



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 7, 2020 – 01:03 AM BST

PDB ID : 6Y1G
Title : Photoconverted HcRed in its optoacoustic state
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Deposited on : 2020-02-12
Resolution : 2.30 Å(reported)

This is a Full wwPDB X-ray Structure Validation 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.13.1
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)	:	2.13.1

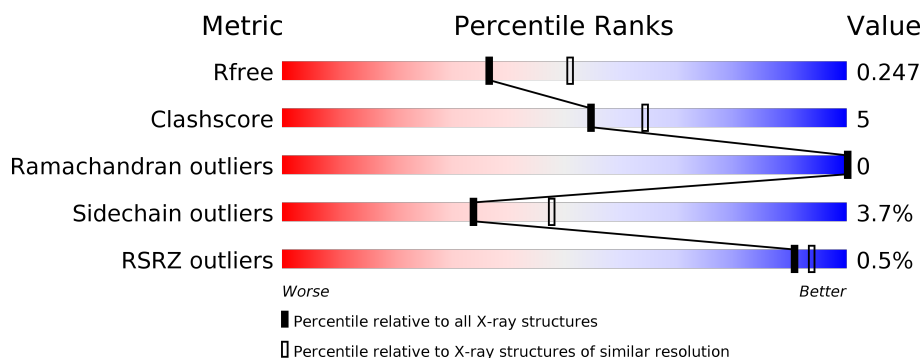
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION





The reported resolution of this entry is 2.30 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	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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 on 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	236	 83% 10% • 6%
1	B	236	 80% 13% • 6%
1	C	236	 78% 14% • 6%
1	D	236	 79% 15% 6%

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	EDO	D	304	-	-	-	X

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7839 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 GFP-like non-fluorescent chromoprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	222	Total	C	N	O	S	0	4	0
			1797	1143	304	335	15			
1	B	222	Total	C	N	O	S	0	4	0
			1795	1143	304	334	14			
1	C	222	Total	C	N	O	S	0	2	0
			1789	1136	305	333	15			
1	D	222	Total	C	N	O	S	0	2	0
			1786	1137	303	331	15			

There are 84 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-9	MET	-	initiating methionine	UNP Q95W85
A	-8	ARG	-	expression tag	UNP Q95W85
A	-7	GLY	-	expression tag	UNP Q95W85
A	-6	SER	-	expression tag	UNP Q95W85
A	-5	HIS	-	expression tag	UNP Q95W85
A	-4	HIS	-	expression tag	UNP Q95W85
A	-3	HIS	-	expression tag	UNP Q95W85
A	-2	HIS	-	expression tag	UNP Q95W85
A	-1	HIS	-	expression tag	UNP Q95W85
A	0	HIS	-	expression tag	UNP Q95W85
A	1	GLY	-	expression tag	UNP Q95W85
A	2	SER	-	expression tag	UNP Q95W85
A	3	SER	ALA	conflict	UNP Q95W85
A	37	ALA	THR	conflict	UNP Q95W85
A	64	CRU	GLU	chromophore	UNP Q95W85
A	64	CRU	TYR	chromophore	UNP Q95W85
A	64	CRU	GLY	chromophore	UNP Q95W85
A	144	SER	CYS	conflict	UNP Q95W85
A	174	HIS	LEU	conflict	UNP Q95W85
A	202	LEU	PRO	conflict	UNP Q95W85
A	205	GLU	LYS	conflict	UNP Q95W85

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-9	MET	-	initiating methionine	UNP Q95W85
B	-8	ARG	-	expression tag	UNP Q95W85
B	-7	GLY	-	expression tag	UNP Q95W85
B	-6	SER	-	expression tag	UNP Q95W85
B	-5	HIS	-	expression tag	UNP Q95W85
B	-4	HIS	-	expression tag	UNP Q95W85
B	-3	HIS	-	expression tag	UNP Q95W85
B	-2	HIS	-	expression tag	UNP Q95W85
B	-1	HIS	-	expression tag	UNP Q95W85
B	0	HIS	-	expression tag	UNP Q95W85
B	1	GLY	-	expression tag	UNP Q95W85
B	2	SER	-	expression tag	UNP Q95W85
B	3	SER	ALA	conflict	UNP Q95W85
B	37	ALA	THR	conflict	UNP Q95W85
B	64	CRU	GLU	chromophore	UNP Q95W85
B	64	CRU	TYR	chromophore	UNP Q95W85
B	64	CRU	GLY	chromophore	UNP Q95W85
B	144	SER	CYS	conflict	UNP Q95W85
B	174	HIS	LEU	conflict	UNP Q95W85
B	202	LEU	PRO	conflict	UNP Q95W85
B	205	GLU	LYS	conflict	UNP Q95W85
C	-9	MET	-	initiating methionine	UNP Q95W85
C	-8	ARG	-	expression tag	UNP Q95W85
C	-7	GLY	-	expression tag	UNP Q95W85
C	-6	SER	-	expression tag	UNP Q95W85
C	-5	HIS	-	expression tag	UNP Q95W85
C	-4	HIS	-	expression tag	UNP Q95W85
C	-3	HIS	-	expression tag	UNP Q95W85
C	-2	HIS	-	expression tag	UNP Q95W85
C	-1	HIS	-	expression tag	UNP Q95W85
C	0	HIS	-	expression tag	UNP Q95W85
C	1	GLY	-	expression tag	UNP Q95W85
C	2	SER	-	expression tag	UNP Q95W85
C	3	SER	ALA	conflict	UNP Q95W85
C	37	ALA	THR	conflict	UNP Q95W85
C	64	CRU	GLU	chromophore	UNP Q95W85
C	64	CRU	TYR	chromophore	UNP Q95W85
C	64	CRU	GLY	chromophore	UNP Q95W85
C	144	SER	CYS	conflict	UNP Q95W85
C	174	HIS	LEU	conflict	UNP Q95W85
C	202	LEU	PRO	conflict	UNP Q95W85
C	205	GLU	LYS	conflict	UNP Q95W85

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-9	MET	-	initiating methionine	UNP Q95W85
D	-8	ARG	-	expression tag	UNP Q95W85
D	-7	GLY	-	expression tag	UNP Q95W85
D	-6	SER	-	expression tag	UNP Q95W85
D	-5	HIS	-	expression tag	UNP Q95W85
D	-4	HIS	-	expression tag	UNP Q95W85
D	-3	HIS	-	expression tag	UNP Q95W85
D	-2	HIS	-	expression tag	UNP Q95W85
D	-1	HIS	-	expression tag	UNP Q95W85
D	0	HIS	-	expression tag	UNP Q95W85
D	1	GLY	-	expression tag	UNP Q95W85
D	2	SER	-	expression tag	UNP Q95W85
D	3	SER	ALA	conflict	UNP Q95W85
D	37	ALA	THR	conflict	UNP Q95W85
D	64	CRU	GLU	chromophore	UNP Q95W85
D	64	CRU	TYR	chromophore	UNP Q95W85
D	64	CRU	GLY	chromophore	UNP Q95W85
D	144	SER	CYS	conflict	UNP Q95W85
D	174	HIS	LEU	conflict	UNP Q95W85
D	202	LEU	PRO	conflict	UNP Q95W85
D	205	GLU	LYS	conflict	UNP Q95W85

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



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	B	1	Total C O 4 2 2	0	0
2	B	1	Total C O 4 2 2	0	0
2	B	1	Total C O 4 2 2	0	0
2	B	1	Total C O 4 2 2	0	0
2	B	1	Total C O 4 2 2	0	0
2	B	1	Total C O 4 2 2	0	0
2	C	1	Total C O 4 2 2	0	0
2	C	1	Total C O 4 2 2	0	0
2	C	1	Total C O 4 2 2	0	0
2	C	1	Total C O 4 2 2	0	0
2	D	1	Total C O 4 2 2	0	0
2	D	1	Total C O 4 2 2	0	0
2	D	1	Total C O 4 2 2	0	0
2	D	1	Total C O 4 2 2	0	0
2	D	1	Total C O 4 2 2	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	161	Total O 161 161	0	0

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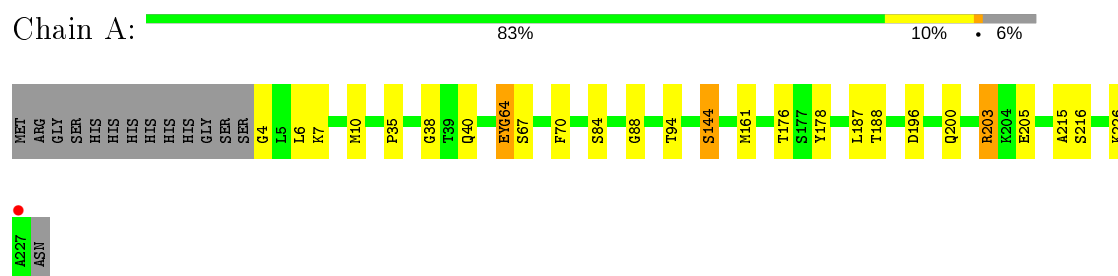
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	157	Total 157	O 157	0	0
3	C	152	Total 152	O 152	0	0
3	D	126	Total 126	O 126	0	0

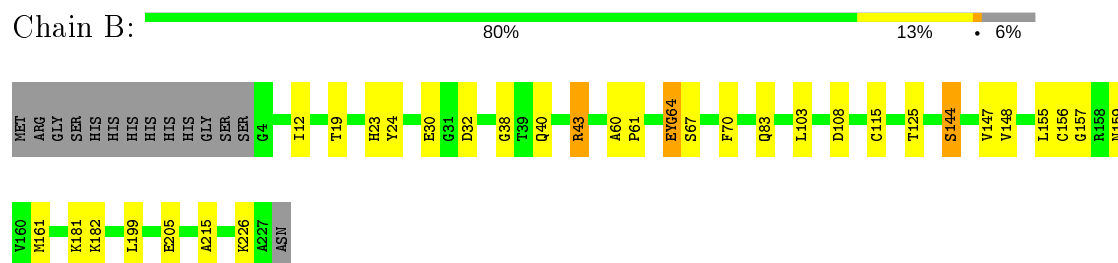
3 Residue-property plots [i](#)

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.

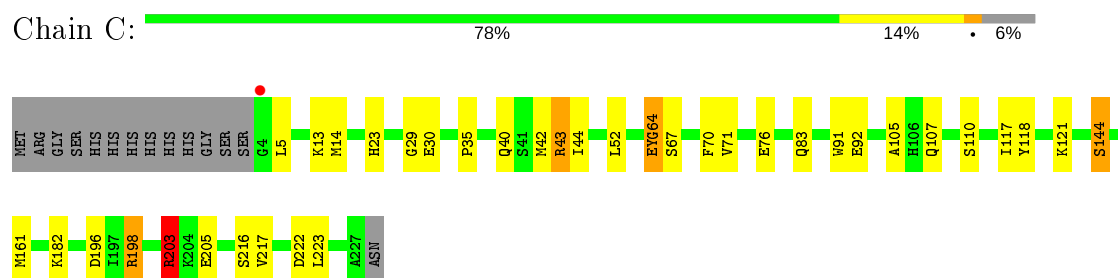
- Molecule 1: GFP-like non-fluorescent chromoprotein



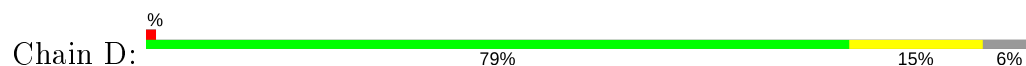
- Molecule 1: GFP-like non-fluorescent chromoprotein

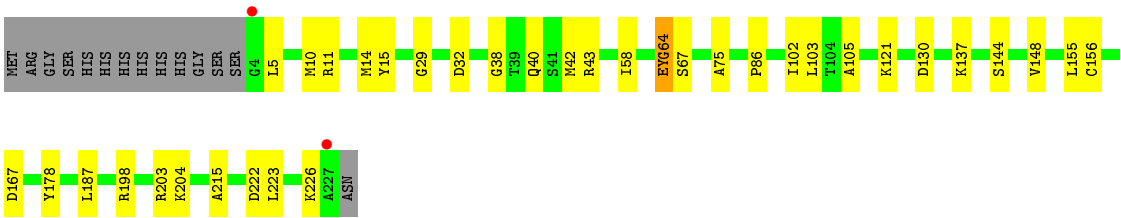


- Molecule 1: GFP-like non-fluorescent chromoprotein



- Molecule 1: GFP-like non-fluorescent chromoprotein





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	50.71 Å 75.20 Å 254.21 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.78 – 2.30 49.73 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.7 (49.78-2.30) 99.7 (49.73-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.02 (at 2.29 Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.168 , 0.247 0.174 , 0.247	Depositor DCC
R_{free} test set	2181 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	30.0	Xtriage
Anisotropy	0.362	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 42.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7839	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.33% of the height of the origin peak. No significant pseudotranslation is detected.*

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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.78	0/1827	0.98	1/2465 (0.0%)
1	B	0.80	0/1825	0.98	1/2464 (0.0%)
1	C	0.79	0/1810	0.97	1/2442 (0.0%)
1	D	0.75	0/1813	0.98	1/2446 (0.0%)
All	All	0.78	0/7275	0.98	4/9817 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	43	ARG	NE-CZ-NH1	5.95	123.27	120.30
1	D	167	ASP	CB-CA-C	5.54	121.47	110.40
1	C	203	ARG	NE-CZ-NH1	5.16	122.88	120.30
1	A	188	THR	CA-CB-OG1	-5.10	98.30	109.00

There are no chirality outliers.

There are no planarity outliers.

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	1797	0	1737	14	0
1	B	1795	0	1739	19	0
1	C	1789	0	1720	25	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	1786	0	1729	21	0
2	A	20	0	30	1	0
2	B	20	0	30	0	0
2	C	16	0	24	0	0
2	D	20	0	30	0	0
3	A	161	0	0	2	0
3	B	157	0	0	2	0
3	C	152	0	0	2	0
3	D	126	0	0	3	0
All	All	7839	0	7039	74	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 5.

All (74) 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:144:SER:OG	1:B:161:MET:HG2	1.80	0.82
1:D:14:MET:SD	1:D:42[A]:MET:HE1	2.32	0.70
1:C:203:ARG:NH2	1:C:205:GLU:OE1	2.25	0.69
1:A:40:GLN:HE22	1:A:67:SER:CB	2.06	0.68
1:A:144:SER:OG	1:A:161:MET:HG2	1.98	0.63
1:D:58:ILE:HD12	1:D:103:LEU:HD12	1.81	0.63
1:C:83:GLN:O	1:C:182:LYS:HE3	1.99	0.63
1:C:198:ARG:NH2	3:C:401:HOH:O	2.31	0.61
1:B:108:ASP:OD1	1:B:181:LYS:HE2	2.01	0.61
1:A:203:ARG:NH2	1:A:205:GLU:OE2	2.35	0.60
1:C:67:SER:HG	1:C:118:TYR:HH	1.47	0.59
1:C:144:SER:OG	1:C:161:MET:HG2	2.02	0.59
1:C:40:GLN:HE22	1:C:67:SER:CB	2.16	0.58
1:A:4:GLY:HA2	3:A:517:HOH:O	2.06	0.56
1:D:40:GLN:HE22	1:D:67:SER:HB3	1.70	0.55
1:C:14:MET:SD	1:C:42:MET:HE1	2.47	0.55
1:D:105:ALA:HA	1:D:121:LYS:O	2.07	0.55
1:C:64:CRU:HD2	1:C:64:CRU:O2	2.08	0.54
1:A:38:GLY:O	1:A:215:ALA:HA	2.09	0.52
1:C:110[B]:SER:OG	1:C:117:ILE:HB	2.10	0.52
1:A:196:ASP:O	1:A:216:SER:HA	2.10	0.51
1:C:42:MET:HE3	1:C:44:ILE:HG13	1.92	0.51
1:D:64:CRU:O2	1:D:64:CRU:HD2	2.11	0.50
1:B:148[A]:VAL:HG12	1:B:155:LEU:HD11	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:125[B]:THR:HG22	1:C:92:GLU:HG3	1.95	0.49
1:D:155:LEU:HB3	1:D:178:TYR:HB2	1.94	0.48
1:D:10:MET:O	1:D:32:ASP:HA	2.13	0.48
1:B:147:VAL:O	1:B:157:GLY:HA2	2.14	0.48
1:B:64:CRU:HE1	1:B:159:ASN:ND2	2.29	0.48
1:A:40:GLN:NE2	3:A:403:HOH:O	2.43	0.47
1:C:83:GLN:O	1:C:182:LYS:CE	2.62	0.47
1:D:148:VAL:HA	1:D:156:CYS:O	2.14	0.47
1:D:38:GLY:O	1:D:215:ALA:HA	2.14	0.47
1:C:217:VAL:HG23	1:D:223:LEU:HD21	1.97	0.46
1:A:94:THR:HG21	1:D:102:ILE:HG21	1.97	0.45
1:A:176:THR:HG21	1:A:178:TYR:CZ	2.51	0.45
1:D:130:ASP:HB2	3:D:464:HOH:O	2.17	0.45
1:A:84[A]:SER:O	1:A:88:GLY:N	2.50	0.44
1:B:43:ARG:HG3	3:B:530:HOH:O	2.17	0.44
1:B:64:CRU:O2	1:B:64:CRU:HD2	2.17	0.44
1:D:40:GLN:NE2	1:D:67:SER:HB3	2.32	0.44
1:B:83:GLN:O	1:B:182:LYS:HE3	2.18	0.44
1:B:40:GLN:HE22	1:B:67:SER:HB3	1.82	0.44
1:D:137:LYS:HE3	3:D:512:HOH:O	2.16	0.44
1:C:35:PRO:HA	1:C:70:PHE:HA	2.00	0.44
1:B:12:ILE:HD11	1:B:70:PHE:CZ	2.53	0.44
1:B:148[B]:VAL:HA	1:B:156:CYS:O	2.18	0.44
1:B:182:LYS:NZ	3:B:406:HOH:O	2.47	0.43
1:D:5:LEU:HD21	1:D:86:PRO:HB3	2.00	0.43
1:D:29:GLY:HA2	1:D:43:ARG:O	2.18	0.43
1:A:6:LEU:HD22	1:A:10[B]:MET:SD	2.59	0.43
1:A:64:CRU:O2	1:A:64:CRU:HD2	2.18	0.43
1:A:35:PRO:HA	1:A:70:PHE:HA	2.01	0.43
1:C:29:GLY:HA2	1:C:43:ARG:O	2.19	0.43
1:C:198:ARG:NH1	1:D:223:LEU:O	2.52	0.42
1:B:19[B]:THR:HA	1:B:23:HIS:O	2.19	0.42
1:C:105:ALA:HA	1:C:121:LYS:O	2.19	0.42
1:C:42:MET:HE2	1:C:44:ILE:HD11	2.01	0.42
1:B:19[A]:THR:HG22	1:B:24:TYR:HA	2.01	0.42
1:B:64:CRU:CE1	1:B:159:ASN:HD22	2.33	0.42
1:D:5:LEU:HD12	1:D:5:LEU:HA	1.91	0.41
1:A:200:GLN:HG2	2:A:302:EDO:H12	2.02	0.41
1:B:60:ALA:HB3	1:B:61:PRO:HD3	2.02	0.41
1:C:196:ASP:O	1:C:216:SER:HA	2.20	0.41
1:D:75:ALA:O	3:D:401:HOH:O	2.22	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:38:GLY:O	1:B:215:ALA:HA	2.20	0.41
1:C:71:VAL:HG22	1:C:216:SER:HB2	2.04	0.40
1:C:5:LEU:HA	1:C:5:LEU:HD12	1.84	0.40
1:D:14:MET:HG2	1:D:15:TYR:N	2.36	0.40
1:C:52:LEU:HD23	1:C:52:LEU:HA	1.95	0.40
1:C:223:LEU:O	1:D:198:ARG:NH1	2.54	0.40
1:B:12:ILE:HD11	1:B:70:PHE:CE1	2.57	0.40
1:C:23:HIS:HB3	3:C:506:HOH:O	2.21	0.40
1:C:91:TRP:CE2	1:C:107:GLN:CB	3.04	0.40

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	223/236 (94%)	217 (97%)	6 (3%)	0	100	100
1	B	223/236 (94%)	215 (96%)	8 (4%)	0	100	100
1	C	221/236 (94%)	214 (97%)	7 (3%)	0	100	100
1	D	221/236 (94%)	213 (96%)	8 (4%)	0	100	100
All	All	888/944 (94%)	859 (97%)	29 (3%)	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	192/200 (96%)	187 (97%)	5 (3%)	46	63
1	B	192/200 (96%)	184 (96%)	8 (4%)	30	42
1	C	190/200 (95%)	182 (96%)	8 (4%)	30	42
1	D	190/200 (95%)	183 (96%)	7 (4%)	34	48
All	All	764/800 (96%)	736 (96%)	28 (4%)	34	48

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	LYS
1	A	144	SER
1	A	187	LEU
1	A	203	ARG
1	A	226	LYS
1	B	30	GLU
1	B	32	ASP
1	B	103	LEU
1	B	115	CYS
1	B	144	SER
1	B	199	LEU
1	B	205	GLU
1	B	226	LYS
1	C	13	LYS
1	C	30	GLU
1	C	43	ARG
1	C	76	GLU
1	C	144	SER
1	C	198	ARG
1	C	203	ARG
1	C	222	ASP
1	D	11	ARG
1	D	144	SER
1	D	187	LEU
1	D	203	ARG
1	D	204	LYS
1	D	222	ASP
1	D	226	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	126	ASN
1	A	174	HIS
1	B	174	HIS
1	C	174	HIS
1	D	126	ASN
1	D	174	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

4 non-standard protein/DNA/RNA residues 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$
1	CRU	A	64	1	21,25,26	3.34	5 (23%)	24,34,36	4.53	11 (45%)
1	CRU	B	64	1	21,25,26	3.50	3 (14%)	24,34,36	4.95	10 (41%)
1	CRU	C	64	1	21,25,26	3.64	6 (28%)	24,34,36	4.58	13 (54%)
1	CRU	D	64	1	21,25,26	3.84	6 (28%)	24,34,36	5.11	12 (50%)

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
1	CRU	A	64	1	-	3/8/32/33	0/2/2/2
1	CRU	B	64	1	-	3/8/32/33	0/2/2/2
1	CRU	C	64	1	-	3/8/32/33	0/2/2/2
1	CRU	D	64	1	-	5/8/32/33	0/2/2/2

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	64	CRU	CB2-CA2	16.09	1.48	1.35
1	C	64	CRU	CB2-CA2	15.60	1.48	1.35
1	B	64	CRU	CB2-CA2	14.89	1.47	1.35
1	A	64	CRU	CB2-CA2	13.64	1.46	1.35
1	D	64	CRU	C2-N3	-3.62	1.31	1.39
1	A	64	CRU	O2-C2	3.42	1.30	1.23
1	B	64	CRU	C2-N3	-3.19	1.32	1.39
1	D	64	CRU	O2-C2	3.14	1.29	1.23
1	A	64	CRU	C2-N3	-2.95	1.32	1.39
1	D	64	CRU	CA2-N2	-2.82	1.32	1.38
1	C	64	CRU	CA2-C2	-2.74	1.45	1.48
1	D	64	CRU	CE1-CZ	2.68	1.44	1.38
1	C	64	CRU	CE1-CZ	2.62	1.43	1.38
1	C	64	CRU	C2-N3	-2.39	1.34	1.39
1	B	64	CRU	CE1-CZ	2.38	1.43	1.38
1	A	64	CRU	CA2-N2	-2.38	1.33	1.38
1	D	64	CRU	CB1-CA1	2.16	1.56	1.50
1	A	64	CRU	CA2-C2	-2.13	1.46	1.48
1	C	64	CRU	CA2-N2	-2.13	1.34	1.38
1	C	64	CRU	O2-C2	2.11	1.27	1.23

All (46) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	64	CRU	CB2-CA2-C2	15.19	140.40	122.28
1	D	64	CRU	CB2-CA2-C2	14.30	139.34	122.28
1	C	64	CRU	CB2-CA2-C2	13.59	138.50	122.28
1	A	64	CRU	CB2-CA2-C2	13.27	138.12	122.28
1	D	64	CRU	CA2-C2-N3	12.48	109.27	103.37
1	B	64	CRU	CA2-C2-N3	12.14	109.11	103.37
1	B	64	CRU	CB2-CA2-N2	-10.98	113.59	128.83
1	C	64	CRU	CB2-CA2-N2	-9.79	115.24	128.83
1	A	64	CRU	CB2-CA2-N2	-9.59	115.53	128.83
1	D	64	CRU	CB2-CA2-N2	-9.40	115.79	128.83
1	C	64	CRU	CA2-C2-N3	9.35	107.79	103.37
1	A	64	CRU	CA2-C2-N3	9.05	107.65	103.37
1	D	64	CRU	CB1-CG1-CD3	6.27	123.18	112.67
1	A	64	CRU	CG2-CB2-CA2	-6.22	122.32	129.94
1	D	64	CRU	C2-CA2-N2	-6.04	104.70	108.93
1	A	64	CRU	O2-C2-CA2	-5.77	127.72	130.96
1	C	64	CRU	O2-C2-CA2	-5.40	127.93	130.96
1	D	64	CRU	O-C-CA3	-5.10	111.00	126.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	64	CRU	C2-CA2-N2	-5.00	105.43	108.93
1	B	64	CRU	CG2-CB2-CA2	-4.76	124.11	129.94
1	C	64	CRU	O-C-CA3	-4.52	112.75	126.39
1	A	64	CRU	O-C-CA3	-4.32	113.35	126.39
1	A	64	CRU	C2-CA2-N2	-4.31	105.91	108.93
1	C	64	CRU	CE1-CD1-CG2	-4.19	115.78	121.25
1	D	64	CRU	CG2-CB2-CA2	-4.10	124.92	129.94
1	C	64	CRU	C2-CA2-N2	-4.07	106.08	108.93
1	D	64	CRU	CA3-N3-C1	3.18	134.45	128.22
1	C	64	CRU	CD2-CG2-CD1	3.04	122.14	117.64
1	C	64	CRU	CG2-CB2-CA2	-2.94	126.34	129.94
1	D	64	CRU	CA2-N2-C1	2.94	109.70	104.33
1	C	64	CRU	N3-C1-N2	-2.90	109.45	113.28
1	D	64	CRU	CE1-CD1-CG2	-2.86	117.52	121.25
1	B	64	CRU	O-C-CA3	-2.82	117.88	126.39
1	B	64	CRU	CD2-CG2-CD1	2.70	121.63	117.64
1	C	64	CRU	CA2-N2-C1	2.62	109.12	104.33
1	D	64	CRU	N3-C1-N2	-2.58	109.87	113.28
1	A	64	CRU	CB1-CG1-CD3	-2.48	108.51	112.67
1	B	64	CRU	CE1-CD1-CG2	-2.44	118.06	121.25
1	A	64	CRU	CD2-CG2-CD1	2.39	121.17	117.64
1	C	64	CRU	CD2-CE2-CZ	-2.28	117.38	119.88
1	D	64	CRU	CA3-N3-C2	-2.20	118.75	123.80
1	C	64	CRU	CD1-CE1-CZ	2.20	122.29	119.88
1	A	64	CRU	CD1-CG2-CB2	-2.19	113.75	121.22
1	A	64	CRU	CA2-N2-C1	2.07	108.12	104.33
1	B	64	CRU	O2-C2-N3	-2.04	120.29	124.35
1	B	64	CRU	CA2-N2-C1	2.01	108.01	104.33

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	64	CRU	C2-CA2-CB2-CG2
1	B	64	CRU	C2-CA2-CB2-CG2
1	C	64	CRU	C1-CA1-CB1-CG1
1	C	64	CRU	C2-CA2-CB2-CG2
1	D	64	CRU	C2-CA2-CB2-CG2
1	D	64	CRU	C-CA3-N3-C2
1	D	64	CRU	CA1-CB1-CG1-CD3
1	A	64	CRU	N2-CA2-CB2-CG2
1	B	64	CRU	N2-CA2-CB2-CG2

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Mol	Chain	Res	Type	Atoms
1	C	64	CRU	N2-CA2-CB2-CG2
1	D	64	CRU	N2-CA2-CB2-CG2
1	D	64	CRU	C-CA3-N3-C1
1	A	64	CRU	C1-CA1-CB1-CG1
1	B	64	CRU	C1-CA1-CB1-CG1

There are no ring outliers.

4 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	64	CRU	1	0
1	B	64	CRU	3	0
1	C	64	CRU	1	0
1	D	64	CRU	1	0

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

19 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	B	303	-	3,3,3	0.09	0	2,2,2	0.27	0
2	EDO	C	301	-	3,3,3	0.18	0	2,2,2	0.33	0
2	EDO	D	303	-	3,3,3	0.42	0	2,2,2	0.58	0
2	EDO	B	302	-	3,3,3	0.07	0	2,2,2	0.10	0
2	EDO	A	302	-	3,3,3	0.37	0	2,2,2	0.66	0
2	EDO	C	302	-	3,3,3	0.10	0	2,2,2	0.16	0
2	EDO	C	303	-	3,3,3	0.03	0	2,2,2	0.15	0
2	EDO	D	302	-	3,3,3	0.07	0	2,2,2	0.20	0
2	EDO	B	301	-	3,3,3	0.10	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	D	305	-	3,3,3	0.08	0	2,2,2	0.16	0
2	EDO	A	303	-	3,3,3	0.13	0	2,2,2	0.22	0
2	EDO	B	305	-	3,3,3	0.10	0	2,2,2	0.34	0
2	EDO	D	304	-	3,3,3	0.12	0	2,2,2	0.32	0
2	EDO	D	301	-	3,3,3	0.08	0	2,2,2	0.08	0
2	EDO	C	304	-	3,3,3	0.15	0	2,2,2	0.36	0
2	EDO	A	301	-	3,3,3	0.22	0	2,2,2	0.39	0
2	EDO	A	304	-	3,3,3	0.07	0	2,2,2	0.27	0
2	EDO	B	304	-	3,3,3	0.12	0	2,2,2	0.22	0
2	EDO	A	305	-	3,3,3	0.09	0	2,2,2	0.16	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	B	303	-	-	1/1/1/1	-
2	EDO	C	301	-	-	1/1/1/1	-
2	EDO	D	303	-	-	0/1/1/1	-
2	EDO	B	302	-	-	1/1/1/1	-
2	EDO	A	302	-	-	0/1/1/1	-
2	EDO	C	302	-	-	1/1/1/1	-
2	EDO	C	303	-	-	1/1/1/1	-
2	EDO	D	302	-	-	1/1/1/1	-
2	EDO	B	301	-	-	1/1/1/1	-
2	EDO	D	305	-	-	1/1/1/1	-
2	EDO	A	303	-	-	1/1/1/1	-
2	EDO	B	305	-	-	1/1/1/1	-
2	EDO	D	304	-	-	1/1/1/1	-
2	EDO	D	301	-	-	1/1/1/1	-
2	EDO	C	304	-	-	1/1/1/1	-
2	EDO	A	301	-	-	1/1/1/1	-
2	EDO	A	304	-	-	1/1/1/1	-
2	EDO	B	304	-	-	1/1/1/1	-
2	EDO	A	305	-	-	0/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (16) torsion outliers are listed below:

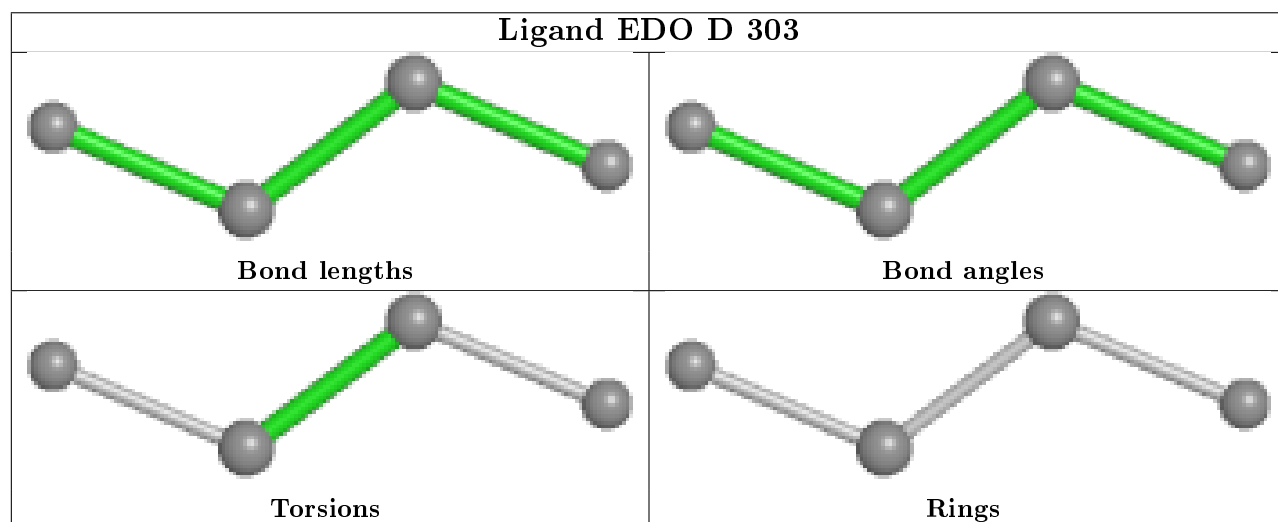
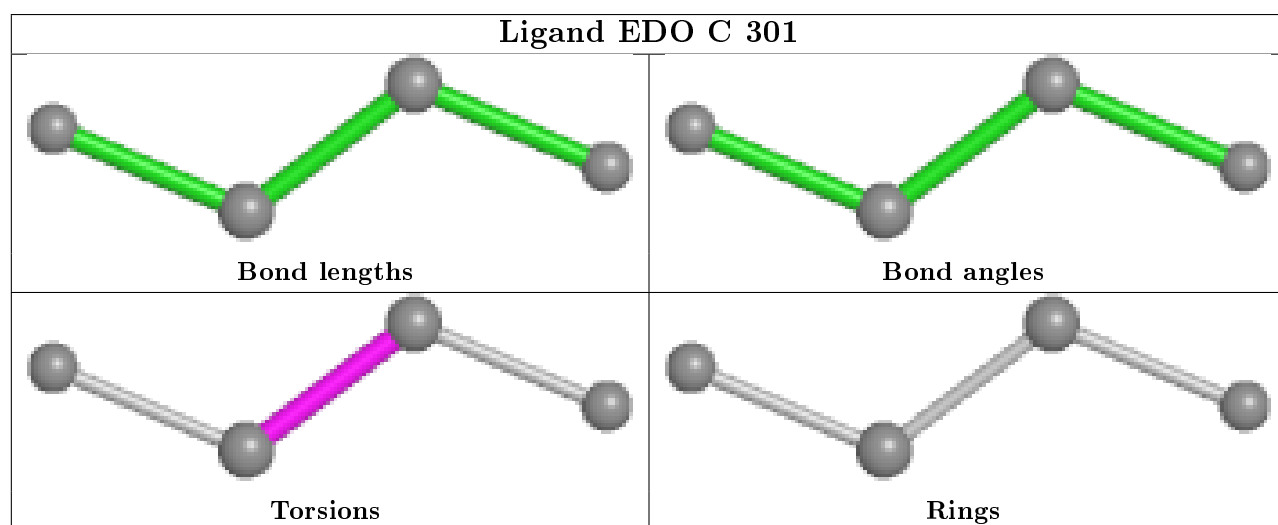
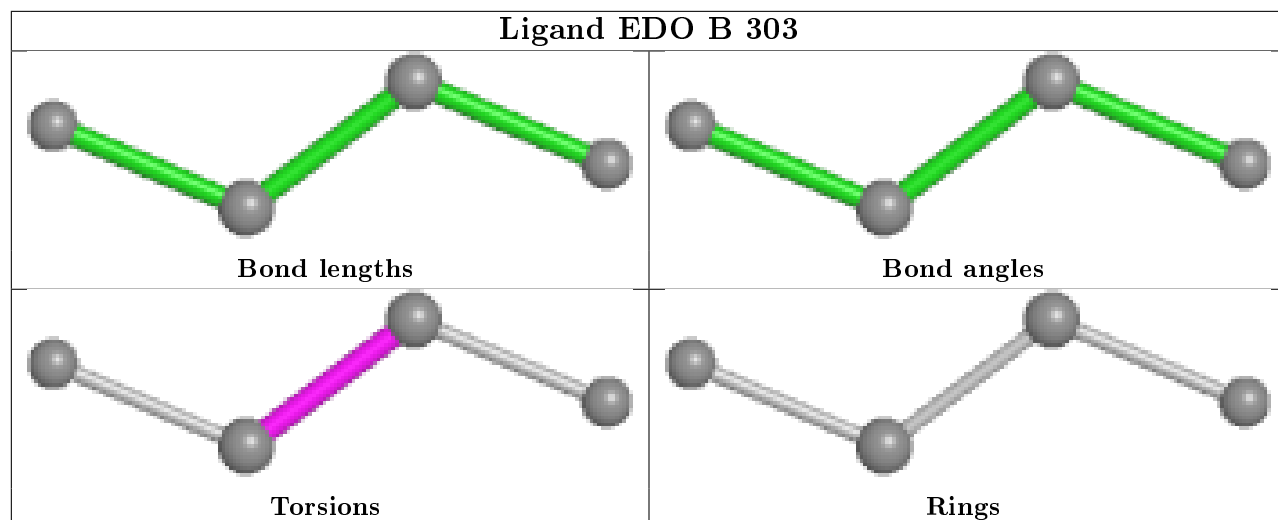
Mol	Chain	Res	Type	Atoms
2	C	301	EDO	O1-C1-C2-O2
2	B	302	EDO	O1-C1-C2-O2
2	C	302	EDO	O1-C1-C2-O2
2	B	301	EDO	O1-C1-C2-O2
2	D	305	EDO	O1-C1-C2-O2
2	A	303	EDO	O1-C1-C2-O2
2	A	301	EDO	O1-C1-C2-O2
2	A	304	EDO	O1-C1-C2-O2
2	D	302	EDO	O1-C1-C2-O2
2	C	303	EDO	O1-C1-C2-O2
2	B	305	EDO	O1-C1-C2-O2
2	D	301	EDO	O1-C1-C2-O2
2	D	304	EDO	O1-C1-C2-O2
2	B	304	EDO	O1-C1-C2-O2
2	B	303	EDO	O1-C1-C2-O2
2	C	304	EDO	O1-C1-C2-O2

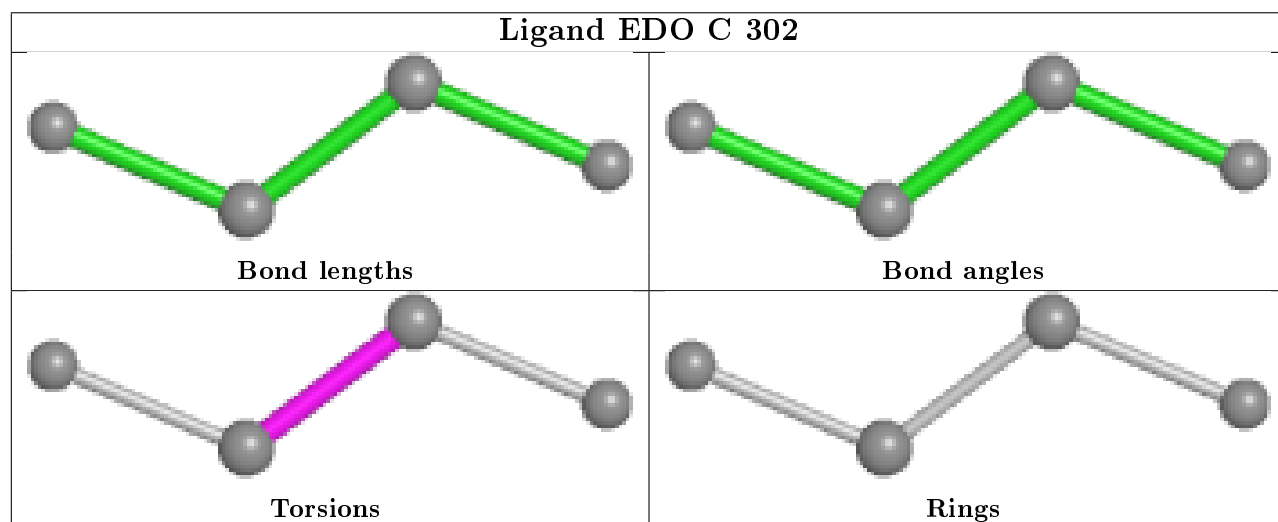
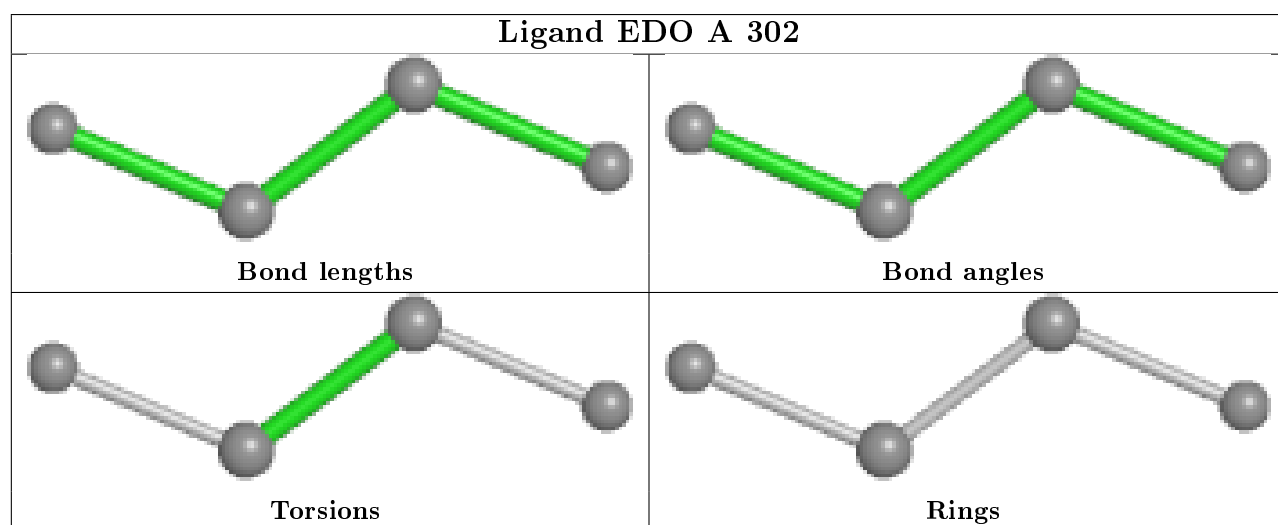
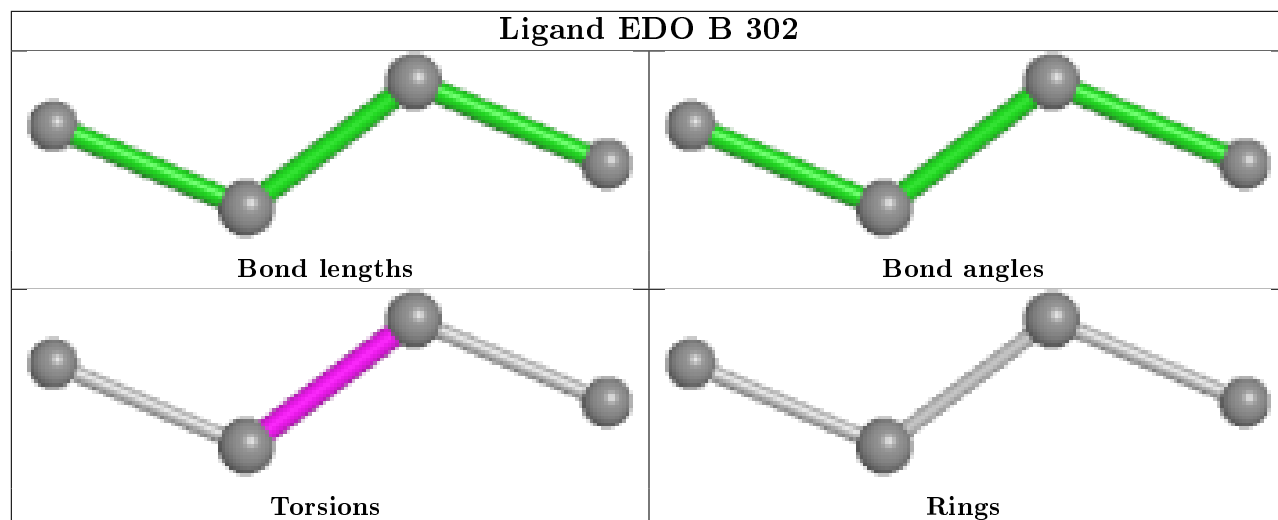
There are no ring outliers.

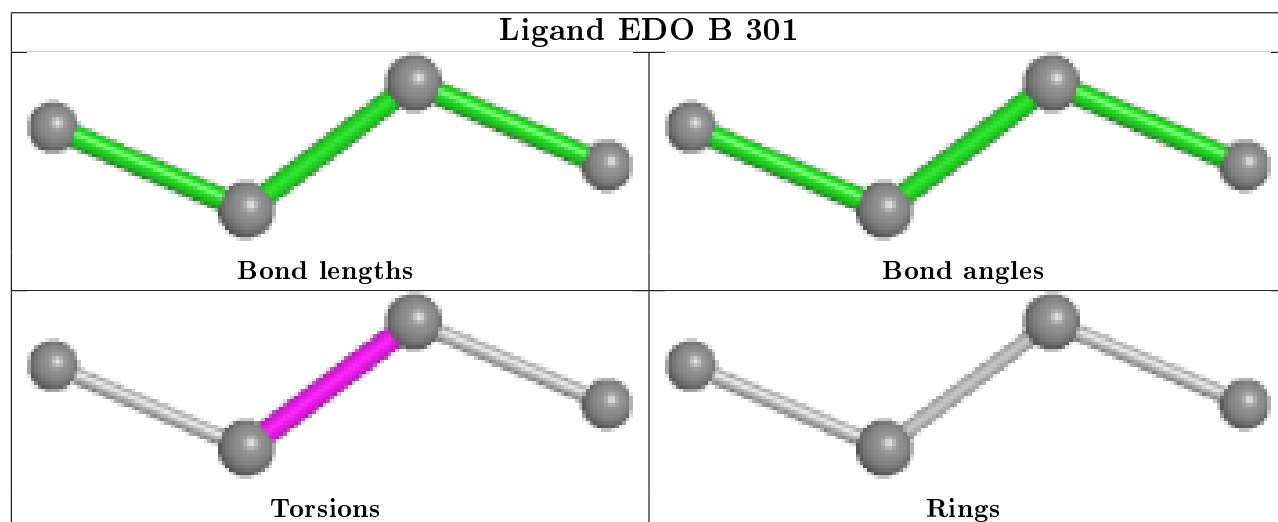
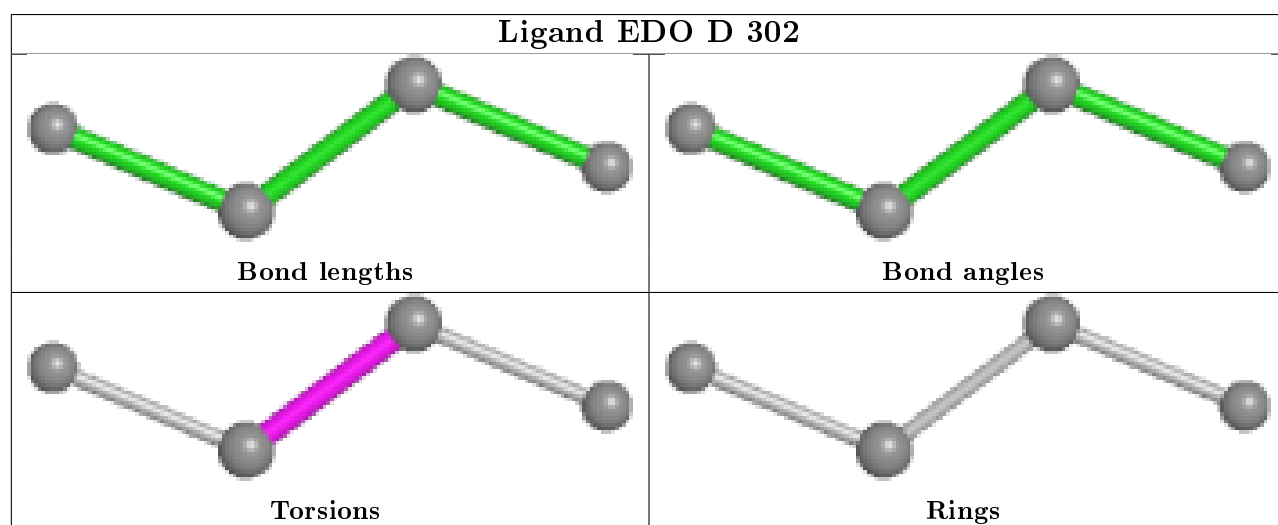
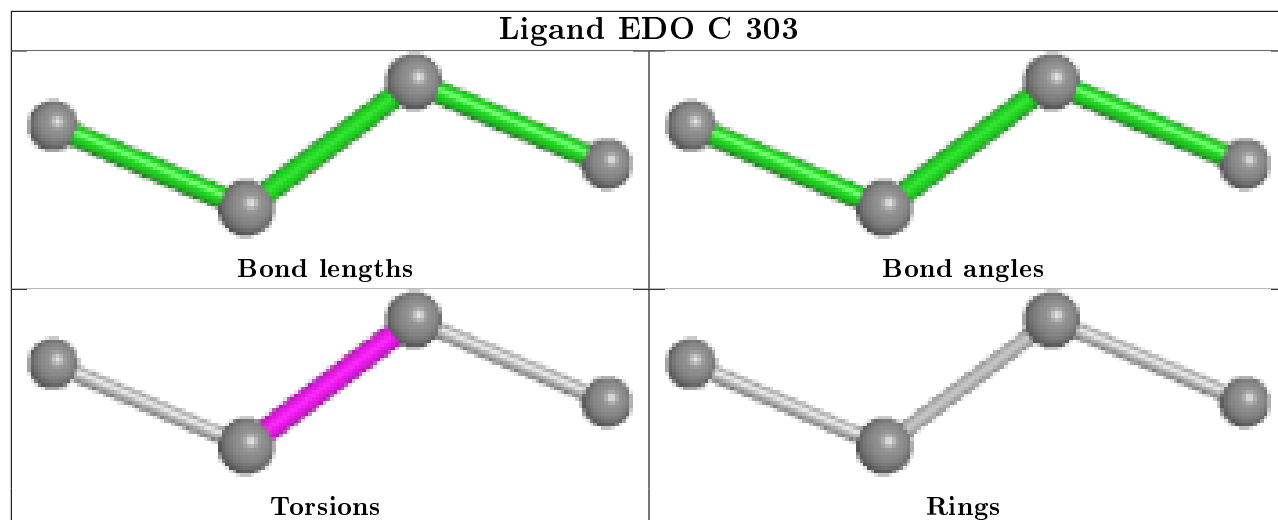
1 monomer is involved in 1 short contact:

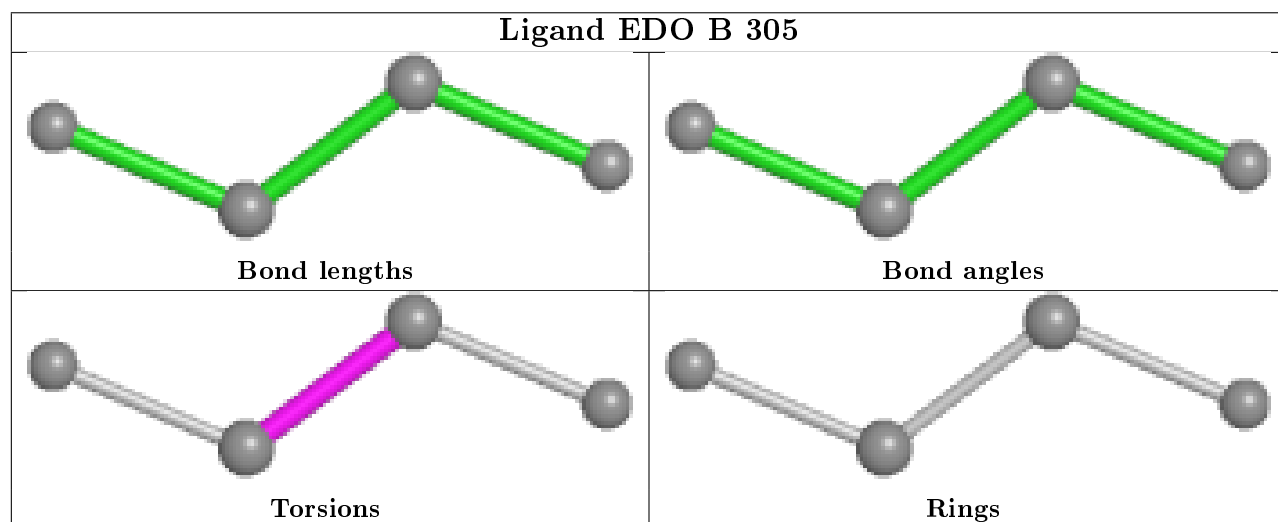
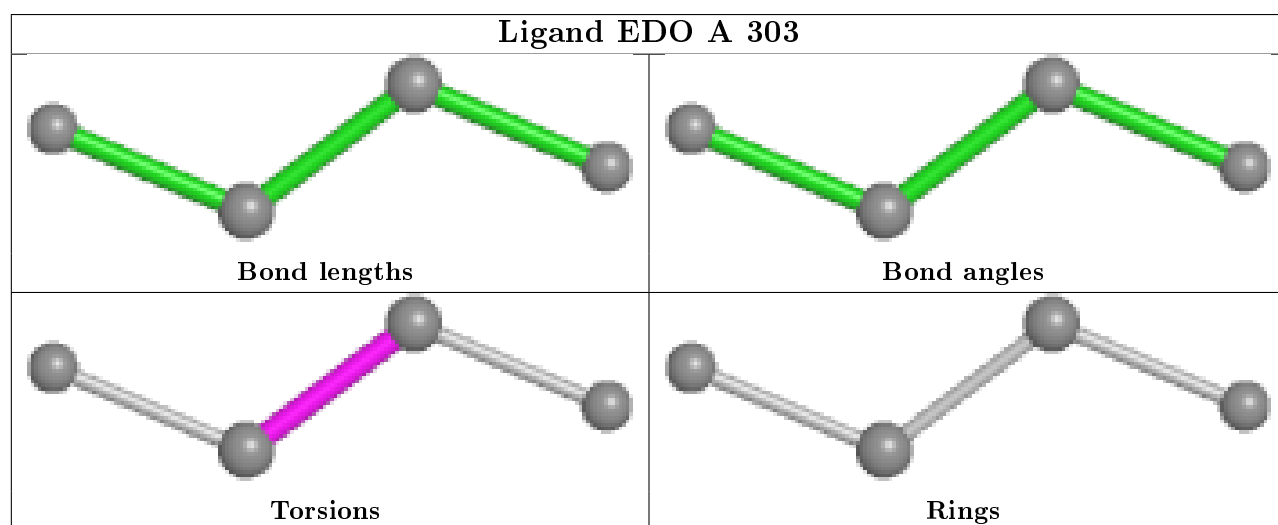
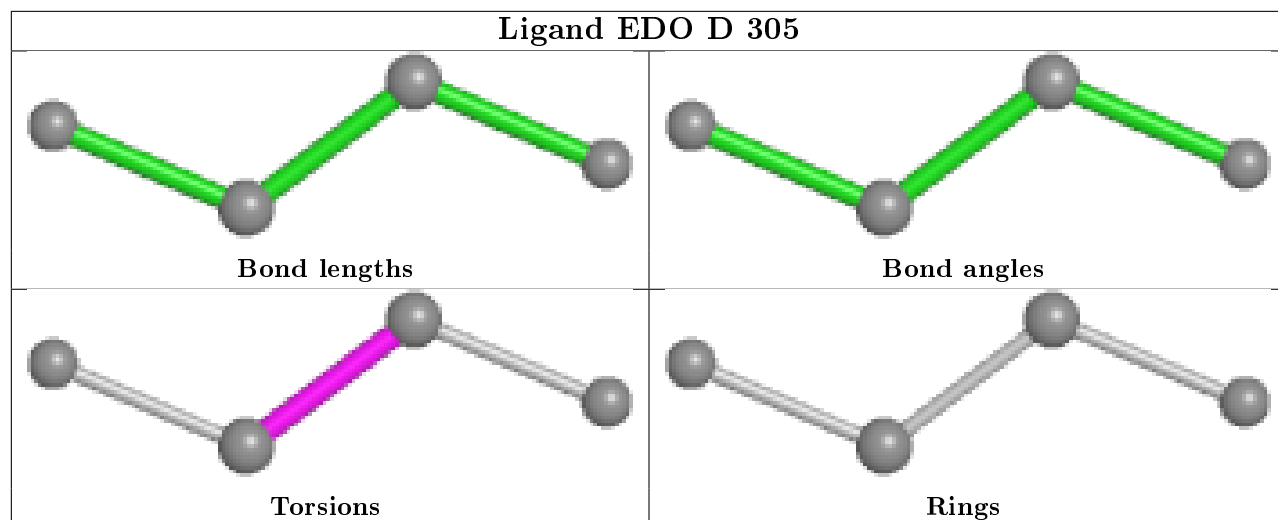
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	302	EDO	1	0

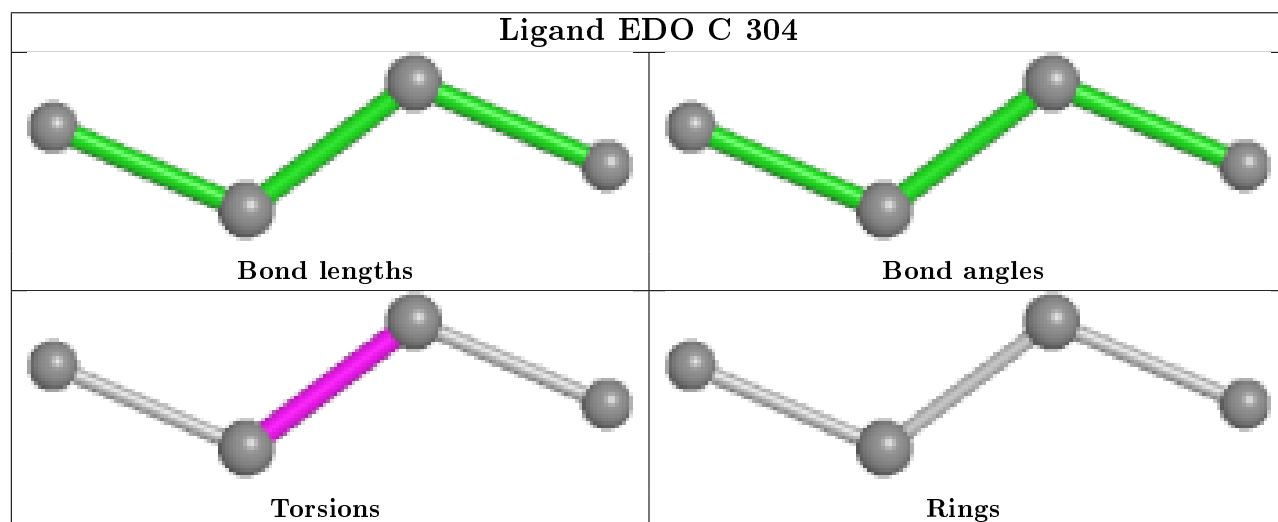
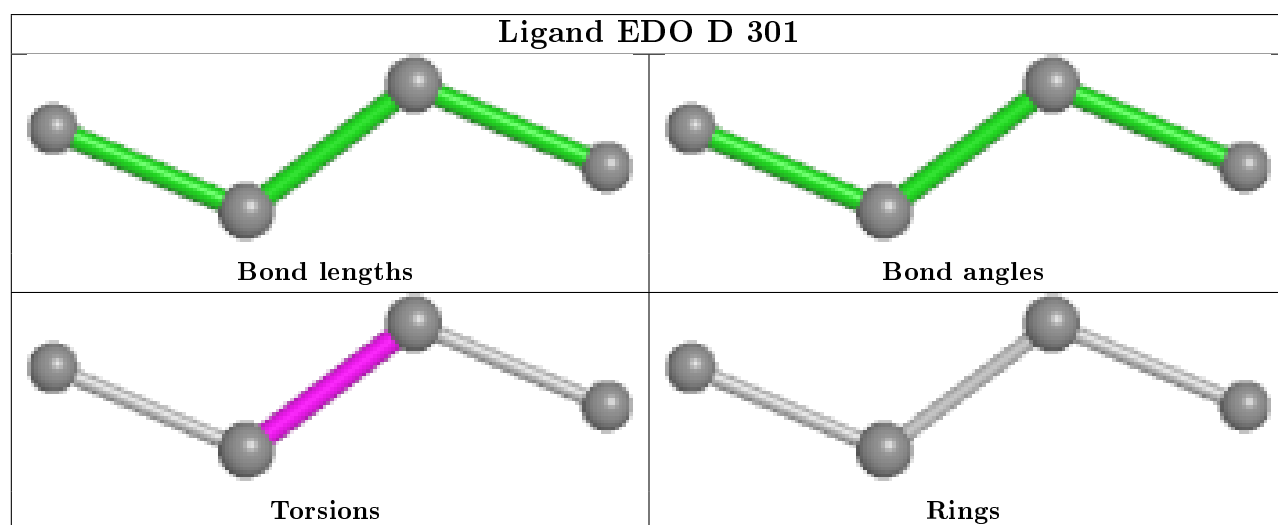
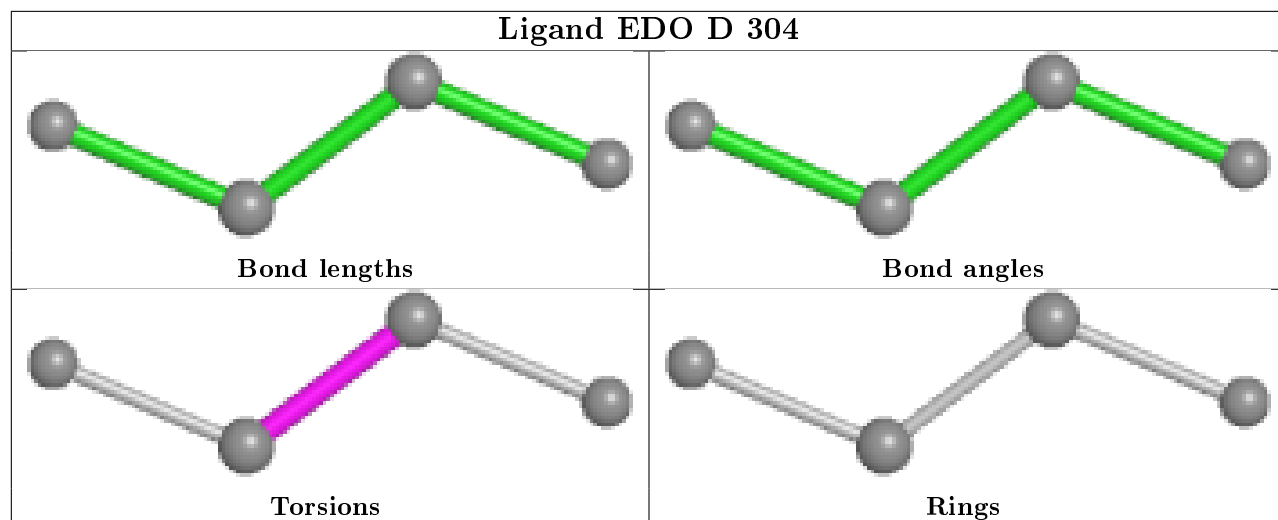
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

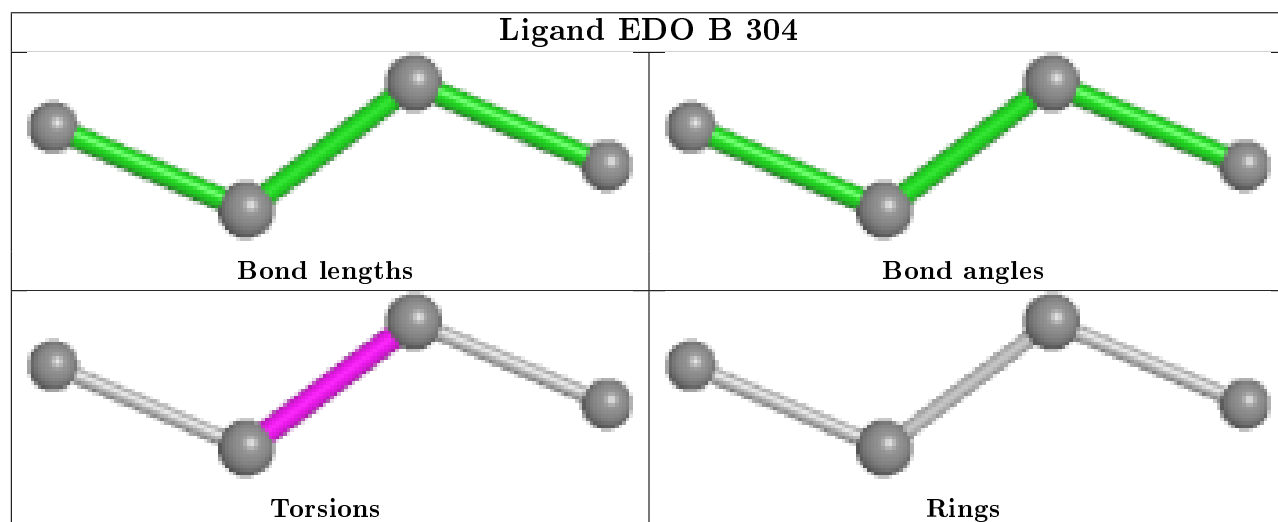
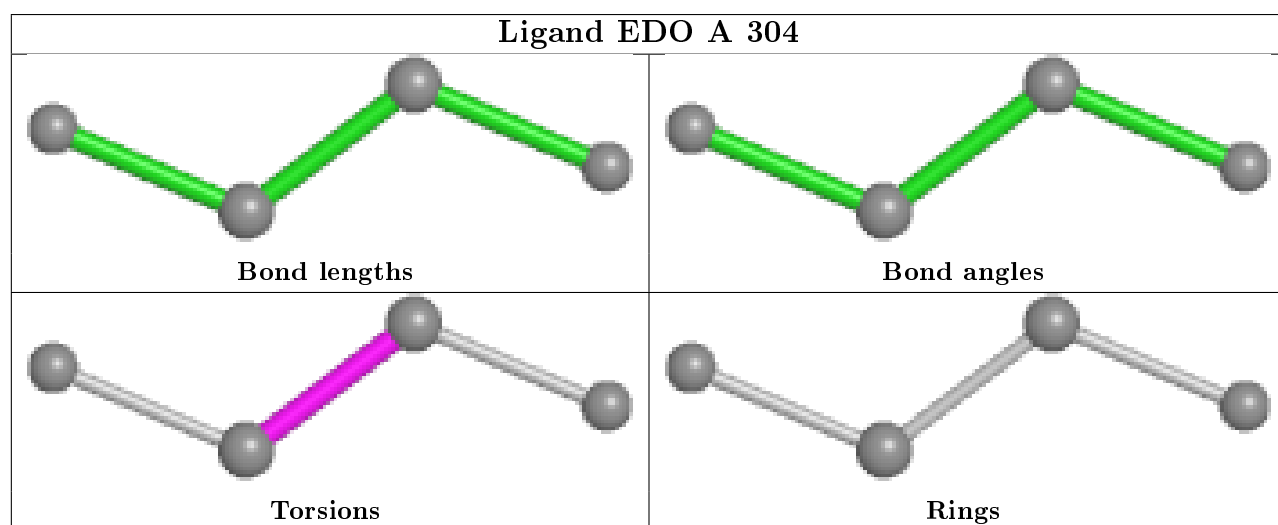
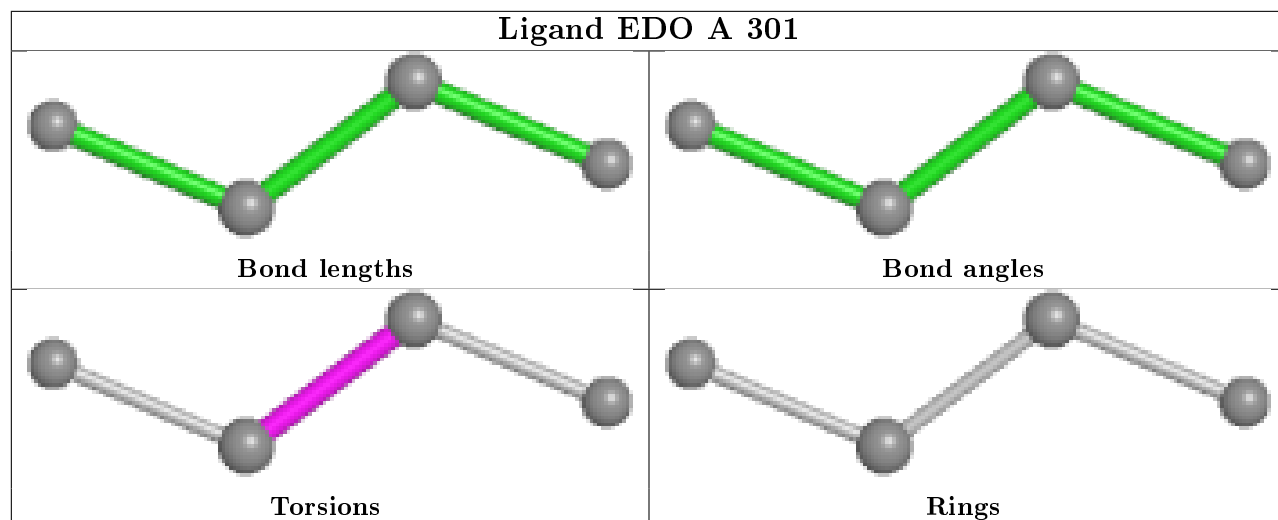


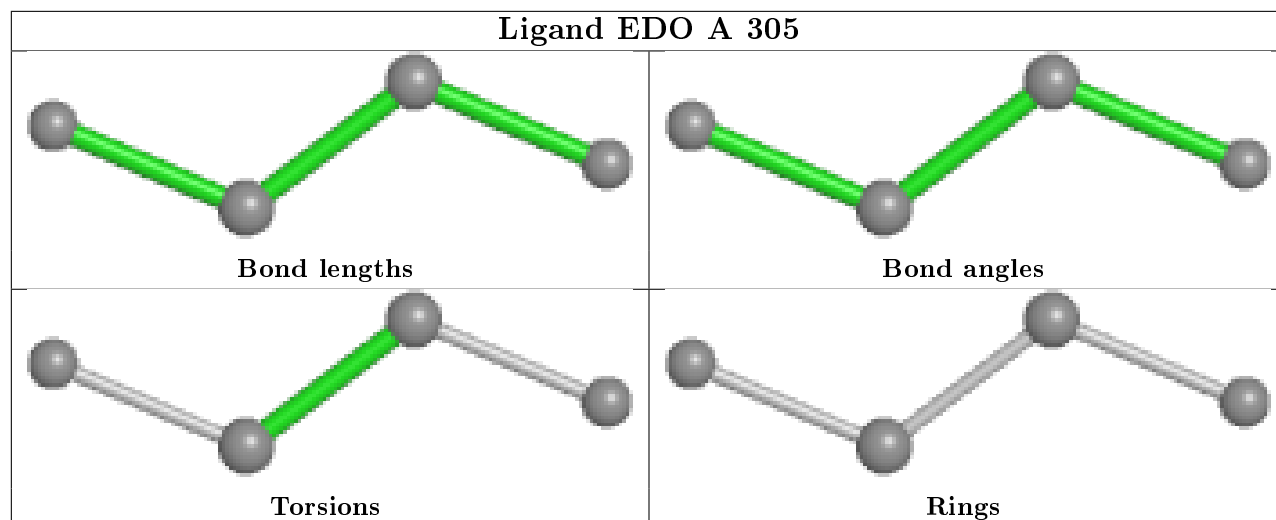












5.7 Other polymers [i](#)

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	221/236 (93%)	-0.42	1 (0%) 91 94	21, 29, 49, 71	1 (0%)
1	B	221/236 (93%)	-0.43	0 100 100	19, 26, 45, 70	3 (1%)
1	C	221/236 (93%)	-0.33	1 (0%) 91 94	19, 29, 49, 72	4 (1%)
1	D	221/236 (93%)	-0.42	2 (0%) 84 88	20, 30, 46, 70	2 (0%)
All	All	884/944 (93%)	-0.40	4 (0%) 91 94	19, 29, 49, 72	10 (1%)

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	4	GLY	3.7
1	D	227	ALA	2.8
1	A	227	ALA	2.2
1	C	4	GLY	2.1

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

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
1	CRU	C	64	24/25	0.92	0.17	30,44,63,72	0
1	CRU	D	64	24/25	0.92	0.16	32,39,51,64	0
1	CRU	A	64	24/25	0.94	0.15	29,40,53,61	0
1	CRU	B	64	24/25	0.94	0.14	26,39,53,59	0

6.3 Carbohydrates ⓘ

There are no monosaccharides in this entry.

6.4 Ligands ⓘ

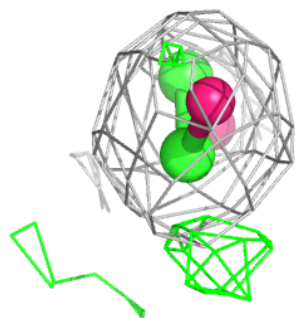
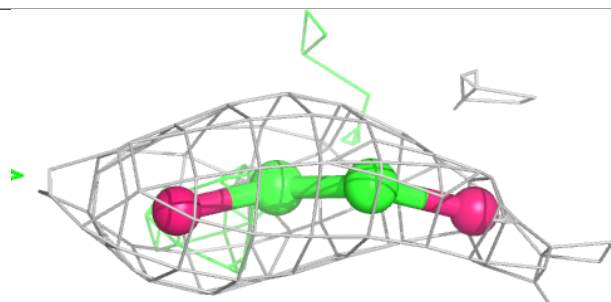
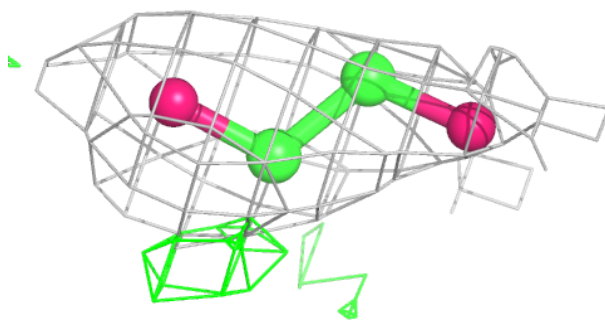
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	EDO	D	304	4/4	0.72	0.40	59,60,61,63	0
2	EDO	A	303	4/4	0.78	0.18	62,66,67,67	0
2	EDO	D	303	4/4	0.79	0.35	45,46,48,49	0
2	EDO	A	304	4/4	0.80	0.16	59,63,64,64	0
2	EDO	A	302	4/4	0.83	0.29	50,55,56,56	0
2	EDO	C	303	4/4	0.84	0.41	58,58,58,61	0
2	EDO	A	301	4/4	0.85	0.43	53,59,62,66	0
2	EDO	C	301	4/4	0.86	0.29	50,52,52,53	0
2	EDO	C	304	4/4	0.87	0.27	52,56,57,60	0
2	EDO	B	302	4/4	0.87	0.13	48,54,57,59	0
2	EDO	D	305	4/4	0.87	0.37	61,65,67,69	0
2	EDO	A	305	4/4	0.87	0.19	50,53,58,61	0
2	EDO	B	304	4/4	0.91	0.16	60,66,66,68	0
2	EDO	B	301	4/4	0.91	0.20	54,57,57,59	0
2	EDO	D	302	4/4	0.92	0.16	48,50,53,56	0
2	EDO	C	302	4/4	0.93	0.17	52,56,58,61	0
2	EDO	B	303	4/4	0.94	0.14	56,56,57,58	0
2	EDO	B	305	4/4	0.95	0.22	52,55,55,55	0
2	EDO	D	301	4/4	0.96	0.13	51,55,55,57	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.

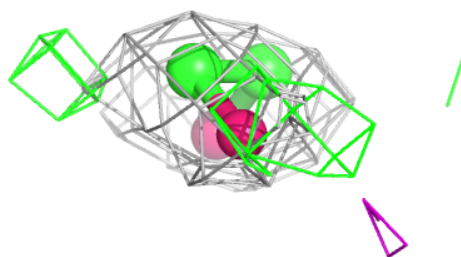
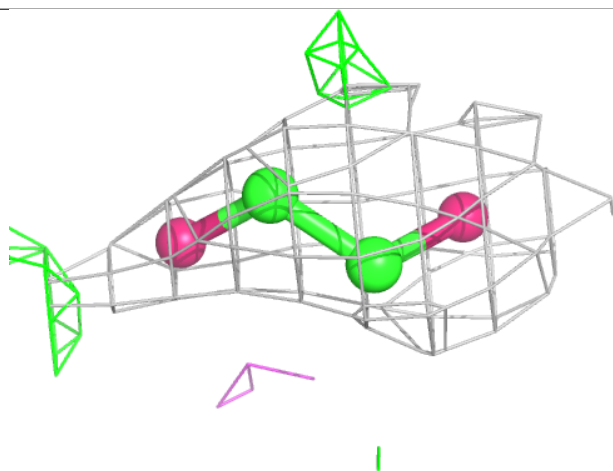
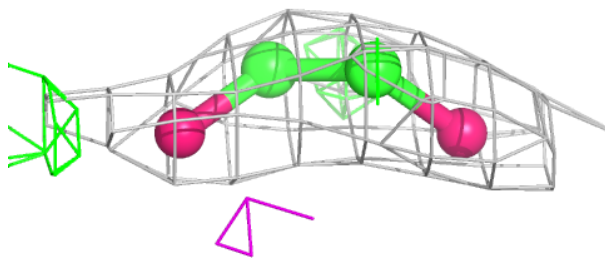
Electron density around EDO D 304:

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



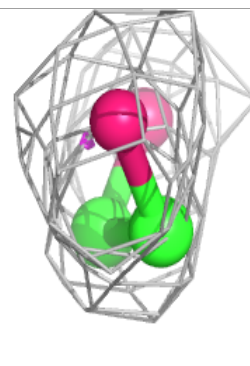
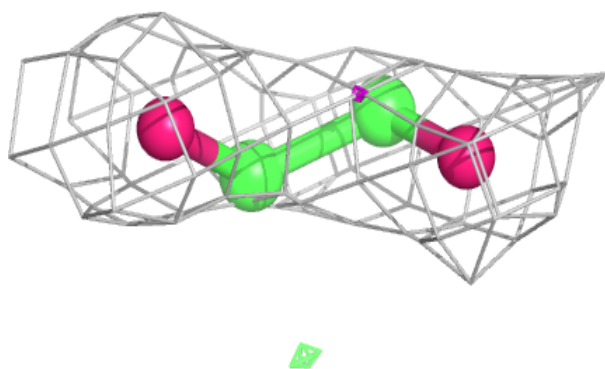
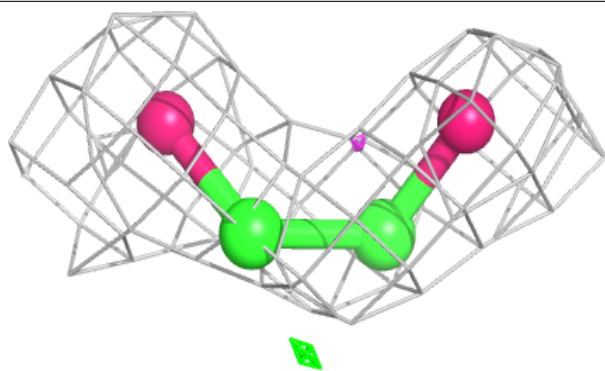
Electron density around EDO A 303:

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

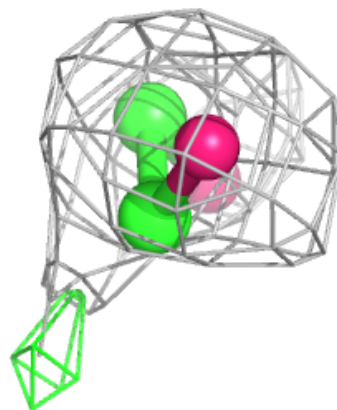
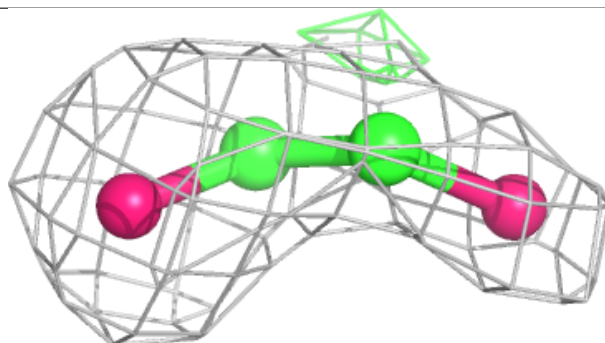
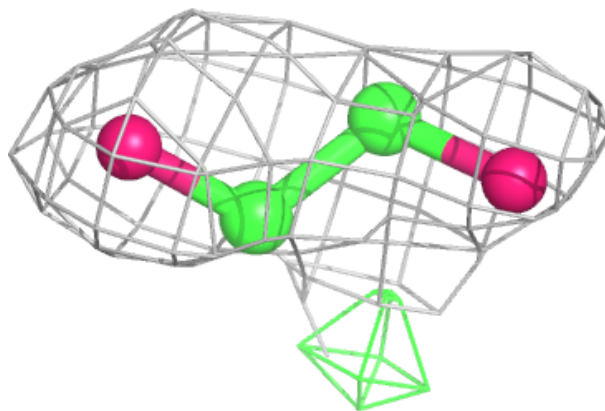


Electron density around EDO D 303:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

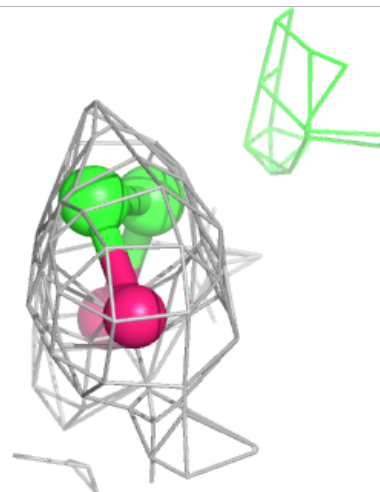
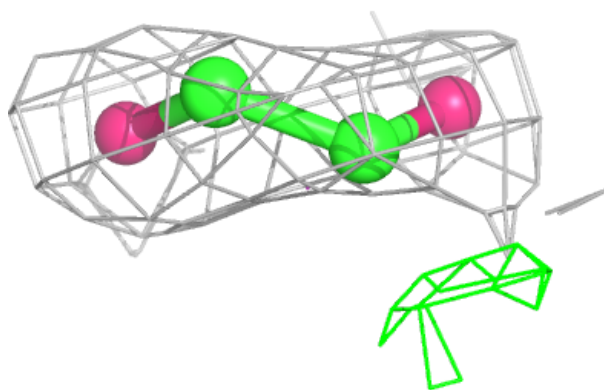
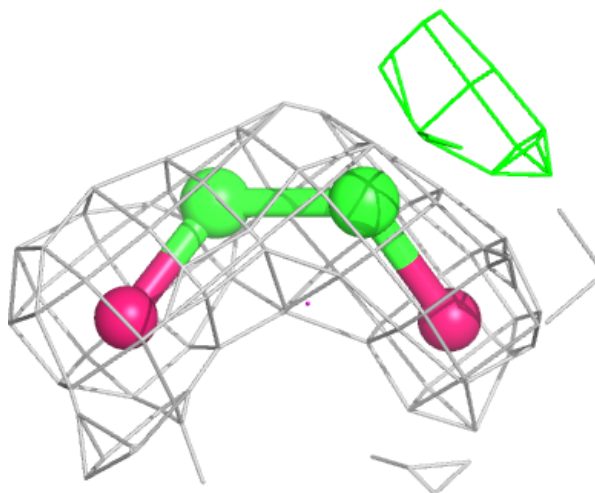
**Electron density around EDO A 304:**

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



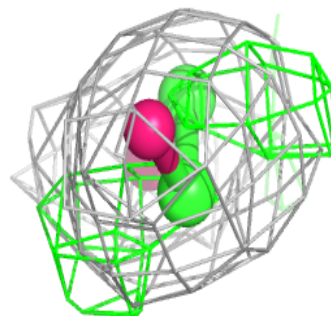
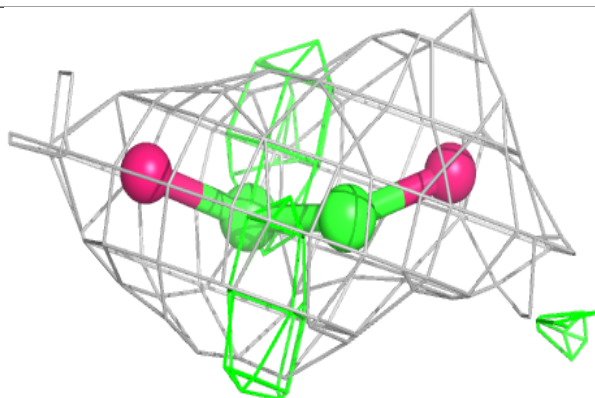
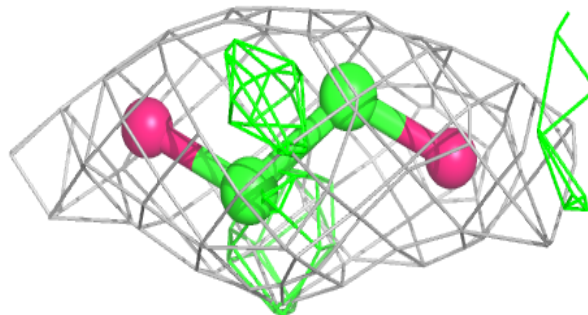
Electron density around EDO A 302:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

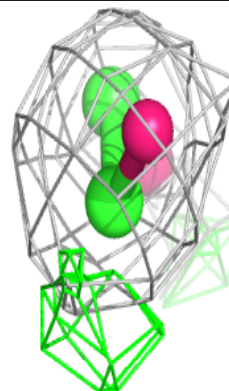
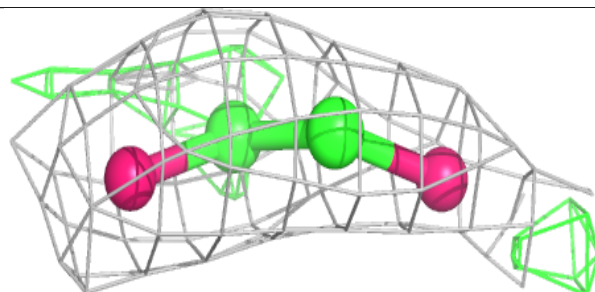
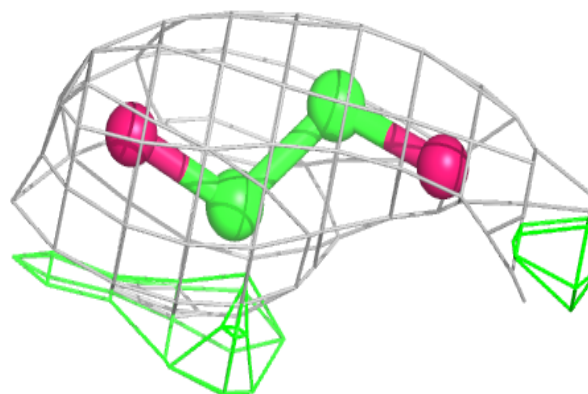


Electron density around EDO C 303:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

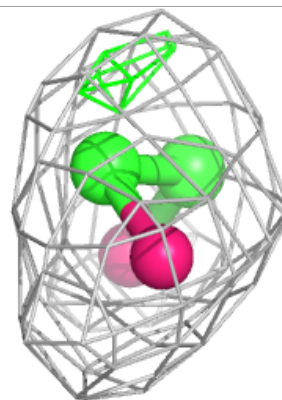
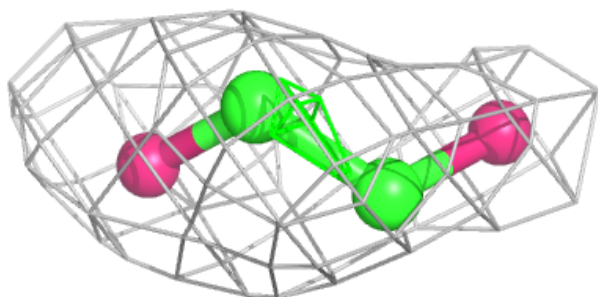
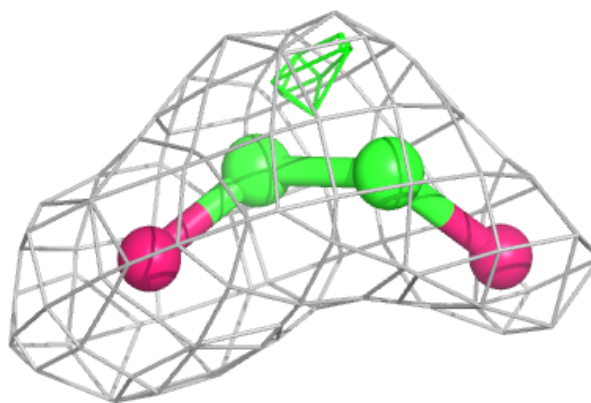
**Electron density around EDO A 301:**

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

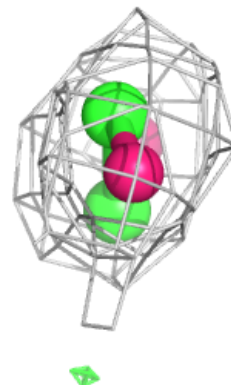
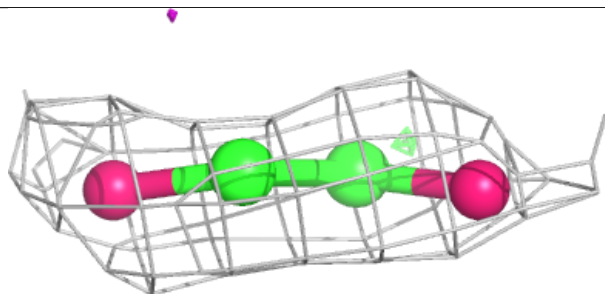
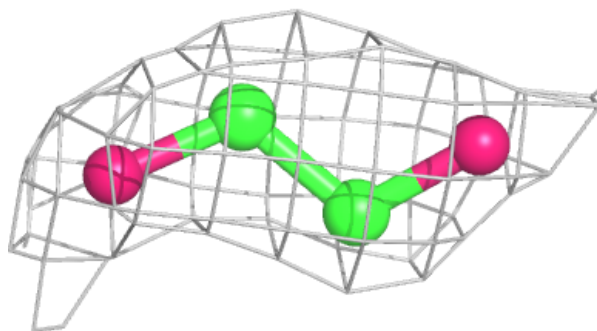


Electron density around EDO C 301:

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

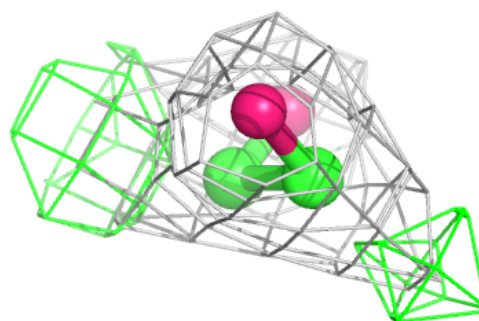
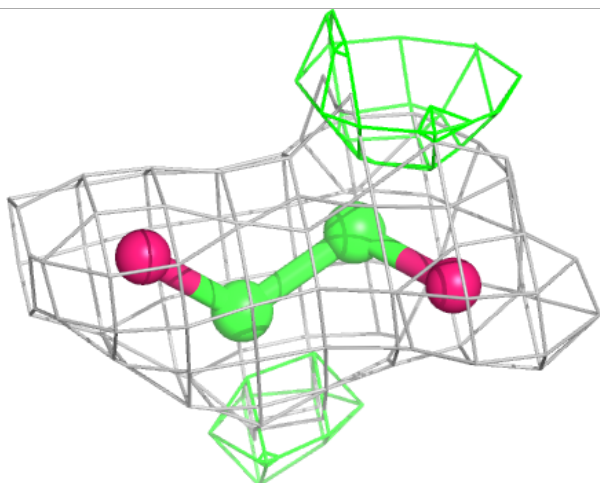
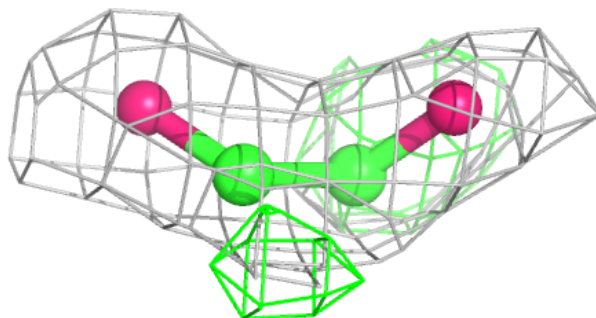
**Electron density around EDO C 304:**

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



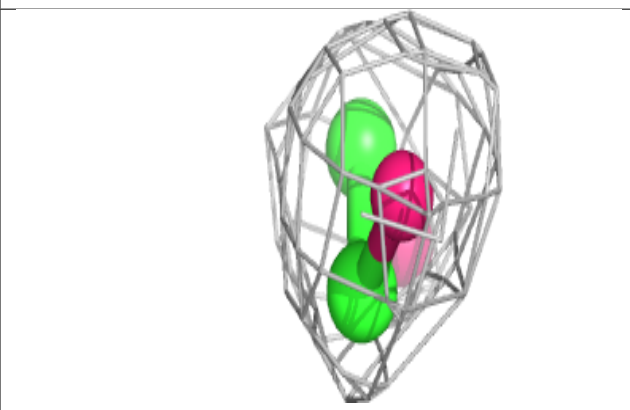
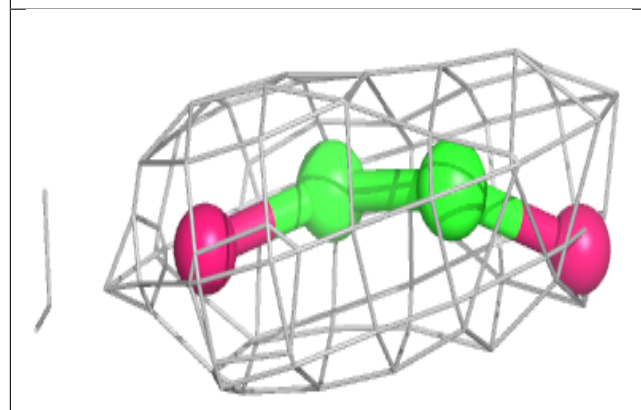
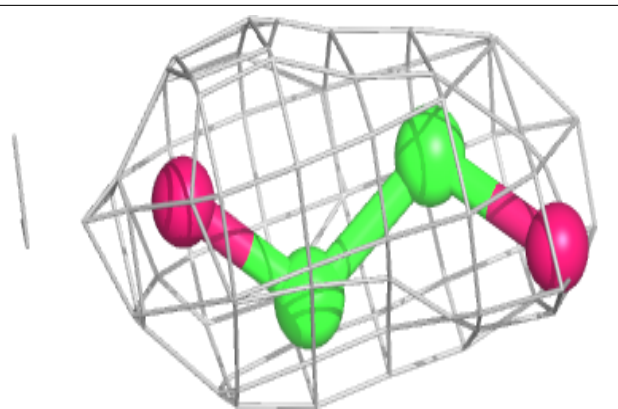
Electron density around EDO B 302:

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



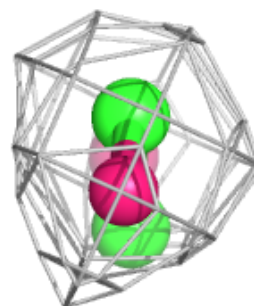
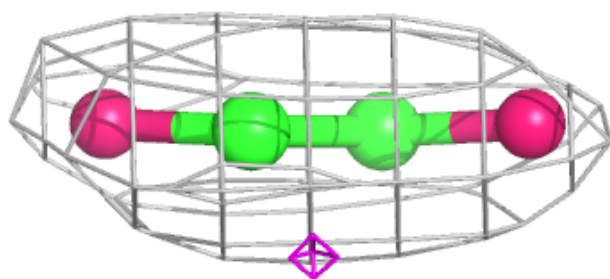
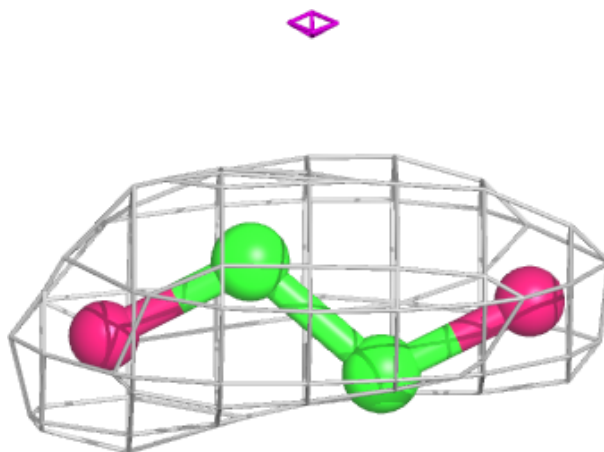
Electron density around EDO D 305:

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



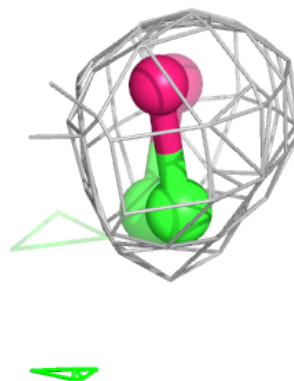
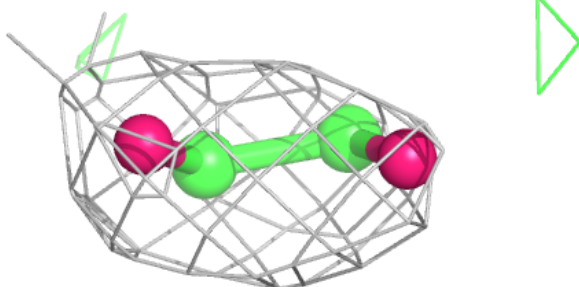
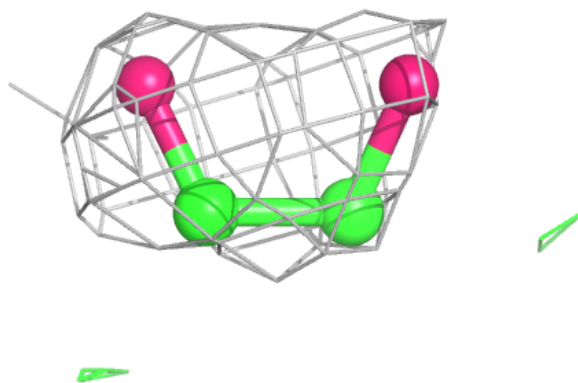
Electron density around EDO A 305:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

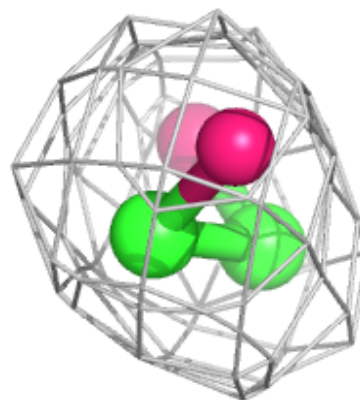
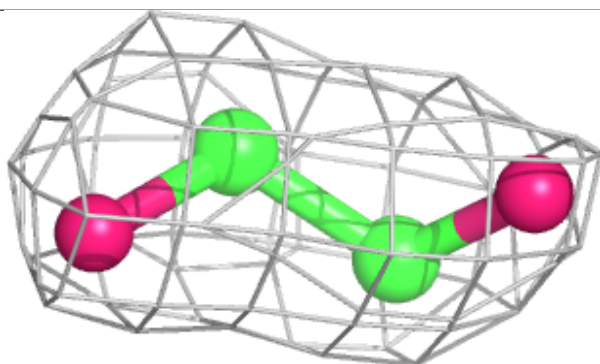
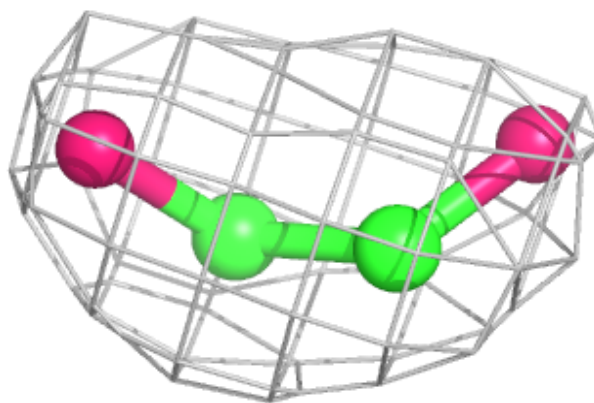


Electron density around EDO B 304:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

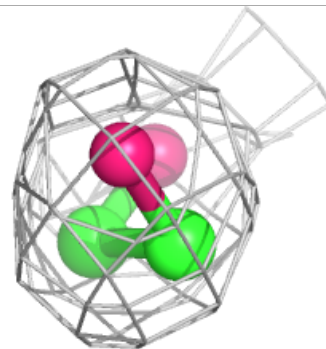
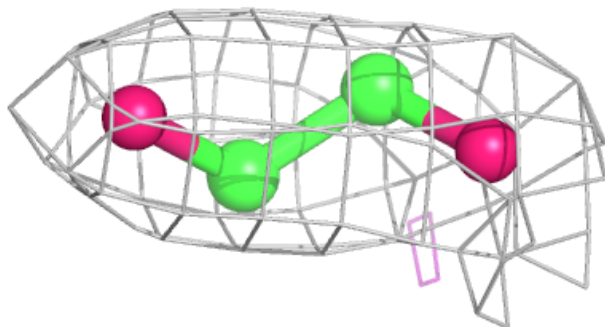
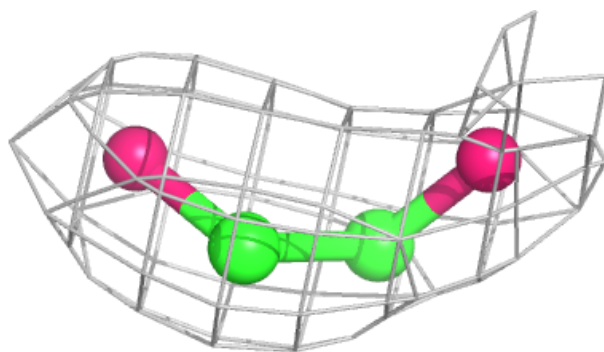
**Electron density around EDO B 301:**

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



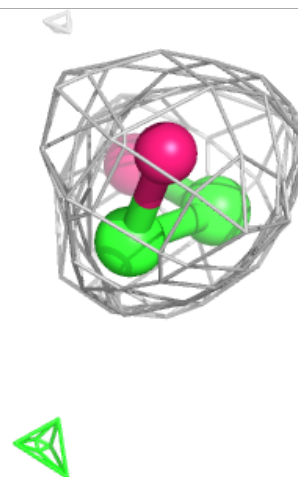
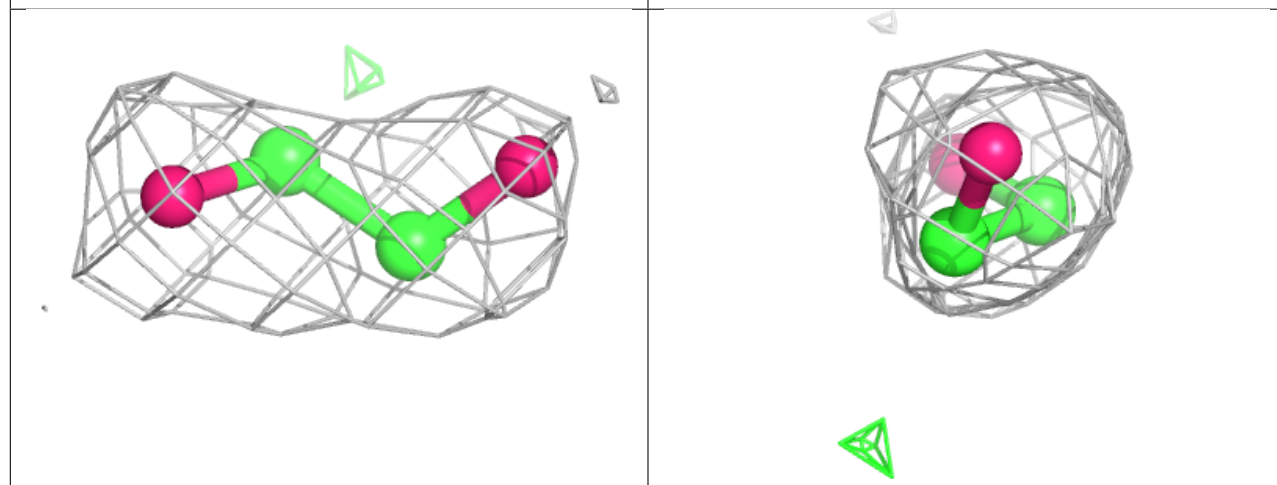
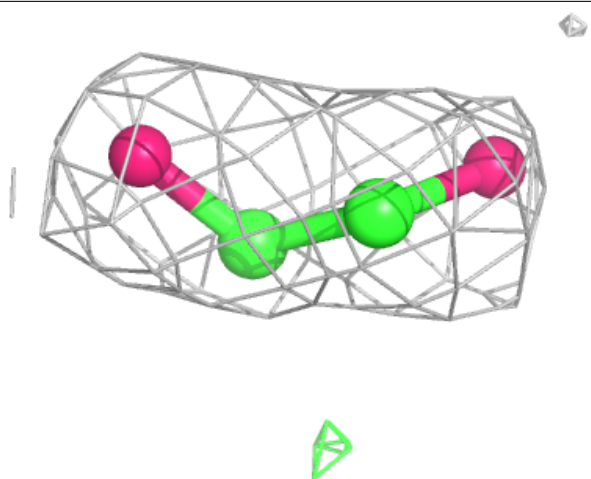
Electron density around EDO D 302:

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



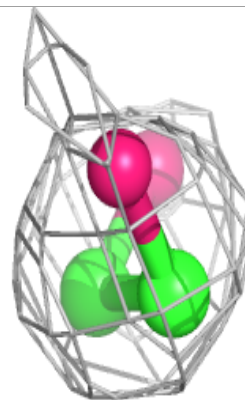
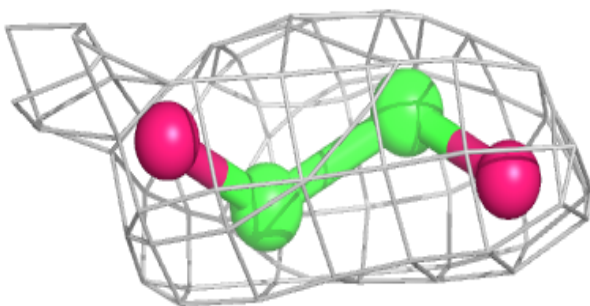
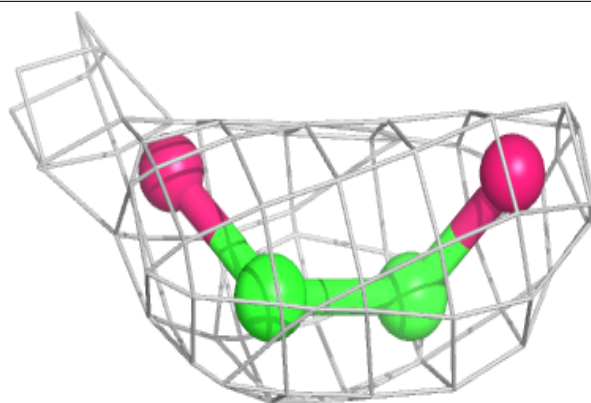
Electron density around EDO C 302:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

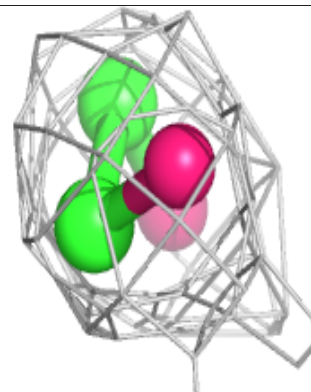
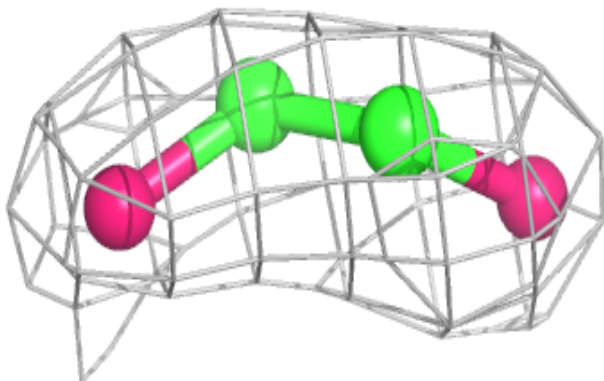
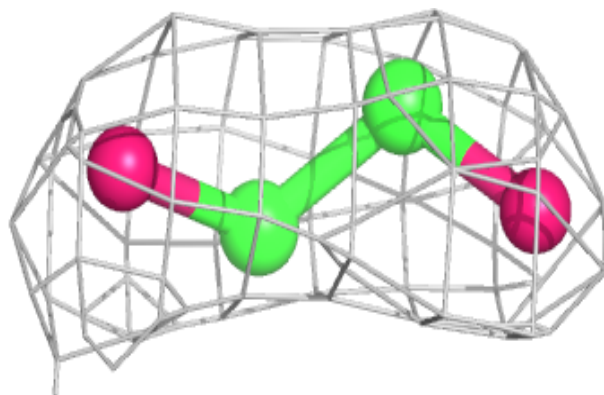


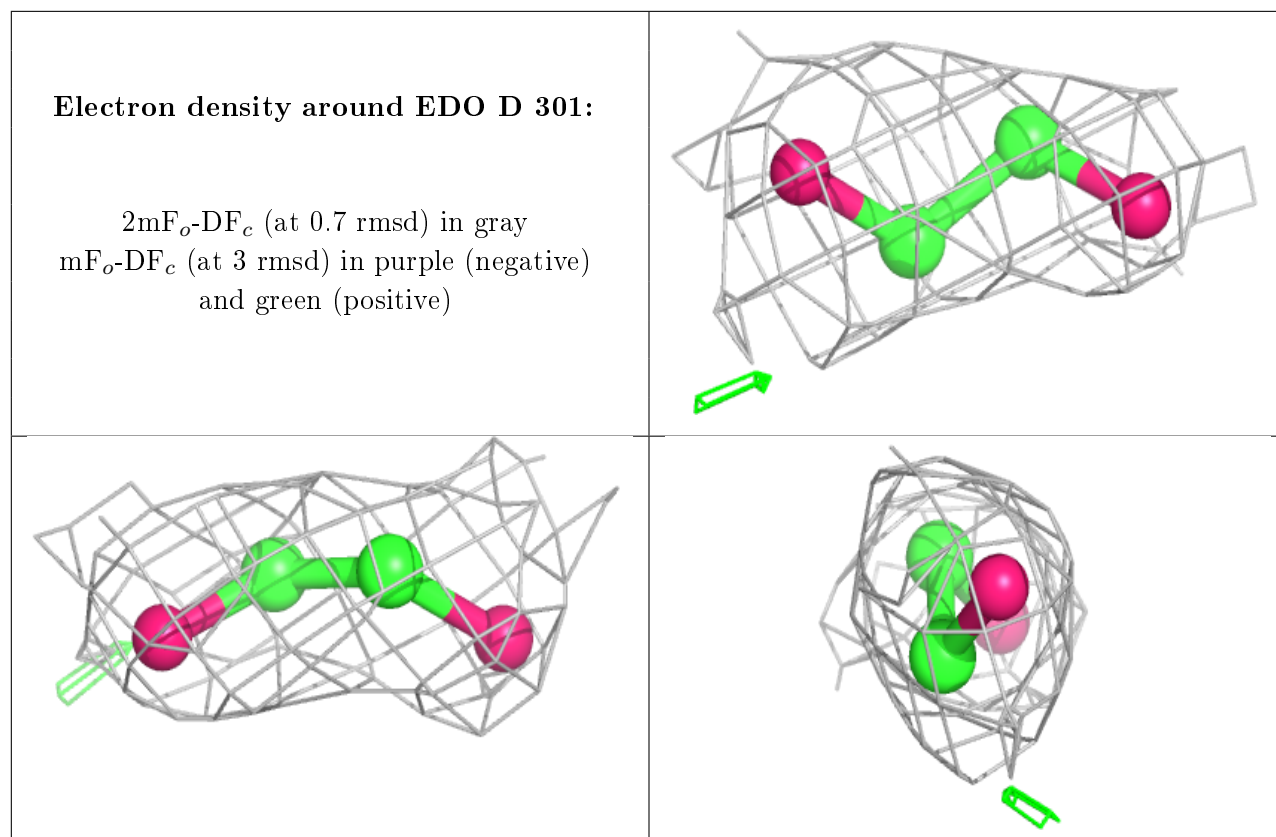
Electron density around EDO B 303:

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

**Electron density around EDO B 305:**

$2mF_o-DF_c$ (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.