



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 09:47 AM UTC

PDB ID : 1SOX / pdb_00001sox
Title : SULFITE OXIDASE FROM CHICKEN LIVER
Authors : Kisker, C.; Schindelin, H.; Rees, D.C.
Deposited on : 1997-12-31
Resolution : 1.90 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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

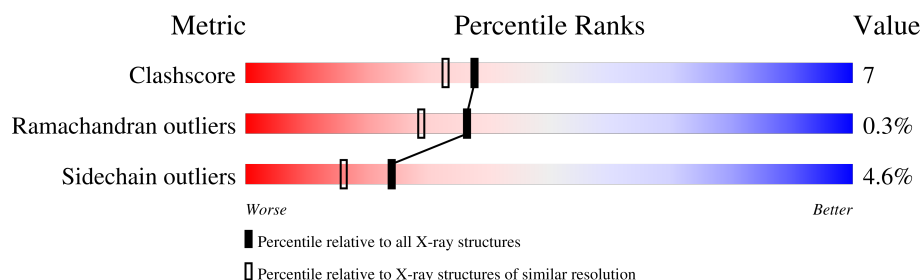
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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))
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	466	 80% 17% ..
1	B	466	 78% 17% ..

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 8189 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 SULFITE OXIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	463	Total	C	N	O	S	0	6	0
			3613	2293	654	659	7			
1	B	458	Total	C	N	O	S	0	6	0
			3577	2271	641	658	7			

There are 16 discrepancies between the modelled and reference sequences:

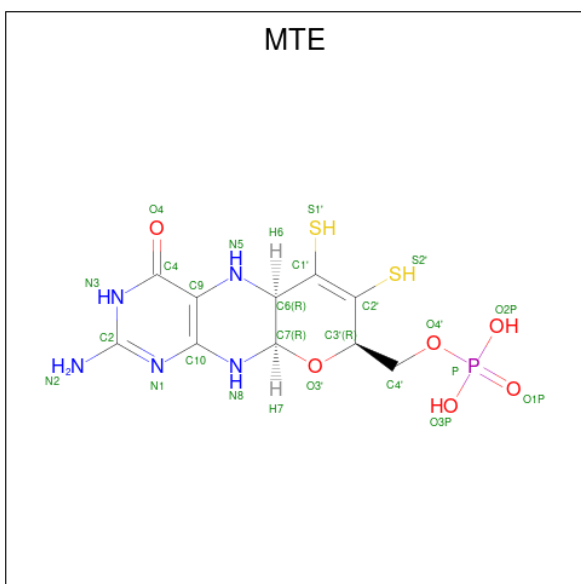
Chain	Residue	Modelled	Actual	Comment	Reference
A	6	GLU	ARG	conflict	UNP P07850
A	?	-	ARG	deletion	UNP P07850
A	161	GLY	ARG	conflict	UNP P07850
A	218	ARG	SER	conflict	UNP P07850
A	235	GLU	GLY	conflict	UNP P07850
A	236	TRP	GLU	conflict	UNP P07850
A	284	GLY	ARG	conflict	UNP P07850
A	316	HIS	ARG	conflict	UNP P07850
B	6	GLU	ARG	conflict	UNP P07850
B	?	-	ARG	deletion	UNP P07850
B	161	GLY	ARG	conflict	UNP P07850
B	218	ARG	SER	conflict	UNP P07850
B	235	GLU	GLY	conflict	UNP P07850
B	236	TRP	GLU	conflict	UNP P07850
B	284	GLY	ARG	conflict	UNP P07850
B	316	HIS	ARG	conflict	UNP P07850

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is PHOSPHONIC ACIDMONO-(2-AMINO-5,6-DIMERCAPTO-4-OXO-3,7,8A, 9,10,10A-HEXAHYDRO-4H-8-OXA-1,3,9,10-TETRAAZA-ANTHRACEN-7-YLMETHYL) ESTER (CCD ID: MTE) (formula: $C_{10}H_{14}N_5O_6PS_2$).

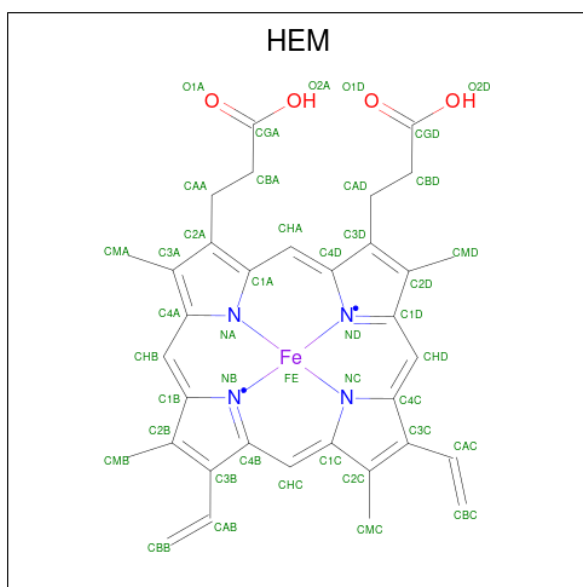


Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	S	0	0
			24	10	5	6	1	2		
3	B	1	Total	C	N	O	P	S	0	0
			24	10	5	6	1	2		

- Molecule 4 is MOLYBDENUM ATOM (CCD ID: MO) (formula: Mo).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Mo	0	0
			1	1		
4	B	1	Total	Mo	0	0
			1	1		

- Molecule 5 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



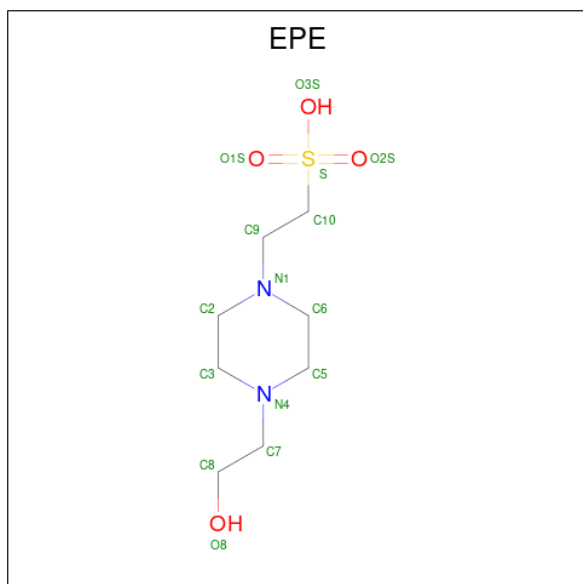
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
5	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 6 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 7 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (CCD ID: EPE) (formula: $C_8H_{18}N_2O_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	B	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	417	Total 417	O 417	0	0
8	B	405	Total 405	O 405	0	0

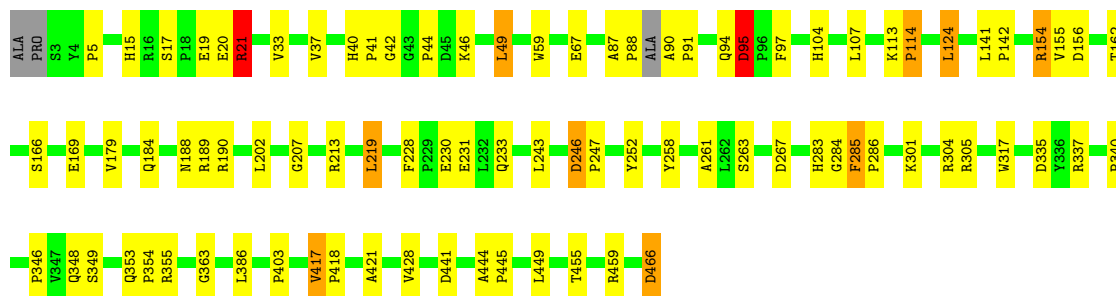
3 Residue-property plots

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. 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. 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.

Note EDS was not executed.

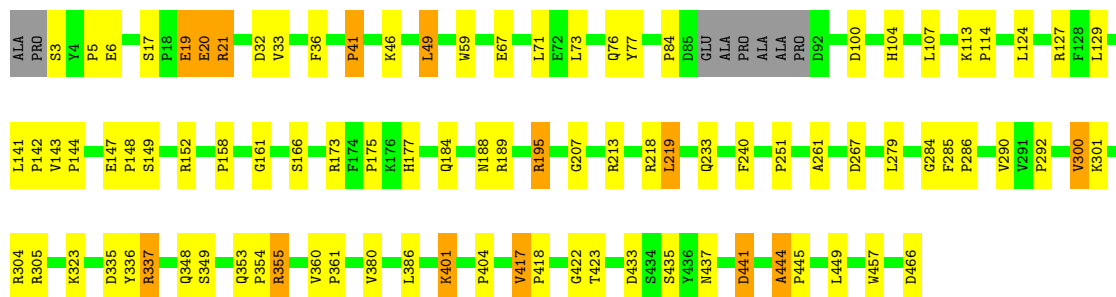
• Molecule 1: SULFITE OXIDASE

Chain A: 



• Molecule 1: SULFITE OXIDASE

Chain B: 



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	55.12Å 197.69Å 56.01Å 90.00° 94.23° 90.00°	Depositor
Resolution (Å)	50.00 – 1.90	Depositor
% Data completeness (in resolution range)	89.2 (50.00-1.90)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
Refinement program	X-PLOR 3.8	Depositor
R, R_{free}	0.175 , 0.220	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	8189	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: HEM, MO, MTE, SO4, EPE, GOL

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	0.61	0/3749	1.39	28/5128 (0.5%)
1	B	0.60	0/3712	1.41	39/5078 (0.8%)
All	All	0.61	0/7461	1.40	67/10206 (0.7%)

There are no bond length outliers.

All (67) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	353	GLN	CA-C-O	-13.37	107.55	119.72
1	B	353	GLN	CA-C-O	-11.98	108.81	119.72
1	B	285	PHE	O-C-N	10.56	133.46	121.32
1	B	113	LYS	CA-C-O	-10.52	107.49	120.74
1	B	113	LYS	O-C-N	9.71	130.15	121.32
1	A	113	LYS	CA-C-O	-9.69	109.03	120.08
1	A	285	PHE	O-C-N	9.20	131.90	121.32
1	B	285	PHE	CA-C-O	-9.17	107.59	120.16
1	B	114	PRO	N-CA-CB	8.31	111.74	102.60
1	B	354	PRO	N-CA-CB	7.83	111.22	102.60
1	A	354	PRO	N-CA-CB	7.77	111.15	102.60
1	A	355	ARG	CD-NE-CZ	7.69	135.17	124.40
1	A	285	PHE	CA-C-O	-7.58	109.78	120.16
1	A	417	VAL	N-CA-CB	-7.04	104.11	111.64
1	B	437	ASN	OD1-CG-ND2	7.01	129.61	122.60
1	B	337	ARG	CD-NE-CZ	6.90	134.06	124.40
1	A	114	PRO	N-CA-CB	6.83	110.11	102.60
1	B	361	PRO	N-CA-CB	6.80	106.76	103.22
1	A	87	ALA	CA-C-O	6.68	129.85	120.83
1	A	353	GLN	O-C-N	6.66	129.64	121.79
1	B	353	GLN	O-C-N	6.58	129.56	121.79
1	B	286	PRO	N-CA-CB	6.51	109.77	102.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	286	PRO	CA-N-CD	-6.43	102.49	111.50
1	A	286	PRO	N-CA-CB	6.37	109.60	102.60
1	A	403	PRO	N-CA-CB	6.37	106.75	103.19
1	B	213	ARG	CD-NE-CZ	6.36	133.31	124.40
1	B	355	ARG	CD-NE-CZ	6.27	133.18	124.40
1	B	360	VAL	CA-C-N	6.24	124.15	119.66
1	B	360	VAL	C-N-CA	6.24	124.15	119.66
1	B	67	GLU	CA-C-N	5.92	125.39	119.24
1	B	67	GLU	C-N-CA	5.92	125.39	119.24
1	B	444	ALA	CA-C-N	5.86	125.48	119.56
1	B	444	ALA	C-N-CA	5.86	125.48	119.56
1	A	286	PRO	N-CD-CG	5.83	110.79	103.80
1	A	113	LYS	O-C-N	5.79	128.19	121.47
1	B	100	ASP	CA-C-N	5.78	123.82	119.66
1	B	100	ASP	C-N-CA	5.78	123.82	119.66
1	B	435[A]	SER	N-CA-C	-5.73	105.88	112.92
1	B	435[B]	SER	N-CA-C	-5.73	105.88	112.92
1	B	104	HIS	CA-C-N	5.66	125.33	119.56
1	B	104	HIS	C-N-CA	5.66	125.33	119.56
1	A	283	HIS	CA-CB-CG	-5.65	108.15	113.80
1	B	158	PRO	N-CA-CB	5.62	108.17	103.17
1	A	104	HIS	CA-C-N	5.62	125.24	119.56
1	A	104	HIS	C-N-CA	5.62	125.24	119.56
1	B	433	ASP	CA-CB-CG	5.53	118.12	112.60
1	B	195	ARG	CA-C-N	5.52	127.63	120.56
1	B	195	ARG	C-N-CA	5.52	127.63	120.56
1	A	5	PRO	N-CA-CB	5.43	108.14	103.31
1	A	286	PRO	CA-N-CD	-5.42	103.91	111.50
1	A	88	PRO	N-CA-CB	5.31	108.84	103.00
1	B	286	PRO	N-CD-CG	5.28	110.13	103.80
1	A	95	ASP	CA-C-N	5.24	125.04	119.28
1	A	95	ASP	C-N-CA	5.24	125.04	119.28
1	A	354	PRO	N-CD-CG	5.21	110.06	103.80
1	B	305	ARG	NE-CZ-NH2	5.20	123.88	119.20
1	B	300	VAL	N-CA-C	5.16	115.89	108.93
1	B	417	VAL	N-CA-CB	-5.16	106.12	111.64
1	B	5	PRO	N-CA-CB	5.14	108.11	103.33
1	B	41	PRO	N-CA-CB	5.14	107.90	103.27
1	B	404	PRO	N-CA-CB	5.12	107.80	103.19
1	B	175	PRO	N-CA-CB	5.07	107.83	103.27
1	A	67	GLU	CA-C-N	5.03	125.05	119.32
1	A	67	GLU	C-N-CA	5.03	125.05	119.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	263	SER	CA-C-N	5.02	124.19	118.97
1	A	263	SER	C-N-CA	5.02	124.19	118.97
1	A	124	LEU	N-CA-C	5.00	117.11	111.11

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	3613	0	3503	52	0
1	B	3577	0	3456	49	0
2	A	10	0	0	0	0
2	B	10	0	0	0	0
3	A	24	0	10	0	0
3	B	24	0	10	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	43	0	30	2	0
5	B	43	0	30	3	0
6	A	6	0	8	0	0
7	B	15	0	18	2	0
8	A	417	0	0	0	0
8	B	405	0	0	0	0
All	All	8189	0	7065	100	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 7.

All (100) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:161:GLY:HA3	7:B:911:EPE:H52	1.57	0.83
1:B:386:LEU:HD11	1:B:417:VAL:HG21	1.67	0.76
1:A:124:LEU:HD11	1:A:207:GLY:HA2	1.69	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:305[B]:ARG:HG3	1:A:305[B]:ARG:HH11	1.51	0.73
1:B:335:ASP:OD1	1:B:337:ARG:HD3	1.89	0.71
1:A:304[B]:ARG:HH11	1:A:304[B]:ARG:HG2	1.55	0.71
1:B:189:ARG:HH11	1:B:348:GLN:HE22	1.41	0.67
1:A:305[B]:ARG:HG3	1:A:305[B]:ARG:NH1	2.08	0.67
1:A:335:ASP:OD1	1:A:337:ARG:HD3	1.96	0.66
1:A:189:ARG:HH11	1:A:348:GLN:HE22	1.45	0.63
1:A:33:VAL:HG11	5:A:502:HEM:HBB1	1.81	0.63
1:A:417:VAL:HG22	1:A:418:PRO:HD2	1.82	0.62
1:B:141:LEU:HB3	1:B:142:PRO:CD	2.30	0.62
1:A:348:GLN:HE21	1:A:349:SER:H	1.48	0.61
1:B:124:LEU:HD11	1:B:207:GLY:HA2	1.83	0.61
1:B:177[B]:HIS:CE1	1:B:218:ARG:HE	2.19	0.60
1:A:15:HIS:HB3	1:A:21:ARG:O	2.02	0.60
1:A:243:LEU:HD21	1:A:304[B]:ARG:HD3	1.83	0.59
1:B:348:GLN:HE21	1:B:349:SER:H	1.50	0.59
1:B:141:LEU:HB3	1:B:142:PRO:HD2	1.85	0.59
1:A:386:LEU:HD11	1:A:417:VAL:HG11	1.85	0.58
1:B:36:PHE:HB2	1:B:77:TYR:CZ	2.38	0.58
1:B:177[A]:HIS:HE1	1:B:267:ASP:OD1	1.86	0.57
1:A:190:ARG:HD2	1:A:202:LEU:O	2.05	0.57
1:B:21:ARG:HD3	1:B:32:ASP:OD2	2.05	0.57
1:A:154:ARG:HG2	1:A:305[B]:ARG:NH1	2.20	0.56
1:A:154:ARG:NH1	1:A:156:ASP:HB3	2.20	0.56
1:A:141:LEU:HB3	1:A:142:PRO:HD2	1.87	0.55
1:B:184:GLN:HE21	1:B:188:ASN:HB3	1.70	0.55
1:A:20:GLU:O	1:A:21:ARG:HB2	2.06	0.54
1:B:177[B]:HIS:CD2	1:B:218:ARG:HH21	2.25	0.54
1:A:17:SER:O	1:A:21:ARG:N	2.40	0.53
1:B:173:ARG:NH1	7:B:911:EPE:O2S	2.43	0.52
1:B:240:PHE:HB2	1:B:300:VAL:HG21	1.93	0.51
1:A:421:ALA:HB1	1:A:466:ASP:HA	1.92	0.51
1:B:189:ARG:HH11	1:B:348:GLN:NE2	2.06	0.51
1:A:189:ARG:HH11	1:A:348:GLN:NE2	2.09	0.51
1:B:19:GLU:H	1:B:19:GLU:CD	2.17	0.51
1:A:154:ARG:NE	1:A:305[B]:ARG:HH12	2.09	0.50
1:A:304[B]:ARG:HH11	1:A:304[B]:ARG:CG	2.24	0.50
1:A:179:VAL:HG21	1:A:267:ASP:HB3	1.92	0.50
1:A:184:GLN:HE21	1:A:188:ASN:HD22	1.60	0.50
1:B:41:PRO:HG2	5:B:502:HEM:C2A	2.47	0.49
1:A:90:ALA:N	1:A:91:PRO:HD3	2.27	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:149:SER:O	1:B:152:ARG:NH2	2.45	0.49
1:A:444:ALA:HB3	1:A:445:PRO:HD3	1.93	0.49
1:B:323:LYS:HD2	1:B:336:TYR:HD2	1.75	0.49
1:A:42:GLY:HA3	5:A:502:HEM:O2D	2.13	0.49
1:A:184:GLN:NE2	1:A:188:ASN:HD22	2.11	0.48
1:B:417:VAL:CG2	1:B:418:PRO:HD2	2.43	0.48
1:A:94:GLN:O	1:A:95:ASP:HB3	2.12	0.48
1:A:417:VAL:CG2	1:A:418:PRO:HD2	2.43	0.48
1:A:154:ARG:HH12	1:A:156:ASP:HB3	1.77	0.48
1:B:33:VAL:HG11	5:B:502:HEM:HBB1	1.95	0.48
1:B:127:ARG:HD3	1:B:129:LEU:O	2.14	0.48
1:B:17:SER:O	1:B:21:ARG:N	2.47	0.47
1:B:251:PRO:HG3	1:B:304:ARG:NH2	2.30	0.47
1:A:246:ASP:HB2	1:A:247:PRO:HD2	1.97	0.47
1:A:252:TYR:HA	1:A:317:TRP:CZ3	2.50	0.47
1:A:428:VAL:HG22	1:A:459:ARG:HG2	1.97	0.47
1:A:46:LYS:O	1:A:49:LEU:HB2	2.15	0.47
1:B:147[A]:GLU:OE2	1:B:148:PRO:HD2	2.15	0.47
1:B:422:GLY:N	1:B:466:ASP:HA	2.30	0.47
1:A:141:LEU:HB3	1:A:142:PRO:CD	2.45	0.46
1:B:386:LEU:HD11	1:B:417:VAL:CG2	2.41	0.46
1:A:166:SER:OG	1:A:169:GLU:HG3	2.17	0.45
1:A:97:PHE:CE2	1:A:114:PRO:HA	2.52	0.45
1:A:154:ARG:CZ	1:A:305[B]:ARG:HH12	2.30	0.45
1:B:417:VAL:HG22	1:B:418:PRO:HD2	1.99	0.45
1:B:441:ASP:HA	1:B:457:TRP:CE2	2.52	0.45
1:A:243:LEU:CD2	1:A:304[B]:ARG:HD3	2.47	0.45
1:B:284:GLY:HA2	1:B:301:LYS:HB3	1.99	0.45
1:A:19:GLU:H	1:A:19:GLU:CD	2.25	0.45
1:B:184:GLN:NE2	1:B:188:ASN:HD22	2.15	0.45
1:A:154:ARG:NH1	1:A:162:THR:HG23	2.33	0.44
1:B:46:LYS:O	1:B:49:LEU:HB2	2.18	0.44
1:A:37:VAL:HG12	1:A:44:PRO:HB3	2.00	0.44
1:B:323:LYS:HD2	1:B:336:TYR:CD2	2.53	0.43
1:A:228:PHE:HB3	1:A:258:TYR:OH	2.18	0.43
1:A:304[B]:ARG:CG	1:A:304[B]:ARG:NH1	2.81	0.43
1:A:346:PRO:HG2	1:A:455:THR:HG21	2.00	0.43
1:B:20:GLU:O	1:B:21:ARG:HB2	2.18	0.43
1:B:290:VAL:O	1:B:292:PRO:HD3	2.19	0.43
1:B:401:LYS:HA	1:B:401:LYS:HE2	2.00	0.42
1:A:284:GLY:HA2	1:A:301:LYS:HB3	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:363:GLY:O	1:A:418:PRO:HA	2.19	0.42
1:B:189:ARG:NH1	1:B:348:GLN:HE22	2.14	0.42
1:B:219:LEU:HD13	1:B:261:ALA:HB1	2.02	0.41
1:B:46:LYS:O	5:B:502:HEM:HBC2	2.20	0.41
1:B:36:PHE:HE1	1:B:73:LEU:HD21	1.85	0.41
1:B:386:LEU:HD11	1:B:417:VAL:HG11	2.02	0.41
1:B:444:ALA:HB3	1:B:445:PRO:HD3	2.03	0.41
1:A:219:LEU:HD13	1:A:261:ALA:HB1	2.03	0.41
1:B:36:PHE:CE1	1:B:73:LEU:HD21	2.55	0.41
1:A:340:PRO:HG3	1:B:380:VAL:HG11	2.03	0.40
1:B:20:GLU:O	1:B:21:ARG:CB	2.69	0.40
1:A:40:HIS:HA	1:A:41:PRO:HD3	1.91	0.40
1:B:152:ARG:HE	1:B:166:SER:HA	1.86	0.40
1:A:284:GLY:O	1:A:285:PHE:C	2.65	0.40
1:B:143:VAL:HA	1:B:144:PRO:HD3	1.92	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	465/466 (100%)	453 (97%)	11 (2%)	1 (0%)	43	36
1	B	460/466 (99%)	448 (97%)	10 (2%)	2 (0%)	30	22
All	All	925/932 (99%)	901 (97%)	21 (2%)	3 (0%)	36	29

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	21	ARG
1	B	21	ARG
1	B	84	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	379/377 (100%)	363 (96%)	16 (4%)	26	19
1	B	377/377 (100%)	359 (95%)	18 (5%)	23	15
All	All	756/754 (100%)	722 (96%)	34 (4%)	24	17

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	21	ARG
1	A	49	LEU
1	A	59	TRP
1	A	95	ASP
1	A	107	LEU
1	A	154	ARG
1	A	155	VAL
1	A	213	ARG
1	A	219	LEU
1	A	230	GLU
1	A	231	GLU
1	A	233	GLN
1	A	246	ASP
1	A	441	ASP
1	A	449	LEU
1	A	466	ASP
1	B	3	SER
1	B	6	GLU
1	B	19	GLU
1	B	20	GLU
1	B	49	LEU
1	B	59	TRP
1	B	71	LEU
1	B	76	GLN
1	B	107	LEU
1	B	195	ARG
1	B	219	LEU

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Mol	Chain	Res	Type
1	B	233	GLN
1	B	279	LEU
1	B	355	ARG
1	B	401	LYS
1	B	423	THR
1	B	441	ASP
1	B	449	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	184	GLN
1	A	233	GLN
1	A	319	GLN
1	A	343	GLN
1	A	348	GLN
1	A	465	GLN
1	B	184	GLN
1	B	319	GLN
1	B	343	GLN
1	B	348	GLN
1	B	353	GLN
1	B	437	ASN
1	B	465	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 12 ligands modelled in this entry, 2 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
5	HEM	B	502	1	50,50,50	1.25	8 (16%)	67,82,82	1.01	3 (4%)
7	EPE	B	911	-	15,15,15	1.81	2 (13%)	19,20,20	2.06	7 (36%)
6	GOL	A	912	-	5,5,5	0.13	0	5,5,5	0.28	0
2	SO4	A	503	-	4,4,4	1.00	0	6,6,6	0.24	0
3	MTE	B	501	4	21,26,26	5.83	10 (47%)	19,40,40	3.32	10 (52%)
2	SO4	B	503	-	4,4,4	0.97	0	6,6,6	0.41	0
3	MTE	A	501	4	21,26,26	5.88	11 (52%)	19,40,40	3.57	10 (52%)
2	SO4	B	504	-	4,4,4	1.04	0	6,6,6	0.27	0
2	SO4	A	504	-	4,4,4	1.06	0	6,6,6	0.21	0
5	HEM	A	502	1	50,50,50	1.28	9 (18%)	67,82,82	1.18	2 (2%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	HEM	B	502	1	-	3/14/54/54	-
3	MTE	A	501	4	-	0/6/34/34	0/3/3/3
5	HEM	A	502	1	-	6/14/54/54	-
7	EPE	B	911	-	-	3/9/19/19	0/1/1/1
3	MTE	B	501	4	-	0/6/34/34	0/3/3/3
6	GOL	A	912	-	-	2/4/4/4	-

All (40) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	501	MTE	C7-C6	15.05	1.65	1.53
3	A	501	MTE	C7-C6	15.00	1.65	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	501	MTE	C9-C10	13.45	1.61	1.38
3	A	501	MTE	C9-C10	13.01	1.60	1.38
3	A	501	MTE	C6-N5	10.51	1.58	1.46
3	B	501	MTE	C6-N5	10.40	1.58	1.46
3	A	501	MTE	C4'-C3'	-7.95	1.40	1.51
3	B	501	MTE	C4'-C3'	-7.29	1.41	1.51
3	A	501	MTE	P-O4'	-7.02	1.38	1.60
3	B	501	MTE	P-O4'	-6.43	1.40	1.60
7	B	911	EPE	O3S-S	5.62	1.68	1.47
3	A	501	MTE	C9-N5	5.50	1.50	1.37
3	B	501	MTE	C9-N5	5.45	1.50	1.37
3	A	501	MTE	P-O3P	-5.30	1.35	1.54
3	B	501	MTE	P-O3P	-4.91	1.36	1.54
3	B	501	MTE	O4-C4	4.86	1.32	1.23
3	A	501	MTE	O4-C4	4.11	1.31	1.23
3	A	501	MTE	C10-N8	3.53	1.39	1.35
7	B	911	EPE	C10-S	3.28	1.82	1.77
5	A	502	HEM	CAB-C3B	3.08	1.55	1.47
5	B	502	HEM	CAB-C3B	2.76	1.54	1.47
5	A	502	HEM	CAC-C3C	2.75	1.54	1.47
5	B	502	HEM	FE-NB	2.69	2.03	1.94
3	B	501	MTE	C2-N1	2.43	1.39	1.33
5	A	502	HEM	FE-ND	2.40	2.02	1.94
5	B	502	HEM	FE-NC	2.32	2.02	1.95
3	A	501	MTE	C2-N1	2.30	1.38	1.33
5	B	502	HEM	CAC-C3C	2.27	1.53	1.47
5	A	502	HEM	CMB-C2B	2.22	1.55	1.50
5	B	502	HEM	FE-ND	2.19	2.01	1.94
5	B	502	HEM	CMA-C3A	2.15	1.55	1.50
3	A	501	MTE	C9-C4	2.09	1.48	1.41
5	A	502	HEM	FE-NC	2.08	2.02	1.95
5	B	502	HEM	CMB-C2B	2.08	1.55	1.50
5	B	502	HEM	CMC-C2C	2.05	1.55	1.50
5	A	502	HEM	CMA-C3A	2.05	1.55	1.50
5	A	502	HEM	FE-NA	2.04	2.01	1.95
3	B	501	MTE	C9-C4	2.02	1.47	1.41
5	A	502	HEM	C2A-C3A	-2.02	1.33	1.38
5	A	502	HEM	CMC-C2C	2.02	1.54	1.50

All (32) bond angle outliers are listed below:

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	501	MTE	C2-N1-C10	8.36	128.12	113.36
3	B	501	MTE	C2-N1-C10	7.69	126.93	113.36
5	A	502	HEM	CBD-CAD-C3D	6.59	130.76	112.53
3	B	501	MTE	O3'-C7-N8	-6.40	102.80	108.61
3	A	501	MTE	O3'-C7-C6	-6.28	104.77	108.96
3	A	501	MTE	O3'-C7-N8	-6.10	103.08	108.61
3	B	501	MTE	O4-C4-C9	-5.07	115.04	127.26
3	A	501	MTE	O4-C4-C9	-5.02	115.15	127.26
3	B	501	MTE	O3'-C7-C6	-4.74	105.80	108.96
7	B	911	EPE	O3S-S-O1S	-4.53	100.06	111.40
3	A	501	MTE	C4-C9-N5	4.12	127.52	116.27
3	B	501	MTE	C4-C9-N5	4.10	127.44	116.27
5	B	502	HEM	CAD-CBD-CGD	3.49	122.92	113.67
3	A	501	MTE	N2-C2-N3	3.44	124.02	116.76
3	A	501	MTE	C9-C4-N3	3.16	120.80	112.13
7	B	911	EPE	O1S-S-C10	3.09	111.40	106.73
7	B	911	EPE	C3-C2-N1	2.95	116.60	110.65
3	A	501	MTE	C7-C6-N5	2.95	110.76	107.87
3	B	501	MTE	O4-C4-N3	2.90	125.58	120.11
7	B	911	EPE	C6-N1-C2	2.79	114.86	108.84
3	B	501	MTE	N3-C2-N1	-2.71	118.36	123.32
3	B	501	MTE	C9-C4-N3	2.63	119.36	112.13
7	B	911	EPE	C2-C3-N4	2.61	115.92	110.65
3	B	501	MTE	N2-C2-N3	2.55	122.14	116.76
3	A	501	MTE	O2P-P-O4'	2.42	112.97	106.67
7	B	911	EPE	C5-C6-N1	2.41	115.51	110.65
5	B	502	HEM	O1A-CGA-CBA	-2.33	115.69	123.09
3	B	501	MTE	O3P-P-O4'	2.33	112.74	106.67
3	A	501	MTE	O4-C4-N3	2.09	124.05	120.11
7	B	911	EPE	O2S-S-C10	2.05	109.82	106.73
5	A	502	HEM	CAD-CBD-CGD	2.02	119.03	113.67
5	B	502	HEM	O2A-CGA-O1A	2.01	128.51	123.33

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	912	GOL	O1-C1-C2-O2
6	A	912	GOL	O1-C1-C2-C3
7	B	911	EPE	C10-C9-N1-C6

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Mol	Chain	Res	Type	Atoms
7	B	911	EPE	C9-C10-S-O2S
5	B	502	HEM	C3D-CAD-CBD-CGD
7	B	911	EPE	C9-C10-S-O3S
5	A	502	HEM	C2B-C3B-CAB-CBB
5	A	502	HEM	C4B-C3B-CAB-CBB
5	A	502	HEM	CAD-CBD-CGD-O1D
5	A	502	HEM	CAD-CBD-CGD-O2D
5	A	502	HEM	CAA-CBA-CGA-O2A
5	A	502	HEM	CAA-CBA-CGA-O1A
5	B	502	HEM	CAA-CBA-CGA-O2A
5	B	502	HEM	CAA-CBA-CGA-O1A

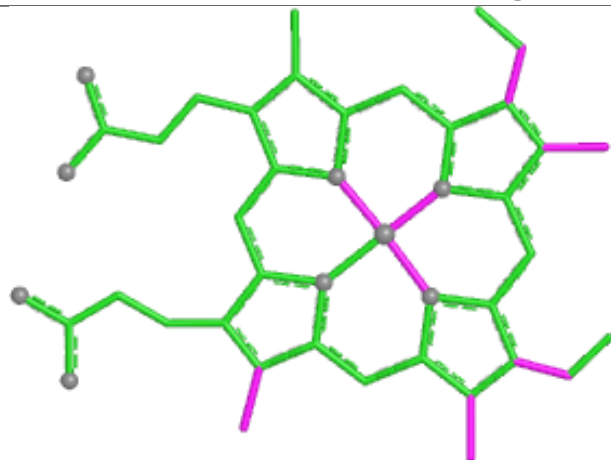
There are no ring outliers.

3 monomers are involved in 7 short contacts:

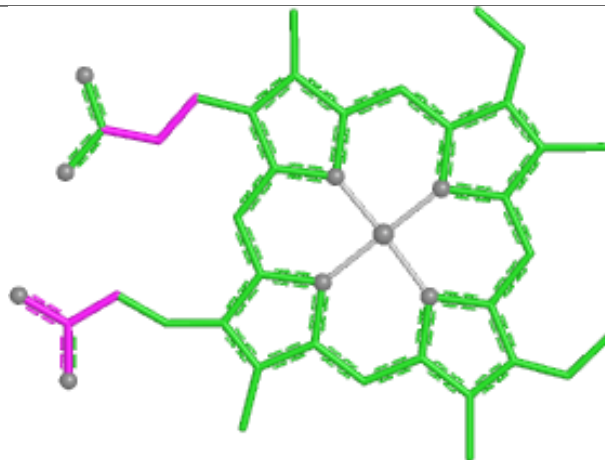
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	502	HEM	3	0
7	B	911	EPE	2	0
5	A	502	HEM	2	0

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.

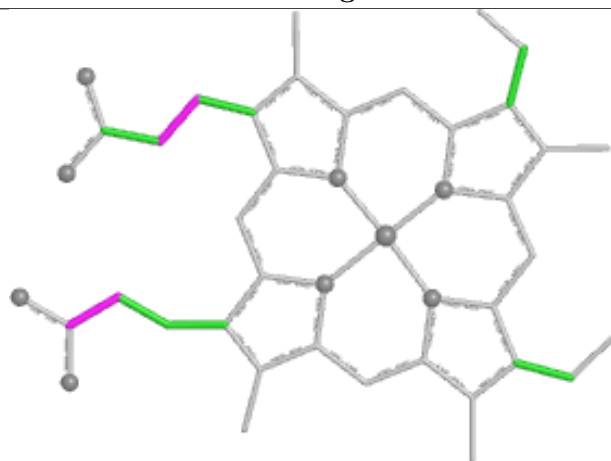
Ligand HEM B 502



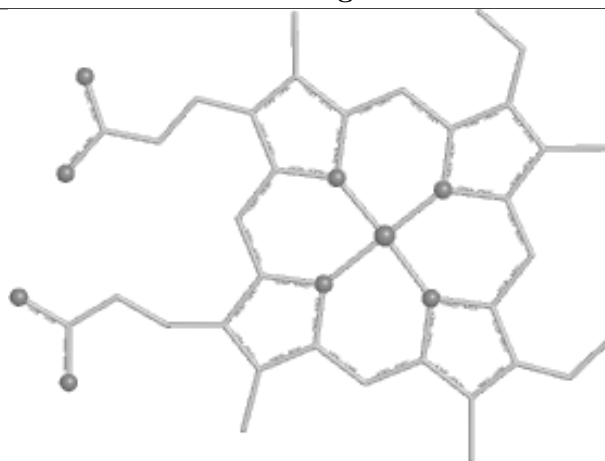
Bond lengths



Bond angles

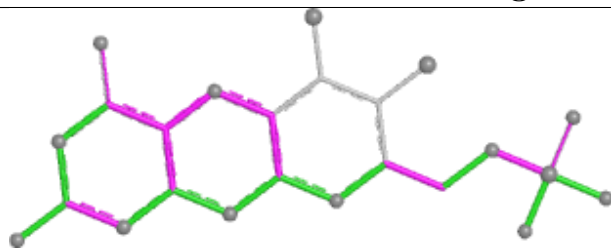


Torsions

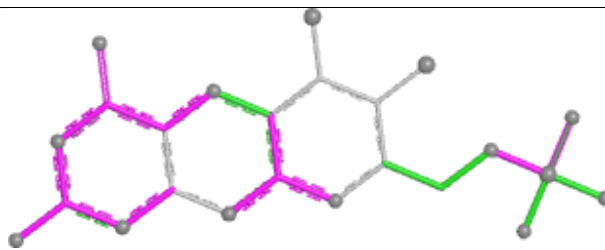


Rings

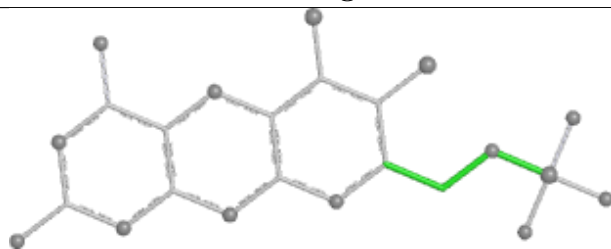
Ligand MTE B 501



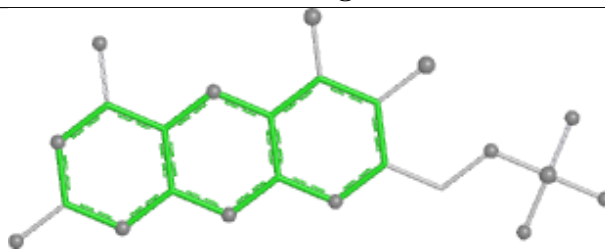
Bond lengths



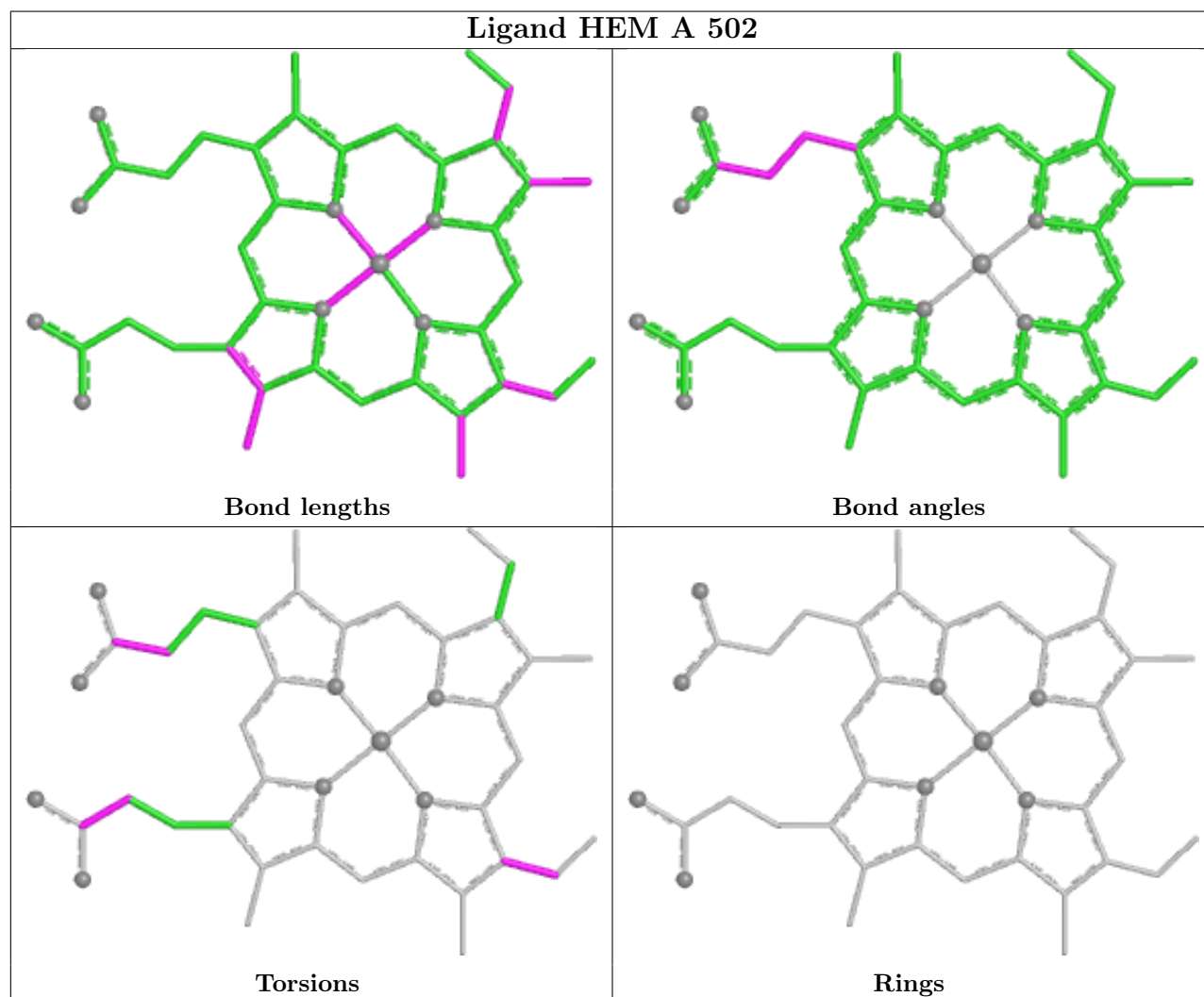
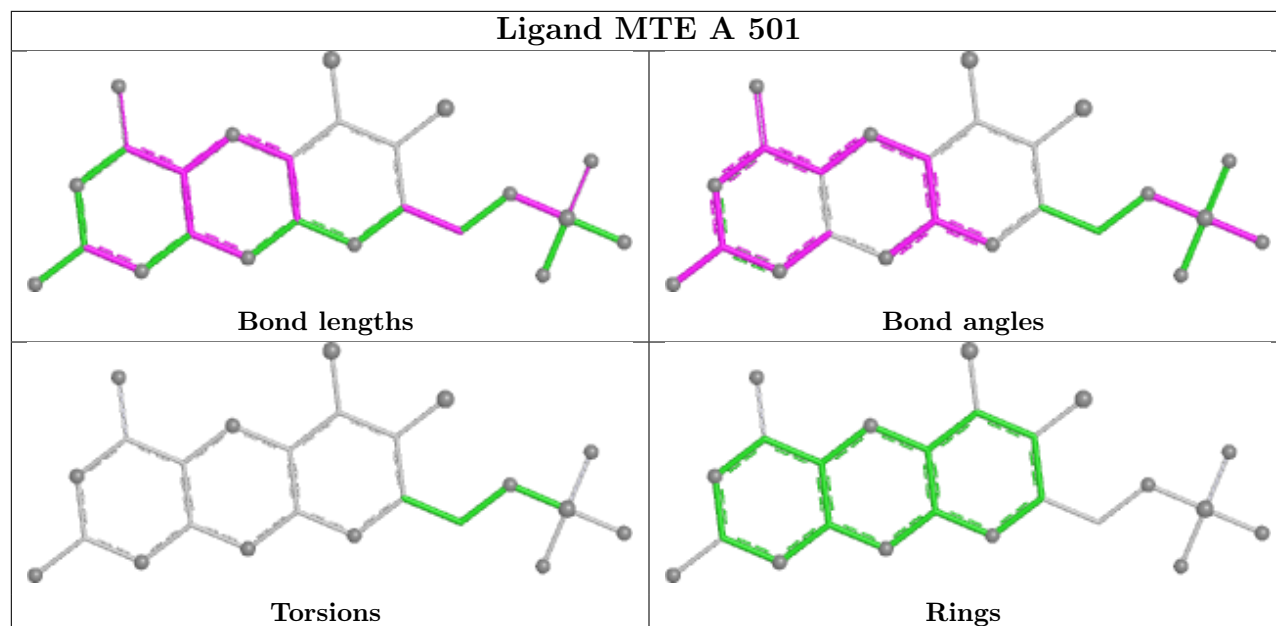
Bond angles



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

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.