Full wwPDB X-ray Structure Validation Report (i)

Jan 28, 2024 – 12:44 AM EST

PDB ID	:	1DMJ
Title	:	BOVINE ENDOTHELIAL NITRIC OXIDE SYNTHASE HEME DOMAIN COMPLEXED WITH 5,6-CYCLIC-TETRAHYDROPTERIDINE
Authors	:	Kotsonis, P.; Frohlich, L.G.; Raman, C.S.; Li, H.; Berg, M.; Gerwig, R.; Groehn, V.; Kang, Y.; Al-Masoudi, N.; Taghavi-Moghadam, S.; Mohr, D.; Munch, U.; Schnabel, J.; Martasek, P.; Masters, B.S.; Strobel, H.; Poulos, T.; Matter, H.; Pfleiderer, W.; Schmidt, H.H.
Deposited on	:	1999-12-14
Resolution	:	2.35 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 (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:

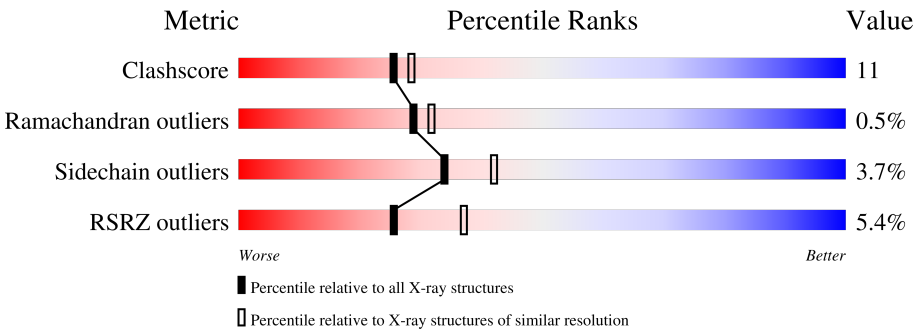
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)

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 2.35 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	141614	1232 (2.36-2.36)
Ramachandran outliers	138981	1211 (2.36-2.36)
Sidechain outliers	138945	1212 (2.36-2.36)
RSRZ outliers	127900	1150 (2.36-2.36)

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	444	6% 75% 17% • 6%
1	B	444	4% 74% 18% • 7%

Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36

2 Entry composition (i)

There are 9 unique types of molecules in this entry. The entry contains 7168 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.

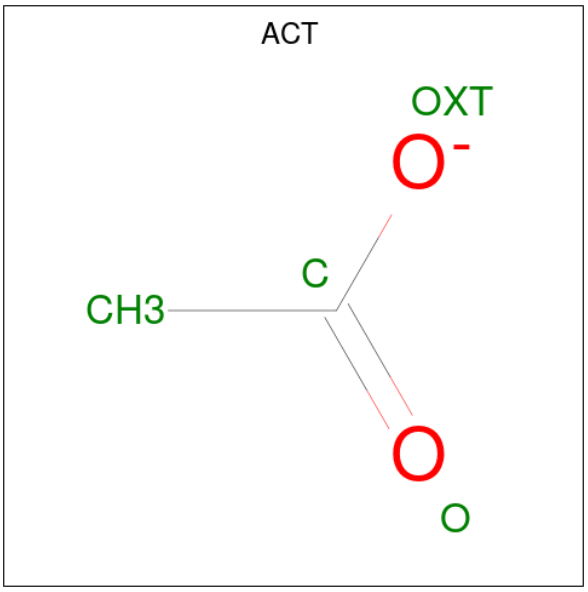
- Molecule 1 is a protein called NITRIC OXIDE SYNTHASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	416	3302	2099	584	603	16	0	0	0
1	B	414	3291	2092	582	601	16	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	100	ARG	CYS	conflict	UNP P29473
B	100	ARG	CYS	conflict	UNP P29473

- Molecule 2 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



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

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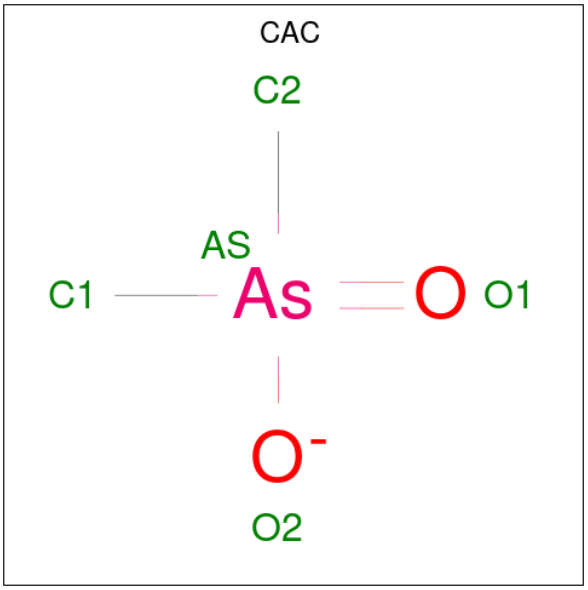
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	A	1	4	2	2	0	0
			Total	C	O		
2	B	1	4	2	2	0	0
			Total	C	O		
2	B	1	4	2	2	0	0

- Molecule 3 is CACODYLATE ION (three-letter code: CAC) (formula: $C_2H_6AsO_2$).

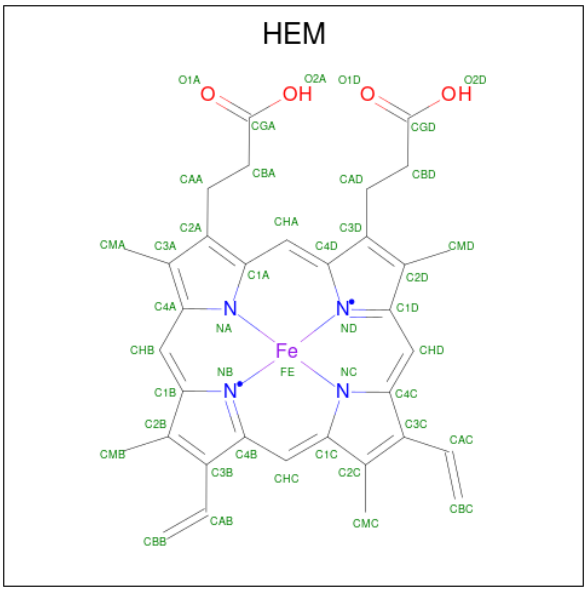


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	As	C		
3	A	1	3	1	2	0	0
			Total	As	C		
3	B	1	3	1	2	0	0

- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

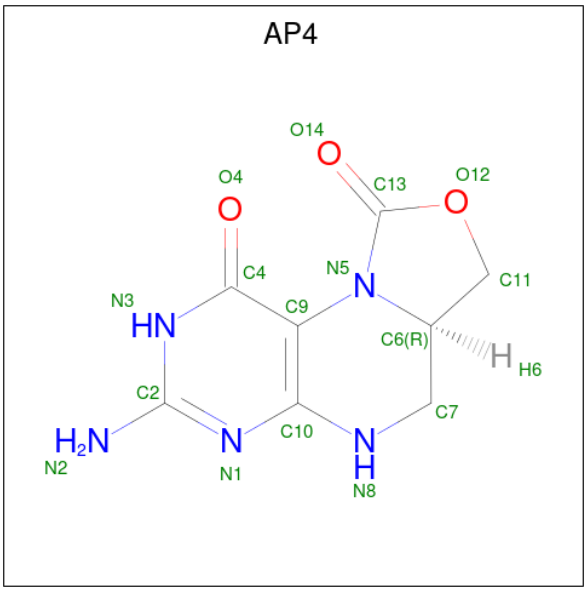
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Zn		
4	A	1	1	1	0	0

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



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

- Molecule 6 is 7-AMINO-3,3A,4,5-TETRAHYDRO-8H-2-OXA-5,6,8,9B-TETRAAZA-CYCLOPENTA[A]NAPHTHALENE-1,9-DIONE (three-letter code: AP4) (formula: $C_8H_9N_5O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
6	A	1	16	8	5	3	0	0

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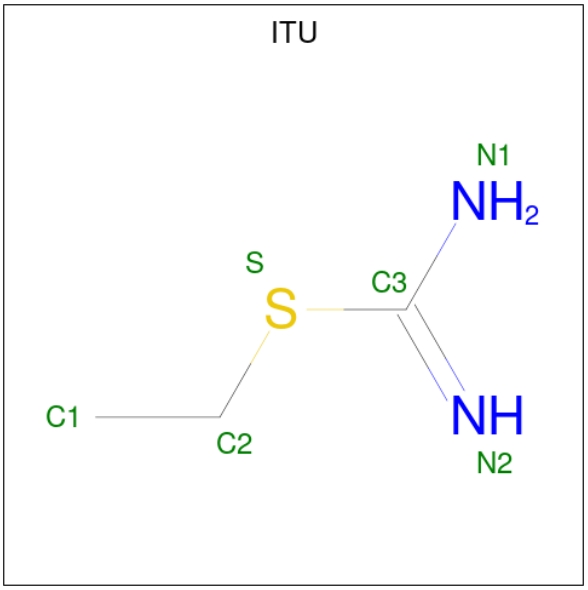
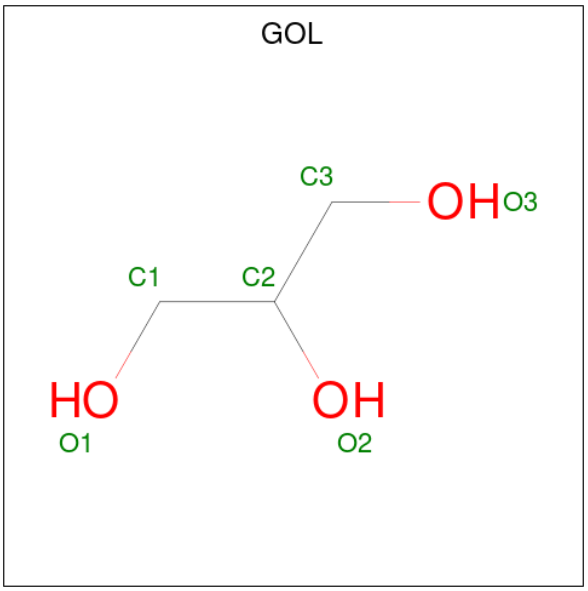
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
6	B	1	16	8	5	3	0	0

- Molecule 7 is ETHYLISOTHIOUREA (three-letter code: ITU) (formula: $C_3H_8N_2S$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	S		
7	A	1	6	3	2	1	0	0
			Total	C	N	S		
7	B	1	6	3	2	1	0	0

- Molecule 8 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 9 is water.

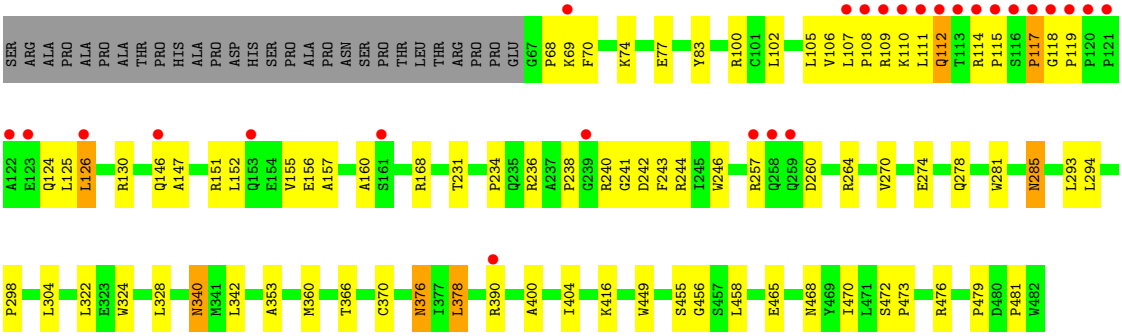
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
9	A	214	214	214	0	0
			Total	O		
9	B	196	196	196	0	0

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.

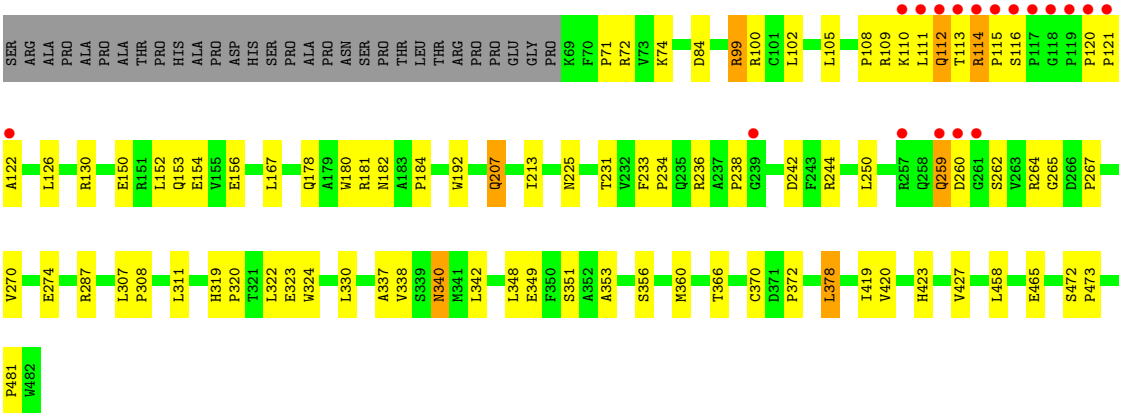
• Molecule 1: NITRIC OXIDE SYNTHASE

Chain A:



• Molecule 1: NITRIC OXIDE SYNTHASE

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	57.90Å 106.49Å 156.55Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.38 – 2.35 48.38 – 2.35	Depositor EDS
% Data completeness (in resolution range)	96.6 (48.38-2.35) 93.7 (48.38-2.35)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	4.43 (at 2.34Å)	Xtriage
Refinement program	CNS 0.9	Depositor
R, R_{free}	0.187 , 0.228 0.187 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å^2)	28.0	Xtriage
Anisotropy	0.382	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 48.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	7168	wwPDB-VP
Average B, all atoms (Å^2)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.38% 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 (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, CAC, AP4, ITU, ACT, HEM, ZN

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	0.35	0/3397	0.62	1/4631 (0.0%)
1	B	0.37	1/3385 (0.0%)	0.63	1/4614 (0.0%)
All	All	0.36	1/6782 (0.0%)	0.62	2/9245 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	122	ALA	CA-CB	-6.71	1.38	1.52

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	360	MET	N-CA-C	-5.38	96.46	111.00
1	A	360	MET	N-CA-C	-5.17	97.05	111.00

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	3302	0	3215	77	0
1	B	3291	0	3205	65	0
2	A	8	0	6	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	8	0	6	0	0
3	A	3	0	0	1	0
3	B	3	0	0	0	0
4	A	1	0	0	0	0
5	A	43	0	30	1	0
5	B	43	0	30	0	0
6	A	16	0	8	1	0
6	B	16	0	8	0	0
7	A	6	0	7	0	0
7	B	6	0	7	1	0
8	A	6	0	8	1	0
8	B	6	0	8	1	0
9	A	214	0	0	4	0
9	B	196	0	0	10	0
All	All	7168	0	6538	141	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 11.

All (141) 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:A:115:PRO:HD3	1:A:479:PRO:HG2	1.16	1.15
1:A:105:LEU:HB2	1:A:108:PRO:HG3	1.47	0.96
1:B:108:PRO:HB3	1:B:111:LEU:HD13	1.54	0.89
1:A:119:PRO:HG3	1:A:238:PRO:HB2	1.57	0.86
1:B:250:LEU:HD22	9:B:2962:HOH:O	1.77	0.84
1:B:116:SER:O	1:B:238:PRO:HA	1.80	0.81
1:A:111:LEU:HD21	1:A:470:ILE:HD13	1.64	0.80
1:A:111:LEU:HD22	1:A:476:ARG:NH1	1.95	0.80
1:B:114:ARG:HA	1:B:114:ARG:HH11	1.46	0.78
1:A:115:PRO:HD3	1:A:479:PRO:CG	2.10	0.72
1:A:119:PRO:HG3	1:A:238:PRO:CB	2.20	0.71
1:A:111:LEU:CD2	1:A:470:ILE:HD13	2.22	0.69
1:B:152:LEU:O	1:B:156:GLU:HG3	1.94	0.67
1:B:366:THR:O	1:B:370:CYS:HB2	1.95	0.67
1:A:240:ARG:HD3	1:A:298:PRO:HB3	1.78	0.65
1:A:257:ARG:HG3	1:A:257:ARG:HH11	1.61	0.65
1:B:338:VAL:HG23	7:B:2800:ITU:H11	1.78	0.65
1:A:69:LYS:O	1:A:69:LYS:HG2	1.95	0.65
1:A:126:LEU:HD11	1:A:156:GLU:HB3	1.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:126:LEU:HD12	1:A:130:ARG:HE	1.61	0.65
1:B:114:ARG:HH11	1:B:115:PRO:HD3	1.61	0.64
1:A:378:LEU:HB2	9:A:2083:HOH:O	1.97	0.64
1:B:236:ARG:HD2	1:B:242:ASP:OD2	1.98	0.64
1:A:146:GLN:NE2	1:A:147:ALA:H	1.96	0.63
1:A:105:LEU:CB	1:A:108:PRO:HG3	2.26	0.63
1:B:423:HIS:O	1:B:427:VAL:HG23	2.00	0.62
1:A:77:GLU:HG3	1:B:372:PRO:HG2	1.81	0.61
1:A:274:GLU:O	1:A:278:GLN:HG3	1.98	0.61
1:B:378:LEU:HB2	9:B:3020:HOH:O	1.99	0.61
1:A:168:ARG:HD2	9:A:2002:HOH:O	2.01	0.60
1:A:111:LEU:CD2	1:A:476:ARG:HH12	2.15	0.60
1:B:113:THR:HG21	1:B:342:LEU:CD1	2.33	0.59
1:A:111:LEU:CD2	1:A:476:ARG:NH1	2.64	0.59
1:B:126:LEU:O	1:B:130:ARG:HG3	2.02	0.59
1:B:114:ARG:HA	1:B:114:ARG:NH1	2.16	0.58
1:B:308:PRO:HD2	1:B:311:LEU:HD12	1.86	0.58
1:A:465:GLU:HB3	1:B:105:LEU:HD22	1.84	0.58
1:B:370:CYS:SG	1:B:378:LEU:HD13	2.44	0.58
2:A:1860:ACT:H2	5:A:1500:HEM:HMB3	1.87	0.56
1:B:114:ARG:HD3	1:B:115:PRO:CD	2.35	0.56
1:A:70:PHE:O	1:B:109:ARG:NH2	2.38	0.56
1:A:115:PRO:O	1:A:117:PRO:HD3	2.07	0.55
1:B:207:GLN:HB2	9:B:3056:HOH:O	2.07	0.55
8:B:2880:GOL:H32	9:B:2968:HOH:O	2.07	0.55
1:A:151:ARG:HD3	1:A:168:ARG:HE	1.72	0.55
1:A:340:ASN:H	1:A:340:ASN:HD22	1.56	0.53
1:A:111:LEU:HD21	1:A:470:ILE:HG21	1.88	0.53
1:B:108:PRO:CB	1:B:111:LEU:HB2	2.39	0.53
1:A:115:PRO:C	1:A:117:PRO:HD3	2.29	0.53
1:B:236:ARG:HD3	1:B:351:SER:HB3	1.91	0.53
1:B:182:ASN:O	1:B:184:PRO:HD3	2.10	0.52
1:A:257:ARG:HG3	1:A:257:ARG:NH1	2.24	0.52
1:A:240:ARG:HD2	1:A:241:GLY:O	2.10	0.52
1:B:114:ARG:NH1	1:B:115:PRO:HD3	2.24	0.52
1:B:72:ARG:NH1	9:B:3010:HOH:O	2.42	0.51
1:B:99:ARG:HB2	1:B:99:ARG:HH11	1.74	0.51
1:A:108:PRO:HG2	9:A:1906:HOH:O	2.10	0.51
1:A:246:TRP:HB2	1:A:294:LEU:HB3	1.93	0.51
1:A:109:ARG:O	1:A:110:LYS:HD2	2.11	0.50
1:A:376:ASN:ND2	9:A:2083:HOH:O	2.34	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:115:PRO:CD	1:A:479:PRO:HG2	2.11	0.50
1:A:152:LEU:O	1:A:156:GLU:HG2	2.11	0.50
1:B:340:ASN:HD22	1:B:340:ASN:H	1.58	0.50
1:A:244:ARG:NH2	1:A:481:PRO:HD3	2.27	0.50
1:A:324:TRP:HB2	3:A:950:CAC:C2	2.41	0.50
1:A:106:VAL:C	1:A:108:PRO:HD3	2.31	0.50
1:B:260:ASP:C	1:B:262:SER:H	2.15	0.49
1:A:270:VAL:O	1:A:274:GLU:HG3	2.12	0.49
1:B:114:ARG:HD3	1:B:115:PRO:HD2	1.93	0.49
1:A:472:SER:HA	1:A:473:PRO:C	2.33	0.49
1:A:111:LEU:HD22	1:A:476:ARG:HH11	1.73	0.48
1:B:231:THR:O	1:B:353:ALA:HA	2.14	0.48
1:A:231:THR:O	1:A:353:ALA:HA	2.14	0.48
1:B:74:LYS:O	1:B:465:GLU:HG3	2.13	0.48
1:B:244:ARG:NH2	1:B:481:PRO:HD3	2.28	0.48
1:A:74:LYS:O	1:A:465:GLU:HG3	2.13	0.48
1:B:337:ALA:HB2	1:B:356:SER:HB3	1.96	0.48
1:B:113:THR:HG21	1:B:342:LEU:HD11	1.96	0.47
1:B:472:SER:HA	1:B:473:PRO:C	2.34	0.47
1:A:342:LEU:C	1:A:342:LEU:HD23	2.35	0.47
1:A:126:LEU:HB3	1:A:130:ARG:NH2	2.29	0.47
1:B:111:LEU:O	1:B:112:GLN:O	2.31	0.47
1:B:322:LEU:HD13	1:B:324:TRP:CZ2	2.50	0.47
1:A:264:ARG:HH11	1:A:264:ARG:HG3	1.79	0.47
1:A:322:LEU:HD13	1:A:324:TRP:CZ2	2.50	0.47
1:B:114:ARG:HD3	1:B:115:PRO:HD3	1.95	0.47
1:B:213:ILE:HG21	9:B:2962:HOH:O	2.14	0.47
1:A:69:LYS:O	1:A:69:LYS:CG	2.63	0.47
1:A:111:LEU:O	1:A:112:GLN:C	2.53	0.46
1:B:167:LEU:HG	1:B:348:LEU:HD12	1.98	0.46
1:B:178:GLN:HE22	1:B:181:ARG:HH11	1.63	0.46
1:A:114:ARG:HD2	1:A:479:PRO:HG3	1.97	0.46
1:B:113:THR:HG21	1:B:342:LEU:HD13	1.97	0.46
1:B:233:PHE:HB3	1:B:234:PRO:CD	2.46	0.45
1:A:146:GLN:H	1:A:146:GLN:CD	2.19	0.45
1:A:400:ALA:O	1:A:404:ILE:HG13	2.17	0.45
1:B:307:LEU:HD11	9:B:2971:HOH:O	2.16	0.45
1:A:340:ASN:HD22	1:A:340:ASN:N	2.14	0.45
1:B:259:GLN:HG3	1:B:260:ASP:H	1.81	0.45
1:B:419:ILE:HG13	1:B:420:VAL:N	2.32	0.45
1:B:270:VAL:O	1:B:274:GLU:HG3	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:113:THR:HG23	1:B:349:GLU:OE2	2.17	0.44
1:A:126:LEU:HB3	1:A:130:ARG:HH21	1.83	0.44
1:B:150:GLU:O	1:B:154:GLU:HG3	2.17	0.44
1:A:236:ARG:HG2	1:A:242:ASP:OD2	2.18	0.44
1:A:370:CYS:SG	1:A:378:LEU:HD13	2.58	0.43
1:A:77:GLU:HG3	1:B:372:PRO:CG	2.46	0.43
1:B:71:PRO:HG2	1:B:84:ASP:HB3	1.99	0.43
1:B:213:ILE:CG2	9:B:2962:HOH:O	2.66	0.43
1:B:108:PRO:HB3	1:B:111:LEU:HB2	2.01	0.43
1:B:264:ARG:HH11	1:B:287:ARG:HG3	1.83	0.43
1:B:180:TRP:CE3	1:B:192:TRP:HA	2.53	0.43
1:A:366:THR:O	1:A:370:CYS:HB2	2.19	0.43
8:A:1880:GOL:H32	9:B:2961:HOH:O	2.17	0.43
1:B:323:GLU:HG2	1:B:324:TRP:N	2.34	0.43
1:A:294:LEU:HD13	1:A:304:LEU:HD13	2.00	0.43
1:B:120:PRO:HA	1:B:121:PRO:HD3	1.93	0.43
1:B:319:HIS:CG	1:B:320:PRO:HD2	2.54	0.42
1:A:111:LEU:HD22	1:A:476:ARG:HH12	1.70	0.42
1:B:340:ASN:HD22	1:B:340:ASN:N	2.13	0.42
1:B:340:ASN:H	1:B:340:ASN:ND2	2.17	0.42
1:A:126:LEU:HD12	1:A:130:ARG:NE	2.30	0.42
1:B:105:LEU:HD23	1:B:105:LEU:N	2.34	0.42
1:B:109:ARG:O	1:B:110:LYS:HG3	2.20	0.42
1:A:109:ARG:HD3	1:A:109:ARG:HA	1.86	0.42
1:A:236:ARG:HH11	1:A:236:ARG:HG3	1.84	0.42
1:A:157:ALA:O	1:A:160:ALA:HB3	2.20	0.41
1:A:68:PRO:HD2	1:A:83:TYR:CZ	2.55	0.41
1:B:233:PHE:HB3	1:B:234:PRO:HD2	2.01	0.41
1:B:265:GLY:O	1:B:267:PRO:HD3	2.20	0.41
1:A:281:TRP:HB2	1:A:304:LEU:HD21	2.01	0.41
1:A:107:LEU:N	1:A:108:PRO:HD3	2.36	0.41
1:A:117:PRO:O	1:A:118:GLY:C	2.59	0.41
1:A:285:ASN:N	1:A:285:ASN:HD22	2.18	0.41
1:A:455:SER:O	1:A:458:LEU:HB2	2.21	0.41
1:A:151:ARG:O	1:A:155:VAL:HG23	2.21	0.40
1:A:234:PRO:HB2	1:A:243:PHE:CE1	2.57	0.40
1:A:236:ARG:HG3	1:A:236:ARG:NH1	2.37	0.40
1:A:449:TRP:HA	6:A:1650:AP4:N1	2.37	0.40
1:A:264:ARG:HG3	1:A:264:ARG:NH1	2.36	0.40
1:A:456:GLY:HA3	9:B:3076:HOH:O	2.21	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	414/444 (93%)	390 (94%)	22 (5%)	2 (0%)	29	32
1	B	412/444 (93%)	384 (93%)	26 (6%)	2 (0%)	29	32
All	All	826/888 (93%)	774 (94%)	48 (6%)	4 (0%)	29	32

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	117	PRO
1	B	112	GLN
1	B	259	GLN
1	A	112	GLN

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	354/377 (94%)	339 (96%)	15 (4%)	30	36
1	B	353/377 (94%)	342 (97%)	11 (3%)	40	48
All	All	707/754 (94%)	681 (96%)	26 (4%)	34	42

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	100	ARG
1	A	102	LEU

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Mol	Chain	Res	Type
1	A	124	GLN
1	A	125	LEU
1	A	126	LEU
1	A	260	ASP
1	A	285	ASN
1	A	293	LEU
1	A	328	LEU
1	A	340	ASN
1	A	376	ASN
1	A	378	LEU
1	A	390	ARG
1	A	416	LYS
1	A	468	ASN
1	B	99	ARG
1	B	100	ARG
1	B	102	LEU
1	B	114	ARG
1	B	153	GLN
1	B	207	GLN
1	B	225	ASN
1	B	330	LEU
1	B	340	ASN
1	B	378	LEU
1	B	458	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	91	GLN
1	A	112	GLN
1	A	146	GLN
1	A	191	GLN
1	A	278	GLN
1	A	285	ASN
1	A	340	ASN
1	A	376	ASN
1	A	413	GLN
1	A	468	ASN
1	B	112	GLN
1	B	178	GLN
1	B	191	GLN
1	B	222	ASN

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Mol	Chain	Res	Type
1	B	225	ASN
1	B	340	ASN
1	B	376	ASN
1	B	405	ASN

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 15 ligands modelled in this entry, 1 is monoatomic - leaving 14 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	$\#\|Z\| > 2$
8	GOL	A	1880	-	5,5,5	0.15	0	5,5,5	0.17	0
6	AP4	B	2650	-	17,18,18	4.36	7 (41%)	14,27,27	3.23	10 (71%)
7	ITU	B	2800	-	4,5,5	0.79	0	4,5,5	2.62	1 (25%)
2	ACT	A	1850	-	3,3,3	1.00	0	3,3,3	0.72	0
3	CAC	A	950	1	0,2,4	-	-	0,1,6	-	-
6	AP4	A	1650	-	17,18,18	4.37	7 (41%)	14,27,27	3.26	10 (71%)
5	HEM	A	1500	1	41,50,50	1.23	2 (4%)	45,82,82	1.21	3 (6%)
8	GOL	B	2880	-	5,5,5	0.20	0	5,5,5	0.18	0
7	ITU	A	1800	-	4,5,5	0.79	0	4,5,5	2.85	1 (25%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	ACT	B	2860	-	3,3,3	0.90	0	3,3,3	0.73	0
2	ACT	A	1860	-	3,3,3	0.85	0	3,3,3	0.83	0
5	HEM	B	2500	1	41,50,50	1.29	4 (9%)	45,82,82	1.25	5 (11%)
3	CAC	B	950	1	0,2,4	-	-	0,1,6	-	-
2	ACT	B	2850	-	3,3,3	0.83	0	3,3,3	0.78	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
7	ITU	B	2800	-	-	0/3/3/3	-
6	AP4	A	1650	-	-	-	0/2/3/3
5	HEM	A	1500	1	-	2/12/54/54	-
8	GOL	B	2880	-	-	0/4/4/4	-
8	GOL	A	1880	-	-	0/4/4/4	-
6	AP4	B	2650	-	-	-	0/2/3/3
5	HEM	B	2500	1	-	2/12/54/54	-
7	ITU	A	1800	-	-	0/3/3/3	-

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	2650	AP4	C13-N5	11.03	1.52	1.36
6	A	1650	AP4	C13-N5	10.90	1.52	1.36
6	A	1650	AP4	C7-C6	-9.44	1.41	1.52
6	B	2650	AP4	C7-C6	-9.32	1.41	1.52
6	B	2650	AP4	O12-C13	6.51	1.43	1.35
6	A	1650	AP4	O12-C13	6.42	1.43	1.35
6	A	1650	AP4	C4-N3	5.47	1.42	1.33
6	B	2650	AP4	C4-N3	5.37	1.42	1.33
6	A	1650	AP4	C6-N5	4.00	1.53	1.48
6	B	2650	AP4	C6-N5	3.44	1.52	1.48
6	A	1650	AP4	C10-N1	3.40	1.41	1.34
6	B	2650	AP4	C10-N1	3.38	1.40	1.34
5	A	1500	HEM	C3C-CAC	-3.34	1.41	1.47
5	B	2500	HEM	C3C-C2C	-2.97	1.36	1.40
5	B	2500	HEM	C3C-CAC	-2.80	1.42	1.47
5	A	1500	HEM	CAB-C3B	-2.56	1.40	1.47
6	B	2650	AP4	O12-C11	-2.45	1.41	1.45
6	A	1650	AP4	O12-C11	-2.31	1.41	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	2500	HEM	C4A-NA	2.08	1.40	1.36
5	B	2500	HEM	CAB-C3B	-2.01	1.42	1.47

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
6	B	2650	AP4	C6-N5-C13	-6.32	105.50	111.12
6	A	1650	AP4	C6-N5-C13	-6.09	105.71	111.12
7	A	1800	ITU	C2-S-C3	5.68	109.15	103.20
7	B	2800	ITU	C2-S-C3	5.21	108.66	103.20
6	B	2650	AP4	C10-C9-C4	4.85	118.23	114.44
6	A	1650	AP4	C10-C9-C4	4.78	118.18	114.44
6	A	1650	AP4	C2-N3-C4	3.94	122.19	115.93
6	B	2650	AP4	O12-C11-C6	3.70	109.31	105.38
6	A	1650	AP4	O12-C11-C6	3.68	109.29	105.38
6	B	2650	AP4	C2-N3-C4	3.63	121.70	115.93
6	B	2650	AP4	O14-C13-N5	3.57	132.73	128.68
6	A	1650	AP4	O14-C13-N5	3.53	132.68	128.68
6	A	1650	AP4	N1-C2-N3	-3.40	120.08	125.42
6	A	1650	AP4	N2-C2-N1	3.31	122.40	117.25
6	B	2650	AP4	N1-C2-N3	-3.25	120.31	125.42
6	B	2650	AP4	N2-C2-N1	3.12	122.11	117.25
5	A	1500	HEM	CMD-C2D-C1D	2.85	129.38	125.04
5	B	2500	HEM	C4B-CHC-C1C	2.85	126.31	122.56
5	B	2500	HEM	C4B-C3B-C2B	-2.80	104.89	107.11
6	A	1650	AP4	C2-N1-C10	2.75	120.71	114.54
6	B	2650	AP4	C2-N1-C10	2.71	120.61	114.54
6	A	1650	AP4	O12-C13-O14	-2.70	119.29	122.53
6	B	2650	AP4	O12-C13-O14	-2.58	119.44	122.53
6	A	1650	AP4	C9-C4-N3	-2.50	118.00	123.14
6	B	2650	AP4	C9-C4-N3	-2.39	118.22	123.14
5	A	1500	HEM	C4B-CHC-C1C	2.34	125.65	122.56
5	B	2500	HEM	C4A-C3A-C2A	-2.27	105.42	107.00
5	B	2500	HEM	C2C-C3C-C4C	-2.19	105.37	106.90
5	A	1500	HEM	C4D-ND-C1D	-2.19	102.81	105.07
5	B	2500	HEM	CBA-CAA-C2A	-2.03	109.16	112.62

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	1500	HEM	CAD-CBD-CGD-O2D

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Mol	Chain	Res	Type	Atoms
5	A	1500	HEM	CAD-CBD-CGD-O1D
5	B	2500	HEM	CAA-CBA-CGA-O2A
5	B	2500	HEM	CAA-CBA-CGA-O1A

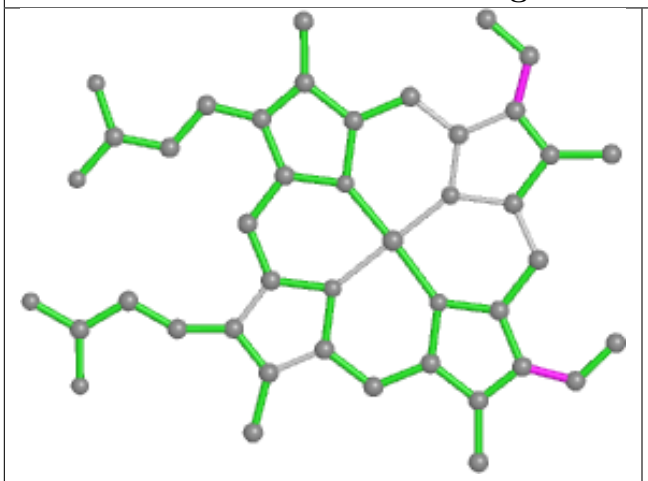
There are no ring outliers.

7 monomers are involved in 6 short contacts:

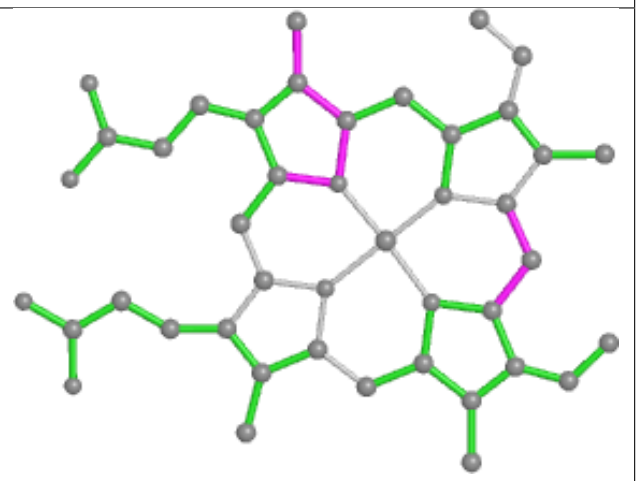
Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	A	1880	GOL	1	0
7	B	2800	ITU	1	0
3	A	950	CAC	1	0
6	A	1650	AP4	1	0
5	A	1500	HEM	1	0
8	B	2880	GOL	1	0
2	A	1860	ACT	1	0

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.

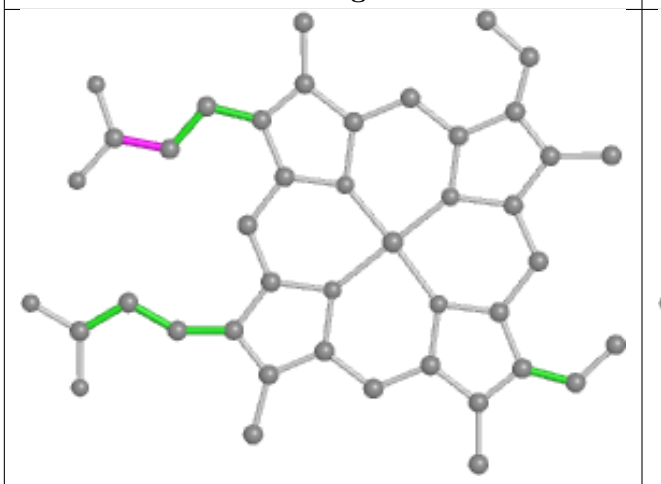
Ligand HEM A 1500



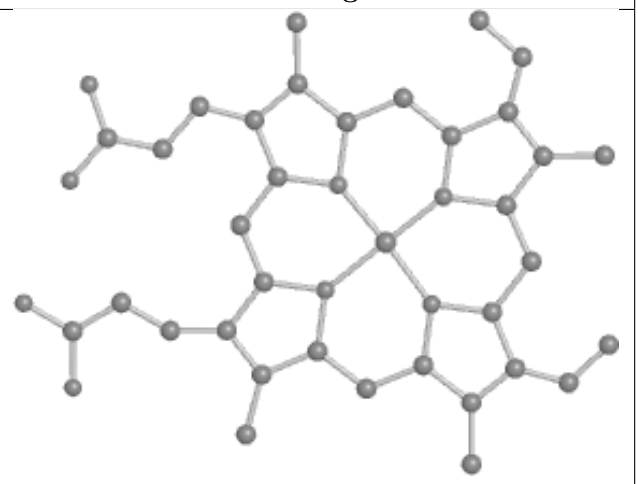
Bond lengths



Bond angles

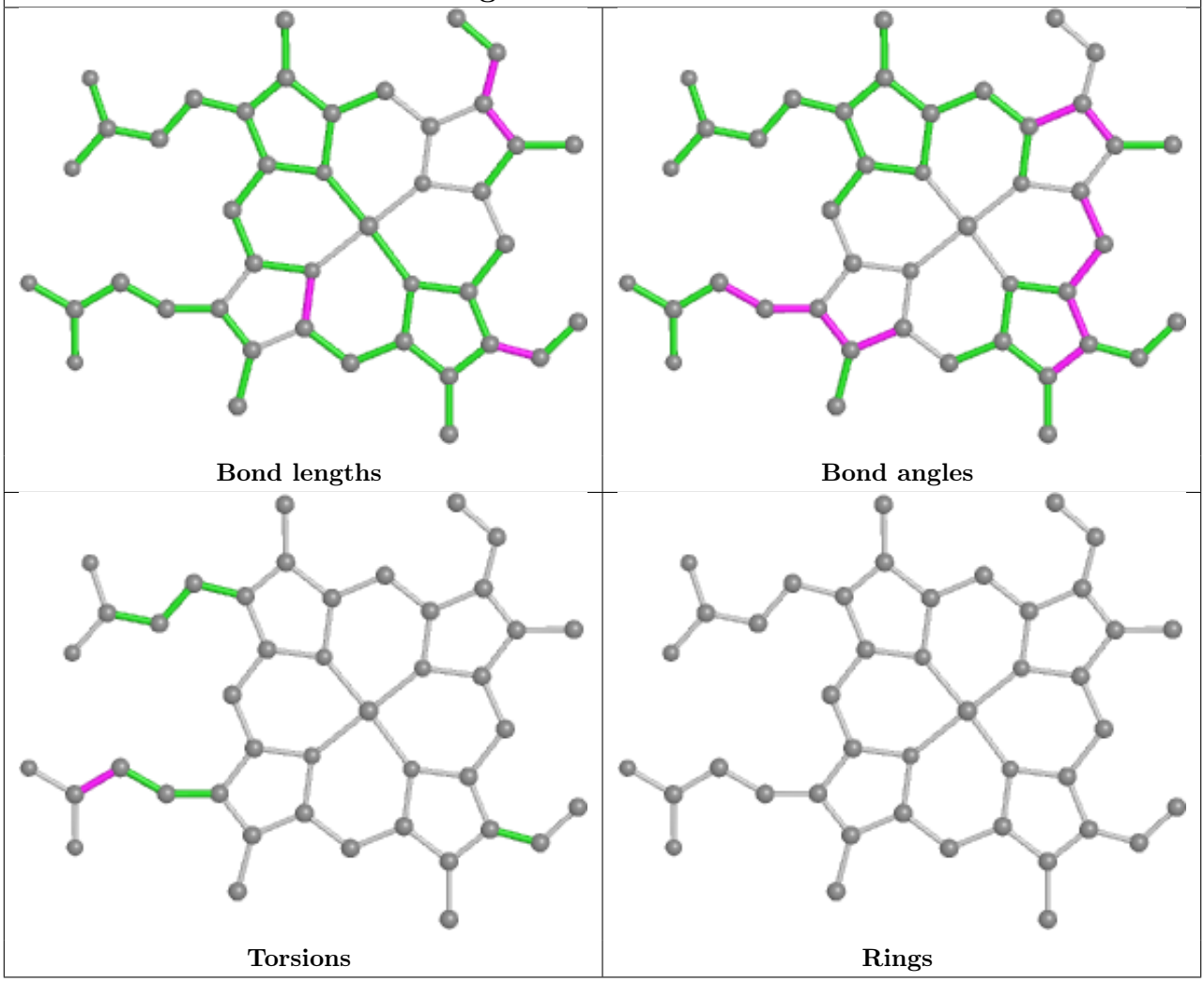


Torsions



Rings

Ligand HEM B 2500



5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	416/444 (93%)	0.27	27 (6%)	18	27	17, 28, 59, 100	0
1	B	414/444 (93%)	0.18	18 (4%)	35	47	16, 30, 63, 100	0
All	All	830/888 (93%)	0.23	45 (5%)	25	37	16, 29, 62, 100	0

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	113	THR	14.4
1	A	119	PRO	13.4
1	A	111	LEU	12.5
1	B	117	PRO	12.3
1	B	119	PRO	11.8
1	A	113	THR	11.7
1	B	118	GLY	11.3
1	A	115	PRO	10.5
1	B	116	SER	9.1
1	B	115	PRO	8.7
1	A	110	LYS	8.3
1	B	114	ARG	8.3
1	B	111	LEU	7.9
1	A	117	PRO	7.7
1	A	118	GLY	7.0
1	A	114	ARG	5.8
1	A	112	GLN	5.7
1	B	110	LYS	5.7
1	B	120	PRO	5.2
1	B	259	GLN	4.9
1	A	116	SER	4.7
1	A	239	GLY	4.3
1	B	112	GLN	4.3
1	A	120	PRO	4.2

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Mol	Chain	Res	Type	RSRZ
1	A	259	GLN	3.8
1	A	109	ARG	3.8
1	B	239	GLY	3.4
1	A	108	PRO	3.2
1	B	261	GLY	3.2
1	A	121	PRO	3.1
1	A	69	LYS	2.9
1	A	153	GLN	2.9
1	A	257	ARG	2.8
1	A	146	GLN	2.7
1	A	123	GLU	2.7
1	B	257	ARG	2.7
1	B	121	PRO	2.6
1	A	122	ALA	2.6
1	A	390	ARG	2.4
1	A	161	SER	2.3
1	B	122	ALA	2.3
1	A	258	GLN	2.2
1	A	126	LEU	2.2
1	B	260	ASP	2.1
1	A	107	LEU	2.0

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

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

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 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
8	GOL	A	1880	6/6	0.85	0.28	51,52,53,55	0
8	GOL	B	2880	6/6	0.93	0.22	33,34,38,44	0

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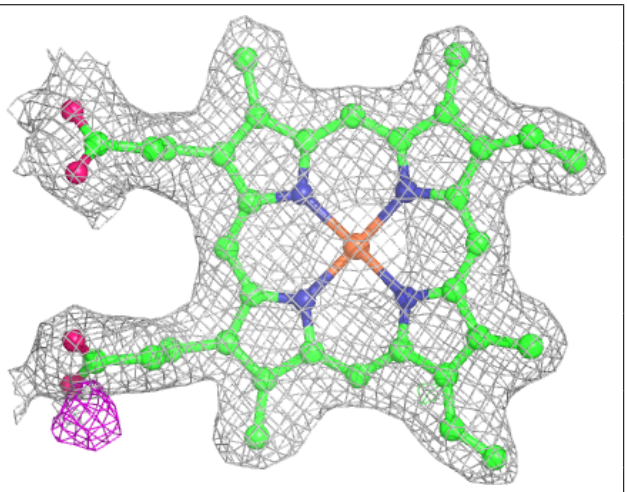
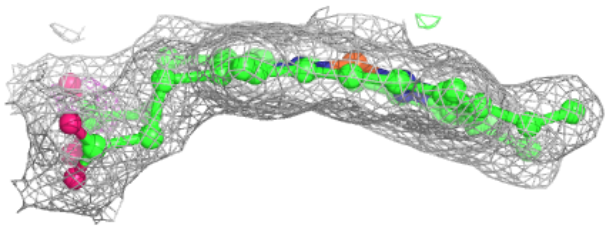
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	CAC	A	950	3/5	0.94	0.16	41,41,47,52	0
2	ACT	B	2860	4/4	0.95	0.10	25,27,28,31	0
6	AP4	A	1650	16/16	0.96	0.17	16,21,23,23	0
2	ACT	A	1850	4/4	0.97	0.09	24,26,29,29	0
6	AP4	B	2650	16/16	0.97	0.17	20,22,26,27	0
7	ITU	A	1800	6/6	0.97	0.19	21,23,24,25	0
2	ACT	B	2850	4/4	0.97	0.10	23,24,25,26	0
5	HEM	B	2500	43/43	0.97	0.13	17,22,28,37	0
2	ACT	A	1860	4/4	0.98	0.14	32,34,34,35	0
3	CAC	B	950	3/5	0.98	0.12	46,46,48,52	0
5	HEM	A	1500	43/43	0.98	0.17	15,19,31,36	0
4	ZN	A	900	1/1	0.99	0.11	23,23,23,23	0
7	ITU	B	2800	6/6	0.99	0.12	16,21,23,23	0

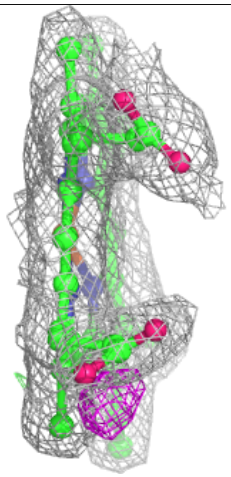
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around HEM B 2500:

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

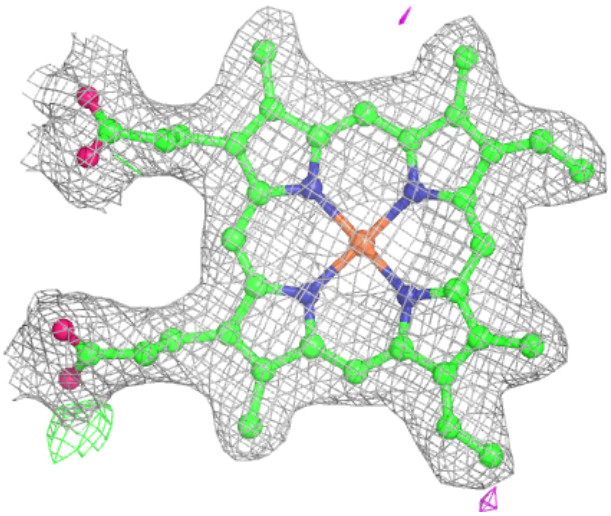


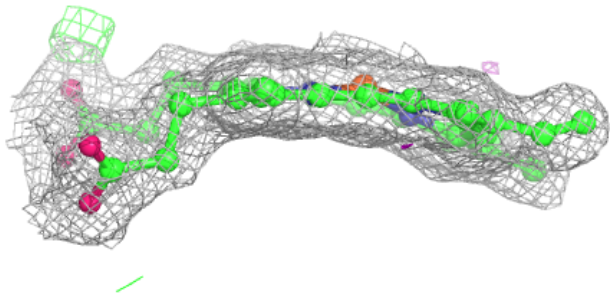


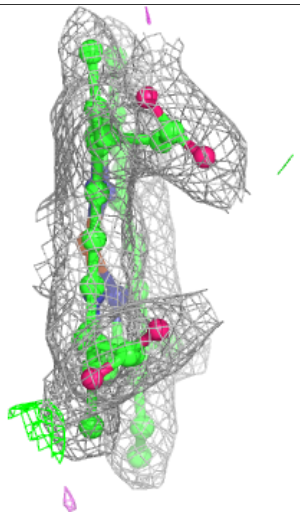


Electron density around HEM A 1500:

$2mF_o$-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.