



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:26 AM UTC

PDB ID : 7SXM / pdb_00007sxm
Title : Structure of Xenon-derivatized Methyl-Coenzyme M Reductase from Methanothermobacter marburgensis
Authors : Chen, P.Y.-T.; Drennan, C.L.
Deposited on : 2021-11-23
Resolution : 2.50 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 types of validation reports are described at

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

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

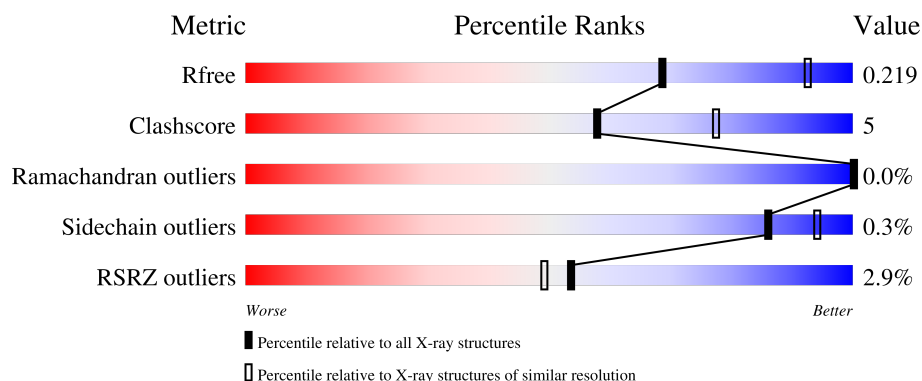
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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

Mol	Chain	Length	Quality of chain
1	A	548	0% 88% 12%
1	D	548	0% 87% 12%
2	B	442	3% 87% 13%
2	E	442	8% 88% 12%
3	C	246	2% 92% 7%

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Mol	Chain	Length	Quality of chain
3	F	246	 A horizontal bar chart showing the quality of chain F. The bar is divided into three segments: a red segment at the beginning labeled '6%', a green segment in the middle labeled '88%', and a yellow segment at the end labeled '10%'. A small black dot is located at the far right end of the bar.

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
1	DYA	A	450	-	X	-	-
10	XE	B	501[A]	-	-	X	-
10	XE	E	501[A]	-	-	X	-

2 Entry composition

There are 11 unique types of molecules in this entry. The entry contains 20235 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 Methyl-coenzyme M reductase I subunit alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	547	Total	C	N	O	S	0	2	0
			4263	2698	717	828	20			
1	D	547	Total	C	N	O	S	0	3	0
			4265	2699	717	829	20			

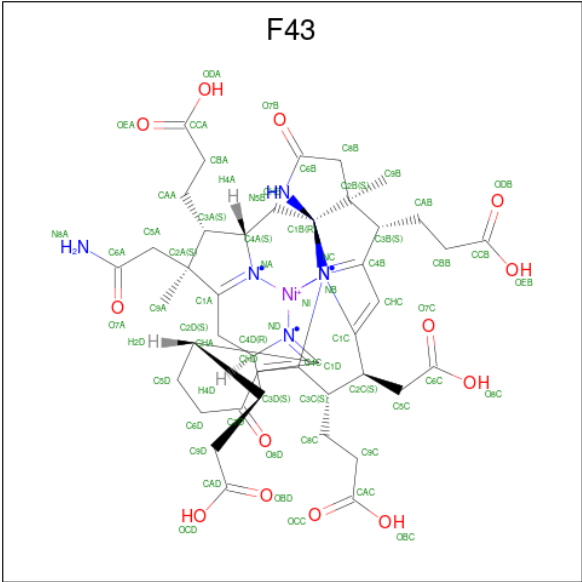
- Molecule 2 is a protein called Methyl-coenzyme M reductase I subunit beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	442	Total	C	N	O	S	0	9	0
			3364	2131	557	658	18			
2	E	442	Total	C	N	O	S	0	9	0
			3364	2131	557	658	18			

- Molecule 3 is a protein called Methyl-coenzyme M reductase I subunit gamma.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	243	Total	C	N	O	S	0	1	0
			1968	1219	350	389	10			
3	F	243	Total	C	N	O	S	0	0	0
			1966	1218	350	388	10			

- Molecule 4 is FACTOR 430 (CCD ID: F43) (formula: C₄₂H₅₁N₆NiO₁₃).

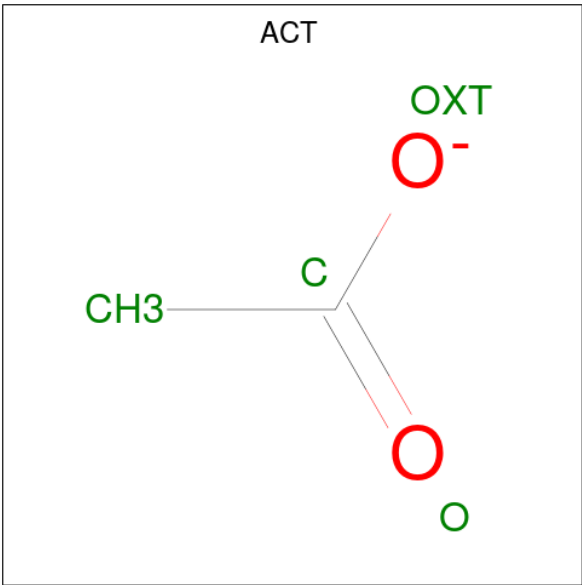


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	Ni	O	0	0
			62	42	6	1	13		
4	A	1	Total	C	N	Ni	O	0	0
			62	42	6	1	13		

- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

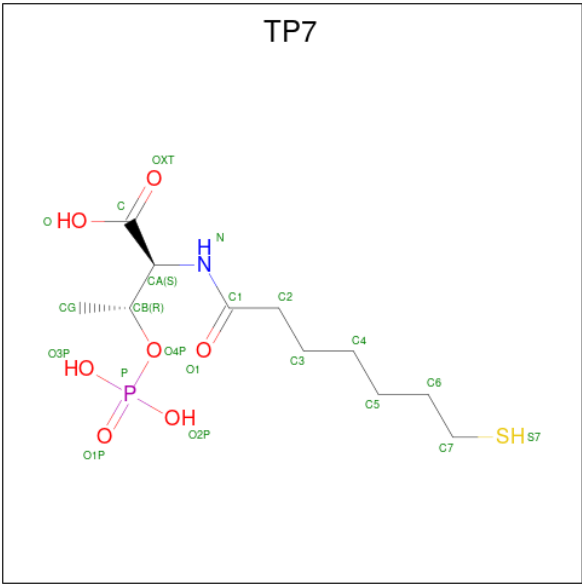
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Mg	0	0
			1	1		
5	D	1	Total	Mg	0	0
			1	1		

- Molecule 6 is ACETATE ION (CCD ID: ACT) (formula: C₂H₃O₂).



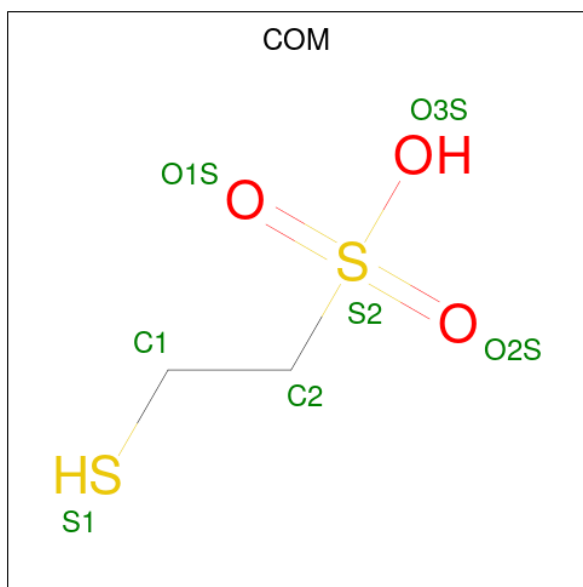
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			4	2	2		
6	A	1	Total	C	O	0	1
			4	2	2		
6	D	1	Total	C	O	0	1
			4	2	2		
6	D	1	Total	C	O	0	0
			4	2	2		

- Molecule 7 is Coenzyme B (CCD ID: TP7) (formula: C₁₁H₂₂NO₇PS).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	A	1	Total	C	N	O	P	S	
			21	11	1	7	1	1	
7	A	1	Total	C	N	O	P	S	
			21	11	1	7	1	1	

- Molecule 8 is 1-THIOETHANESULFONIC ACID (CCD ID: COM) (formula: $C_2H_6O_3S_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	A	1	Total	C	O	S		
			7	2	3	2		
8	D	1	Total	C	O	S		
			7	2	3	2		

- Molecule 9 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	1	Total	Na		
			1	1		

- Molecule 10 is XENON (CCD ID: XE) (formula: Xe) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	1	Total	Xe		
			1	1		
10	B	3	Total	Xe		
			3	3		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	D	1	Total 1	Xe 1	0	0
10	E	3	Total 3	Xe 3	0	2

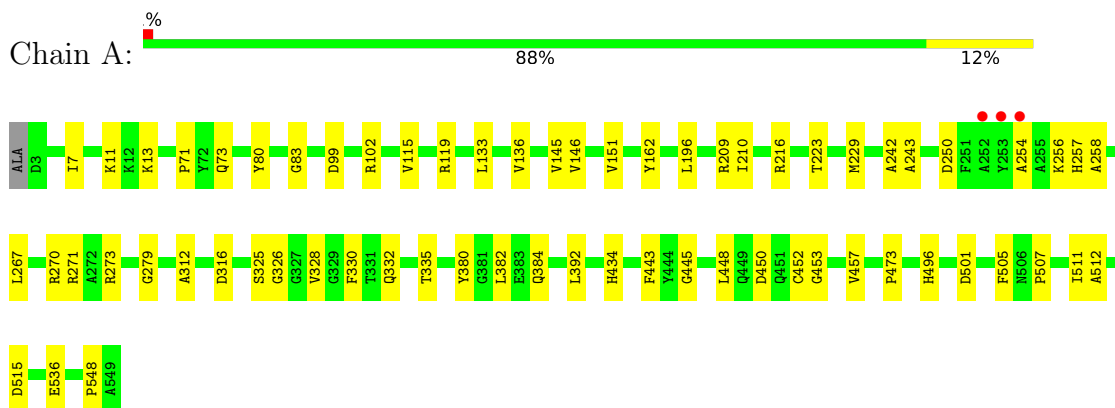
- Molecule 11 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	A	236	Total 236	O 236	0	8
11	B	129	Total 129	O 129	0	1
11	C	67	Total 67	O 67	0	0
11	D	210	Total 210	O 210	0	9
11	E	124	Total 124	O 124	0	1
11	F	72	Total 72	O 72	0	0

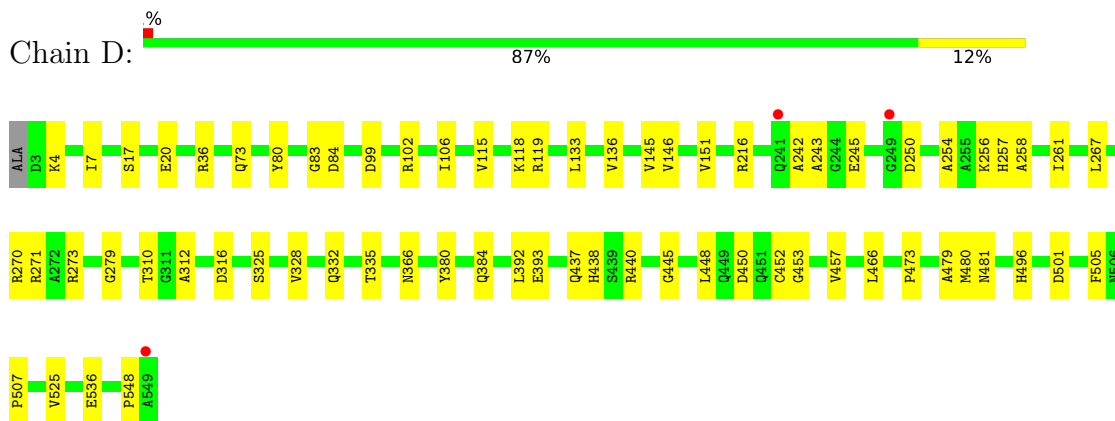
3 Residue-property plots

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

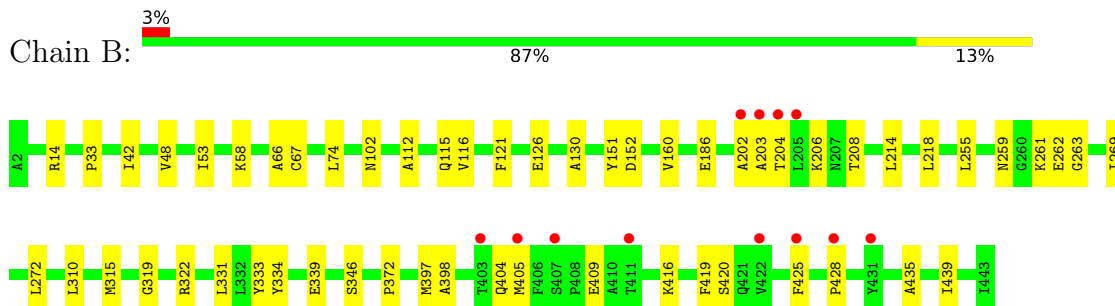
- Molecule 1: Methyl-coenzyme M reductase I subunit alpha



- Molecule 1: Methyl-coenzyme M reductase I subunit alpha

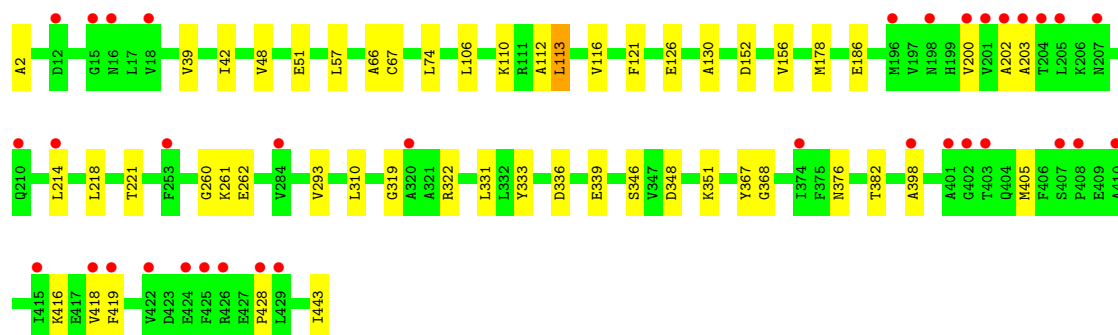


- Molecule 2: Methyl-coenzyme M reductase I subunit beta



- Molecule 2: Methyl-coenzyme M reductase I subunit beta

Chain E: 



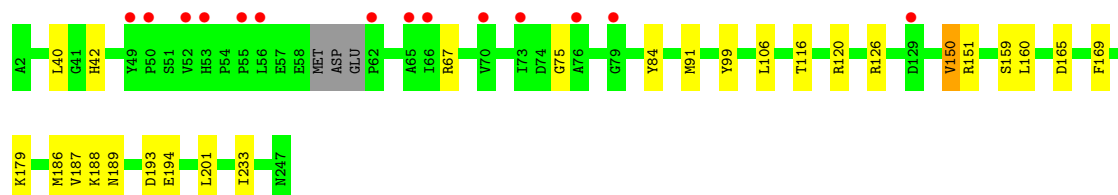
- Molecule 3: Methyl-coenzyme M reductase I subunit gamma

Chain C: 



- Molecule 3: Methyl-coenzyme M reductase I subunit gamma

Chain F: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	81.96Å 115.74Å 123.40Å 90.00° 92.53° 90.00°	Depositor
Resolution (Å)	69.63 – 2.50 69.63 – 2.50	Depositor EDS
% Data completeness (in resolution range)	97.0 (69.63-2.50) 91.7 (69.63-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.19 (at 2.51Å)	Xtriage
Refinement program	PHENIX 1.18_3845	Depositor
R, R_{free}	0.178 , 0.218 0.179 , 0.219	Depositor DCC
R_{free} test set	3842 reflections (2.73%)	wwPDB-VP
Wilson B-factor (Å ²)	32.5	Xtriage
Anisotropy	0.376	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 44.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.026 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	20235	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.67% 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: GL3, F43, AGM, ACT, MGN, NA, MG, SMC, MHS, XE, TP7, COM, DYA

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	0.12	0/4307	0.31	0/5842
1	D	0.11	0/4313	0.31	0/5850
2	B	0.11	0/3421	0.29	0/4635
2	E	0.11	0/3421	0.27	0/4635
3	C	0.12	0/2007	0.33	0/2702
3	F	0.11	0/2001	0.31	0/2694
All	All	0.11	0/19470	0.30	0/26358

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [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	4263	0	4071	42	0
1	D	4265	0	4072	49	0
2	B	3364	0	3347	42	0
2	E	3364	0	3347	35	0
3	C	1968	0	1901	12	0
3	F	1966	0	1900	21	0
4	A	124	0	86	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	1	0	0	0	0
5	D	1	0	0	0	0
6	A	8	0	6	1	0
6	D	8	0	4	1	0
7	A	42	0	36	7	0
8	A	7	0	4	2	0
8	D	7	0	4	2	0
9	A	1	0	0	0	0
10	A	1	0	0	0	0
10	B	3	0	0	3	0
10	D	1	0	0	0	0
10	E	3	0	0	2	0
11	A	236	0	0	4	0
11	B	129	0	0	4	0
11	C	67	0	0	0	0
11	D	210	0	0	3	0
11	E	124	0	0	3	0
11	F	72	0	0	2	0
All	All	20235	0	18778	173	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 (173) 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)
2:E:202:ALA:HB1	2:E:416:LYS:HB3	1.63	0.81
2:B:405:MET:HE3	1:D:115:VAL:HA	1.66	0.77
2:B:405:MET:HE2	1:D:119:ARG:HB2	1.66	0.75
1:D:254:ALA:HA	1:D:258:ALA:HB3	1.67	0.75
1:A:254:ALA:HA	1:A:258:ALA:HB3	1.69	0.74
2:B:214:LEU:HB2	2:B:428:PRO:HG3	1.70	0.74
2:B:42:ILE:HG12	2:B:419[B]:PHE:HZ	1.54	0.72
1:A:243:ALA:HB1	4:A:601:F43:H9B1	1.77	0.67
1:A:145:VAL:HG23	1:A:146:VAL:HG23	1.78	0.66
1:D:270[B]:ARG:NH2	11:D:702:HOH:O	2.26	0.65
2:B:405:MET:HG3	1:D:115:VAL:HG22	1.77	0.65
2:E:214:LEU:HB2	2:E:428:PRO:HG3	1.79	0.65
2:E:351:LYS:NZ	11:E:604:HOH:O	2.25	0.65
1:A:328:VAL:HB	4:A:608:F43:H9A1	1.77	0.64
2:B:214:LEU:HD21	10:B:501[A]:XE:XE	2.76	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:405:MET:HE1	1:D:118:LYS:HB2	1.79	0.64
3:F:165:ASP:OD1	3:F:179:LYS:NZ	2.32	0.63
1:A:270[B]:ARG:NH2	11:A:707:HOH:O	2.33	0.62
4:A:601:F43:H9A1	1:D:328:VAL:HB	1.81	0.61
2:B:202:ALA:HB1	2:B:416:LYS:HB2	1.82	0.61
1:A:270[A]:ARG:NH1	7:A:604[A]:TP7:O1P	2.34	0.60
1:D:145:VAL:HG23	1:D:146:VAL:HG23	1.83	0.60
2:B:203:ALA:HB3	10:B:501[A]:XE:XE	2.79	0.60
2:E:319:GLY:O	2:E:322:ARG:NH1	2.31	0.60
2:E:398:ALA:O	3:F:67:ARG:NH2	2.30	0.59
2:B:310:LEU:HD11	2:B:331:LEU:HD23	1.84	0.59
2:B:116:VAL:HG21	2:B:130:ALA:HA	1.85	0.58
2:E:203:ALA:HB3	10:E:501[A]:XE:XE	2.81	0.58
2:B:206:LYS:NZ	11:B:611:HOH:O	2.35	0.58
2:B:48:VAL:HB	2:B:112:ALA:HB3	1.86	0.57
2:B:255:LEU:HD11	2:B:272:LEU:HD23	1.86	0.57
8:D:602:COM:O2S	3:F:120:ARG:NH1	2.35	0.57
2:B:74:LEU:HD11	2:B:152:ASP:HB3	1.86	0.56
1:D:496:HIS:HB3	1:D:501:ASP:HB2	1.86	0.56
2:E:116:VAL:HG21	2:E:130:ALA:HA	1.88	0.56
3:C:64:ASP:HB3	3:C:67:ARG:HB2	1.87	0.55
2:B:14:ARG:HH12	3:C:63:GLU:HG2	1.71	0.55
1:A:256:LYS:NZ	6:A:606[B]:ACT:O	2.41	0.54
1:A:7:ILE:HG22	1:A:11:LYS:HE2	1.91	0.53
3:C:117:LEU:HD13	3:C:120:ARG:HG3	1.90	0.53
1:A:151:VAL:HG11	1:D:83:GLY:HA3	1.89	0.53
2:B:204:THR:O	11:B:601:HOH:O	2.18	0.53
2:E:42:ILE:HG12	2:E:419[B]:PHE:HZ	1.72	0.53
1:D:106:ILE:HB	1:D:261:ILE:HB	1.91	0.53
1:A:270[B]:ARG:NH1	2:B:186:GLU:OE2	2.41	0.53
1:A:133:LEU:HA	1:A:136:VAL:HG12	1.91	0.53
2:E:348:ASP:HB3	2:E:351:LYS:HB2	1.91	0.52
1:D:4:LYS:HB2	1:D:7:ILE:HG12	1.91	0.52
1:D:270[B]:ARG:NH2	11:D:718:HOH:O	2.41	0.52
1:D:256:LYS:NZ	6:D:601[B]:ACT:O	2.43	0.52
3:F:189:ASN:HD21	3:F:193:ASP:HB2	1.74	0.52
1:A:119:ARG:NH2	1:A:250:ASP:OD2	2.43	0.51
1:D:73:GLN:HB2	1:D:80:TYR:CE2	2.45	0.51
2:B:261:LYS:HG3	2:B:262:GLU:HG3	1.92	0.51
1:D:36:ARG:HG2	1:D:84:ASP:HB2	1.92	0.51
2:E:48:VAL:HB	2:E:112:ALA:HB3	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:33:PRO:HD2	2:B:435:ALA:HB1	1.91	0.51
1:A:13:LYS:NZ	1:A:71:PRO:O	2.34	0.51
7:A:604[A]:TP7:O	11:A:701:HOH:O	2.19	0.51
2:B:339:GLU:HB3	3:C:237:ILE:HD13	1.93	0.51
1:A:453:GLY:O	1:A:457:VAL:HG23	2.11	0.51
2:E:200:VAL:HA	10:E:501[A]:XE:XE	2.89	0.50
1:A:115:VAL:HG22	2:E:405:MET:HG3	1.92	0.50
4:A:601:F43:NB	8:D:602:COM:H22	2.26	0.50
2:E:106:LEU:HD12	2:E:418:VAL:HG22	1.93	0.50
1:D:453:GLY:O	1:D:457:VAL:HG23	2.11	0.50
2:E:310:LEU:HD11	2:E:331:LEU:HD23	1.93	0.50
3:F:75:GLY:HA3	3:F:126:ARG:HB3	1.93	0.50
8:A:605:COM:H22	4:A:608:F43:NB	2.28	0.49
1:A:270[B]:ARG:NH2	11:A:717:HOH:O	2.44	0.49
1:D:133:LEU:HA	1:D:136:VAL:HG12	1.94	0.49
1:D:119:ARG:NH2	1:D:250:ASP:OD2	2.46	0.49
1:A:83:GLY:HA3	1:D:151:VAL:HG11	1.94	0.48
1:A:536:GLU:HA	1:D:548:PRO:HG3	1.94	0.48
1:A:443:PHE:HB2	8:A:605:COM:O1S	2.13	0.48
2:B:66:ALA:O	1:D:507:PRO:HD2	2.14	0.48
2:E:336:ASP:OD1	11:E:601:HOH:O	2.20	0.48
7:A:604[A]:TP7:OXT	11:A:702:HOH:O	2.20	0.48
1:A:548:PRO:HG3	1:D:536:GLU:HA	1.96	0.48
3:C:84:TYR:CZ	1:D:242:ALA:HB2	2.49	0.47
1:A:229:MET:HE1	2:E:367[A]:TYR:HA	1.96	0.47
2:E:261:LYS:HG3	2:E:262:GLU:HG3	1.97	0.47
2:B:151:TYR:CZ	1:D:366:ASN:HA	2.49	0.47
1:D:270[B]:ARG:NH1	2:E:186:GLU:OE2	2.46	0.47
1:D:448:LEU:O	1:D:452:SMC:HCS3	2.15	0.47
2:E:42:ILE:HD12	2:E:218:LEU:HD21	1.97	0.47
3:F:40:LEU:HB3	3:F:42:HIS:CD2	2.49	0.47
1:A:242:ALA:HB2	3:F:84:TYR:CZ	2.50	0.47
1:A:392:LEU:HG	3:C:159:SER:HB2	1.96	0.47
2:B:208:THR:HG22	2:B:315:MET:HE2	1.96	0.47
3:C:174:ARG:HG2	3:C:188:LYS:HB2	1.97	0.47
2:E:333:TYR:HB2	3:F:106:LEU:HD22	1.95	0.47
3:C:84:TYR:CE1	1:D:242:ALA:HB2	2.50	0.47
1:A:209:ARG:NH2	1:A:512:ALA:O	2.38	0.46
2:E:178:MET:HG3	2:E:419[B]:PHE:CZ	2.49	0.46
7:A:609[A]:TP7:H52C	1:D:325:SER:HB3	1.97	0.46
2:B:102:ASN:HB3	2:B:115:GLN:HB2	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:2:ALA:HB2	2:E:443:ILE:HG22	1.96	0.46
2:B:33:PRO:HD3	2:B:439:ILE:HD13	1.98	0.46
1:D:99:ASP:OD1	1:D:102:ARG:NH2	2.49	0.46
1:A:332:GLN:HA	1:A:335:THR:OG1	2.16	0.46
2:E:39:VAL:HG21	2:E:221:THR:HG21	1.96	0.46
2:B:425:PHE:HB3	10:B:501[A]:XE:XE	2.94	0.45
2:E:113:LEU:HB2	2:E:418:VAL:HG13	1.97	0.45
1:A:448:LEU:O	1:A:452:SMC:HCS3	2.16	0.45
7:A:609[A]:TP7:H32C	1:D:480:MET:O	2.16	0.45
1:D:267:LEU:HD12	1:D:273:ARG:HB2	1.98	0.45
3:F:187:VAL:HG23	3:F:188:LYS:HG3	1.98	0.45
1:D:279:GLY:HA2	1:D:473:PRO:HB2	1.98	0.45
2:E:178:MET:HE3	2:E:178:MET:HB3	1.72	0.45
4:A:608:F43:H9B1	1:D:243:ALA:HB1	1.98	0.45
1:D:332:GLN:HA	1:D:335:THR:OG1	2.17	0.45
2:E:57:LEU:HD21	2:E:156:VAL:HA	1.98	0.45
1:A:73:GLN:HB2	1:A:80:TYR:CE2	2.52	0.45
1:A:507:PRO:HD2	2:E:66:ALA:O	2.16	0.45
1:A:216:ARG:HB3	1:D:216:ARG:HB3	1.99	0.45
1:A:505:PHE:CE1	2:E:67:CYS:HB3	2.53	0.44
2:B:404:GLN:OE1	11:B:602:HOH:O	2.21	0.44
1:D:392:LEU:HG	3:F:159:SER:HB2	1.99	0.44
3:F:42:HIS:N	11:F:307:HOH:O	2.42	0.44
7:A:609[A]:TP7:S7	1:D:481:ASN:HA	2.58	0.44
3:F:193:ASP:OD2	11:F:301:HOH:O	2.20	0.44
1:A:279:GLY:HA2	1:A:473:PRO:HB2	1.99	0.44
1:D:393:GLU:HG3	3:F:160:LEU:HG	2.00	0.44
1:A:382:LEU:HD11	1:A:434:HIS:HA	2.00	0.43
2:B:372[B]:PRO:HG2	2:B:397:MET:SD	2.57	0.43
1:D:440:ARG:NH1	3:F:91:MET:HG2	2.34	0.43
2:E:339:GLU:HG3	2:E:346:SER:HB3	2.00	0.43
1:A:210:ILE:HD13	1:A:223:THR:HA	2.01	0.43
2:B:398:ALA:O	3:C:67:ARG:NH1	2.45	0.43
2:E:74:LEU:HD11	2:E:152:ASP:HB3	2.01	0.43
1:A:99:ASP:OD1	1:A:102:ARG:NH2	2.52	0.43
1:A:496:HIS:HB3	1:A:501:ASP:HB2	2.00	0.43
3:C:85:ILE:HG22	3:C:130:LEU:HD11	2.01	0.43
2:B:42:ILE:HD12	2:B:218:LEU:HD21	1.99	0.43
1:A:267:LEU:HD12	1:A:273:ARG:HB2	2.01	0.43
1:A:325:SER:OG	1:A:326:GLY:N	2.52	0.42
2:B:333:TYR:HB2	3:C:106:LEU:HD22	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:409:GLU:OE1	11:B:603:HOH:O	2.21	0.42
1:A:196:LEU:HB3	1:A:511:ILE:HD11	2.01	0.42
2:B:339:GLU:HG3	2:B:346:SER:HB3	2.02	0.42
1:A:162:TYR:OH	1:A:515:ASP:OD2	2.34	0.42
3:C:187:VAL:HG23	3:C:188:LYS:HG3	2.01	0.42
2:B:67:CYS:HB3	1:D:505:PHE:CE1	2.55	0.42
2:B:121:PHE:HZ	2:B:130:ALA:HB2	1.84	0.42
2:B:319:GLY:O	2:B:322:ARG:NH1	2.41	0.42
1:D:380:TYR:O	1:D:384:GLN:HG2	2.20	0.42
1:A:312:ALA:O	1:A:316:ASP:HB2	2.20	0.41
3:F:151:ARG:NH2	3:F:169:PHE:O	2.46	0.41
3:F:186:MET:HB2	3:F:201:LEU:HD11	2.02	0.41
1:A:380:TYR:O	1:A:384:GLN:HG2	2.21	0.41
1:D:17:SER:HB3	1:D:20:GLU:HG3	2.01	0.41
1:D:245:GLU:OE1	11:D:701:HOH:O	2.22	0.41
2:E:376:ASN:O	2:E:382:THR:OG1	2.38	0.41
3:F:150:VAL:HB	3:F:186:MET:HE1	2.02	0.41
3:F:188:LYS:HG2	3:F:194:GLU:HA	2.01	0.41
2:B:58:LYS:HB2	2:B:58:LYS:HE3	1.86	0.41
3:F:189:ASN:ND2	3:F:193:ASP:HB2	2.34	0.41
2:B:203:ALA:HA	2:B:420:SER:HB3	2.01	0.41
1:A:330:PHE:CE2	7:A:604[A]:TP7:H71C	2.55	0.41
3:F:99:TYR:CG	3:F:116:THR:HG21	2.56	0.41
2:B:53:ILE:HD13	2:B:160:VAL:HG22	2.03	0.41
2:B:259:ASN:HB3	2:B:263:GLY:HA3	2.03	0.41
2:E:121:PHE:HZ	2:E:130:ALA:HB2	1.85	0.41
2:E:260:GLY:N	11:E:618:HOH:O	2.53	0.41
2:E:293:VAL:HG22	3:F:233:ILE:HD11	2.02	0.41
1:D:312:ALA:O	1:D:316:ASP:HB2	2.21	0.40
1:D:473:PRO:O	1:D:479:ALA:HA	2.20	0.40
2:B:269:ILE:HD11	2:B:334:TYR:HA	2.02	0.40
1:D:310:THR:OG1	1:D:525:VAL:HG21	2.22	0.40
1:D:437:GLN:HG2	1:D:438:HIS:CD2	2.57	0.40
2:E:51:GLU:CD	2:E:110:LYS:HD2	2.47	0.40
1:D:466:LEU:HD23	1:D:466:LEU:HA	1.97	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	541/548 (99%)	522 (96%)	19 (4%)	0	100	100
1	D	542/548 (99%)	522 (96%)	20 (4%)	0	100	100
2	B	449/442 (102%)	435 (97%)	14 (3%)	0	100	100
2	E	449/442 (102%)	434 (97%)	13 (3%)	2 (0%)	30	49
3	C	240/246 (98%)	236 (98%)	4 (2%)	0	100	100
3	F	239/246 (97%)	236 (99%)	3 (1%)	0	100	100
All	All	2460/2472 (100%)	2385 (97%)	73 (3%)	2 (0%)	100	68

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	E	368[A]	GLY
2	E	368[B]	GLY

5.3.2 Protein sidechains

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	434/432 (100%)	434 (100%)	0	100	100
1	D	435/432 (101%)	435 (100%)	0	100	100
2	B	346/341 (102%)	345 (100%)	1 (0%)	86	94
2	E	346/341 (102%)	344 (99%)	2 (1%)	78	91
3	C	212/214 (99%)	210 (99%)	2 (1%)	70	87

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
3	F	211/214 (99%)	210 (100%)	1 (0%)	81 92
All	All	1984/1974 (100%)	1978 (100%)	6 (0%)	86 94

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	126	GLU
3	C	3	GLN
3	C	150	VAL
2	E	113	LEU
2	E	126	GLU
3	F	150	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	296	ASN
1	D	131	HIS
2	E	318	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

12 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	AGM	D	271	1	10,11,12	1.33	1 (10%)	7,13,15	1.14	0
1	MHS	D	257	1	10,11,12	1.85	3 (30%)	5,14,16	1.12	1 (20%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	MGN	D	400	1	6,9,10	0.84	0	7,12,14	1.05	0
1	SMC	D	452	1	5,6,7	0.97	0	3,6,8	1.33	0
1	GL3	D	445	1	2,3,4	1.46	1 (50%)	1,2,4	0.42	0
1	MGN	A	400	1	6,9,10	0.85	0	7,12,14	0.98	0
1	MHS	A	257	1	10,11,12	1.82	3 (30%)	5,14,16	1.16	1 (20%)
1	SMC	A	452	1	5,6,7	1.00	0	3,6,8	1.40	0
1	DYA	A	450	1	6,7,8	1.93	2 (33%)	6,8,10	2.74	5 (83%)
1	DYA	D	450	1	6,7,8	1.51	1 (16%)	6,8,10	2.49	3 (50%)
1	AGM	A	271	1	10,11,12	1.39	1 (10%)	7,13,15	1.28	0
1	GL3	A	445	1	2,3,4	1.45	1 (50%)	1,2,4	0.42	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
1	AGM	D	271	1	-	3/10/11/13	-
1	MHS	D	257	1	-	0/5/6/8	0/1/1/1
1	MGN	D	400	1	-	0/7/9/12	-
1	SMC	D	452	1	-	1/3/5/7	-
1	GL3	D	445	1	-	1/1/1/2	-
1	MGN	A	400	1	-	0/7/9/12	-
1	MHS	A	257	1	-	0/5/6/8	0/1/1/1
1	SMC	A	452	1	-	1/3/5/7	-
1	DYA	A	450	1	-	4/4/6/8	-
1	DYA	D	450	1	-	4/4/6/8	-
1	AGM	A	271	1	-	2/10/11/13	-
1	GL3	A	445	1	-	1/1/1/2	-

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	257	MHS	CD2-CG	3.90	1.41	1.36
1	A	257	MHS	CD2-CG	3.87	1.41	1.36
1	A	450	DYA	OD1-CG	-3.23	1.22	1.30
1	A	271	AGM	CZ-NE1	3.21	1.38	1.33
1	D	271	AGM	CZ-NE1	3.00	1.38	1.33
1	D	450	DYA	C-CA	2.38	1.49	1.45
1	A	450	DYA	C-CA	2.20	1.48	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	257	MHS	CG-ND1	2.15	1.41	1.38
1	A	257	MHS	CG-ND1	2.11	1.41	1.38
1	D	445	GL3	C-S	-2.03	1.72	1.80
1	D	257	MHS	CE1-ND1	2.02	1.40	1.35
1	A	445	GL3	C-S	-2.01	1.72	1.80
1	A	257	MHS	CE1-ND1	2.01	1.40	1.35

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	450	DYA	CB-CA-N	-3.99	116.00	123.25
1	A	450	DYA	CB-CA-N	-3.99	116.00	123.25
1	A	450	DYA	O-C-CA	-3.79	120.64	125.39
1	D	450	DYA	O-C-CA	-3.76	120.68	125.39
1	A	450	DYA	OD2-CG-CB	-2.33	116.89	124.02
1	A	450	DYA	OD1-CG-CB	2.26	120.11	113.40
1	A	257	MHS	ND1-CE1-NE2	-2.16	111.66	112.94
1	A	450	DYA	CA-CB-CG	-2.07	119.69	126.27
1	D	257	MHS	ND1-CE1-NE2	-2.07	111.71	112.94
1	D	450	DYA	CA-CB-CG	-2.04	119.81	126.27

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	450	DYA	N-CA-CB-CG
1	A	450	DYA	O-C-CA-CB
1	A	452	SMC	CA-CB-SG-CS
1	D	450	DYA	N-CA-CB-CG
1	D	450	DYA	O-C-CA-CB
1	D	452	SMC	CA-CB-SG-CS
1	D	450	DYA	CA-CB-CG-OD1
1	D	450	DYA	CA-CB-CG-OD2
1	A	450	DYA	CA-CB-CG-OD1
1	A	450	DYA	CA-CB-CG-OD2
1	A	271	AGM	CE2-CD-NE1-CZ
1	D	445	GL3	S-C-CA-N
1	D	271	AGM	NE1-CD-CG-CB
1	D	271	AGM	CE2-CD-NE1-CZ
1	A	271	AGM	N-CA-CB-CG
1	A	445	GL3	S-C-CA-N
1	D	271	AGM	N-CA-CB-CG

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	D	452	SMC	1	0
1	A	452	SMC	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 21 ligands modelled in this entry, 11 are monoatomic - leaving 10 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	F43	A	608	1,8	63,71,71	2.54	13 (20%)	73,118,118	1.64	18 (24%)
4	F43	A	601	1,8	63,71,71	2.57	13 (20%)	73,118,118	1.64	18 (24%)
6	ACT	D	601[B]	-	3,3,3	1.35	0	3,3,3	1.51	0
6	ACT	A	606[B]	-	3,3,3	1.36	0	3,3,3	1.50	0
6	ACT	D	605	5	3,3,3	1.37	0	3,3,3	1.37	0
8	COM	A	605	4	6,6,6	1.46	1 (16%)	8,8,8	0.74	0
7	TP7	A	609[A]	-	19,20,20	0.89	0	24,26,26	1.06	1 (4%)
6	ACT	A	603	5	3,3,3	1.40	1 (33%)	3,3,3	1.54	0
7	TP7	A	604[A]	-	19,20,20	0.96	1 (5%)	24,26,26	0.90	0
8	COM	D	602	4	6,6,6	1.53	1 (16%)	8,8,8	0.90	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	F43	A	608	1,8	-	9/28/185/185	-
8	COM	A	605	4	-	3/4/4/4	-
4	F43	A	601	1,8	-	10/28/185/185	-
7	TP7	A	609[A]	-	-	2/24/24/24	-
7	TP7	A	604[A]	-	-	1/24/24/24	-
8	COM	D	602	4	-	3/4/4/4	-

All (30) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	601	F43	NI-NB	9.37	2.12	1.89
4	A	608	F43	NI-NA	9.30	2.11	1.89
4	A	608	F43	NI-NB	9.25	2.11	1.89
4	A	601	F43	NI-NA	9.14	2.11	1.89
4	A	601	F43	NI-ND	7.06	2.06	1.89
4	A	608	F43	NI-ND	6.98	2.06	1.89
4	A	601	F43	CHB-C1B	6.32	1.57	1.53
4	A	608	F43	CHB-C1B	6.21	1.57	1.53
4	A	601	F43	CHA-C4D	5.84	1.58	1.53
4	A	608	F43	CHA-C4D	5.34	1.58	1.53
4	A	608	F43	CHC-C4B	3.54	1.49	1.39
4	A	601	F43	CHC-C4B	3.52	1.49	1.39
4	A	601	F43	C3C-C4C	3.31	1.55	1.50
4	A	601	F43	CHB-C4A	2.90	1.55	1.51
4	A	608	F43	C3C-C4C	2.88	1.55	1.50
8	D	602	COM	C2-S2	-2.79	1.73	1.77
4	A	608	F43	CHB-C4A	2.73	1.55	1.51
8	A	605	COM	C2-S2	-2.66	1.73	1.77
4	A	608	F43	OEB-CCB	-2.53	1.22	1.30
4	A	601	F43	OEB-CCB	-2.47	1.22	1.30
7	A	604[A]	TP7	P-O3P	-2.38	1.46	1.54
4	A	601	F43	OCD-CAD	-2.27	1.23	1.30
4	A	608	F43	C4A-NA	2.25	1.52	1.49
4	A	601	F43	C4C-NC	2.24	1.39	1.35
4	A	608	F43	OCD-CAD	-2.23	1.23	1.30
4	A	601	F43	CBB-CCB	2.11	1.55	1.50
4	A	608	F43	C4C-NC	2.07	1.39	1.35
6	A	603	ACT	CH3-C	2.04	1.57	1.49
4	A	601	F43	C4A-NA	2.01	1.52	1.49
4	A	608	F43	CBB-CCB	2.00	1.55	1.50

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	601	F43	C5D-C2D-C1D	3.51	115.01	110.43
4	A	601	F43	O8D-C7D-C6D	-3.49	115.15	120.87
4	A	608	F43	O8D-C7D-C6D	-3.47	115.17	120.87
4	A	608	F43	C5D-C2D-C1D	3.43	114.91	110.43
4	A	608	F43	OBC-CAC-C9C	3.40	124.73	114.00
4	A	601	F43	OBC-CAC-C9C	3.39	124.72	114.00
4	A	608	F43	C3D-C4D-ND	3.19	107.29	102.34
4	A	608	F43	OCC-CAC-C9C	-3.11	113.24	123.09
4	A	601	F43	C3D-C4D-ND	3.09	107.14	102.34
4	A	601	F43	OCC-CAC-C9C	-3.09	113.30	123.09
4	A	608	F43	C2C-C5C-C6C	-2.94	108.45	113.95
4	A	601	F43	C2C-C5C-C6C	-2.93	108.47	113.95
4	A	608	F43	CAB-C3B-C2B	-2.74	112.95	119.00
4	A	608	F43	C2B-C3B-C4B	-2.64	98.69	101.64
4	A	601	F43	C1B-C2B-C3B	2.56	105.24	101.51
4	A	601	F43	ODB-CCB-CBB	-2.55	115.00	123.09
4	A	608	F43	ODB-CCB-CBB	-2.54	115.04	123.09
4	A	601	F43	CAB-C3B-C2B	-2.53	113.42	119.00
4	A	601	F43	O7C-C6C-C5C	-2.44	115.35	122.84
4	A	608	F43	O7C-C6C-C5C	-2.42	115.41	122.84
4	A	608	F43	C9A-C2A-C1A	2.41	113.32	107.54
4	A	608	F43	C1B-C2B-C3B	2.38	104.98	101.51
4	A	601	F43	C9A-C2A-C1A	2.38	113.24	107.54
4	A	608	F43	OCD-CAD-C9D	2.37	121.37	114.00
4	A	601	F43	OCD-CAD-C9D	2.33	121.23	114.00
4	A	601	F43	C2B-C3B-C4B	-2.28	99.10	101.64
4	A	601	F43	C4B-CHC-C1C	2.21	129.41	125.84
4	A	601	F43	C3B-C4B-CHC	-2.20	118.65	123.33
4	A	608	F43	C4B-CHC-C1C	2.19	129.39	125.84
7	A	609[A]	TP7	O3P-P-O2P	2.19	116.00	107.80
4	A	608	F43	C3B-C4B-CHC	-2.16	118.73	123.33
4	A	601	F43	O8C-C6C-C5C	2.16	120.70	114.00
4	A	608	F43	O8C-C6C-C5C	2.13	120.64	114.00
4	A	601	F43	C9B-C2B-C8B	-2.10	105.31	110.61
4	A	608	F43	OBD-CAD-C9D	-2.07	116.50	122.84
4	A	608	F43	C9B-C2B-C8B	-2.07	105.41	110.61
4	A	601	F43	OBD-CAD-C9D	-2.04	116.59	122.84

There are no chirality outliers.

All (28) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
8	A	605	COM	C1-C2-S2-O1S

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Mol	Chain	Res	Type	Atoms
8	A	605	COM	C1-C2-S2-O2S
8	D	602	COM	C1-C2-S2-O1S
8	D	602	COM	C1-C2-S2-O2S
8	D	602	COM	C1-C2-S2-O3S
4	A	608	F43	C3A-CAA-CBA-CCA
8	A	605	COM	C1-C2-S2-O3S
4	A	601	F43	C3A-CAA-CBA-CCA
4	A	608	F43	C2C-C5C-C6C-O8C
4	A	608	F43	C2C-C5C-C6C-O7C
7	A	604[A]	TP7	C2-C3-C4-C5
4	A	601	F43	CAB-CBB-CCB-OEB
4	A	601	F43	C2C-C5C-C6C-O7C
4	A	601	F43	C2C-C5C-C6C-O8C
4	A	601	F43	CAB-CBB-CCB-ODB
4	A	601	F43	CAA-CBA-CCA-OEA
7	A	609[A]	TP7	C2-C3-C4-C5
4	A	608	F43	CAA-CBA-CCA-OEA
4	A	608	F43	CAB-CBB-CCB-OEB
4	A	601	F43	CAA-CBA-CCA-ODA
4	A	608	F43	CAA-CBA-CCA-ODA
4	A	608	F43	C3D-C9D-CAD-OC'D
7	A	609[A]	TP7	C1-C2-C3-C4
4	A	608	F43	CAB-CBB-CCB-ODB
4	A	601	F43	C3D-C9D-CAD-OC'D
4	A	608	F43	C3D-C9D-CAD-OB'D
4	A	601	F43	C4B-C3B-CAB-CBB
4	A	601	F43	C3D-C9D-CAD-OB'D

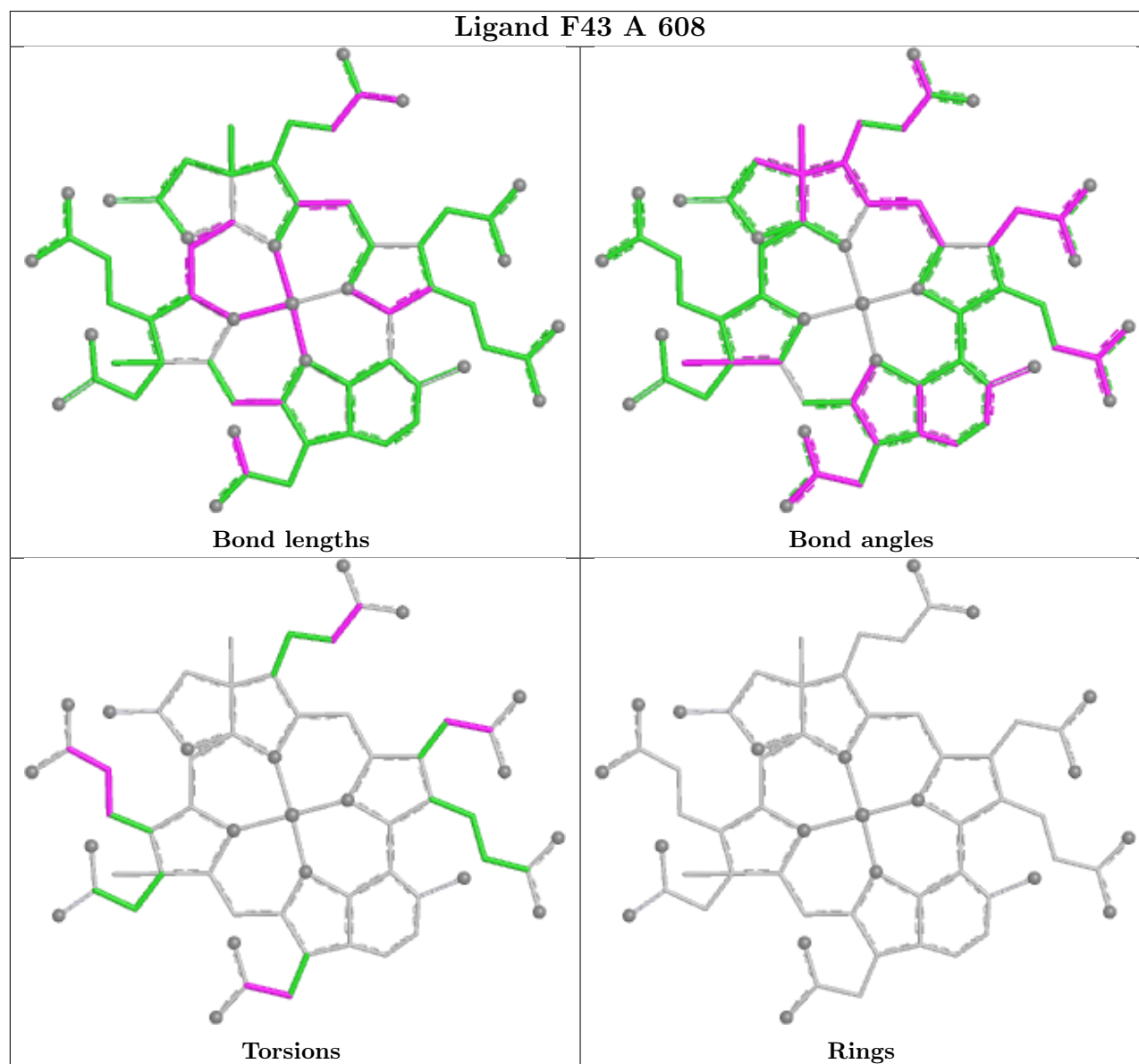
There are no ring outliers.

8 monomers are involved in 17 short contacts:

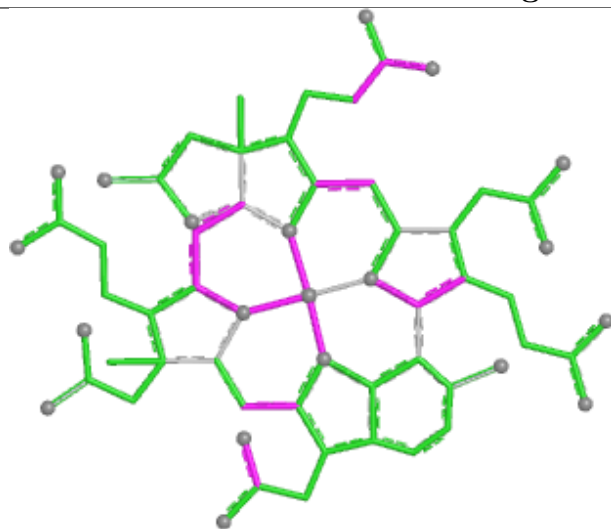
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	608	F43	3	0
4	A	601	F43	3	0
6	D	601[B]	ACT	1	0
6	A	606[B]	ACT	1	0
8	A	605	COM	2	0
7	A	609[A]	TP7	3	0
7	A	604[A]	TP7	4	0
8	D	602	COM	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths,

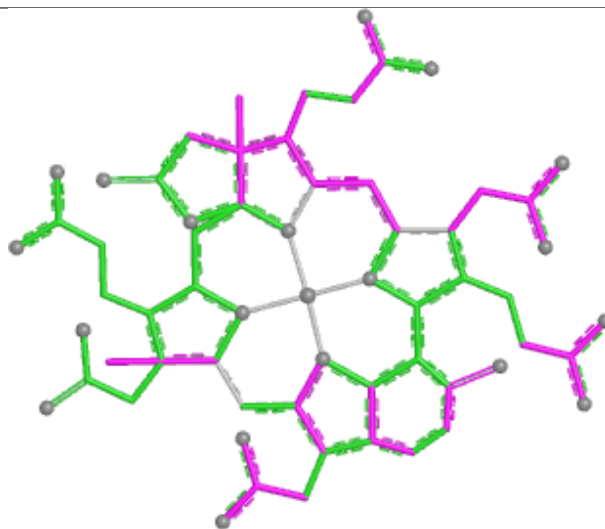
bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within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



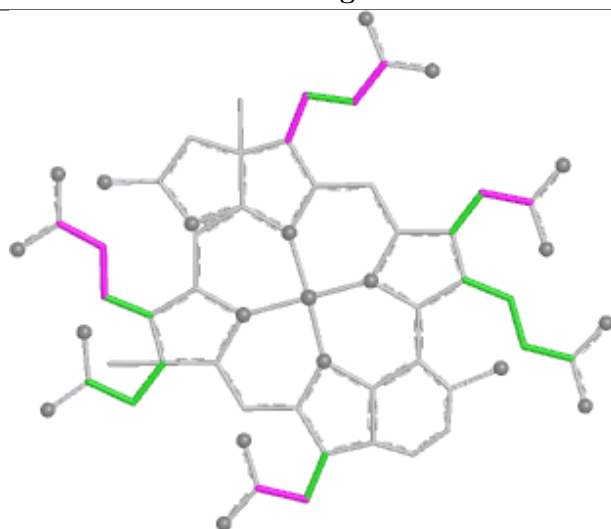
Ligand F43 A 601



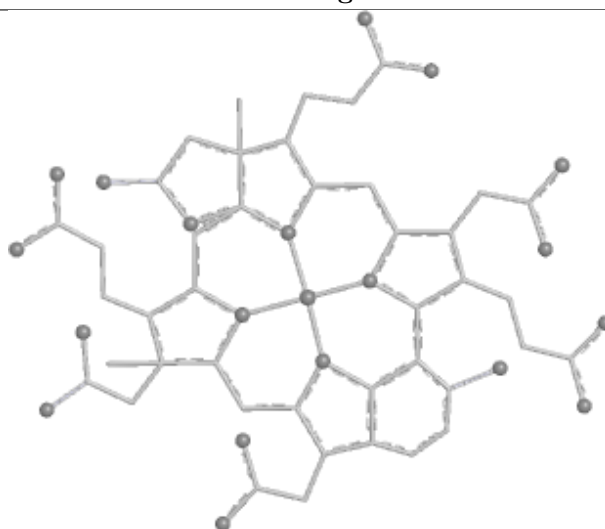
Bond lengths



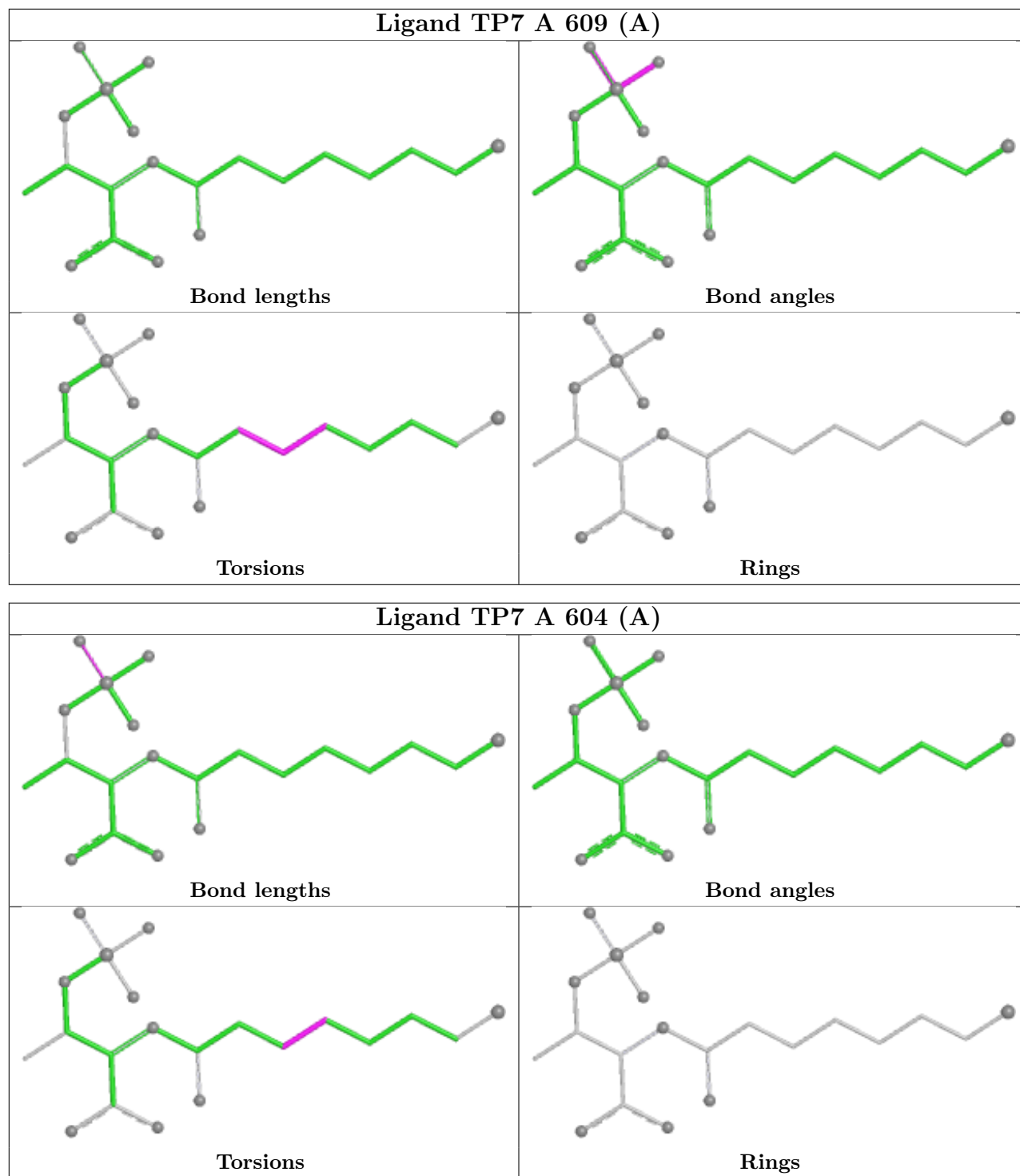
Bond angles



Torsions



Rings



5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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	541/548 (98%)	-0.28	3 (0%) 85 83	20, 31, 50, 69	2 (0%)
1	D	541/548 (98%)	-0.33	3 (0%) 85 83	17, 30, 46, 66	3 (0%)
2	B	442/442 (100%)	0.11	12 (2%) 56 51	17, 39, 57, 71	9 (2%)
2	E	442/442 (100%)	0.42	35 (7%) 18 16	19, 44, 71, 82	9 (2%)
3	C	243/246 (98%)	0.13	4 (1%) 70 67	27, 41, 64, 77	1 (0%)
3	F	243/246 (98%)	0.47	14 (5%) 29 25	30, 49, 76, 88	0
All	All	2452/2472 (99%)	0.02	71 (2%) 53 49	17, 37, 65, 88	24 (0%)

All (71) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	E	202	ALA	4.4
2	E	205	LEU	4.4
2	B	203	ALA	3.9
2	B	202	ALA	3.8
2	E	422	VAL	3.7
3	F	52	VAL	3.6
3	F	76	ALA	3.5
2	E	407	SER	3.5
2	E	374	ILE	3.4
3	F	62	PRO	3.3
2	E	424	GLU	3.3
2	E	425	PHE	3.2
3	F	66	ILE	3.2
2	E	398	ALA	3.2
2	E	429	LEU	3.2
3	F	70	VAL	3.2
3	C	76	ALA	3.2
2	E	418	VAL	3.0
2	B	422	VAL	2.9

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Mol	Chain	Res	Type	RSRZ
1	D	549	ALA	2.9
2	E	320	ALA	2.9
2	E	210	GLN	2.9
2	E	428	PRO	2.8
3	F	56	LEU	2.8
1	A	253	TYR	2.8
2	E	410	ALA	2.7
2	B	425	PHE	2.7
2	E	419[A]	PHE	2.7
3	C	62	PRO	2.7
2	E	200	VAL	2.6
2	E	401	ALA	2.6
2	E	203	ALA	2.6
2	B	405	MET	2.6
2	E	253	PHE	2.6
2	E	214	LEU	2.6
3	F	79	GLY	2.5
2	E	201	VAL	2.5
1	A	254	ALA	2.5
2	E	426	ARG	2.5
2	E	408	PRO	2.4
2	E	15	GLY	2.4
3	F	129	ASP	2.4
2	B	205	LEU	2.3
2	E	12	ASP	2.3
3	F	53	HIS	2.3
3	F	73	ILE	2.3
2	E	196	MET	2.3
2	B	431	TYR	2.3
2	B	204	THR	2.3
1	D	241	GLN	2.3
2	E	415	ILE	2.3
2	E	204	THR	2.2
2	B	411	THR	2.2
1	A	252	ALA	2.2
2	E	16	ASN	2.2
3	F	65	ALA	2.2
2	B	407	SER	2.2
2	B	428	PRO	2.2
3	F	50	PRO	2.1
2	E	403	THR	2.1
2	E	18	VAL	2.1

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Mol	Chain	Res	Type	RSRZ
2	E	198	ASN	2.1
1	D	249	GLY	2.1
2	E	402	GLY	2.1
2	E	207	ASN	2.1
3	F	49	TYR	2.1
3	F	55	PRO	2.1
2	E	284	VAL	2.0
3	C	56	LEU	2.0
2	B	403	THR	2.0
3	C	54	PRO	2.0

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	MHS	A	257	11/12	0.81	0.15	49,51,55,57	0
1	MHS	D	257	11/12	0.85	0.15	44,46,50,50	0
1	AGM	A	271	12/13	0.92	0.09	27,32,35,36	0
1	AGM	D	271	12/13	0.92	0.10	27,30,32,32	0
1	MGN	A	400	10/11	0.94	0.07	21,25,28,28	0
1	MGN	D	400	10/11	0.94	0.07	23,27,29,33	0
1	DYA	D	450	8/9	0.96	0.07	24,29,30,31	0
1	DYA	A	450	8/9	0.98	0.05	24,27,29,29	0
1	SMC	A	452	7/8	0.98	0.06	27,28,30,30	0
1	GL3	A	445	4/5	0.98	0.06	23,24,24,26	0
1	SMC	D	452	7/8	0.98	0.06	27,29,31,31	0
1	GL3	D	445	4/5	0.99	0.05	26,27,27,29	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides 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,

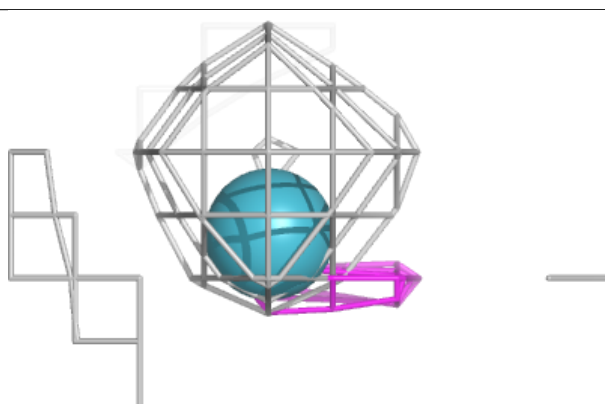
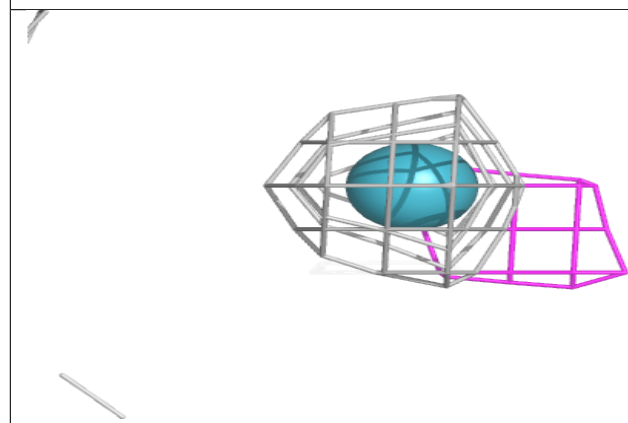
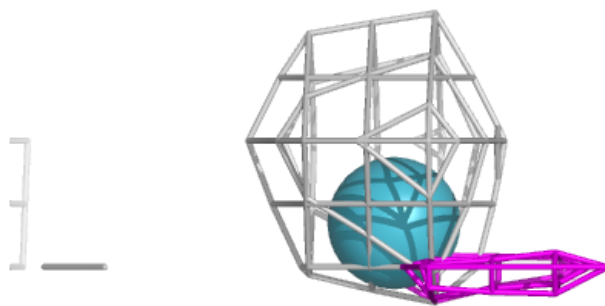
median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
10	XE	B	503	1/1	0.84	0.29	40,40,40,40	1
7	TP7	A	609[A]	21/21	0.91	0.12	28,34,42,43	21
6	ACT	D	601[B]	4/4	0.92	0.15	35,35,35,36	4
10	XE	E	501[A]	1/1	0.93	0.07	45,45,45,45	1
6	ACT	A	606[B]	4/4	0.94	0.17	37,38,39,40	4
7	TP7	A	604[A]	21/21	0.95	0.08	25,33,38,39	21
10	XE	B	501[A]	1/1	0.95	0.07	51,51,51,51	1
8	COM	D	602	7/7	0.96	0.12	35,37,44,45	0
8	COM	A	605	7/7	0.96	0.16	31,34,40,42	0
6	ACT	D	605	4/4	0.97	0.10	32,32,33,35	0
6	ACT	A	603	4/4	0.97	0.06	33,35,37,37	0
10	XE	B	502[B]	1/1	0.97	0.05	47,47,47,47	1
4	F43	A	601	62/62	0.97	0.07	26,30,35,39	0
10	XE	D	603	1/1	0.97	0.05	49,49,49,49	1
4	F43	A	608	62/62	0.97	0.07	21,26,31,33	0
10	XE	E	503	1/1	0.97	0.25	40,40,40,40	1
10	XE	E	502[B]	1/1	0.98	0.04	47,47,47,47	1
10	XE	A	610	1/1	0.98	0.05	46,46,46,46	1
9	NA	A	607	1/1	0.99	0.04	20,20,20,20	0
5	MG	A	602	1/1	0.99	0.02	37,37,37,37	0
5	MG	D	604	1/1	0.99	0.04	34,34,34,34	0

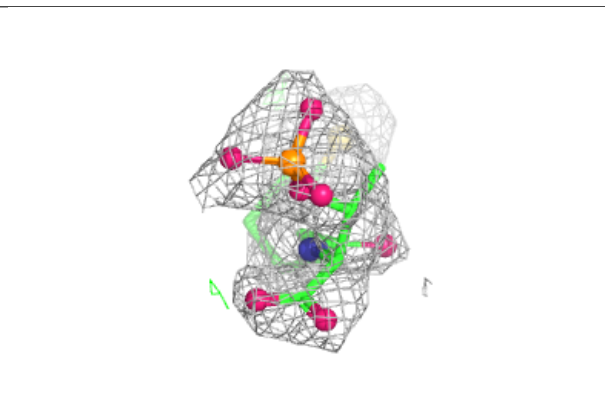
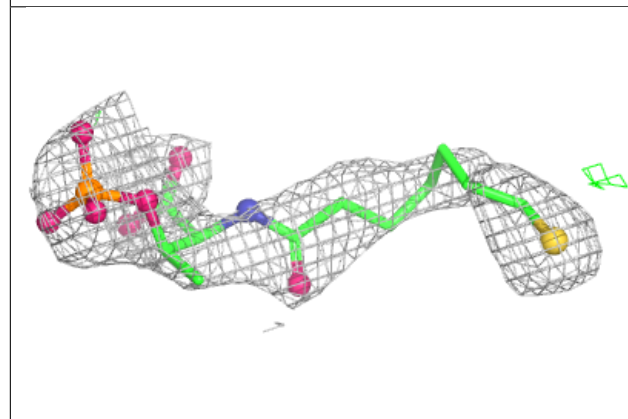
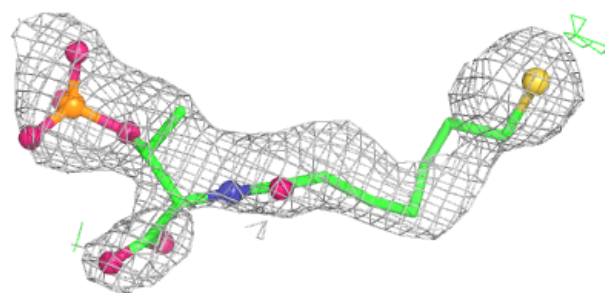
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around XE B 503:

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

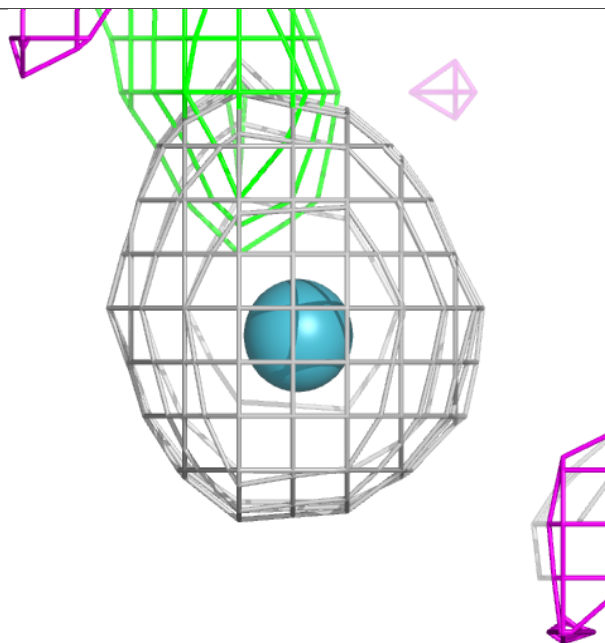
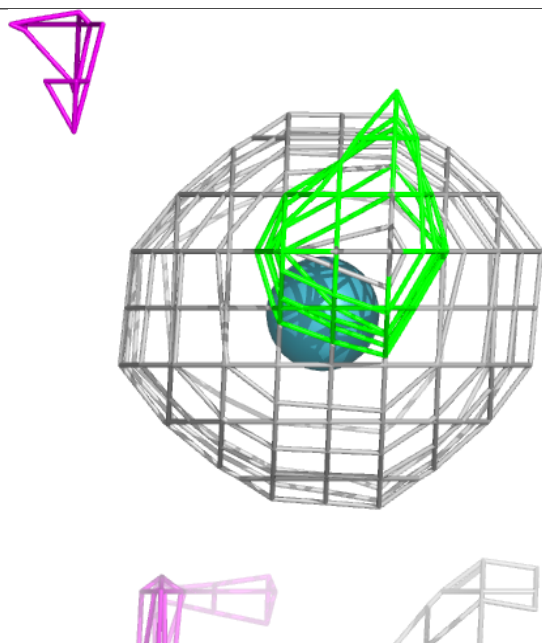
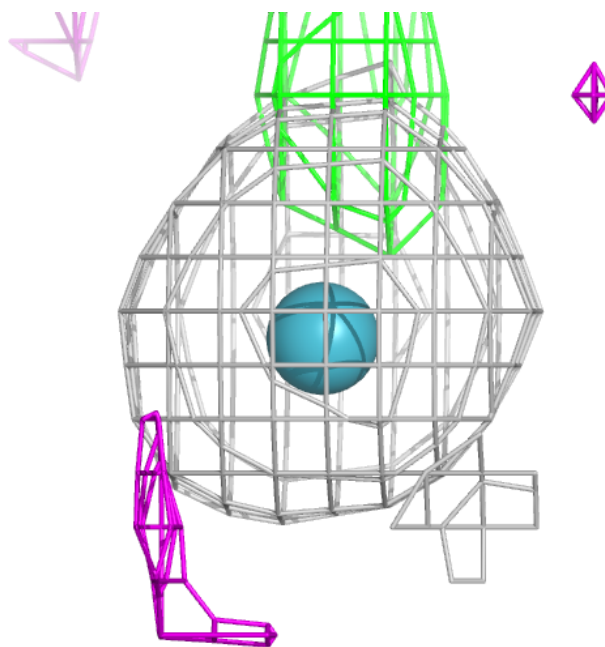
**Electron density around TP7 A 609 (A):**

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



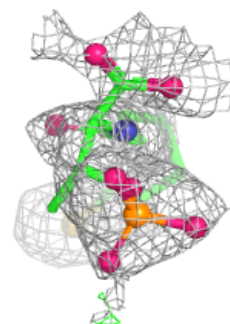
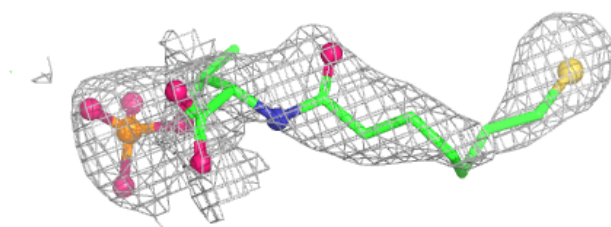
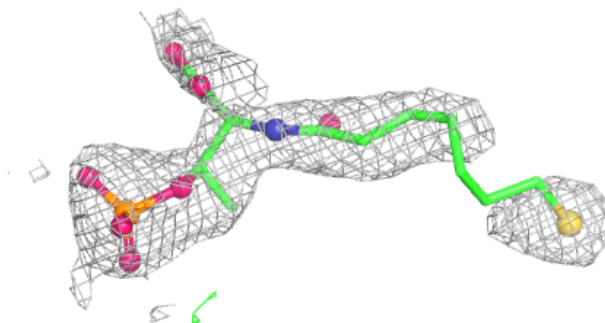
Electron density around XE E 501 (A):

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



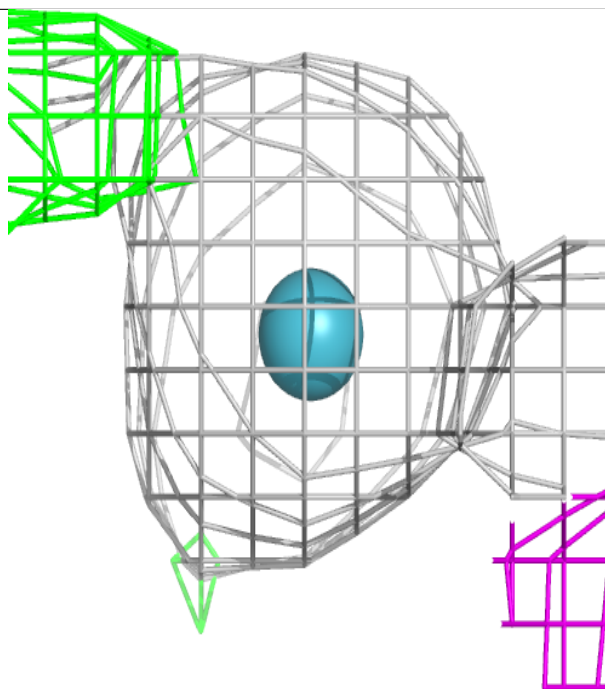
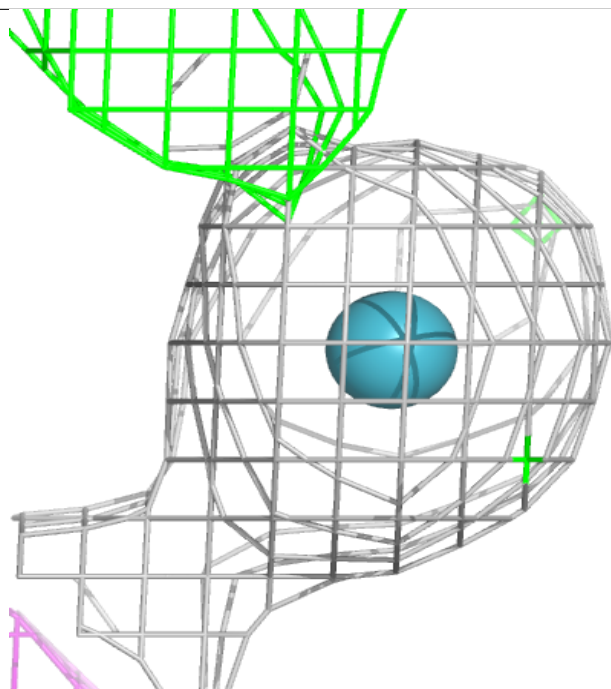
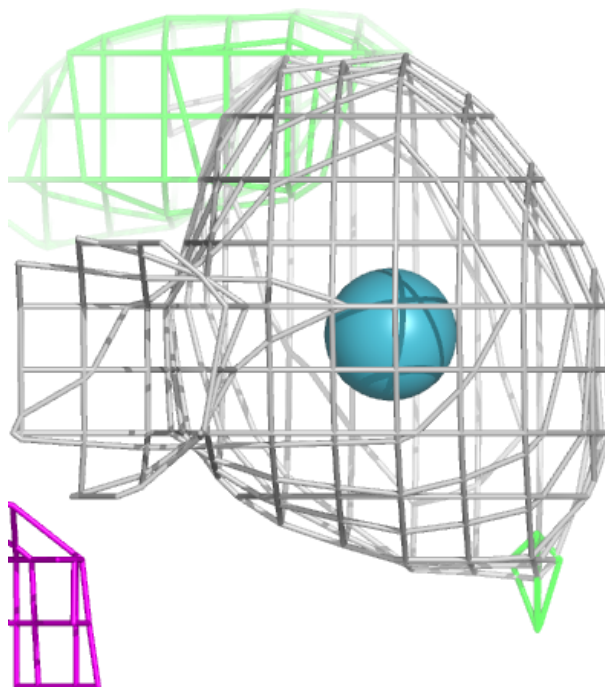
Electron density around TP7 A 604 (A):

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



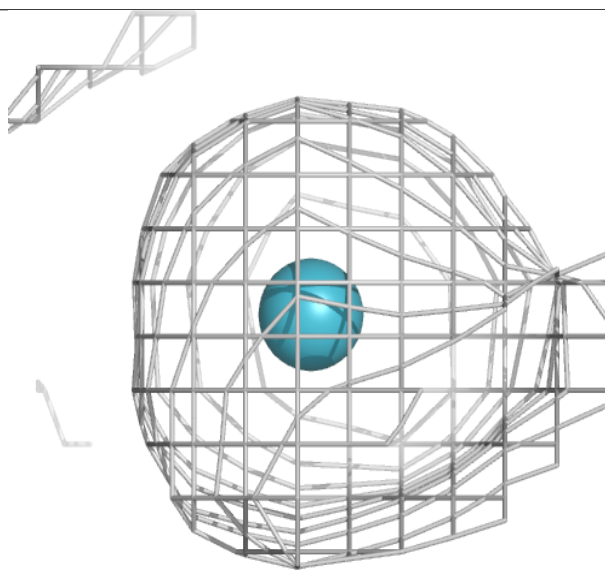
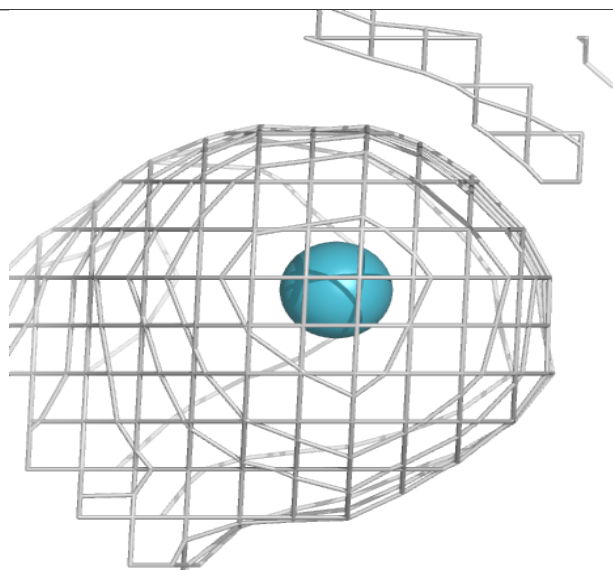
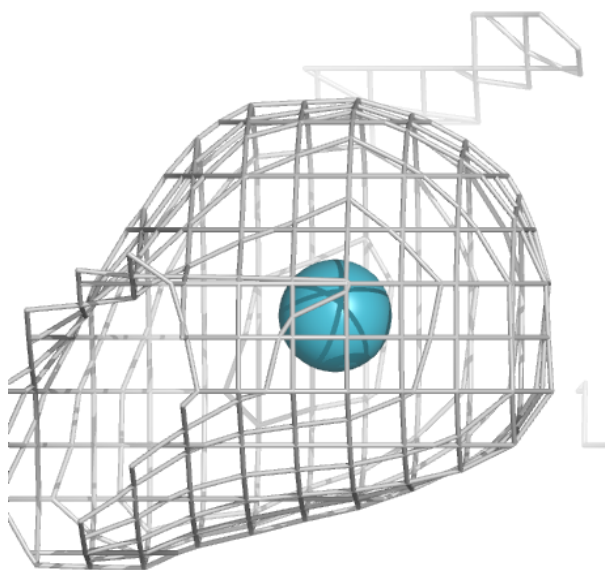
Electron density around XE B 501 (A):

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



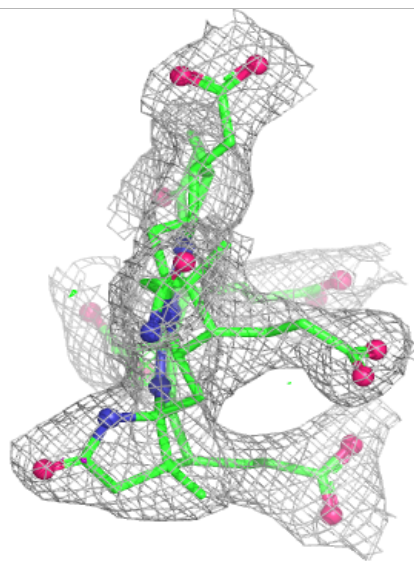
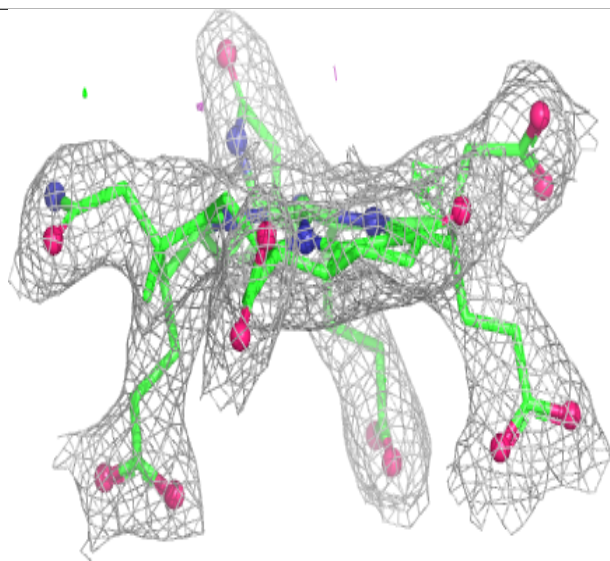
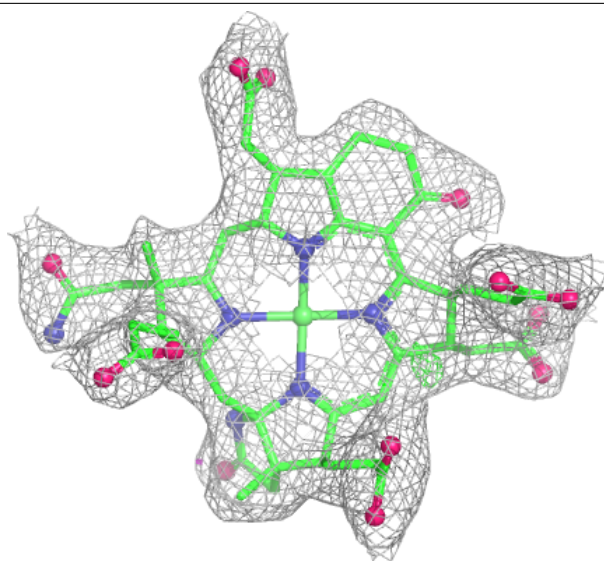
Electron density around XE B 502 (B):

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



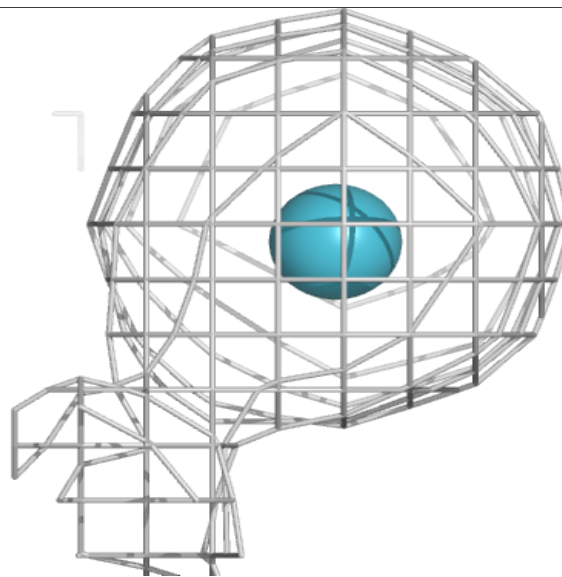
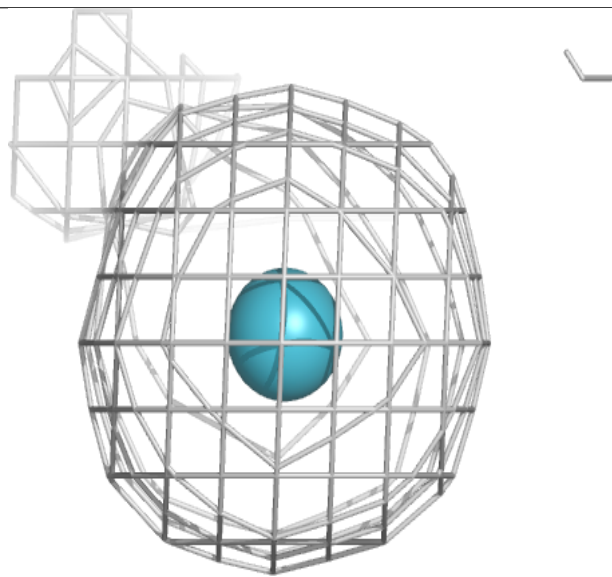
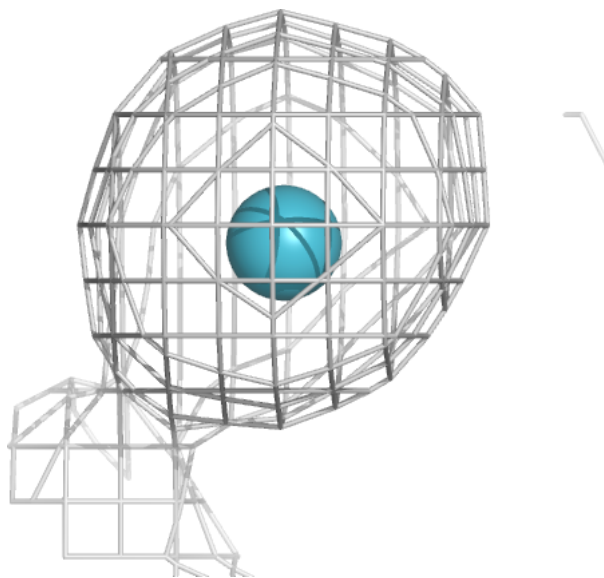
Electron density around F43 A 601:

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



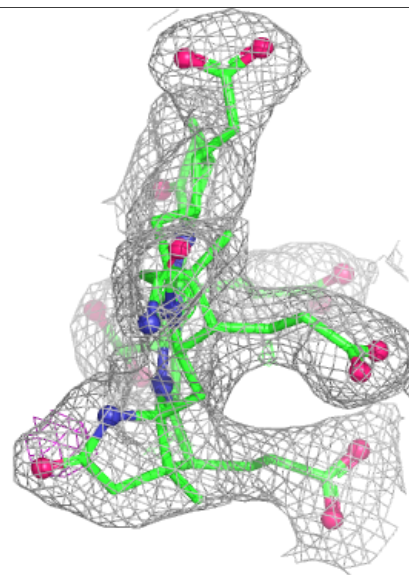
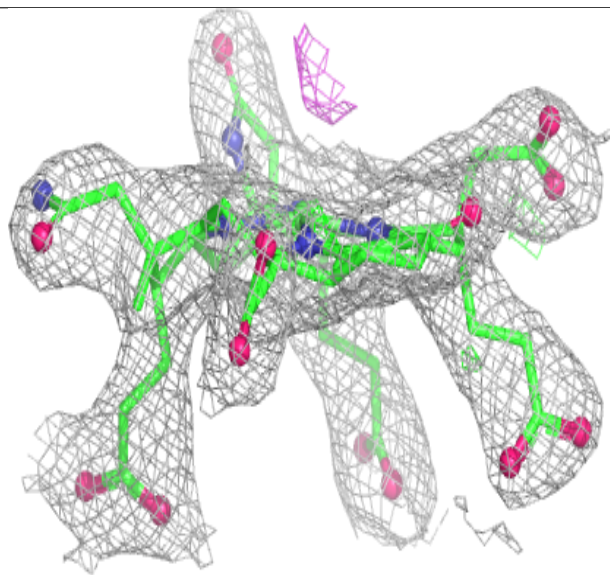
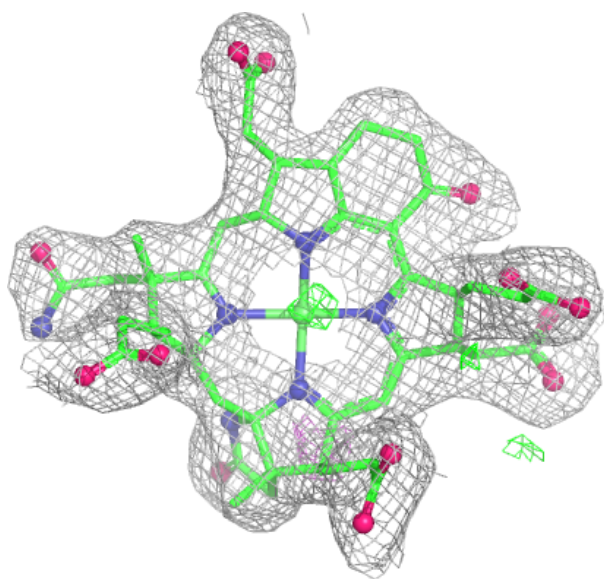
Electron density around XE D 603:

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



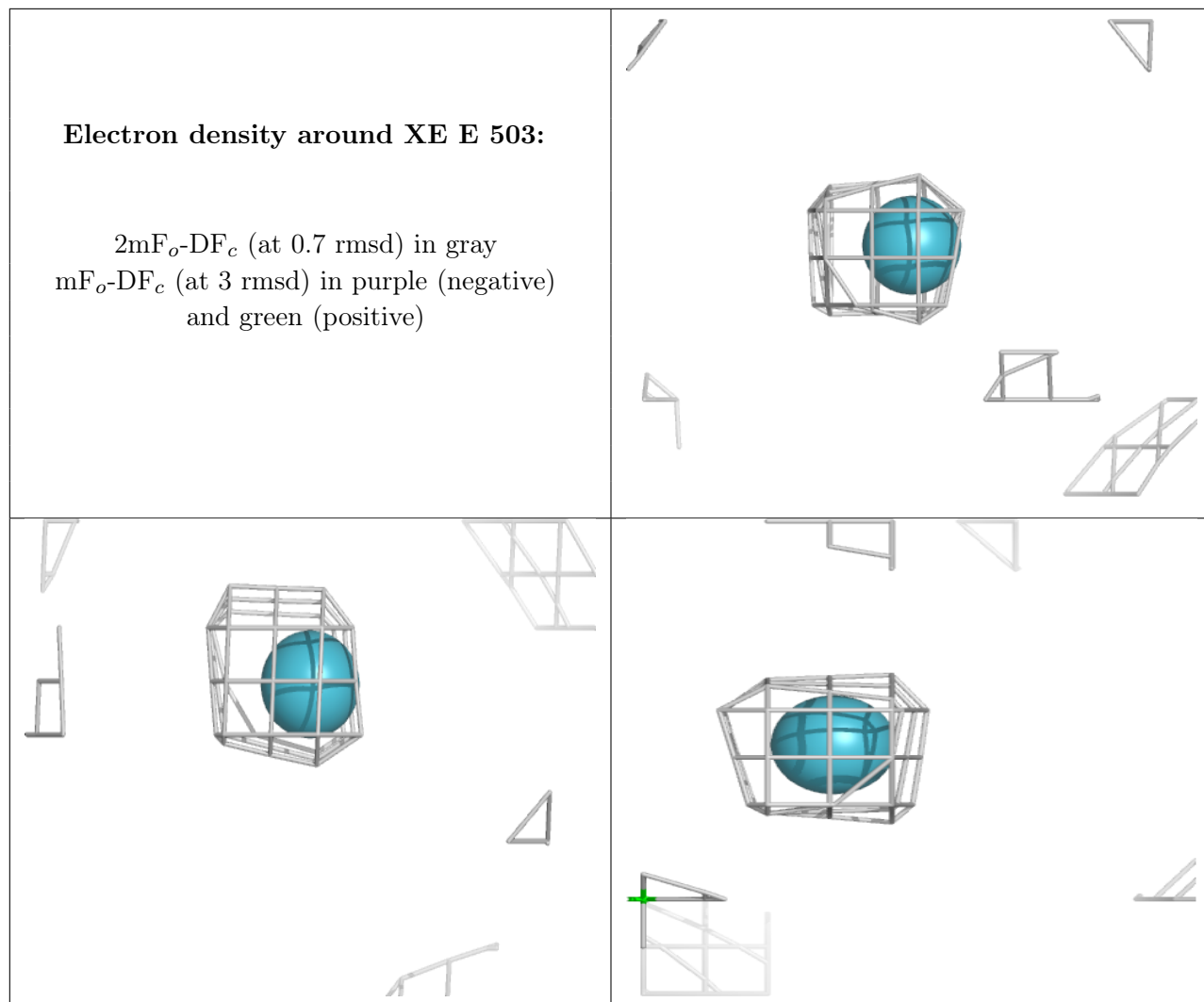
Electron density around F43 A 608:

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



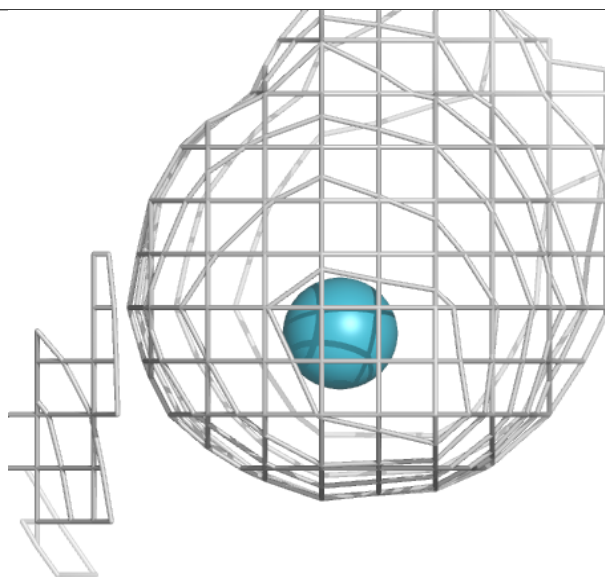
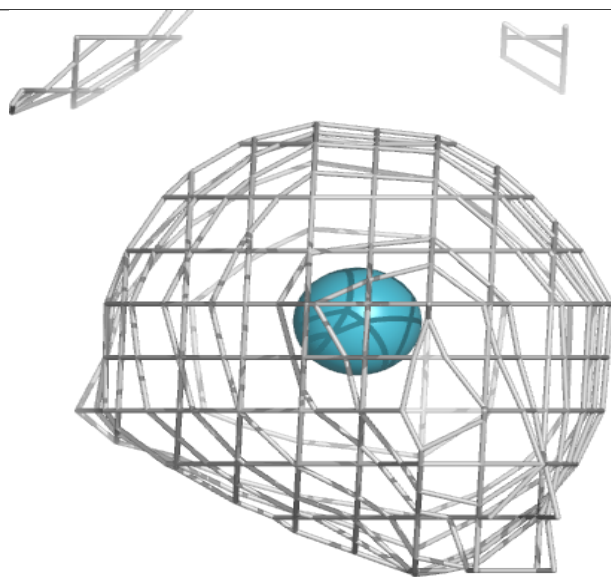
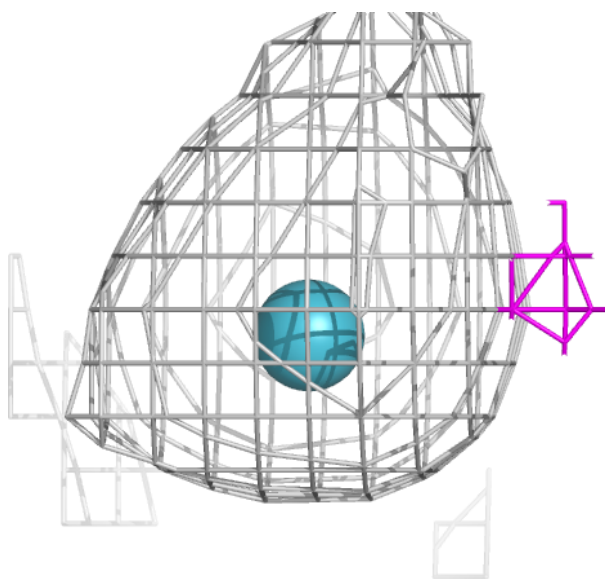
Electron density around XE E 503:

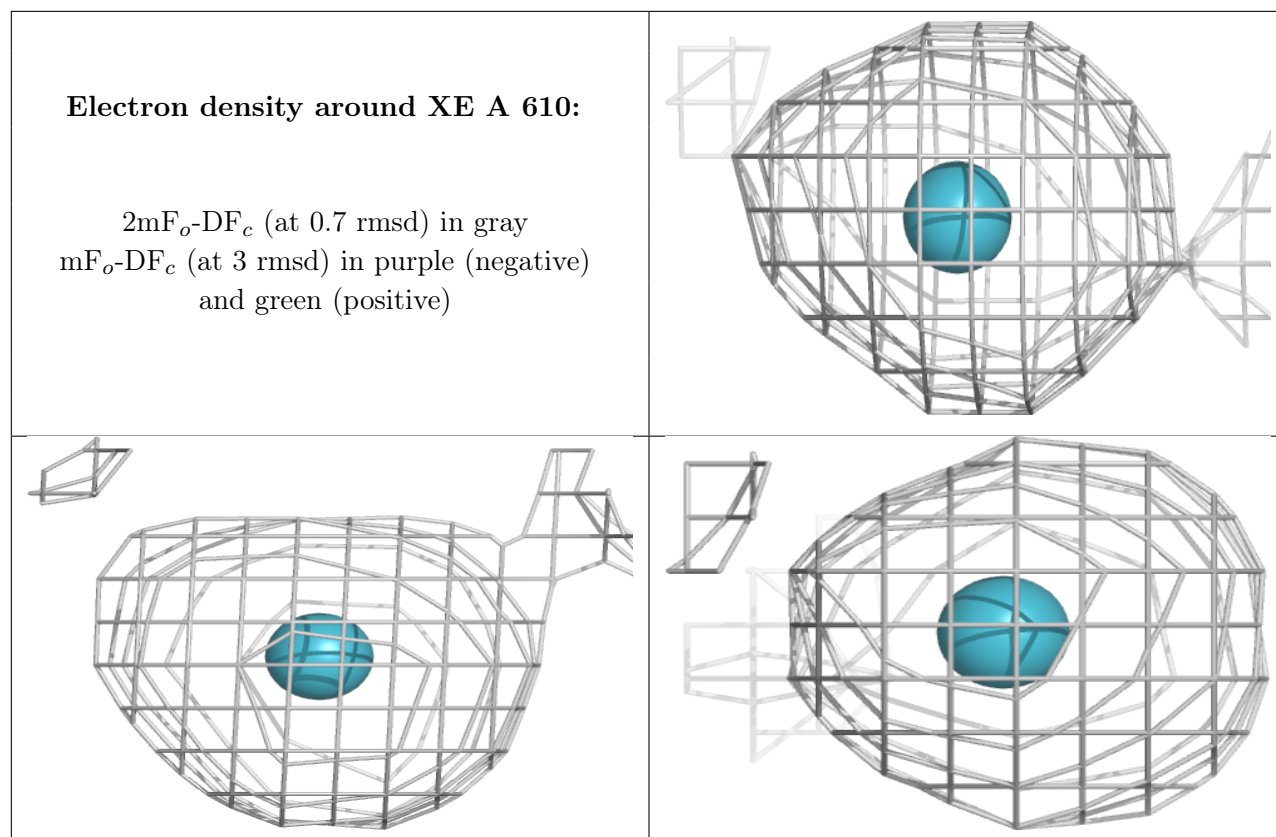
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around XE E 502 (B):

$2mF_o-DF_c$ (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.