



# wwPDB X-ray Structure Validation Summary Report ⓘ

May 26, 2026 – 02:22 PM JST

PDB ID : 26JL / pdb\_000026jl  
Title : Crystal structure of the AAA+ domain of Vibrio cholerae FlrA  
Authors : Dasgupta, J.; Sen, U.; Das, R.  
Deposited on : 2026-05-01  
Resolution : 3.20 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 2.0  
EDS : 3.0  
Buster-report : wwPDB partial adaption of 1.1.7 (2018)  
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)  
CCP4 : 9.0.010 (Gargrove)  
Density-Fitness : 1.0.12  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.49

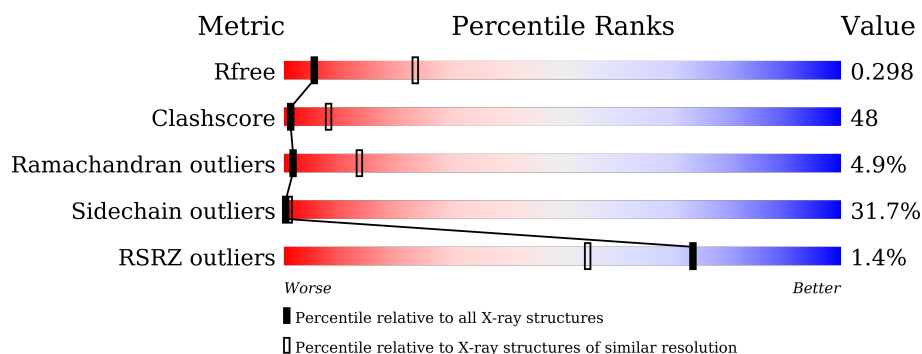
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.20 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	180053	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	259	<div> <div>2%</div> <div>15% 58% 23% 5%</div> </div>
1	B	259	<div> <div>%</div> <div>18% 56% 21% 5%</div> </div>
1	C	259	<div> <div>2%</div> <div>16% 55% 25% .</div> </div>
1	D	259	<div> <div>%</div> <div>19% 53% 25% .</div> </div>
1	E	259	<div> <div>2%</div> <div>18% 53% 25% 5%</div> </div>
1	F	259	<div> <div>%</div> <div>18% 51% 27% 5%</div> </div>



The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	EDO	B	440	-	X	-	-
2	EDO	C	404	-	-	X	-
2	EDO	D	405	-	-	X	-

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 13833 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Flagellar regulatory protein A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	259	Total	C	N	O	S	0	0	0
			2035	1276	367	377	15			
1	B	259	Total	C	N	O	S	0	0	0
			2035	1276	367	377	15			
1	C	259	Total	C	N	O	S	0	0	0
			2035	1276	367	377	15			
1	D	259	Total	C	N	O	S	0	0	0
			2035	1276	367	377	15			
1	E	259	Total	C	N	O	S	0	0	0
			2035	1276	367	377	15			
1	F	259	Total	C	N	O	S	0	0	0
			2035	1276	367	377	15			

There are 66 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	127	ALA	-	expression tag	UNP A0A0H3AM86
A	128	SER	-	expression tag	UNP A0A0H3AM86
A	129	GLY	-	expression tag	UNP A0A0H3AM86
A	130	LEU	-	expression tag	UNP A0A0H3AM86
A	131	VAL	-	expression tag	UNP A0A0H3AM86
A	132	PRO	-	expression tag	UNP A0A0H3AM86
A	133	ARG	-	expression tag	UNP A0A0H3AM86
A	134	GLY	-	expression tag	UNP A0A0H3AM86
A	135	SER	-	expression tag	UNP A0A0H3AM86
A	136	HIS	-	expression tag	UNP A0A0H3AM86
A	137	MET	-	expression tag	UNP A0A0H3AM86
B	127	ALA	-	expression tag	UNP A0A0H3AM86
B	128	SER	-	expression tag	UNP A0A0H3AM86
B	129	GLY	-	expression tag	UNP A0A0H3AM86
B	130	LEU	-	expression tag	UNP A0A0H3AM86
B	131	VAL	-	expression tag	UNP A0A0H3AM86
B	132	PRO	-	expression tag	UNP A0A0H3AM86

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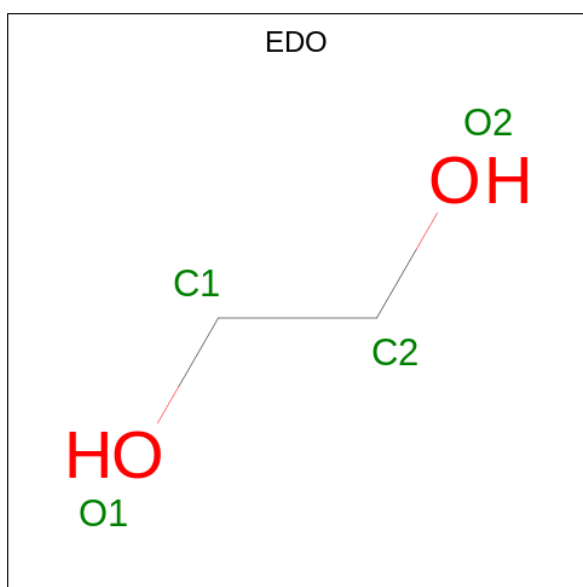
Chain	Residue	Modelled	Actual	Comment	Reference
B	133	ARG	-	expression tag	UNP A0A0H3AM86
B	134	GLY	-	expression tag	UNP A0A0H3AM86
B	135	SER	-	expression tag	UNP A0A0H3AM86
B	136	HIS	-	expression tag	UNP A0A0H3AM86
B	137	MET	-	expression tag	UNP A0A0H3AM86
C	127	ALA	-	expression tag	UNP A0A0H3AM86
C	128	SER	-	expression tag	UNP A0A0H3AM86
C	129	GLY	-	expression tag	UNP A0A0H3AM86
C	130	LEU	-	expression tag	UNP A0A0H3AM86
C	131	VAL	-	expression tag	UNP A0A0H3AM86
C	132	PRO	-	expression tag	UNP A0A0H3AM86
C	133	ARG	-	expression tag	UNP A0A0H3AM86
C	134	GLY	-	expression tag	UNP A0A0H3AM86
C	135	SER	-	expression tag	UNP A0A0H3AM86
C	136	HIS	-	expression tag	UNP A0A0H3AM86
C	137	MET	-	expression tag	UNP A0A0H3AM86
D	127	ALA	-	expression tag	UNP A0A0H3AM86
D	128	SER	-	expression tag	UNP A0A0H3AM86
D	129	GLY	-	expression tag	UNP A0A0H3AM86
D	130	LEU	-	expression tag	UNP A0A0H3AM86
D	131	VAL	-	expression tag	UNP A0A0H3AM86
D	132	PRO	-	expression tag	UNP A0A0H3AM86
D	133	ARG	-	expression tag	UNP A0A0H3AM86
D	134	GLY	-	expression tag	UNP A0A0H3AM86
D	135	SER	-	expression tag	UNP A0A0H3AM86
D	136	HIS	-	expression tag	UNP A0A0H3AM86
D	137	MET	-	expression tag	UNP A0A0H3AM86
E	127	ALA	-	expression tag	UNP A0A0H3AM86
E	128	SER	-	expression tag	UNP A0A0H3AM86
E	129	GLY	-	expression tag	UNP A0A0H3AM86
E	130	LEU	-	expression tag	UNP A0A0H3AM86
E	131	VAL	-	expression tag	UNP A0A0H3AM86
E	132	PRO	-	expression tag	UNP A0A0H3AM86
E	133	ARG	-	expression tag	UNP A0A0H3AM86
E	134	GLY	-	expression tag	UNP A0A0H3AM86
E	135	SER	-	expression tag	UNP A0A0H3AM86
E	136	HIS	-	expression tag	UNP A0A0H3AM86
E	137	MET	-	expression tag	UNP A0A0H3AM86
F	127	ALA	-	expression tag	UNP A0A0H3AM86
F	128	SER	-	expression tag	UNP A0A0H3AM86
F	129	GLY	-	expression tag	UNP A0A0H3AM86
F	130	LEU	-	expression tag	UNP A0A0H3AM86

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Chain	Residue	Modelled	Actual	Comment	Reference
F	131	VAL	-	expression tag	UNP A0A0H3AM86
F	132	PRO	-	expression tag	UNP A0A0H3AM86
F	133	ARG	-	expression tag	UNP A0A0H3AM86
F	134	GLY	-	expression tag	UNP A0A0H3AM86
F	135	SER	-	expression tag	UNP A0A0H3AM86
F	136	HIS	-	expression tag	UNP A0A0H3AM86
F	137	MET	-	expression tag	UNP A0A0H3AM86

- Molecule 2 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C<sub>2</sub>H<sub>6</sub>O<sub>2</sub>) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			4	2	2		
2	A	1	Total	C	O	0	0
			4	2	2		
2	A	1	Total	C	O	0	0
			4	2	2		
2	A	1	Total	C	O	0	0
			4	2	2		
2	A	1	Total	C	O	0	0
			4	2	2		
2	A	1	Total	C	O	0	0
			4	2	2		
2	A	1	Total	C	O	0	0
			4	2	2		
2	A	1	Total	C	O	0	0
			4	2	2		
2	A	1	Total	C	O	0	0
			4	2	2		
2	A	1	Total	C	O	0	0
			4	2	2		
2	A	1	Total	C	O	0	0
			4	2	2		
2	A	1	Total	C	O	0	0
			4	2	2		
2	A	1	Total	C	O	0	0
			4	2	2		
2	A	1	Total	C	O	0	0
			4	2	2		
2	A	1	Total	C	O	0	0
			4	2	2		
2	A	1	Total	C	O	0	0
			4	2	2		
2	A	1	Total	C	O	0	0
			4	2	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total 4	C 2	O 2	0	0
2	A	1	Total 4	C 2	O 2	0	0
2	A	1	Total 4	C 2	O 2	0	0
2	A	1	Total 4	C 2	O 2	0	0
2	A	1	Total 4	C 2	O 2	0	0
2	A	1	Total 4	C 2	O 2	0	0
2	A	1	Total 4	C 2	O 2	0	0
2	A	1	Total 4	C 2	O 2	0	0
2	A	1	Total 4	C 2	O 2	0	0
2	A	1	Total 4	C 2	O 2	0	0
2	A	1	Total 4	C 2	O 2	0	0
2	A	1	Total 4	C 2	O 2	0	0
2	A	1	Total 4	C 2	O 2	0	0
2	A	1	Total 4	C 2	O 2	0	0
2	A	1	Total 4	C 2	O 2	0	0
2	B	1	Total 4	C 2	O 2	0	0
2	B	1	Total 4	C 2	O 2	0	0
2	B	1	Total 4	C 2	O 2	0	0
2	B	1	Total 4	C 2	O 2	0	0
2	B	1	Total 4	C 2	O 2	0	0
2	B	1	Total 4	C 2	O 2	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		
2	C	1	Total	C	O	0	0
			4	2	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	C	O	0	0
			4	2	2		
2	C	1	Total	C	O	0	0
			4	2	2		
2	C	1	Total	C	O	0	0
			4	2	2		
2	C	1	Total	C	O	0	0
			4	2	2		
2	C	1	Total	C	O	0	0
			4	2	2		
2	C	1	Total	C	O	0	0
			4	2	2		
2	C	1	Total	C	O	0	0
			4	2	2		
2	C	1	Total	C	O	0	0
			4	2	2		
2	C	1	Total	C	O	0	0
			4	2	2		
2	C	1	Total	C	O	0	0
			4	2	2		
2	C	1	Total	C	O	0	0
			4	2	2		
2	C	1	Total	C	O	0	0
			4	2	2		
2	C	1	Total	C	O	0	0
			4	2	2		
2	C	1	Total	C	O	0	0
			4	2	2		
2	C	1	Total	C	O	0	0
			4	2	2		
2	C	1	Total	C	O	0	0
			4	2	2		
2	C	1	Total	C	O	0	0
			4	2	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	C	O	0	0
			4	2	2		
2	C	1	Total	C	O	0	0
			4	2	2		
2	C	1	Total	C	O	0	0
			4	2	2		
2	C	1	Total	C	O	0	0
			4	2	2		
2	C	1	Total	C	O	0	0
			4	2	2		
2	C	1	Total	C	O	0	0
			4	2	2		
2	C	1	Total	C	O	0	0
			4	2	2		
2	C	1	Total	C	O	0	0
			4	2	2		
2	C	1	Total	C	O	0	0
			4	2	2		
2	C	1	Total	C	O	0	0
			4	2	2		
2	C	1	Total	C	O	0	0
			4	2	2		
2	C	1	Total	C	O	0	0
			4	2	2		
2	D	1	Total	C	O	0	0
			4	2	2		
2	D	1	Total	C	O	0	0
			4	2	2		
2	D	1	Total	C	O	0	0
			4	2	2		
2	D	1	Total	C	O	0	0
			4	2	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	D	1	Total	C	O	0	0
			4	2	2		
2	D	1	Total	C	O	0	0
			4	2	2		
2	D	1	Total	C	O	0	0
			4	2	2		
2	D	1	Total	C	O	0	0
			4	2	2		
2	D	1	Total	C	O	0	0
			4	2	2		
2	D	1	Total	C	O	0	0
			4	2	2		
2	D	1	Total	C	O	0	0
			4	2	2		
2	D	1	Total	C	O	0	0
			4	2	2		
2	D	1	Total	C	O	0	0
			4	2	2		
2	D	1	Total	C	O	0	0
			4	2	2		
2	D	1	Total	C	O	0	0
			4	2	2		
2	D	1	Total	C	O	0	0
			4	2	2		
2	D	1	Total	C	O	0	0
			4	2	2		
2	D	1	Total	C	O	0	0
			4	2	2		
2	D	1	Total	C	O	0	0
			4	2	2		
2	D	1	Total	C	O	0	0
			4	2	2		
2	D	1	Total	C	O	0	0
			4	2	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	D	1	Total	C	O	0	0
			4	2	2		
2	D	1	Total	C	O	0	0
			4	2	2		
2	D	1	Total	C	O	0	0
			4	2	2		
2	D	1	Total	C	O	0	0
			4	2	2		
2	D	1	Total	C	O	0	0
			4	2	2		
2	D	1	Total	C	O	0	0
			4	2	2		
2	D	1	Total	C	O	0	0
			4	2	2		
2	D	1	Total	C	O	0	0
			4	2	2		
2	D	1	Total	C	O	0	0
			4	2	2		
2	D	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total	C	O	0	0
			4	2	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	E	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total	C	O	0	0
			4	2	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	E	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total	C	O	0	0
			4	2	2		
2	F	1	Total	C	O	0	0
			4	2	2		
2	F	1	Total	C	O	0	0
			4	2	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	F	1	Total	C	O	0	0
			4	2	2		
2	F	1	Total	C	O	0	0
			4	2	2		
2	F	1	Total	C	O	0	0
			4	2	2		
2	F	1	Total	C	O	0	0
			4	2	2		
2	F	1	Total	C	O	0	0
			4	2	2		
2	F	1	Total	C	O	0	0
			4	2	2		
2	F	1	Total	C	O	0	0
			4	2	2		
2	F	1	Total	C	O	0	0
			4	2	2		
2	F	1	Total	C	O	0	0
			4	2	2		
2	F	1	Total	C	O	0	0
			4	2	2		
2	F	1	Total	C	O	0	0
			4	2	2		
2	F	1	Total	C	O	0	0
			4	2	2		
2	F	1	Total	C	O	0	0
			4	2	2		
2	F	1	Total	C	O	0	0
			4	2	2		
2	F	1	Total	C	O	0	0
			4	2	2		
2	F	1	Total	C	O	0	0
			4	2	2		
2	F	1	Total	C	O	0	0
			4	2	2		
2	F	1	Total	C	O	0	0
			4	2	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	F	1	Total 4	C 2	O 2	0	0
2	F	1	Total 4	C 2	O 2	0	0
2	F	1	Total 4	C 2	O 2	0	0
2	F	1	Total 4	C 2	O 2	0	0
2	F	1	Total 4	C 2	O 2	0	0
2	F	1	Total 4	C 2	O 2	0	0
2	F	1	Total 4	C 2	O 2	0	0
2	F	1	Total 4	C 2	O 2	0	0
2	F	1	Total 4	C 2	O 2	0	0
2	F	1	Total 4	C 2	O 2	0	0
2	F	1	Total 4	C 2	O 2	0	0
2	F	1	Total 4	C 2	O 2	0	0
2	F	1	Total 4	C 2	O 2	0	0
2	F	1	Total 4	C 2	O 2	0	0
2	F	1	Total 4	C 2	O 2	0	0
2	F	1	Total 4	C 2	O 2	0	0
2	F	1	Total 4	C 2	O 2	0	0
2	F	1	Total 4	C 2	O 2	0	0
2	F	1	Total 4	C 2	O 2	0	0
2	F	1	Total 4	C 2	O 2	0	0
2	F	1	Total 4	C 2	O 2	0	0
2	F	1	Total 4	C 2	O 2	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	F	1	Total	C	O	0	0
			4	2	2		

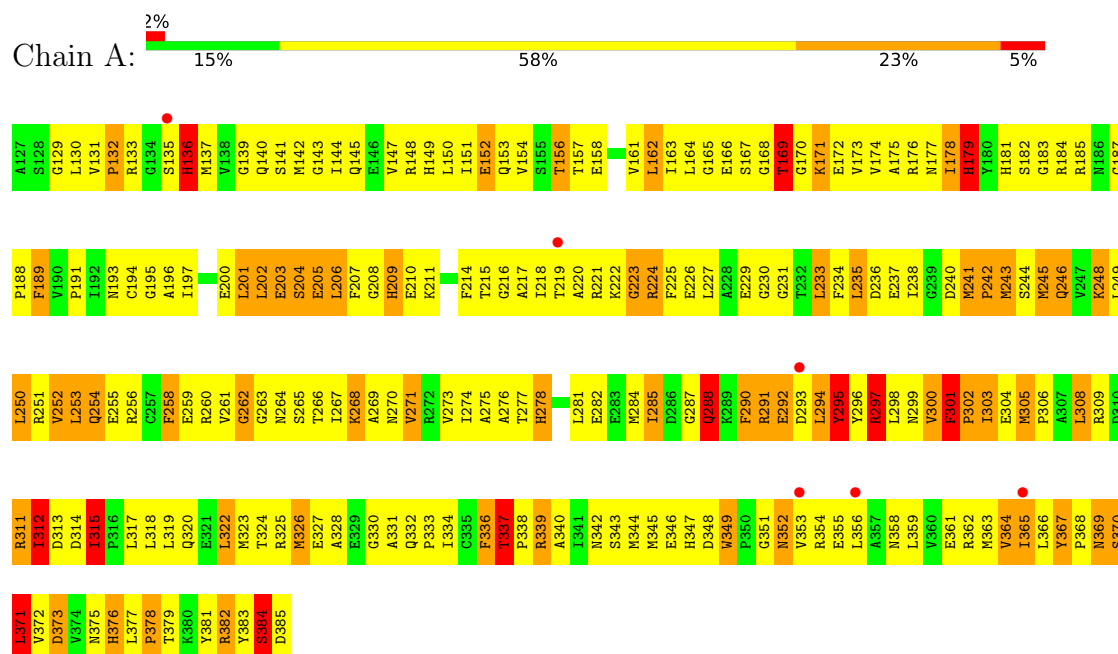
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	96	Total	O	0	0
			96	96		
3	B	91	Total	O	0	0
			91	91		
3	C	100	Total	O	0	0
			100	100		
3	D	98	Total	O	0	0
			98	98		
3	E	97	Total	O	0	0
			97	97		
3	F	101	Total	O	0	0
			101	101		

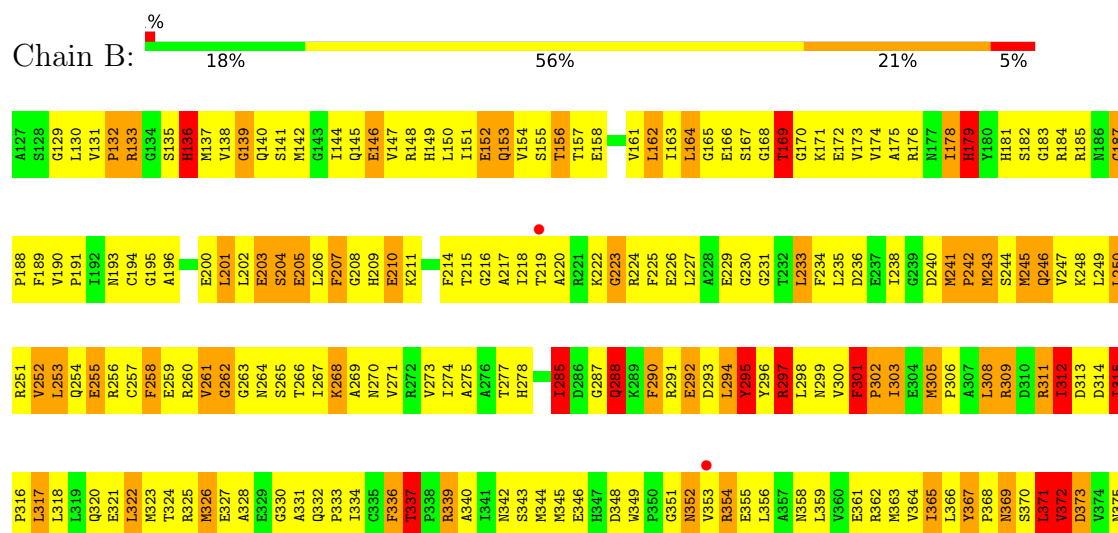
### 3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

#### • Molecule 1: Flagellar regulatory protein A

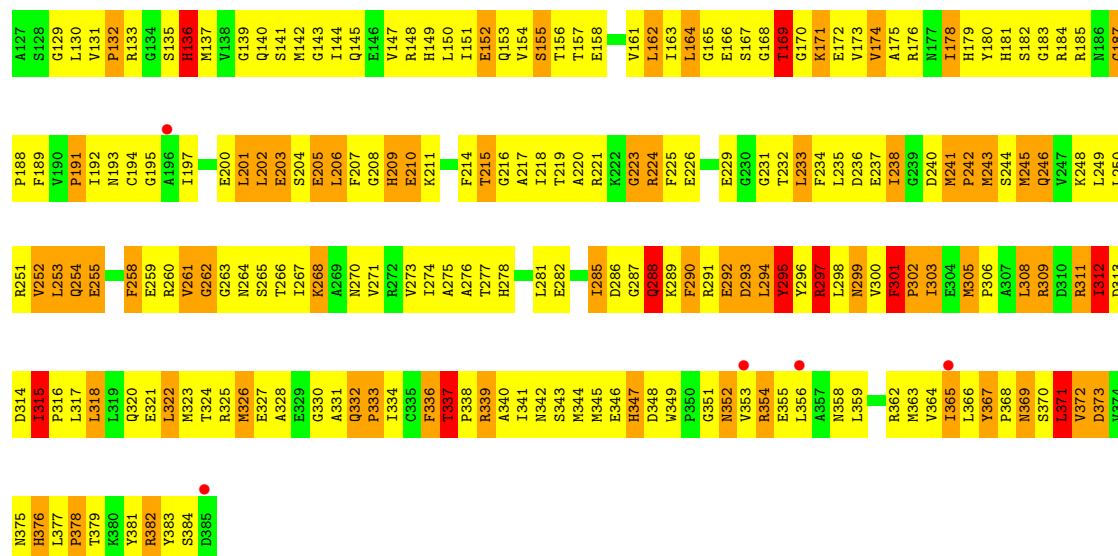
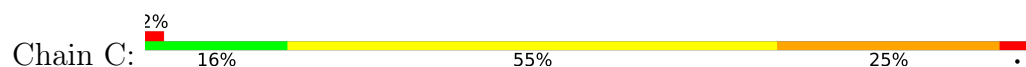


#### • Molecule 1: Flagellar regulatory protein A

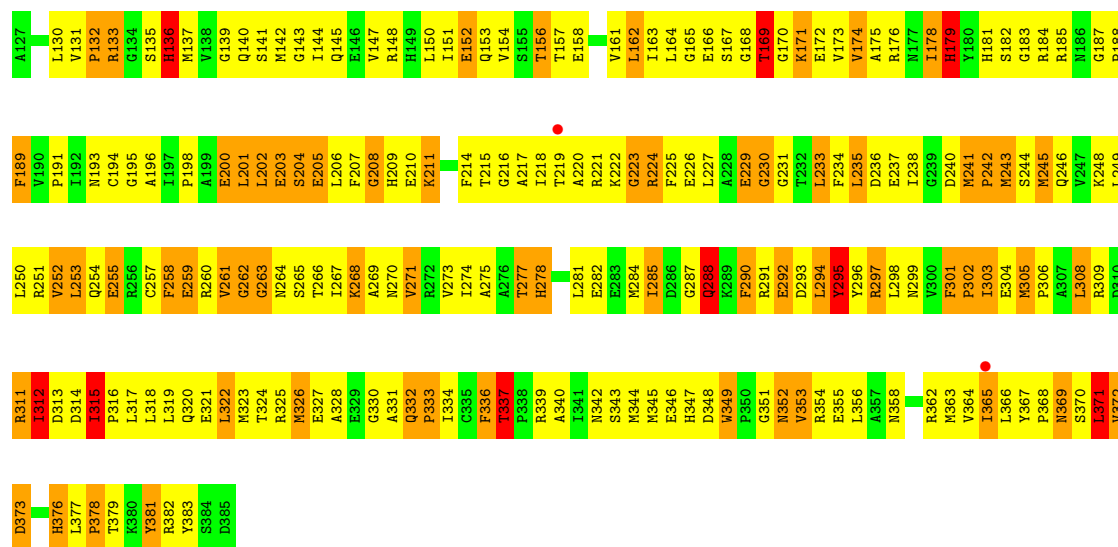
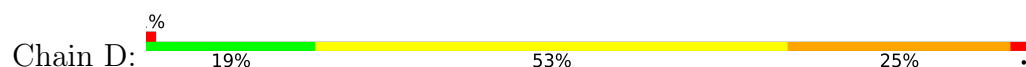


H376  
L377  
P378  
T379  
K380  
Y381  
R382  
S384  
D385

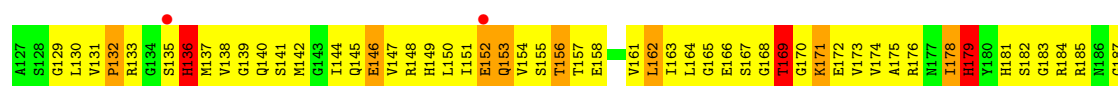
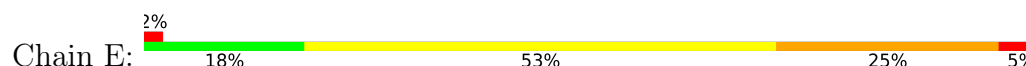
• Molecule 1: Flagellar regulatory protein A

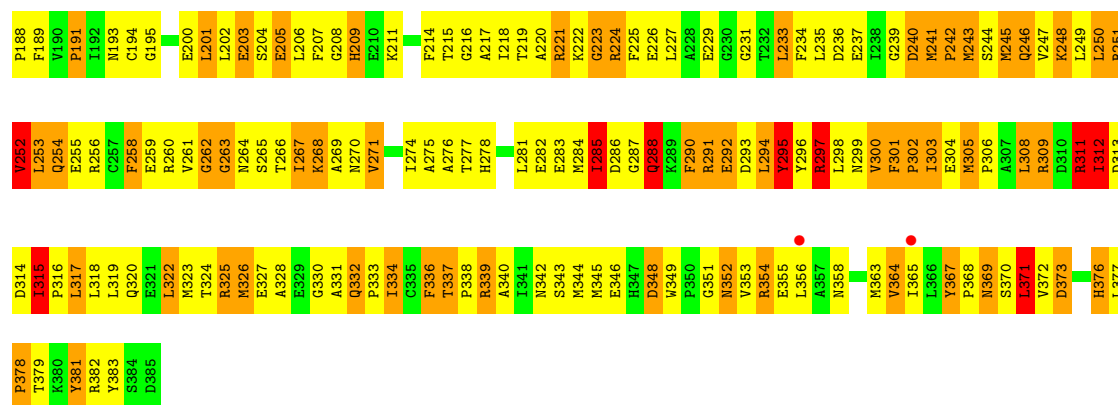


• Molecule 1: Flagellar regulatory protein A

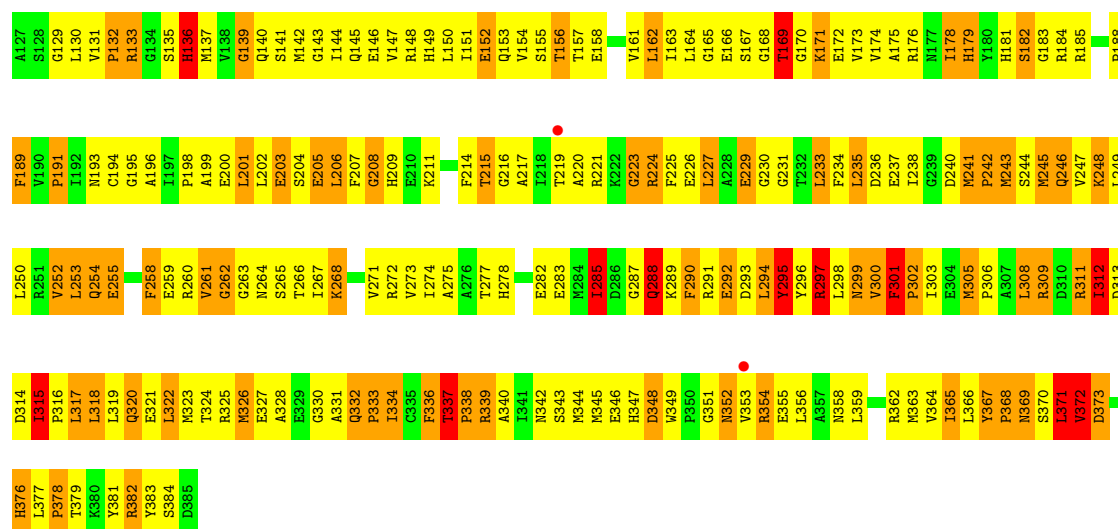
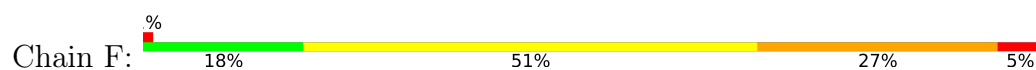


• Molecule 1: Flagellar regulatory protein A





● Molecule 1: Flagellar regulatory protein A



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	101.17Å 101.17Å 49.91Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	9.98 – 3.20 9.98 – 3.20	Depositor EDS
% Data completeness (in resolution range)	99.1 (9.98-3.20) 99.2 (9.98-3.20)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.18 (at 3.12Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, $R_{free}$	0.220 , 0.295 0.261 , 0.298	Depositor DCC
$R_{free}$ test set	2047 reflections (6.64%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	52.0	Xtriage
Anisotropy	0.370	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.16 , 45.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.499 for h+k,-h,l 0.499 for -k,h+k,l 0.499 for k,-h-k,l 0.499 for -h-k,h,l 0.499 for -h,-k,l 0.418 for h,-h-k,-l 0.418 for -h-k,k,-l 0.418 for k,h,-l 0.418 for -k,-h,-l 0.418 for -h,h+k,-l 0.418 for h+k,-k,-l	Xtriage
$F_o, F_c$ correlation	0.83	EDS
Total number of atoms	13833	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.52% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: EDO

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z  > 5$	RMSZ	$\# Z  > 5$
1	A	0.94	1/2072 (0.0%)	1.55	45/2798 (1.6%)
1	B	0.95	2/2072 (0.1%)	1.56	38/2798 (1.4%)
1	C	0.98	4/2072 (0.2%)	1.59	38/2798 (1.4%)
1	D	0.93	1/2072 (0.0%)	1.56	38/2798 (1.4%)
1	E	0.97	2/2072 (0.1%)	1.57	45/2798 (1.6%)
1	F	0.93	1/2072 (0.0%)	1.61	51/2798 (1.8%)
All	All	0.95	11/12432 (0.1%)	1.57	255/16788 (1.5%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	4
1	B	0	7
1	C	0	4
1	D	0	5
1	E	0	6
1	F	0	6
All	All	0	32

The worst 5 of 11 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	347	HIS	CG-CD2	-7.40	1.27	1.35
1	C	209	HIS	CG-CD2	-7.06	1.28	1.35
1	C	352	ASN	CA-C	-6.33	1.50	1.53
1	E	352	ASN	CA-C	-6.30	1.50	1.53
1	B	352	ASN	CA-C	-6.20	1.50	1.53

The worst 5 of 255 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	262	GLY	N-CA-C	11.75	126.83	112.73
1	B	351	GLY	N-CA-C	-10.45	92.55	111.25
1	E	351	GLY	N-CA-C	-10.32	91.20	112.04
1	F	351	GLY	N-CA-C	-10.26	93.02	112.51
1	B	262	GLY	N-CA-C	10.01	124.74	112.73

There are no chirality outliers.

5 of 32 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	169	THR	Peptide
1	A	223	GLY	Peptide
1	A	294	LEU	Peptide
1	A	297	ARG	Peptide
1	B	133	ARG	Peptide

## 5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2035	0	2036	217	0
1	B	2035	0	2036	217	0
1	C	2035	0	2036	232	0
1	D	2035	0	2036	218	0
1	E	2035	0	2036	209	0
1	F	2035	0	2036	219	0
2	A	172	0	258	8	0
2	B	188	0	281	8	0
2	C	152	0	228	9	0
2	D	156	0	234	9	0
2	E	192	0	288	8	0
2	F	180	0	270	5	0
3	A	96	0	0	18	0
3	B	91	0	0	16	0
3	C	100	0	0	20	0
3	D	98	0	0	18	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	E	97	0	0	16	0
3	F	101	0	0	18	0
All	All	13833	0	13775	1294	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 48.

The worst 5 of 1294 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:154:VAL:HG21	1:B:181:HIS:HB3	1.30	1.10
1:F:154:VAL:HG21	1:F:181:HIS:HB3	1.33	1.10
1:C:208:GLY:HA3	1:C:260:ARG:HG3	1.08	1.08
1:F:208:GLY:HA3	1:F:260:ARG:HG3	1.10	1.07
1:D:208:GLY:HA3	1:D:260:ARG:HG3	1.10	1.06

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	257/259 (99%)	213 (83%)	32 (12%)	12 (5%)	2	14
1	B	257/259 (99%)	215 (84%)	29 (11%)	13 (5%)	1	13
1	C	257/259 (99%)	214 (83%)	31 (12%)	12 (5%)	2	14
1	D	257/259 (99%)	214 (83%)	31 (12%)	12 (5%)	2	14
1	E	257/259 (99%)	215 (84%)	30 (12%)	12 (5%)	2	14
1	F	257/259 (99%)	217 (84%)	25 (10%)	15 (6%)	1	10
All	All	1542/1554 (99%)	1288 (84%)	178 (12%)	76 (5%)	1	13



5 of 76 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	191	PRO
1	A	263	GLY
1	A	288	GLN
1	A	328	ALA
1	B	263	GLY

### 5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	220/220 (100%)	147 (67%)	73 (33%)	0	1
1	B	220/220 (100%)	150 (68%)	70 (32%)	0	1
1	C	220/220 (100%)	149 (68%)	71 (32%)	0	1
1	D	220/220 (100%)	155 (70%)	65 (30%)	0	1
1	E	220/220 (100%)	150 (68%)	70 (32%)	0	1
1	F	220/220 (100%)	150 (68%)	70 (32%)	0	1
All	All	1320/1320 (100%)	901 (68%)	419 (32%)	0	1

5 of 419 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	204	SER
1	E	156	THR
1	F	309	ARG
1	D	235	LEU
1	D	305	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 53 such sidechains are listed below:

Mol	Chain	Res	Type
1	D	160	ASN
1	E	140	GLN

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Mol	Chain	Res	Type
1	F	181	HIS
1	D	177	ASN
1	D	270	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

### 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

### 5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

### 5.6 Ligand geometry [i](#)

260 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
2	EDO	F	405	-	3,3,3	0.42	0	2,2,2	0.37	0
2	EDO	F	428	-	3,3,3	0.55	0	2,2,2	0.17	0
2	EDO	E	434	-	3,3,3	0.77	0	2,2,2	0.30	0
2	EDO	B	447	-	3,3,3	0.42	0	2,2,2	0.37	0
2	EDO	F	444	-	3,3,3	0.42	0	2,2,2	0.40	0
2	EDO	E	445	-	3,3,3	0.55	0	2,2,2	0.21	0
2	EDO	A	440	-	3,3,3	0.49	0	2,2,2	0.30	0
2	EDO	B	408	-	3,3,3	0.56	0	2,2,2	0.02	0
2	EDO	F	423	-	3,3,3	0.64	0	2,2,2	0.24	0
2	EDO	B	443	-	3,3,3	0.50	0	2,2,2	0.21	0
2	EDO	A	438	-	3,3,3	0.42	0	2,2,2	0.40	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	EDO	A	408	-	3,3,3	0.46	0	2,2,2	0.74	0
2	EDO	C	419	-	3,3,3	0.52	0	2,2,2	0.18	0
2	EDO	A	416	-	3,3,3	0.48	0	2,2,2	0.26	0
2	EDO	A	423	-	3,3,3	0.58	0	2,2,2	0.13	0
2	EDO	E	402	-	3,3,3	0.44	0	2,2,2	0.37	0
2	EDO	C	436	-	3,3,3	0.45	0	2,2,2	0.47	0
2	EDO	E	426	-	3,3,3	0.44	0	2,2,2	0.43	0
2	EDO	F	419	-	3,3,3	0.50	0	2,2,2	0.15	0
2	EDO	A	424	-	3,3,3	0.44	0	2,2,2	0.53	0
2	EDO	D	433	-	3,3,3	0.41	0	2,2,2	0.40	0
2	EDO	B	445	-	3,3,3	0.58	0	2,2,2	0.11	0
2	EDO	E	431	-	3,3,3	0.49	0	2,2,2	0.28	0
2	EDO	F	413	-	3,3,3	0.54	0	2,2,2	0.12	0
2	EDO	A	413	-	3,3,3	0.61	0	2,2,2	0.18	0
2	EDO	D	417	-	3,3,3	0.54	0	2,2,2	0.19	0
2	EDO	E	432	-	3,3,3	0.38	0	2,2,2	0.44	0
2	EDO	A	418	-	3,3,3	0.54	0	2,2,2	0.20	0
2	EDO	B	435	-	3,3,3	0.38	0	2,2,2	0.50	0
2	EDO	C	408	-	3,3,3	0.48	0	2,2,2	0.27	0
2	EDO	C	438	-	3,3,3	0.58	0	2,2,2	0.40	0
2	EDO	E	403	-	3,3,3	0.48	0	2,2,2	0.29	0
2	EDO	E	436	-	3,3,3	0.44	0	2,2,2	0.47	0
2	EDO	E	437	-	3,3,3	0.52	0	2,2,2	0.17	0
2	EDO	C	424	-	3,3,3	0.54	0	2,2,2	0.21	0
2	EDO	E	447	-	3,3,3	0.52	0	2,2,2	0.23	0
2	EDO	A	412	-	3,3,3	0.43	0	2,2,2	0.41	0
2	EDO	C	413	-	3,3,3	0.49	0	2,2,2	0.14	0
2	EDO	B	430	-	3,3,3	0.57	0	2,2,2	0.08	0
2	EDO	E	420	-	3,3,3	0.43	0	2,2,2	0.50	0
2	EDO	D	410	-	3,3,3	0.37	0	2,2,2	0.54	0
2	EDO	E	413	-	3,3,3	0.46	0	2,2,2	0.39	0
2	EDO	E	438	-	3,3,3	0.72	0	2,2,2	0.15	0
2	EDO	E	418	-	3,3,3	0.58	0	2,2,2	0.08	0
2	EDO	F	414	-	3,3,3	0.42	0	2,2,2	0.33	0
2	EDO	A	427	-	3,3,3	0.37	0	2,2,2	0.54	0
2	EDO	C	410	-	3,3,3	0.49	0	2,2,2	0.65	0
2	EDO	F	436	-	3,3,3	0.56	0	2,2,2	0.26	0
2	EDO	A	431	-	3,3,3	0.49	0	2,2,2	0.28	0
2	EDO	C	430	-	3,3,3	0.50	0	2,2,2	0.16	0
2	EDO	F	416	-	3,3,3	0.40	0	2,2,2	0.83	0
2	EDO	D	439	-	3,3,3	0.53	0	2,2,2	0.30	0
2	EDO	E	446	-	3,3,3	0.54	0	2,2,2	0.17	0
2	EDO	F	422	-	3,3,3	0.58	0	2,2,2	0.14	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	EDO	B	428	-	3,3,3	0.45	0	2,2,2	0.28	0
2	EDO	A	422	-	3,3,3	0.45	0	2,2,2	0.40	0
2	EDO	E	444	-	3,3,3	0.51	0	2,2,2	0.19	0
2	EDO	B	444	-	3,3,3	0.52	0	2,2,2	0.26	0
2	EDO	A	409	-	3,3,3	0.40	0	2,2,2	0.33	0
2	EDO	C	416	-	3,3,3	0.59	0	2,2,2	0.26	0
2	EDO	C	431	-	3,3,3	0.48	0	2,2,2	0.27	0
2	EDO	A	429	-	3,3,3	0.47	0	2,2,2	0.27	0
2	EDO	F	438	-	3,3,3	0.49	0	2,2,2	0.10	0
2	EDO	D	422	-	3,3,3	0.52	0	2,2,2	0.12	0
2	EDO	A	437	-	3,3,3	0.51	0	2,2,2	0.20	0
2	EDO	C	422	-	3,3,3	0.39	0	2,2,2	0.68	0
2	EDO	C	432	-	3,3,3	0.55	0	2,2,2	0.18	0
2	EDO	D	428	-	3,3,3	0.47	0	2,2,2	0.31	0
2	EDO	C	425	-	3,3,3	0.43	0	2,2,2	0.47	0
2	EDO	C	429	-	3,3,3	0.65	0	2,2,2	0.42	0
2	EDO	C	437	-	3,3,3	0.57	0	2,2,2	0.09	0
2	EDO	C	412	-	3,3,3	0.46	0	2,2,2	0.41	0
2	EDO	B	404	-	3,3,3	0.42	0	2,2,2	0.44	0
2	EDO	B	431	-	3,3,3	0.50	0	2,2,2	0.19	0
2	EDO	F	442	-	3,3,3	0.51	0	2,2,2	0.32	0
2	EDO	B	438	-	3,3,3	0.54	0	2,2,2	0.07	0
2	EDO	B	422	-	3,3,3	0.53	0	2,2,2	0.24	0
2	EDO	F	410	-	3,3,3	0.43	0	2,2,2	0.41	0
2	EDO	F	441	-	3,3,3	0.49	0	2,2,2	0.29	0
2	EDO	A	404	-	3,3,3	0.61	0	2,2,2	0.11	0
2	EDO	A	410	-	3,3,3	0.46	0	2,2,2	0.48	0
2	EDO	B	409	-	3,3,3	0.32	0	2,2,2	0.65	0
2	EDO	B	423	-	3,3,3	0.44	0	2,2,2	0.57	0
2	EDO	D	416	-	3,3,3	0.65	0	2,2,2	0.25	0
2	EDO	D	411	-	3,3,3	0.45	0	2,2,2	0.77	0
2	EDO	E	401	-	3,3,3	0.45	0	2,2,2	0.32	0
2	EDO	D	406	-	3,3,3	0.56	0	2,2,2	0.19	0
2	EDO	B	442	-	3,3,3	0.47	0	2,2,2	0.30	0
2	EDO	C	427	-	3,3,3	0.45	0	2,2,2	0.36	0
2	EDO	D	404	-	3,3,3	0.50	0	2,2,2	0.24	0
2	EDO	E	435	-	3,3,3	0.49	0	2,2,2	0.21	0
2	EDO	B	418	-	3,3,3	0.48	0	2,2,2	0.27	0
2	EDO	C	404	-	3,3,3	0.41	0	2,2,2	0.50	0
2	EDO	F	408	-	3,3,3	0.53	0	2,2,2	0.24	0
2	EDO	A	426	-	3,3,3	0.47	0	2,2,2	0.47	0
2	EDO	F	439	-	3,3,3	0.59	0	2,2,2	0.12	0
2	EDO	A	402	-	3,3,3	0.39	0	2,2,2	0.51	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	EDO	D	438	-	3,3,3	0.47	0	2,2,2	0.16	0
2	EDO	E	414	-	3,3,3	0.60	0	2,2,2	0.11	0
2	EDO	A	443	-	3,3,3	0.55	0	2,2,2	0.12	0
2	EDO	F	404	-	3,3,3	0.71	0	2,2,2	0.17	0
2	EDO	D	423	-	3,3,3	0.57	0	2,2,2	0.20	0
2	EDO	F	418	-	3,3,3	0.50	0	2,2,2	0.32	0
2	EDO	F	443	-	3,3,3	0.50	0	2,2,2	0.23	0
2	EDO	F	432	-	3,3,3	0.44	0	2,2,2	0.53	0
2	EDO	A	421	-	3,3,3	0.43	0	2,2,2	0.60	0
2	EDO	A	430	-	3,3,3	0.44	0	2,2,2	0.50	0
2	EDO	A	432	-	3,3,3	0.76	0	2,2,2	0.26	0
2	EDO	F	403	-	3,3,3	0.56	0	2,2,2	0.09	0
2	EDO	E	440	-	3,3,3	0.48	0	2,2,2	0.15	0
2	EDO	D	402	-	3,3,3	0.51	0	2,2,2	0.25	0
2	EDO	E	411	-	3,3,3	0.50	0	2,2,2	0.27	0
2	EDO	E	415	-	3,3,3	0.52	0	2,2,2	0.26	0
2	EDO	C	402	-	3,3,3	0.55	0	2,2,2	0.14	0
2	EDO	F	412	-	3,3,3	0.52	0	2,2,2	0.18	0
2	EDO	B	439	-	3,3,3	0.46	0	2,2,2	0.21	0
2	EDO	B	424	-	3,3,3	0.44	0	2,2,2	0.43	0
2	EDO	A	417	-	3,3,3	0.41	0	2,2,2	0.41	0
2	EDO	F	417	-	3,3,3	0.42	0	2,2,2	0.29	0
2	EDO	B	413	-	3,3,3	0.60	0	2,2,2	0.05	0
2	EDO	F	401	-	3,3,3	0.44	0	2,2,2	0.32	0
2	EDO	B	433	-	3,3,3	0.56	0	2,2,2	0.11	0
2	EDO	F	420	-	3,3,3	0.47	0	2,2,2	0.31	0
2	EDO	A	439	-	3,3,3	0.58	0	2,2,2	0.06	0
2	EDO	D	427	-	3,3,3	0.43	0	2,2,2	0.54	0
2	EDO	E	433	-	3,3,3	0.55	0	2,2,2	0.04	0
2	EDO	F	434	-	3,3,3	0.47	0	2,2,2	0.22	0
2	EDO	B	426	-	3,3,3	0.48	0	2,2,2	0.46	0
2	EDO	B	402	-	3,3,3	0.53	0	2,2,2	0.29	0
2	EDO	E	441	-	3,3,3	0.39	0	2,2,2	0.66	0
2	EDO	C	421	-	3,3,3	0.59	0	2,2,2	0.18	0
2	EDO	A	403	-	3,3,3	0.53	0	2,2,2	0.35	0
2	EDO	D	425	-	3,3,3	0.43	0	2,2,2	0.41	0
2	EDO	E	409	-	3,3,3	0.33	0	2,2,2	0.67	0
2	EDO	F	425	-	3,3,3	0.46	0	2,2,2	0.33	0
2	EDO	B	432	-	3,3,3	0.49	0	2,2,2	0.41	0
2	EDO	D	429	-	3,3,3	0.46	0	2,2,2	0.33	0
2	EDO	E	408	-	3,3,3	0.56	0	2,2,2	0.05	0
2	EDO	E	439	-	3,3,3	0.39	0	2,2,2	0.48	0
2	EDO	F	402	-	3,3,3	0.44	0	2,2,2	0.31	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	EDO	B	403	-	3,3,3	0.51	0	2,2,2	0.28	0
2	EDO	D	436	-	3,3,3	0.52	0	2,2,2	0.18	0
2	EDO	C	423	-	3,3,3	0.45	0	2,2,2	0.41	0
2	EDO	E	427	-	3,3,3	0.64	0	2,2,2	0.01	0
2	EDO	D	403	-	3,3,3	0.52	0	2,2,2	0.04	0
2	EDO	E	442	-	3,3,3	0.52	0	2,2,2	0.18	0
2	EDO	B	427	-	3,3,3	0.46	0	2,2,2	0.31	0
2	EDO	B	412	-	3,3,3	0.45	0	2,2,2	0.44	0
2	EDO	E	443	-	3,3,3	0.54	0	2,2,2	0.27	0
2	EDO	F	424	-	3,3,3	0.46	0	2,2,2	0.28	0
2	EDO	C	426	-	3,3,3	0.46	0	2,2,2	0.29	0
2	EDO	A	435	-	3,3,3	0.55	0	2,2,2	0.14	0
2	EDO	E	419	-	3,3,3	0.54	0	2,2,2	0.15	0
2	EDO	B	420	-	3,3,3	0.50	0	2,2,2	0.26	0
2	EDO	D	407	-	3,3,3	0.59	0	2,2,2	0.08	0
2	EDO	E	425	-	3,3,3	0.42	0	2,2,2	0.66	0
2	EDO	D	413	-	3,3,3	0.44	0	2,2,2	0.50	0
2	EDO	A	420	-	3,3,3	0.59	0	2,2,2	0.16	0
2	EDO	F	421	-	3,3,3	0.59	0	2,2,2	0.18	0
2	EDO	B	429	-	3,3,3	0.38	0	2,2,2	0.49	0
2	EDO	A	442	-	3,3,3	0.42	0	2,2,2	0.57	0
2	EDO	A	414	-	3,3,3	0.53	0	2,2,2	0.17	0
2	EDO	B	415	-	3,3,3	0.55	0	2,2,2	0.01	0
2	EDO	D	434	-	3,3,3	0.46	0	2,2,2	0.42	0
2	EDO	F	415	-	3,3,3	0.33	0	2,2,2	0.55	0
2	EDO	B	446	-	3,3,3	0.50	0	2,2,2	0.35	0
2	EDO	B	416	-	3,3,3	0.59	0	2,2,2	0.07	0
2	EDO	A	415	-	3,3,3	0.55	0	2,2,2	0.08	0
2	EDO	E	416	-	3,3,3	0.61	0	2,2,2	0.16	0
2	EDO	D	408	-	3,3,3	0.52	0	2,2,2	0.17	0
2	EDO	D	430	-	3,3,3	0.34	0	2,2,2	0.64	0
2	EDO	E	428	-	3,3,3	0.49	0	2,2,2	0.21	0
2	EDO	D	424	-	3,3,3	0.43	0	2,2,2	0.63	0
2	EDO	C	403	-	3,3,3	0.50	0	2,2,2	0.30	0
2	EDO	E	404	-	3,3,3	0.33	0	2,2,2	0.71	0
2	EDO	C	411	-	3,3,3	0.49	0	2,2,2	0.21	0
2	EDO	C	417	-	3,3,3	0.54	0	2,2,2	0.14	0
2	EDO	B	441	-	3,3,3	0.49	0	2,2,2	0.37	0
2	EDO	C	409	-	3,3,3	0.35	0	2,2,2	0.67	0
2	EDO	B	419	-	3,3,3	0.50	0	2,2,2	0.28	0
2	EDO	B	405	-	3,3,3	0.47	0	2,2,2	0.27	0
2	EDO	A	441	-	3,3,3	0.35	0	2,2,2	0.59	0
2	EDO	B	414	-	3,3,3	0.49	0	2,2,2	0.28	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	EDO	F	409	-	3,3,3	0.51	0	2,2,2	0.33	0
2	EDO	A	419	-	3,3,3	0.56	0	2,2,2	0.02	0
2	EDO	E	424	-	3,3,3	0.63	0	2,2,2	0.07	0
2	EDO	B	407	-	3,3,3	0.57	0	2,2,2	0.17	0
2	EDO	F	431	-	3,3,3	0.48	0	2,2,2	0.26	0
2	EDO	F	437	-	3,3,3	0.40	0	2,2,2	0.68	0
2	EDO	E	406	-	3,3,3	0.58	0	2,2,2	0.04	0
2	EDO	F	407	-	3,3,3	0.56	0	2,2,2	0.24	0
2	EDO	C	420	-	3,3,3	0.52	0	2,2,2	0.18	0
2	EDO	A	407	-	3,3,3	0.37	0	2,2,2	0.56	0
2	EDO	D	415	-	3,3,3	0.38	0	2,2,2	0.49	0
2	EDO	E	412	-	3,3,3	0.43	0	2,2,2	0.26	0
2	EDO	F	433	-	3,3,3	0.43	0	2,2,2	0.35	0
2	EDO	B	401	-	3,3,3	0.56	0	2,2,2	0.16	0
2	EDO	B	417	-	3,3,3	0.53	0	2,2,2	0.10	0
2	EDO	C	414	-	3,3,3	0.43	0	2,2,2	0.36	0
2	EDO	A	433	-	3,3,3	0.42	0	2,2,2	0.46	0
2	EDO	A	401	-	3,3,3	0.53	0	2,2,2	0.30	0
2	EDO	C	415	-	3,3,3	0.61	0	2,2,2	0.16	0
2	EDO	D	432	-	3,3,3	0.47	0	2,2,2	0.36	0
2	EDO	F	430	-	3,3,3	0.45	0	2,2,2	0.43	0
2	EDO	F	411	-	3,3,3	0.46	0	2,2,2	0.40	0
2	EDO	B	406	-	3,3,3	0.61	0	2,2,2	0.05	0
2	EDO	A	411	-	3,3,3	0.56	0	2,2,2	0.13	0
2	EDO	F	406	-	3,3,3	0.57	0	2,2,2	0.17	0
2	EDO	B	440	-	3,3,3	2.11	2 (66%)	2,2,2	1.17	0
2	EDO	F	445	-	3,3,3	0.51	0	2,2,2	0.25	0
2	EDO	D	401	-	3,3,3	0.56	0	2,2,2	0.20	0
2	EDO	D	412	-	3,3,3	0.46	0	2,2,2	0.15	0
2	EDO	E	417	-	3,3,3	0.56	0	2,2,2	0.18	0
2	EDO	D	420	-	3,3,3	0.54	0	2,2,2	0.09	0
2	EDO	E	423	-	3,3,3	0.55	0	2,2,2	0.10	0
2	EDO	D	414	-	3,3,3	0.51	0	2,2,2	0.16	0
2	EDO	A	425	-	3,3,3	0.46	0	2,2,2	0.31	0
2	EDO	E	429	-	3,3,3	0.42	0	2,2,2	0.54	0
2	EDO	A	434	-	3,3,3	0.59	0	2,2,2	0.15	0
2	EDO	E	410	-	3,3,3	0.42	0	2,2,2	0.75	0
2	EDO	B	411	-	3,3,3	0.43	0	2,2,2	0.26	0
2	EDO	C	407	-	3,3,3	0.47	0	2,2,2	0.31	0
2	EDO	F	427	-	3,3,3	0.53	0	2,2,2	0.14	0
2	EDO	D	409	-	3,3,3	0.51	0	2,2,2	0.20	0
2	EDO	C	434	-	3,3,3	0.46	0	2,2,2	0.28	0
2	EDO	B	421	-	3,3,3	0.51	0	2,2,2	0.19	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	EDO	D	405	-	3,3,3	0.39	0	2,2,2	0.46	0
2	EDO	F	426	-	3,3,3	0.49	0	2,2,2	0.24	0
2	EDO	C	433	-	3,3,3	0.80	0	2,2,2	0.29	0
2	EDO	D	418	-	3,3,3	0.56	0	2,2,2	0.15	0
2	EDO	F	440	-	3,3,3	0.49	0	2,2,2	0.20	0
2	EDO	D	437	-	3,3,3	0.58	0	2,2,2	0.25	0
2	EDO	E	407	-	3,3,3	0.49	0	2,2,2	0.26	0
2	EDO	A	428	-	3,3,3	0.59	0	2,2,2	0.08	0
2	EDO	C	401	-	3,3,3	0.60	0	2,2,2	0.24	0
2	EDO	B	436	-	3,3,3	0.53	0	2,2,2	0.20	0
2	EDO	C	435	-	3,3,3	0.44	0	2,2,2	0.43	0
2	EDO	D	421	-	3,3,3	0.51	0	2,2,2	0.25	0
2	EDO	F	429	-	3,3,3	0.42	0	2,2,2	0.64	0
2	EDO	F	435	-	3,3,3	0.33	0	2,2,2	0.56	0
2	EDO	D	435	-	3,3,3	0.59	0	2,2,2	0.17	0
2	EDO	B	425	-	3,3,3	0.53	0	2,2,2	0.24	0
2	EDO	A	436	-	3,3,3	0.43	0	2,2,2	0.39	0
2	EDO	E	422	-	3,3,3	0.46	0	2,2,2	0.31	0
2	EDO	C	406	-	3,3,3	0.54	0	2,2,2	0.05	0
2	EDO	D	431	-	3,3,3	0.68	0	2,2,2	0.49	0
2	EDO	E	405	-	3,3,3	0.52	0	2,2,2	0.22	0
2	EDO	D	419	-	3,3,3	0.61	0	2,2,2	0.04	0
2	EDO	E	430	-	3,3,3	0.44	0	2,2,2	0.33	0
2	EDO	C	428	-	3,3,3	0.34	0	2,2,2	0.56	0
2	EDO	E	421	-	3,3,3	0.44	0	2,2,2	0.33	0
2	EDO	A	405	-	3,3,3	0.57	0	2,2,2	0.03	0
2	EDO	A	406	-	3,3,3	0.47	0	2,2,2	0.24	0
2	EDO	B	434	-	3,3,3	0.73	0	2,2,2	0.20	0
2	EDO	B	437	-	3,3,3	0.52	0	2,2,2	0.13	0
2	EDO	D	426	-	3,3,3	0.56	0	2,2,2	0.12	0
2	EDO	C	405	-	3,3,3	0.53	0	2,2,2	0.25	0
2	EDO	C	418	-	3,3,3	0.47	0	2,2,2	0.33	0
2	EDO	E	448	-	3,3,3	0.26	0	2,2,2	0.84	0
2	EDO	B	410	-	3,3,3	0.51	0	2,2,2	0.73	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	EDO	F	405	-	-	0/1/1/1	-
2	EDO	F	428	-	-	0/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	EDO	E	434	-	-	0/1/1/1	-
2	EDO	B	447	-	-	1/1/1/1	-
2	EDO	F	444	-	-	0/1/1/1	-
2	EDO	E	445	-	-	1/1/1/1	-
2	EDO	A	440	-	-	0/1/1/1	-
2	EDO	B	408	-	-	1/1/1/1	-
2	EDO	F	423	-	-	1/1/1/1	-
2	EDO	B	443	-	-	1/1/1/1	-
2	EDO	A	438	-	-	1/1/1/1	-
2	EDO	A	408	-	-	0/1/1/1	-
2	EDO	C	419	-	-	0/1/1/1	-
2	EDO	A	416	-	-	1/1/1/1	-
2	EDO	A	423	-	-	1/1/1/1	-
2	EDO	E	402	-	-	1/1/1/1	-
2	EDO	C	436	-	-	0/1/1/1	-
2	EDO	E	426	-	-	0/1/1/1	-
2	EDO	F	419	-	-	0/1/1/1	-
2	EDO	A	424	-	-	1/1/1/1	-
2	EDO	D	433	-	-	1/1/1/1	-
2	EDO	B	445	-	-	1/1/1/1	-
2	EDO	E	431	-	-	0/1/1/1	-
2	EDO	F	413	-	-	0/1/1/1	-
2	EDO	A	413	-	-	0/1/1/1	-
2	EDO	D	417	-	-	0/1/1/1	-
2	EDO	E	432	-	-	1/1/1/1	-
2	EDO	A	418	-	-	1/1/1/1	-
2	EDO	B	435	-	-	0/1/1/1	-
2	EDO	C	408	-	-	1/1/1/1	-
2	EDO	C	438	-	-	0/1/1/1	-
2	EDO	E	403	-	-	1/1/1/1	-
2	EDO	E	436	-	-	0/1/1/1	-
2	EDO	E	437	-	-	1/1/1/1	-
2	EDO	C	424	-	-	1/1/1/1	-
2	EDO	E	447	-	-	1/1/1/1	-
2	EDO	A	412	-	-	0/1/1/1	-
2	EDO	C	413	-	-	1/1/1/1	-
2	EDO	B	430	-	-	0/1/1/1	-
2	EDO	E	420	-	-	1/1/1/1	-
2	EDO	D	410	-	-	0/1/1/1	-
2	EDO	E	413	-	-	0/1/1/1	-
2	EDO	E	438	-	-	1/1/1/1	-
2	EDO	E	418	-	-	1/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	EDO	F	414	-	-	1/1/1/1	-
2	EDO	A	427	-	-	1/1/1/1	-
2	EDO	C	410	-	-	0/1/1/1	-
2	EDO	F	436	-	-	1/1/1/1	-
2	EDO	A	431	-	-	1/1/1/1	-
2	EDO	C	430	-	-	0/1/1/1	-
2	EDO	F	416	-	-	0/1/1/1	-
2	EDO	D	439	-	-	1/1/1/1	-
2	EDO	E	446	-	-	1/1/1/1	-
2	EDO	F	422	-	-	0/1/1/1	-
2	EDO	B	428	-	-	0/1/1/1	-
2	EDO	A	422	-	-	0/1/1/1	-
2	EDO	E	444	-	-	1/1/1/1	-
2	EDO	B	444	-	-	1/1/1/1	-
2	EDO	A	409	-	-	0/1/1/1	-
2	EDO	C	416	-	-	1/1/1/1	-
2	EDO	C	431	-	-	1/1/1/1	-
2	EDO	A	429	-	-	0/1/1/1	-
2	EDO	F	438	-	-	1/1/1/1	-
2	EDO	D	422	-	-	0/1/1/1	-
2	EDO	A	437	-	-	0/1/1/1	-
2	EDO	C	422	-	-	1/1/1/1	-
2	EDO	C	432	-	-	1/1/1/1	-
2	EDO	D	428	-	-	1/1/1/1	-
2	EDO	C	425	-	-	1/1/1/1	-
2	EDO	C	429	-	-	1/1/1/1	-
2	EDO	C	437	-	-	1/1/1/1	-
2	EDO	C	412	-	-	0/1/1/1	-
2	EDO	B	404	-	-	1/1/1/1	-
2	EDO	B	431	-	-	1/1/1/1	-
2	EDO	F	442	-	-	1/1/1/1	-
2	EDO	B	438	-	-	1/1/1/1	-
2	EDO	B	422	-	-	0/1/1/1	-
2	EDO	F	410	-	-	0/1/1/1	-
2	EDO	F	441	-	-	0/1/1/1	-
2	EDO	A	404	-	-	0/1/1/1	-
2	EDO	A	410	-	-	0/1/1/1	-
2	EDO	B	409	-	-	0/1/1/1	-
2	EDO	B	423	-	-	1/1/1/1	-
2	EDO	D	416	-	-	0/1/1/1	-
2	EDO	D	411	-	-	1/1/1/1	-
2	EDO	E	401	-	-	1/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	EDO	D	406	-	-	0/1/1/1	-
2	EDO	B	442	-	-	1/1/1/1	-
2	EDO	C	427	-	-	1/1/1/1	-
2	EDO	D	404	-	-	0/1/1/1	-
2	EDO	E	435	-	-	0/1/1/1	-
2	EDO	B	418	-	-	1/1/1/1	-
2	EDO	C	404	-	-	1/1/1/1	-
2	EDO	F	408	-	-	0/1/1/1	-
2	EDO	A	426	-	-	0/1/1/1	-
2	EDO	F	439	-	-	0/1/1/1	-
2	EDO	A	402	-	-	1/1/1/1	-
2	EDO	D	438	-	-	0/1/1/1	-
2	EDO	E	414	-	-	1/1/1/1	-
2	EDO	A	443	-	-	1/1/1/1	-
2	EDO	F	404	-	-	0/1/1/1	-
2	EDO	D	423	-	-	0/1/1/1	-
2	EDO	F	418	-	-	0/1/1/1	-
2	EDO	F	443	-	-	1/1/1/1	-
2	EDO	F	432	-	-	1/1/1/1	-
2	EDO	A	421	-	-	1/1/1/1	-
2	EDO	A	430	-	-	1/1/1/1	-
2	EDO	A	432	-	-	1/1/1/1	-
2	EDO	F	403	-	-	1/1/1/1	-
2	EDO	E	440	-	-	1/1/1/1	-
2	EDO	D	402	-	-	0/1/1/1	-
2	EDO	E	411	-	-	1/1/1/1	-
2	EDO	E	415	-	-	0/1/1/1	-
2	EDO	C	402	-	-	0/1/1/1	-
2	EDO	F	412	-	-	0/1/1/1	-
2	EDO	B	439	-	-	1/1/1/1	-
2	EDO	B	424	-	-	0/1/1/1	-
2	EDO	A	417	-	-	1/1/1/1	-
2	EDO	F	417	-	-	0/1/1/1	-
2	EDO	B	413	-	-	0/1/1/1	-
2	EDO	F	401	-	-	0/1/1/1	-
2	EDO	B	433	-	-	1/1/1/1	-
2	EDO	F	420	-	-	0/1/1/1	-
2	EDO	A	439	-	-	0/1/1/1	-
2	EDO	D	427	-	-	1/1/1/1	-
2	EDO	E	433	-	-	0/1/1/1	-
2	EDO	F	434	-	-	0/1/1/1	-
2	EDO	B	426	-	-	1/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	EDO	B	402	-	-	0/1/1/1	-
2	EDO	E	441	-	-	1/1/1/1	-
2	EDO	C	421	-	-	0/1/1/1	-
2	EDO	A	403	-	-	0/1/1/1	-
2	EDO	D	425	-	-	0/1/1/1	-
2	EDO	E	409	-	-	0/1/1/1	-
2	EDO	F	425	-	-	1/1/1/1	-
2	EDO	B	432	-	-	0/1/1/1	-
2	EDO	D	429	-	-	0/1/1/1	-
2	EDO	E	408	-	-	1/1/1/1	-
2	EDO	E	439	-	-	1/1/1/1	-
2	EDO	F	402	-	-	1/1/1/1	-
2	EDO	B	403	-	-	1/1/1/1	-
2	EDO	D	436	-	-	0/1/1/1	-
2	EDO	C	423	-	-	1/1/1/1	-
2	EDO	E	427	-	-	1/1/1/1	-
2	EDO	D	403	-	-	1/1/1/1	-
2	EDO	E	442	-	-	0/1/1/1	-
2	EDO	B	427	-	-	1/1/1/1	-
2	EDO	B	412	-	-	0/1/1/1	-
2	EDO	E	443	-	-	1/1/1/1	-
2	EDO	F	424	-	-	0/1/1/1	-
2	EDO	C	426	-	-	1/1/1/1	-
2	EDO	A	435	-	-	0/1/1/1	-
2	EDO	E	419	-	-	1/1/1/1	-
2	EDO	B	420	-	-	0/1/1/1	-
2	EDO	D	407	-	-	0/1/1/1	-
2	EDO	E	425	-	-	1/1/1/1	-
2	EDO	D	413	-	-	0/1/1/1	-
2	EDO	A	420	-	-	0/1/1/1	-
2	EDO	F	421	-	-	0/1/1/1	-
2	EDO	B	429	-	-	1/1/1/1	-
2	EDO	A	442	-	-	1/1/1/1	-
2	EDO	A	414	-	-	0/1/1/1	-
2	EDO	B	415	-	-	0/1/1/1	-
2	EDO	D	434	-	-	0/1/1/1	-
2	EDO	F	415	-	-	0/1/1/1	-
2	EDO	B	446	-	-	0/1/1/1	-
2	EDO	B	416	-	-	1/1/1/1	-
2	EDO	A	415	-	-	1/1/1/1	-
2	EDO	E	416	-	-	1/1/1/1	-
2	EDO	D	408	-	-	0/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	EDO	D	430	-	-	1/1/1/1	-
2	EDO	E	428	-	-	1/1/1/1	-
2	EDO	D	424	-	-	1/1/1/1	-
2	EDO	C	403	-	-	0/1/1/1	-
2	EDO	E	404	-	-	1/1/1/1	-
2	EDO	C	411	-	-	0/1/1/1	-
2	EDO	C	417	-	-	1/1/1/1	-
2	EDO	B	441	-	-	1/1/1/1	-
2	EDO	C	409	-	-	1/1/1/1	-
2	EDO	B	419	-	-	1/1/1/1	-
2	EDO	B	405	-	-	1/1/1/1	-
2	EDO	A	441	-	-	1/1/1/1	-
2	EDO	B	414	-	-	0/1/1/1	-
2	EDO	F	409	-	-	0/1/1/1	-
2	EDO	A	419	-	-	0/1/1/1	-
2	EDO	E	424	-	-	0/1/1/1	-
2	EDO	B	407	-	-	0/1/1/1	-
2	EDO	F	431	-	-	1/1/1/1	-
2	EDO	F	437	-	-	1/1/1/1	-
2	EDO	E	406	-	-	0/1/1/1	-
2	EDO	F	407	-	-	0/1/1/1	-
2	EDO	C	420	-	-	0/1/1/1	-
2	EDO	A	407	-	-	0/1/1/1	-
2	EDO	D	415	-	-	0/1/1/1	-
2	EDO	E	412	-	-	1/1/1/1	-
2	EDO	F	433	-	-	0/1/1/1	-
2	EDO	B	401	-	-	0/1/1/1	-
2	EDO	B	417	-	-	1/1/1/1	-
2	EDO	C	414	-	-	0/1/1/1	-
2	EDO	A	433	-	-	0/1/1/1	-
2	EDO	A	401	-	-	0/1/1/1	-
2	EDO	C	415	-	-	1/1/1/1	-
2	EDO	D	432	-	-	0/1/1/1	-
2	EDO	F	430	-	-	1/1/1/1	-
2	EDO	F	411	-	-	0/1/1/1	-
2	EDO	B	406	-	-	0/1/1/1	-
2	EDO	A	411	-	-	1/1/1/1	-
2	EDO	F	406	-	-	1/1/1/1	-
2	EDO	B	440	-	-	1/1/1/1	-
2	EDO	F	445	-	-	0/1/1/1	-
2	EDO	D	401	-	-	0/1/1/1	-
2	EDO	D	412	-	-	0/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	EDO	E	417	-	-	0/1/1/1	-
2	EDO	D	420	-	-	1/1/1/1	-
2	EDO	E	423	-	-	0/1/1/1	-
2	EDO	D	414	-	-	1/1/1/1	-
2	EDO	A	425	-	-	1/1/1/1	-
2	EDO	E	429	-	-	1/1/1/1	-
2	EDO	A	434	-	-	1/1/1/1	-
2	EDO	E	410	-	-	0/1/1/1	-
2	EDO	B	411	-	-	0/1/1/1	-
2	EDO	C	407	-	-	0/1/1/1	-
2	EDO	F	427	-	-	1/1/1/1	-
2	EDO	D	409	-	-	1/1/1/1	-
2	EDO	C	434	-	-	1/1/1/1	-
2	EDO	B	421	-	-	0/1/1/1	-
2	EDO	D	405	-	-	1/1/1/1	-
2	EDO	F	426	-	-	1/1/1/1	-
2	EDO	C	433	-	-	0/1/1/1	-
2	EDO	D	418	-	-	0/1/1/1	-
2	EDO	F	440	-	-	1/1/1/1	-
2	EDO	D	437	-	-	0/1/1/1	-
2	EDO	E	407	-	-	0/1/1/1	-
2	EDO	A	428	-	-	0/1/1/1	-
2	EDO	C	401	-	-	0/1/1/1	-
2	EDO	B	436	-	-	1/1/1/1	-
2	EDO	C	435	-	-	1/1/1/1	-
2	EDO	D	421	-	-	1/1/1/1	-
2	EDO	F	429	-	-	1/1/1/1	-
2	EDO	F	435	-	-	1/1/1/1	-
2	EDO	D	435	-	-	1/1/1/1	-
2	EDO	B	425	-	-	1/1/1/1	-
2	EDO	A	436	-	-	1/1/1/1	-
2	EDO	E	422	-	-	0/1/1/1	-
2	EDO	C	406	-	-	1/1/1/1	-
2	EDO	D	431	-	-	1/1/1/1	-
2	EDO	E	405	-	-	0/1/1/1	-
2	EDO	D	419	-	-	0/1/1/1	-
2	EDO	E	430	-	-	0/1/1/1	-
2	EDO	C	428	-	-	1/1/1/1	-
2	EDO	E	421	-	-	1/1/1/1	-
2	EDO	A	405	-	-	0/1/1/1	-
2	EDO	A	406	-	-	1/1/1/1	-
2	EDO	B	434	-	-	1/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	EDO	B	437	-	-	1/1/1/1	-
2	EDO	D	426	-	-	1/1/1/1	-
2	EDO	C	405	-	-	0/1/1/1	-
2	EDO	C	418	-	-	1/1/1/1	-
2	EDO	E	448	-	-	1/1/1/1	-
2	EDO	B	410	-	-	0/1/1/1	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	440	EDO	O1-C1	-2.82	1.27	1.42
2	B	440	EDO	O2-C2	2.27	1.53	1.42

There are no bond angle outliers.

There are no chirality outliers.

5 of 134 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	429	EDO	O1-C1-C2-O2
2	E	444	EDO	O1-C1-C2-O2
2	A	406	EDO	O1-C1-C2-O2
2	A	417	EDO	O1-C1-C2-O2
2	A	421	EDO	O1-C1-C2-O2

There are no ring outliers.

34 monomers are involved in 46 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	F	444	EDO	1	0
2	A	408	EDO	2	0
2	C	419	EDO	1	0
2	A	412	EDO	1	0
2	D	410	EDO	1	0
2	E	413	EDO	1	0
2	C	410	EDO	2	0
2	F	436	EDO	1	0
2	F	416	EDO	2	0
2	C	412	EDO	1	0
2	B	404	EDO	3	0
2	F	410	EDO	1	0
2	A	410	EDO	1	0

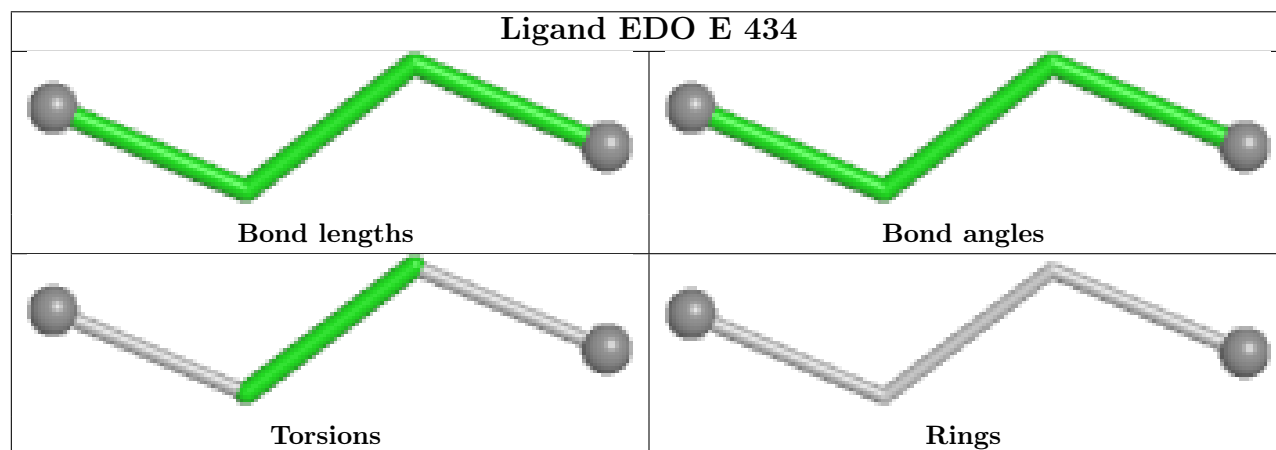
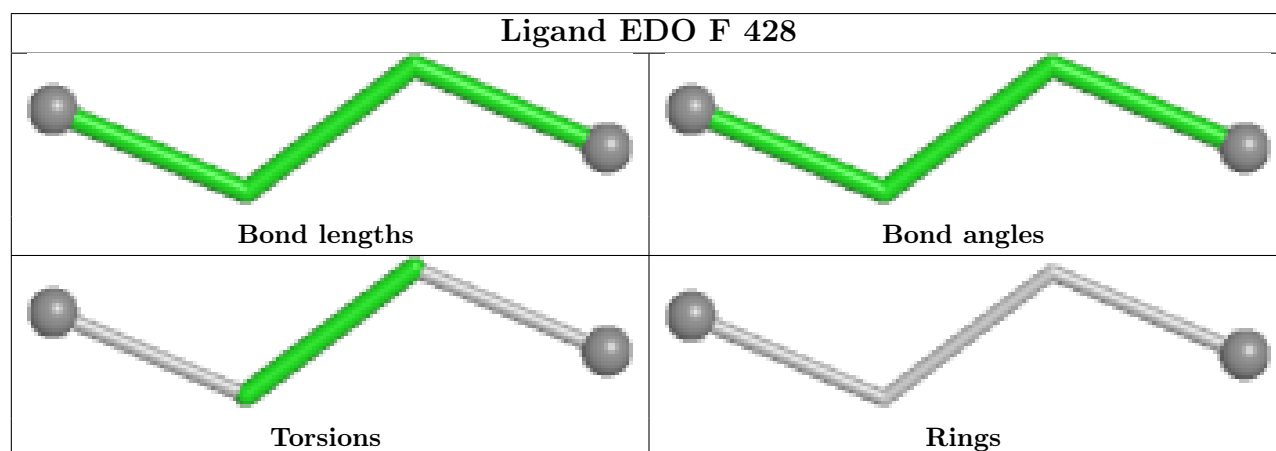
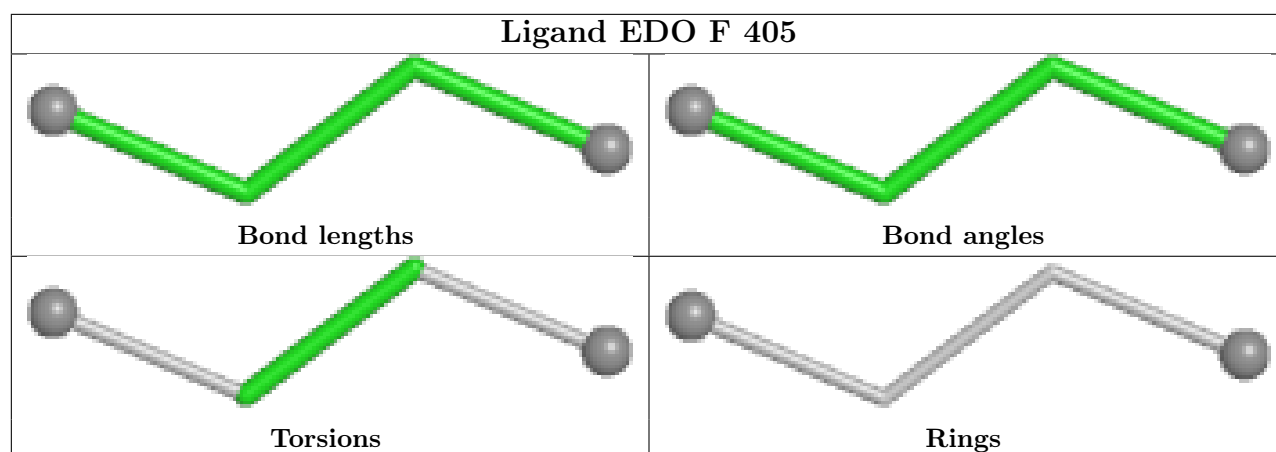
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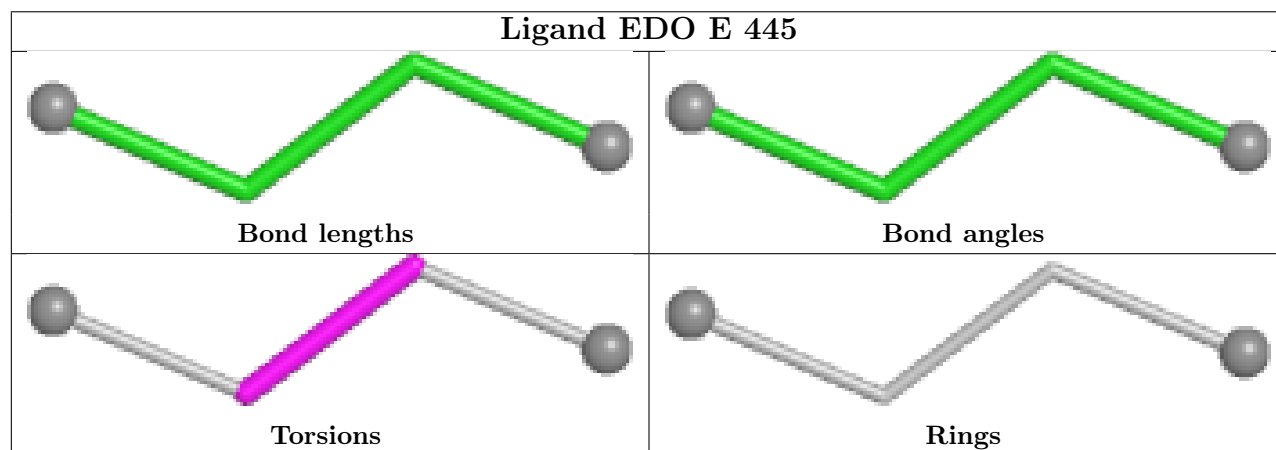
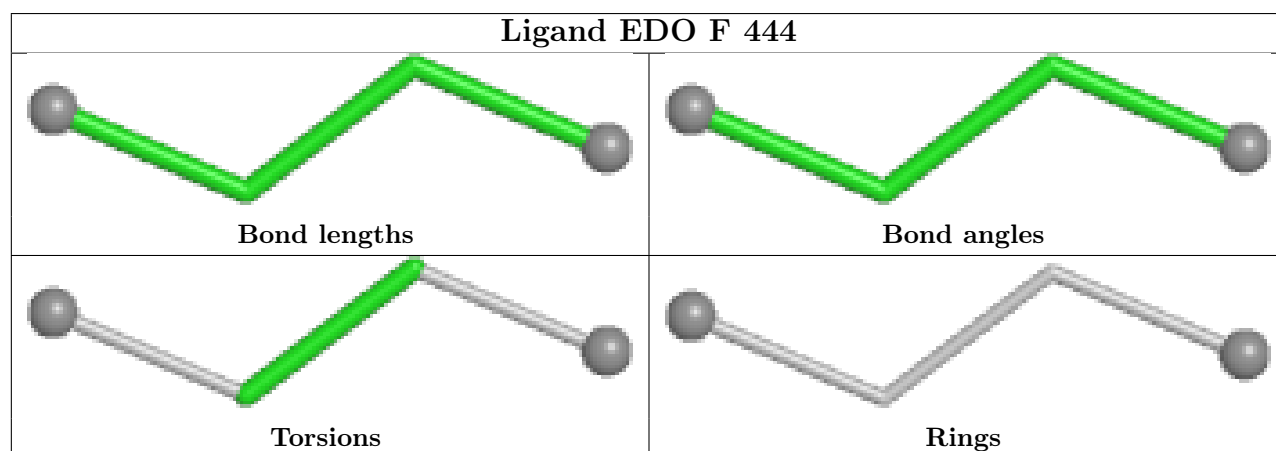
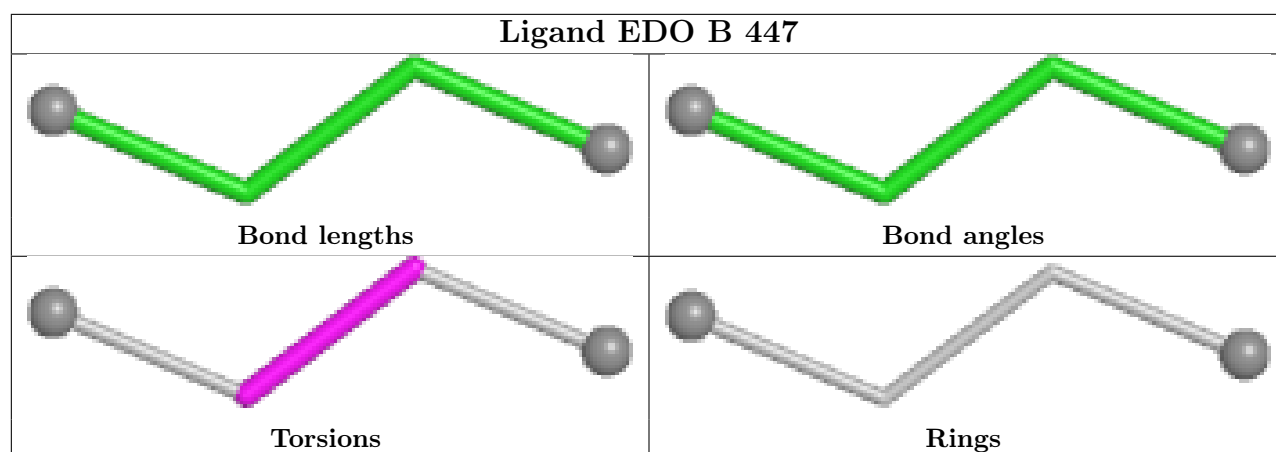
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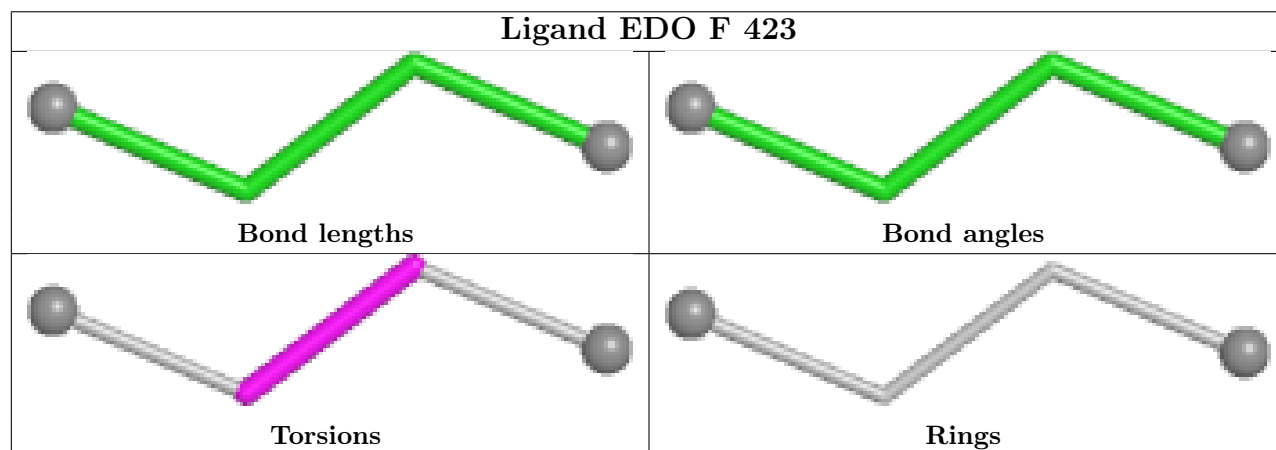
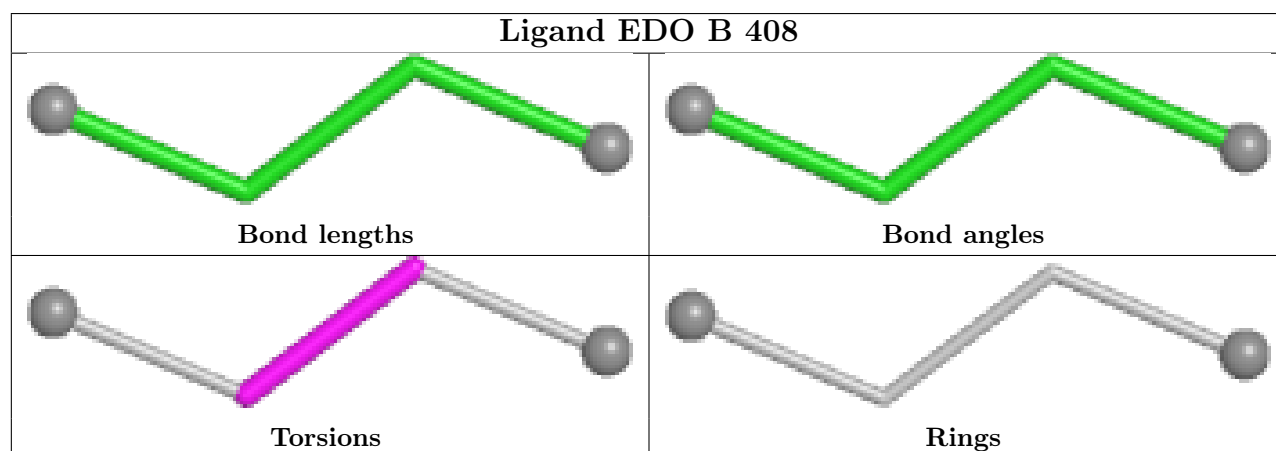
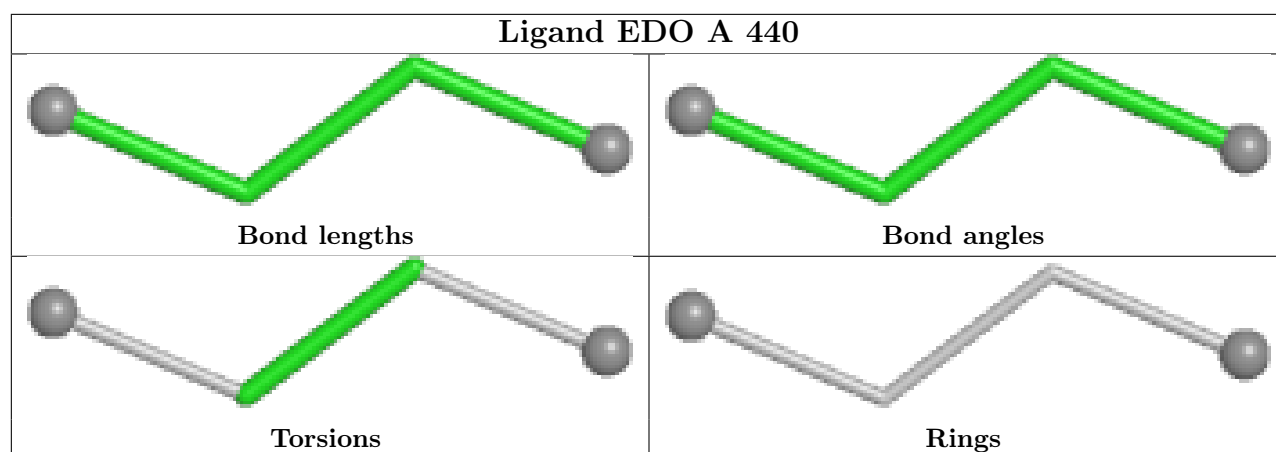
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	409	EDO	1	0
2	D	411	EDO	2	0
2	C	404	EDO	4	0
2	A	402	EDO	3	0
2	A	443	EDO	1	0
2	E	415	EDO	1	0
2	E	409	EDO	1	0
2	B	412	EDO	1	0
2	B	420	EDO	1	0
2	D	413	EDO	2	0
2	F	415	EDO	1	0
2	E	404	EDO	3	0
2	C	409	EDO	1	0
2	B	414	EDO	1	0
2	A	407	EDO	1	0
2	D	415	EDO	1	0
2	C	414	EDO	1	0
2	E	410	EDO	2	0
2	D	405	EDO	4	0
2	E	422	EDO	1	0
2	B	410	EDO	2	0

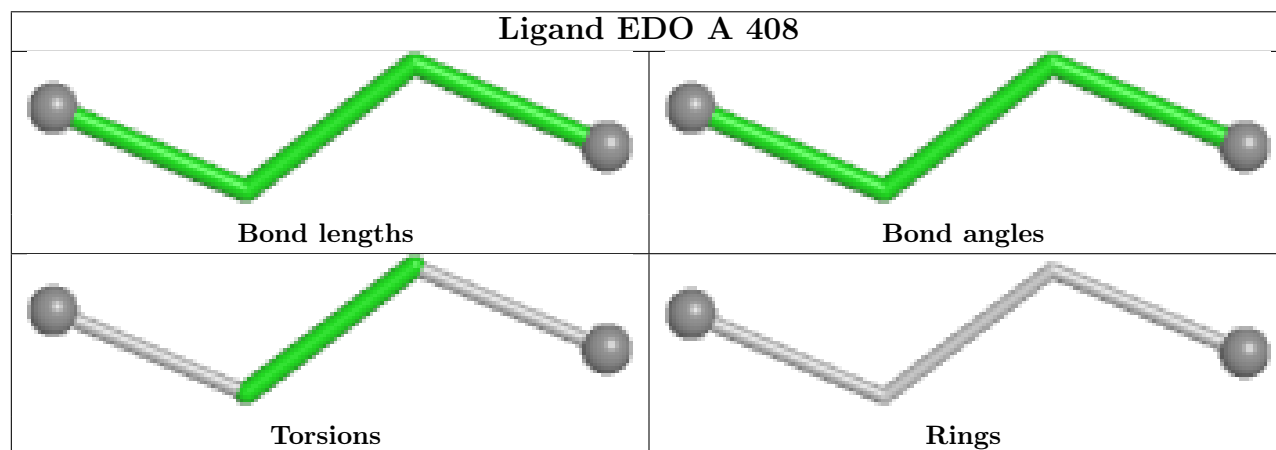
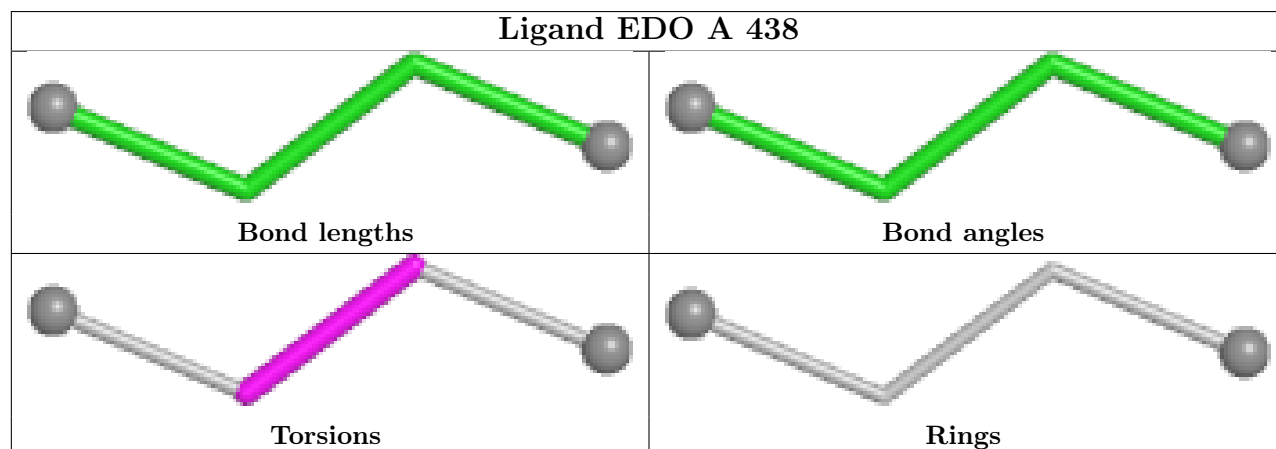
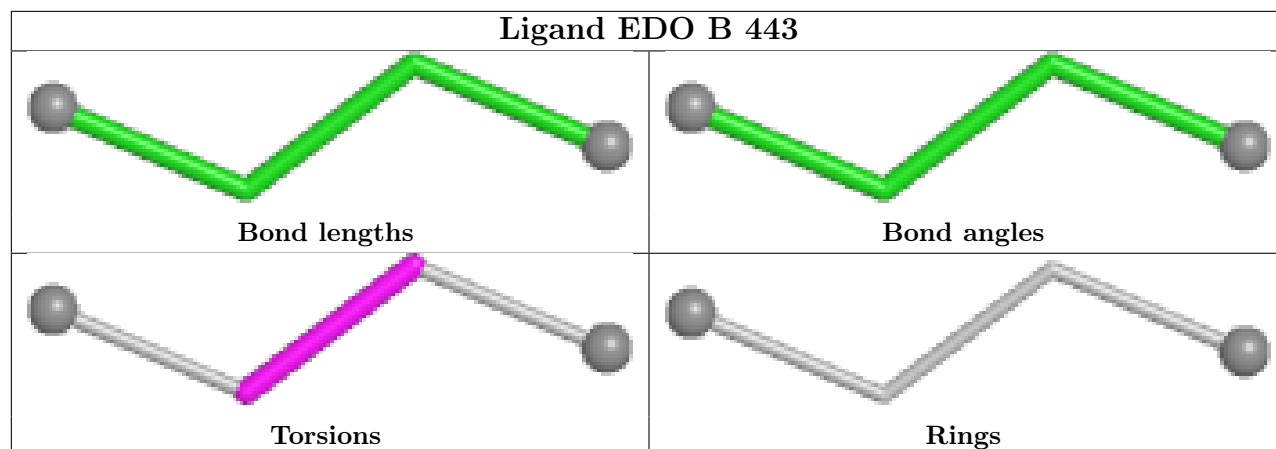
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

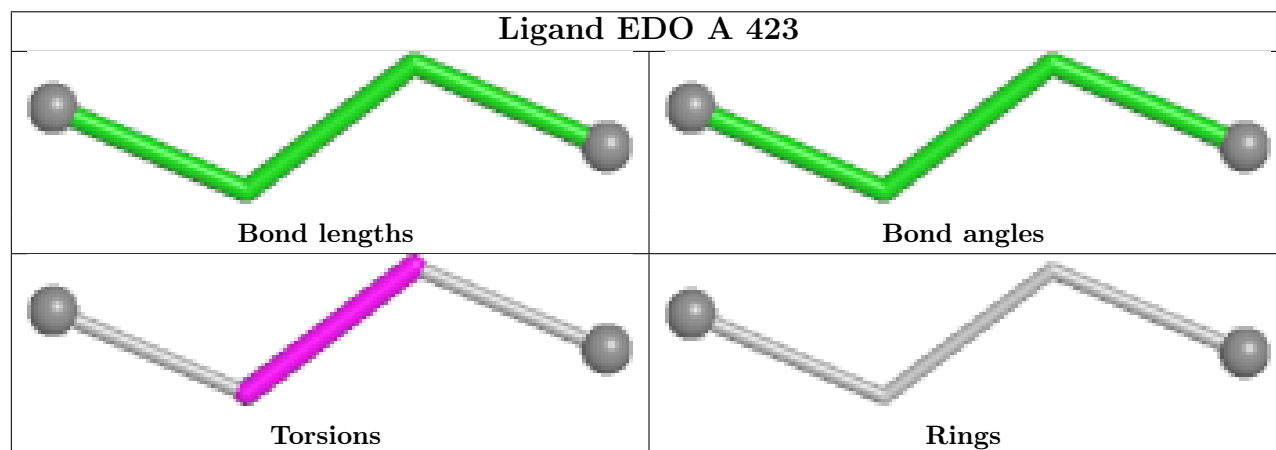
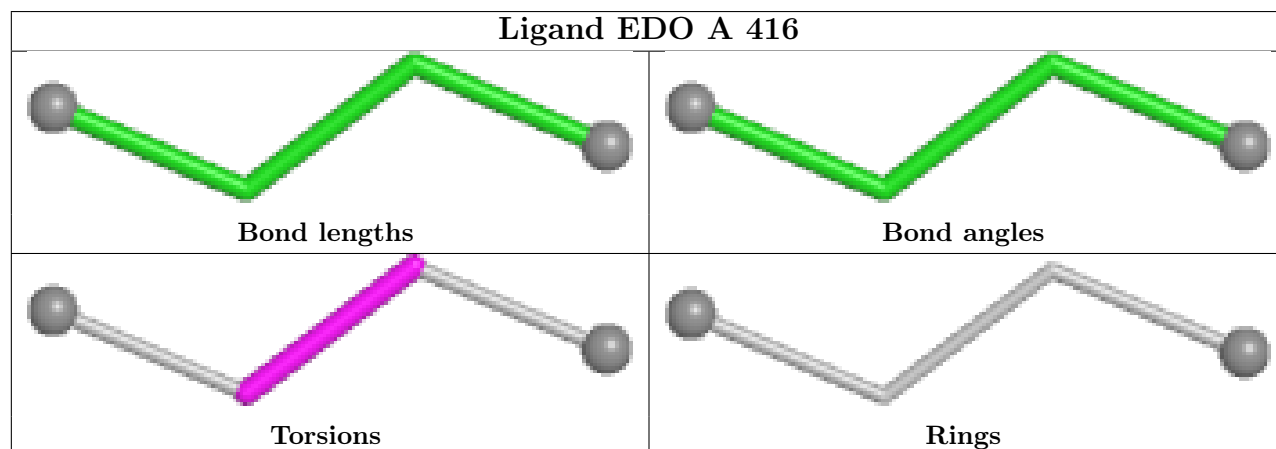
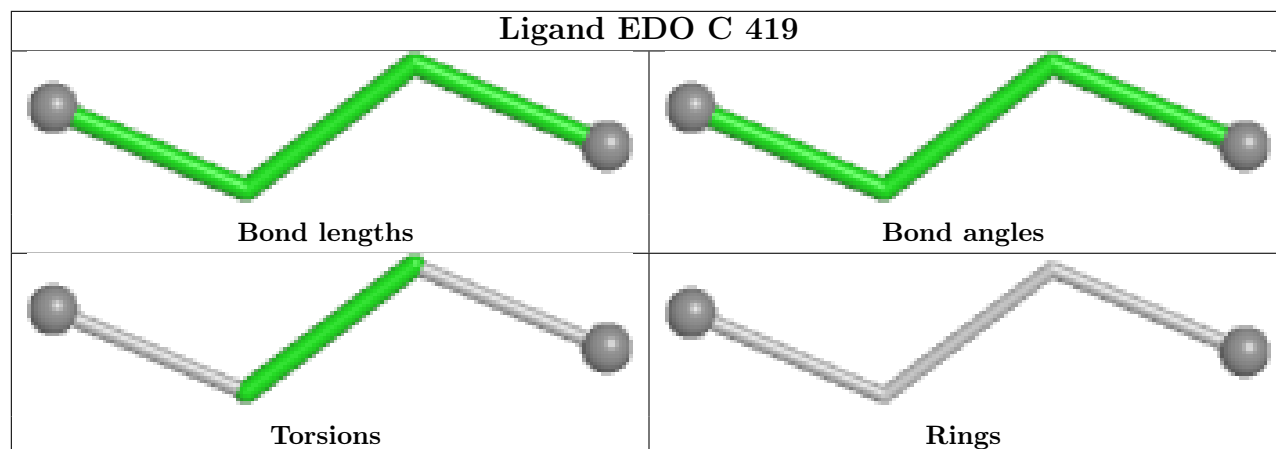


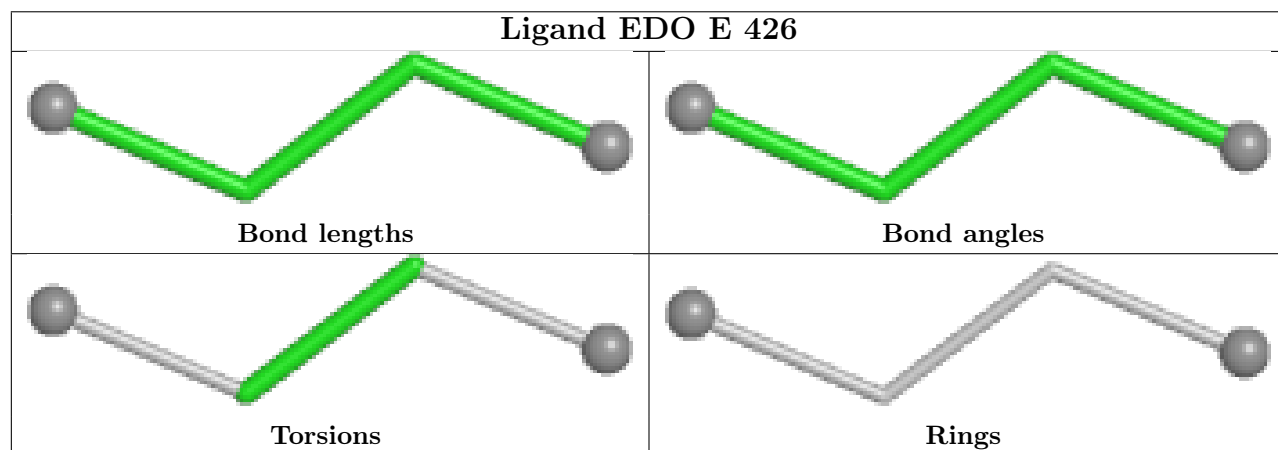
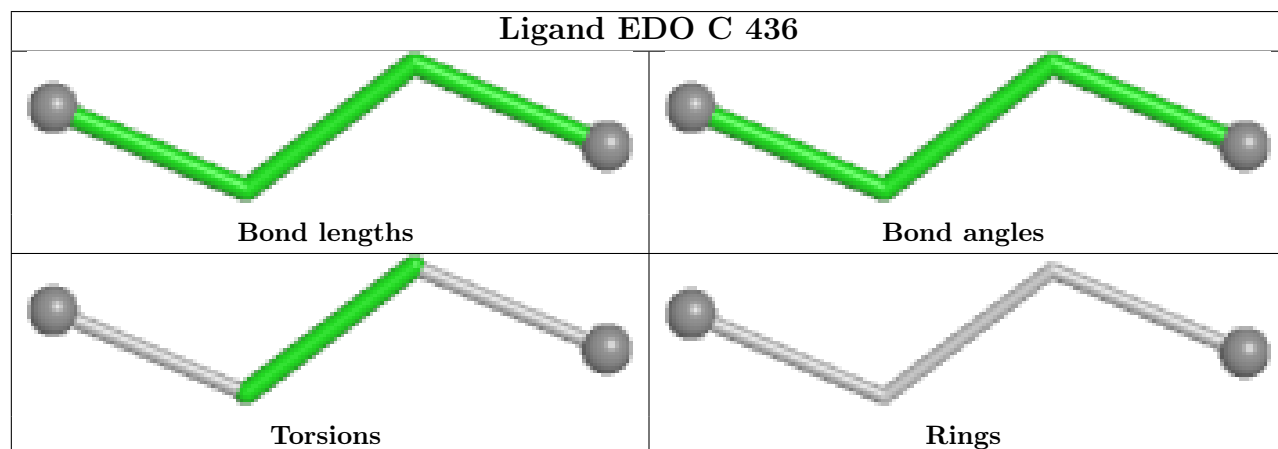
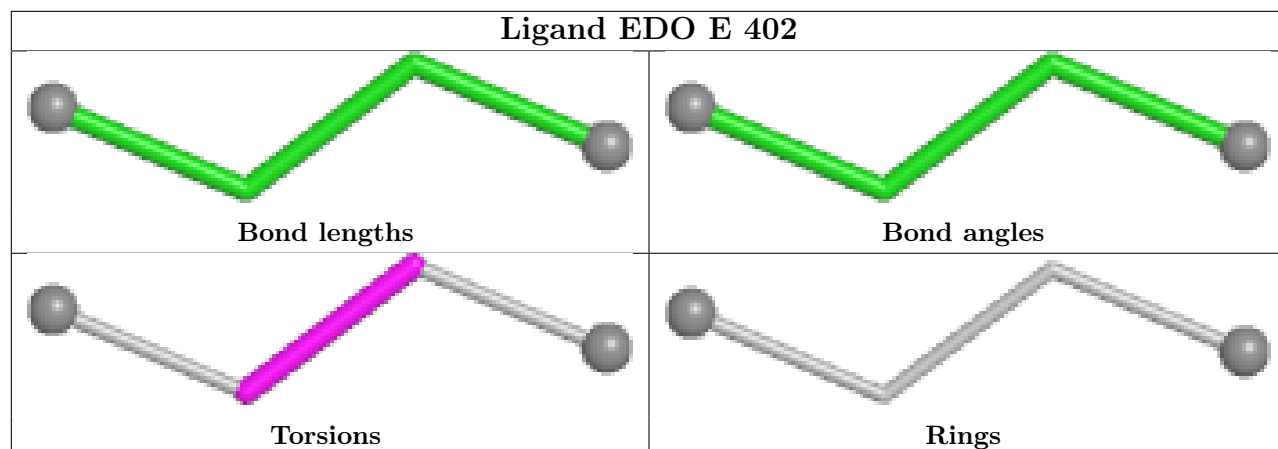


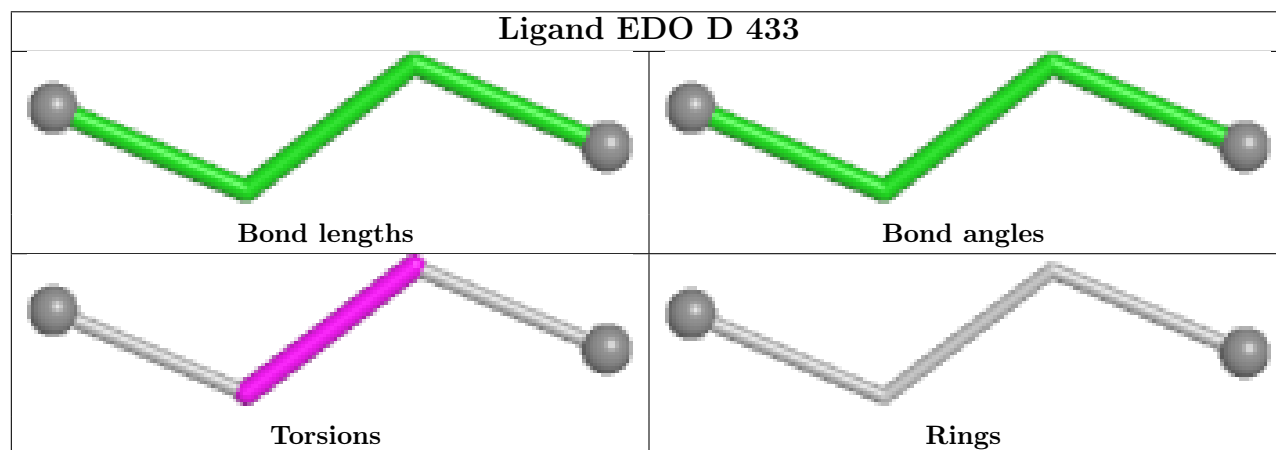
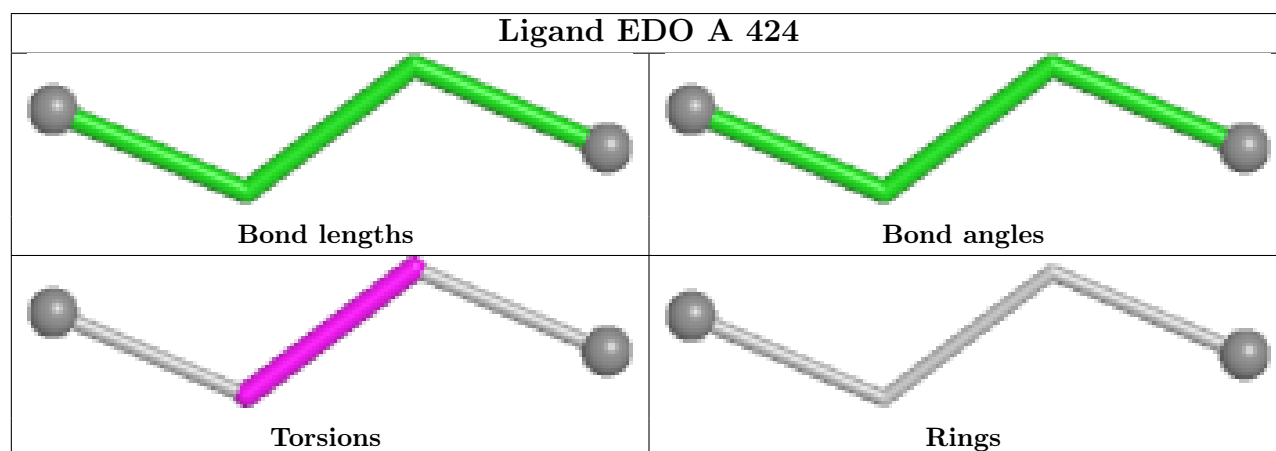
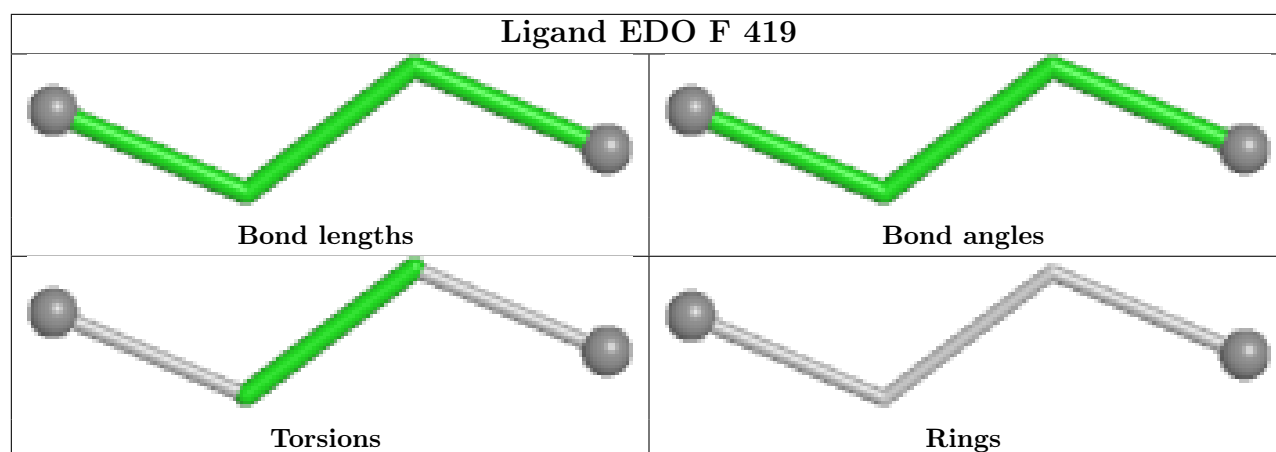


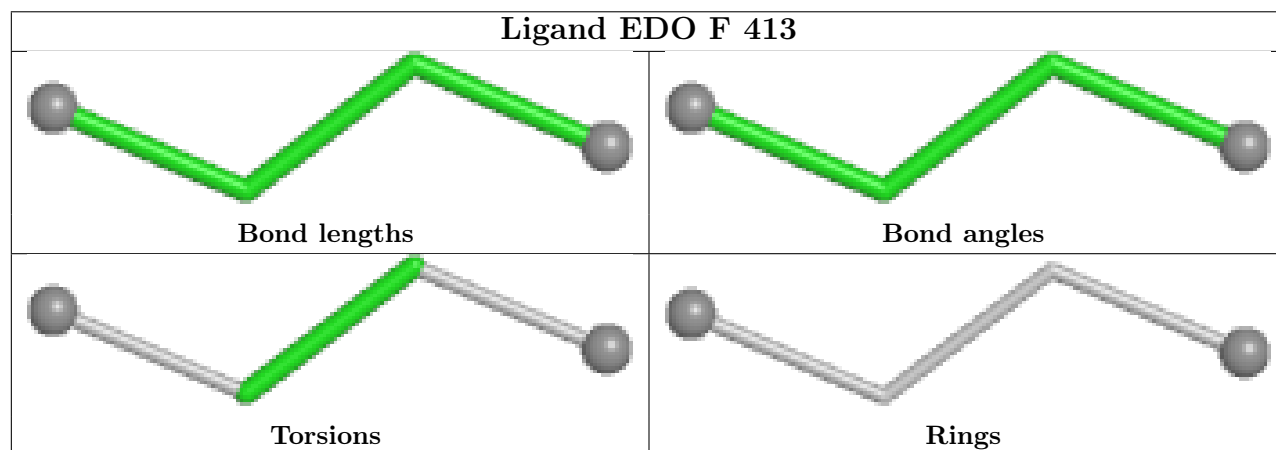
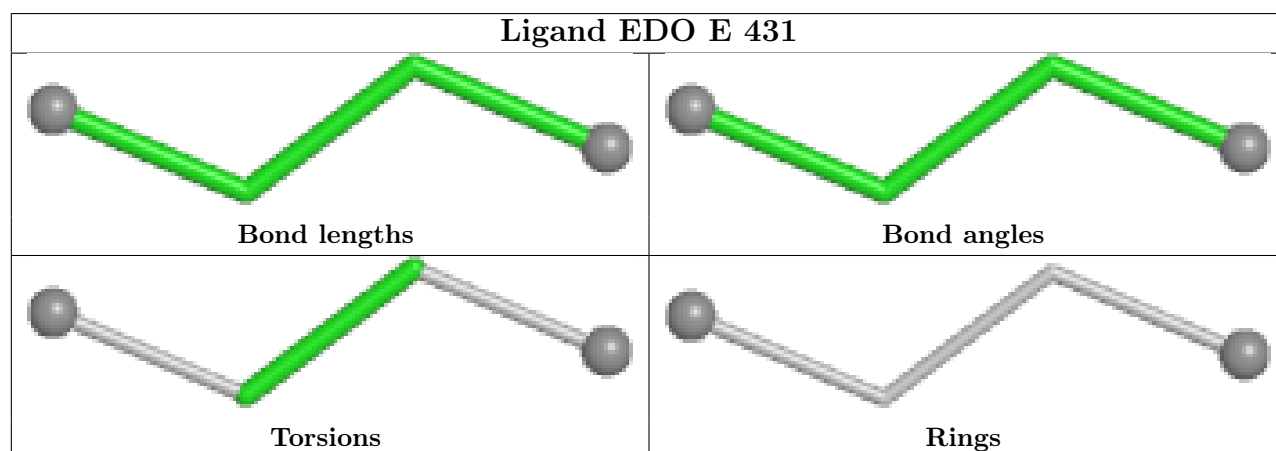
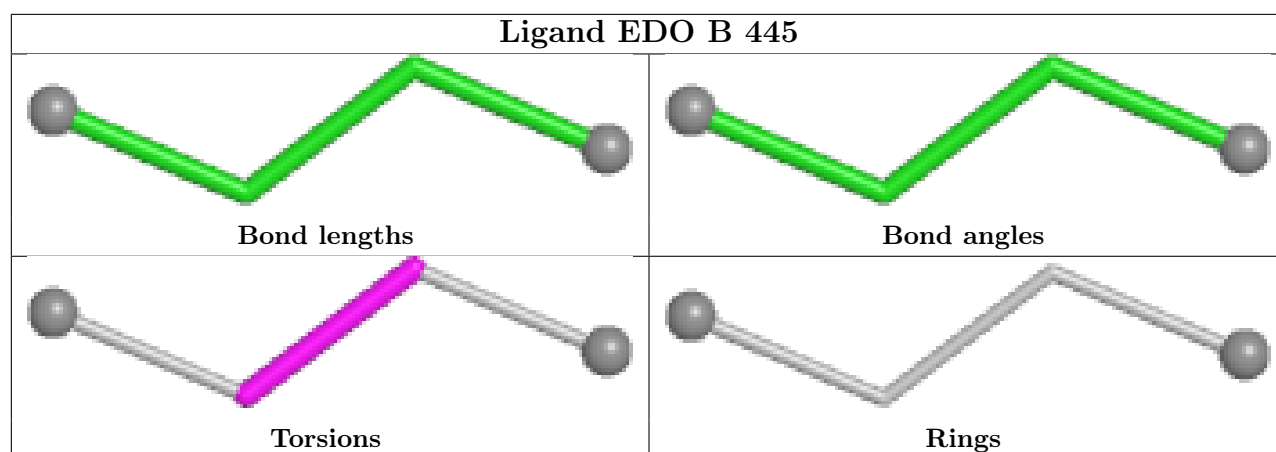




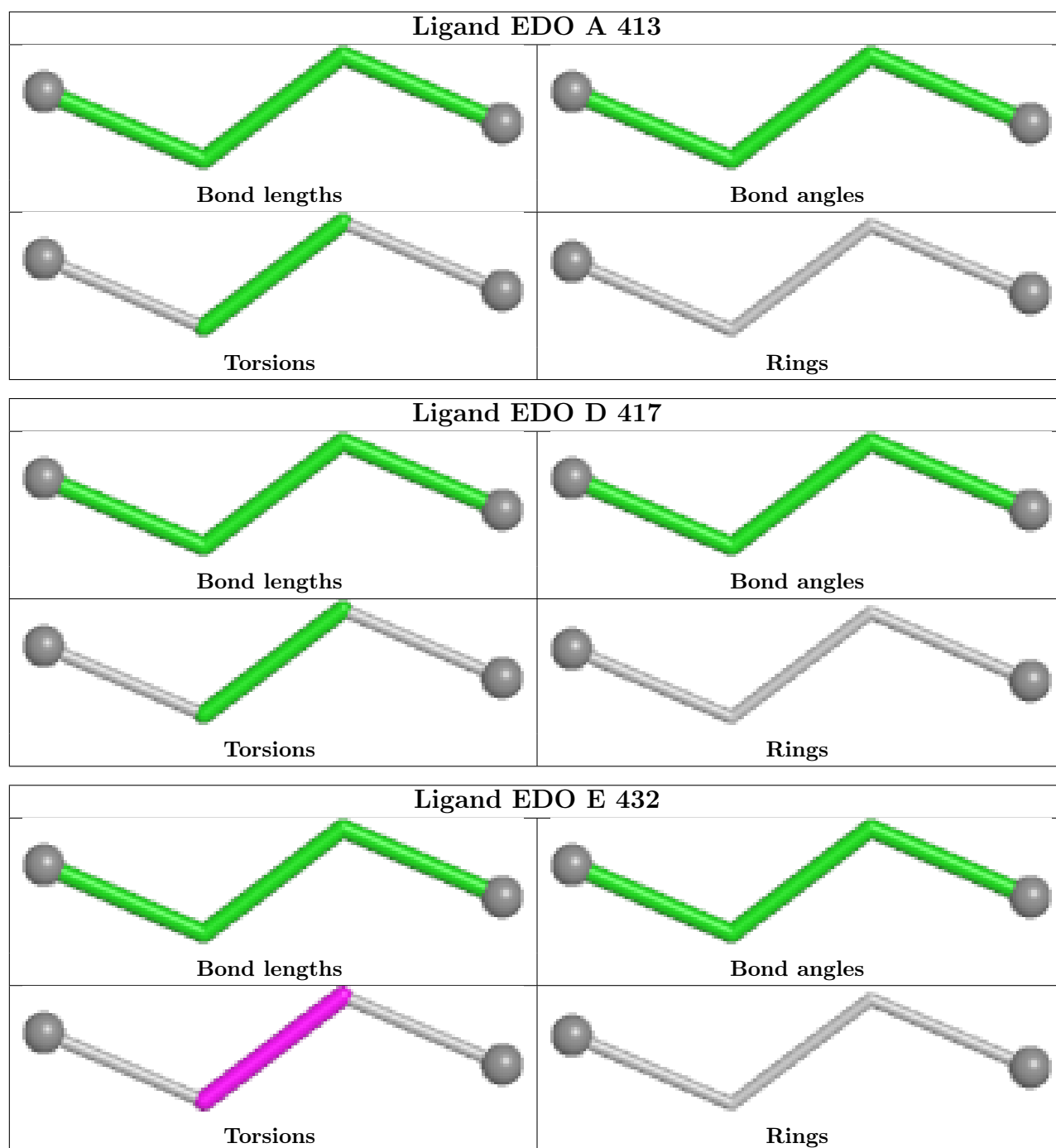


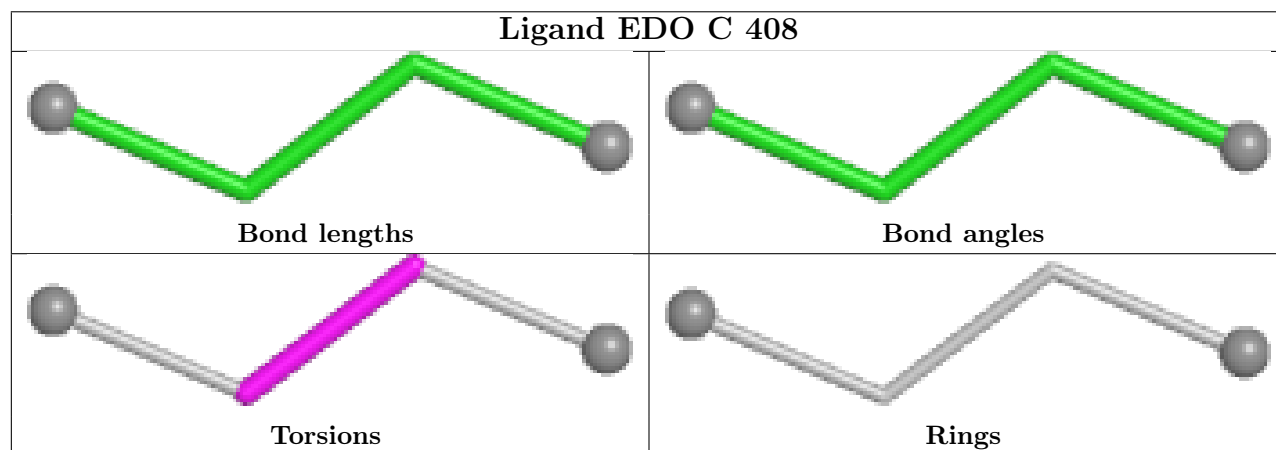
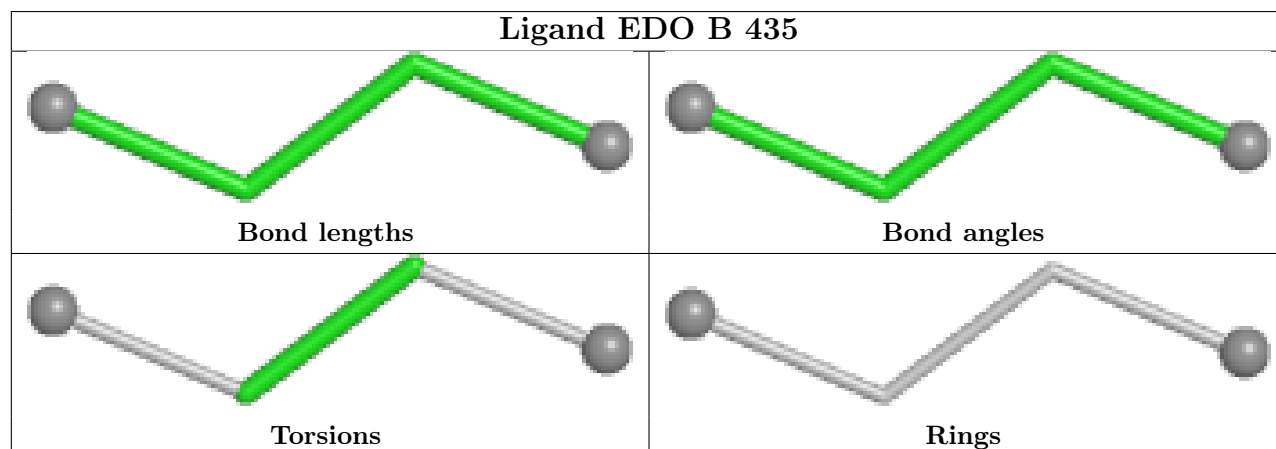
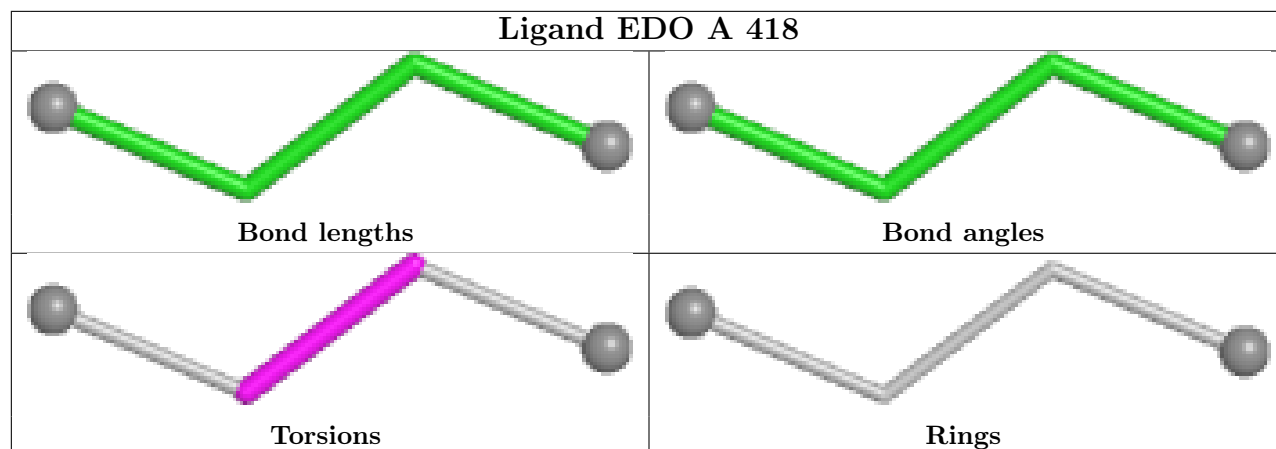


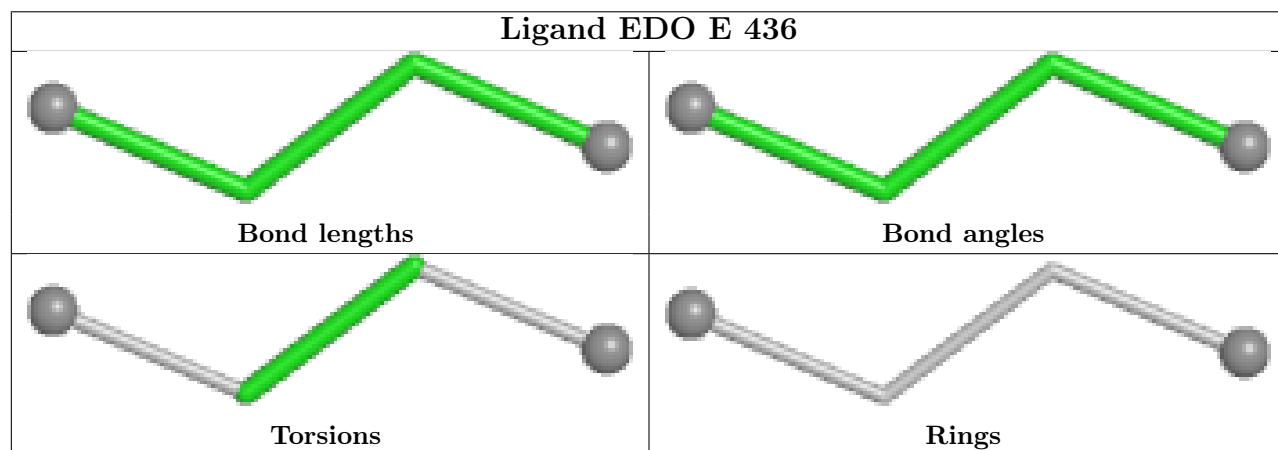
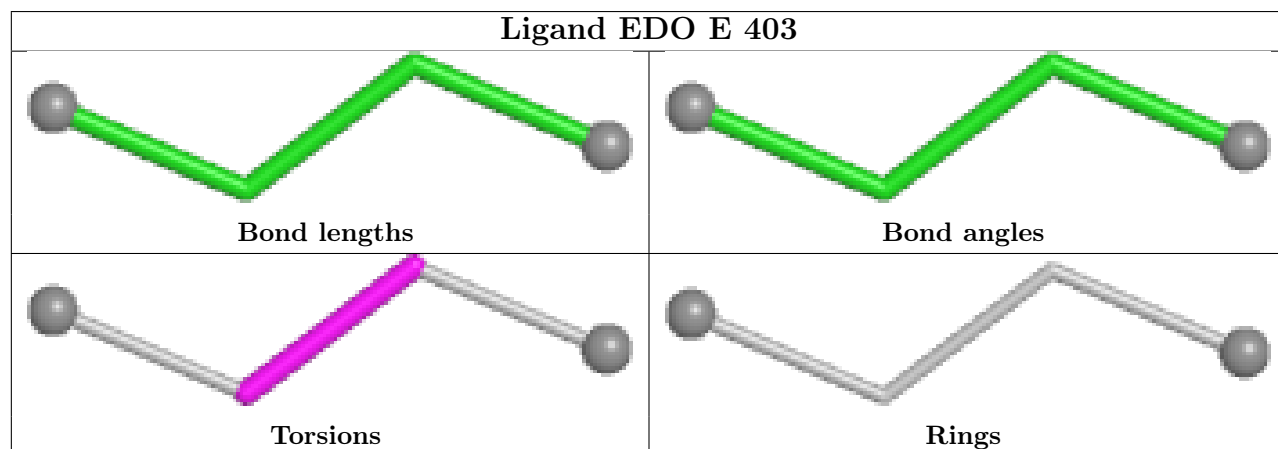
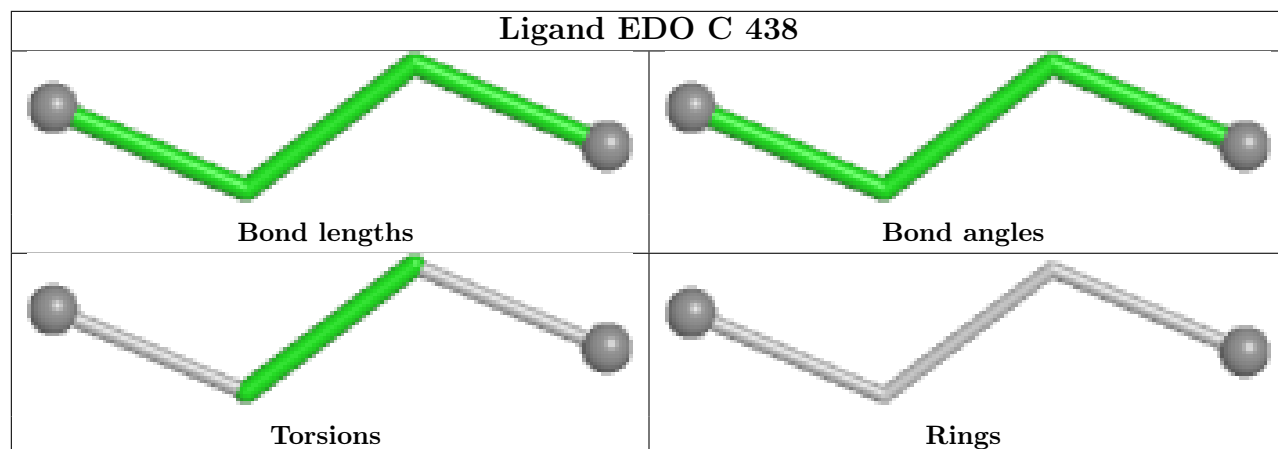


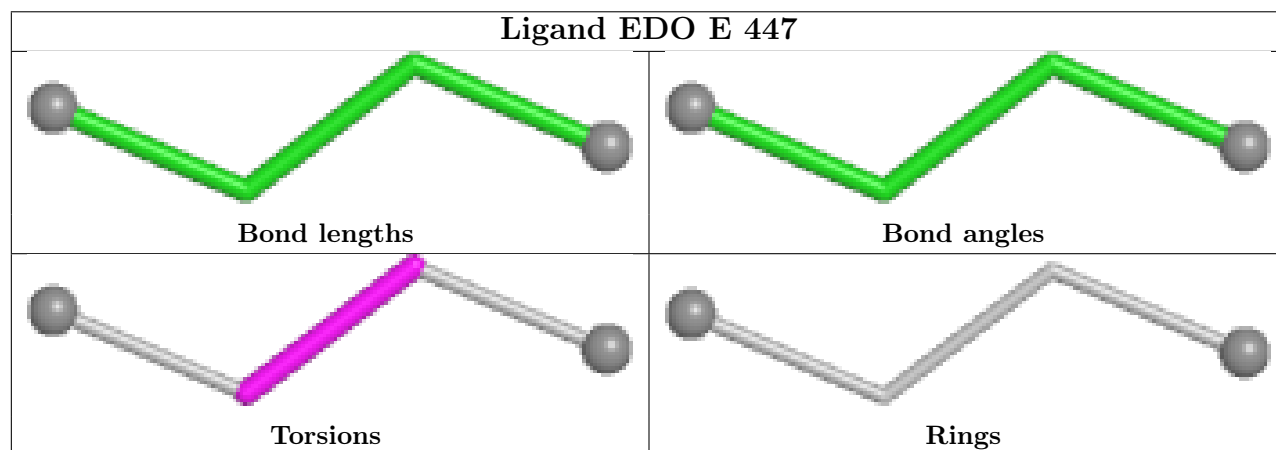
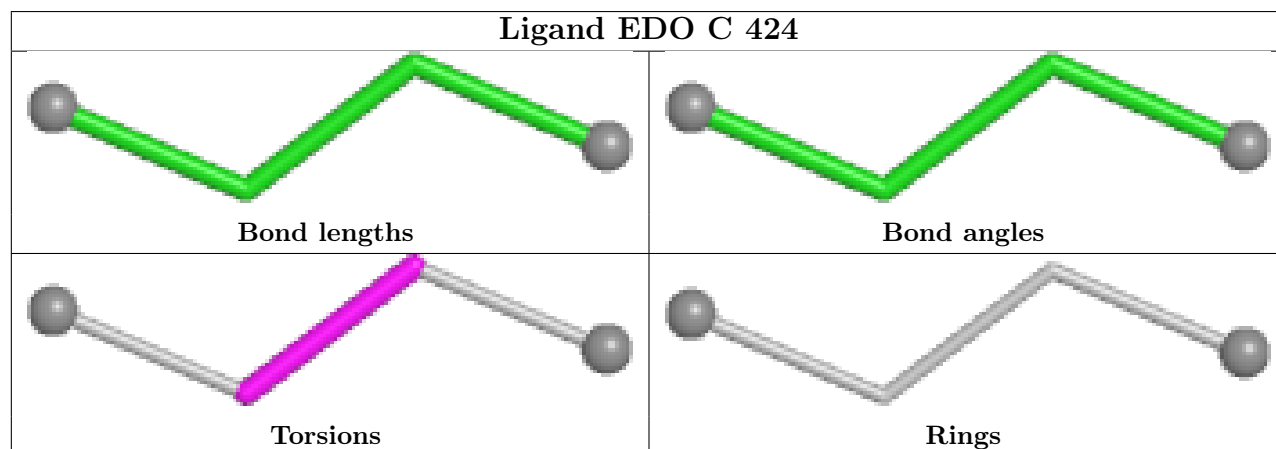
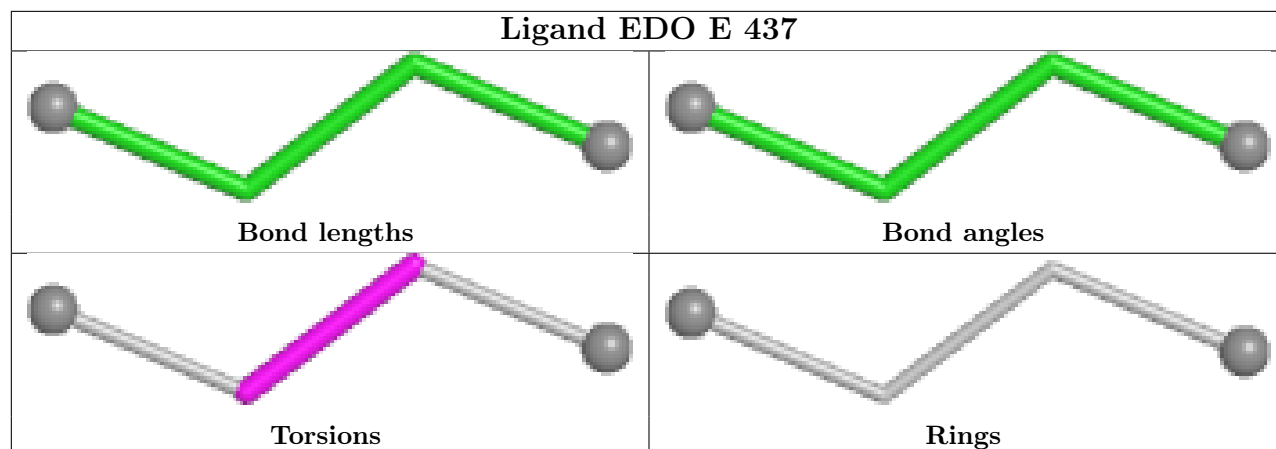


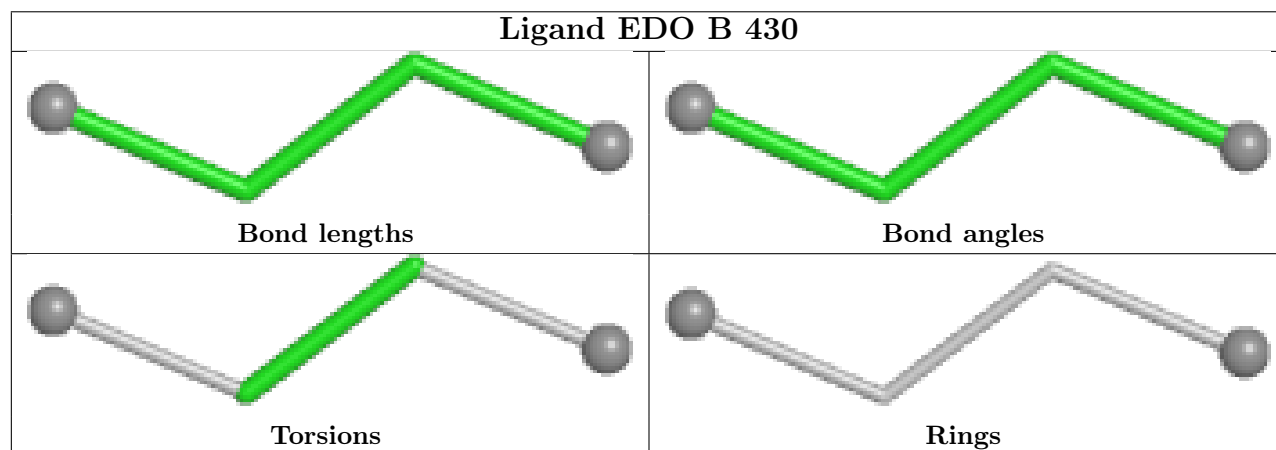
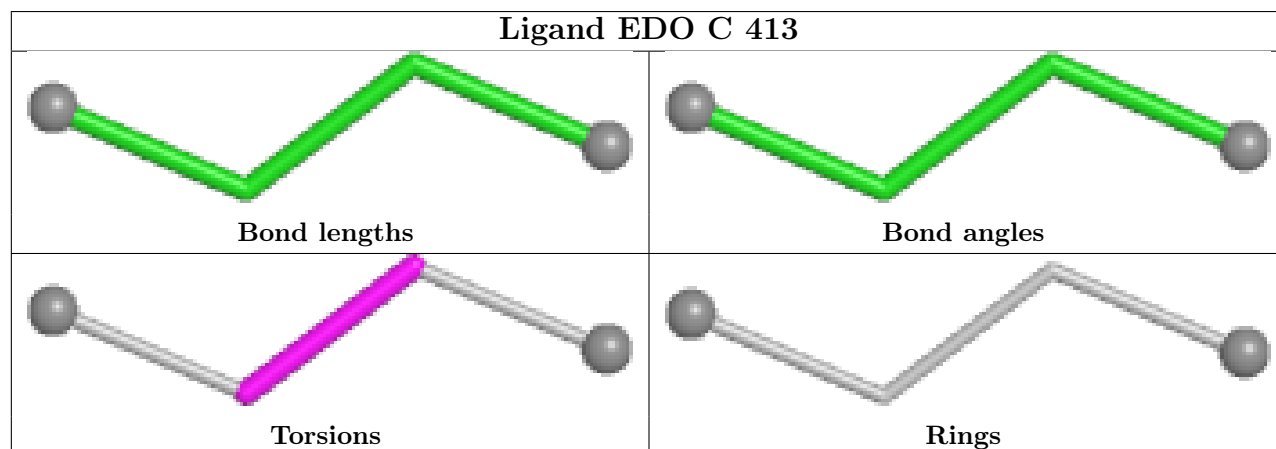
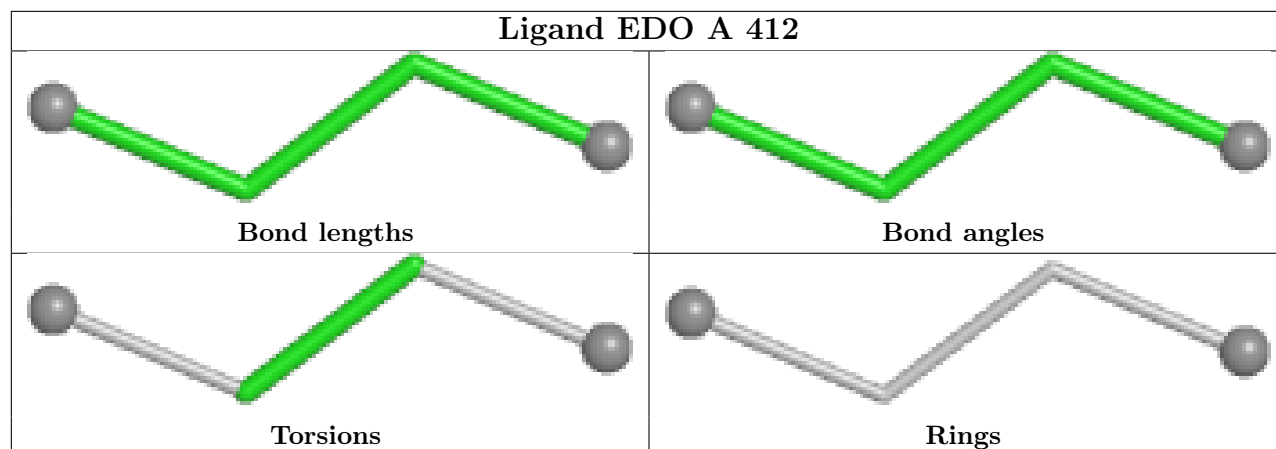


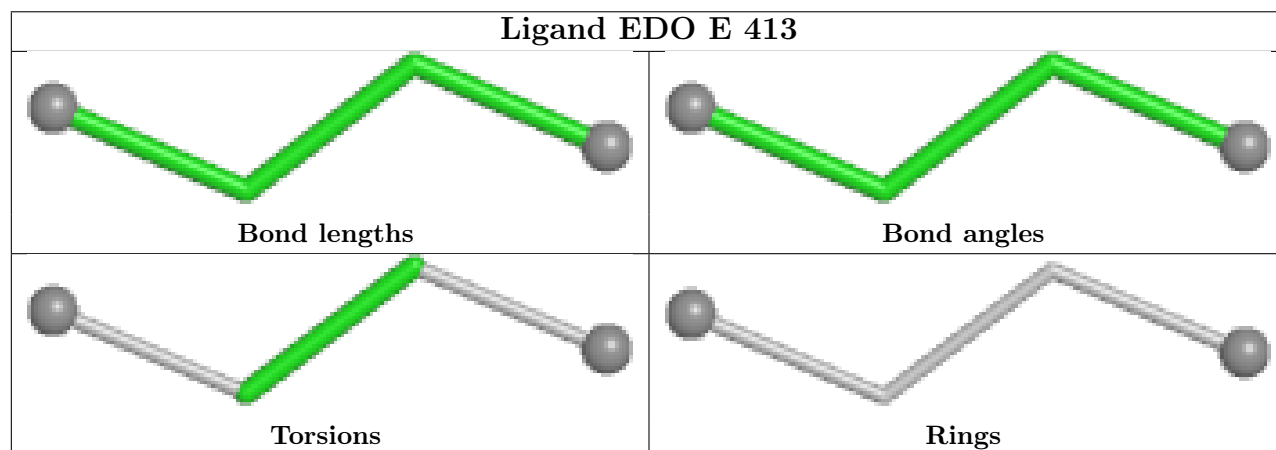
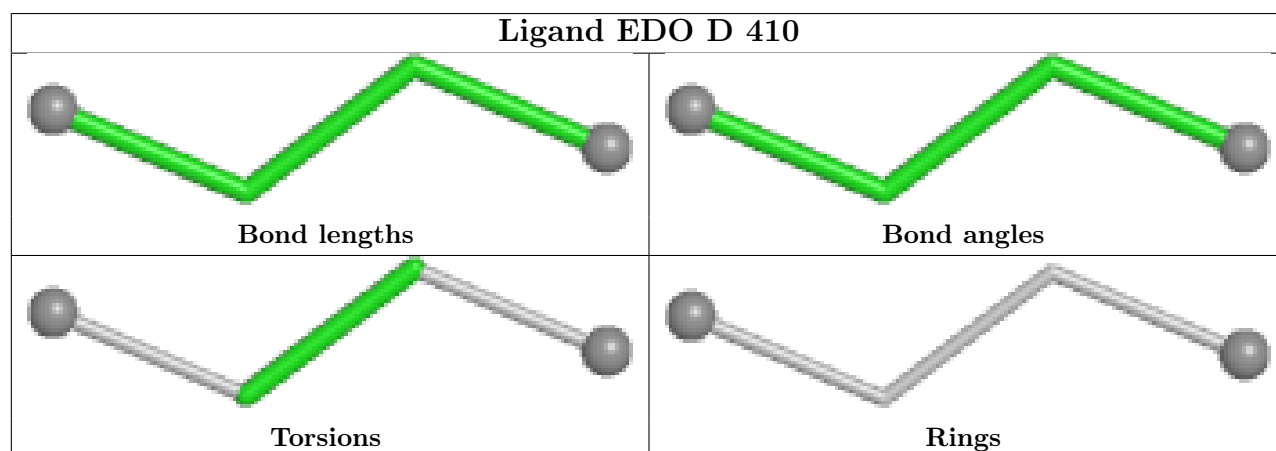
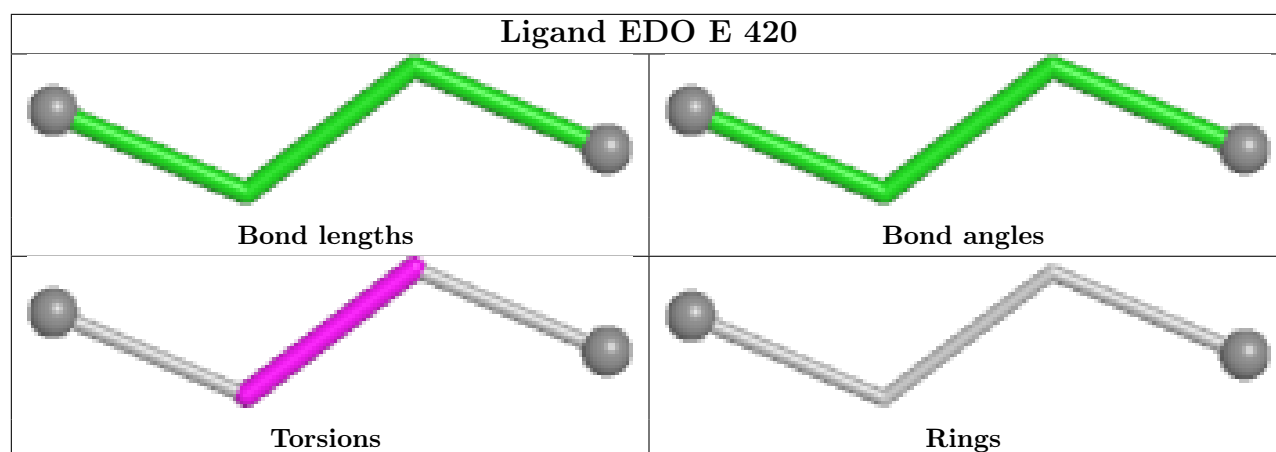


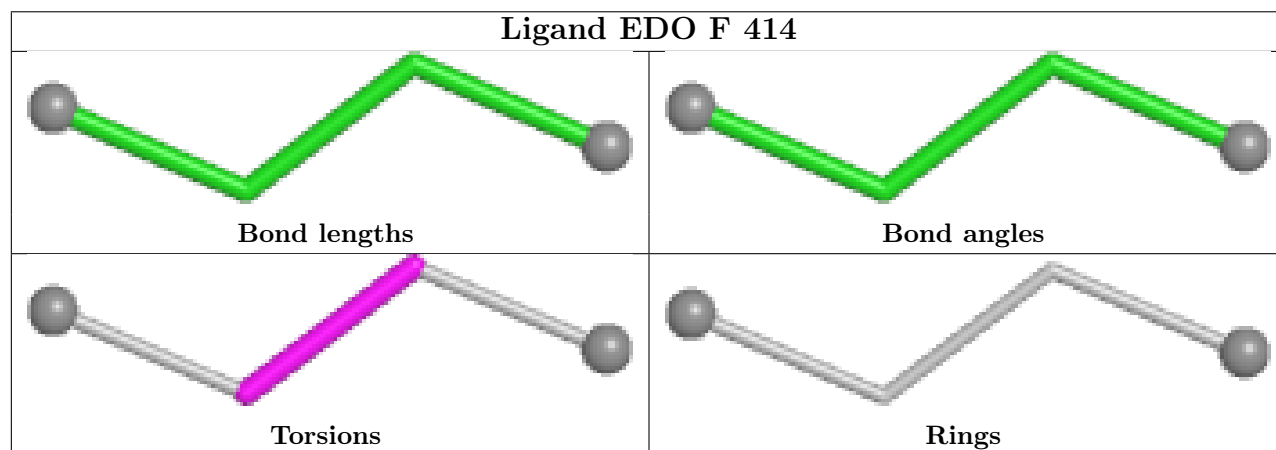
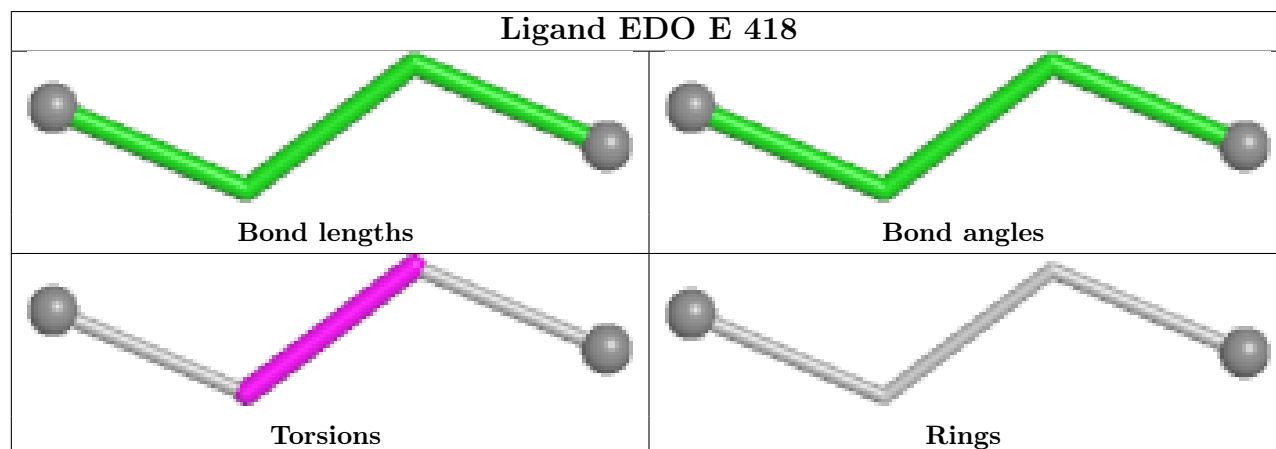
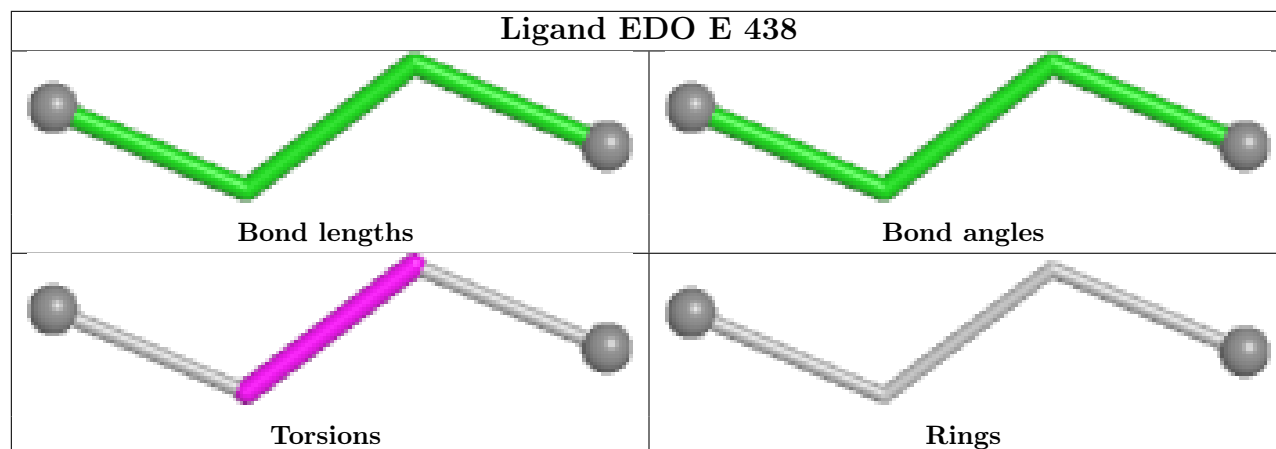


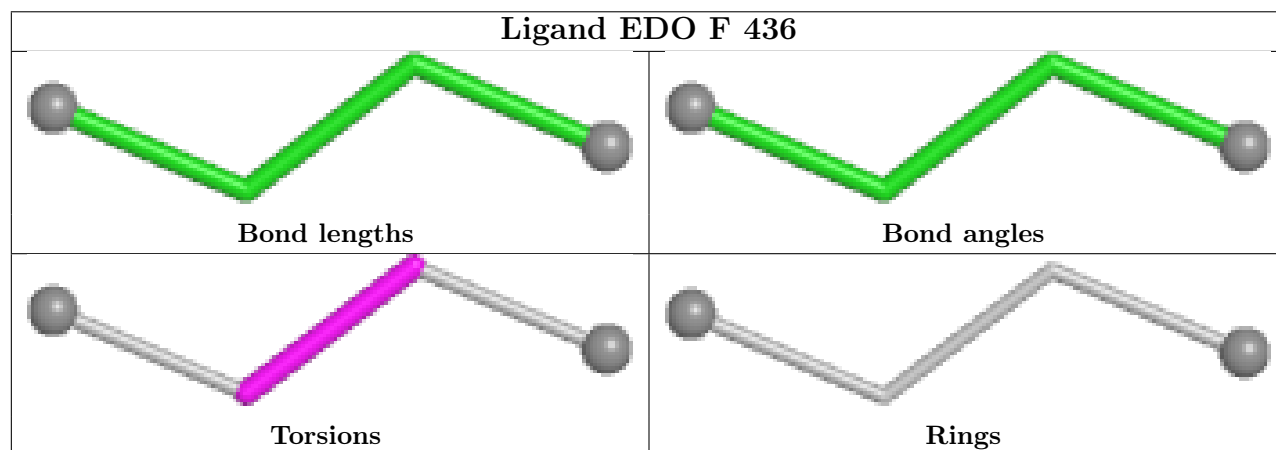
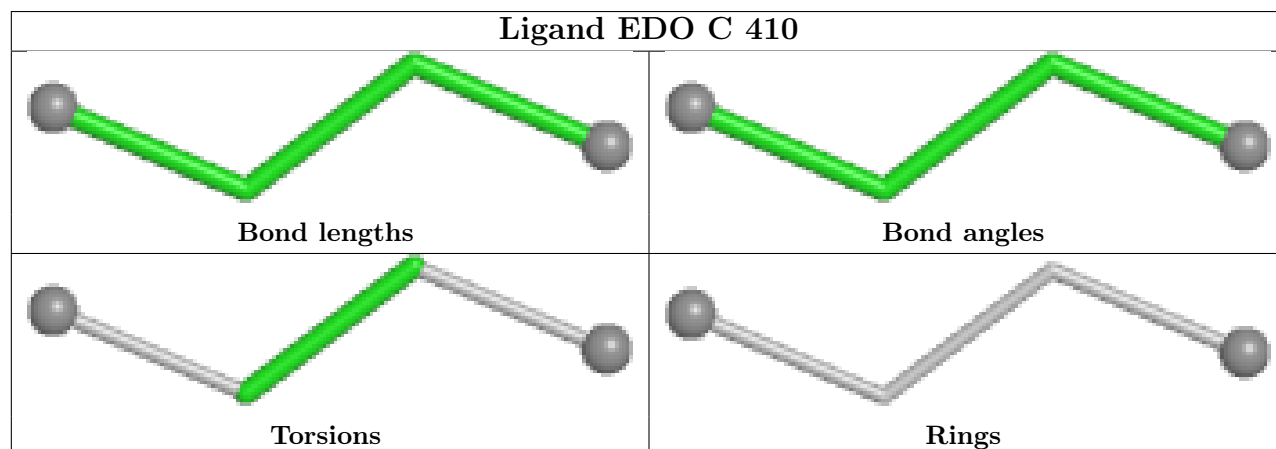
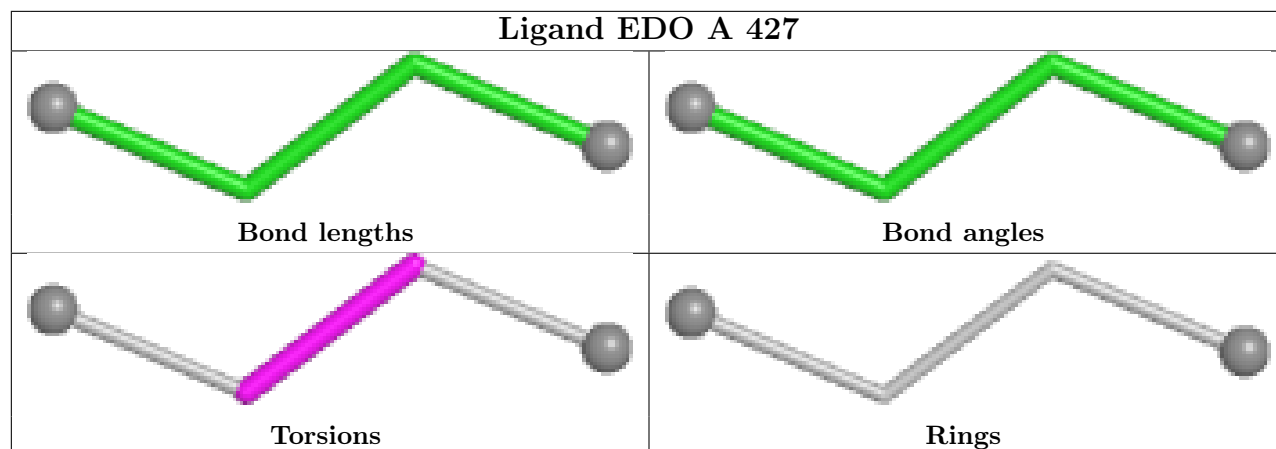




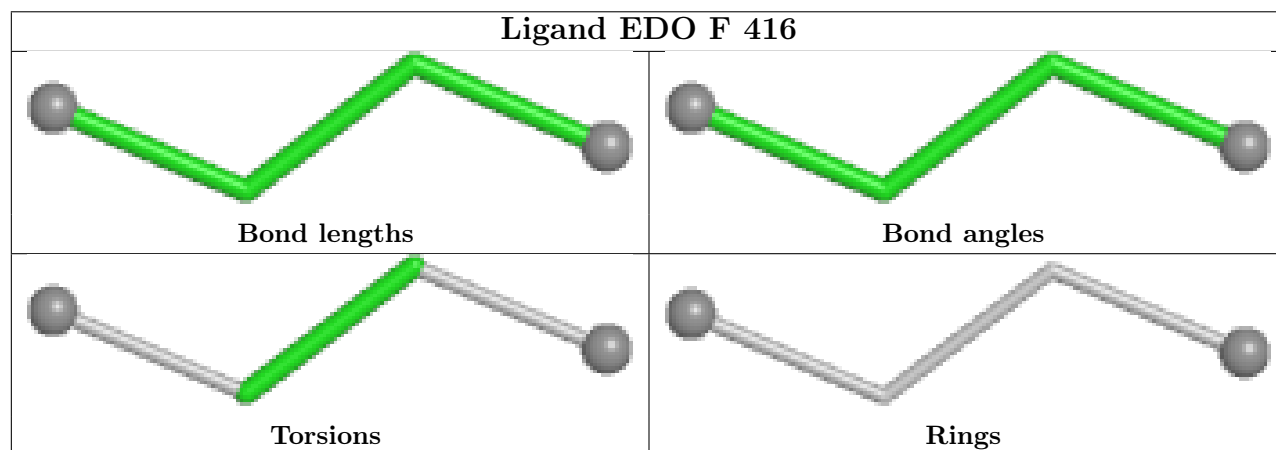
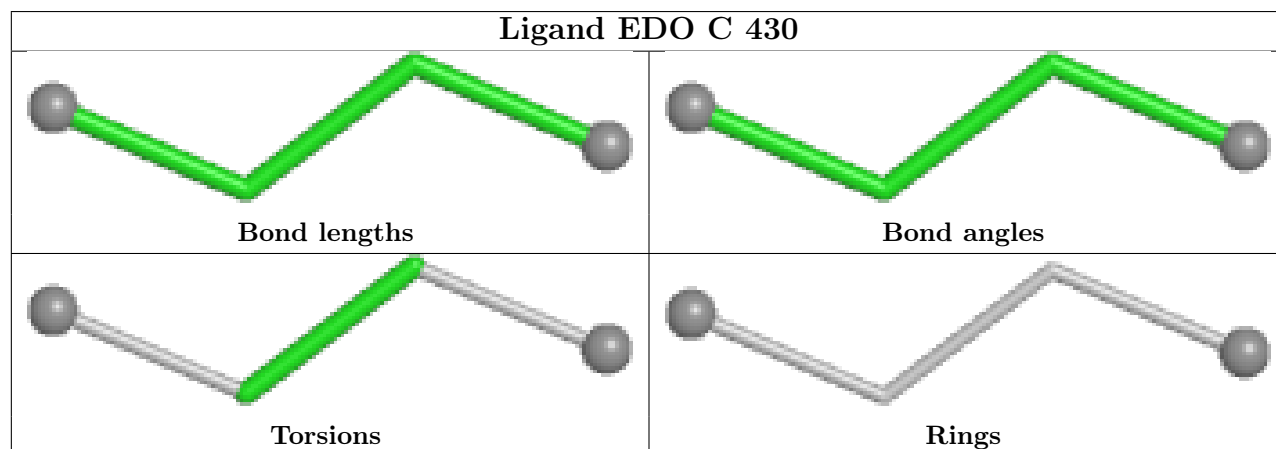
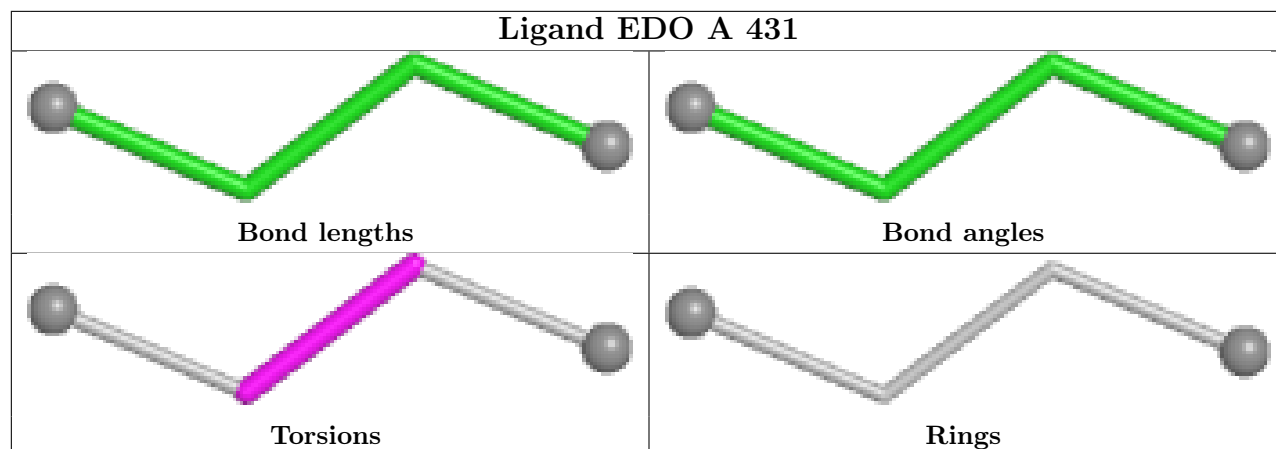


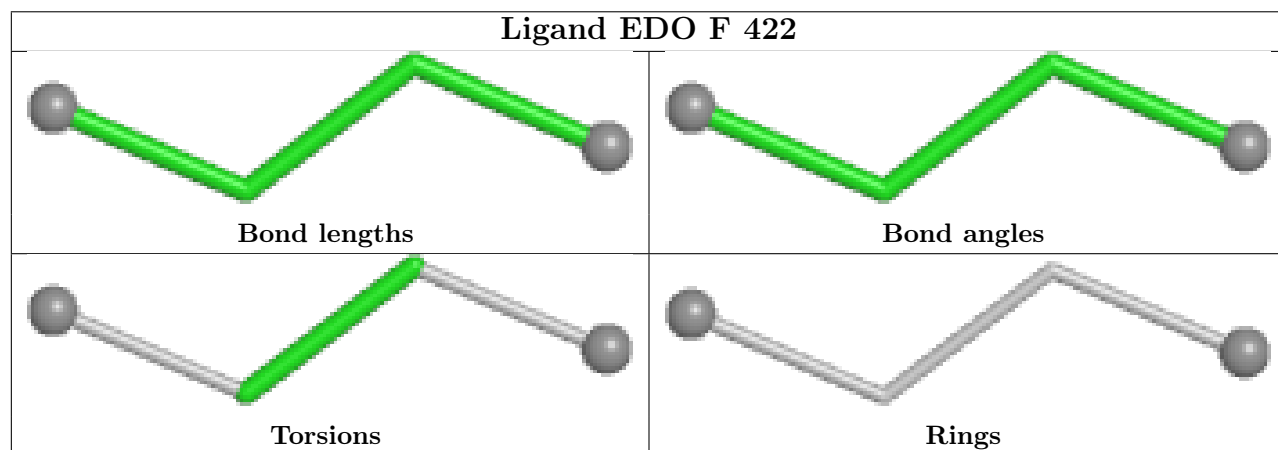
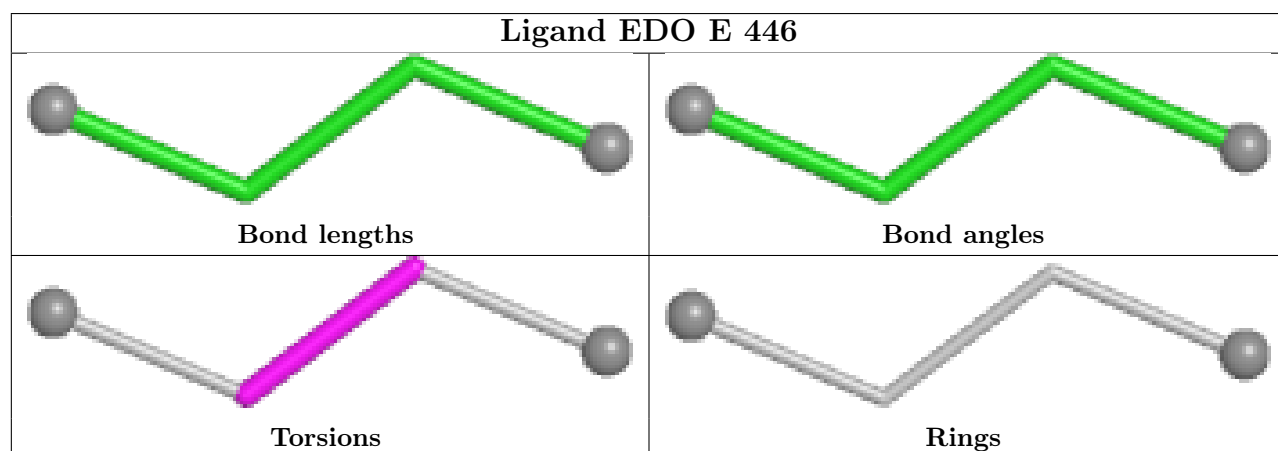
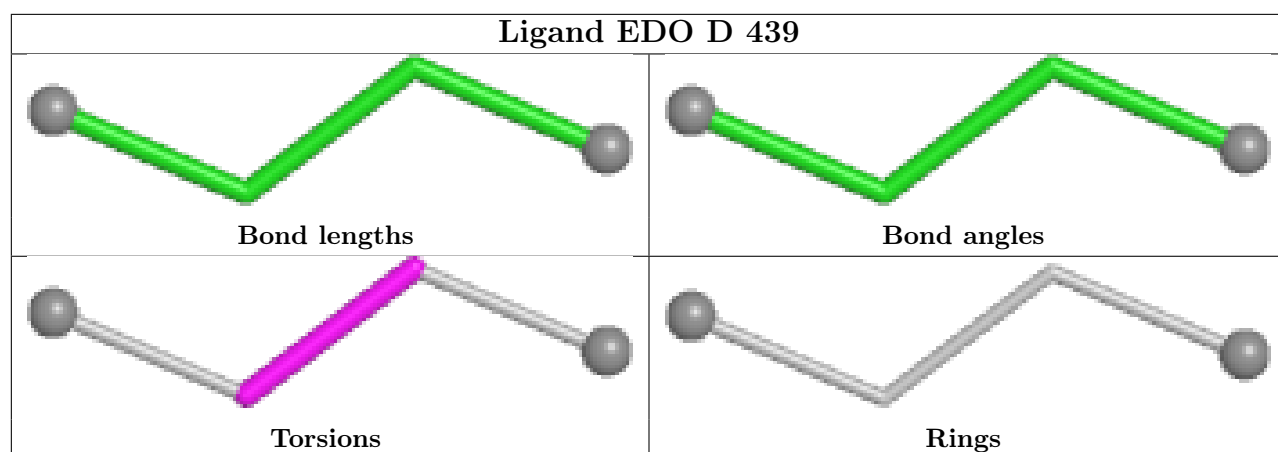


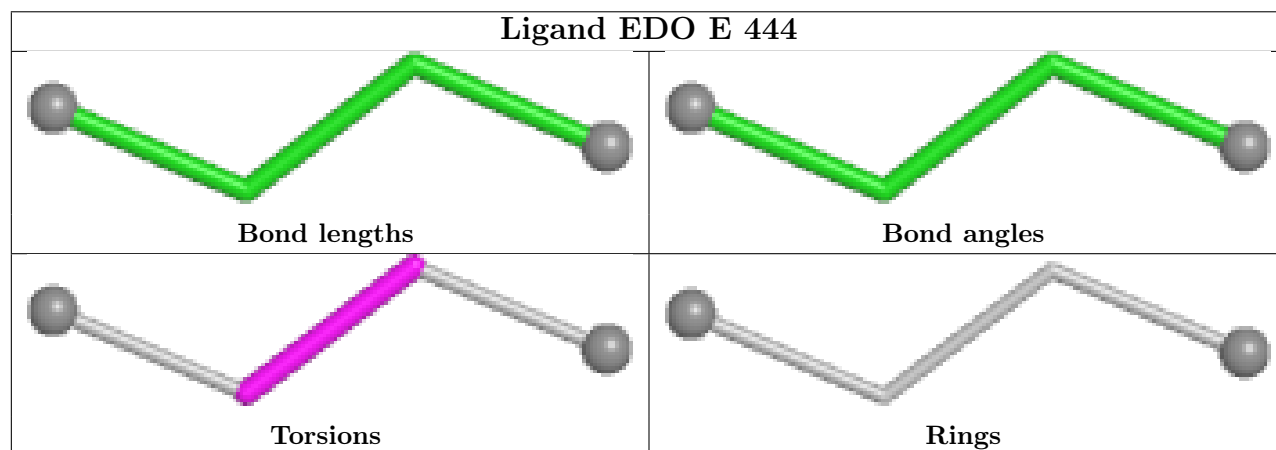
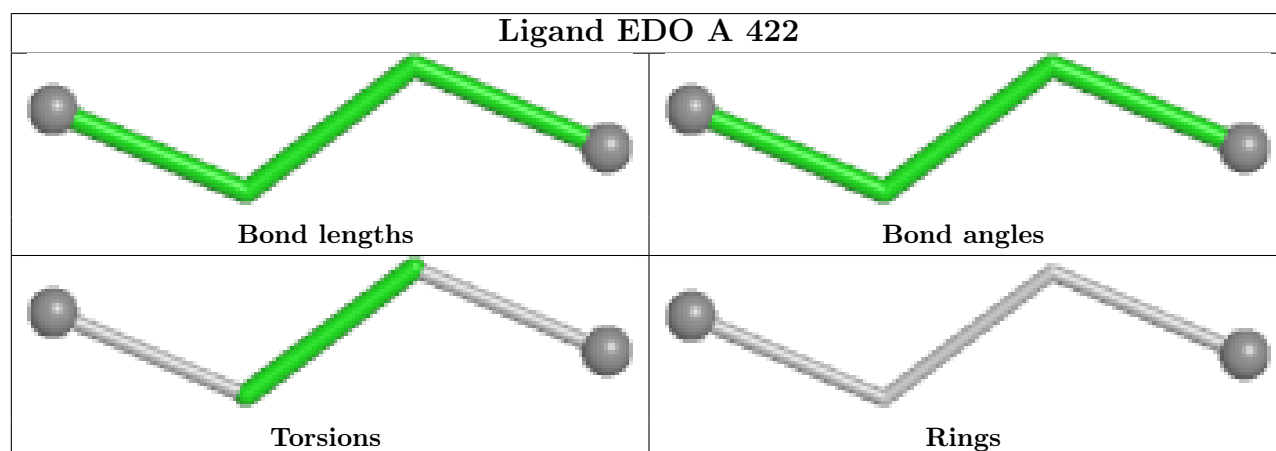
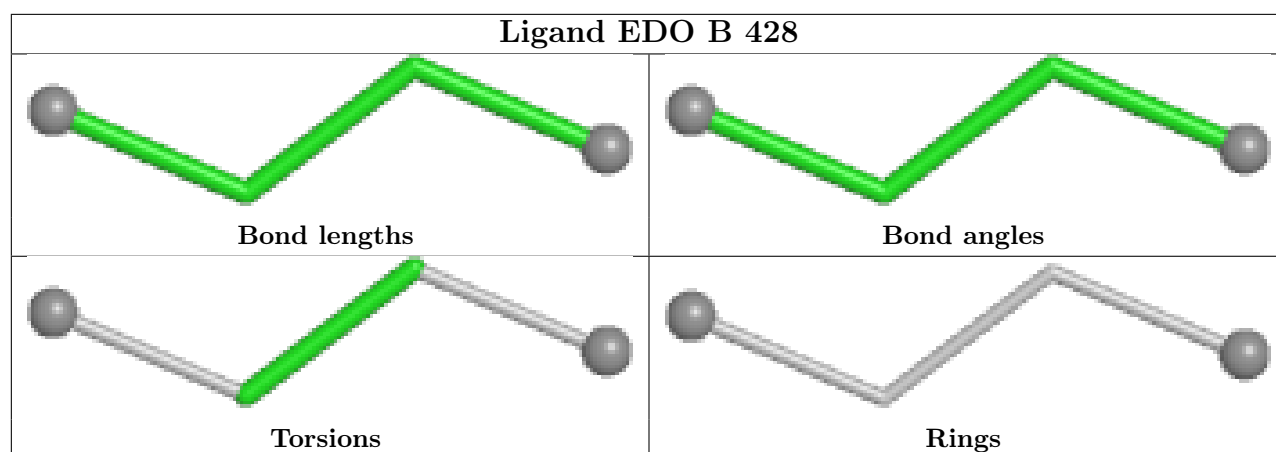


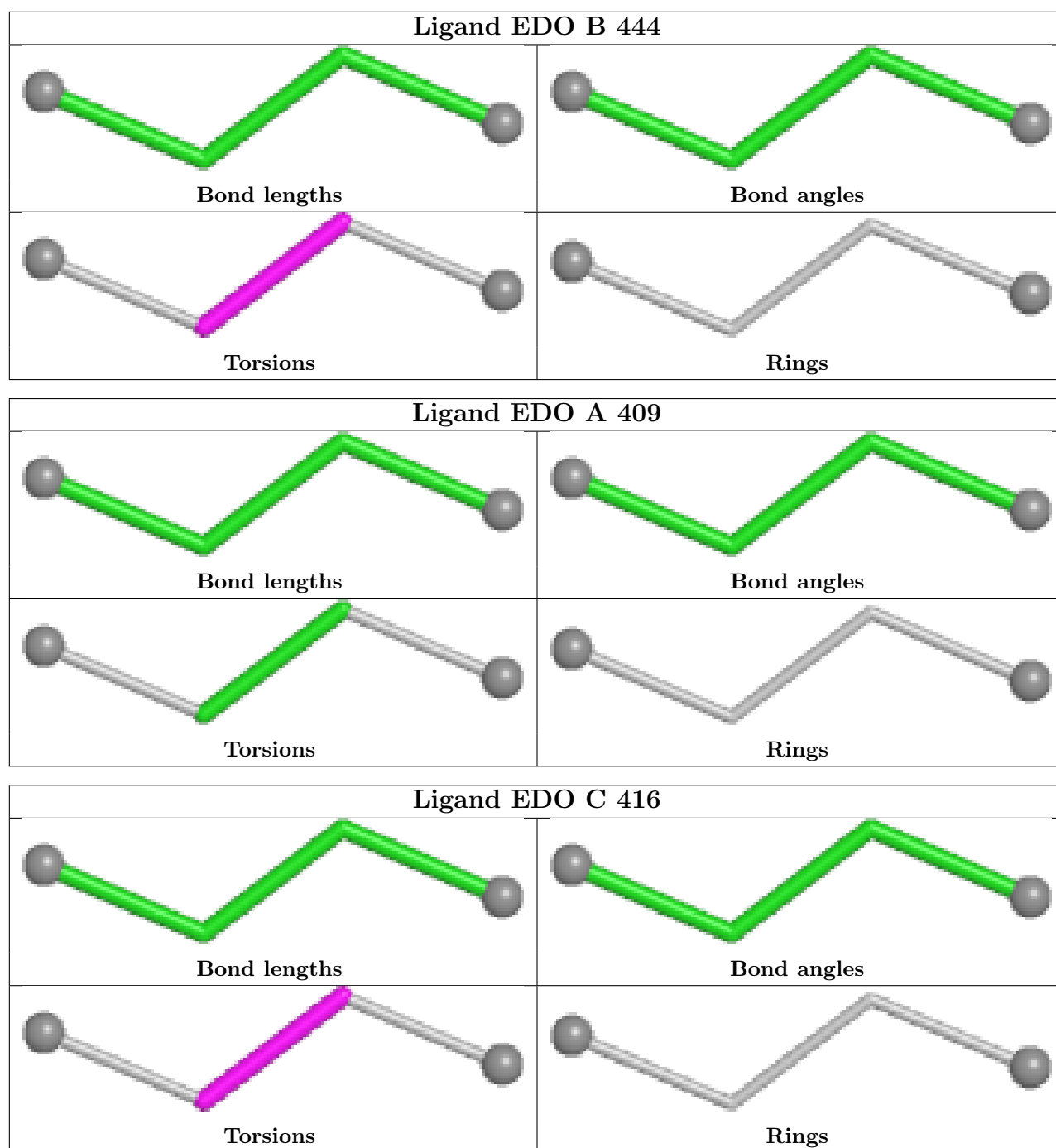


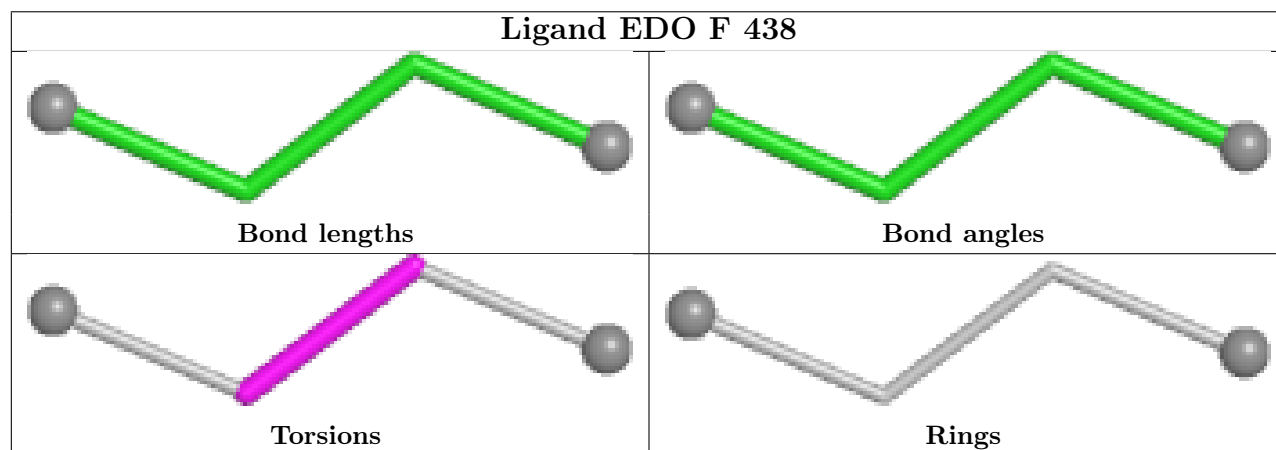
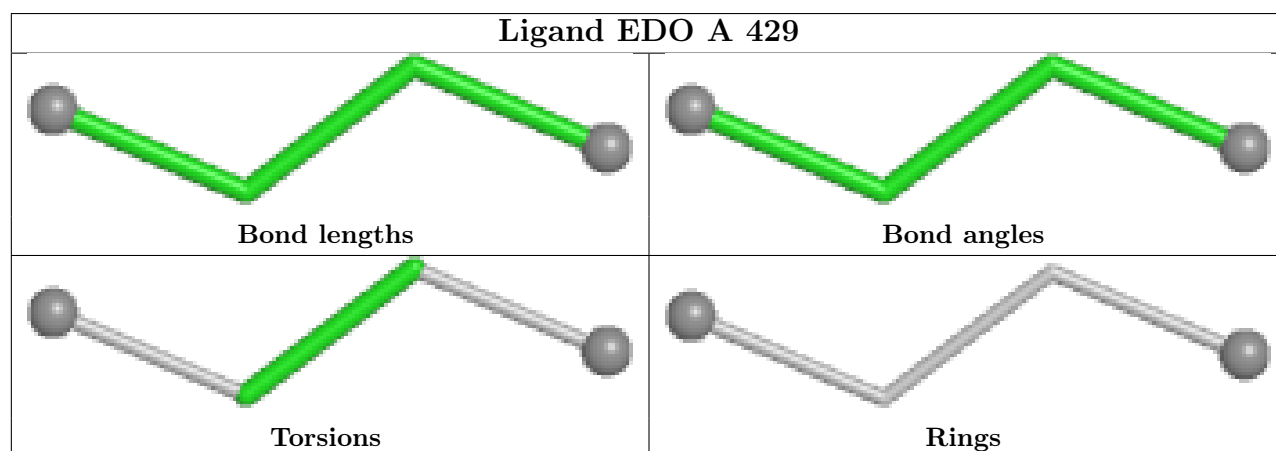
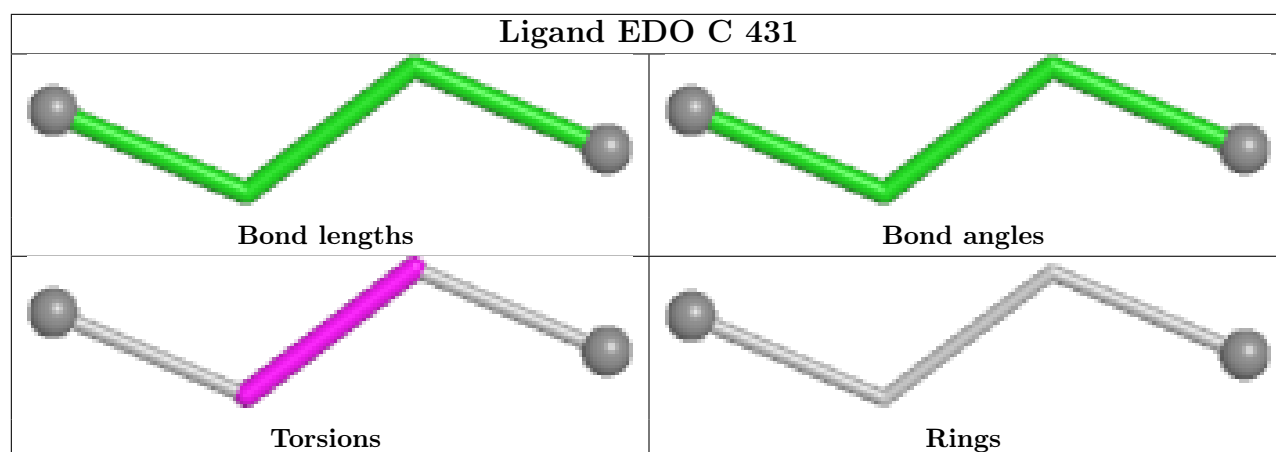


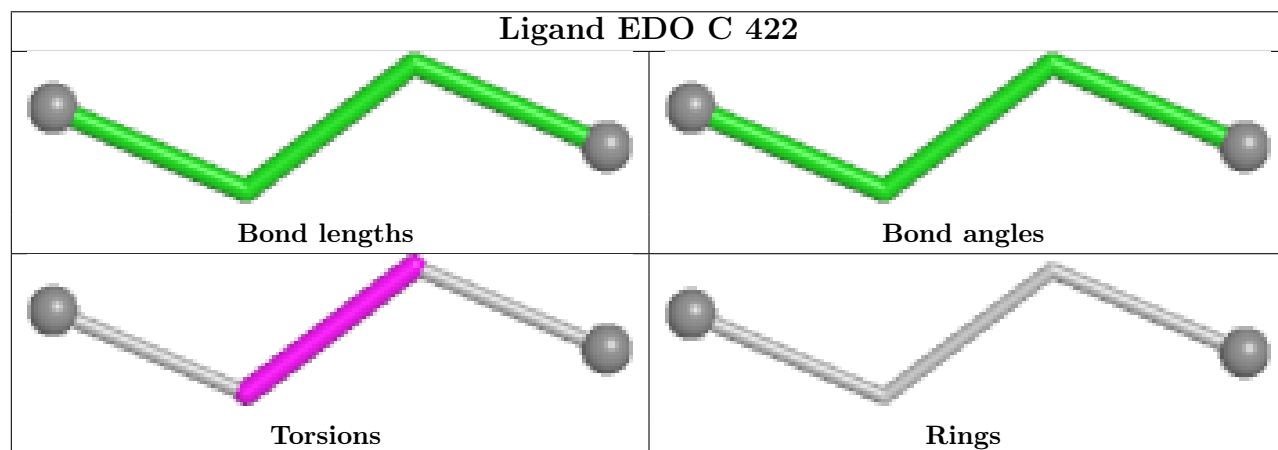
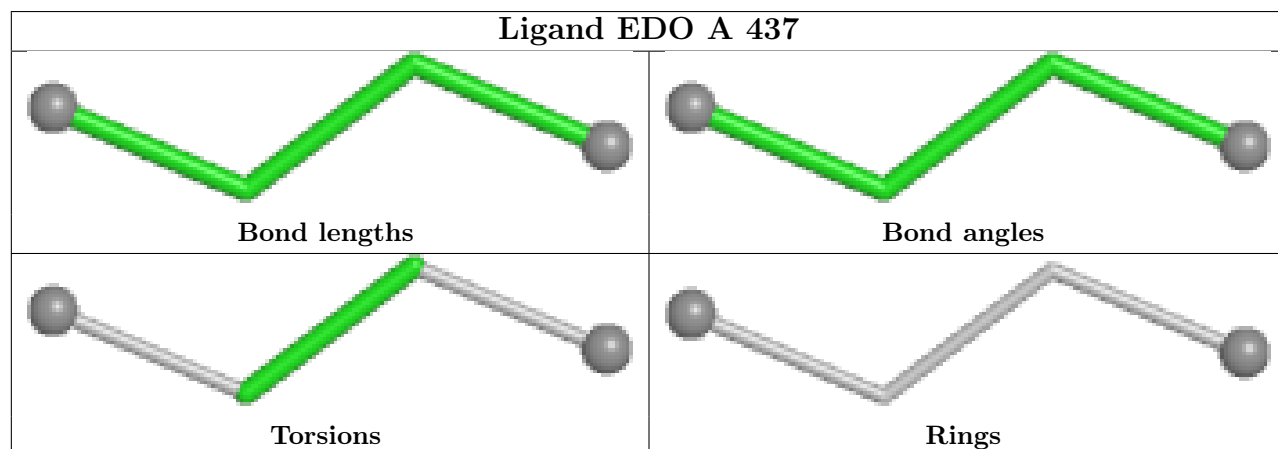
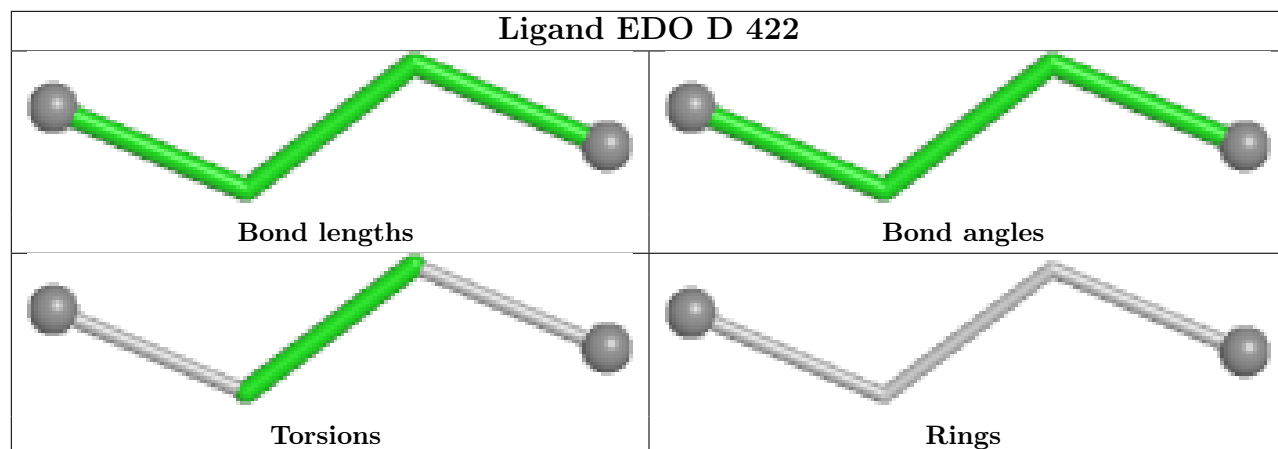


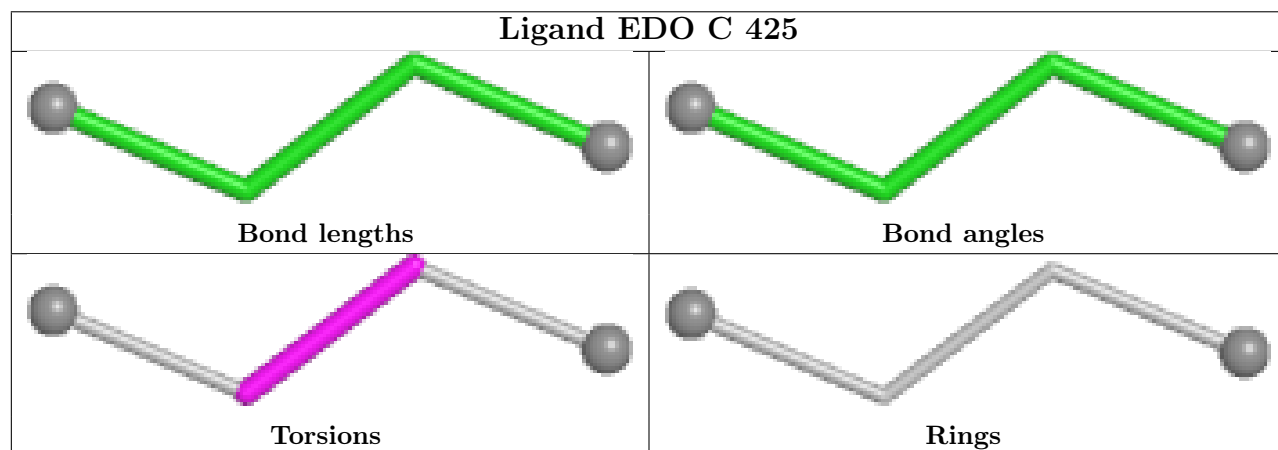
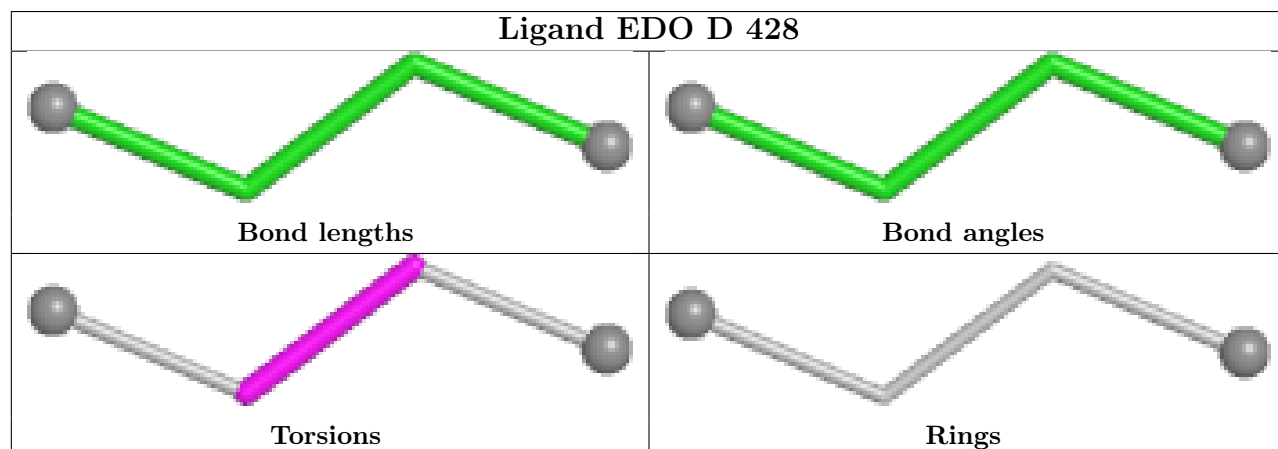
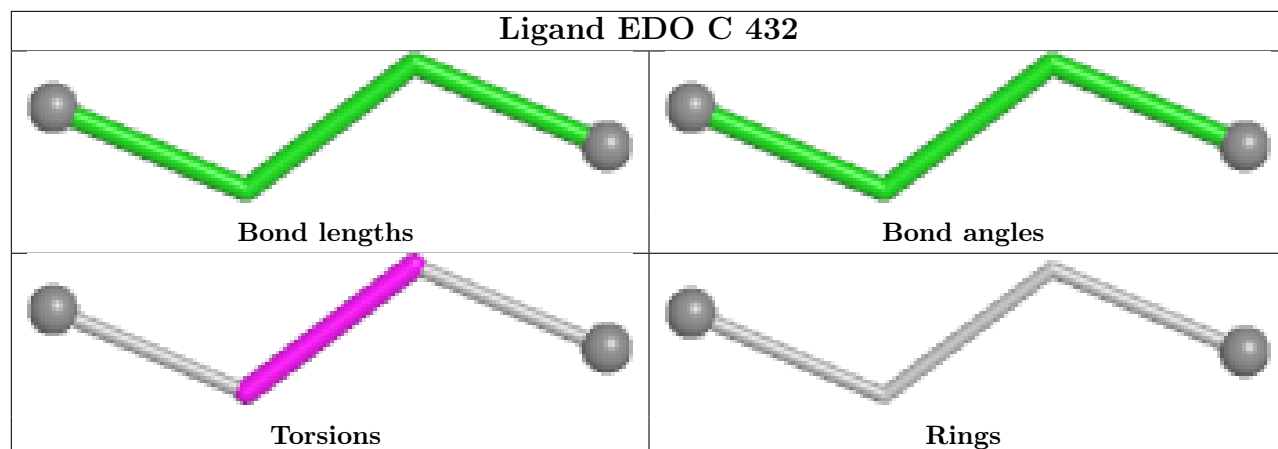


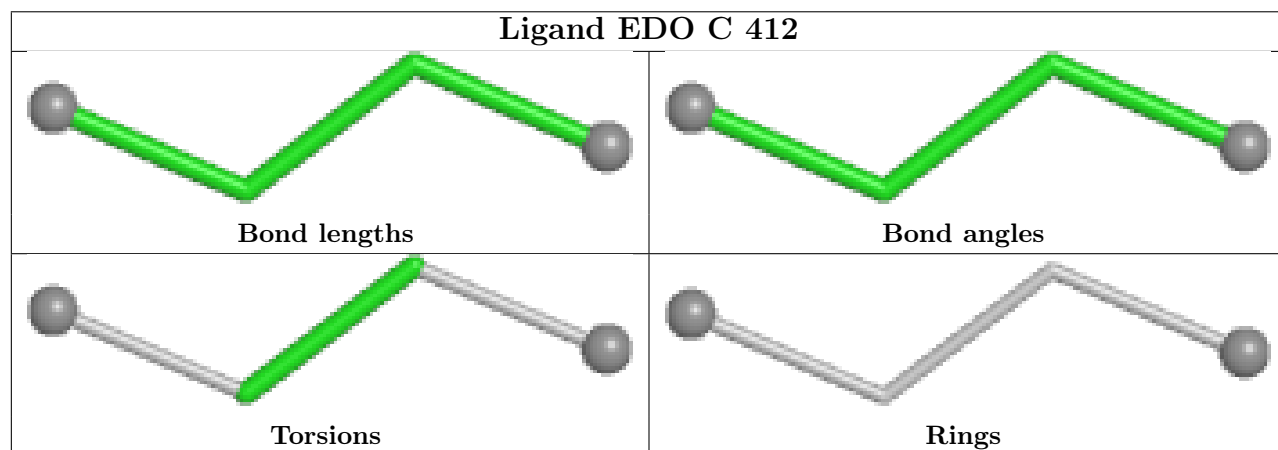
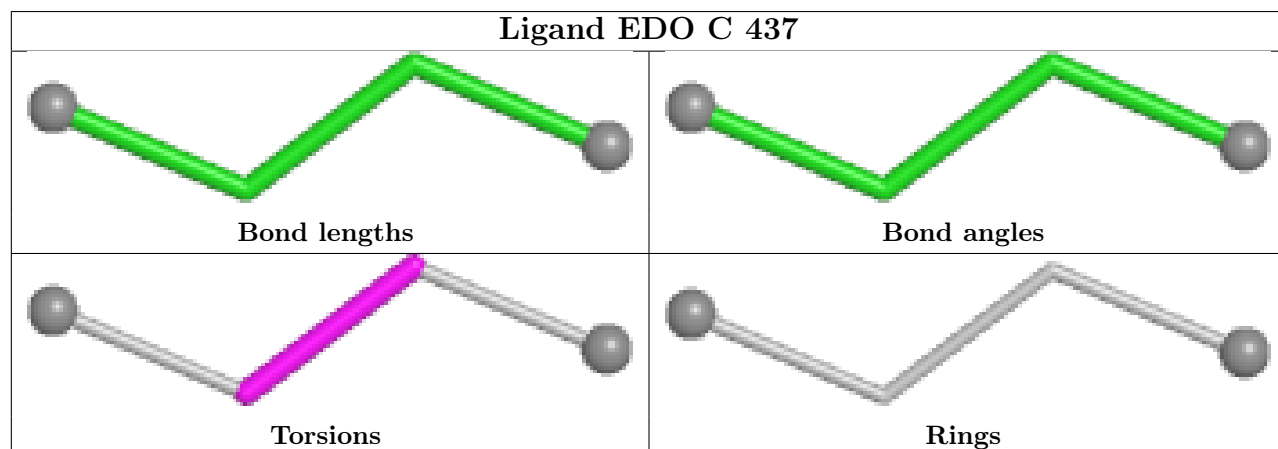
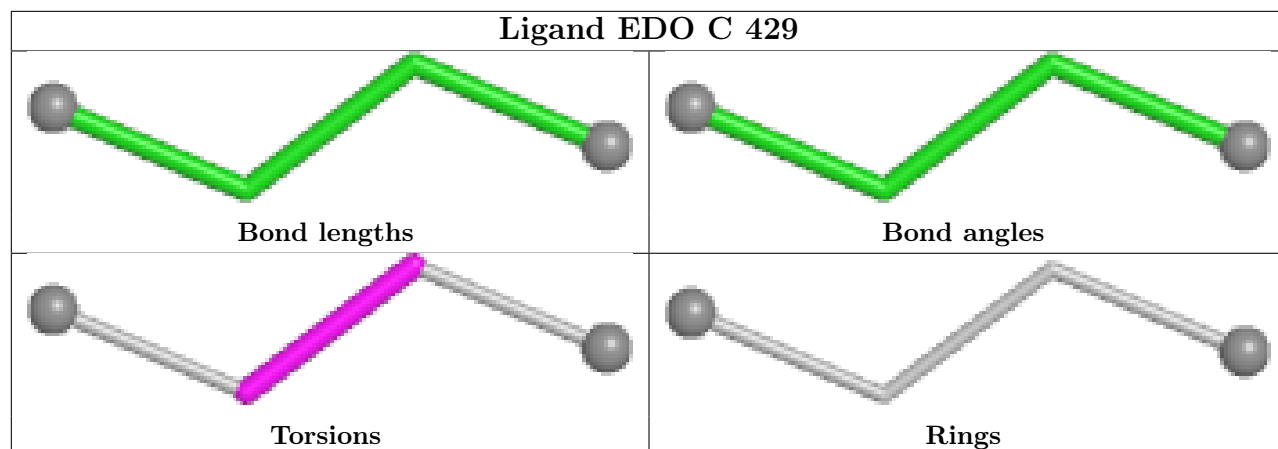




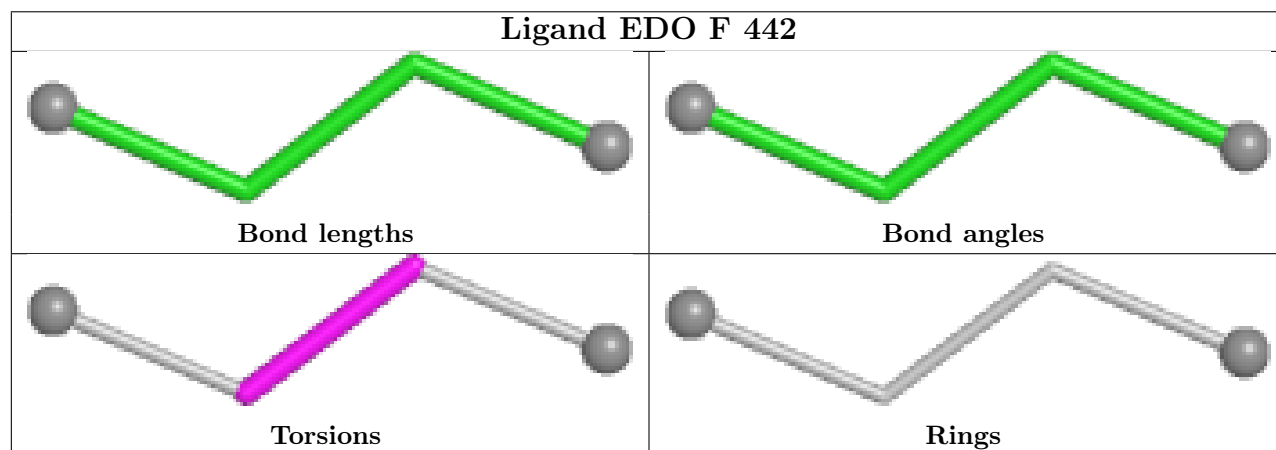
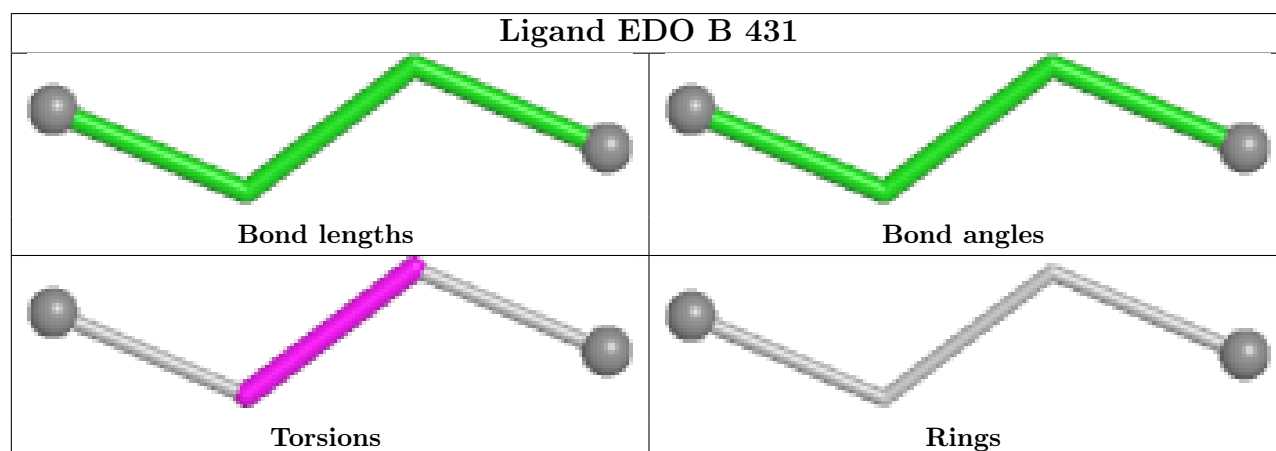
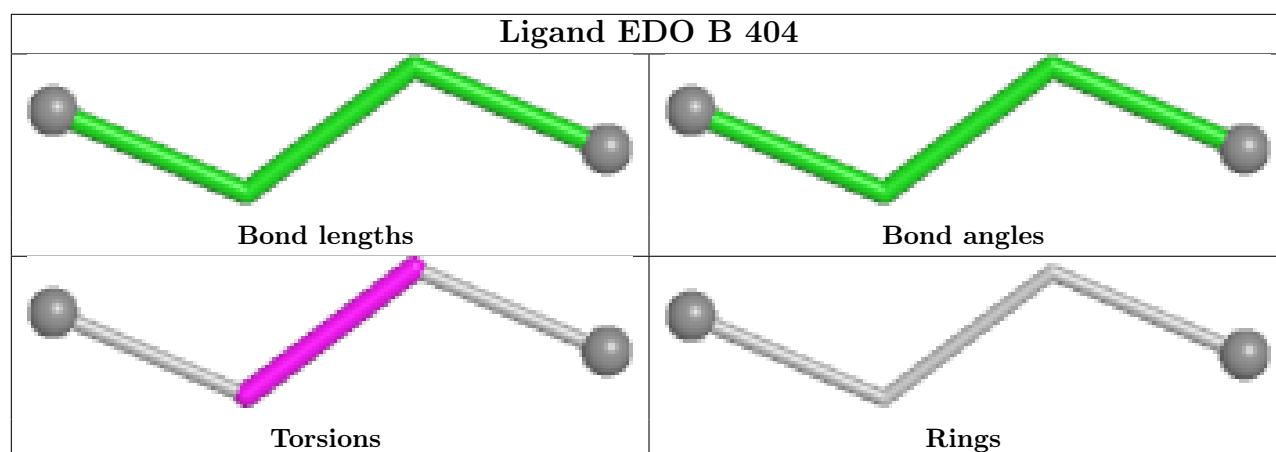


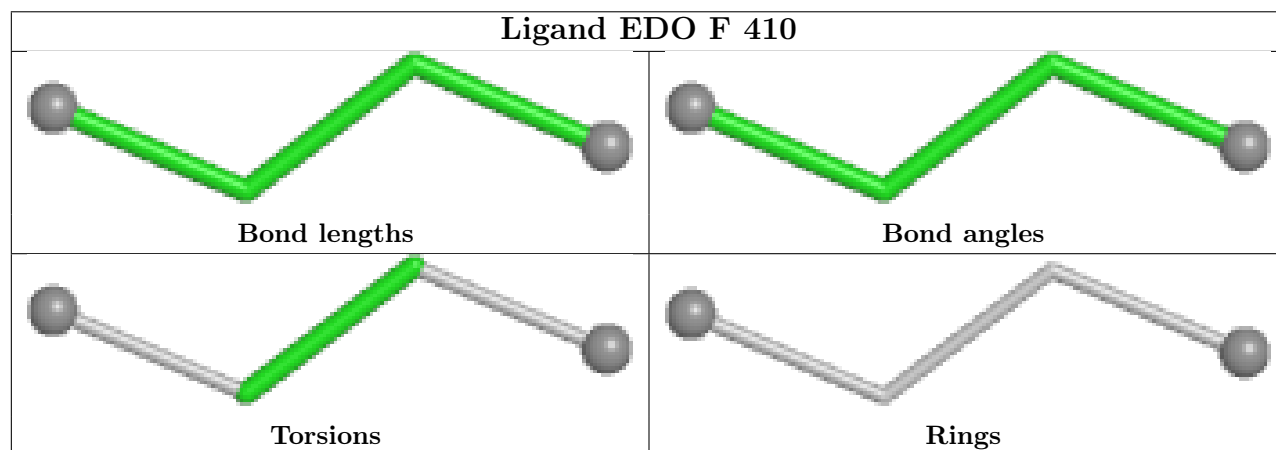
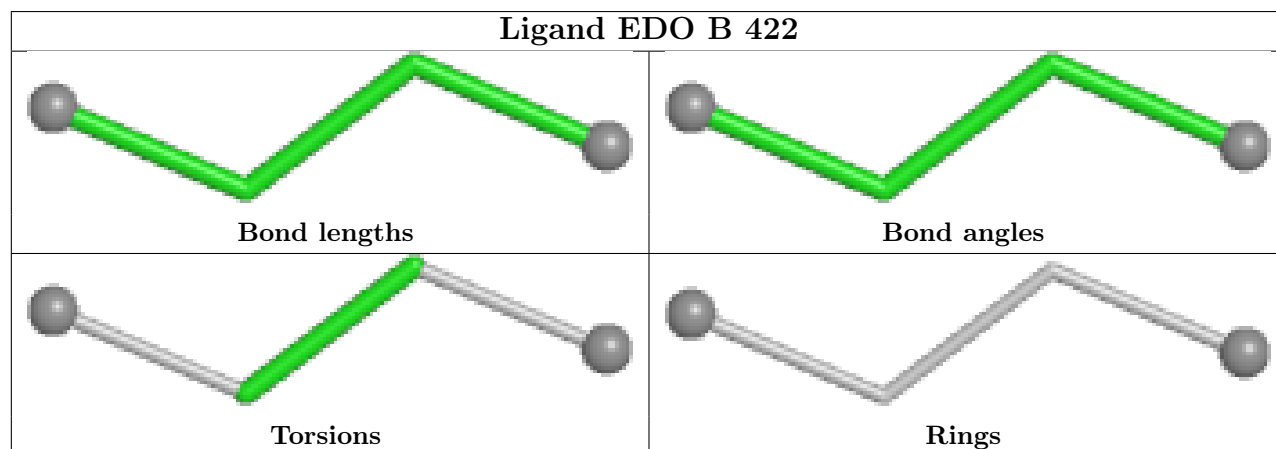
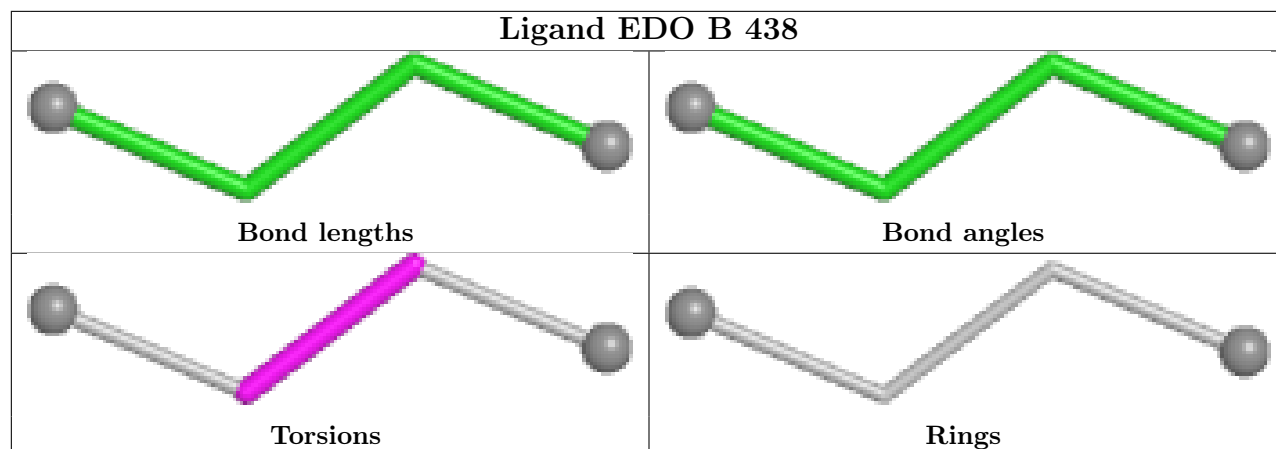


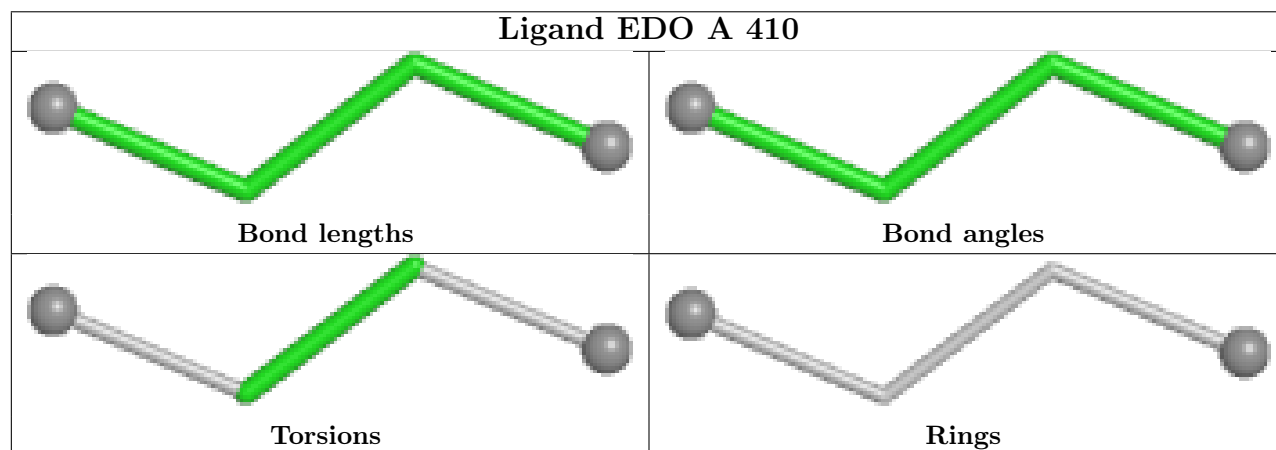
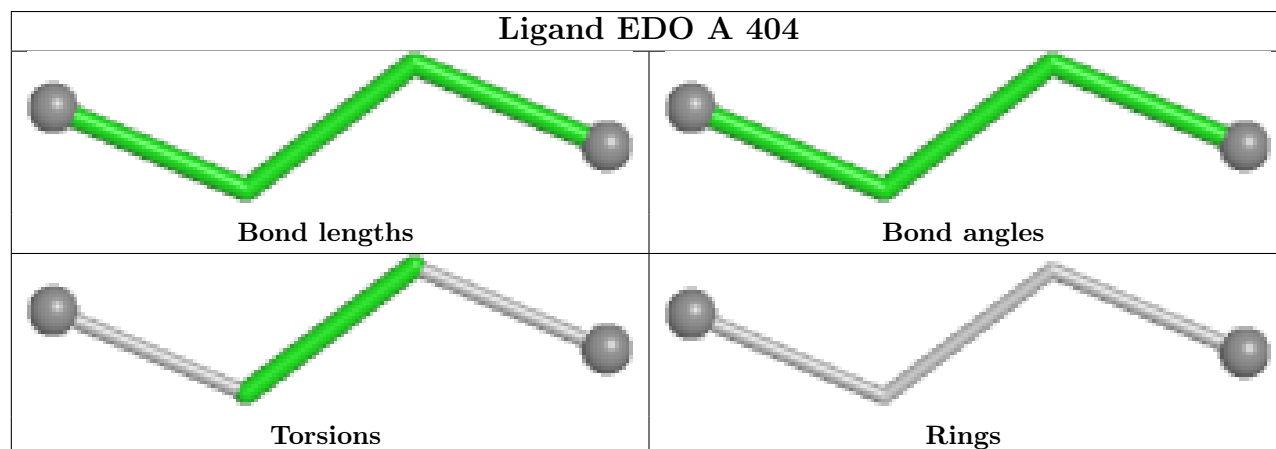
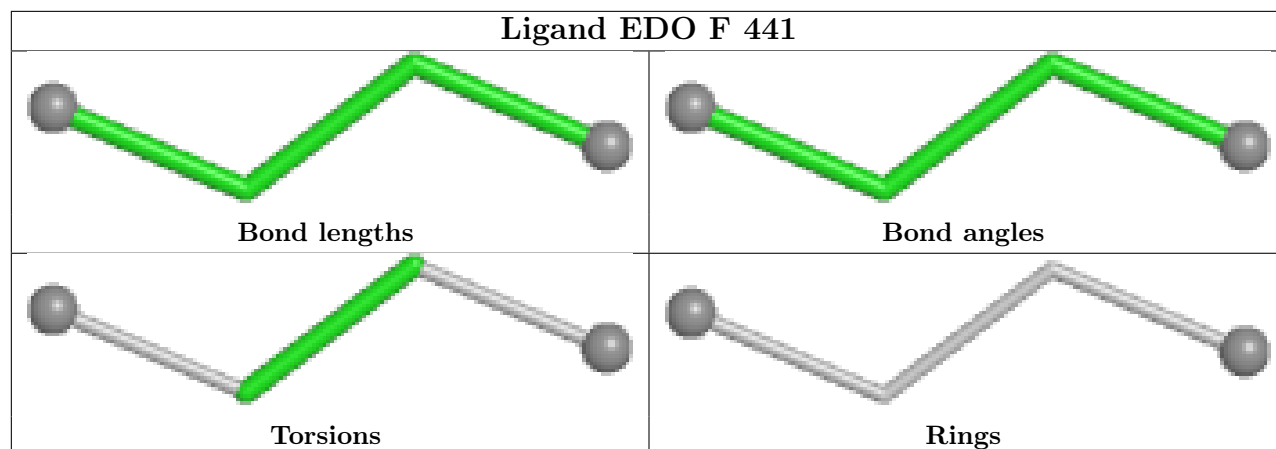


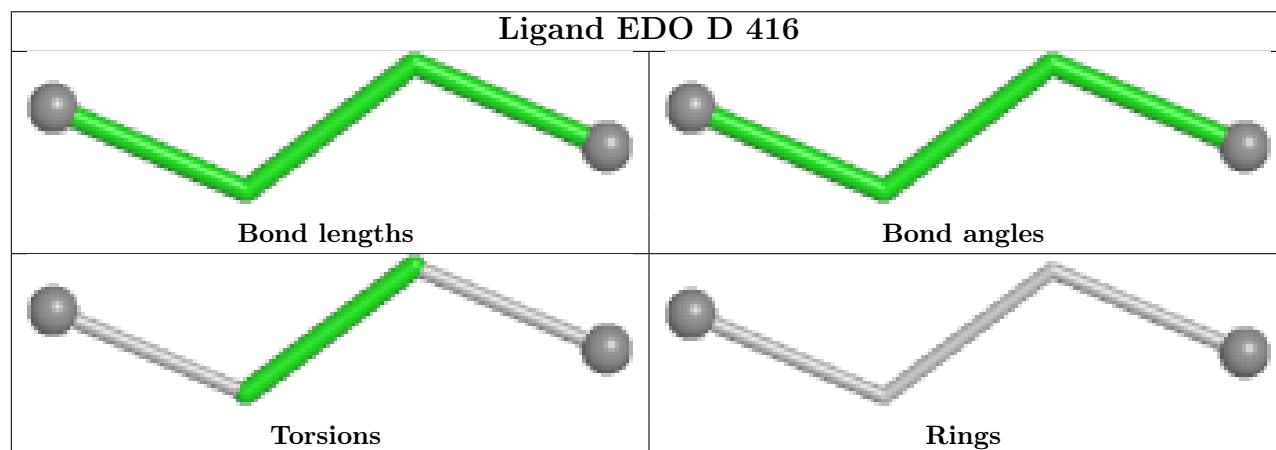
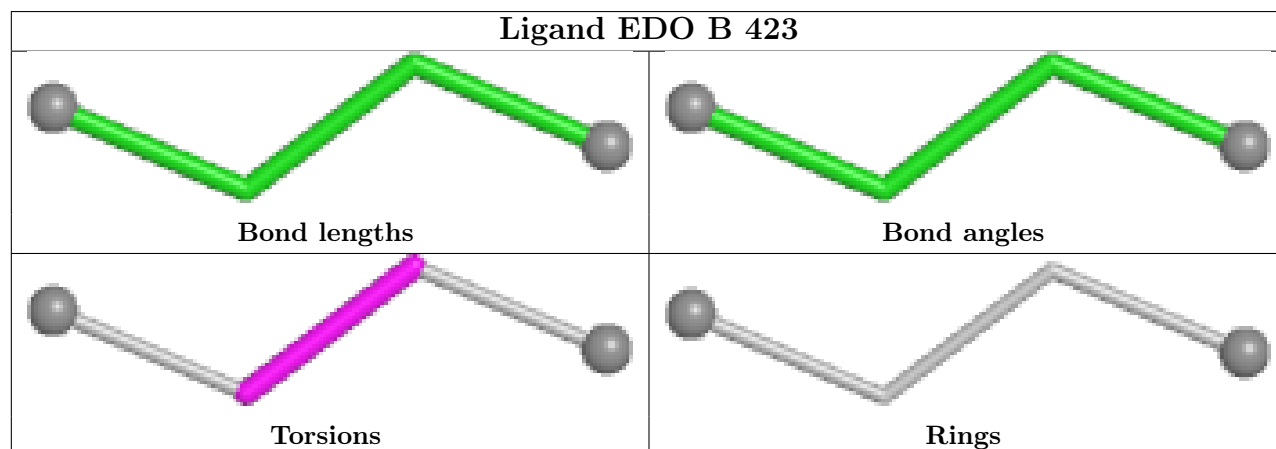
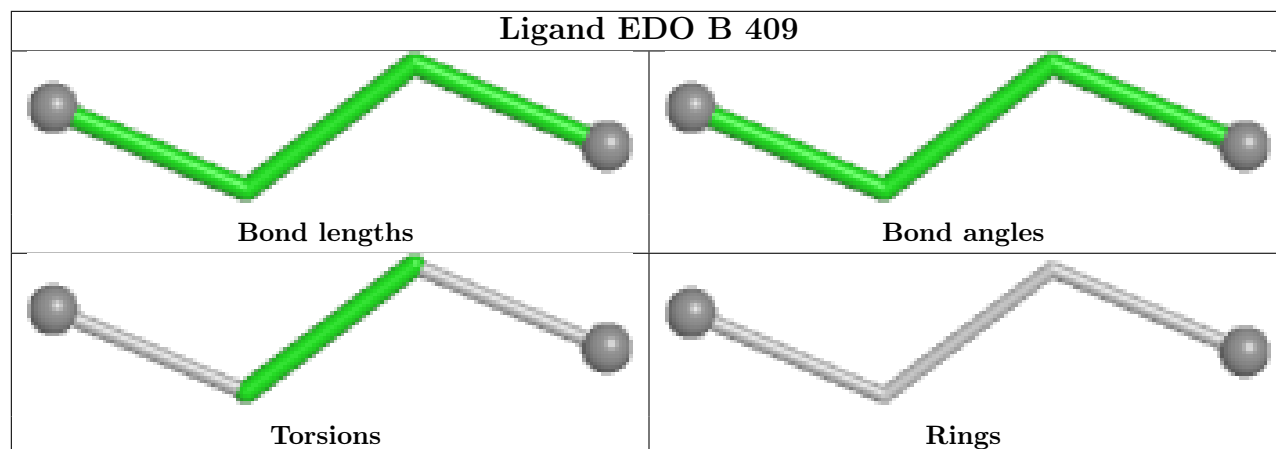


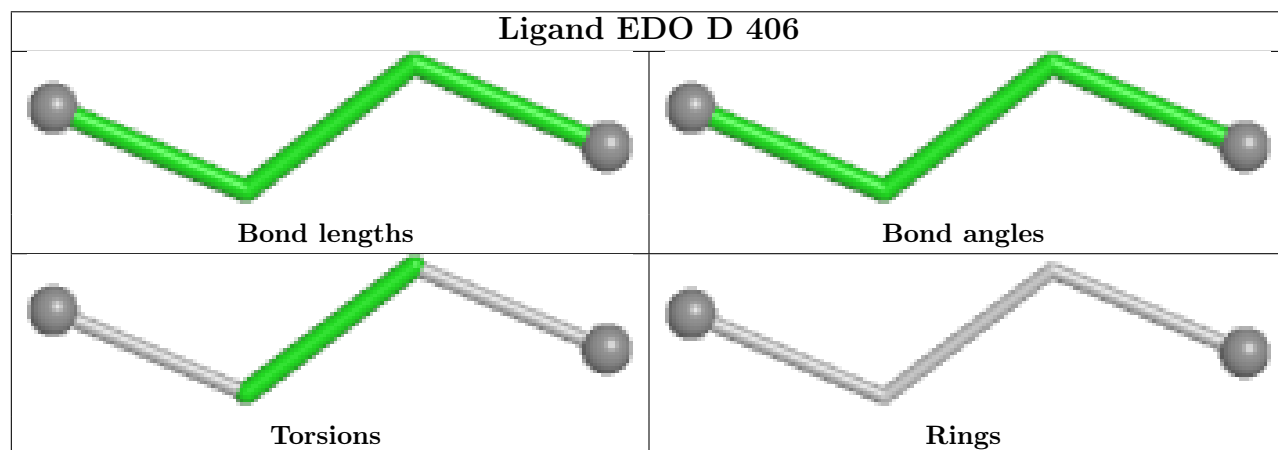
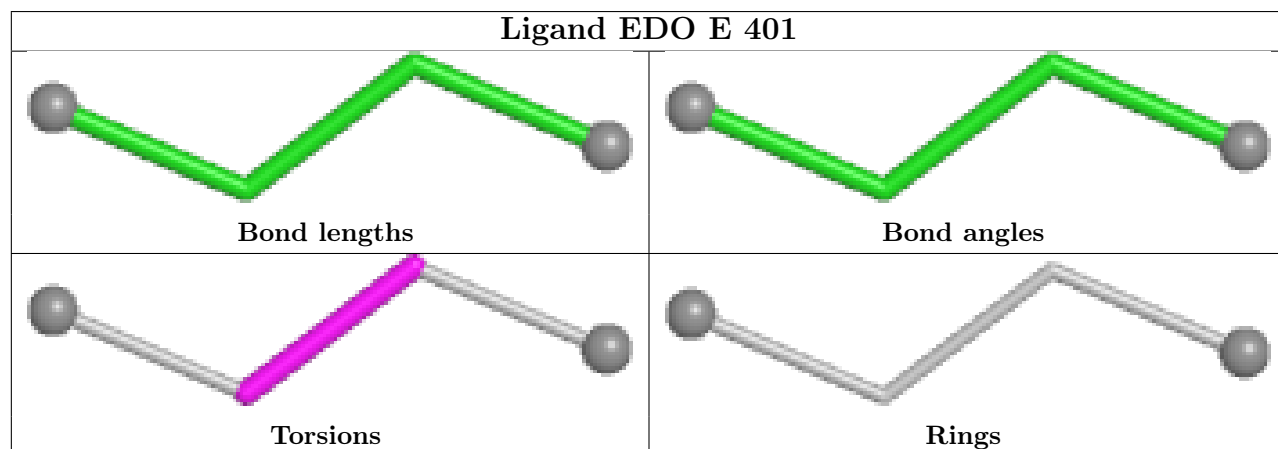
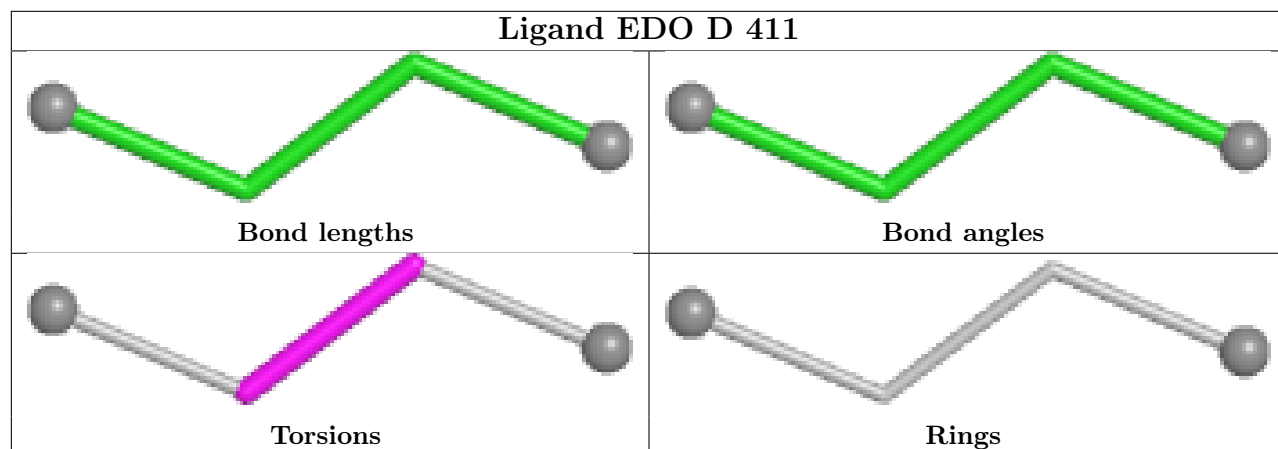


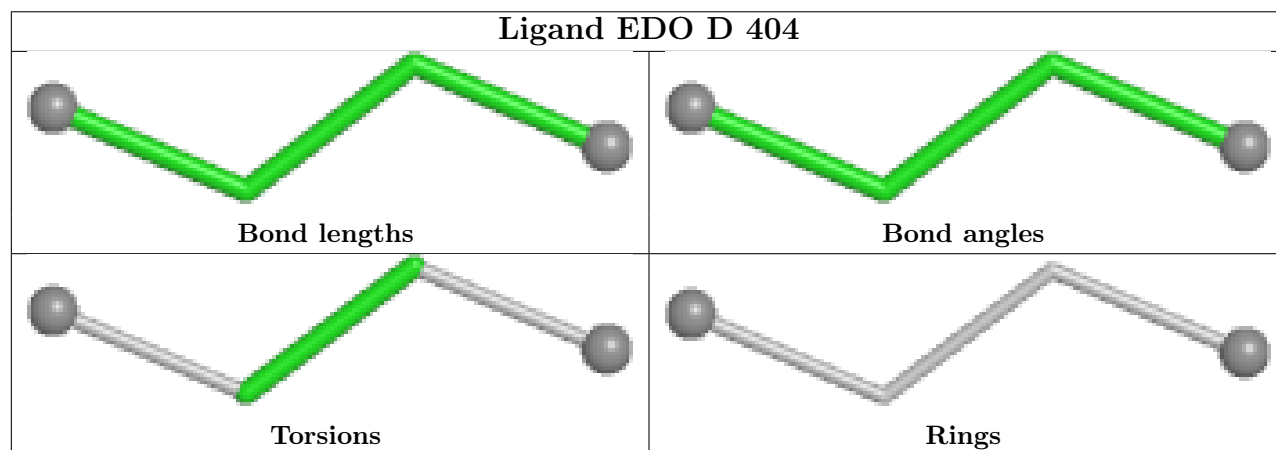
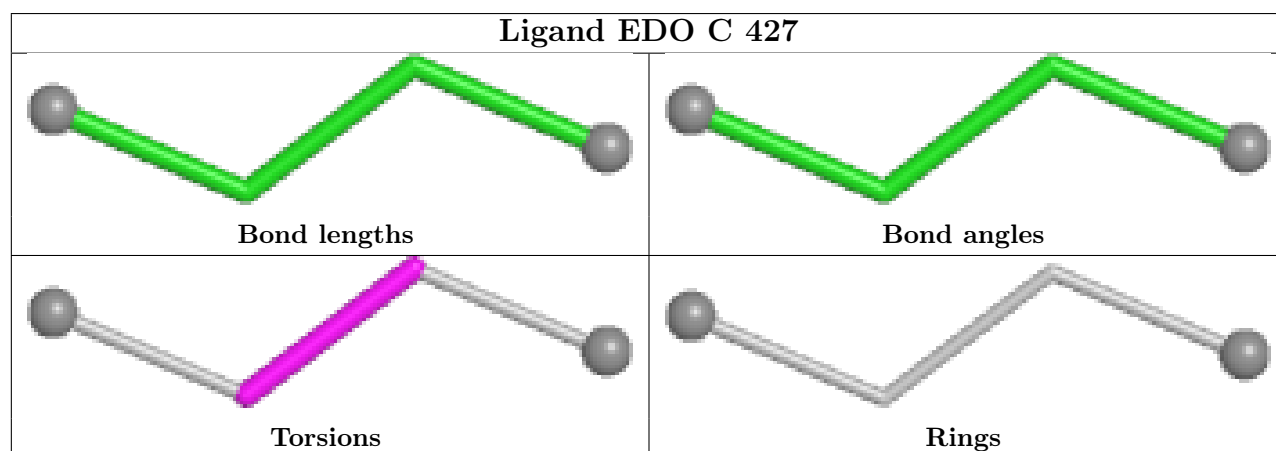
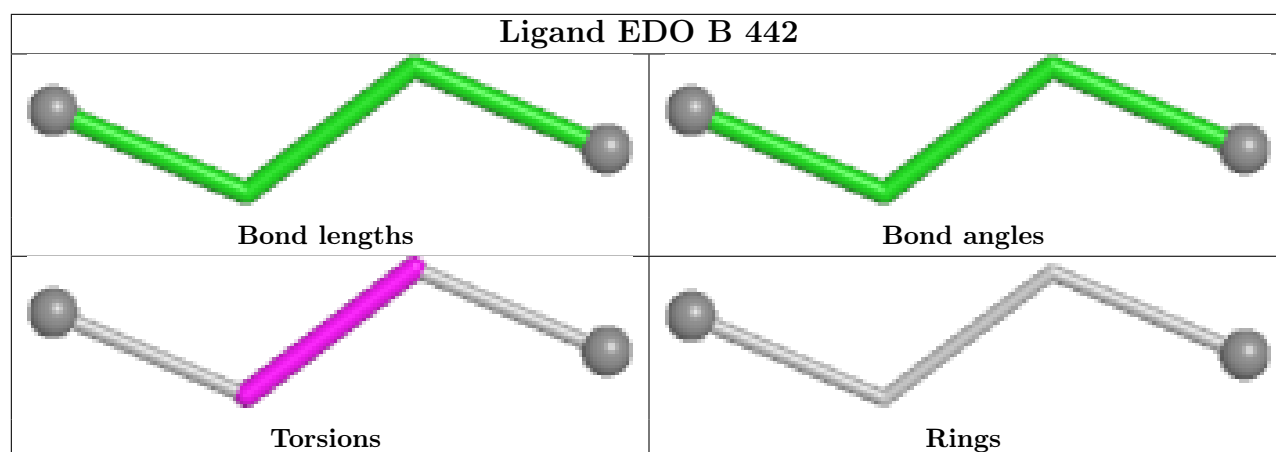


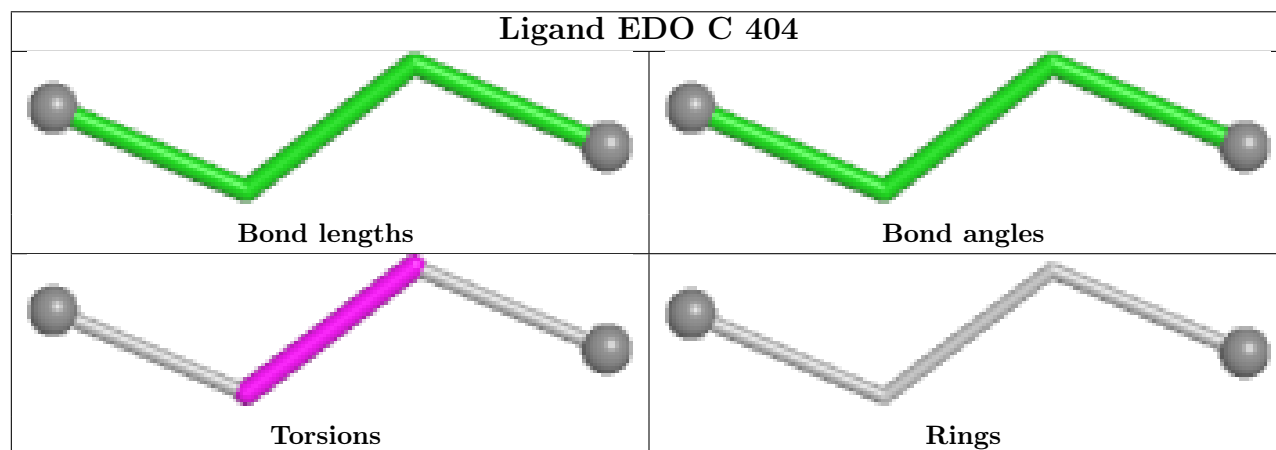
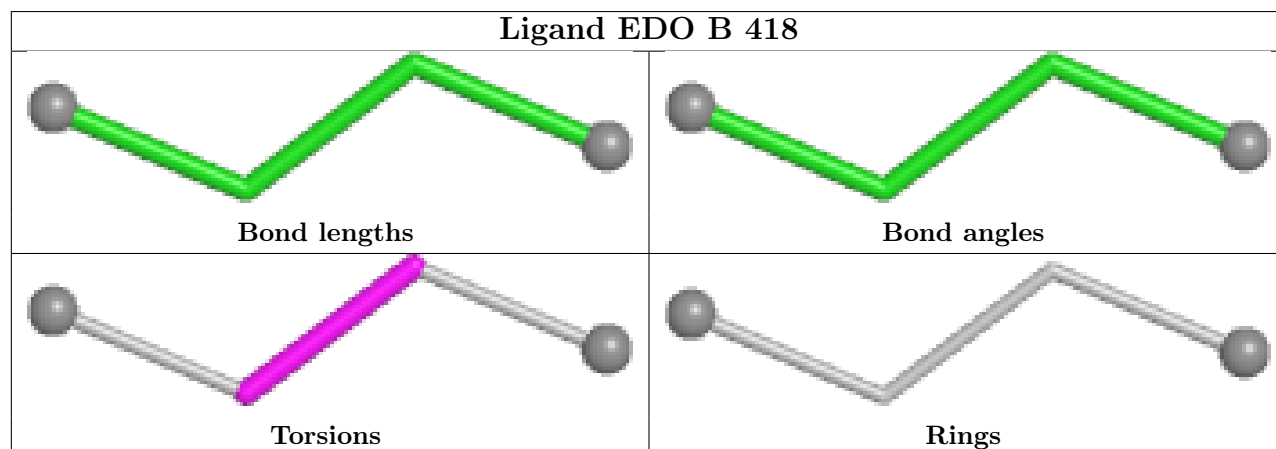
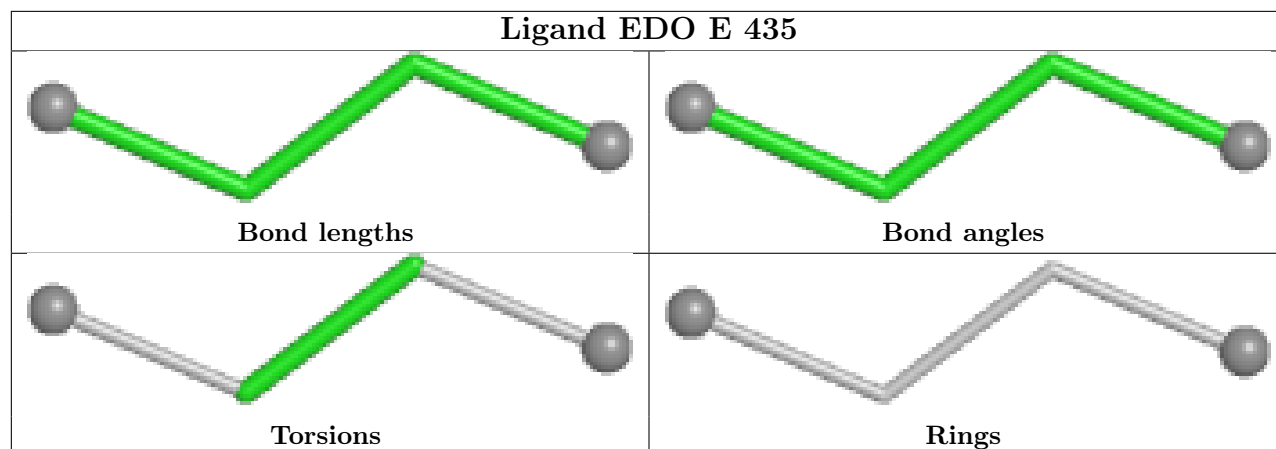


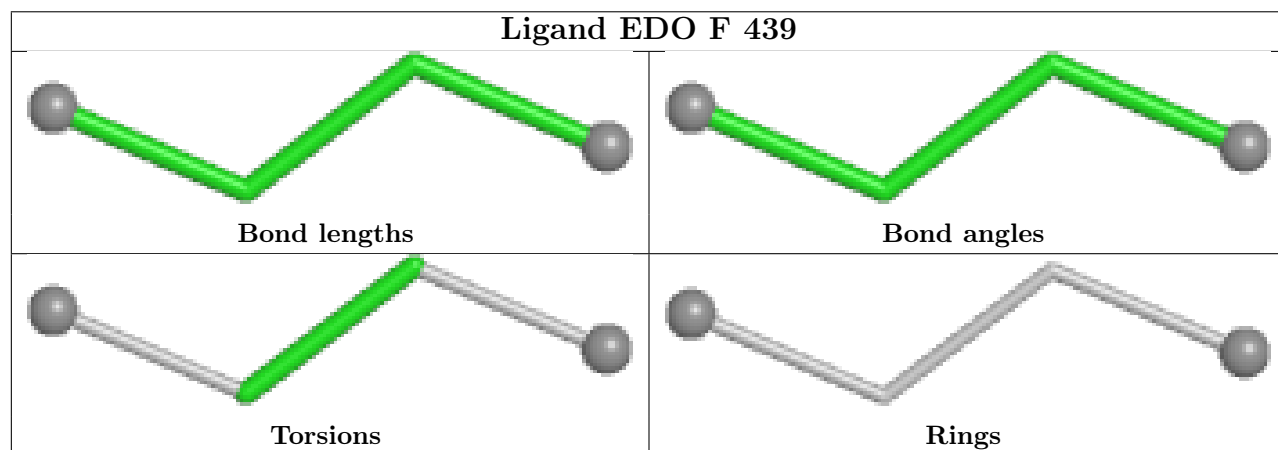
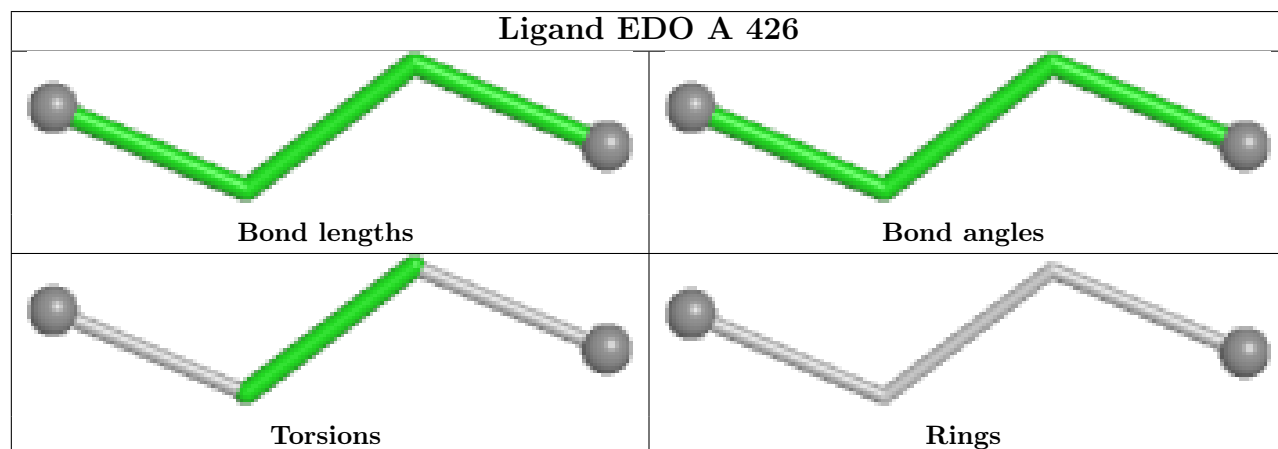
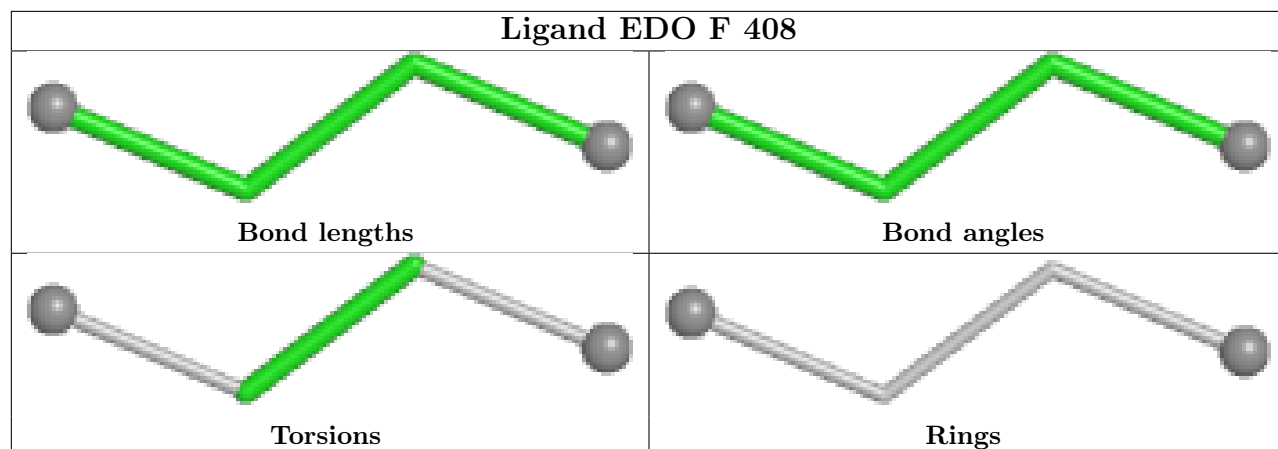




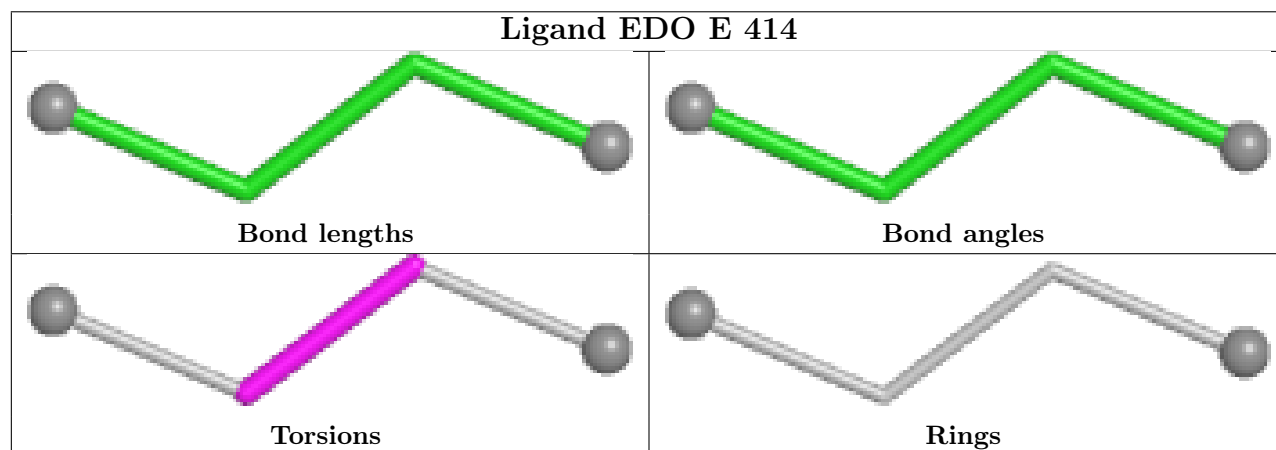
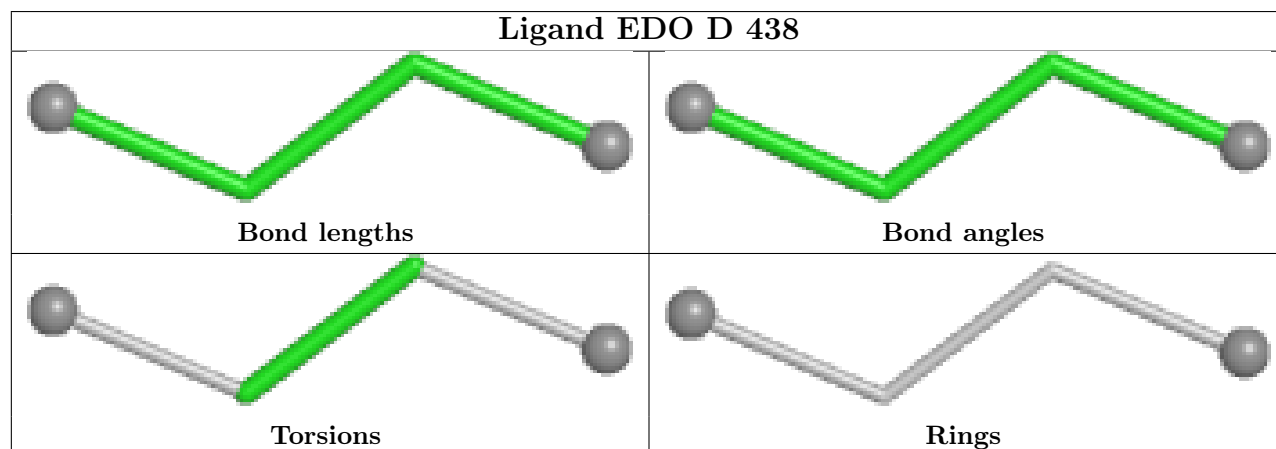
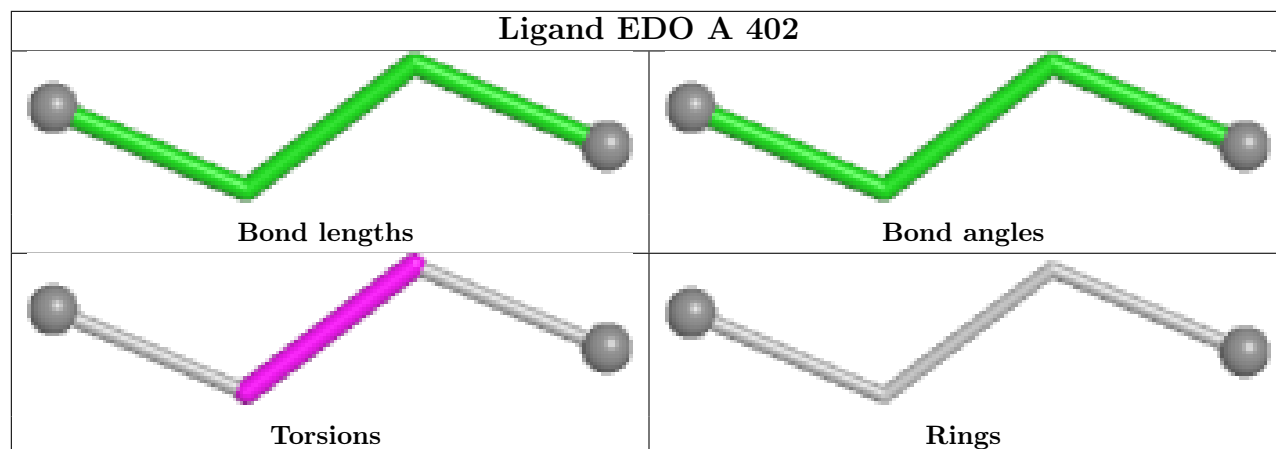


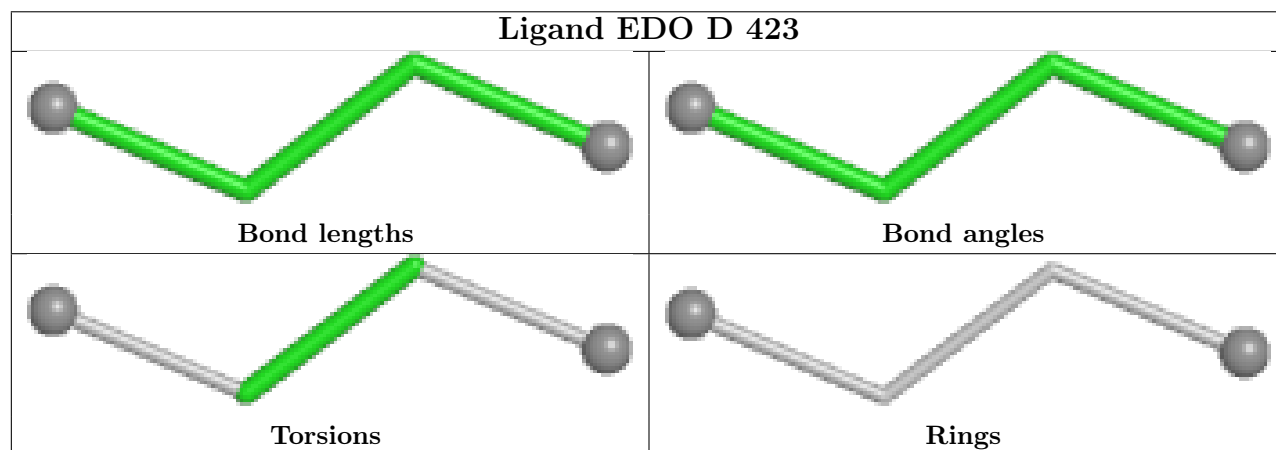
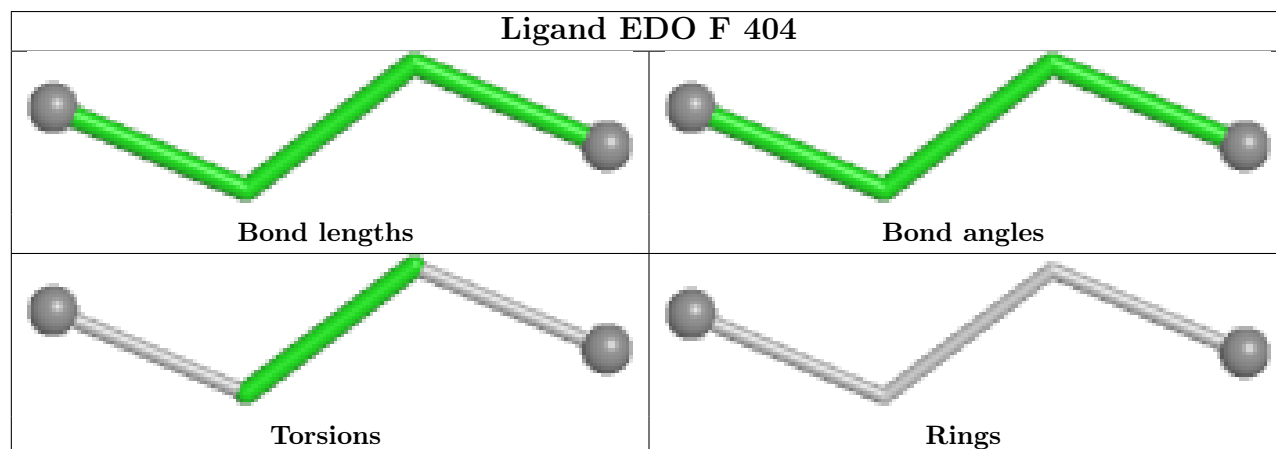
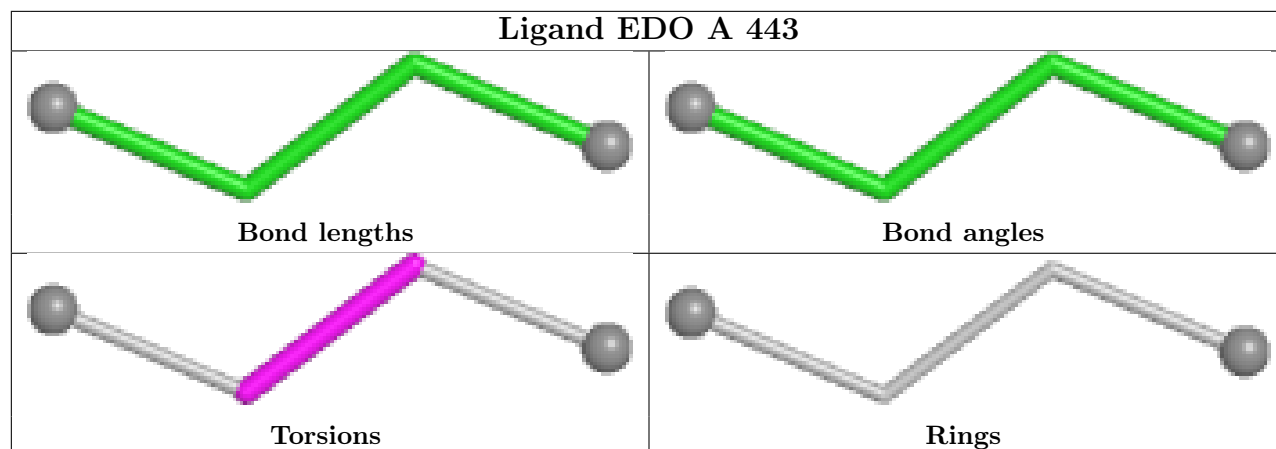


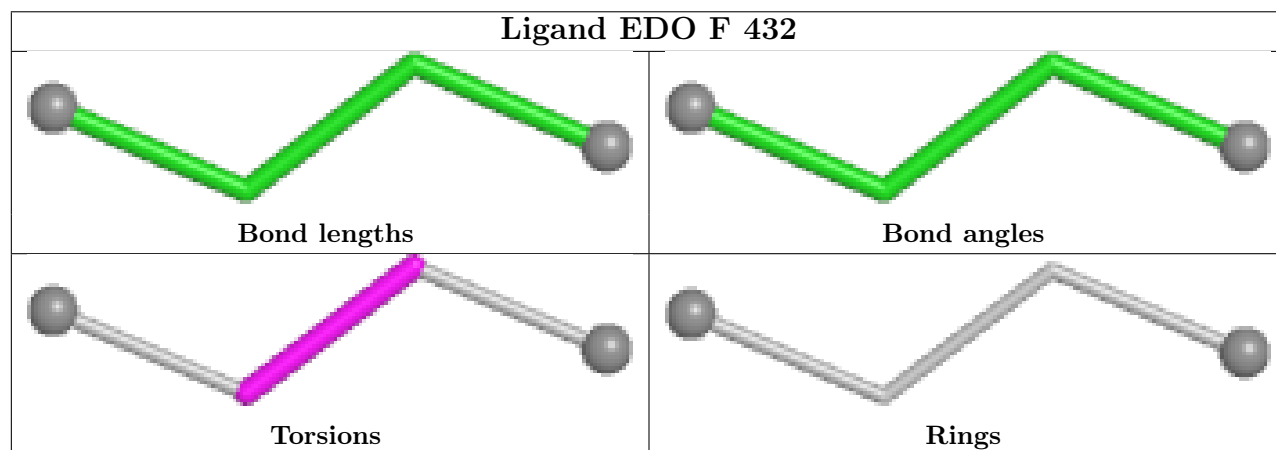
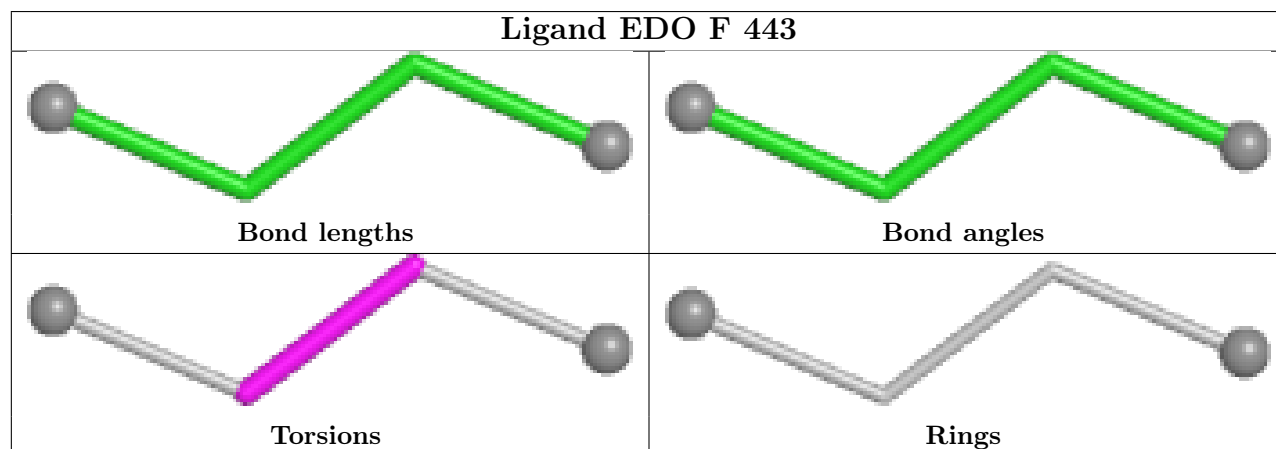
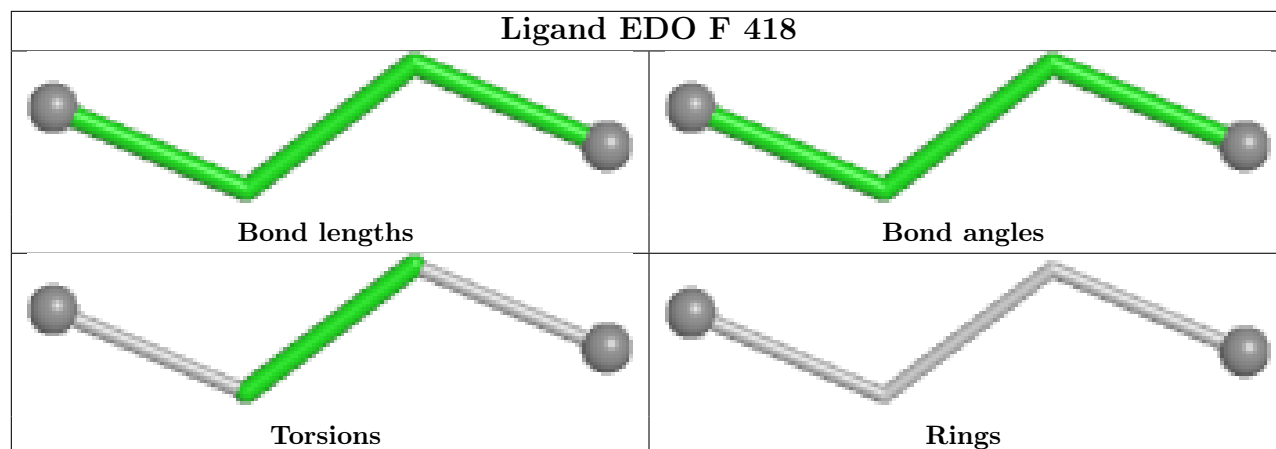


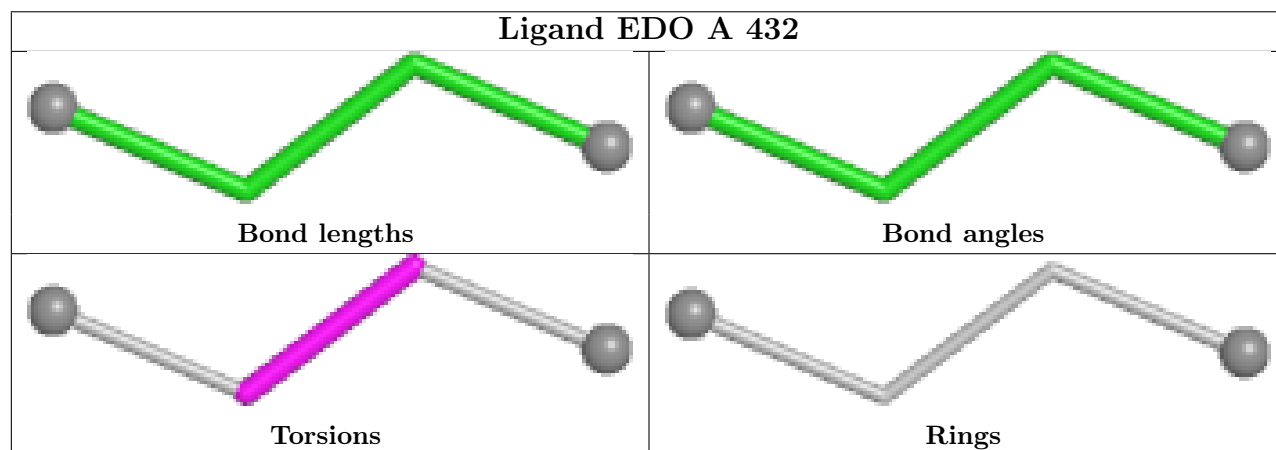
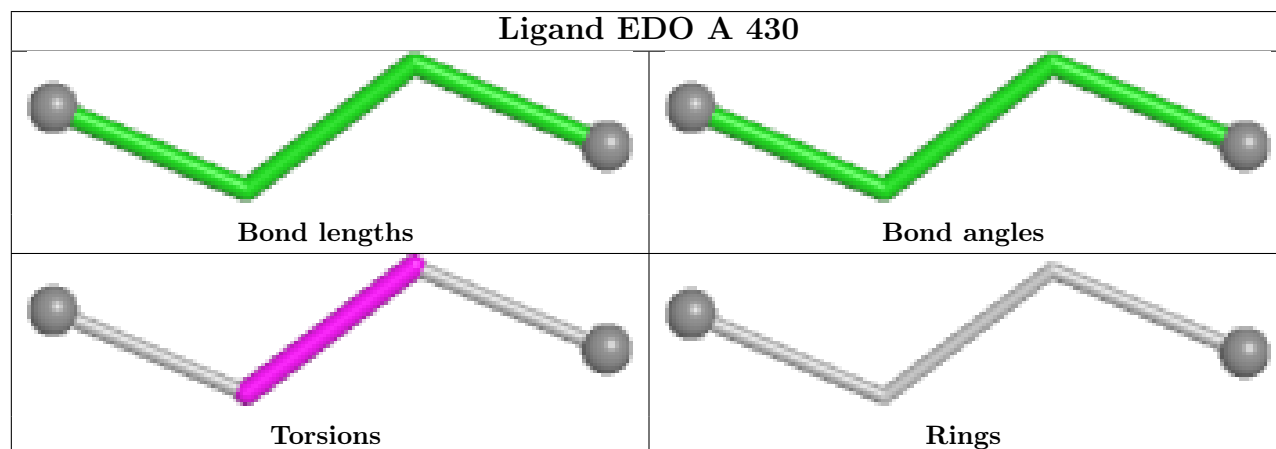
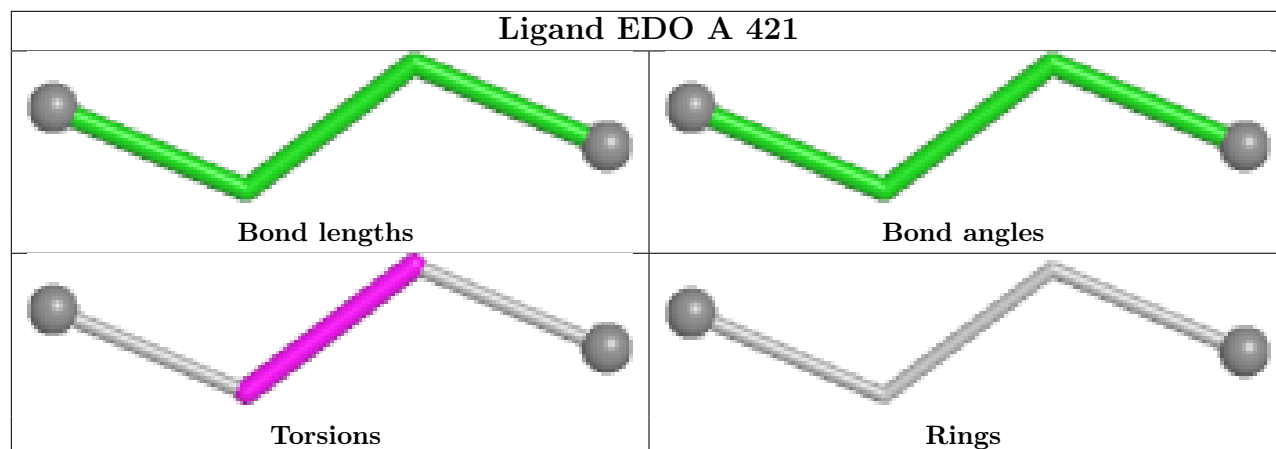


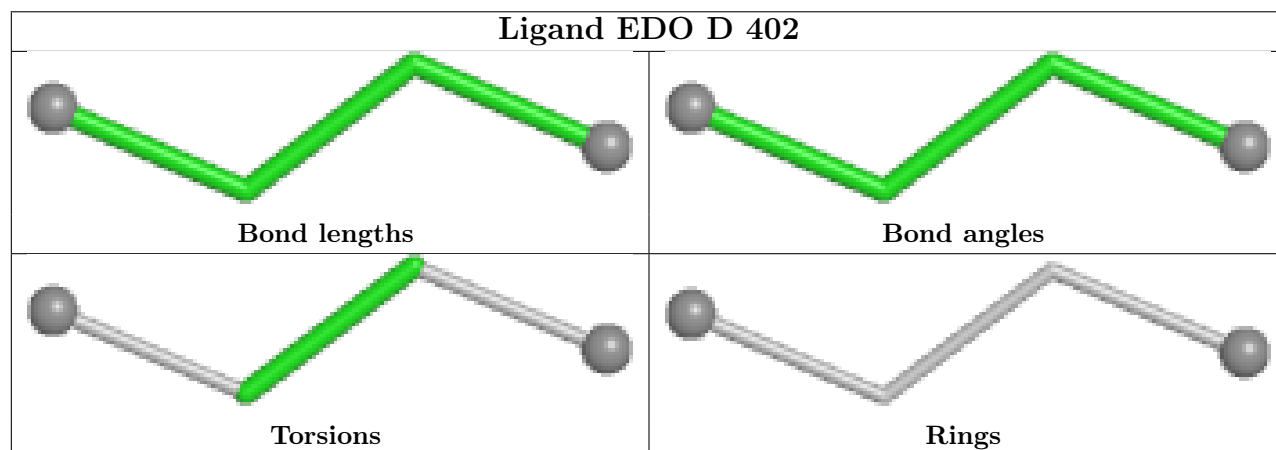
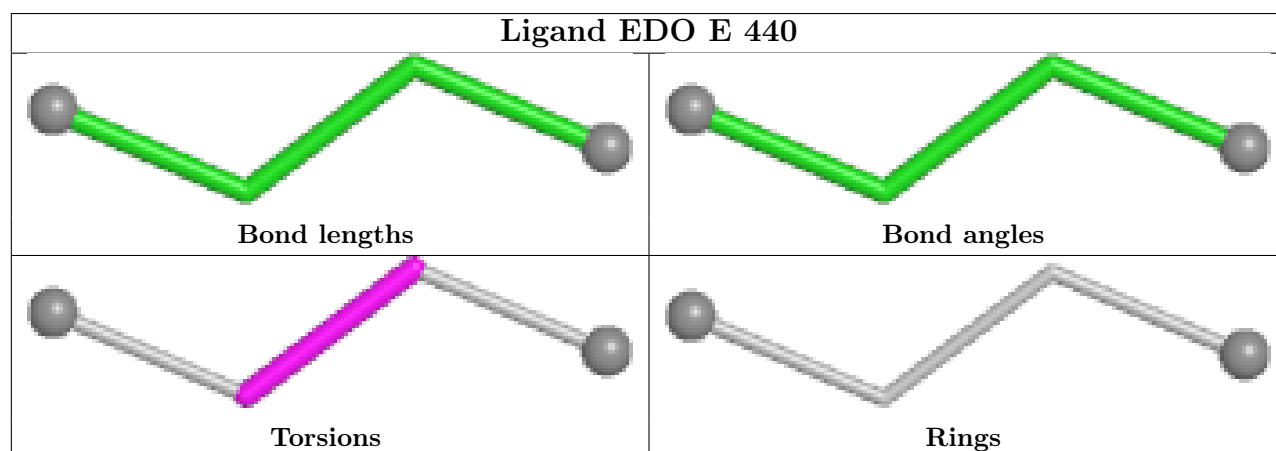
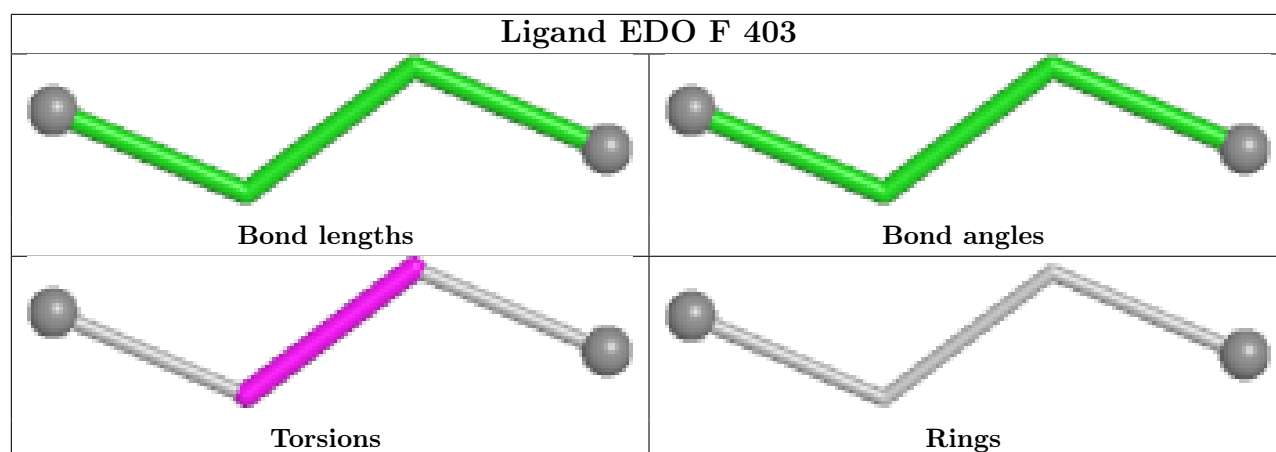


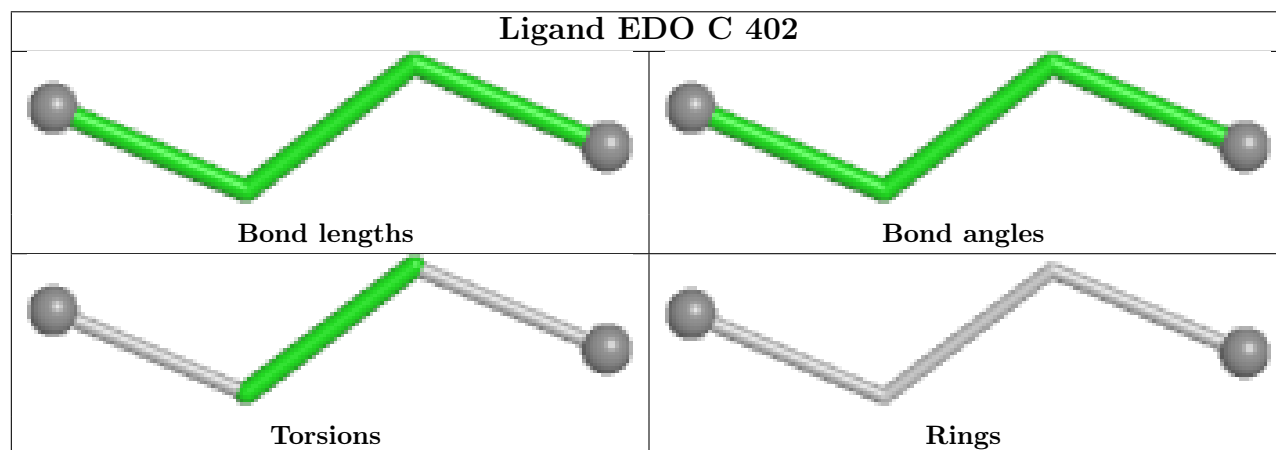
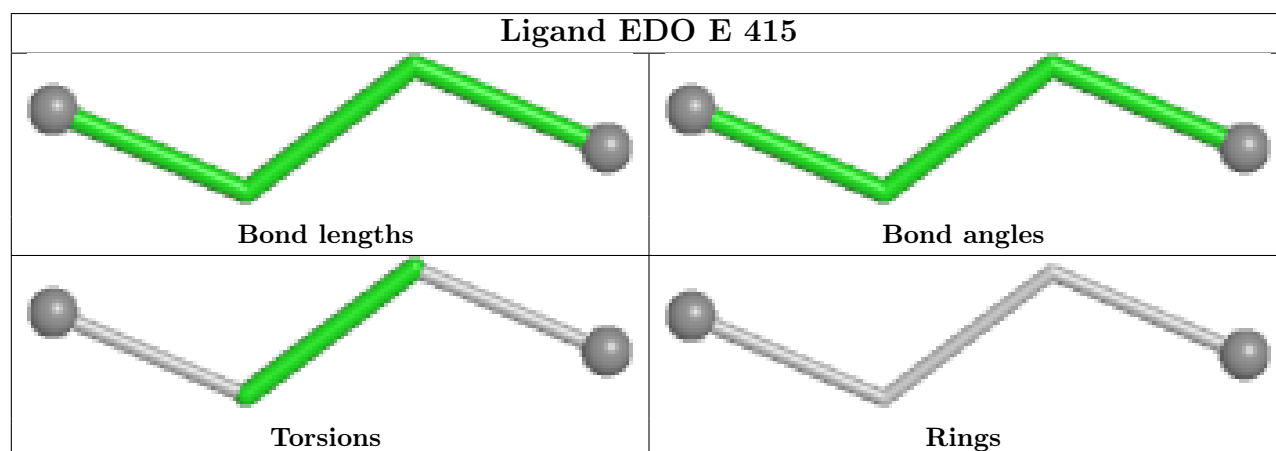
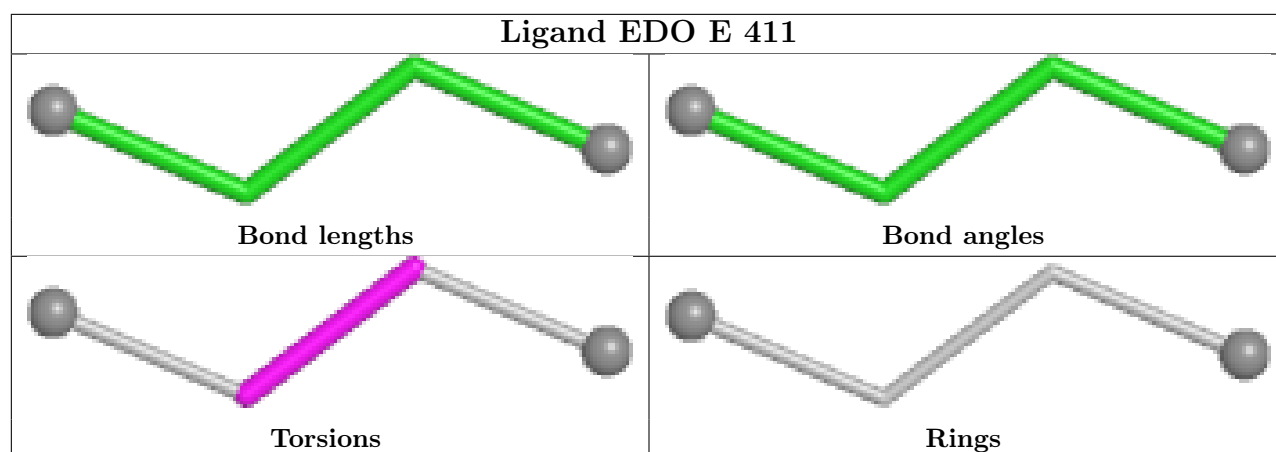


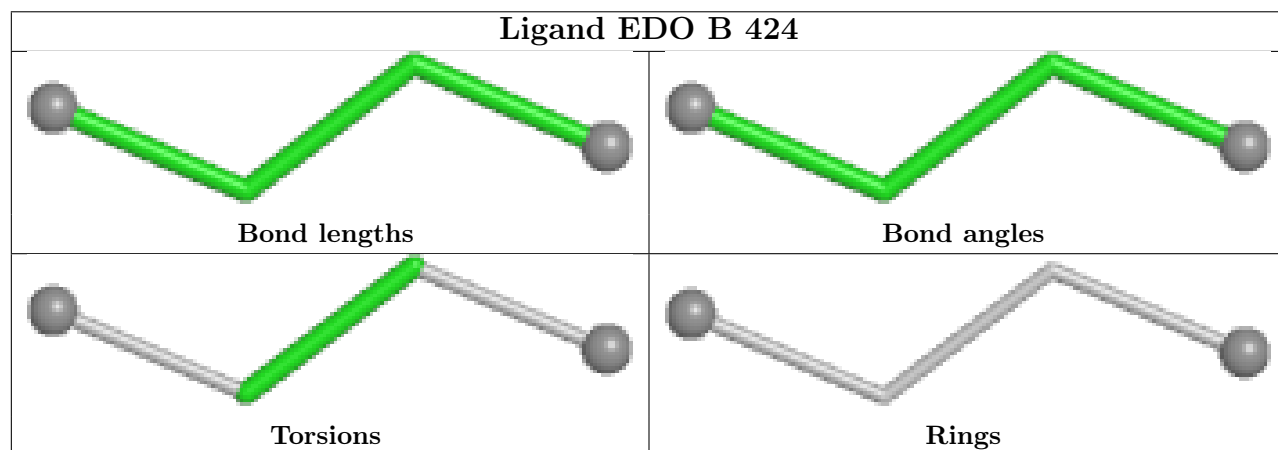
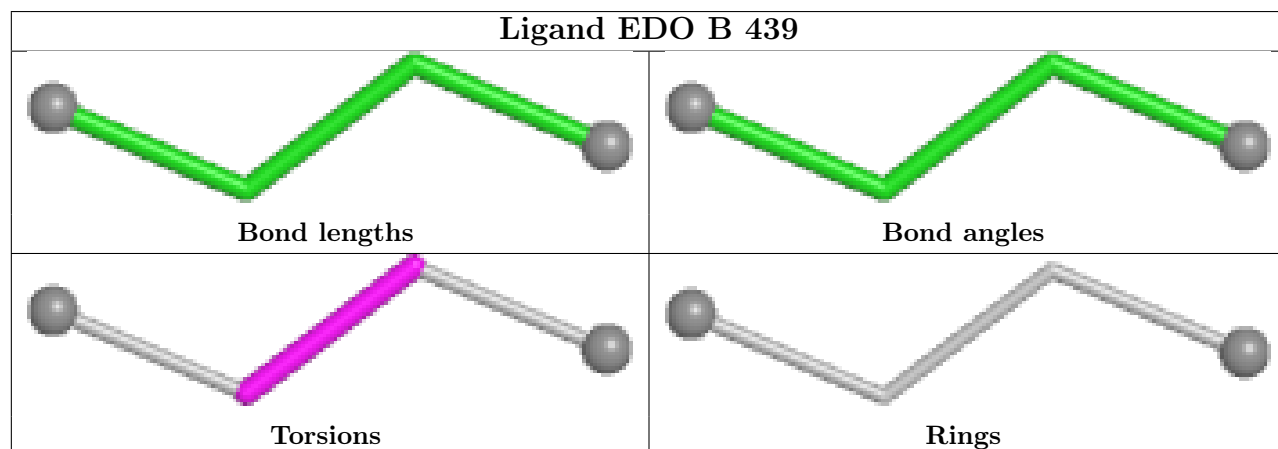
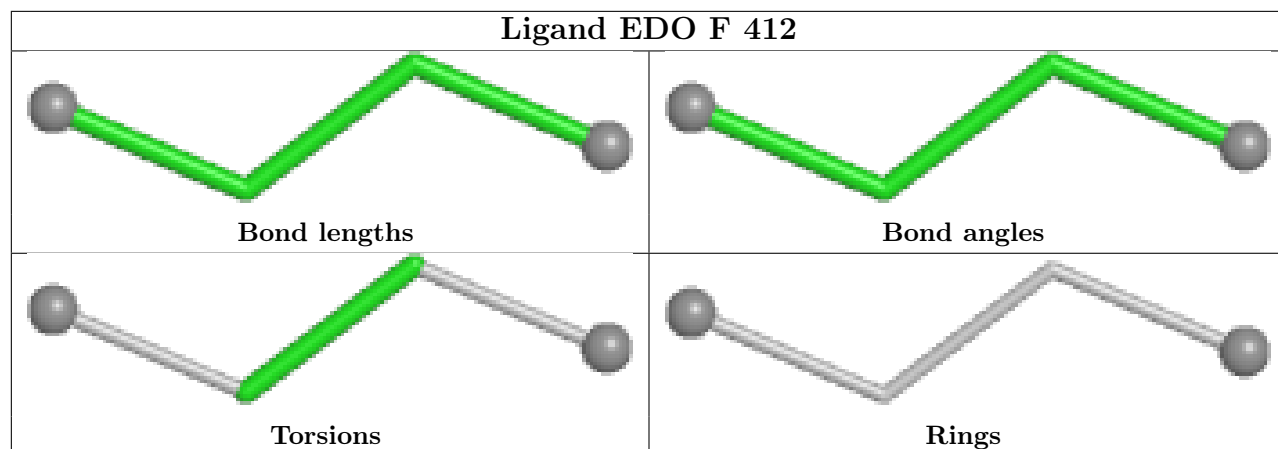


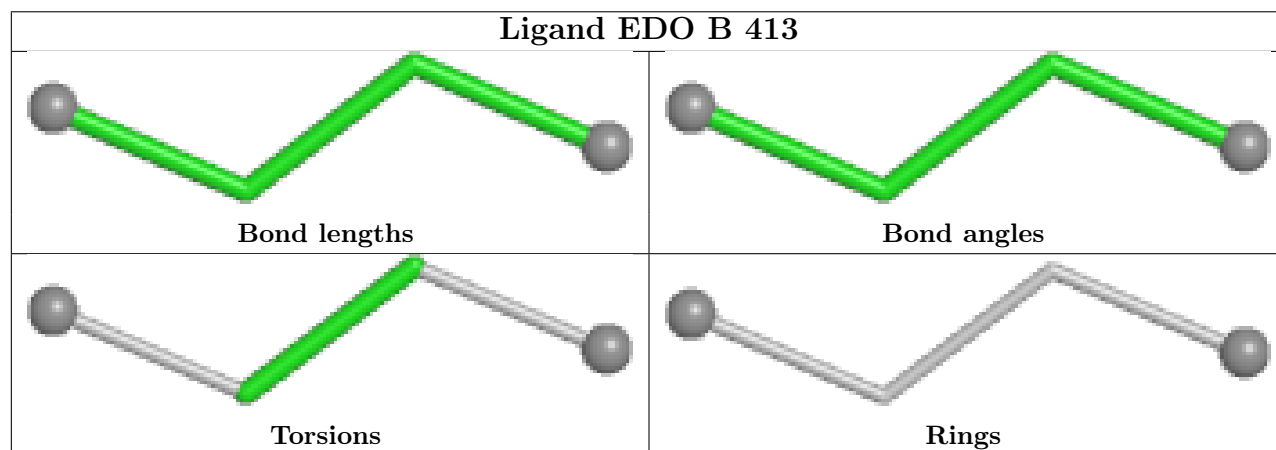
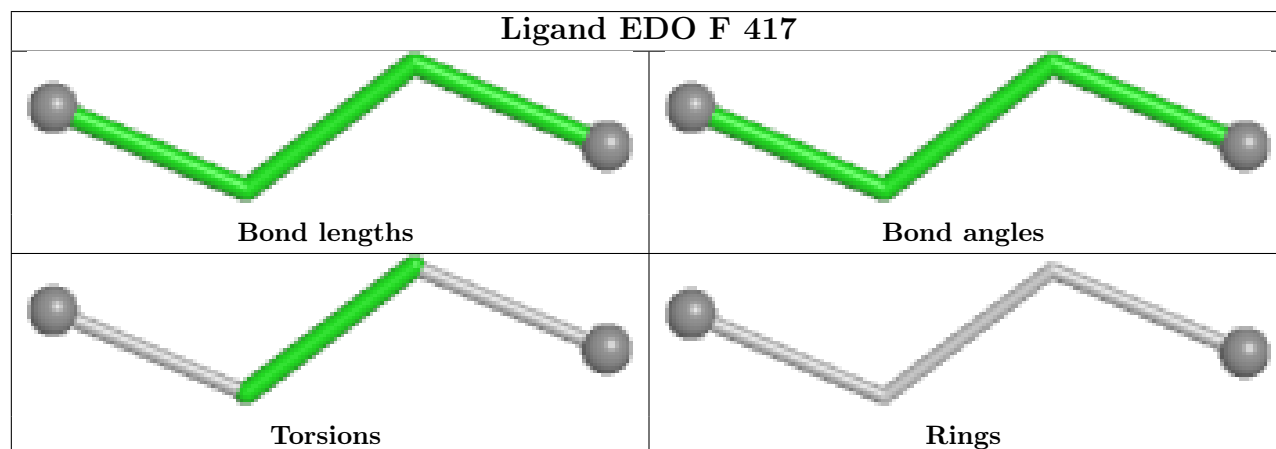
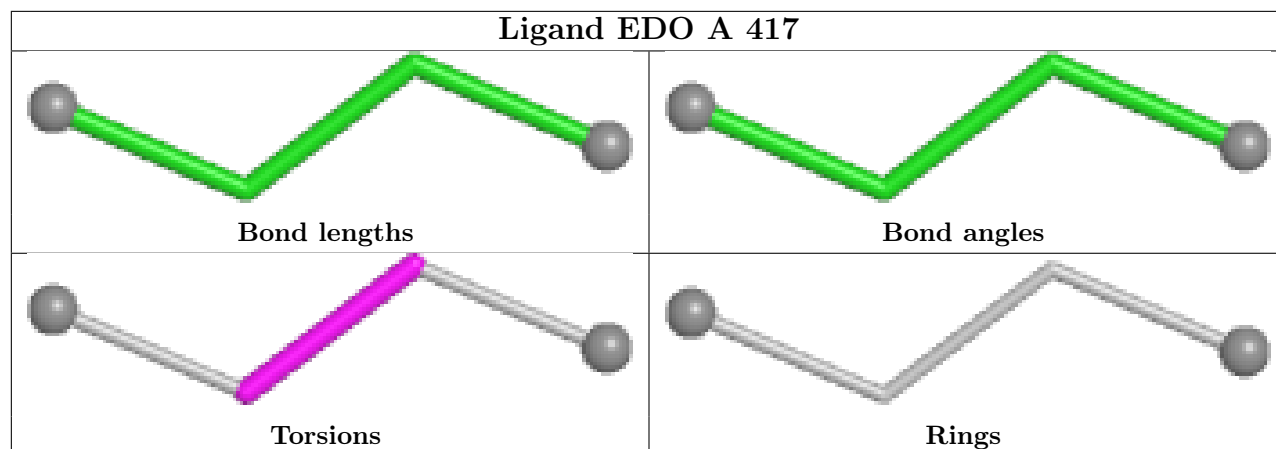




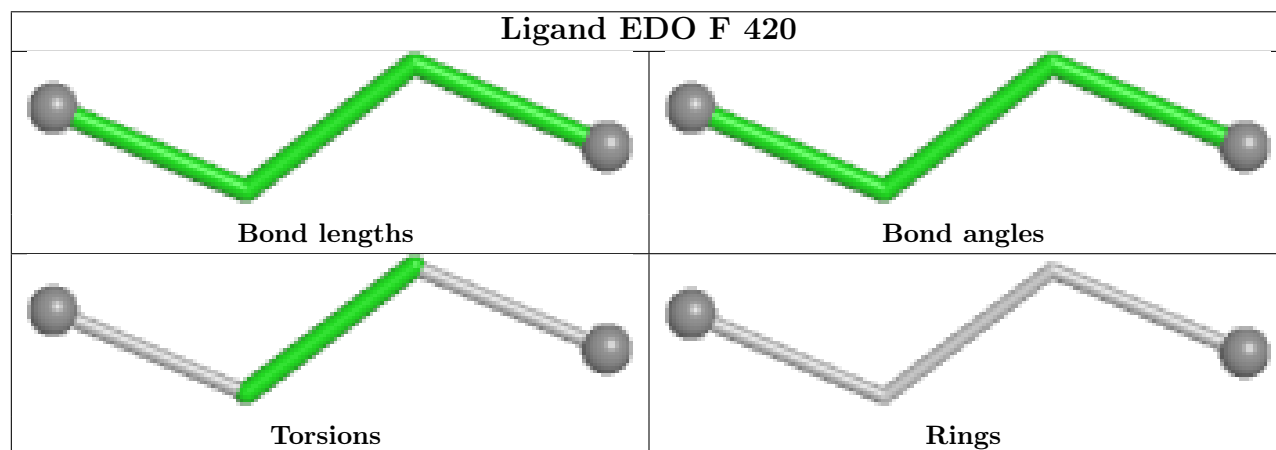
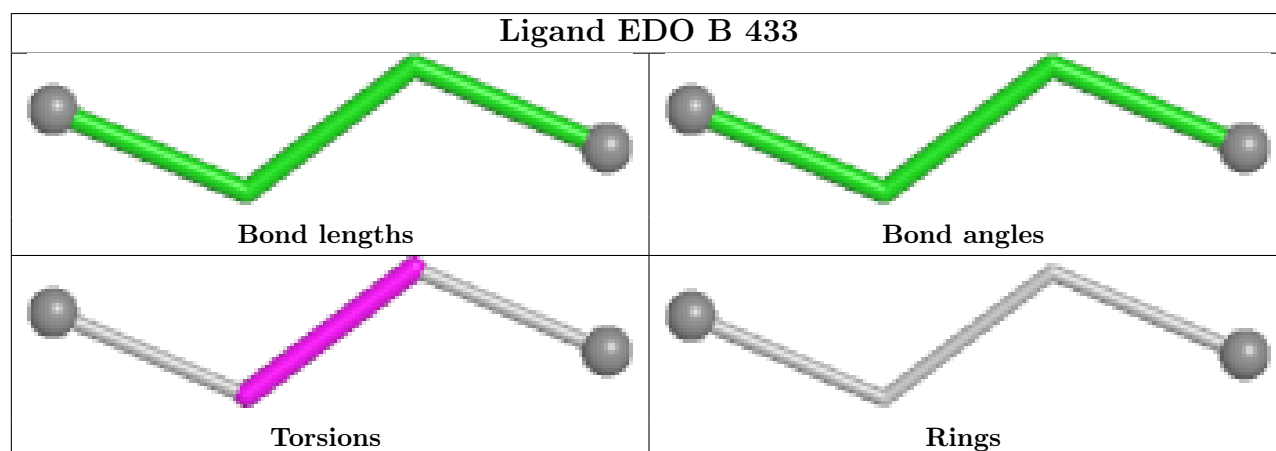
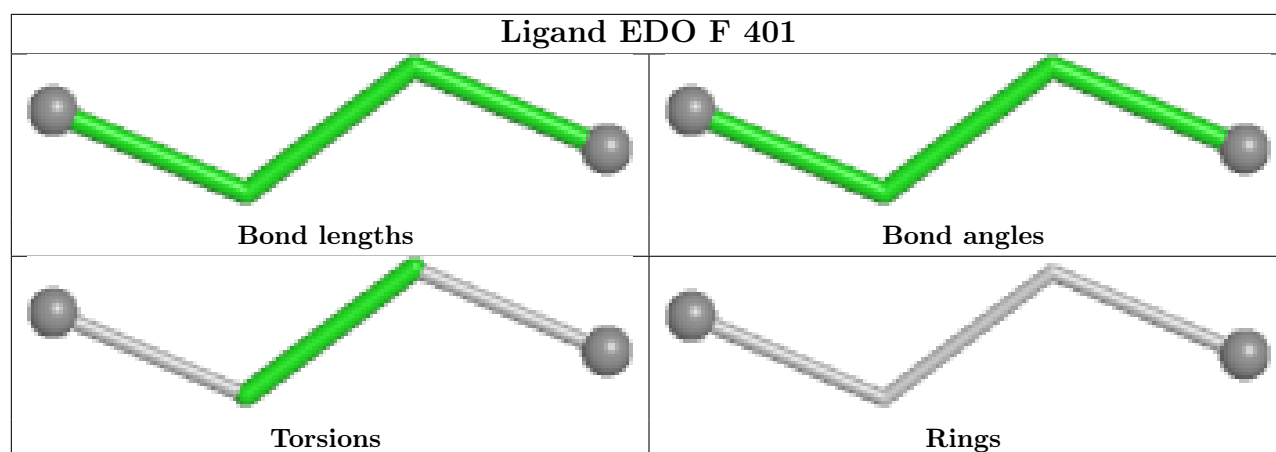


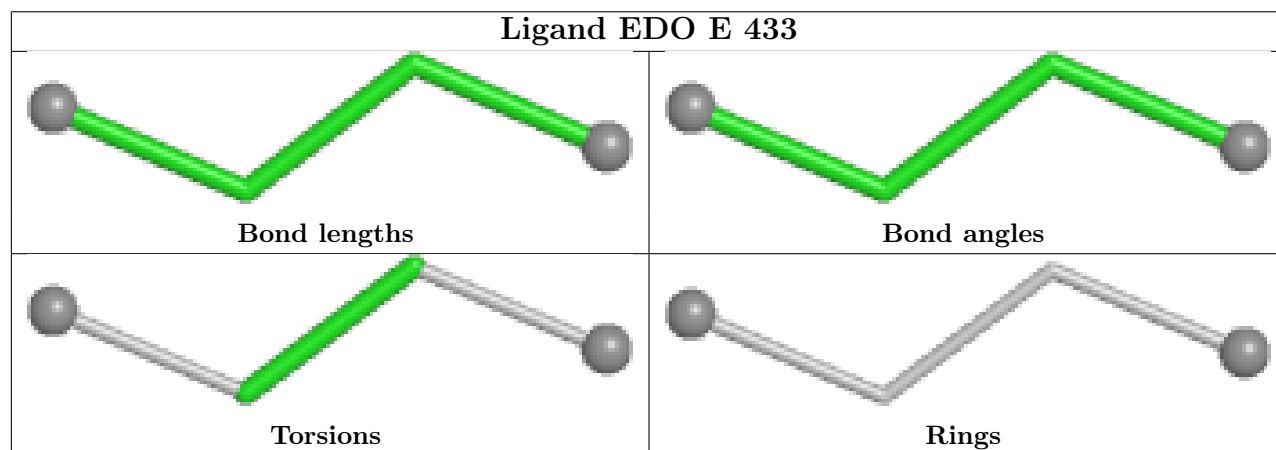
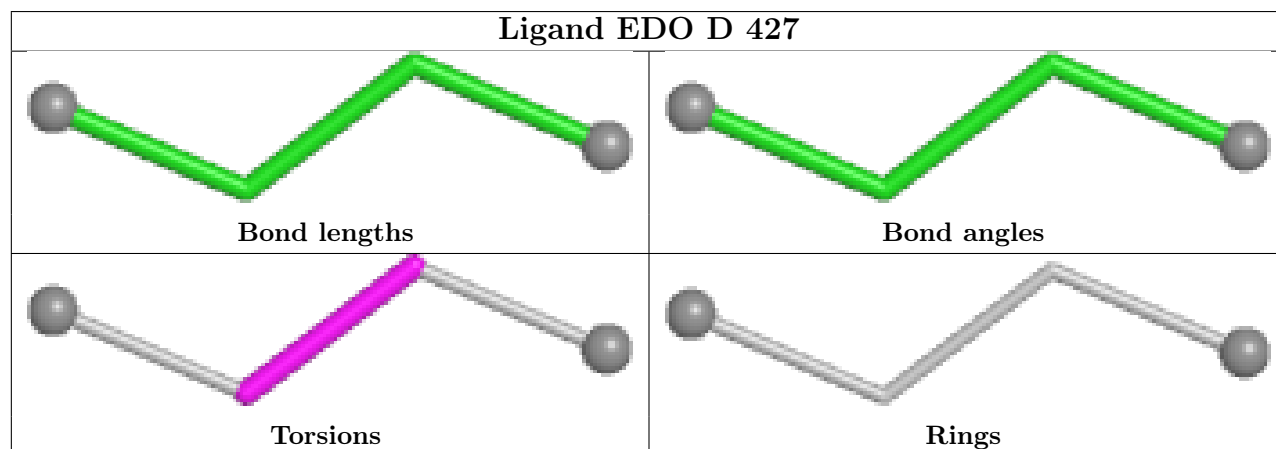
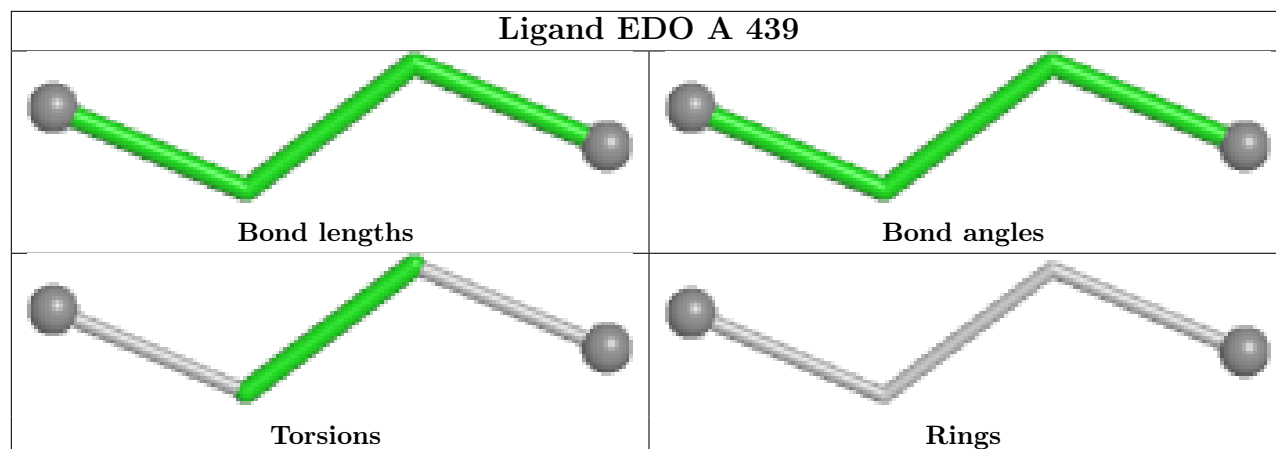


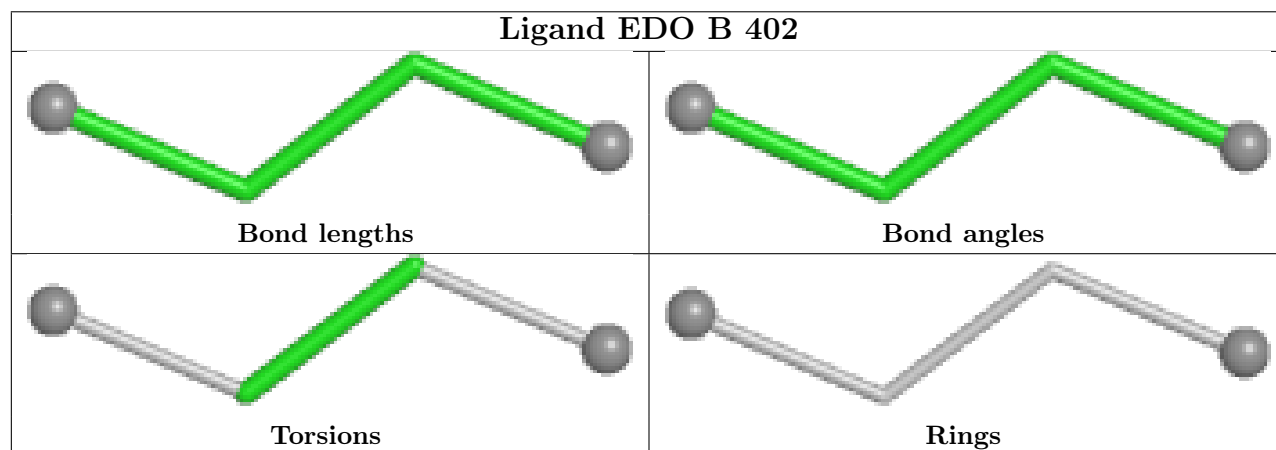
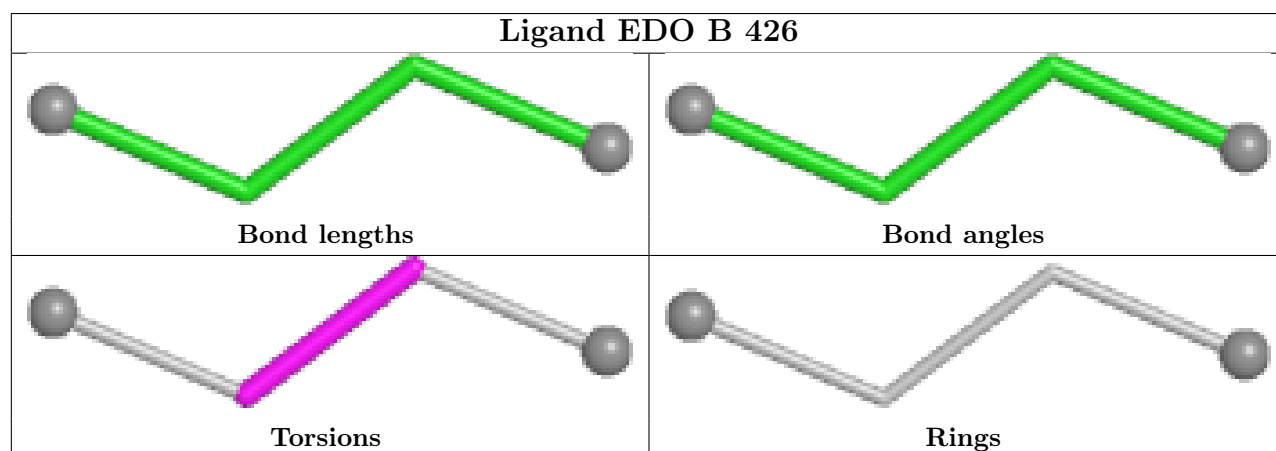
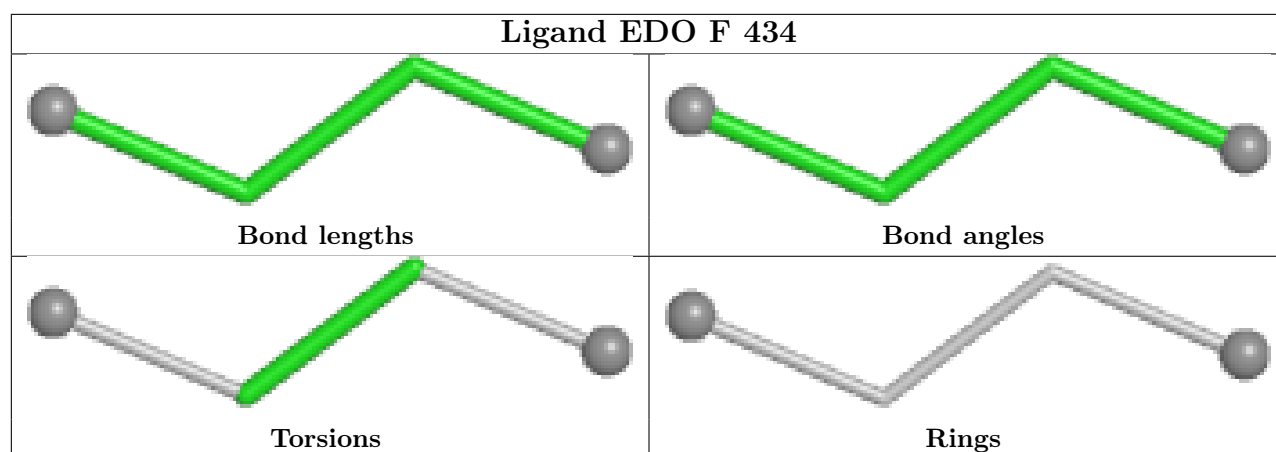


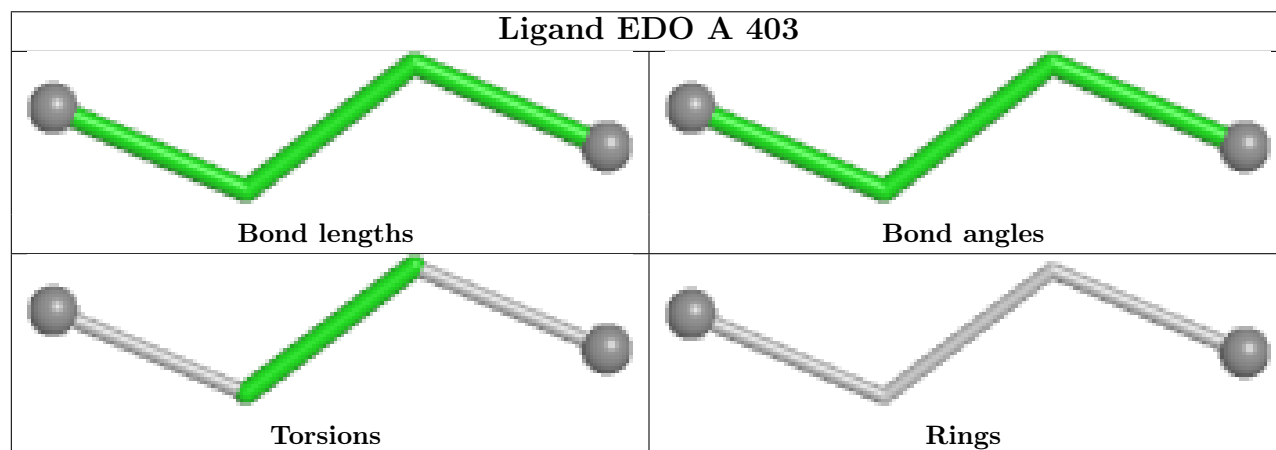
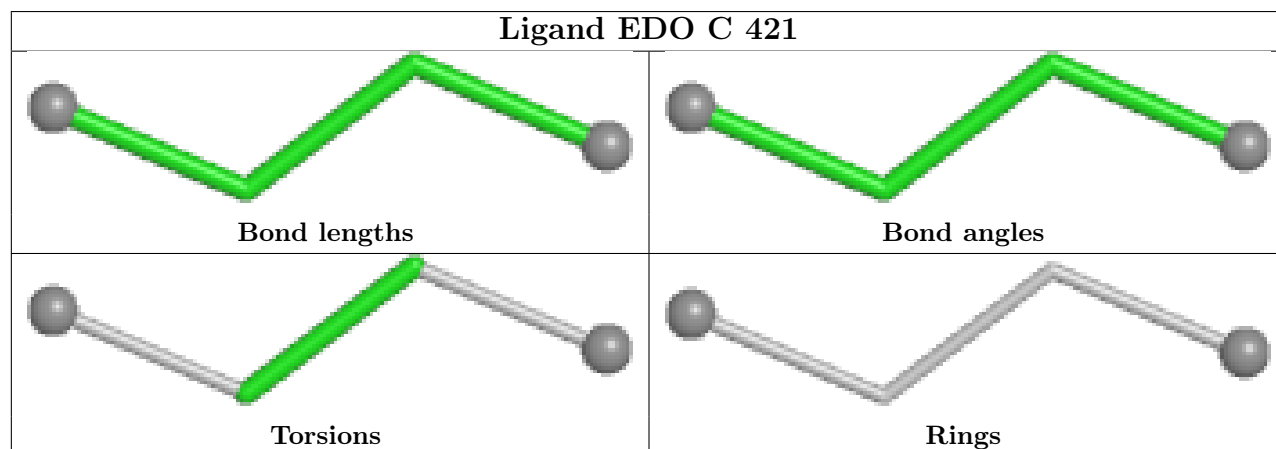
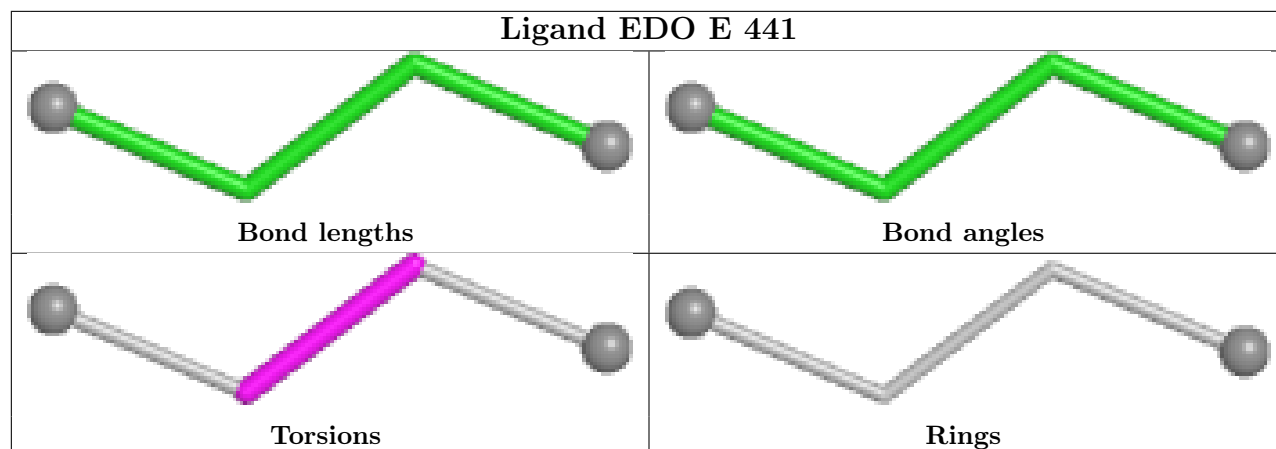


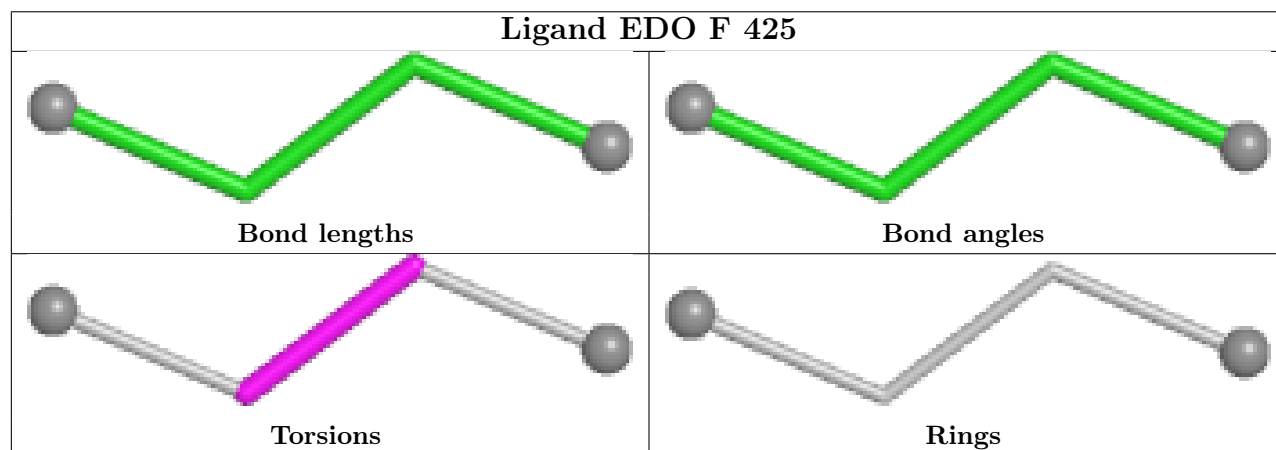
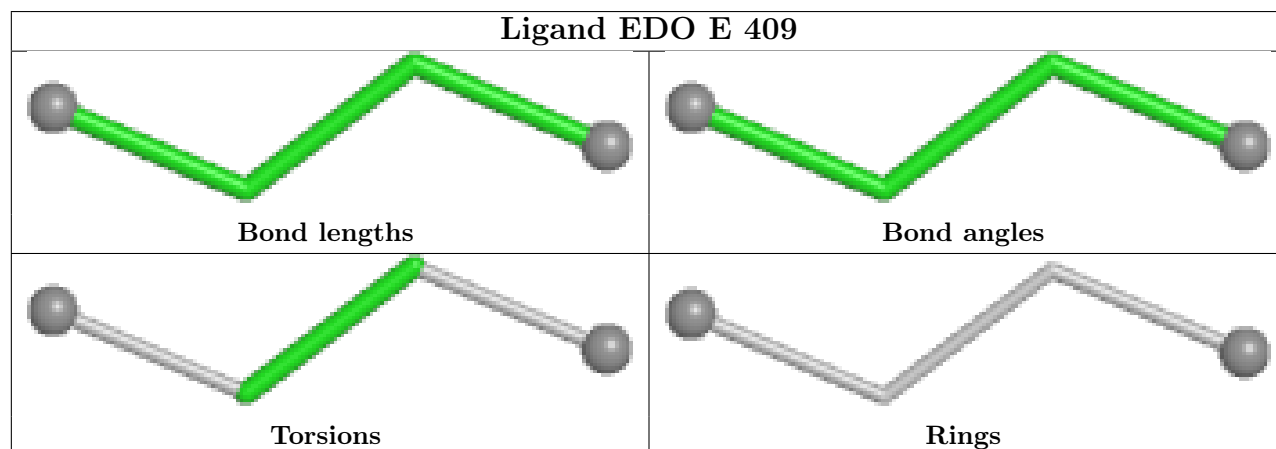
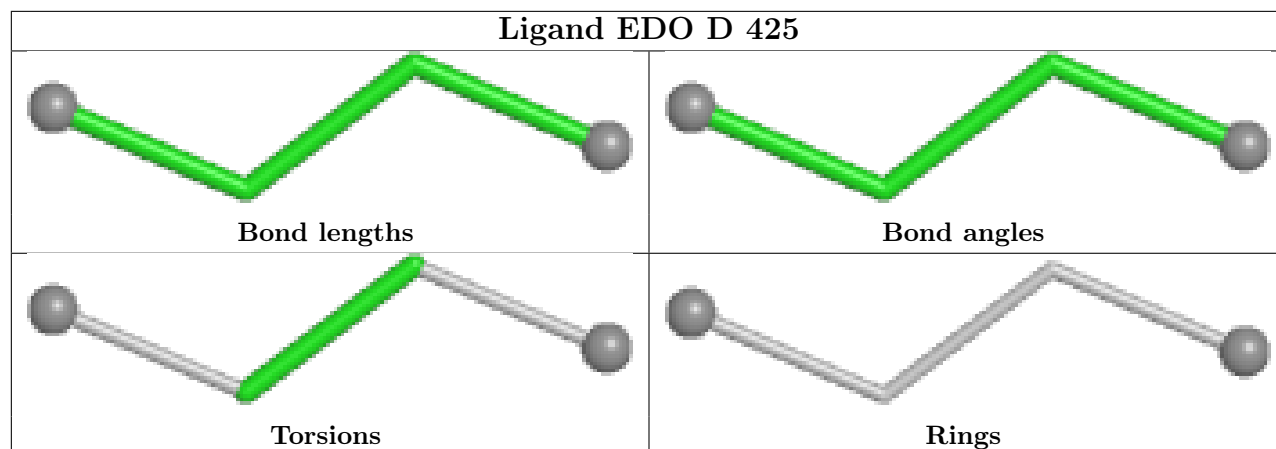


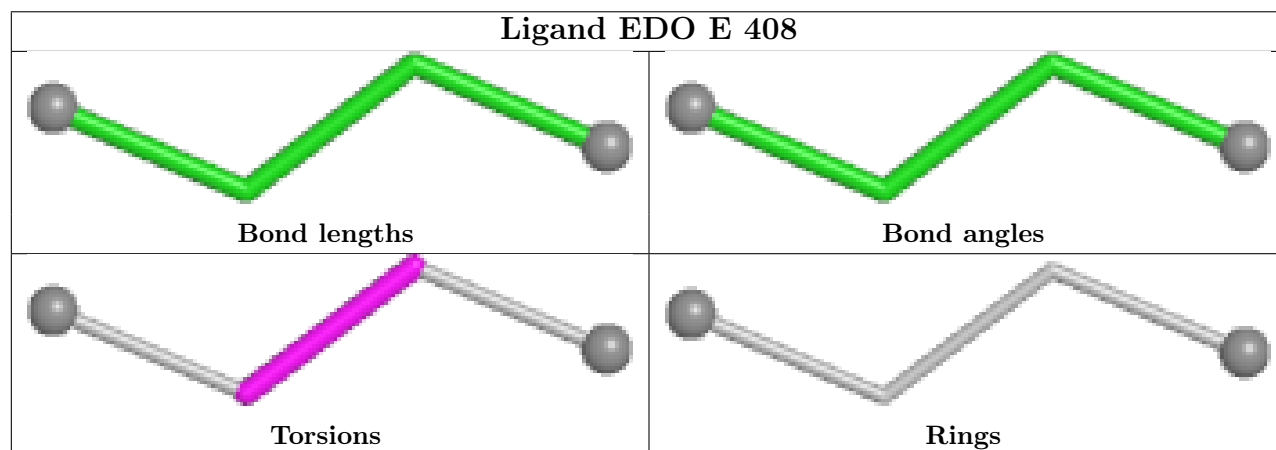
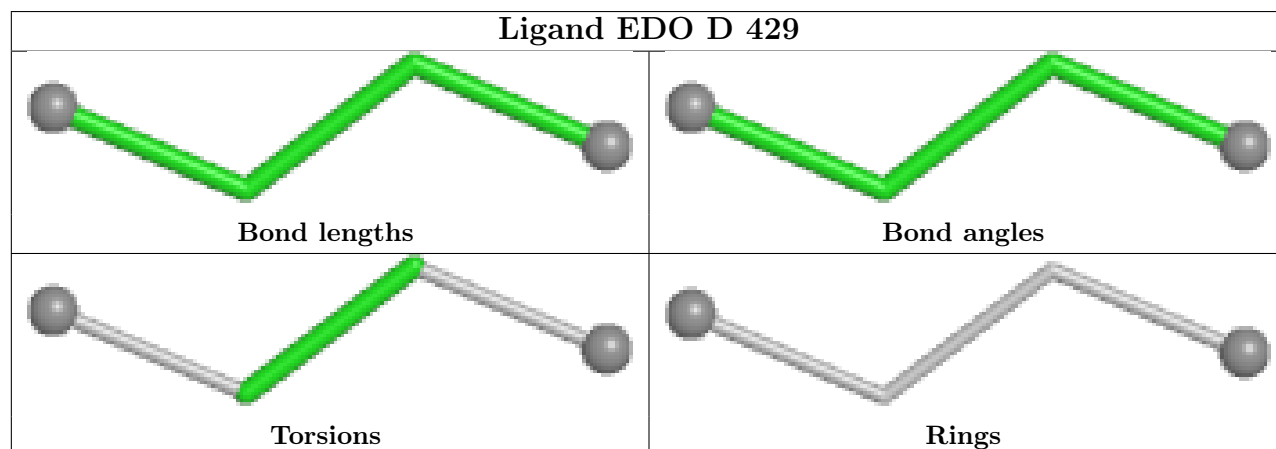
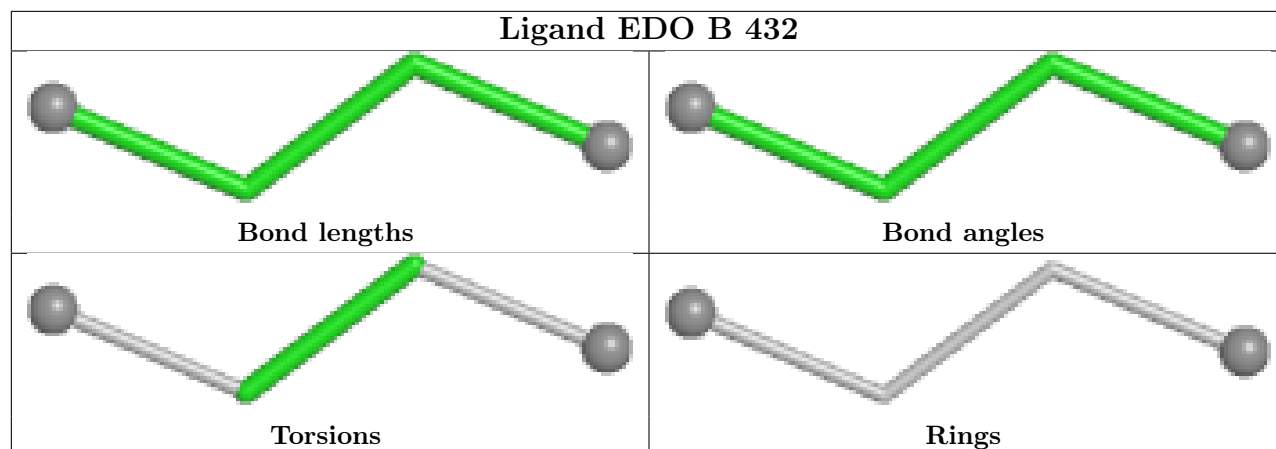


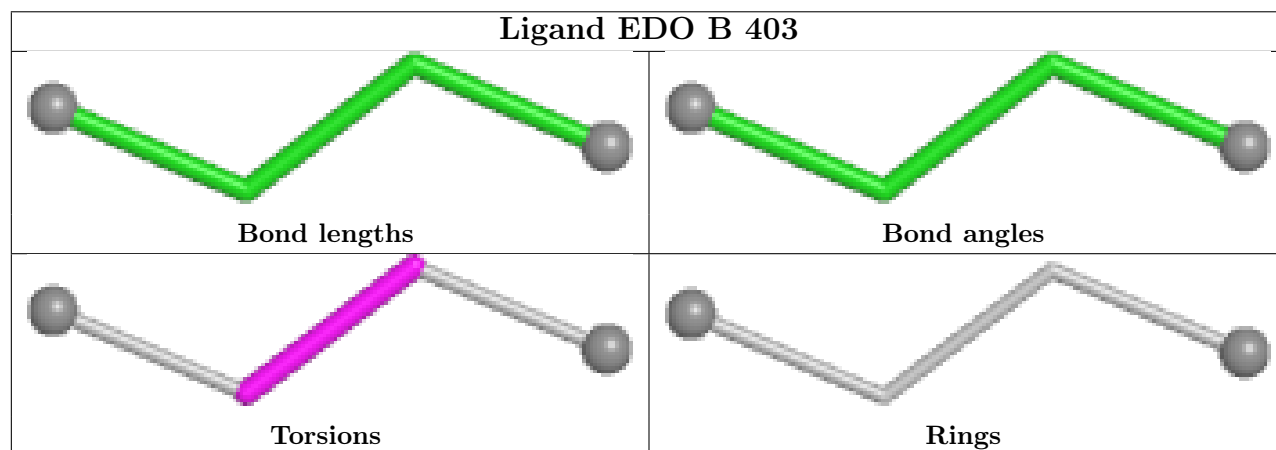
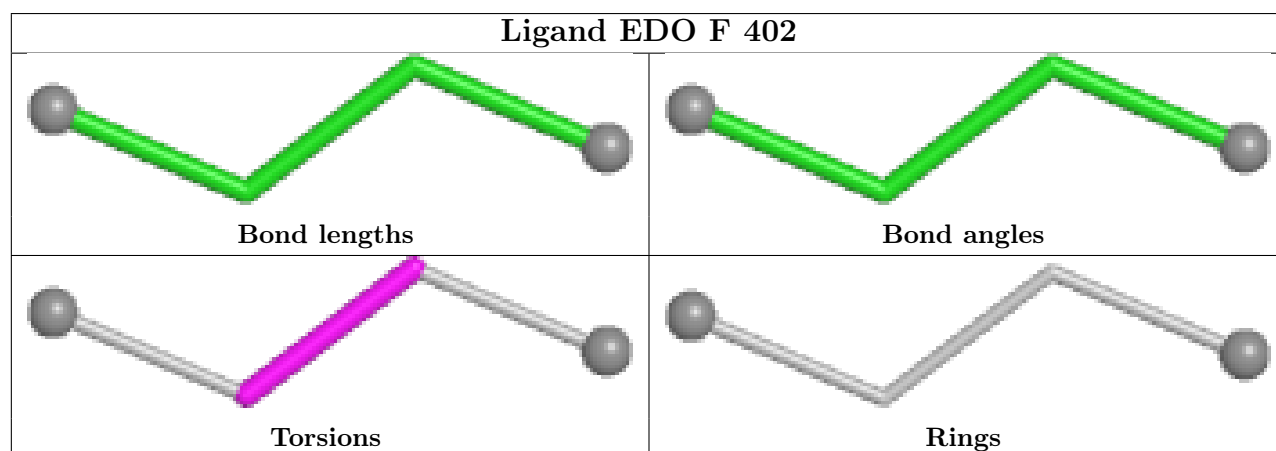
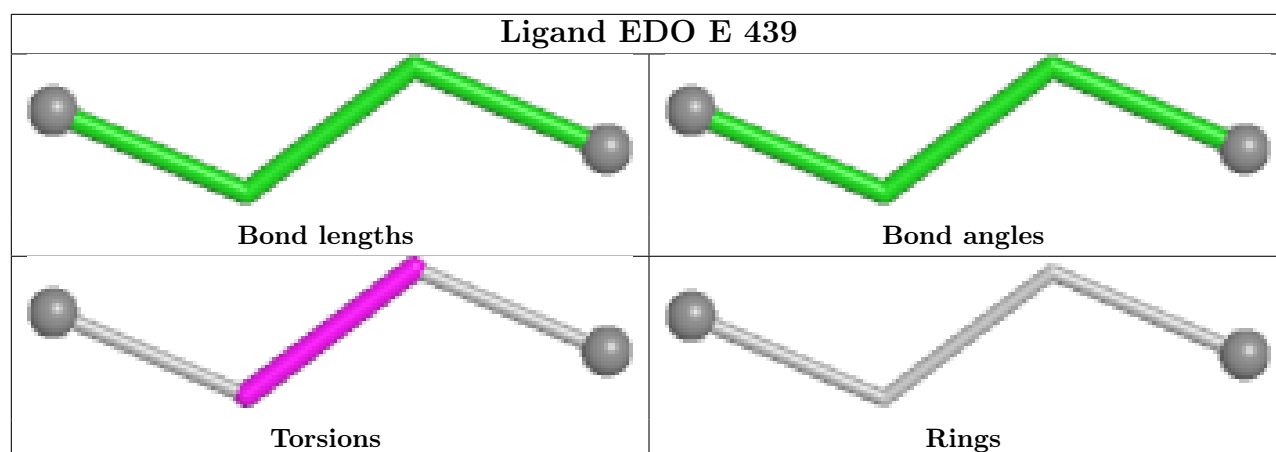


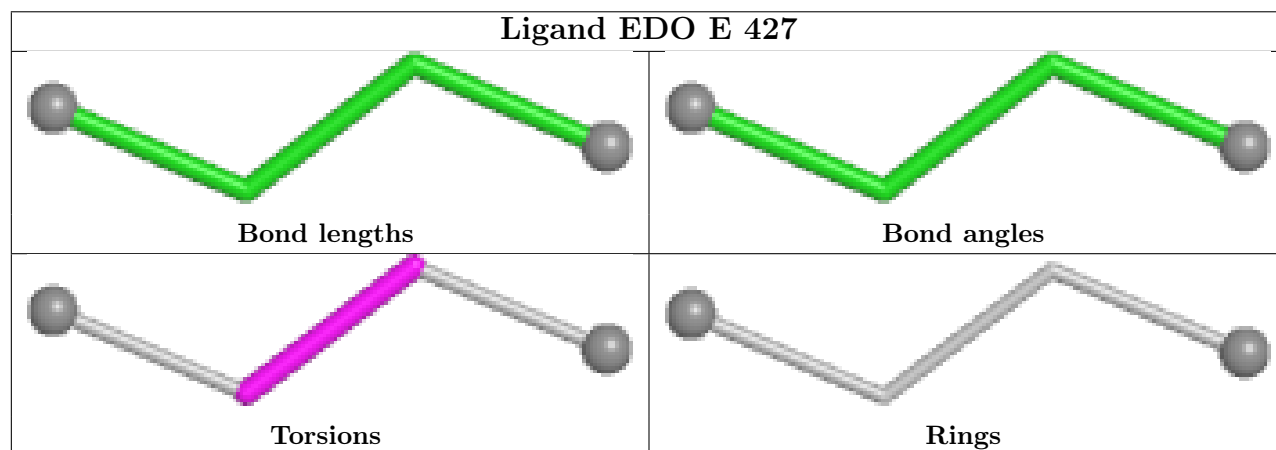
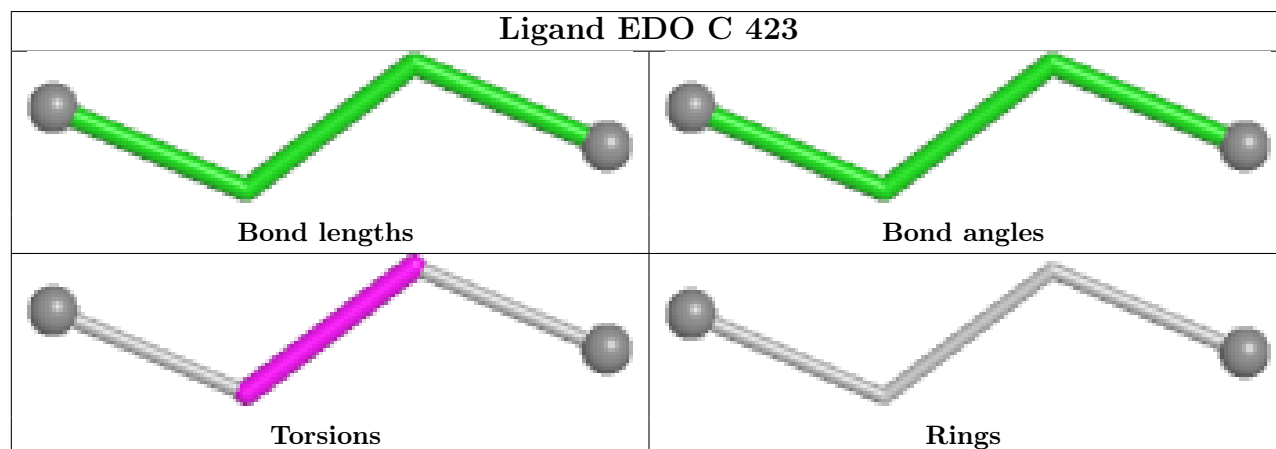
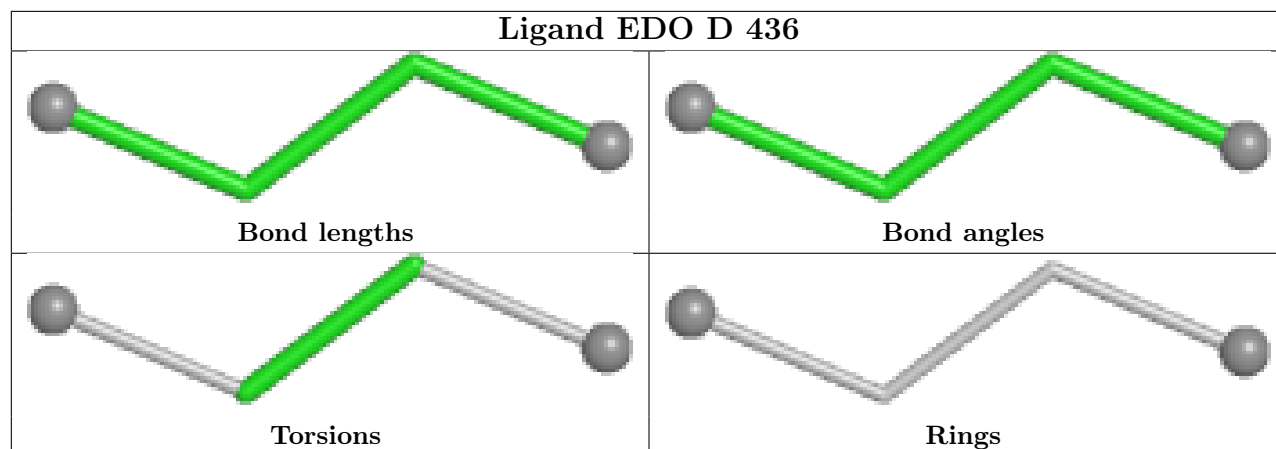




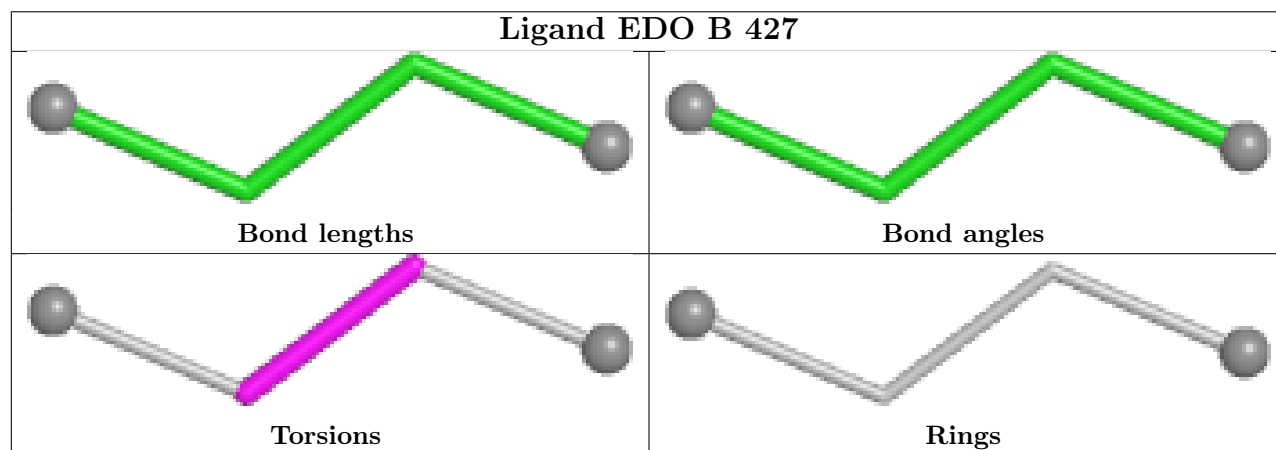
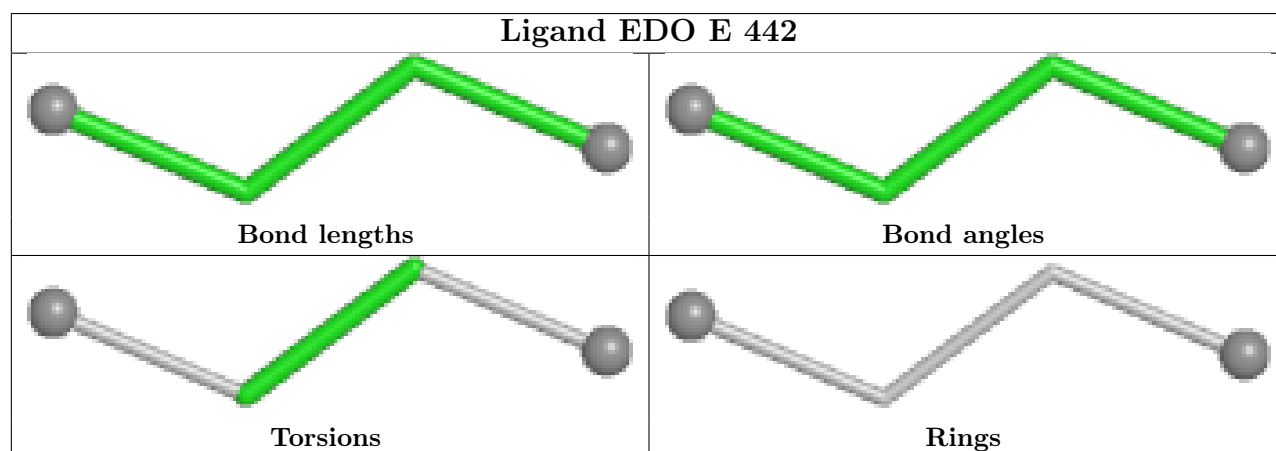
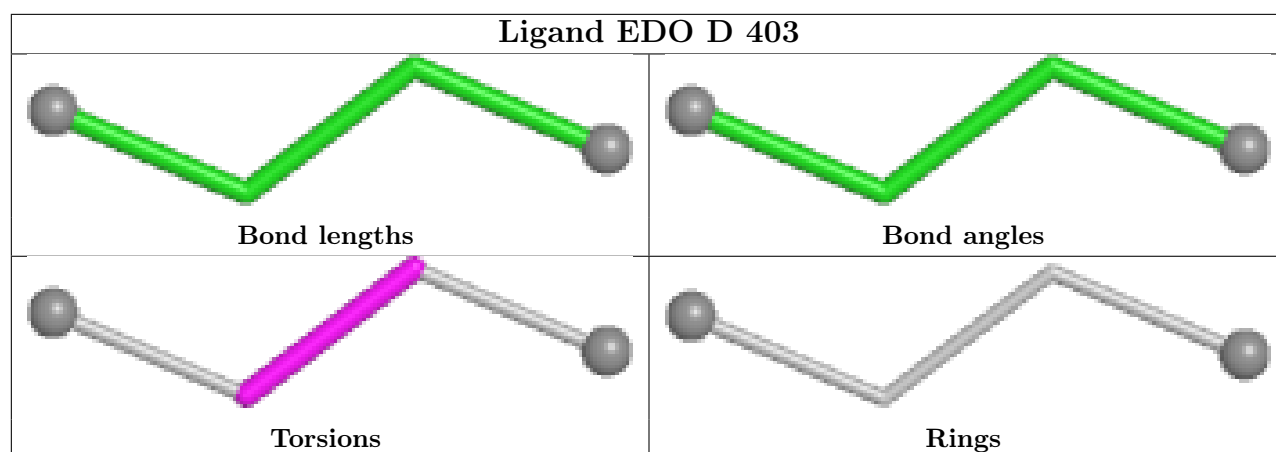


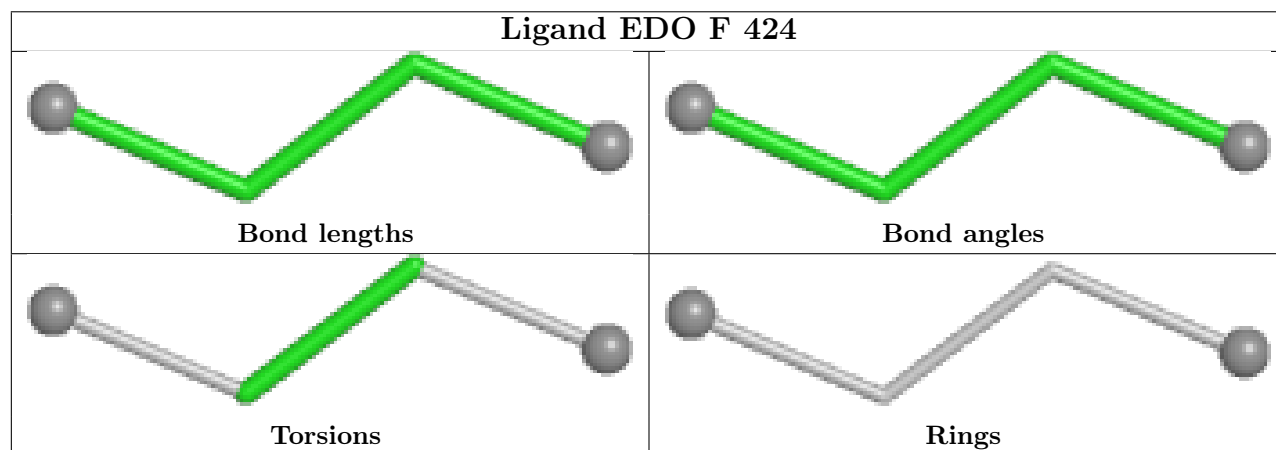
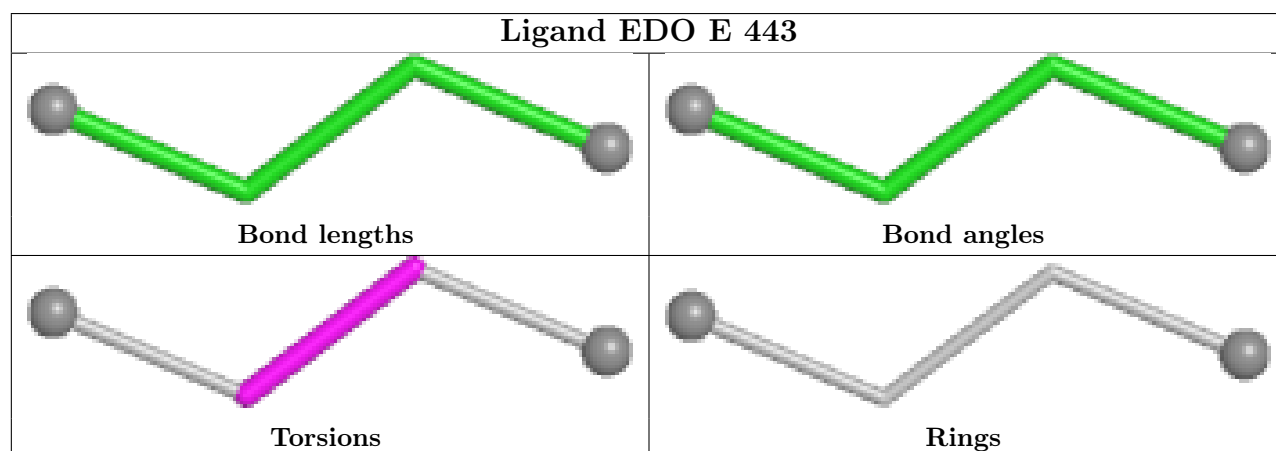
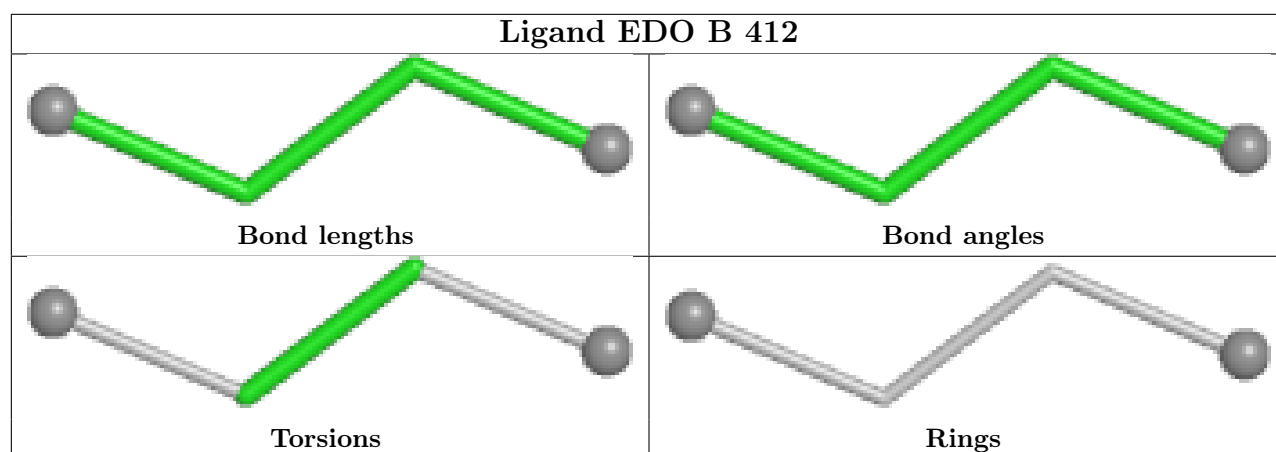


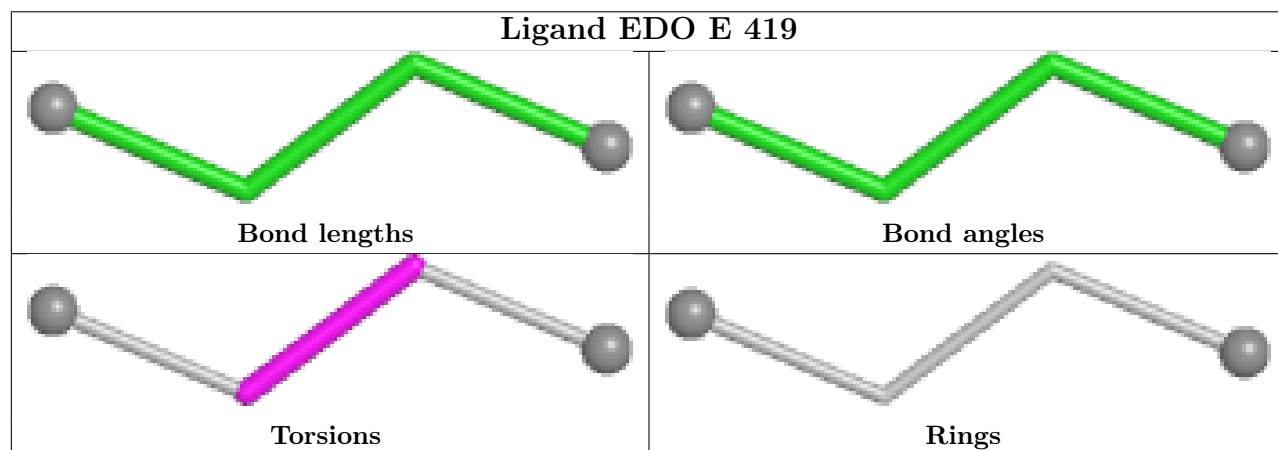
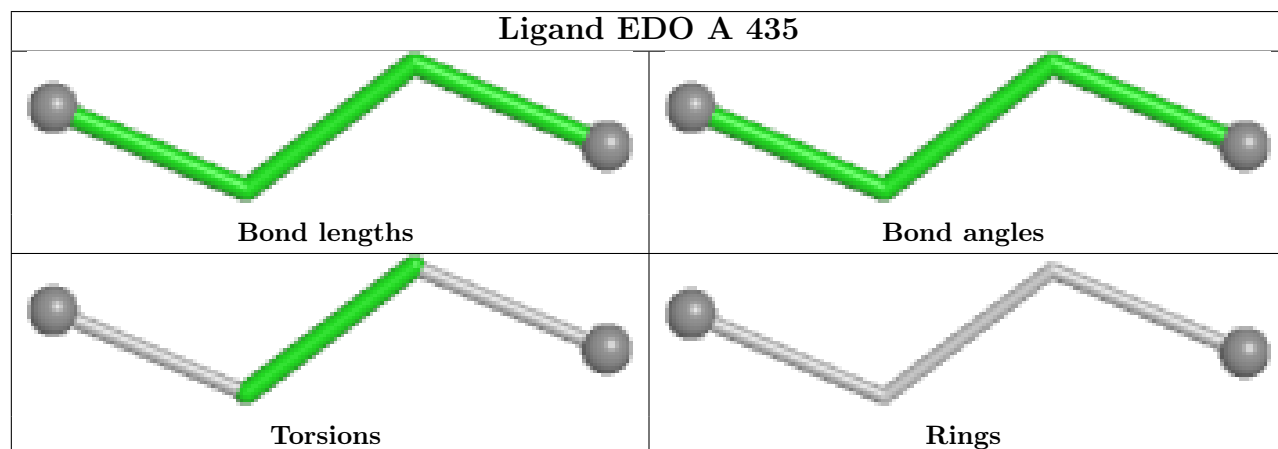
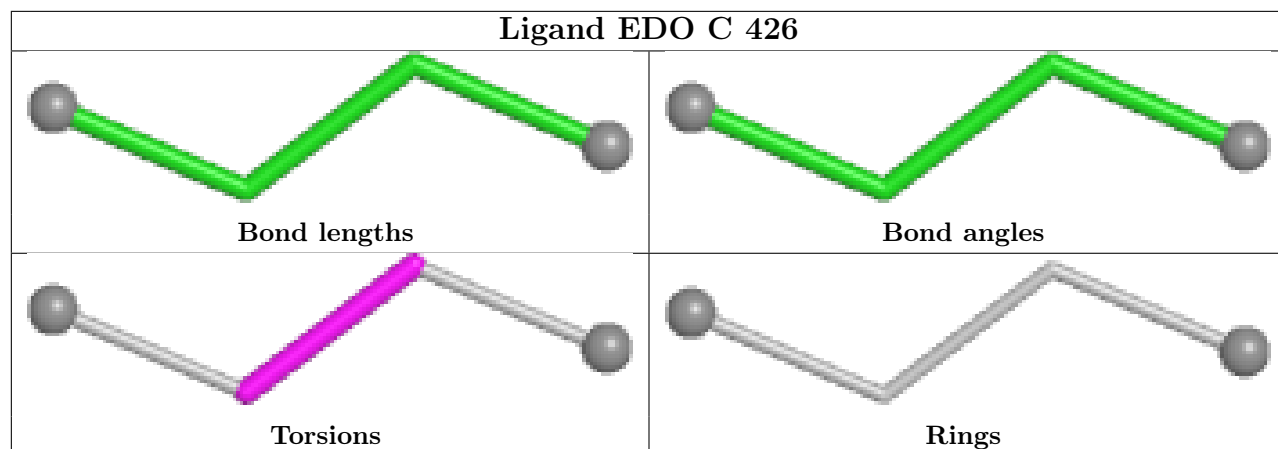


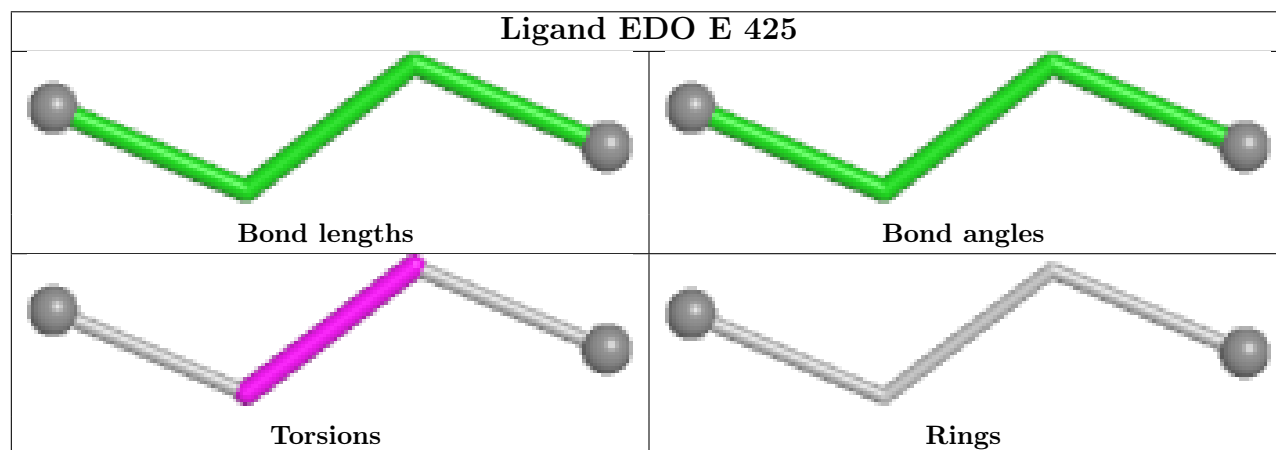
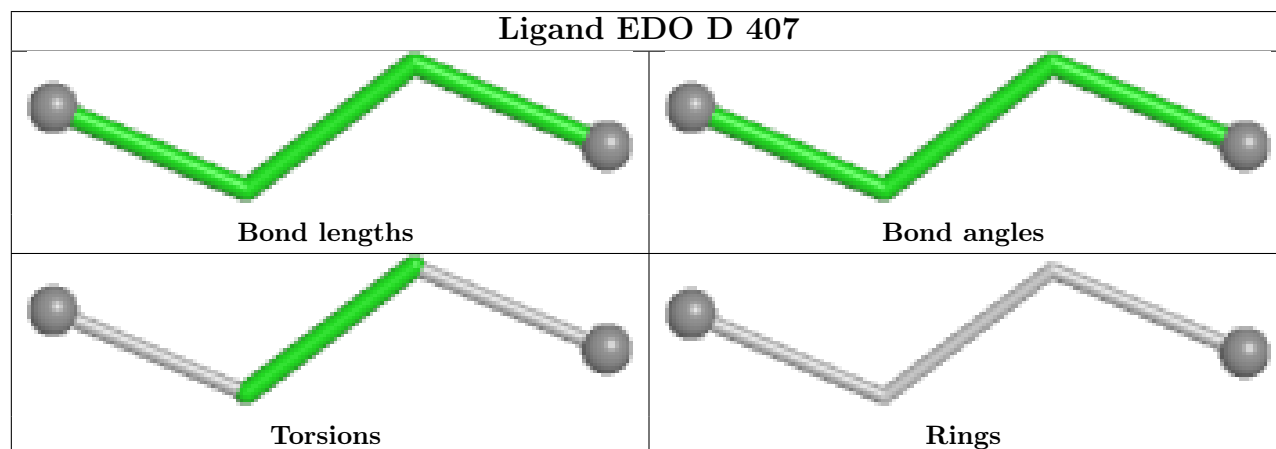
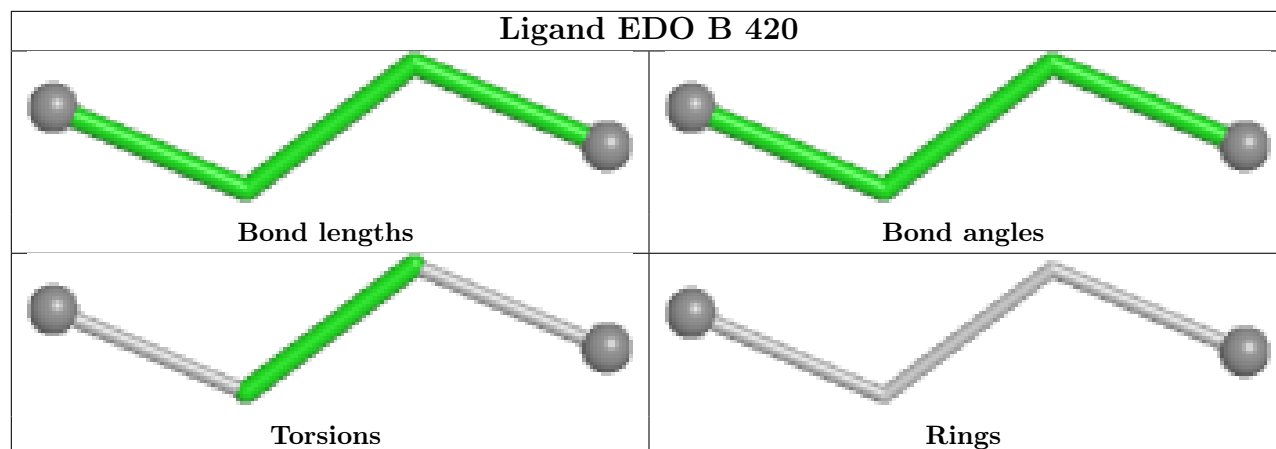


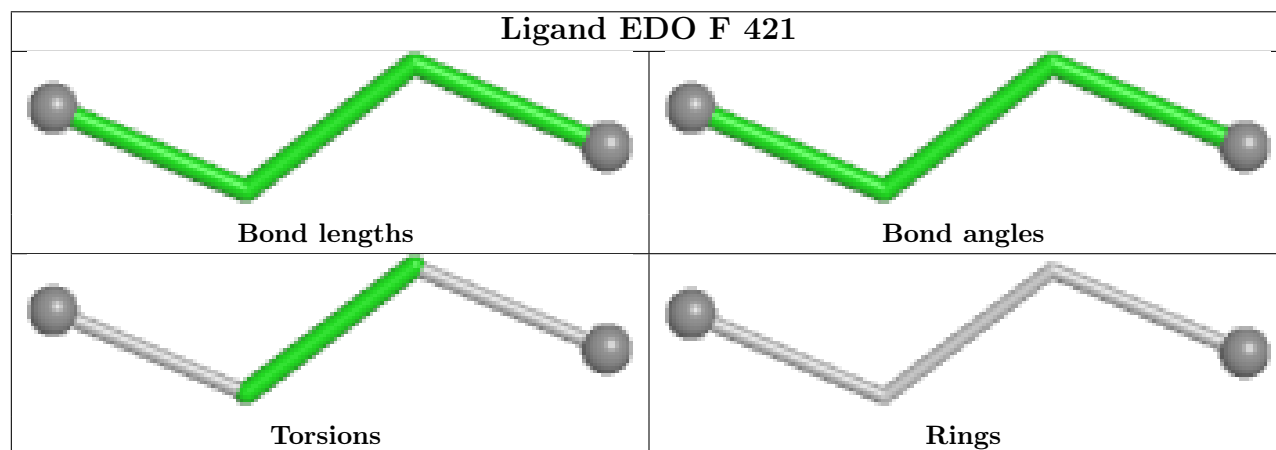
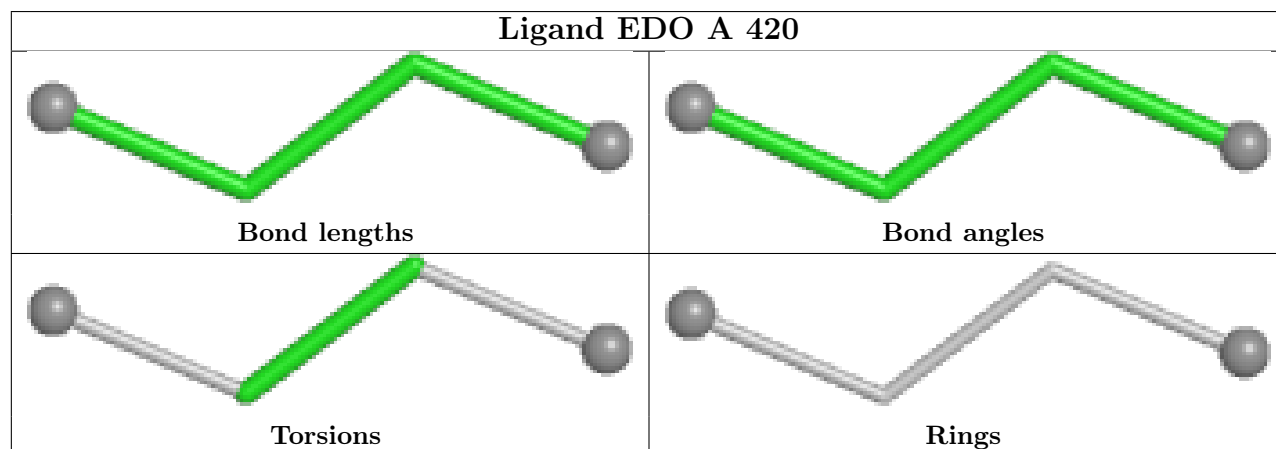
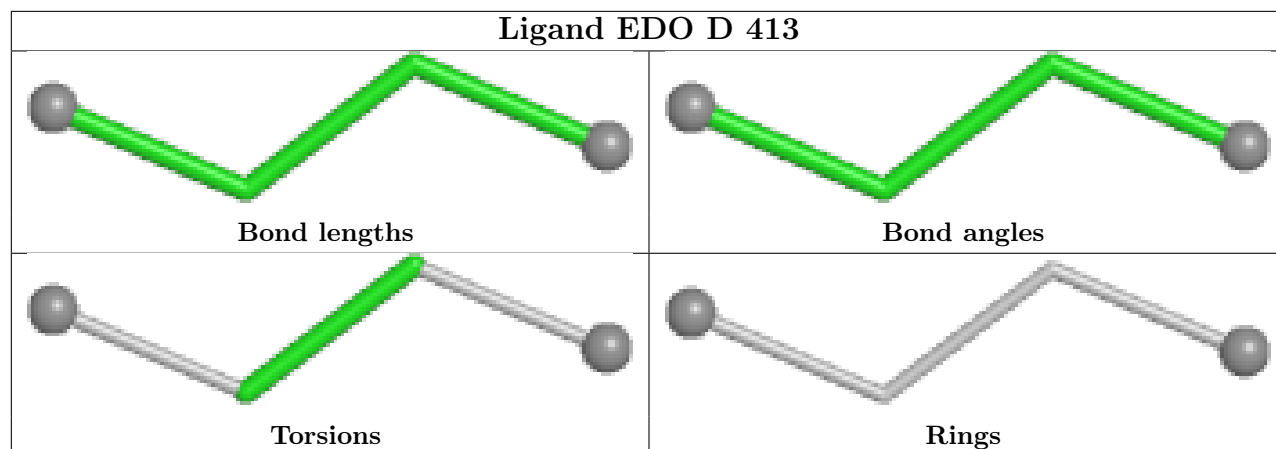


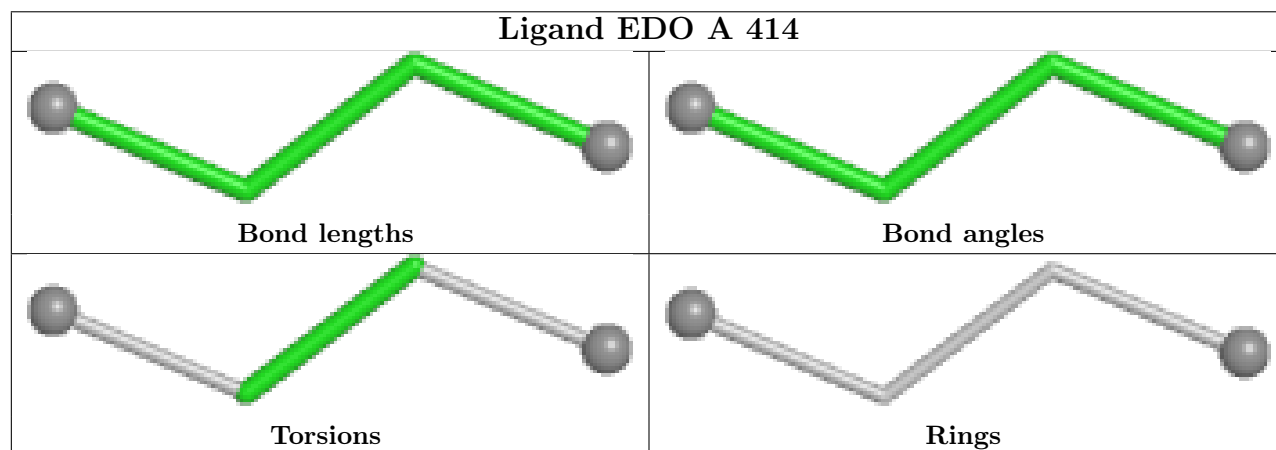
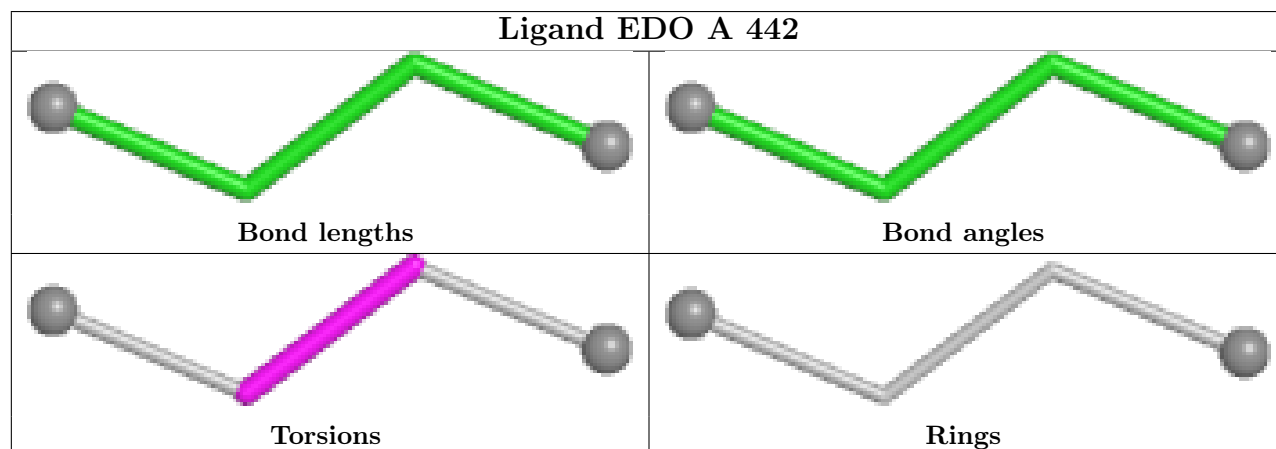
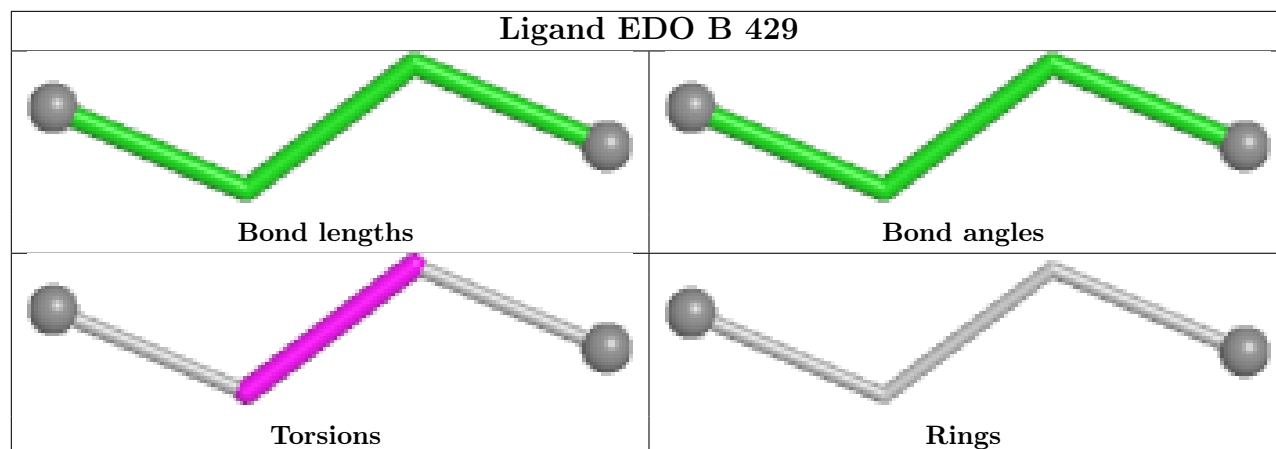


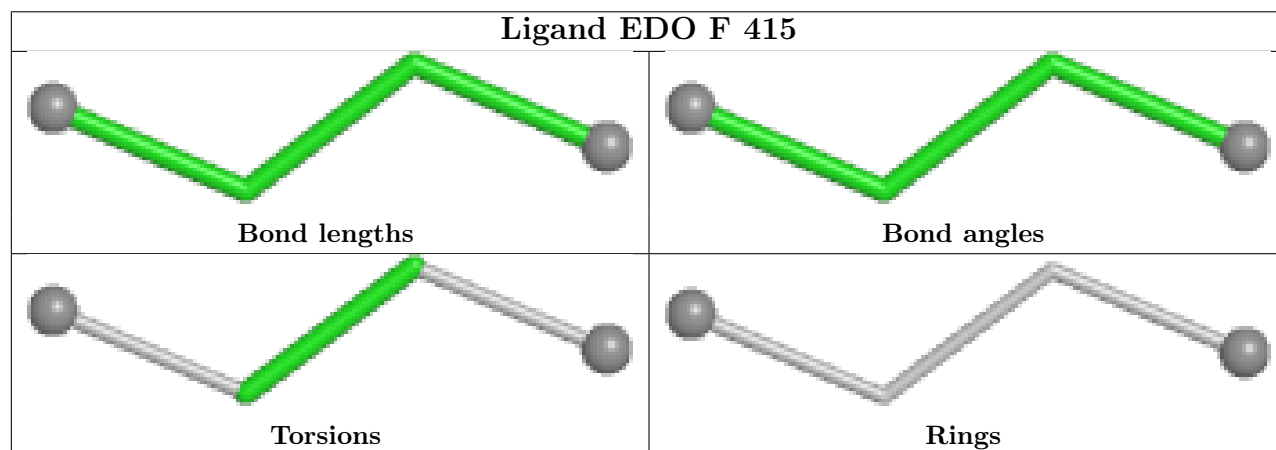
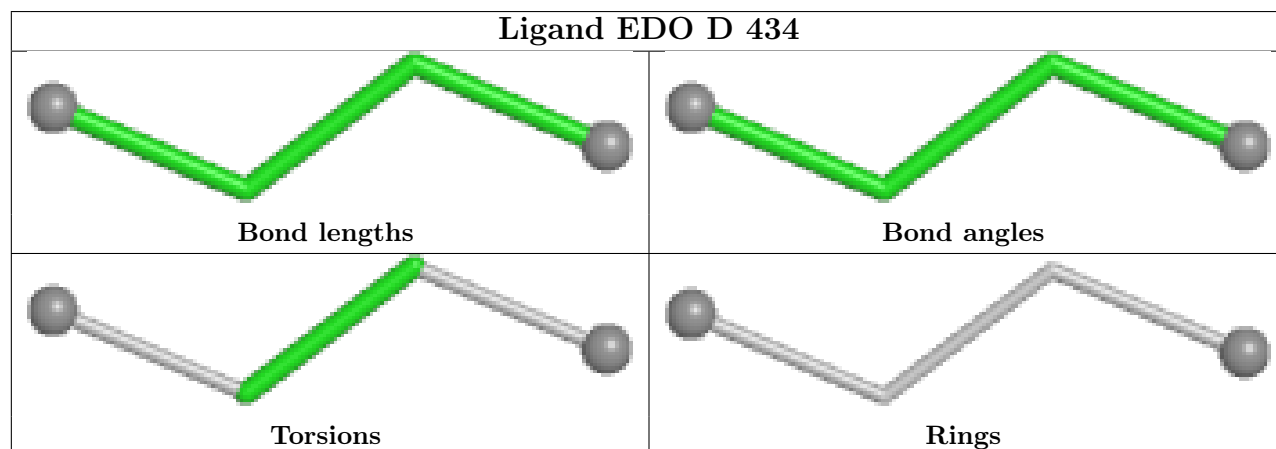
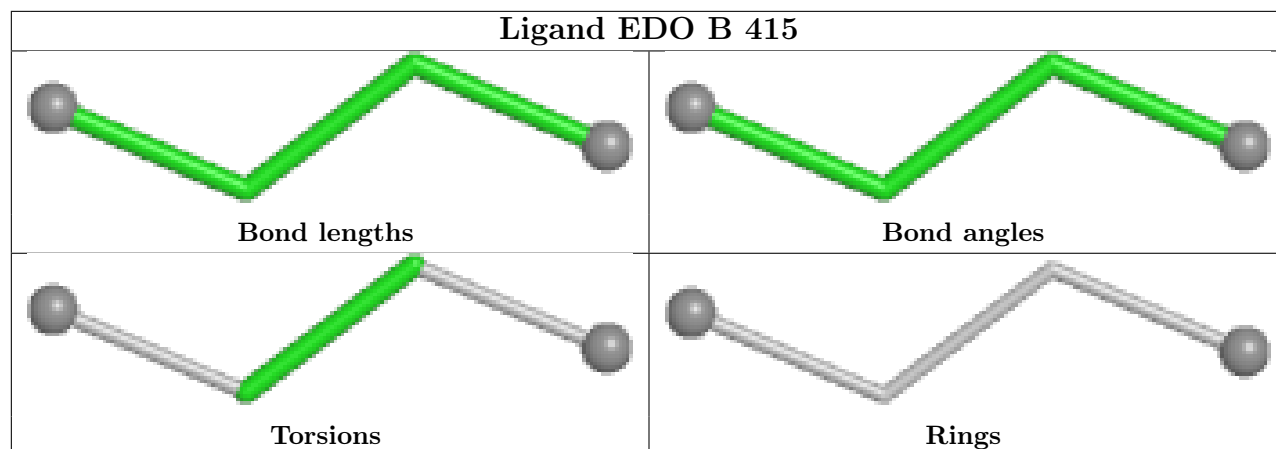


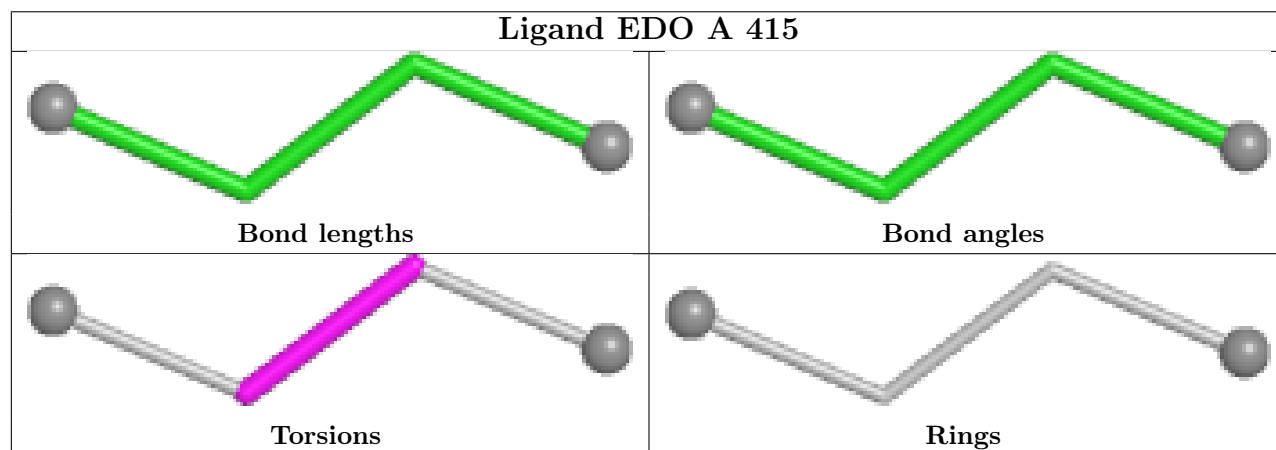
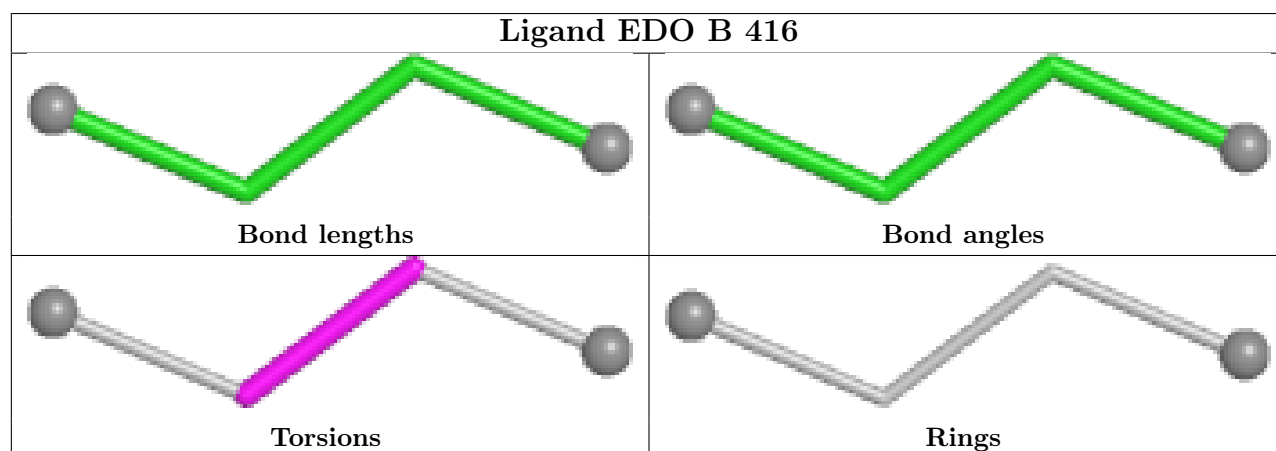
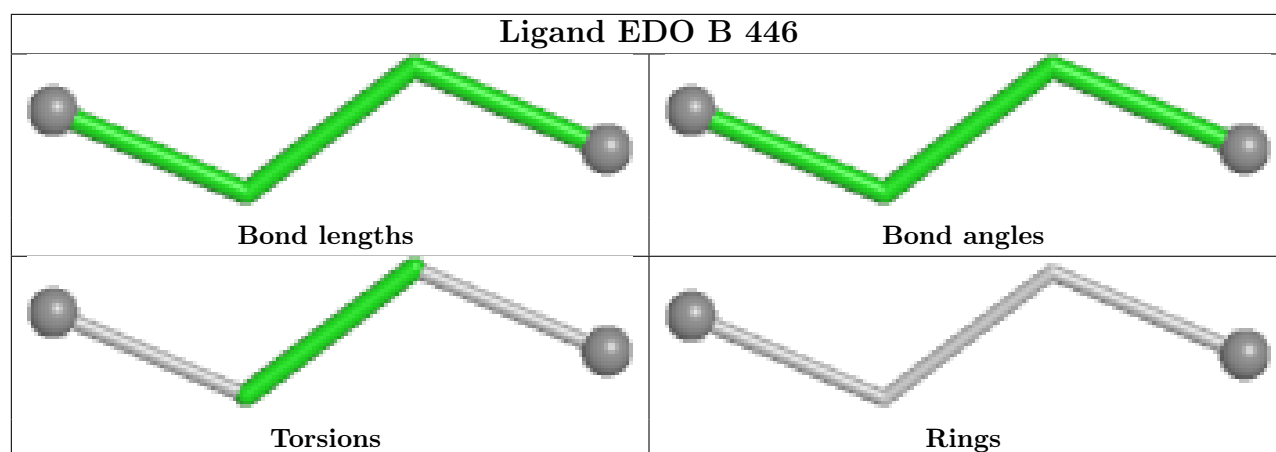




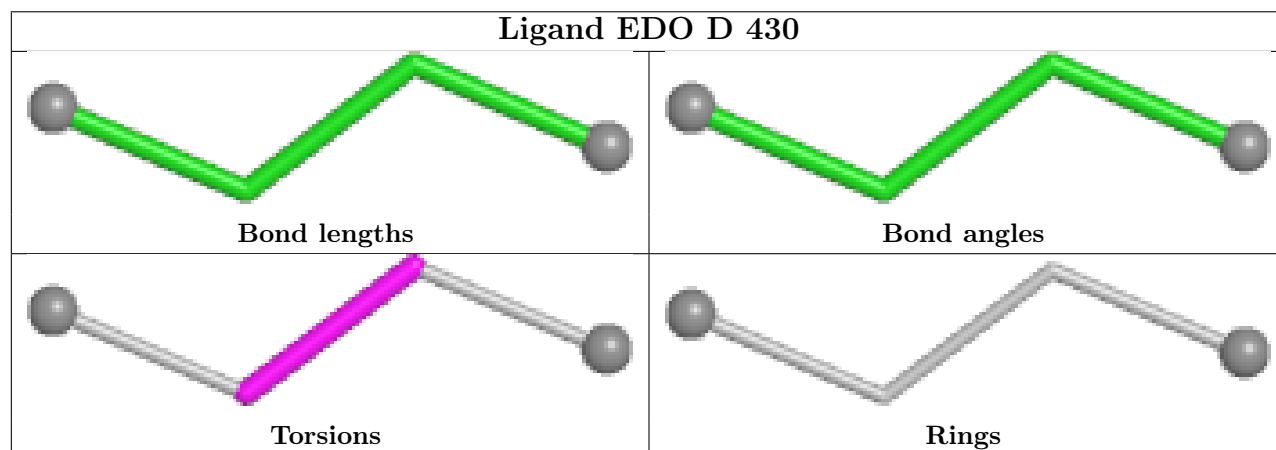
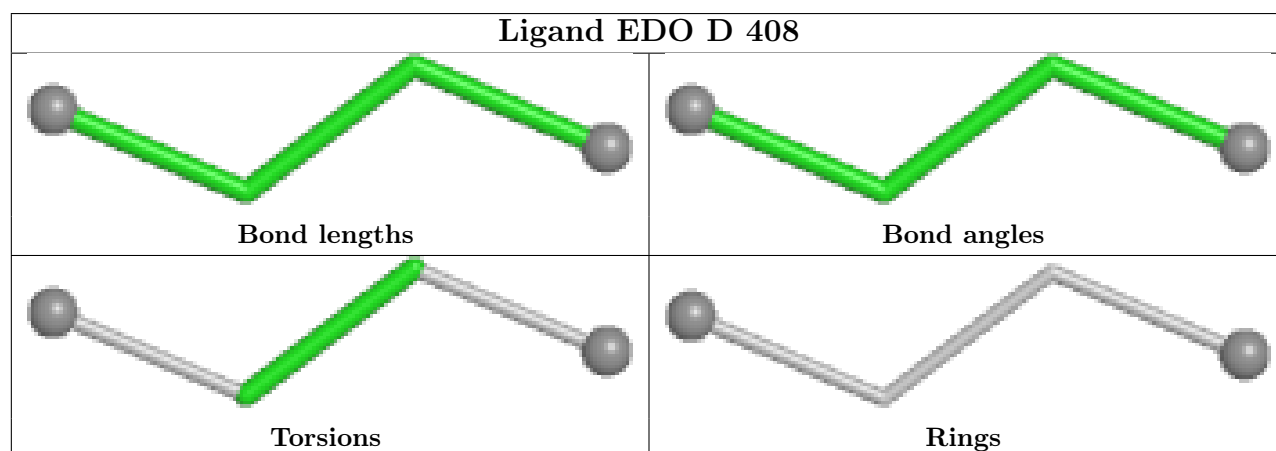
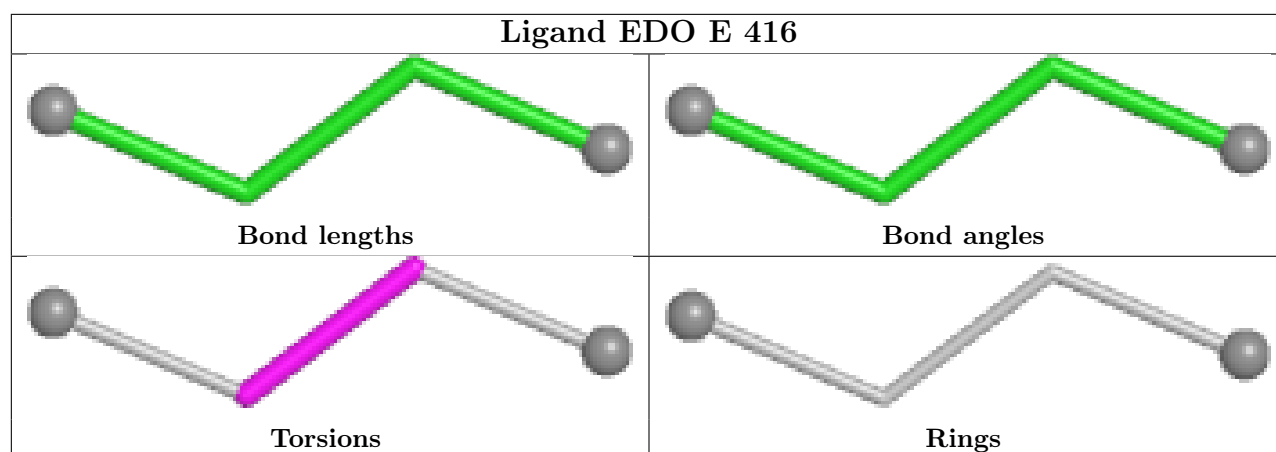




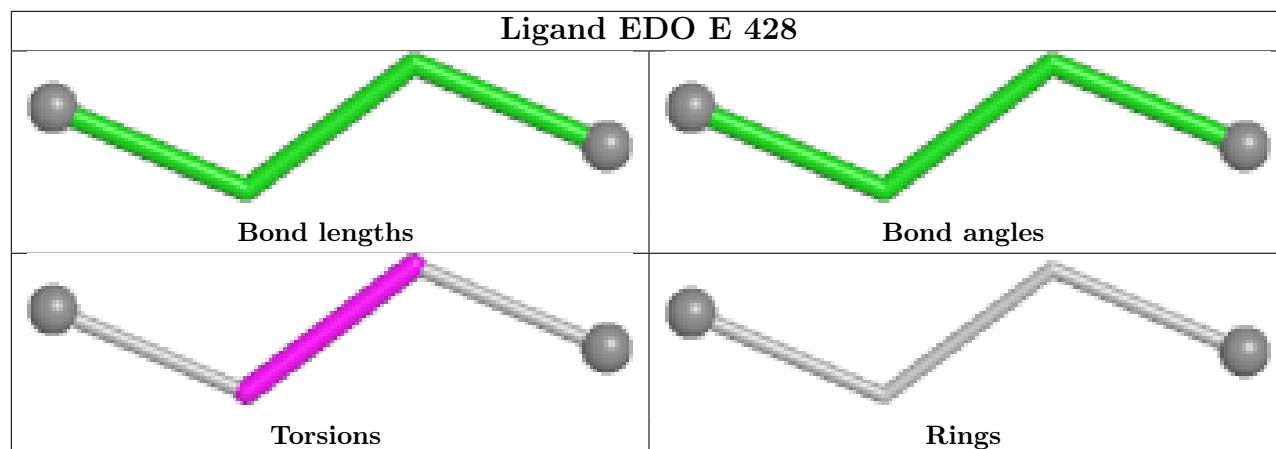




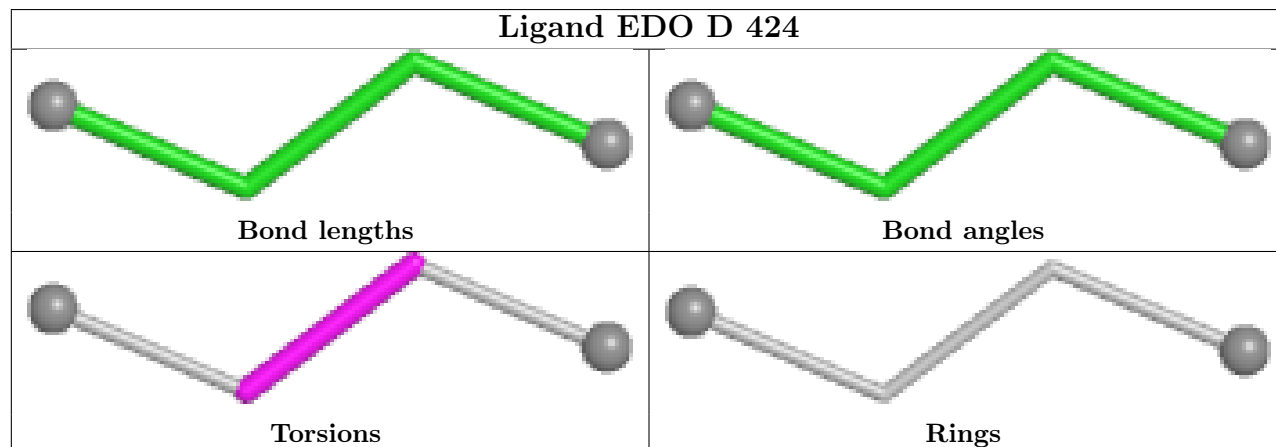




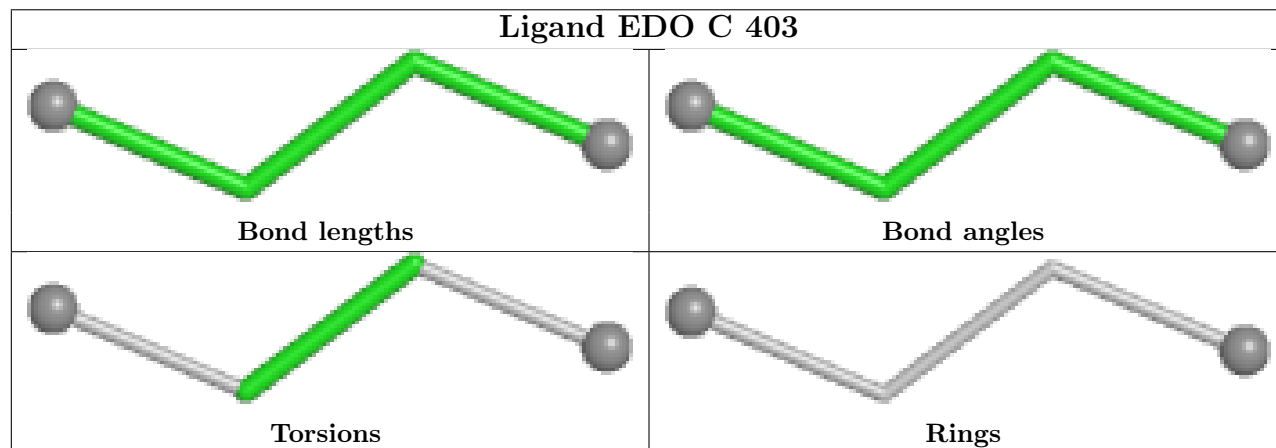
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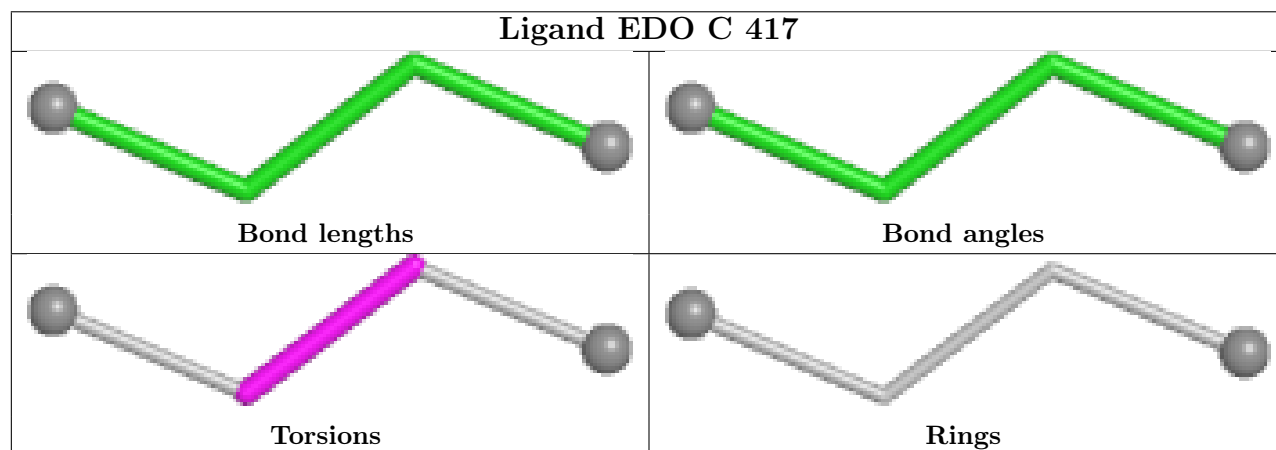
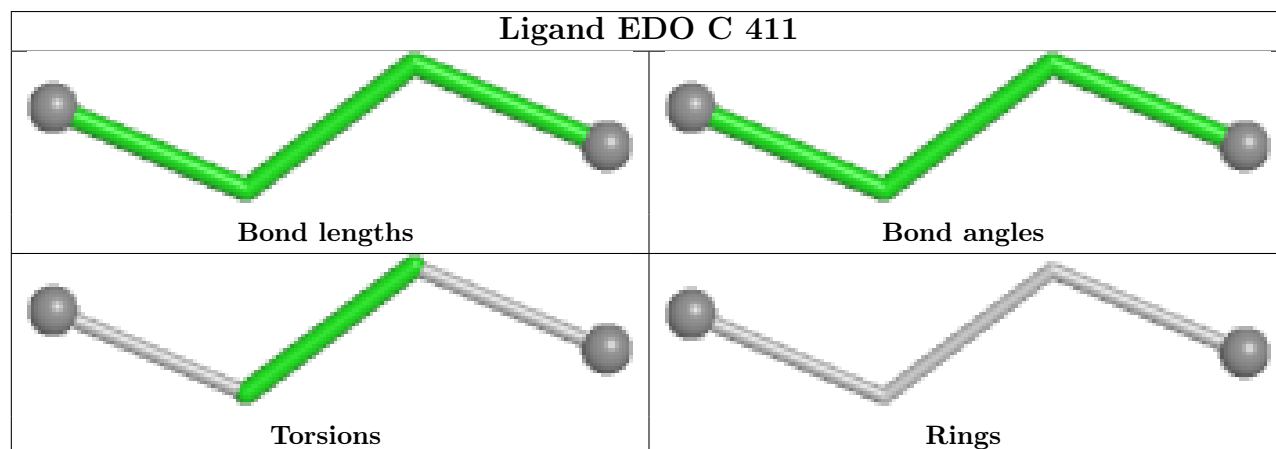
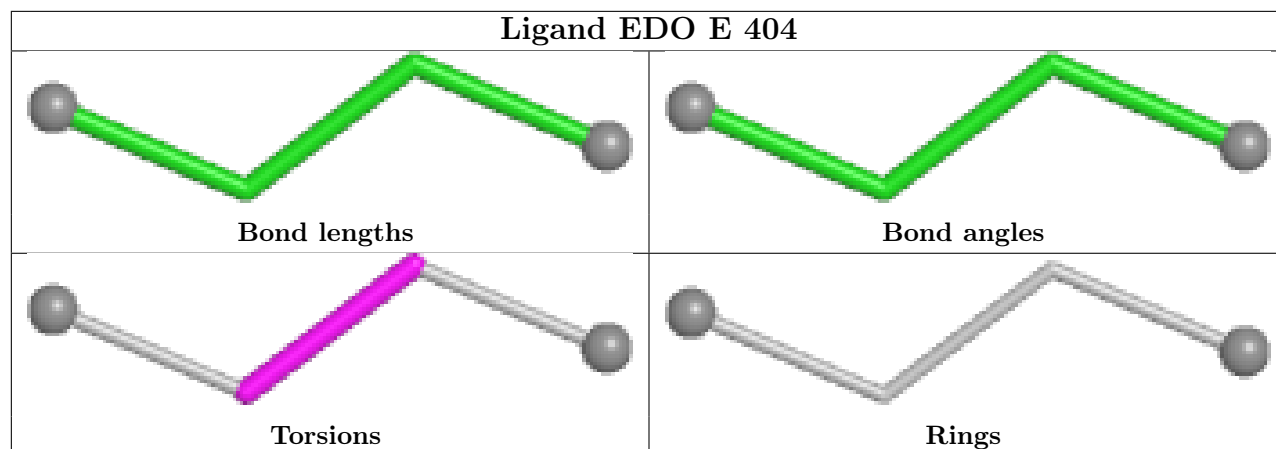


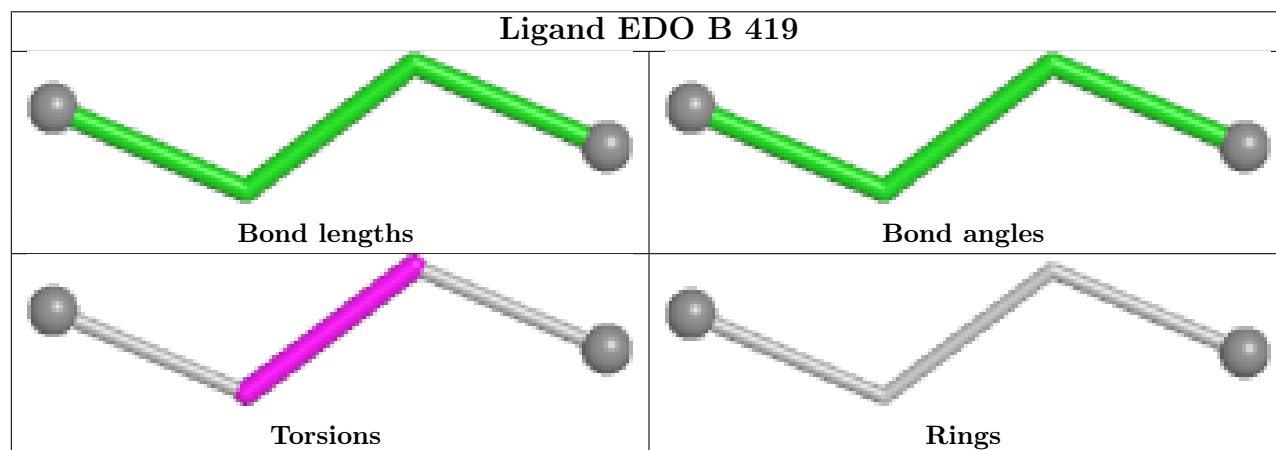
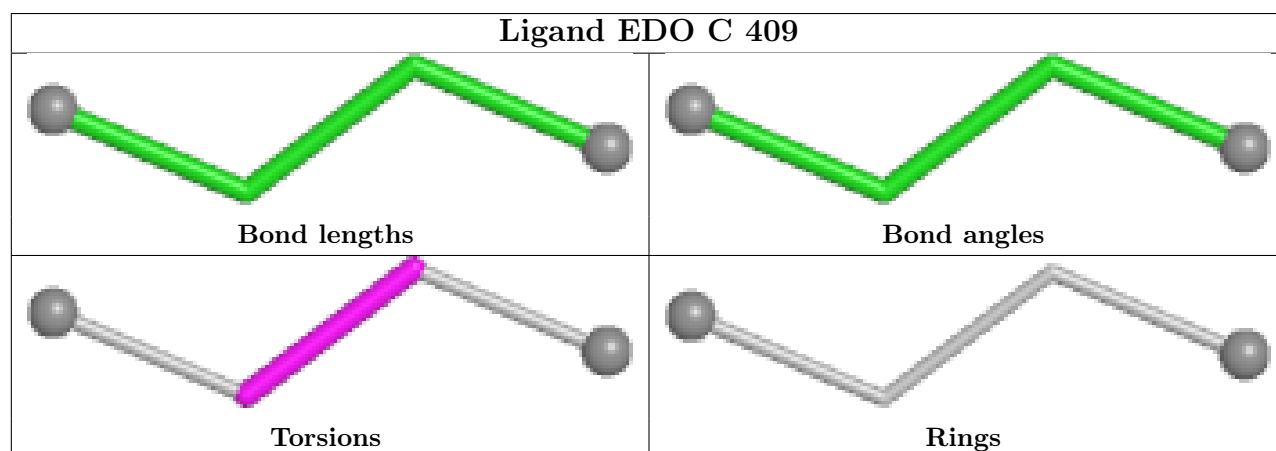
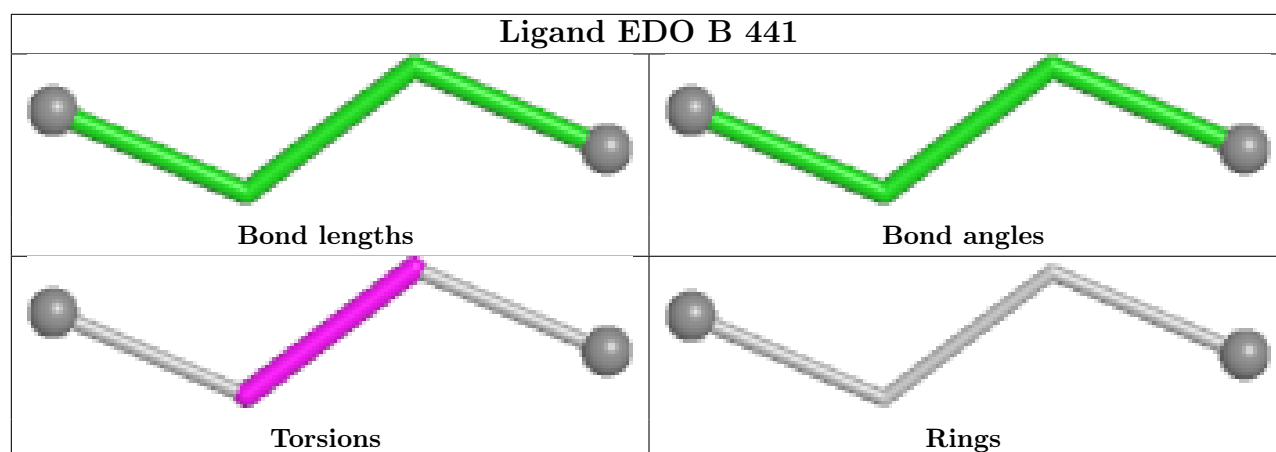
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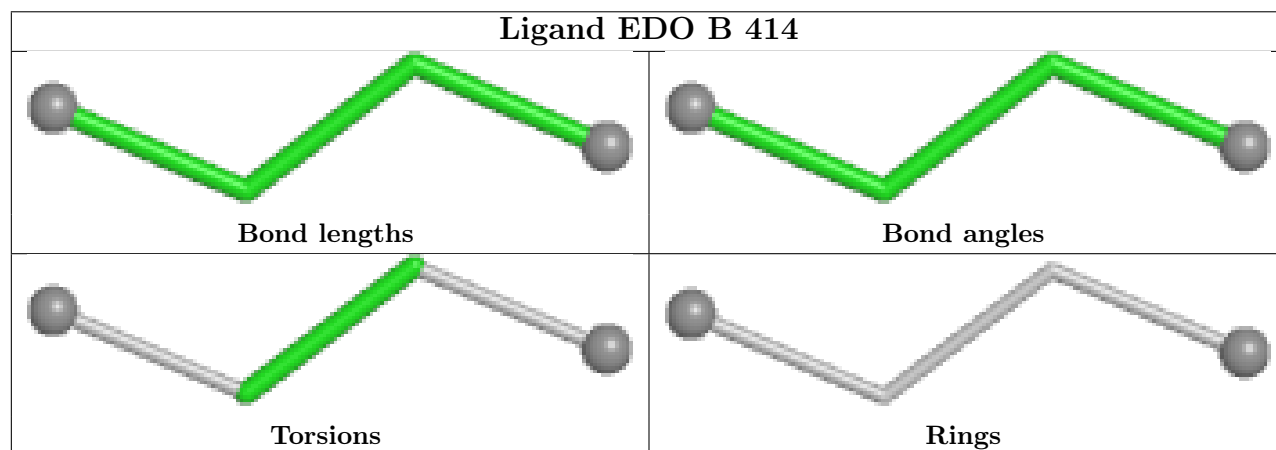
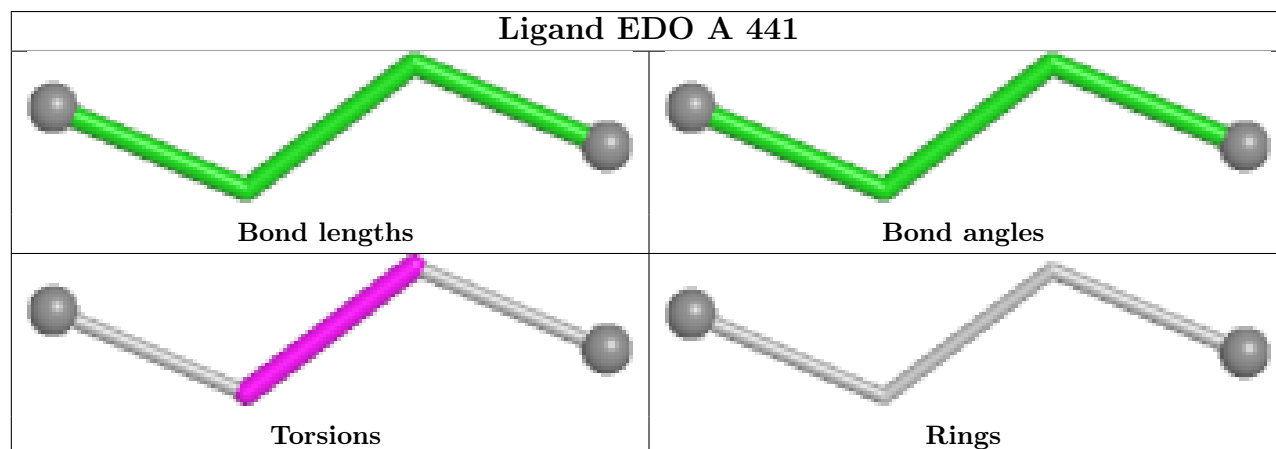
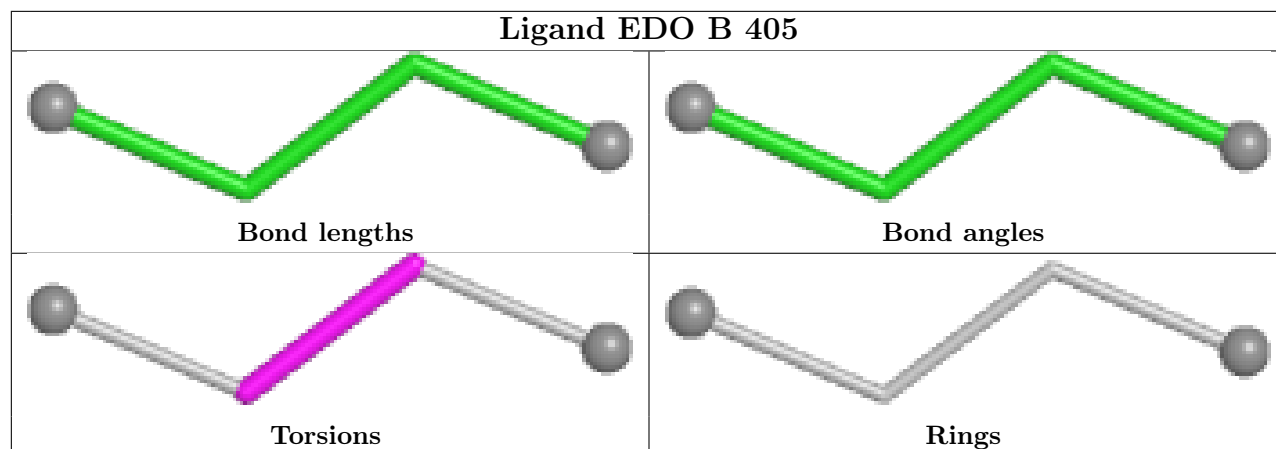


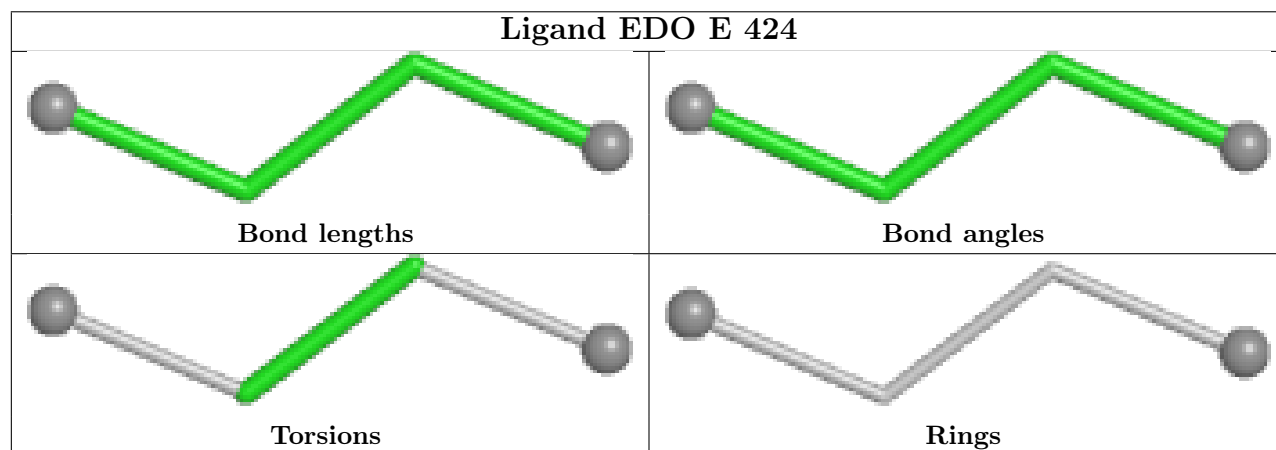
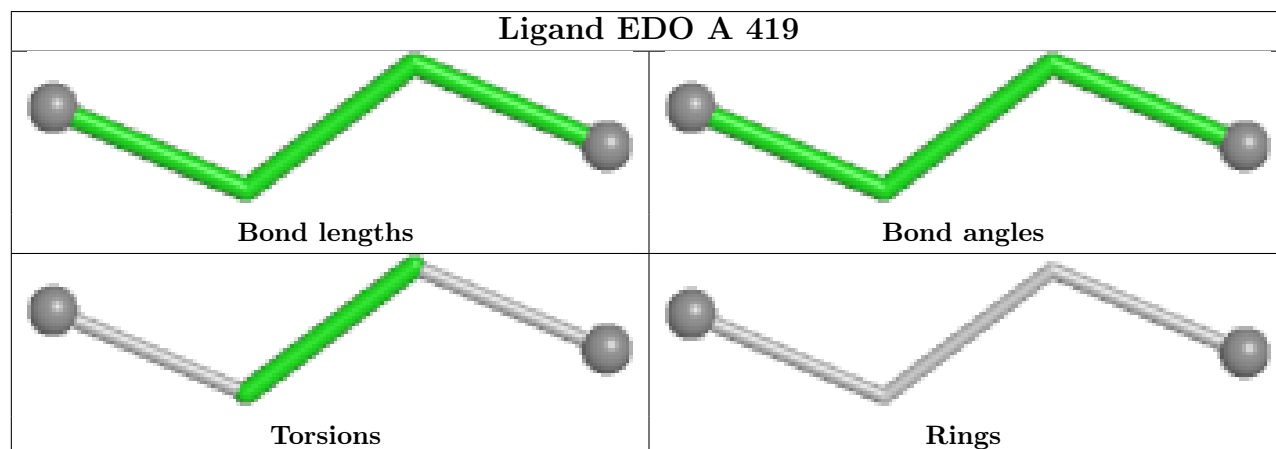
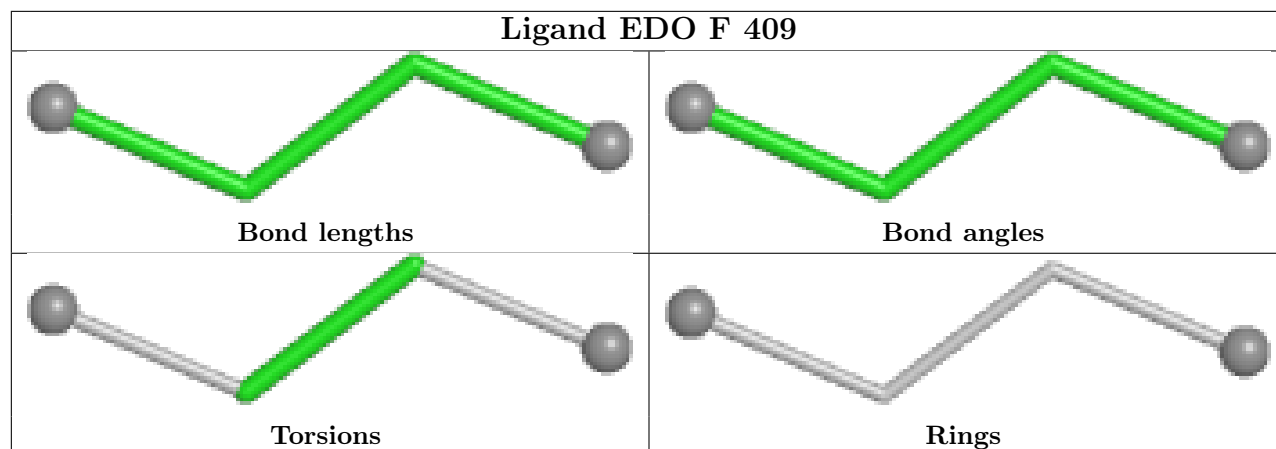
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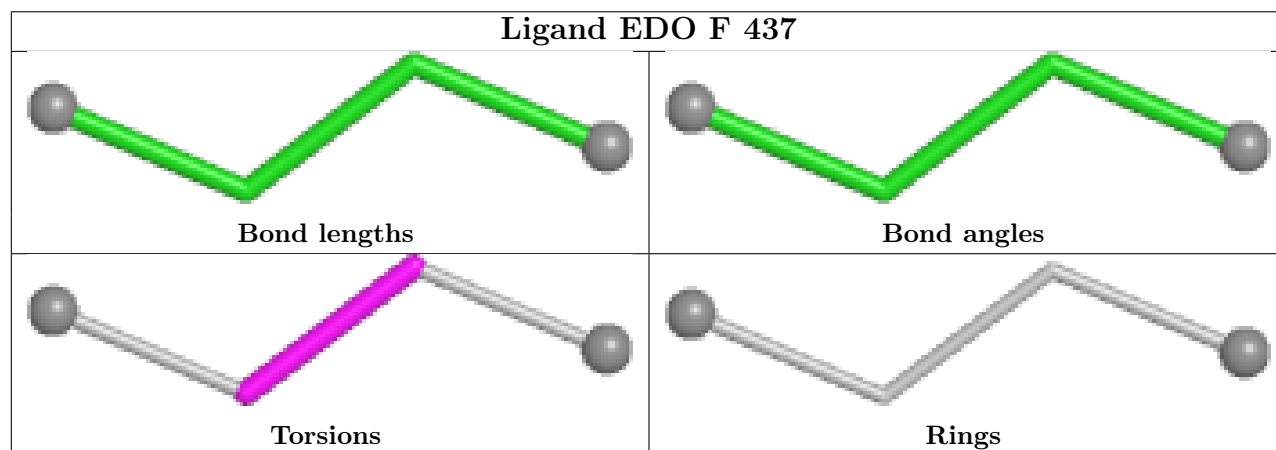
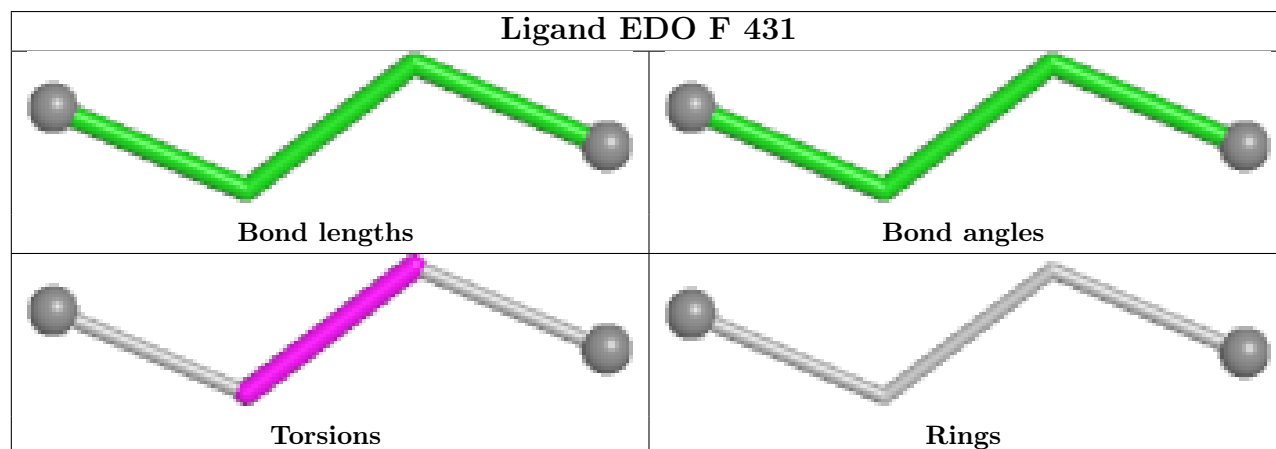
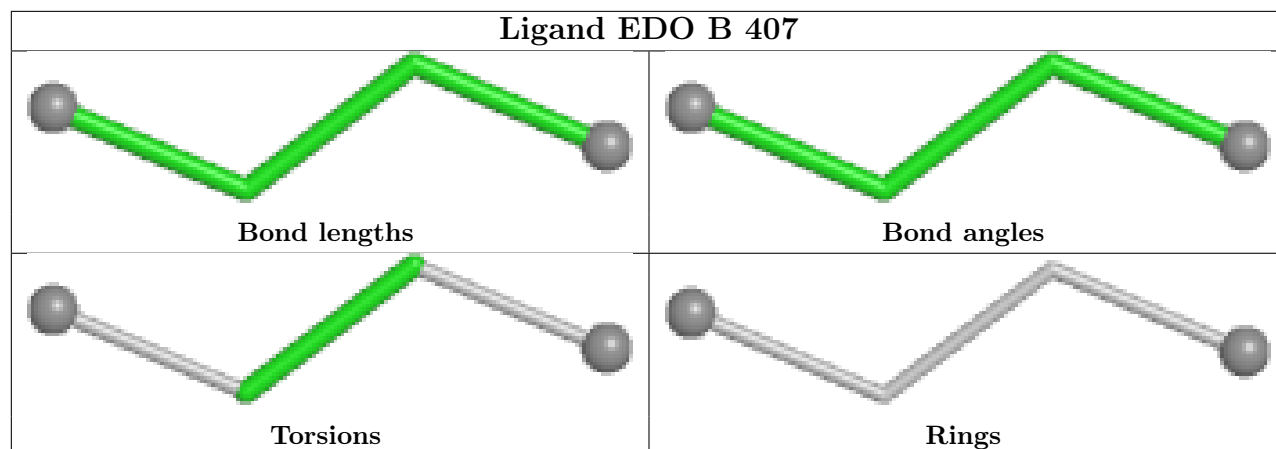


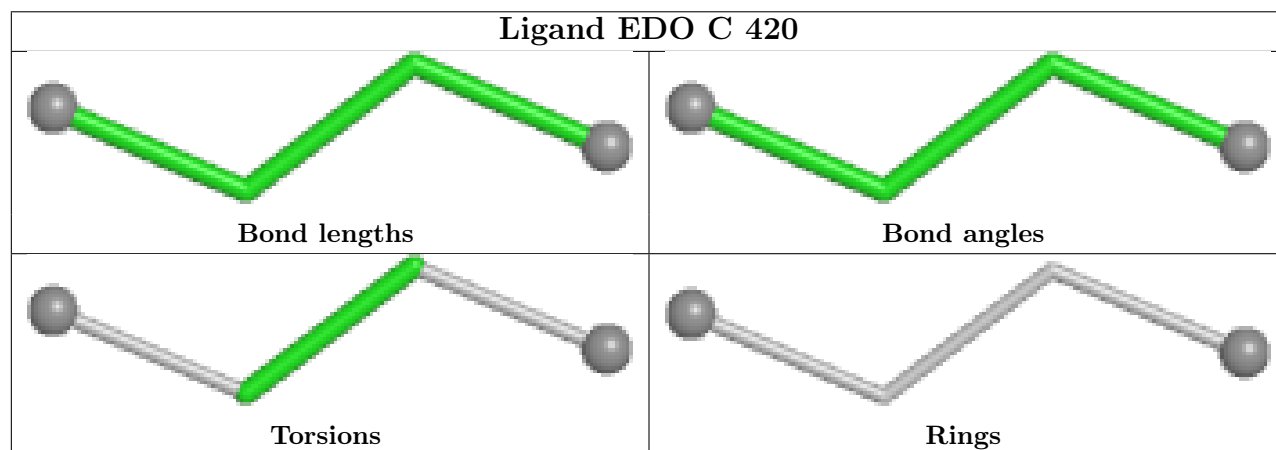
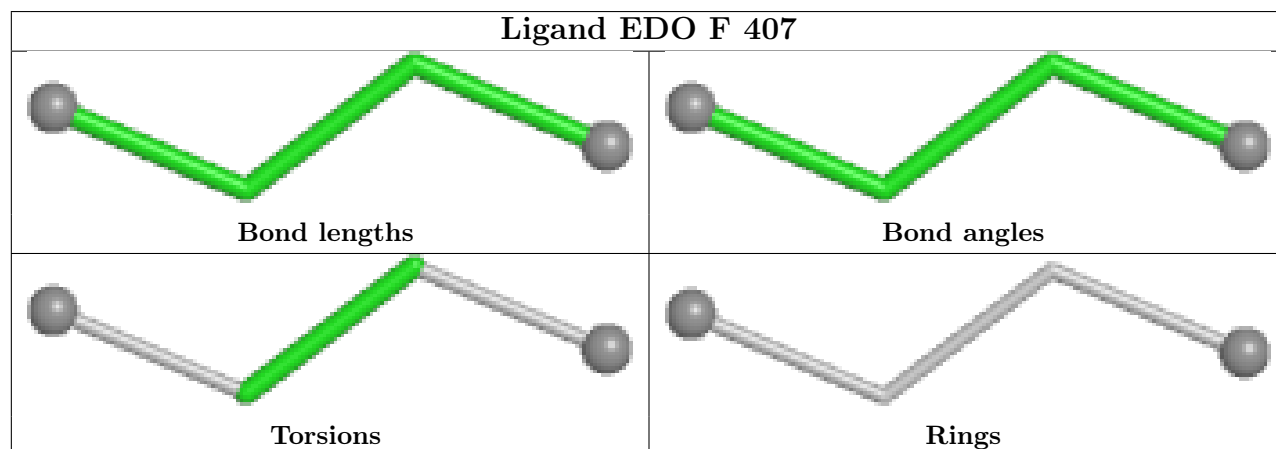
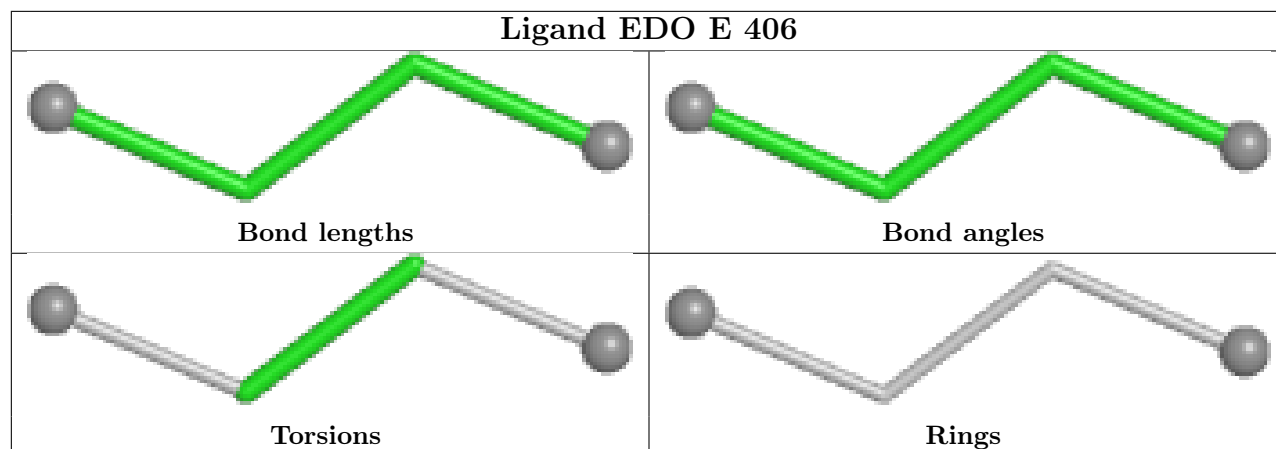




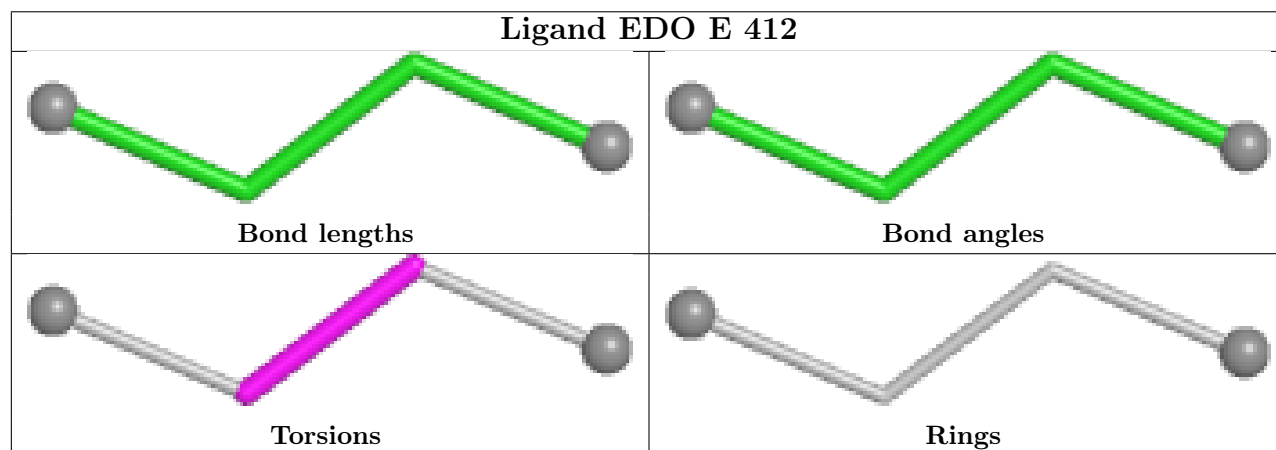
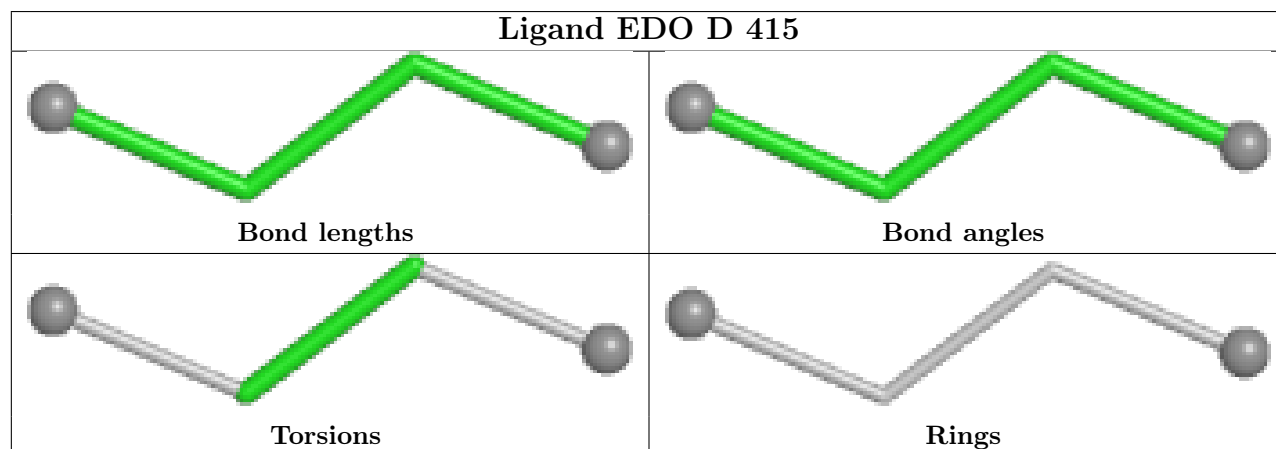
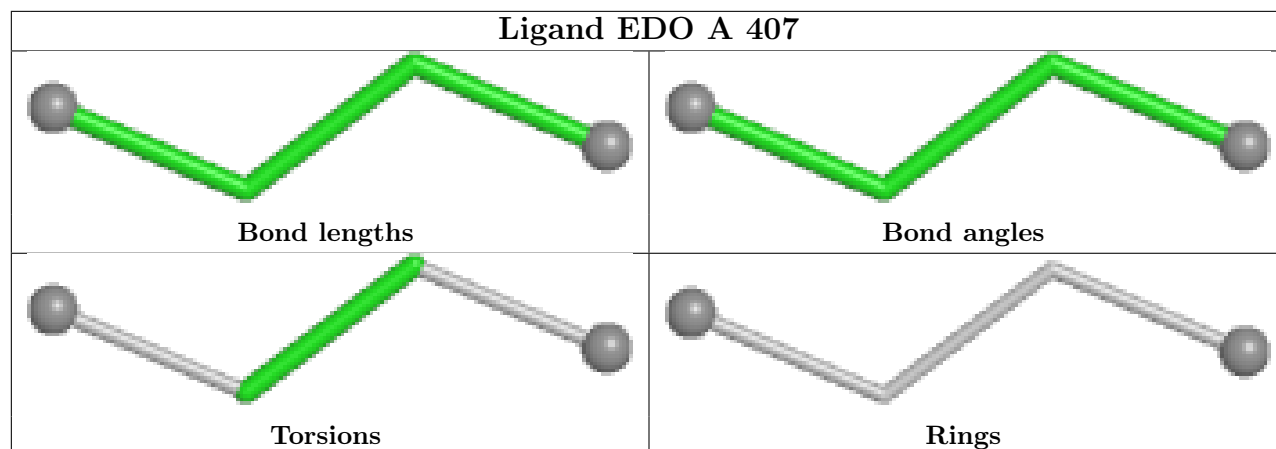


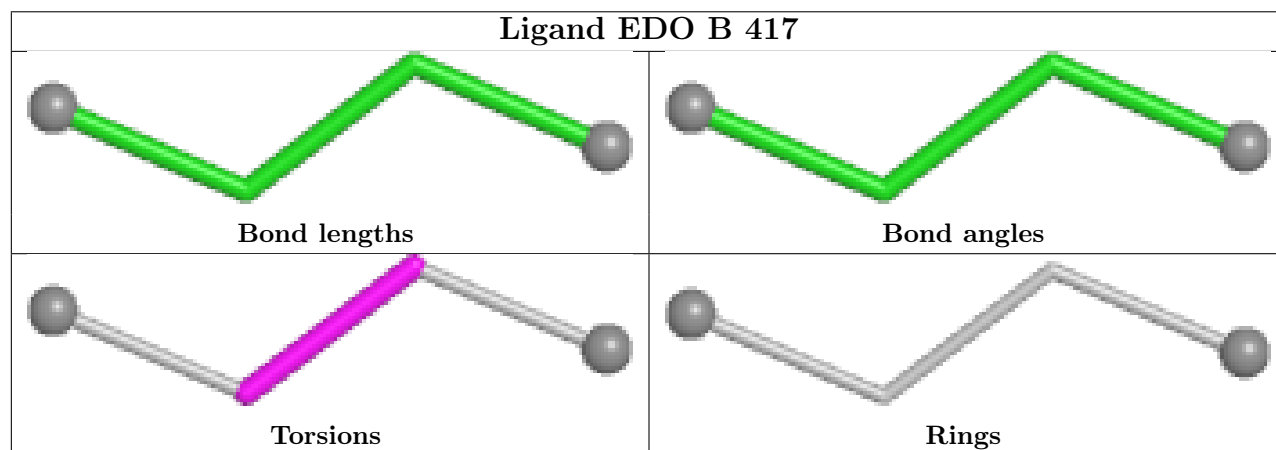
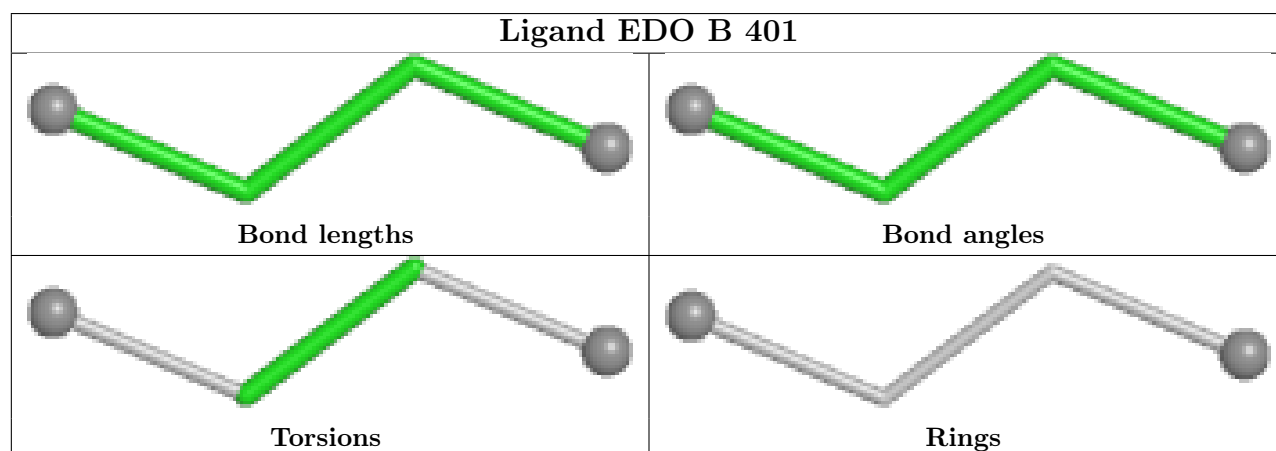
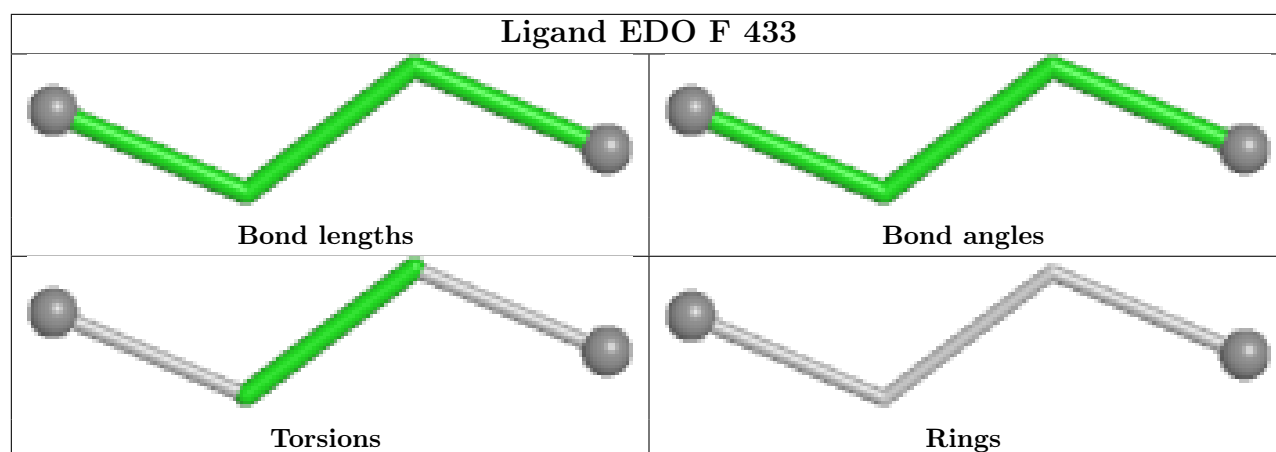


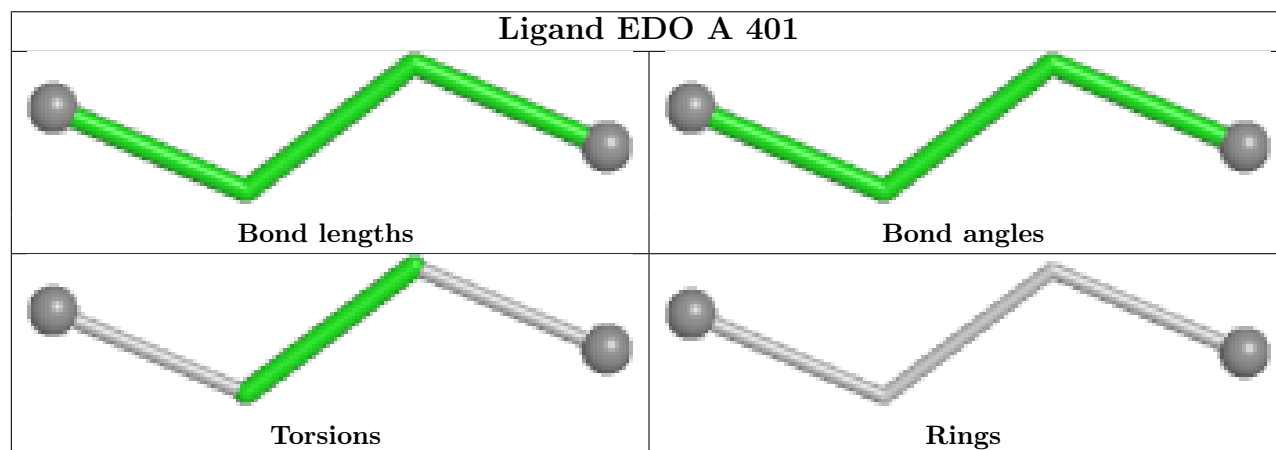
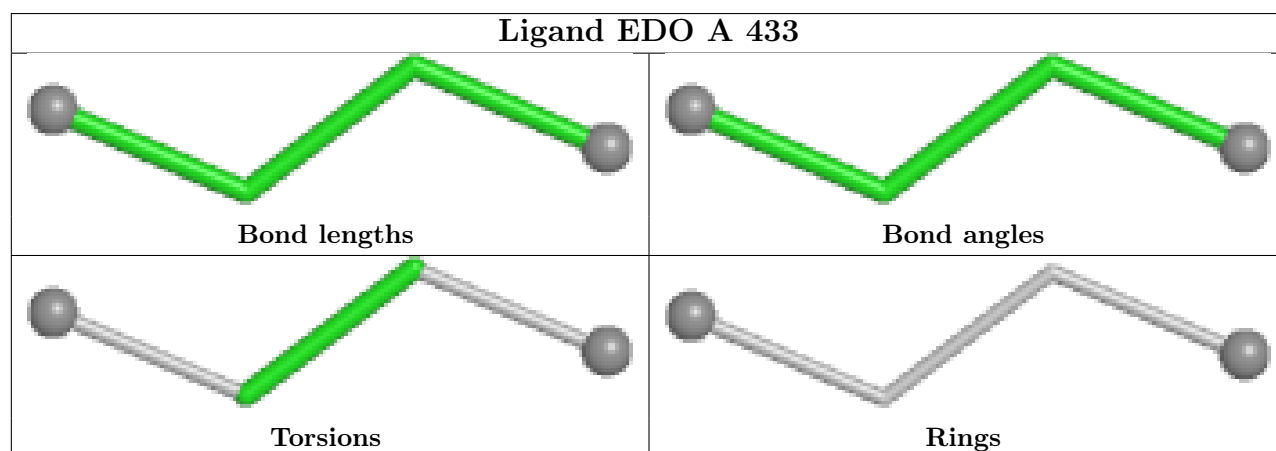
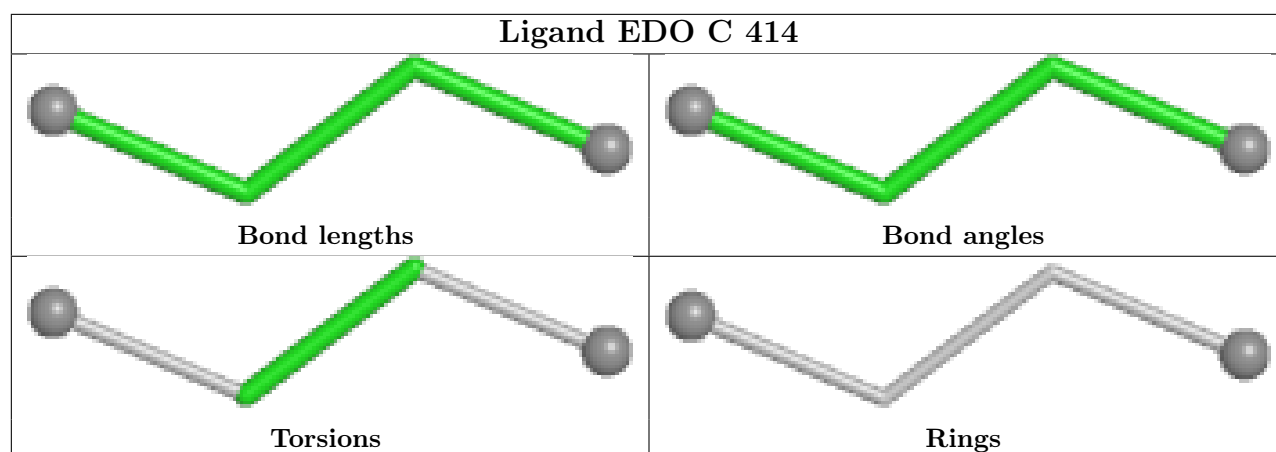


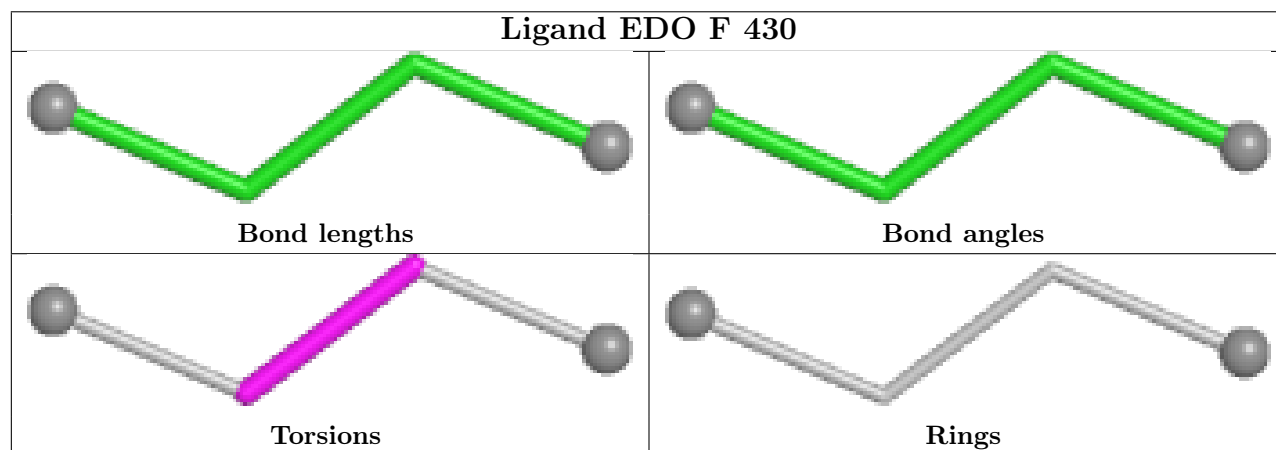
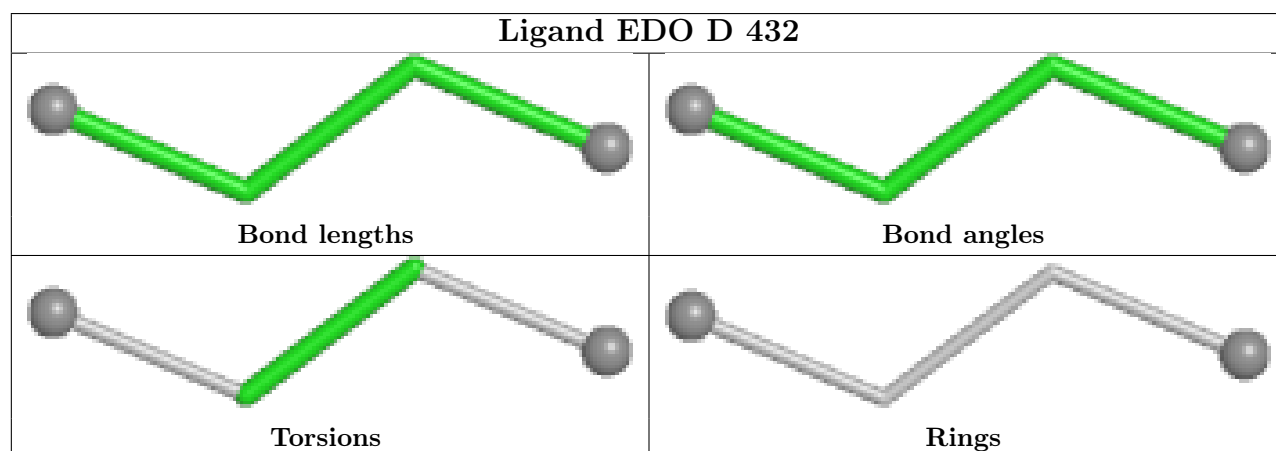
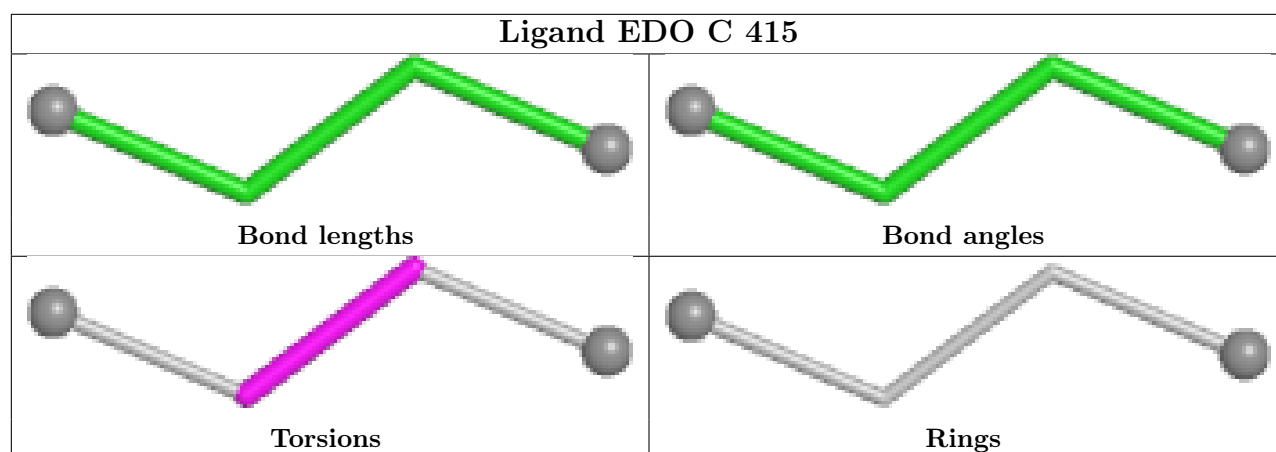


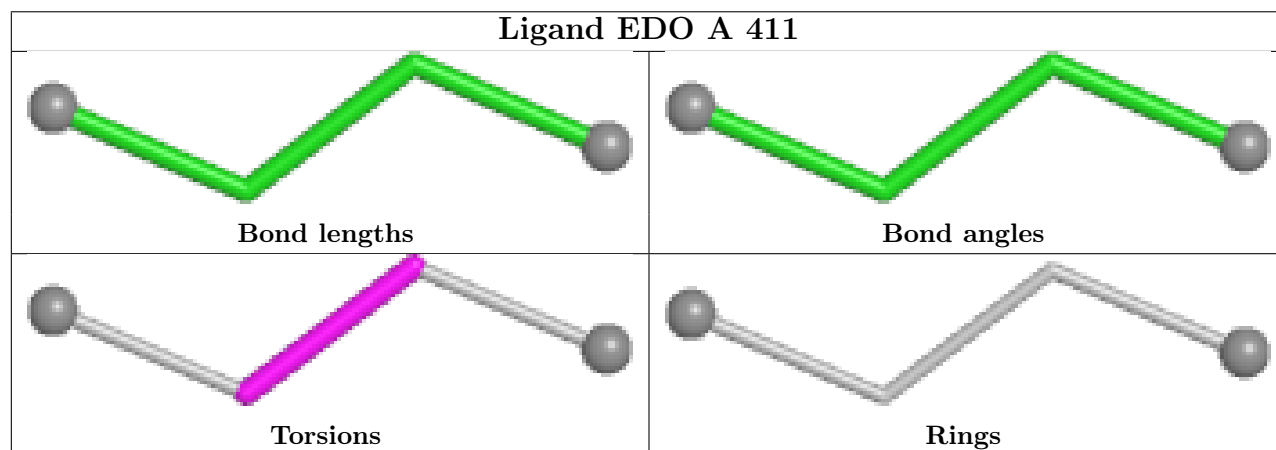
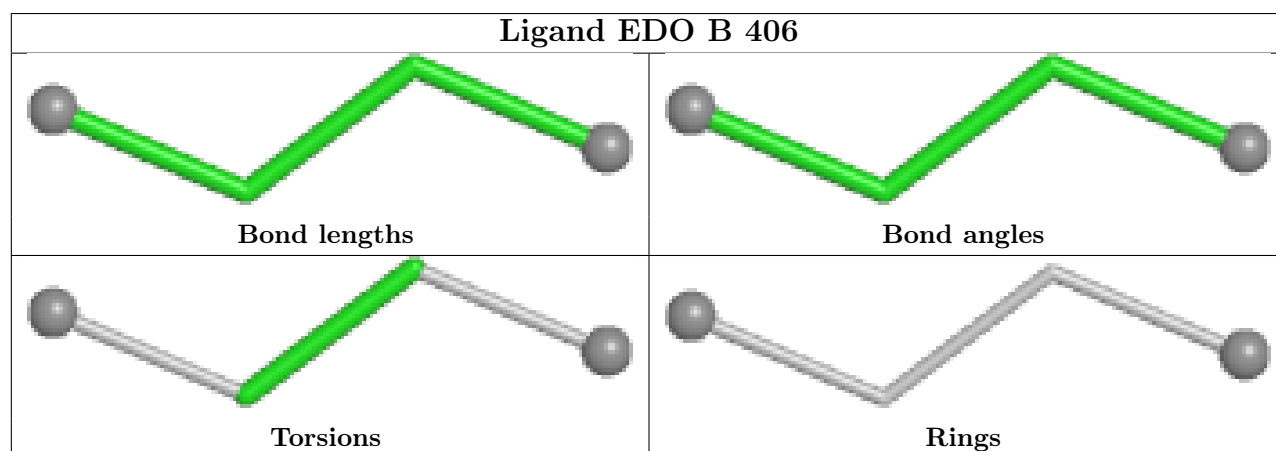
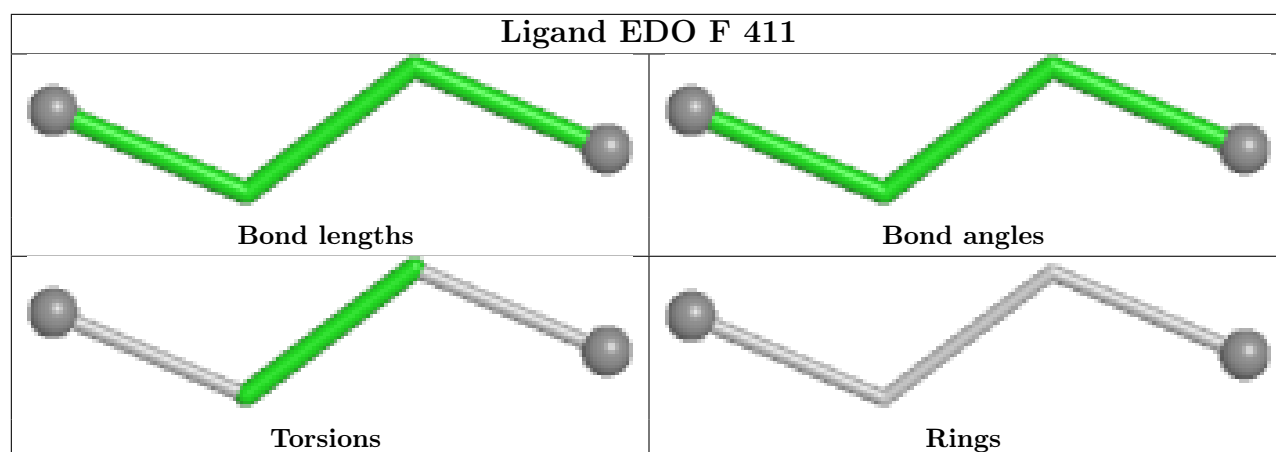


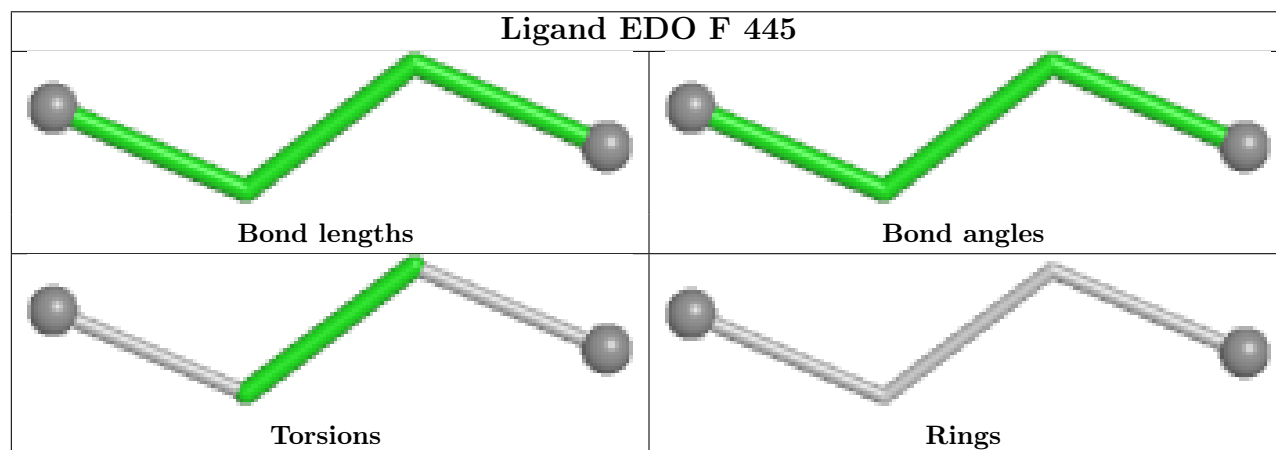
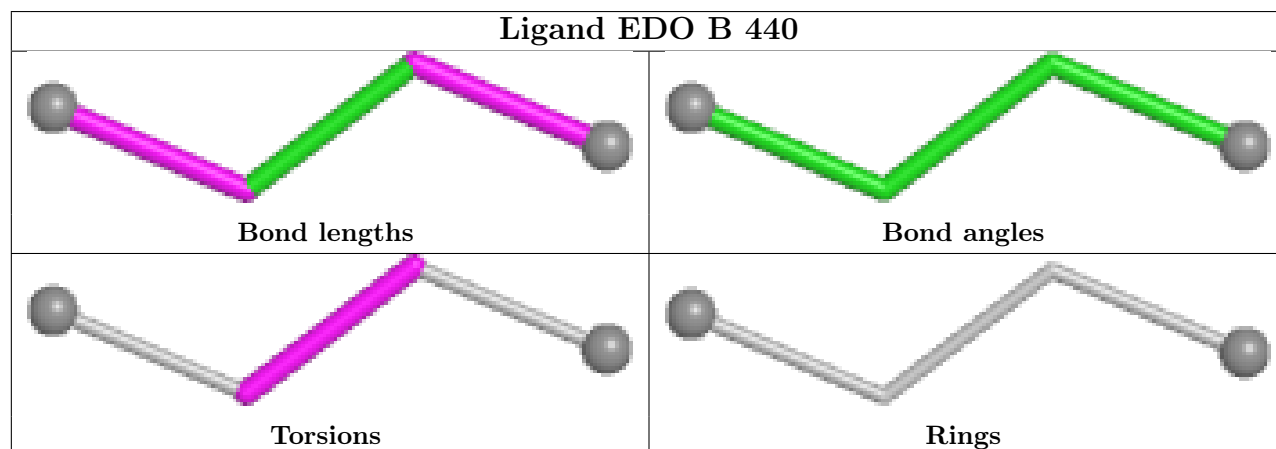
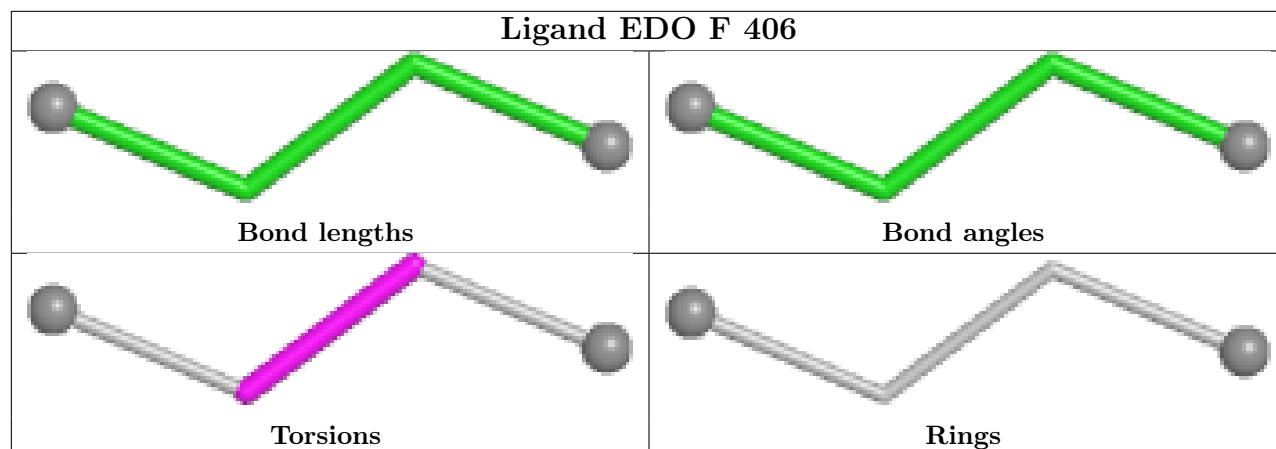


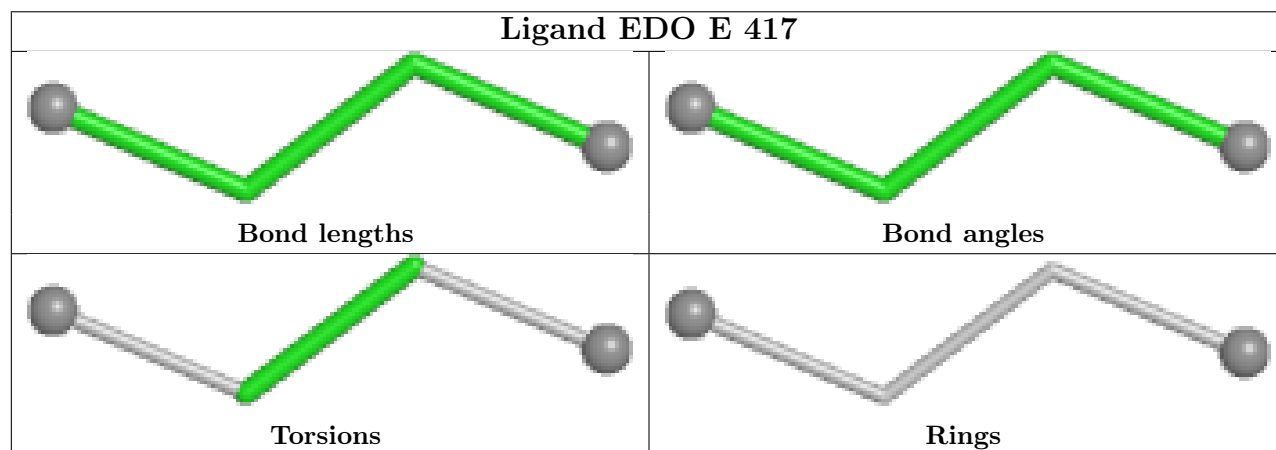
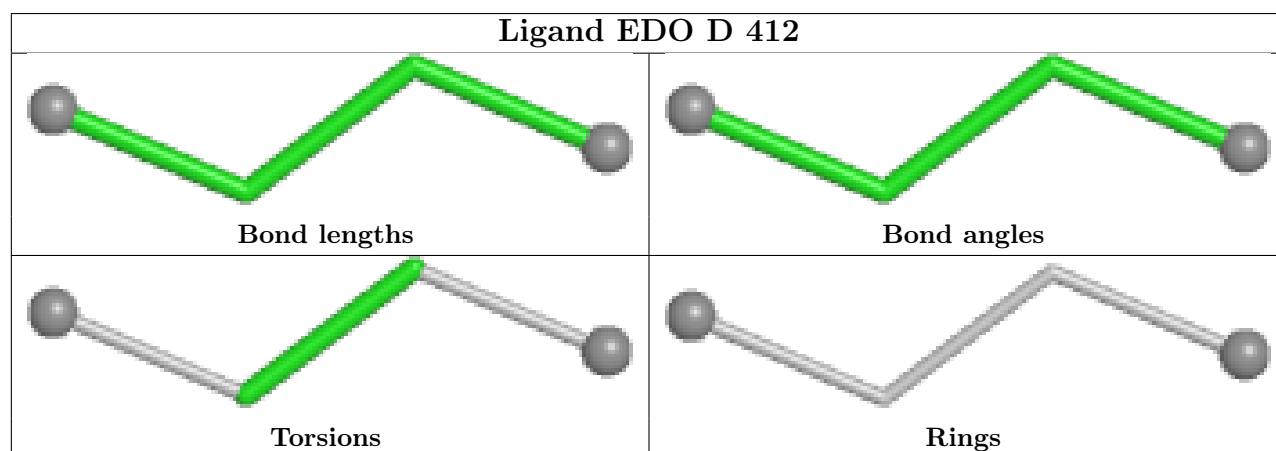
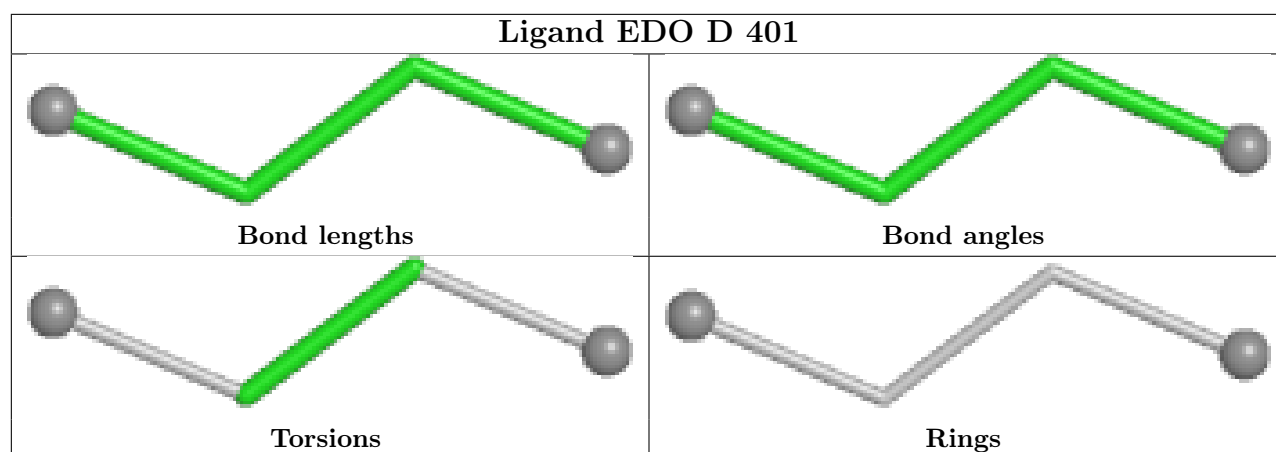


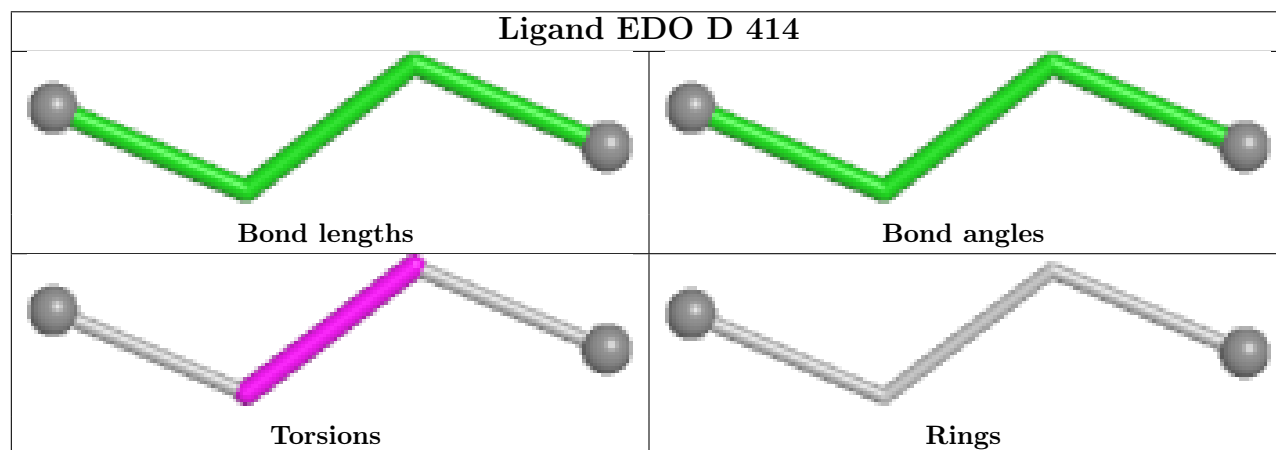
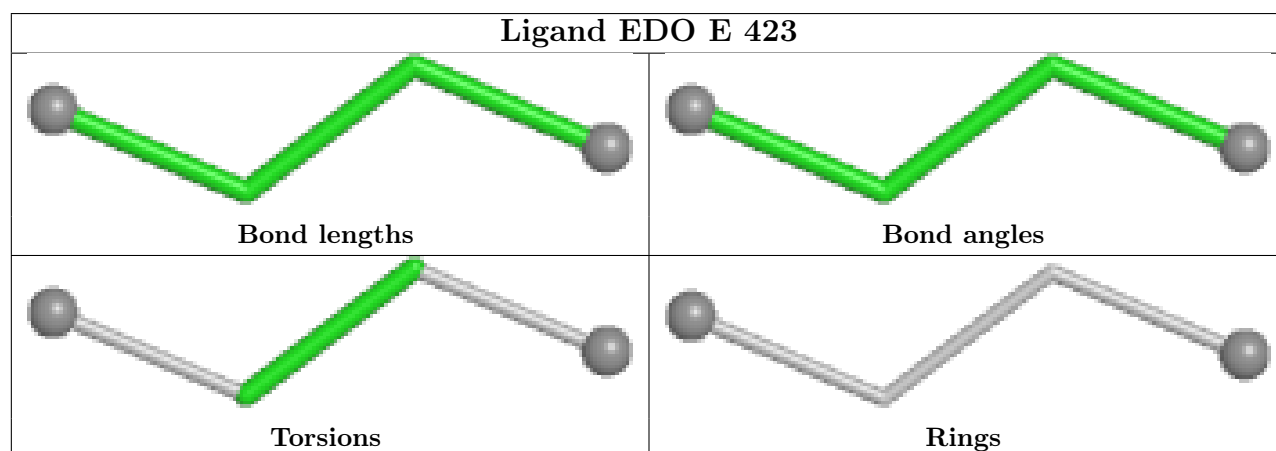
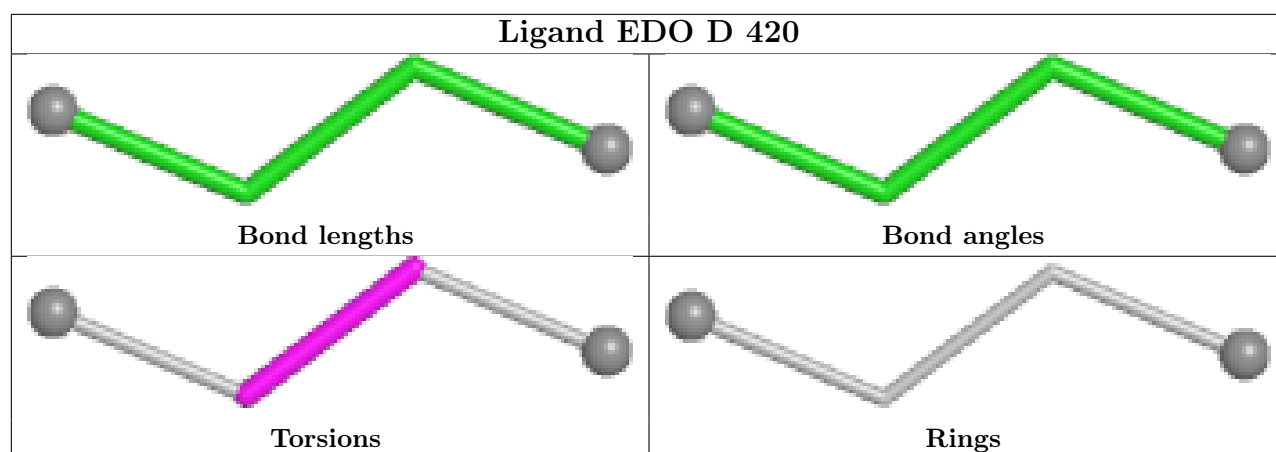




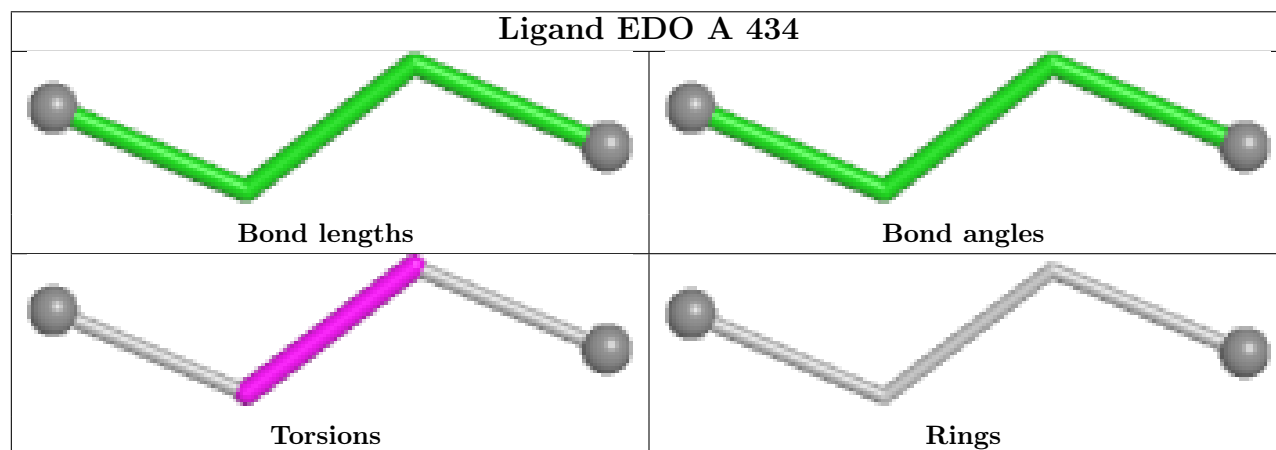
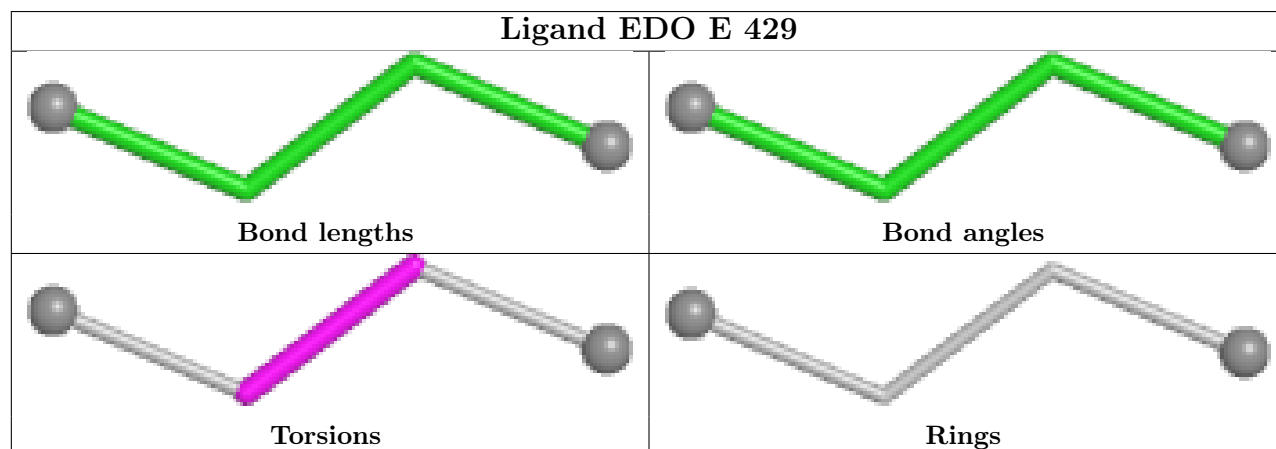
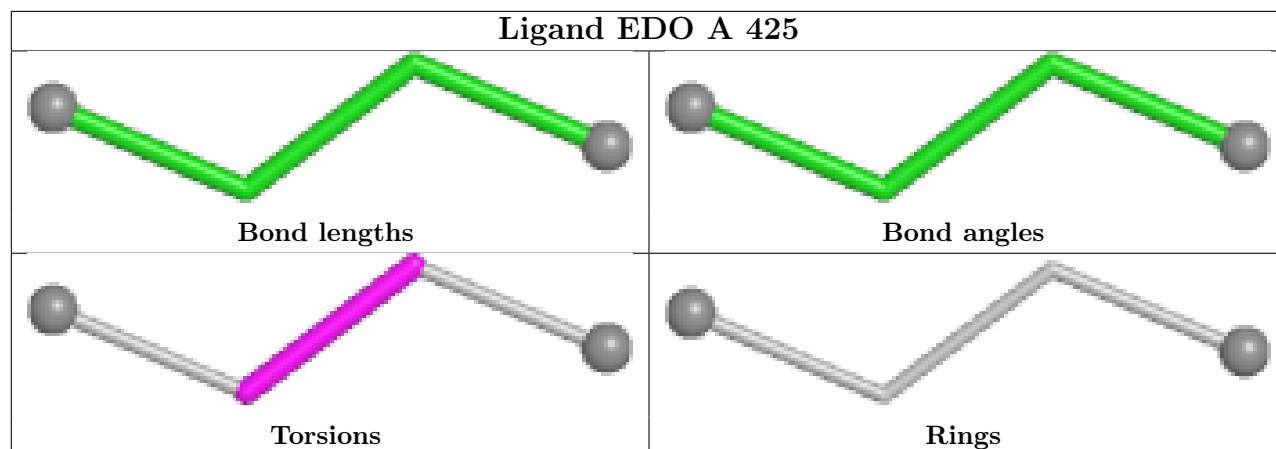


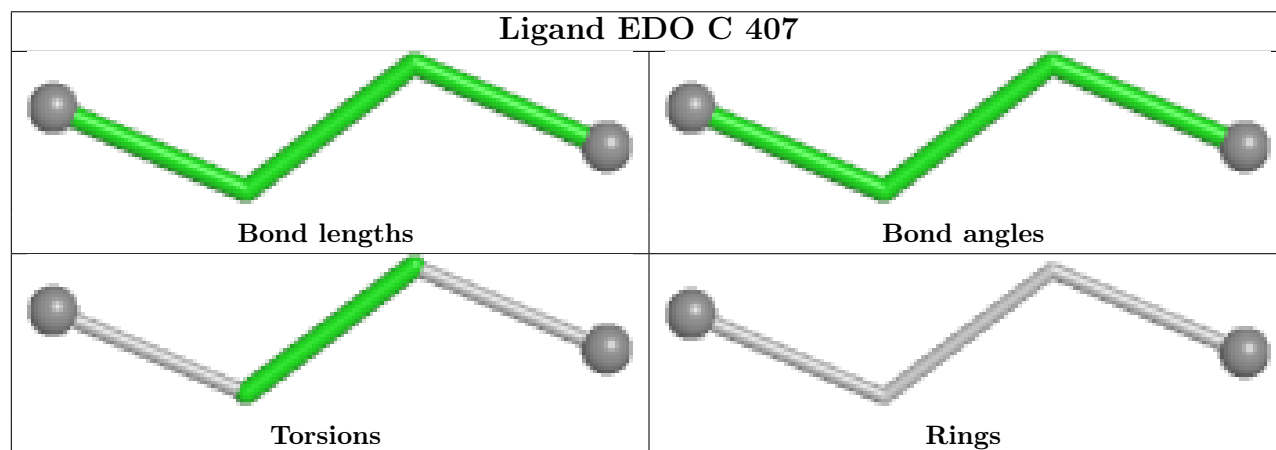
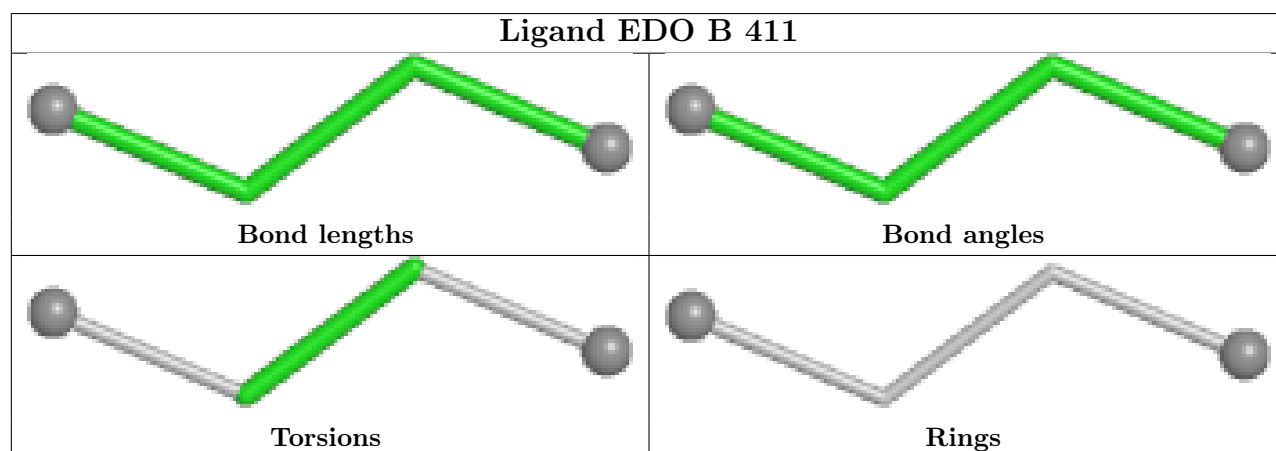
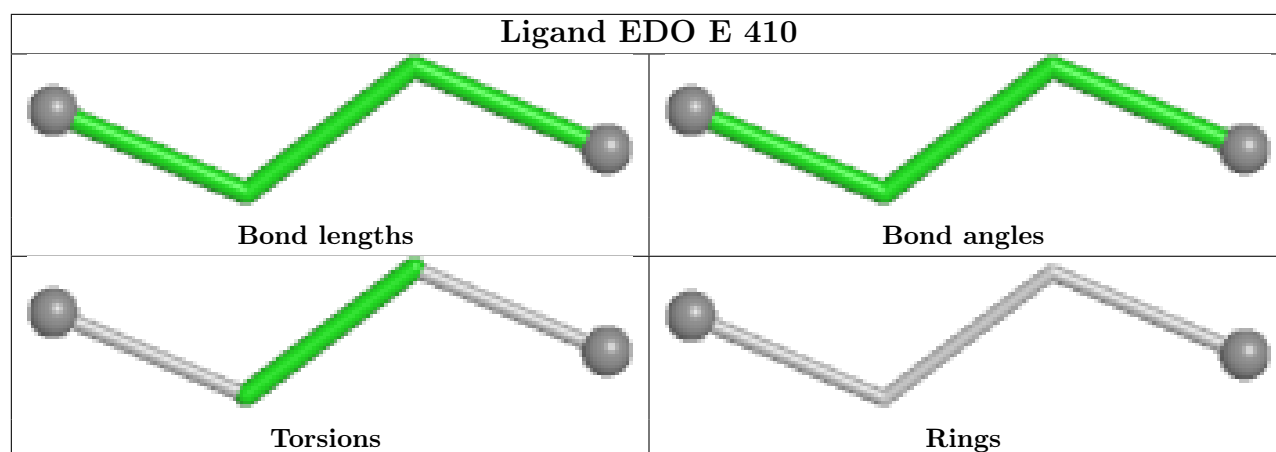


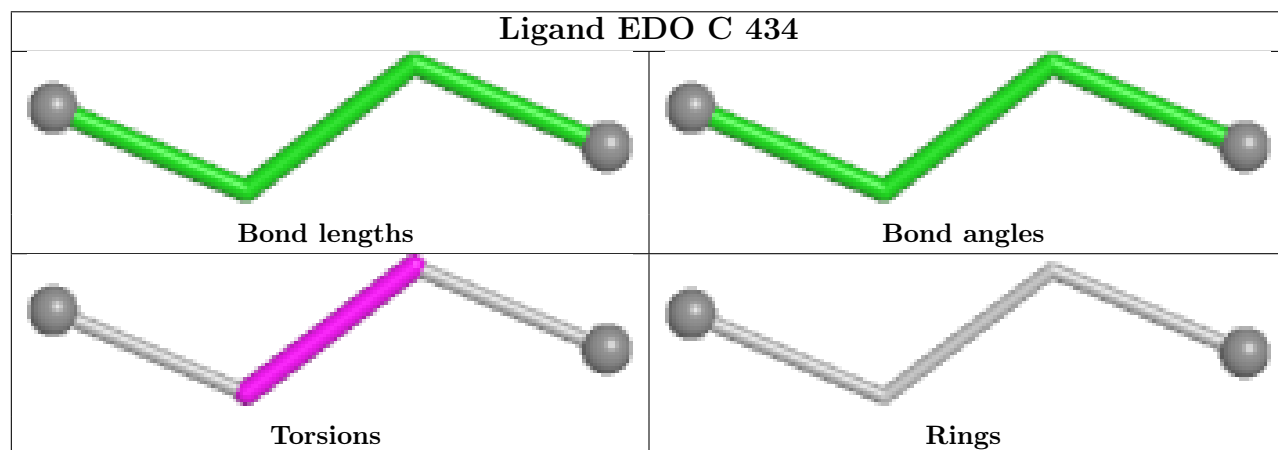
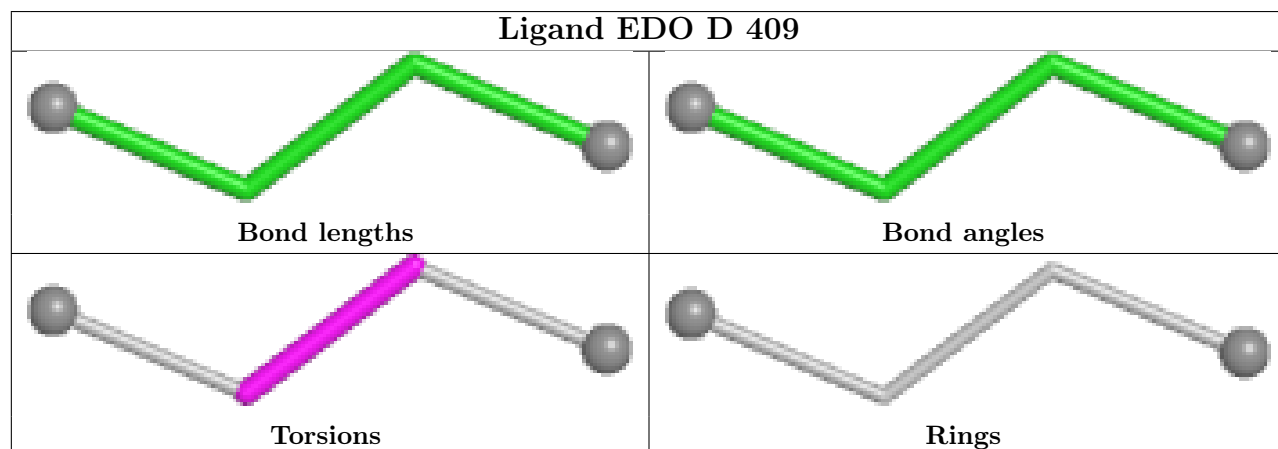
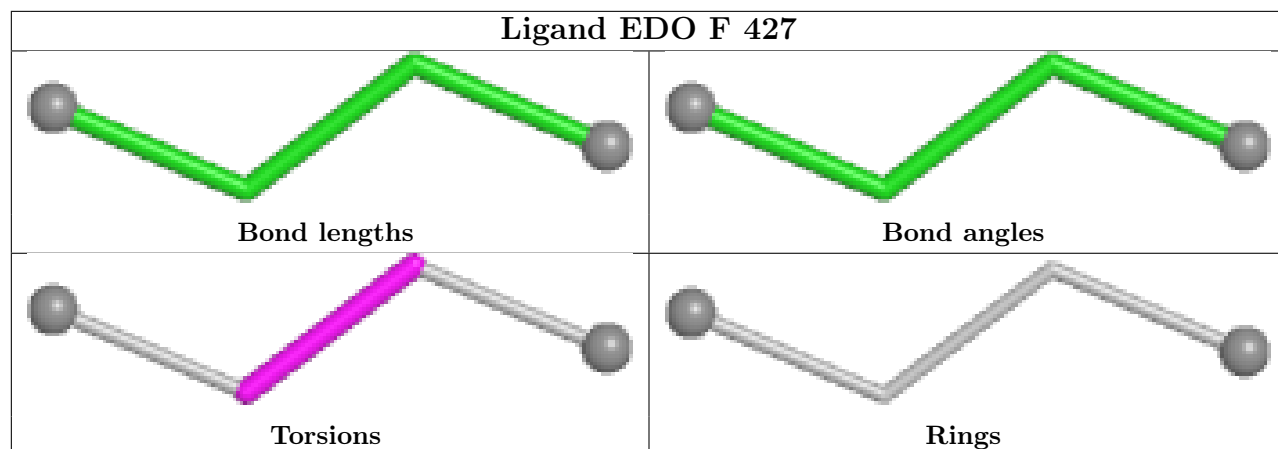


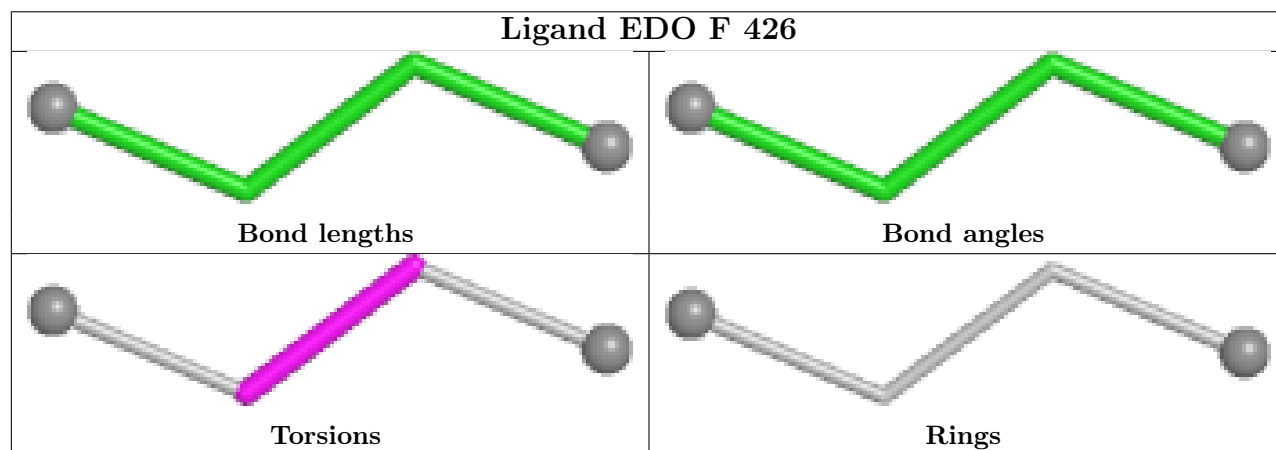
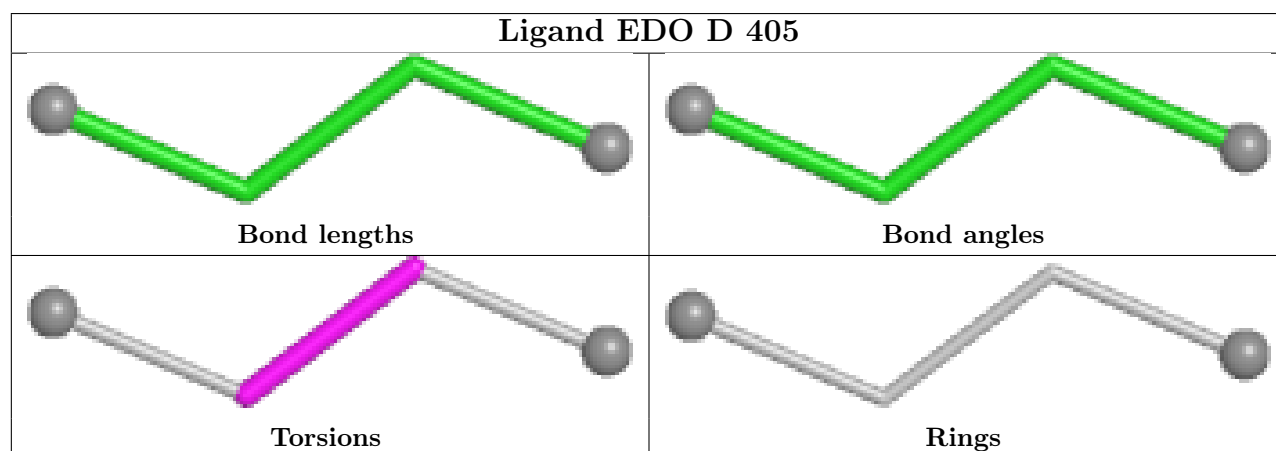
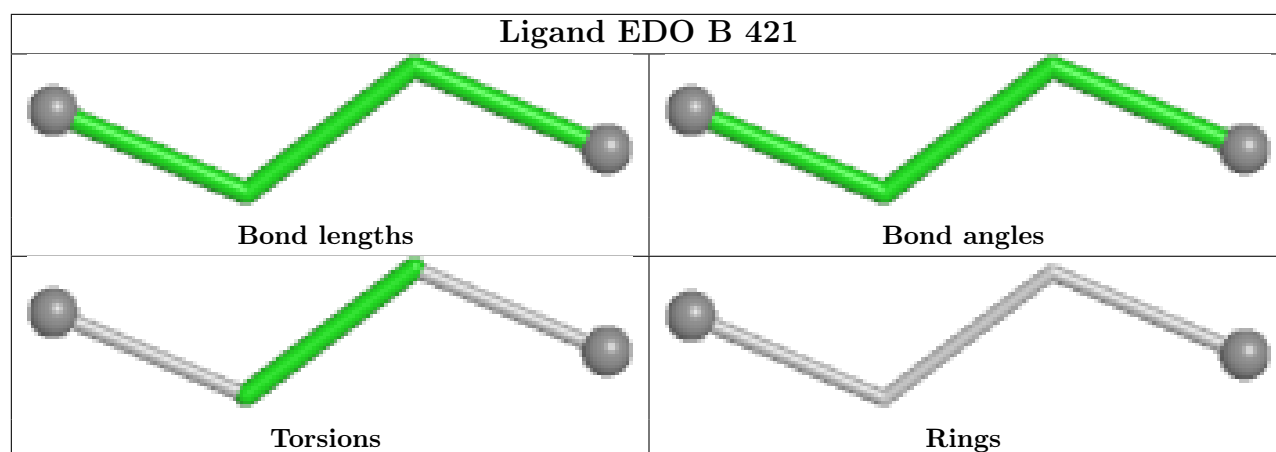


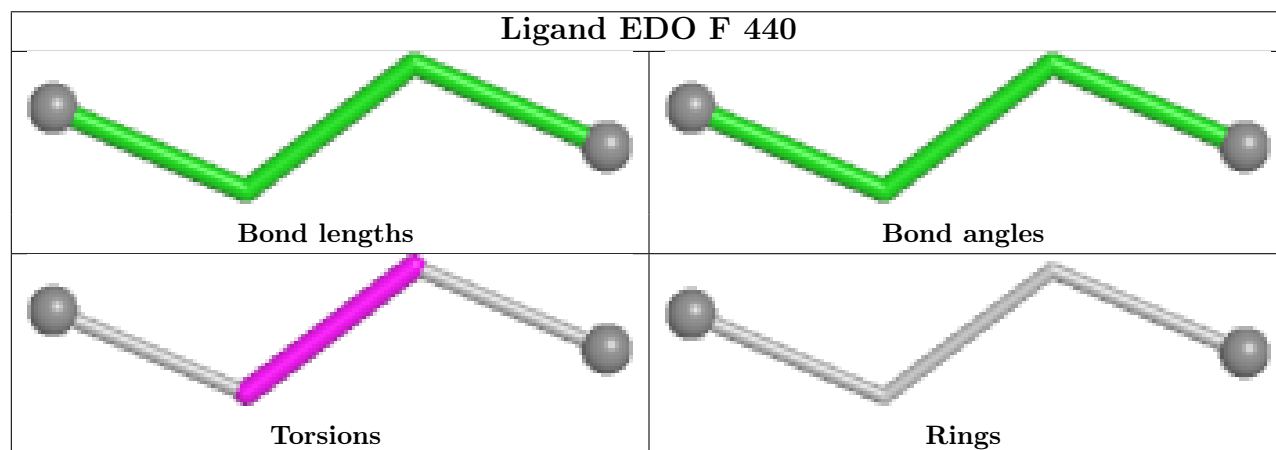
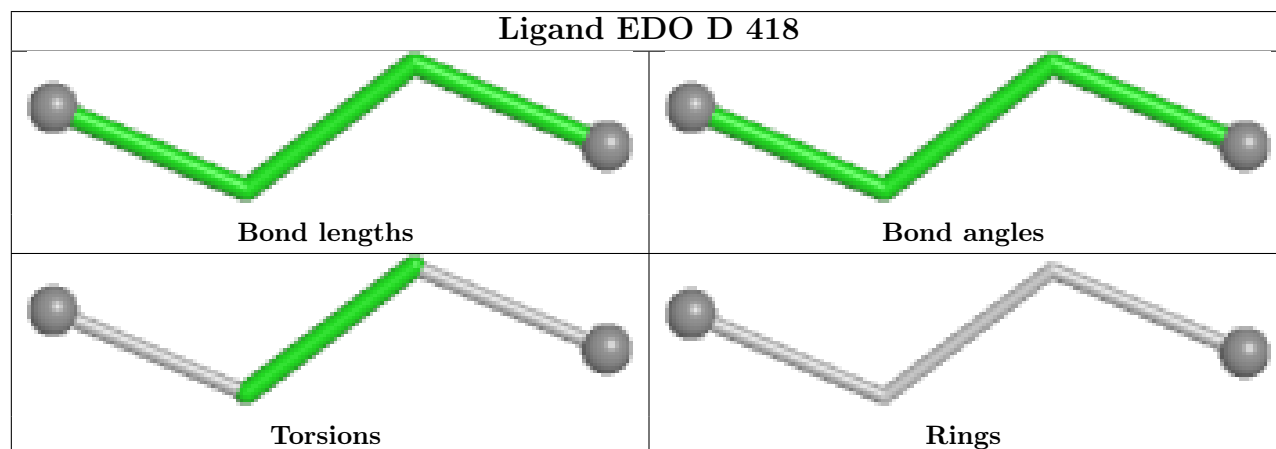
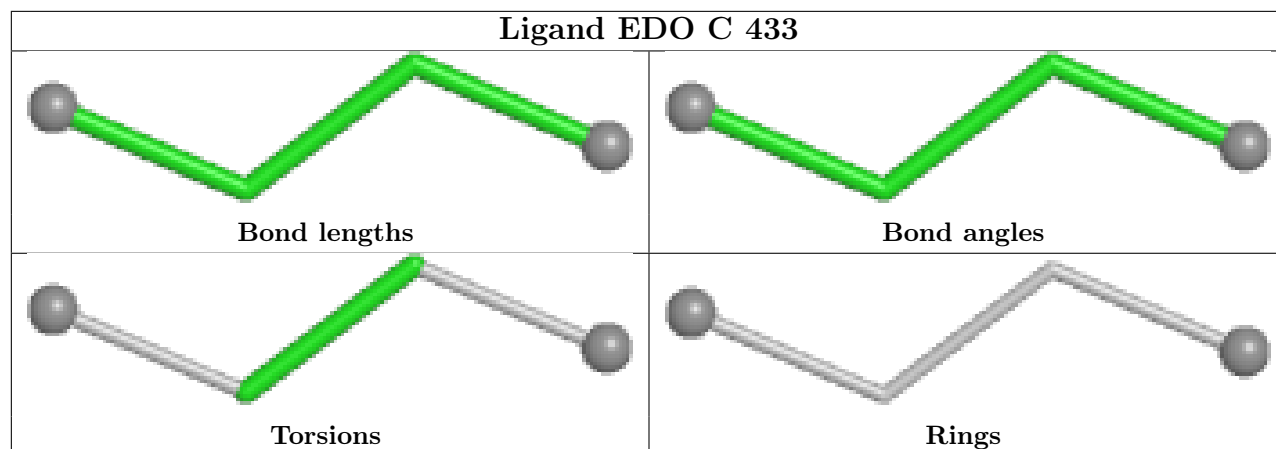


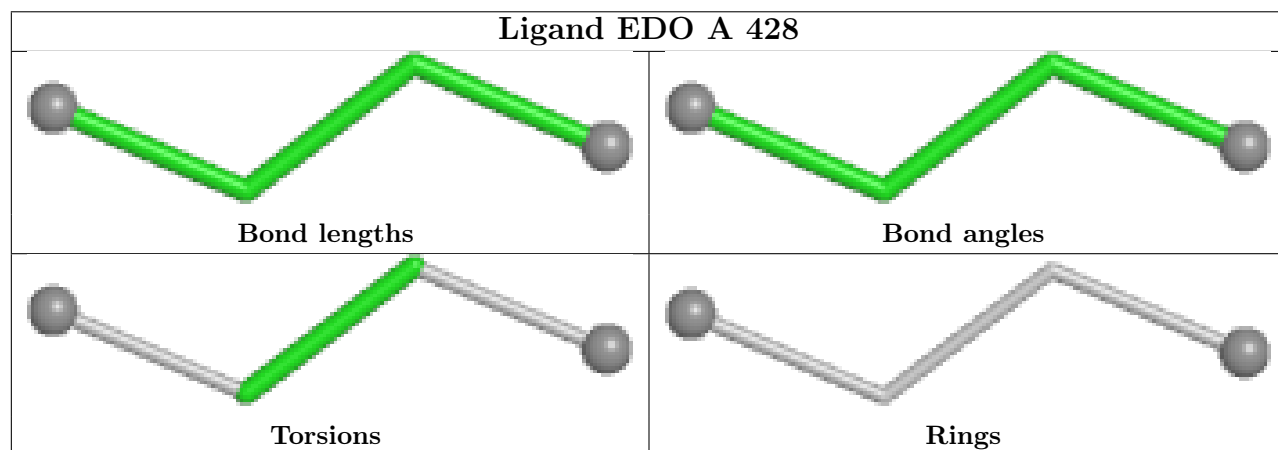
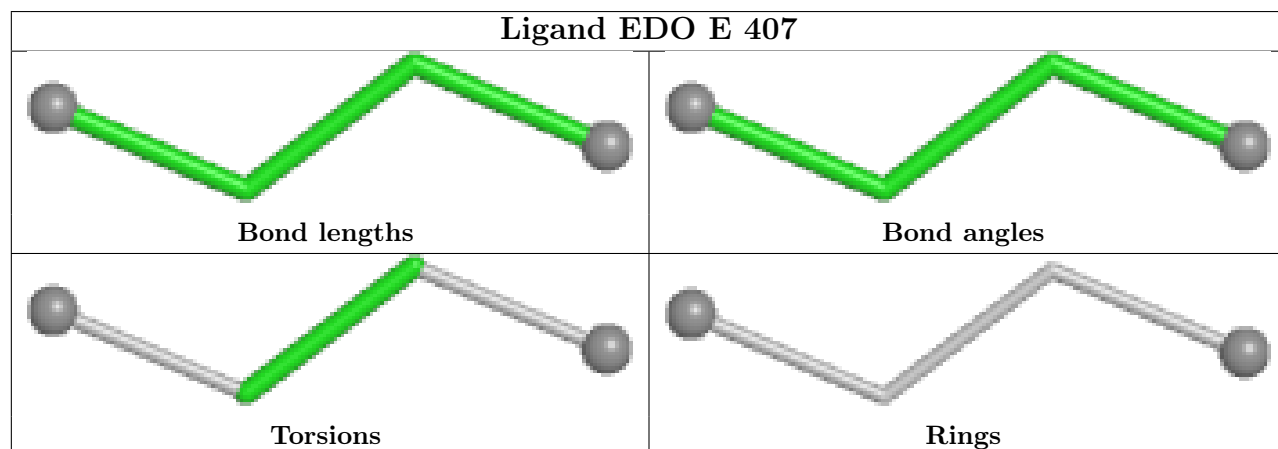
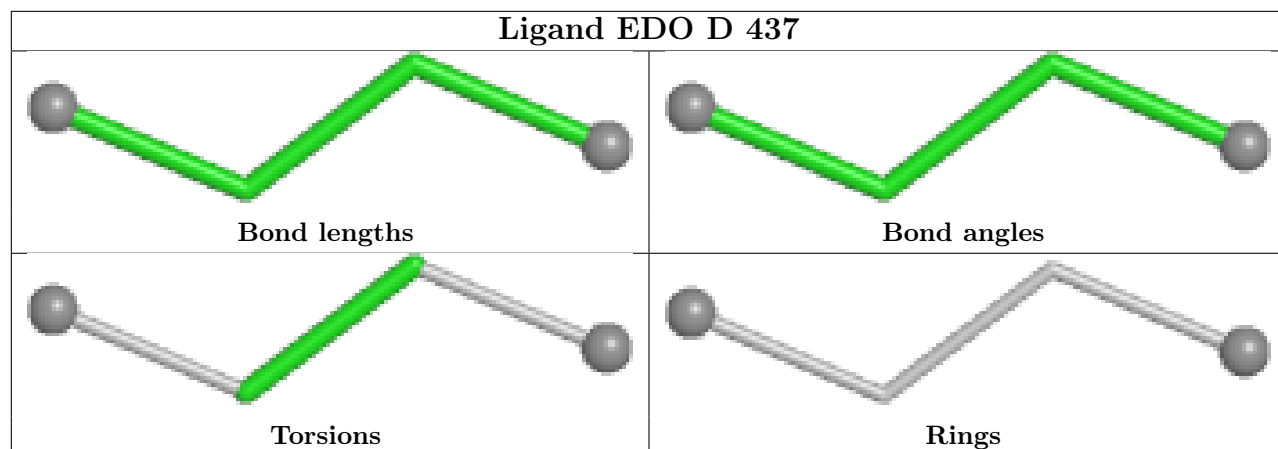


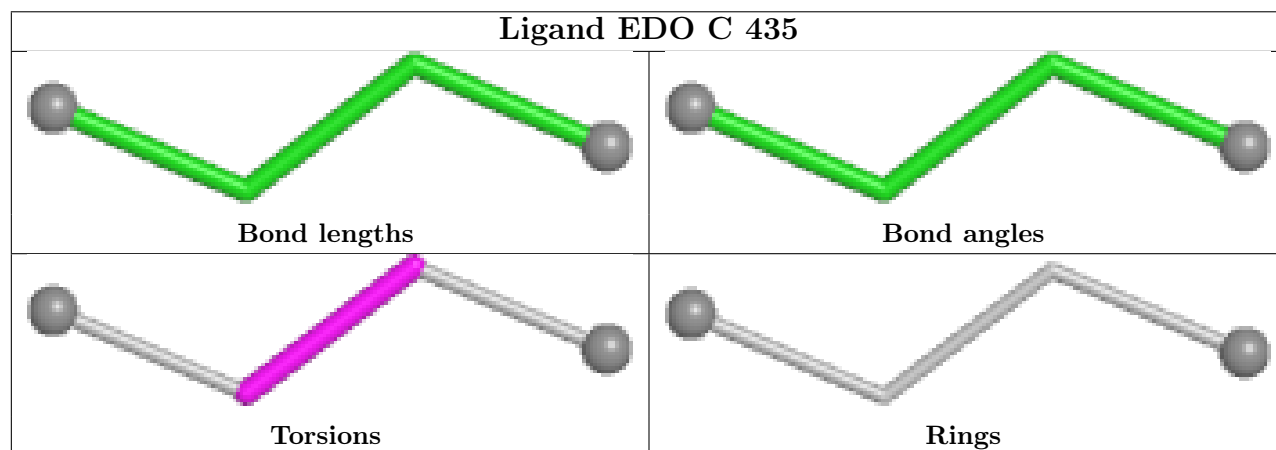
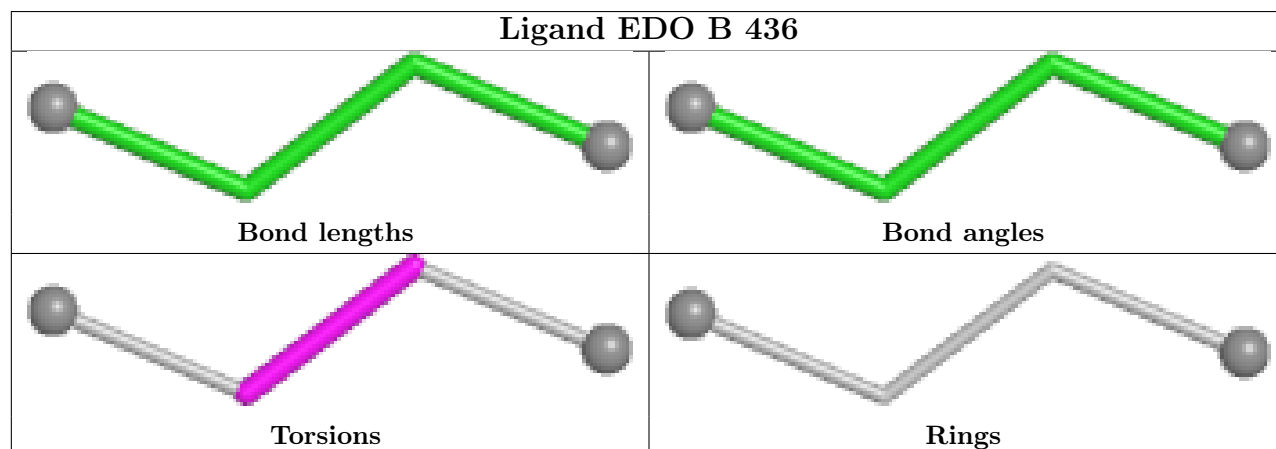
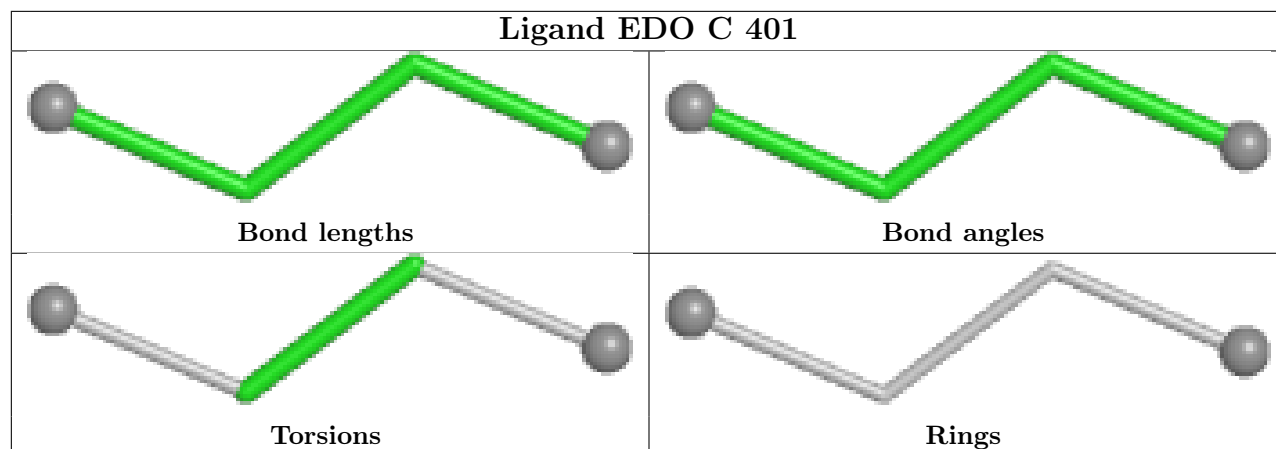


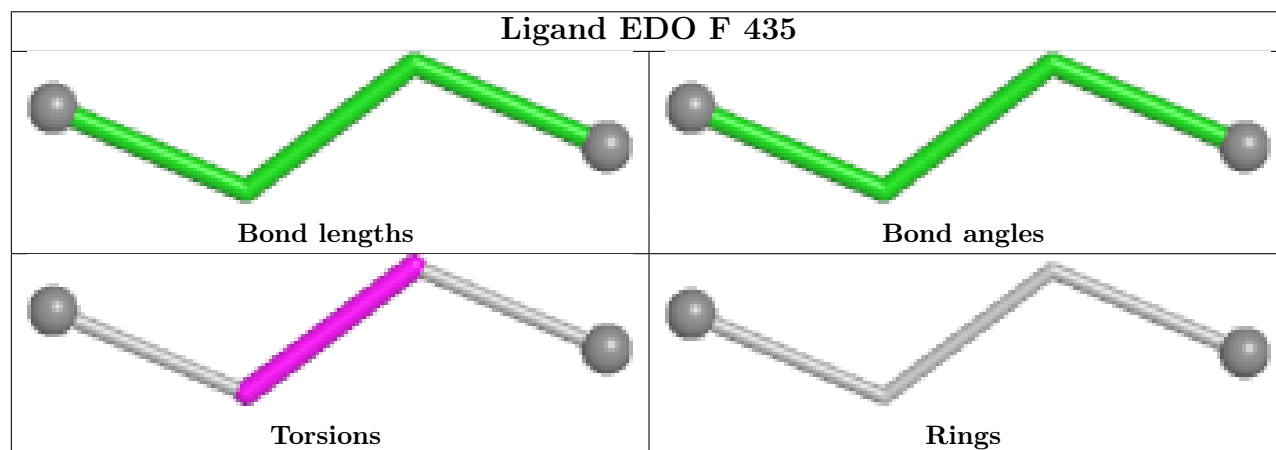
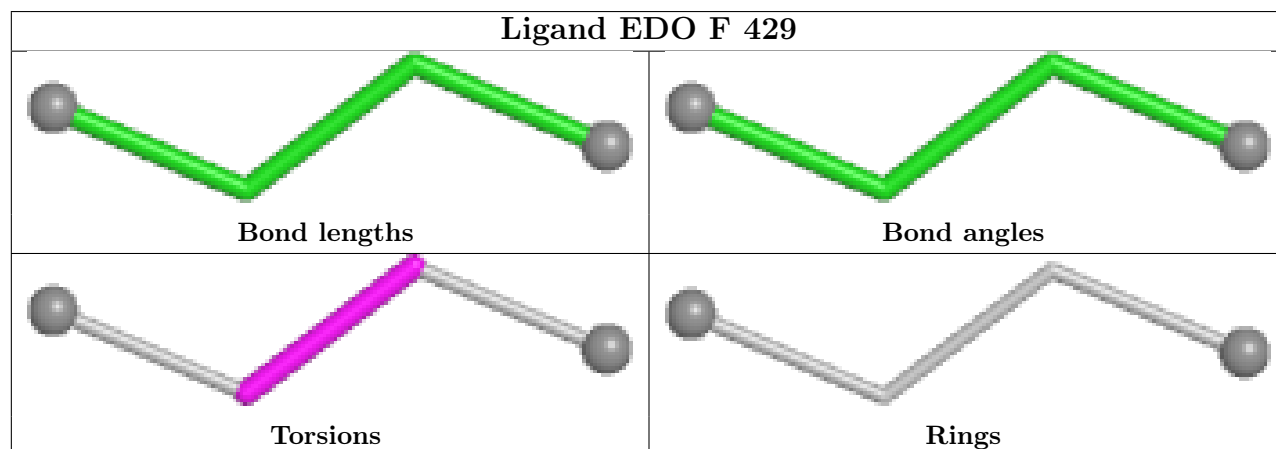
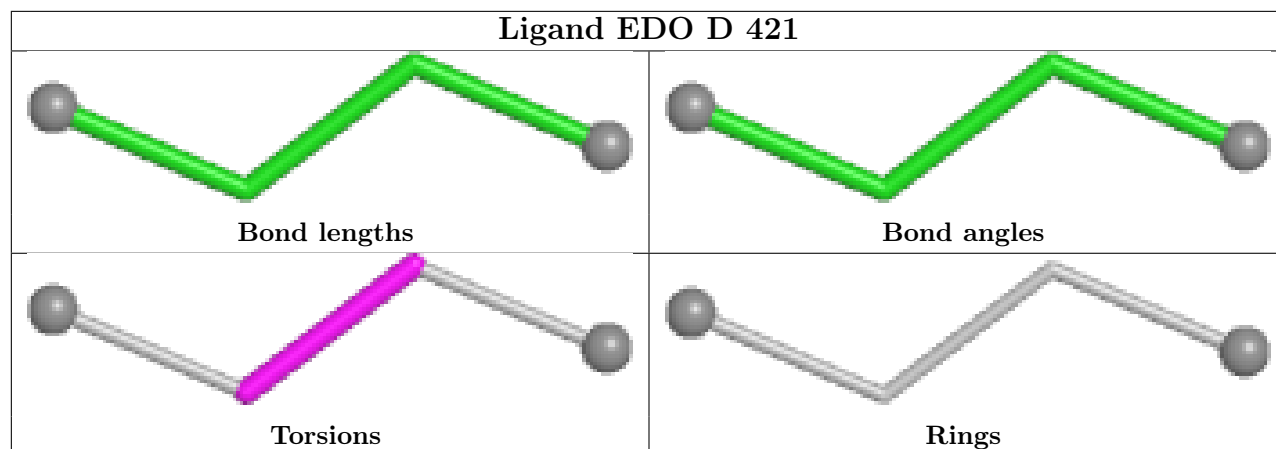




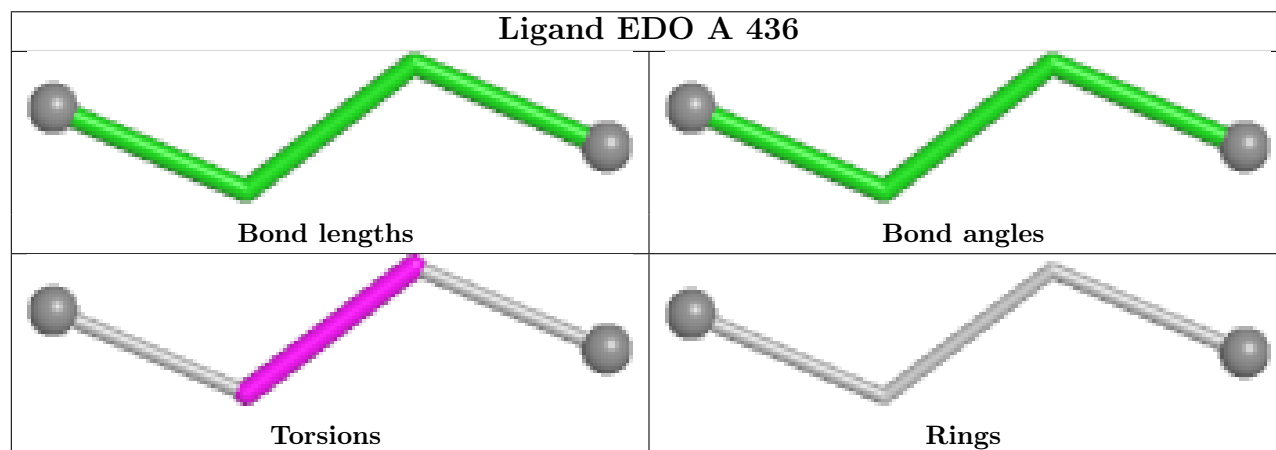
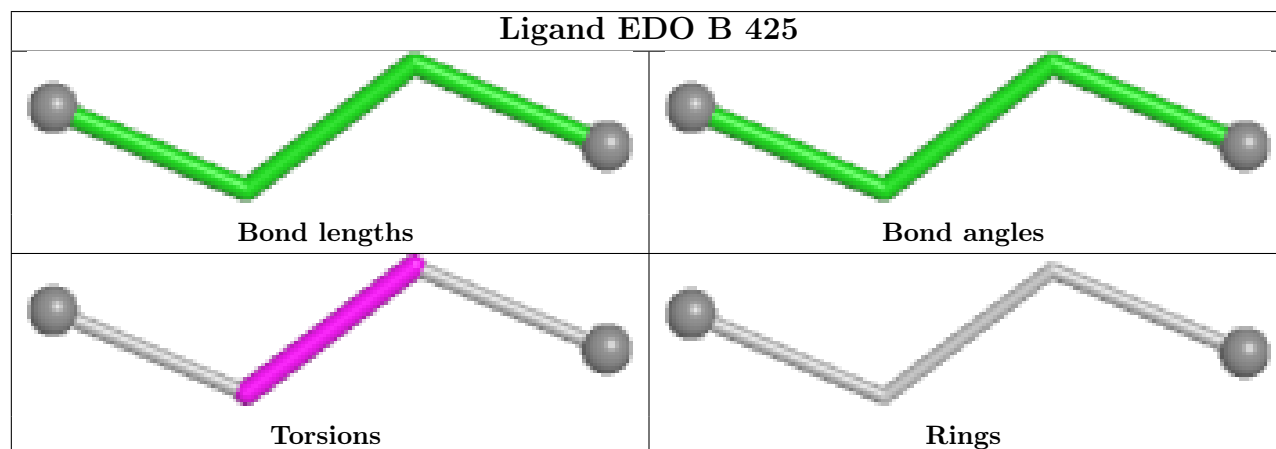
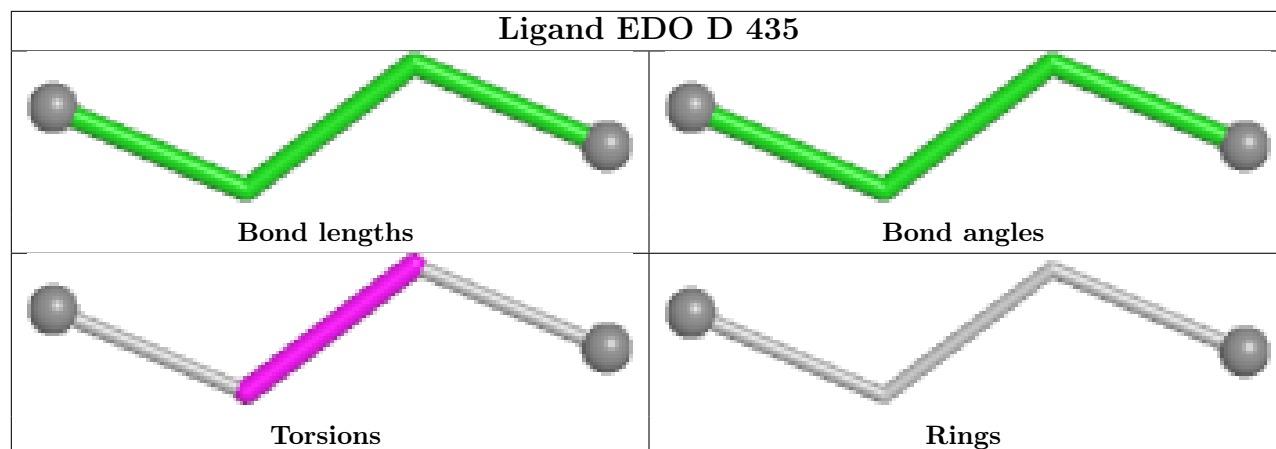


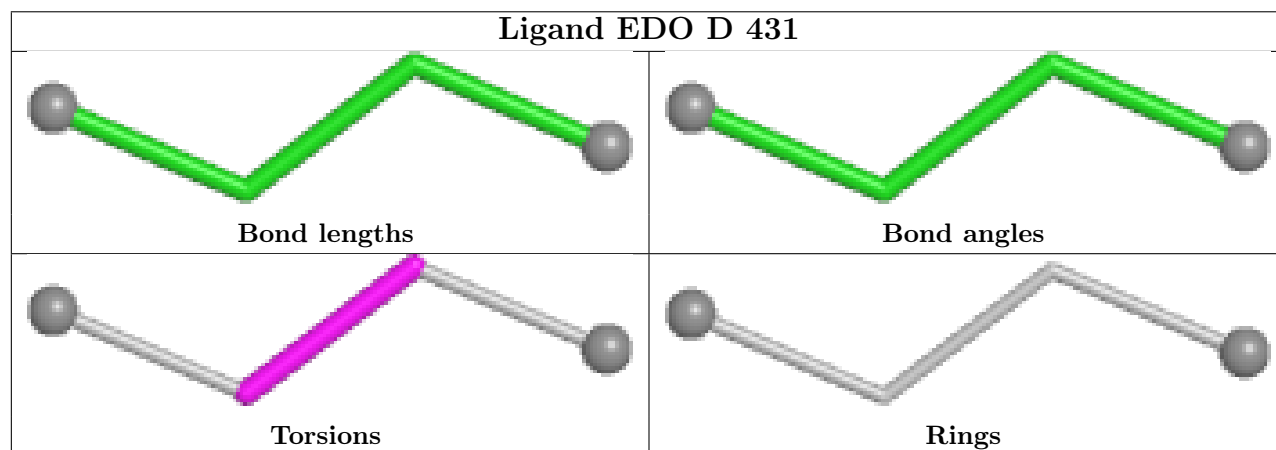
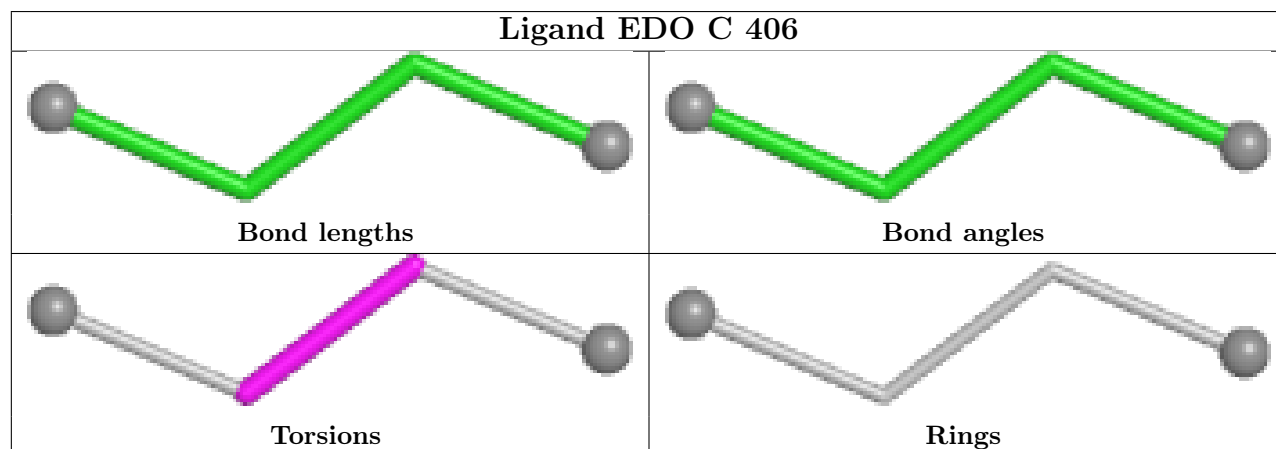
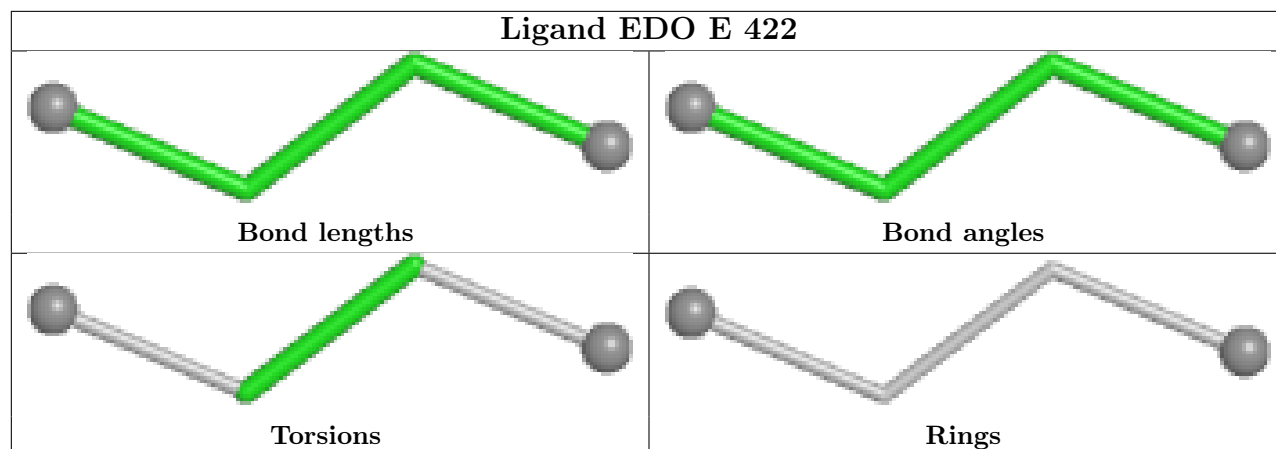


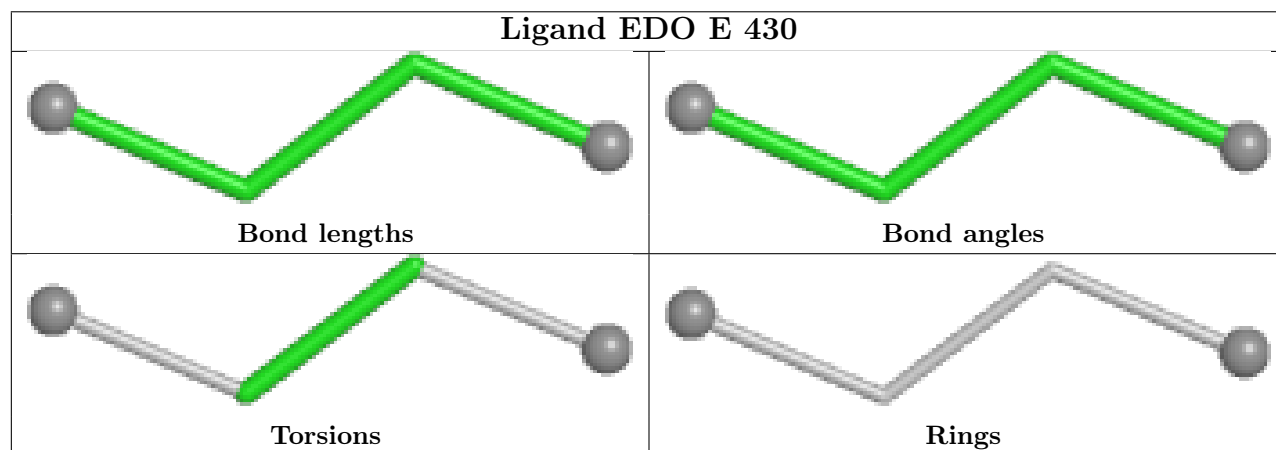
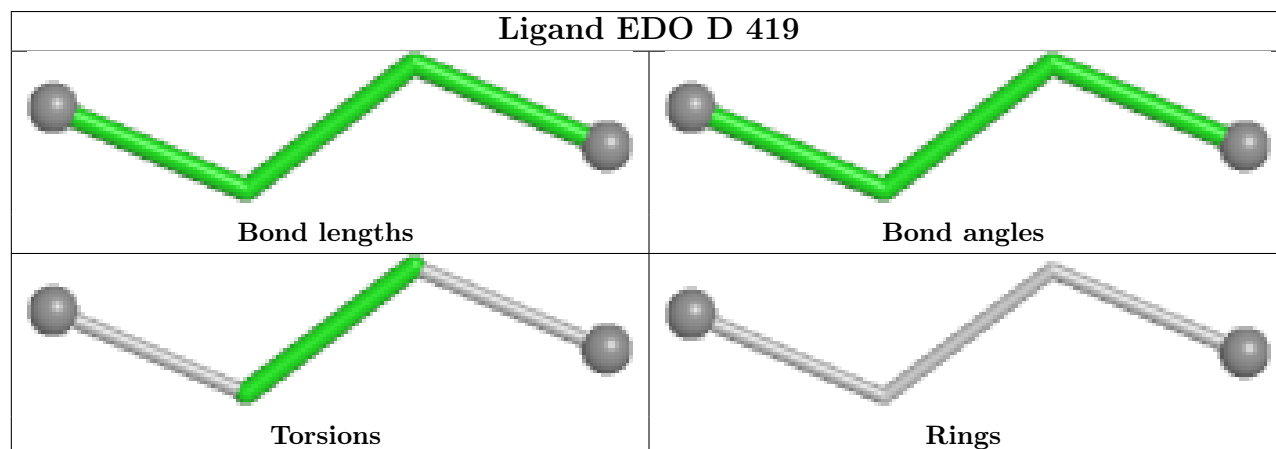
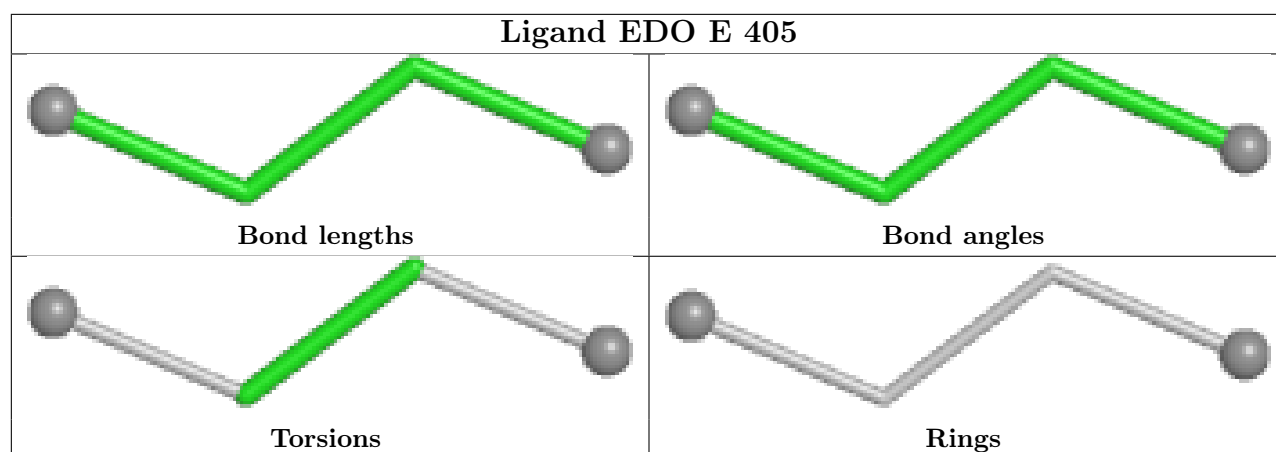


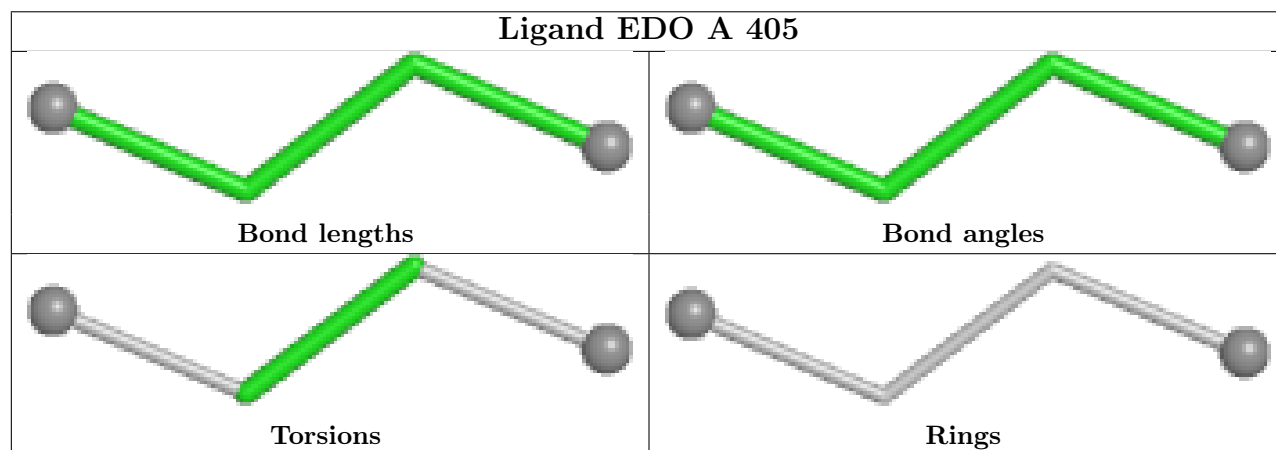
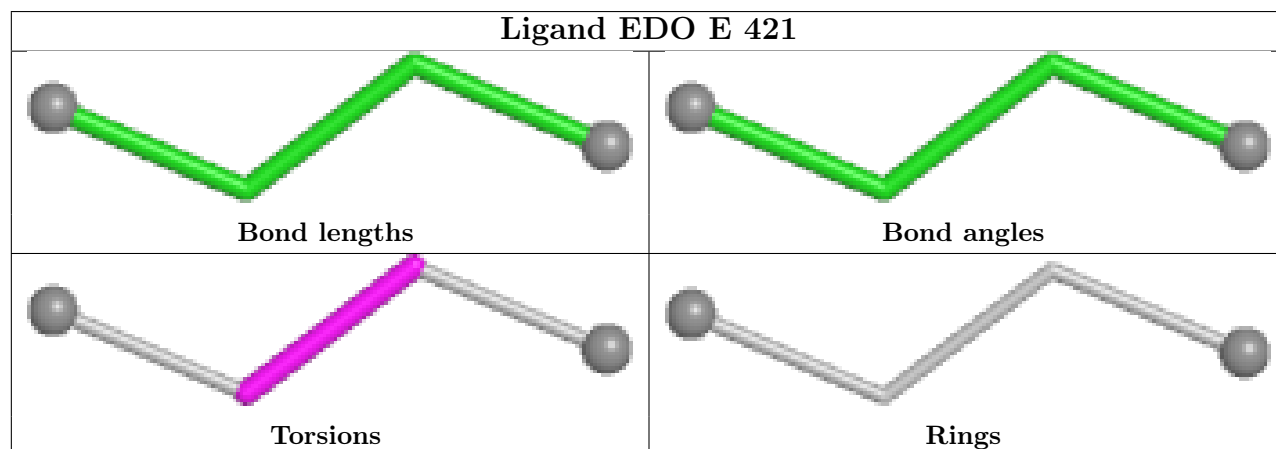
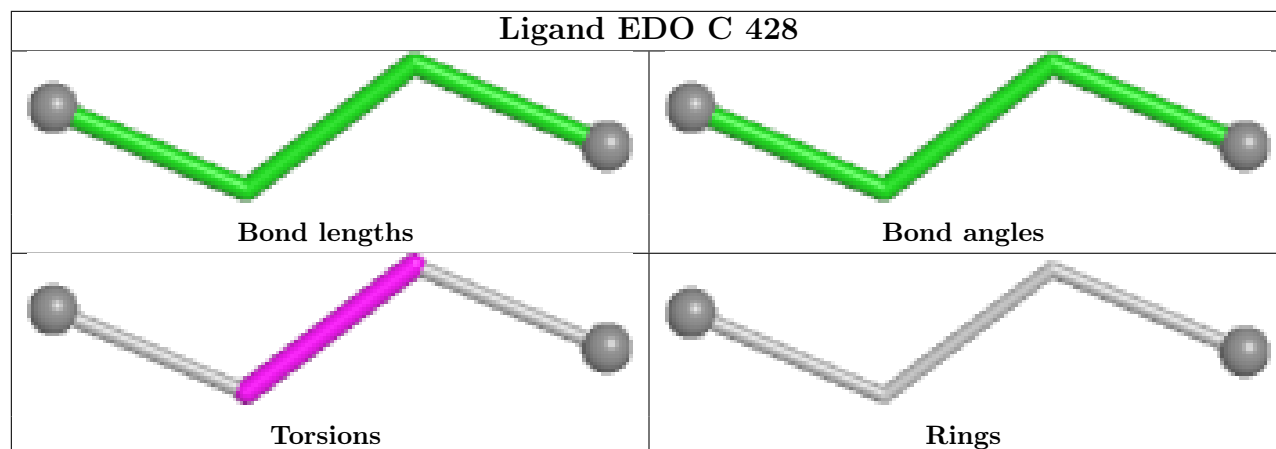


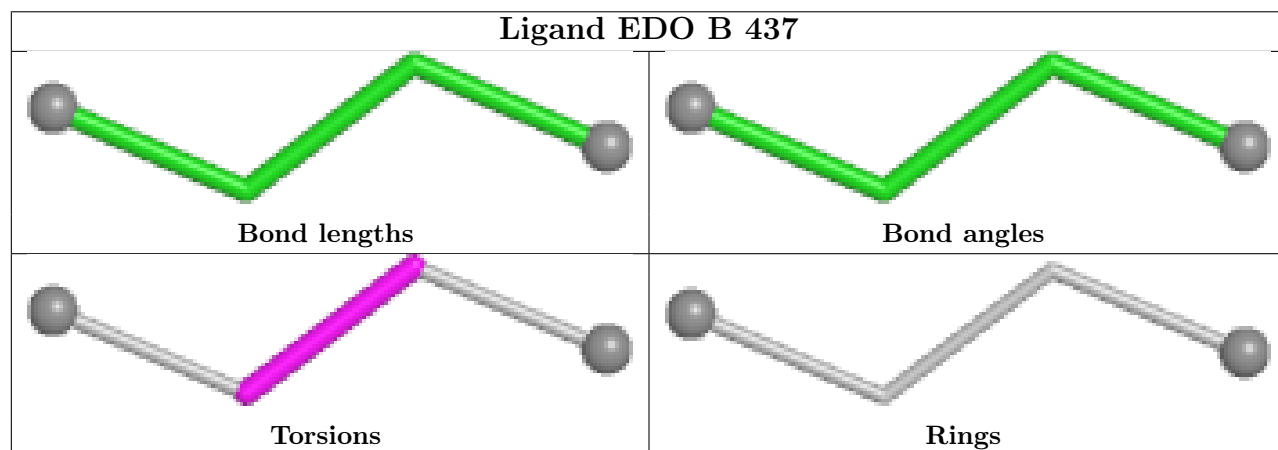
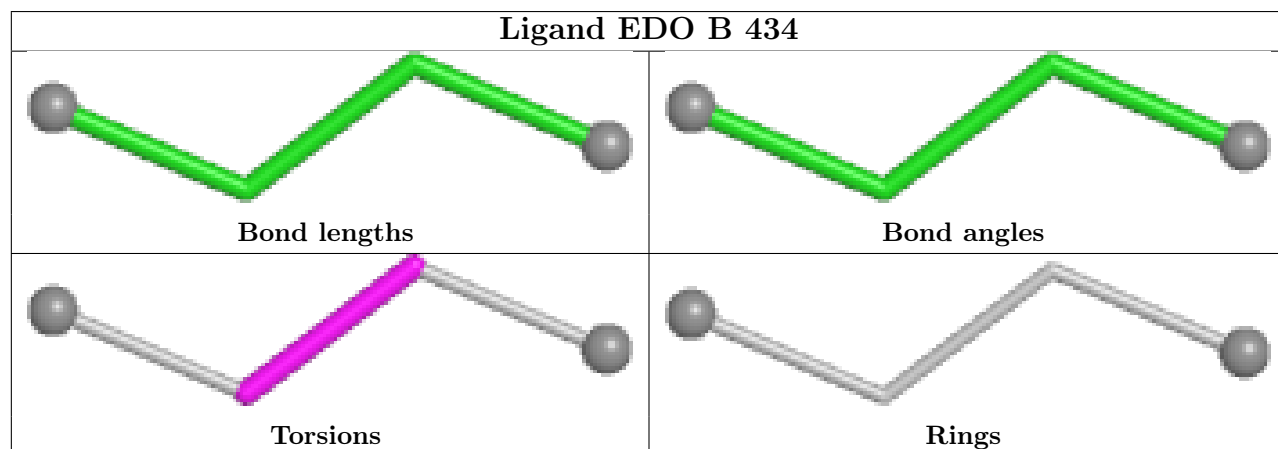
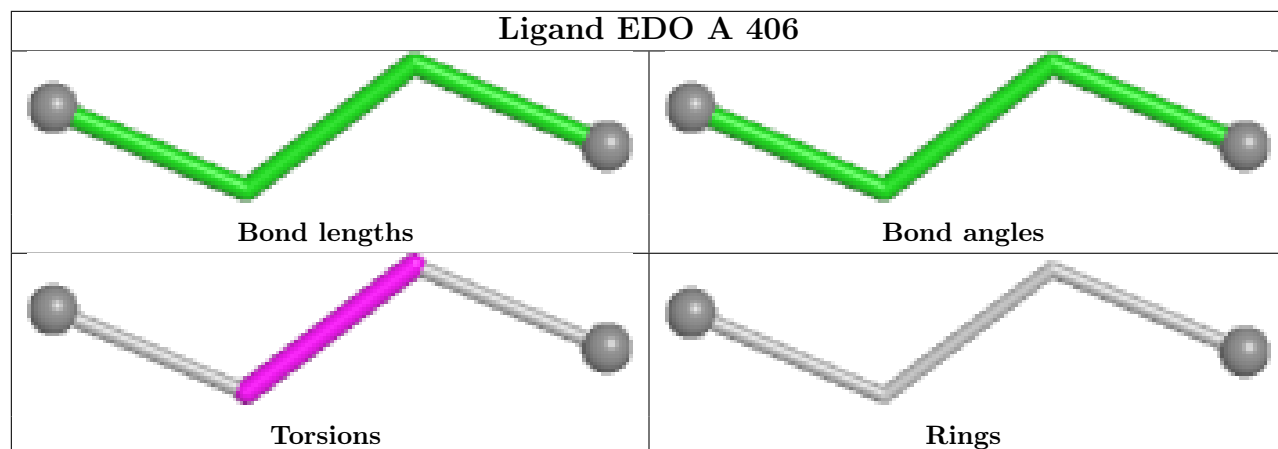


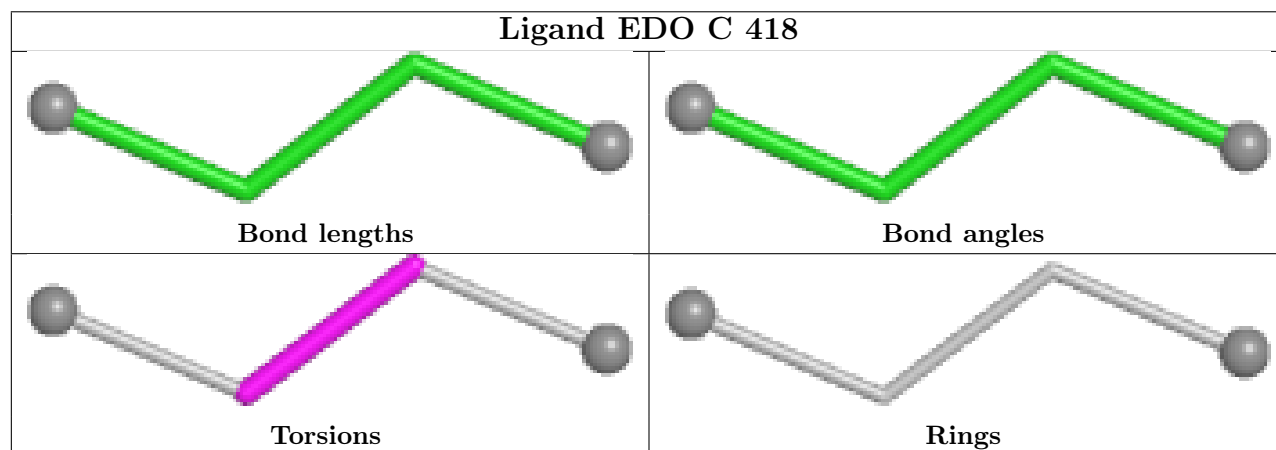
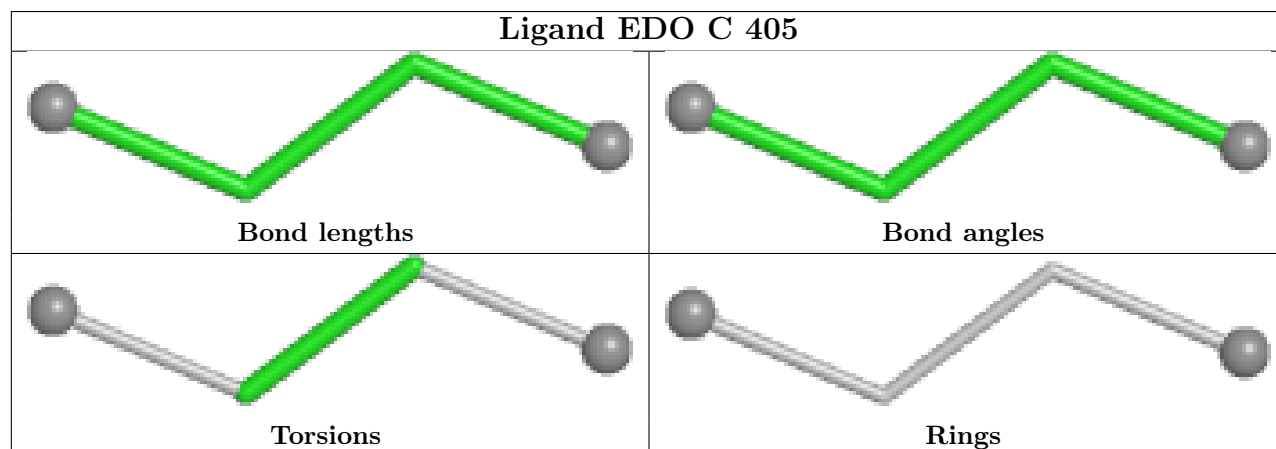
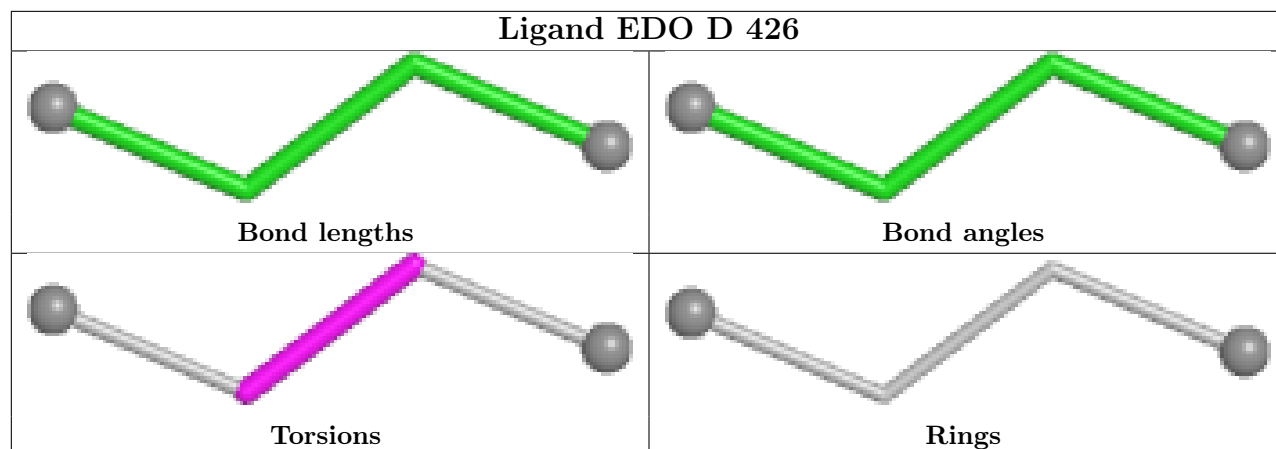


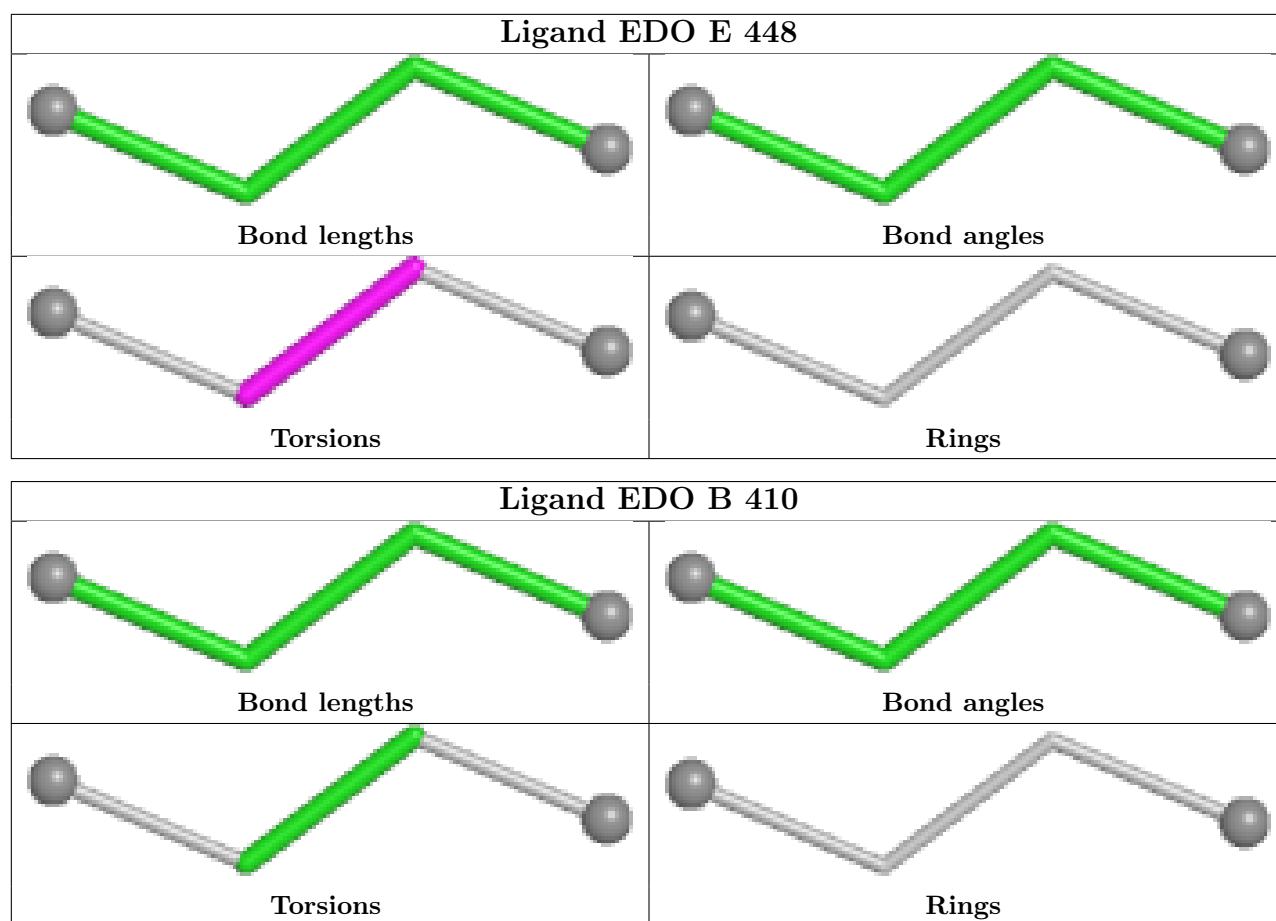












## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	259/259 (100%)	-0.58	6 (2%) 61 41	13, 32, 52, 74	0
1	B	259/259 (100%)	-0.59	2 (0%) 82 67	15, 32, 51, 73	0
1	C	259/259 (100%)	-0.51	5 (1%) 66 46	13, 32, 53, 71	0
1	D	259/259 (100%)	-0.59	2 (0%) 82 67	13, 31, 52, 73	0
1	E	259/259 (100%)	-0.56	4 (1%) 72 52	14, 32, 52, 73	0
1	F	259/259 (100%)	-0.59	2 (0%) 82 67	12, 32, 51, 71	0
All	All	1554/1554 (100%)	-0.57	21 (1%) 73 54	12, 32, 52, 74	0

The worst 5 of 21 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	219	THR	4.3
1	A	353	VAL	3.6
1	C	353	VAL	3.6
1	D	365	ILE	3.4
1	F	353	VAL	3.4

### 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

### 6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.



## 6.4 Ligands

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	EDO	E	446	4/4	0.89	0.16	48,48,48,48	0
2	EDO	C	427	4/4	0.90	0.26	26,26,26,26	0
2	EDO	E	431	4/4	0.90	0.29	26,26,26,26	0
2	EDO	B	428	4/4	0.90	0.20	22,22,22,22	0
2	EDO	F	444	4/4	0.91	0.09	37,37,37,37	0
2	EDO	A	426	4/4	0.92	0.30	27,27,27,27	0
2	EDO	E	445	4/4	0.92	0.33	40,40,40,40	0
2	EDO	B	441	4/4	0.92	0.09	54,54,54,54	0
2	EDO	F	418	4/4	0.92	0.06	19,25,34,34	0
2	EDO	F	434	4/4	0.92	0.29	25,25,25,25	0
2	EDO	B	414	4/4	0.92	0.07	15,20,23,24	0
2	EDO	A	443	4/4	0.93	0.06	41,41,41,41	0
2	EDO	E	432	4/4	0.93	0.20	32,32,32,32	0
2	EDO	B	403	4/4	0.93	0.17	7,8,11,16	0
2	EDO	A	423	4/4	0.93	0.06	17,17,17,17	0
2	EDO	C	436	4/4	0.93	0.11	36,36,36,36	0
2	EDO	D	425	4/4	0.93	0.09	72,72,72,72	0
2	EDO	E	426	4/4	0.93	0.08	70,70,70,70	0
2	EDO	D	403	4/4	0.94	0.24	54,61,66,68	0
2	EDO	D	420	4/4	0.94	0.13	61,68,73,75	0
2	EDO	A	428	4/4	0.94	0.09	39,39,39,39	0
2	EDO	D	429	4/4	0.94	0.17	20,20,20,20	0
2	EDO	B	434	4/4	0.94	0.09	26,26,26,26	0
2	EDO	B	437	4/4	0.94	0.10	23,23,23,23	0
2	EDO	A	401	4/4	0.94	0.28	4,6,9,14	0
2	EDO	E	433	4/4	0.94	0.09	39,39,39,39	0
2	EDO	B	446	4/4	0.94	0.14	56,56,56,56	0
2	EDO	C	403	4/4	0.94	0.20	5,7,10,15	0
2	EDO	F	403	4/4	0.94	0.05	29,29,29,29	0
2	EDO	F	404	4/4	0.94	0.09	21,21,21,21	0
2	EDO	A	416	4/4	0.94	0.07	42,49,54,56	0
2	EDO	C	433	4/4	0.94	0.09	21,21,21,21	0
2	EDO	F	443	4/4	0.94	0.30	37,37,37,37	0
2	EDO	A	427	4/4	0.94	0.27	37,37,37,37	0
2	EDO	F	445	4/4	0.94	0.07	46,46,46,46	0
2	EDO	B	426	4/4	0.95	0.10	11,11,11,11	0
2	EDO	E	441	4/4	0.95	0.16	35,35,35,35	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	EDO	D	404	4/4	0.95	0.22	6,8,11,15	0
2	EDO	C	418	4/4	0.95	0.06	24,27,28,29	0
2	EDO	B	438	4/4	0.95	0.21	37,37,37,37	0
2	EDO	C	432	4/4	0.95	0.05	34,34,34,34	0
2	EDO	F	406	4/4	0.95	0.06	22,22,22,22	0
2	EDO	D	439	4/4	0.95	0.05	40,40,40,40	0
2	EDO	F	420	4/4	0.95	0.06	13,18,21,22	0
2	EDO	F	424	4/4	0.95	0.06	42,50,55,56	0
2	EDO	A	431	4/4	0.95	0.05	30,30,30,30	0
2	EDO	F	436	4/4	0.95	0.13	38,38,38,38	0
2	EDO	F	440	4/4	0.95	0.09	26,26,26,26	0
2	EDO	F	442	4/4	0.95	0.12	34,34,34,34	0
2	EDO	E	429	4/4	0.95	0.10	11,11,11,11	0
2	EDO	A	421	4/4	0.95	0.06	21,21,21,21	0
2	EDO	D	402	4/4	0.95	0.11	17,23,26,26	0
2	EDO	D	424	4/4	0.96	0.07	23,23,23,23	0
2	EDO	A	440	4/4	0.96	0.12	19,19,19,19	0
2	EDO	D	426	4/4	0.96	0.05	19,19,19,19	0
2	EDO	B	444	4/4	0.96	0.10	30,30,30,30	0
2	EDO	D	432	4/4	0.96	0.07	39,39,39,39	0
2	EDO	D	433	4/4	0.96	0.08	22,22,22,22	0
2	EDO	D	435	4/4	0.96	0.04	34,34,34,34	0
2	EDO	D	436	4/4	0.96	0.08	36,36,36,36	0
2	EDO	A	442	4/4	0.96	0.10	28,28,28,28	0
2	EDO	E	403	4/4	0.96	0.24	4,6,9,14	0
2	EDO	E	404	4/4	0.96	0.10	15,19,19,37	0
2	EDO	E	411	4/4	0.96	0.06	13,28,35,40	0
2	EDO	E	413	4/4	0.96	0.05	18,24,33,33	0
2	EDO	E	419	4/4	0.96	0.06	42,50,54,56	0
2	EDO	E	421	4/4	0.96	0.05	19,21,23,24	0
2	EDO	B	447	4/4	0.96	0.05	32,32,32,32	0
2	EDO	E	428	4/4	0.96	0.04	32,32,32,32	0
2	EDO	C	402	4/4	0.96	0.08	14,19,22,23	0
2	EDO	A	417	4/4	0.96	0.05	13,16,17,18	0
2	EDO	A	412	4/4	0.96	0.06	10,15,17,19	0
2	EDO	C	420	4/4	0.96	0.09	15,15,15,15	0
2	EDO	E	434	4/4	0.96	0.08	20,20,20,20	0
2	EDO	C	424	4/4	0.96	0.04	30,30,30,30	0
2	EDO	E	443	4/4	0.96	0.11	18,18,18,18	0
2	EDO	C	425	4/4	0.96	0.08	15,15,15,15	0
2	EDO	A	422	4/4	0.96	0.07	73,73,73,73	0
2	EDO	C	428	4/4	0.96	0.18	37,37,37,37	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	EDO	B	424	4/4	0.96	0.08	68,68,68,68	0
2	EDO	A	403	4/4	0.96	0.10	19,21,23,28	0
2	EDO	F	408	4/4	0.96	0.10	9,15,18,18	0
2	EDO	F	409	4/4	0.96	0.14	7,9,12,17	0
2	EDO	F	416	4/4	0.96	0.13	2,3,4,22	0
2	EDO	C	434	4/4	0.96	0.07	35,35,35,35	0
2	EDO	A	434	4/4	0.96	0.05	19,19,19,19	0
2	EDO	C	438	4/4	0.96	0.08	40,40,40,40	0
2	EDO	F	425	4/4	0.96	0.07	24,26,28,29	0
2	EDO	F	427	4/4	0.96	0.09	18,18,18,18	0
2	EDO	F	429	4/4	0.96	0.07	26,26,26,26	0
2	EDO	F	430	4/4	0.96	0.05	58,58,58,58	0
2	EDO	B	430	4/4	0.96	0.09	30,30,30,30	0
2	EDO	A	436	4/4	0.96	0.07	40,40,40,40	0
2	EDO	F	439	4/4	0.96	0.07	24,24,24,24	0
2	EDO	A	438	4/4	0.96	0.06	30,30,30,30	0
2	EDO	F	441	4/4	0.96	0.11	55,55,55,55	0
2	EDO	D	406	4/4	0.96	0.11	29,31,33,37	0
2	EDO	D	407	4/4	0.96	0.07	4,17,20,28	0
2	EDO	A	439	4/4	0.96	0.09	30,30,30,30	0
2	EDO	D	422	4/4	0.96	0.08	18,18,18,18	0
2	EDO	D	434	4/4	0.97	0.08	38,38,38,38	0
2	EDO	B	405	4/4	0.97	0.09	21,23,25,30	0
2	EDO	C	406	4/4	0.97	0.06	5,17,20,28	0
2	EDO	C	410	4/4	0.97	0.12	4,5,5,24	0
2	EDO	E	401	4/4	0.97	0.04	25,25,25,25	0
2	EDO	E	402	4/4	0.97	0.05	26,28,30,31	0
2	EDO	C	414	4/4	0.97	0.05	14,19,21,23	0
2	EDO	B	410	4/4	0.97	0.12	3,4,4,23	0
2	EDO	E	406	4/4	0.97	0.04	8,21,23,31	0
2	EDO	E	410	4/4	0.97	0.12	5,6,6,24	0
2	EDO	B	412	4/4	0.97	0.03	16,22,31,31	0
2	EDO	C	422	4/4	0.97	0.07	18,18,18,18	0
2	EDO	A	429	4/4	0.97	0.05	14,28,35,41	0
2	EDO	B	417	4/4	0.97	0.04	2,2,3,8	0
2	EDO	E	423	4/4	0.97	0.08	14,14,14,14	0
2	EDO	E	425	4/4	0.97	0.06	21,21,21,21	0
2	EDO	B	418	4/4	0.97	0.05	53,61,66,67	0
2	EDO	E	427	4/4	0.97	0.04	20,20,20,20	0
2	EDO	B	419	4/4	0.97	0.04	18,21,22,23	0
2	EDO	C	429	4/4	0.97	0.10	33,33,33,33	0
2	EDO	C	431	4/4	0.97	0.05	27,28,31,36	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	EDO	B	423	4/4	0.97	0.06	22,22,22,22	0
2	EDO	A	407	4/4	0.97	0.12	6,7,14,17	0
2	EDO	A	432	4/4	0.97	0.08	15,15,15,15	0
2	EDO	E	437	4/4	0.97	0.04	20,20,20,20	0
2	EDO	E	438	4/4	0.97	0.11	26,26,26,26	0
2	EDO	E	439	4/4	0.97	0.09	31,31,31,31	0
2	EDO	C	435	4/4	0.97	0.06	31,36,40,40	0
2	EDO	A	433	4/4	0.97	0.07	33,33,33,33	0
2	EDO	C	437	4/4	0.97	0.15	43,43,43,43	0
2	EDO	B	429	4/4	0.97	0.17	36,36,36,36	0
2	EDO	E	447	4/4	0.97	0.05	42,42,42,42	0
2	EDO	D	401	4/4	0.97	0.05	4,5,5,6	0
2	EDO	A	419	4/4	0.97	0.09	16,16,16,16	0
2	EDO	B	431	4/4	0.97	0.06	11,25,32,38	0
2	EDO	F	407	4/4	0.97	0.05	6,6,9,11	0
2	EDO	A	408	4/4	0.97	0.14	4,5,6,24	0
2	EDO	B	435	4/4	0.97	0.05	26,26,26,26	0
2	EDO	F	411	4/4	0.97	0.10	22,24,26,30	0
2	EDO	A	410	4/4	0.97	0.05	17,24,32,32	0
2	EDO	D	408	4/4	0.97	0.07	13,14,17,18	0
2	EDO	D	411	4/4	0.97	0.12	5,6,6,24	0
2	EDO	D	415	4/4	0.97	0.05	8,13,15,17	0
2	EDO	D	418	4/4	0.97	0.06	4,5,5,10	0
2	EDO	D	419	4/4	0.97	0.06	7,7,10,11	0
2	EDO	A	411	4/4	0.97	0.05	16,19,23,27	0
2	EDO	B	439	4/4	0.97	0.05	42,42,42,42	0
2	EDO	F	431	4/4	0.97	0.04	14,14,14,14	0
2	EDO	F	432	4/4	0.97	0.09	15,15,15,15	0
2	EDO	A	424	4/4	0.97	0.06	15,15,15,15	0
2	EDO	F	435	4/4	0.97	0.18	31,31,31,31	0
2	EDO	B	442	4/4	0.97	0.05	32,32,32,32	0
2	EDO	F	437	4/4	0.97	0.07	21,21,21,21	0
2	EDO	F	438	4/4	0.97	0.07	48,48,48,48	0
2	EDO	A	404	4/4	0.97	0.06	10,22,25,33	0
2	EDO	D	427	4/4	0.97	0.09	14,14,14,14	0
2	EDO	A	413	4/4	0.97	0.10	10,11,14,16	0
2	EDO	D	430	4/4	0.97	0.23	36,36,36,36	0
2	EDO	D	431	4/4	0.97	0.12	39,39,39,39	0
2	EDO	B	402	4/4	0.97	0.07	19,24,28,28	0
2	EDO	A	406	4/4	0.97	0.06	6,9,12,13	0
2	EDO	E	414	4/4	0.98	0.05	21,24,28,32	0
2	EDO	E	415	4/4	0.98	0.05	15,20,22,24	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	EDO	E	416	4/4	0.98	0.10	11,12,15,17	0
2	EDO	E	417	4/4	0.98	0.05	39,41,47,50	0
2	EDO	E	418	4/4	0.98	0.03	4,4,4,9	0
2	EDO	B	421	4/4	0.98	0.07	15,15,15,15	0
2	EDO	E	420	4/4	0.98	0.05	22,23,26,31	0
2	EDO	B	445	4/4	0.98	0.07	38,38,38,38	0
2	EDO	E	422	4/4	0.98	0.07	12,17,33,40	0
2	EDO	A	402	4/4	0.98	0.07	13,16,17,34	0
2	EDO	B	406	4/4	0.98	0.06	7,20,23,31	0
2	EDO	C	401	4/4	0.98	0.04	6,6,9,11	0
2	EDO	B	425	4/4	0.98	0.04	21,21,21,21	0
2	EDO	B	407	4/4	0.98	0.06	9,11,14,15	0
2	EDO	C	404	4/4	0.98	0.07	12,16,17,34	0
2	EDO	E	430	4/4	0.98	0.08	70,70,70,70	0
2	EDO	C	405	4/4	0.98	0.10	18,20,22,26	0
2	EDO	D	410	4/4	0.98	0.11	7,8,15,18	0
2	EDO	B	427	4/4	0.98	0.07	74,74,74,74	0
2	EDO	D	412	4/4	0.98	0.05	8,9,11,12	0
2	EDO	E	435	4/4	0.98	0.05	10,24,31,37	0
2	EDO	D	413	4/4	0.98	0.03	17,23,32,32	0
2	EDO	C	407	4/4	0.98	0.07	14,15,19,20	0
2	EDO	D	416	4/4	0.98	0.10	13,14,17,19	0
2	EDO	C	408	4/4	0.98	0.07	8,11,13,15	0
2	EDO	E	442	4/4	0.98	0.04	33,33,33,33	0
2	EDO	C	409	4/4	0.98	0.09	7,8,15,18	0
2	EDO	E	444	4/4	0.98	0.05	41,41,41,41	0
2	EDO	B	408	4/4	0.98	0.07	5,8,10,12	0
2	EDO	D	421	4/4	0.98	0.07	13,18,34,41	0
2	EDO	C	411	4/4	0.98	0.04	3,5,8,9	0
2	EDO	F	402	4/4	0.98	0.04	27,27,30,35	0
2	EDO	C	412	4/4	0.98	0.04	19,25,34,34	0
2	EDO	B	409	4/4	0.98	0.10	5,6,13,16	0
2	EDO	F	405	4/4	0.98	0.04	35,35,35,35	0
2	EDO	C	415	4/4	0.98	0.11	9,10,13,15	0
2	EDO	C	417	4/4	0.98	0.04	4,4,4,9	0
2	EDO	D	428	4/4	0.98	0.06	74,74,74,74	0
2	EDO	A	430	4/4	0.98	0.05	27,27,30,35	0
2	EDO	C	419	4/4	0.98	0.06	11,16,32,39	0
2	EDO	F	412	4/4	0.98	0.06	6,19,22,30	0
2	EDO	F	413	4/4	0.98	0.06	18,19,23,24	0
2	EDO	F	415	4/4	0.98	0.11	5,6,14,16	0
2	EDO	A	435	4/4	0.98	0.04	2,2,5,7	0

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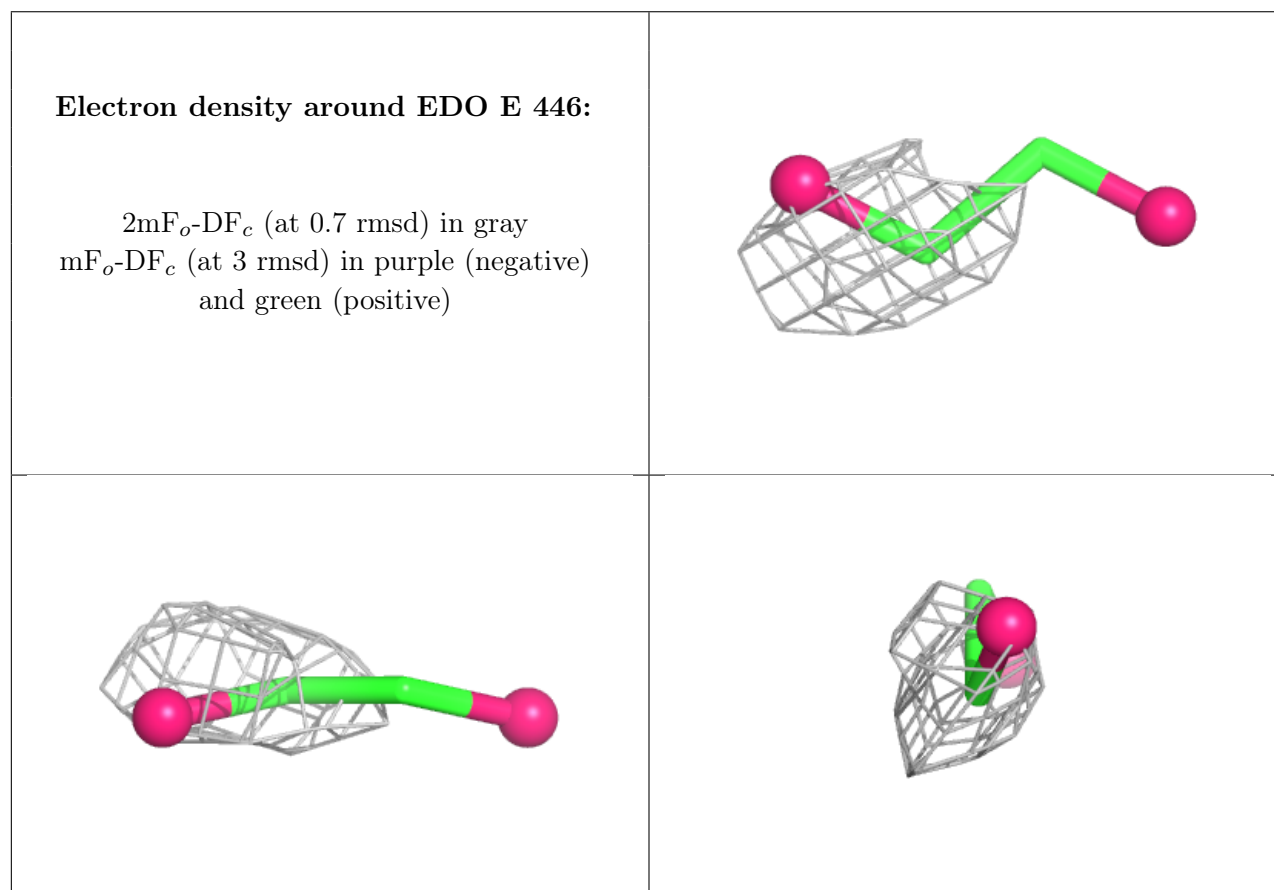
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	EDO	B	432	4/4	0.98	0.04	25,25,28,33	0
2	EDO	F	419	4/4	0.98	0.06	19,22,26,30	0
2	EDO	C	423	4/4	0.98	0.04	62,62,62,62	0
2	EDO	F	421	4/4	0.98	0.11	14,15,18,20	0
2	EDO	F	423	4/4	0.98	0.03	2,2,2,7	0
2	EDO	B	433	4/4	0.98	0.03	34,34,34,34	0
2	EDO	B	413	4/4	0.98	0.07	17,19,24,28	0
2	EDO	F	426	4/4	0.98	0.07	12,17,33,40	0
2	EDO	C	426	4/4	0.98	0.07	66,66,66,66	0
2	EDO	D	437	4/4	0.98	0.08	38,38,38,38	0
2	EDO	D	438	4/4	0.98	0.04	32,32,32,32	0
2	EDO	A	405	4/4	0.98	0.07	16,17,21,22	0
2	EDO	B	436	4/4	0.98	0.04	20,20,20,20	0
2	EDO	F	433	4/4	0.98	0.08	65,65,65,65	0
2	EDO	B	415	4/4	0.98	0.10	9,10,13,15	0
2	EDO	B	416	4/4	0.98	0.07	43,44,50,53	0
2	EDO	B	401	4/4	0.98	0.05	7,7,10,12	0
2	EDO	E	405	4/4	0.98	0.10	18,20,22,26	0
2	EDO	B	440	4/4	0.98	0.07	34,34,34,34	0
2	EDO	E	407	4/4	0.98	0.07	13,14,18,19	0
2	EDO	E	408	4/4	0.98	0.06	5,8,11,12	0
2	EDO	E	409	4/4	0.98	0.10	5,6,13,16	0
2	EDO	A	437	4/4	0.98	0.04	24,24,24,24	0
2	EDO	A	425	4/4	0.98	0.06	70,70,70,70	0
2	EDO	E	412	4/4	0.98	0.03	1,2,4,7	0
2	EDO	B	443	4/4	0.98	0.07	39,39,39,39	0
2	EDO	F	417	4/4	0.99	0.05	5,5,7,10	0
2	EDO	B	411	4/4	0.99	0.03	2,3,4,5	0
2	EDO	E	440	4/4	0.99	0.04	16,16,16,16	0
2	EDO	C	413	4/4	0.99	0.04	11,14,18,22	0
2	EDO	B	420	4/4	0.99	0.05	10,15,31,38	0
2	EDO	F	422	4/4	0.99	0.05	38,39,45,49	0
2	EDO	B	404	4/4	0.99	0.08	16,19,20,37	0
2	EDO	D	423	4/4	0.99	0.03	16,16,16,16	0
2	EDO	D	405	4/4	0.99	0.09	12,15,16,33	0
2	EDO	E	424	4/4	0.99	0.04	19,19,19,19	0
2	EDO	C	416	4/4	0.99	0.05	45,46,52,56	0
2	EDO	F	428	4/4	0.99	0.03	13,13,13,13	0
2	EDO	E	448	4/4	0.99	0.04	29,29,29,29	0
2	EDO	F	401	4/4	0.99	0.04	12,26,33,39	0
2	EDO	C	430	4/4	0.99	0.04	11,25,32,38	0
2	EDO	B	422	4/4	0.99	0.04	18,18,18,18	0

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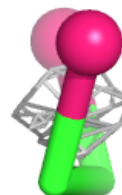
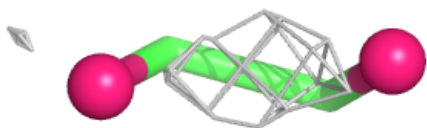
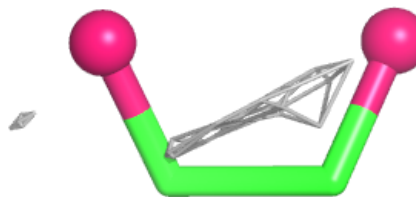
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	EDO	D	409	4/4	0.99	0.07	7,10,13,14	0
2	EDO	A	441	4/4	0.99	0.04	17,17,17,17	0
2	EDO	A	409	4/4	0.99	0.04	3,4,5,6	0
2	EDO	A	414	4/4	0.99	0.06	36,37,44,47	0
2	EDO	C	421	4/4	0.99	0.04	18,18,18,18	0
2	EDO	D	414	4/4	0.99	0.05	12,14,19,23	0
2	EDO	F	410	4/4	0.99	0.05	14,18,19,36	0
2	EDO	A	418	4/4	0.99	0.06	13,18,34,41	0
2	EDO	A	415	4/4	0.99	0.04	3,3,3,8	0
2	EDO	E	436	4/4	0.99	0.04	20,20,23,28	0
2	EDO	F	414	4/4	0.99	0.06	5,8,11,12	0
2	EDO	D	417	4/4	0.99	0.07	36,37,43,46	0
2	EDO	A	420	4/4	0.99	0.04	20,20,20,20	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

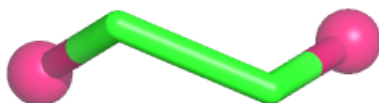
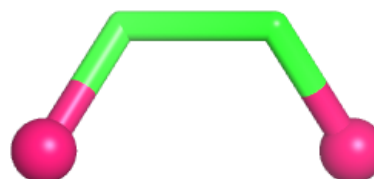


**Electron density around EDO C 427:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around EDO E 431:**

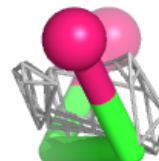
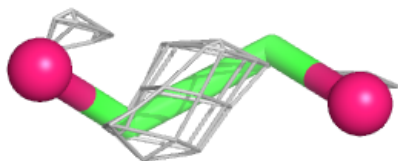
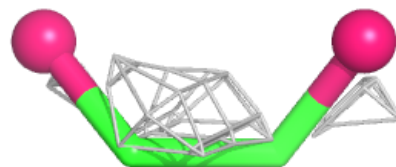
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



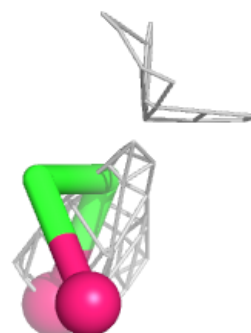
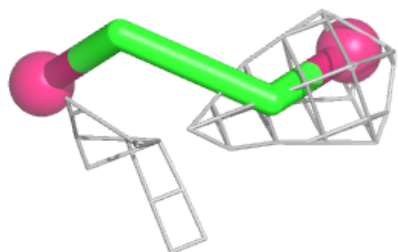
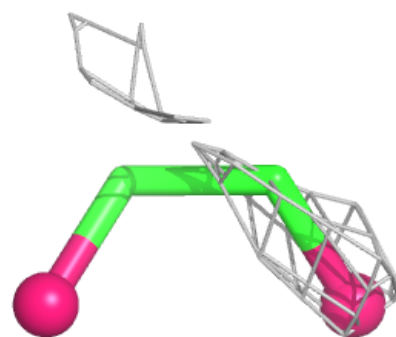


**Electron density around EDO B 428:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

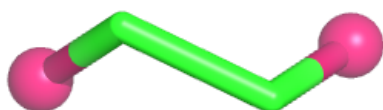
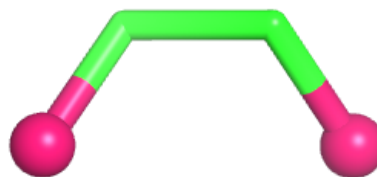
**Electron density around EDO F 444:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

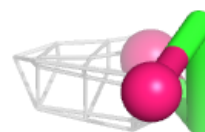
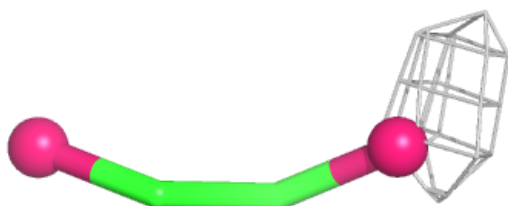


**Electron density around EDO A 426:**

$2mF_o - DF_c$  (at 0.7 rmsd) in gray  
 $mF_o - DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

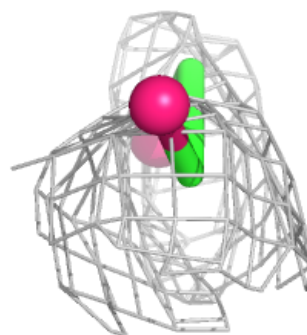
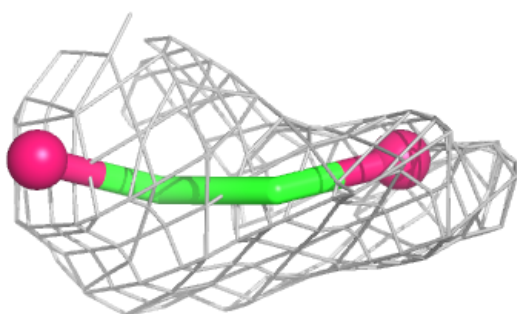
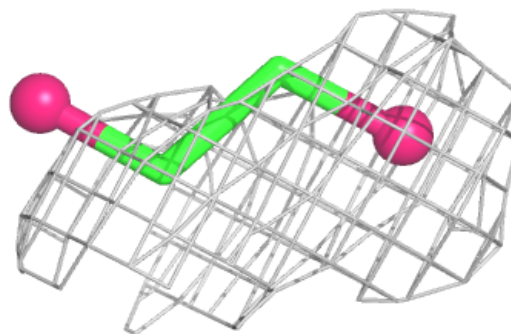
**Electron density around EDO E 445:**

$2mF_o - DF_c$  (at 0.7 rmsd) in gray  
 $mF_o - DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

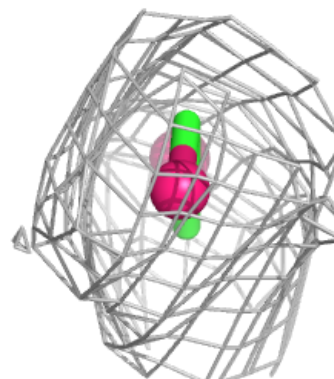
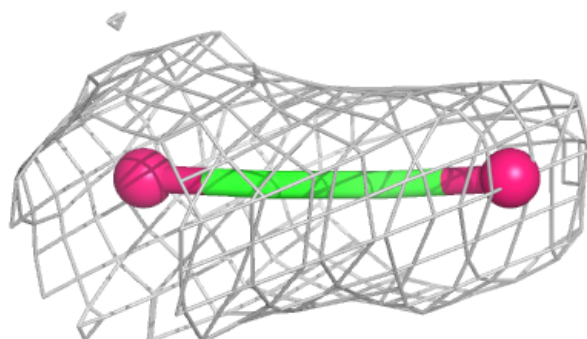
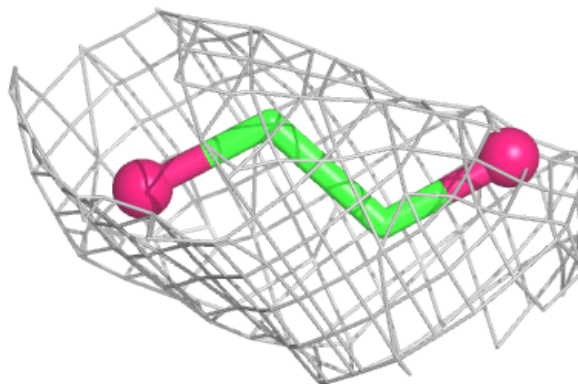


**Electron density around EDO B 441:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

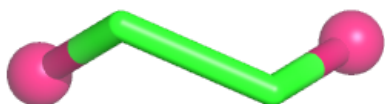
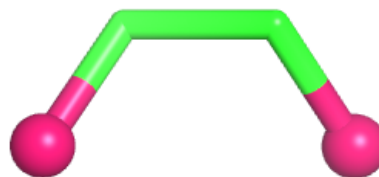
**Electron density around EDO F 418:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

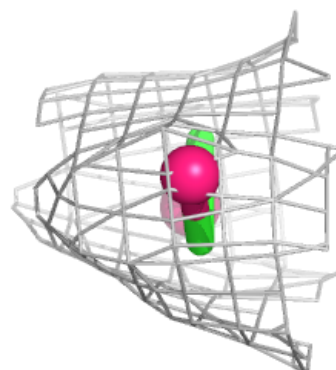
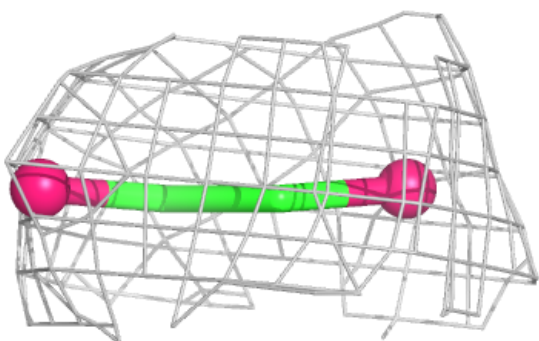
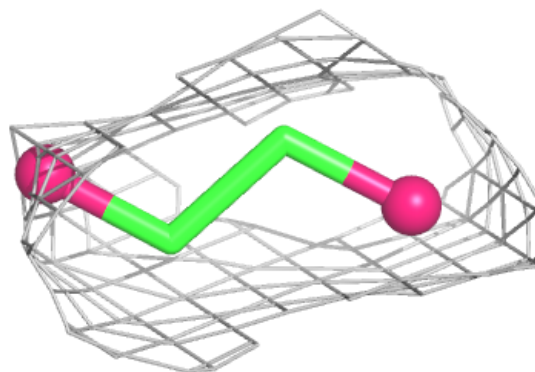


**Electron density around EDO F 434:**

$2mF_o - DF_c$  (at 0.7 rmsd) in gray  
 $mF_o - DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

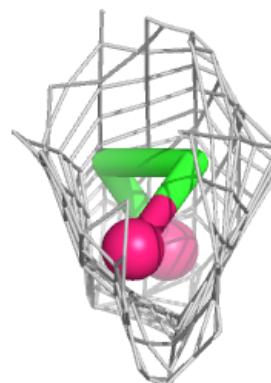
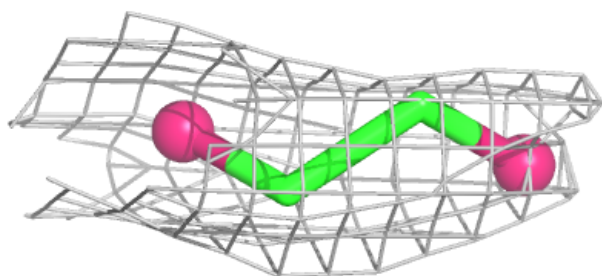
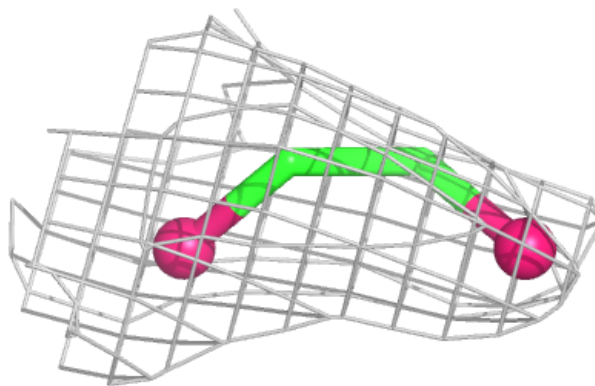
**Electron density around EDO B 414:**

$2mF_o - DF_c$  (at 0.7 rmsd) in gray  
 $mF_o - DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

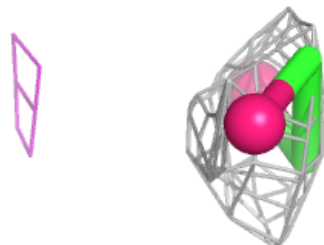
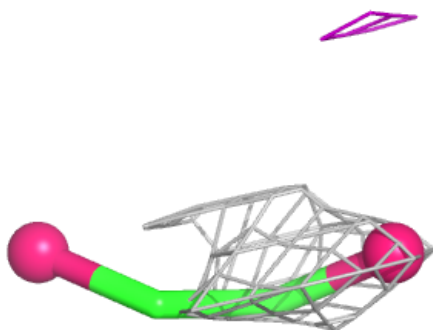
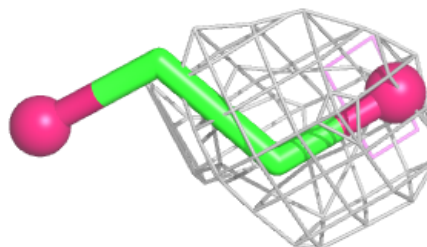


**Electron density around EDO A 443:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

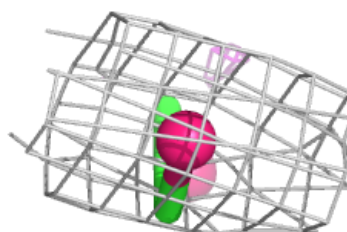
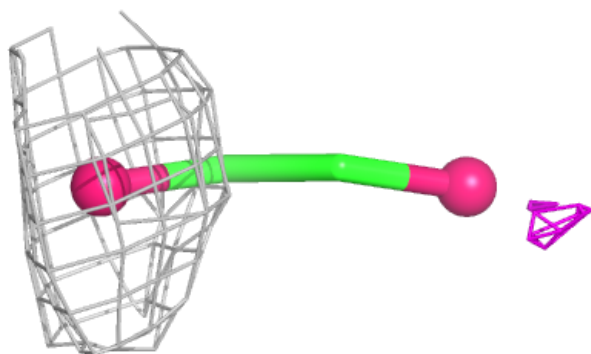
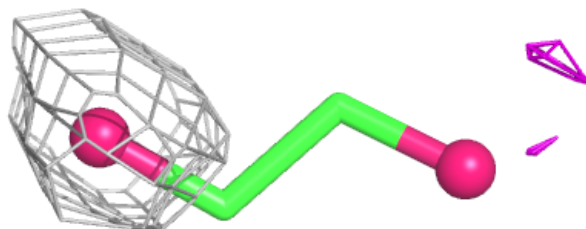
**Electron density around EDO E 432:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

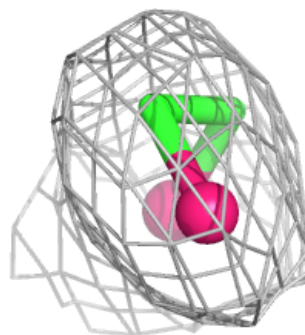
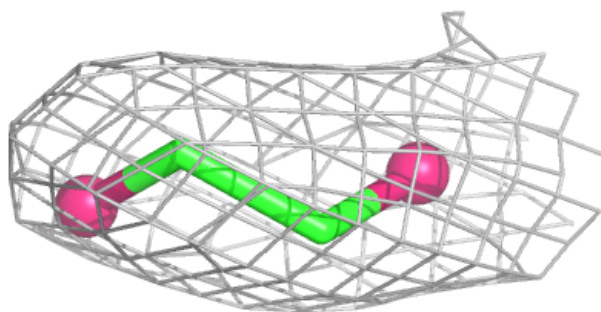
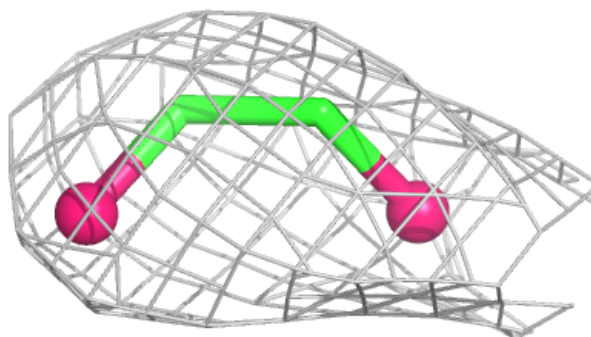


**Electron density around EDO B 403:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around EDO A 423:**

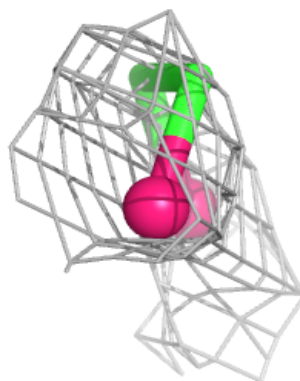
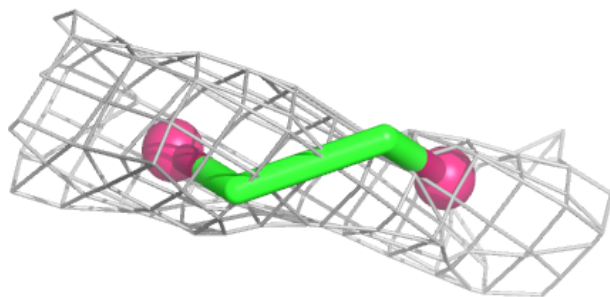
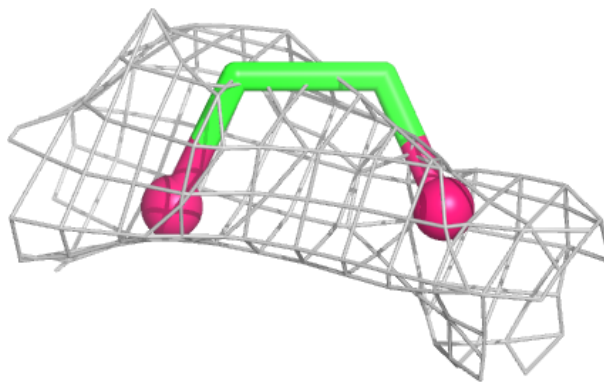
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



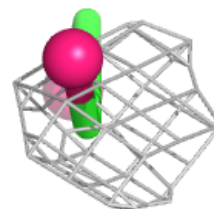
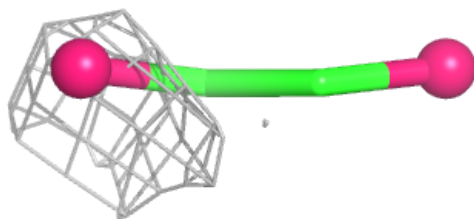
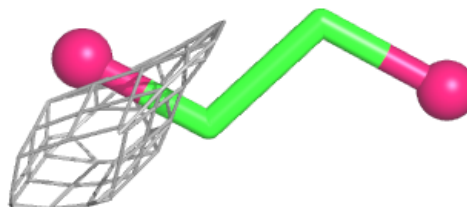


**Electron density around EDO C 436:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

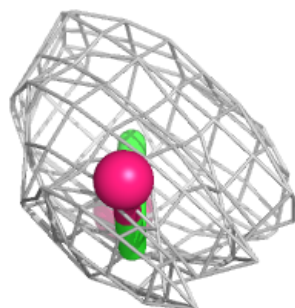
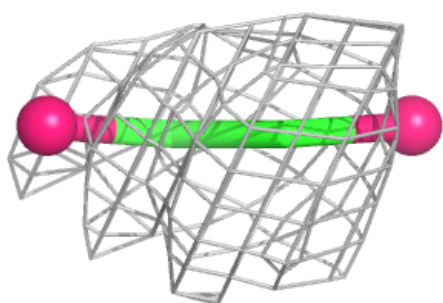
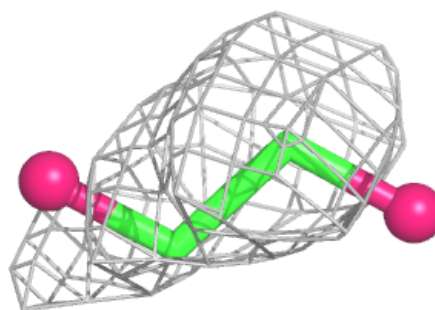
**Electron density around EDO D 425:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

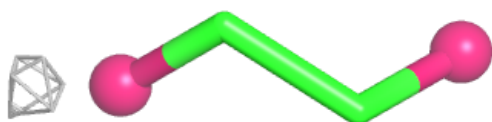


**Electron density around EDO E 426:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around EDO D 403:**

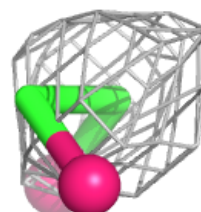
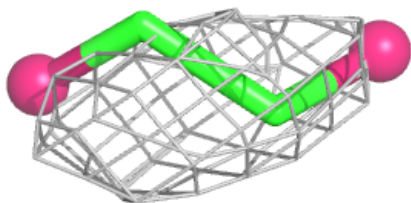
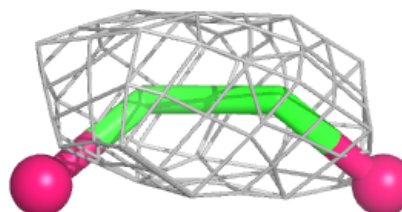
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



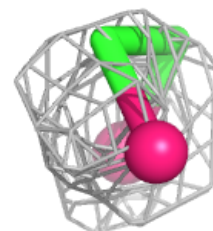
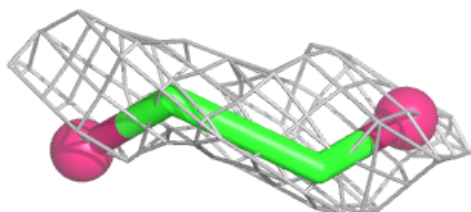
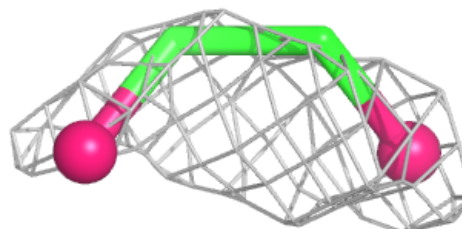


**Electron density around EDO D 420:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

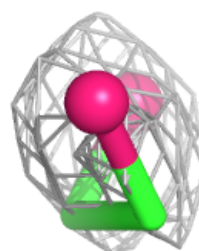
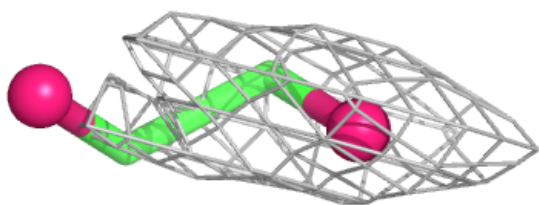
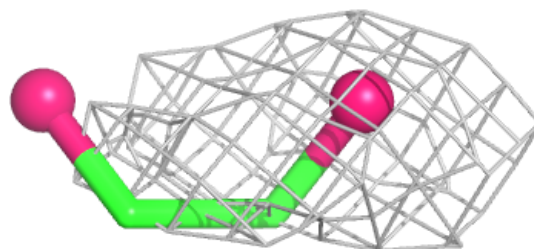
**Electron density around EDO A 428:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

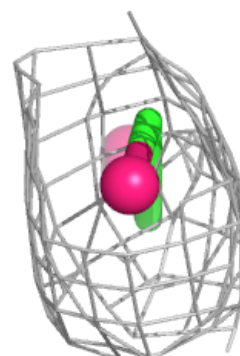
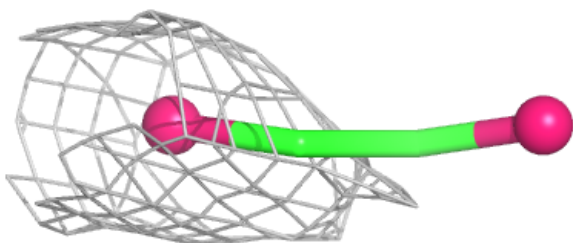
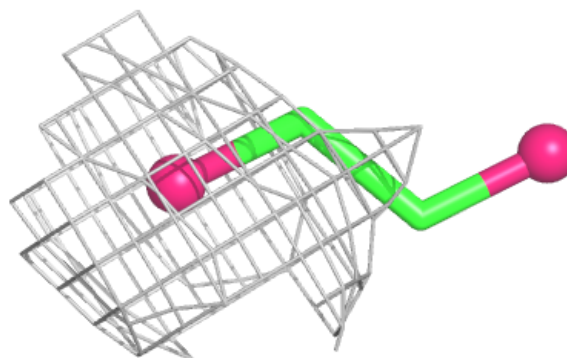


**Electron density around EDO D 429:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

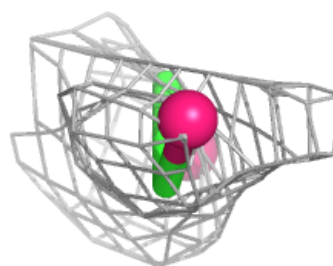
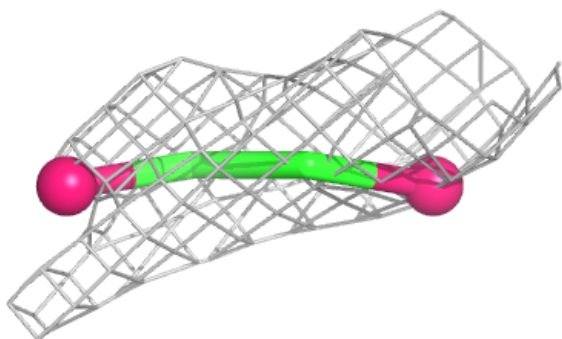
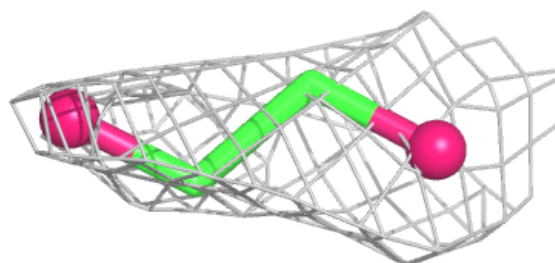
**Electron density around EDO B 434:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



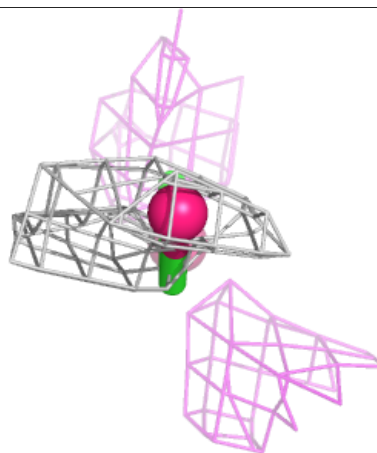
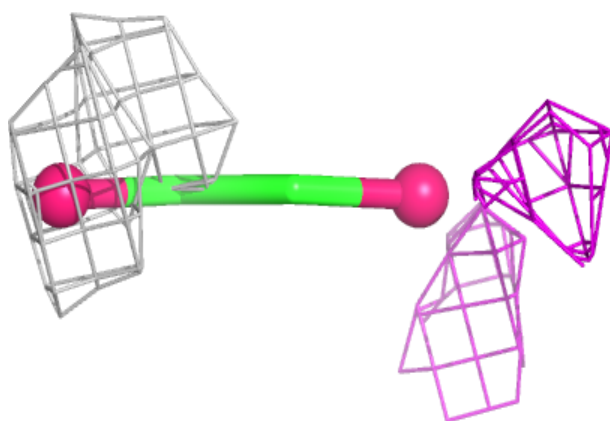
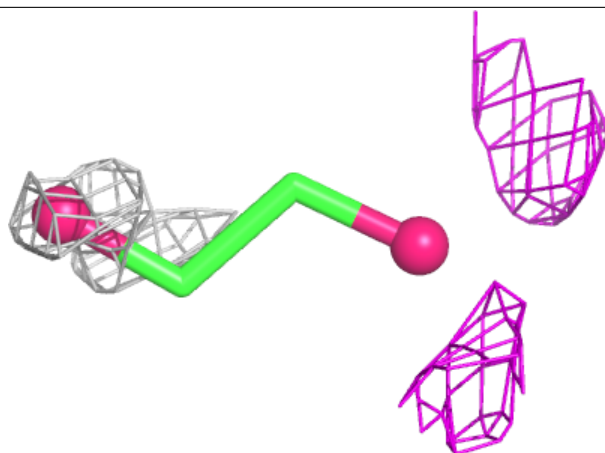
**Electron density around EDO B 437:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



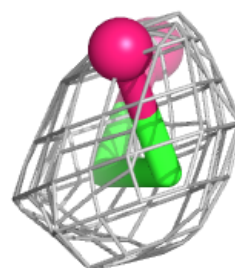
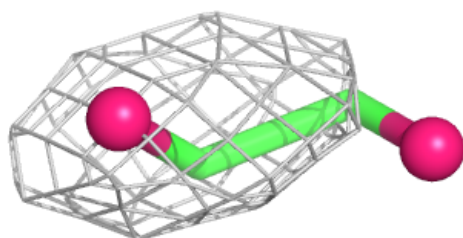
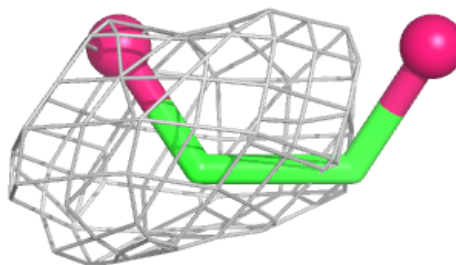
**Electron density around EDO A 401:**

$2mF_o - DF_c$  (at 0.7 rmsd) in gray  
 $mF_o - DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

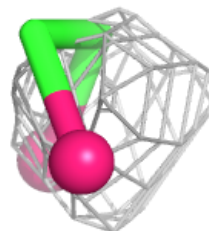
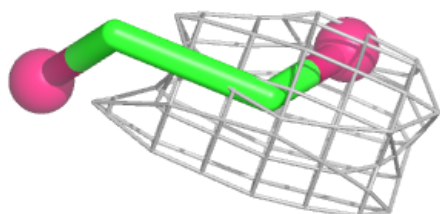
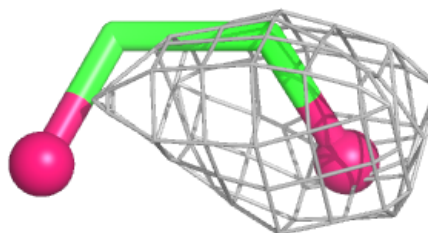


**Electron density around EDO E 433:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

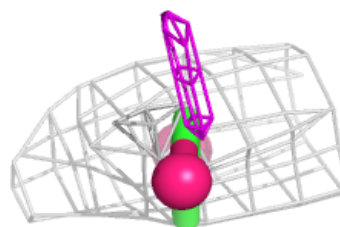
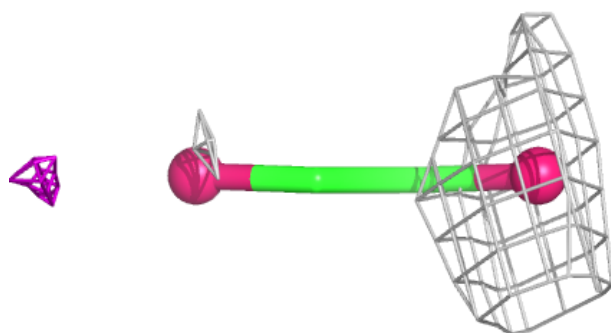
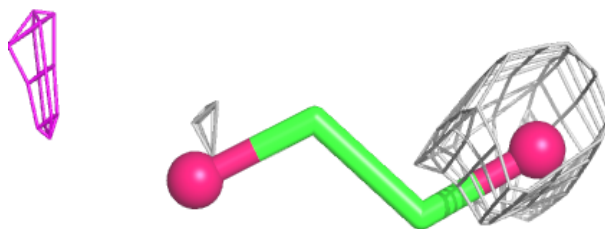
**Electron density around EDO B 446:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

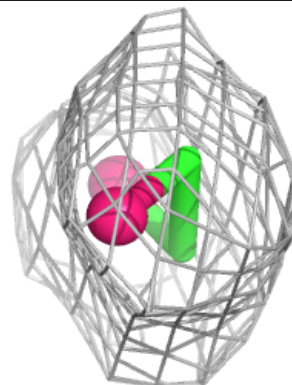
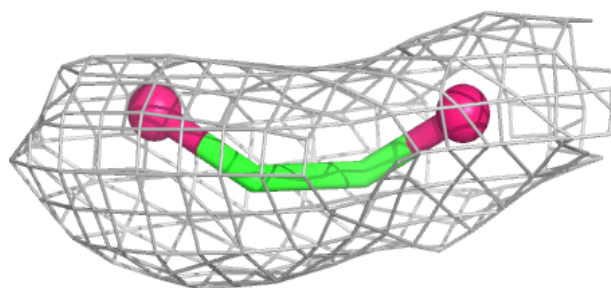
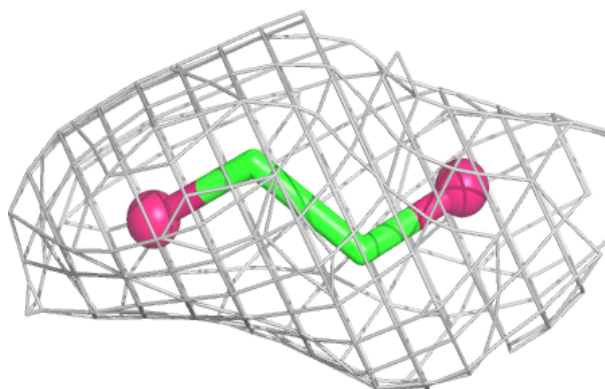


**Electron density around EDO C 403:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around EDO F 403:**

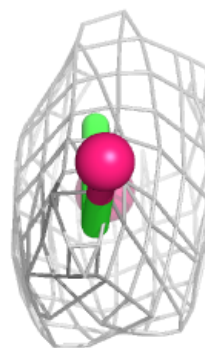
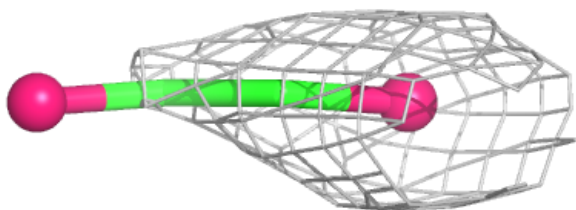
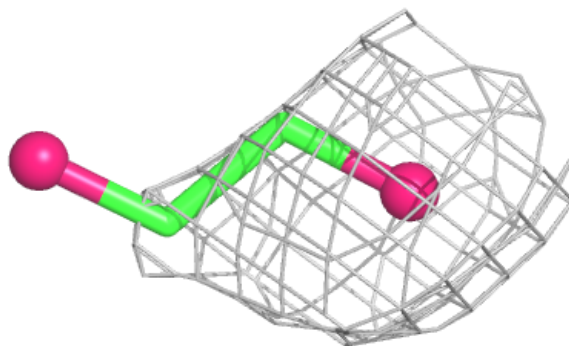
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



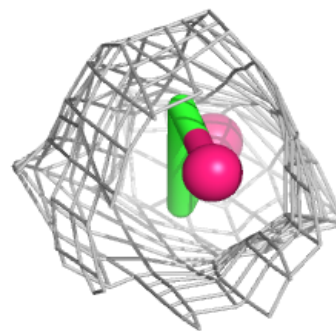
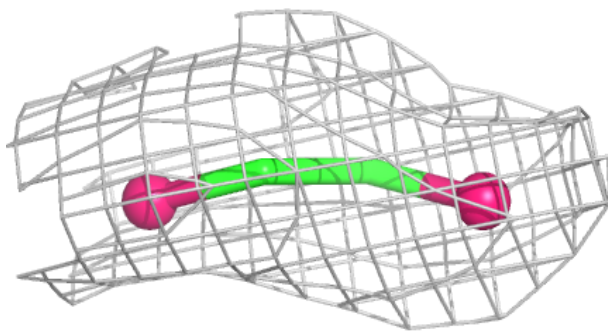
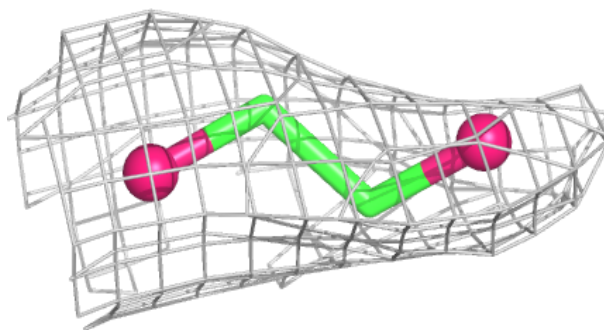


**Electron density around EDO F 404:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

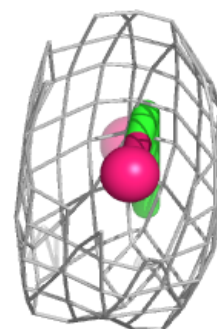
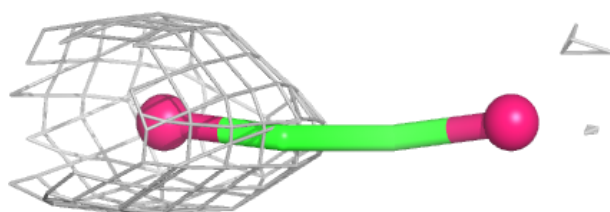
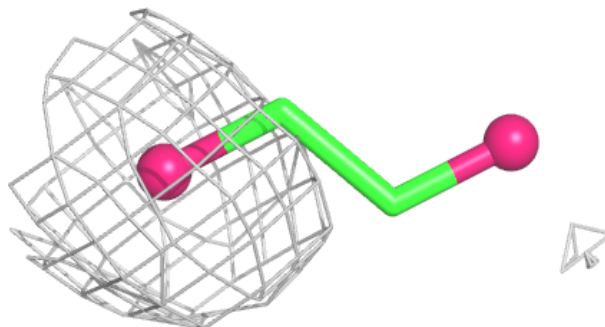
**Electron density around EDO A 416:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



**Electron density around EDO C 433:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around EDO F 443:**

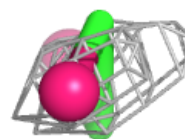
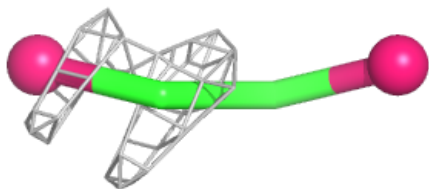
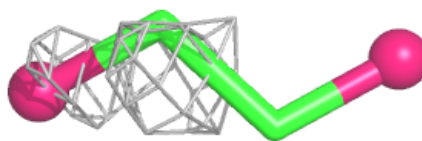
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



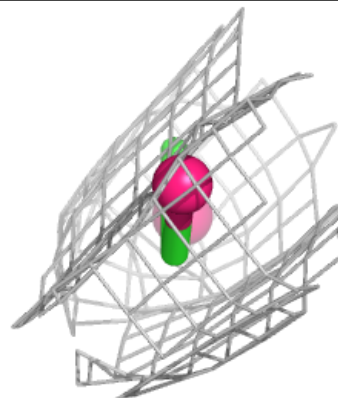
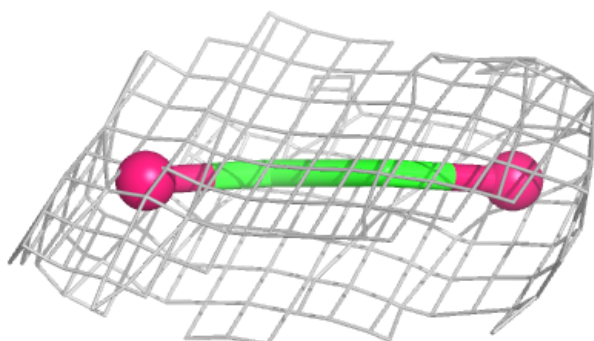
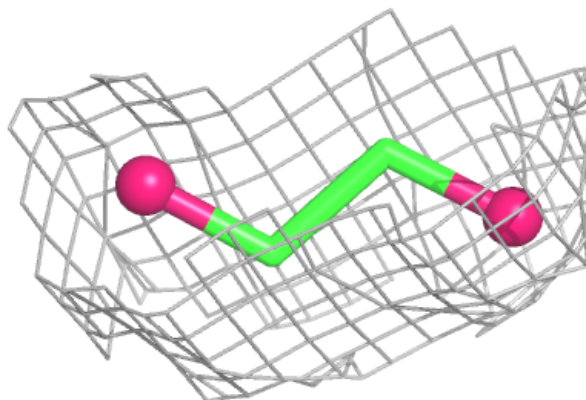


**Electron density around EDO A 427:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

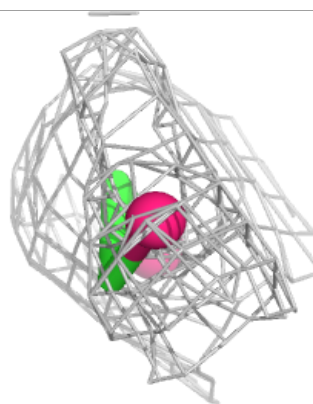
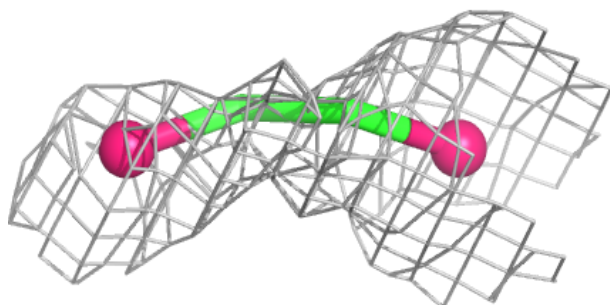
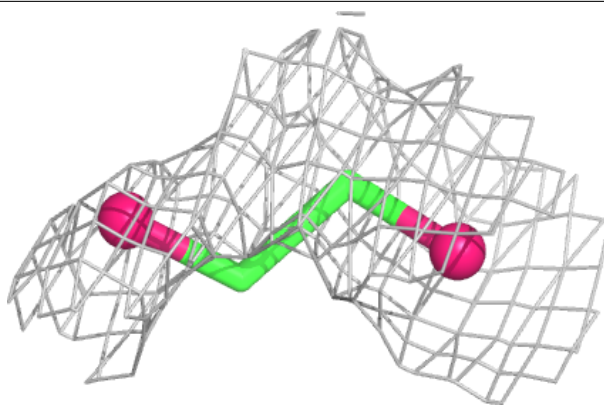
**Electron density around EDO F 445:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

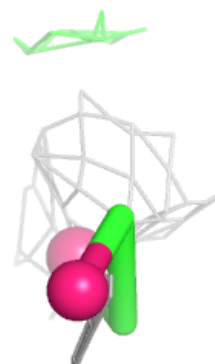
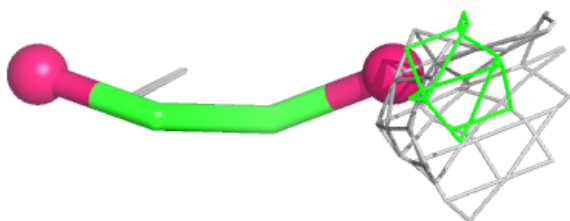
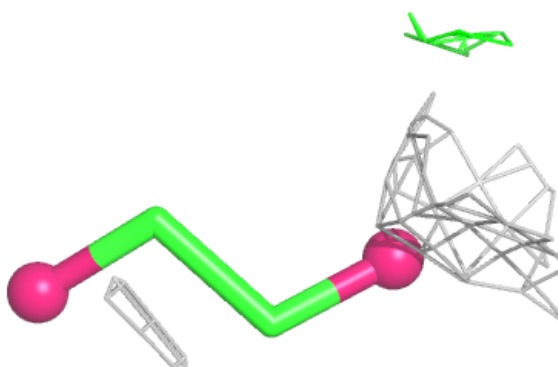


**Electron density around EDO B 426:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

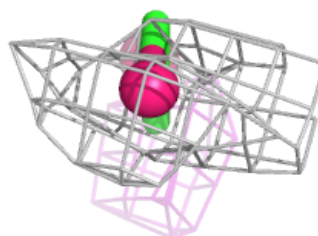
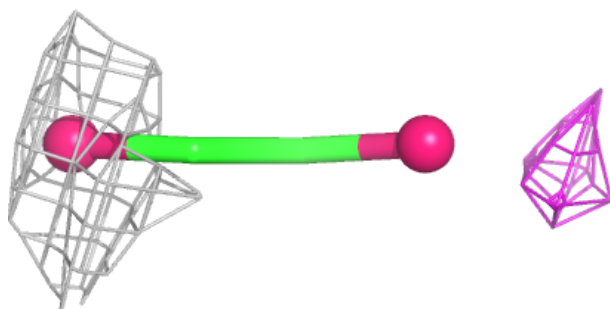
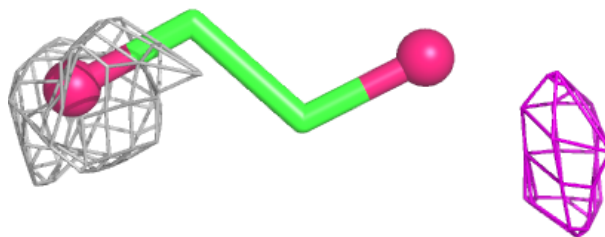
**Electron density around EDO E 441:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

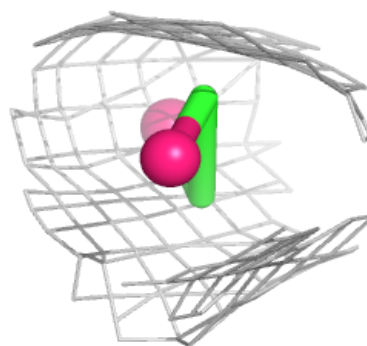
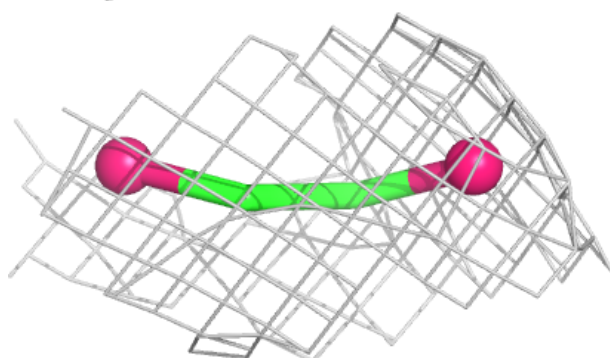
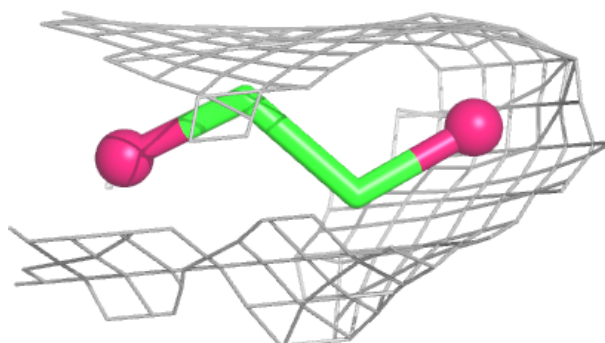


**Electron density around EDO D 404:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

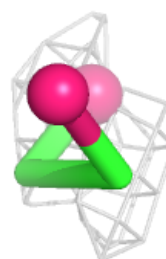
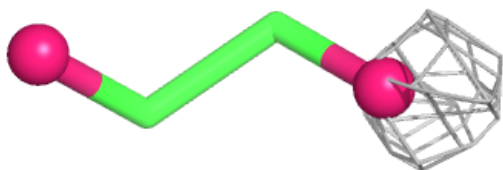
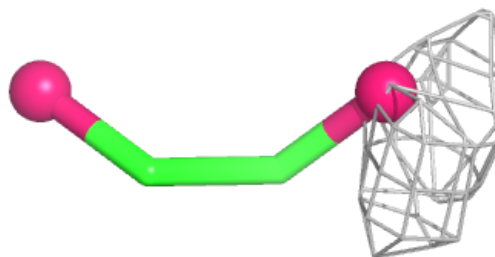
**Electron density around EDO C 418:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

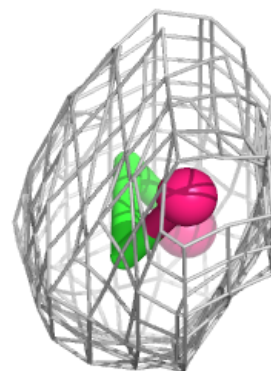
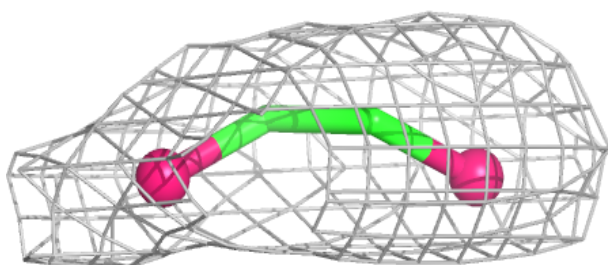
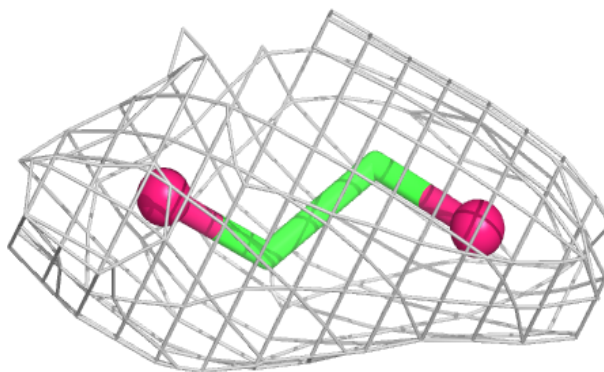


**Electron density around EDO B 438:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

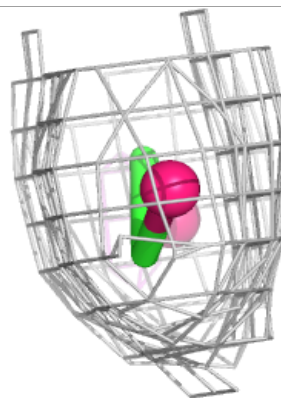
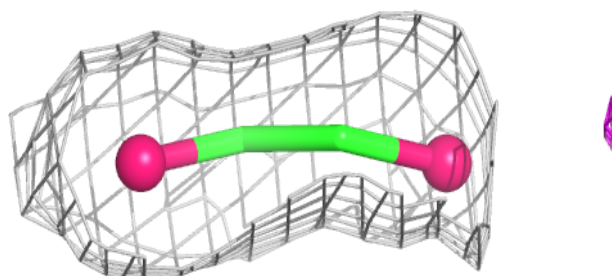
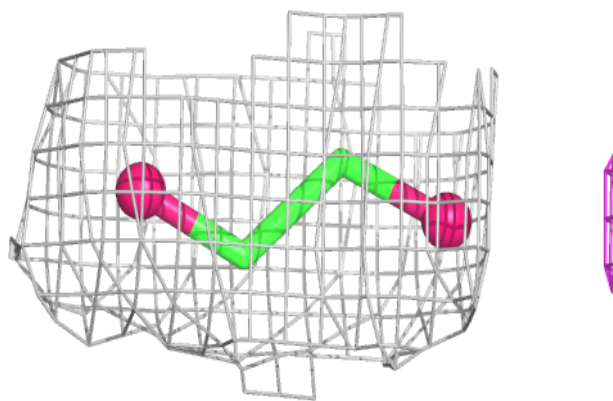
**Electron density around EDO C 432:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

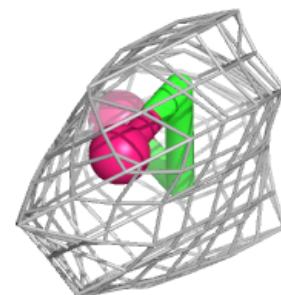
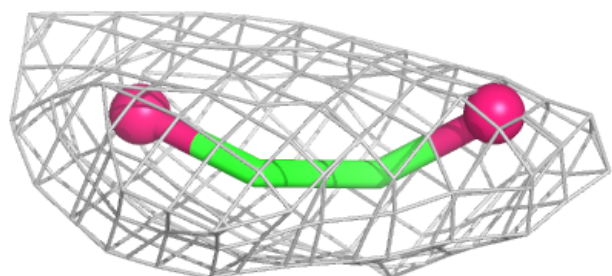
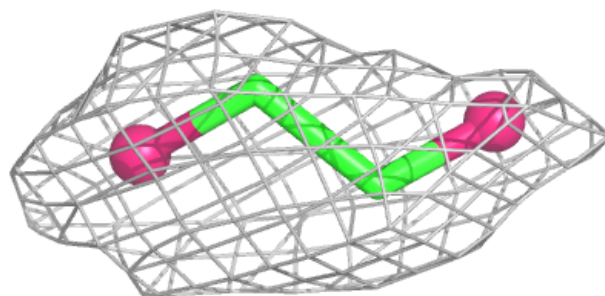


**Electron density around EDO F 406:**

$2mF_o - DF_c$  (at 0.7 rmsd) in gray  
 $mF_o - DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around EDO D 439:**

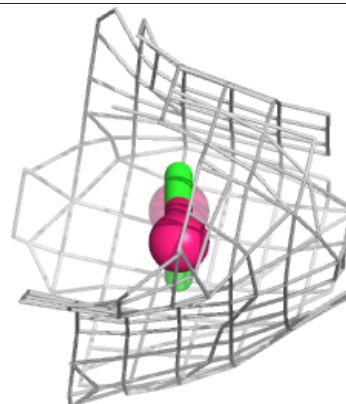
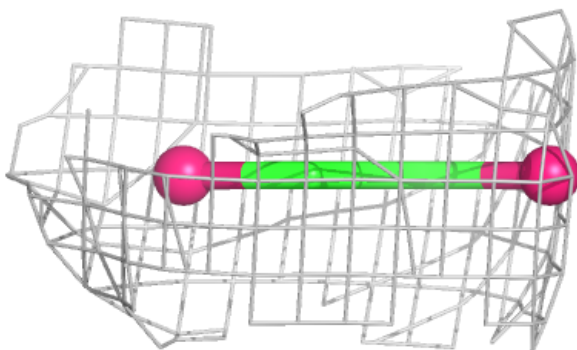
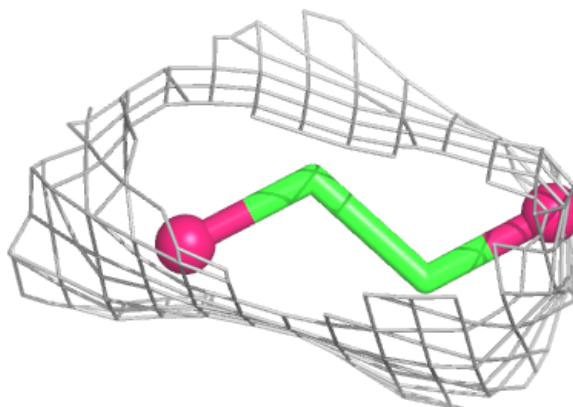
$2mF_o - DF_c$  (at 0.7 rmsd) in gray  
 $mF_o - DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



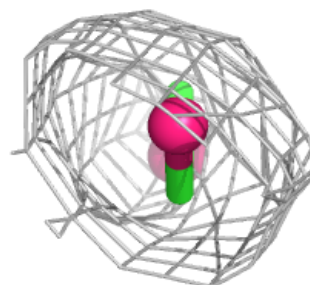
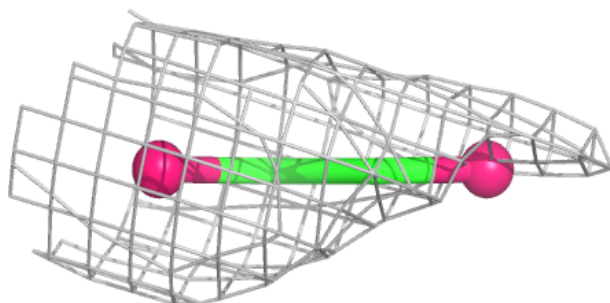
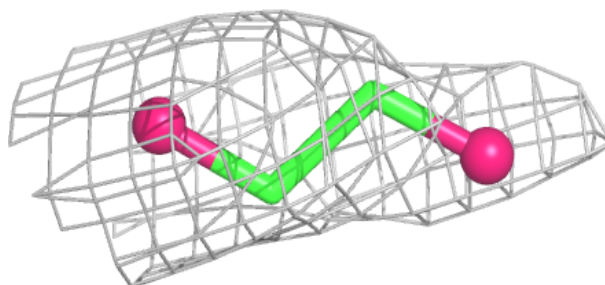


**Electron density around EDO F 420:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

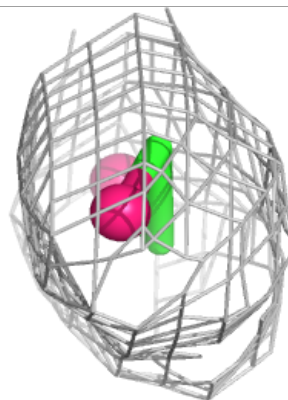
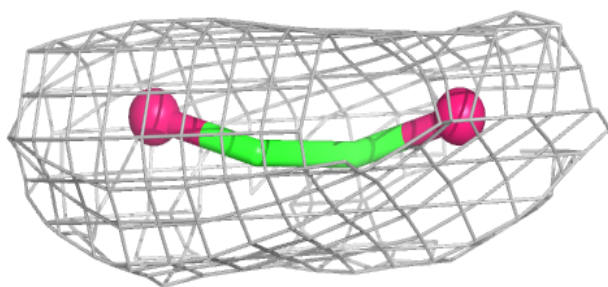
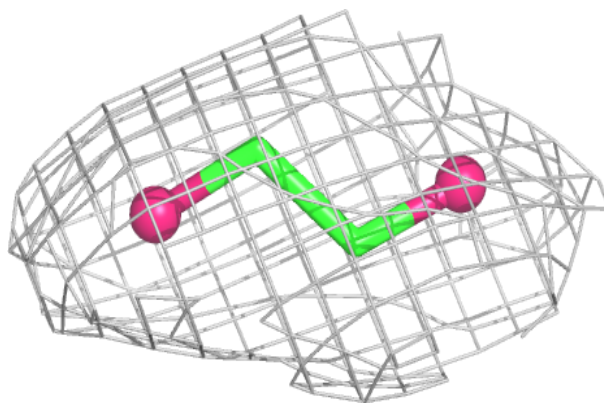
**Electron density around EDO F 424:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

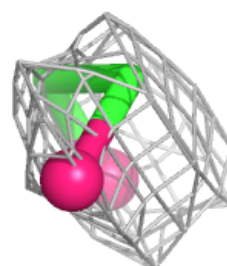
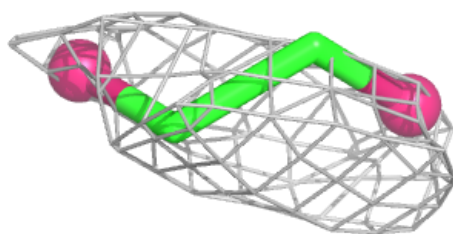
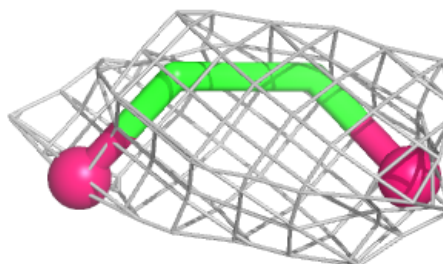


**Electron density around EDO A 431:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

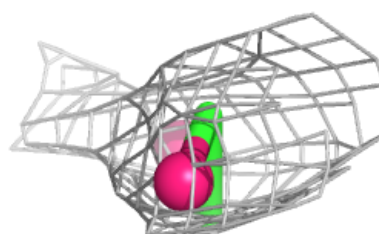
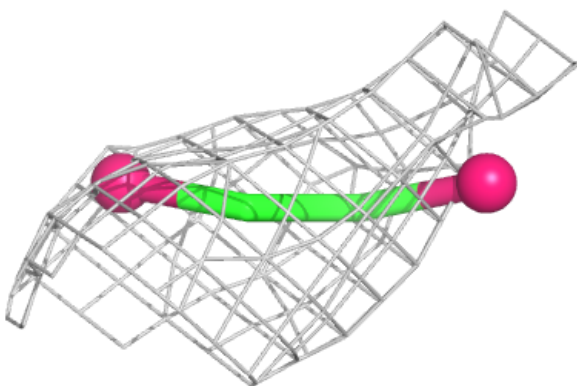
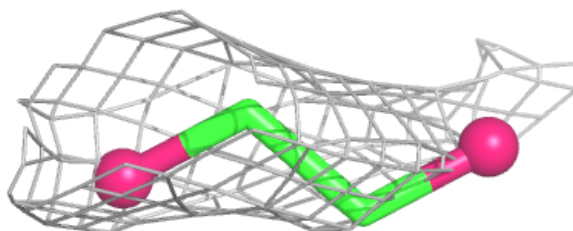
**Electron density around EDO F 436:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

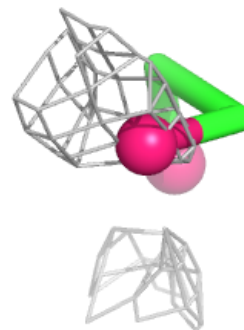
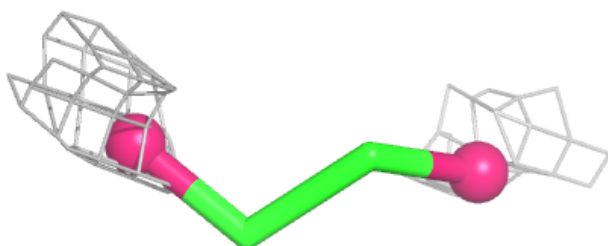
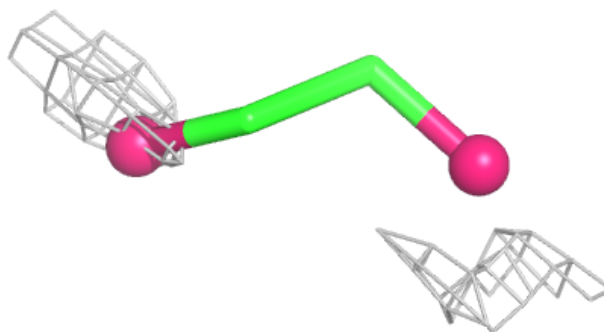


**Electron density around EDO F 440:**

$2mF_o - DF_c$  (at 0.7 rmsd) in gray  
 $mF_o - DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around EDO F 442:**

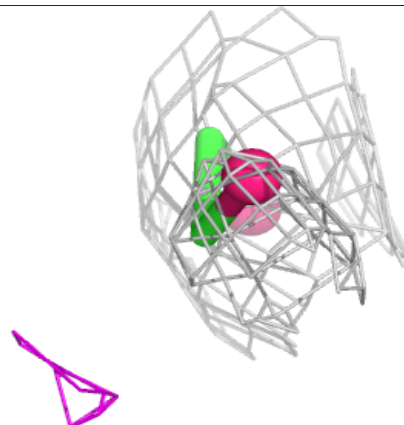
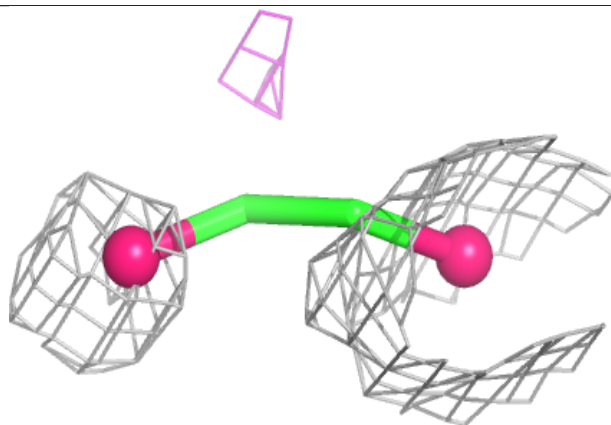
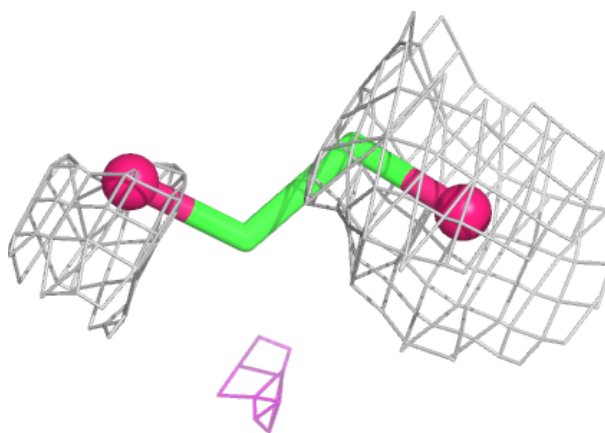
$2mF_o - DF_c$  (at 0.7 rmsd) in gray  
 $mF_o - DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



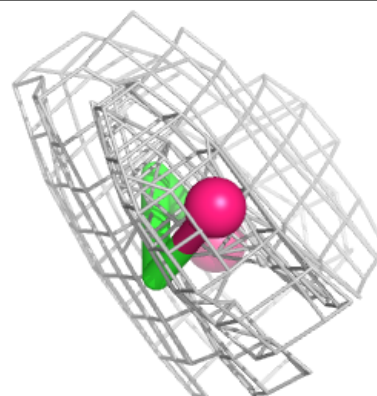
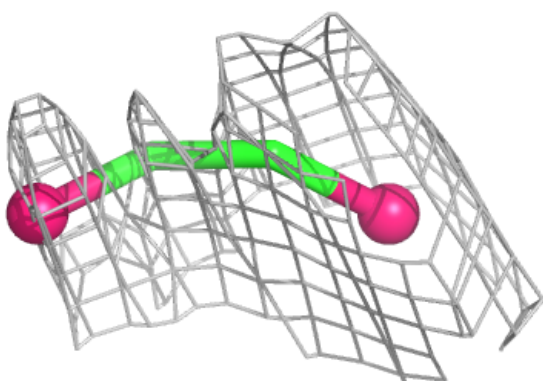
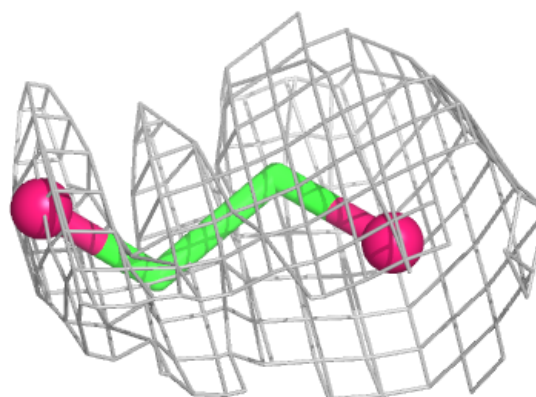


**Electron density around EDO E 429:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

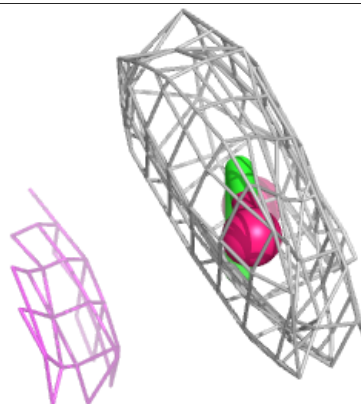
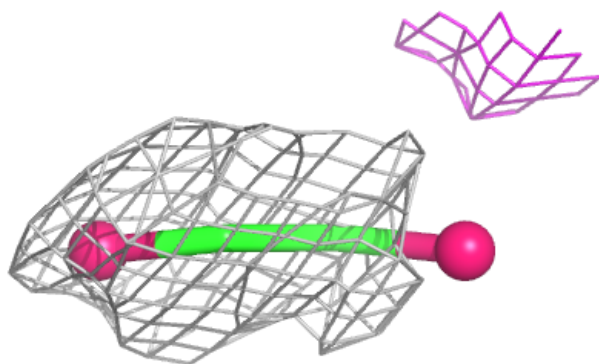
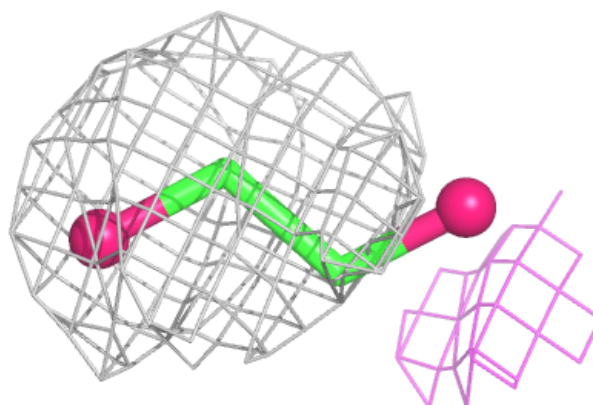
**Electron density around EDO A 421:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

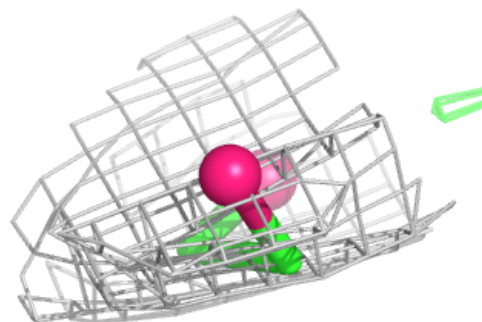
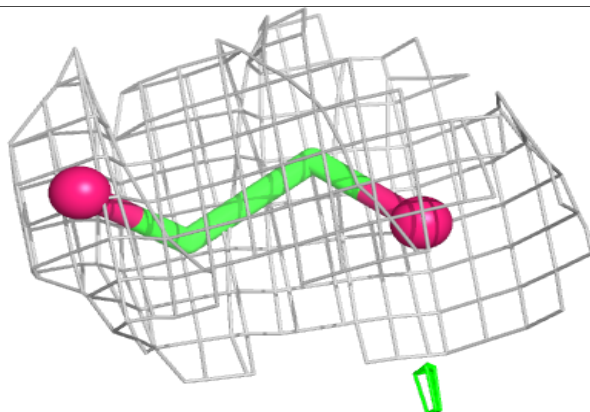
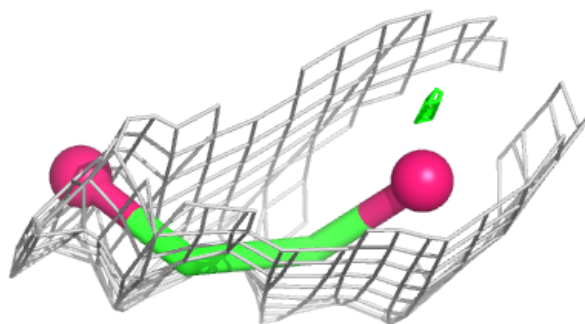


**Electron density around EDO D 402:**

$2mF_o - DF_c$  (at 0.7 rmsd) in gray  
 $mF_o - DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

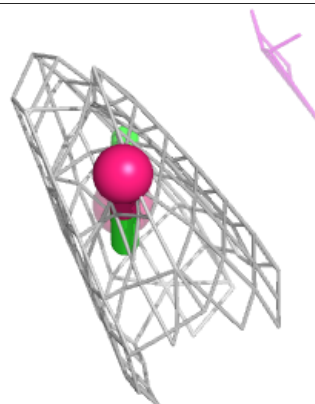
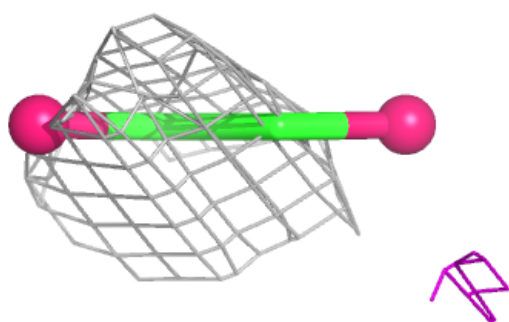
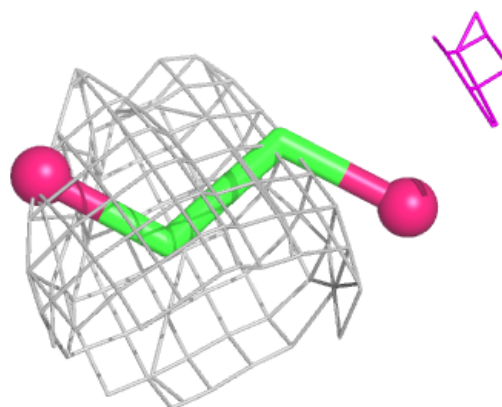
**Electron density around EDO D 424:**

$2mF_o - DF_c$  (at 0.7 rmsd) in gray  
 $mF_o - DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

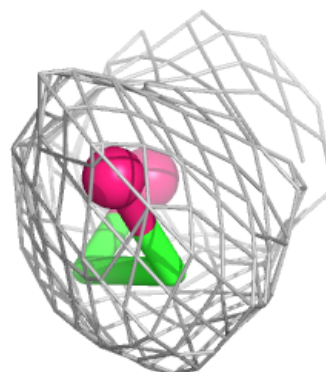
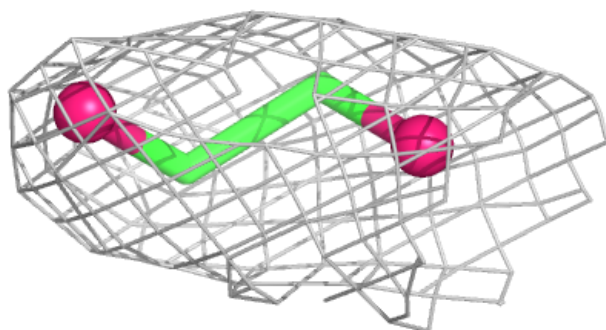
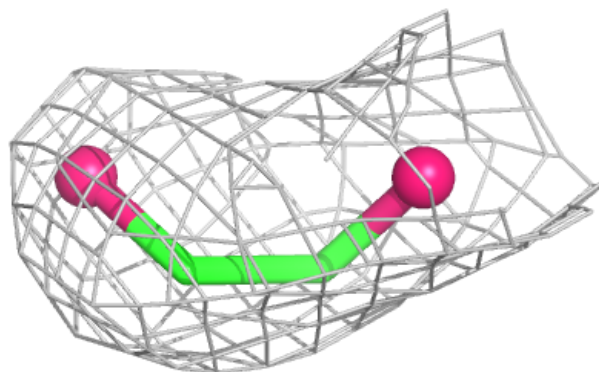


**Electron density around EDO A 440:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

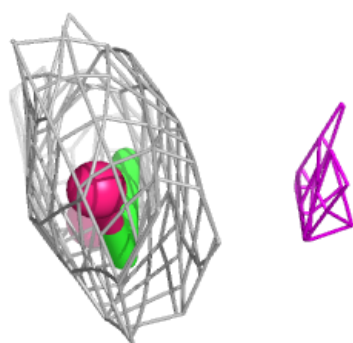
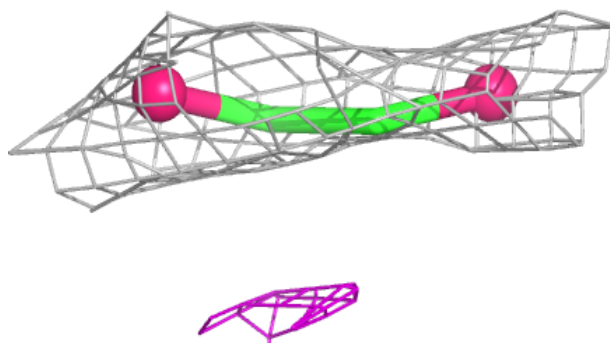
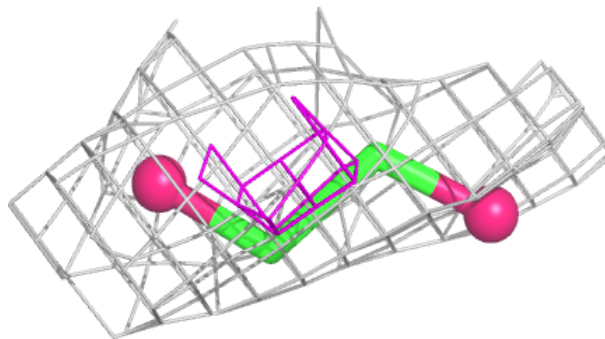
**Electron density around EDO D 426:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

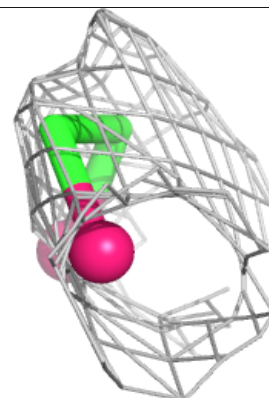
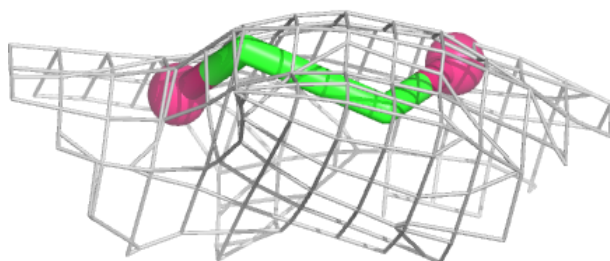


**Electron density around EDO B 444:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around EDO D 432:**

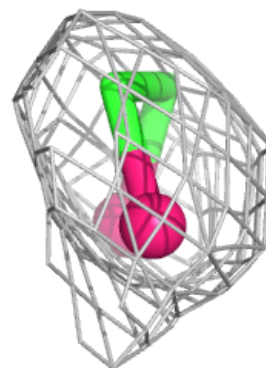
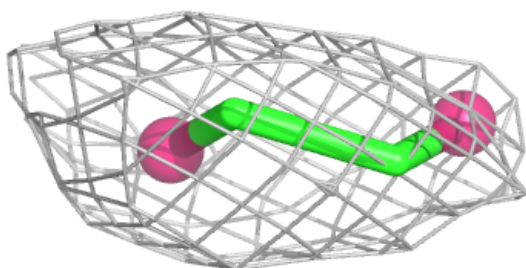
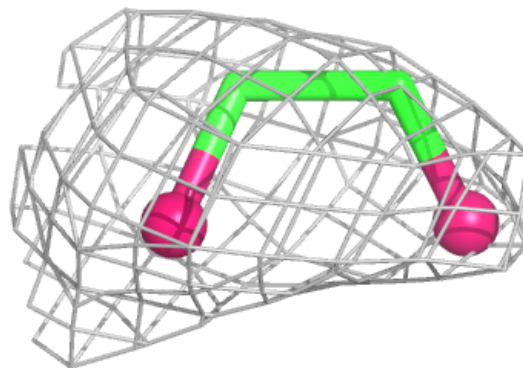
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



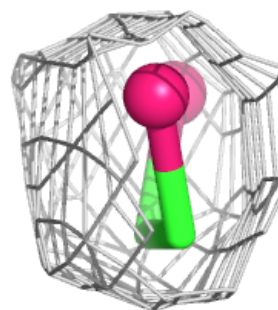
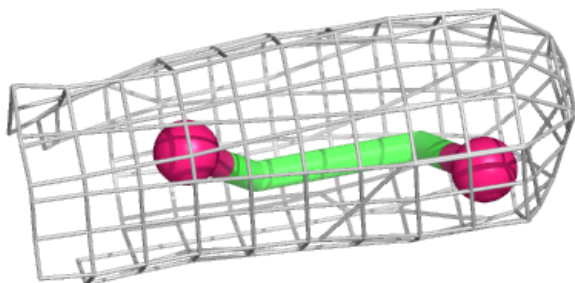
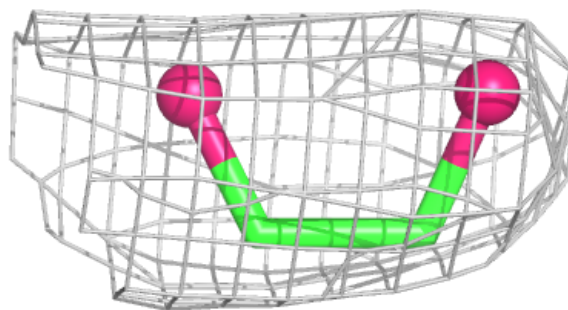


**Electron density around EDO D 433:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

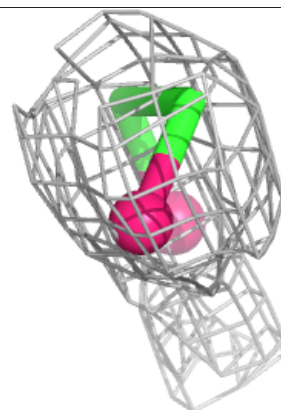
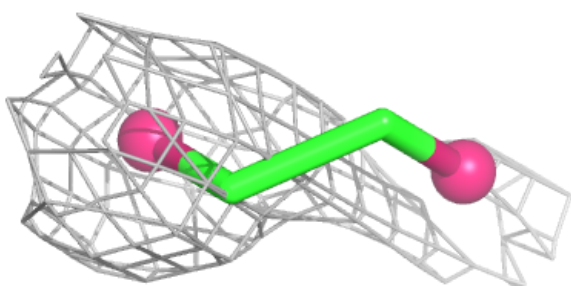
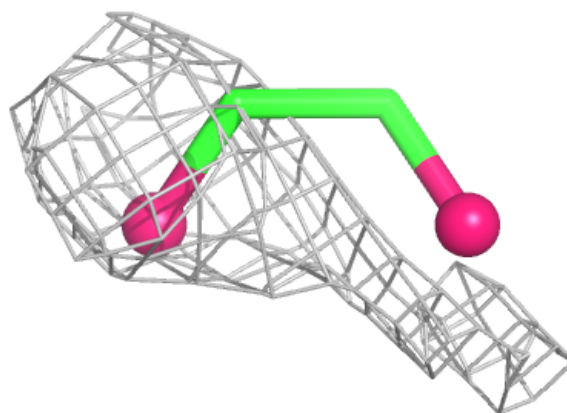
**Electron density around EDO D 435:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

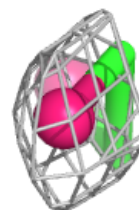
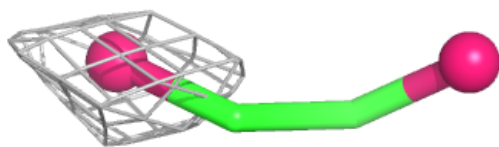
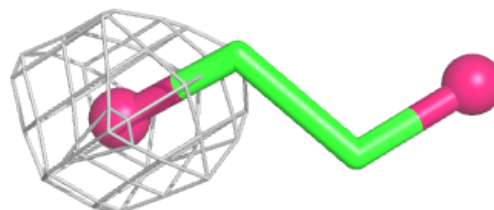


**Electron density around EDO D 436:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

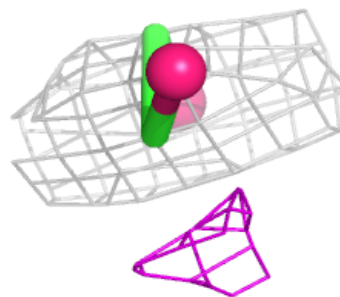
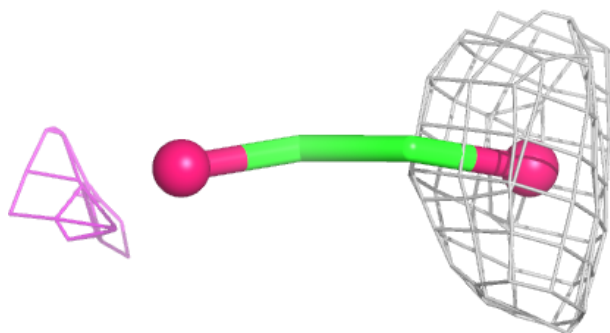
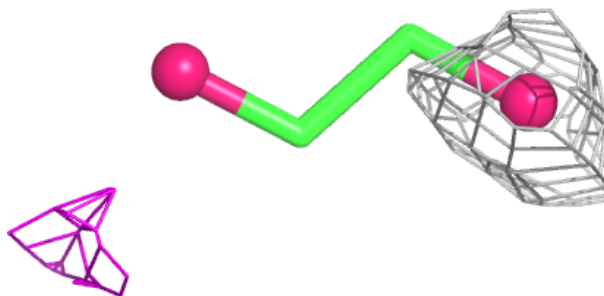
**Electron density around EDO A 442:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

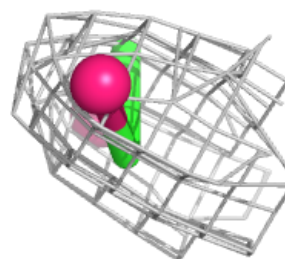
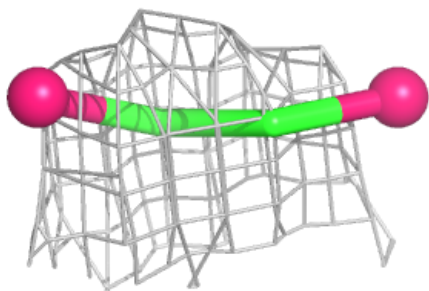
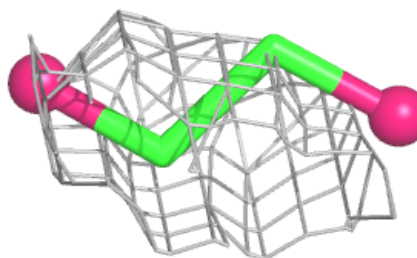


**Electron density around EDO E 403:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

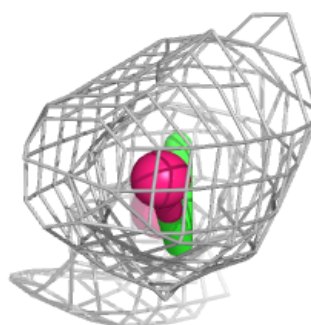
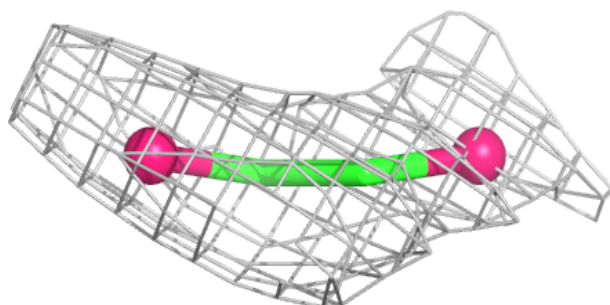
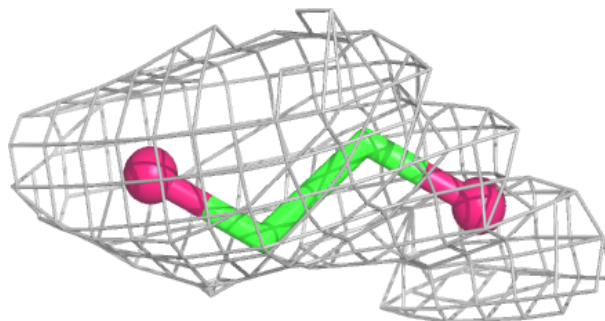
**Electron density around EDO E 404:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

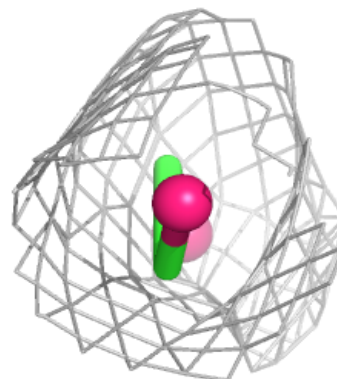
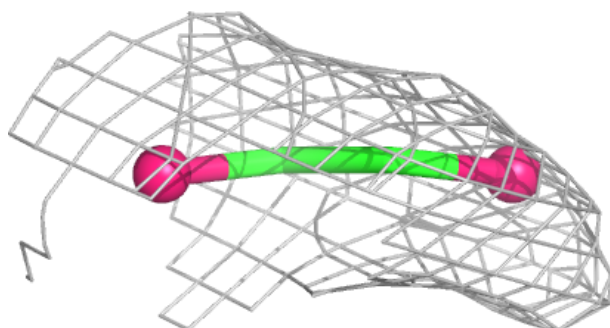
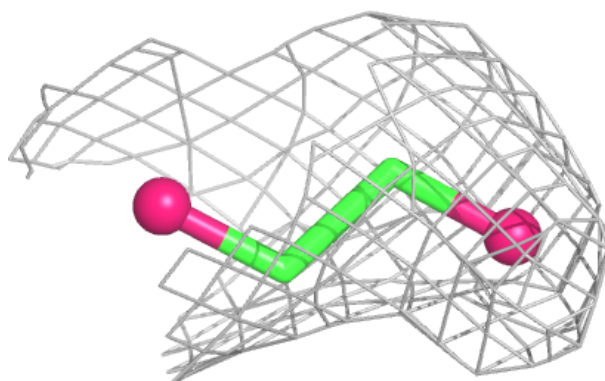


**Electron density around EDO E 411:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around EDO E 413:**

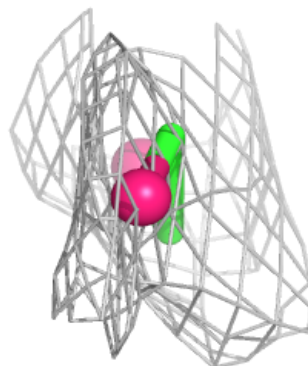
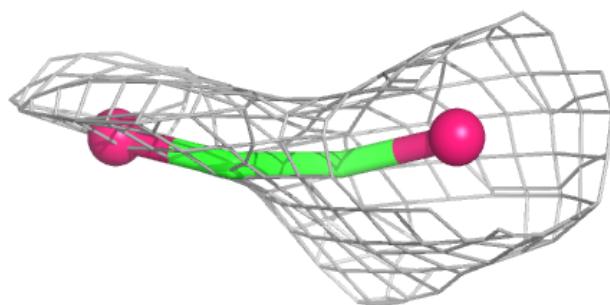
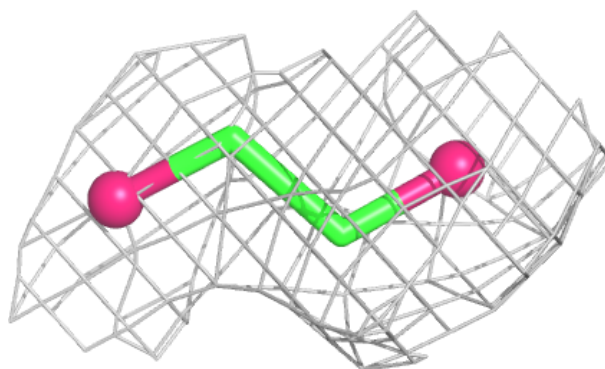
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



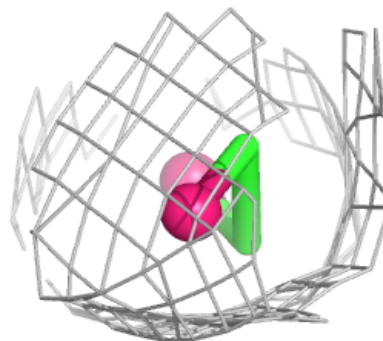
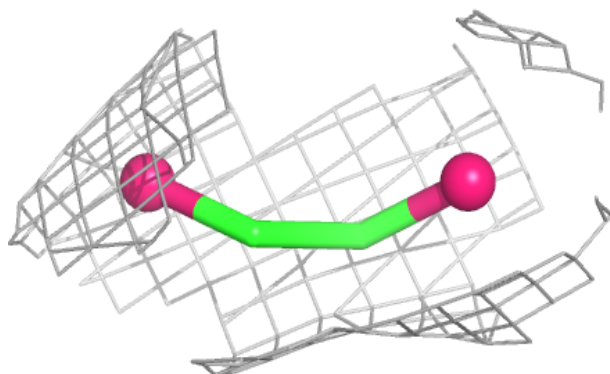
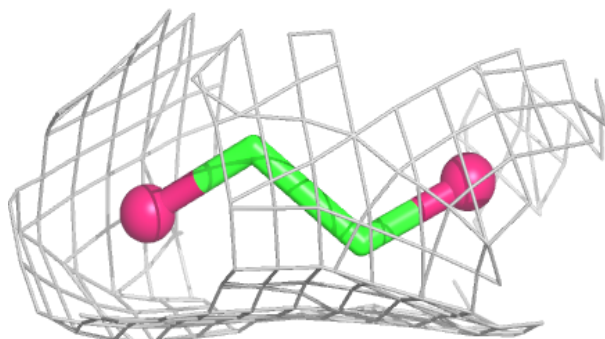


**Electron density around EDO E 419:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

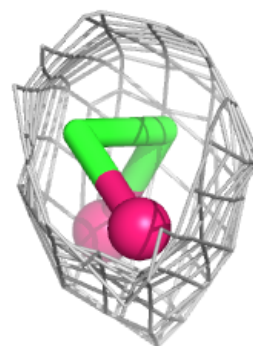
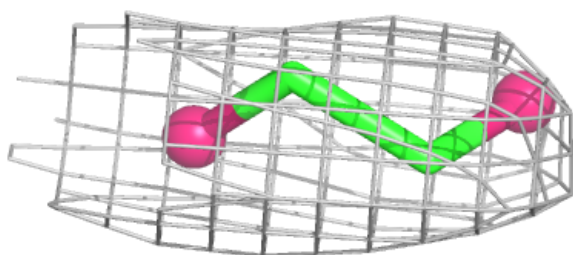
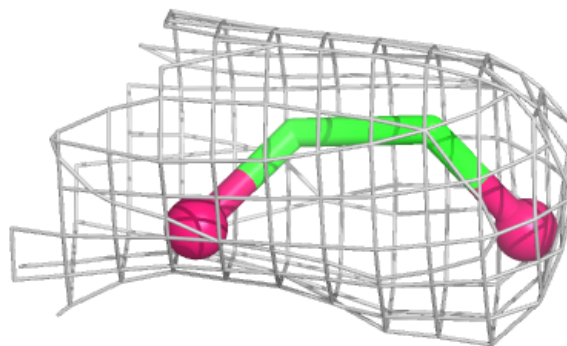
**Electron density around EDO E 421:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

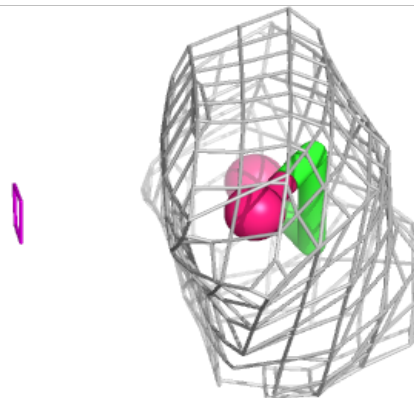
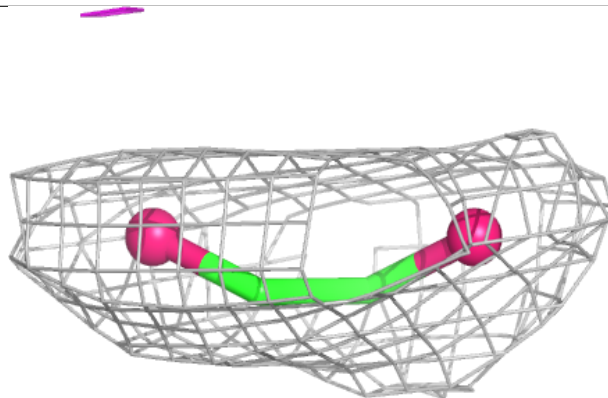
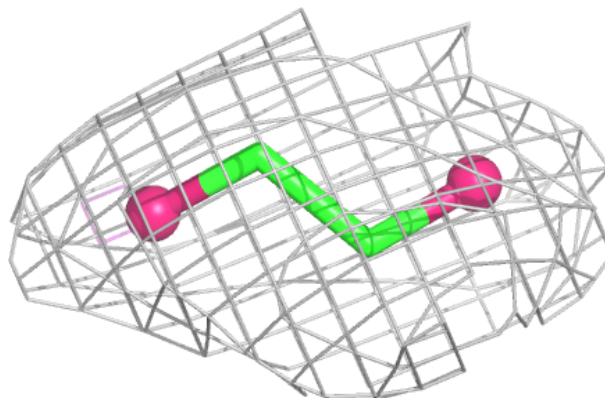


**Electron density around EDO B 447:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

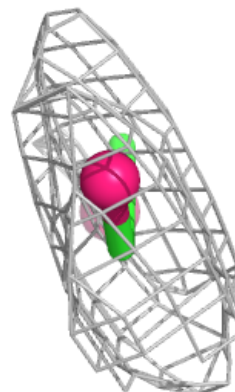
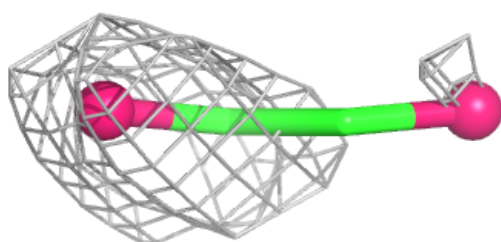
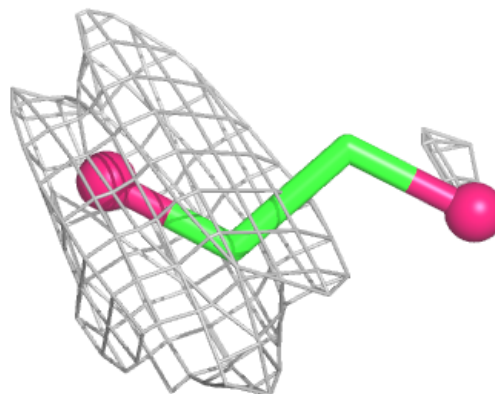
**Electron density around EDO E 428:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

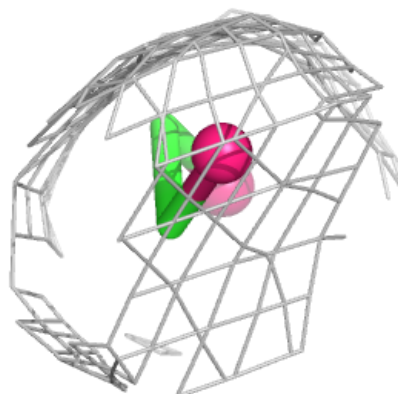
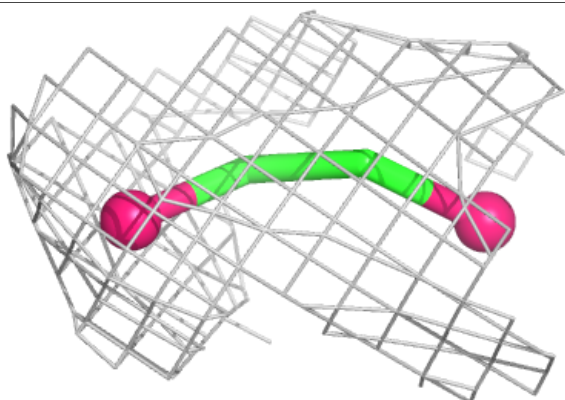
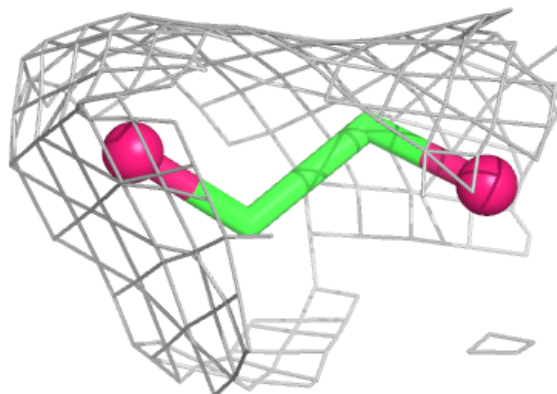


**Electron density around EDO C 402:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

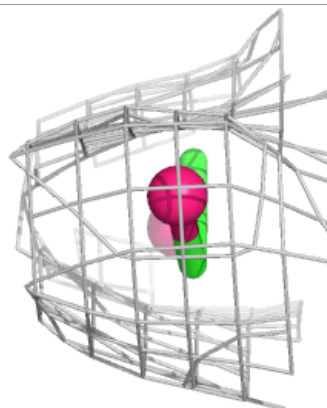
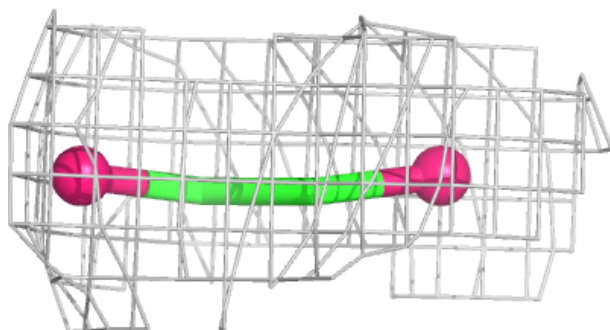
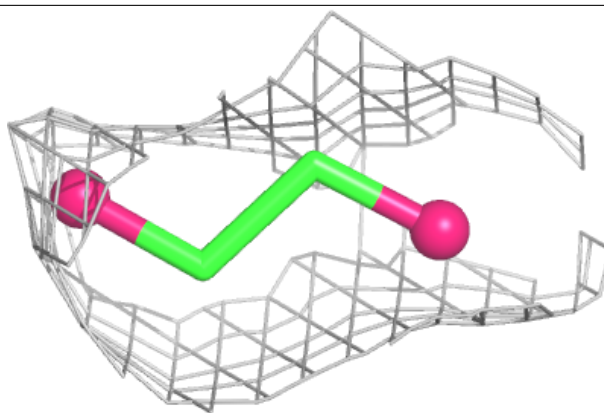
**Electron density around EDO A 417:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

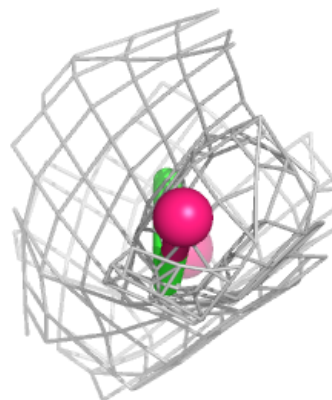
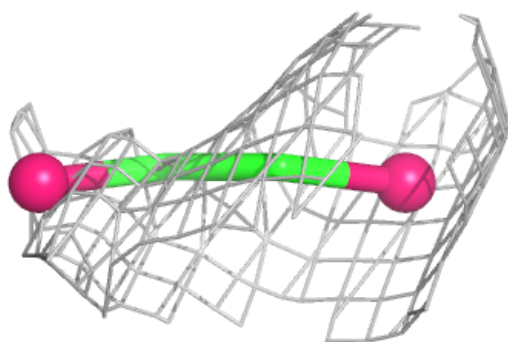
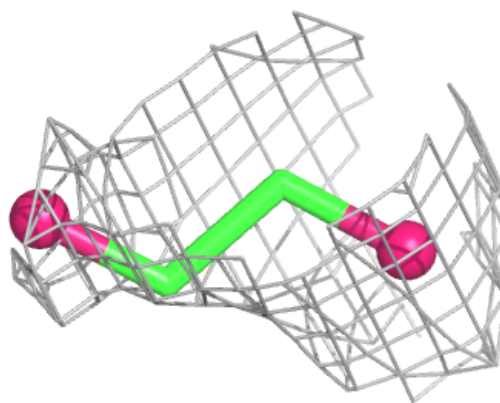


**Electron density around EDO A 412:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around EDO C 420:**

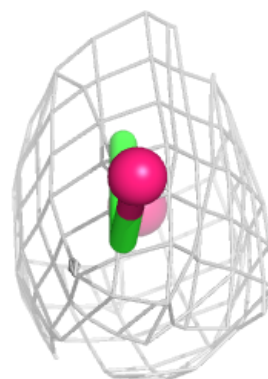
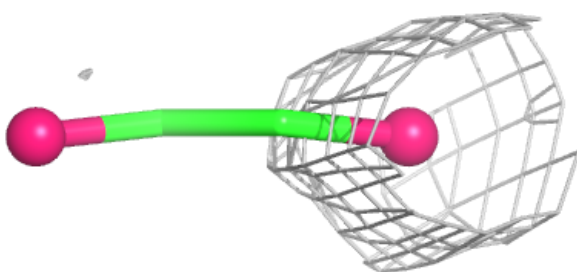
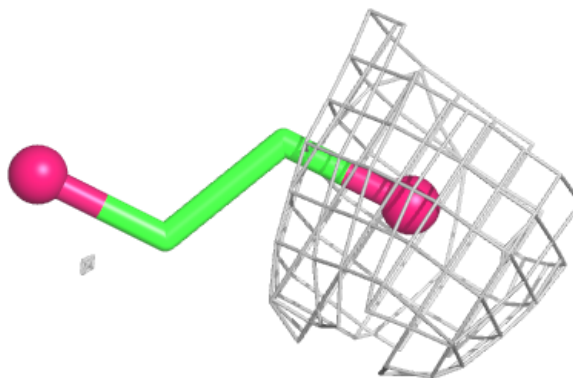
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



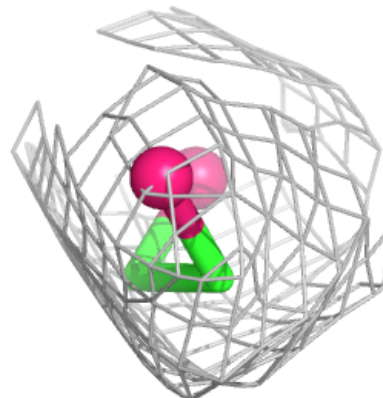
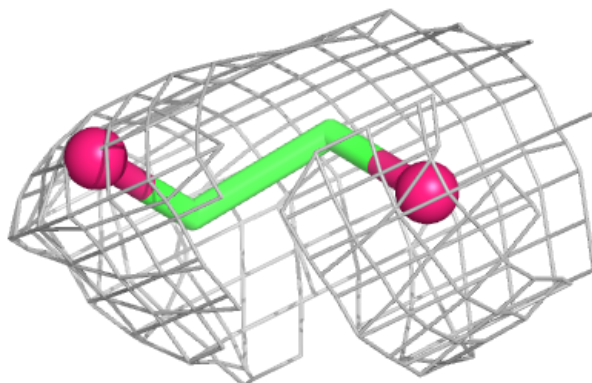
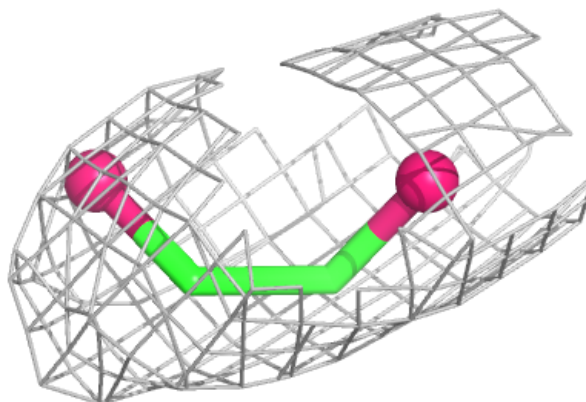


**Electron density around EDO E 434:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

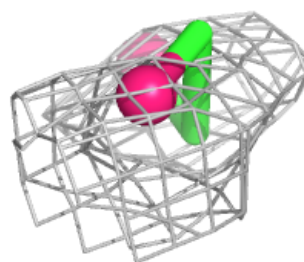
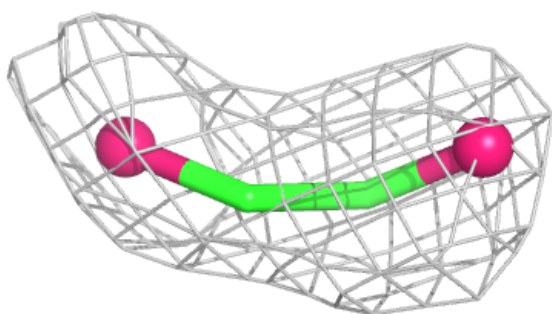
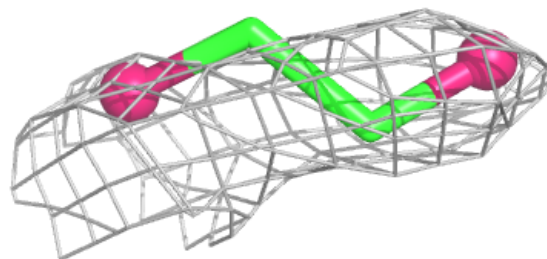
**Electron density around EDO C 424:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

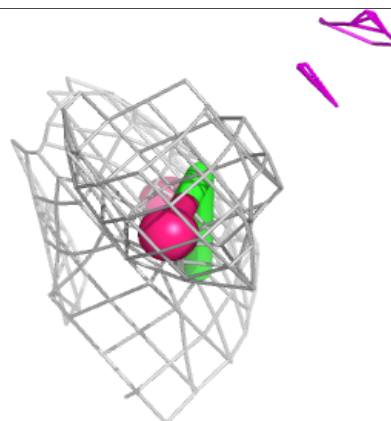
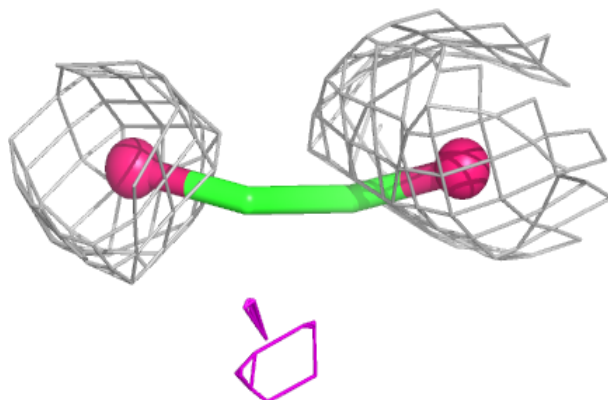
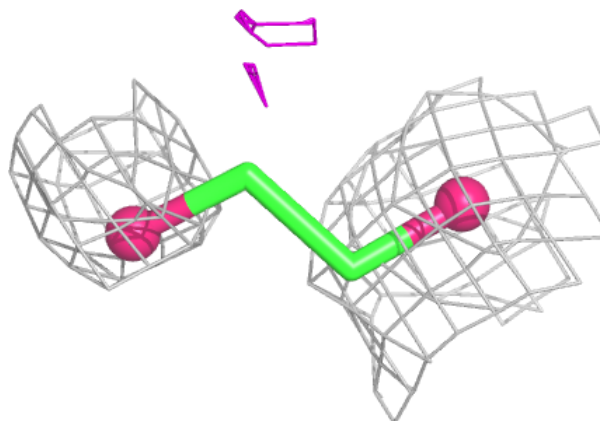


**Electron density around EDO E 443:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

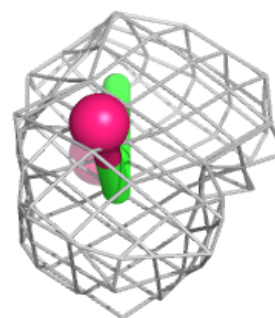
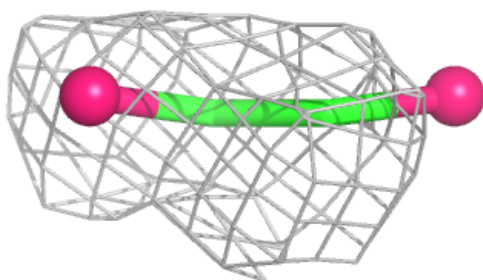
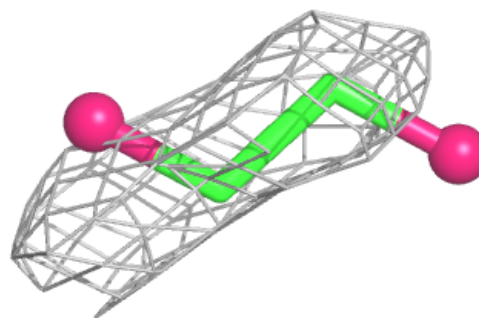
**Electron density around EDO C 425:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

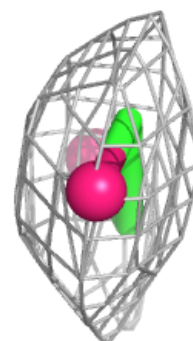
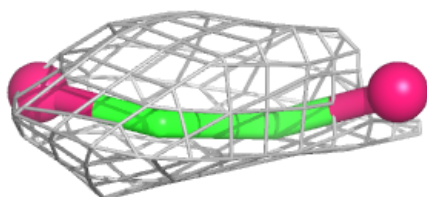
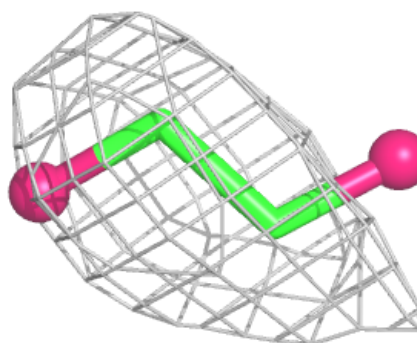


**Electron density around EDO A 422:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

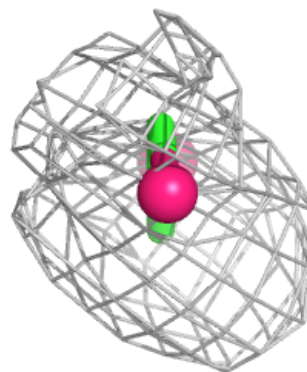
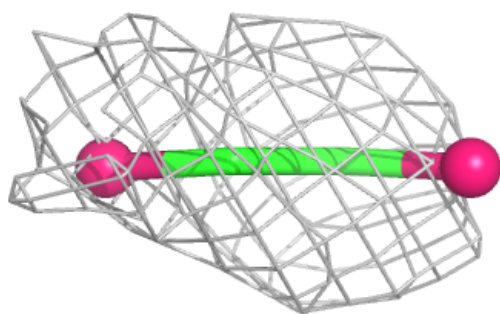
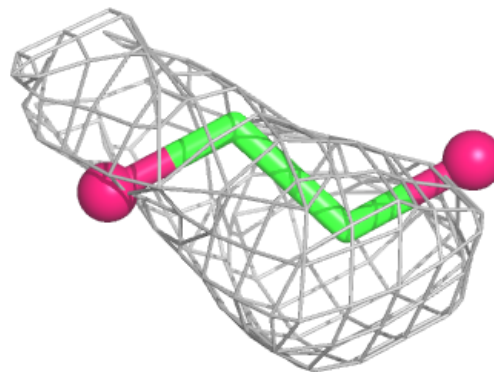
**Electron density around EDO C 428:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



**Electron density around EDO B 424:**

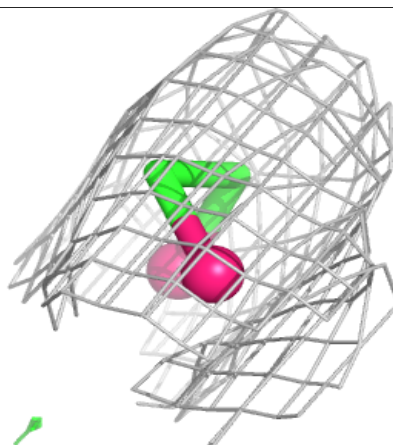
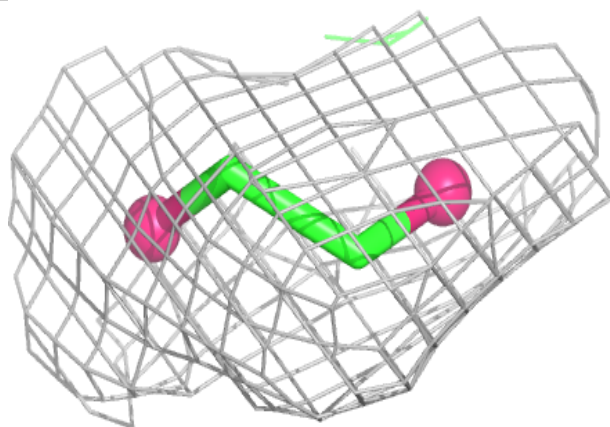
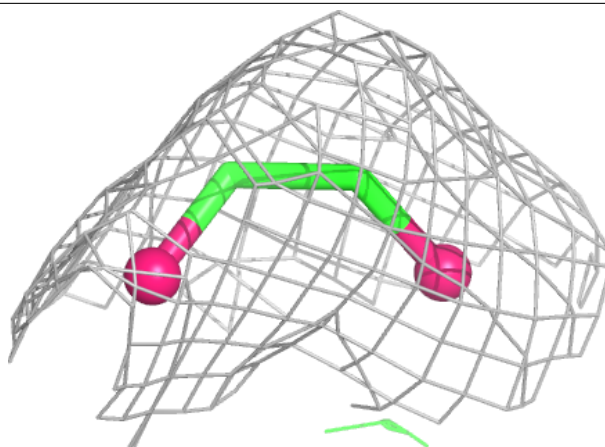
$2mF_o - DF_c$  (at 0.7 rmsd) in gray  
 $mF_o - DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





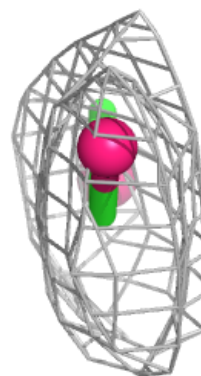
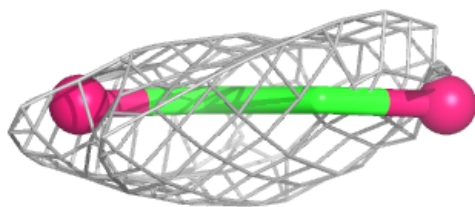
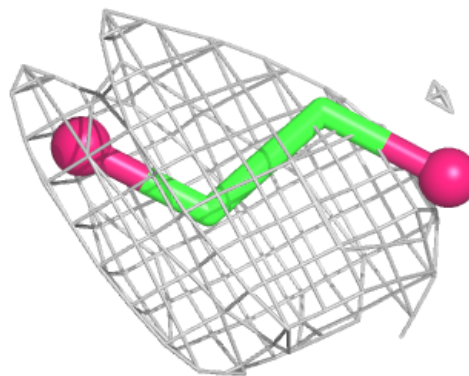
**Electron density around EDO A 403:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

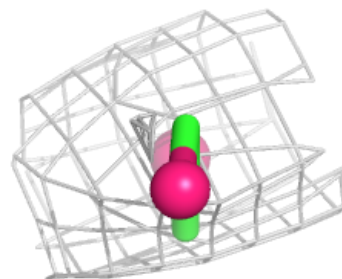
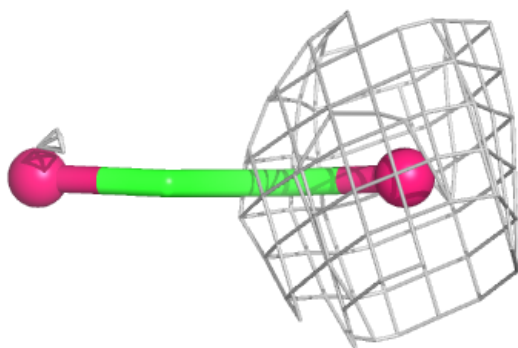
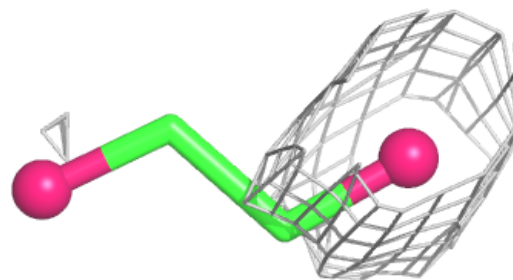


**Electron density around EDO F 408:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

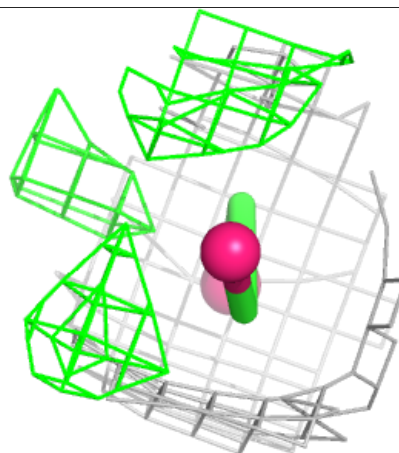
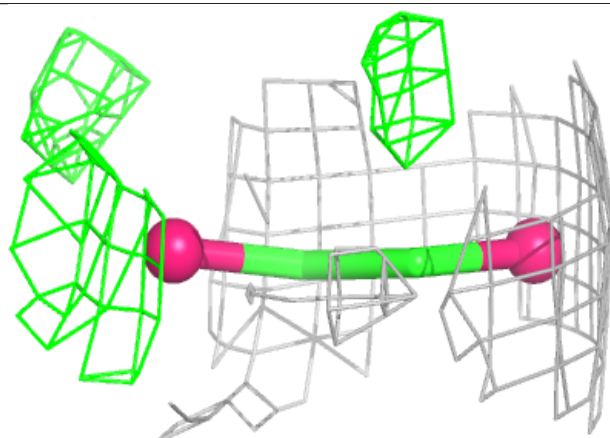
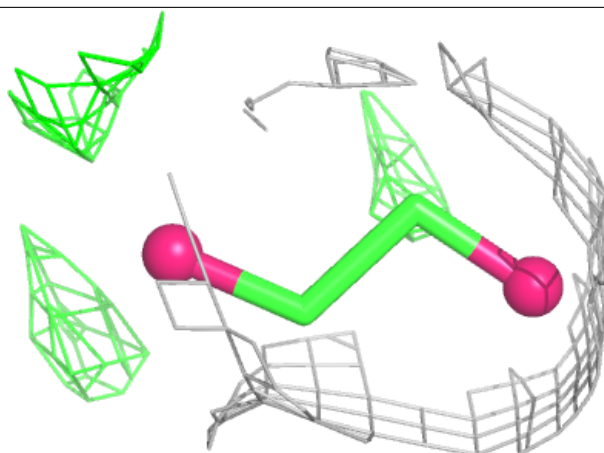
**Electron density around EDO F 409:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

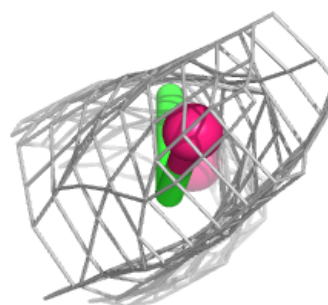
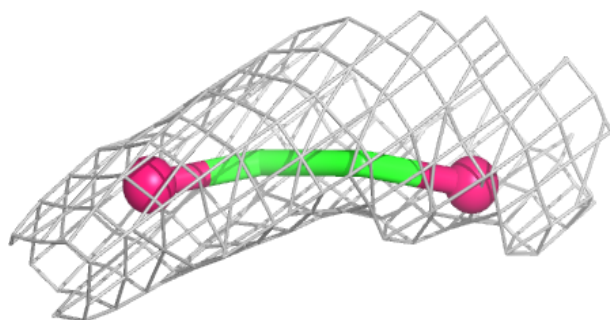
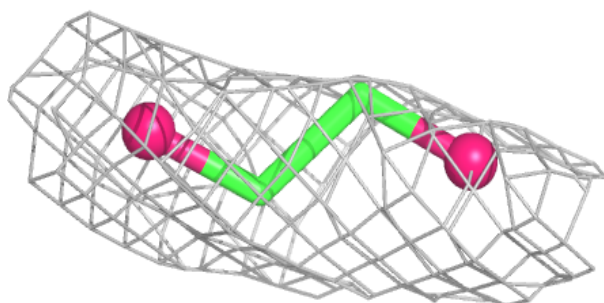


**Electron density around EDO F 416:**

$2mF_o - DF_c$  (at 0.7 rmsd) in gray  
 $mF_o - DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

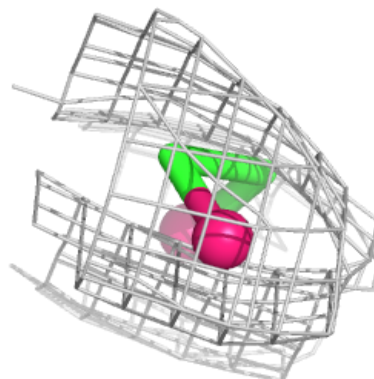
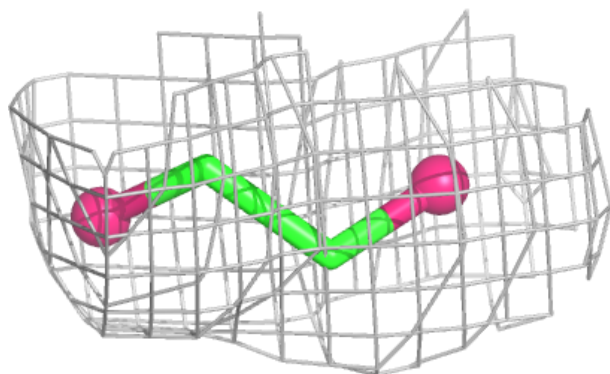
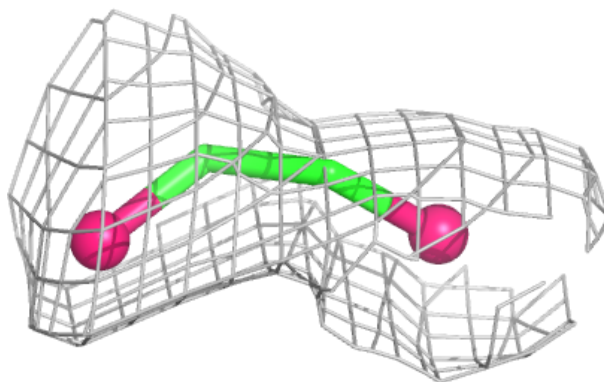
**Electron density around EDO C 434:**

$2mF_o - DF_c$  (at 0.7 rmsd) in gray  
 $mF_o - DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

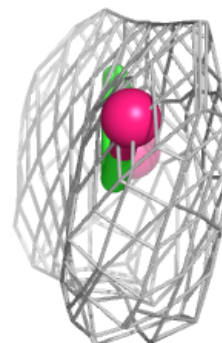
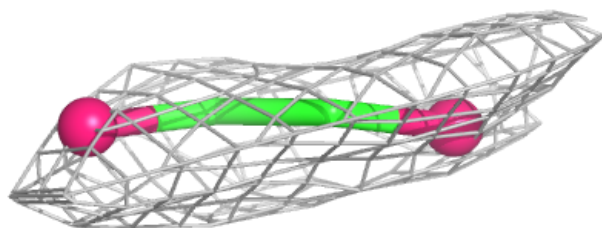
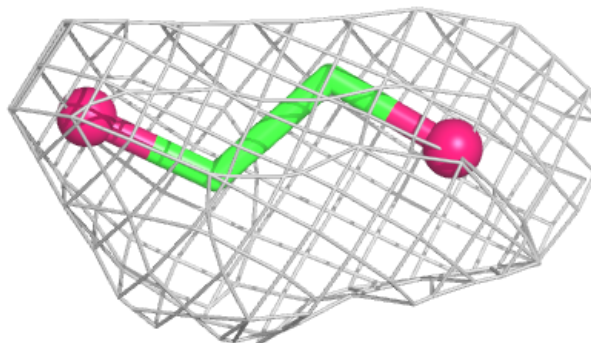


**Electron density around EDO A 434:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around EDO C 438:**

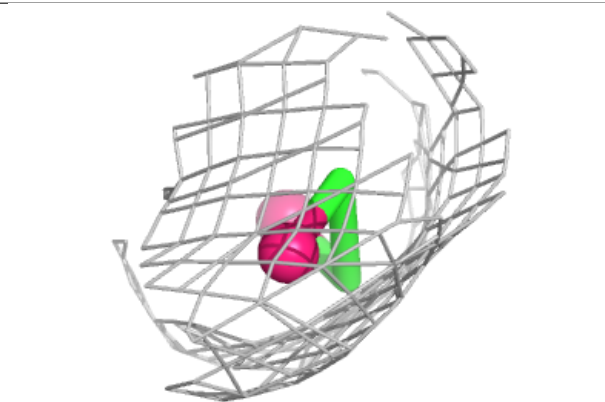
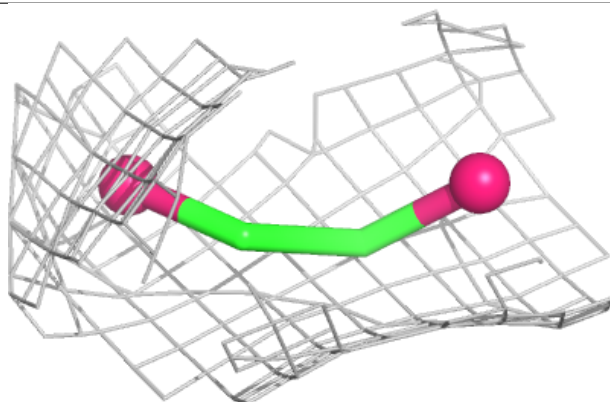
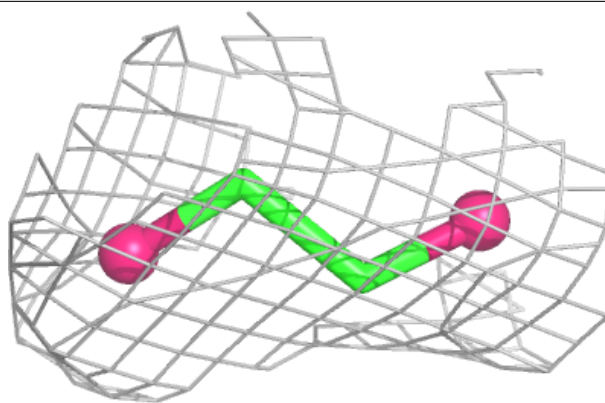
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



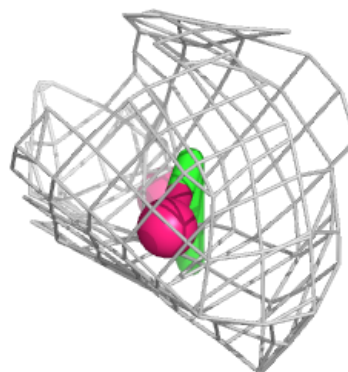
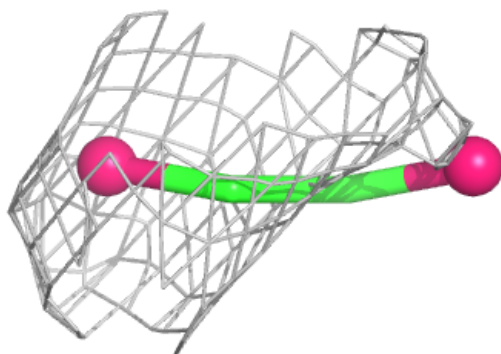
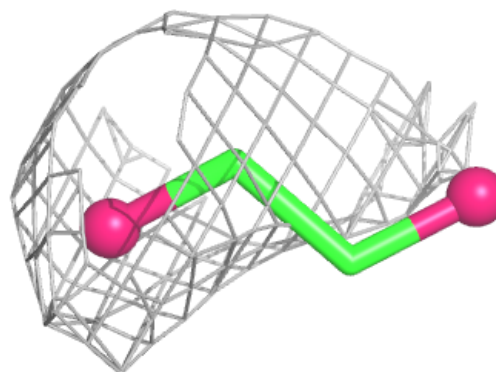


**Electron density around EDO F 425:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

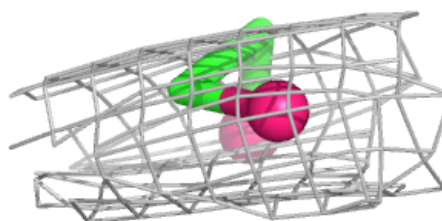
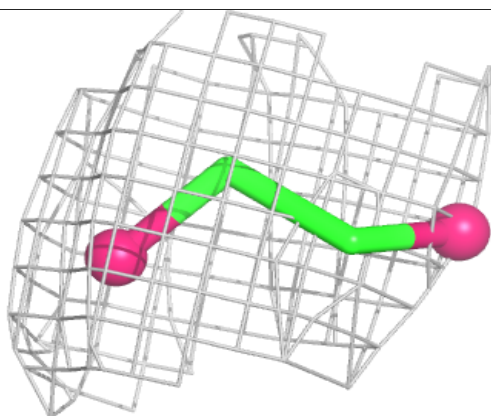
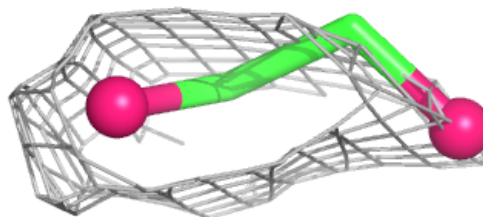
**Electron density around EDO F 427:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

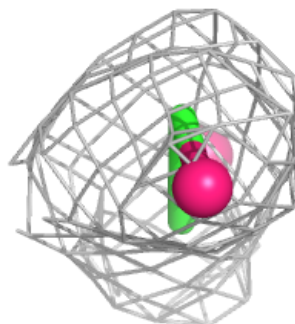
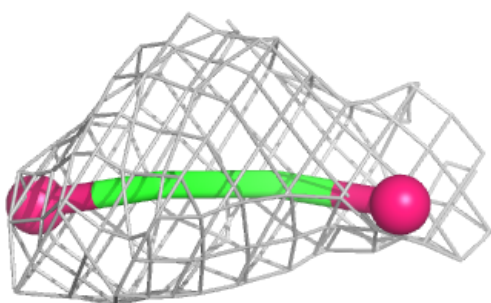
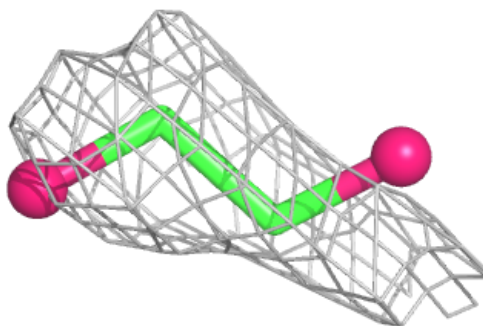


**Electron density around EDO F 429:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

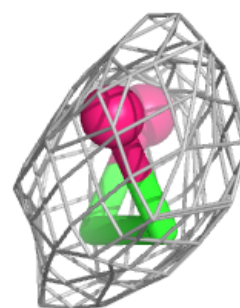
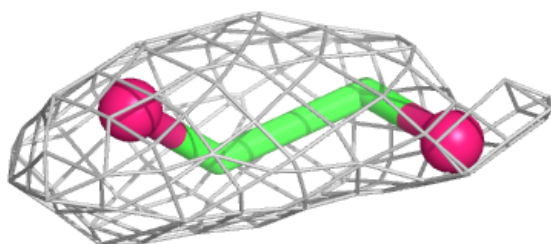
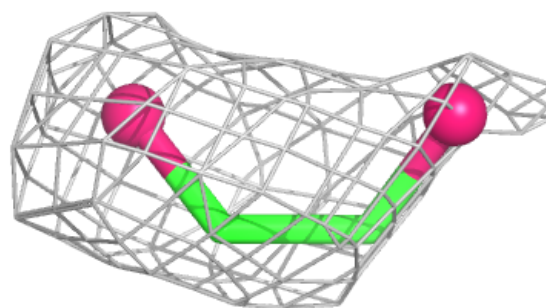
**Electron density around EDO F 430:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

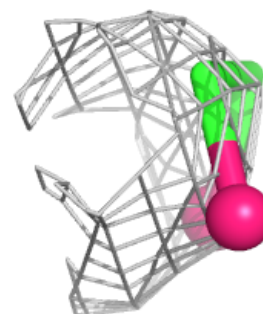
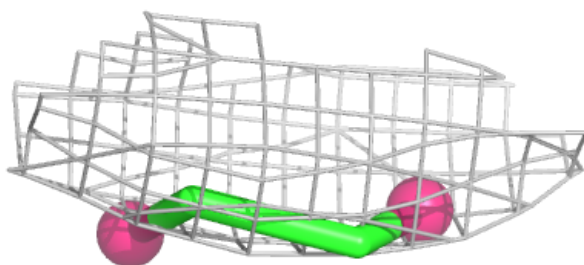
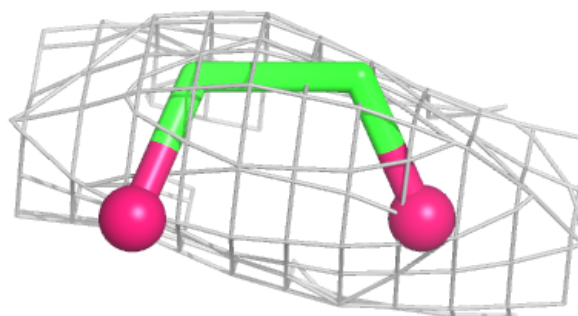


**Electron density around EDO B 430:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

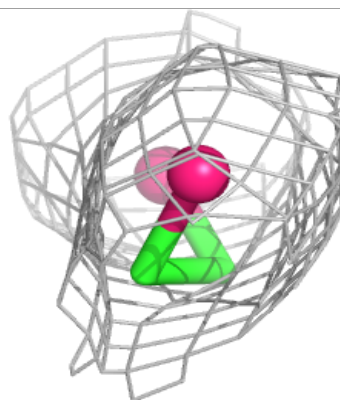
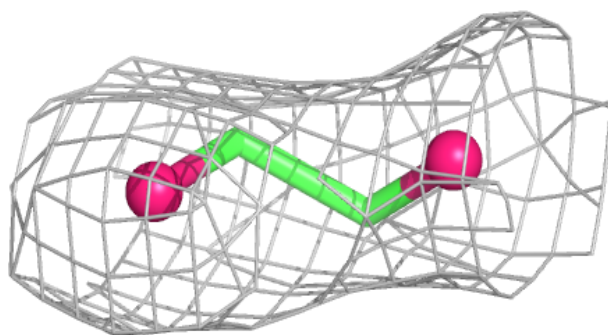
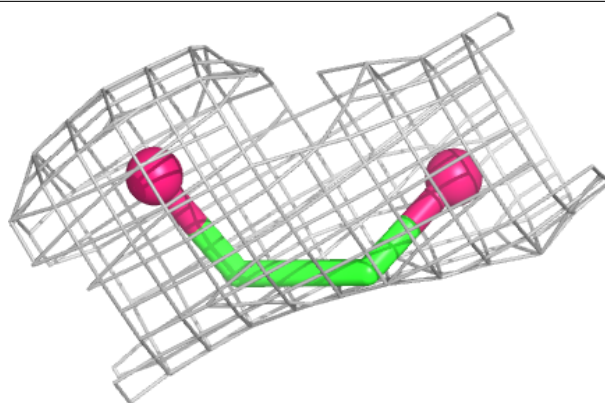
**Electron density around EDO A 436:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

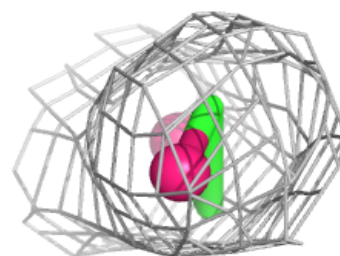
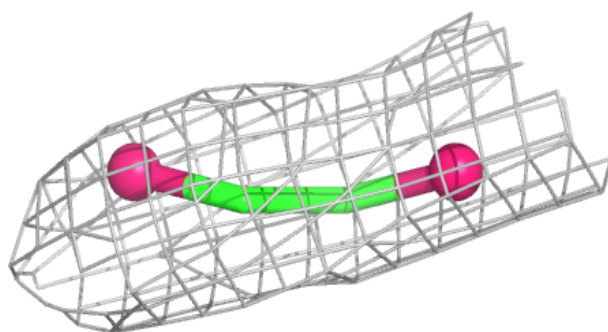
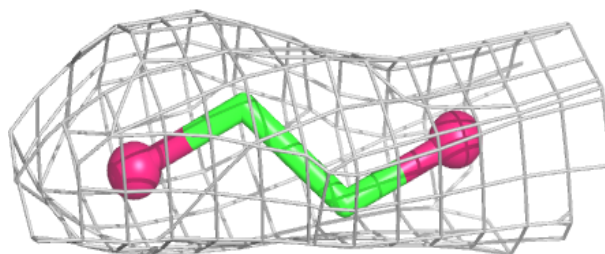


**Electron density around EDO F 439:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around EDO A 438:**

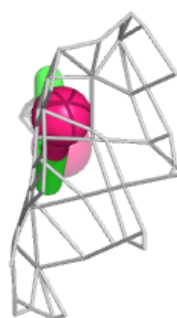
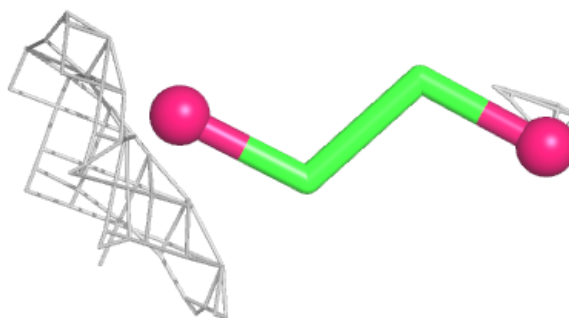
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





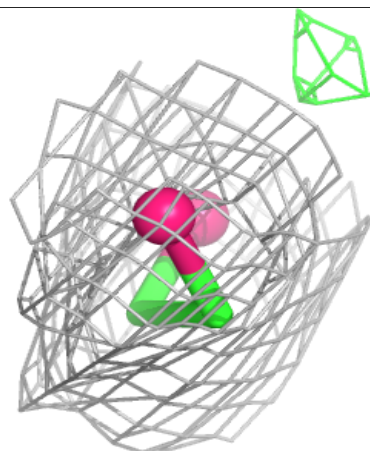
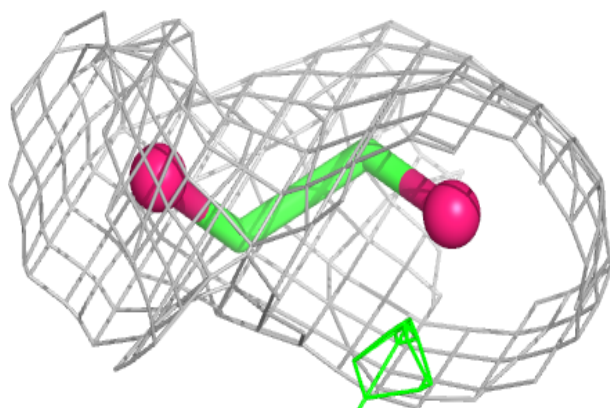
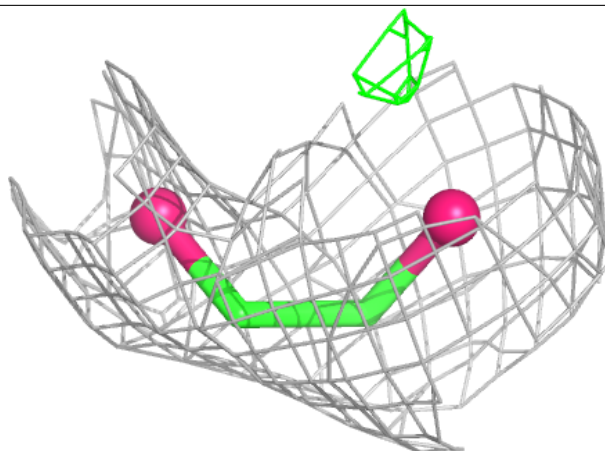
**Electron density around EDO F 441:**

$2mF_o - DF_c$  (at 0.7 rmsd) in gray  
 $mF_o - DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

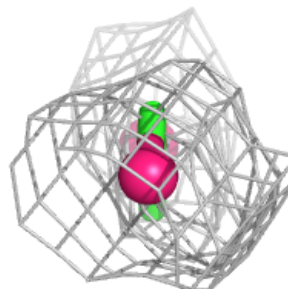
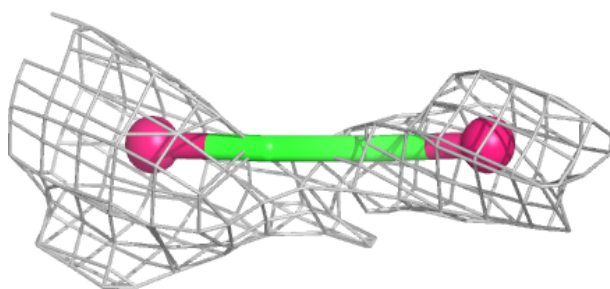
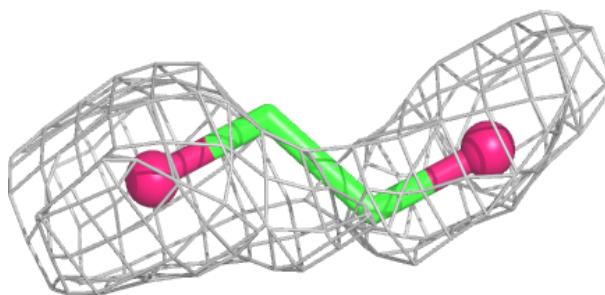


**Electron density around EDO D 406:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

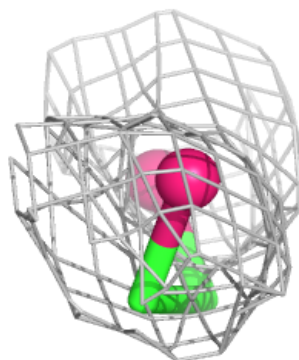
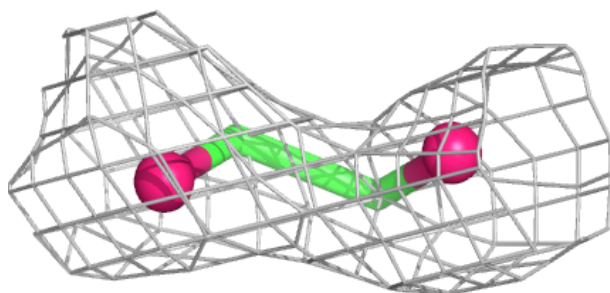
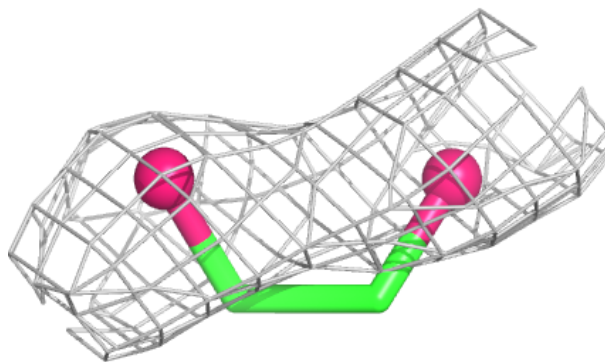
**Electron density around EDO D 407:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

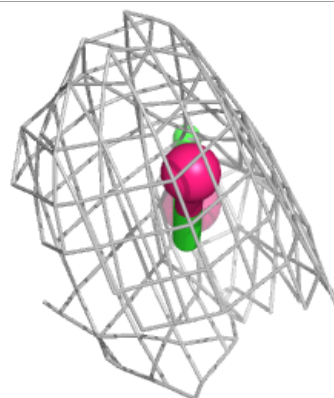
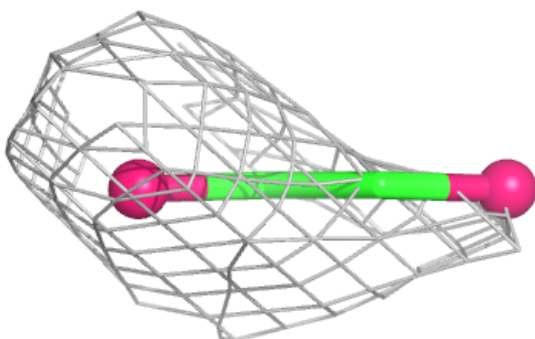
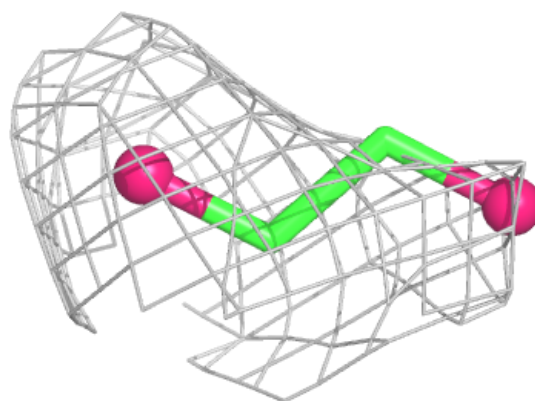


**Electron density around EDO A 439:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

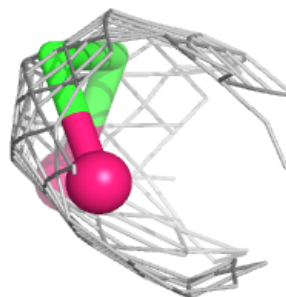
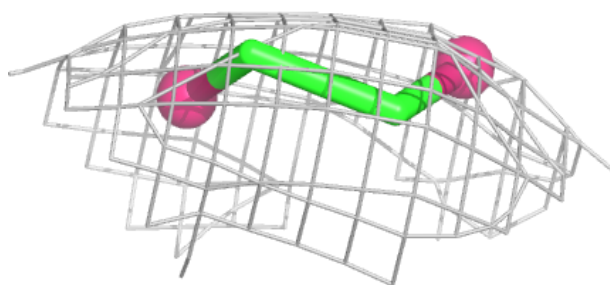
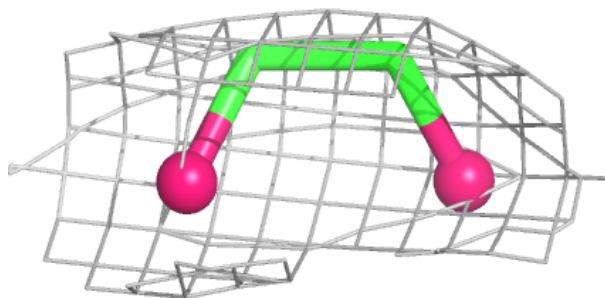
**Electron density around EDO D 422:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

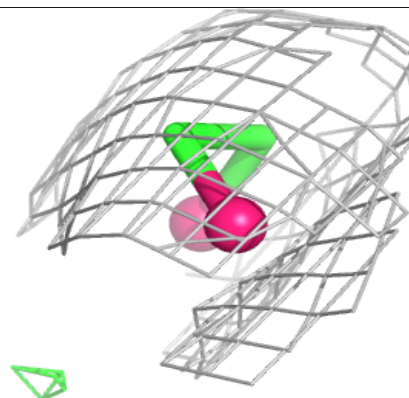
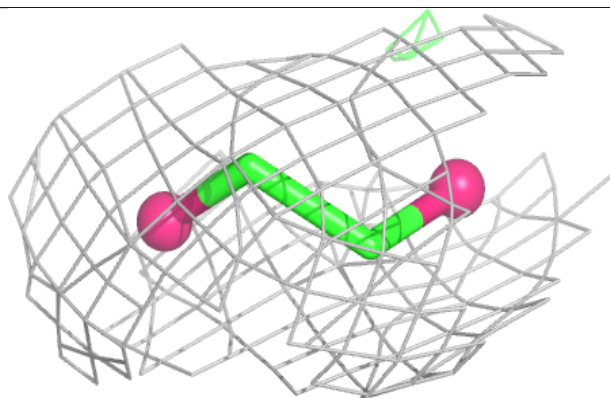
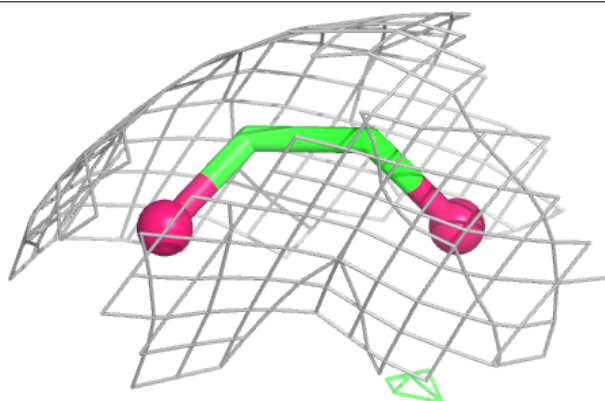


**Electron density around EDO D 434:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

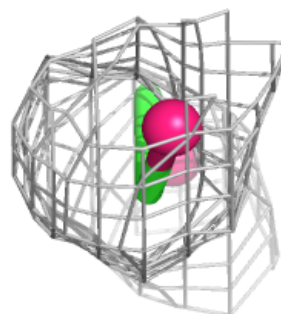
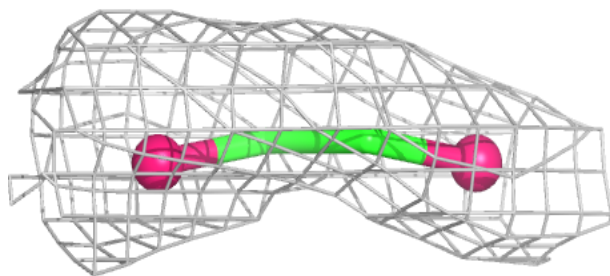
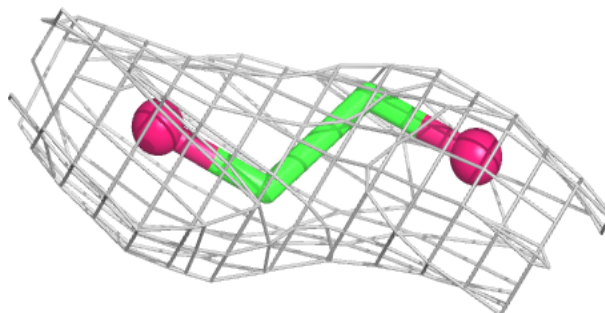
**Electron density around EDO B 405:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

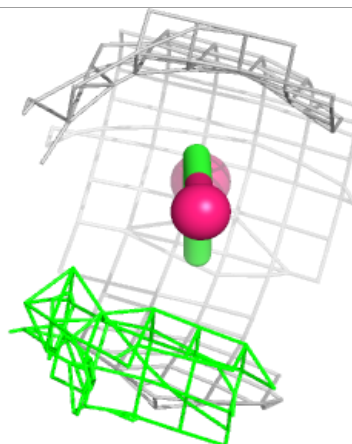
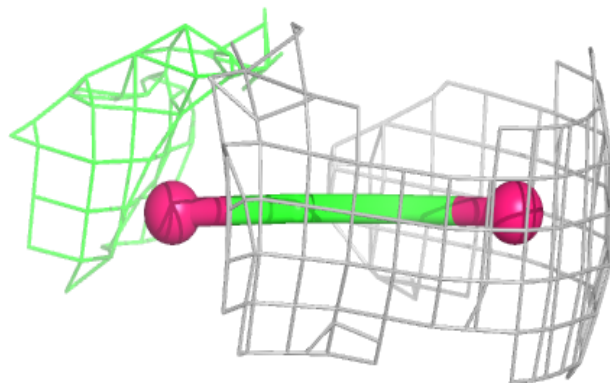
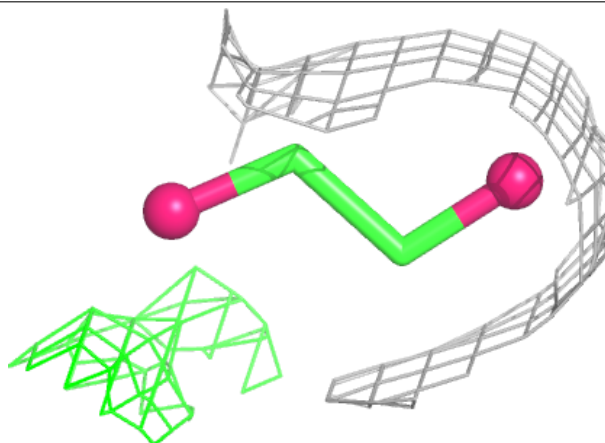


**Electron density around EDO C 406:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around EDO C 410:**

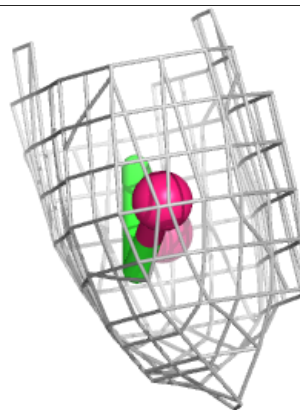
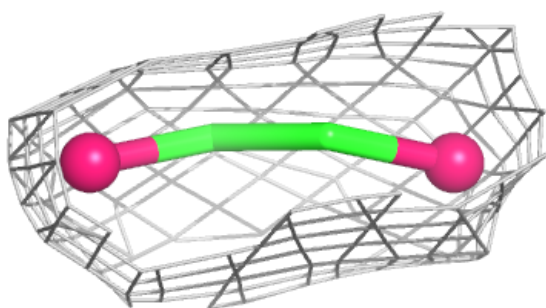
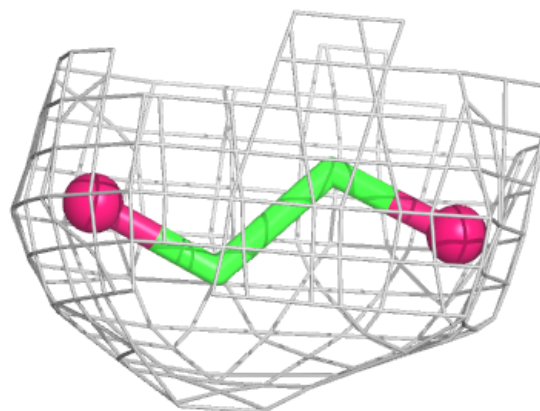
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



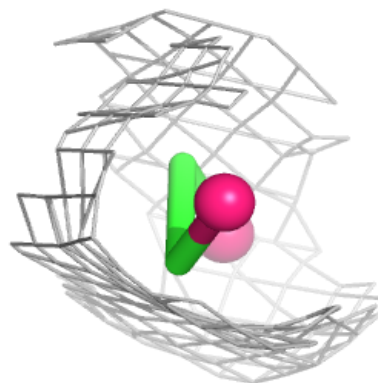
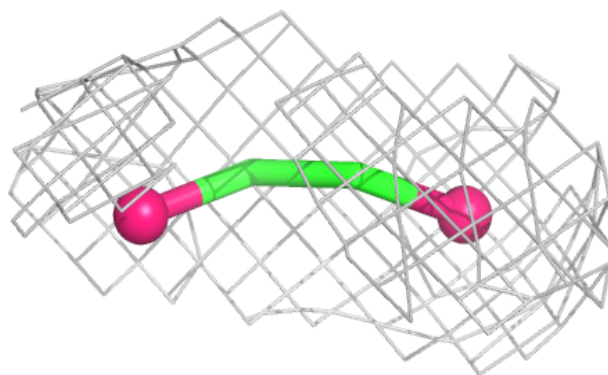
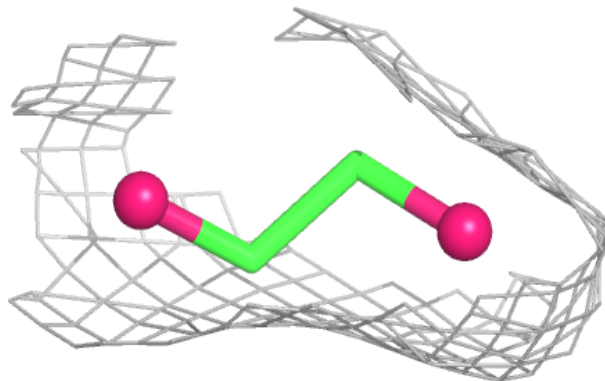


**Electron density around EDO E 401:**

$2mF_o - DF_c$  (at 0.7 rmsd) in gray  
 $mF_o - DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

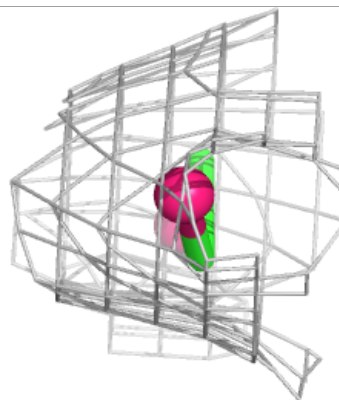
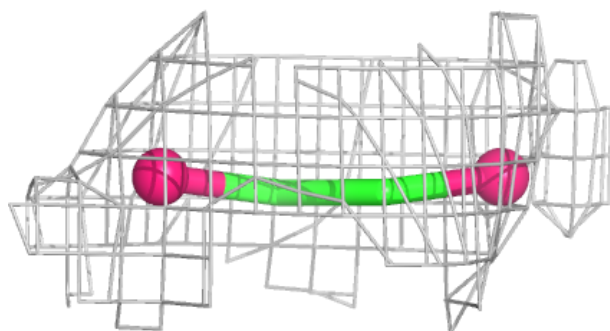
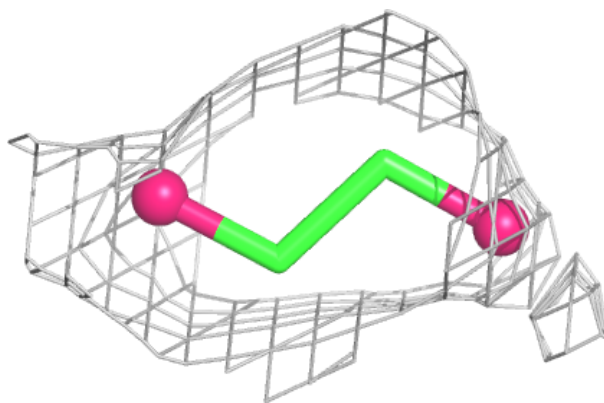
**Electron density around EDO E 402:**

$2mF_o - DF_c$  (at 0.7 rmsd) in gray  
 $mF_o - DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

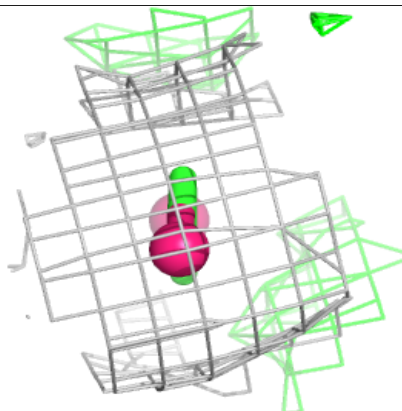
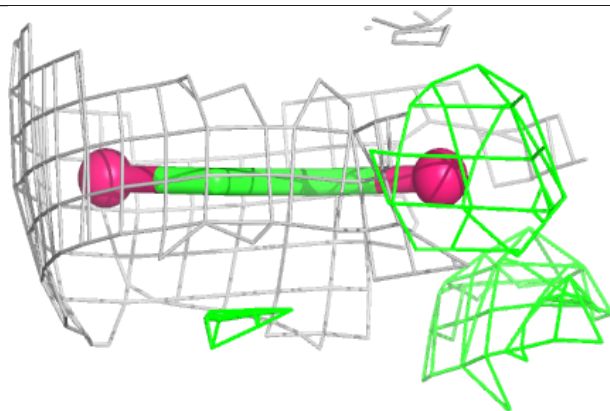
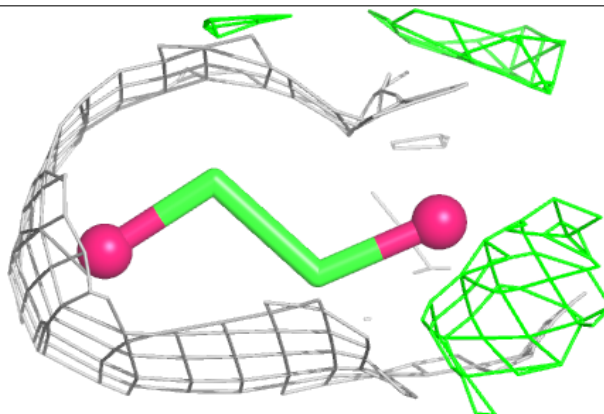


**Electron density around EDO C 414:**

$2mF_o - DF_c$  (at 0.7 rmsd) in gray  
 $mF_o - DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

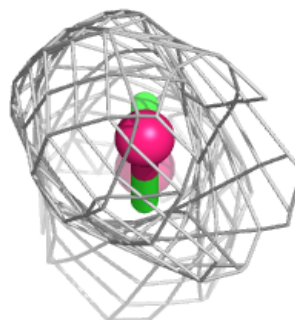
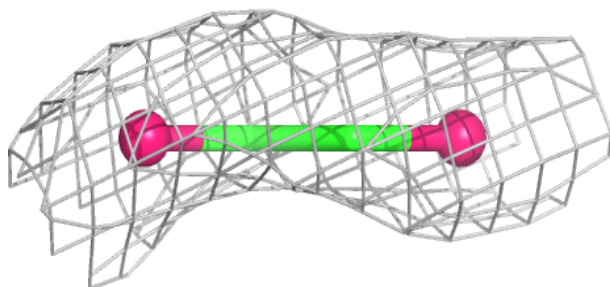
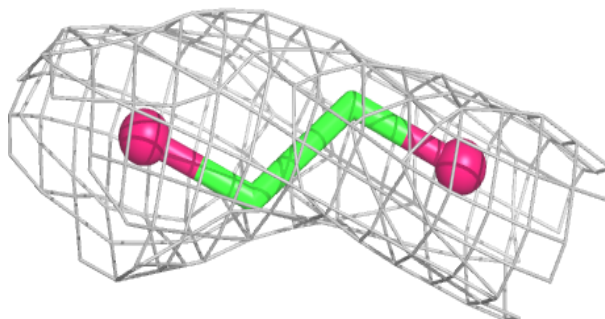
**Electron density around EDO B 410:**

$2mF_o - DF_c$  (at 0.7 rmsd) in gray  
 $mF_o - DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

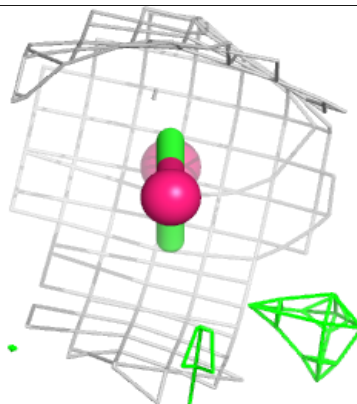
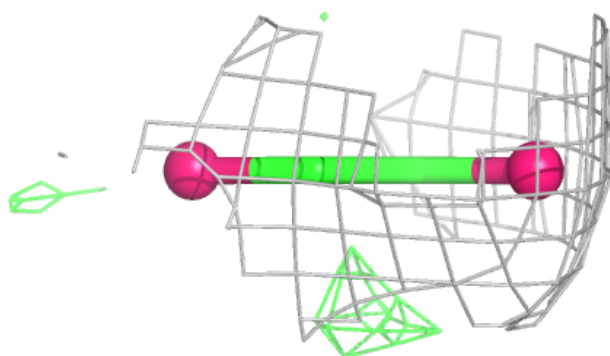
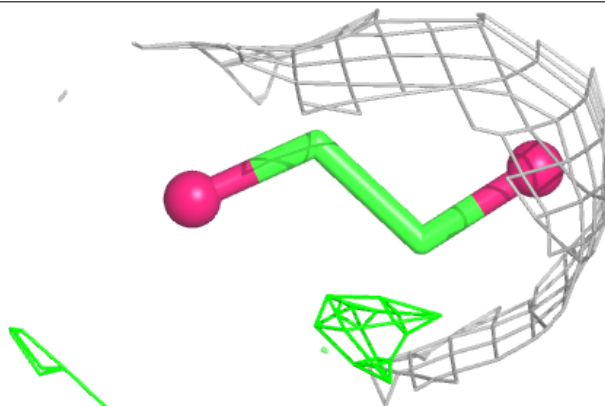


**Electron density around EDO E 406:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around EDO E 410:**

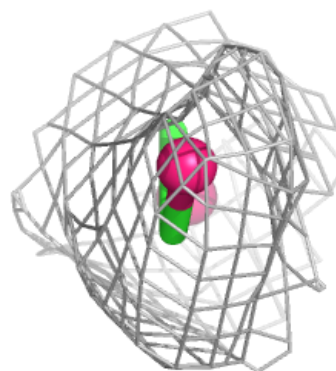
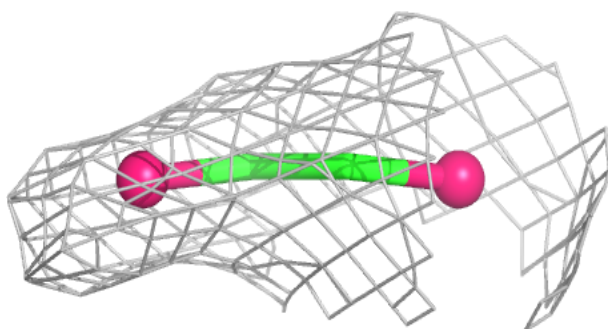
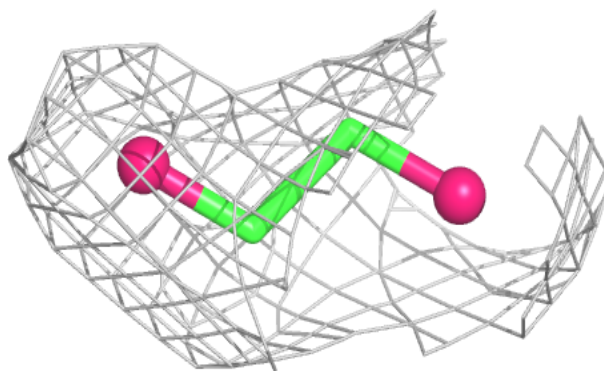
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



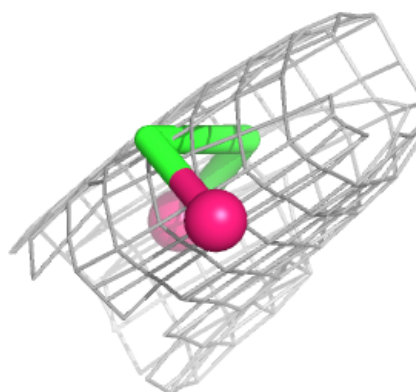
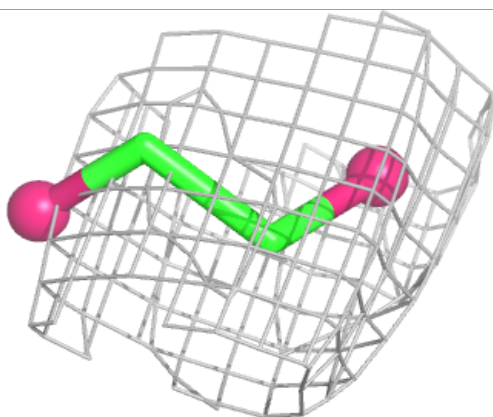
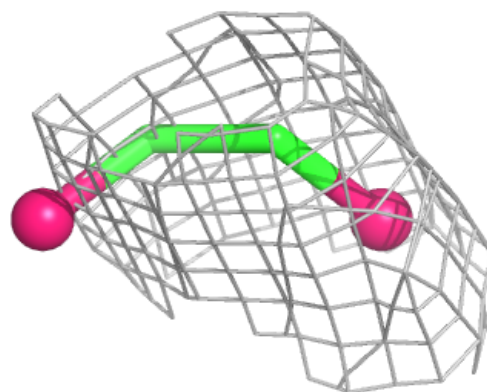


**Electron density around EDO B 412:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

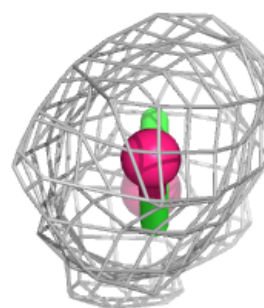
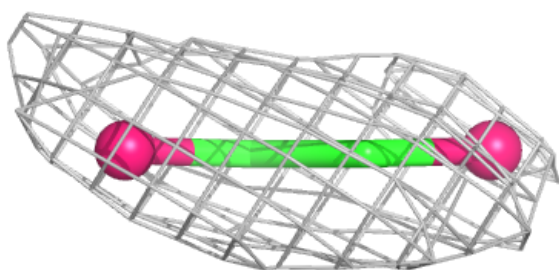
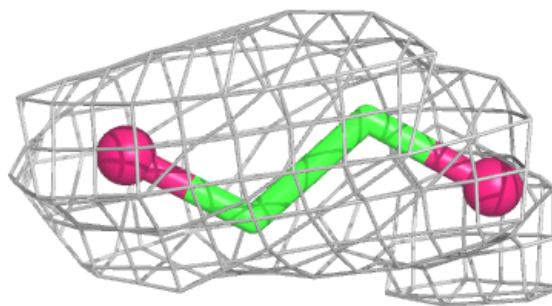
**Electron density around EDO C 422:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

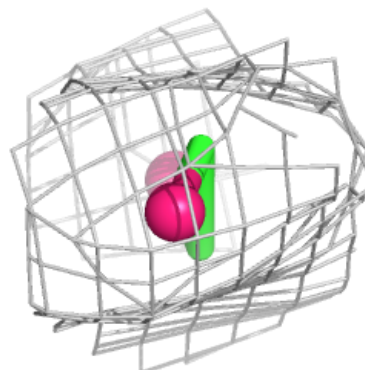
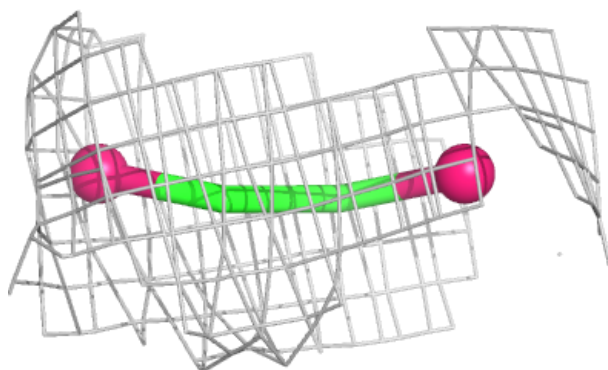
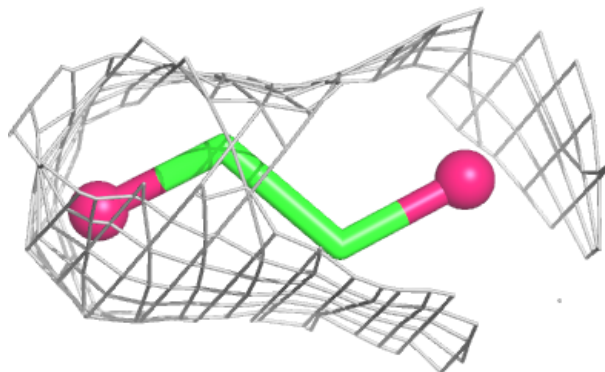


**Electron density around EDO A 429:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

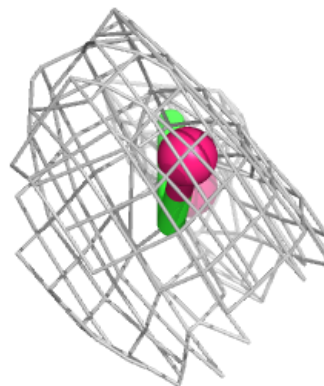
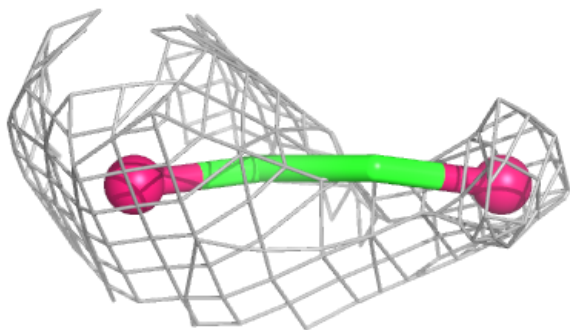
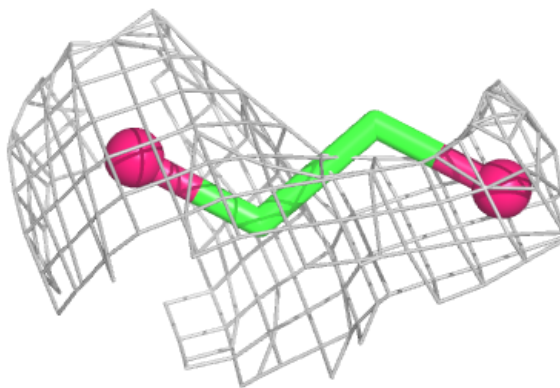
**Electron density around EDO B 417:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

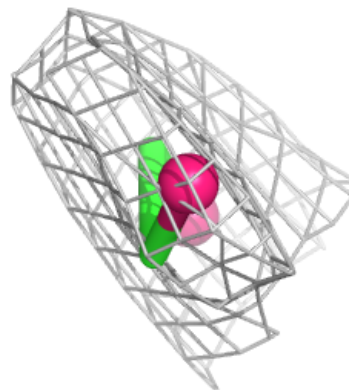
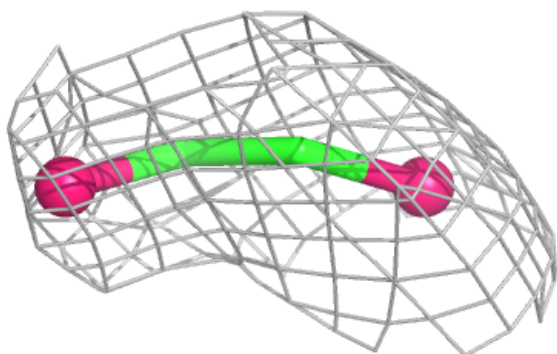
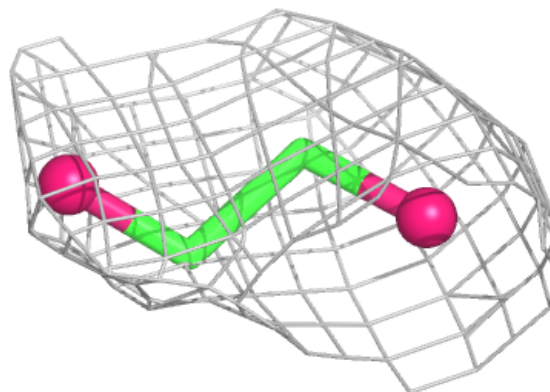


**Electron density around EDO E 423:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

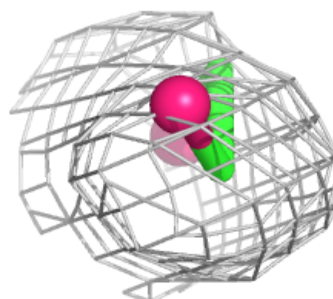
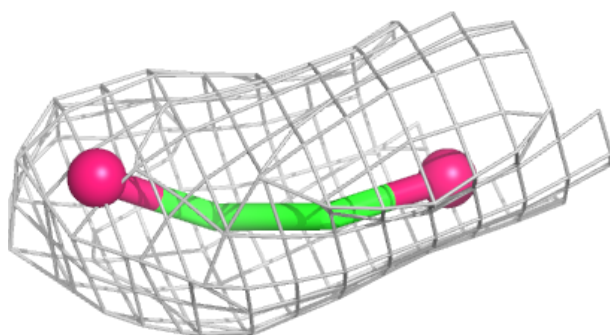
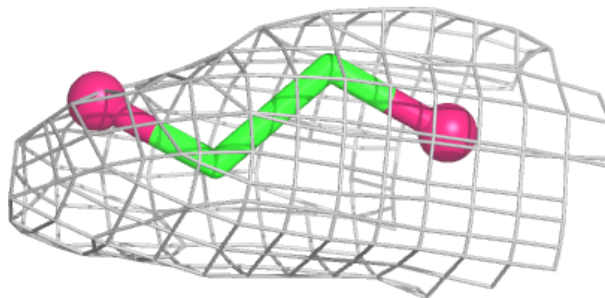
**Electron density around EDO E 425:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

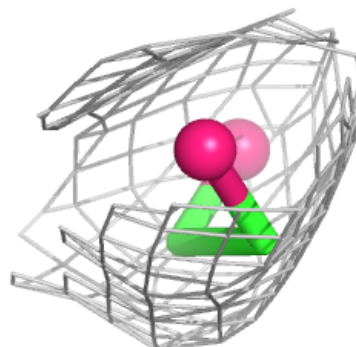
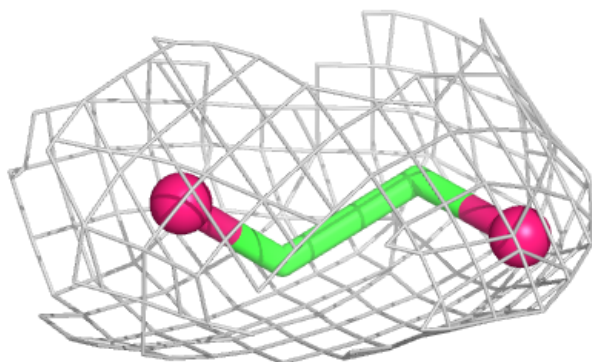
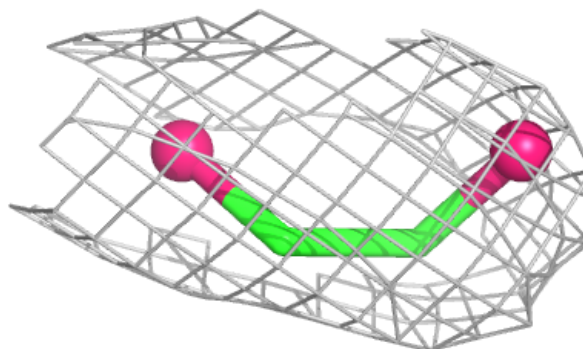


**Electron density around EDO B 418:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around EDO E 427:**

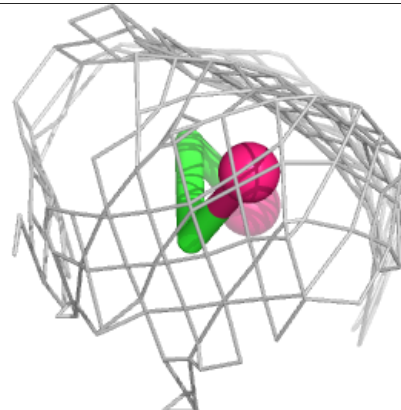
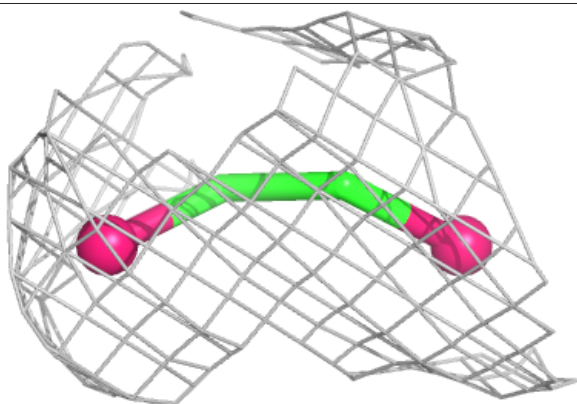
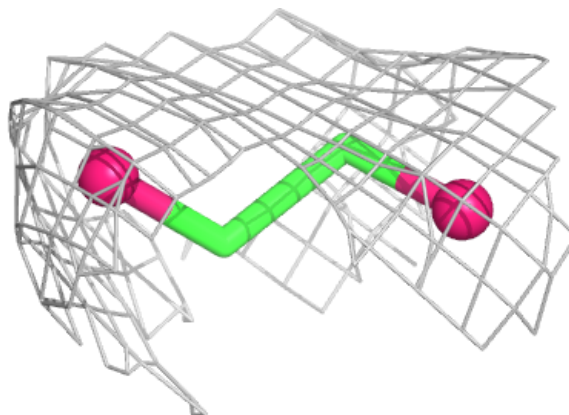
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



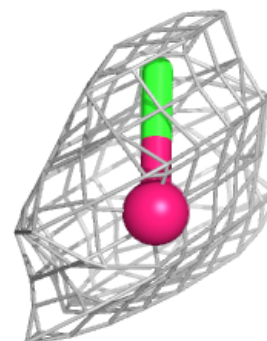
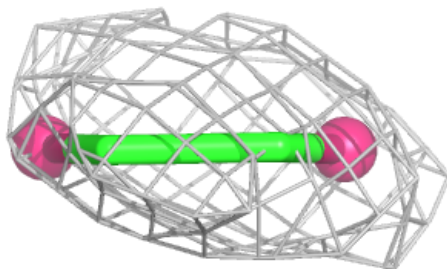
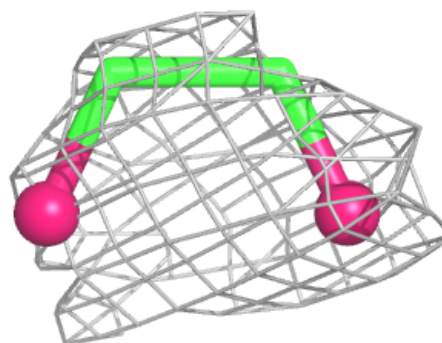


**Electron density around EDO B 419:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

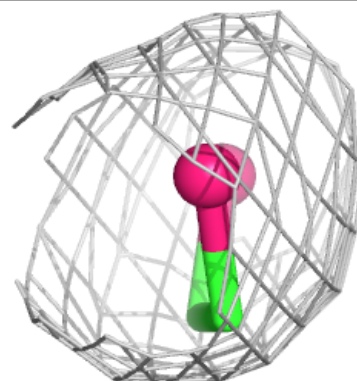
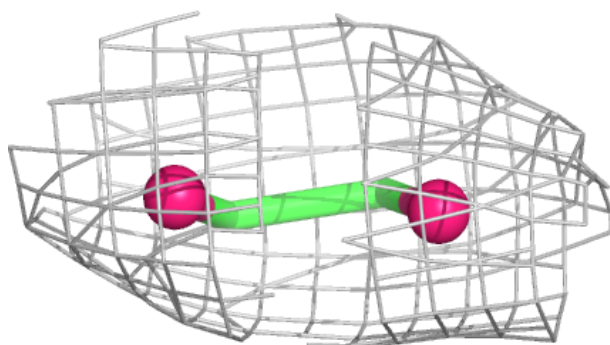
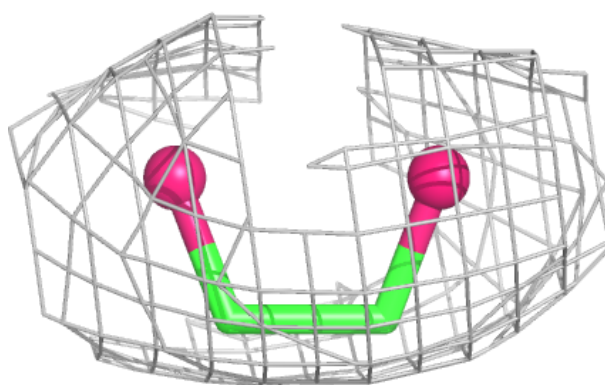
**Electron density around EDO C 429:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

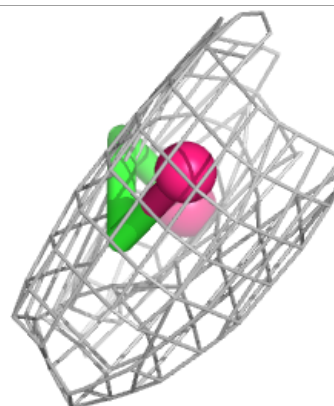
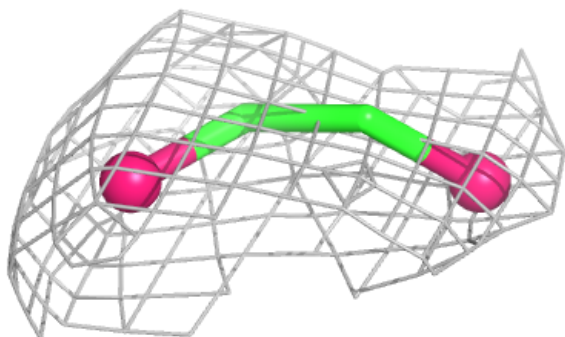
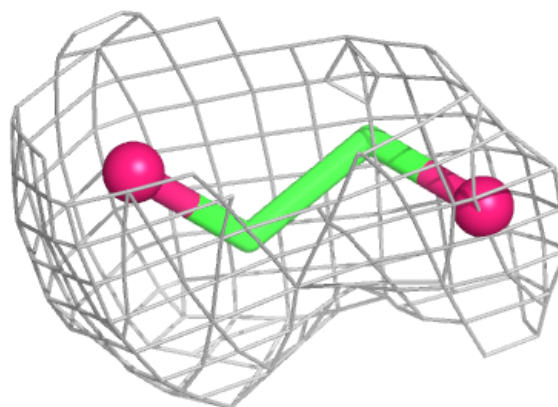


**Electron density around EDO C 431:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

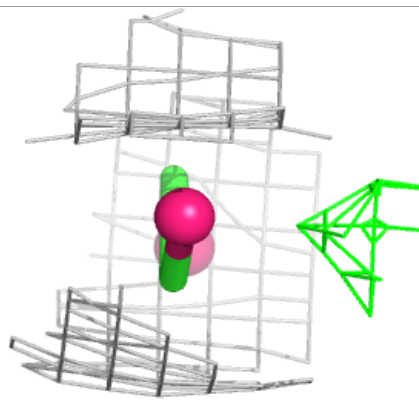
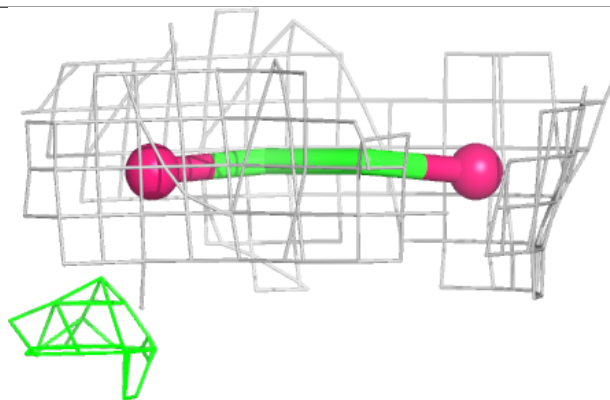
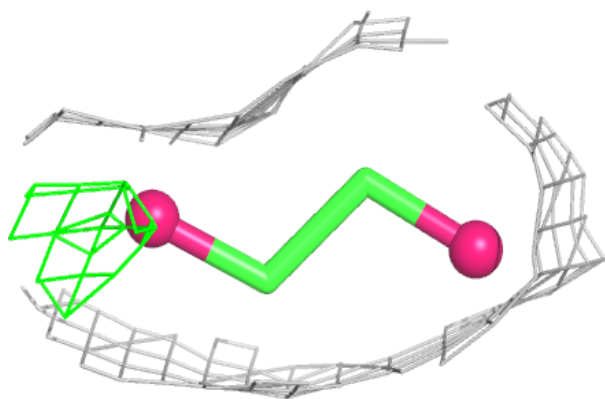
**Electron density around EDO B 423:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

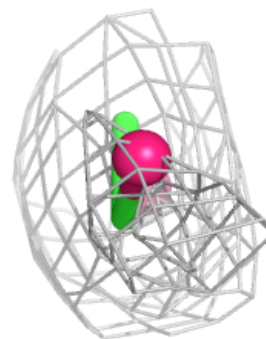
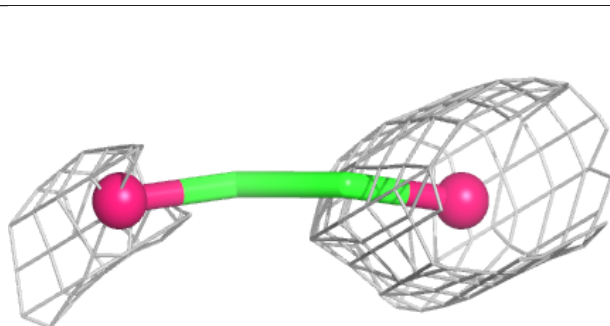
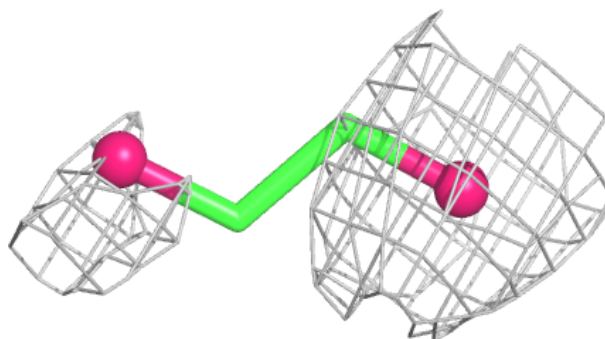


**Electron density around EDO A 407:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

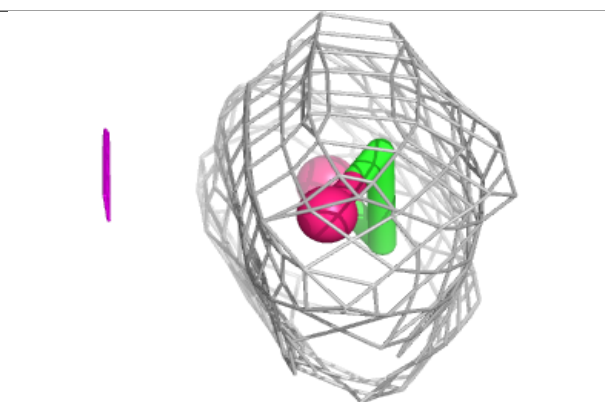
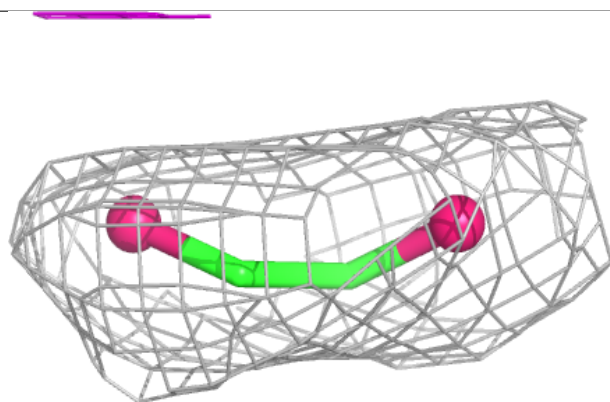
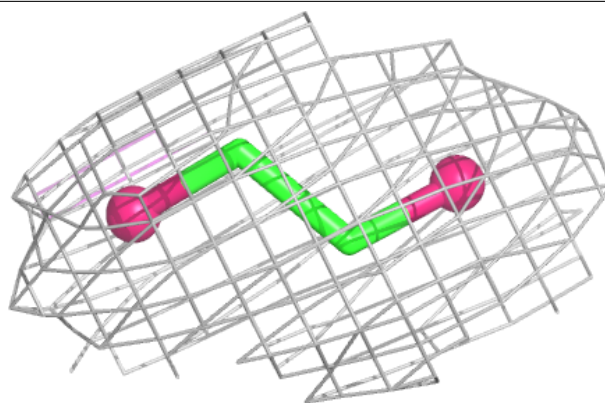
**Electron density around EDO A 432:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

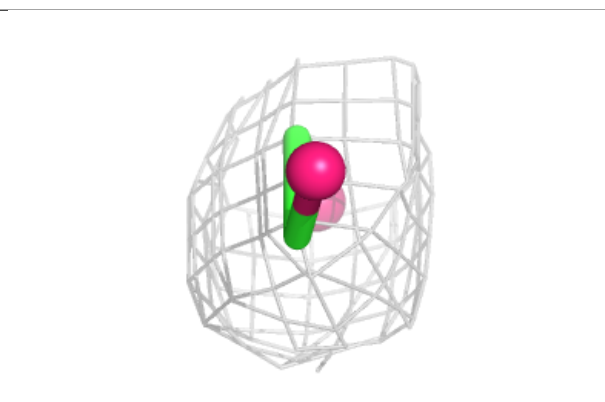
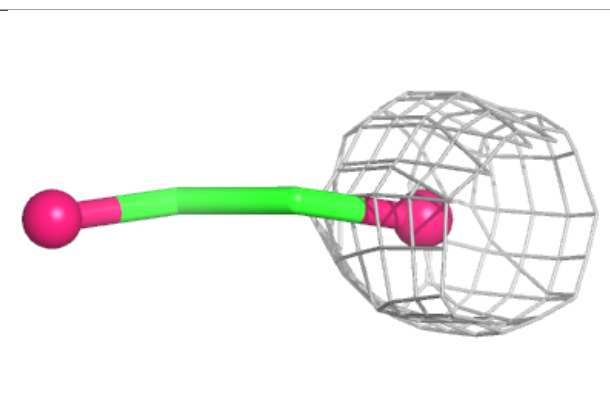
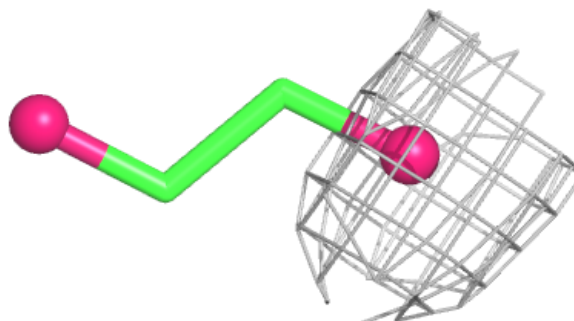


**Electron density around EDO E 437:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around EDO E 438:**

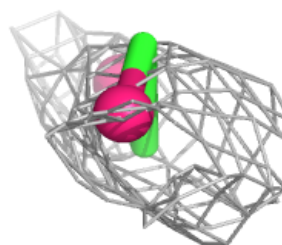
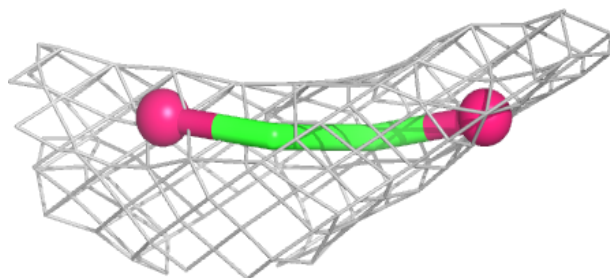
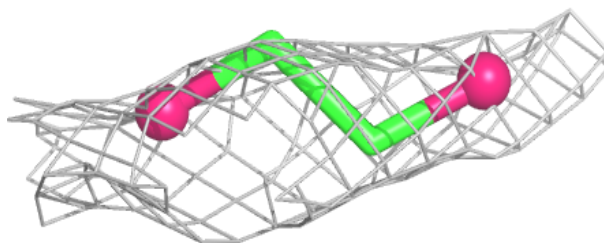
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



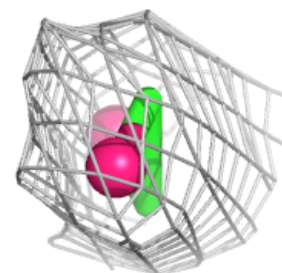
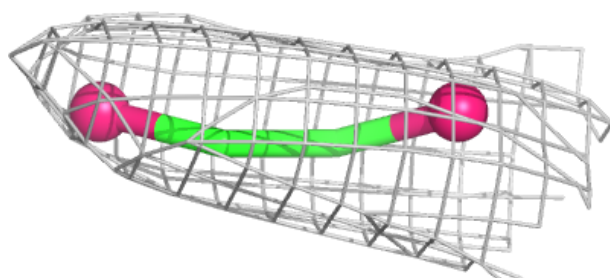
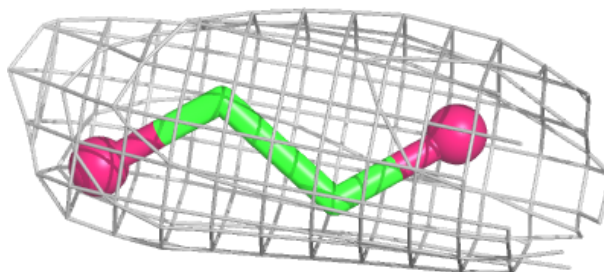


**Electron density around EDO E 439:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

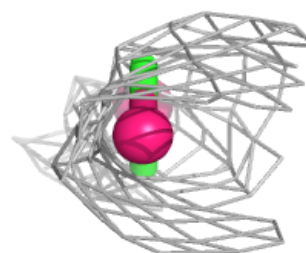
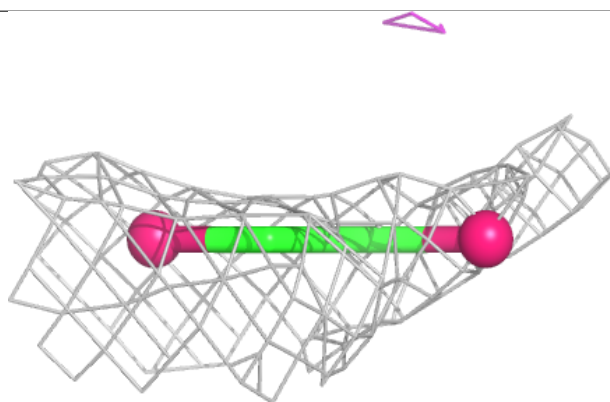
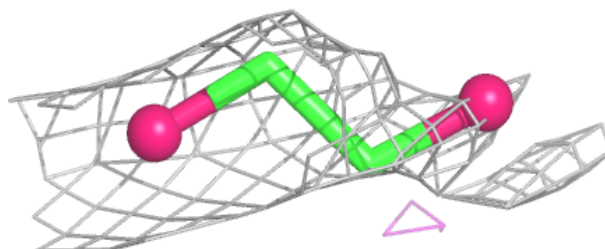
**Electron density around EDO C 435:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

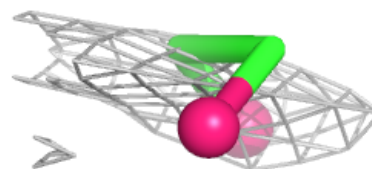
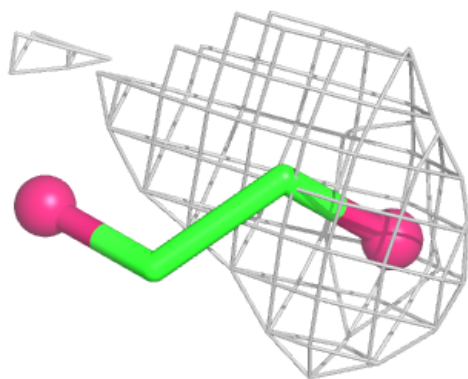
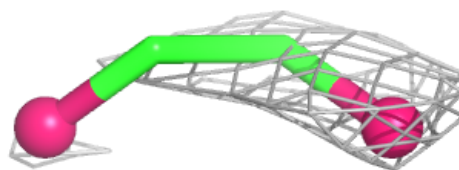


**Electron density around EDO A 433:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

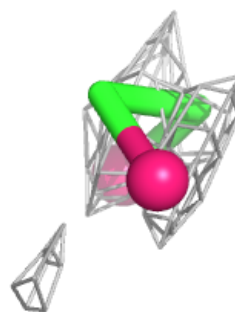
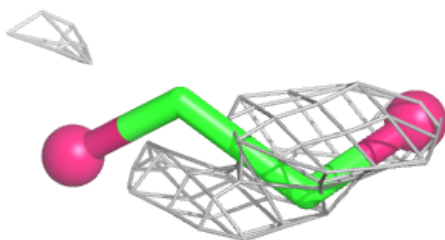
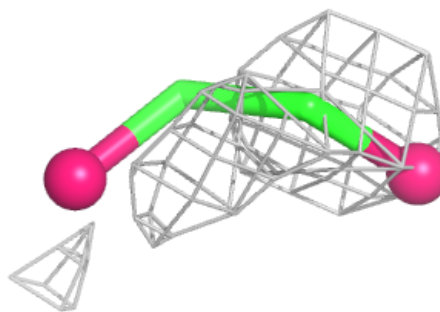
**Electron density around EDO C 437:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



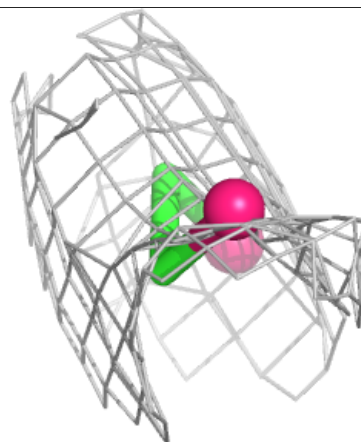
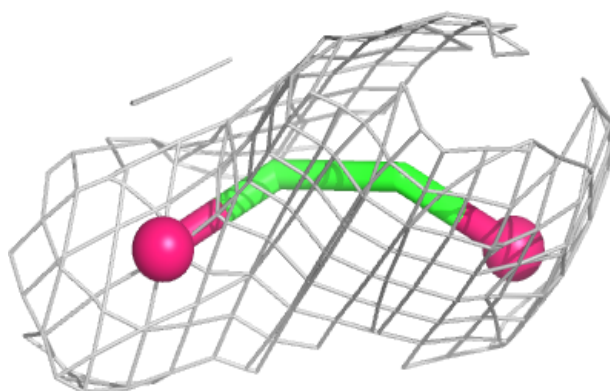
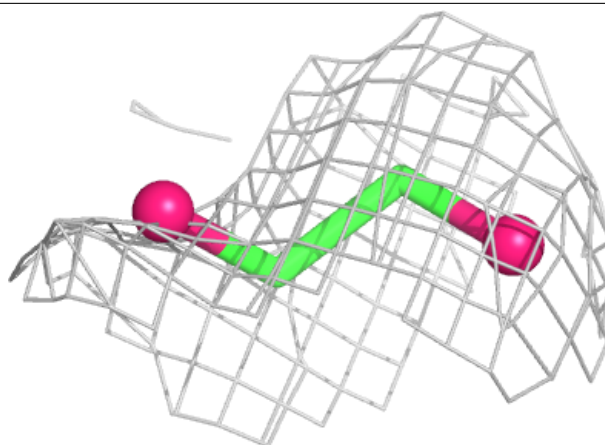
**Electron density around EDO B 429:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

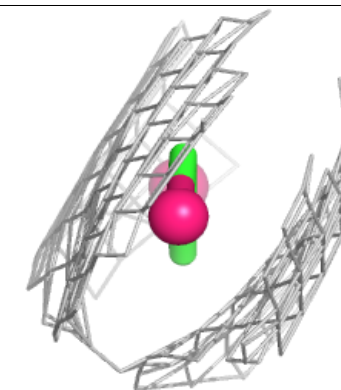
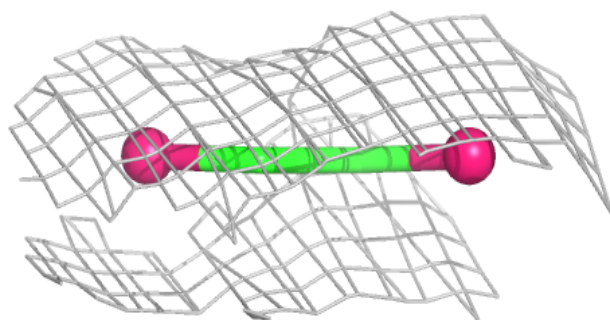
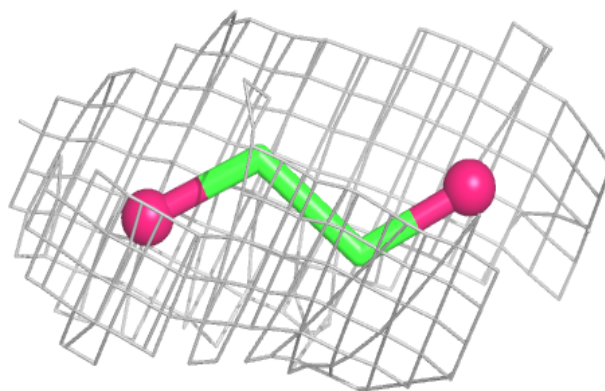


**Electron density around EDO E 447:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

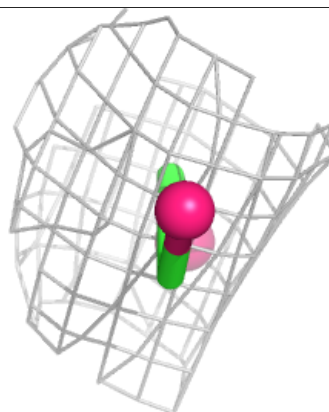
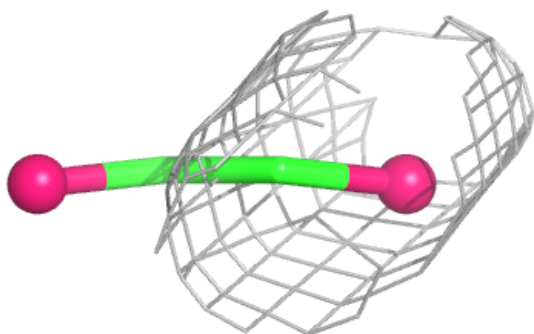
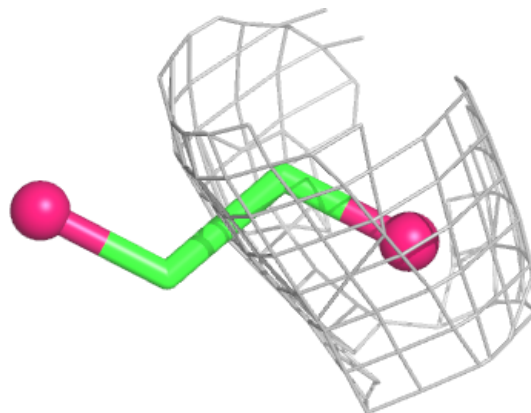
**Electron density around EDO D 401:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

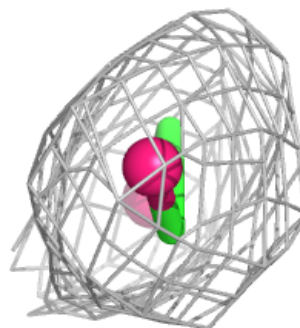
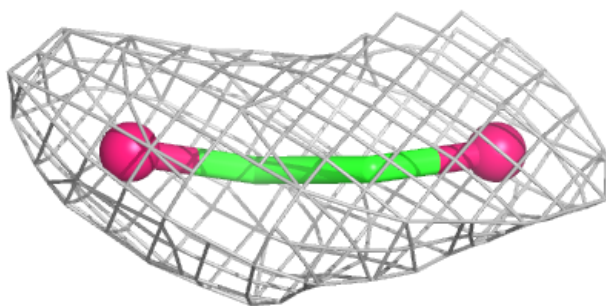
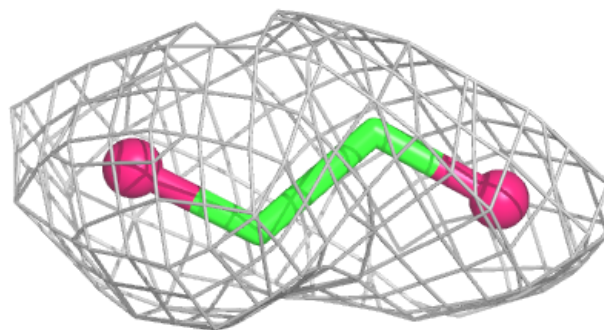


**Electron density around EDO A 419:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around EDO B 431:**

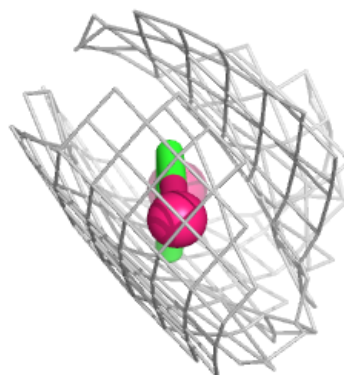
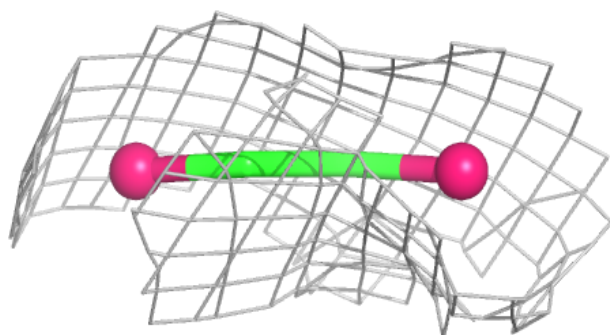
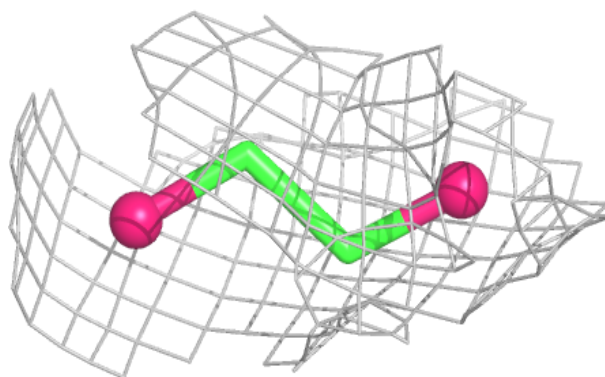
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



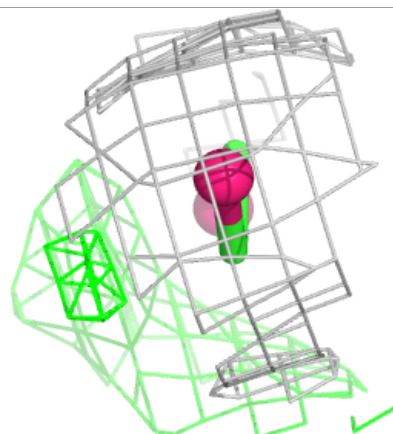
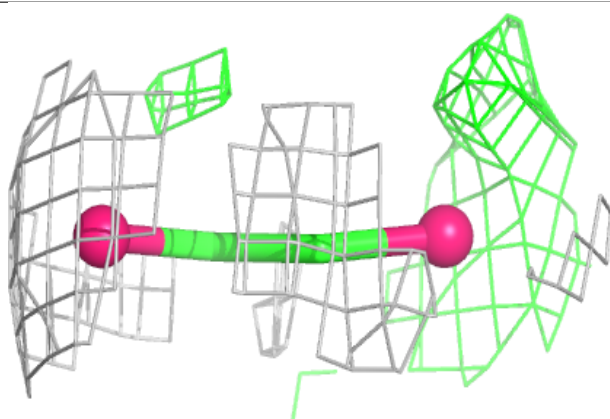
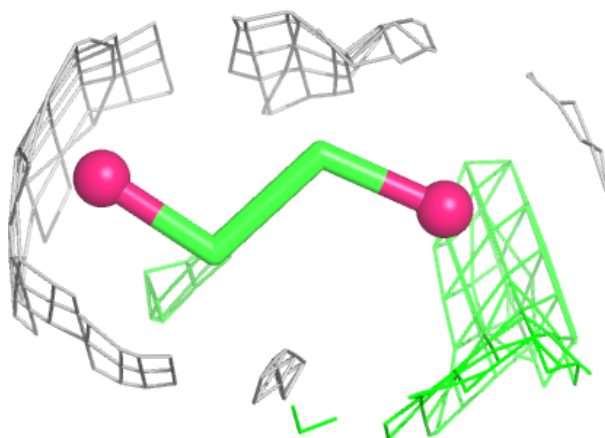


**Electron density around EDO F 407:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

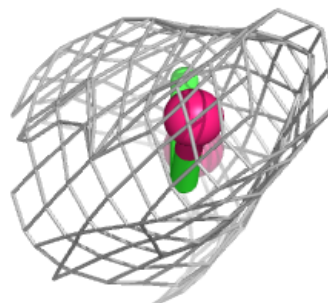
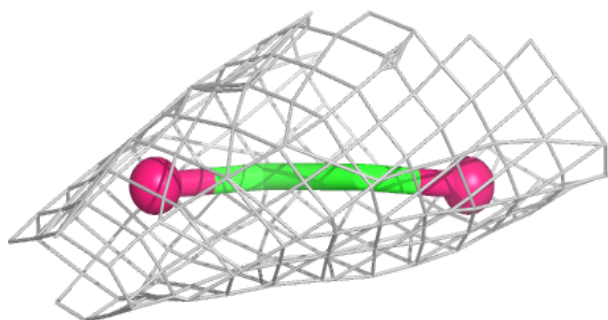
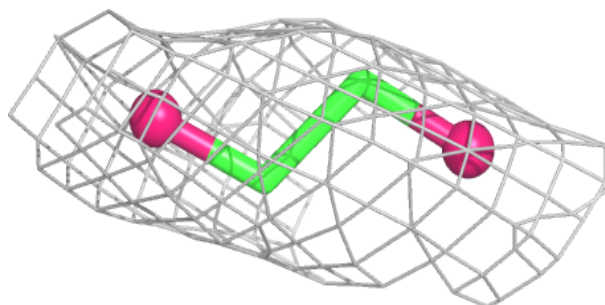
**Electron density around EDO A 408:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



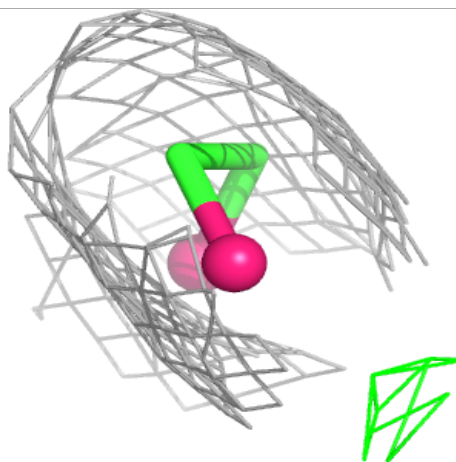
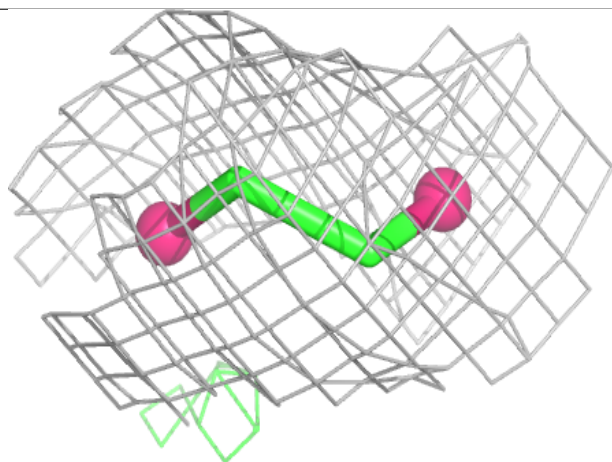
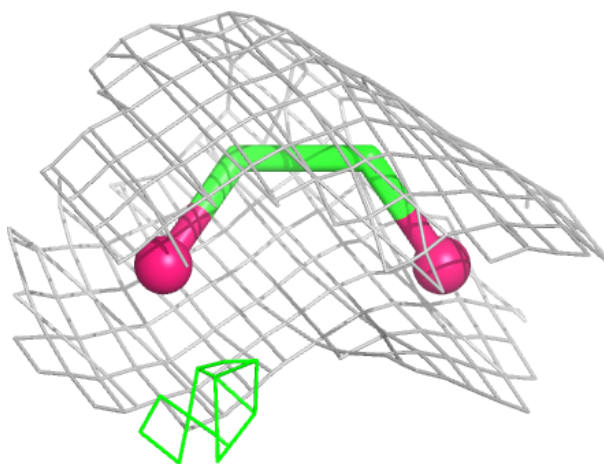
**Electron density around EDO B 435:**

$2mF_o - DF_c$  (at 0.7 rmsd) in gray  
 $mF_o - DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



**Electron density around EDO F 411:**

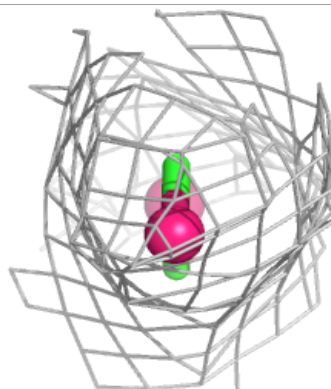
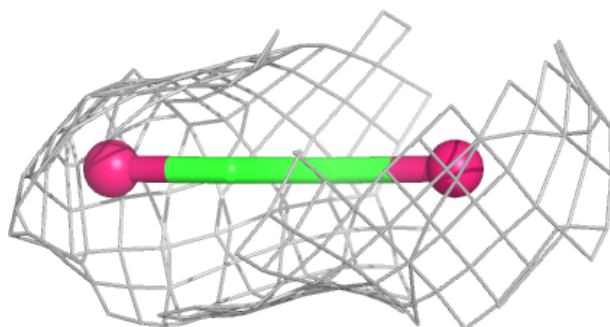
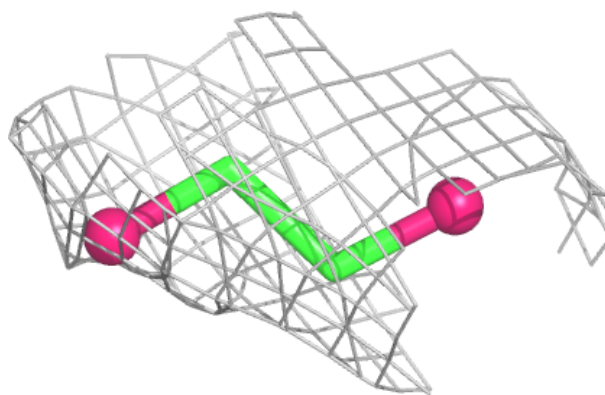
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



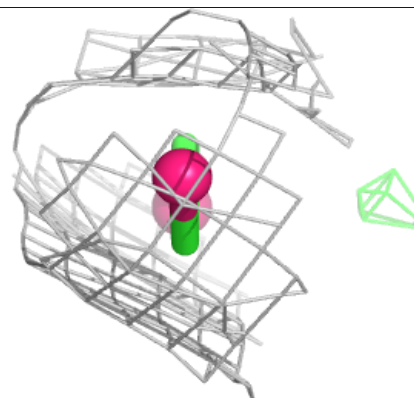
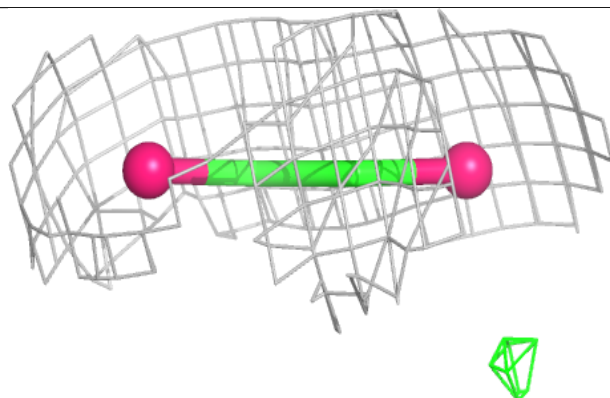
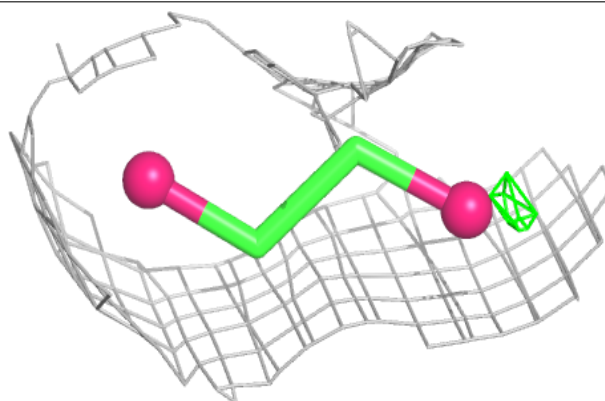


**Electron density around EDO A 410:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

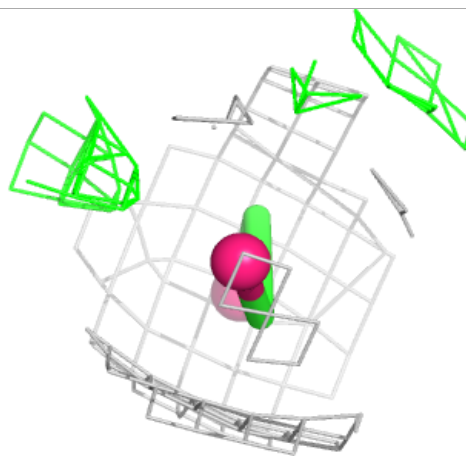
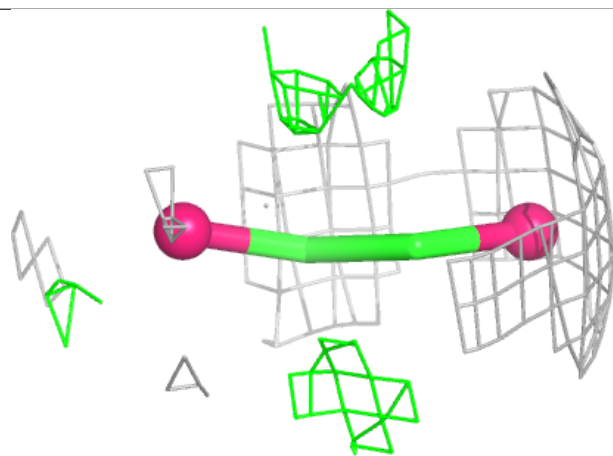
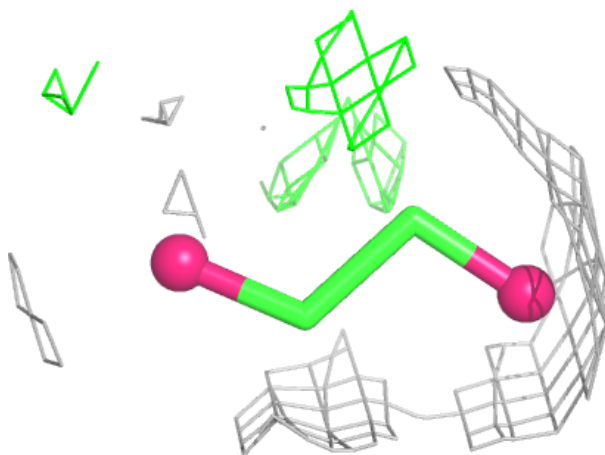
**Electron density around EDO D 408:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



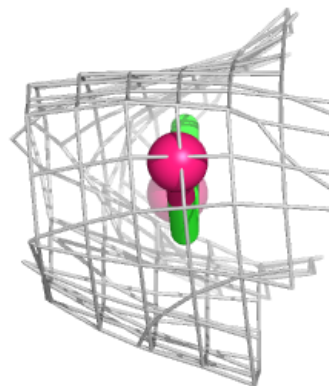
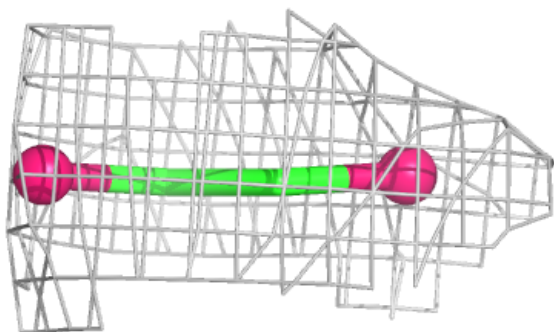
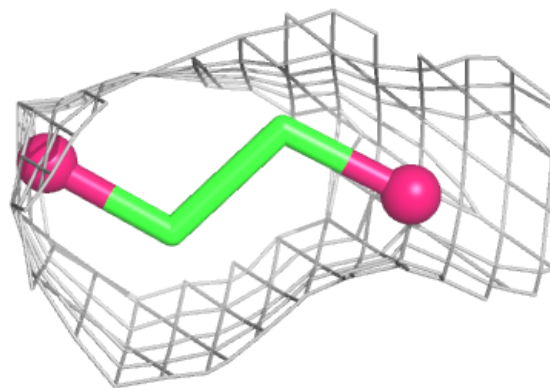
**Electron density around EDO D 411:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

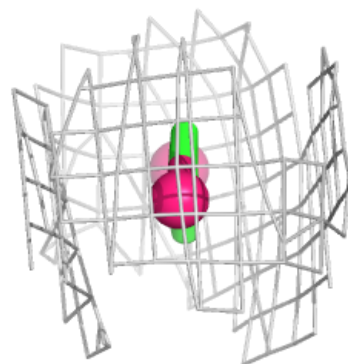
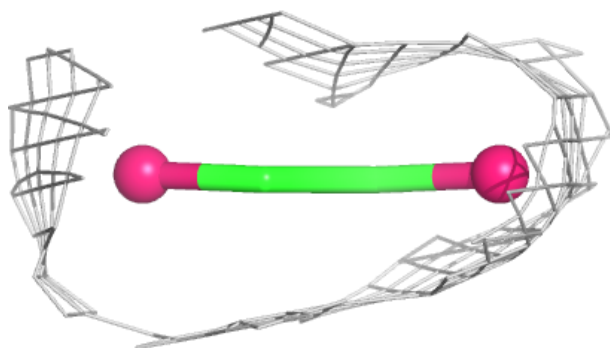
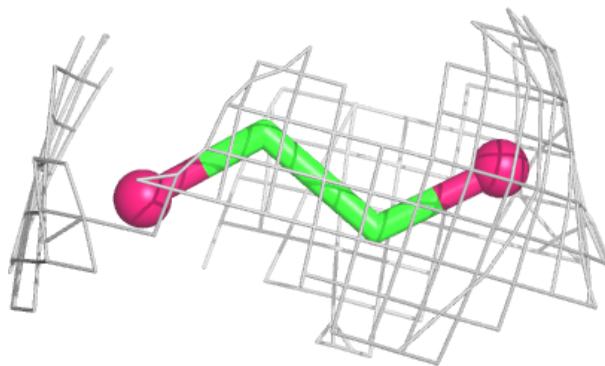


**Electron density around EDO D 415:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

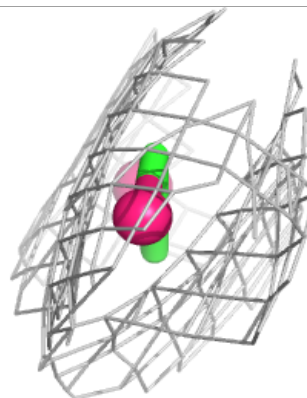
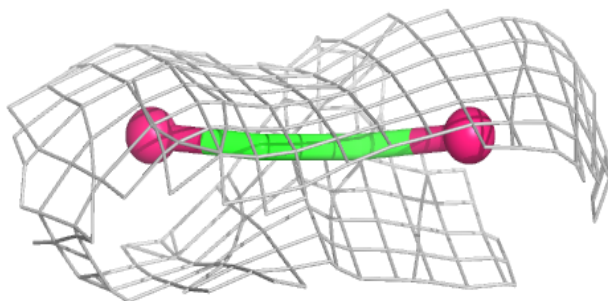
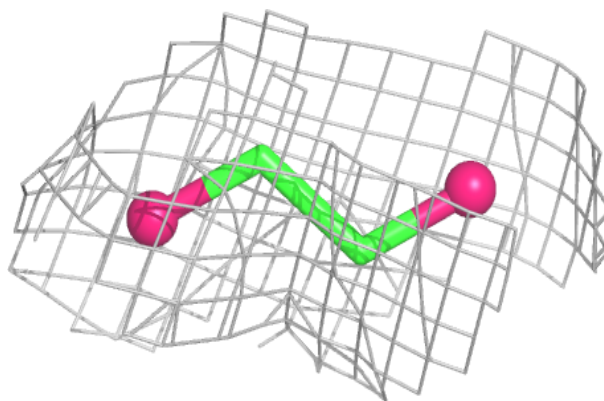
**Electron density around EDO D 418:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

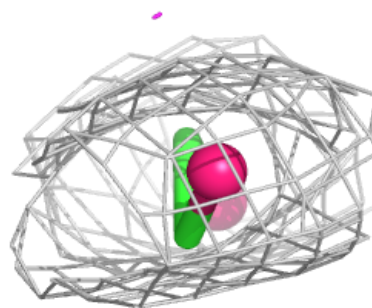
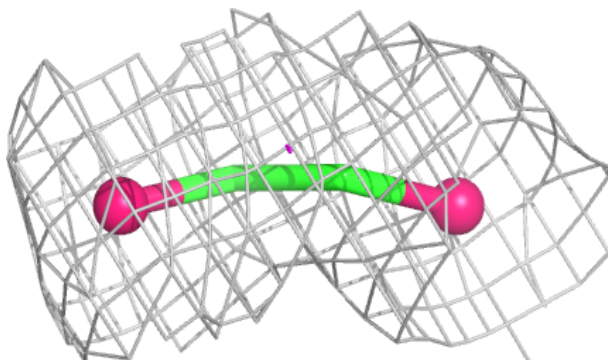
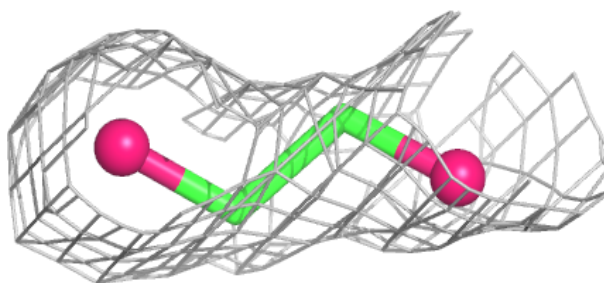


**Electron density around EDO D 419:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

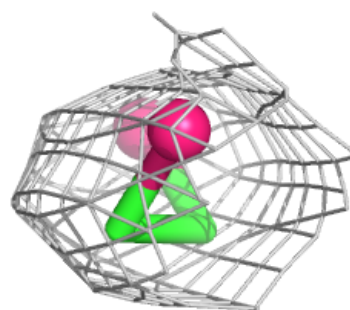
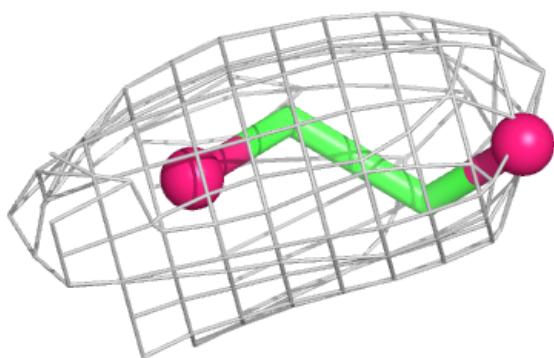
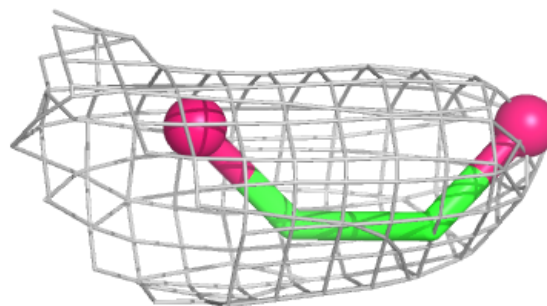
**Electron density around EDO A 411:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

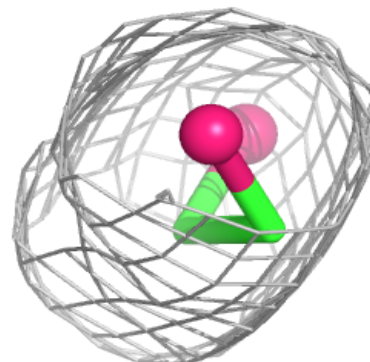
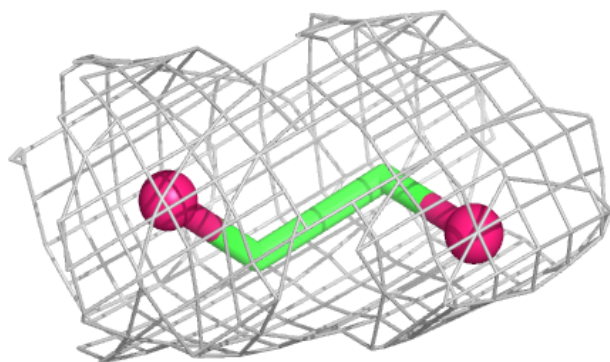
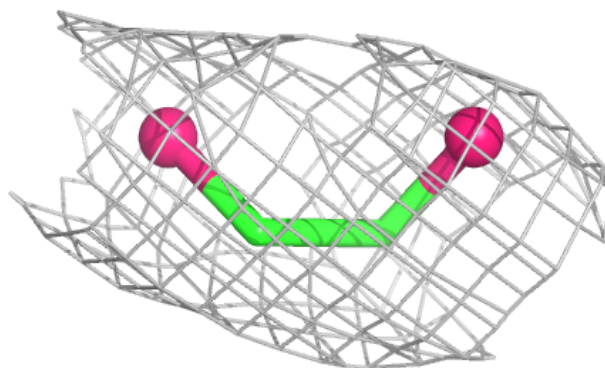


**Electron density around EDO B 439:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around EDO F 431:**

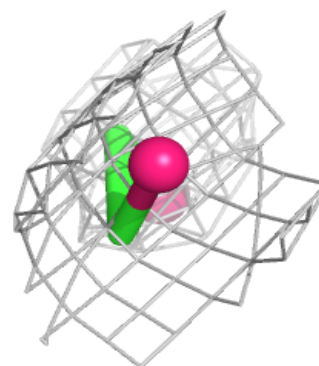
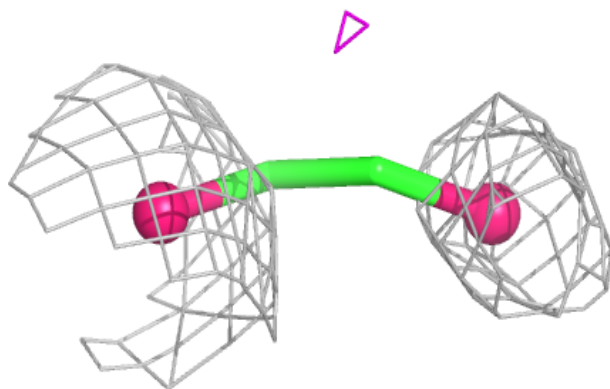
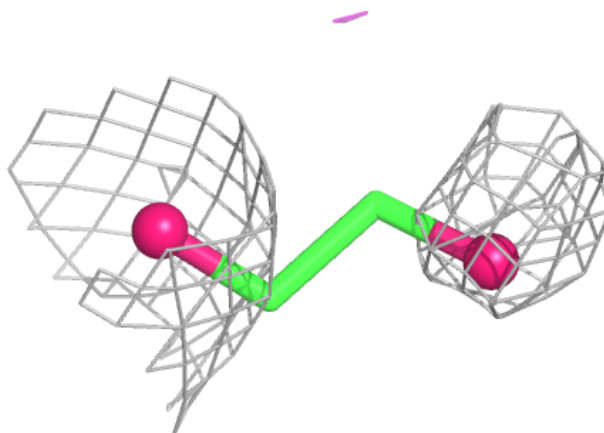
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





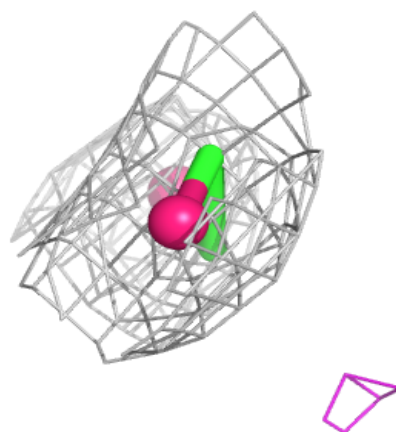
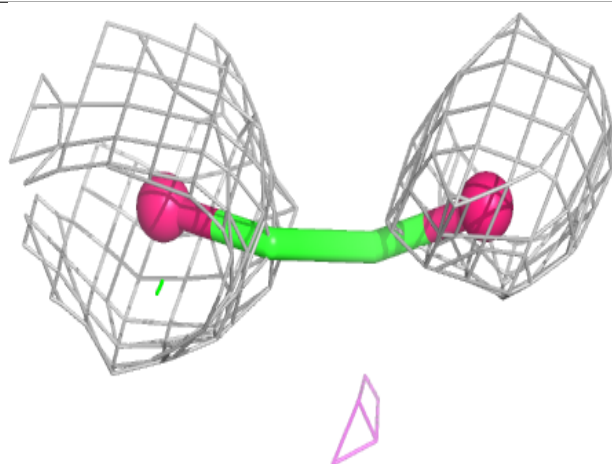
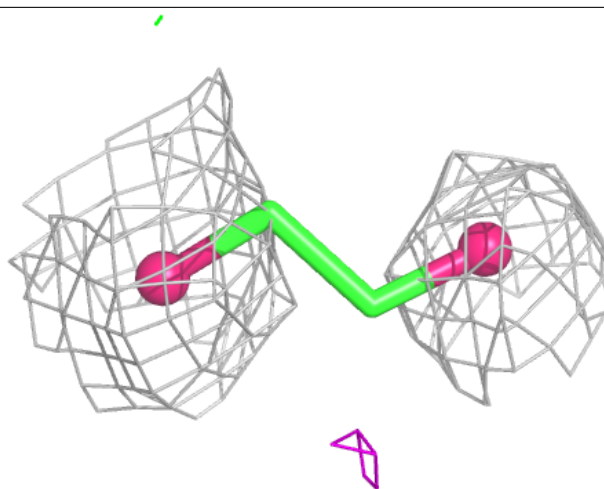
**Electron density around EDO F 432:**

$2mF_o - DF_c$  (at 0.7 rmsd) in gray  
 $mF_o - DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



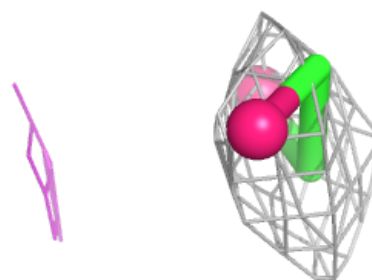
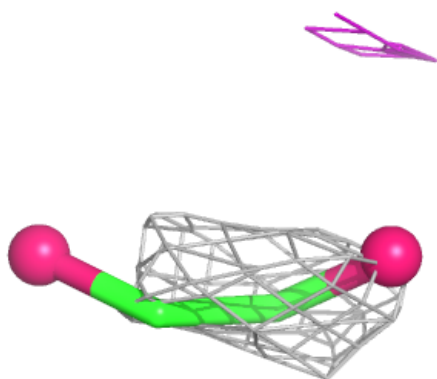
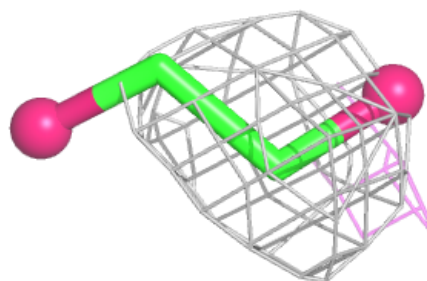
**Electron density around EDO A 424:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

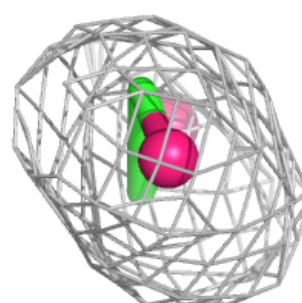
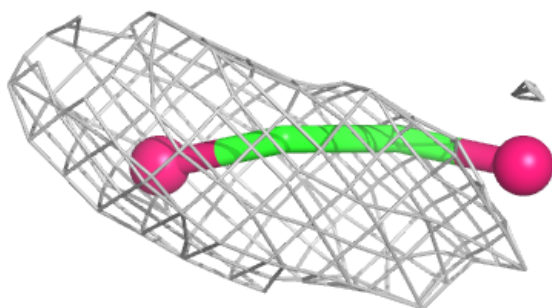
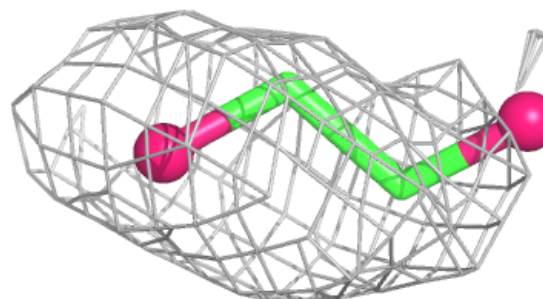


**Electron density around EDO F 435:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around EDO B 442:**

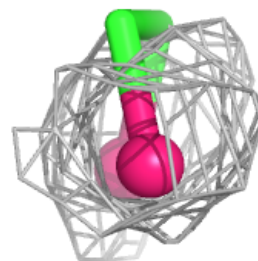
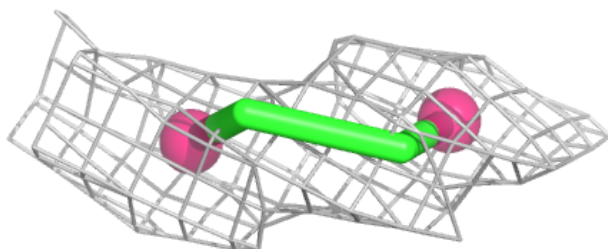
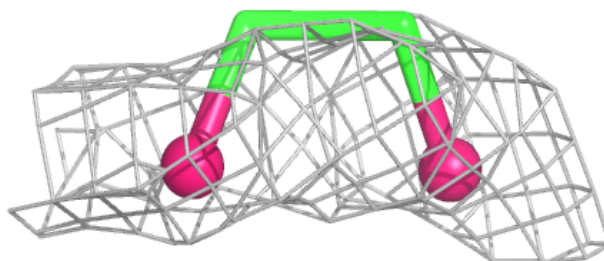
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



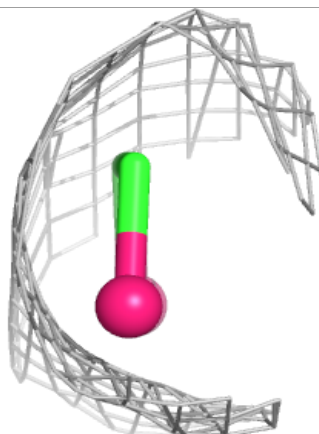
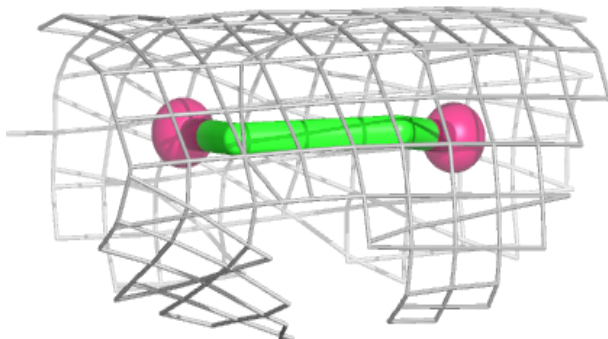
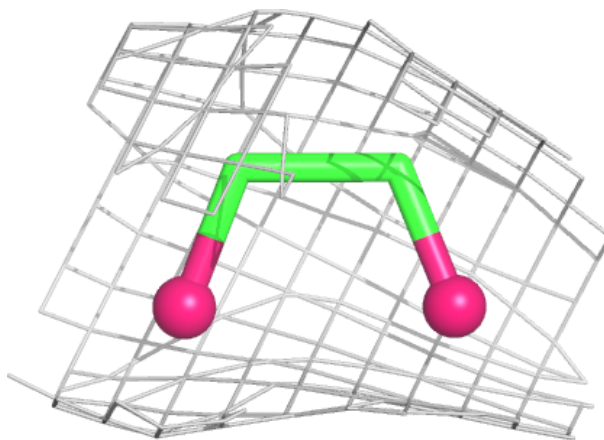


**Electron density around EDO F 437:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

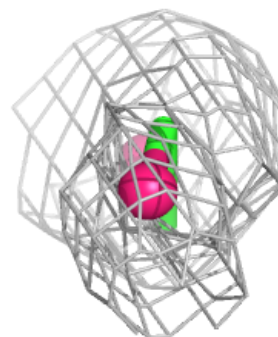
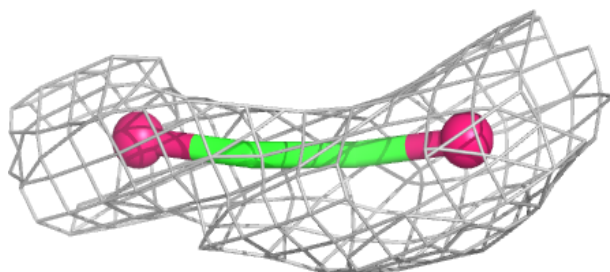
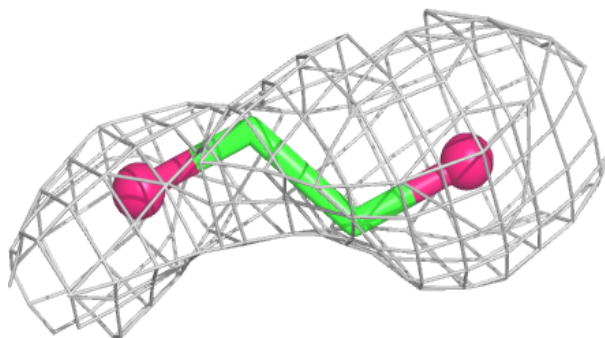
**Electron density around EDO F 438:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

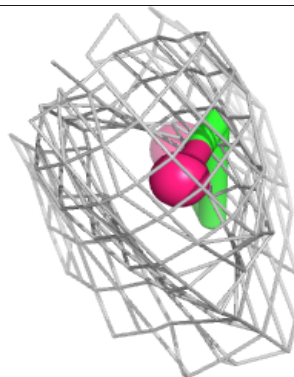
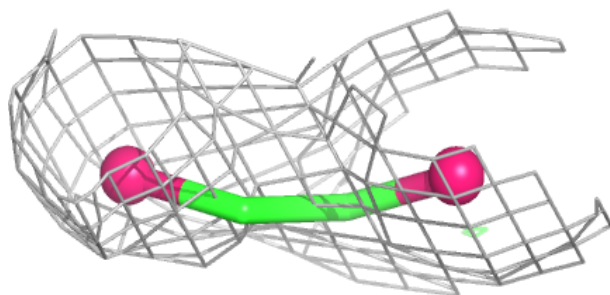
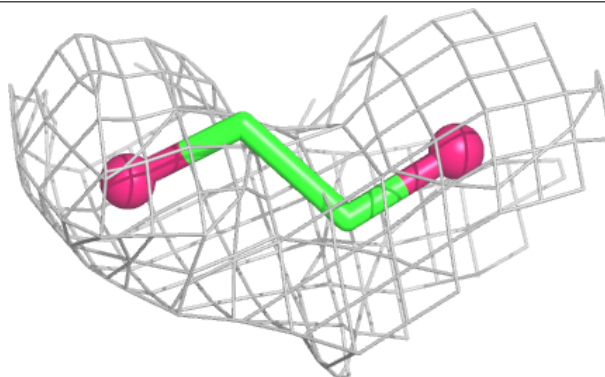


**Electron density around EDO A 404:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

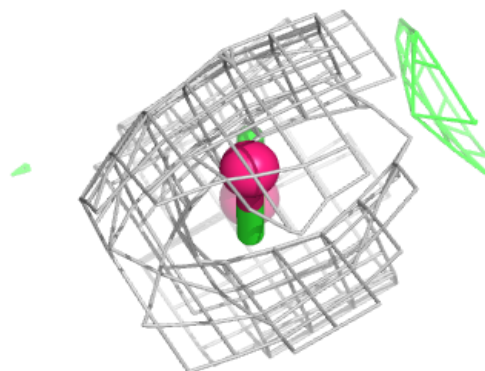
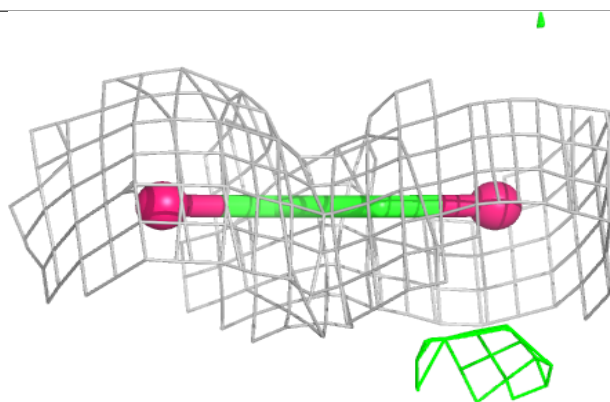
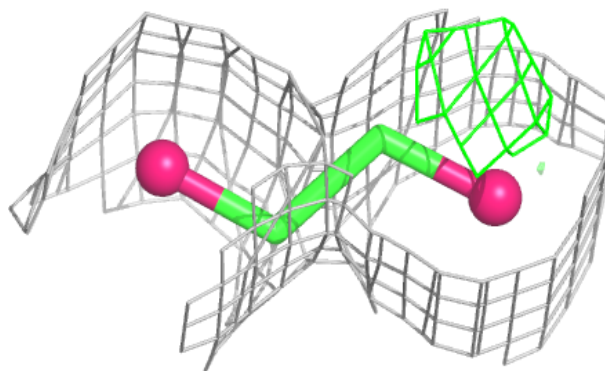
**Electron density around EDO D 427:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

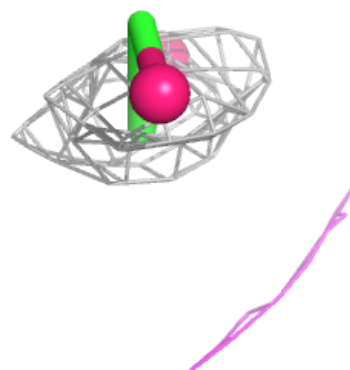
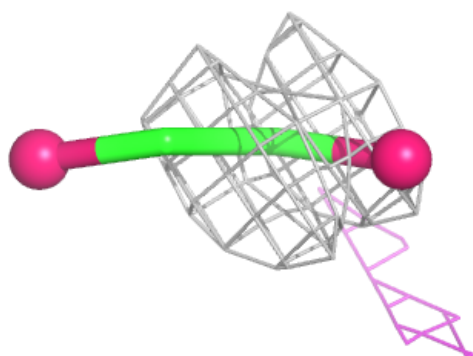
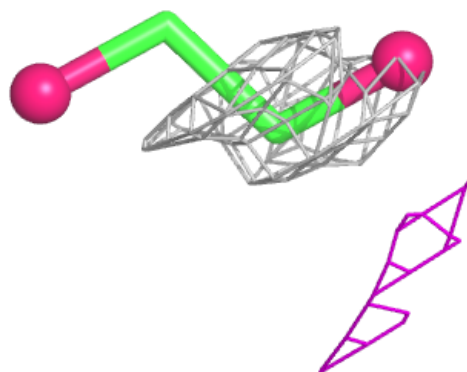


**Electron density around EDO A 413:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

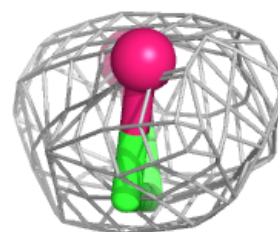
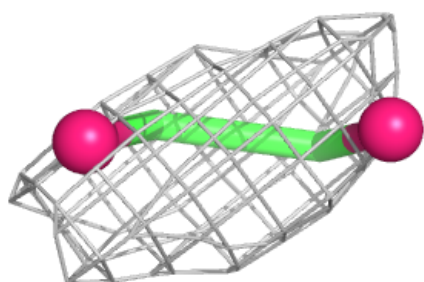
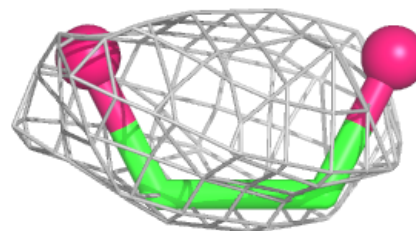
**Electron density around EDO D 430:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

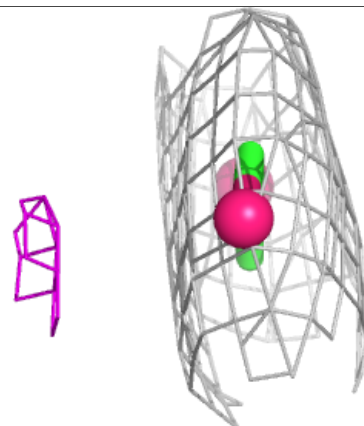
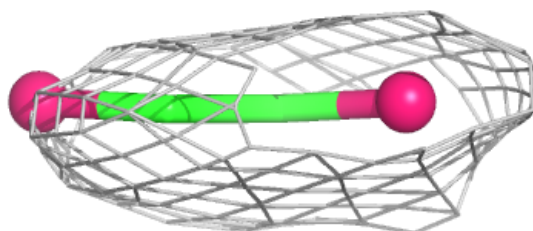
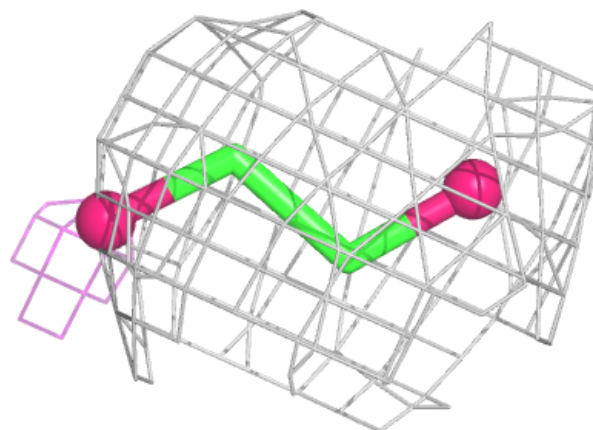


**Electron density around EDO D 431:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around EDO B 402:**

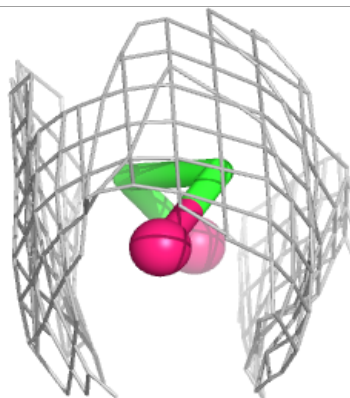
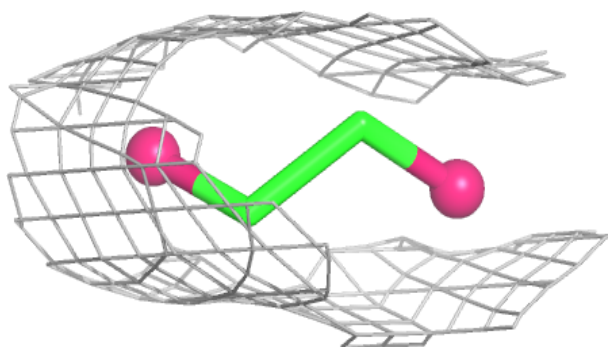
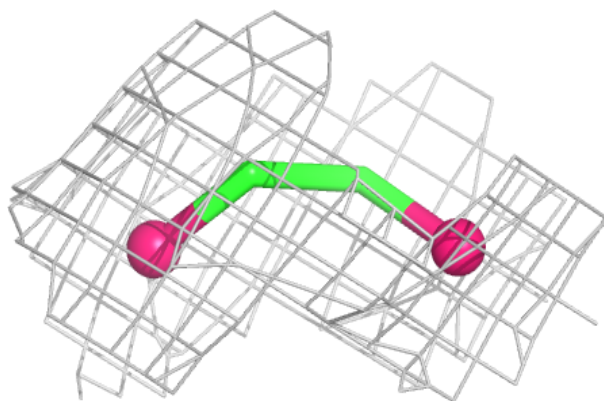
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



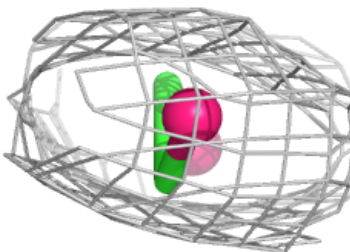
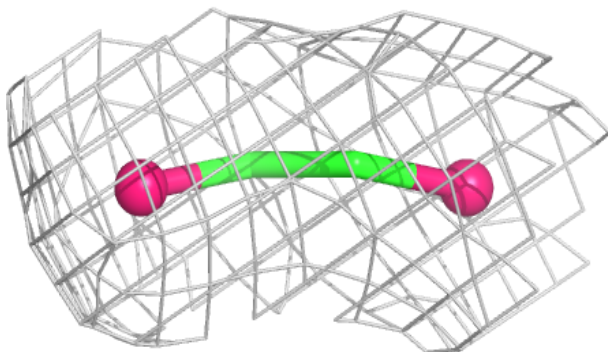
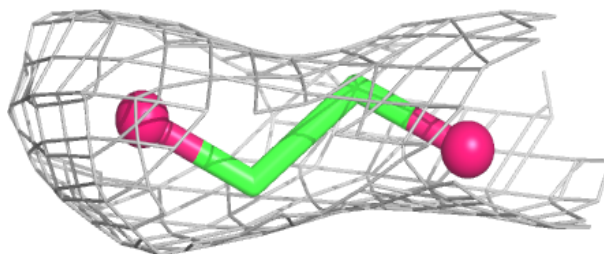


**Electron density around EDO A 406:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

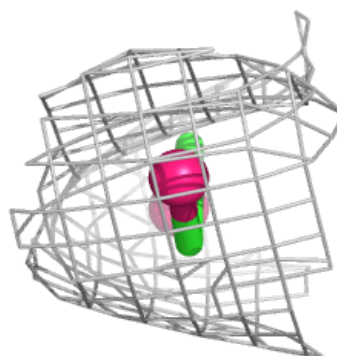
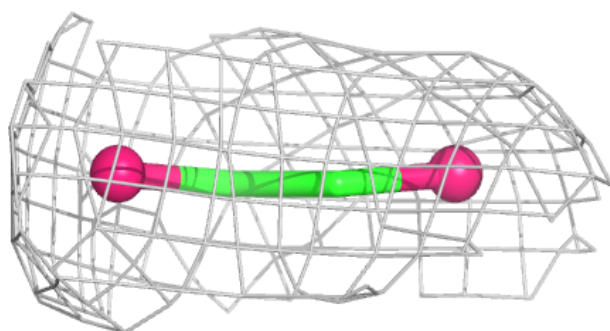
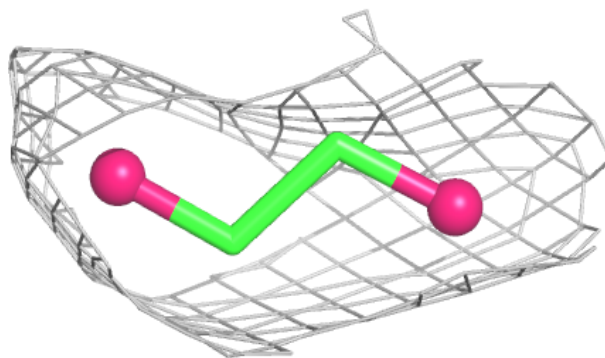
**Electron density around EDO E 414:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

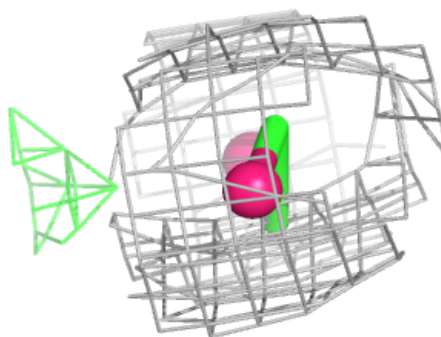
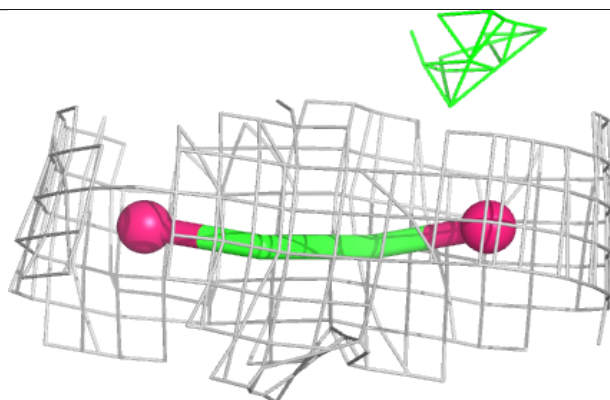
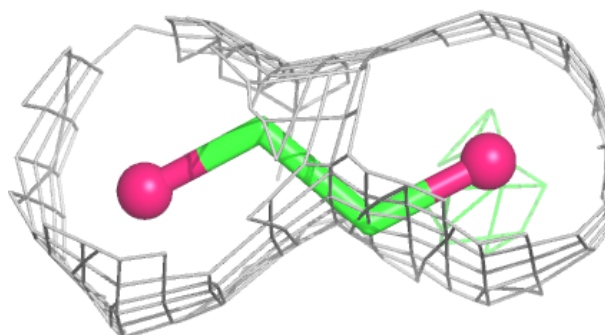


**Electron density around EDO E 415:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

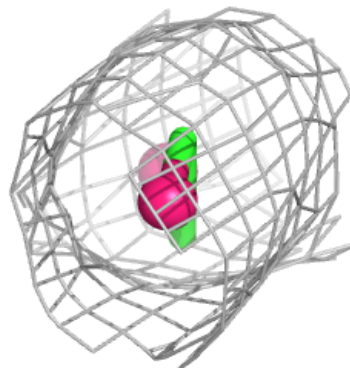
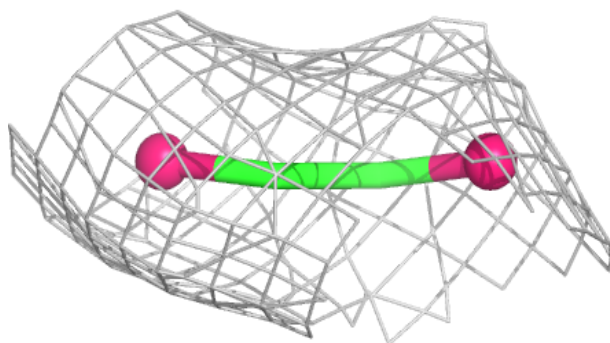
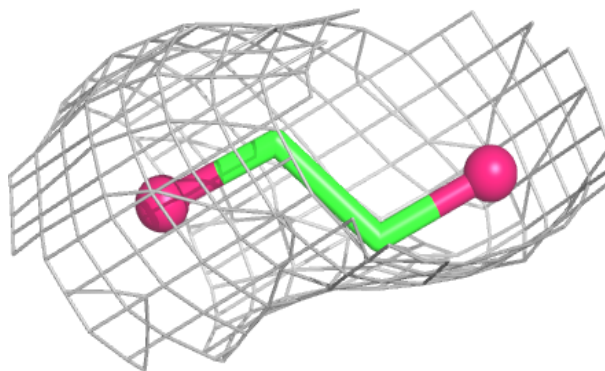
**Electron density around EDO E 416:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

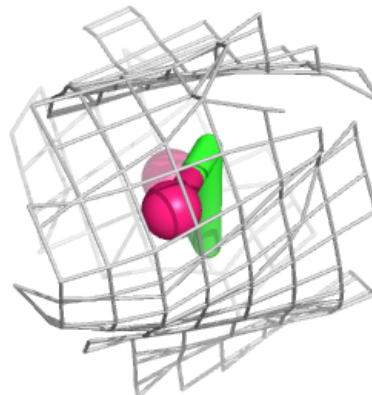
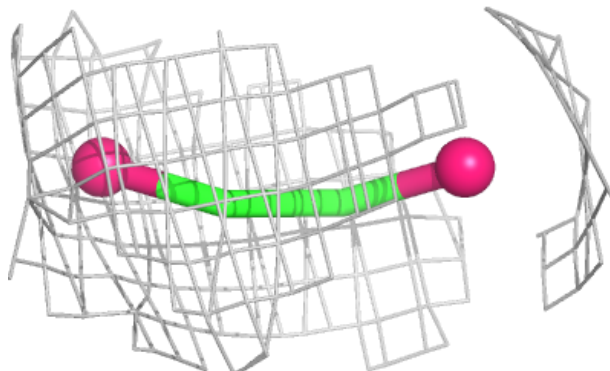
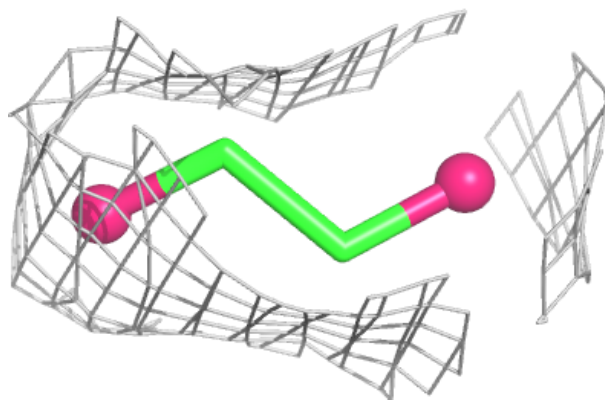


**Electron density around EDO E 417:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

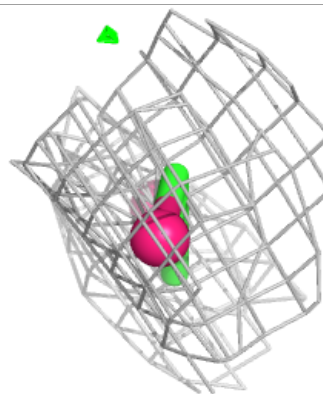
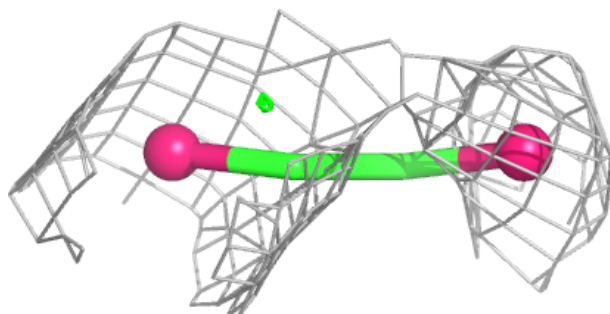
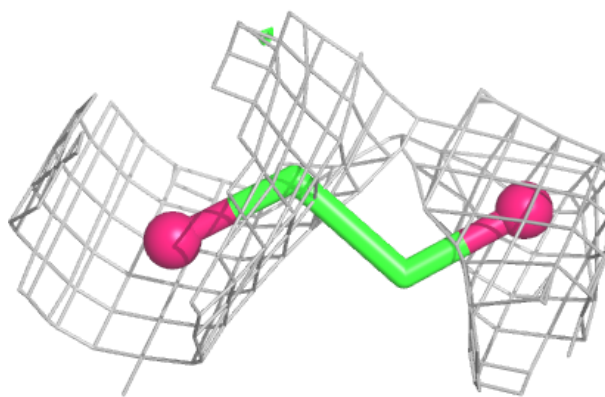
**Electron density around EDO E 418:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

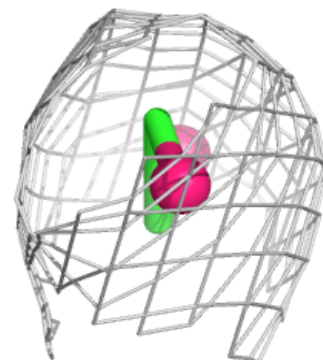
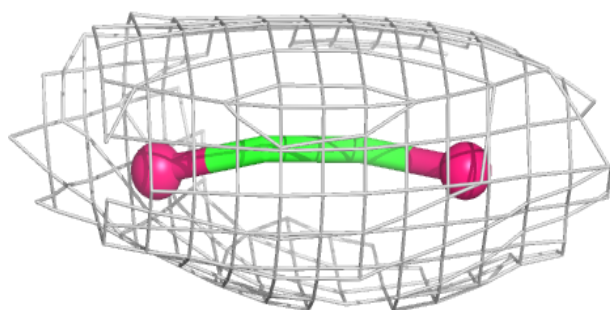
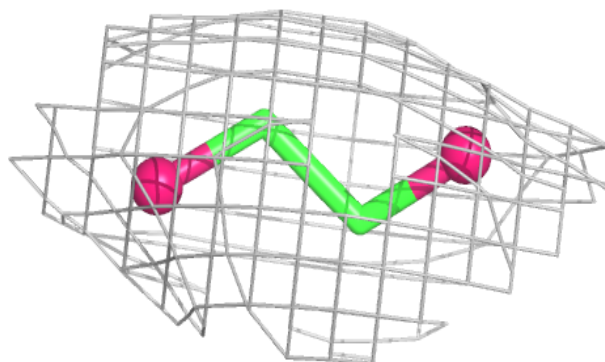


**Electron density around EDO B 421:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around EDO E 420:**

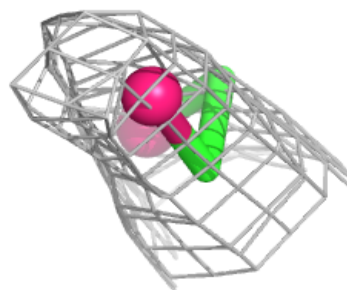
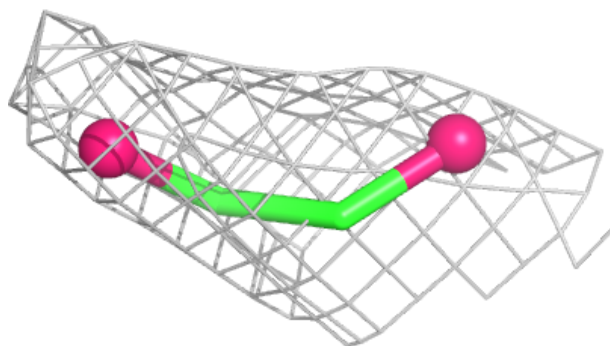
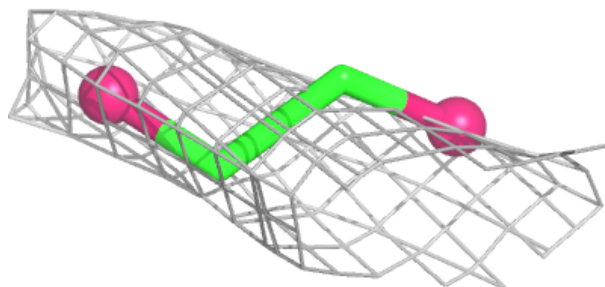
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



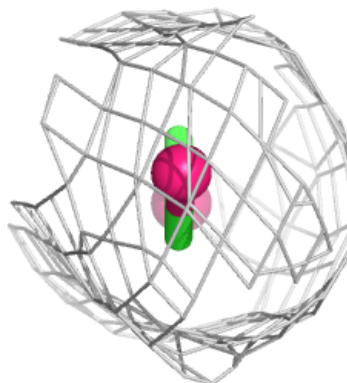
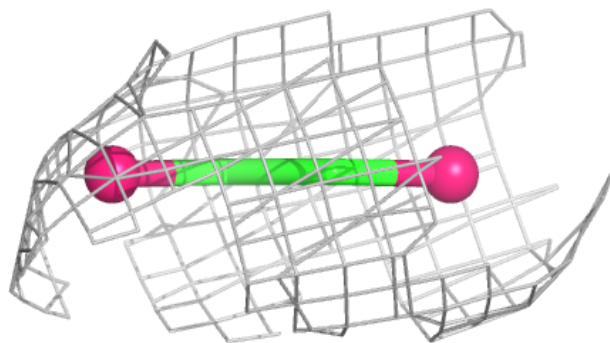
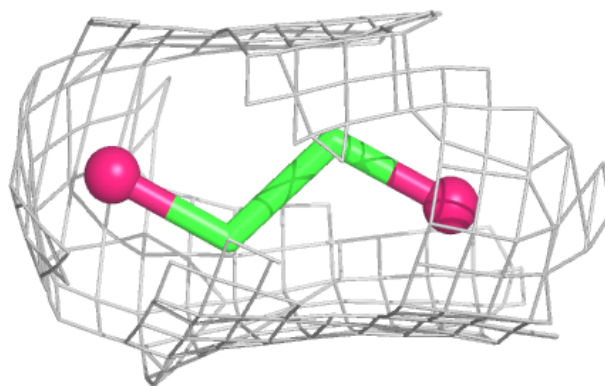


**Electron density around EDO B 445:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

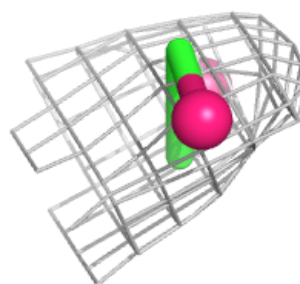
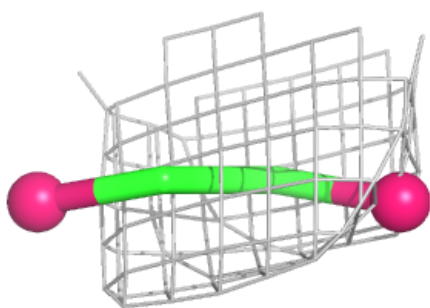
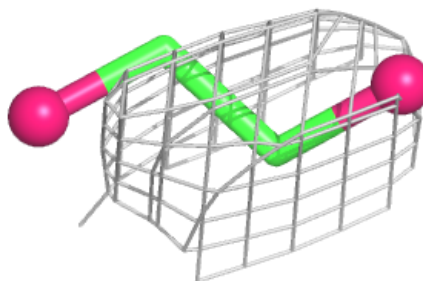
**Electron density around EDO E 422:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

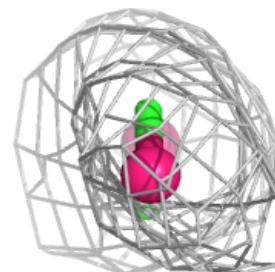
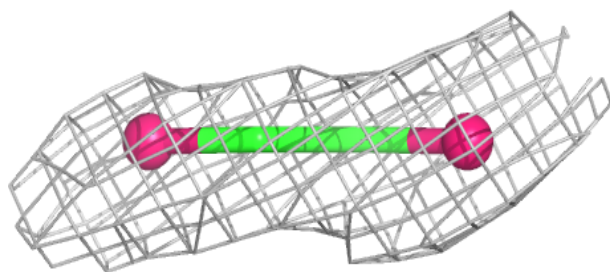
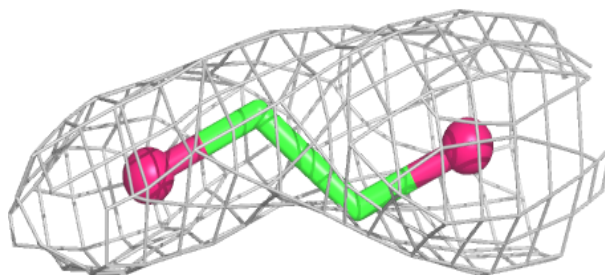


**Electron density around EDO A 402:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

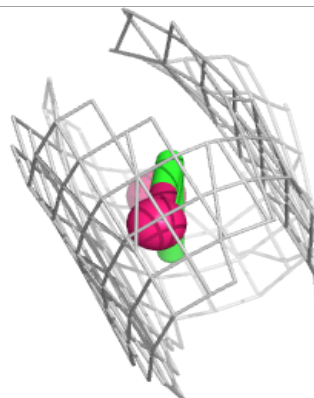
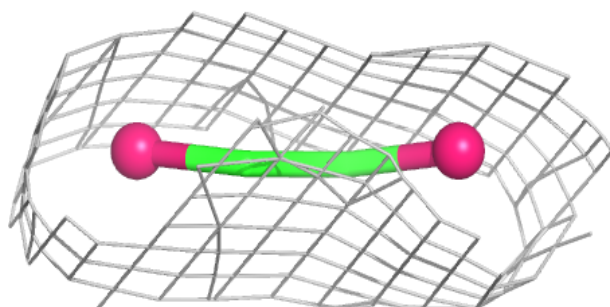
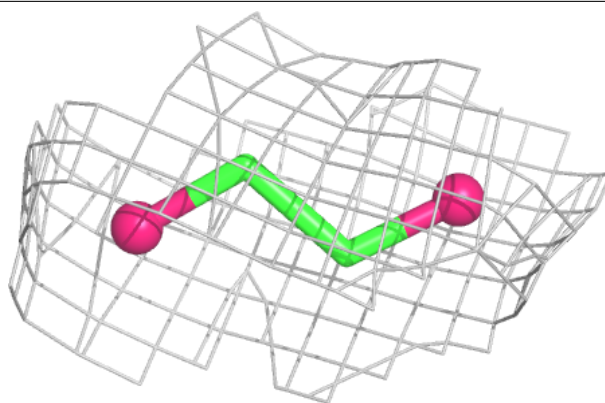
**Electron density around EDO B 406:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

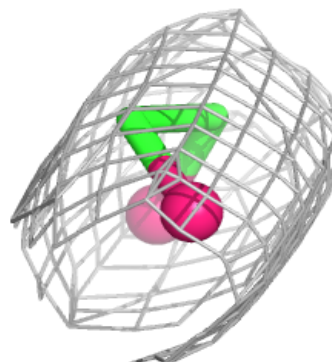
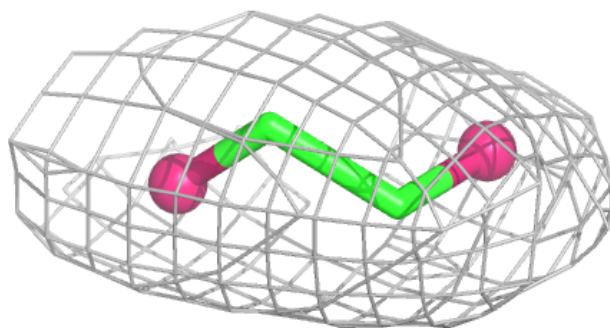
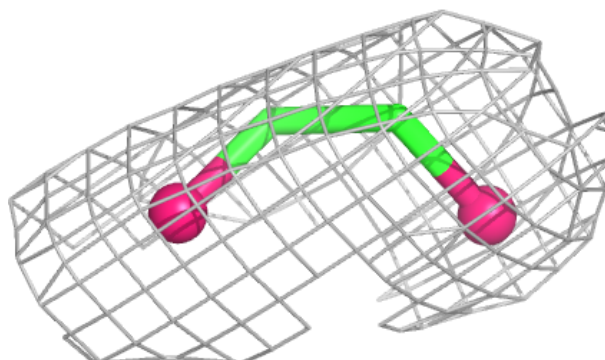


**Electron density around EDO C 401:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

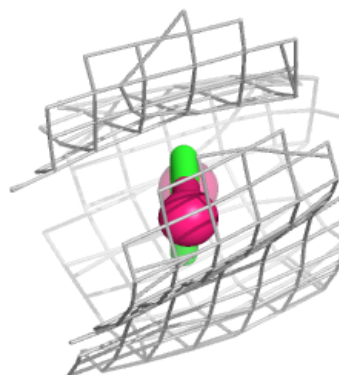
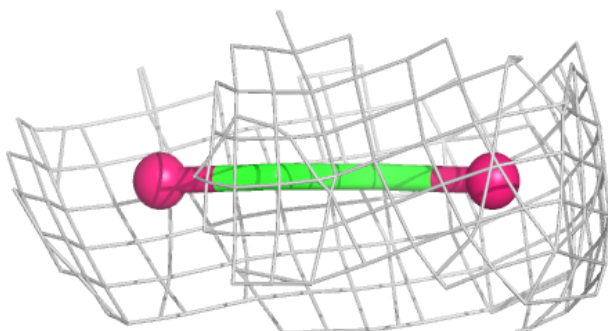
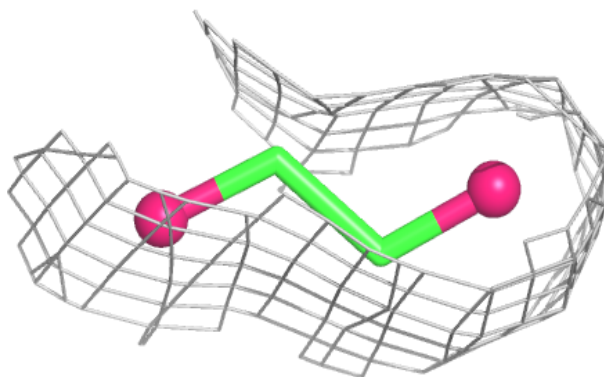
**Electron density around EDO B 425:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

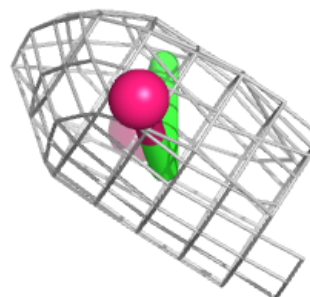
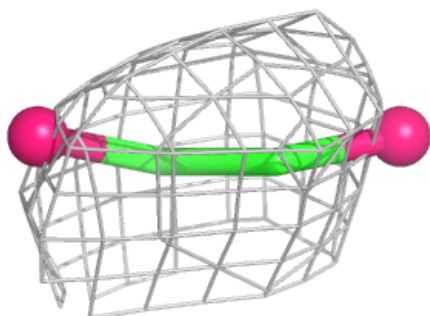
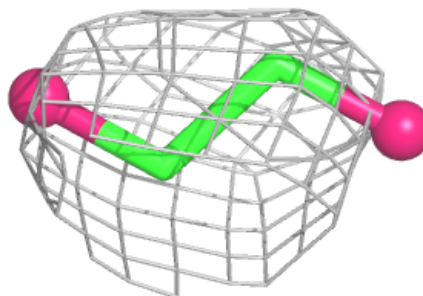


**Electron density around EDO B 407:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around EDO C 404:**

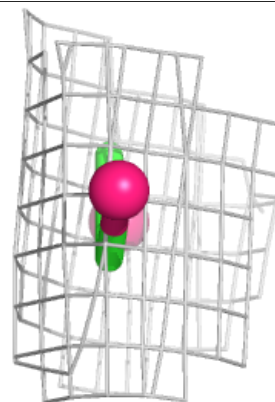
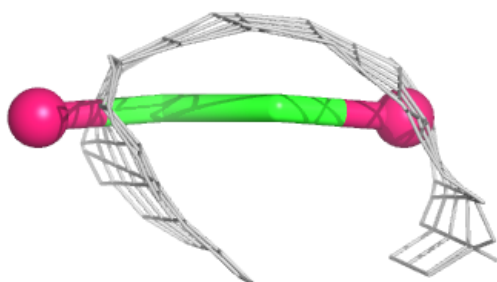
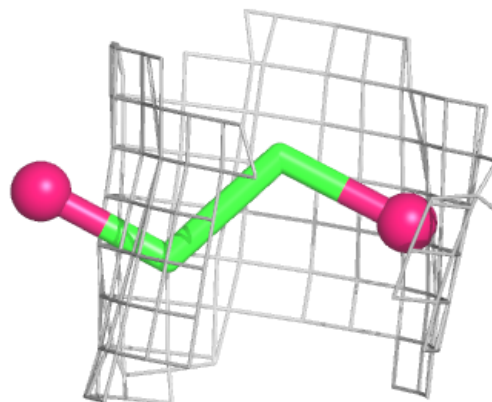
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



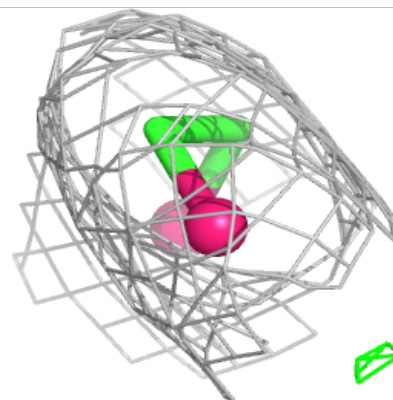
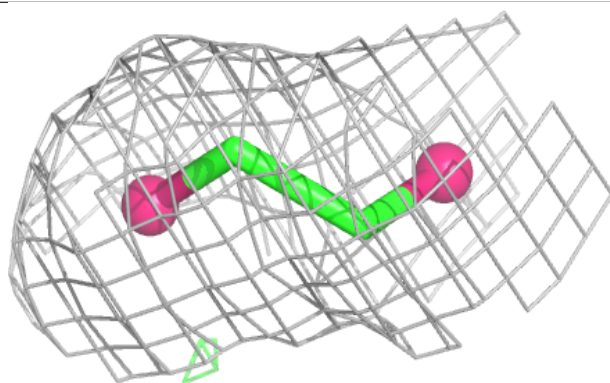
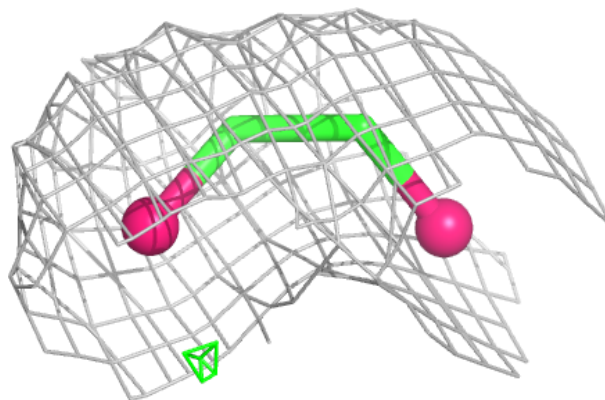


**Electron density around EDO E 430:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

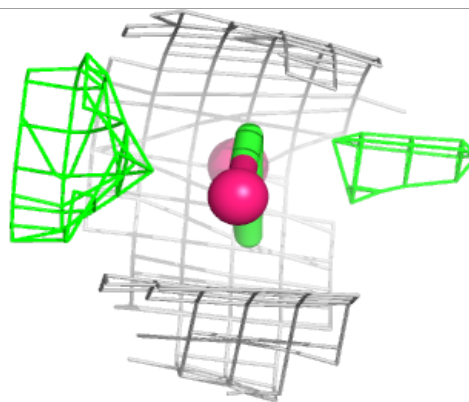
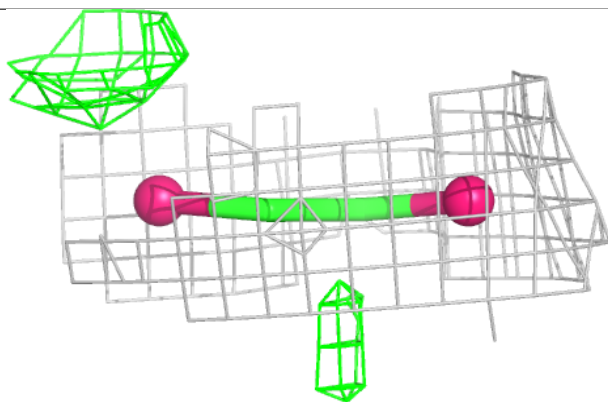
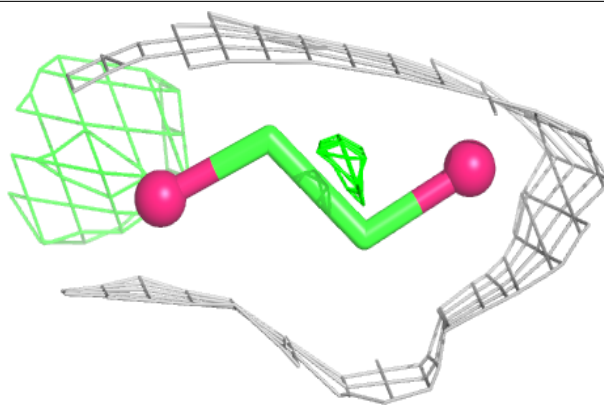
**Electron density around EDO C 405:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

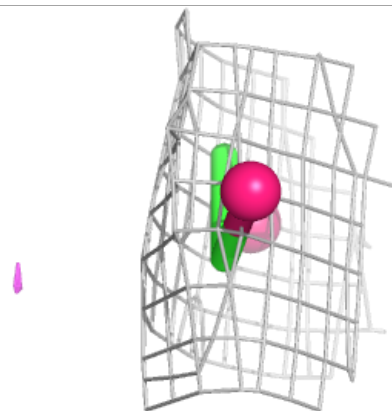
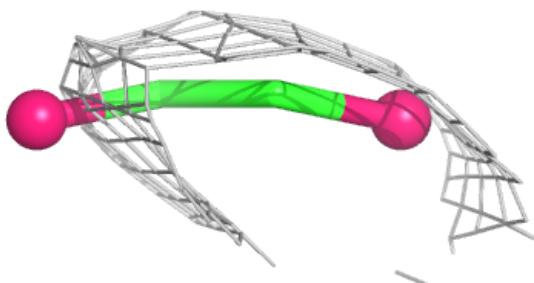
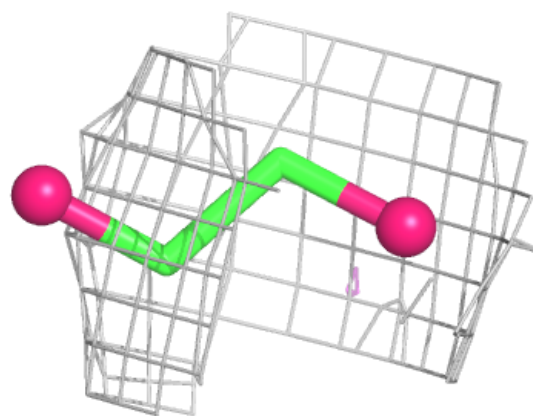


**Electron density around EDO D 410:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

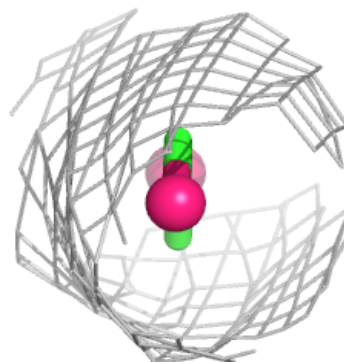
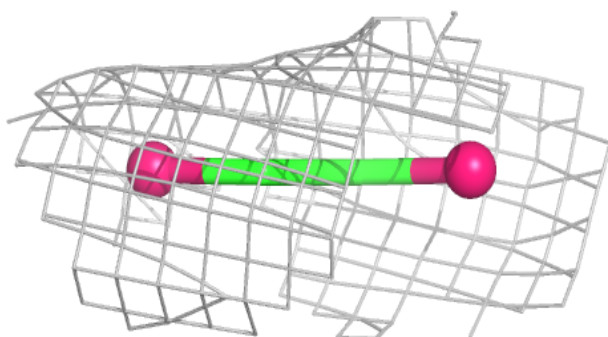
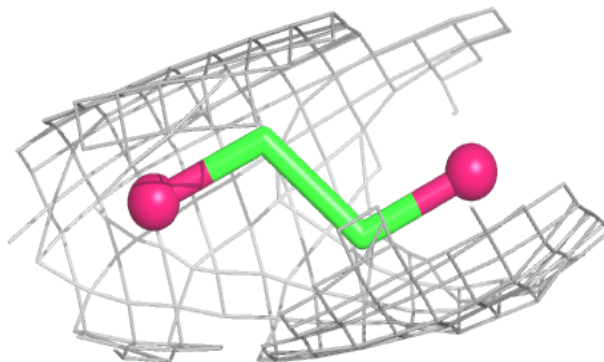
**Electron density around EDO B 427:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

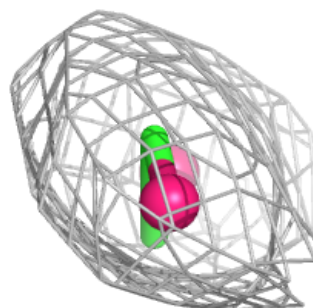
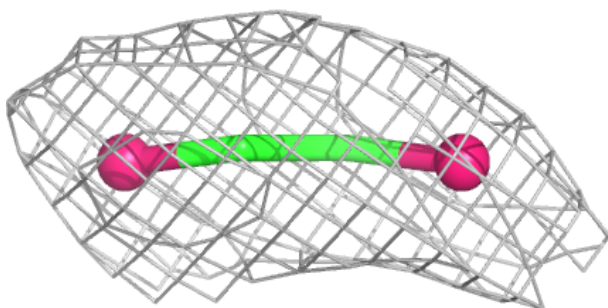
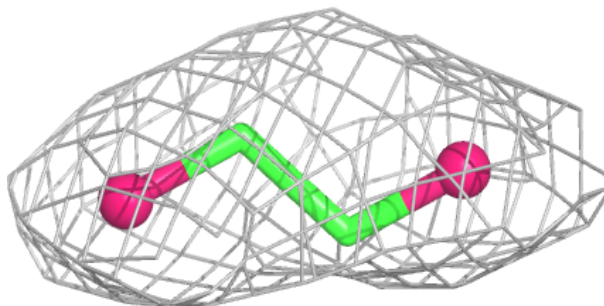


**Electron density around EDO D 412:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

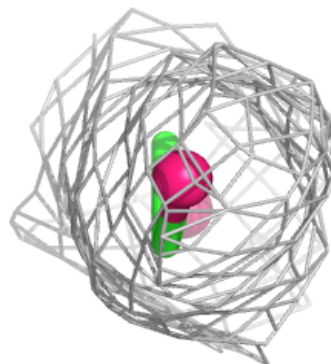
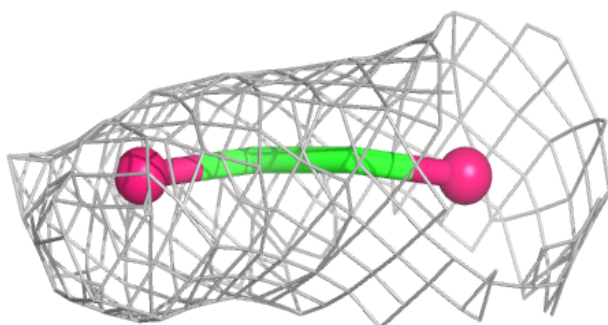
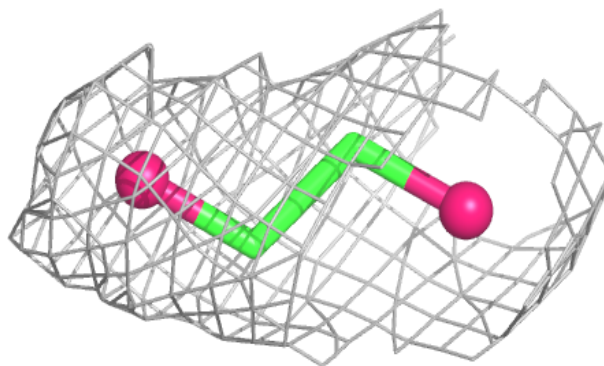
**Electron density around EDO E 435:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

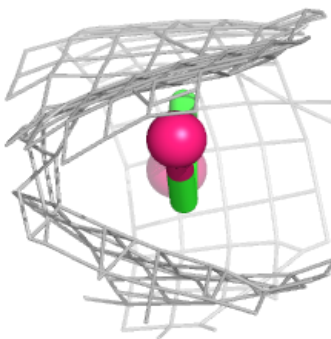
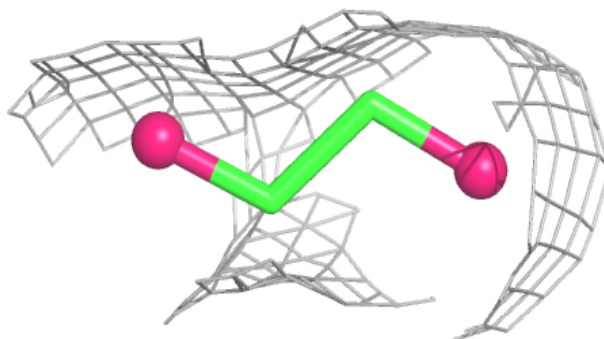


**Electron density around EDO D 413:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around EDO C 407:**

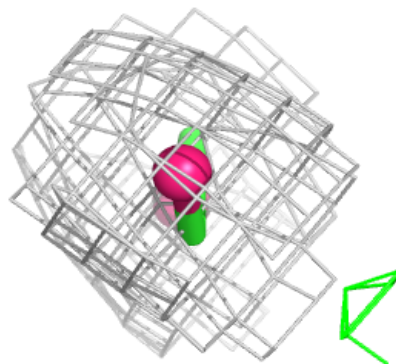
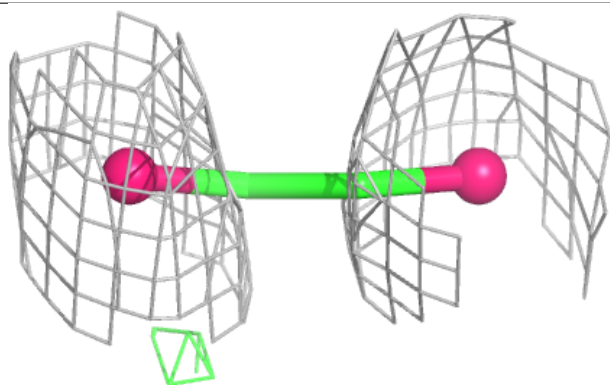
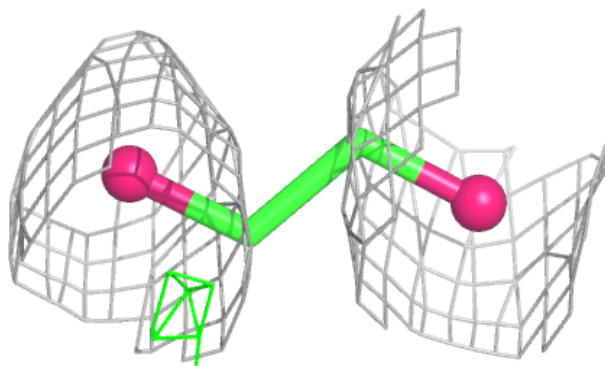
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



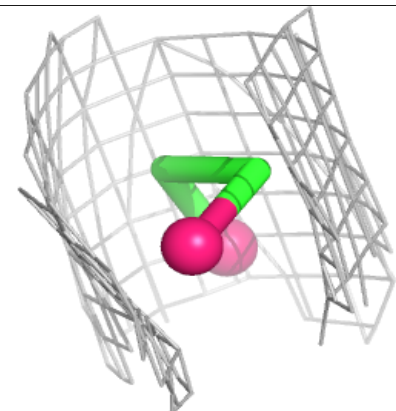
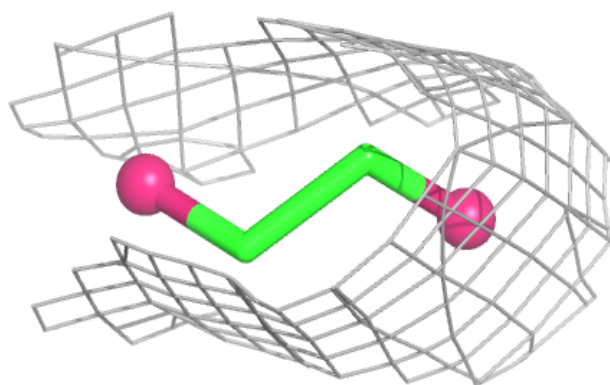
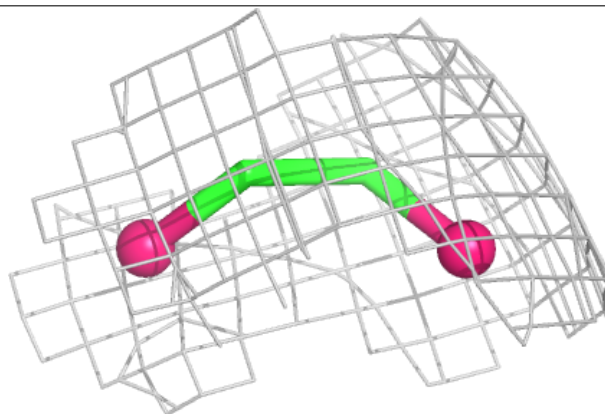


**Electron density around EDO D 416:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

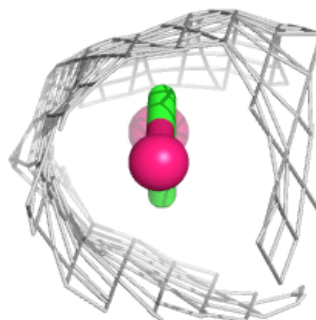
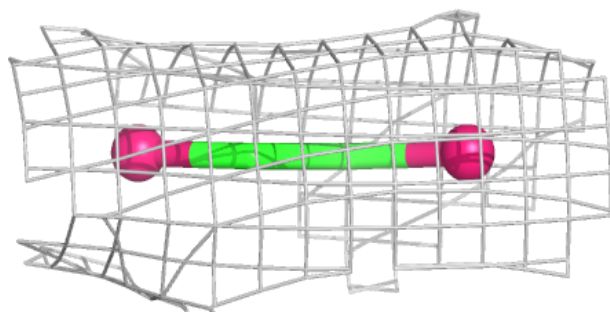
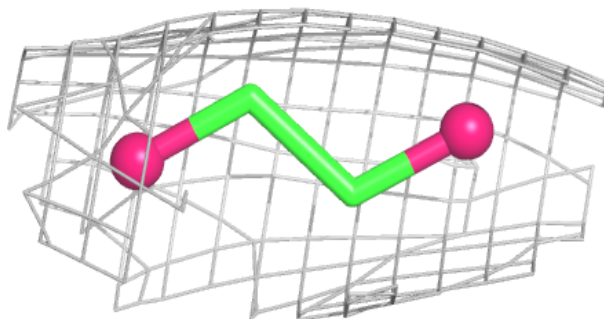
**Electron density around EDO C 408:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

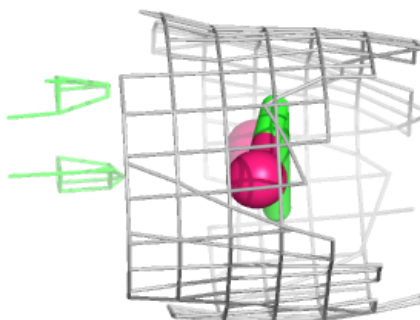
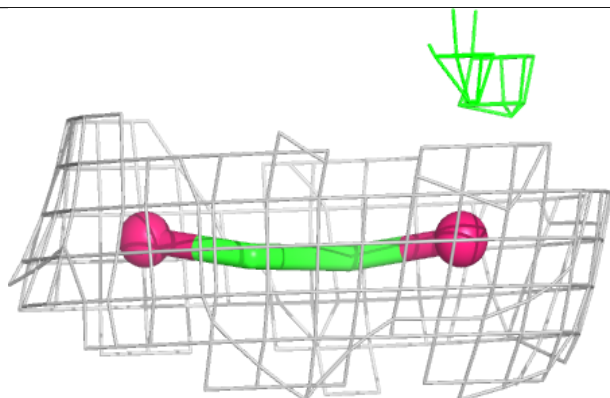
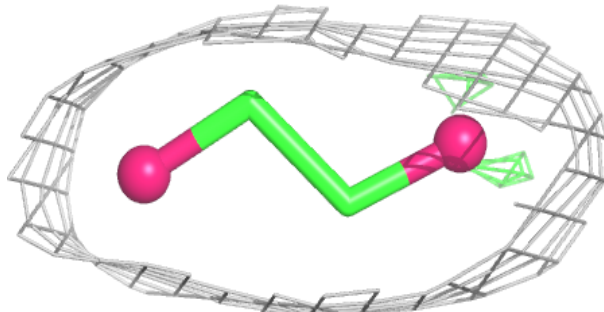


**Electron density around EDO E 442:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

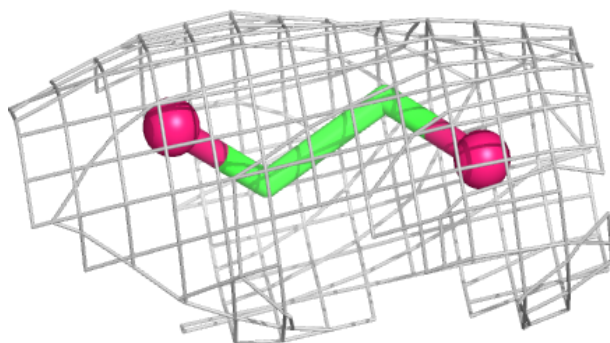
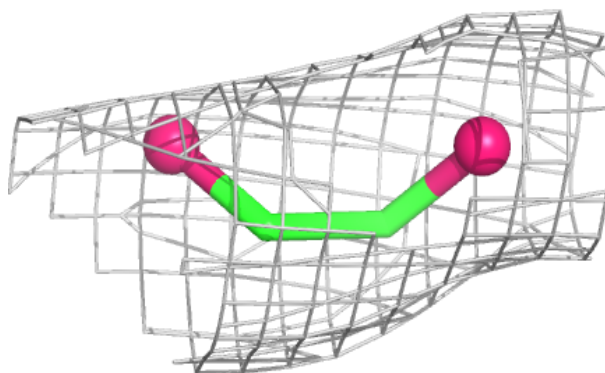
**Electron density around EDO C 409:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

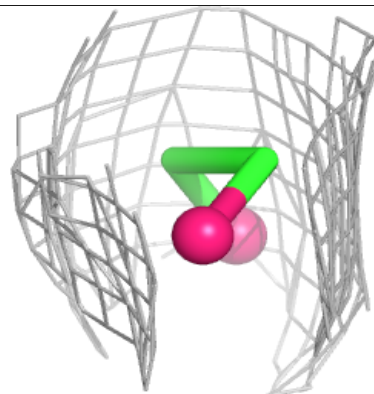
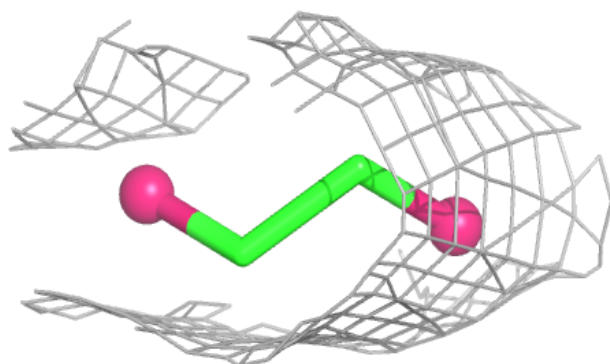
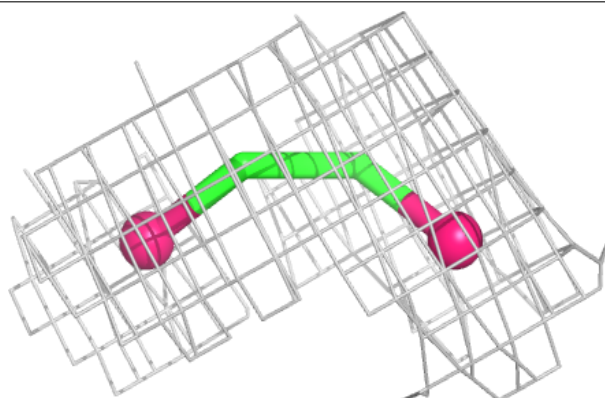


**Electron density around EDO E 444:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

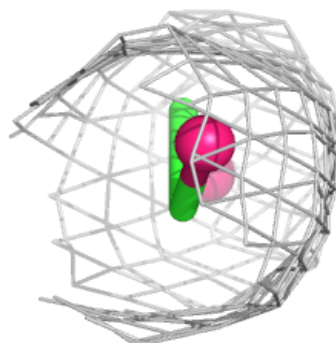
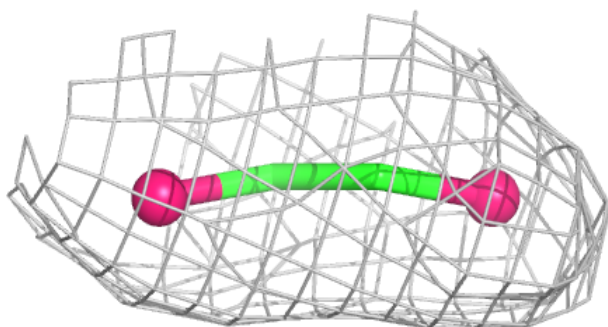
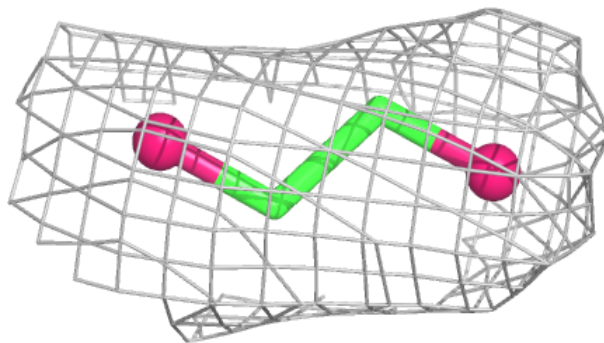
**Electron density around EDO B 408:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

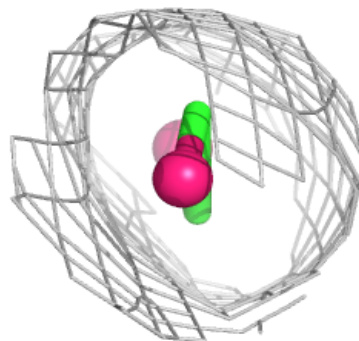
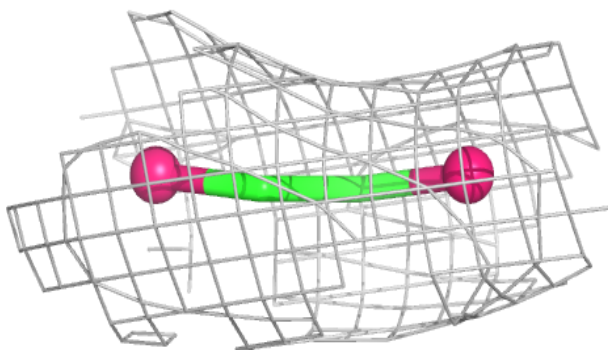
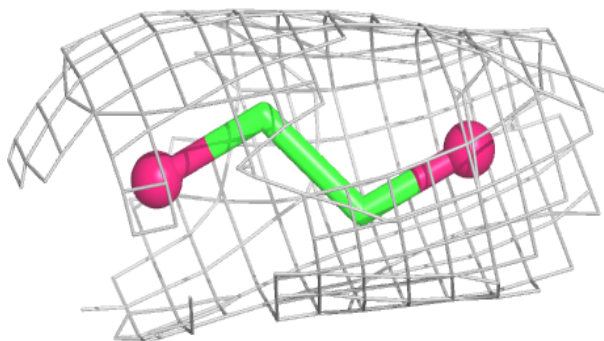


**Electron density around EDO D 421:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around EDO C 411:**

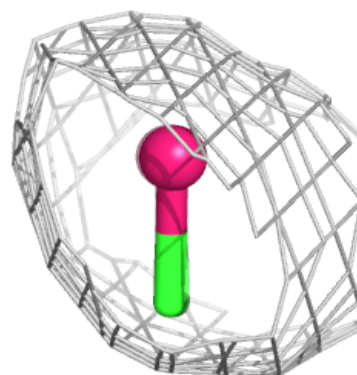
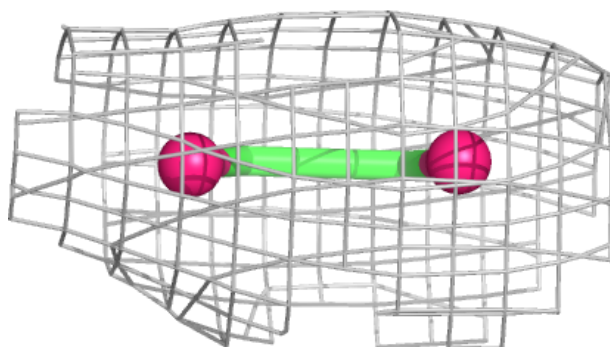
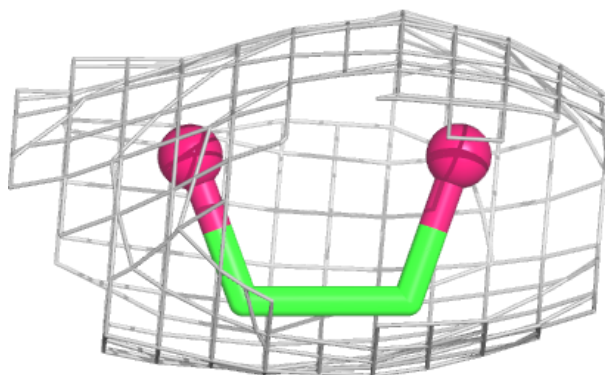
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



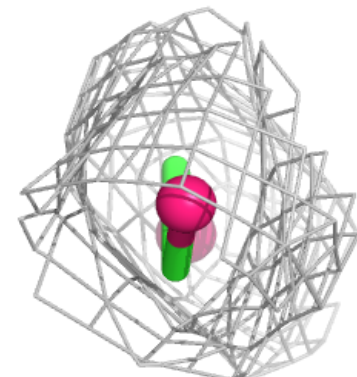
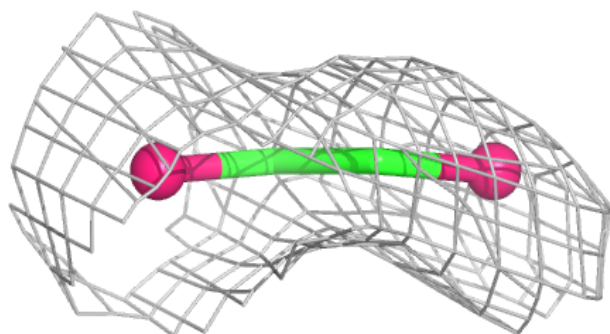
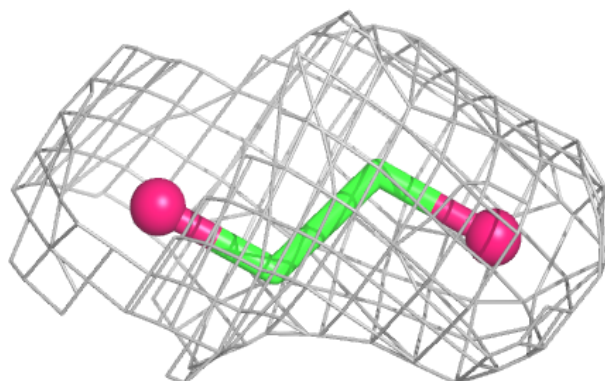


**Electron density around EDO F 402:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

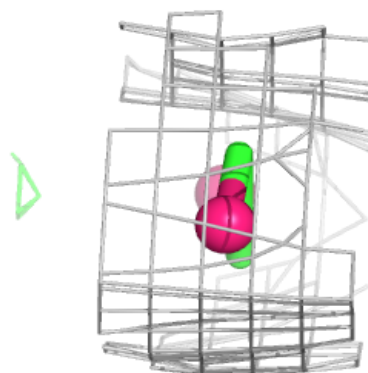
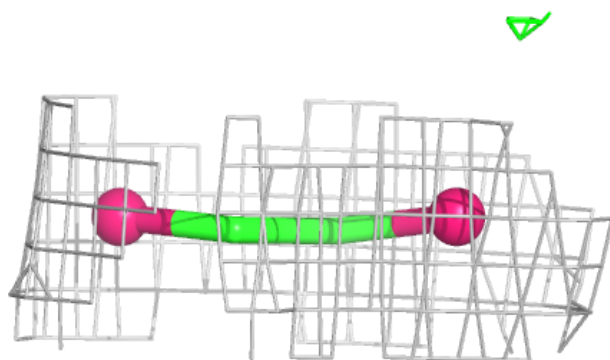
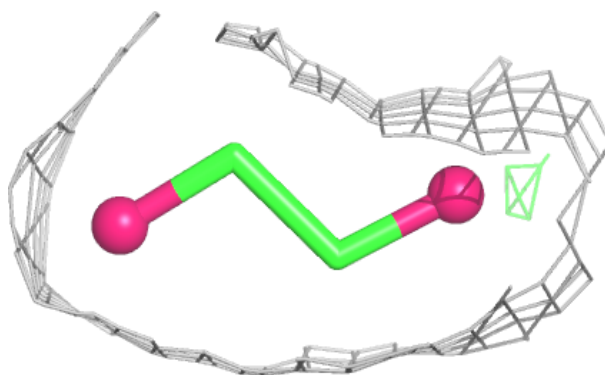
**Electron density around EDO C 412:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

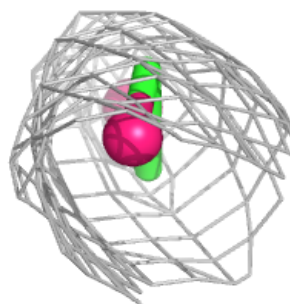
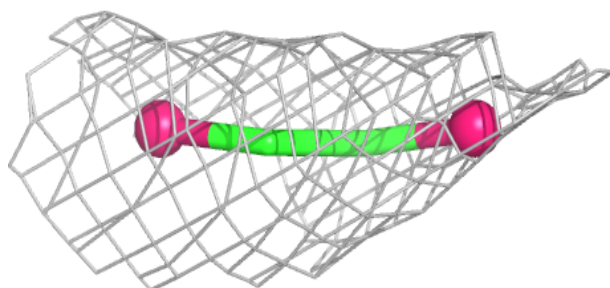
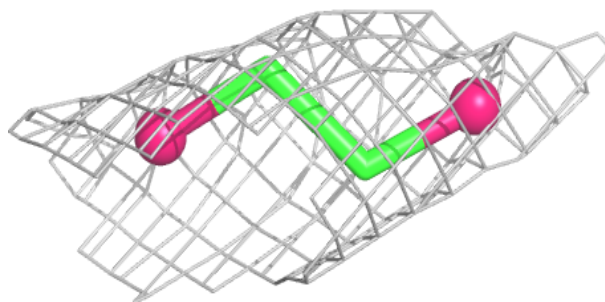


**Electron density around EDO B 409:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

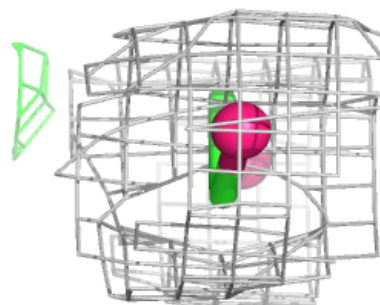
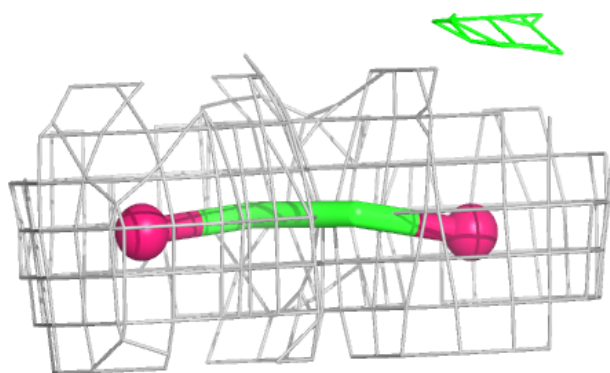
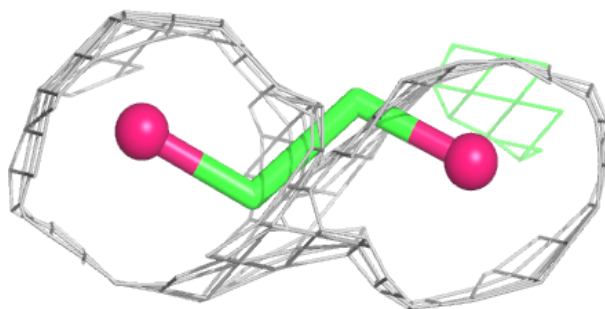
**Electron density around EDO F 405:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

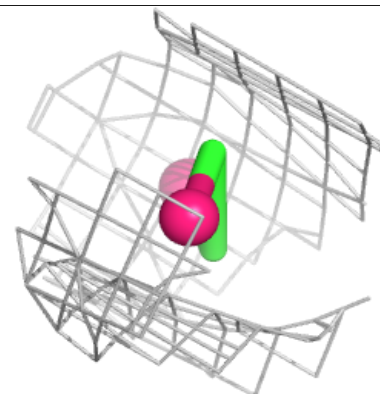
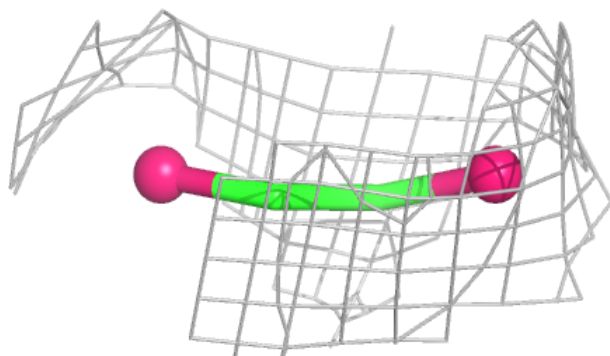
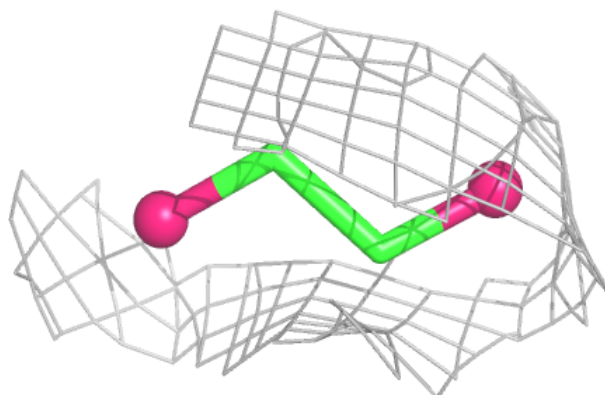


**Electron density around EDO C 415:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

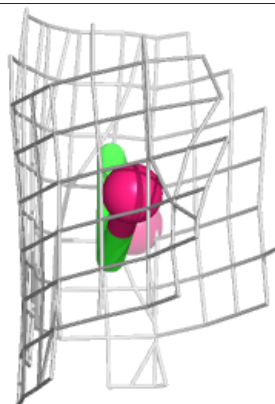
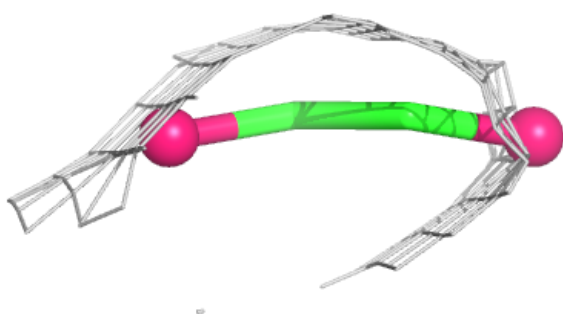
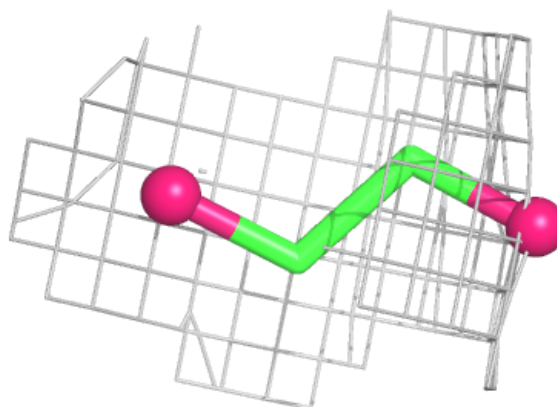
**Electron density around EDO C 417:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

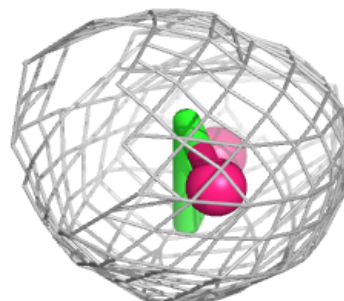
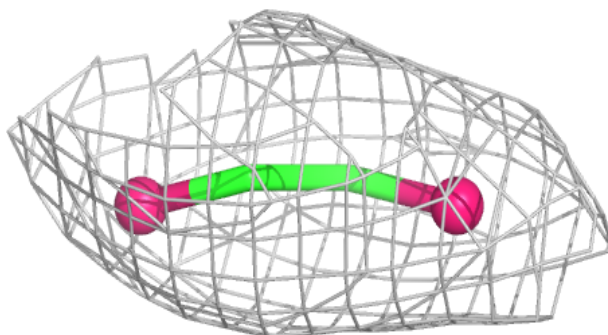
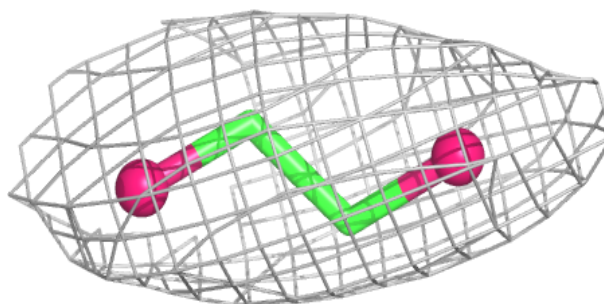


**Electron density around EDO D 428:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around EDO A 430:**

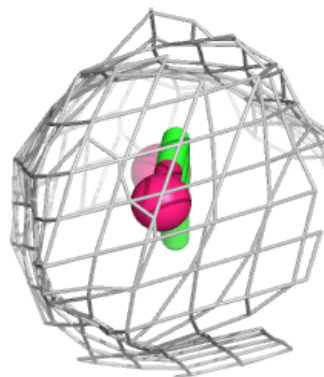
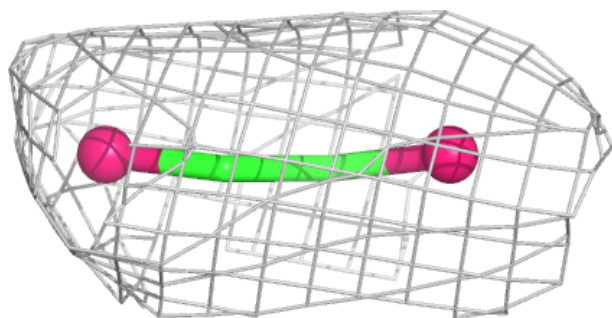
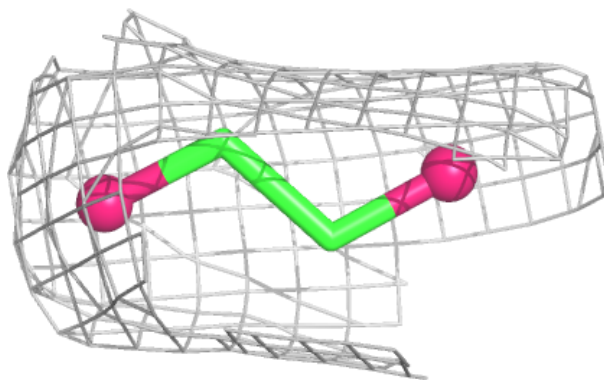
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



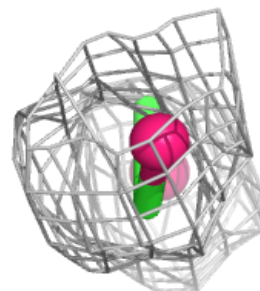
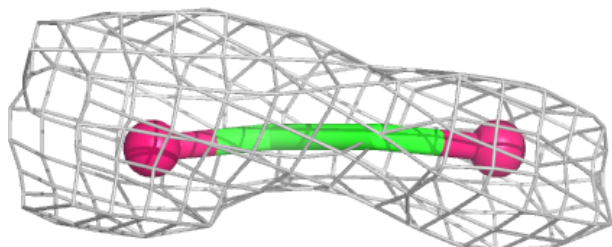
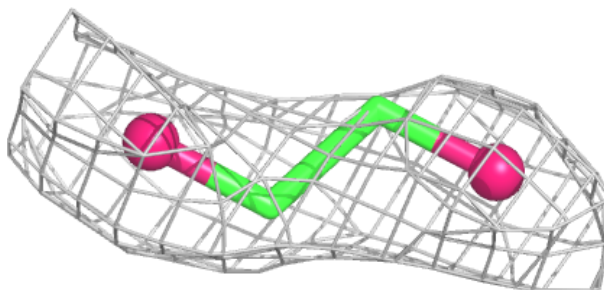


**Electron density around EDO C 419:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

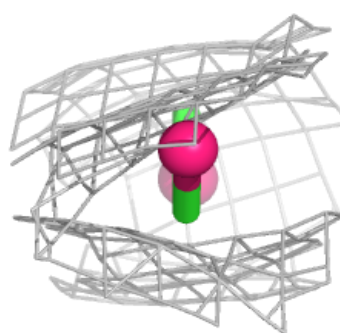
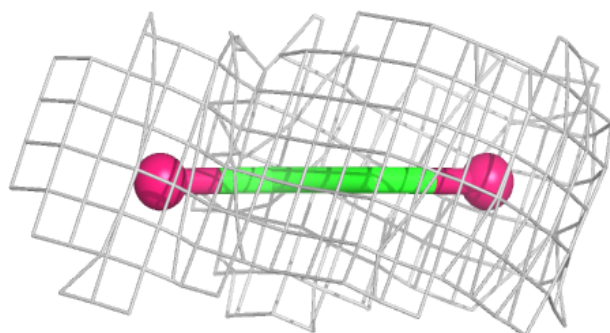
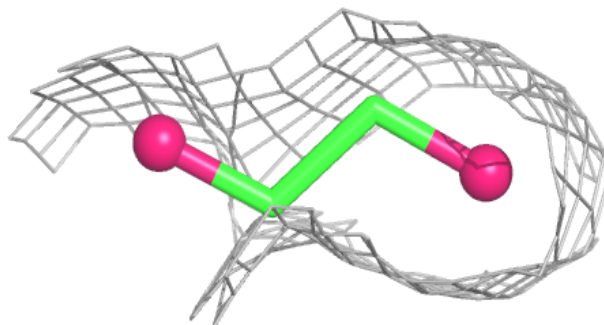
**Electron density around EDO F 412:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

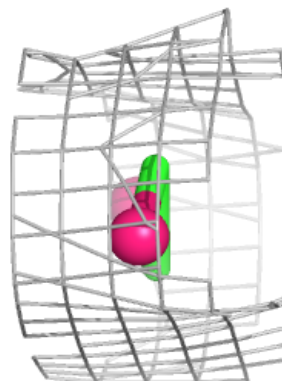
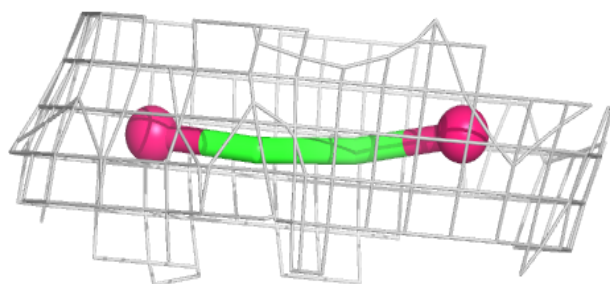
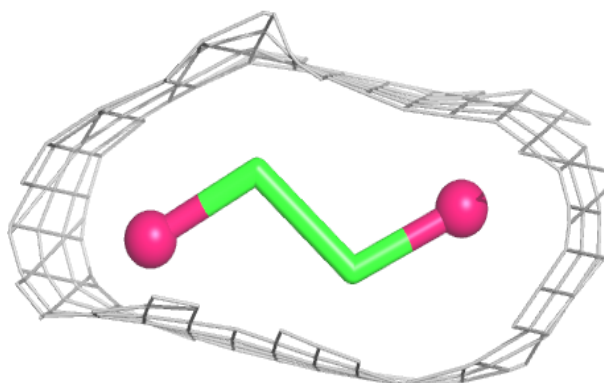


**Electron density around EDO F 413:**

$2mF_o - DF_c$  (at 0.7 rmsd) in gray  
 $mF_o - DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

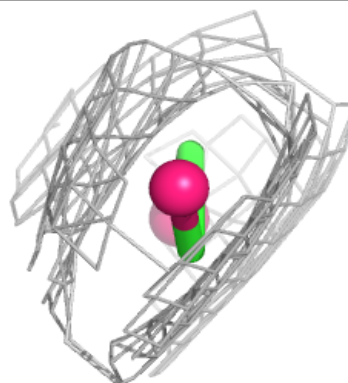
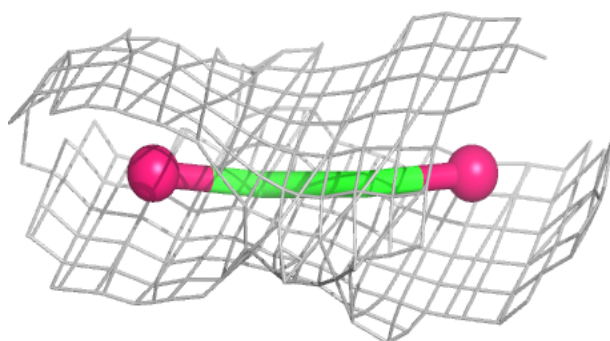
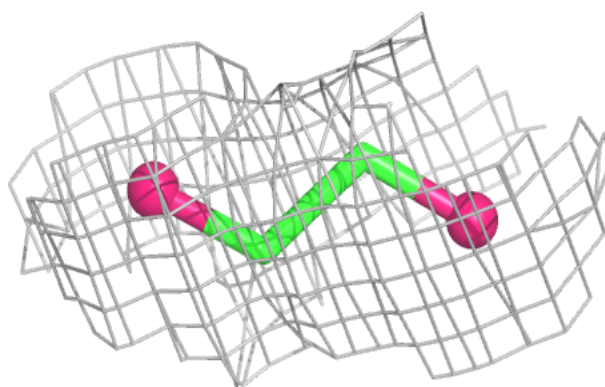
**Electron density around EDO F 415:**

$2mF_o - DF_c$  (at 0.7 rmsd) in gray  
 $mF_o - DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

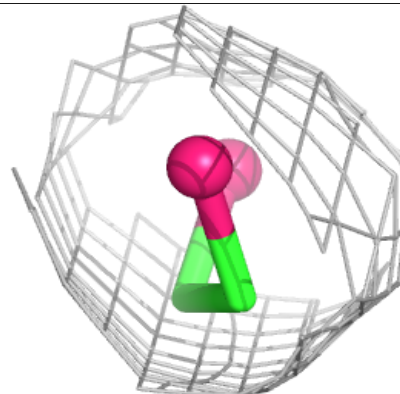
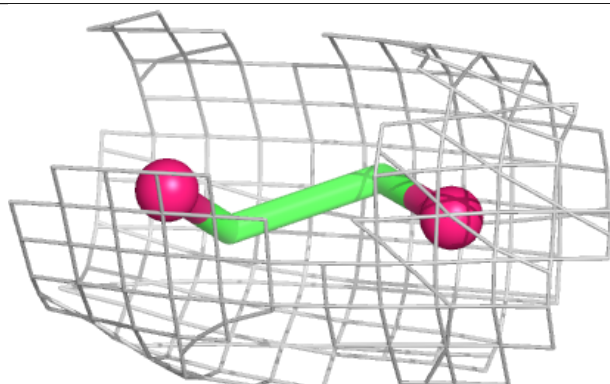


**Electron density around EDO A 435:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

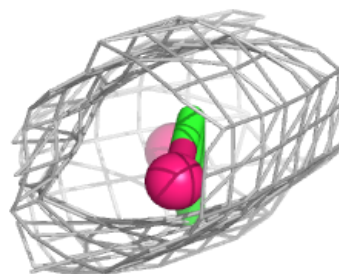
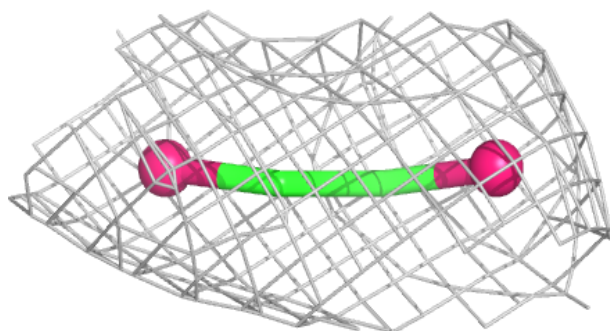
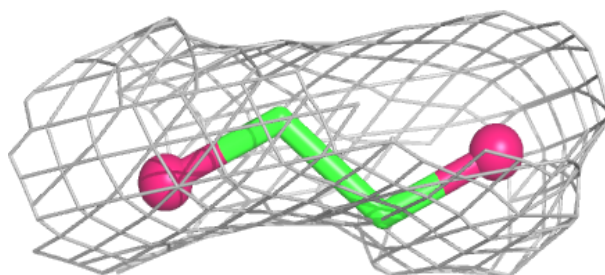
**Electron density around EDO B 432:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

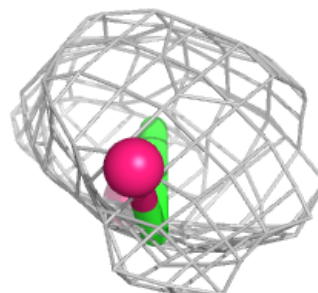
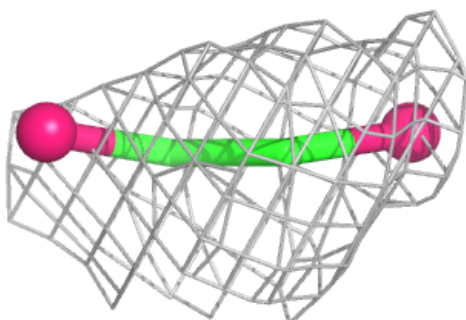
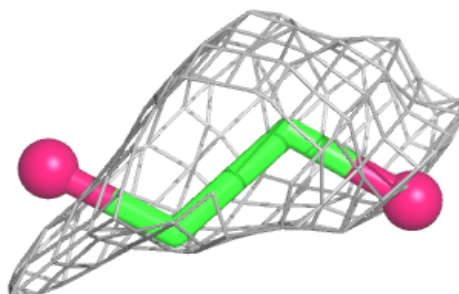


**Electron density around EDO F 419:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around EDO C 423:**

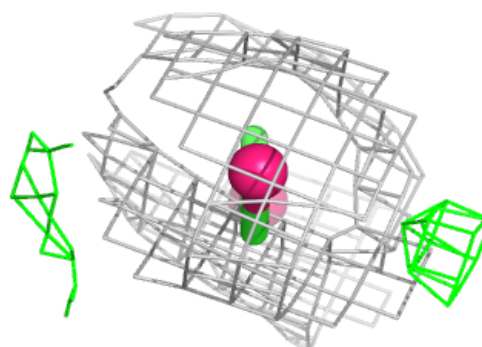
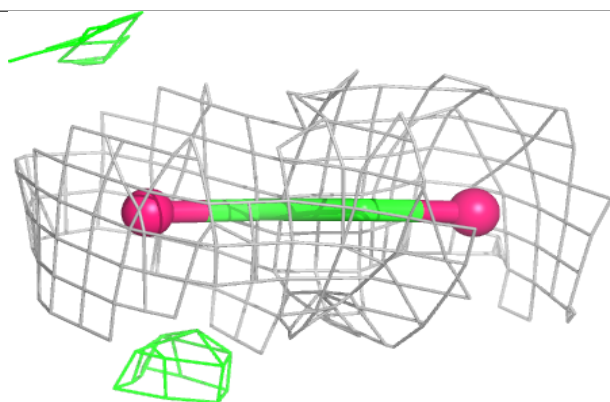
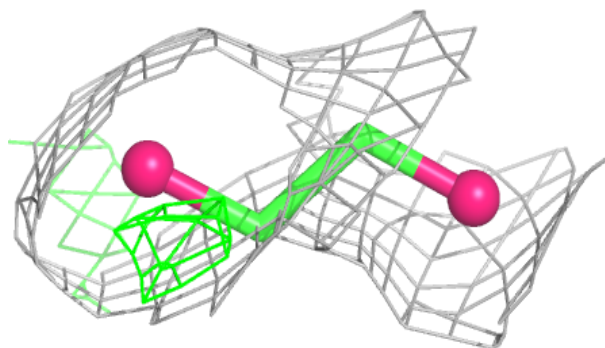
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



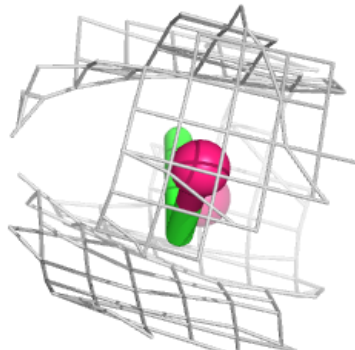
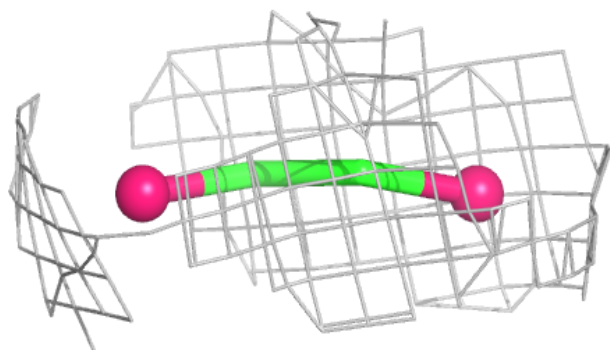
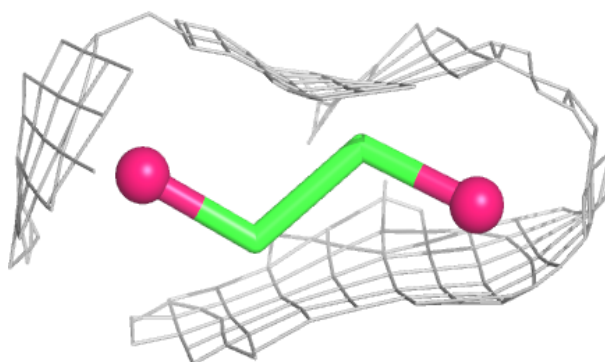


**Electron density around EDO F 421:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

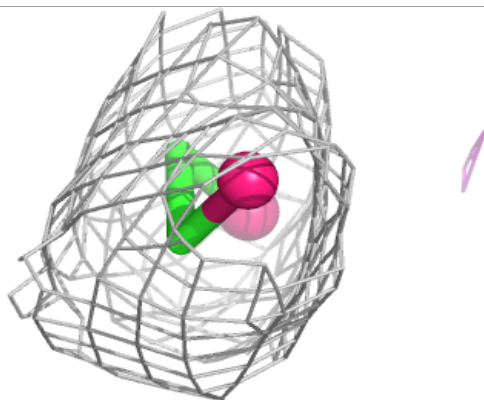
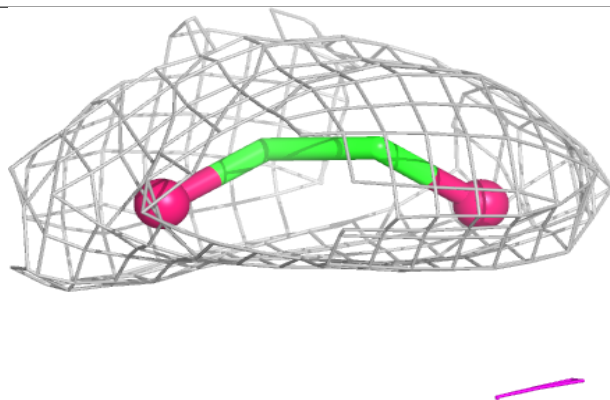
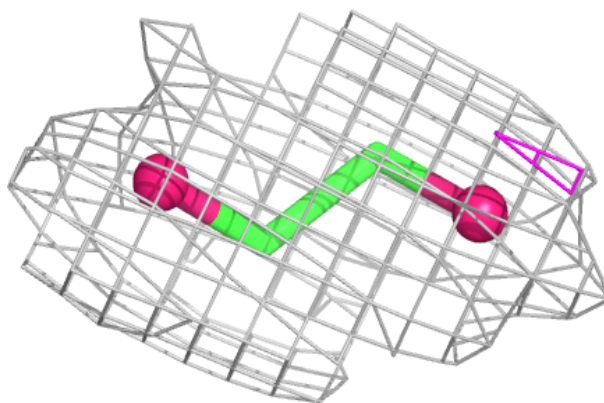
**Electron density around EDO F 423:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

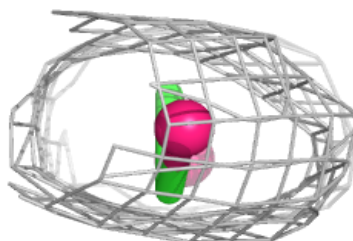
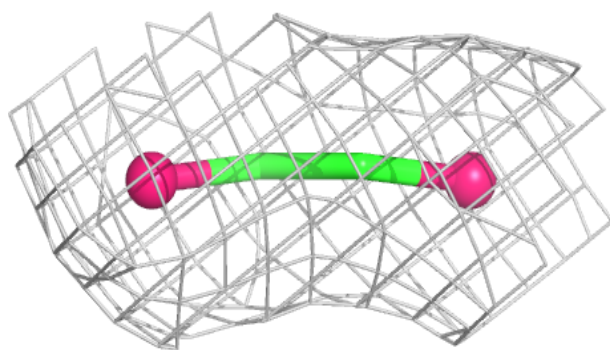
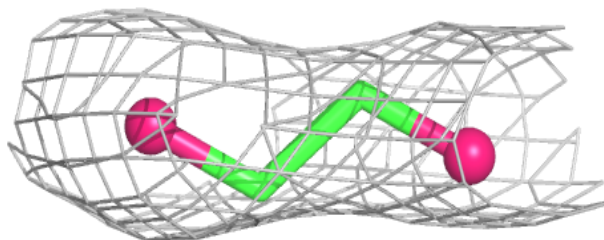


**Electron density around EDO B 433:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

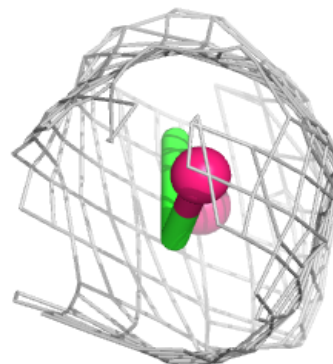
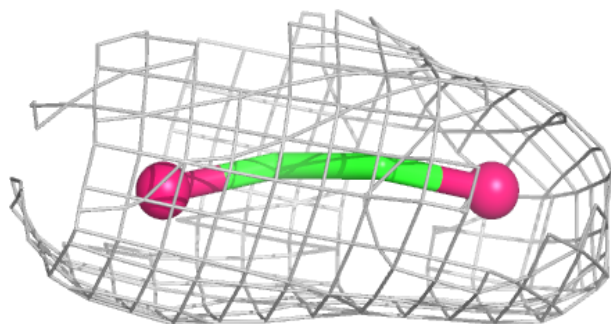
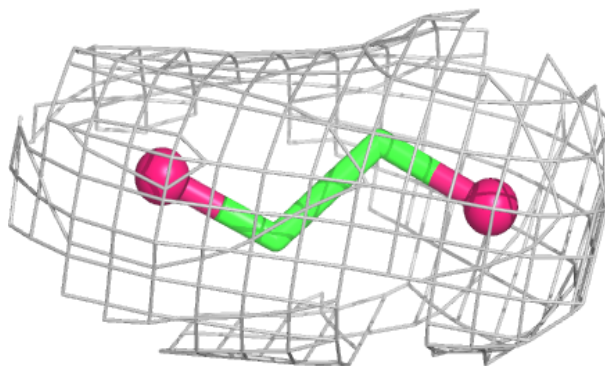
**Electron density around EDO B 413:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

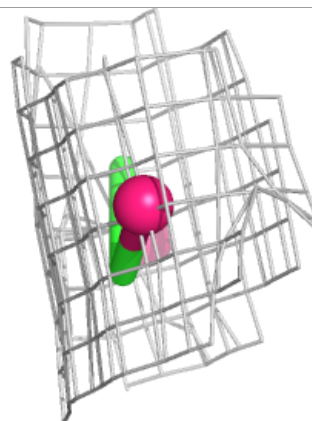
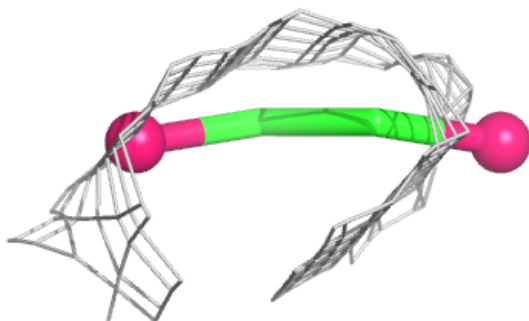
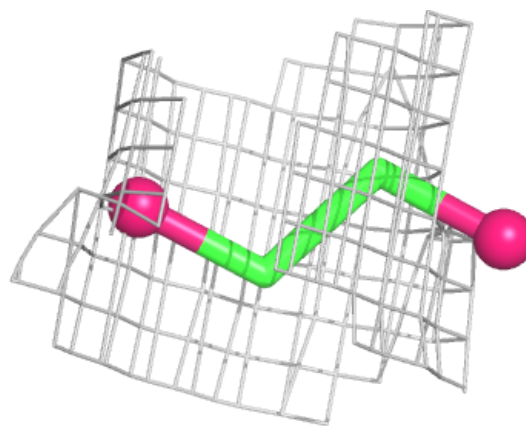


**Electron density around EDO F 426:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

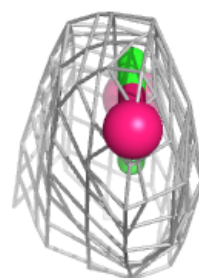
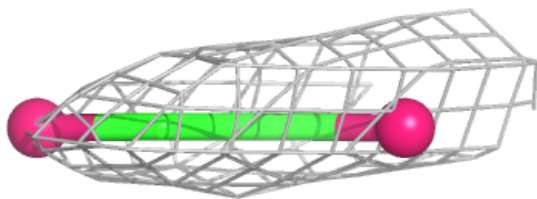
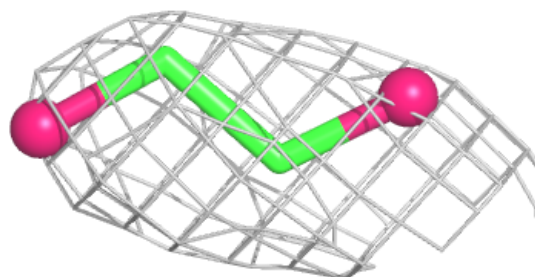
**Electron density around EDO C 426:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

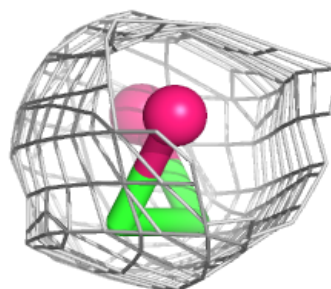
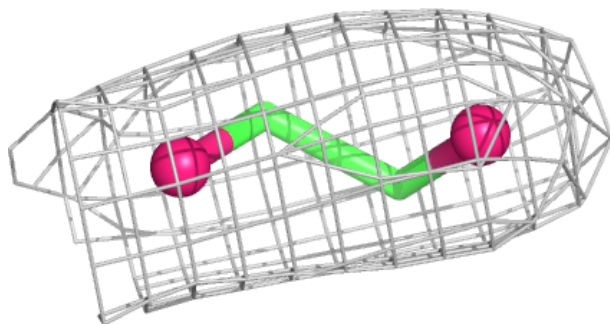
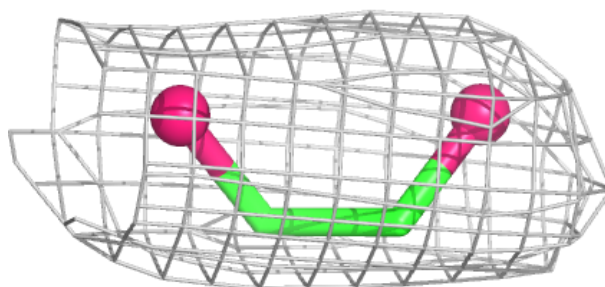


**Electron density around EDO D 437:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around EDO D 438:**

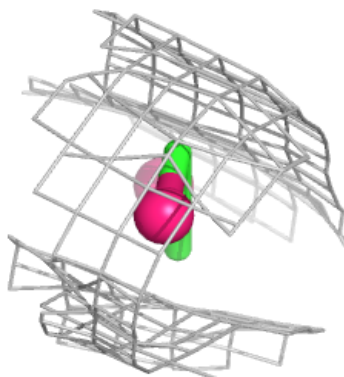
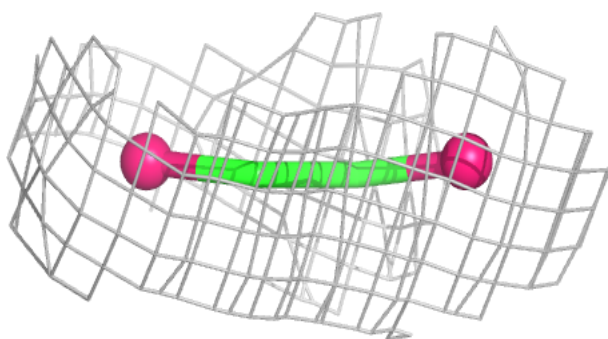
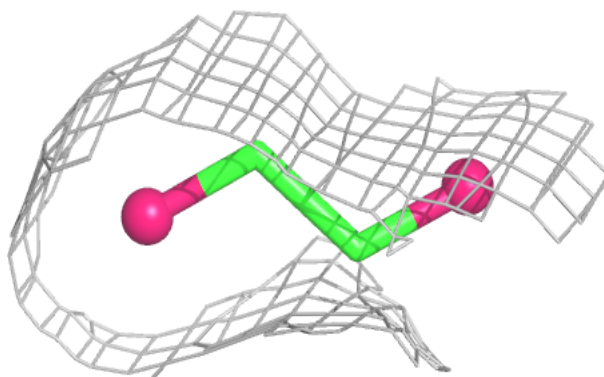
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



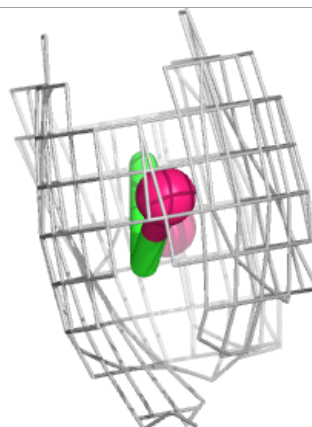
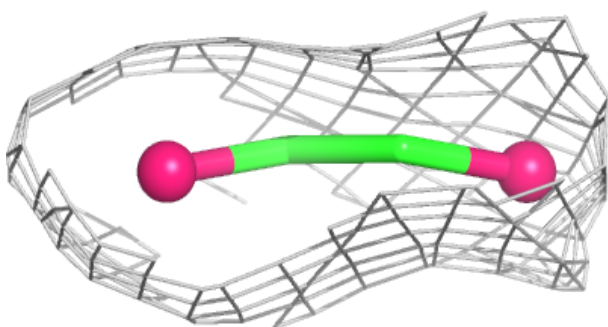
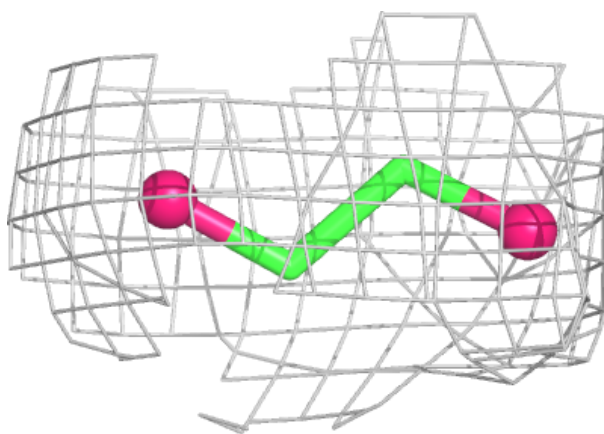


**Electron density around EDO A 405:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

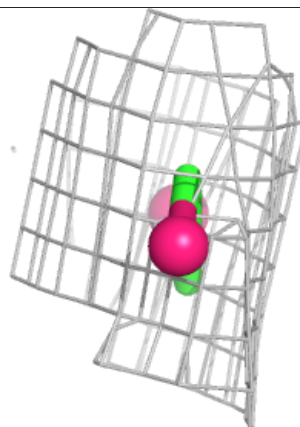
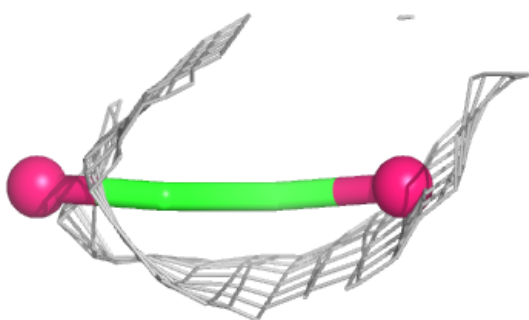
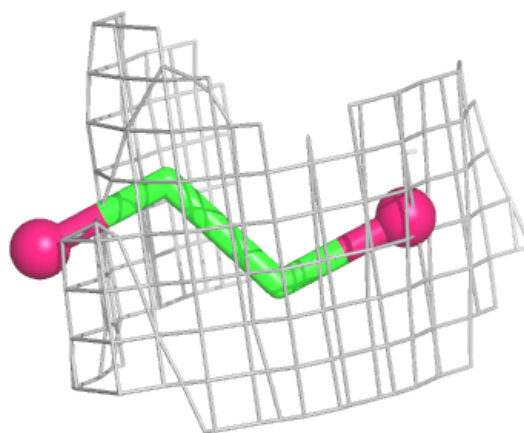
**Electron density around EDO B 436:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

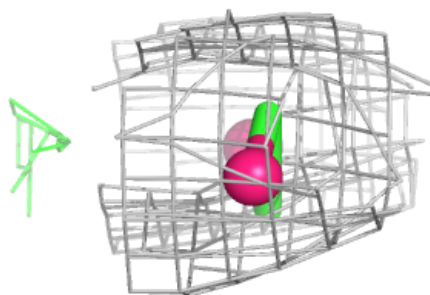
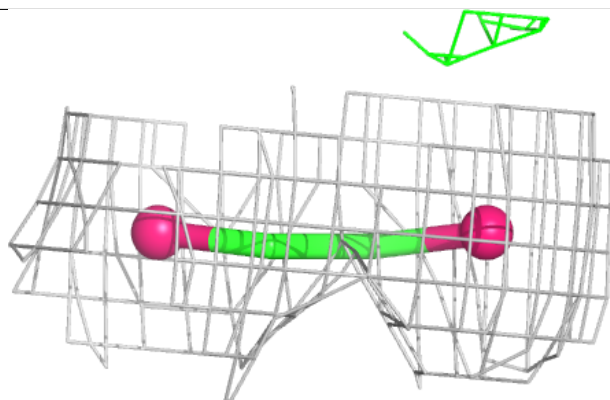
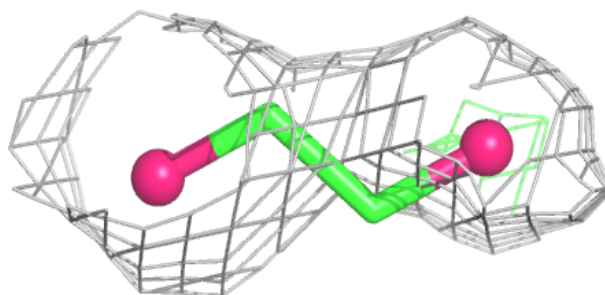


**Electron density around EDO F 433:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

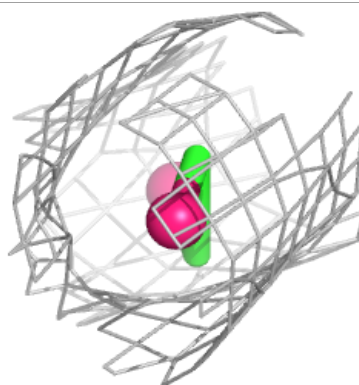
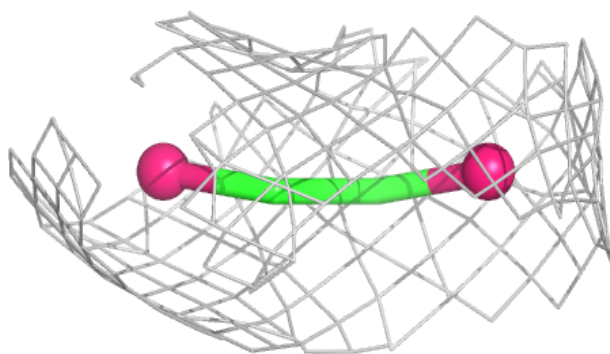
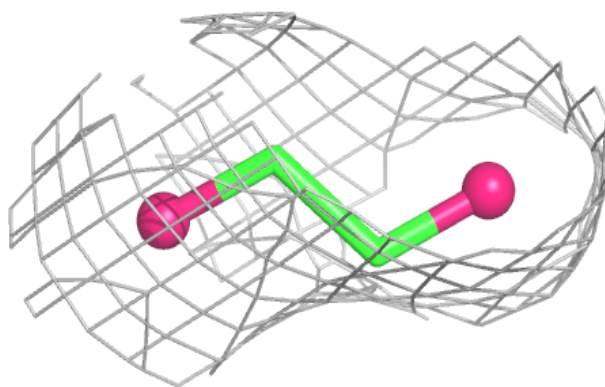
**Electron density around EDO B 415:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

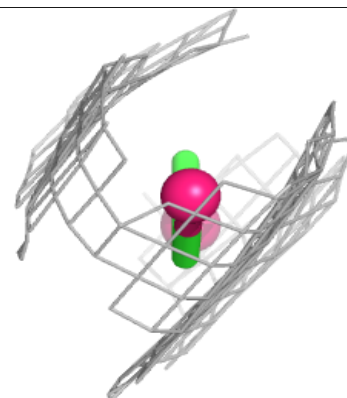
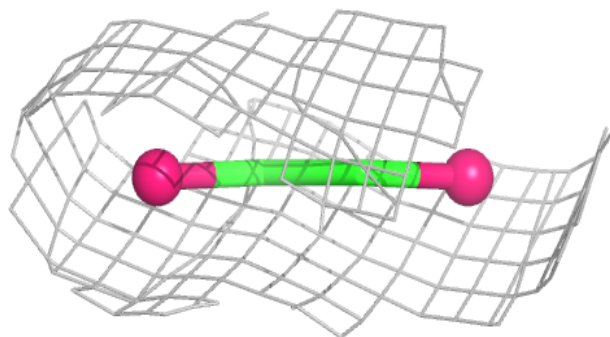
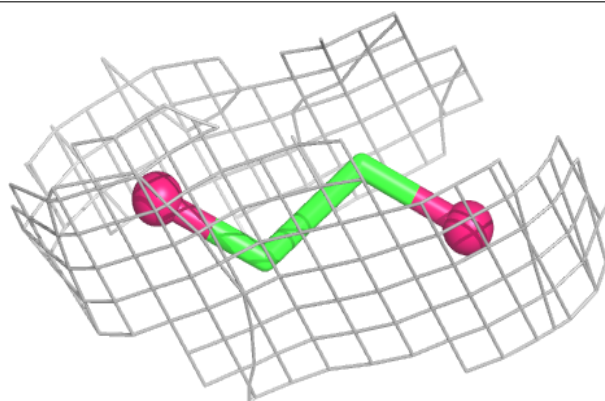


**Electron density around EDO B 416:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

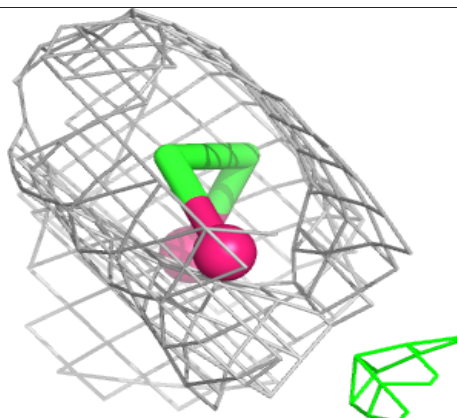
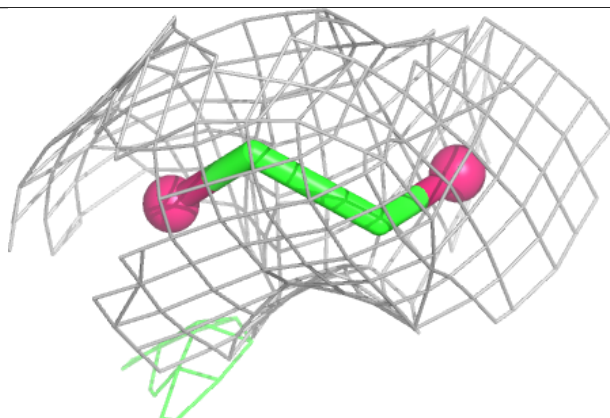
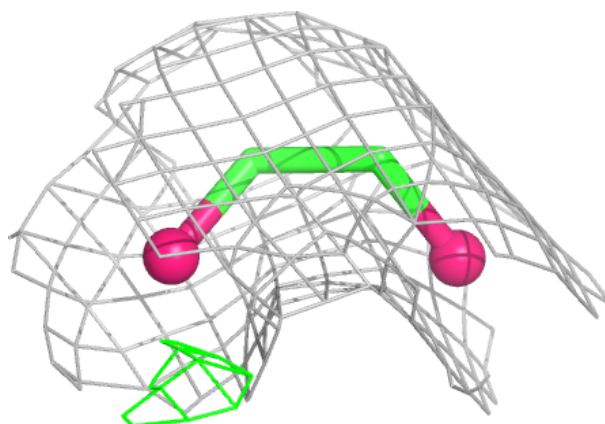
**Electron density around EDO B 401:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

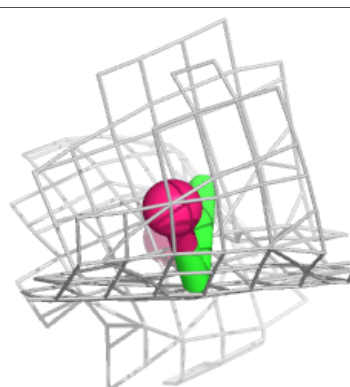
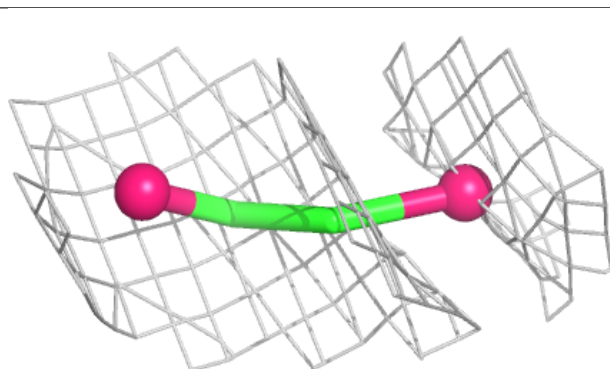
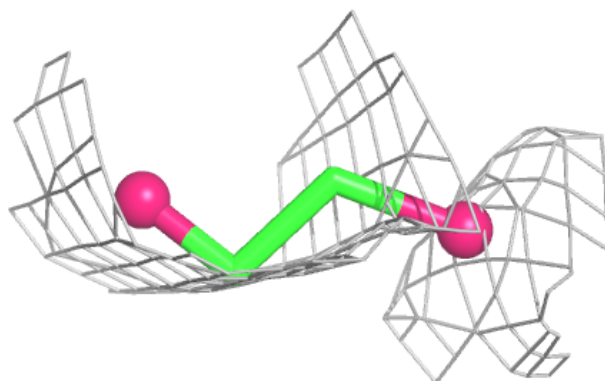


**Electron density around EDO E 405:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around EDO B 440:**

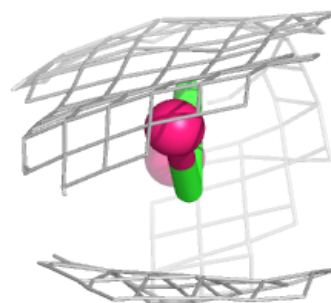
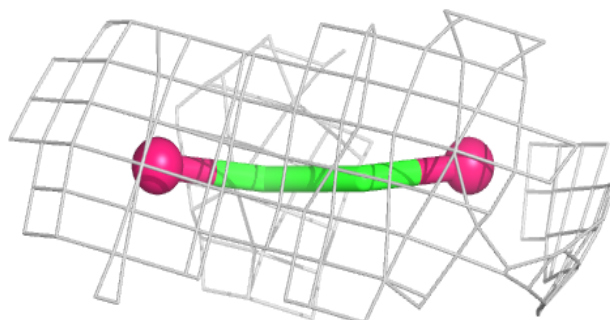
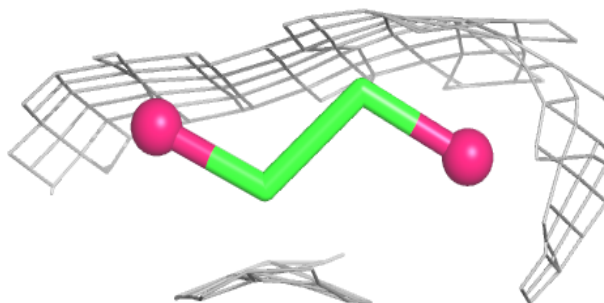
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



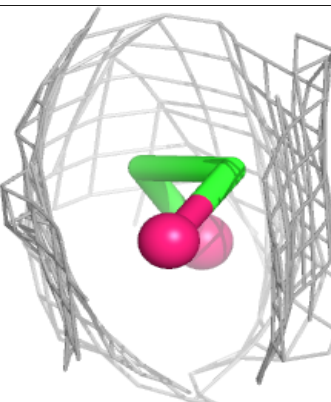
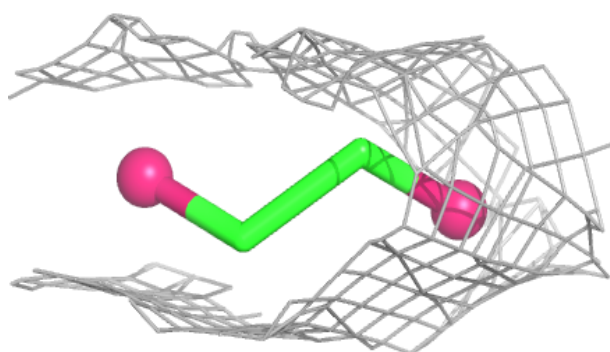
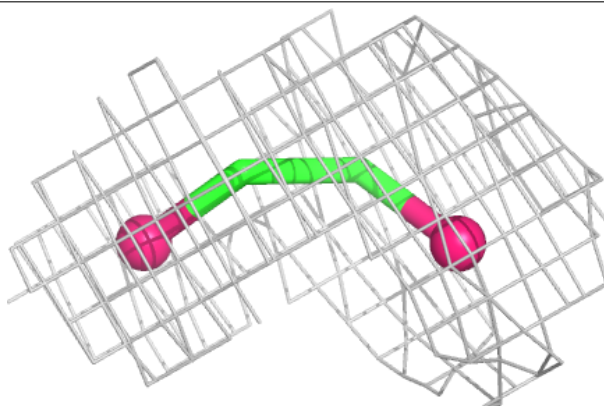


**Electron density around EDO E 407:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

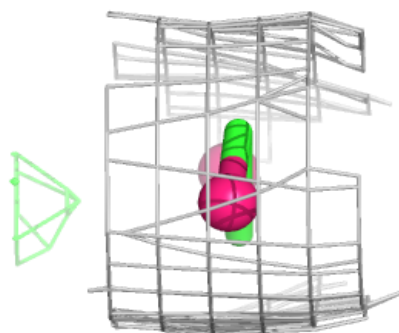
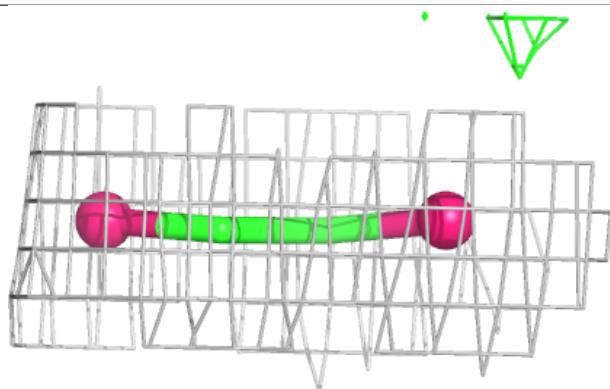
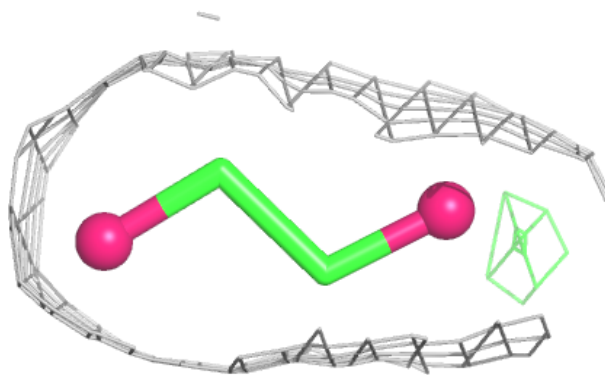
**Electron density around EDO E 408:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



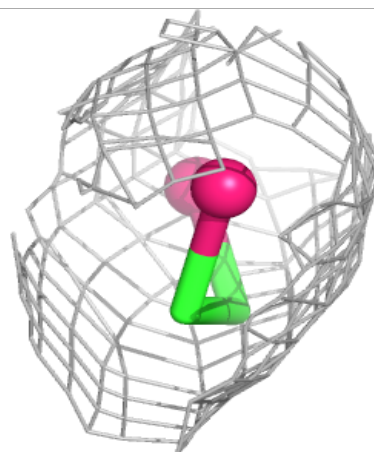
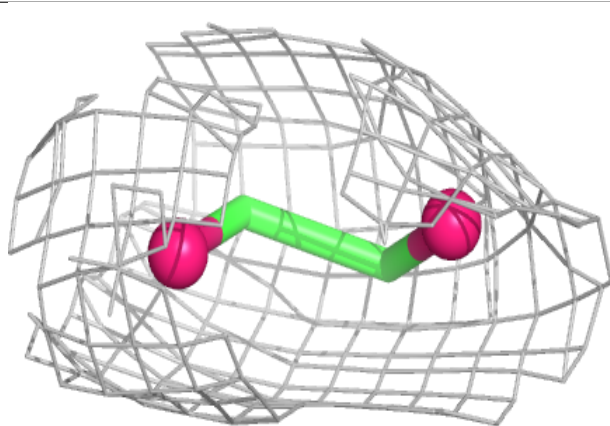
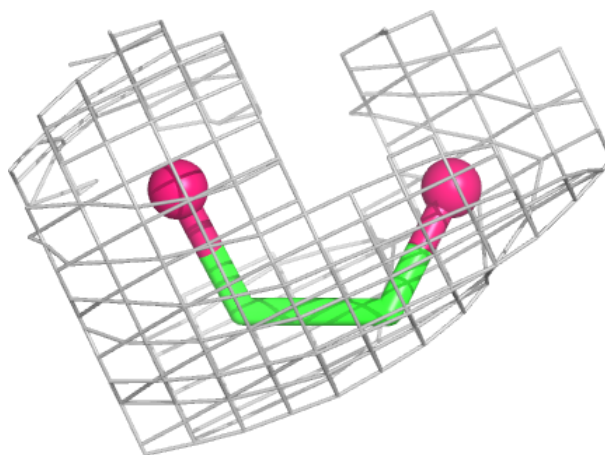
**Electron density around EDO E 409:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



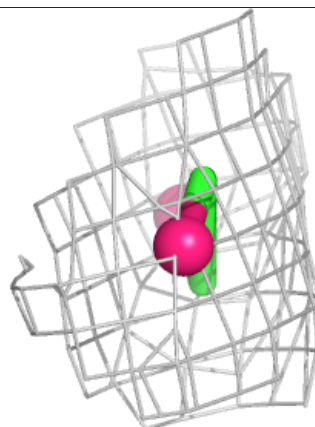
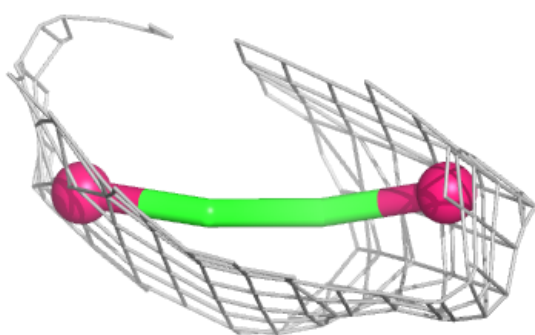
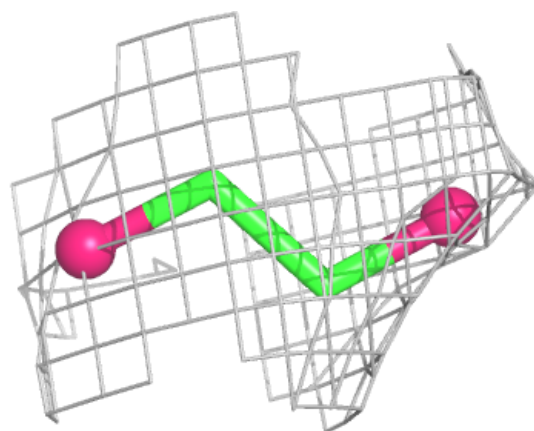
**Electron density around EDO A 437:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

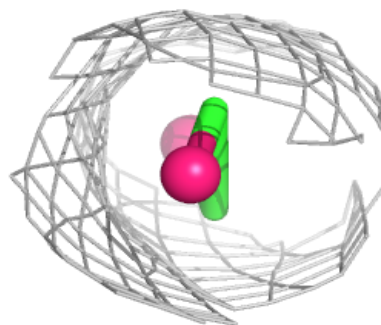
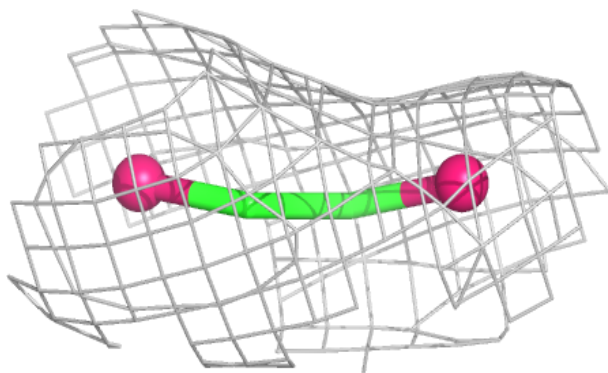
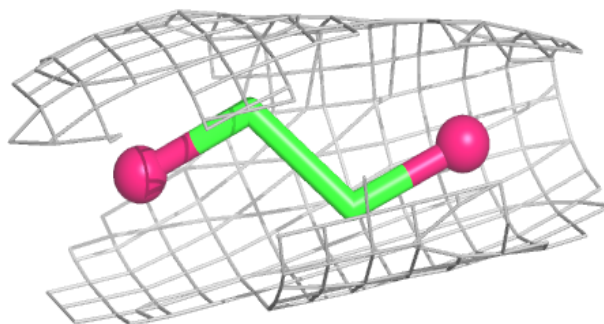


**Electron density around EDO A 425:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around EDO E 412:**

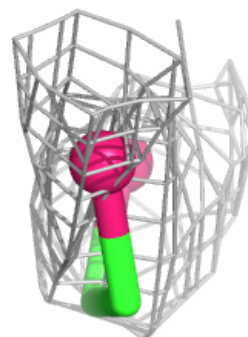
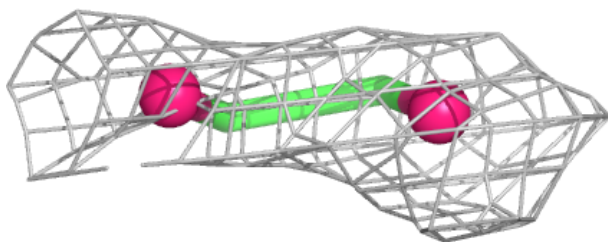
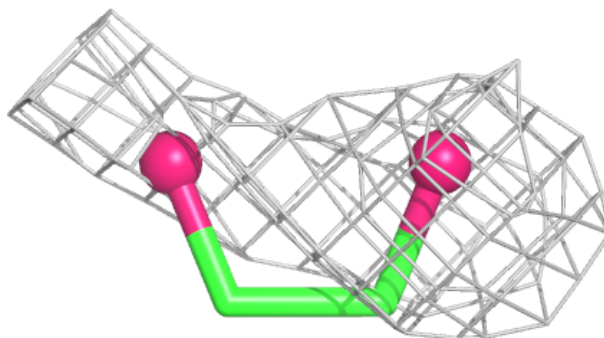
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



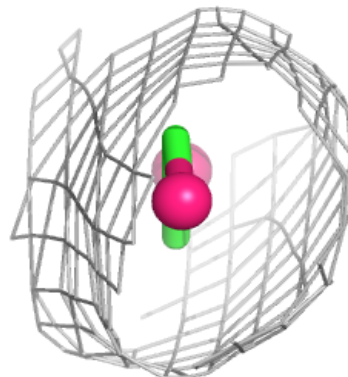
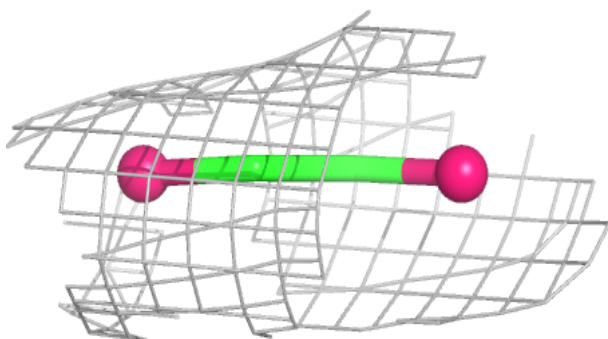
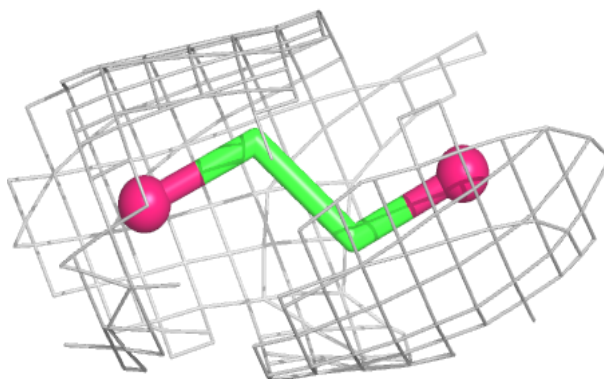


**Electron density around EDO B 443:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

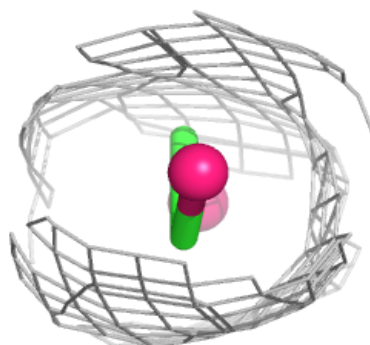
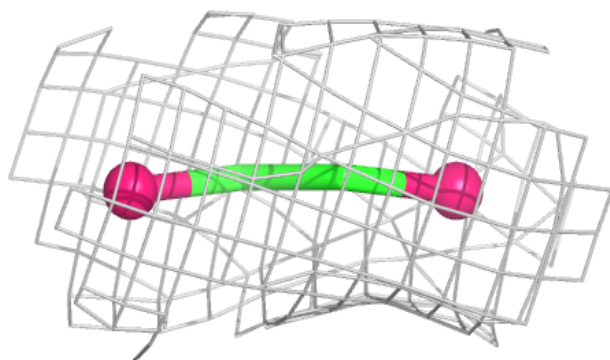
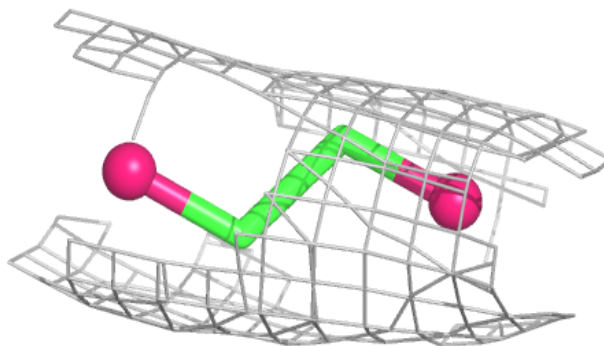
**Electron density around EDO F 417:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

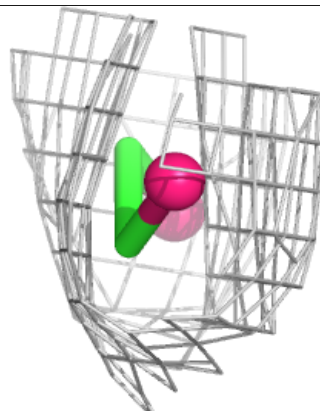
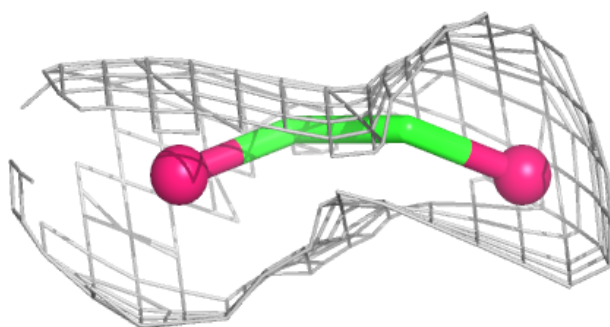
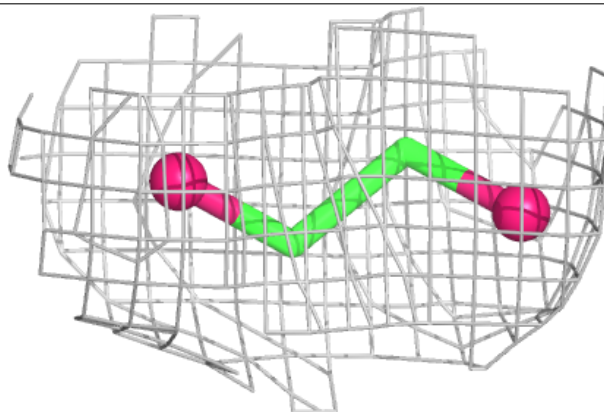


**Electron density around EDO B 411:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

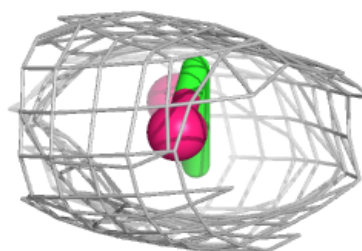
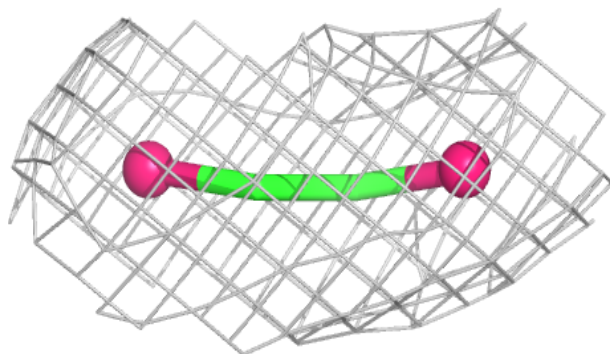
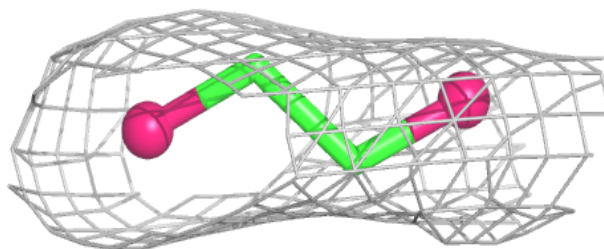
**Electron density around EDO E 440:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

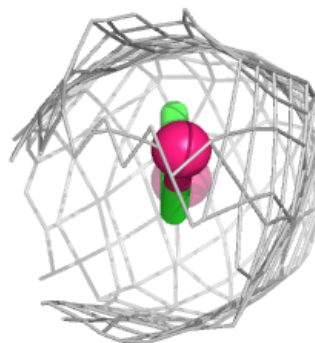
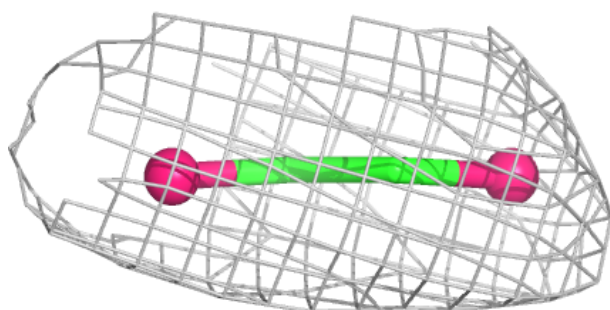
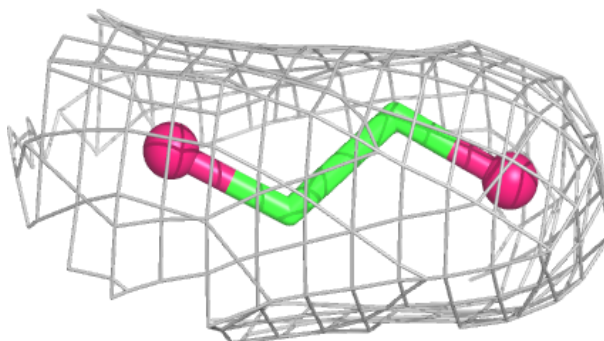


**Electron density around EDO C 413:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

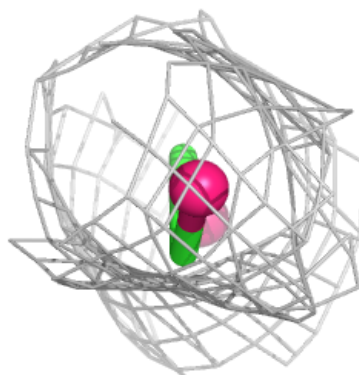
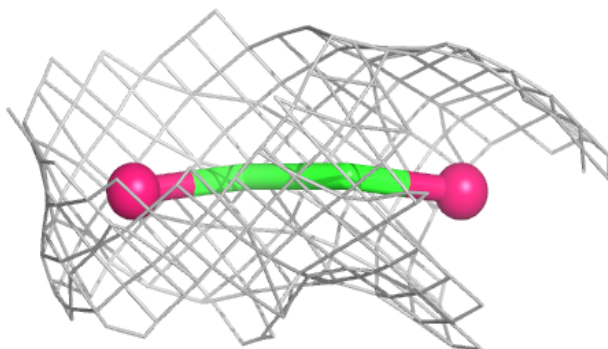
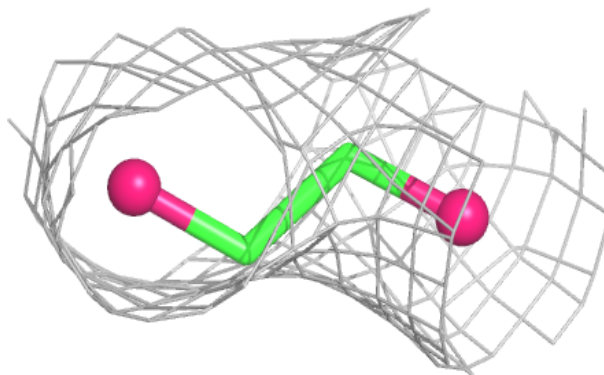
**Electron density around EDO B 420:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

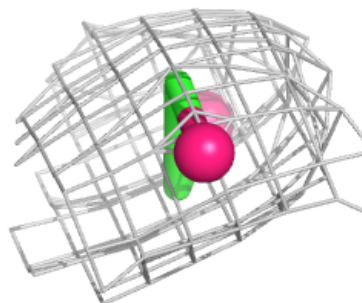
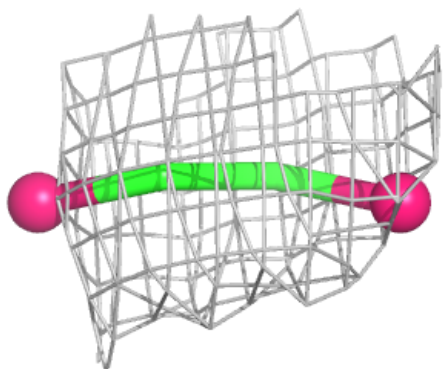
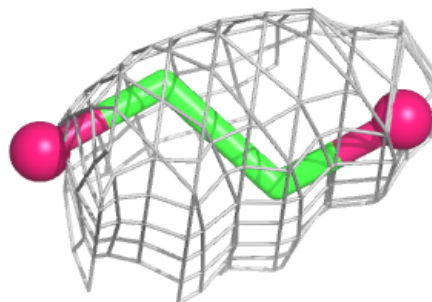


**Electron density around EDO F 422:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around EDO B 404:**

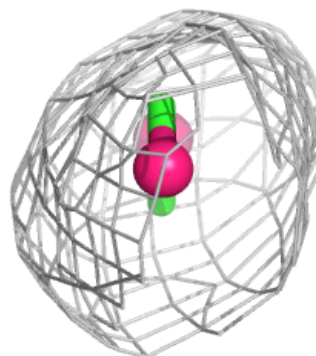
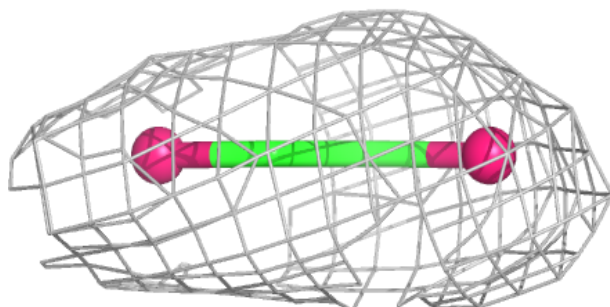
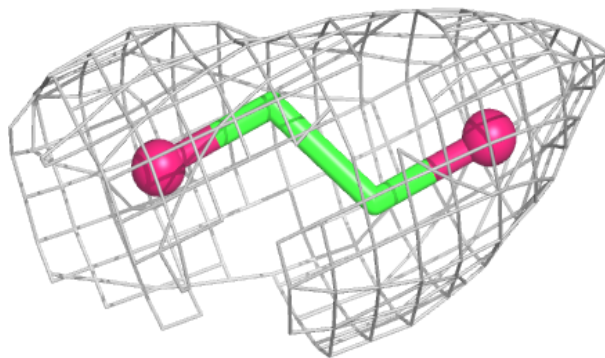
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



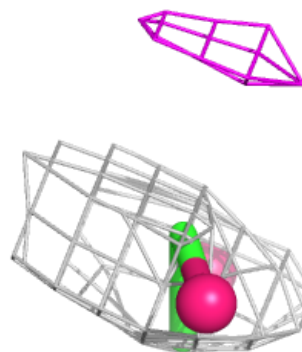
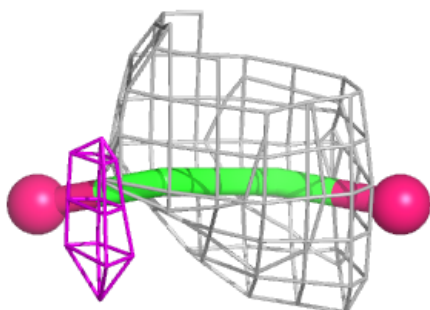
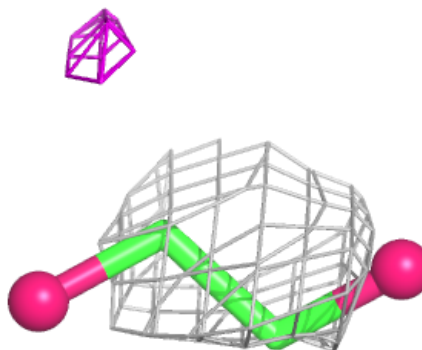


**Electron density around EDO D 423:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

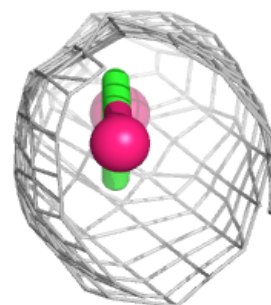
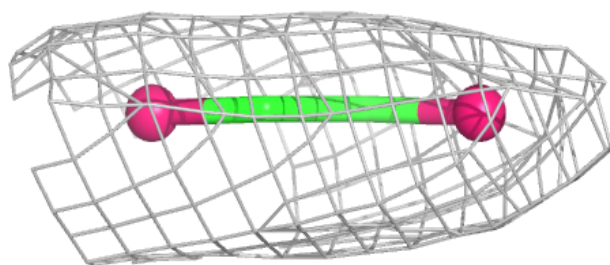
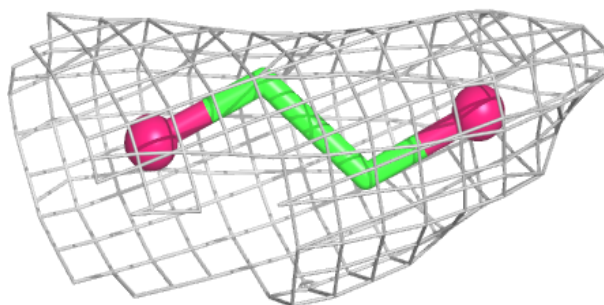
**Electron density around EDO D 405:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

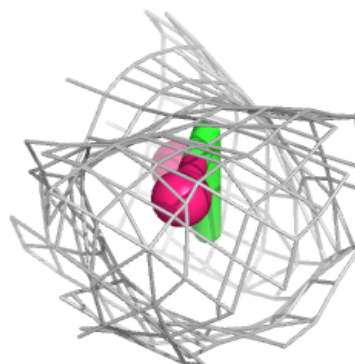
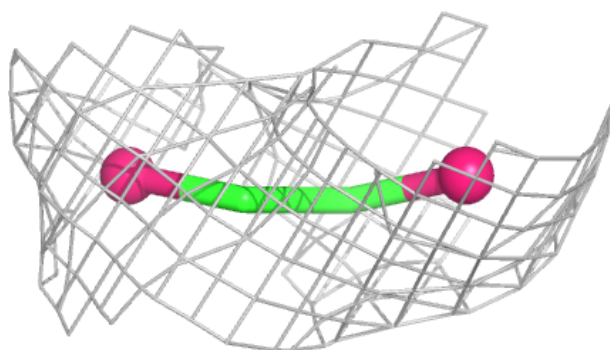
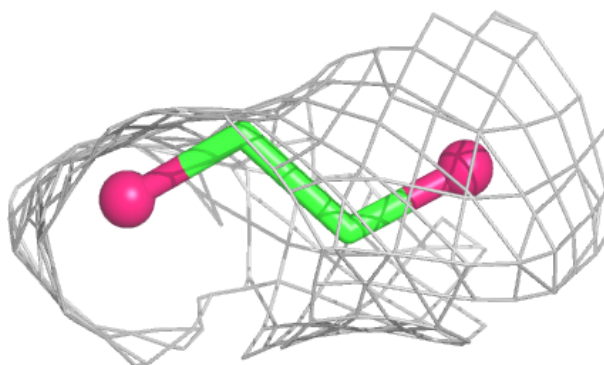


**Electron density around EDO E 424:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

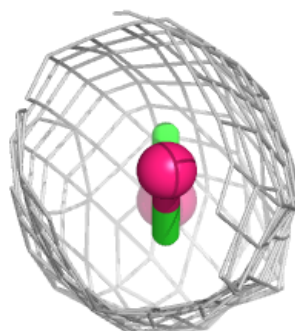
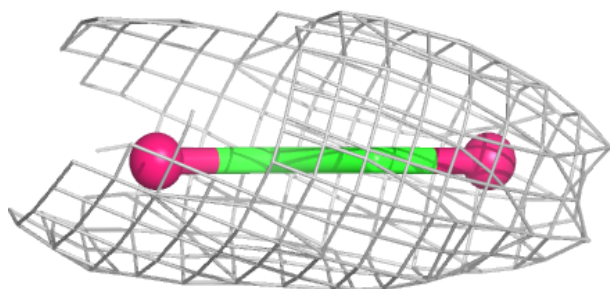
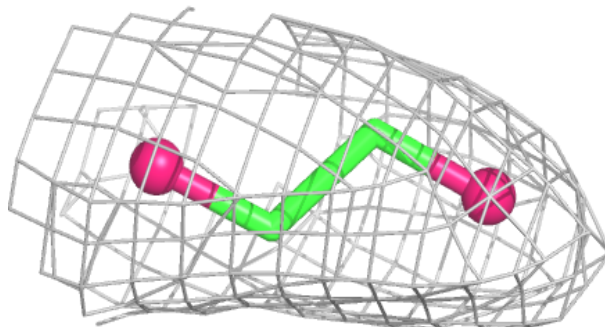
**Electron density around EDO C 416:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

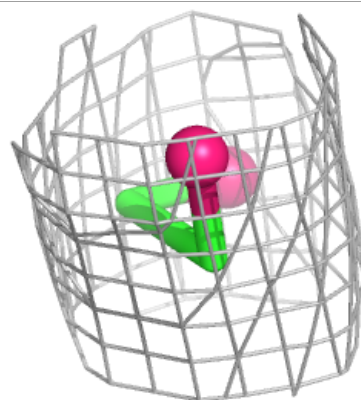
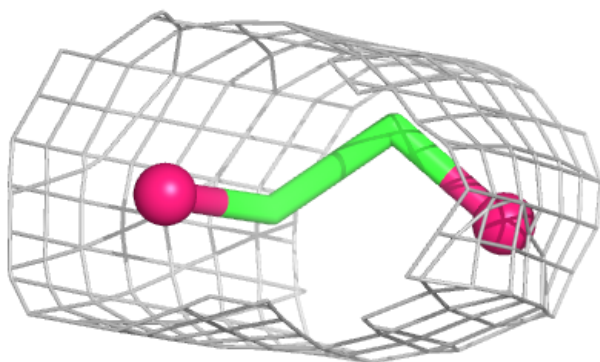
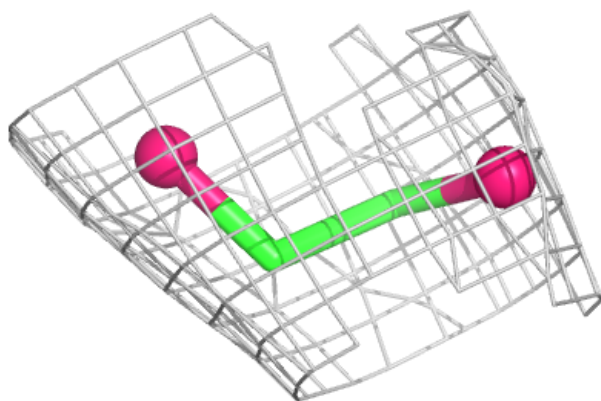


**Electron density around EDO F 428:**

$2mF_o - DF_c$  (at 0.7 rmsd) in gray  
 $mF_o - DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

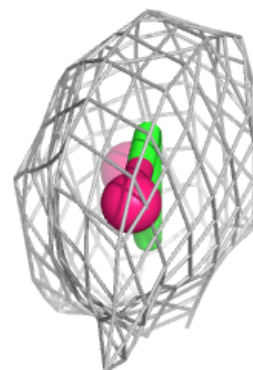
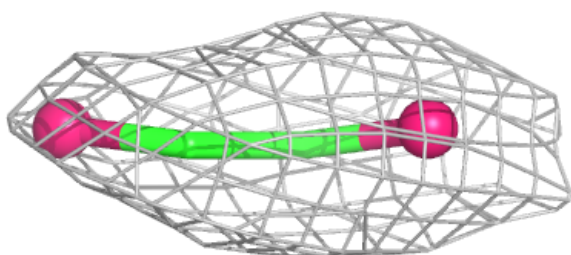
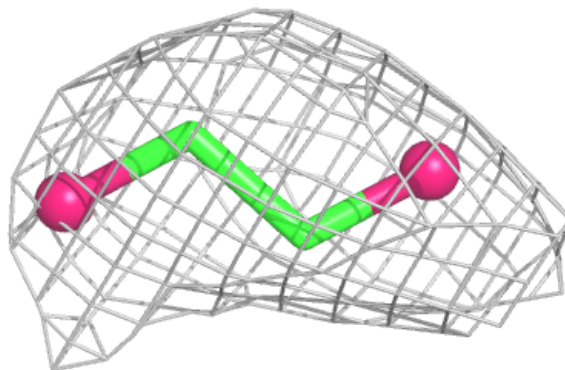
**Electron density around EDO E 448:**

$2mF_o - DF_c$  (at 0.7 rmsd) in gray  
 $mF_o - DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

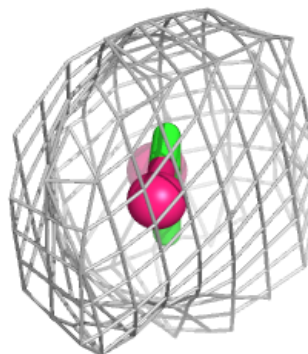
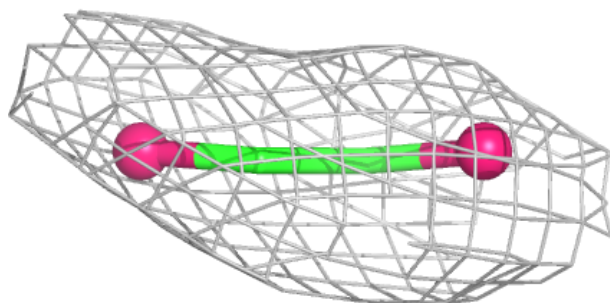
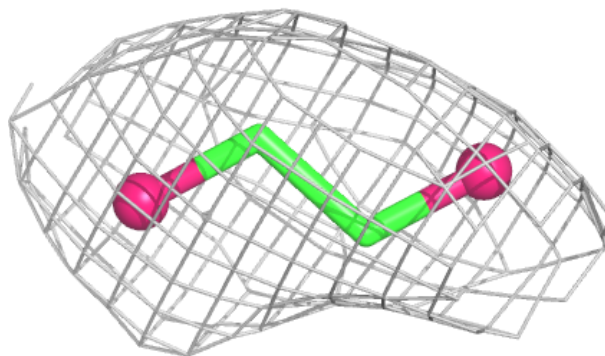


**Electron density around EDO F 401:**

$2mF_o - DF_c$  (at 0.7 rmsd) in gray  
 $mF_o - DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around EDO C 430:**

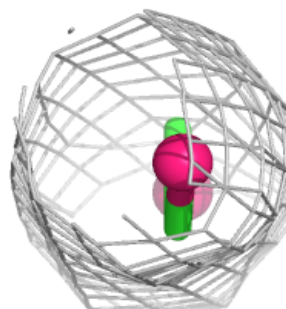
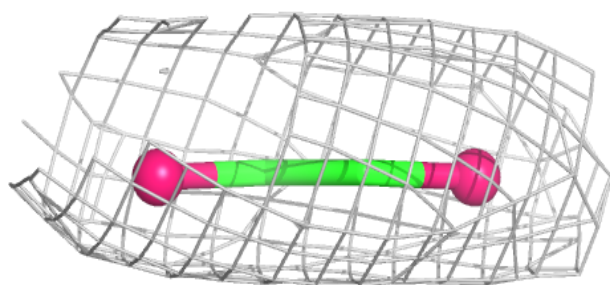
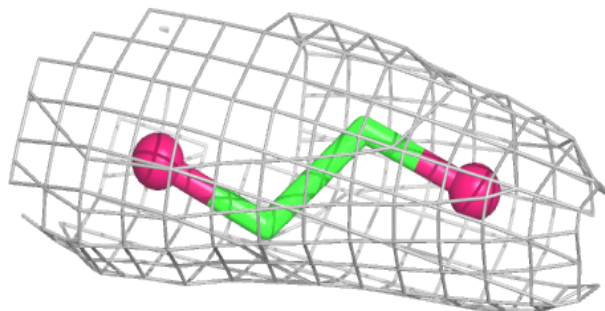
$2mF_o - DF_c$  (at 0.7 rmsd) in gray  
 $mF_o - DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



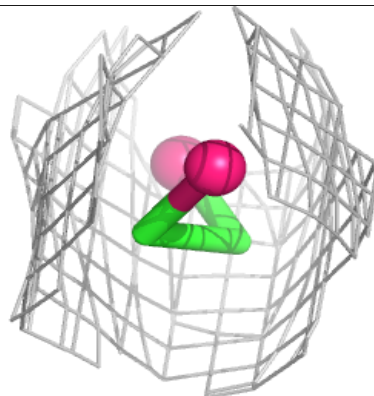
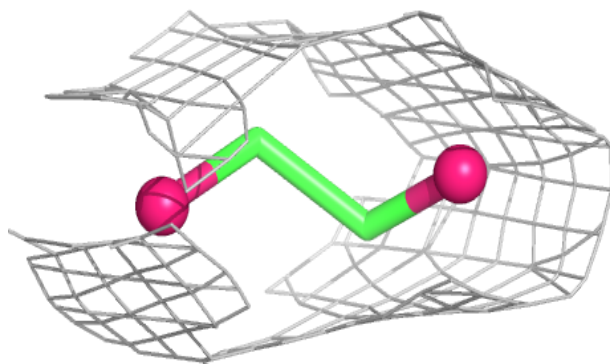
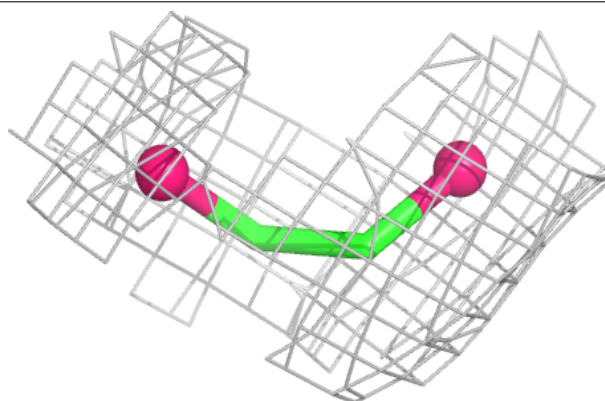


**Electron density around EDO B 422:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

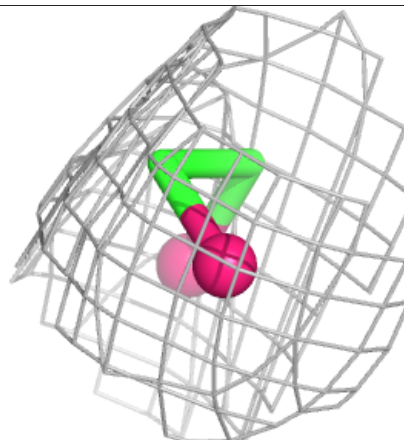
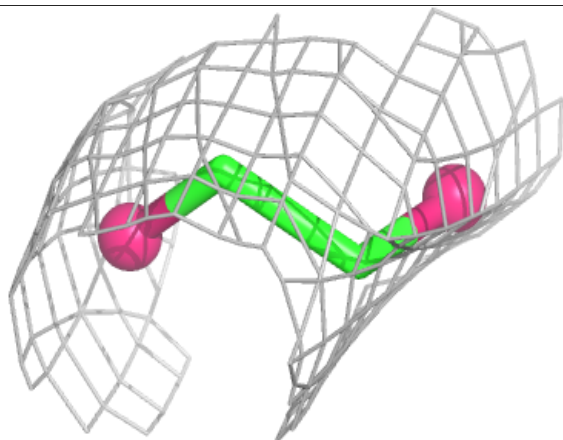
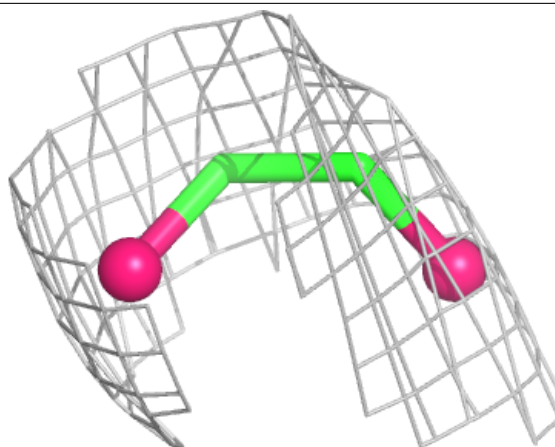
**Electron density around EDO D 409:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

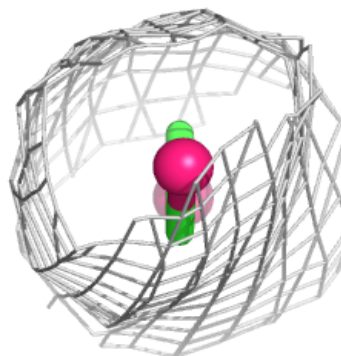
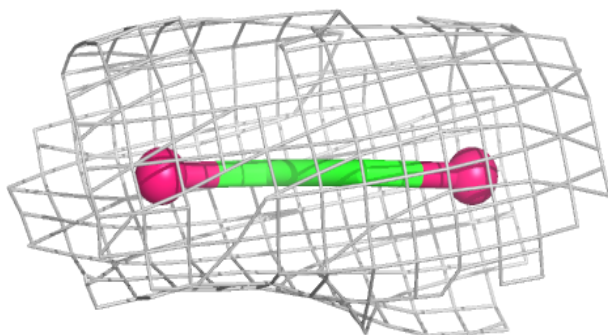
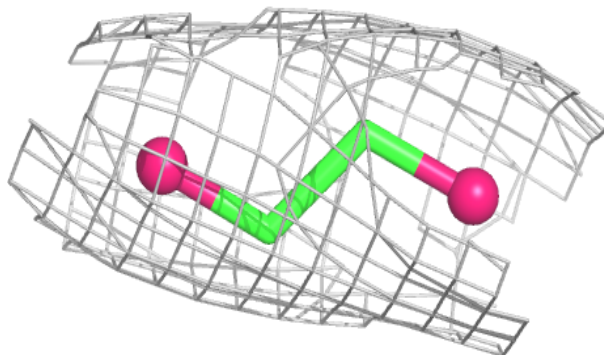


**Electron density around EDO A 441:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

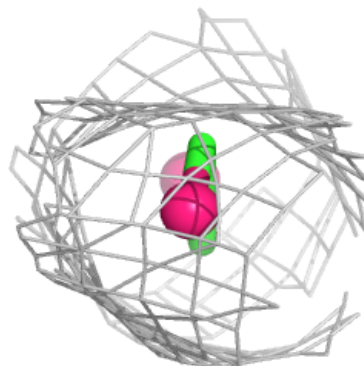
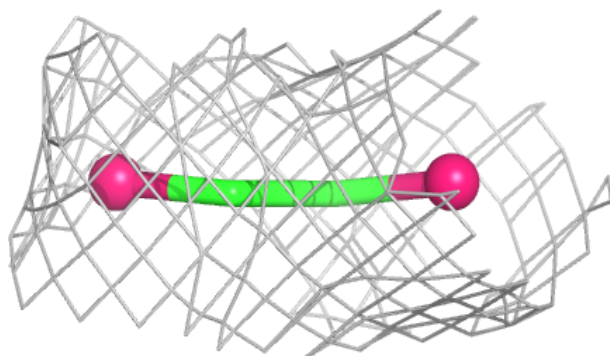
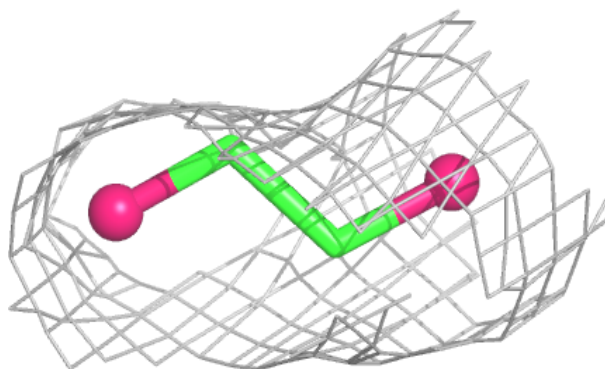
**Electron density around EDO A 409:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

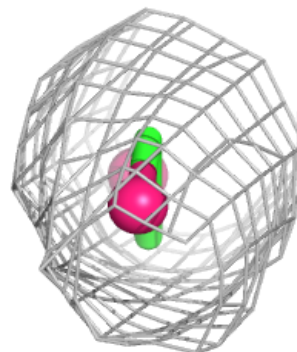
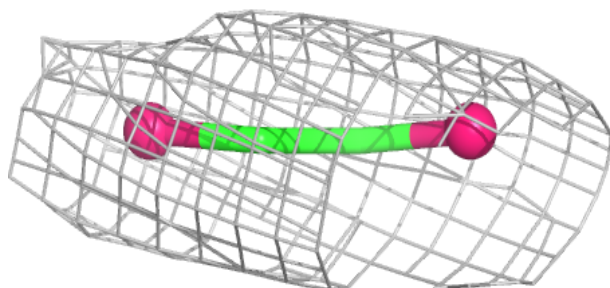
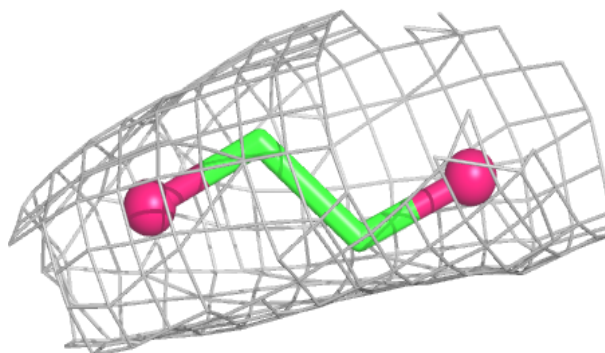


**Electron density around EDO A 414:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

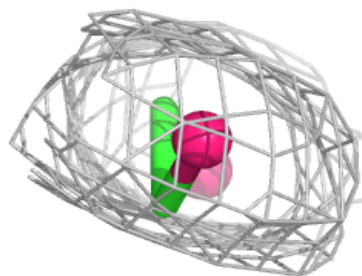
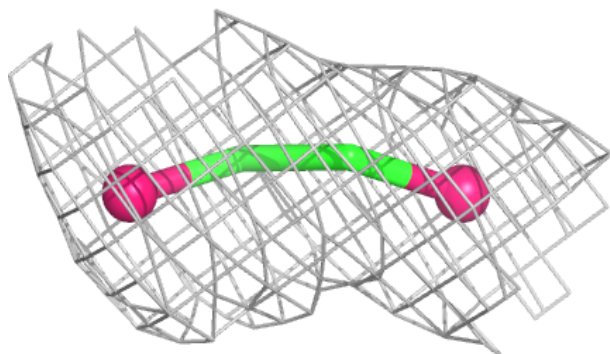
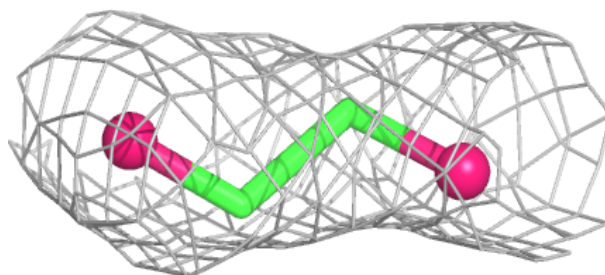
**Electron density around EDO C 421:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

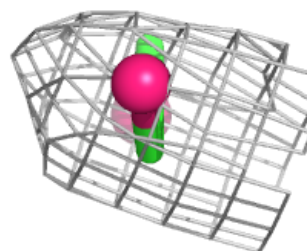
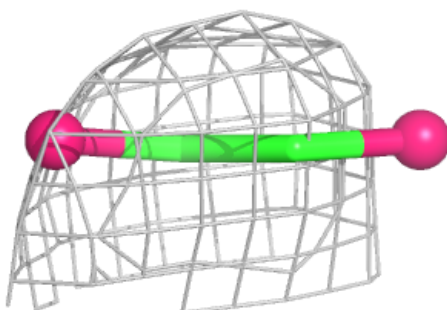
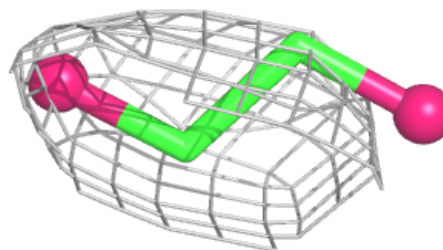


**Electron density around EDO D 414:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around EDO F 410:**

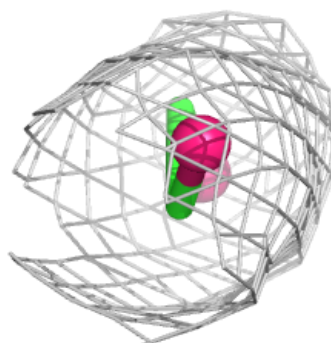
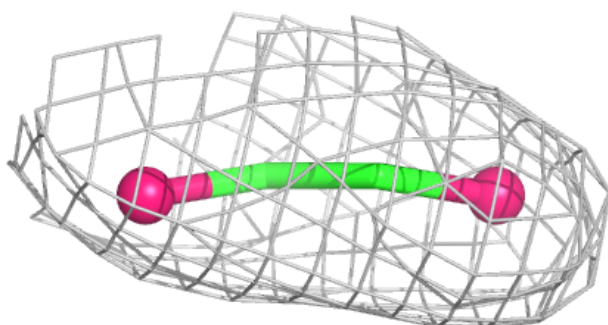
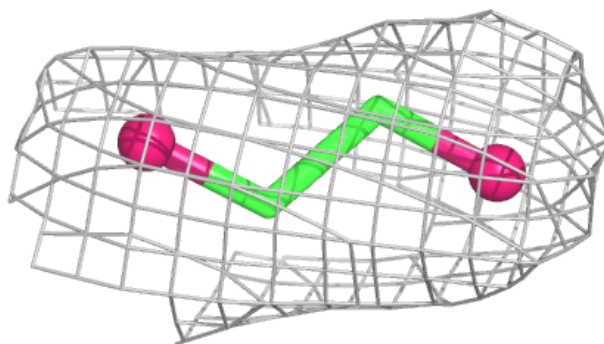
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



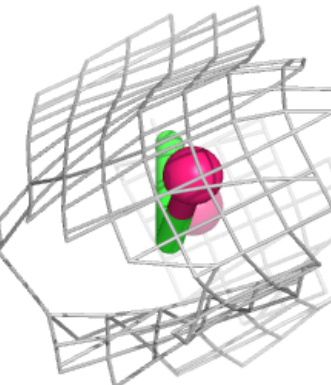
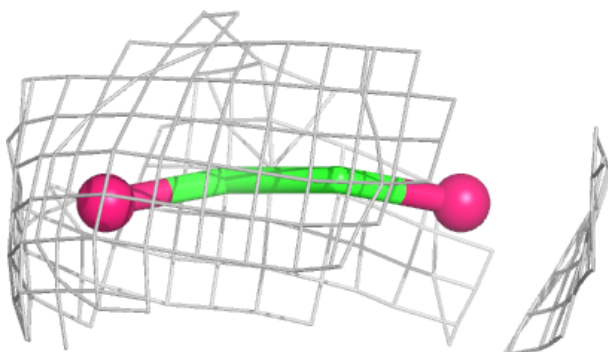
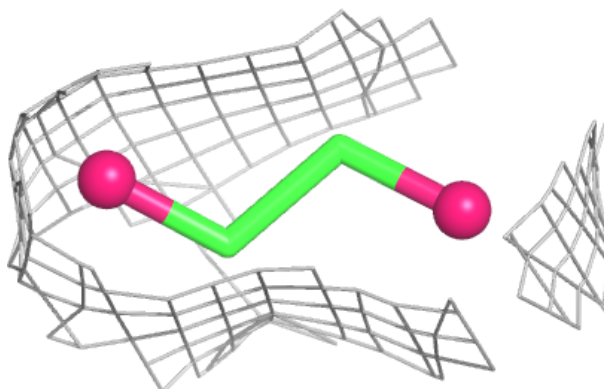


**Electron density around EDO A 418:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

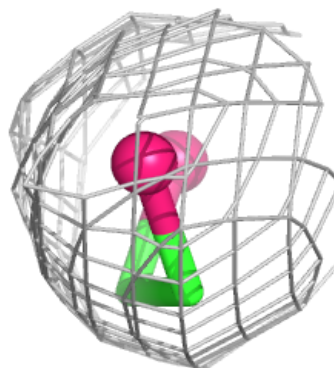
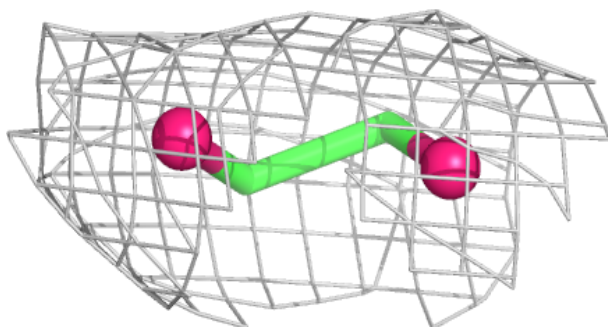
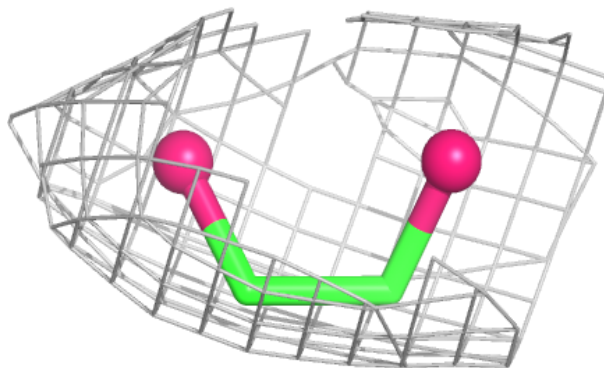
**Electron density around EDO A 415:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

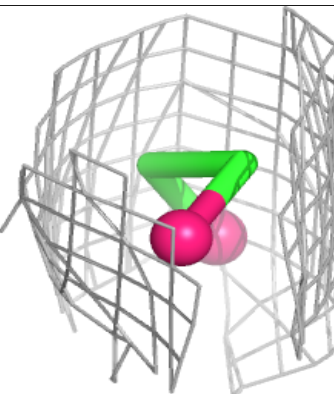
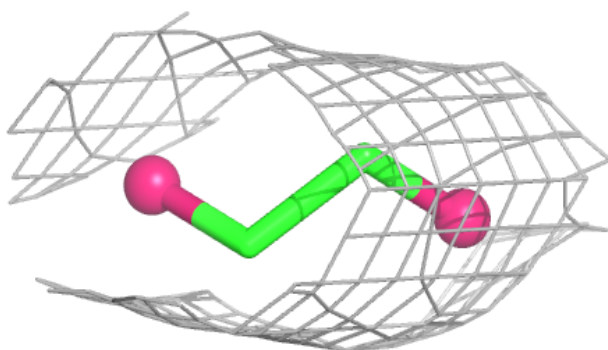
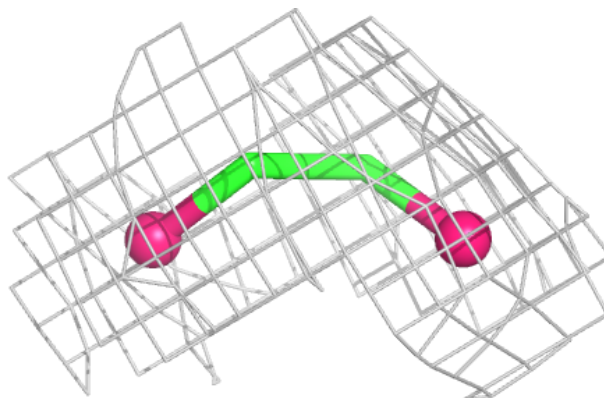


**Electron density around EDO E 436:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

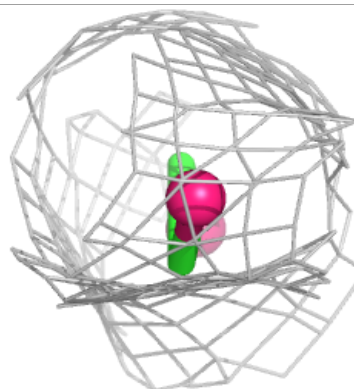
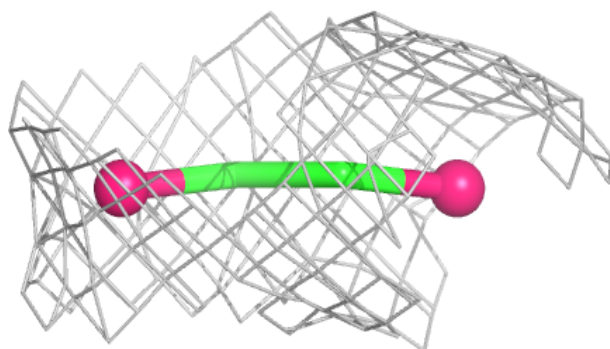
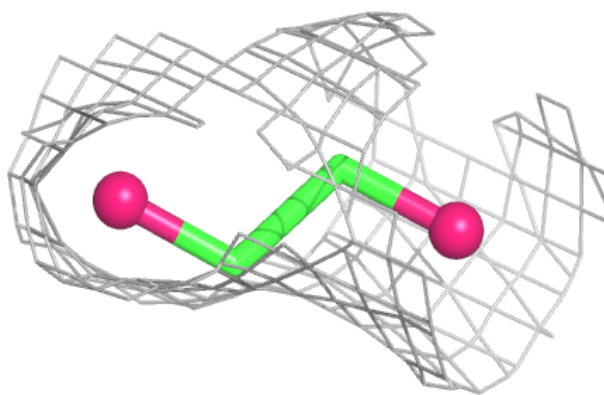
**Electron density around EDO F 414:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

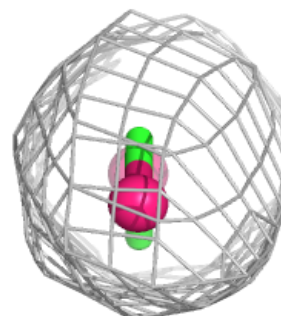
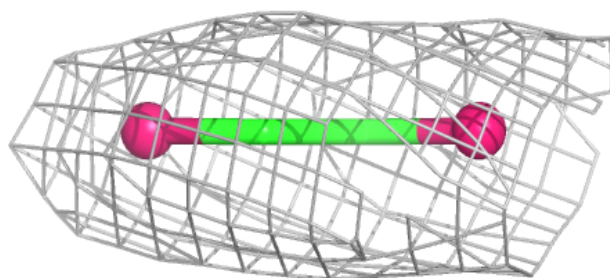
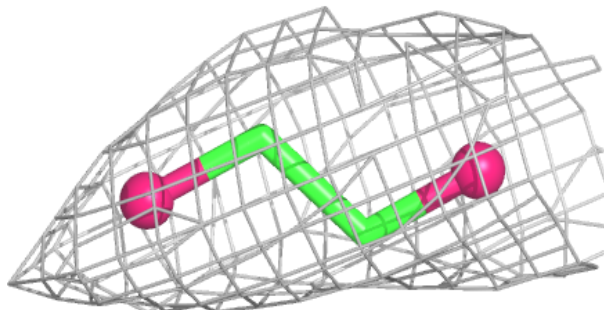


**Electron density around EDO D 417:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around EDO A 420:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



## 6.5 Other polymers [i](#)

There are no such residues in this entry.