



Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 06:36 am BST

PDB ID : 5A8K
Title : METHYL-COENZYME M REDUCTASE FROM METHANOTHERMOBACTER WOLFEII AT 1.4 Å RESOLUTION
Authors : Wagner, T.; Ermler, U.
Deposited on : 2015-07-16
Resolution : 1.41 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

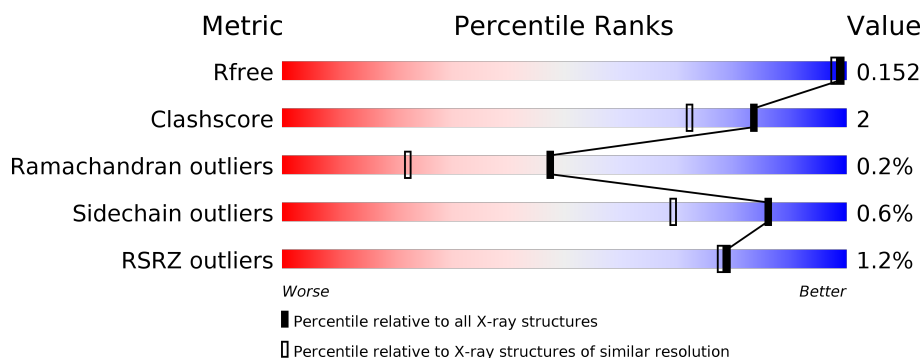
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.41 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	2579 (1.44-1.40)
Clashscore	141614	2696 (1.44-1.40)
Ramachandran outliers	138981	2632 (1.44-1.40)
Sidechain outliers	138945	2631 (1.44-1.40)
RSRZ outliers	127900	2528 (1.44-1.40)

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

Mol	Chain	Length	Quality of chain
1	A	550	<div> <div>%</div> <div> <div></div> <div>95%</div> <div>.</div> </div> </div>
1	D	550	<div> <div>%</div> <div> <div></div> <div>95%</div> <div>5%</div> </div> </div>
2	B	443	<div> <div>%</div> <div> <div></div> <div>94%</div> <div>6%</div> </div> </div>
2	E	443	<div> <div></div> <div> <div></div> <div>95%</div> <div>.</div> </div> </div>
3	C	249	<div> <div>4%</div> <div> <div></div> <div>95%</div> <div>5%</div> </div> </div>
3	F	249	<div> <div>2%</div> <div> <div></div> <div>98%</div> <div>.</div> </div> </div>

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
9	ETX	A	1551	-	-	X	-
9	ETX	D	1551	-	-	X	-

2 Entry composition

There are 11 unique types of molecules in this entry. The entry contains 22338 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.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	548	Total	C	N	O	S	0	15	0
			4327	2751	716	840	20			
1	D	548	Total	C	N	O	S	0	16	0
			4332	2757	716	839	20			

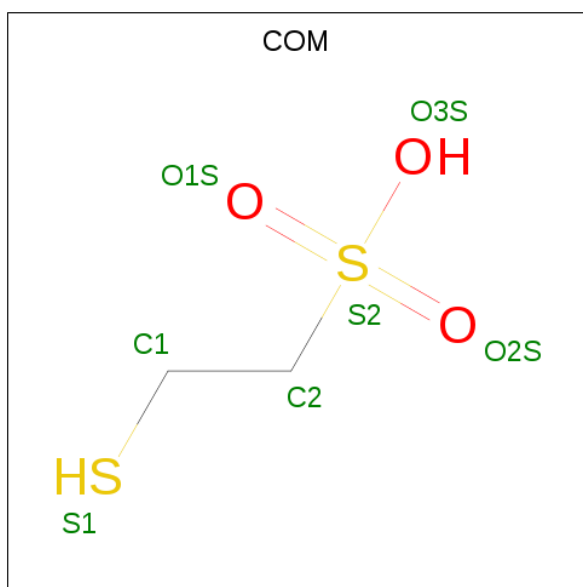
- Molecule 2 is a protein called METHYL-COENZYME M REDUCTASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	442	Total	C	N	O	S	0	11	0
			3371	2140	553	656	22			
2	E	442	Total	C	N	O	S	0	15	0
			3391	2151	555	666	19			

- Molecule 3 is a protein called METHYL-COENZYME M REDUCTASE.

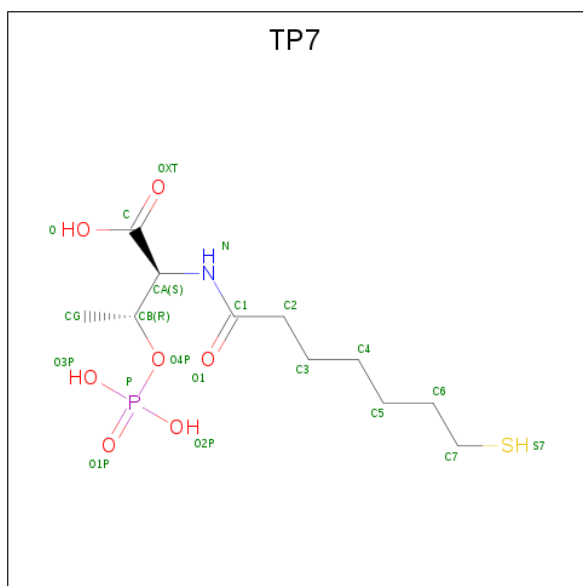
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	248	Total	C	N	O	S	0	8	0
			2049	1274	358	405	12			
3	F	248	Total	C	N	O	S	0	8	0
			2052	1276	358	406	12			

- Molecule 4 is 1-THIOETHANESULFONIC ACID (three-letter code: COM) (formula: C₂H₆O₃S₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	O	S	0	0
			7	2	3	2		
4	D	1	Total	C	O	S	0	0
			7	2	3	2		

- Molecule 5 is Coenzyme B (three-letter code: TP7) (formula: $C_{11}H_{22}NO_7PS$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
5	A	1	Total	C	N	O	P	S	0	0
			21	11	1	7	1	1		
5	D	1	Total	C	N	O	P	S	0	0
			21	11	1	7	1	1		

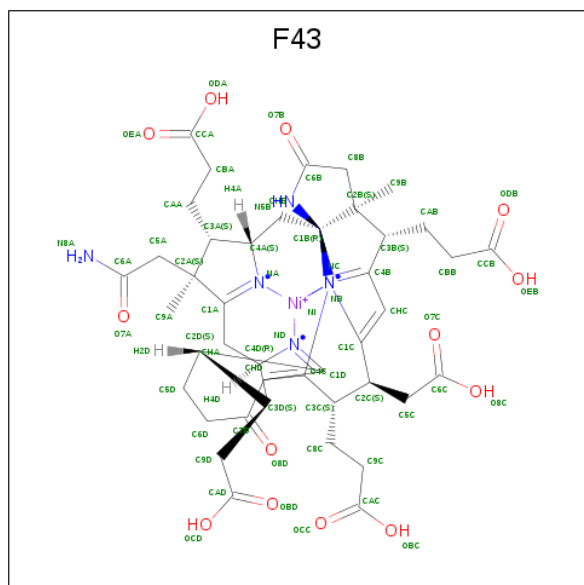
- Molecule 6 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	2	Total K 2 2	0	0
6	D	1	Total K 1 1	0	0
6	C	1	Total K 1 1	0	0
6	E	1	Total K 1 1	0	0

- Molecule 7 is CALCIUM ION (three-letter code: CA) (formula: Ca).

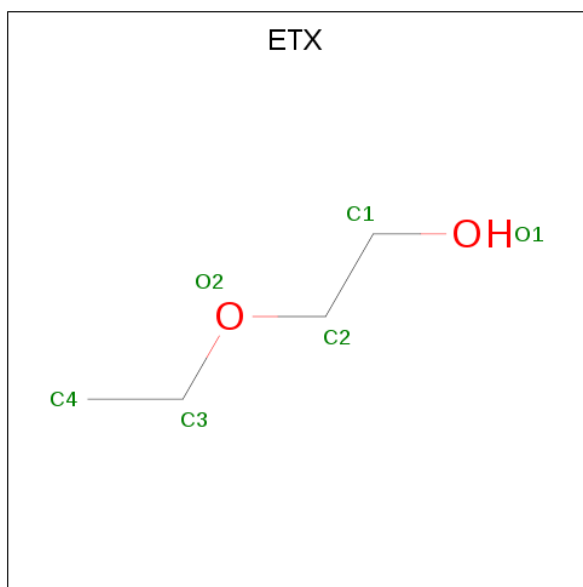
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	B	1	Total Ca 1 1	0	0
7	A	4	Total Ca 4 4	0	0
7	D	5	Total Ca 5 5	0	0
7	F	1	Total Ca 1 1	0	0
7	E	1	Total Ca 1 1	0	0

- Molecule 8 is FACTOR 430 (three-letter code: F43) (formula: $C_{42}H_{51}N_6NiO_{13}$).



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

- Molecule 9 is 2-ETHOXYETHANOL (three-letter code: ETX) (formula: C₄H₁₀O₂).



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

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	C	1	Total C O 6 4 2	0	0
9	D	1	Total C O 6 4 2	0	0
9	D	1	Total C O 6 4 2	0	0
9	D	1	Total C O 6 4 2	0	0
9	E	1	Total C O 6 4 2	0	0
9	E	1	Total C O 6 4 2	0	0
9	E	1	Total C O 6 4 2	0	0
9	E	1	Total C O 6 4 2	0	0
9	E	1	Total C O 6 4 2	0	0
9	E	1	Total C O 6 4 2	0	0
9	E	1	Total C O 6 4 2	0	0
9	E	1	Total C O 6 4 2	0	0
9	E	1	Total C O 6 4 2	0	0
9	F	1	Total C O 6 4 2	0	0
9	F	1	Total C O 6 4 2	0	0
9	F	1	Total C O 6 4 2	0	0

- Molecule 10 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

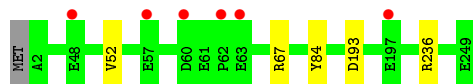
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
10	F	1	Total Mg 1 1	0	0

- Molecule 11 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	A	586	Total 586	O 586	0	0
11	B	445	Total 445	O 445	0	0
11	C	273	Total 273	O 273	0	0
11	D	489	Total 489	O 489	0	0
11	E	418	Total 418	O 418	0	0
11	F	251	Total 251	O 251	0	0

- Molecule 3: METHYL-COENZYME M REDUCTASE

Chain F: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	84.14Å 150.54Å 187.09Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.47 – 1.41 48.47 – 1.41	Depositor EDS
% Data completeness (in resolution range)	99.9 (48.47-1.41) 99.9 (48.47-1.41)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.44 (at 1.41Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.114 , 0.140 0.130 , 0.152	Depositor DCC
R_{free} test set	22742 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	10.4	Xtriage
Anisotropy	0.302	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 40.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	22338	wwPDB-VP
Average B, all atoms (Å ²)	13.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.51% 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: MG, COM, ETX, AGM, F43, MGN, TP7, K, GL3, SMC, MHS, CA

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	0.47	0/4420	0.66	0/6004
1	D	0.47	0/4430	0.67	1/6017 (0.0%)
2	B	0.44	0/3458	0.63	0/4678
2	E	0.45	0/3487	0.64	0/4718
3	C	0.47	0/2110	0.70	1/2841 (0.0%)
3	F	0.46	0/2114	0.68	0/2846
All	All	0.46	0/20019	0.66	2/27104 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	120	ARG	NE-CZ-NH1	5.85	123.23	120.30
1	D	36	ARG	NE-CZ-NH2	-5.17	117.72	120.30

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	4327	0	4180	18	0
1	D	4332	0	4183	21	0
2	B	3371	0	3402	29	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	E	3391	0	3415	18	0
3	C	2049	0	1995	8	0
3	F	2052	0	1999	7	0
4	A	7	0	5	1	0
4	D	7	0	5	1	0
5	A	21	0	19	0	0
5	D	21	0	19	0	0
6	A	2	0	0	0	0
6	C	1	0	0	0	0
6	D	1	0	0	0	0
6	E	1	0	0	0	0
7	A	4	0	0	0	0
7	B	1	0	0	0	0
7	D	5	0	0	0	0
7	E	1	0	0	0	0
7	F	1	0	0	0	0
8	A	62	0	43	1	0
8	D	62	0	43	1	0
9	A	30	0	50	6	0
9	B	30	0	50	6	0
9	C	6	0	10	0	0
9	D	18	0	29	7	0
9	E	54	0	90	4	0
9	F	18	0	30	1	0
10	F	1	0	0	0	0
11	A	586	0	0	10	0
11	B	445	0	0	10	0
11	C	273	0	0	3	0
11	D	489	0	0	9	1
11	E	418	0	0	6	1
11	F	251	0	0	1	0
All	All	22338	0	19567	96	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 2.

All (96) 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)
11:D:2436:HOH:O	2:E:351[B]:LYS:NZ	1.88	1.07
11:A:2511:HOH:O	2:B:351[B]:LYS:NZ	1.88	1.01

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:14:ARG:NH1	3:C:63:GLU:OE2	2.01	0.93
2:E:97[A]:GLU:OE1	11:E:2161:HOH:O	1.87	0.92
2:B:40[C]:GLN:NE2	11:B:2076:HOH:O	2.03	0.91
1:A:4[B]:LYS:NZ	11:A:2012:HOH:O	2.04	0.91
2:E:147[B]:ASP:OD1	11:E:2206:HOH:O	1.99	0.80
2:E:274:GLU:O	2:E:278[A]:GLU:HG3	1.81	0.79
1:D:42[B]:ASN:ND2	11:D:2120:HOH:O	2.14	0.78
1:D:241[B]:GLN:NE2	11:D:2298:HOH:O	2.17	0.77
2:E:91[B]:GLU:OE1	11:E:2163:HOH:O	2.02	0.77
1:D:522:PHE:H	9:D:1551:ETX:H41	1.54	0.72
1:D:123:GLU:OE1	11:D:2196:HOH:O	2.07	0.71
3:C:27:LYS:NZ	3:C:30:GLU:OE1	2.24	0.71
11:D:2435:HOH:O	2:E:351[A]:LYS:NZ	2.13	0.71
1:A:67:ARG:NH2	11:A:2151:HOH:O	2.25	0.68
11:A:2510:HOH:O	2:B:351[A]:LYS:NZ	2.06	0.66
2:E:91[B]:GLU:OE2	11:E:2162:HOH:O	2.15	0.64
3:C:60:ASP:OD1	11:C:2102:HOH:O	2.15	0.64
3:C:83:ARG:NH2	11:C:2148:HOH:O	2.31	0.63
1:A:42[B]:ASN:OD1	11:A:2103:HOH:O	2.16	0.62
1:A:521:ASP:HA	9:A:1551:ETX:H31	1.83	0.60
1:A:45:LYS:NZ	11:A:2109:HOH:O	2.35	0.59
2:B:40[A]:GLN:OE1	11:B:2080:HOH:O	2.16	0.58
2:B:40[B]:GLN:OE1	11:B:2080:HOH:O	2.17	0.58
2:E:77:ASP:OD1	11:E:2096:HOH:O	2.17	0.57
3:F:193:ASP:OD2	11:F:2196:HOH:O	2.17	0.57
9:D:1551:ETX:H32	11:D:2331:HOH:O	2.03	0.57
1:D:521:ASP:HA	9:D:1551:ETX:H31	1.87	0.57
1:A:352[A]:GLU:HG2	11:A:2427:HOH:O	2.05	0.56
11:A:2559:HOH:O	1:D:548:PRO:O	2.18	0.55
2:B:27:LEU:HD22	2:B:246[B]:MET:SD	2.47	0.54
2:B:405:MET:SD	9:B:1445:ETX:H43	2.48	0.54
1:A:522:PHE:H	9:A:1551:ETX:H41	1.73	0.53
9:E:1447:ETX:H42	3:F:52:VAL:CG2	2.38	0.53
9:A:1551:ETX:H32	11:A:2394:HOH:O	2.07	0.53
2:B:77:ASP:OD1	11:B:2103:HOH:O	2.18	0.52
1:A:240:LYS:HE2	3:F:193:ASP:OD2	2.10	0.51
2:B:86:ALA:HB3	9:B:1447:ETX:H43	1.93	0.51
2:B:246[B]:MET:HE1	2:B:432:VAL:HG12	1.92	0.51
2:E:322:ARG:NH2	3:F:67[B]:ARG:HH11	2.10	0.50
1:A:328:VAL:HB	8:A:1550:F43:H9A1	1.93	0.50
2:B:121:PHE:CB	9:B:1444:ETX:H22	2.42	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:147:ARG:NH2	3:C:207:GLU:OE2	2.46	0.49
2:B:65:PRO:HB3	9:D:1553:ETX:H42	1.93	0.49
9:B:1444:ETX:H21	11:B:2079:HOH:O	2.12	0.49
3:C:84:TYR:CE1	1:D:242:ALA:HB2	2.48	0.49
2:B:121:PHE:HB3	9:B:1444:ETX:H22	1.94	0.48
2:B:438:GLU:OE2	11:B:2440:HOH:O	2.20	0.48
3:C:185:GLU:OE2	11:C:2207:HOH:O	2.20	0.48
1:D:522:PHE:H	9:D:1551:ETX:C4	2.24	0.48
2:B:230:VAL:HG23	11:B:2314:HOH:O	2.13	0.48
2:B:224:PHE:HA	11:B:2314:HOH:O	2.14	0.48
1:A:528:GLU:OE2	9:A:1555:ETX:H43	2.15	0.47
1:D:328:VAL:HB	8:D:1550:F43:H9A1	1.96	0.47
1:A:242:ALA:HB2	3:F:84:TYR:CE1	2.51	0.45
2:B:188:PRO:HD3	11:E:2226:HOH:O	2.17	0.45
1:D:268[B]:PRO:HG3	11:D:2448:HOH:O	2.16	0.45
1:D:182:ILE:HD11	1:D:207[B]:VAL:HG11	1.97	0.45
3:C:27:LYS:CE	3:C:30:GLU:OE1	2.66	0.44
2:B:36:LYS:O	2:B:40[C]:GLN:HG2	2.18	0.44
1:D:352:GLU:HG2	11:D:2363:HOH:O	2.17	0.44
2:B:27:LEU:HD22	2:B:246[B]:MET:CE	2.49	0.43
11:B:2261:HOH:O	2:E:188:PRO:HD3	2.18	0.43
2:E:246:MET:HE1	2:E:432:VAL:HB	2.01	0.43
1:A:531:LYS:HZ3	9:A:1555:ETX:C3	2.31	0.43
1:A:522:PHE:HB2	9:A:1551:ETX:H41	2.00	0.43
1:D:123:GLU:N	1:D:123:GLU:OE1	2.48	0.43
1:D:268[B]:PRO:CG	11:D:2448:HOH:O	2.66	0.42
2:B:178[B]:MET:HG2	2:B:195:ILE:HG23	2.00	0.42
2:B:232:ALA:HB2	9:E:1452:ETX:H12	2.01	0.42
1:D:443:PHE:HB2	4:D:555:COM:H11	2.02	0.42
1:A:332:GLN:HA	1:A:335:THR:OG1	2.19	0.42
2:B:126:GLU:HB3	2:E:126:GLU:HB3	2.01	0.42
1:D:521:ASP:OD1	9:D:1551:ETX:H31	2.19	0.42
1:D:97:LEU:C	1:D:97:LEU:HD23	2.40	0.42
1:D:453:GLY:O	1:D:457:VAL:HG23	2.20	0.41
1:A:182:ILE:HD11	1:A:207[B]:VAL:HG11	2.01	0.41
2:B:246[A]:MET:CE	2:B:429:LEU:HD12	2.49	0.41
2:B:362:PHE:O	2:B:369:GLY:HA3	2.20	0.41
1:A:507:PRO:HD2	2:E:66:ALA:O	2.20	0.41
1:A:127:GLU:HG3	11:A:2215:HOH:O	2.20	0.41
1:D:522:PHE:HB2	9:D:1551:ETX:H41	2.02	0.41
3:F:236:ARG:HG3	9:F:1252:ETX:H42	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:21:GLN:OE1	9:B:1448:ETX:H43	2.21	0.41
1:D:154:HIS:NE2	1:D:545[A]:LEU:HD21	2.36	0.41
2:E:140:GLN:OE1	9:E:1452:ETX:H11	2.21	0.41
2:E:214:LEU:HB2	2:E:428:PRO:HG3	2.02	0.41
9:E:1452:ETX:H42	9:E:1452:ETX:H11	2.03	0.40
1:D:73:GLN:HB2	1:D:80:TYR:CE2	2.57	0.40
2:B:123:VAL:HG12	2:E:36:LYS:HA	2.04	0.40
1:A:443:PHE:HB2	4:A:556:COM:H11	2.04	0.40
2:B:113:LEU:C	2:B:113:LEU:HD23	2.42	0.40
2:B:58:LYS:HE2	11:B:2104:HOH:O	2.21	0.40
2:E:322:ARG:CZ	3:F:67[A]:ARG:HD2	2.52	0.40
2:E:362:PHE:O	2:E:369:GLY:HA3	2.22	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:D:2251:HOH:O	11:E:2389:HOH:O[1_655]	2.19	0.01

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	556/550 (101%)	536 (96%)	18 (3%)	2 (0%)	34	12
1	D	557/550 (101%)	538 (97%)	17 (3%)	2 (0%)	34	12
2	B	452/443 (102%)	445 (98%)	7 (2%)	0	100	100
2	E	455/443 (103%)	448 (98%)	7 (2%)	0	100	100
3	C	253/249 (102%)	245 (97%)	8 (3%)	0	100	100
3	F	254/249 (102%)	249 (98%)	5 (2%)	0	100	100
All	All	2527/2484 (102%)	2461 (97%)	62 (2%)	4 (0%)	47	22

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	325	SER
1	D	325	SER
1	D	63	PRO
1	A	63	PRO

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	449/436 (103%)	446 (99%)	3 (1%)	84	65
1	D	450/436 (103%)	446 (99%)	4 (1%)	78	56
2	B	354/343 (103%)	353 (100%)	1 (0%)	92	81
2	E	357/343 (104%)	355 (99%)	2 (1%)	86	69
3	C	224/217 (103%)	221 (99%)	3 (1%)	69	41
3	F	224/217 (103%)	224 (100%)	0	100	100
All	All	2058/1992 (103%)	2045 (99%)	13 (1%)	86	69

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	LYS
1	A	444	TYR
1	A	543	ARG
2	B	126	GLU
3	C	19[A]	CYS
3	C	19[B]	CYS
3	C	211	MET
1	D	360[A]	LEU
1	D	360[B]	LEU
1	D	444	TYR
1	D	543	ARG
2	E	126	GLU
2	E	178	MET

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

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

10 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	MGN	A	400	1	6,9,10	0.67	0	5,12,14	0.17	0
1	MGN	D	400	1	6,9,10	0.82	0	5,12,14	0.19	0
1	AGM	D	271	1	10,11,12	1.14	1 (10%)	6,13,15	1.07	0
1	AGM	A	271	1	10,11,12	1.54	3 (30%)	6,13,15	0.80	0
1	SMC	A	452	1	5,6,7	0.92	0	2,6,8	1.56	1 (50%)
1	MHS	D	257	1	7,11,12	1.11	0	6,14,16	1.51	1 (16%)
1	MHS	A	257	1	7,11,12	1.20	0	6,14,16	1.15	0
1	GL3	A	445	1	2,3,4	3.61	1 (50%)	1,2,4	0.24	0
1	GL3	D	445	1	2,3,4	4.16	1 (50%)	1,2,4	0.45	0
1	SMC	D	452	1	5,6,7	0.61	0	2,6,8	1.54	1 (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	MGN	A	400	1	-	0/7/9/12	-
1	MGN	D	400	1	-	0/7/9/12	-
1	AGM	D	271	1	-	3/10/11/13	-
1	AGM	A	271	1	-	2/10/11/13	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	SMC	A	452	1	-	1/3/5/7	-
1	MHS	D	257	1	-	0/5/6/8	0/1/1/1
1	MHS	A	257	1	-	0/5/6/8	0/1/1/1
1	GL3	A	445	1	-	0/1/1/2	-
1	GL3	D	445	1	-	0/1/1/2	-
1	SMC	D	452	1	-	1/3/5/7	-

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	445	GL3	C-S	-5.88	1.60	1.80
1	A	445	GL3	C-S	-5.10	1.63	1.80
1	A	271	AGM	CG-CD	2.64	1.57	1.53
1	A	271	AGM	CB-CA	2.50	1.56	1.53
1	A	271	AGM	CZ-NE1	2.45	1.37	1.33
1	D	271	AGM	CZ-NE1	2.12	1.36	1.33

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	257	MHS	CB-CA-C	-2.43	106.91	111.47
1	A	452	SMC	CA-CB-SG	-2.17	110.53	114.04
1	D	452	SMC	CA-CB-SG	-2.08	110.67	114.04

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	452	SMC	CA-CB-SG-CS
1	D	452	SMC	CA-CB-SG-CS
1	A	271	AGM	CE2-CD-NE1-CZ
1	D	271	AGM	CE2-CD-NE1-CZ
1	D	271	AGM	CE2-CD-CG-CB
1	D	271	AGM	NE1-CD-CG-CB
1	A	271	AGM	NE1-CD-CG-CB

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

Of 50 ligands modelled in this entry, 18 are monoatomic - leaving 32 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
9	ETX	F	1251	-	5,5,5	0.39	0	4,4,4	0.39	0
9	ETX	E	1447	-	5,5,5	0.36	0	4,4,4	0.41	0
9	ETX	E	1450	-	5,5,5	0.39	0	4,4,4	0.56	0
9	ETX	A	1554	-	5,5,5	0.60	0	4,4,4	0.58	0
9	ETX	D	1552	6	5,5,5	0.45	0	4,4,4	0.36	0
9	ETX	E	1448	-	5,5,5	0.38	0	4,4,4	0.77	0
4	COM	D	555	8	6,6,6	1.39	1 (16%)	7,8,8	2.11	1 (14%)
9	ETX	D	1551	-	5,5,5	0.62	0	4,4,4	0.53	0
9	ETX	B	1445	-	5,5,5	0.40	0	4,4,4	0.09	0
9	ETX	D	1553	-	5,5,5	0.41	0	4,4,4	0.42	0
9	ETX	F	1250	-	5,5,5	0.33	0	4,4,4	0.34	0
5	TP7	D	554	-	16,20,20	0.54	0	18,26,26	0.77	0
9	ETX	B	1447	-	5,5,5	0.36	0	4,4,4	0.57	0
4	COM	A	556	8	6,6,6	1.70	2 (33%)	7,8,8	2.87	3 (42%)
9	ETX	B	1448	-	5,5,5	0.52	0	4,4,4	0.61	0
9	ETX	A	1552	-	5,5,5	0.44	0	4,4,4	0.59	0
9	ETX	B	1446	-	5,5,5	0.36	0	4,4,4	0.61	0
9	ETX	C	1250	-	5,5,5	0.39	0	4,4,4	0.44	0
9	ETX	A	1555	-	5,5,5	0.30	0	4,4,4	1.07	0
9	ETX	E	1452	-	5,5,5	0.37	0	4,4,4	0.70	0
9	ETX	A	1551	-	5,5,5	0.55	0	4,4,4	0.53	0
8	F43	D	1550	1,4	46,71,71	2.48	5 (10%)	48,118,118	1.18	4 (8%)
9	ETX	F	1252	-	5,5,5	0.41	0	4,4,4	0.57	0
9	ETX	E	1451	-	5,5,5	0.44	0	4,4,4	0.36	0
9	ETX	E	1449	-	5,5,5	0.41	0	4,4,4	0.36	0
9	ETX	B	1444	-	5,5,5	0.44	0	4,4,4	0.74	0
9	ETX	A	1553	-	5,5,5	0.48	0	4,4,4	0.59	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
9	ETX	E	1445	-	5,5,5	0.36	0	4,4,4	0.37	0
5	TP7	A	558	-	16,20,20	0.51	0	18,26,26	0.69	0
9	ETX	E	1444	-	5,5,5	0.35	0	4,4,4	0.61	0
8	F43	A	1550	1,4	46,71,71	2.37	5 (10%)	48,118,118	1.15	4 (8%)
9	ETX	E	1446	-	5,5,5	0.45	0	4,4,4	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
9	ETX	F	1251	-	-	1/3/3/3	-
9	ETX	E	1447	-	-	2/3/3/3	-
9	ETX	E	1450	-	-	1/3/3/3	-
9	ETX	A	1554	-	-	2/3/3/3	-
9	ETX	D	1552	6	-	1/3/3/3	-
9	ETX	E	1448	-	-	0/3/3/3	-
4	COM	D	555	8	-	0/4/4/4	-
9	ETX	D	1551	-	-	1/3/3/3	-
9	ETX	B	1445	-	-	3/3/3/3	-
9	ETX	D	1553	-	-	2/3/3/3	-
9	ETX	F	1250	-	-	2/3/3/3	-
5	TP7	D	554	-	-	0/20/24/24	-
9	ETX	B	1447	-	-	3/3/3/3	-
4	COM	A	556	8	-	0/4/4/4	-
9	ETX	B	1448	-	-	1/3/3/3	-
9	ETX	A	1552	-	-	1/3/3/3	-
9	ETX	B	1446	-	-	0/3/3/3	-
9	ETX	C	1250	-	-	0/3/3/3	-
9	ETX	A	1555	-	-	3/3/3/3	-
9	ETX	E	1452	-	-	2/3/3/3	-
9	ETX	A	1551	-	-	1/3/3/3	-
8	F43	D	1550	1,4	-	1/18/185/185	-
9	ETX	F	1252	-	-	1/3/3/3	-
9	ETX	E	1451	-	-	3/3/3/3	-
9	ETX	E	1449	-	-	1/3/3/3	-
9	ETX	B	1444	-	-	2/3/3/3	-
9	ETX	A	1553	-	-	0/3/3/3	-
9	ETX	E	1445	-	-	3/3/3/3	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	TP7	A	558	-	-	1/20/24/24	-
9	ETX	E	1444	-	-	0/3/3/3	-
8	F43	A	1550	1,4	-	1/18/185/185	-
9	ETX	E	1446	-	-	1/3/3/3	-

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	D	1550	F43	NI-NA	9.81	2.10	1.89
8	A	1550	F43	NI-NA	9.03	2.09	1.89
8	D	1550	F43	NI-ND	8.12	2.07	1.89
8	D	1550	F43	NI-NB	8.08	2.06	1.89
8	A	1550	F43	NI-NB	7.52	2.05	1.89
8	A	1550	F43	NI-ND	7.51	2.05	1.89
8	A	1550	F43	CHD-C1D	-5.31	1.36	1.43
8	D	1550	F43	CHD-C1D	-4.49	1.37	1.43
4	A	556	COM	O1S-S2	2.99	1.53	1.45
8	D	1550	F43	CHC-C4B	2.90	1.47	1.39
8	A	1550	F43	CHC-C4B	2.59	1.46	1.39
4	A	556	COM	O2S-S2	2.46	1.52	1.45
4	D	555	COM	O2S-S2	2.17	1.51	1.45

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	556	COM	O2S-S2-C2	5.59	113.64	106.92
4	D	555	COM	O2S-S2-C2	4.38	112.19	106.92
8	D	1550	F43	C2B-C1B-NB	2.92	106.21	101.84
4	A	556	COM	C2-C1-S1	-2.86	105.86	113.10
4	A	556	COM	O3S-S2-O2S	-2.67	104.75	111.27
8	D	1550	F43	O8D-C7D-C6D	-2.60	116.61	120.86
8	A	1550	F43	CAB-C3B-C2B	-2.49	113.83	119.09
8	D	1550	F43	CAB-C3B-C2B	-2.45	113.91	119.09
8	A	1550	F43	C2B-C1B-NB	2.43	105.47	101.84
8	D	1550	F43	C4D-ND-C1D	2.35	111.60	108.51
8	A	1550	F43	C4A-NA-C1A	-2.08	106.44	108.97
8	A	1550	F43	C6D-C5D-C2D	-2.04	107.35	111.46

There are no chirality outliers.

All (40) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
9	E	1451	ETX	O1-C1-C2-O2
9	B	1444	ETX	O1-C1-C2-O2
9	E	1445	ETX	O1-C1-C2-O2
8	D	1550	F43	C3A-CAA-CBA-CCA
8	A	1550	F43	C3A-CAA-CBA-CCA
9	B	1444	ETX	C4-C3-O2-C2
9	D	1553	ETX	O1-C1-C2-O2
9	F	1250	ETX	O1-C1-C2-O2
9	D	1553	ETX	C4-C3-O2-C2
9	F	1251	ETX	O1-C1-C2-O2
9	B	1448	ETX	O1-C1-C2-O2
9	E	1452	ETX	O1-C1-C2-O2
9	A	1554	ETX	O1-C1-C2-O2
9	B	1447	ETX	O1-C1-C2-O2
9	E	1451	ETX	C1-C2-O2-C3
9	E	1452	ETX	C4-C3-O2-C2
9	F	1250	ETX	C4-C3-O2-C2
9	E	1446	ETX	C4-C3-O2-C2
9	E	1445	ETX	C4-C3-O2-C2
9	E	1447	ETX	C4-C3-O2-C2
9	A	1554	ETX	C4-C3-O2-C2
9	B	1447	ETX	C1-C2-O2-C3
9	B	1445	ETX	C4-C3-O2-C2
9	E	1445	ETX	C1-C2-O2-C3
9	E	1449	ETX	C4-C3-O2-C2
9	A	1555	ETX	O1-C1-C2-O2
9	B	1445	ETX	O1-C1-C2-O2
9	A	1552	ETX	O1-C1-C2-O2
9	B	1447	ETX	C4-C3-O2-C2
9	A	1555	ETX	C4-C3-O2-C2
9	E	1447	ETX	C1-C2-O2-C3
9	F	1252	ETX	O1-C1-C2-O2
9	E	1451	ETX	C4-C3-O2-C2
9	D	1551	ETX	C4-C3-O2-C2
9	A	1551	ETX	C4-C3-O2-C2
9	E	1450	ETX	C4-C3-O2-C2
5	A	558	TP7	CB-O4P-P-O3P
9	A	1555	ETX	C1-C2-O2-C3
9	D	1552	ETX	O1-C1-C2-O2
9	B	1445	ETX	C1-C2-O2-C3

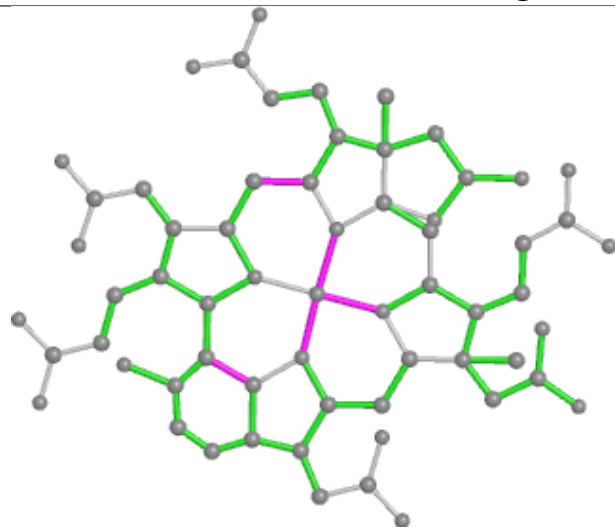
There are no ring outliers.

15 monomers are involved in 28 short contacts:

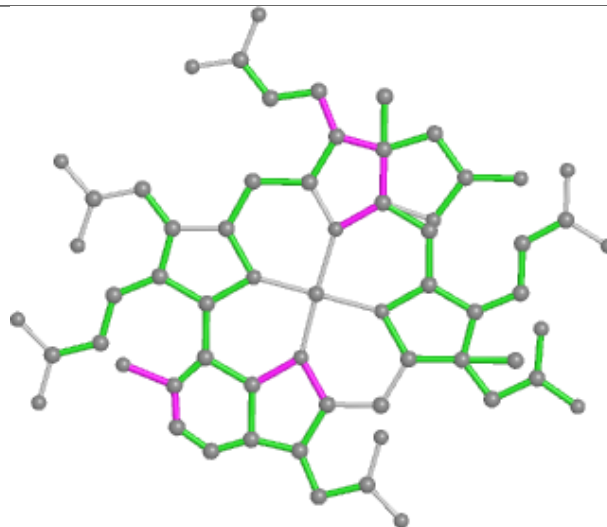
Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	E	1447	ETX	1	0
4	D	555	COM	1	0
9	D	1551	ETX	6	0
9	B	1445	ETX	1	0
9	D	1553	ETX	1	0
9	B	1447	ETX	1	0
4	A	556	COM	1	0
9	B	1448	ETX	1	0
9	A	1555	ETX	2	0
9	E	1452	ETX	3	0
9	A	1551	ETX	4	0
8	D	1550	F43	1	0
9	F	1252	ETX	1	0
9	B	1444	ETX	3	0
8	A	1550	F43	1	0

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

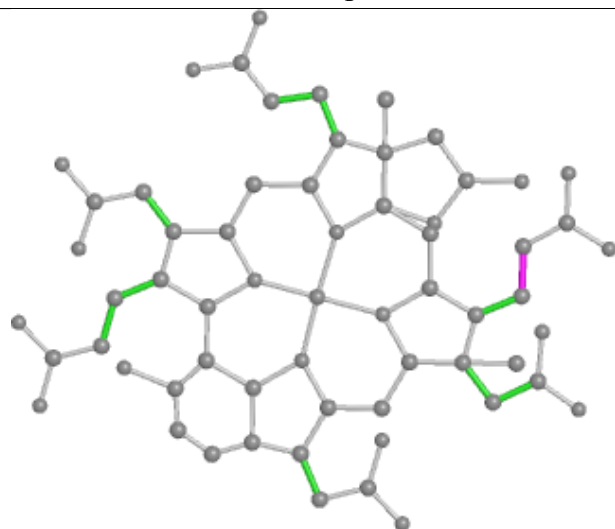
Ligand F43 D 1550



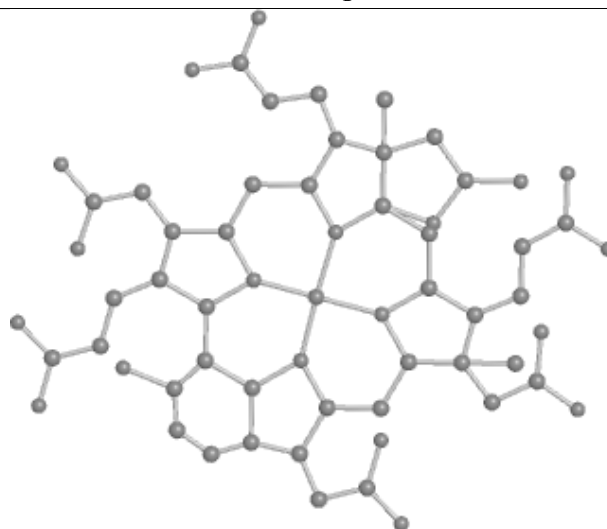
Bond lengths



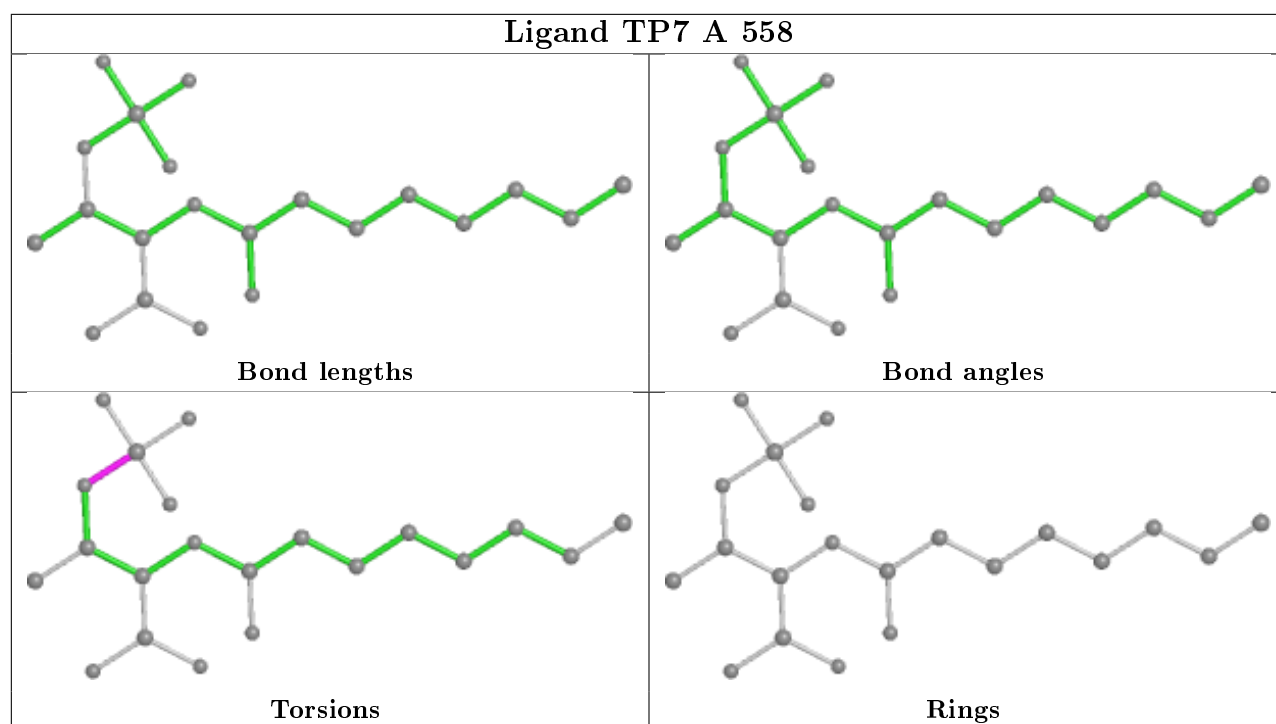
Bond angles

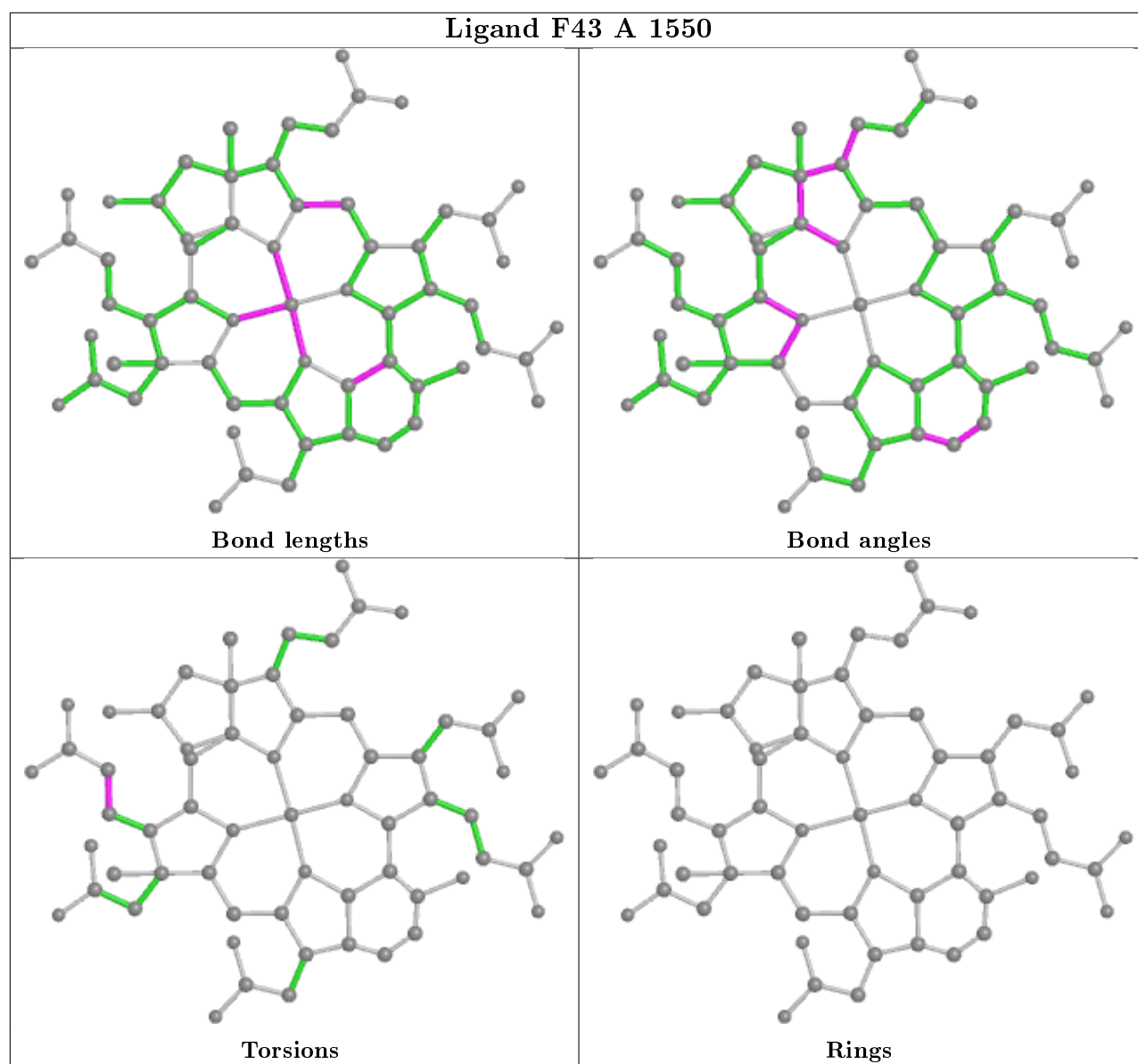


Torsions



Rings





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 ⓘ

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	543/550 (98%)	-0.37	4 (0%) 87 86	5, 9, 20, 43	0
1	D	543/550 (98%)	-0.35	3 (0%) 89 88	5, 8, 18, 39	1 (0%)
2	B	442/443 (99%)	-0.40	6 (1%) 75 74	6, 10, 21, 38	0
2	E	442/443 (99%)	-0.44	0 100 100	6, 9, 18, 31	0
3	C	248/249 (99%)	-0.26	10 (4%) 38 38	7, 13, 26, 34	0
3	F	248/249 (99%)	-0.27	6 (2%) 59 58	7, 11, 27, 50	0
All	All	2466/2484 (99%)	-0.36	29 (1%) 79 77	5, 10, 21, 50	1 (0%)

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	549	ALA	8.3
1	D	549	ALA	8.0
3	C	45	PRO	3.7
3	F	60	ASP	3.6
3	C	48	GLU	3.4
2	B	2	ALA	3.2
3	C	59	MET	3.2
1	A	240	LYS	3.0
2	B	443	ILE	2.9
3	C	46	GLY	2.9
3	C	197	GLU	2.8
1	A	171	GLU	2.7
2	B	441	ASN	2.7
3	C	63	GLU	2.6
1	A	2	ALA	2.6
3	F	63	GLU	2.6
3	F	48	GLU	2.5
3	C	47	GLU	2.4
2	B	297	ASP	2.4

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Mol	Chain	Res	Type	RSRZ
2	B	442	GLU	2.3
3	F	57	GLU	2.3
3	F	197	GLU	2.3
3	C	196	ASP	2.3
3	C	60	ASP	2.3
1	D	123	GLU	2.2
2	B	438	GLU	2.2
3	F	62	PRO	2.2
3	C	44	ALA	2.2
1	D	241[A]	GLN	2.1

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

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	AGM	D	271	12/13	0.97	0.11	4,6,6,7	0
1	MHS	D	257	11/12	0.97	0.06	9,10,13,15	0
1	MHS	A	257	11/12	0.97	0.07	10,11,14,16	0
1	MGN	D	400	10/11	0.98	0.10	5,6,6,7	0
1	AGM	A	271	12/13	0.98	0.10	5,6,7,8	0
1	MGN	A	400	10/11	0.99	0.07	5,6,6,7	0
1	SMC	A	452	7/8	0.99	0.07	5,6,8,8	0
1	GL3	A	445	4/5	0.99	0.07	5,6,6,6	0
1	SMC	D	452	7/8	0.99	0.08	5,5,7,8	0
1	GL3	D	445	4/5	1.00	0.09	4,5,5,5	0

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
9	ETX	A	1554	6/6	0.46	0.27	32,34,35,35	0
9	ETX	E	1446	6/6	0.66	0.21	36,36,37,38	0
9	ETX	D	1551	6/6	0.69	0.20	24,27,28,28	0
9	ETX	F	1250	6/6	0.70	0.23	33,36,37,37	0
9	ETX	C	1250	6/6	0.70	0.25	31,31,32,32	0
9	ETX	B	1447	6/6	0.70	0.23	29,30,32,35	0
9	ETX	F	1252	6/6	0.71	0.20	28,29,29,30	0
9	ETX	A	1551	6/6	0.71	0.17	22,25,26,27	0
9	ETX	A	1555	6/6	0.72	0.24	28,29,30,31	0
9	ETX	E	1449	6/6	0.73	0.26	33,35,35,36	0
9	ETX	E	1451	6/6	0.73	0.34	33,33,35,36	0
9	ETX	B	1446	6/6	0.75	0.15	28,28,29,31	0
9	ETX	E	1448	6/6	0.75	0.24	29,30,31,31	0
9	ETX	E	1447	6/6	0.78	0.26	36,37,38,40	0
9	ETX	E	1452	6/6	0.78	0.21	25,26,28,28	0
9	ETX	B	1448	6/6	0.79	0.38	32,33,33,34	0
9	ETX	D	1552	6/6	0.79	0.17	30,30,31,31	0
9	ETX	A	1553	6/6	0.80	0.18	24,26,28,28	0
9	ETX	E	1450	6/6	0.82	0.23	34,35,35,36	0
9	ETX	F	1251	6/6	0.82	0.21	26,28,31,32	0
9	ETX	B	1445	6/6	0.83	0.18	36,38,41,41	0
9	ETX	D	1553	6/6	0.85	0.38	31,32,34,35	0
9	ETX	E	1445	6/6	0.87	0.12	31,31,32,33	0
9	ETX	A	1552	6/6	0.88	0.17	32,33,35,37	0
9	ETX	E	1444	6/6	0.88	0.13	17,20,25,27	0
9	ETX	B	1444	6/6	0.89	0.17	17,22,28,31	0
4	COM	A	556	7/7	0.94	0.15	9,15,18,19	7
4	COM	D	555	7/7	0.94	0.16	8,13,16,16	7
7	CA	E	444	1/1	0.96	0.21	49,49,49,49	0
5	TP7	D	554	21/21	0.98	0.08	5,7,8,9	0
5	TP7	A	558	21/21	0.98	0.07	6,7,8,9	0
7	CA	B	444	1/1	0.98	0.24	42,42,42,42	0
7	CA	D	559	1/1	0.99	0.22	30,30,30,30	0
8	F43	D	1550	62/62	0.99	0.08	4,6,9,12	0
8	F43	A	1550	62/62	0.99	0.07	5,6,9,12	0
6	K	A	566	1/1	0.99	0.03	18,18,18,18	1
7	CA	A	565	1/1	1.00	0.11	23,23,23,23	0
10	MG	F	251	1/1	1.00	0.07	14,14,14,14	0
6	K	C	250	1/1	1.00	0.09	15,15,15,15	0
7	CA	D	558	1/1	1.00	0.09	25,25,25,25	0
7	CA	D	557	1/1	1.00	0.04	8,8,8,8	0
6	K	A	561	1/1	1.00	0.04	11,11,11,11	0
7	CA	A	562	1/1	1.00	0.10	23,23,23,23	0

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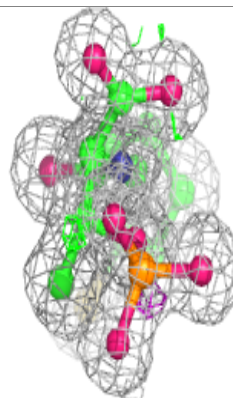
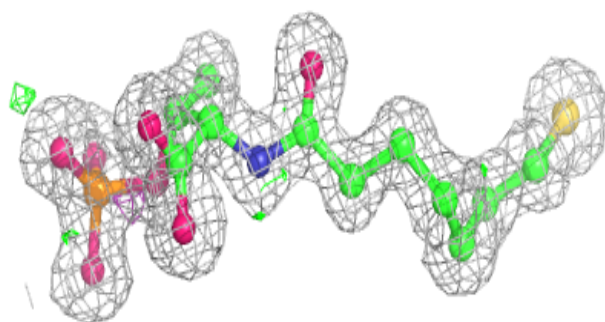
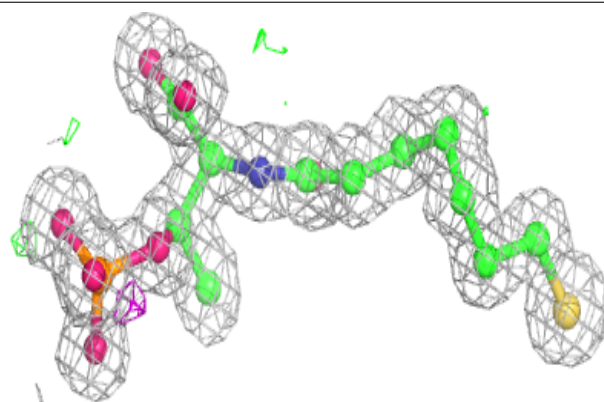
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
7	CA	F	250	1/1	1.00	0.04	15,15,15,15	0
7	CA	D	556	1/1	1.00	0.04	11,11,11,11	0
6	K	E	445	1/1	1.00	0.18	20,20,20,20	0
6	K	D	561	1/1	1.00	0.05	13,13,13,13	1
7	CA	D	560	1/1	1.00	0.10	16,16,16,16	0
7	CA	A	563	1/1	1.00	0.05	8,8,8,8	0
7	CA	A	564	1/1	1.00	0.09	24,24,24,24	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.

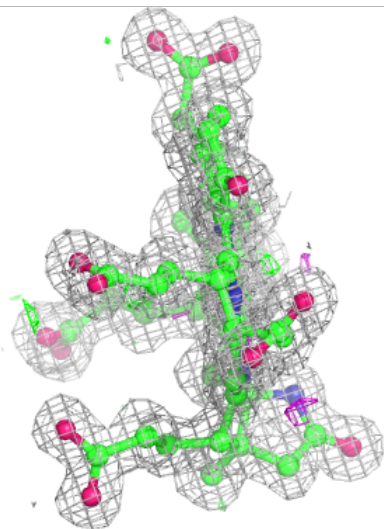
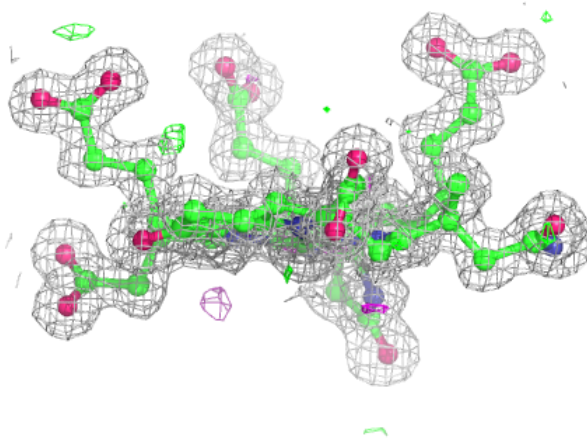
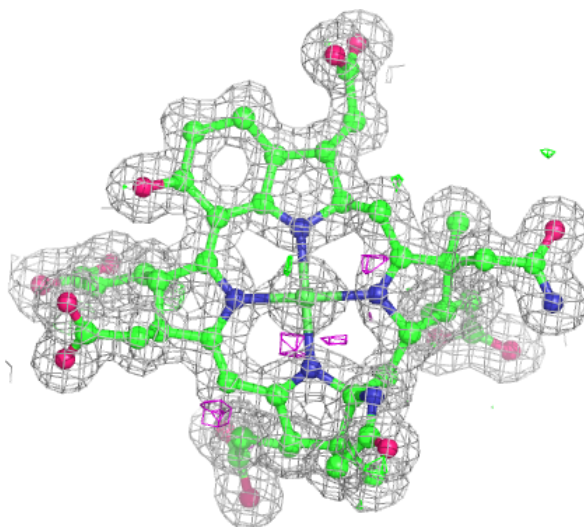
Electron density around TP7 A 558:

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



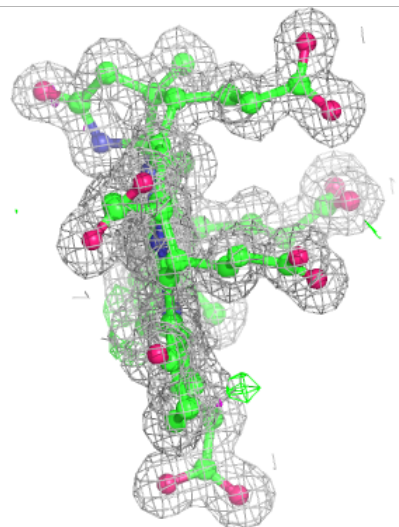
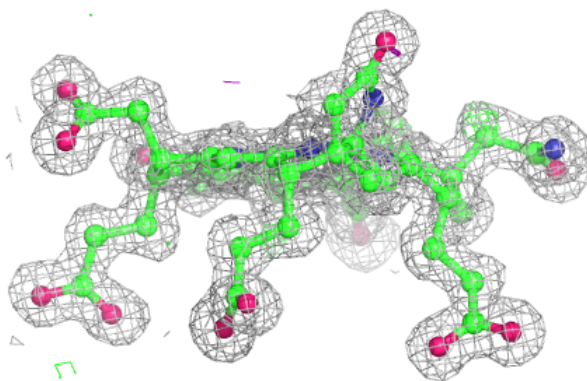
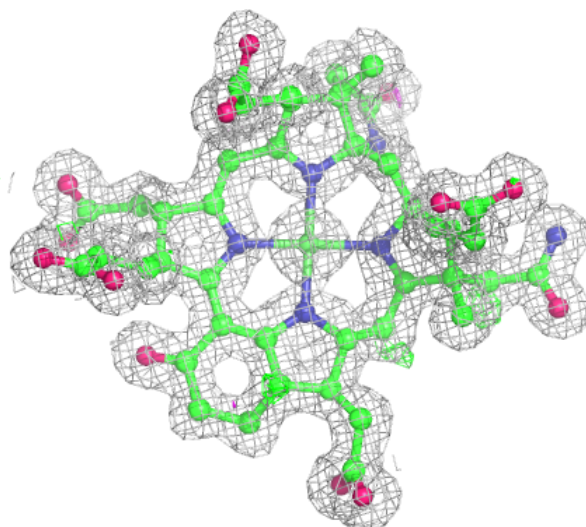
Electron density around F43 D 1550:

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



Electron density around F43 A 1550:

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



6.5 Other polymers ⓘ

There are no such residues in this entry.