

wwPDB X-ray Structure Validation Summary Report (i)

Oct 10, 2023 – 03:36 PM EDT

PDB ID : 7KXG

Title: Dihydrodipicolinate synthase (DHDPS) from C.jejuni with pyruvate bound in

the active site and L-histidine bound at the allosteric site

Authors: Saran, S.; Sanders, D.A.R.

Deposited on : 2020-12-03

Resolution : 2.28 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 (i) 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 (1)) 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 : 2.35.1

buster-report : 1.1.7 (2018)

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

 $Refmac \quad : \quad 5.8.0158$

CCP4 : 7.0.044 (Gargrove)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

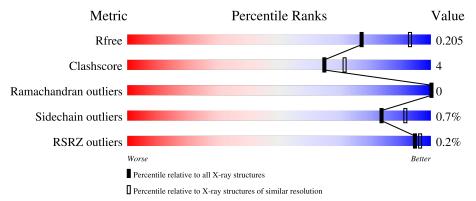
Validation Pipeline (wwPDB-VP) : 2.35.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: X-RAY DIFFRACTION

The reported resolution of this entry is 2.28 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 $(\# \mathrm{Entries})$	Similar resolution $(\# \text{Entries, resolution range}(\text{\AA}))$
R_{free}	130704	6980 (2.30-2.26)
Clashscore	141614	7711 (2.30-2.26)
Ramachandran outliers	138981	7597 (2.30-2.26)
Sidechain outliers	138945	7598 (2.30-2.26)
RSRZ outliers	127900	6849 (2.30-2.26)

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 <=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	310	85% 11%	5%
1	В	310	85% 10%	• 5%
1	С	310	87% 8%	. •
1	D	310	88% 79	6 5%
1	Е	310	87% 8%	• 5%



Continued from previous page...

Mol	Chain	Length	Quality of chain					
-	Б	210						
1	F'	310	87%	8%	•			

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
7	ACT	В	307	-	-	X	X



2 Entry composition (i)

There are 9 unique types of molecules in this entry. The entry contains 14653 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 4-hydroxy-tetrahydrodipicolinate synthase.

Mol	Chain	Residues		At	oms			ZeroOcc	AltConf	Trace	
1	A	296	Total	С	Ν	О	S	0	1	0	
1	Λ	290	2281	1450	378	440	13	U	1		
1	В	296	Total	С	N	О	S	0	0	0	
1	Ъ	290	2269	1444	375	437	13	U	0	U	
1	С	297	Total	С	N	О	S	0	0	0	
1		291	2278	1448	377	440	13	0			
1	D	296	Total	С	N	О	S	0	1	0	
1	D	290	2285	1453	380	439	13	U	1		
1	E	296	Total	С	N	О	S	0	0	0	
1	l L	290	2273	1446	377	437	13	U	0		
1	F	297	Total	С	N	О	S	0	0	0	
1	I'	291	2275	1446	378	438	13	U	U		

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-11	MET	-	expression tag	UNP Q9PPB4
A	-10	ARG	-	expression tag	UNP Q9PPB4
A	-9	GLY	-	expression tag	UNP Q9PPB4
A	-8	SER	-	expression tag	UNP Q9PPB4
A	-7	HIS	-	expression tag	UNP Q9PPB4
A	-6	HIS	-	expression tag	UNP Q9PPB4
A	-5	HIS	-	expression tag	UNP Q9PPB4
A	-4	HIS	-	expression tag	UNP Q9PPB4
A	-3	HIS	-	expression tag	UNP Q9PPB4
A	-2	HIS	-	expression tag	UNP Q9PPB4
A	-1	GLY	-	expression tag	UNP Q9PPB4
A	0	SER	-	expression tag	UNP Q9PPB4
В	-11	MET	-	expression tag	UNP Q9PPB4
В	-10	ARG	-	expression tag	UNP Q9PPB4
В	-9	GLY	-	expression tag	UNP Q9PPB4
В	-8	SER	-	expression tag	UNP Q9PPB4
В	-7	HIS	-	expression tag	UNP Q9PPB4



 $Continued\ from\ previous\ page...$

Chain	Residue	Modelled Modelled	Actual	Comment	Reference
В	-6	HIS	-	expression tag	UNP Q9PPB4
В	-5	HIS	-	expression tag	UNP Q9PPB4
В	-4	HIS	-	expression tag	UNP Q9PPB4
В	-3	HIS	-	expression tag	UNP Q9PPB4
В	-2	HIS	-	expression tag	UNP Q9PPB4
В	-1	GLY	-	expression tag	UNP Q9PPB4
В	0	SER	-	expression tag	UNP Q9PPB4
С	-11	MET	-	expression tag	UNP Q9PPB4
С	-10	ARG	-	expression tag	UNP Q9PPB4
С	-9	GLY	-	expression tag	UNP Q9PPB4
С	-8	SER	-	expression tag	UNP Q9PPB4
С	-7	HIS	-	expression tag	UNP Q9PPB4
С	-6	HIS	-	expression tag	UNP Q9PPB4
С	-5	HIS	_	expression tag	UNP Q9PPB4
С	-4	HIS	-	expression tag	UNP Q9PPB4
С	-3	HIS	_	expression tag	UNP Q9PPB4
С	-2	HIS	-	expression tag	UNP Q9PPB4
С	-1	GLY	-	expression tag	UNP Q9PPB4
С	0	SER	-	expression tag	UNP Q9PPB4
D	-11	MET	-	expression tag	UNP Q9PPB4
D	-10	ARG	_	expression tag	UNP Q9PPB4
D	-9	GLY	-	expression tag	UNP Q9PPB4
D	-8	SER	-	expression tag	UNP Q9PPB4
D	-7	HIS	-	expression tag	UNP Q9PPB4
D	-6	HIS	-	expression tag	UNP Q9PPB4
D	-5	HIS	-	expression tag	UNP Q9PPB4
D	-4	HIS	-	expression tag	UNP Q9PPB4
D	-3	HIS	-	expression tag	UNP Q9PPB4
D	-2	HIS	-	expression tag	UNP Q9PPB4
D	-1	GLY	-	expression tag	UNP Q9PPB4
D	0	SER	-	expression tag	UNP Q9PPB4
Е	-11	MET	-	expression tag	UNP Q9PPB4
Е	-10	ARG	-	expression tag	UNP Q9PPB4
Е	-9	GLY	-	expression tag	UNP Q9PPB4
Е	-8	SER	_	expression tag	UNP Q9PPB4
Е	-7	HIS	-	expression tag	UNP Q9PPB4
Е	-6	HIS	-	expression tag	UNP Q9PPB4
Е	-5	HIS	-	expression tag	UNP Q9PPB4
Е	-4	HIS	-	expression tag	UNP Q9PPB4
Е	-3	HIS	-	expression tag	UNP Q9PPB4
E	-2	HIS	-	expression tag	UNP Q9PPB4
E	-1	GLY	_	expression tag	UNP Q9PPB4



Continued from previous page...

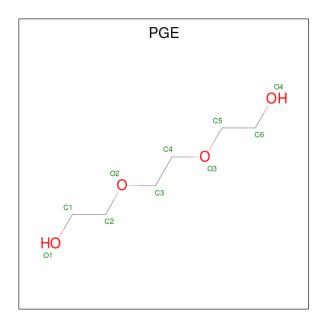
Chain	Residue	Modelled	Actual Comment		Reference
Е	0	SER	-	expression tag	UNP Q9PPB4
F	-11	MET	-	expression tag	UNP Q9PPB4
F	-10	ARG	-	expression tag	UNP Q9PPB4
F	-9	GLY	-	expression tag	UNP Q9PPB4
F	-8	SER	-	expression tag	UNP Q9PPB4
F	-7	HIS	-	expression tag	UNP Q9PPB4
F	-6	HIS	-	expression tag	UNP Q9PPB4
F	-5	HIS	-	expression tag	UNP Q9PPB4
F	-4	HIS	-	expression tag	UNP Q9PPB4
F	-3	HIS	-	expression tag	UNP Q9PPB4
F	-2	HIS	-	expression tag	UNP Q9PPB4
F	-1	GLY	-	expression tag	UNP Q9PPB4
F	0	SER	-	expression tag	UNP Q9PPB4

• Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	2	Total Mg 2 2	0	0
2	В	4	Total Mg 4 4	0	0
2	С	3	Total Mg 3 3	0	0
2	D	4	Total Mg 4 4	0	0
2	Ε	1	Total Mg 1 1	0	0
2	F	3	Total Mg 3 3	0	0

• Molecule 3 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: $C_6H_{14}O_4$) (labeled as "Ligand of Interest" by depositor).

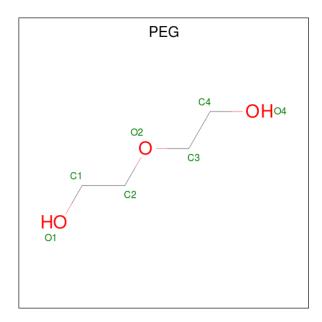




Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 10 6 4	0	0
3	В	1	Total C O 10 6 4	0	0
3	С	1	Total C O 10 6 4	0	0
3	D	1	Total C O 10 6 4	0	0
3	Е	1	Total C O 10 6 4	0	0
3	E	1	Total C O 10 6 4	0	0
3	F	1	Total C O 10 6 4	0	0

• Molecule 4 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: $C_4H_{10}O_3$) (labeled as "Ligand of Interest" by depositor).

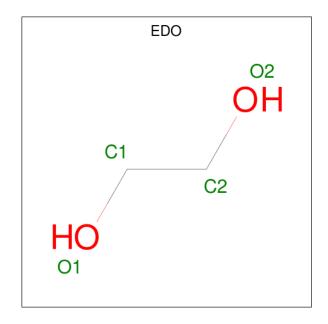




Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 7 4 3	0	0
4	В	1	Total C O 7 4 3	0	0
4	С	1	Total C O 7 4 3	0	0
4	D	1	Total C O 7 4 3	0	0
4	F	1	Total C O 7 4 3	0	0

 \bullet Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2)$ (labeled as "Ligand of Interest" by depositor).

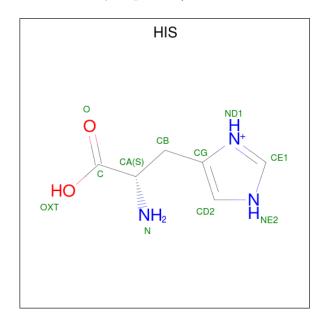




Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total C O 4 2 2	0	0
5	A	1	Total C O 4 2 2	0	0
5	A	1	Total C O 4 2 2	0	0
5	В	1	Total C O 4 2 2	0	0
5	В	1	Total C O 4 2 2	0	0
5	В	1	Total C O 4 2 2	0	0
5	С	1	Total C O 4 2 2	0	0
5	С	1	Total C O 4 2 2	0	0
5	D	1	Total C O 4 2 2	0	0
5	D	1	Total C O 4 2 2	0	0
5	Е	1	Total C O 4 2 2	0	0
5	F	1	Total C O 4 2 2	0	0
5	F	1	Total C O 4 2 2	0	0
5	F	1	Total C O 4 2 2	0	0



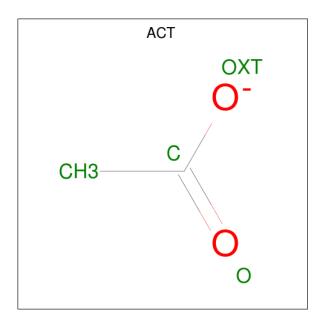
• Molecule 6 is HISTIDINE (three-letter code: HIS) (formula: $C_6H_{10}N_3O_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total C N O 11 6 3 2	0	0
6	В	1	Total C N O 11 6 3 2	0	0
6	С	1	Total C N O 11 6 3 2	0	0
6	D	1	Total C N O 11 6 3 2	0	0
6	E	1	Total C N O 11 6 3 2	0	0
6	F	1	Total C N O 11 6 3 2	0	0

• Molecule 7 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$) (labeled as "Ligand of Interest" by depositor).

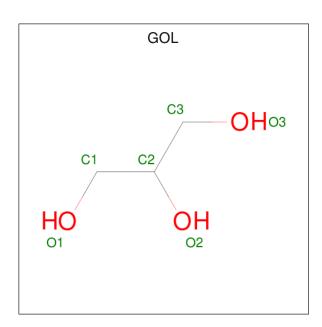




Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	В	1	Total C O 4 2 2	0	0
7	D	1	Total C O 4 2 2	0	0
7	E	1	Total C O 4 2 2	0	0
7	E	1	Total C O 4 2 2	0	0
7	F	1	Total C O 4 2 2	0	0

• Molecule 8 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$) (labeled as "Ligand of Interest" by depositor).





Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	F	1	Total C O 6 3 3	0	0

• Molecule 9 is water.

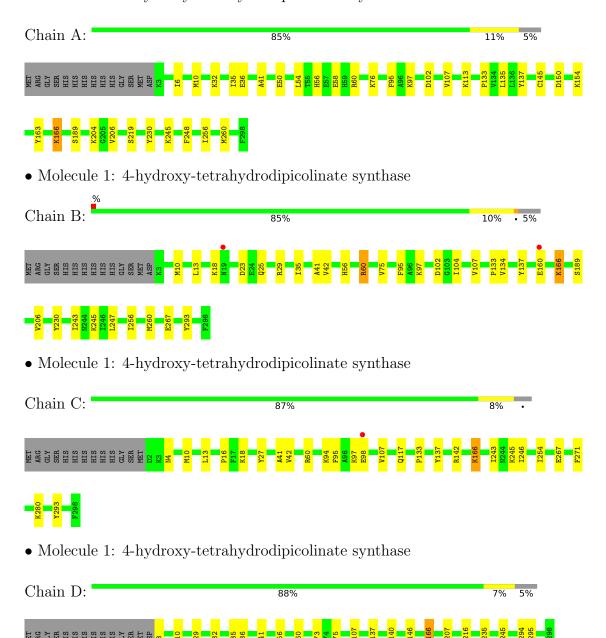
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	A	139	Total O 139 139	0	0
9	В	131	Total O 131 131	0	0
9	С	133	Total O 133 133	0	0
9	D	126	Total O 126 126	0	0
9	Е	96	Total O 96 96	0	0
9	F	97	Total O 97 97	0	0



3 Residue-property plots (i)

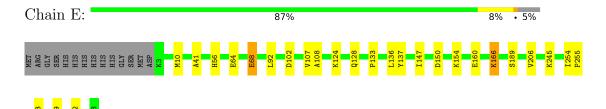
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: 4-hydroxy-tetrahydrodipicolinate synthase

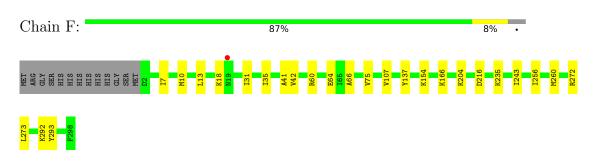




• Molecule 1: 4-hydroxy-tetrahydrodipicolinate synthase



• Molecule 1: 4-hydroxy-tetrahydrodipicolinate synthase





4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants	85.31Å 231.72Å 202.35Å	Donositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.36 - 2.28	Depositor
Resolution (A)	46.36 - 2.28	EDS
% Data completeness	98.7 (46.36-2.28)	Depositor
(in resolution range)	95.6 (46.36-2.28)	EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	2.74 (at 2.29Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
D D.	0.170 , 0.208	Depositor
R, R_{free}	0.170 , 0.205	DCC
R_{free} test set	4518 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	31.2	Xtriage
Anisotropy	0.271	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.35, 39.4	EDS
L-test for twinning ²	$ < L >=0.49, < L^2>=0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	14653	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 2.55% of the height of the origin peak. No significant pseudotranslation is detected.

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



¹Intensities estimated from amplitudes.

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: KPI, PGE, EDO, MG, GOL, ACT, 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	Bo	nd lengths	Bond angles	
MIOI	Chain	RMSZ	# Z > 5	RMSZ	# Z >5
1	A	0.58	2/2305~(0.1%)	0.51	0/3118
1	В	0.40	0/2293	0.50	0/3102
1	С	0.38	0/2302	0.55	0/3115
1	D	0.41	0/2309	0.51	0/3121
1	Е	0.38	0/2297	0.49	0/3107
1	F	0.44	0/2299	0.53	0/3111
All	All	0.44	$2/13805 \ (0.0\%)$	0.51	0/18674

All (2) bond length outliers are listed below:

Mo	Chain	Res	Type	Atoms	\mathbf{Z}	$\operatorname{Observed}(\operatorname{\AA})$	$Ideal(\AA)$
1	A	58	GLU	CD-OE2	-5.67	1.19	1.25
1	A	50	GLU	CD-OE2	-5.51	1.19	1.25

There are no bond angle outliers.

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	2281	0	2304	23	0
1	В	2269	0	2292	25	0
1	С	2278	0	2296	20	0



 $Continued\ from\ previous\ page...$

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	2285	0	2323	15	0
1	Е	2273	0	2300	19	0
1	F	2275	0	2296	17	0
2	A	2	0	0	0	0
2	В	4	0	0	0	0
2	С	3	0	0	0	0
2	D	4	0	0	0	0
2	Е	1	0	0	0	0
2	F	3	0	0	0	0
3	A	10	0	14	3	0
3	В	10	0	14	0	0
3	С	10	0	14	0	0
3	D	10	0	14	0	0
3	Е	20	0	28	5	0
3	F	10	0	14	0	0
4	A	7	0	10	2	0
4	В	7	0	10	1	0
4	С	7	0	10	0	0
4	D	7	0	10	2	0
4	F	7	0	10	1	0
5	A	12	0	18	2	0
5	В	12	0	18	0	0
5	С	8	0	12	0	0
5	D	8	0	12	0	0
5	Е	4	0	6	0	0
5	F	12	0	18	3	0
6	A	11	0	6	1	0
6	В	11	0	6	0	0
6	С	11	0	6	0	0
6	D	11	0	6	1	0
6	E	11	0	6	0	0
6	F	11	0	6	0	0
7	В	4	0	3	3	0
7	D	4	0	3	0	0
7	Е	8	0	6	0	0
7	F	4	0	3	0	0
8	F	6	0	8	0	0
9	A	139	0	0	0	0
9	В	131	0	0	2	0
9	С	133	0	0	1	0
9	D	126	0	0	1	0
9	Е	96	0	0	1	0



Continued from previous page...

Mol	Chain	Non-H	H(model)	$\mathbf{H}(\mathbf{added})$	Clashes	Symm-Clashes
9	F	97	0	0	0	0
All	All	14653	0	14102	116	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 4.

The worst 5 of 116 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	$egin{aligned} ext{Clash} \ ext{overlap } (ext{Å}) \end{aligned}$
1:E:64:GLU:O	1:E:68:GLU:HG2	1.72	0.90
1:B:25:GLN:H	7:B:307:ACT:H3	1.44	0.81
1:C:60:ARG:HB2	1:C:95:PHE:HZ	1.44	0.81
1:E:124:LYS:HE2	1:E:128:GLN:OE1	1.82	0.78
1:C:60:ARG:HB2	1:C:95:PHE:CZ	2.18	0.78

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	Percentile	\mathbf{s}
1	A	294/310~(95%)	288 (98%)	6 (2%)	0	100 100	
1	В	293/310 (94%)	287 (98%)	6 (2%)	0	100 100	
1	С	$294/310 \ (95\%)$	288 (98%)	6 (2%)	0	100 100	
1	D	294/310 (95%)	288 (98%)	6 (2%)	0	100 100	
1	E	293/310 (94%)	287 (98%)	6 (2%)	0	100 100	
1	F	294/310 (95%)	288 (98%)	6 (2%)	0	100 100	
All	All	1762/1860~(95%)	1726 (98%)	36 (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	Perce	ntiles
1	A	247/260 (95%)	244 (99%)	3 (1%)	71	82
1	В	246/260 (95%)	244 (99%)	2 (1%)	81	90
1	C	246/260 (95%)	245 (100%)	1 (0%)	91	95
1	D	249/260 (96%)	246 (99%)	3 (1%)	71	82
1	E	246/260 (95%)	245 (100%)	1 (0%)	91	95
1	F	246/260 (95%)	246 (100%)	0	100	100
All	All	1480/1560 (95%)	1470 (99%)	10 (1%)	84	91

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

Mol	Chain	Res	Type
1	D	235	LYS
1	D	294	LYS
1	Е	68	GLU
1	В	60	ARG
1	В	97	LYS

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)

6 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	Mol Type Cl		Res	Link	Bo	ond leng	$ ag{ths}$	В	ond ang	les
IVIOI	туре	Chain	rtes	LILIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	KPI	С	166	1	11,13,14	2.21	3 (27%)	10,15,17	4.02	6 (60%)
1	KPI	F	166	1	11,13,14	1.15	2 (18%)	10,15,17	2.35	4 (40%)
1	KPI	D	166	1	11,13,14	1.49	2 (18%)	10,15,17	3.42	6 (60%)
1	KPI	Е	166	1	11,13,14	0.91	0	10,15,17	2.41	5 (50%)
1	KPI	A	166	1	11,13,14	1.52	2 (18%)	10,15,17	3.64	6 (60%)
1	KPI	В	166	1	11,13,14	0.94	0	10,15,17	3.03	5 (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	KPI	С	166	1	-	3/13/14/16	-
1	KPI	F	166	1	-	3/13/14/16	-
1	KPI	D	166	1	-	3/13/14/16	-
1	KPI	Ε	166	1	-	0/13/14/16	-
1	KPI	A	166	1	-	6/13/14/16	-
1	KPI	В	166	1	-	9/13/14/16	-

The worst 5 of 9 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	${f Z}$	$\operatorname{Observed}(\text{\AA})$	$\operatorname{Ideal}(ext{\AA})$
1	С	166	KPI	O2-CX2	5.32	1.36	1.22
1	A	166	KPI	O-C	4.13	1.36	1.19
1	С	166	KPI	O-C	4.10	1.36	1.19
1	D	166	KPI	O-C	4.07	1.36	1.19
1	С	166	KPI	O1-CX2	-2.27	1.24	1.30

The worst 5 of 32 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^o)$	$\mathrm{Ideal}(^{o})$
1	С	166	KPI	C1-CX1-CX2	-9.15	109.28	118.17
1	A	166	KPI	C1-CX1-CX2	-8.95	109.47	118.17
1	D	166	KPI	C1-CX1-CX2	-8.18	110.21	118.17
1	В	166	KPI	CX2-CX1-NZ	6.08	129.78	114.98



Continued from previous page...

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\mathbf{Observed}(^{o})$	$Ideal(^{o})$
1	A	166	KPI	O2-CX2-CX1	5.29	128.13	121.38

There are no chirality outliers.

5 of 24 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	166	KPI	C1-CX1-NZ-CE
1	A	166	KPI	NZ-CX1-CX2-O1
1	A	166	KPI	NZ-CX1-CX2-O2
1	A	166	KPI	C1-CX1-CX2-O1
1	A	166	KPI	C1-CX1-CX2-O2

There are no ring outliers.

5 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	С	166	KPI	2	0
1	D	166	KPI	3	0
1	Е	166	KPI	1	0
1	A	166	KPI	1	0
1	В	166	KPI	2	0

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 55 ligands modelled in this entry, 17 are monoatomic - leaving 38 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 Chai	Chain	Chain Res		Pog I;	Res Link	Bond lengths				Bond angles		
		туре	Chain	nes	Lilik	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2		
	7	ACT	F	306	-	3,3,3	1.16	0	3,3,3	1.54	0		



					B	ond leng	orths	P	ond ang	rles
Mol	Type	Chain	Res	Link	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	EDO	В	309	_	3,3,3	0.45	0	2,2,2	0.40	0
$\frac{}{6}$	HIS	A	308	_	6,11,11	0.97	1 (16%)	7,14,14	1.56	2 (28%)
6	HIS	В	311	_	6,11,11	0.97	1 (16%)	7,14,14	1.57	2 (28%)
4	PEG	D	306	_	6,6,6	0.50	0	5,5,5	0.22	0
3	PGE	A	303	_	9,9,9	0.30	0	8,8,8	0.22	0
3	PGE	E	303	-	9,9,9	0.30	0	8,8,8	0.28	0
5	EDO	C	306	_	3,3,3	0.30	0	2,2,2	0.27	0
$\frac{5}{6}$	HIS	D	310							
				-	6,11,11	0.99		7,14,14	1.58	(' ' ' ')
4	PEG	A	304	_	6,6,6	0.48	0	5,5,5	0.32	0
6	HIS	Е	307	-	6,11,11	0.96	1 (16%)	7,14,14	1.55	3 (42%)
7	ACT	В	307	-	3,3,3	0.94	0	3,3,3	0.74	0
5	EDO	F	309	-	3,3,3	0.46	0	2,2,2	0.32	0
5	EDO	F	307	-	3,3,3	0.46	0	2,2,2	0.37	0
5	EDO	A	306	-	3,3,3	0.45	0	2,2,2	0.38	0
3	PGE	D	305	-	9,9,9	0.30	0	8,8,8	0.29	0
5	EDO	E	306	_	3,3,3	0.45	0	2,2,2	0.28	0
5	EDO	С	307	_	3,3,3	0.46	0	2,2,2	0.33	0
5	EDO	D	309	-	3,3,3	0.46	0	2,2,2	0.29	0
6	HIS	F	310	-	6,11,11	0.98	1 (16%)	7,14,14	1.56	2 (28%)
8	GOL	F	311	-	5,5,5	1.04	0	5,5,5	0.88	0
4	PEG	В	306	-	6,6,6	0.49	0	5,5,5	0.27	0
7	ACT	D	307	-	3,3,3	1.31	0	3,3,3	1.52	0
3	PGE	С	304	-	9,9,9	0.30	0	8,8,8	0.34	0
5	EDO	F	308	-	3,3,3	0.45	0	2,2,2	0.33	0
7	ACT	Е	305	-	3,3,3	1.31	0	3,3,3	1.53	0
3	PGE	F	304	-	9,9,9	0.31	0	8,8,8	0.30	0
5	EDO	A	305	-	3,3,3	0.44	0	2,2,2	0.41	0
4	PEG	F	305	-	6,6,6	0.48	0	5,5,5	0.27	0
6	HIS	С	308	-	6,11,11	0.98	1 (16%)	7,14,14	1.56	3 (42%)
5	EDO	В	308	-	3,3,3	0.44	0	2,2,2	0.46	0
4	PEG	С	305	-	6,6,6	0.48	0	5,5,5	0.29	0
3	PGE	В	305	-	9,9,9	0.29	0	8,8,8	0.33	0
5	EDO	D	308	-	3,3,3	0.46	0	2,2,2	0.30	0
5	EDO	A	307	-	3,3,3	0.45	0	2,2,2	0.28	0
7	ACT	E	304	-	3,3,3	1.29	0	3,3,3	1.51	0
3	PGE	Е	302	-	9,9,9	0.29	0	8,8,8	0.31	0
5	EDO	В	310	-	3,3,3	0.46	0	2,2,2	0.31	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
5	EDO	В	309	_	-	0/1/1/1	-
6	HIS	A	308	-	-	1/8/8/8	0/1/1/1
6	HIS	В	311	-	-	0/8/8/8	0/1/1/1
4	PEG	D	306	-	-	0/4/4/4	-
3	PGE	A	303	-	-	4/7/7/7	-
3	PGE	Е	303	-	-	4/7/7/7	-
5	EDO	С	306	-	-	0/1/1/1	-
6	HIS	D	310	-	-	1/8/8/8	0/1/1/1
4	PEG	A	304	-	-	1/4/4/4	-
6	HIS	E	307	_	-	1/8/8/8	0/1/1/1
5	EDO	F	309	_	-	0/1/1/1	-
5	EDO	F	307	_	-	0/1/1/1	-
5	EDO	A	306	_	-	0/1/1/1	-
3	PGE	D	305	_	-	0/7/7/7	-
5	EDO	E	306	-	-	1/1/1/1	-
5	EDO	С	307	_	-	0/1/1/1	-
5	EDO	D	309	_	-	0/1/1/1	-
6	HIS	F	310	-	-	1/8/8/8	0/1/1/1
8	GOL	F	311	-	-	3/4/4/4	-
4	PEG	В	306	_	-	2/4/4/4	-
3	PGE	С	304	-	-	0/7/7/7	-
5	EDO	F	308	-	-	0/1/1/1	-
3	PGE	F	304	-	-	0/7/7/7	-
5	EDO	A	305	_	-	1/1/1/1	-
4	PEG	F	305	-	-	1/4/4/4	-
6	HIS	С	308	-	-	2/8/8/8	0/1/1/1
5	EDO	В	308	_	-	1/1/1/1	-
4	PEG	С	305	-	-	4/4/4/4	-
3	PGE	В	305	-	-	0/7/7/7	-
5	EDO	D	308	-	-	0/1/1/1	-
5	EDO	A	307	-	-	0/1/1/1	-
3	PGE	Е	302	-	-	5/7/7/7	-
5	EDO	В	310	-	-	0/1/1/1	-

The worst 5 of 6 bond length outliers are listed below:

Mol	Chain	Res	<i>U</i> 1	Atoms	\mathbf{Z}	$\operatorname{Observed}(\text{\AA})$	Ideal(Å)
6	D	310	HIS	OXT-C	-2.16	1.23	1.30
6	F	310	HIS	OXT-C	-2.14	1.23	1.30
6	С	308	HIS	OXT-C	-2.13	1.23	1.30
6	A	308	HIS	OXT-C	-2.12	1.23	1.30
6	В	311	HIS	OXT-C	-2.09	1.23	1.30



The worst	5	of	14	bond	angle	outliers	are	listed	below:
I IIC WOIDU	$\mathbf{\mathcal{I}}$	\circ		Ollu	WII SIC	Outiloid	COL C	IIDUCA	DOIOW.

Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$Ideal(^{o})$
6	D	310	HIS	OXT-C-O	-2.72	117.91	124.09
6	F	310	HIS	OXT-C-O	-2.68	118.01	124.09
6	В	311	HIS	OXT-C-O	-2.63	118.11	124.09
6	С	308	HIS	OXT-C-O	-2.63	118.13	124.09
6	A	308	HIS	OXT-C-O	-2.62	118.14	124.09

There are no chirality outliers.

5 of 33 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	308	HIS	CA-CB-CG-ND1
6	С	308	HIS	CA-CB-CG-ND1
6	D	310	HIS	CA-CB-CG-ND1
6	F	310	HIS	CA-CB-CG-ND1
8	F	311	GOL	O1-C1-C2-C3

There are no ring outliers.

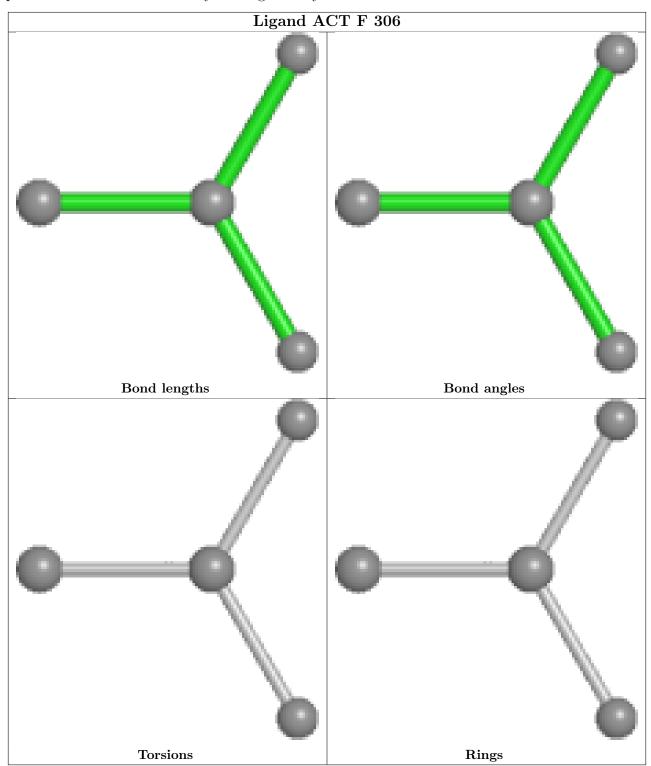
13 monomers are involved in 23 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	308	HIS	1	0
4	D	306	PEG	2	0
3	A	303	PGE	3	0
3	Е	303	PGE	2	0
6	D	310	HIS	1	0
4	A	304	PEG	2	0
7	В	307	ACT	3	0
4	В	306	PEG	1	0
5	F	308	EDO	3	0
5	A	305	EDO	1	0
4	F	305	PEG	1	0
5	A	307	EDO	1	0
3	Е	302	PGE	3	0

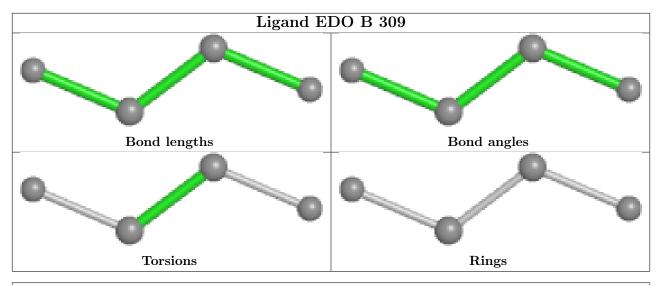
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

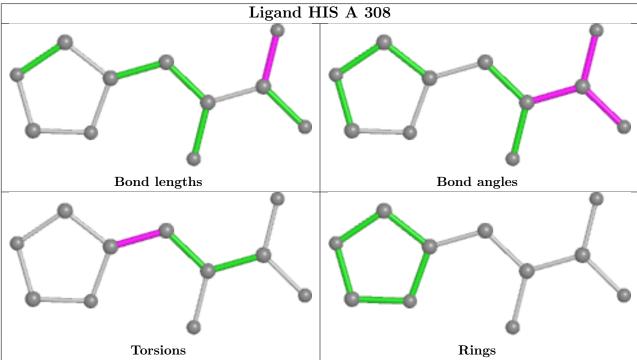


in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

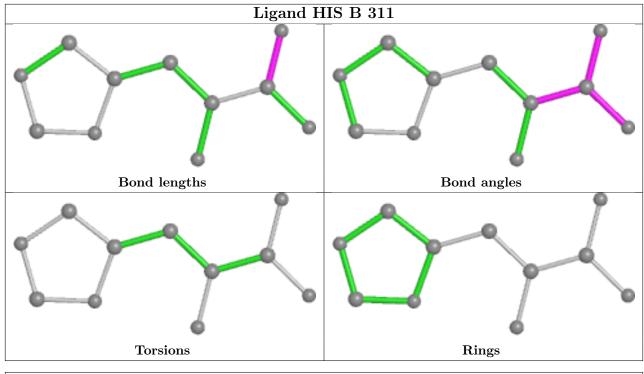


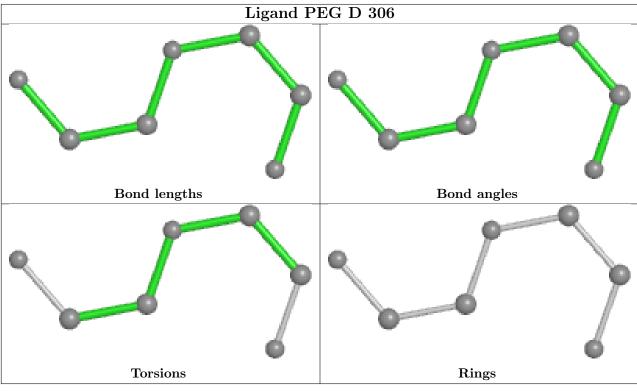


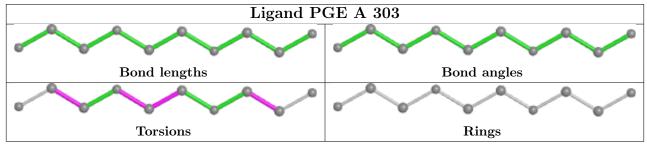




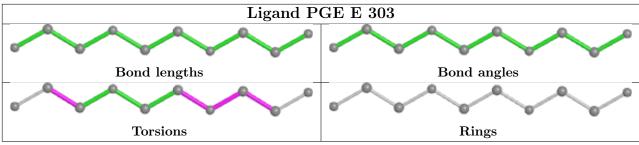


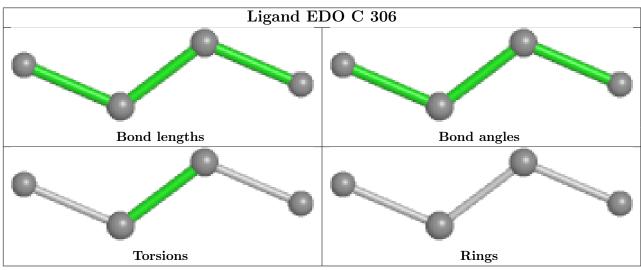


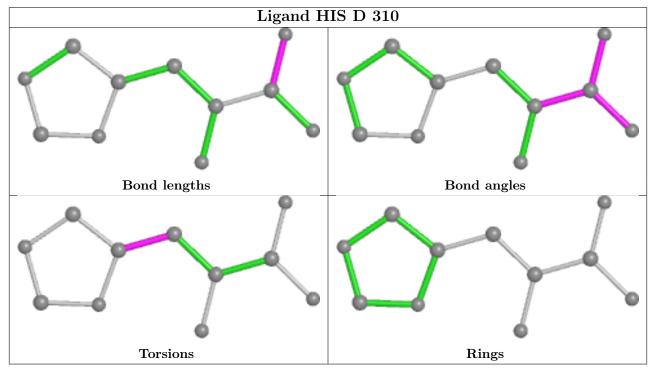




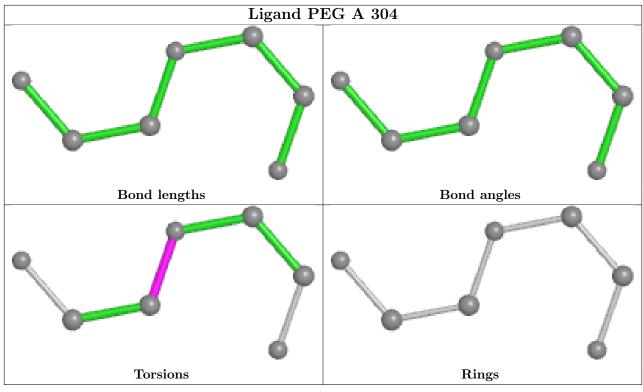


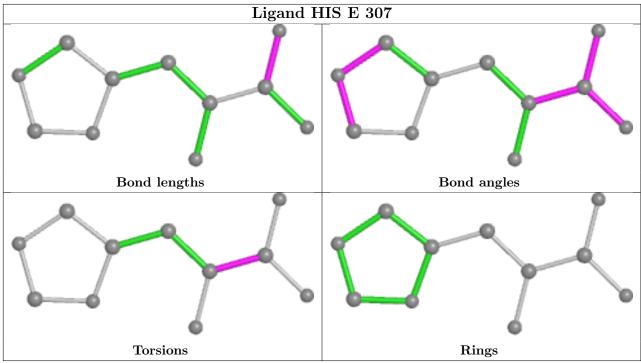




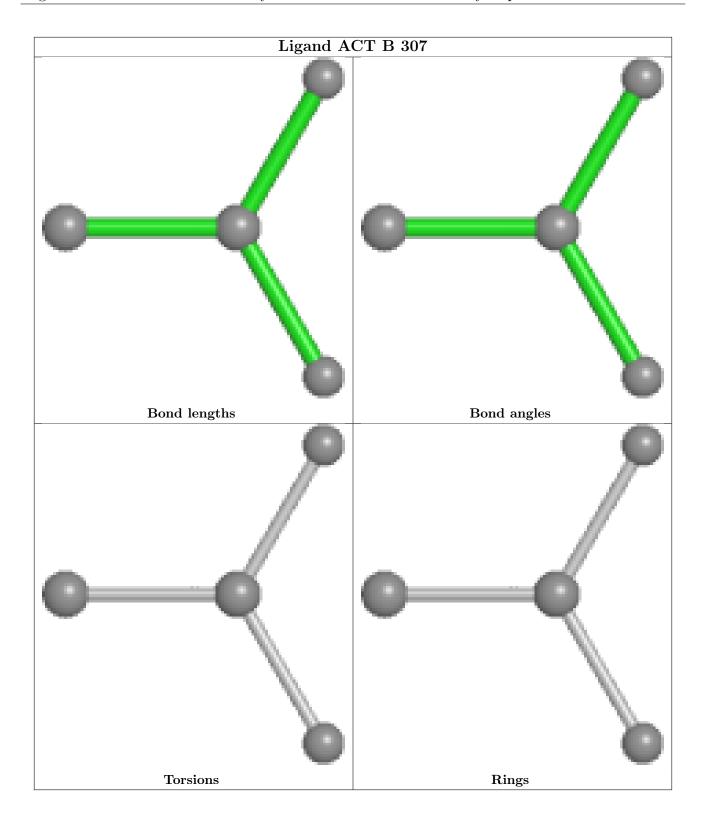




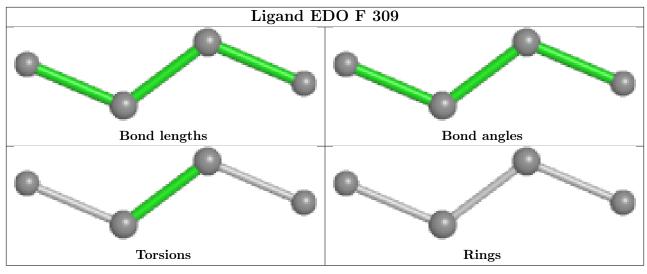


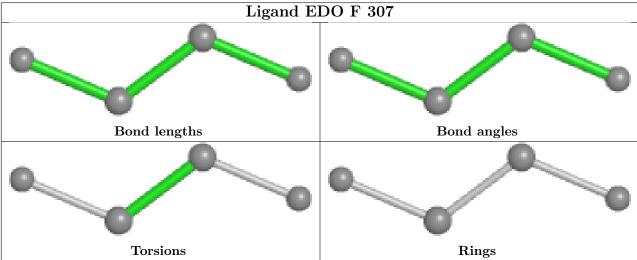


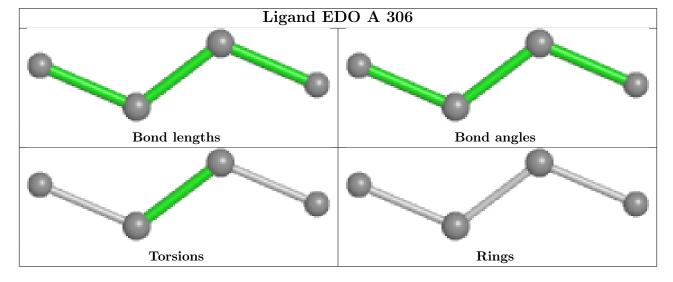




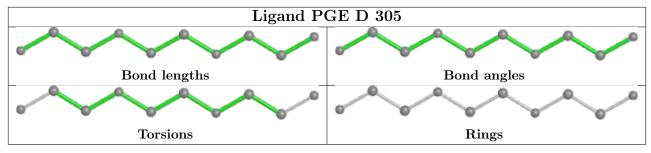


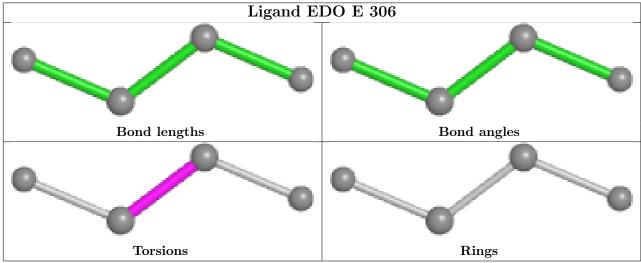


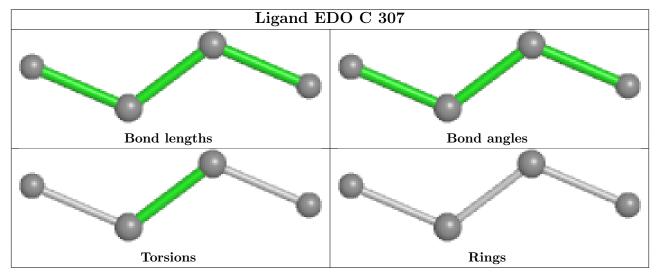




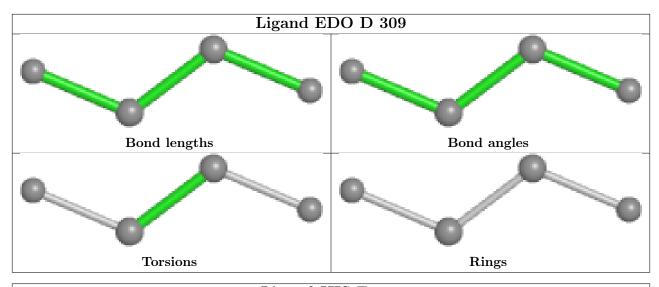


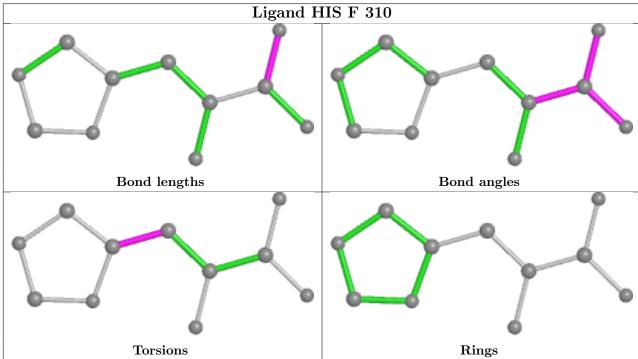




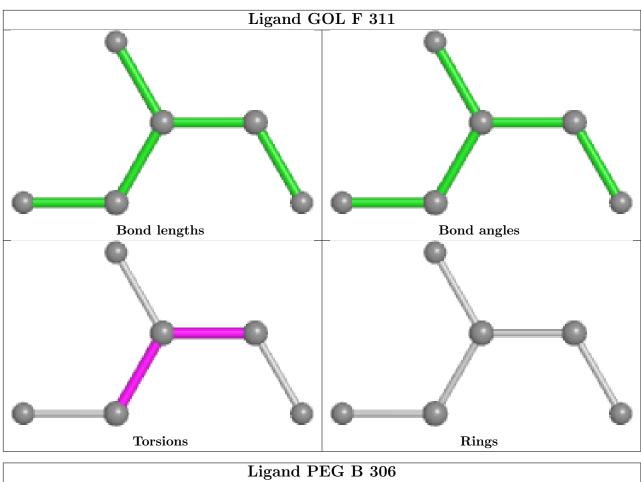


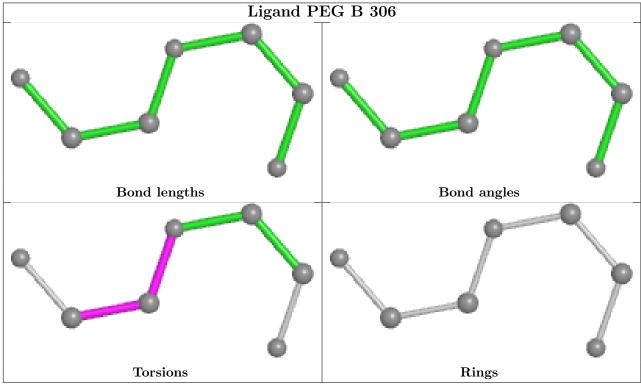




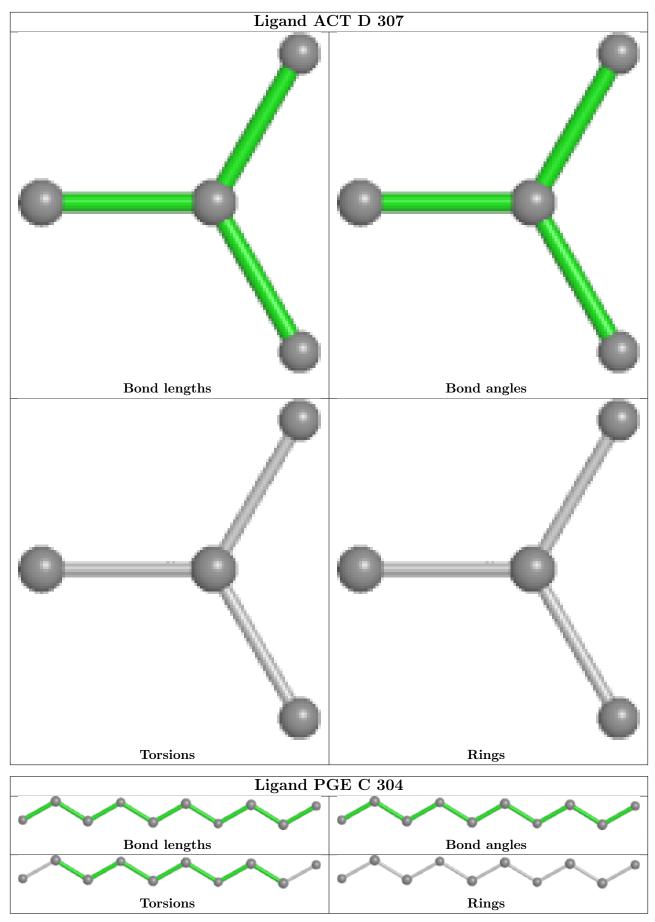




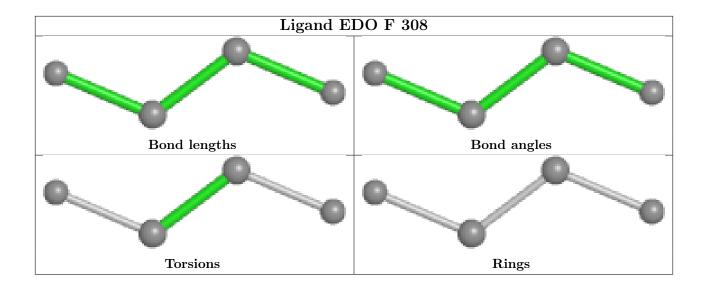




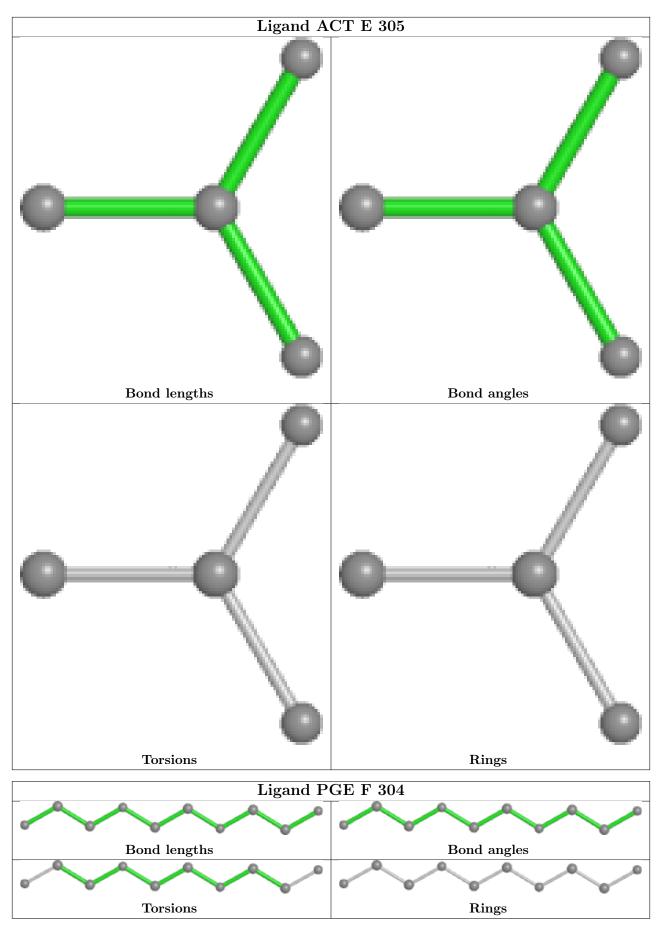




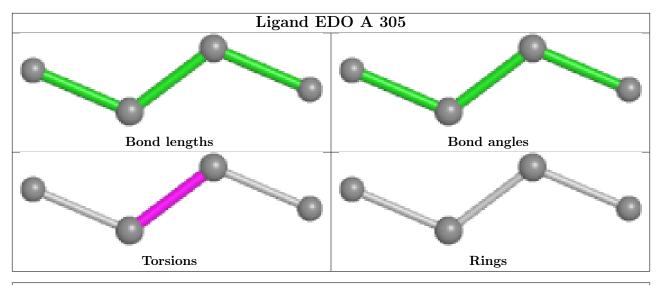


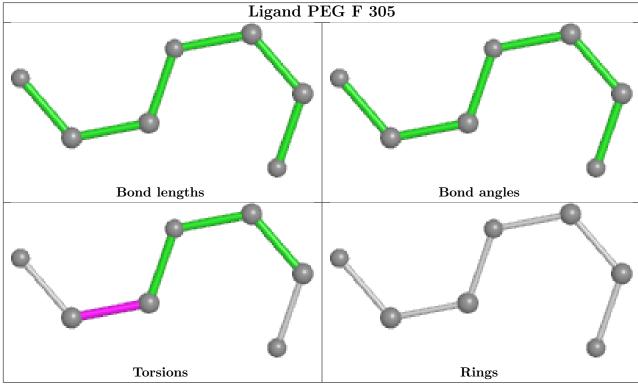




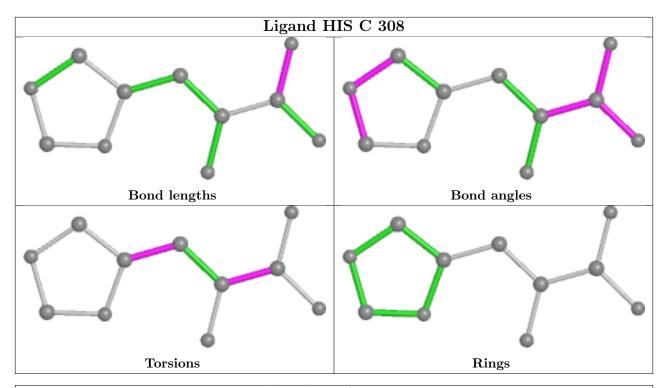


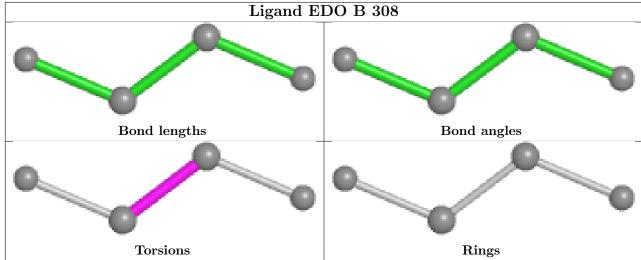




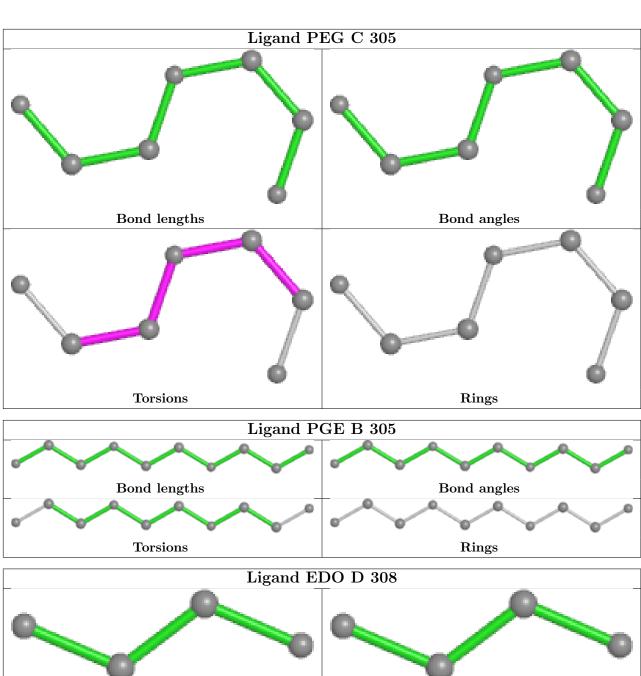


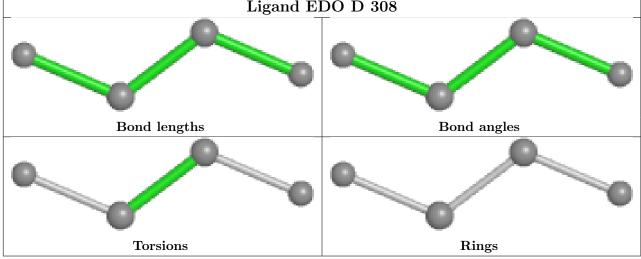




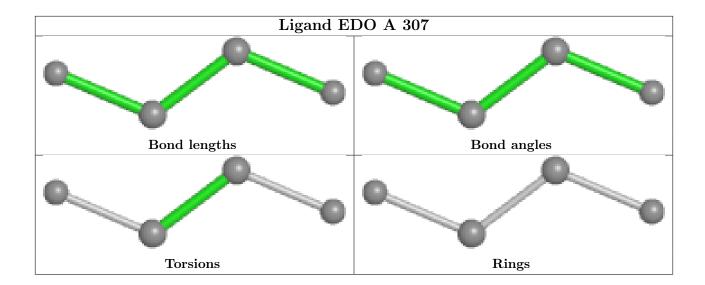




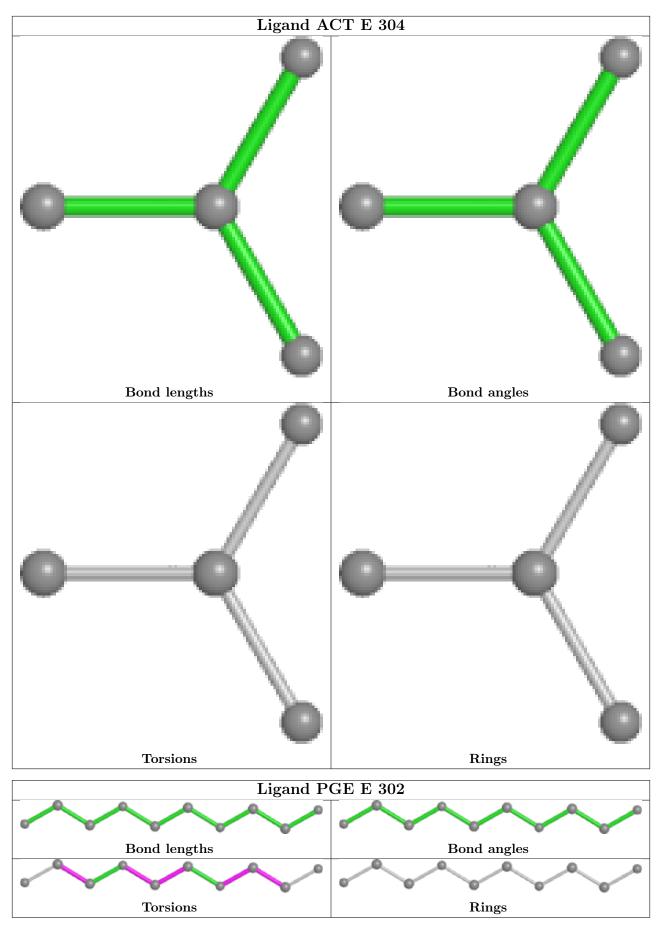




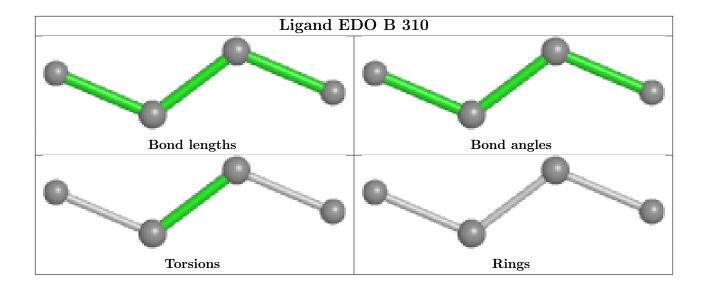












5.7 Other polymers (i)

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, 95^{th} 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>	$\# \mathrm{RSRZ}{>}2$	$OWAB(A^2)$	Q < 0.9
1	A	295/310~(95%)	-0.34	0 100 100	25, 31, 50, 67	0
1	В	295/310~(95%)	-0.36	2 (0%) 87 90	24, 32, 51, 62	0
1	С	296/310 (95%)	-0.31	1 (0%) 94 95	26, 32, 51, 69	0
1	D	295/310~(95%)	-0.28	0 100 100	24, 34, 52, 68	0
1	E	295/310~(95%)	-0.34	0 100 100	25, 38, 56, 74	0
1	F	296/310~(95%)	-0.16	1 (0%) 94 95	27, 40, 62, 87	0
All	All	1772/1860 (95%)	-0.30	4 (0%) 95 96	24, 35, 56, 87	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	В	19	ASN	2.5
1	F	19	ASN	2.4
1	В	160	GLU	2.3
1	С	98	GLU	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, 95^{th} 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	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
1	KPI	В	166	14/15	0.79	0.21	31,37,47,48	0
1	KPI	Ε	166	14/15	0.80	0.20	30,35,58,59	0
1	KPI	A	166	14/15	0.81	0.23	23,27,50,53	0
1	KPI	С	166	14/15	0.85	0.19	26,31,45,48	0
1	KPI	F	166	14/15	0.87	0.20	32,37,57,65	0

Continued on next page...



Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q < 0.9
1	KPI	D	166	14/15	0.89	0.21	25,31,57,57	0

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (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, 95^{th} 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	$\operatorname{B-factors}(\mathring{\mathbf{A}}^2)$	Q < 0.9
5	EDO	F	309	4/4	0.73	0.17	55,58,62,63	0
5	EDO	В	309	4/4	0.78	0.25	60,60,61,62	0
4	PEG	A	304	7/7	0.78	0.18	56,56,62,62	0
4	PEG	В	306	7/7	0.80	0.21	53,59,66,67	0
7	ACT	В	307	4/4	0.80	0.44	20,20,20,20	0
5	EDO	D	309	4/4	0.81	0.60	48,54,58,61	0
5	EDO	В	308	4/4	0.81	0.45	53,54,57,57	0
5	EDO	A	306	4/4	0.81	0.28	45,54,54,57	0
7	ACT	D	307	4/4	0.81	0.22	51,51,54,60	0
7	ACT	Е	305	4/4	0.81	0.23	57,62,62,64	0
5	EDO	В	310	4/4	0.82	0.26	47,53,56,60	0
7	ACT	F	306	4/4	0.82	0.24	44,52,55,59	0
5	EDO	A	307	4/4	0.83	0.41	45,51,54,57	0
4	PEG	С	305	7/7	0.85	0.23	65,68,69,71	0
3	PGE	Е	302	10/10	0.85	0.19	57,59,61,61	0
3	PGE	E	303	10/10	0.85	0.15	58,59,63,63	0
5	EDO	С	306	4/4	0.86	0.18	43,48,48,48	0
5	EDO	С	307	4/4	0.86	0.14	52,52,54,55	0
2	MG	E	301	1/1	0.86	0.15	55,55,55,55	0
5	EDO	F	307	4/4	0.86	0.12	61,61,61,62	0
5	EDO	F	308	4/4	0.86	0.21	52,54,56,59	0
4	PEG	D	306	7/7	0.87	0.18	45,49,57,59	0
2	MG	D	304	1/1	0.87	0.10	61,61,61,61	0
3	PGE	F	304	10/10	0.88	0.18	40,51,52,56	0
5	EDO	Е	306	4/4	0.88	0.27	44,48,48,53	0
2	MG	В	303	1/1	0.88	0.06	40,40,40,40	0
3	PGE	D	305	10/10	0.89	0.18	42,47,50,51	0

Continued on next page...



Continued from previous page...

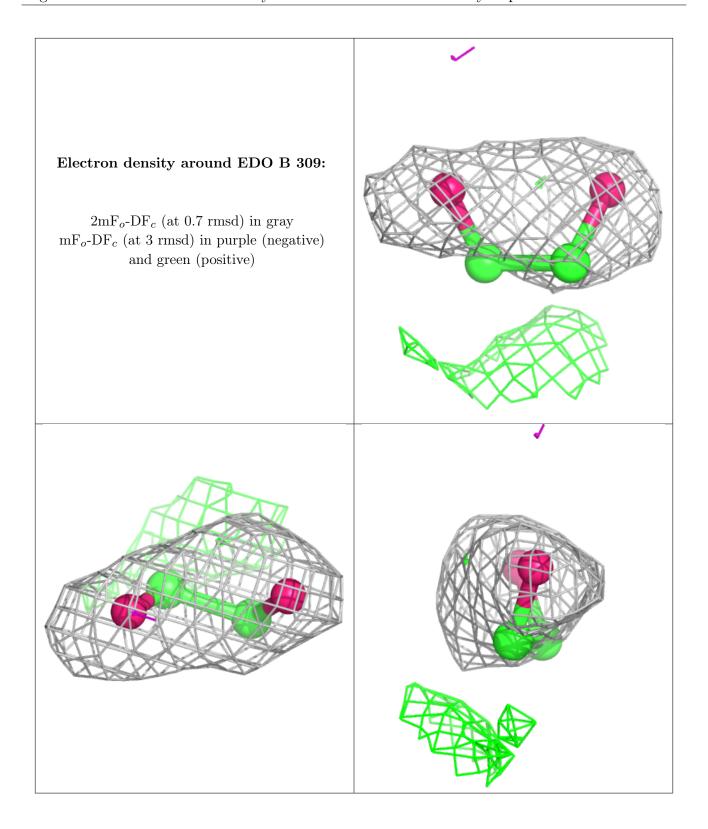
Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q < 0.9
3	PGE	С	304	10/10	0.89	0.14	52,53,59,60	0
2	MG	D	302	1/1	0.90	0.51	53,53,53,53	0
3	PGE	A	303	10/10	0.90	0.15	48,53,56,58	0
2	MG	В	301	1/1	0.90	0.12	40,40,40,40	0
2	MG	С	302	1/1	0.91	0.13	48,48,48,48	0
5	EDO	A	305	4/4	0.91	0.25	44,44,46,47	0
2	MG	В	302	1/1	0.92	0.19	61,61,61,61	0
2	MG	A	301	1/1	0.92	0.16	37,37,37,37	0
5	EDO	D	308	4/4	0.92	0.17	56,57,57,59	0
7	ACT	Е	304	4/4	0.93	0.23	52,55,56,58	0
2	MG	С	303	1/1	0.93	0.16	47,47,47,47	0
2	MG	F	301	1/1	0.93	0.49	54,54,54,54	0
2	MG	F	303	1/1	0.94	0.24	55,55,55,55	0
4	PEG	F	305	7/7	0.94	0.13	53,55,55,58	0
3	PGE	В	305	10/10	0.94	0.17	45,50,55,56	0
8	GOL	F	311	6/6	0.94	0.21	35,44,47,51	0
2	MG	С	301	1/1	0.95	0.18	35,35,35,35	0
2	MG	D	303	1/1	0.95	0.20	37,37,37,37	0
2	MG	D	301	1/1	0.96	0.22	34,34,34,34	0
6	HIS	A	308	11/11	0.96	0.12	30,32,36,38	0
6	HIS	В	311	11/11	0.97	0.11	31,32,36,36	0
6	HIS	С	308	11/11	0.97	0.12	29,31,35,37	0
6	HIS	Е	307	11/11	0.97	0.11	29,32,34,35	0
2	MG	В	304	1/1	0.98	0.13	34,34,34,34	0
6	HIS	F	310	11/11	0.98	0.09	29,32,34,34	0
2	MG	F	302	1/1	0.98	0.16	41,41,41,41	0
6	HIS	D	310	11/11	0.98	0.12	32,32,34,36	0
2	MG	A	302	1/1	0.99	0.12	28,28,28,28	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 F 309: $2 \mathrm{mF}_o\text{-}\mathrm{DF}_c$ (at 0.7 rmsd) in gray ${ m mF}_o{ m -DF}_c$ (at 3 rmsd) in purple (negative) and green (positive)



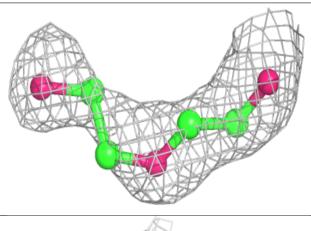


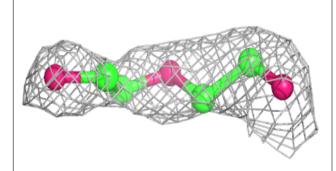


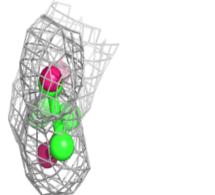
Electron density around PEG A 304: 2mF_o-DF_c (at 0.7 rmsd) in gray mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

Electron density around PEG B 306:

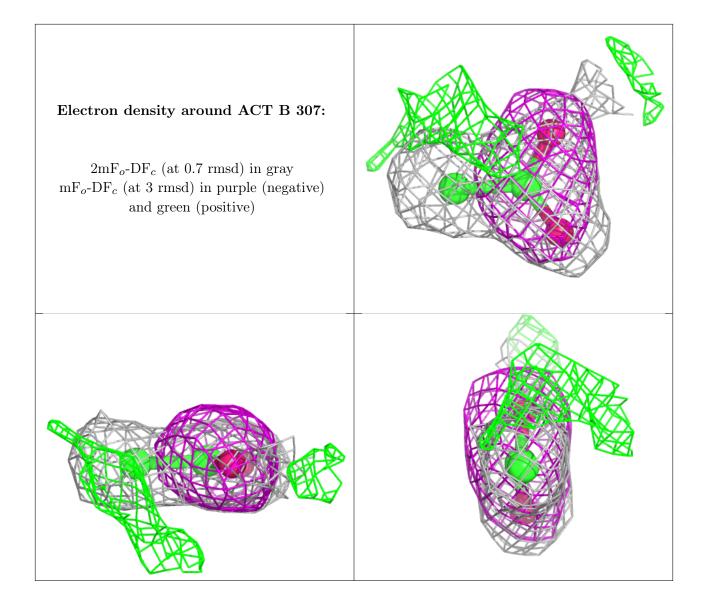
 $2 {\rm mF}_o\text{-}{\rm DF}_c$ (at 0.7 rmsd) in gray ${\rm mF}_o\text{-}{\rm DF}_c$ (at 3 rmsd) in purple (negative) and green (positive)



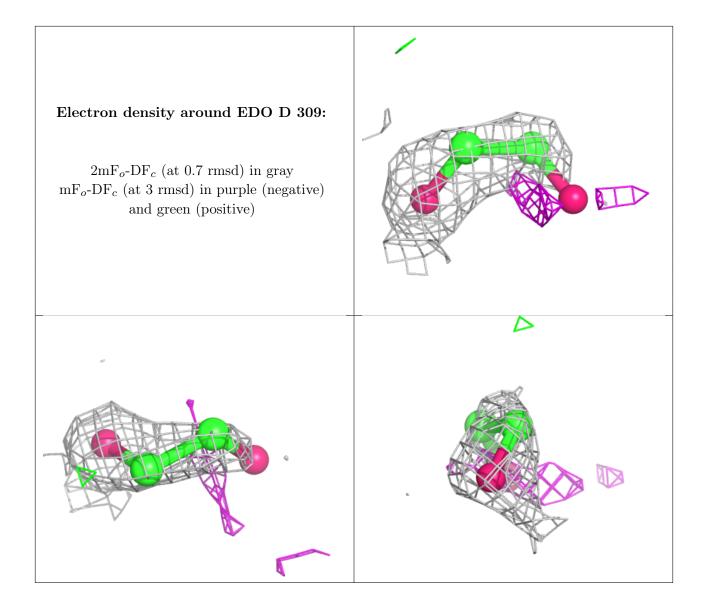














Electron density around EDO B 308: $2 \mathrm{mF}_o\text{-}\mathrm{DF}_c$ (at 0.7 rmsd) in gray ${\rm mF}_o\text{-}{\rm DF}_c$ (at 3 rmsd) in purple (negative) and green (positive)



Electron density around EDO A 306: $2 \mathrm{mF}_o\text{-}\mathrm{DF}_c$ (at 0.7 rmsd) in gray $\mathrm{mF}_o\text{-}\mathrm{DF}_c$ (at 3 rmsd) in purple (negative) and green (positive)

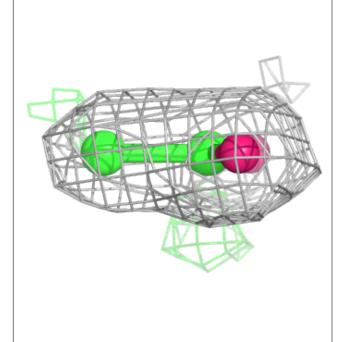


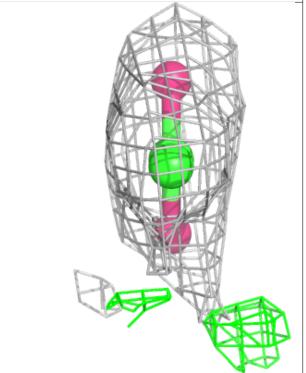


Electron density around ACT E 305:

 $2 {
m mF}_o {
m -DF}_c$ (at 0.7 rmsd) in gray ${
m mF}_o {
m -DF}_c$ (at 3 rmsd) in purple (negative) and green (positive)



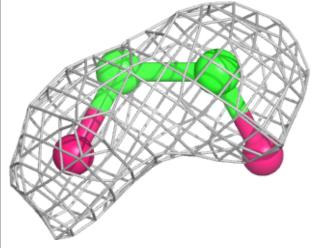


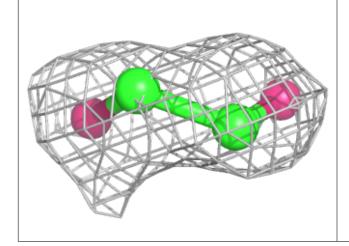


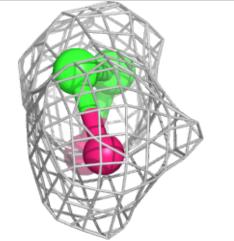


Electron density around EDO B 310:

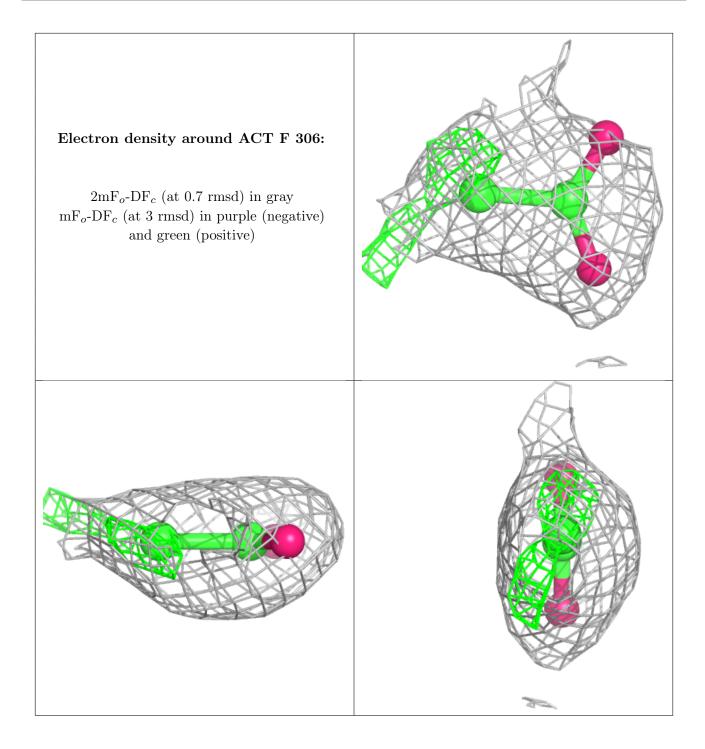
 $2 \text{mF}_o\text{-DF}_c$ (at 0.7 rmsd) in gray $\text{mF}_o\text{-DF}_c$ (at 3 rmsd) in purple (negative) and green (positive)



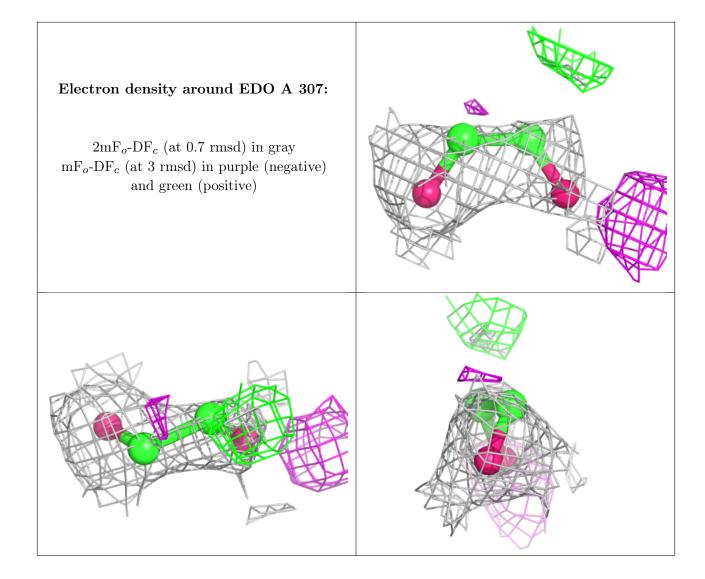




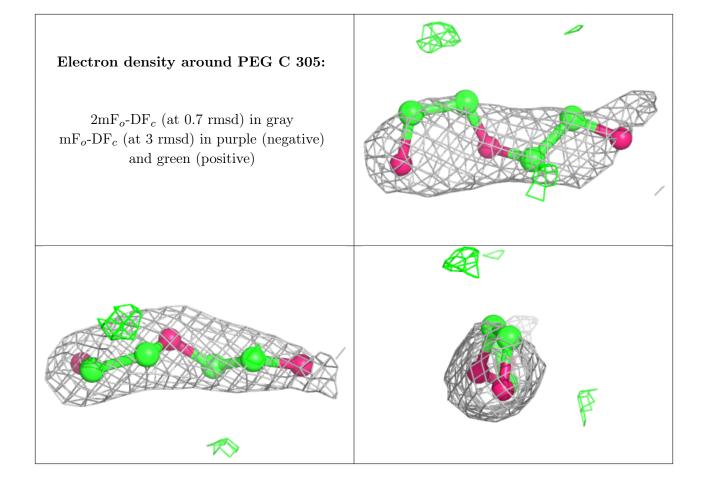




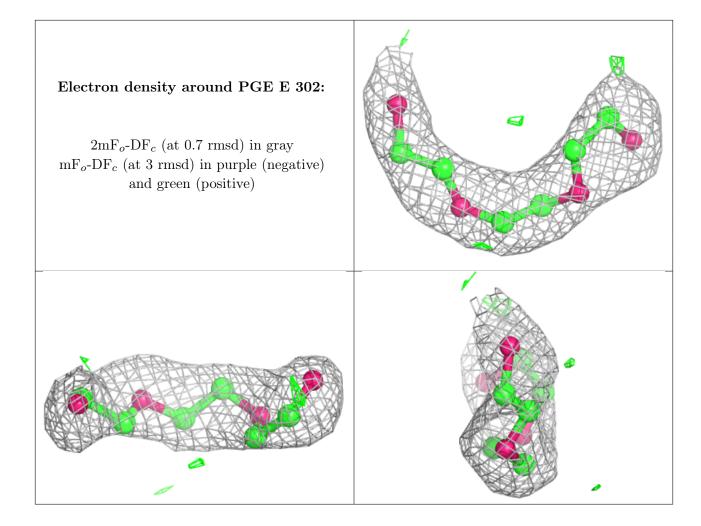




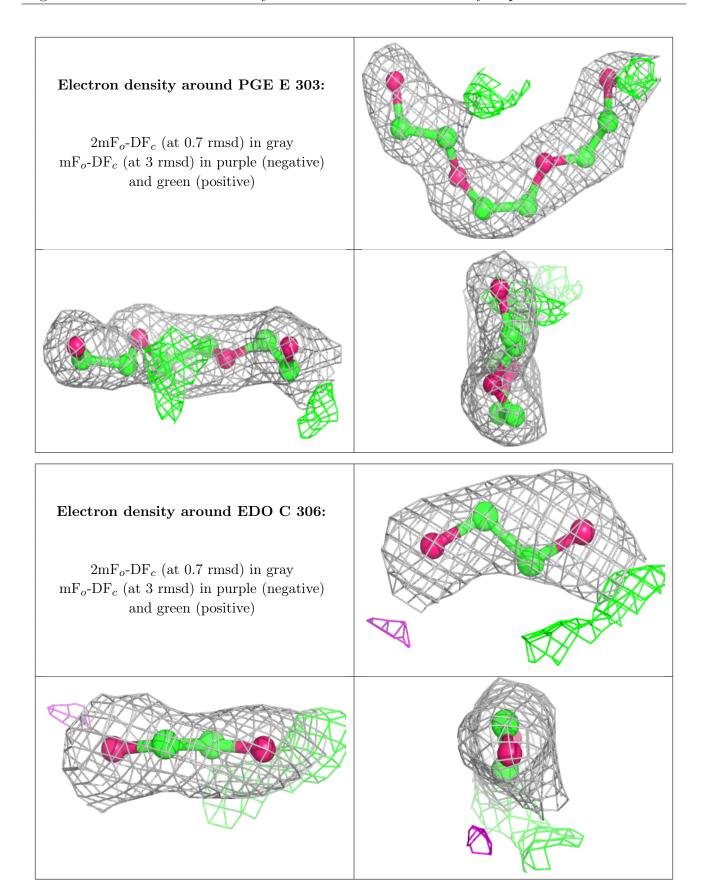




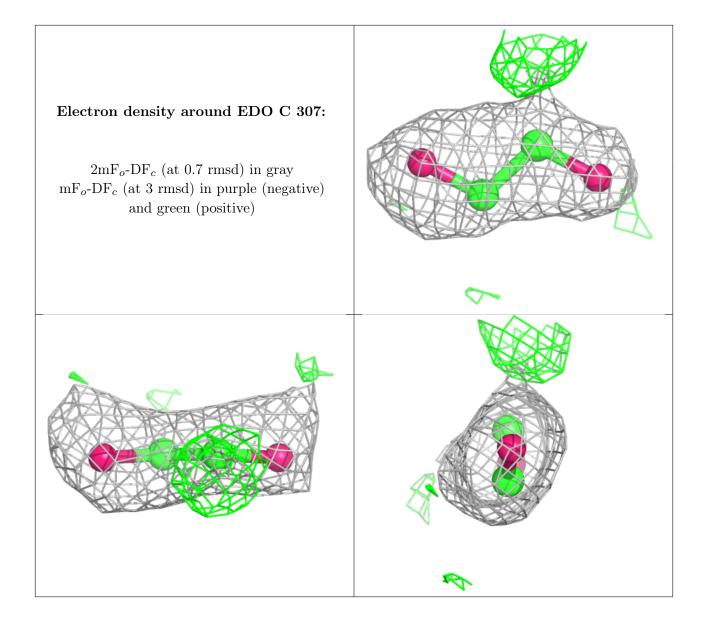




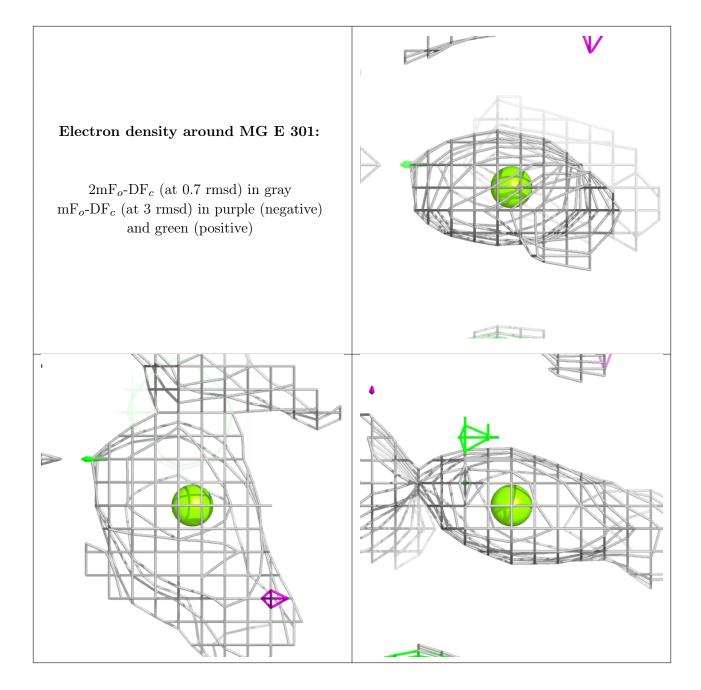








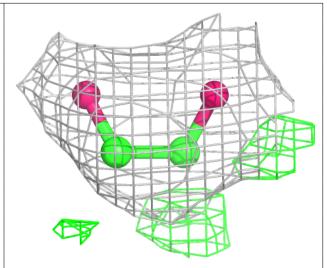


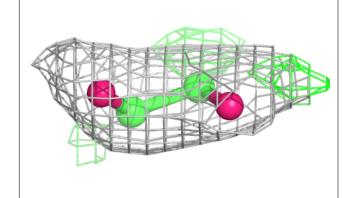


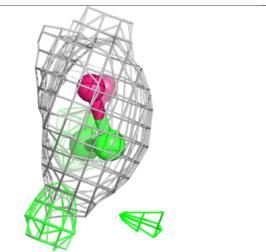


Electron density around EDO F 307:

 $2 {
m mF}_o {
m -DF}_c$ (at 0.7 rmsd) in gray ${
m mF}_o {
m -DF}_c$ (at 3 rmsd) in purple (negative) and green (positive)





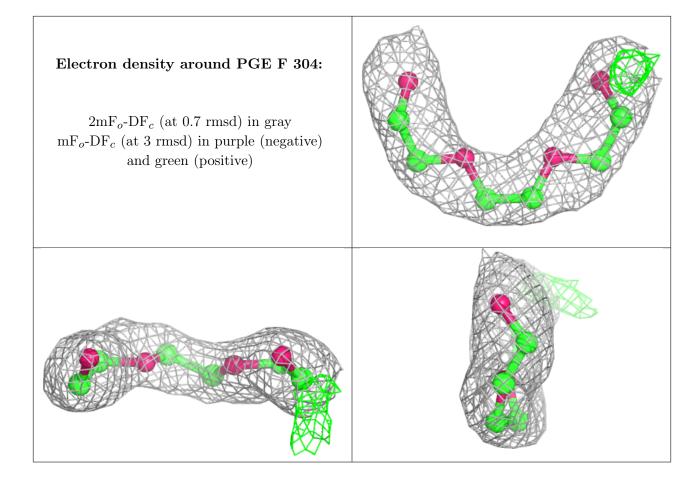




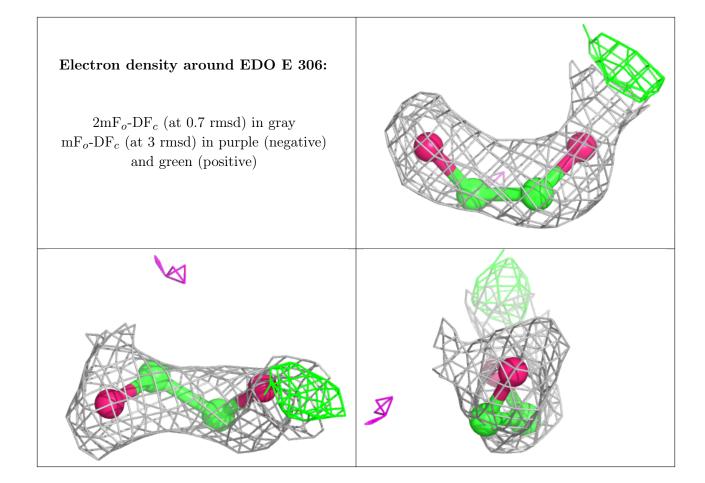
Electron density around EDO F 308: $2 \mathrm{mF}_o\text{-}\mathrm{DF}_c$ (at 0.7 rmsd) in gray ${ m mF}_o{ m -DF}_c$ (at 3 rmsd) in purple (negative) and green (positive) Electron density around PEG D 306: $2 \mathrm{mF}_o\text{-}\mathrm{DF}_c$ (at 0.7 rmsd) in gray mF_o -DF_c (at 3 rmsd) in purple (negative) and green (positive)











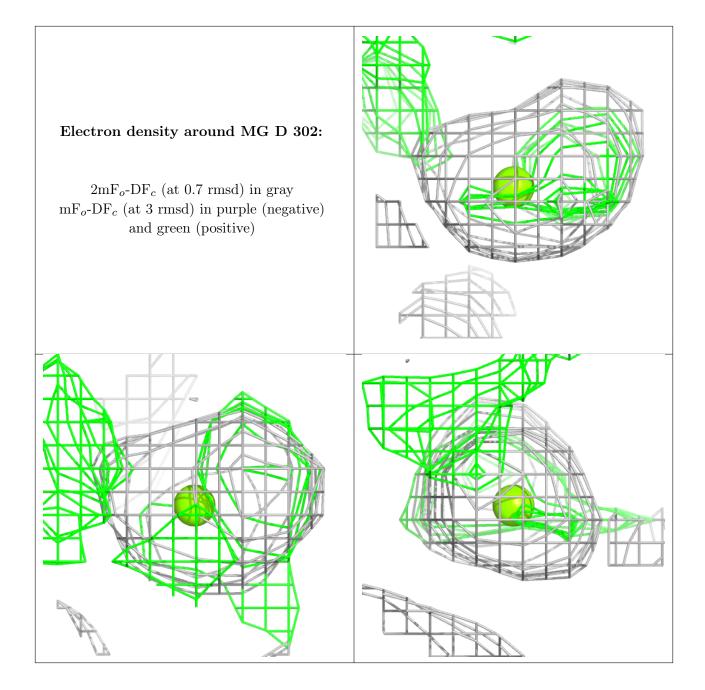


Electron density around MG B 303: $2 \mathrm{mF}_o\text{-}\mathrm{DF}_c$ (at 0.7 rmsd) in gray ${\rm mF}_o\text{-}{\rm DF}_c$ (at 3 rmsd) in purple (negative) and green (positive)

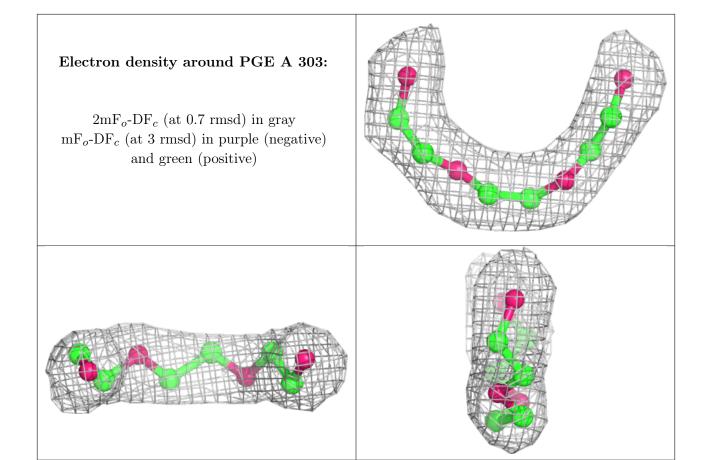


Electron density around PGE D 305: $2 \mathrm{mF}_o\text{-}\mathrm{DF}_c$ (at 0.7 rmsd) in gray ${ m mF}_o{ m -DF}_c$ (at 3 rmsd) in purple (negative) and green (positive) Electron density around PGE C 304: $2 \mathrm{mF}_o\text{-}\mathrm{DF}_c$ (at 0.7 rmsd) in gray ${ m mF}_o{ m -DF}_c$ (at 3 rmsd) in purple (negative) and green (positive)

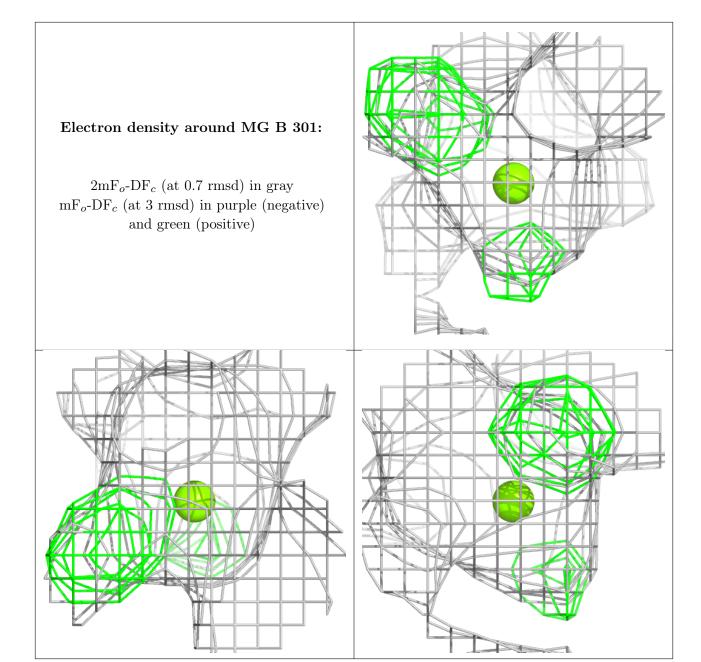




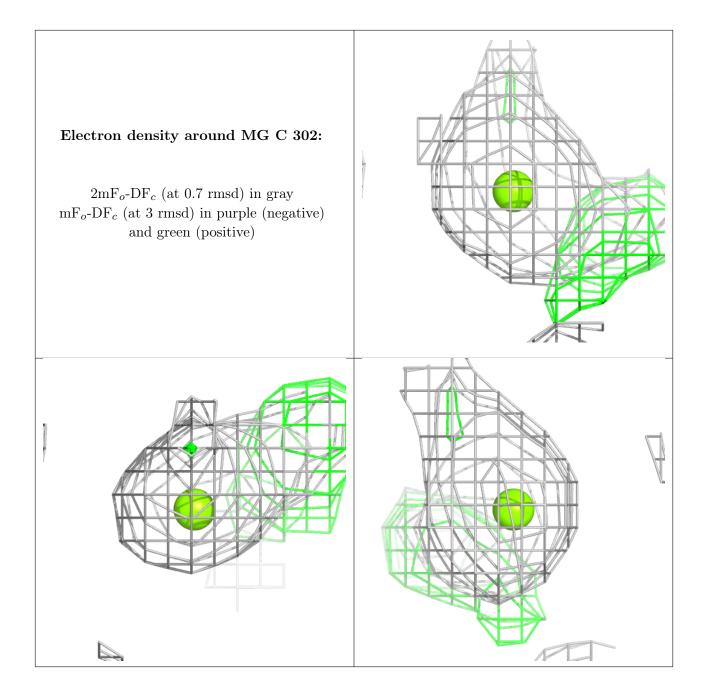




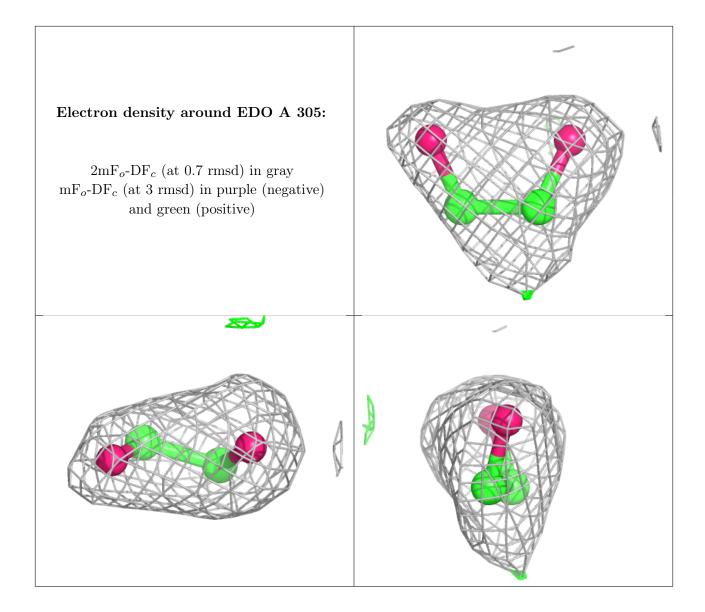








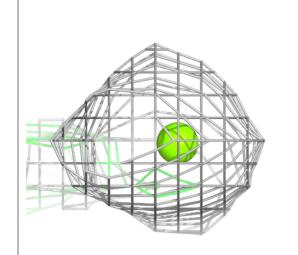


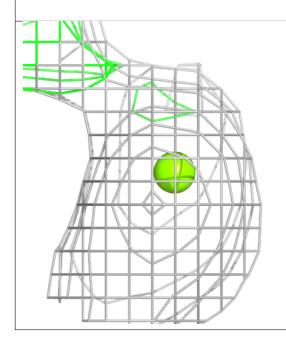


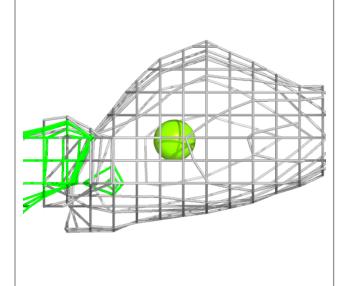


Electron density around MG B 302:

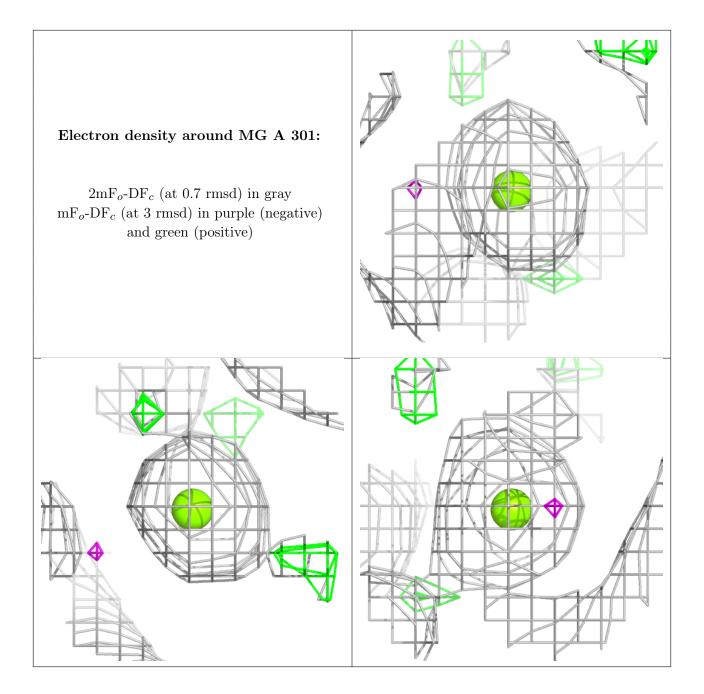
 $2 \mathrm{mF}_o\text{-}\mathrm{DF}_c$ (at 0.7 rmsd) in gray $\mathrm{mF}_o\text{-}\mathrm{DF}_c$ (at 3 rmsd) in purple (negative) and green (positive)







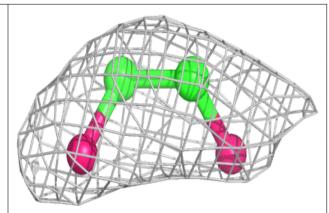


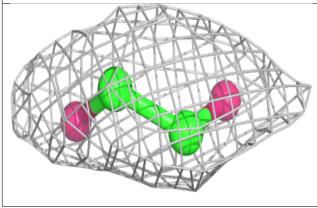


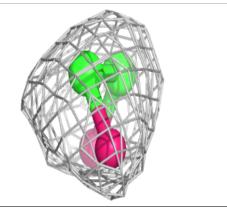


Electron density around EDO D 308:

 $2 \mathrm{mF}_o\text{-}\mathrm{DF}_c$ (at 0.7 rmsd) in gray $\mathrm{mF}_o\text{-}\mathrm{DF}_c$ (at 3 rmsd) in purple (negative) and green (positive)

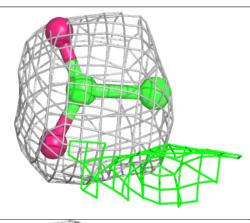


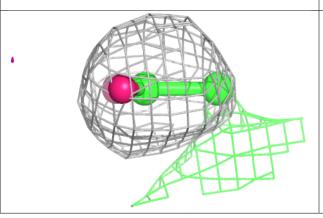


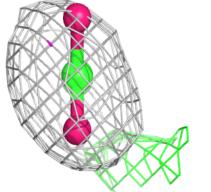


Electron density around ACT E 304:

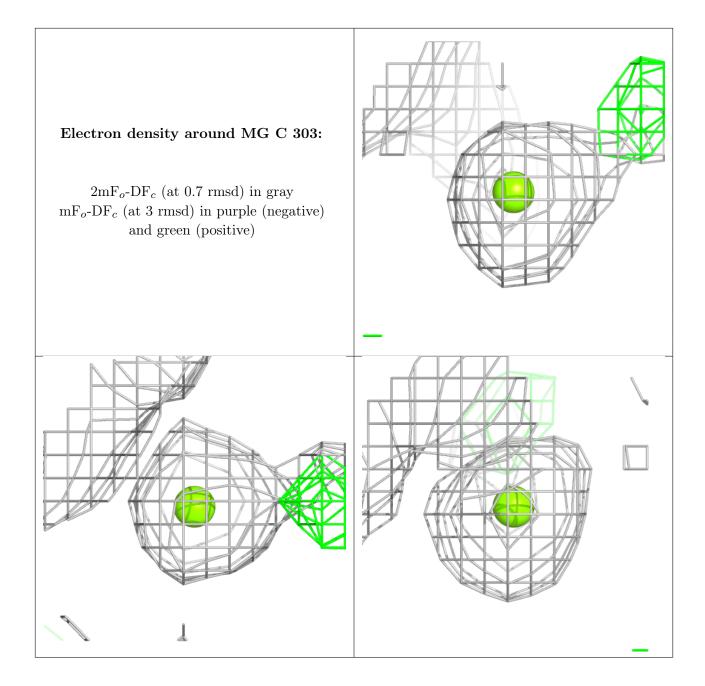
 $2 {
m mF}_o {
m -DF}_c$ (at 0.7 rmsd) in gray ${
m mF}_o {
m -DF}_c$ (at 3 rmsd) in purple (negative) and green (positive)







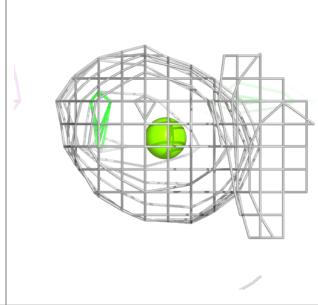


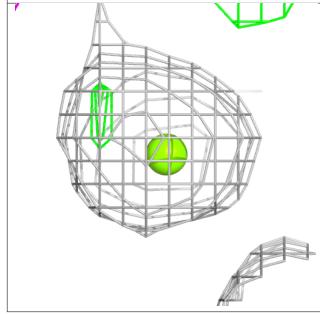


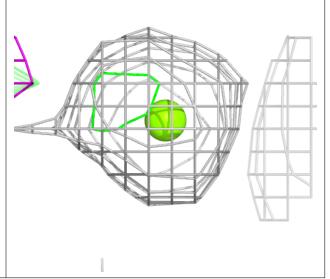


Electron density around MG F 301:

 $2 {
m mF}_o {
m -DF}_c$ (at 0.7 rmsd) in gray ${
m mF}_o {
m -DF}_c$ (at 3 rmsd) in purple (negative) and green (positive)

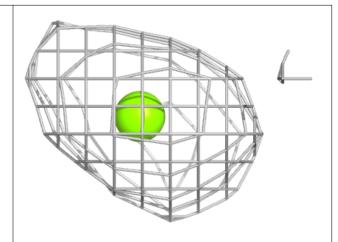


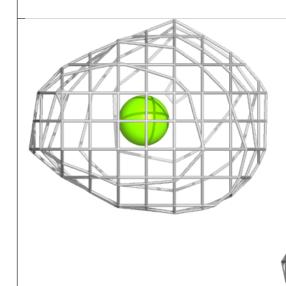


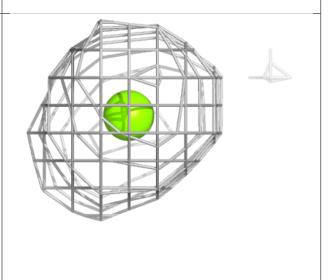


Electron density around MG F 303:

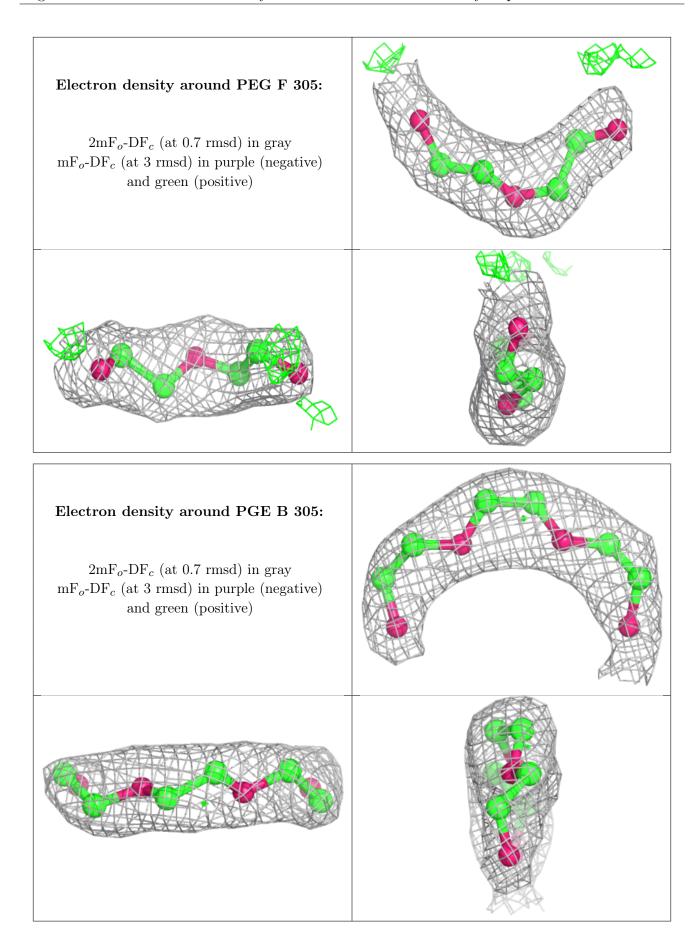
 $2 {
m mF}_o {
m -DF}_c$ (at 0.7 rmsd) in gray ${
m mF}_o {
m -DF}_c$ (at 3 rmsd) in purple (negative) and green (positive)



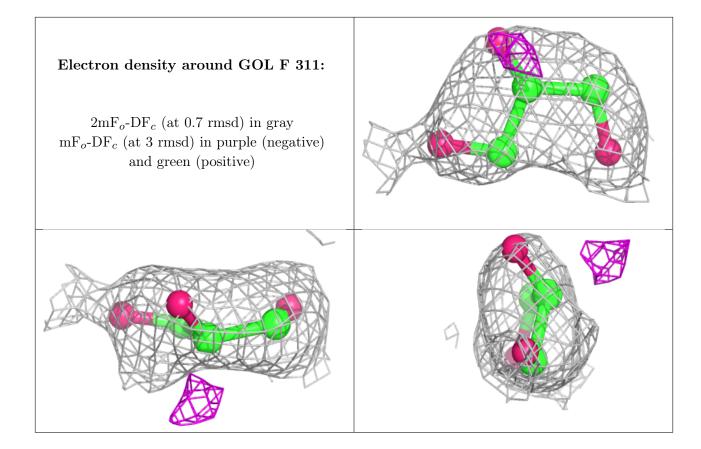




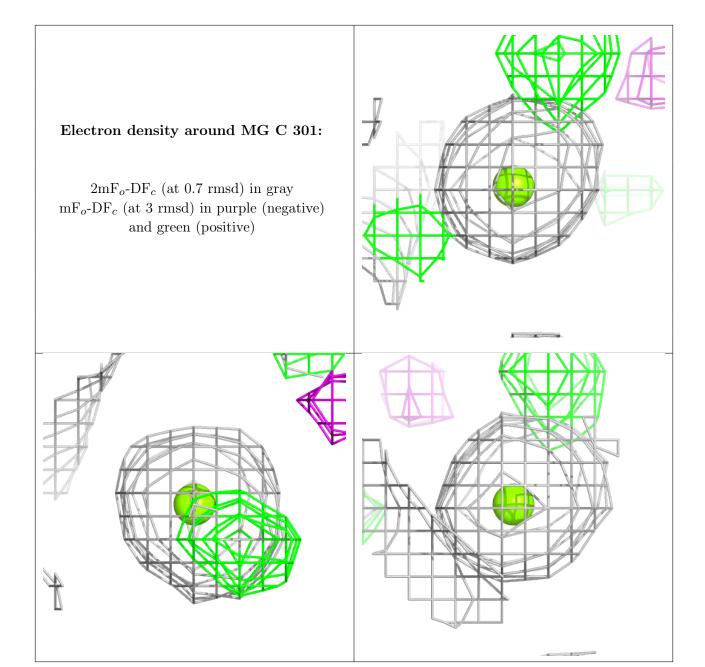












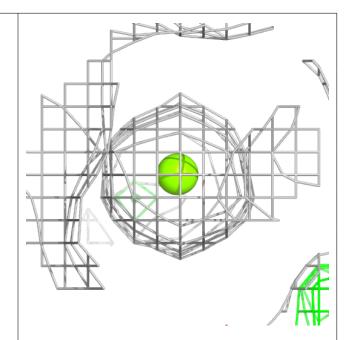


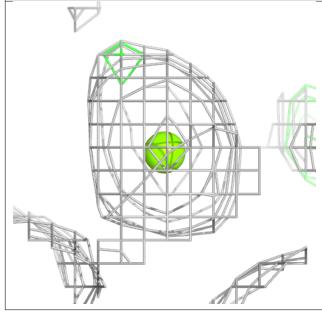
Electron density around MG D 303: $2 \mathrm{mF}_o\text{-}\mathrm{DF}_c$ (at 0.7 rmsd) in gray ${\rm mF}_o\text{-DF}_c$ (at 3 rmsd) in purple (negative) and green (positive)

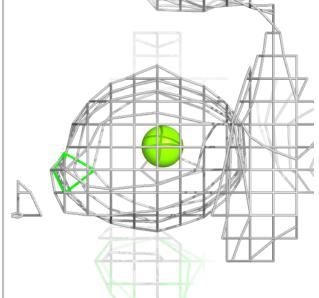


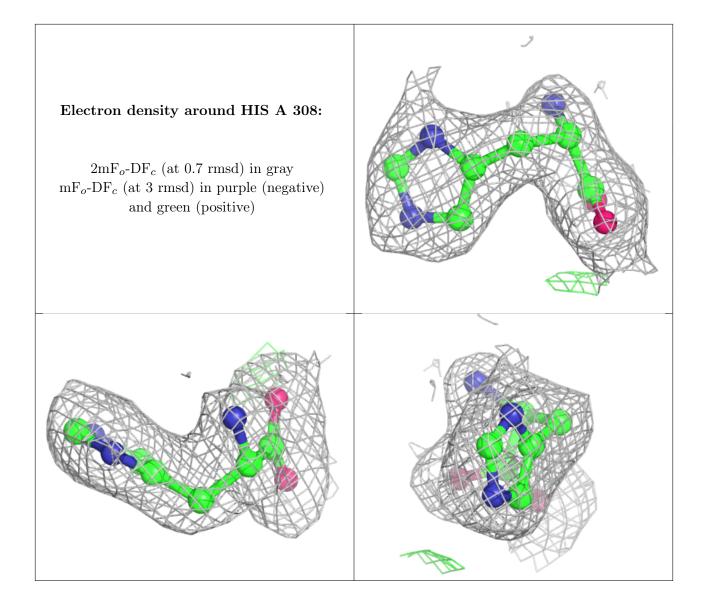
Electron density around MG D 301: $2 \mathrm{mF}_o\text{-DF}_c \text{ (at } 0.7 \text{ rmsd) in gray} \\ \mathrm{mF}_o\text{-DF}_c \text{ (at } 3 \text{ rmsd) in purple (negative)}$

and green (positive)





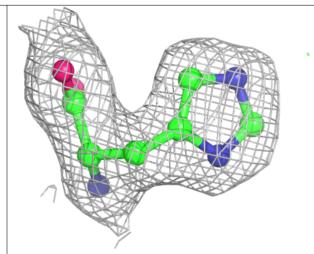


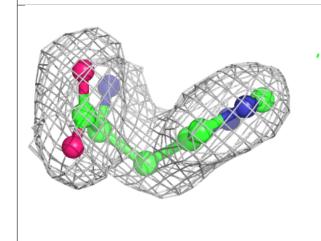


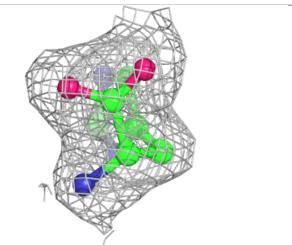


Electron density around HIS B 311:

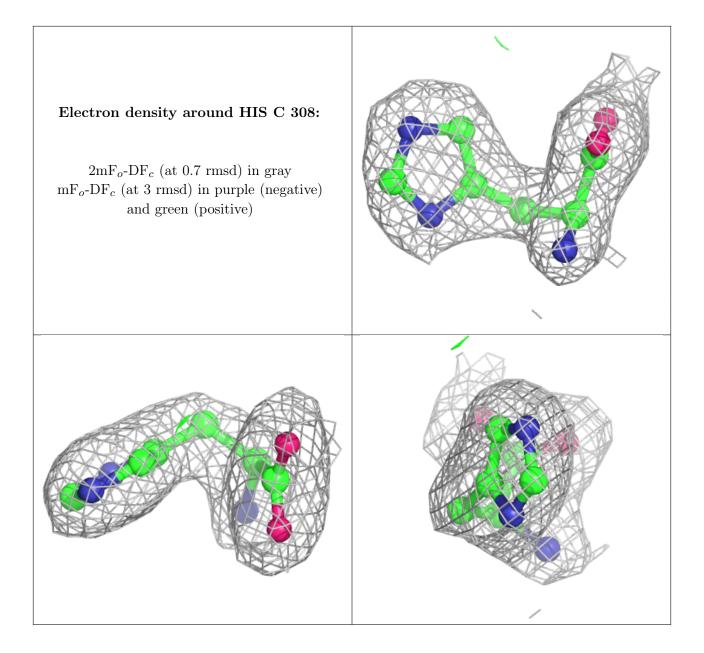
 $2 \text{mF}_o\text{-DF}_c$ (at 0.7 rmsd) in gray $\text{mF}_o\text{-DF}_c$ (at 3 rmsd) in purple (negative) and green (positive)



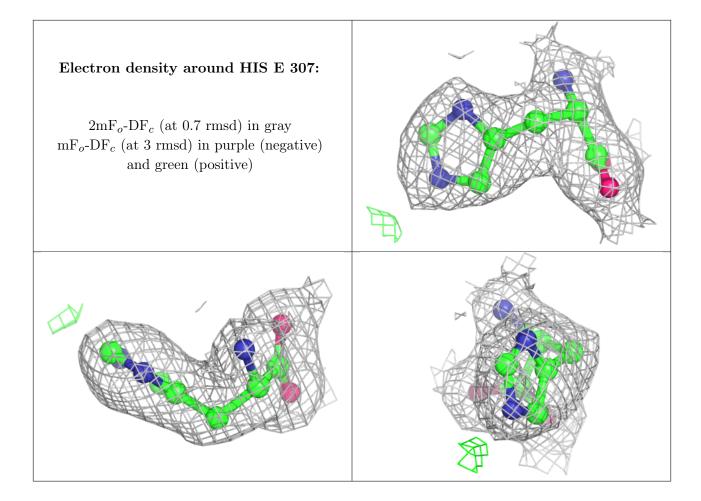




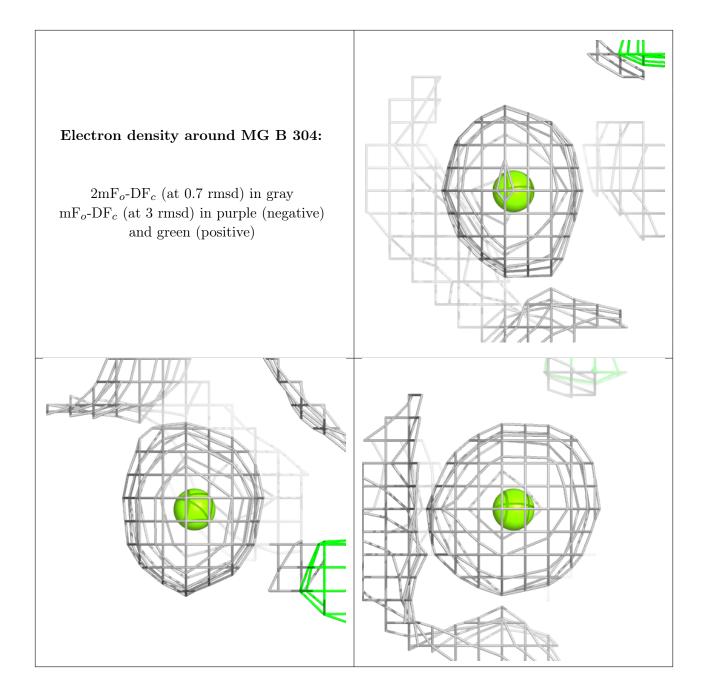




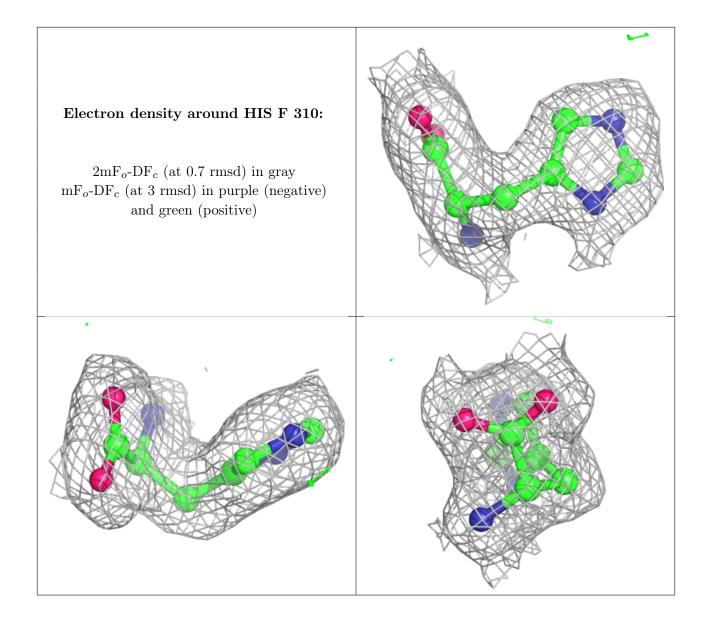








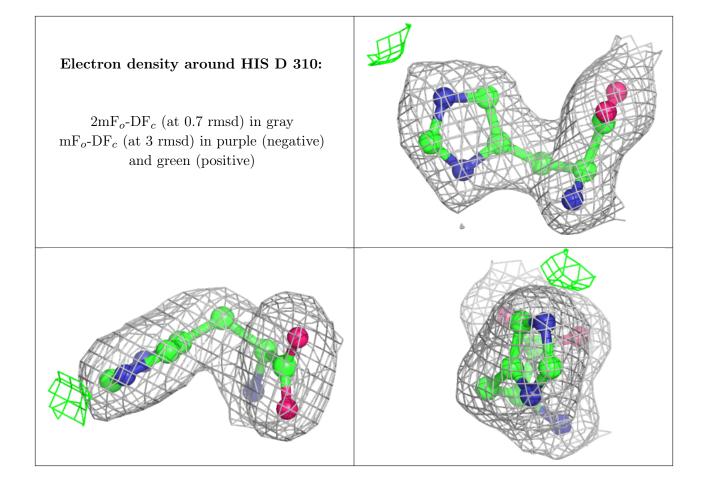




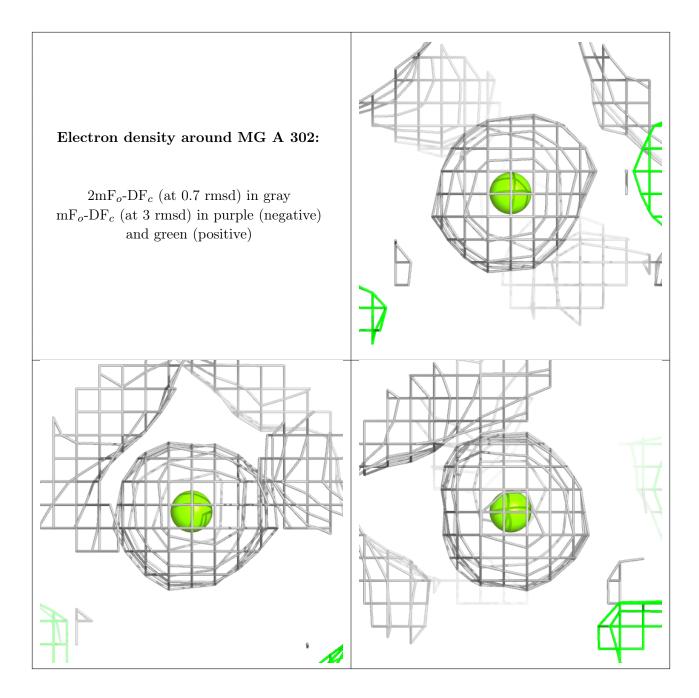


Electron density around MG F 302: $2 \mathrm{mF}_o\text{-}\mathrm{DF}_c$ (at 0.7 rmsd) in gray ${\rm mF}_o\text{-}{\rm DF}_c$ (at 3 rmsd) in purple (negative) and green (positive)









6.5 Other polymers (i)

There are no such residues in this entry.

