wwPDB X-ray Structure Validation Summary Report (i)

Sep 10, 2026 – 05:04 pm BST

PDB ID	:	5ODH / pdb_00005odh
Title	:	Heterodisulfide reductase / [NiFe]-hydrogenase complex from Methanothermococcus thermolithotrophicus soaked with heterodisulfide for 3.5 minutes
Authors	:	Wagner, T.; Koch, J.; Ermler, U.; Shima, S.
Deposited on	:	2017-07-05
Resolution	:	2.20 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 (i)) were used in the production of this report:

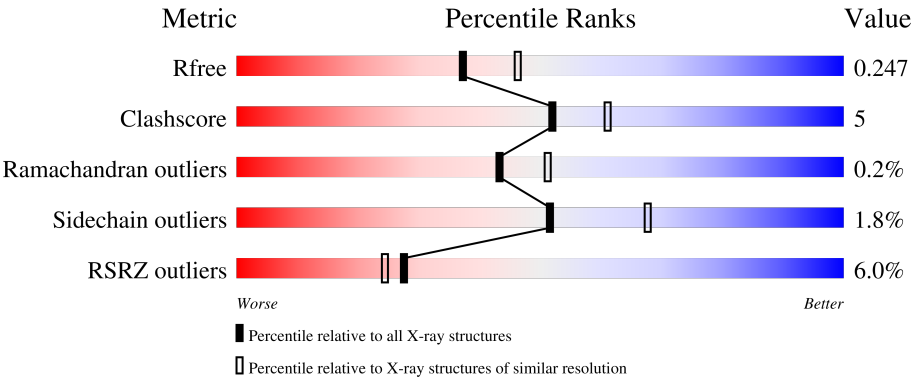
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

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 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	654	10% 88% 11%
1	G	654	5% 88% 11%
2	B	291	2% 87% 13%
2	H	291	27% 73% 27%
3	C	184	3% 84% 15%

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Mol	Chain	Length	Quality of chain
3	I	184	8% 79% 20%
4	D	140	84% 14%
4	J	140	2% 89% 9%
5	E	299	5% 88% 10%
5	K	299	2% 89% 10%
6	F	473	% 85% 9% 5%
6	L	473	2% 87% 7% 5%

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
12	COM	B	303	-	X	-	-

2 Entry composition (i)

There are 20 unique types of molecules in this entry. The entry contains 32106 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 Heterodisulfide reductase, subunit A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	652	Total	C	N	O	S	0	0	0
			4977	3140	845	943	49			
1	G	653	Total	C	N	O	S	0	0	0
			4984	3145	846	944	49			

- Molecule 2 is a protein called Heterodisulfide reductase, subunit B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	291	Total	C	N	O	S	0	0	0
			2236	1420	379	413	24			
2	H	291	Total	C	N	O	S	0	0	0
			2232	1417	378	413	24			

- Molecule 3 is a protein called Heterodisulfide reductase, subunit C.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	184	Total	C	N	O	S	0	0	0
			1421	887	246	274	14			
3	I	183	Total	C	N	O	S	0	0	0
			1416	885	246	271	14			

- Molecule 4 is a protein called Methyl-viologen reducing hydrogenase, subunit D.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	137	Total	C	N	O	S	0	0	0
			1097	698	187	200	12			
4	J	138	Total	C	N	O	S	0	0	0
			1106	703	188	203	12			

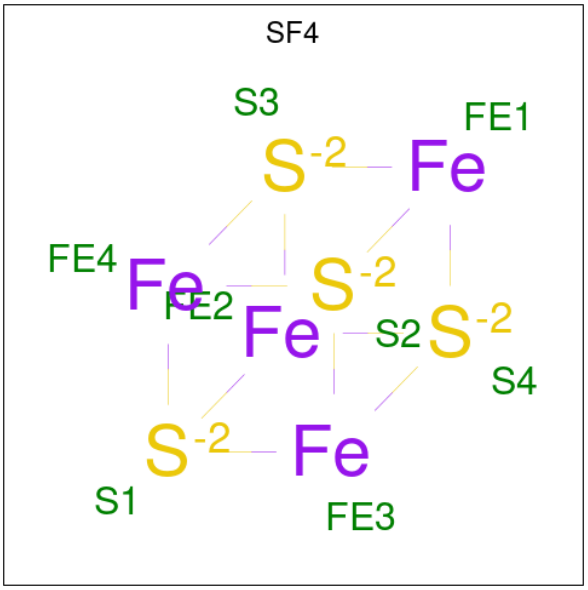
- Molecule 5 is a protein called Methyl-viologen reducing hydrogenase, subunit G.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	298	Total 2258	C 1425	N 369	O 445	S 19	0	0	0
5	K	298	Total 2258	C 1425	N 369	O 445	S 19	0	0	0

- Molecule 6 is a protein called Methyl-viologen reducing hydrogenase, subunit A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	F	447	Total 3521	C 2230	N 600	O 672	S 19	0	0	0
6	L	447	Total 3521	C 2230	N 600	O 672	S 19	0	0	0

- Molecule 7 is IRON/SULFUR CLUSTER (CCD ID: SF4) (formula: Fe_4S_4).



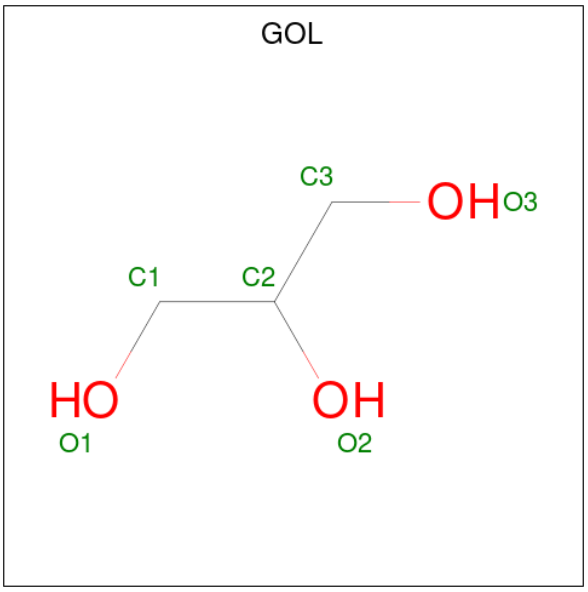
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	1	Total 8	Fe 4	S 4		0	0
7	A	1	Total 8	Fe 4	S 4		0	0
7	A	1	Total 8	Fe 4	S 4		0	0
7	A	1	Total 8	Fe 4	S 4		0	0
7	A	1	Total 8	Fe 4	S 4		0	0
7	A	1	Total 8	Fe 4	S 4		0	0

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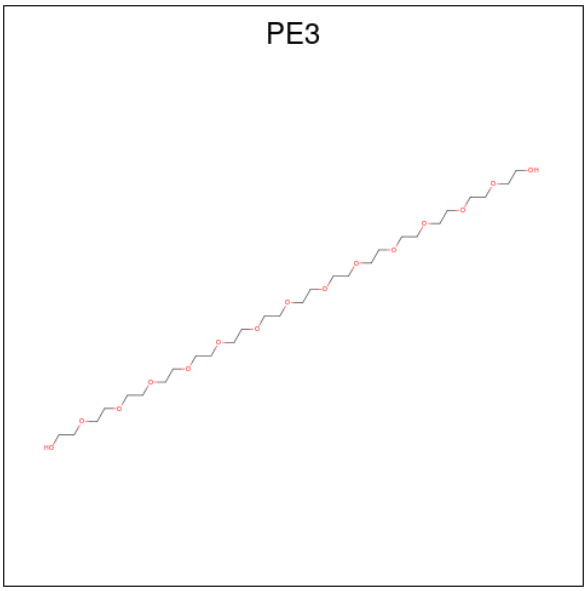
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	C	1	Total Fe S 8 4 4	0	0
7	C	1	Total Fe S 8 4 4	0	0
7	E	1	Total Fe S 8 4 4	0	0
7	E	1	Total Fe S 8 4 4	0	0
7	E	1	Total Fe S 8 4 4	0	0
7	G	1	Total Fe S 8 4 4	0	0
7	G	1	Total Fe S 8 4 4	0	0
7	G	1	Total Fe S 8 4 4	0	0
7	G	1	Total Fe S 8 4 4	0	0
7	G	1	Total Fe S 8 4 4	0	0
7	G	1	Total Fe S 8 4 4	0	0
7	I	1	Total Fe S 8 4 4	0	0
7	I	1	Total Fe S 8 4 4	0	0
7	K	1	Total Fe S 8 4 4	0	0
7	K	1	Total Fe S 8 4 4	0	0
7	K	1	Total Fe S 8 4 4	0	0

- Molecule 8 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



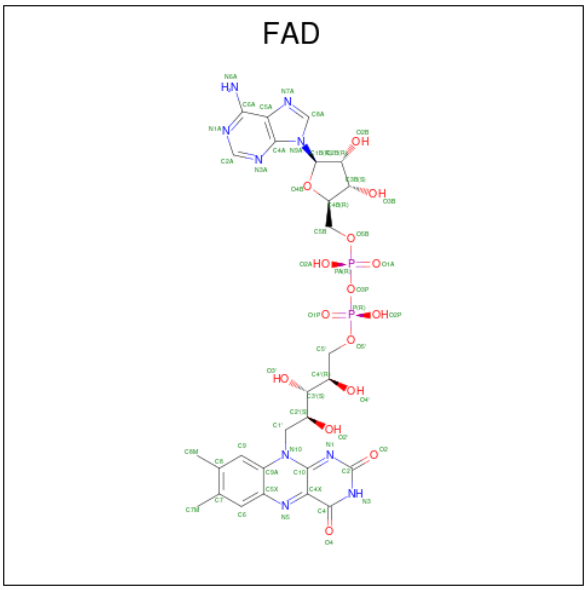
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	1	Total C O 6 3 3	0	0
8	B	1	Total C O 6 3 3	0	0
8	B	1	Total C O 6 3 3	0	0
8	E	1	Total C O 6 3 3	0	0
8	G	1	Total C O 6 3 3	0	0
8	K	1	Total C O 6 3 3	0	0

- Molecule 9 is 3,6,9,12,15,18,21,24,27,30,33,36,39-TRIDECAOXAHENTETRACONTANE-1 ,41-DIOL (CCD ID: PE3) (formula: $C_{28}H_{58}O_{15}$).



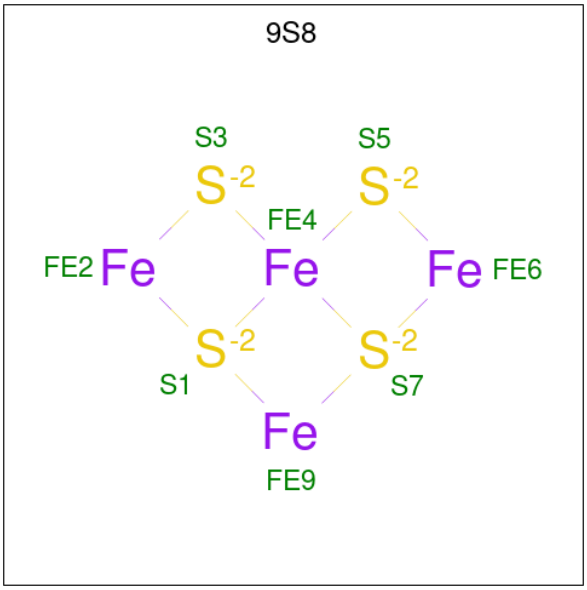
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	A	1	Total C O 14 9 5	0	0
9	F	1	Total C O 9 6 3	0	0
9	F	1	Total C O 9 6 3	0	0
9	L	1	Total C O 9 6 3	0	0

- Molecule 10 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



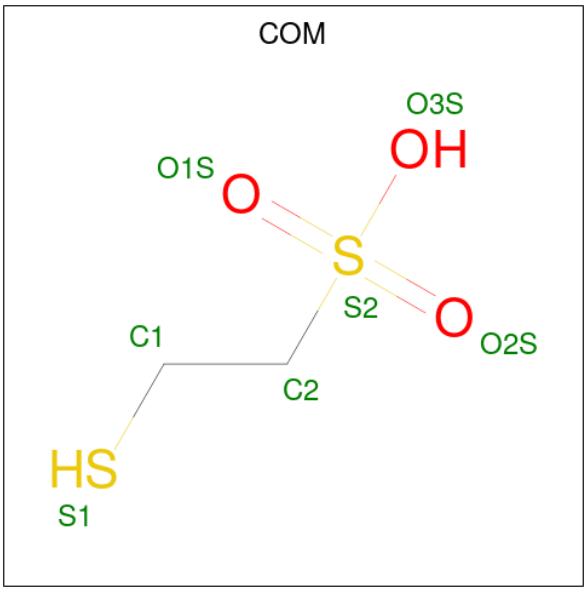
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
10	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
10	G	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 11 is Non-cubane [4Fe-4S]-cluster (CCD ID: 9S8) (formula: Fe_4S_4).



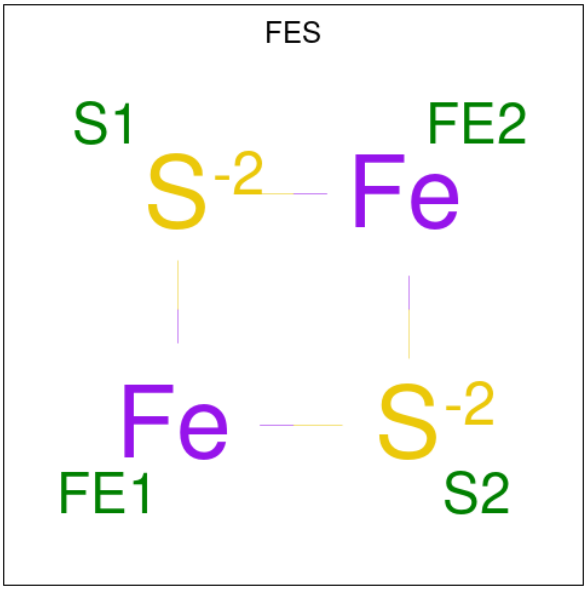
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
11	B	1	Total	Fe	S	0	0
			8	4	4		
11	B	1	Total	Fe	S	0	0
			8	4	4		
11	H	1	Total	Fe	S	0	0
			8	4	4		
11	H	1	Total	Fe	S	0	0
			8	4	4		

- Molecule 12 is 1-THIOETHANESULFONIC ACID (CCD ID: COM) (formula: $C_2H_6O_3S_2$).



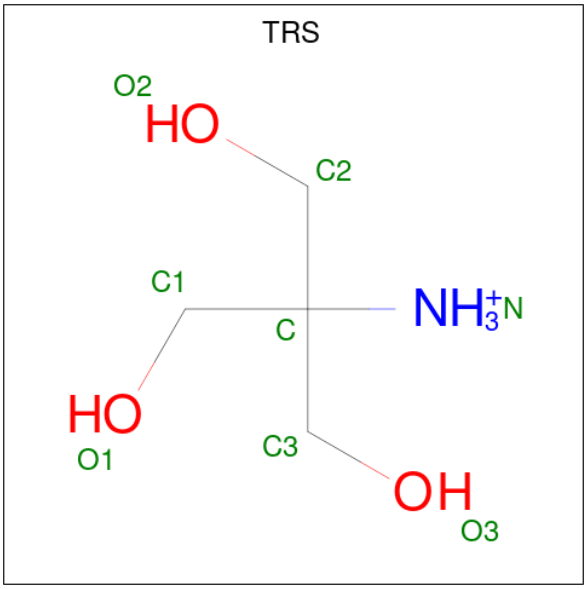
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
12	B	1	Total C O S 7 2 3 2	0	0
12	H	1	Total C O S 7 2 3 2	0	0

- Molecule 13 is FE2/S2 (INORGANIC) CLUSTER (CCD ID: FES) (formula: Fe_2S_2).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
13	D	1	Total Fe S 4 2 2	0	0
13	J	1	Total Fe S 4 2 2	0	0

- Molecule 14 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (CCD ID: TRS) (formula: $C_4H_{12}NO_3$).

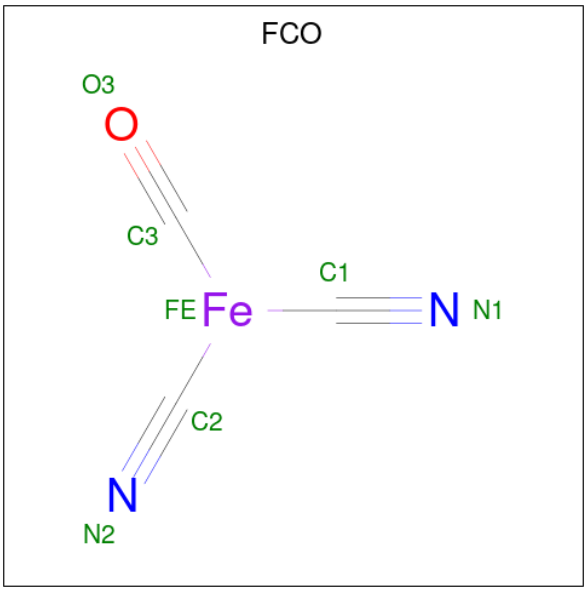


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
14	D	1	Total C N O 8 4 1 3	0	0

- Molecule 15 is FE (III) ION (CCD ID: FE) (formula: Fe).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
15	F	1	Total Fe 1 1	0	0
15	G	1	Total Fe 1 1	0	0
15	L	1	Total Fe 1 1	0	0

- Molecule 16 is CARBONMONOXIDE-(DICYANO) IRON (CCD ID: FCO) (formula: C_3FeN_2O).

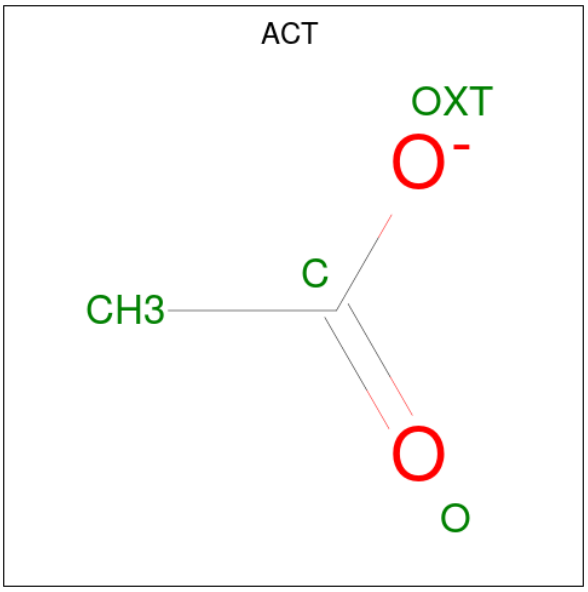


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
16	F	1	Total C Fe N O 7 3 1 2 1	0	0
16	L	1	Total C Fe N O 7 3 1 2 1	0	0

- Molecule 17 is NICKEL (II) ION (CCD ID: NI) (formula: Ni).

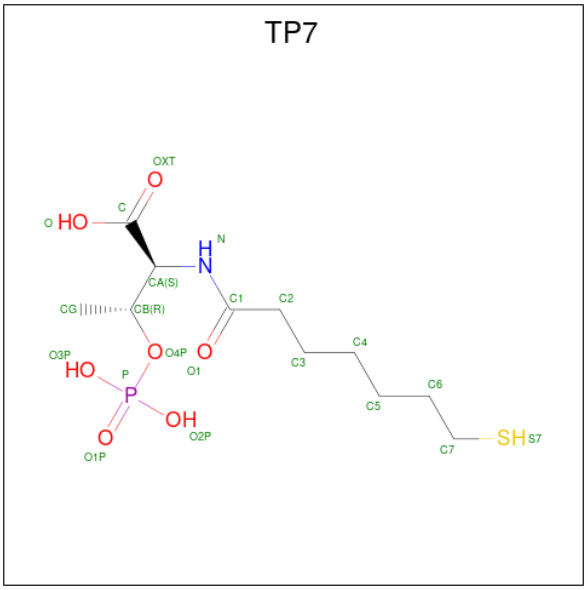
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
17	F	1	Total Ni 1 1	0	0
17	L	1	Total Ni 1 1	0	0

- Molecule 18 is ACETATE ION (CCD ID: ACT) (formula: $C_2H_3O_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
18	G	1	Total C O 4 2 2	0	0
18	G	1	Total C O 4 2 2	0	0

- Molecule 19 is Coenzyme B (CCD ID: TP7) (formula: $C_{11}H_{22}NO_7PS$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
19	H	1	Total C N O P S 21 11 1 7 1 1	0	0

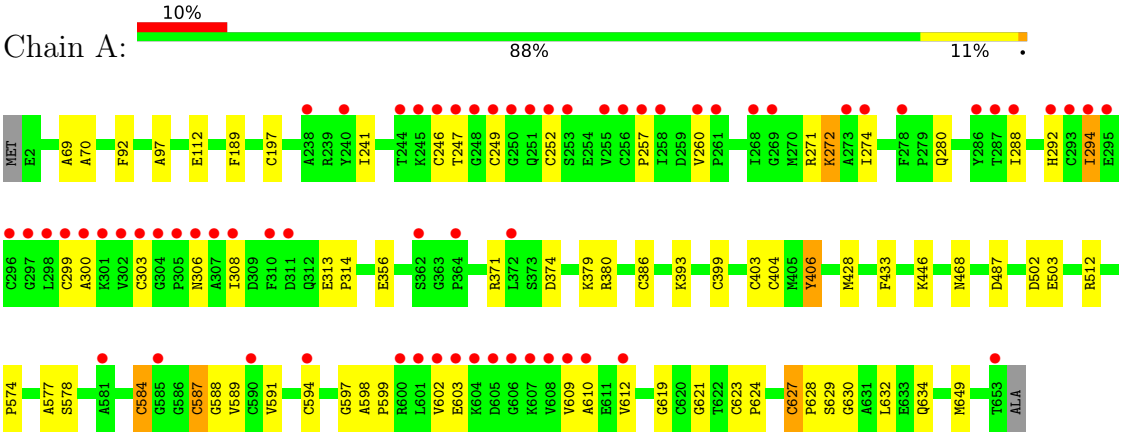
- Molecule 20 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
20	A	111	Total O 111 111	0	0
20	B	27	Total O 27 27	0	0
20	C	25	Total O 25 25	0	0
20	D	29	Total O 29 29	0	0
20	E	37	Total O 37 37	0	0
20	F	67	Total O 67 67	0	0
20	G	134	Total O 134 134	0	0
20	H	5	Total O 5 5	0	0
20	I	14	Total O 14 14	0	0
20	J	31	Total O 31 31	0	0
20	K	51	Total O 51 51	0	0
20	L	79	Total O 79 79	0	0

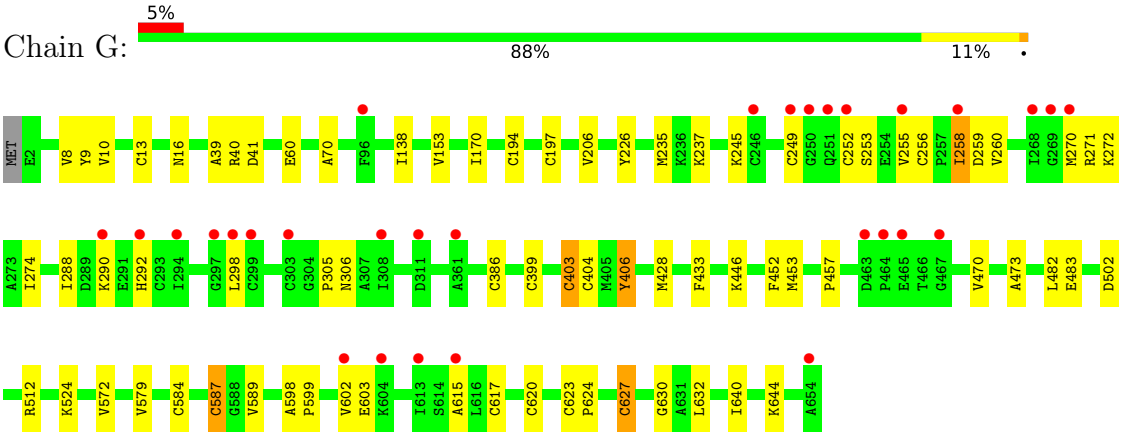
3 Residue-property plots (i)

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.

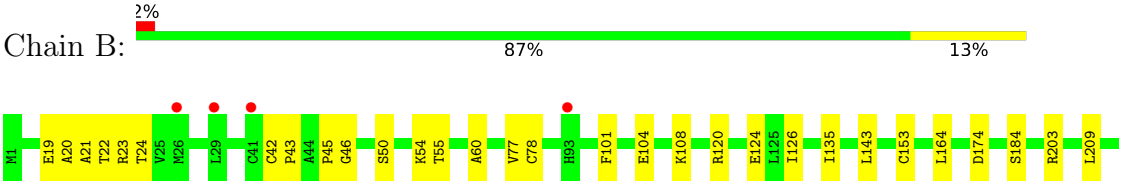
• Molecule 1: Heterodisulfide reductase, subunit A

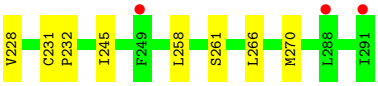


• Molecule 1: Heterodisulfide reductase, subunit A



• Molecule 2: Heterodisulfide reductase, subunit B

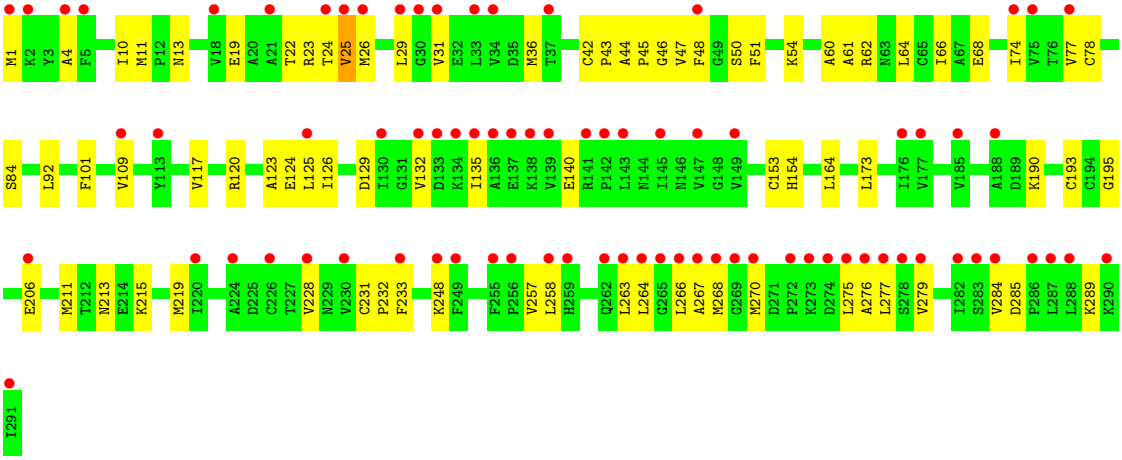




• Molecule 2: Heterodisulfide reductase, subunit B

Chain H:





• Molecule 3: Heterodisulfide reductase, subunit C

Chain C:

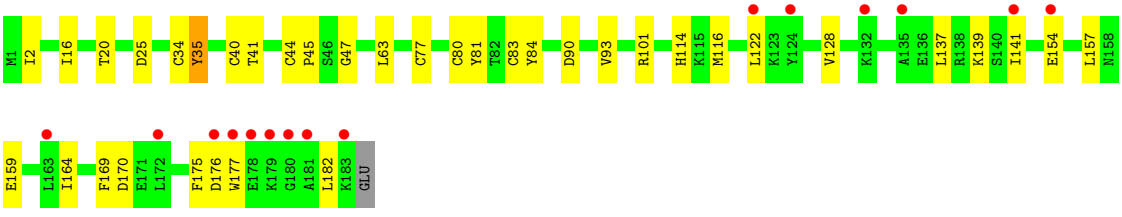




• Molecule 3: Heterodisulfide reductase, subunit C

Chain I:

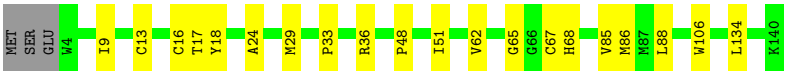




• Molecule 4: Methyl-viologen reducing hydrogenase, subunit D

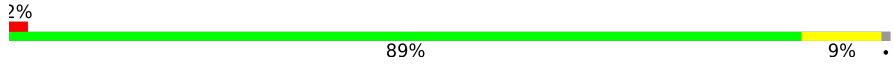
Chain D:

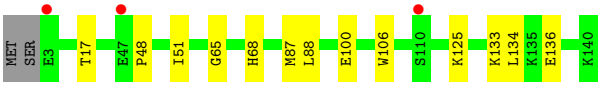




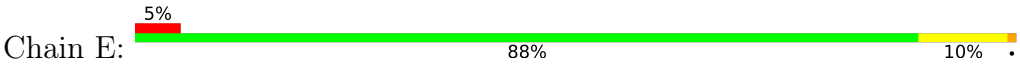
• Molecule 4: Methyl-viologen reducing hydrogenase, subunit D

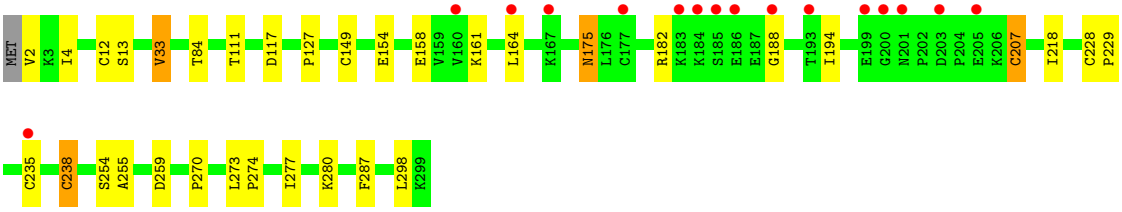
Chain J:



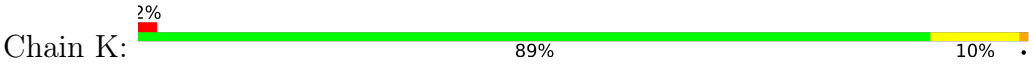


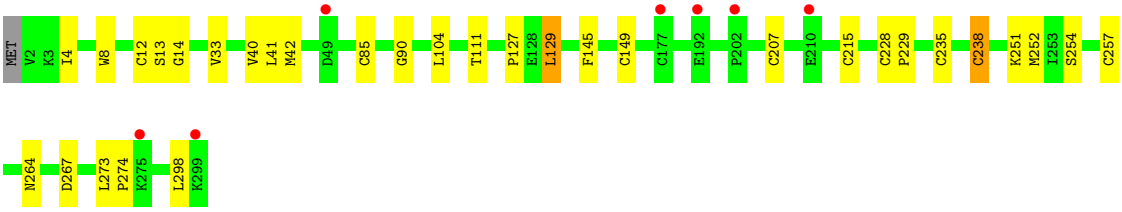
• Molecule 5: Methyl-viologen reducing hydrogenase, subunit G



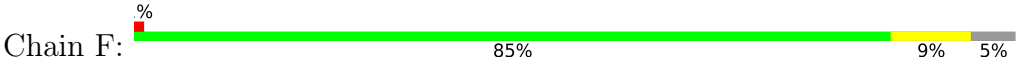


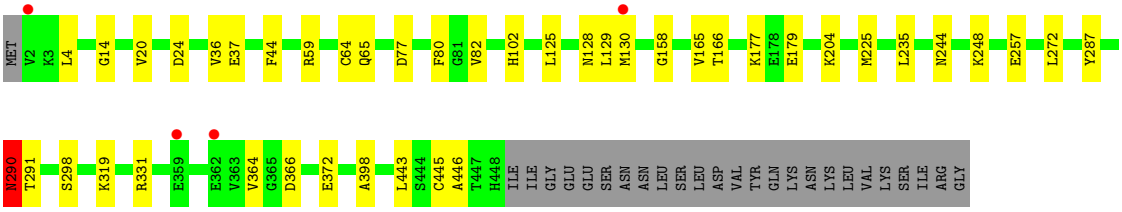
• Molecule 5: Methyl-viologen reducing hydrogenase, subunit G



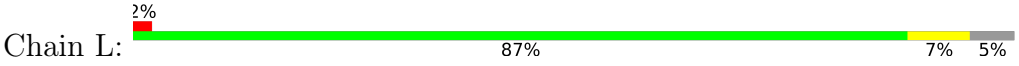


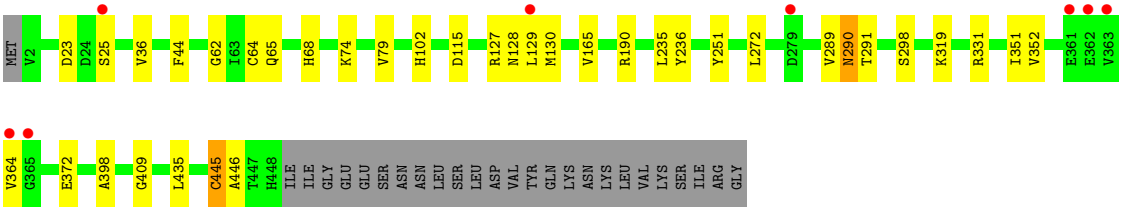
• Molecule 6: Methyl-viologen reducing hydrogenase, subunit A





• Molecule 6: Methyl-viologen reducing hydrogenase, subunit A





4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	366.38Å 97.10Å 133.97Å 90.00° 108.43° 90.00°	Depositor
Resolution (Å)	48.55 – 2.20 48.55 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.8 (48.55-2.20) 99.8 (48.55-2.20)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.86 (at 2.20Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155: ???)	Depositor
R, R_{free}	0.214 , 0.245 0.218 , 0.247	Depositor DCC
R_{free} test set	11273 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å^2)	28.7	Xtriage
Anisotropy	0.224	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 34.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.45$, $< L^2 > = 0.28$	Xtriage
Estimated twinning fraction	0.024 for -h-2*l,-k,l	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	32106	wwPDB-VP
Average B, all atoms (Å^2)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.71% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical 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.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: FES, FCO, SF4, FE, 9S8, FAD, NI, TRS, TP7, ACT, COM, GOL, PE3

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	0.19	0/5069	0.36	0/6850
1	G	0.18	0/5076	0.36	0/6860
2	B	0.20	0/2277	0.43	0/3070
2	H	0.20	0/2273	0.47	0/3066
3	C	0.21	0/1442	0.42	0/1942
3	I	0.21	0/1437	0.46	0/1934
4	D	0.20	0/1123	0.39	0/1508
4	J	0.18	0/1132	0.36	0/1520
5	E	0.20	0/2297	0.41	0/3113
5	K	0.19	0/2297	0.40	0/3113
6	F	0.18	0/3590	0.39	0/4853
6	L	0.19	0/3590	0.39	0/4853
All	All	0.19	0/31603	0.40	0/42682

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4977	0	4968	57	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	G	4984	0	4979	45	0
2	B	2236	0	2243	28	0
2	H	2232	0	2234	64	0
3	C	1421	0	1431	18	0
3	I	1416	0	1435	28	0
4	D	1097	0	1068	14	0
4	J	1106	0	1074	9	0
5	E	2258	0	2265	26	0
5	K	2258	0	2265	21	0
6	F	3521	0	3506	26	1
6	L	3521	0	3505	22	1
7	A	48	0	0	1	0
7	C	16	0	0	0	0
7	E	24	0	0	0	0
7	G	48	0	0	1	0
7	I	16	0	0	0	0
7	K	24	0	0	0	0
8	A	6	0	8	0	0
8	B	12	0	16	0	0
8	E	6	0	8	0	0
8	G	6	0	8	2	0
8	K	6	0	8	1	0
9	A	14	0	16	0	0
9	F	18	0	18	1	0
9	L	9	0	10	0	0
10	A	53	0	31	0	0
10	G	53	0	31	0	0
11	B	16	0	0	1	0
11	H	16	0	0	1	0
12	B	7	0	6	1	0
12	H	7	0	6	2	0
13	D	4	0	0	0	0
13	J	4	0	0	0	0
14	D	8	0	12	0	0
15	F	1	0	0	0	0
15	G	1	0	0	0	0
15	L	1	0	0	0	0
16	F	7	0	0	1	0
16	L	7	0	0	1	0
17	F	1	0	0	0	0
17	L	1	0	0	0	0
18	G	8	0	6	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
19	H	21	0	19	3	0
20	A	111	0	0	2	0
20	B	27	0	0	1	0
20	C	25	0	0	0	0
20	D	29	0	0	0	0
20	E	37	0	0	0	0
20	F	67	0	0	1	0
20	G	134	0	0	0	0
20	H	5	0	0	0	0
20	I	14	0	0	0	0
20	J	31	0	0	0	0
20	K	51	0	0	0	0
20	L	79	0	0	0	0
All	All	32106	0	31176	318	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 5.

The worst 5 of 318 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)
6:F:290:ASN:OD1	6:F:291:THR:N	2.09	0.85
1:G:587:CYS:SG	1:G:589:VAL:HG23	2.24	0.77
1:A:197:CYS:SG	20:A:910:HOH:O	2.49	0.71
2:B:104:GLU:OE2	20:B:401:HOH:O	2.10	0.67
6:L:289:VAL:O	6:L:290:ASN:HB2	1.95	0.67

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)
6:F:24:ASP:OD2	6:L:190:ARG:NH1[3_546]	1.99	0.21

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries

of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	650/654 (99%)	624 (96%)	25 (4%)	1 (0%)	43 51
1	G	651/654 (100%)	615 (94%)	35 (5%)	1 (0%)	43 51
2	B	289/291 (99%)	281 (97%)	8 (3%)	0	100 100
2	H	289/291 (99%)	278 (96%)	11 (4%)	0	100 100
3	C	182/184 (99%)	177 (97%)	5 (3%)	0	100 100
3	I	181/184 (98%)	177 (98%)	4 (2%)	0	100 100
4	D	135/140 (96%)	132 (98%)	3 (2%)	0	100 100
4	J	136/140 (97%)	133 (98%)	3 (2%)	0	100 100
5	E	296/299 (99%)	286 (97%)	10 (3%)	0	100 100
5	K	296/299 (99%)	287 (97%)	9 (3%)	0	100 100
6	F	445/473 (94%)	437 (98%)	6 (1%)	2 (0%)	30 34
6	L	445/473 (94%)	439 (99%)	4 (1%)	2 (0%)	30 34
All	All	3995/4082 (98%)	3866 (97%)	123 (3%)	6 (0%)	43 51

5 of 6 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
6	L	290	ASN
6	F	290	ASN
6	L	398	ALA
1	A	70	ALA
6	F	398	ALA

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	539/541 (100%)	530 (98%)	9 (2%)	53 69

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	G	540/541 (100%)	526 (97%)	14 (3%)	40 55
2	B	242/242 (100%)	242 (100%)	0	100 100
2	H	241/242 (100%)	239 (99%)	2 (1%)	73 85
3	C	156/157 (99%)	150 (96%)	6 (4%)	29 40
3	I	156/157 (99%)	151 (97%)	5 (3%)	34 47
4	D	116/119 (98%)	116 (100%)	0	100 100
4	J	117/119 (98%)	117 (100%)	0	100 100
5	E	255/256 (100%)	247 (97%)	8 (3%)	35 48
5	K	255/256 (100%)	249 (98%)	6 (2%)	43 58
6	F	386/410 (94%)	380 (98%)	6 (2%)	55 71
6	L	386/410 (94%)	381 (99%)	5 (1%)	61 76
All	All	3389/3450 (98%)	3328 (98%)	61 (2%)	51 68

5 of 61 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
6	F	445	CYS
5	K	238	CYS
1	G	404	CYS
5	K	235	CYS
6	L	251	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 20 such sidechains are listed below:

Mol	Chain	Res	Type
2	H	262	GLN
4	J	124	ASN
6	L	394	ASN
6	L	97	GLN
1	G	61	HIS

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates (i)

There are no oligosaccharides in this entry.

5.6 Ligand geometry (i)

Of 53 ligands modelled in this entry, 5 are monoatomic - leaving 48 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	$\#\|Z\| > 2$
7	SF4	G	703	1	0,12,12	-	-	-		
10	FAD	A	709	-	56,58,58	1.43	9 (16%)	81,89,89	1.64	17 (20%)
18	ACT	G	710	-	3,3,3	0.76	0	3,3,3	1.44	0
7	SF4	I	201	3	0,12,12	-	-	-		
7	SF4	A	702	-	0,12,12	-	-	-		
7	SF4	G	704	1	0,12,12	-	-	-		
11	9S8	H	302	2	2,10,10	1.14	0	-		
7	SF4	E	303	5	0,12,12	-	-	-		
12	COM	B	303	-	6,6,6	1.33	1 (16%)	7,8,8	2.99	4 (57%)
11	9S8	B	301	2	2,10,10	1.55	1 (50%)	-		
7	SF4	K	303	5	0,12,12	-	-	-		
19	TP7	H	304	-	19,20,20	0.64	0	24,26,26	1.05	0
7	SF4	C	202	3	0,12,12	-	-	-		
11	9S8	H	301	2	2,10,10	1.13	0	-		
9	PE3	F	501	-	8,8,42	0.45	0	7,7,41	0.31	0
8	GOL	B	304	-	5,5,5	0.38	0	5,5,5	0.19	0
8	GOL	G	707	-	5,5,5	0.35	0	5,5,5	0.13	0
18	ACT	G	709	-	3,3,3	0.80	0	3,3,3	1.35	0
9	PE3	F	502	-	8,8,42	0.46	0	7,7,41	0.22	0
7	SF4	G	702	1	0,12,12	-	-	-		
8	GOL	E	304	-	5,5,5	0.34	0	5,5,5	0.22	0
9	PE3	A	708	-	13,13,42	0.49	0	12,12,41	0.19	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
7	SF4	A	701	1	0,12,12	-	-	-		
7	SF4	E	302	5	0,12,12	-	-	-		
7	SF4	E	301	5	0,12,12	-	-	-		
9	PE3	L	501	-	8,8,42	0.50	0	7,7,41	0.21	0
7	SF4	G	705	1	0,12,12	-	-	-		
16	FCO	F	505	6	1,6,6	3.98	1 (100%)	-		
7	SF4	A	705	1	0,12,12	-	-	-		
7	SF4	A	704	1	0,12,12	-	-	-		
7	SF4	A	706	1	0,12,12	-	-	-		
7	SF4	K	301	5	0,12,12	-	-	-		
14	TRS	D	202	-	7,7,7	0.27	0	9,9,9	0.39	0
8	GOL	B	305	-	5,5,5	0.30	0	5,5,5	0.25	0
7	SF4	C	201	3	0,12,12	-	-	-		
7	SF4	I	202	3	0,12,12	-	-	-		
7	SF4	G	701	1,15	0,12,12	-	-	-		
7	SF4	G	706	1	0,12,12	-	-	-		
7	SF4	A	703	1	0,12,12	-	-	-		
8	GOL	A	707	-	5,5,5	0.31	0	5,5,5	0.35	0
10	FAD	G	711	-	56,58,58	1.45	9 (16%)	81,89,89	1.61	12 (14%)
13	FES	D	201	4	0,4,4	-	-	-		
12	COM	H	303	-	6,6,6	1.42	2 (33%)	7,8,8	2.67	4 (57%)
11	9S8	B	302	2	2,10,10	1.40	0	-		
7	SF4	K	302	5	0,12,12	-	-	-		
8	GOL	K	304	-	5,5,5	0.35	0	5,5,5	0.22	0
16	FCO	L	504	6	1,6,6	4.21	1 (100%)	-		
13	FES	J	201	4	0,4,4	-	-	-		

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
7	SF4	G	703	1	-	-	0/6/5/5
10	FAD	A	709	-	-	2/34/50/50	0/6/6/6
7	SF4	I	201	3	-	-	0/6/5/5
7	SF4	A	702	-	-	-	0/6/5/5
7	SF4	G	704	1	-	-	0/6/5/5
11	9S8	H	302	2	-	-	0/3/3/3
7	SF4	E	303	5	-	-	0/6/5/5
12	COM	B	303	-	-	4/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
11	9S8	B	301	2	-	-	0/3/3/3
7	SF4	K	303	5	-	-	0/6/5/5
19	TP7	H	304	-	-	10/24/24/24	-
7	SF4	C	202	3	-	-	0/6/5/5
11	9S8	H	301	2	-	-	0/3/3/3
9	PE3	F	501	-	-	3/6/6/40	-
8	GOL	B	304	-	-	4/4/4/4	-
8	GOL	G	707	-	-	2/4/4/4	-
9	PE3	F	502	-	-	5/6/6/40	-
7	SF4	G	702	1	-	-	0/6/5/5
8	GOL	E	304	-	-	1/4/4/4	-
9	PE3	A	708	-	-	5/11/11/40	-
7	SF4	A	701	1	-	-	0/6/5/5
9	PE3	L	501	-	-	2/6/6/40	-
7	SF4	E	301	5	-	-	0/6/5/5
7	SF4	E	302	5	-	-	0/6/5/5
7	SF4	G	705	1	-	-	0/6/5/5
7	SF4	A	704	1	-	-	0/6/5/5
7	SF4	A	705	1	-	-	0/6/5/5
7	SF4	A	706	1	-	-	0/6/5/5
7	SF4	K	301	5	-	-	0/6/5/5
14	TRS	D	202	-	-	8/9/9/9	-
8	GOL	B	305	-	-	0/4/4/4	-
7	SF4	C	201	3	-	-	0/6/5/5
7	SF4	I	202	3	-	-	0/6/5/5
7	SF4	G	701	1,15	-	-	0/6/5/5
7	SF4	G	706	1	-	-	0/6/5/5
8	GOL	A	707	-	-	2/4/4/4	-
7	SF4	A	703	1	-	-	0/6/5/5
10	FAD	G	711	-	-	1/34/50/50	0/6/6/6
13	FES	D	201	4	-	-	0/1/1/1
12	COM	H	303	-	-	1/4/4/4	-
11	9S8	B	302	2	-	-	0/3/3/3
7	SF4	K	302	5	-	-	0/6/5/5
8	GOL	K	304	-	-	2/4/4/4	-
13	FES	J	201	4	-	-	0/1/1/1

The worst 5 of 24 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
10	G	711	FAD	C9A-C5X	5.35	1.50	1.41
10	A	709	FAD	C9A-C5X	5.01	1.49	1.41
10	A	709	FAD	C5A-C4A	4.51	1.47	1.39
10	G	711	FAD	C5A-C4A	4.33	1.47	1.39
16	L	504	FCO	O3-C3	4.21	1.21	1.16

The worst 5 of 37 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
10	G	711	FAD	C5A-C4A-N3A	-6.08	118.82	126.75
10	A	709	FAD	C5A-C4A-N3A	-5.86	119.11	126.75
10	G	711	FAD	N3A-C4A-N9A	4.84	135.06	127.08
12	B	303	COM	O3S-S2-O2S	-4.42	100.48	111.27
10	A	709	FAD	N3A-C4A-N9A	4.34	134.24	127.08

There are no chirality outliers.

5 of 52 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
8	B	304	GOL	O1-C1-C2-C3
8	B	304	GOL	C1-C2-C3-O3
12	B	303	COM	S1-C1-C2-S2
12	B	303	COM	C1-C2-S2-O1S
12	B	303	COM	C1-C2-S2-O2S

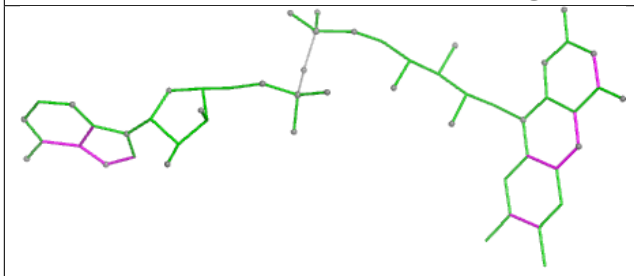
There are no ring outliers.

13 monomers are involved in 16 short contacts:

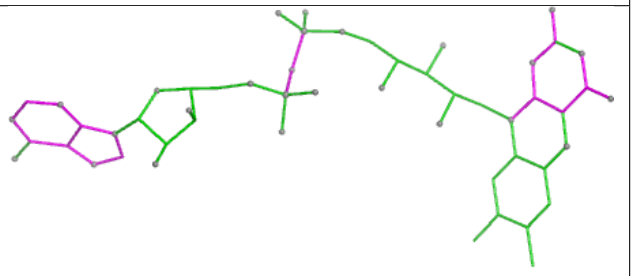
Mol	Chain	Res	Type	Clashes	Symm-Clashes
18	G	710	ACT	1	0
12	B	303	COM	1	0
19	H	304	TP7	3	0
11	H	301	9S8	1	0
8	G	707	GOL	2	0
9	F	502	PE3	1	0
7	G	702	SF4	1	0
16	F	505	FCO	1	0
7	A	706	SF4	1	0
12	H	303	COM	2	0
11	B	302	9S8	1	0
8	K	304	GOL	1	0
16	L	504	FCO	1	0

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 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.

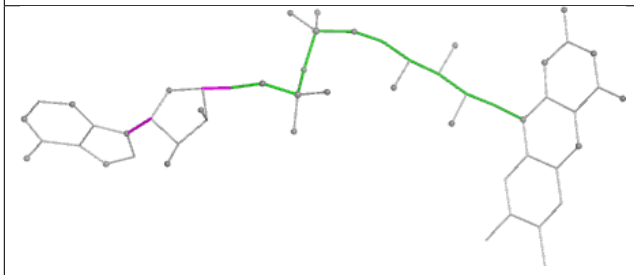
Ligand FAD A 709



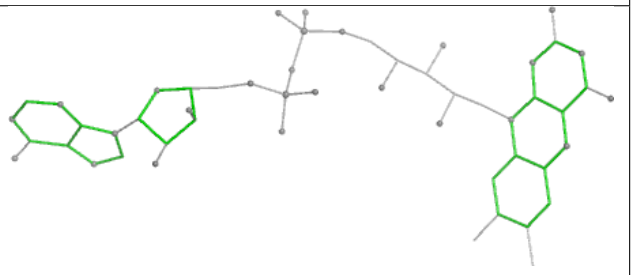
Bond lengths



Bond angles

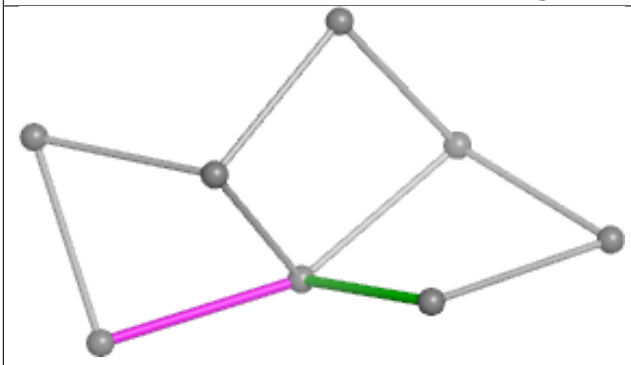


Torsions

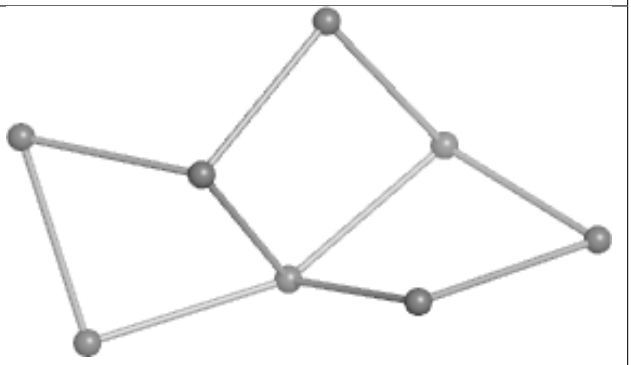


Rings

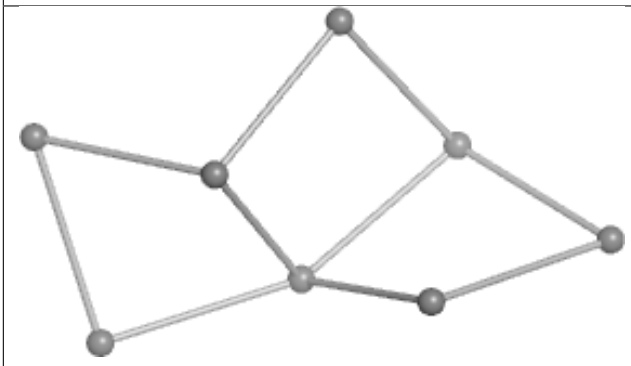
Ligand 9S8 B 301



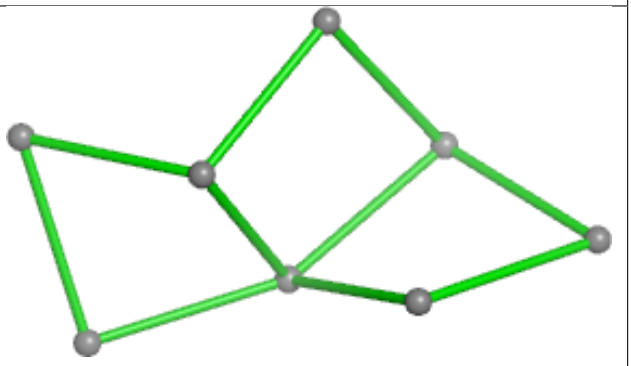
Bond lengths



Bond angles

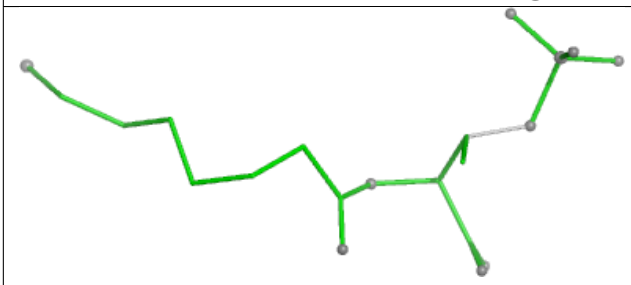


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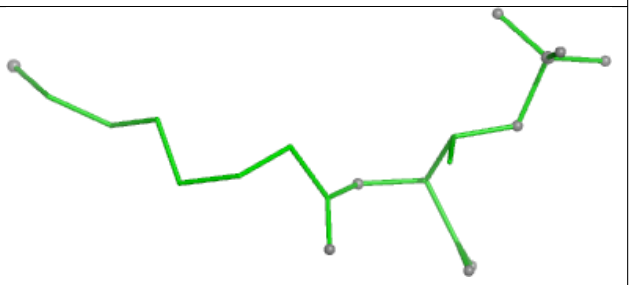


Rings

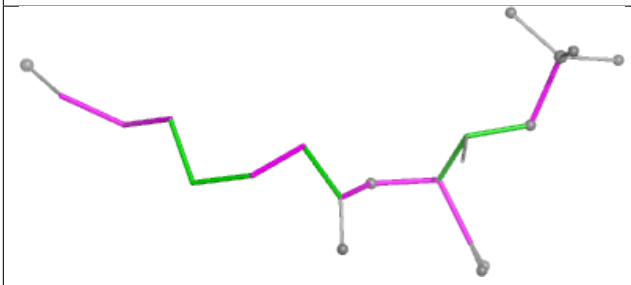
Ligand TP7 H 304



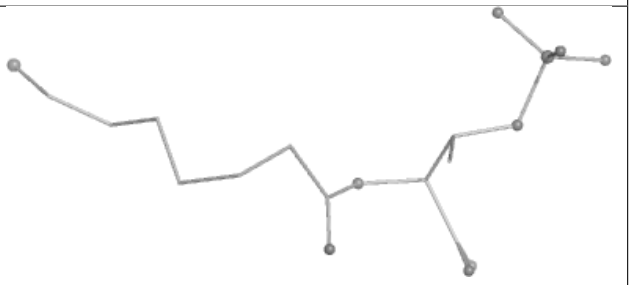
Bond lengths



Bond angles

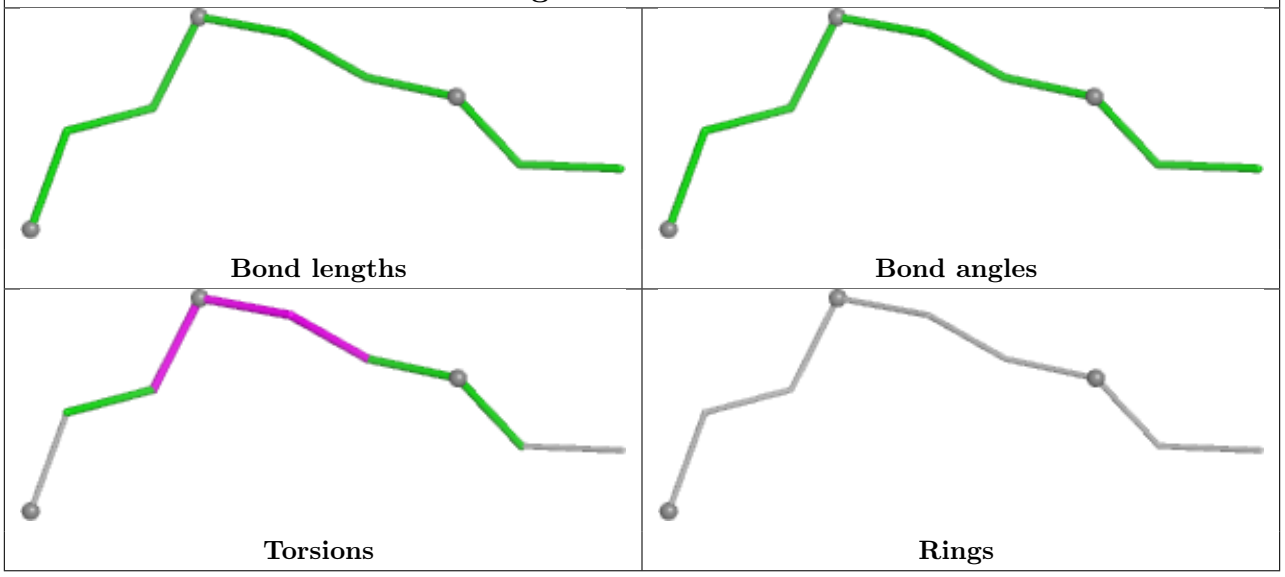


Torsions

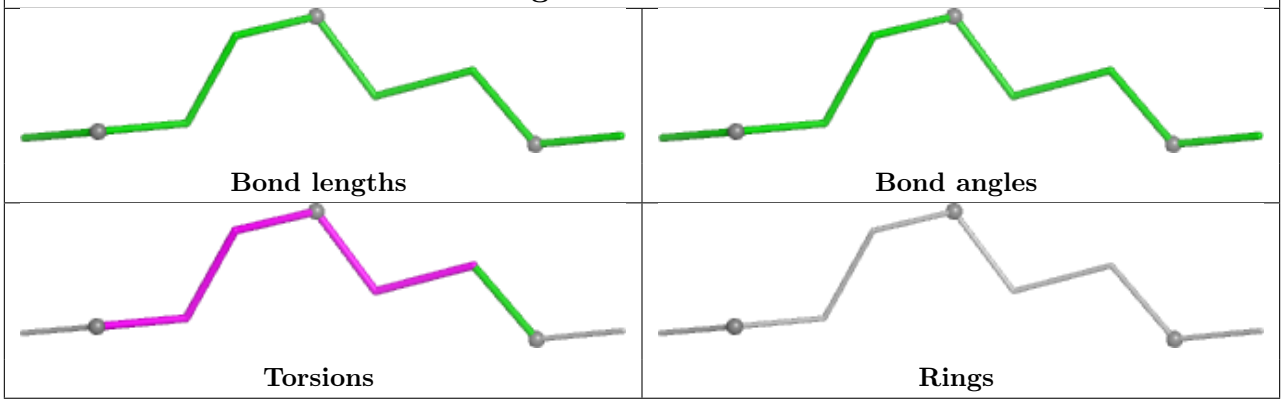


Rings

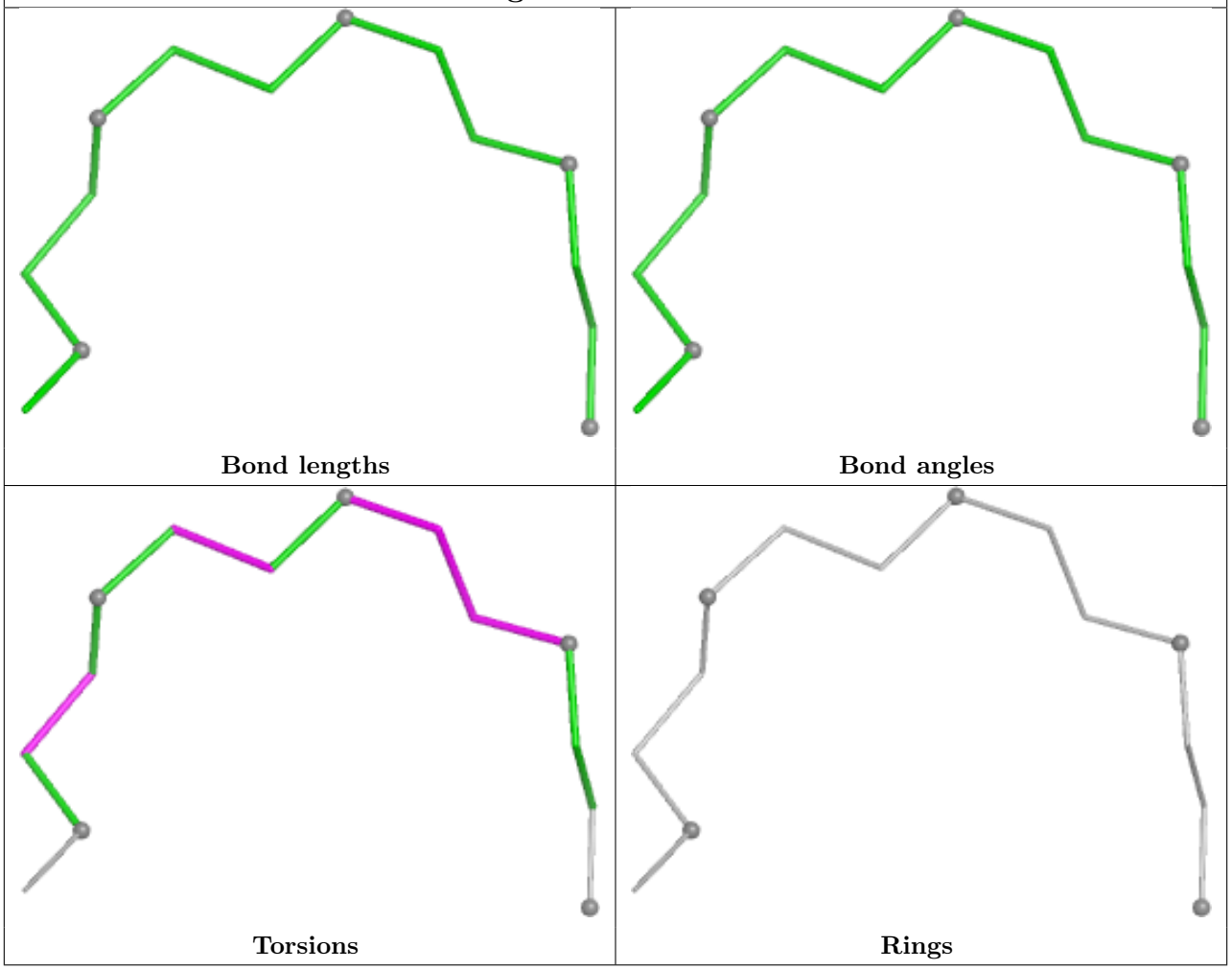
Ligand PE3 F 501



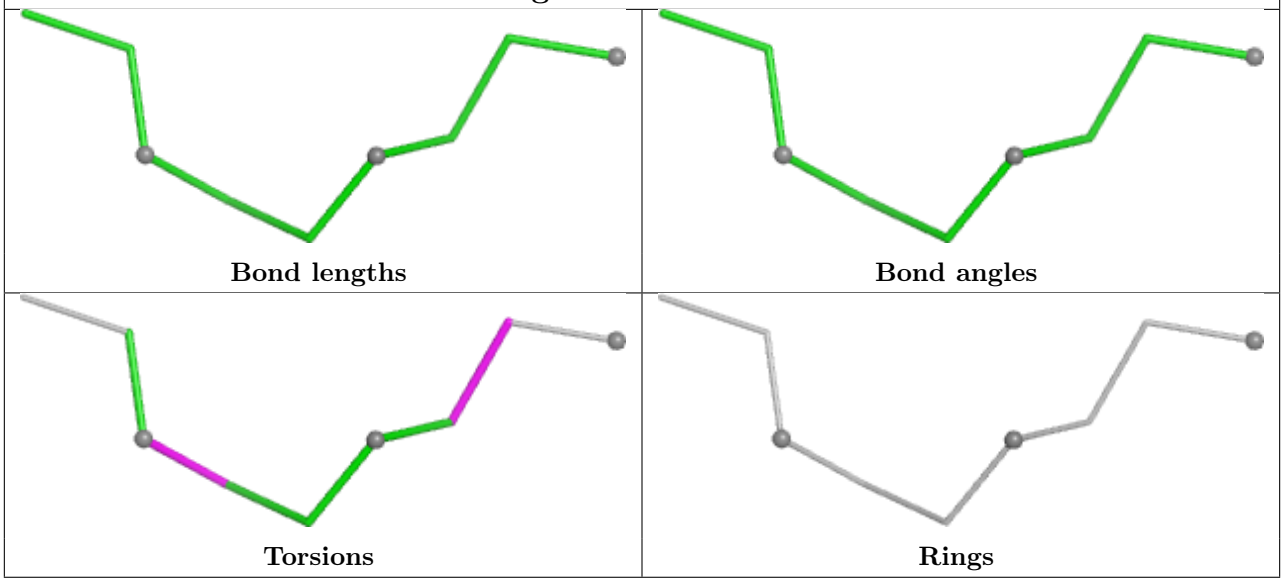
Ligand PE3 F 502



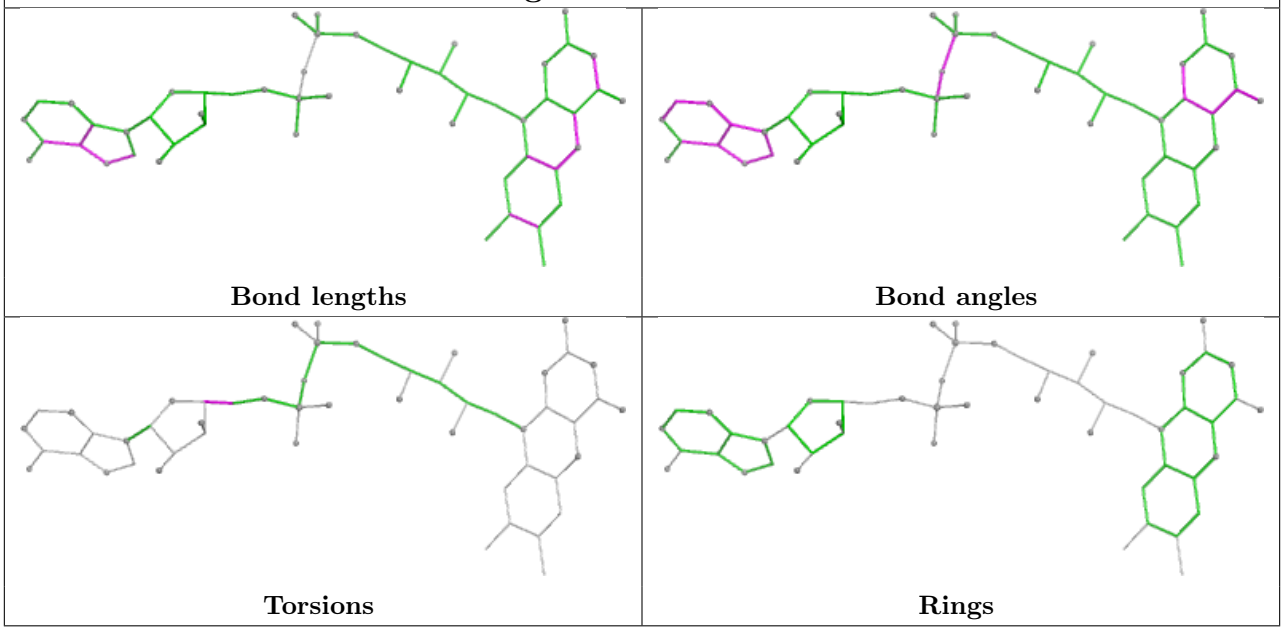
Ligand PE3 A 708



Ligand PE3 L 501



Ligand FAD G 711



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, 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>2	OWAB(Å^2)	Q<0.9
1	A	652/654 (99%)	0.46	65 (9%) 12 10	13, 31, 81, 112	0
1	G	653/654 (99%)	0.34	30 (4%) 37 34	16, 31, 71, 90	0
2	B	291/291 (100%)	0.54	7 (2%) 59 56	21, 40, 55, 70	0
2	H	291/291 (100%)	1.42	79 (27%) 1 1	31, 54, 98, 116	0
3	C	184/184 (100%)	0.41	6 (3%) 49 46	21, 35, 53, 74	0
3	I	183/184 (99%)	0.80	15 (8%) 17 15	19, 43, 71, 98	0
4	D	137/140 (97%)	0.04	0 100 100	19, 28, 40, 54	0
4	J	138/140 (98%)	0.25	3 (2%) 62 59	21, 30, 43, 76	0
5	E	298/299 (99%)	0.72	16 (5%) 31 28	18, 34, 62, 81	0
5	K	298/299 (99%)	0.26	7 (2%) 61 58	16, 29, 50, 66	0
6	F	447/473 (94%)	0.24	4 (0%) 81 79	15, 30, 52, 75	0
6	L	447/473 (94%)	0.11	8 (1%) 67 64	15, 28, 49, 106	0
All	All	4019/4082 (98%)	0.45	240 (5%) 27 24	13, 33, 67, 116	0

The worst 5 of 240 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	610	ALA	5.4
1	A	602	VAL	5.4
2	H	25	VAL	5.2
1	G	258	ILE	4.7
3	I	179	LYS	4.6

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

There are no non-standard protein/DNA/RNA residues in this entry.

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, 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	B-factors(Å^2)	Q<0.9
18	ACT	G	709	4/4	0.74	0.14	38,38,38,47	0
19	TP7	H	304	21/21	0.75	0.16	39,61,74,80	0
18	ACT	G	710	4/4	0.77	0.17	31,39,39,47	0
9	PE3	A	708	14/43	0.82	0.14	30,41,56,57	0
8	GOL	B	304	6/6	0.82	0.16	31,38,41,48	0
8	GOL	E	304	6/6	0.82	0.14	38,42,50,53	0
8	GOL	K	304	6/6	0.82	0.16	29,36,41,49	0
9	PE3	L	501	9/43	0.83	0.13	28,39,44,50	0
12	COM	B	303	7/7	0.86	0.12	34,50,61,71	0
9	PE3	F	501	9/43	0.87	0.13	31,33,43,46	0
7	SF4	A	702	8/8	0.87	0.10	101,125,137,146	0
8	GOL	A	707	6/6	0.88	0.12	33,38,40,47	0
8	GOL	B	305	6/6	0.89	0.10	28,39,41,49	0
8	GOL	G	707	6/6	0.90	0.15	26,36,45,46	0
11	9S8	H	301	8/8	0.90	0.10	38,54,72,77	0
9	PE3	F	502	9/43	0.91	0.11	32,41,47,47	0
14	TRS	D	202	8/8	0.92	0.09	25,31,38,40	0
7	SF4	A	703	8/8	0.92	0.08	72,84,107,111	0
11	9S8	H	302	8/8	0.92	0.07	30,44,52,57	0
11	9S8	B	302	8/8	0.92	0.08	21,31,36,37	0
11	9S8	B	301	8/8	0.93	0.10	30,40,47,58	0
7	SF4	G	703	8/8	0.93	0.07	66,69,73,95	0
7	SF4	G	704	8/8	0.93	0.08	73,87,106,125	0
7	SF4	A	706	8/8	0.94	0.05	19,34,42,43	0
12	COM	H	303	7/7	0.94	0.09	34,36,45,53	0
7	SF4	A	704	8/8	0.95	0.06	27,30,37,37	0
7	SF4	C	202	8/8	0.95	0.05	21,29,31,32	0
17	NI	F	506	1/1	0.95	0.10	23,23,23,23	0
7	SF4	G	705	8/8	0.95	0.06	27,30,38,41	0
7	SF4	I	201	8/8	0.95	0.06	19,24,25,28	0
7	SF4	K	303	8/8	0.95	0.06	27,29,36,37	0
7	SF4	E	302	8/8	0.96	0.05	17,31,38,39	0

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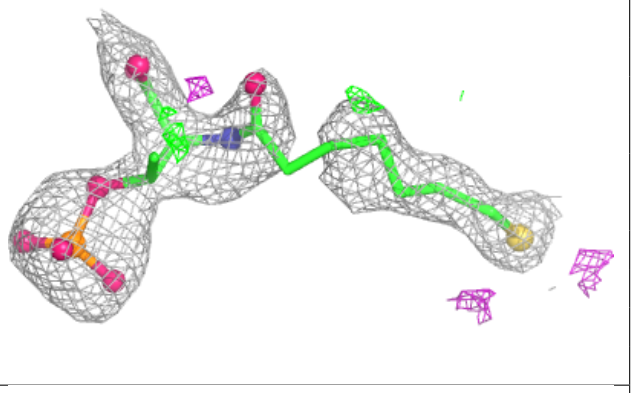
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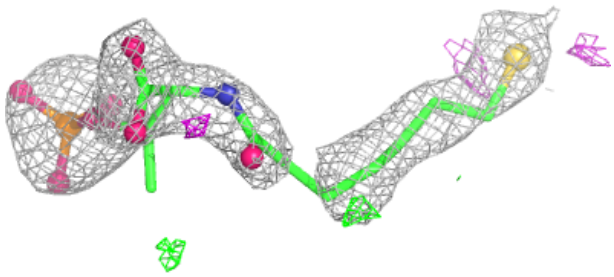
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
7	SF4	K	301	8/8	0.96	0.05	15,21,23,23	0
7	SF4	K	302	8/8	0.96	0.06	23,34,36,37	0
7	SF4	E	303	8/8	0.96	0.06	28,39,49,63	0
7	SF4	C	201	8/8	0.96	0.06	20,25,27,32	0
7	SF4	A	705	8/8	0.96	0.05	27,31,42,43	0
10	FAD	A	709	53/53	0.96	0.07	10,19,34,37	0
7	SF4	E	301	8/8	0.96	0.05	20,27,29,34	0
7	SF4	G	706	8/8	0.96	0.05	21,30,36,40	0
7	SF4	I	202	8/8	0.97	0.04	21,23,31,33	0
16	FCO	L	504	7/7	0.97	0.10	20,22,25,29	0
7	SF4	G	702	8/8	0.97	0.05	19,21,24,26	0
7	SF4	A	701	8/8	0.97	0.04	14,20,22,25	0
10	FAD	G	711	53/53	0.97	0.06	10,20,27,28	0
7	SF4	G	701	8/8	0.97	0.05	18,28,37,40	0
17	NI	L	505	1/1	0.98	0.09	25,25,25,25	0
16	FCO	F	505	7/7	0.98	0.08	14,21,23,24	0
13	FES	D	201	4/4	0.98	0.04	24,24,25,27	0
15	FE	F	504	1/1	0.98	0.03	23,23,23,23	0
15	FE	G	708	1/1	0.99	0.03	32,32,32,32	1
15	FE	L	503	1/1	0.99	0.03	20,20,20,20	0
13	FES	J	201	4/4	0.99	0.04	20,21,26,29	0

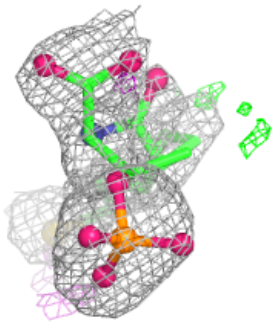
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around TP7 H 304:

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

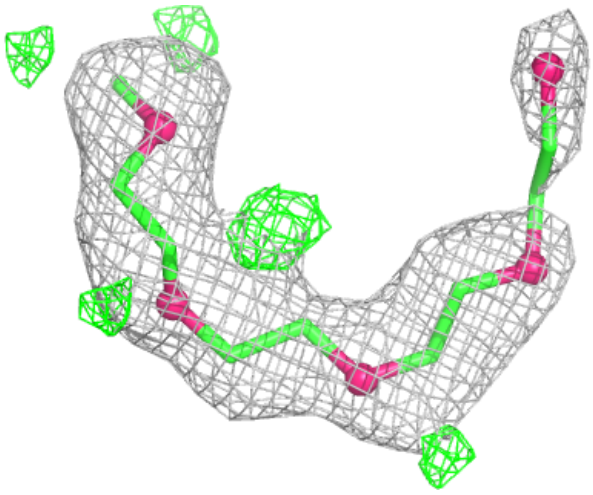


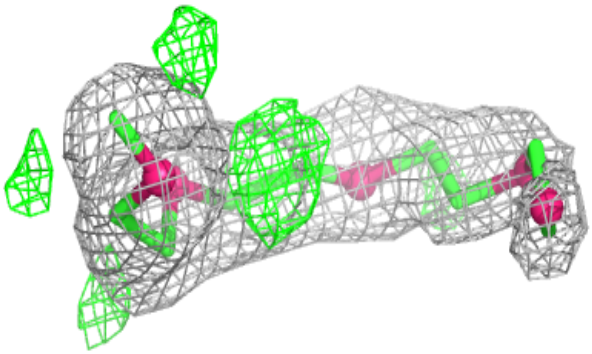


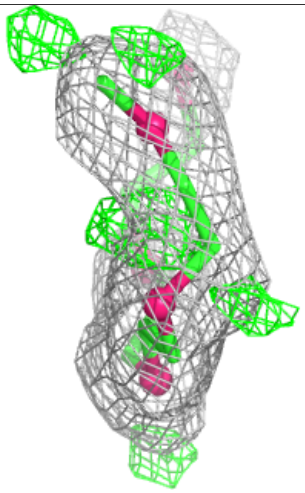


Electron density around PE3 A 708:

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

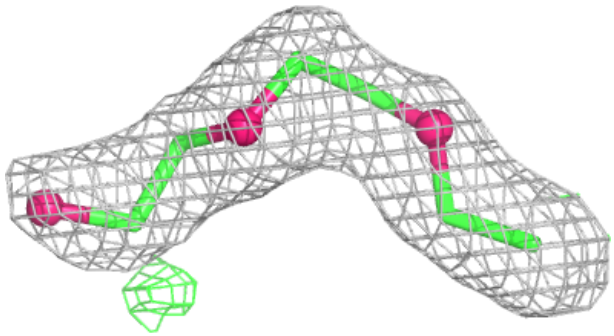


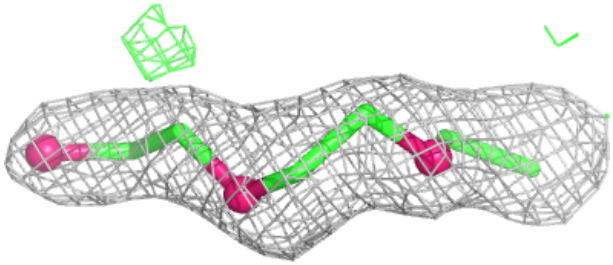


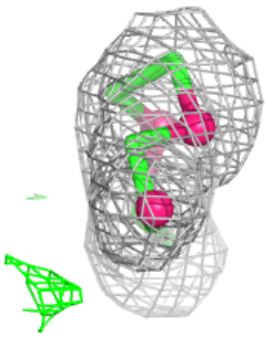


Electron density around PE3 L 501:

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

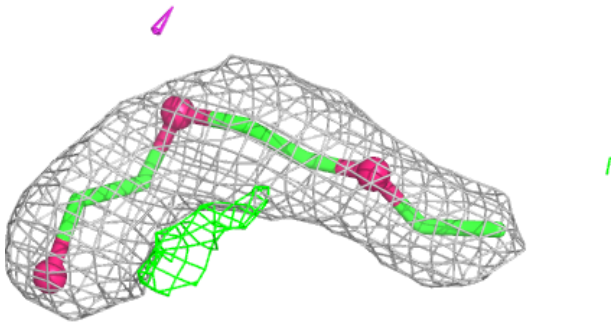


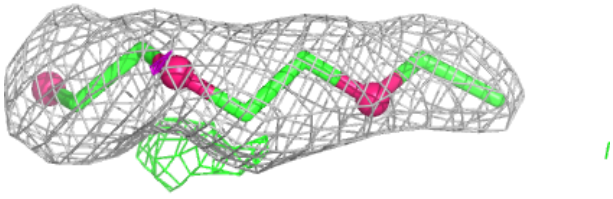


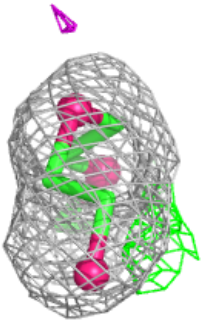


Electron density around PE3 F 501:

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

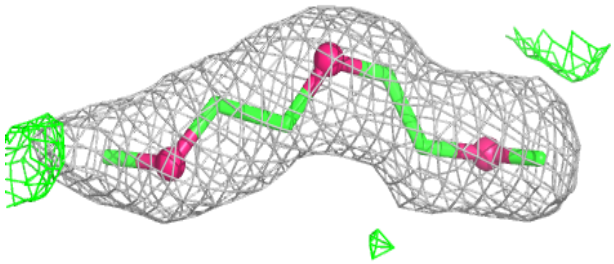


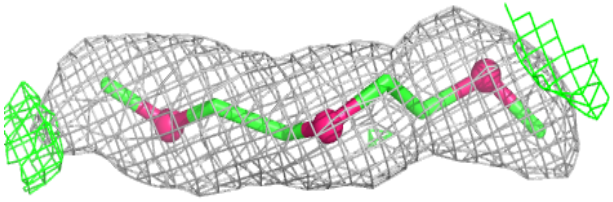


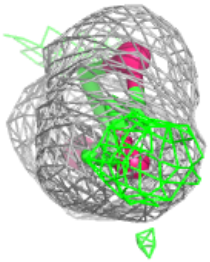


Electron density around PE3 F 502:

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

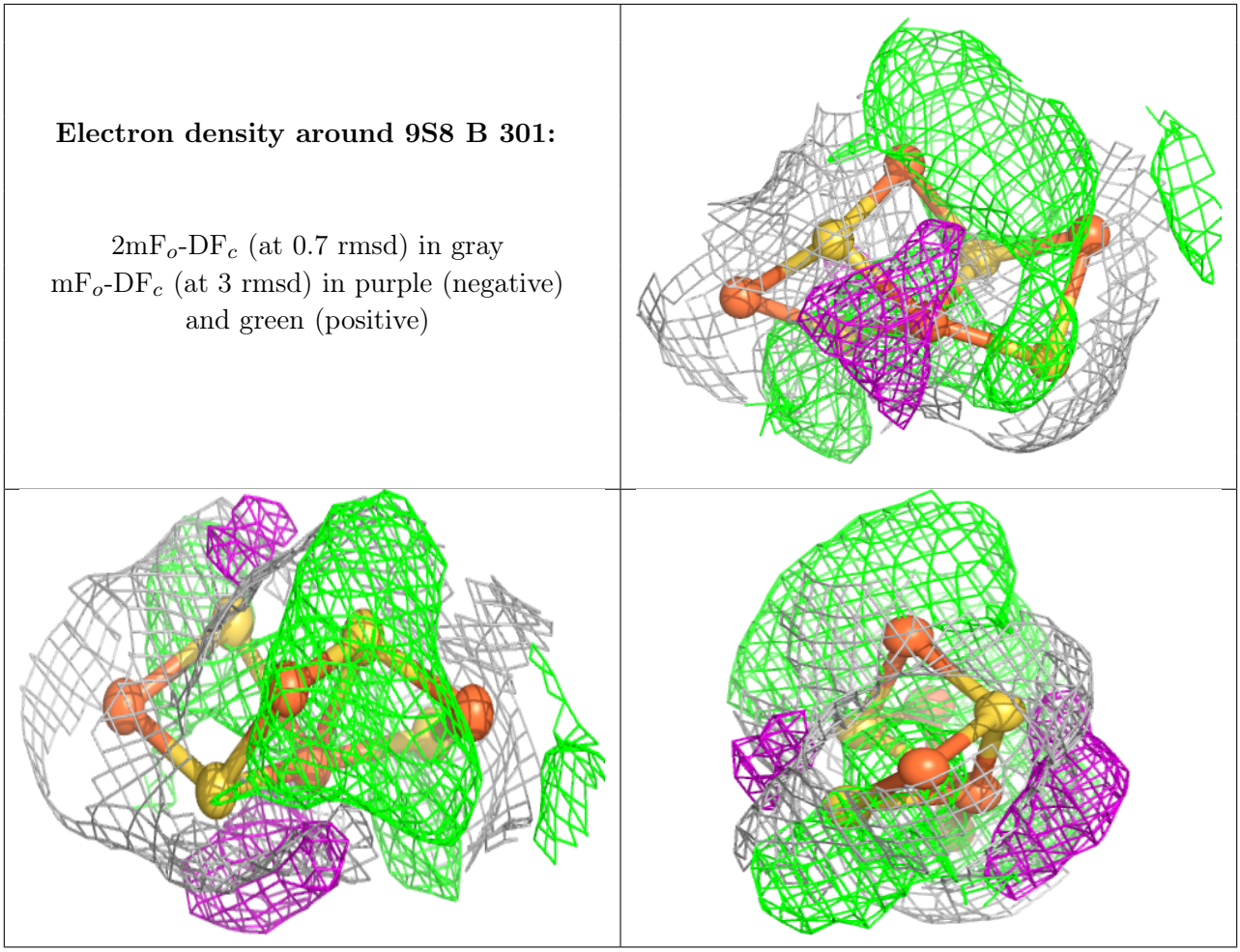






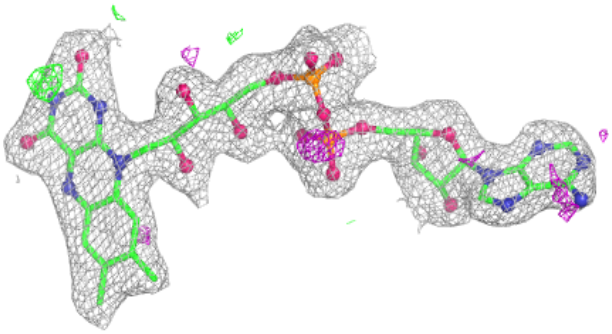
Electron density around 9S8 B 301:

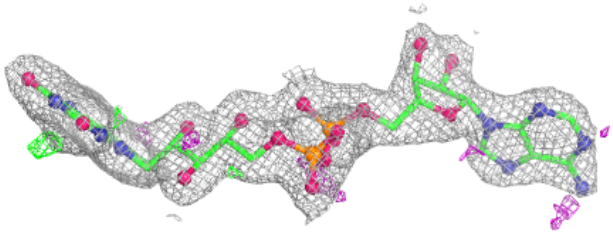
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

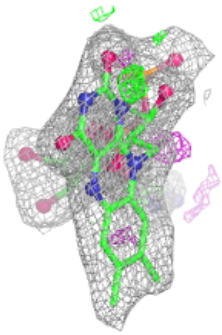


Electron density around FAD A 709:

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

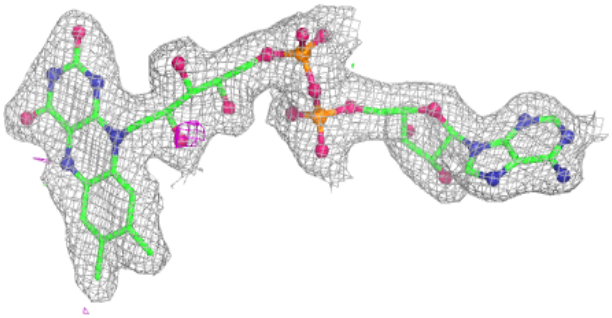


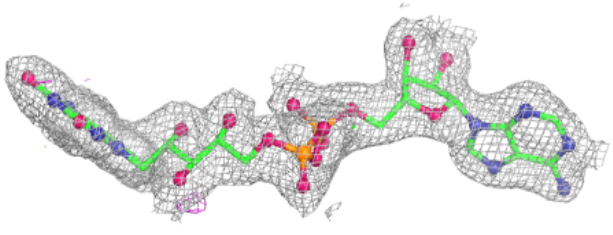


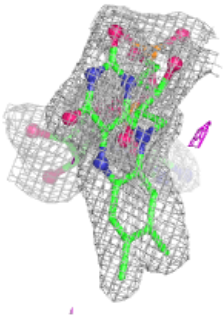


Electron density around FAD G 711:

$2mF_o$-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.