



wwPDB X-ray Structure Validation Summary Report ⓘ

May 6, 2024 – 11:53 am BST

PDB ID : 8RHQ
Title : Crystal structure of HLA-A*11:01 in complex with SVLNDIFSRL, an 10-mer epitope from SARS-CoV-2 Spike (S975-984)
Authors : Ahn, Y.M.; Maddumage, J.C.; Szeto, C.; Gras, S.
Deposited on : 2023-12-16
Resolution : 2.00 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

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>.

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

MolProbity : 4.02b-467
Mogul : 1.8.4, CSD as541be (2020)
Xtrriage (Phenix) : 1.13
EDS : 2.36.2
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36.2

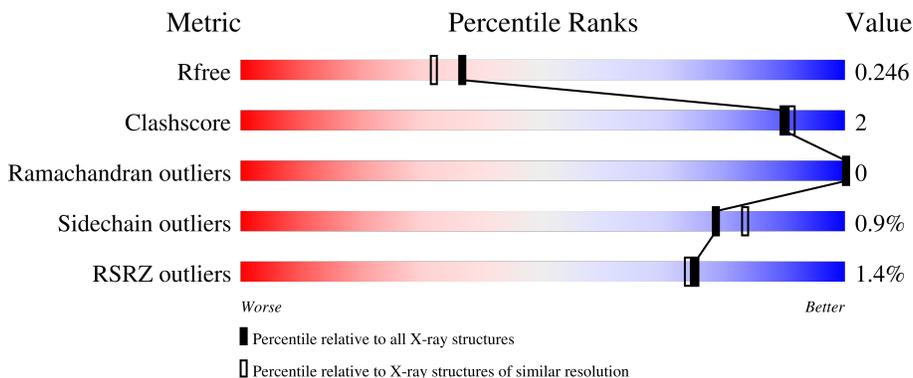
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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.



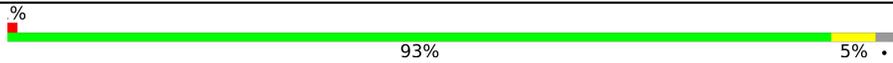
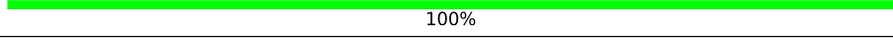
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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 ≥ 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	365	
1	C	365	
1	E	365	
2	B	100	
2	D	100	

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Mol	Chain	Length	Quality of chain
2	F	100	 <p>% 93% 5%</p>
3	G	10	 <p>100%</p>
3	H	10	 <p>90% 10%</p>
3	I	10	 <p>100%</p>

2 Entry composition [i](#)

There are 9 unique types of molecules in this entry. The entry contains 10187 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 HLA class I histocompatibility antigen.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	273	Total 2255	C 1403	N 408	O 433	S 11	0	3	0
1	C	274	Total 2274	C 1411	N 419	O 434	S 10	0	6	0
1	E	273	Total 2245	C 1395	N 407	O 433	S 10	0	3	0

- Molecule 2 is a protein called Beta-2-microglobulin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	99	Total 845	C 539	N 143	O 160	S 3	0	4	0
2	D	100	Total 852	C 544	N 142	O 162	S 4	0	4	0
2	F	98	Total 836	C 534	N 142	O 158	S 2	0	4	0

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	0	MET	-	initiating methionine	UNP P61769
D	0	MET	-	initiating methionine	UNP P61769
F	0	MET	-	initiating methionine	UNP P61769

- Molecule 3 is a protein called Spike protein S2'.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
3	G	10	Total 82	C 52	N 14	O 16	0	0	0
3	H	10	Total 82	C 52	N 14	O 16	0	0	0

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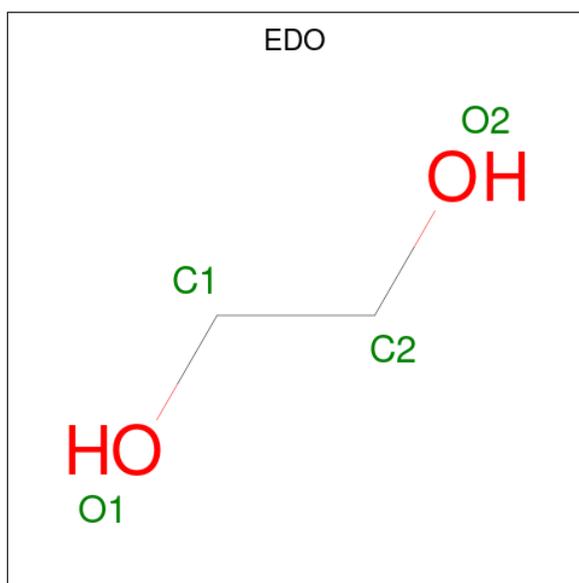
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
3	I	10	88	55	15	18	0	1	0

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
G	7	PHE	LEU	variant	UNP P0DTC2
H	7	PHE	LEU	variant	UNP P0DTC2
I	7	PHE	LEU	variant	UNP P0DTC2

- Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂) (labeled as "Ligand of Interest" by depositor).



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

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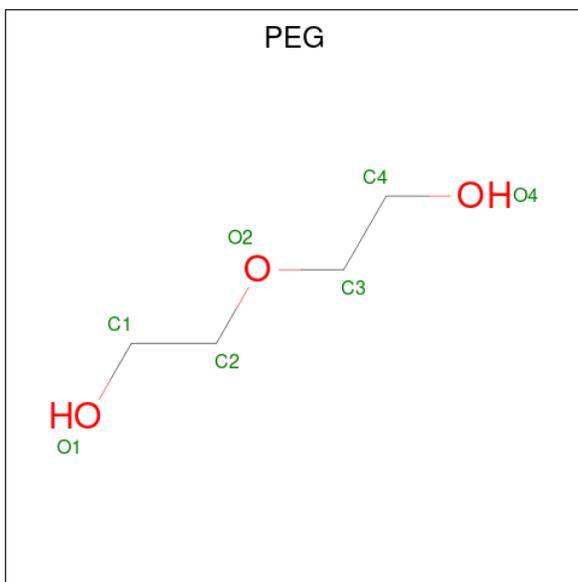
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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total C O 4 2 2	0	0
4	C	1	Total C O 4 2 2	0	0
4	C	1	Total C O 4 2 2	0	0
4	C	1	Total C O 4 2 2	0	0
4	C	1	Total C O 4 2 2	0	0
4	D	1	Total C O 4 2 2	0	0
4	D	1	Total C O 4 2 2	0	0
4	D	1	Total C O 4 2 2	0	0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

- Molecule 5 is SODIUM ION (three-letter code: NA) (formula: Na) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	B	1	Total Na 1 1	0	0
5	E	1	Total Na 1 1	0	0
5	F	1	Total Na 1 1	0	0

- Molecule 6 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C₄H₁₀O₃) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	C	1	Total	C O	0	0
			7	4 3		

- Molecule 7 is CHLORIDE ION (three-letter code: CL) (formula: Cl) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	C	1	Total	Cl	0	0
			1	1		
7	E	2	Total	Cl	0	0
			2	2		

- Molecule 8 is MAGNESIUM ION (three-letter code: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	D	1	Total	Mg	0	0
			1	1		

- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	121	Total	O	0	0
			121	121		
9	B	38	Total	O	0	0
			38	38		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	C	144	Total 144	O 144	0	0
9	D	36	Total 36	O 36	0	0
9	E	143	Total 143	O 143	0	0
9	F	37	Total 37	O 37	0	0
9	G	9	Total 9	O 9	0	0
9	H	6	Total 6	O 6	0	0
9	I	12	Total 12	O 12	0	0



- Molecule 2: Beta-2-microglobulin



- Molecule 2: Beta-2-microglobulin



- Molecule 3: Spike protein S2'



There are no outlier residues recorded for this chain.

- Molecule 3: Spike protein S2'



- Molecule 3: Spike protein S2'



There are no outlier residues recorded for this chain.

4 Data and refinement statistics i

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	85.49Å 85.49Å 312.03Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	42.75 – 2.00 42.75 – 2.00	Depositor EDS
% Data completeness (in resolution range)	100.0 (42.75-2.00) 100.0 (42.75-2.00)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.30 (at 2.00Å)	Xtrriage
Refinement program	PHENIX 1.21rc1_5127	Depositor
R, R_{free}	0.198 , 0.244 0.199 , 0.246	Depositor DCC
R_{free} test set	4539 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	25.4	Xtrriage
Anisotropy	0.088	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 43.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtrriage
Estimated twinning fraction	0.036 for -h,-k,l	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	10187	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 23.88 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 4.2775e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CL, NA, MG, EDO, PEG

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	0.29	0/2320	0.58	0/3147
1	C	0.30	0/2352	0.59	0/3189
1	E	0.30	0/2314	0.56	0/3139
2	B	0.29	0/882	0.52	0/1192
2	D	0.30	0/886	0.53	0/1198
2	F	0.29	0/873	0.53	0/1182
3	G	0.31	0/82	0.64	0/108
3	H	0.35	0/82	0.63	0/108
3	I	0.33	0/88	0.62	0/116
All	All	0.30	0/9879	0.56	0/13379

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	2255	0	2107	12	0
1	C	2274	0	2132	9	0
1	E	2245	0	2092	2	0
2	B	845	0	818	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	852	0	822	4	0
2	F	836	0	809	5	0
3	G	82	0	86	0	0
3	H	82	0	86	1	0
3	I	88	0	90	0	0
4	A	20	0	30	0	0
4	B	8	0	12	2	0
4	C	16	0	24	2	0
4	D	12	0	18	0	0
4	E	4	0	6	0	0
4	F	8	0	12	0	0
5	B	1	0	0	0	0
5	E	1	0	0	0	0
5	F	1	0	0	0	0
6	C	7	0	10	0	0
7	C	1	0	0	0	0
7	E	2	0	0	0	0
8	D	1	0	0	0	0
9	A	121	0	0	0	0
9	B	38	0	0	0	0
9	C	144	0	0	0	0
9	D	36	0	0	0	0
9	E	143	0	0	0	0
9	F	37	0	0	0	0
9	G	9	0	0	0	0
9	H	6	0	0	0	0
9	I	12	0	0	0	0
All	All	10187	0	9154	33	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 2.

The worst 5 of 33 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:C:163:ARG:NH1	3:H:4:ASN:OD1	2.23	0.72
1:A:21:ARG:NH1	1:A:39:ASP:OD2	2.32	0.62
1:A:44:ARG:NH1	1:A:61:ASP:OD1	2.34	0.61
2:B:16:GLU:HB3	2:B:19:LYS:HD2	1.83	0.59
1:C:258:THR:HG22	1:C:273:ARG:HG2	1.87	0.57

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	272/365 (74%)	265 (97%)	7 (3%)	0	100	100
1	C	278/365 (76%)	274 (99%)	4 (1%)	0	100	100
1	E	272/365 (74%)	269 (99%)	3 (1%)	0	100	100
2	B	101/100 (101%)	100 (99%)	1 (1%)	0	100	100
2	D	102/100 (102%)	100 (98%)	2 (2%)	0	100	100
2	F	100/100 (100%)	98 (98%)	2 (2%)	0	100	100
3	G	8/10 (80%)	8 (100%)	0	0	100	100
3	H	8/10 (80%)	8 (100%)	0	0	100	100
3	I	9/10 (90%)	9 (100%)	0	0	100	100
All	All	1150/1425 (81%)	1131 (98%)	19 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	235/300 (78%)	234 (100%)	1 (0%)	91	93
1	C	237/300 (79%)	235 (99%)	2 (1%)	81	86
1	E	234/300 (78%)	233 (100%)	1 (0%)	91	93

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	B	98/95 (103%)	95 (97%)	3 (3%)	40	40
2	D	99/95 (104%)	97 (98%)	2 (2%)	55	58
2	F	97/95 (102%)	96 (99%)	1 (1%)	76	81
3	G	10/10 (100%)	10 (100%)	0	100	100
3	H	10/10 (100%)	10 (100%)	0	100	100
3	I	11/10 (110%)	11 (100%)	0	100	100
All	All	1031/1215 (85%)	1021 (99%)	10 (1%)	78	81

5 of 10 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	D	75	LYS
1	E	231	VAL
2	F	70	PHE
2	B	88[B]	SER
1	C	180	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 25 ligands modelled in this entry, 7 are monoatomic - leaving 18 for Mogul analysis.

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
4	EDO	C	402	-	3,3,3	0.28	0	2,2,2	0.14	0
4	EDO	A	404	-	3,3,3	0.32	0	2,2,2	0.05	0
4	EDO	D	102	-	3,3,3	0.26	0	2,2,2	0.33	0
6	PEG	C	405	-	6,6,6	0.25	0	5,5,5	0.16	0
4	EDO	C	401	-	3,3,3	0.27	0	2,2,2	0.27	0
4	EDO	B	101	-	3,3,3	0.27	0	2,2,2	0.22	0
4	EDO	A	402	-	3,3,3	0.29	0	2,2,2	0.36	0
4	EDO	A	403	-	3,3,3	0.28	0	2,2,2	0.21	0
4	EDO	F	102	-	3,3,3	0.29	0	2,2,2	0.33	0
4	EDO	A	401	-	3,3,3	0.25	0	2,2,2	0.29	0
4	EDO	F	101	-	3,3,3	0.28	0	2,2,2	0.35	0
4	EDO	A	405	-	3,3,3	0.28	0	2,2,2	0.25	0
4	EDO	B	102	-	3,3,3	0.29	0	2,2,2	0.15	0
4	EDO	C	403	-	3,3,3	0.28	0	2,2,2	0.05	0
4	EDO	D	101	-	3,3,3	0.29	0	2,2,2	0.57	0
4	EDO	E	401	-	3,3,3	0.29	0	2,2,2	0.10	0
4	EDO	D	103	-	3,3,3	0.27	0	2,2,2	0.38	0
4	EDO	C	404	-	3,3,3	0.28	0	2,2,2	0.29	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
4	EDO	C	402	-	-	0/1/1/1	-
4	EDO	A	404	-	-	1/1/1/1	-
4	EDO	D	102	-	-	0/1/1/1	-
6	PEG	C	405	-	-	2/4/4/4	-
4	EDO	C	401	-	-	0/1/1/1	-
4	EDO	B	101	-	-	0/1/1/1	-
4	EDO	A	402	-	-	1/1/1/1	-
4	EDO	A	403	-	-	0/1/1/1	-
4	EDO	F	102	-	-	1/1/1/1	-
4	EDO	A	401	-	-	1/1/1/1	-
4	EDO	F	101	-	-	1/1/1/1	-
4	EDO	A	405	-	-	0/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	EDO	B	102	-	-	1/1/1/1	-
4	EDO	C	403	-	-	0/1/1/1	-
4	EDO	D	101	-	-	0/1/1/1	-
4	EDO	E	401	-	-	1/1/1/1	-
4	EDO	D	103	-	-	1/1/1/1	-
4	EDO	C	404	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

5 of 11 torsion outliers are listed below:

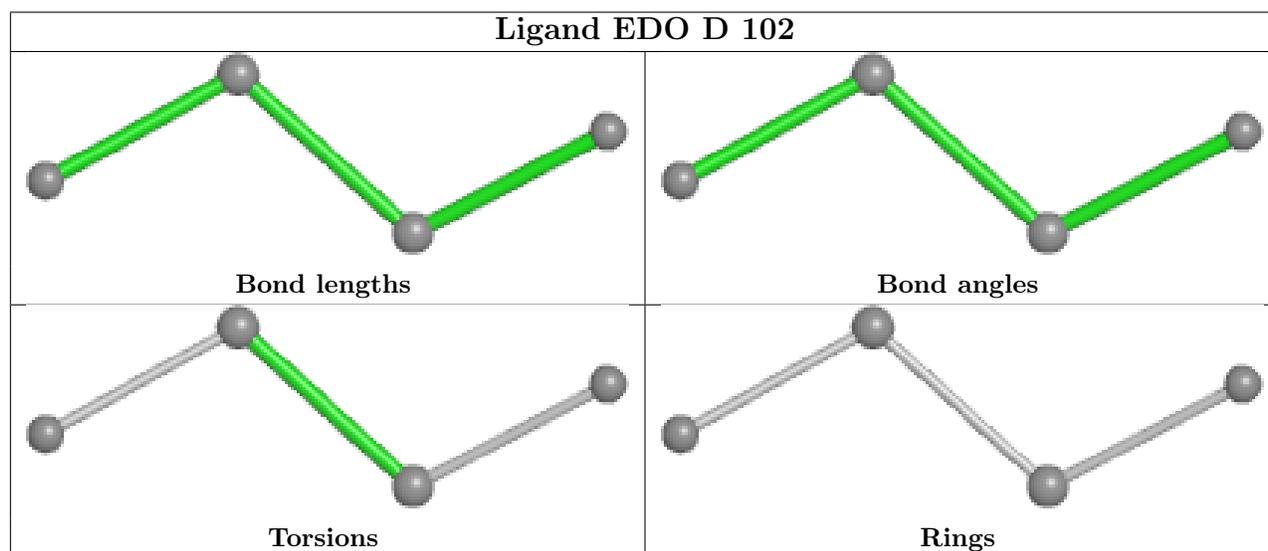
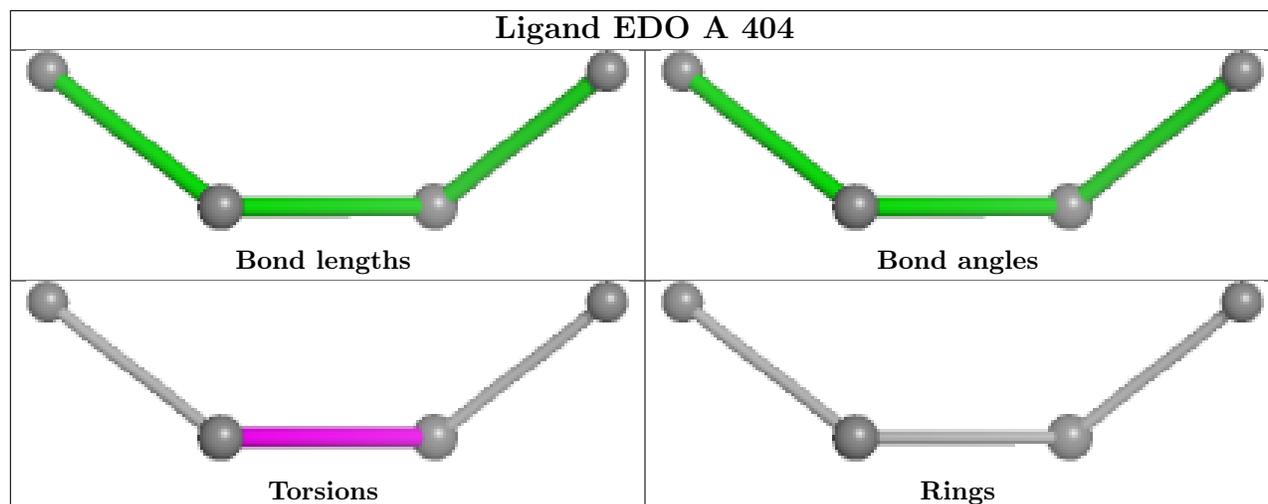
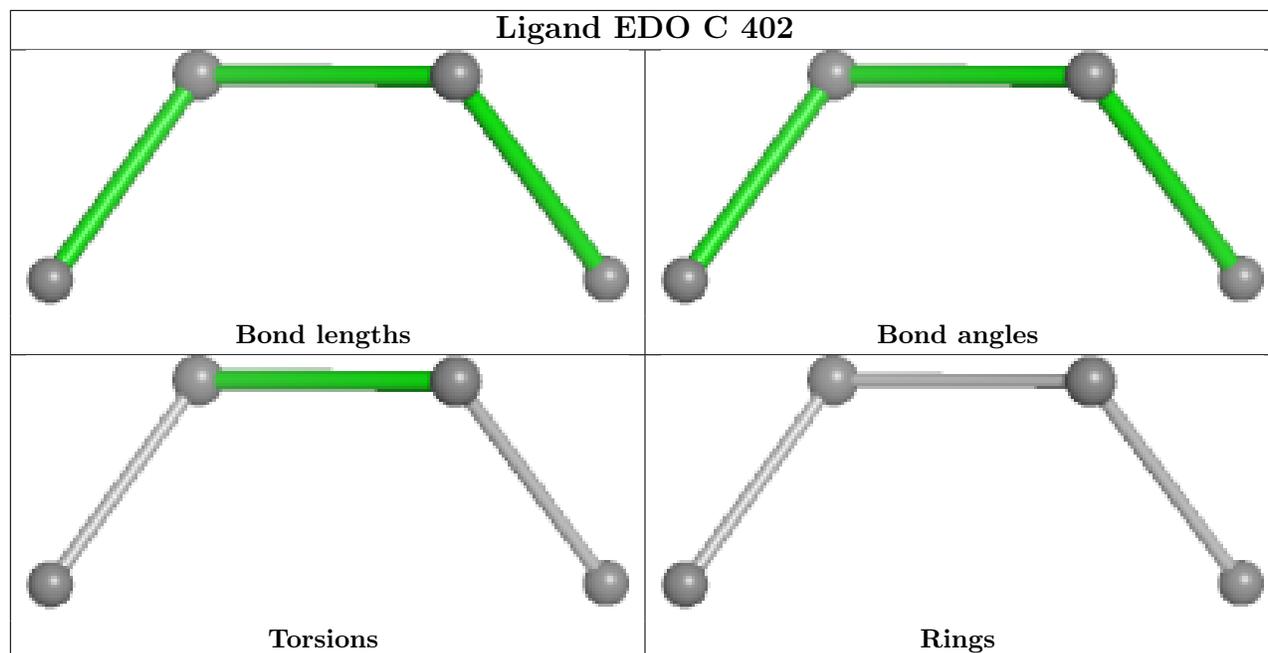
Mol	Chain	Res	Type	Atoms
6	C	405	PEG	O2-C3-C4-O4
4	A	404	EDO	O1-C1-C2-O2
6	C	405	PEG	O1-C1-C2-O2
4	D	103	EDO	O1-C1-C2-O2
4	F	102	EDO	O1-C1-C2-O2

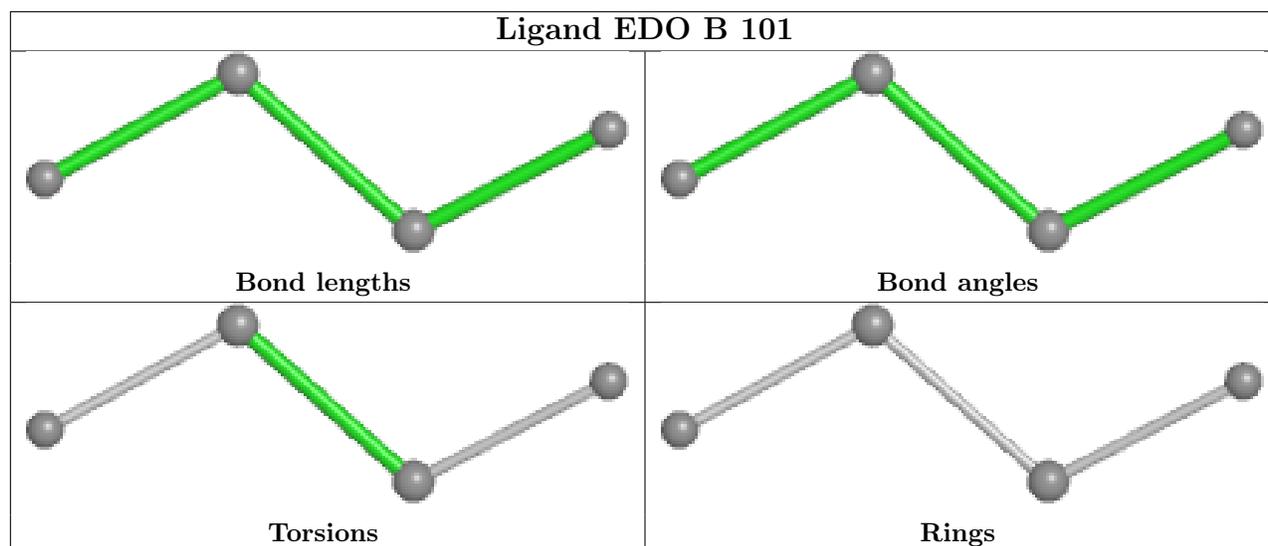
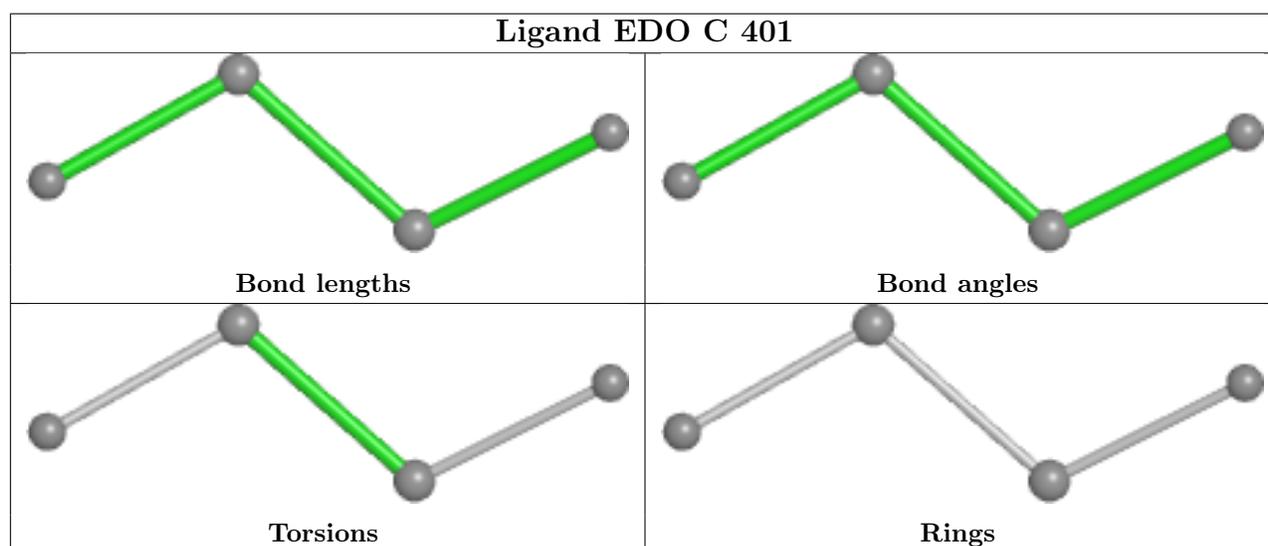
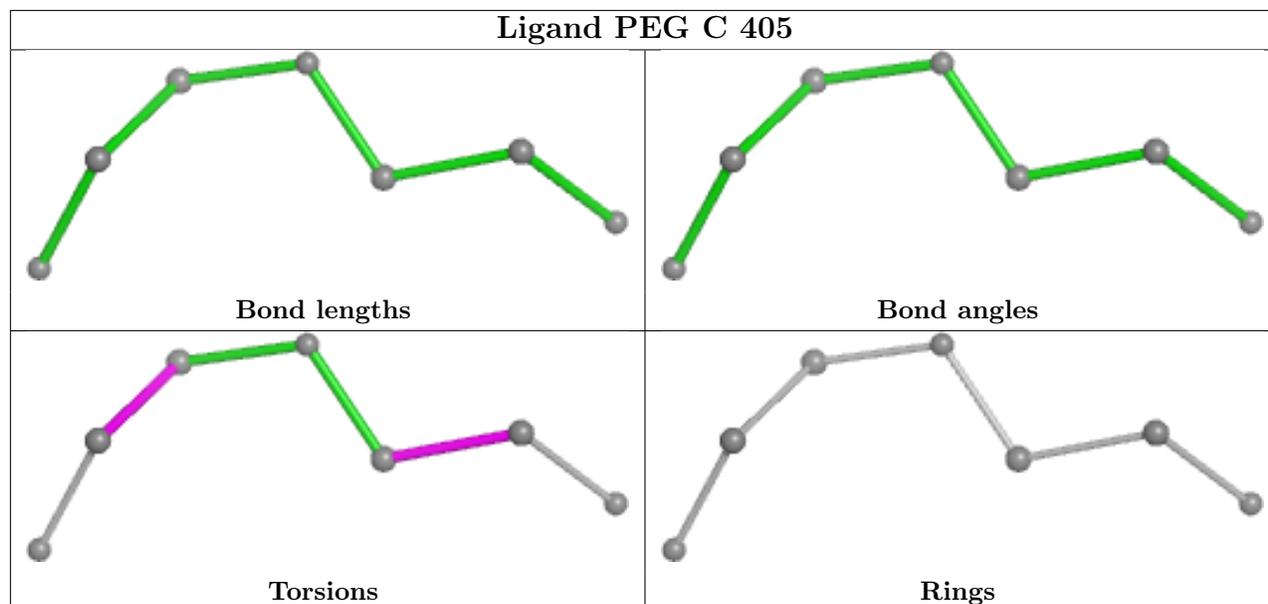
There are no ring outliers.

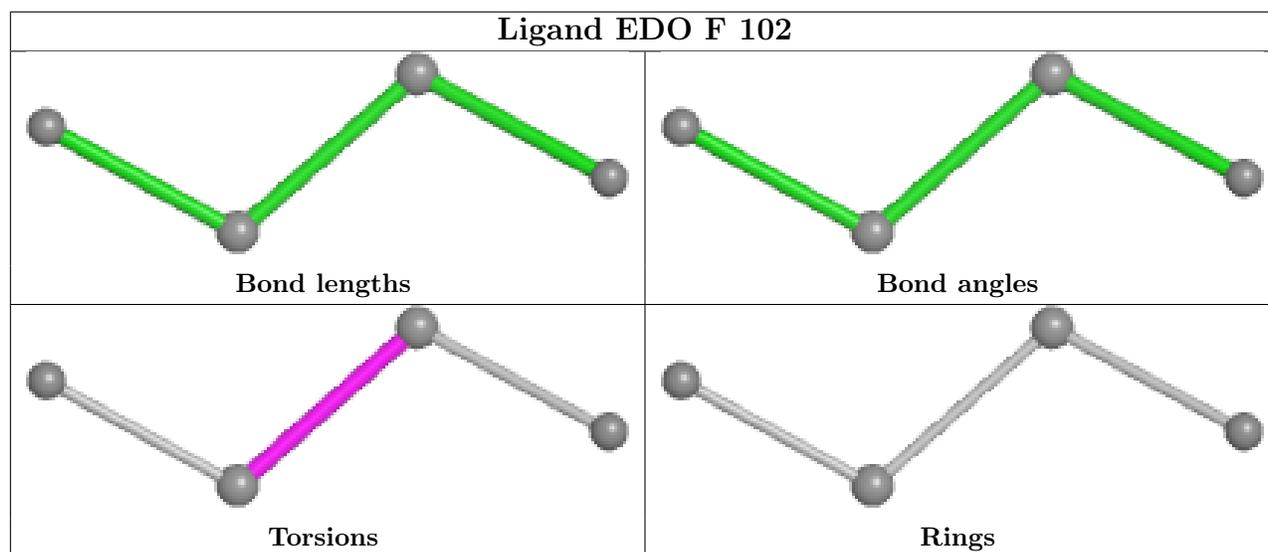
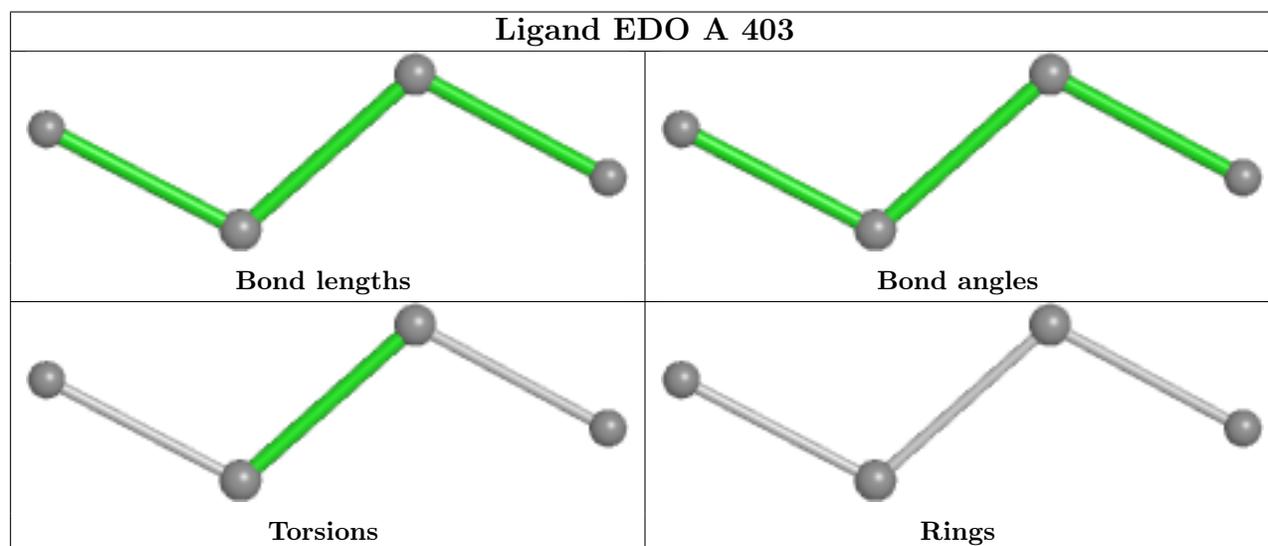
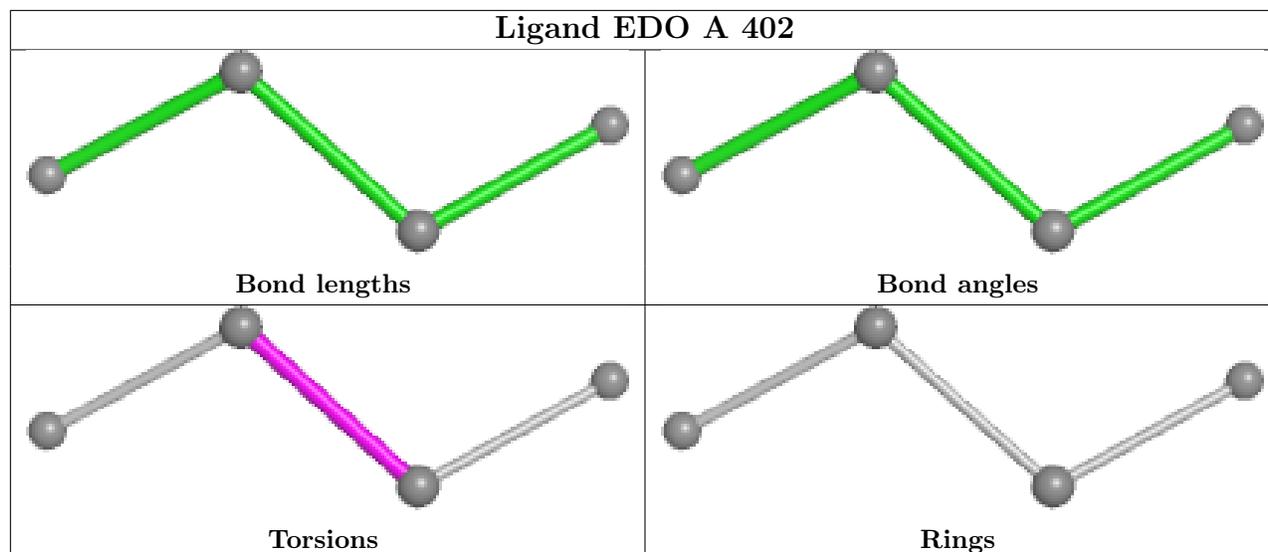
3 monomers are involved in 4 short contacts:

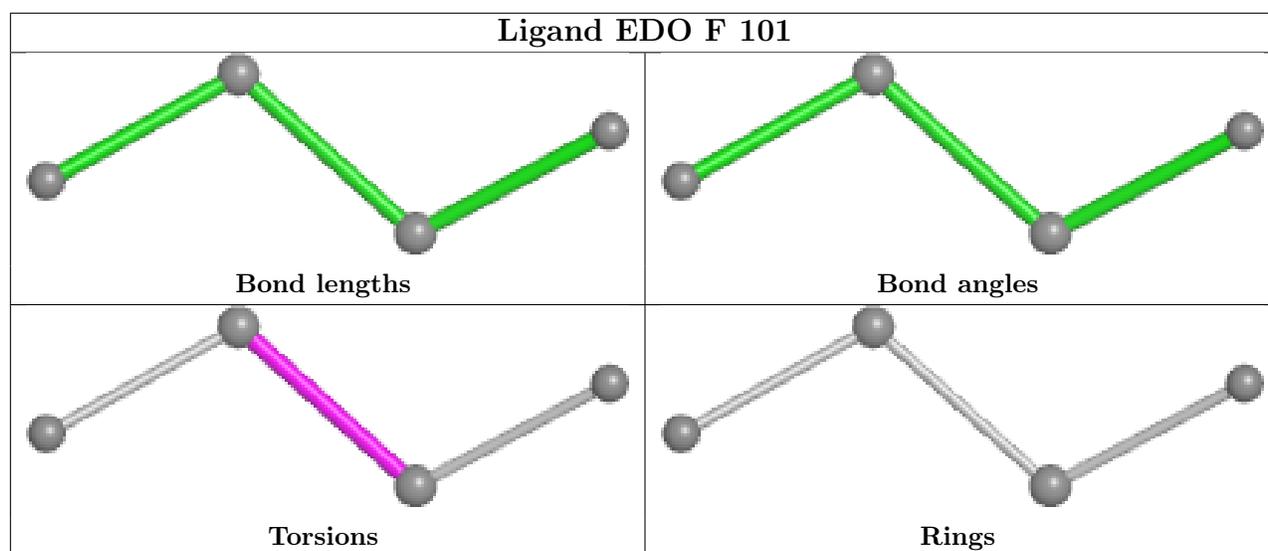
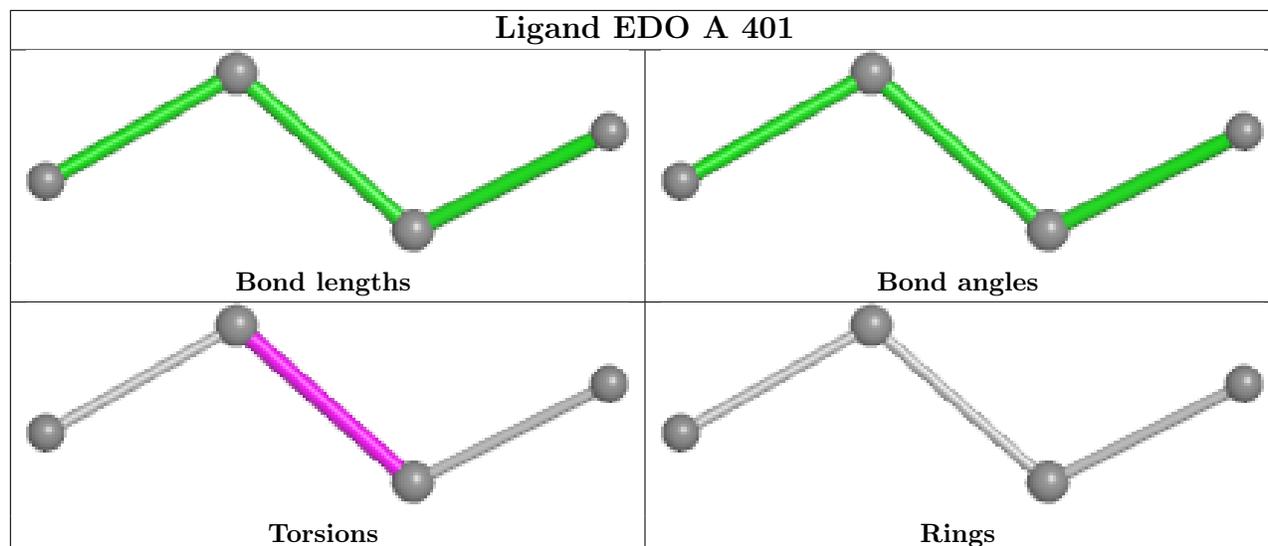
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	401	EDO	1	0
4	B	102	EDO	2	0
4	C	403	EDO	1	0

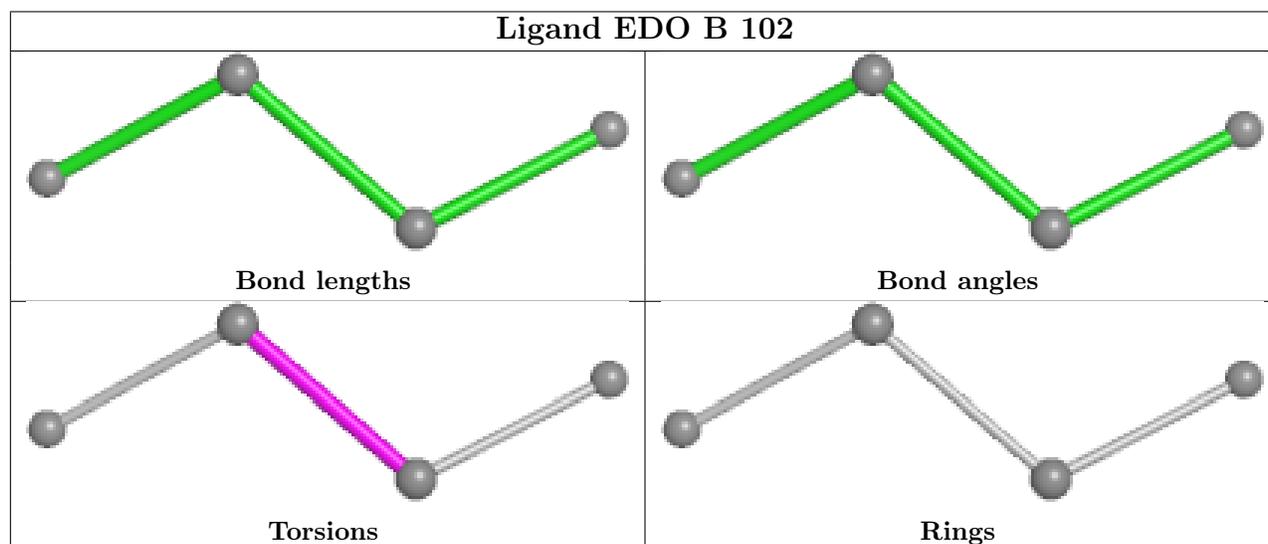
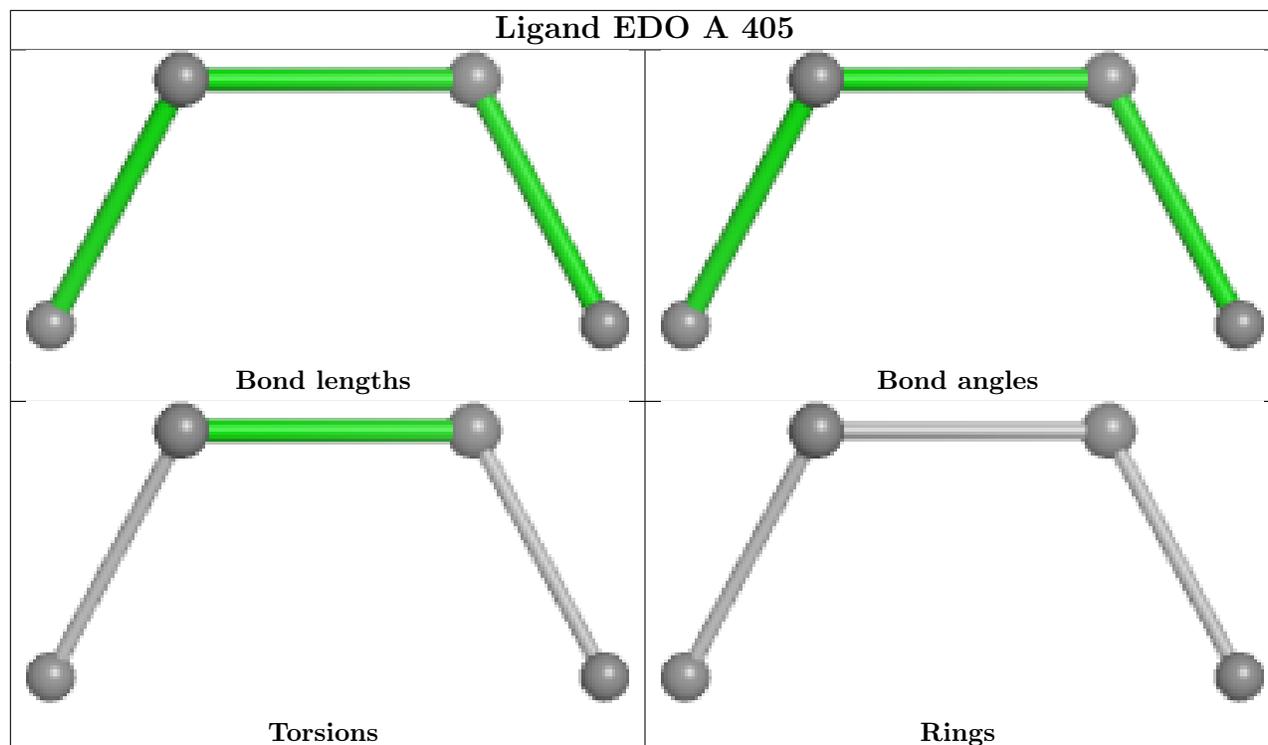
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

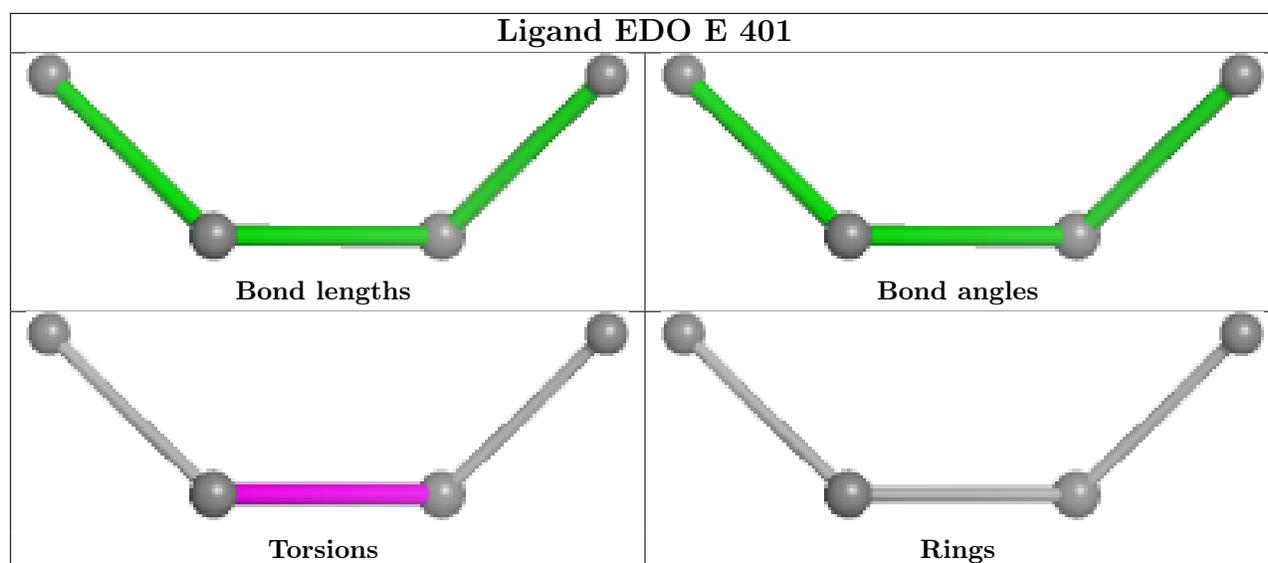
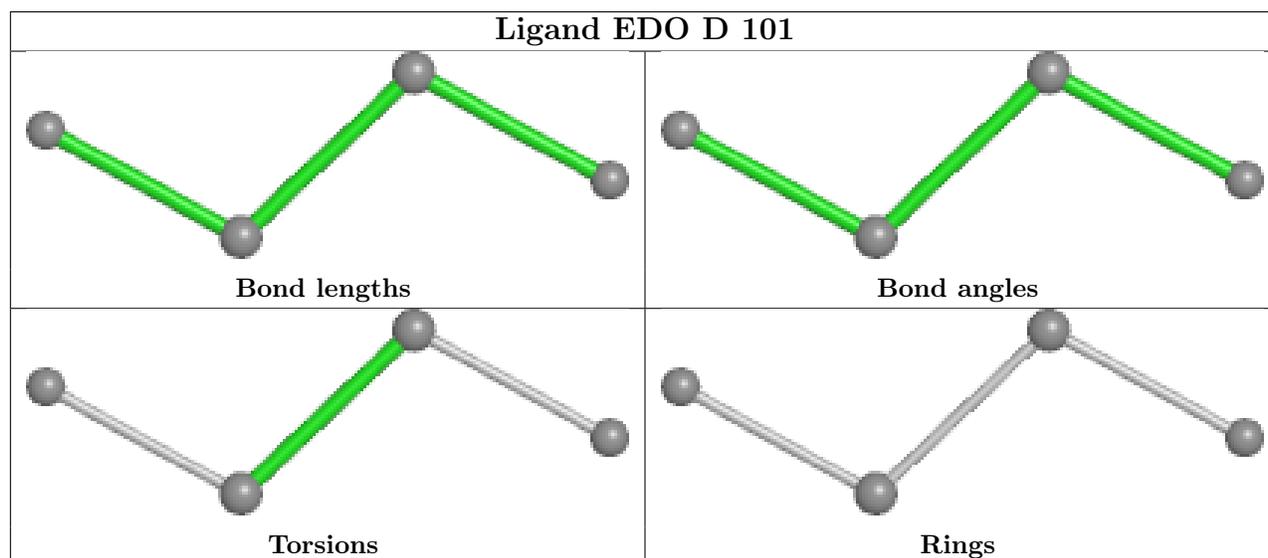
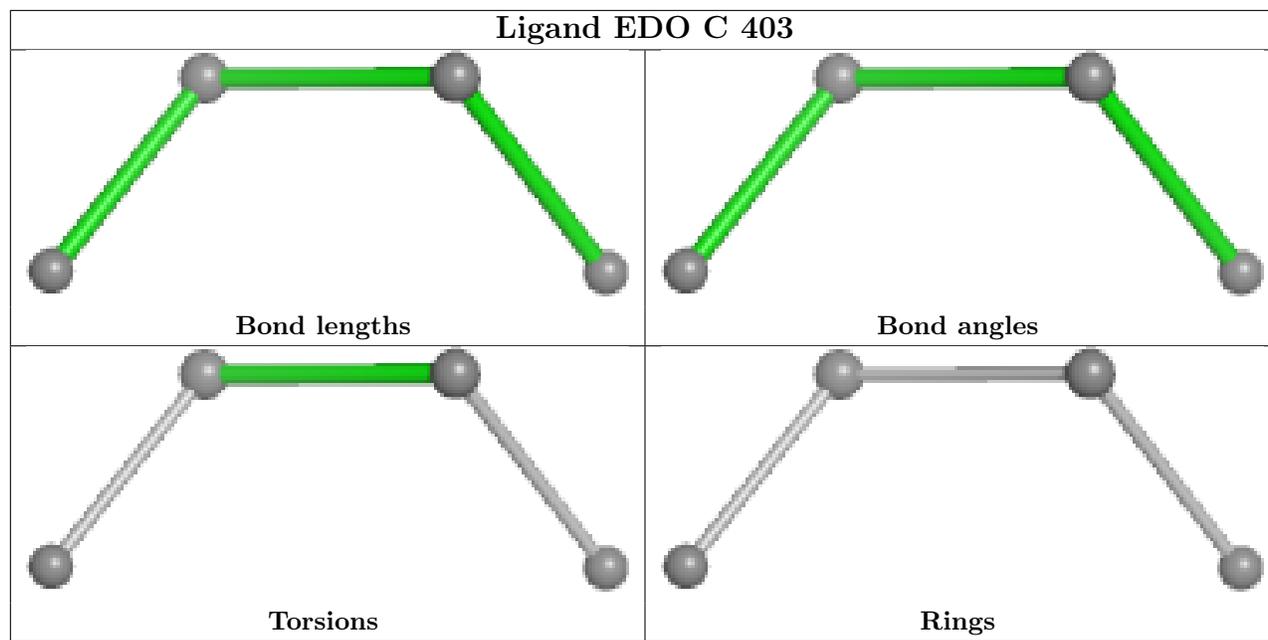


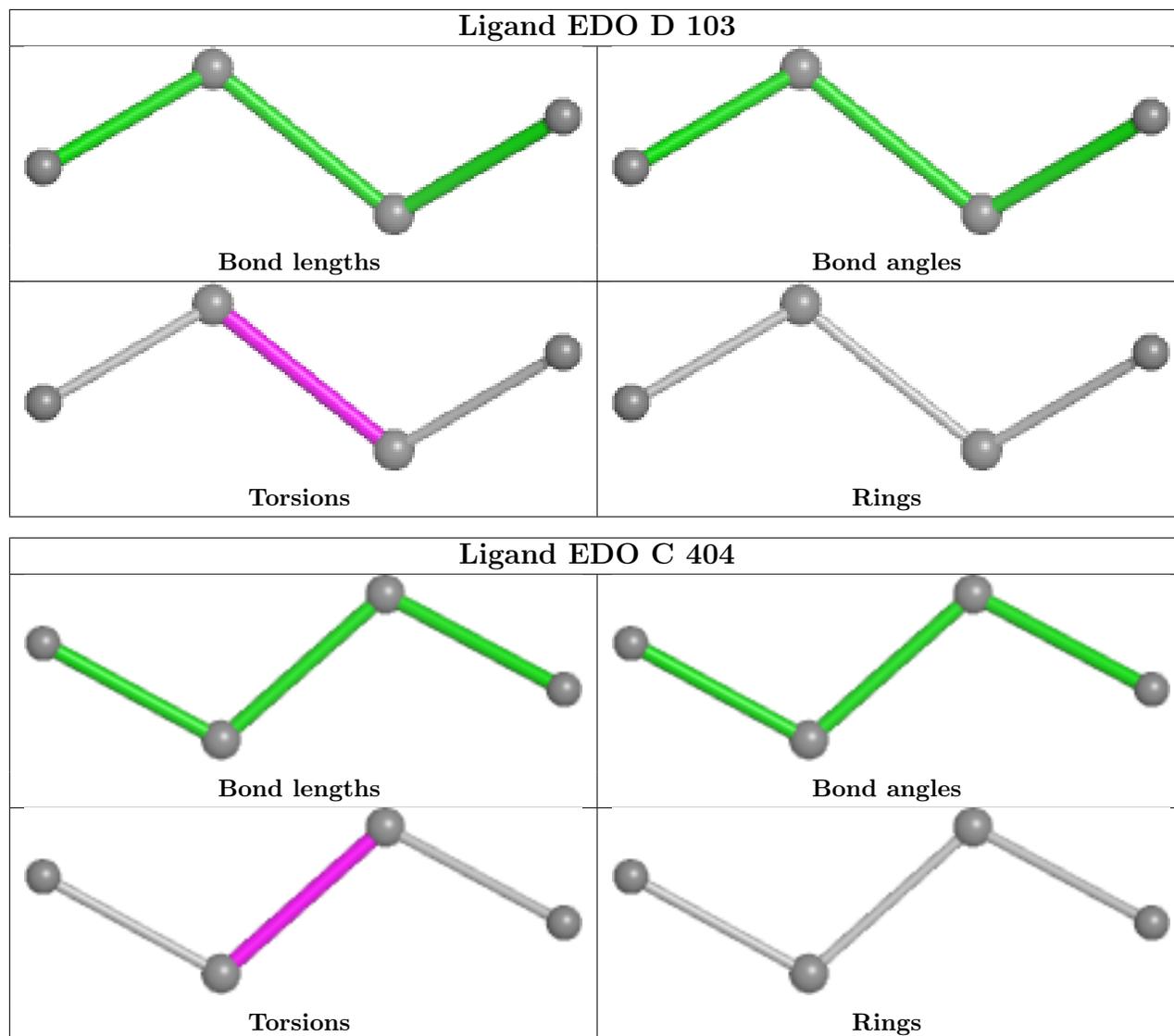












5.7 Other polymers [i](#)

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, 95th 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 ²)	Q<0.9
1	A	273/365 (74%)	-0.17	5 (1%) 68 66	13, 28, 50, 75	0
1	C	274/365 (75%)	-0.22	3 (1%) 80 79	12, 26, 47, 78	0
1	E	273/365 (74%)	-0.27	3 (1%) 80 79	12, 28, 46, 66	0
2	B	99/100 (99%)	-0.03	3 (3%) 50 49	17, 31, 57, 67	0
2	D	100/100 (100%)	-0.18	1 (1%) 82 81	14, 27, 48, 61	0
2	F	98/100 (98%)	-0.12	1 (1%) 82 81	17, 30, 52, 61	0
3	G	10/10 (100%)	-0.44	0 100 100	18, 23, 28, 29	0
3	H	10/10 (100%)	-0.42	0 100 100	14, 18, 22, 25	0
3	I	10/10 (100%)	-0.38	0 100 100	14, 18, 21, 26	0
All	All	1147/1425 (80%)	-0.19	16 (1%) 75 74	12, 28, 49, 78	0

The worst 5 of 16 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	18	GLY	4.1
1	E	90	ASP	3.8
1	A	41	ALA	3.8
2	B	75	LYS	2.9
1	E	197	HIS	2.8

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](#)

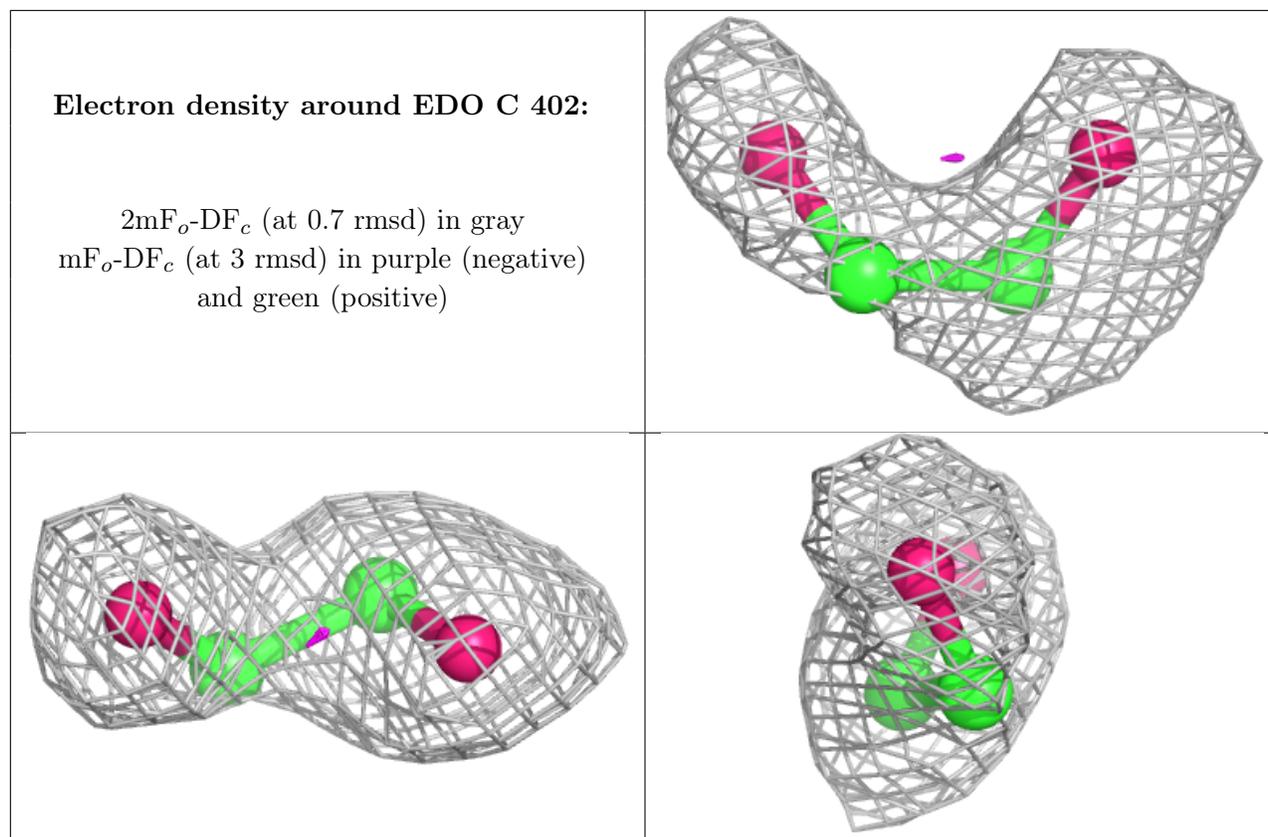
There are no monosaccharides in this entry.

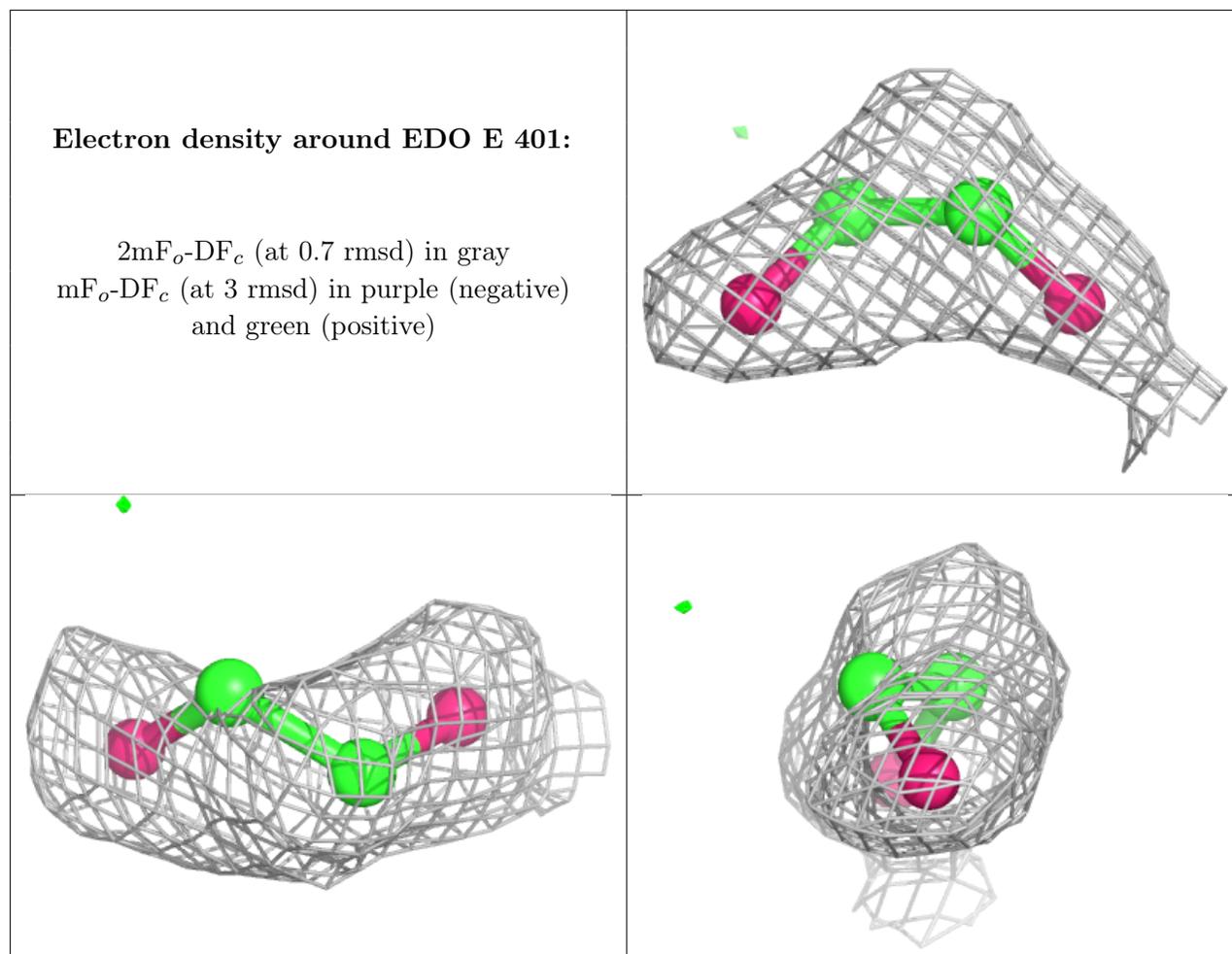
6.4 Ligands

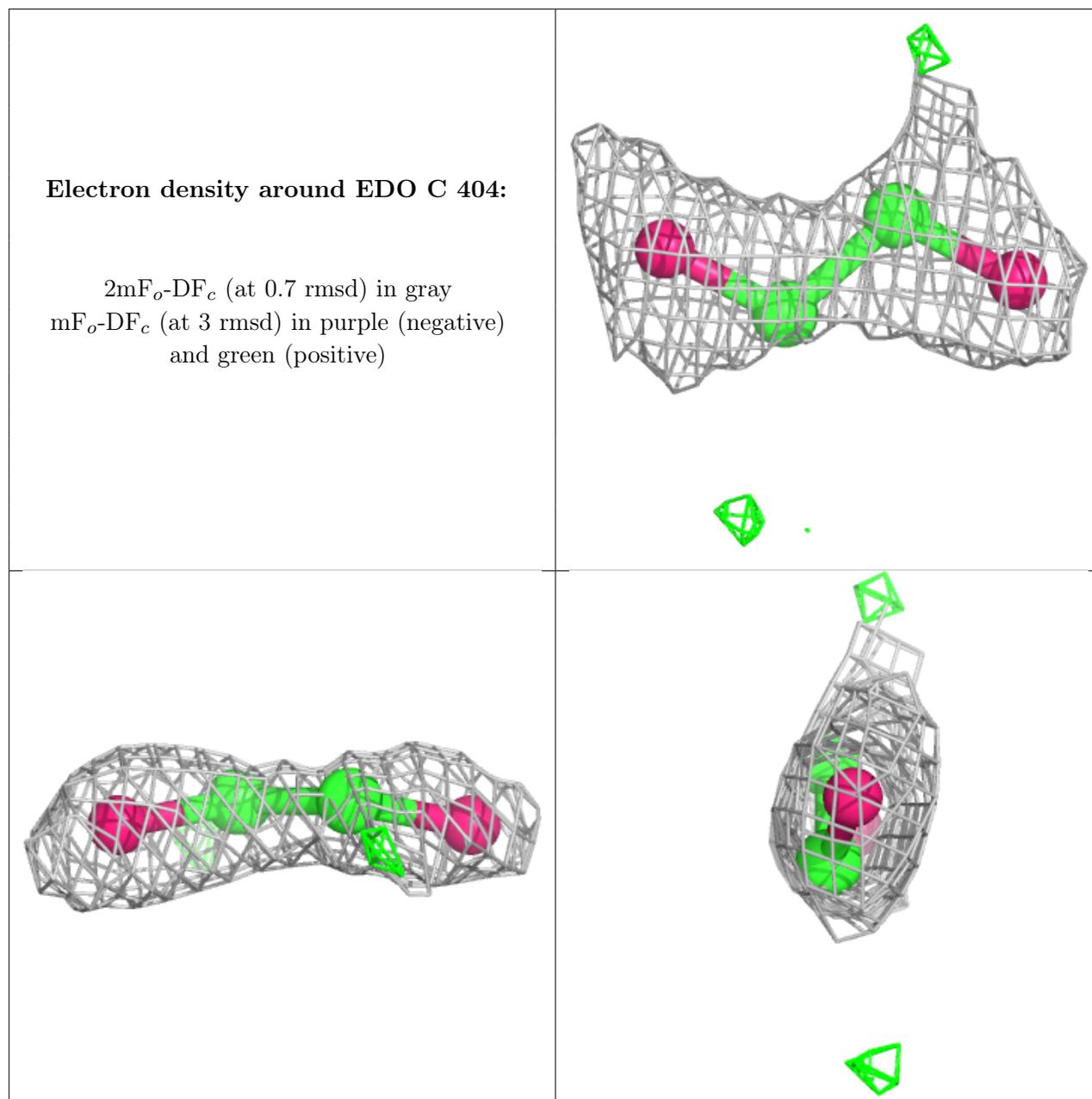
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	EDO	C	402	4/4	0.77	0.20	26,33,38,38	0
4	EDO	E	401	4/4	0.81	0.15	41,42,42,56	0
4	EDO	C	404	4/4	0.82	0.23	46,47,48,49	0
4	EDO	A	404	4/4	0.82	0.19	36,37,44,44	0
8	MG	D	104	1/1	0.82	0.15	29,29,29,29	0
4	EDO	B	102	4/4	0.85	0.34	26,39,46,51	0
4	EDO	B	101	4/4	0.86	0.37	34,45,45,58	0
6	PEG	C	405	7/7	0.87	0.20	22,32,52,53	0
4	EDO	A	402	4/4	0.87	0.29	28,35,37,39	0
5	NA	F	103	1/1	0.89	0.10	32,32,32,32	0
4	EDO	A	401	4/4	0.89	0.17	34,34,39,44	0
4	EDO	A	403	4/4	0.89	0.29	32,35,39,44	0
4	EDO	C	401	4/4	0.90	0.29	31,32,40,42	0
4	EDO	A	405	4/4	0.91	0.21	32,36,41,45	0
4	EDO	F	102	4/4	0.92	0.19	30,39,48,53	0
4	EDO	D	102	4/4	0.93	0.18	23,38,48,54	0
4	EDO	F	101	4/4	0.93	0.12	29,32,32,35	0
5	NA	B	103	1/1	0.94	0.06	33,33,33,33	0
4	EDO	C	403	4/4	0.94	0.27	25,29,40,43	0
4	EDO	D	103	4/4	0.95	0.18	24,32,34,39	0
7	CL	E	403	1/1	0.97	0.16	32,32,32,32	0
4	EDO	D	101	4/4	0.98	0.17	22,24,26,36	0
5	NA	E	404	1/1	0.99	0.09	17,17,17,17	0
7	CL	E	402	1/1	0.99	0.09	31,31,31,31	0
7	CL	C	406	1/1	1.00	0.09	32,32,32,32	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.

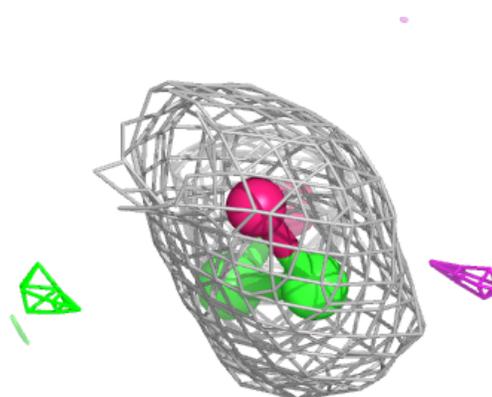
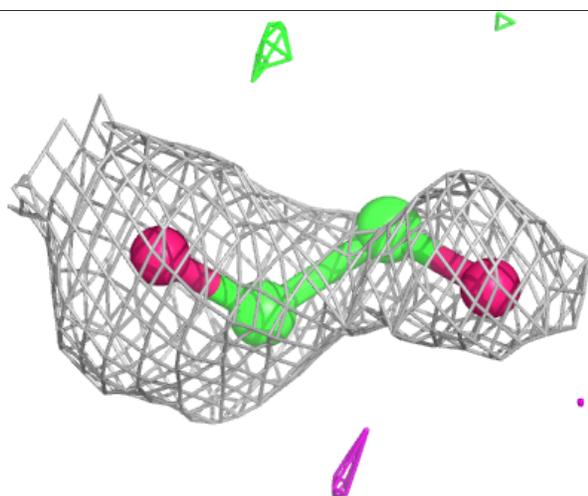
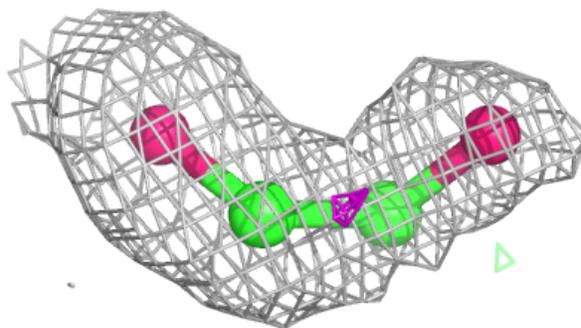






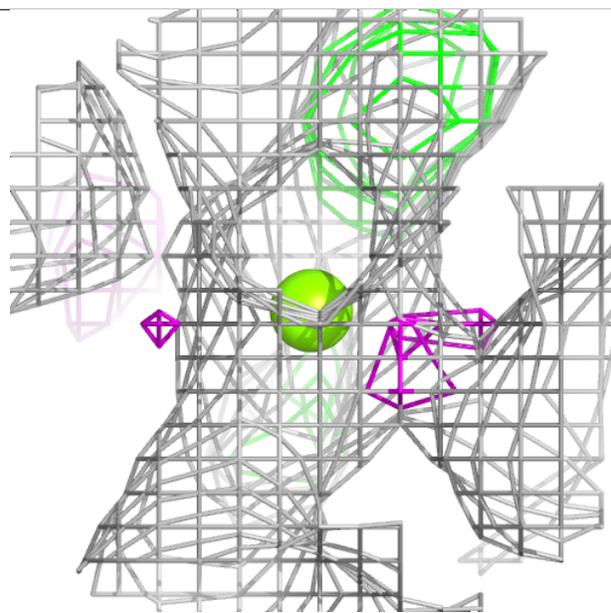
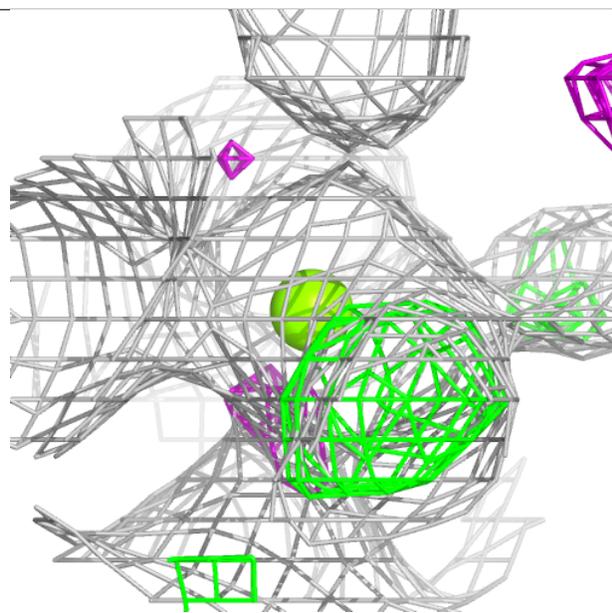
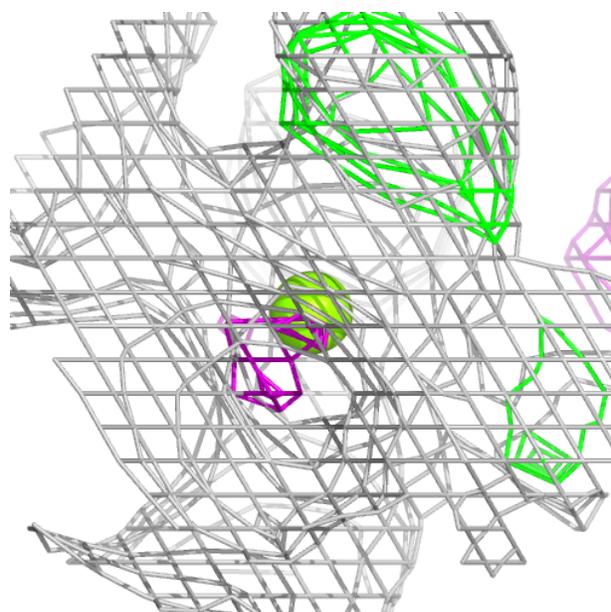
Electron density around EDO A 404:

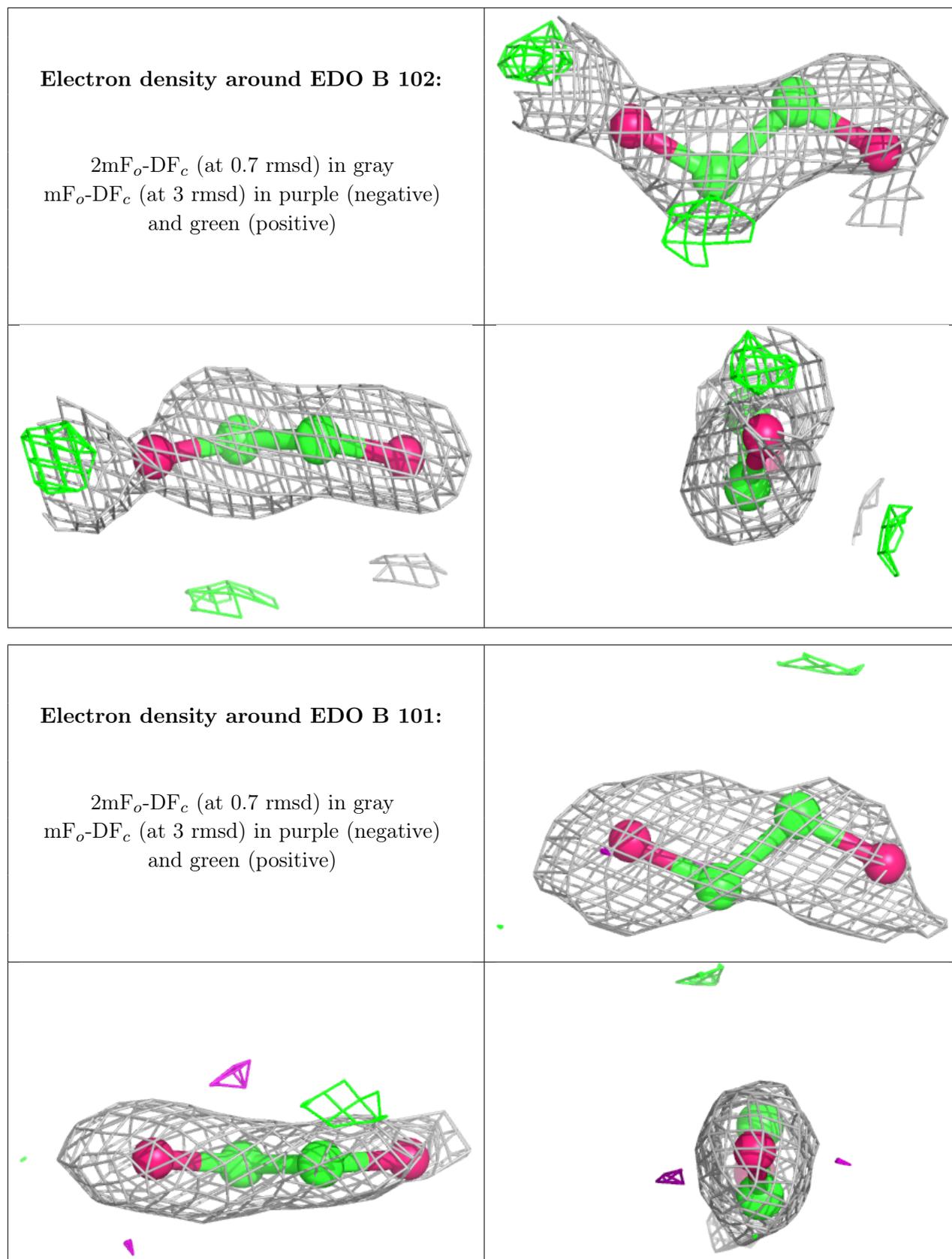
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

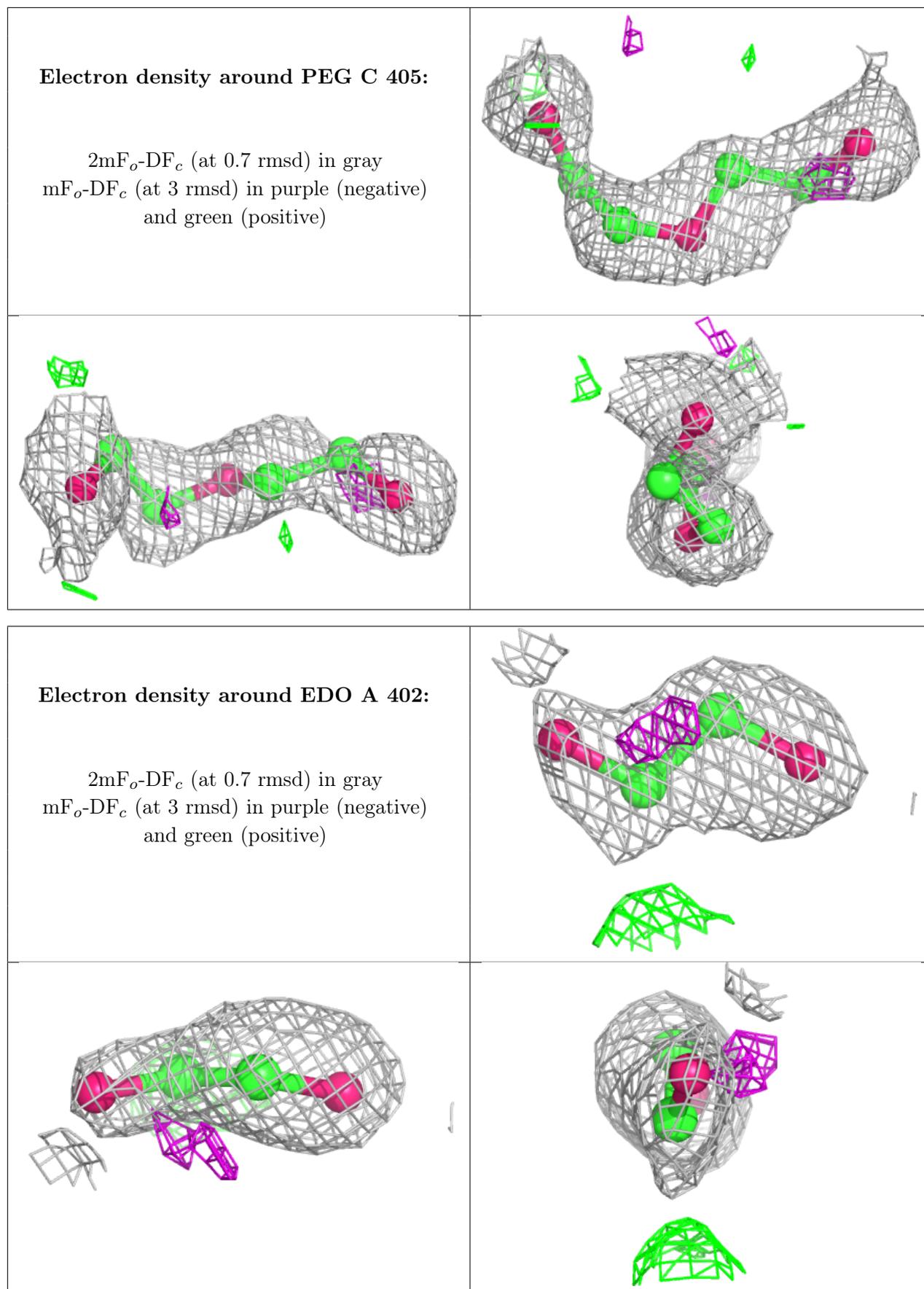


Electron density around MG D 104:

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

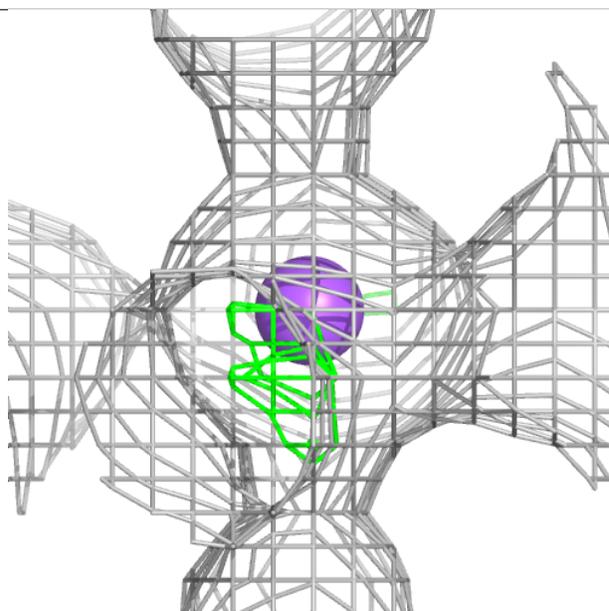
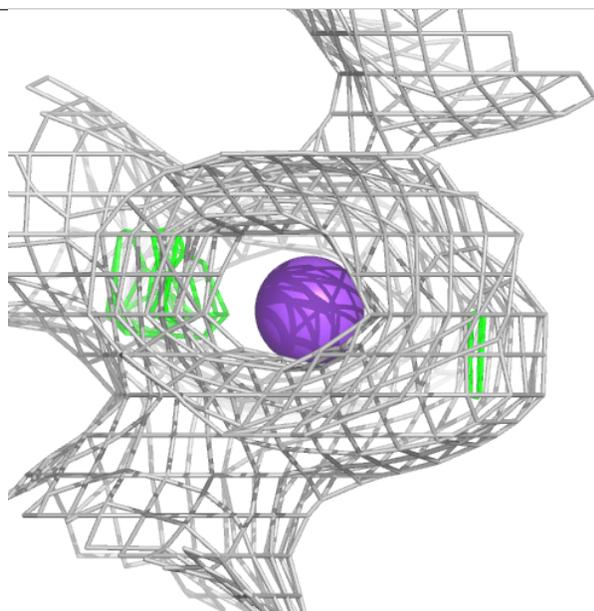
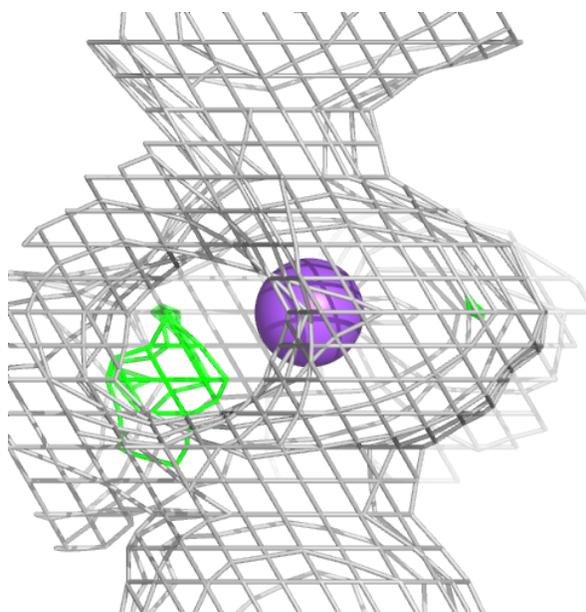


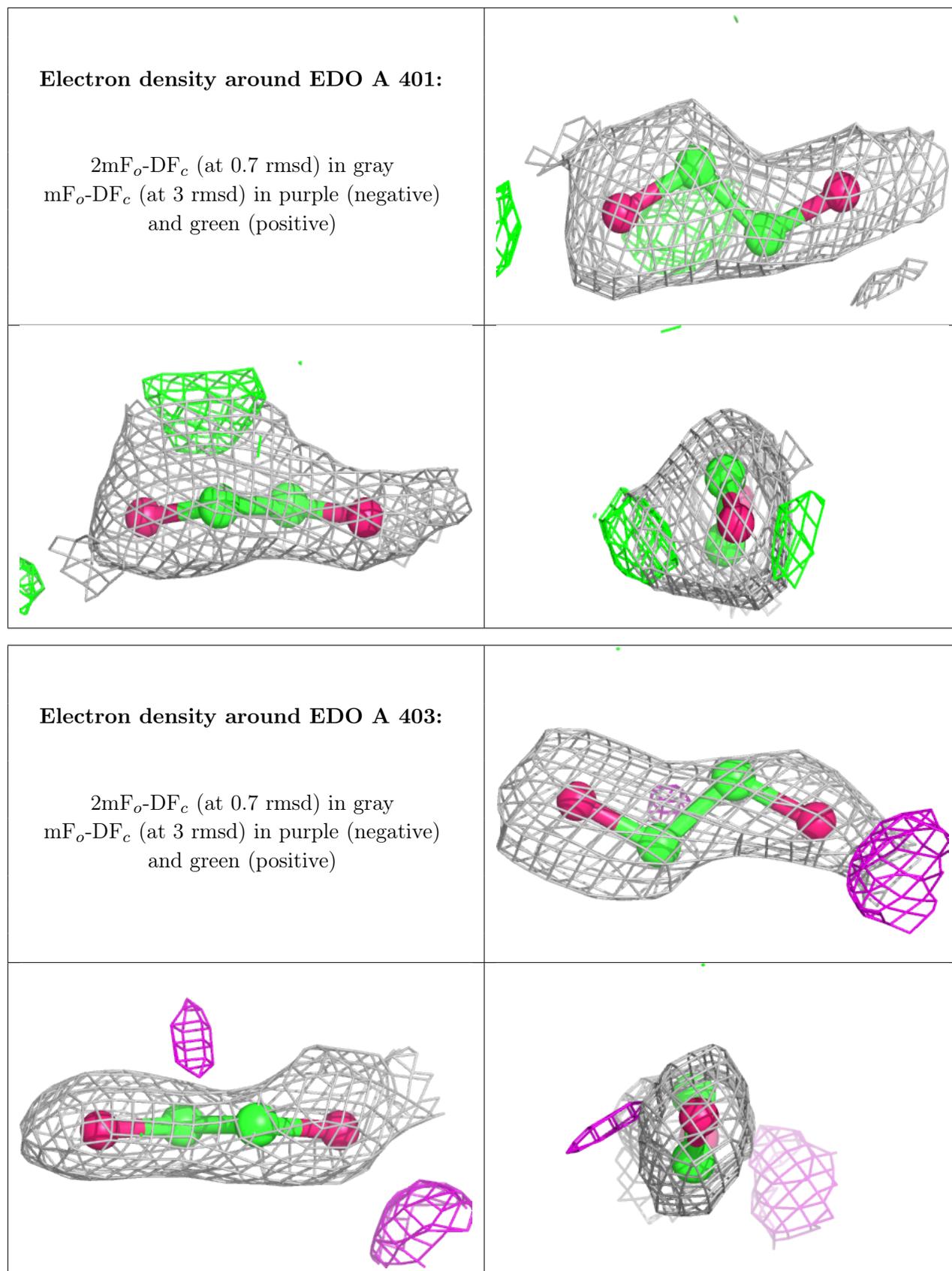


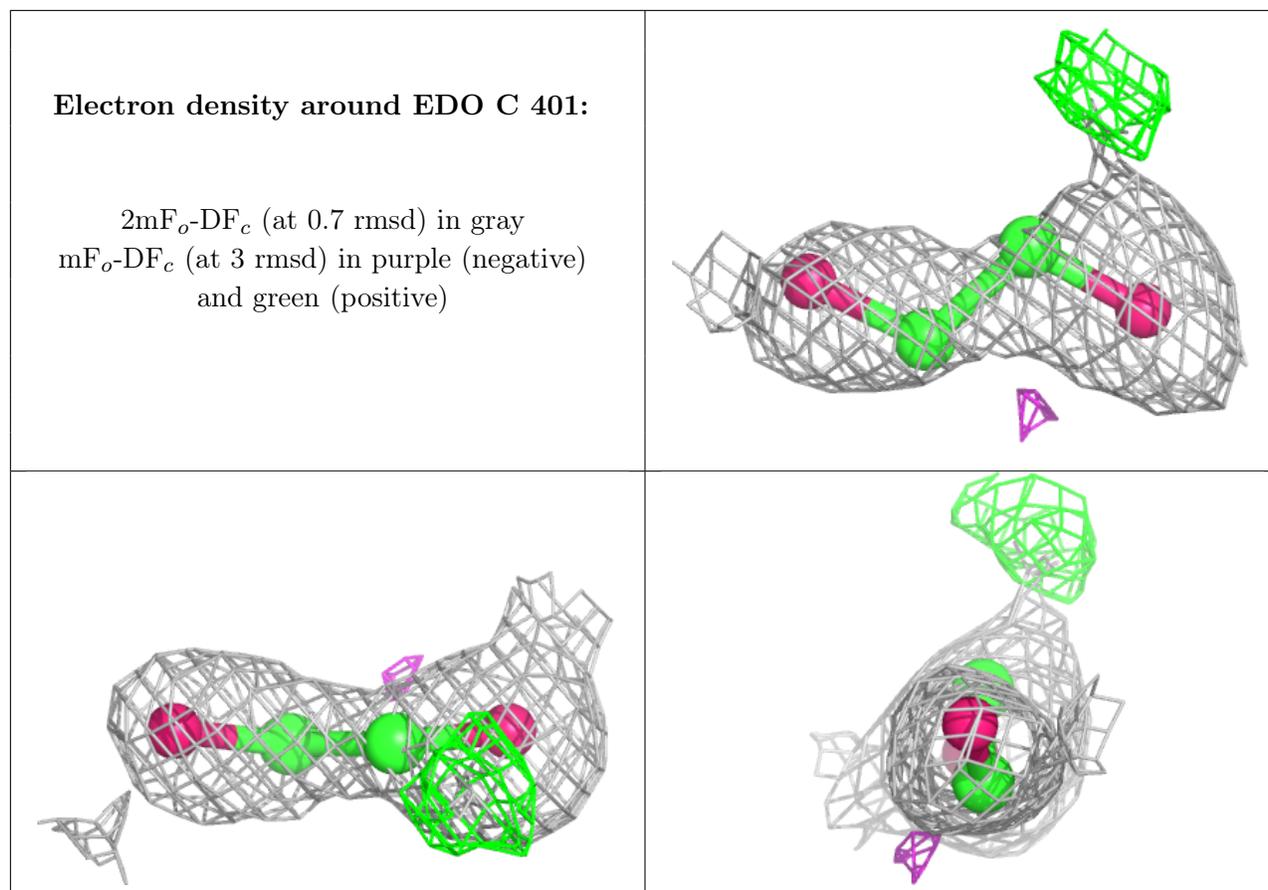


Electron density around NA F 103:

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

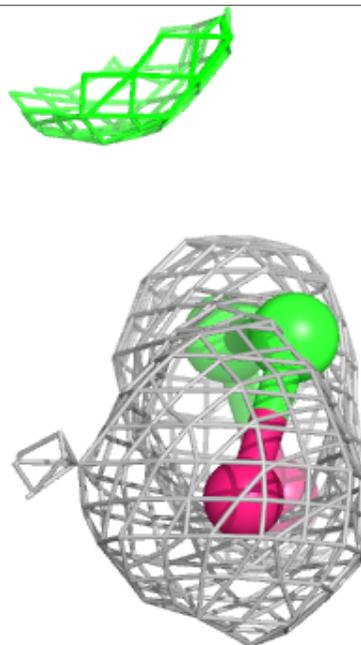
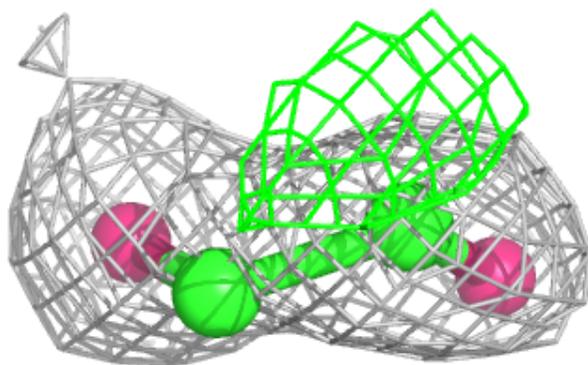
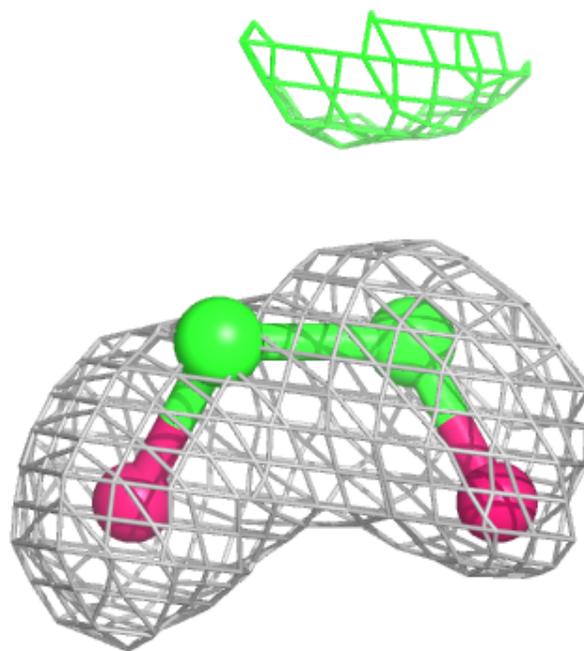


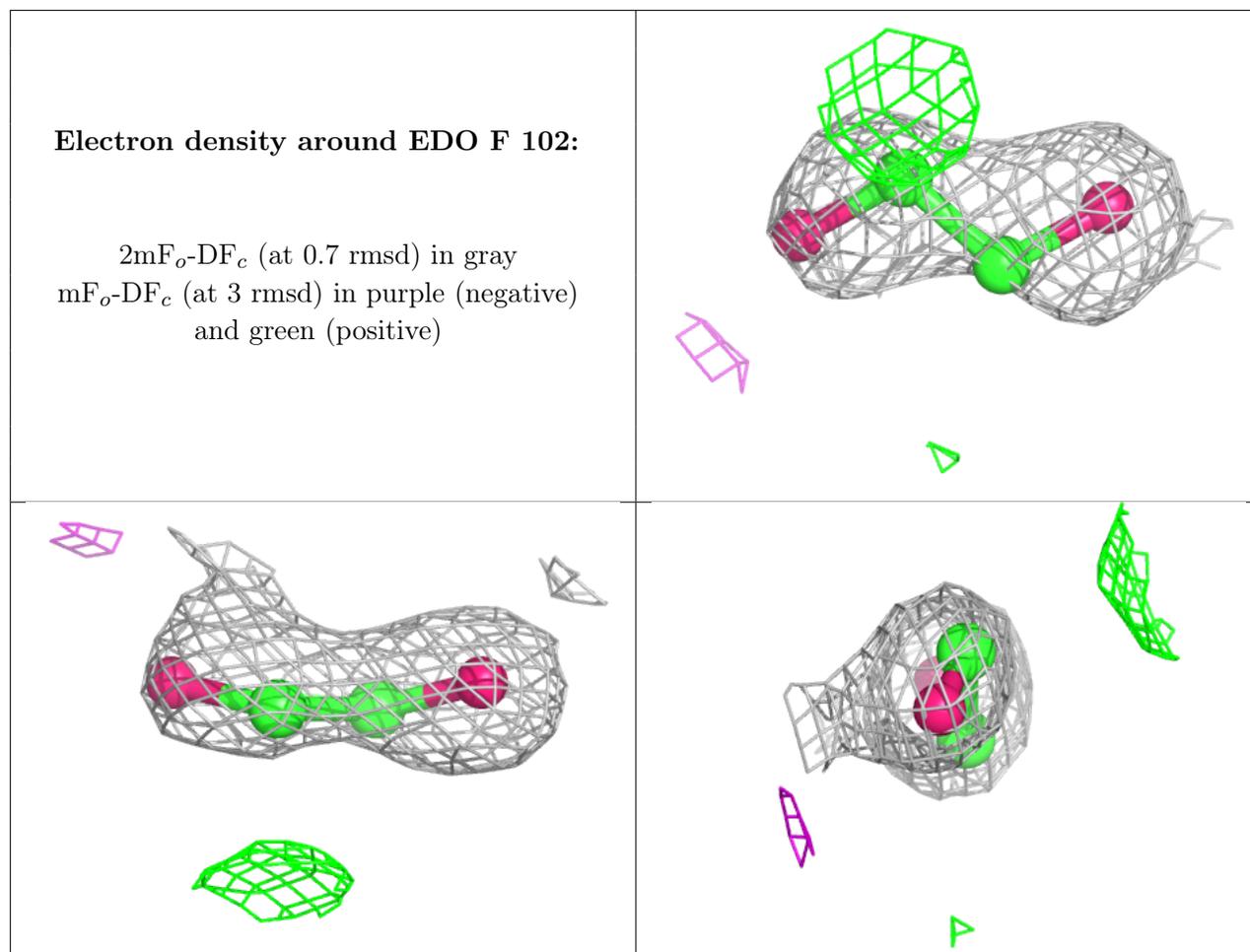


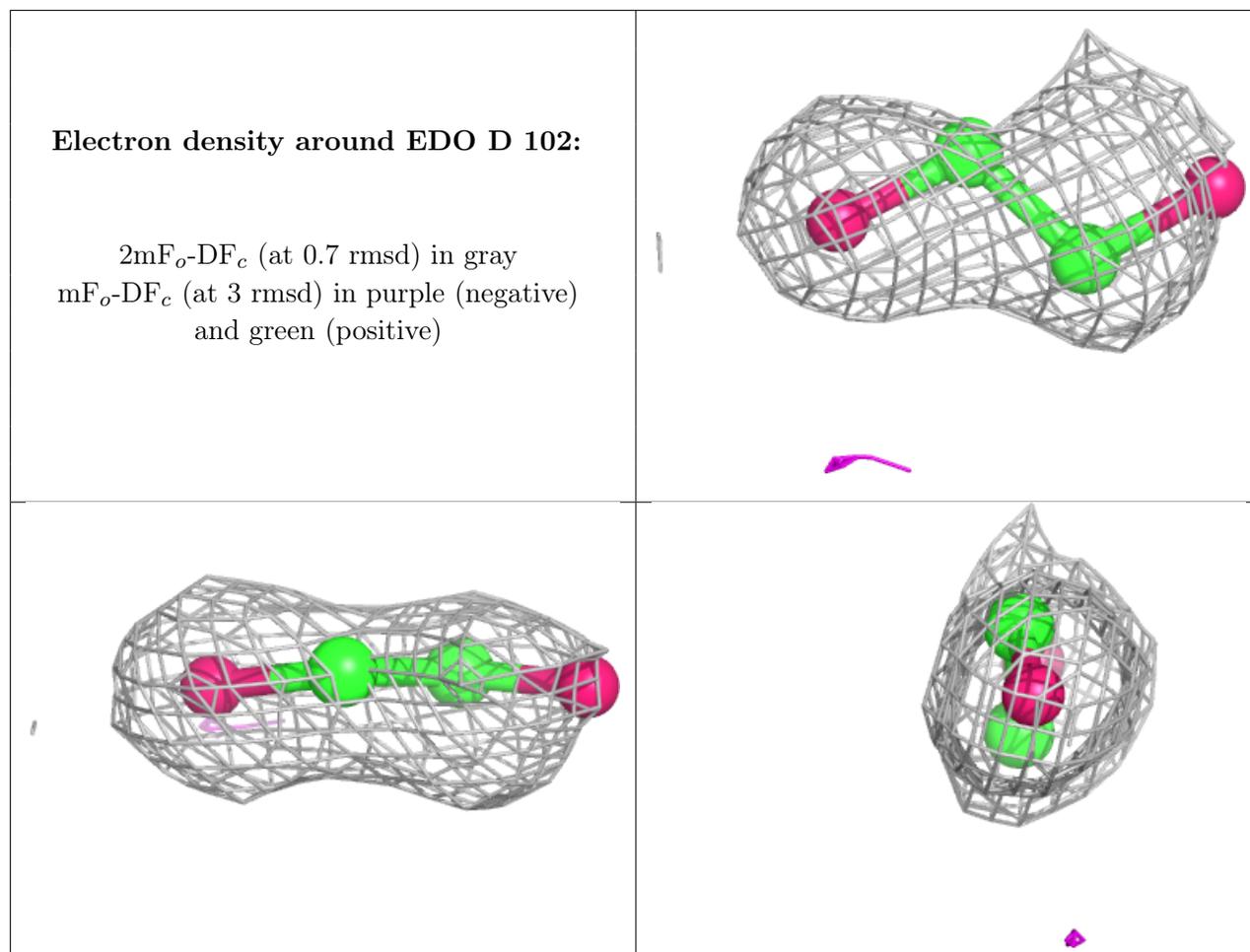


Electron density around EDO A 405:

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

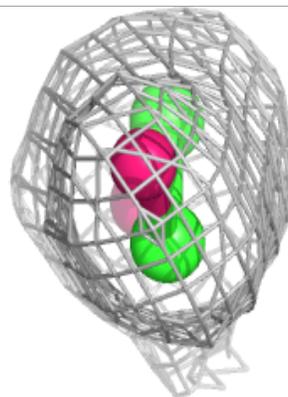
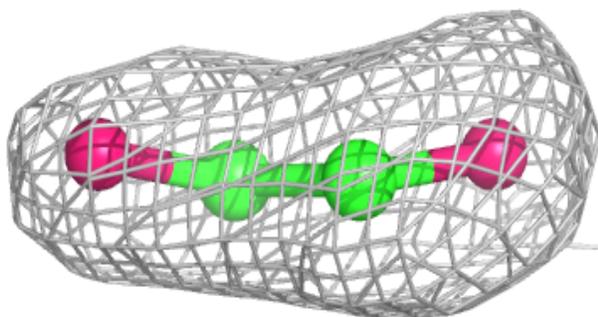
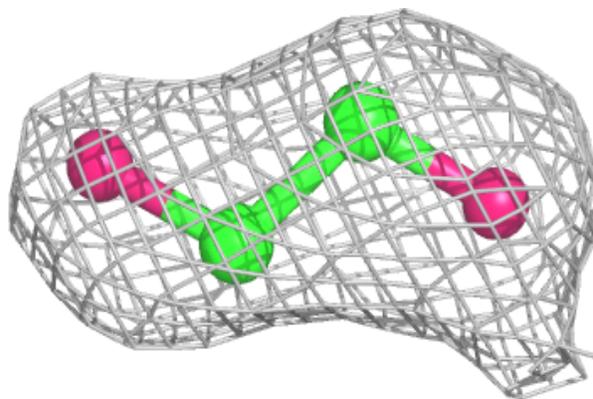






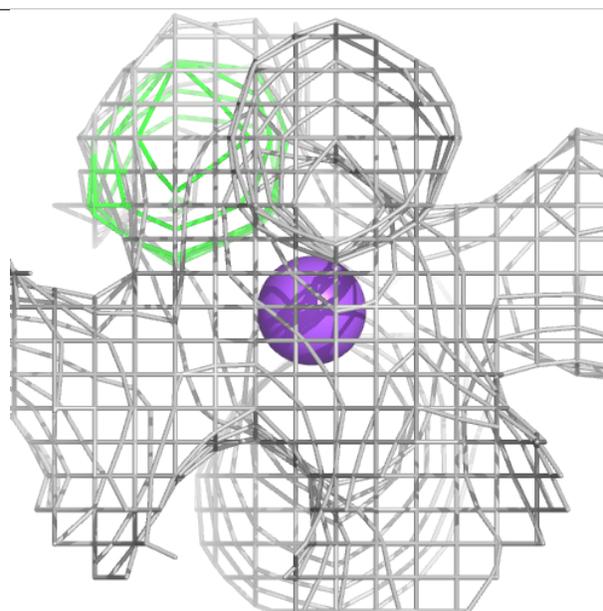
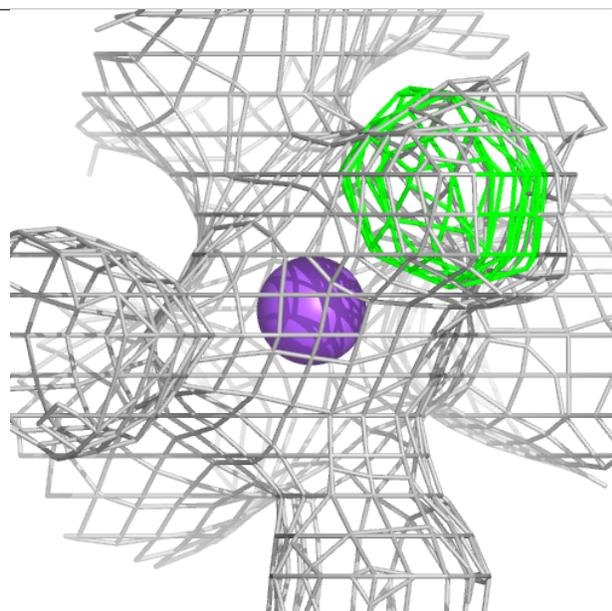
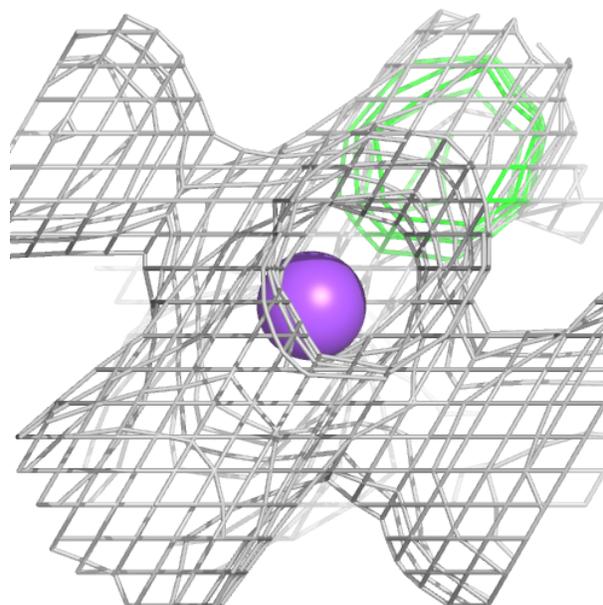
Electron density around EDO F 101:

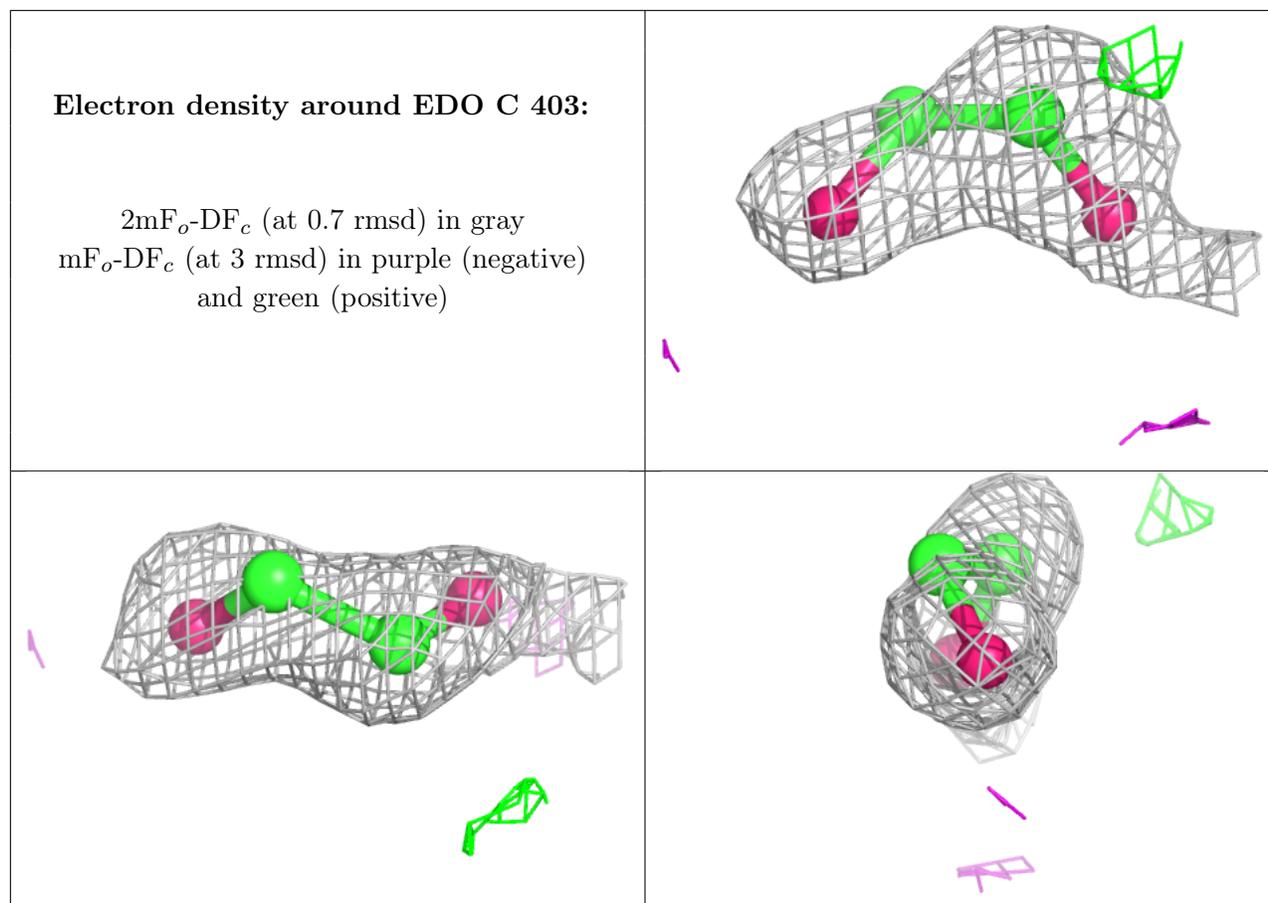
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around NA B 103:

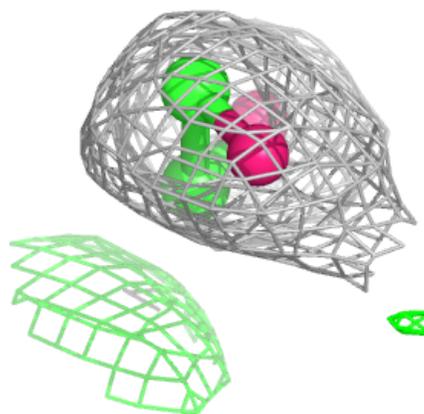
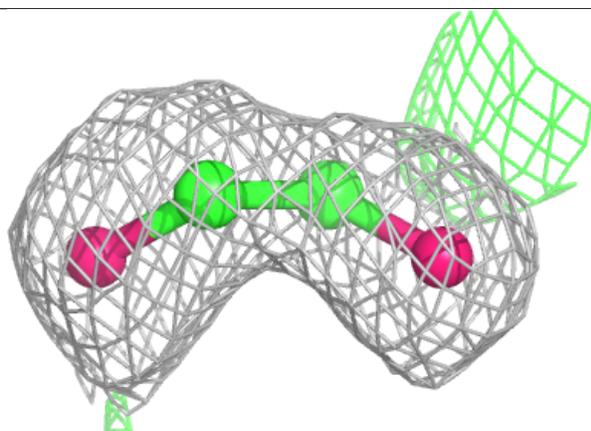
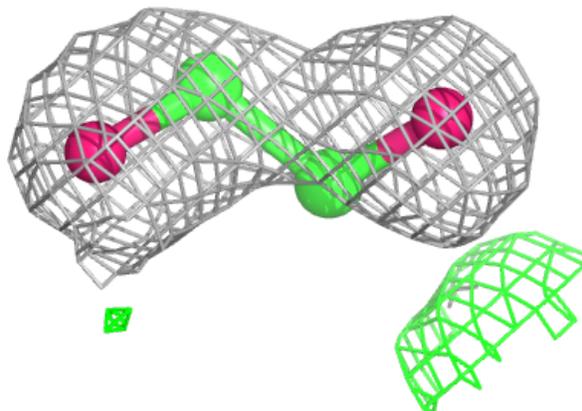
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

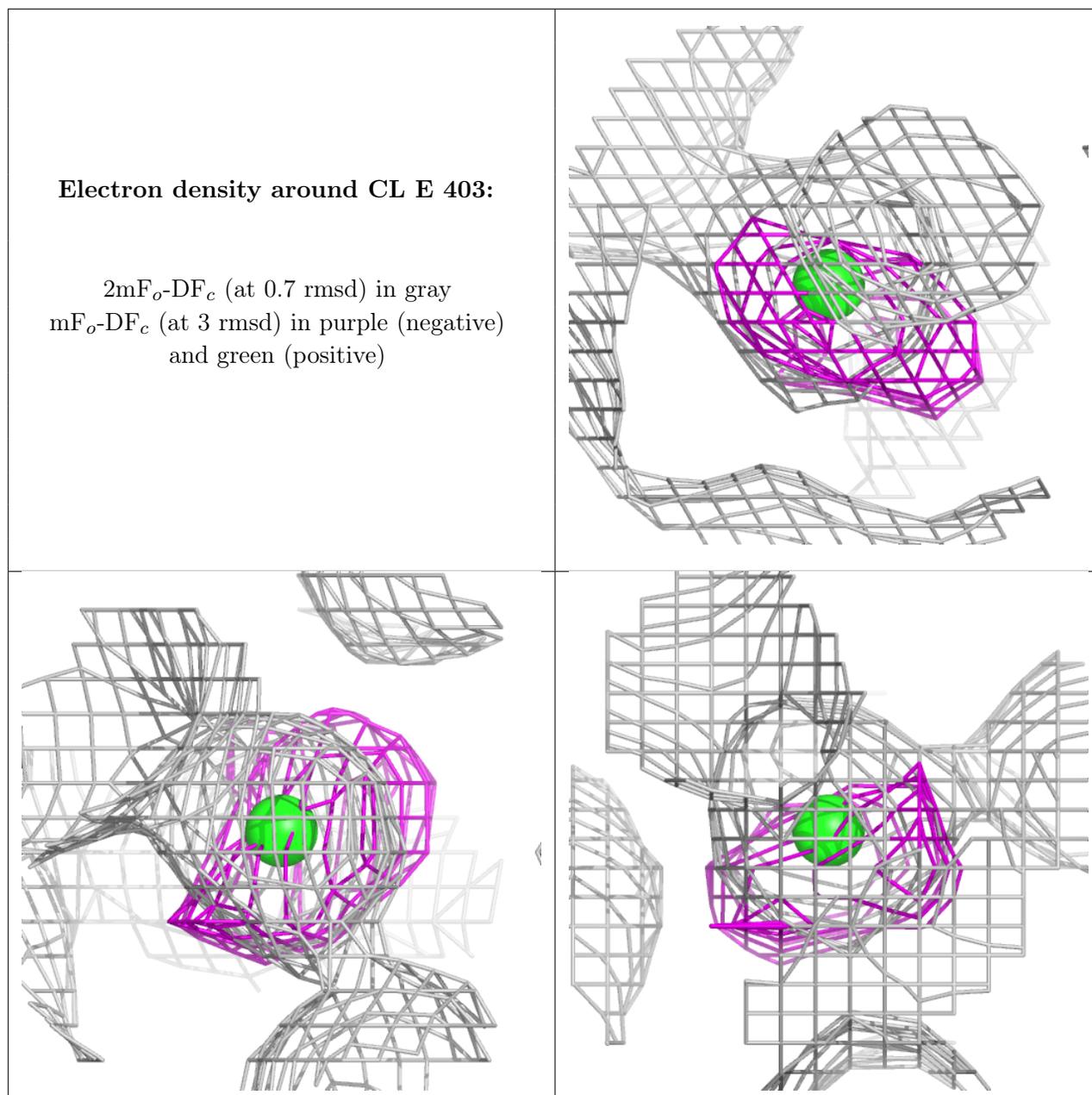


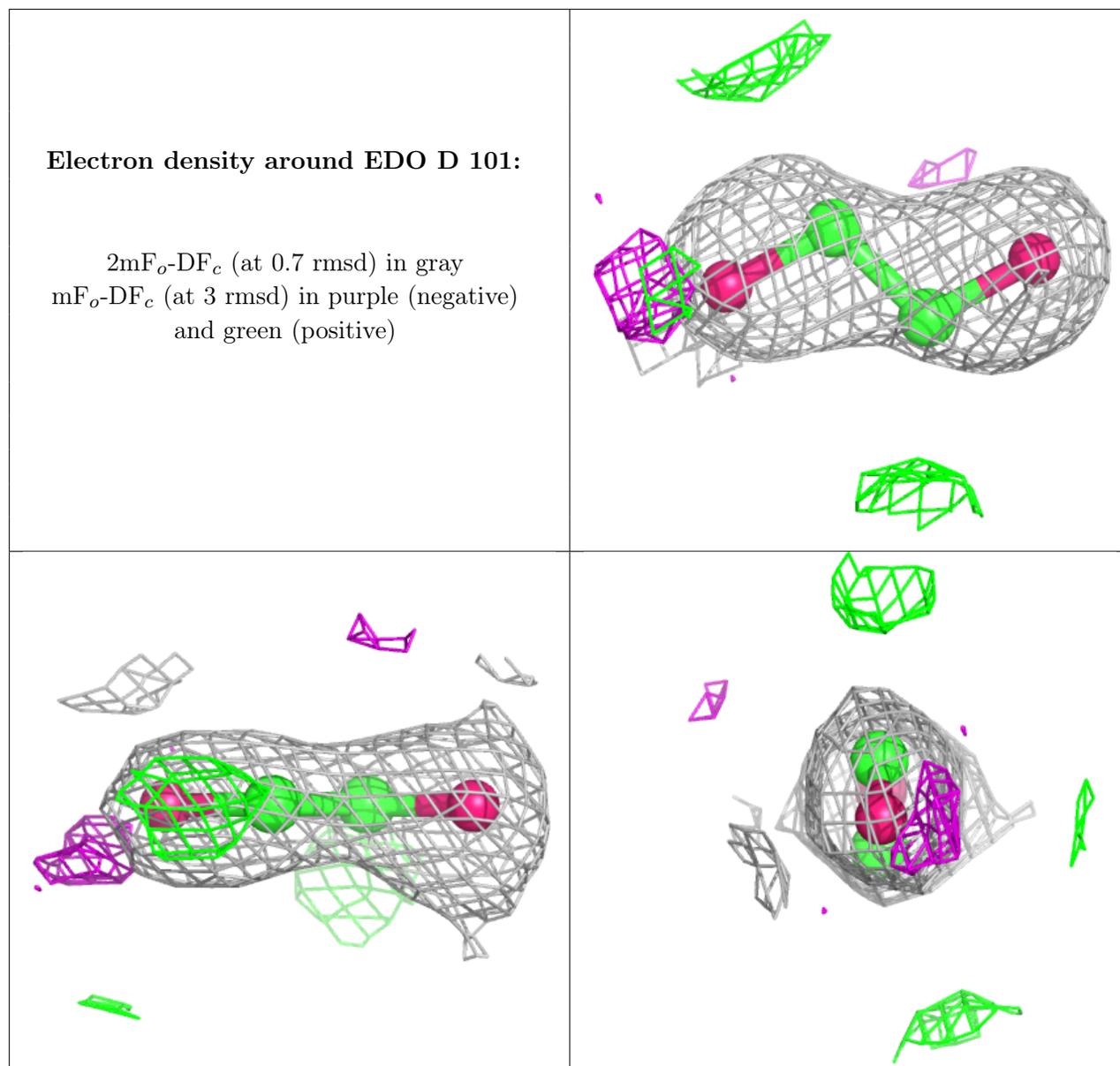


Electron density around EDO D 103:

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

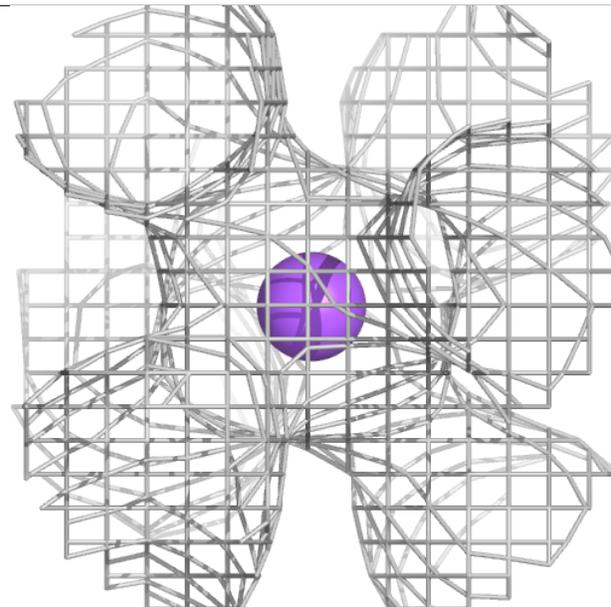
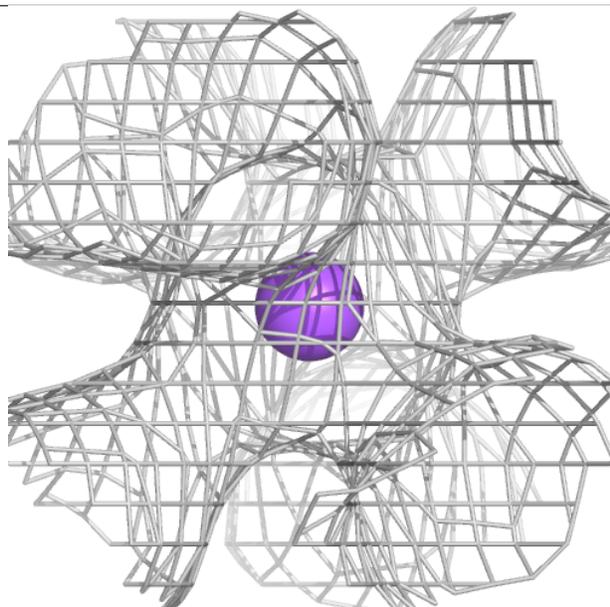
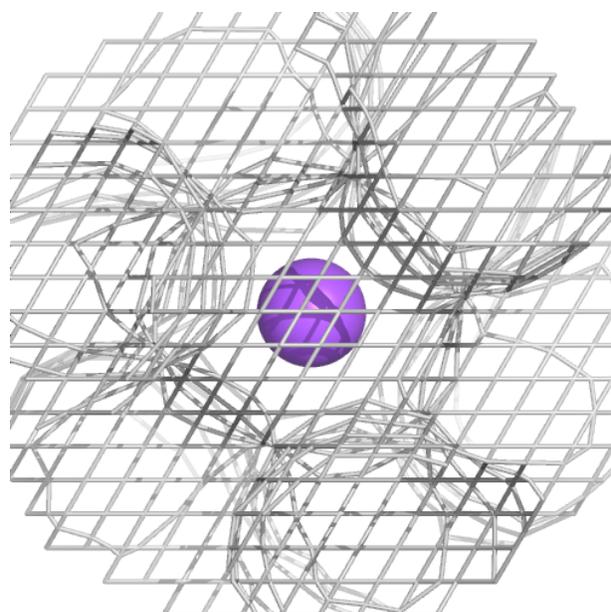


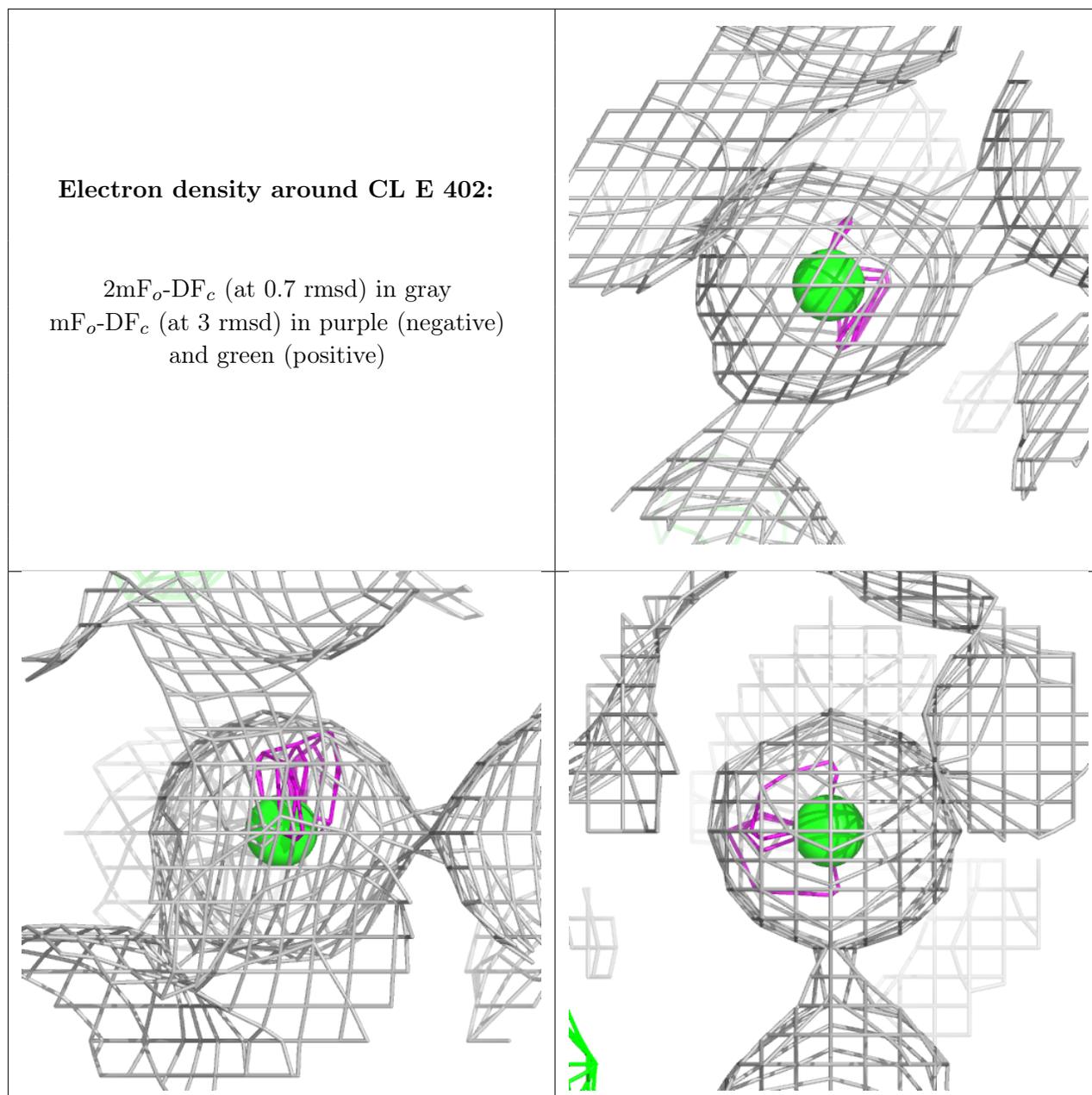


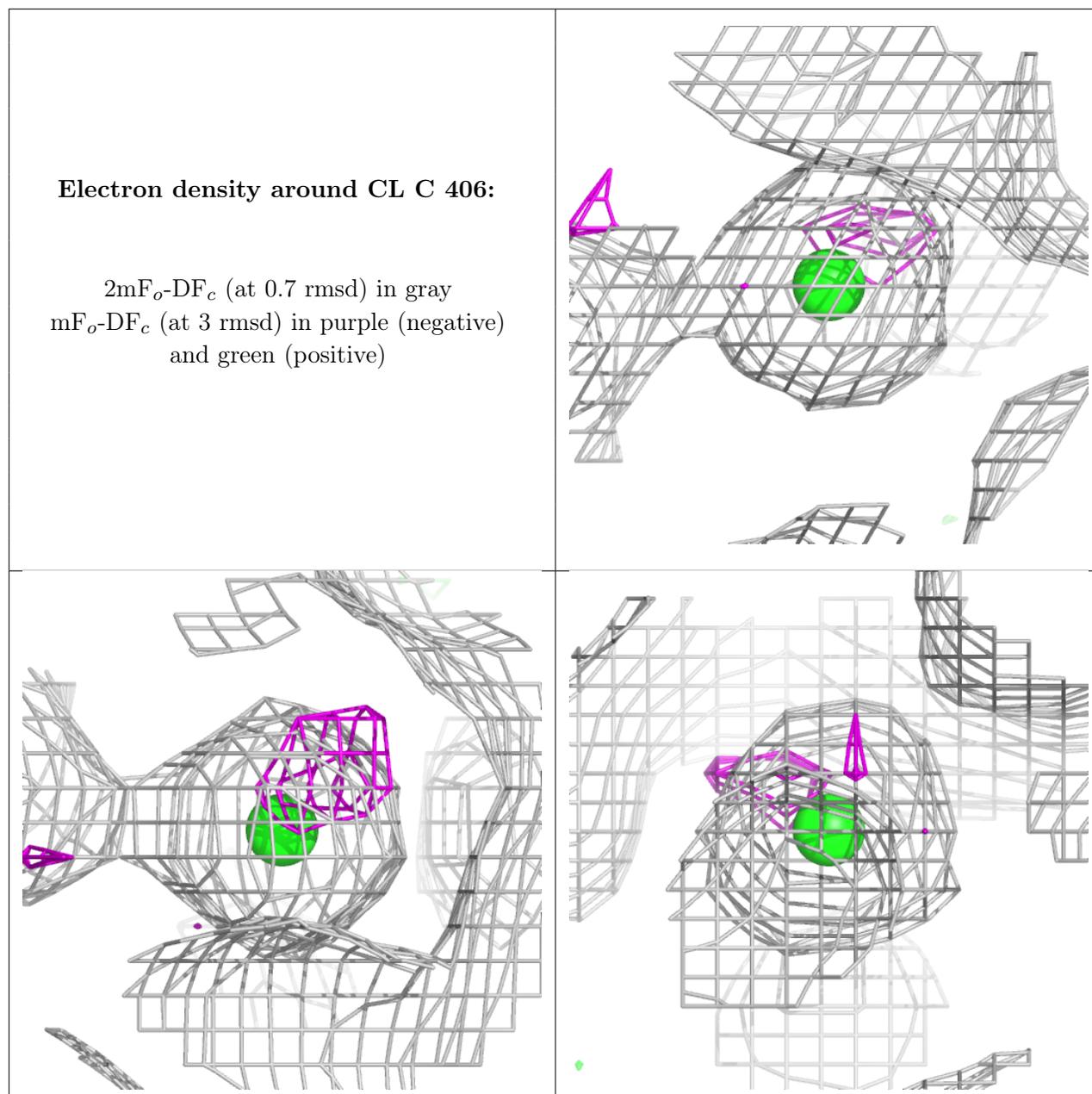


Electron density around NA E 404:

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







6.5 Other polymers [i](#)

There are no such residues in this entry.