



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 3, 2026 – 11:05 PM EDT

PDB ID : 1DMI / pdb_00001dmi
Title : BOVINE ENDOTHELIAL NITRIC OXIDE SYNTHASE HEME DOMAIN
COMPLEXED WITH 6S-H4B
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Deposited on : 1999-12-14
Resolution : 2.00 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (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

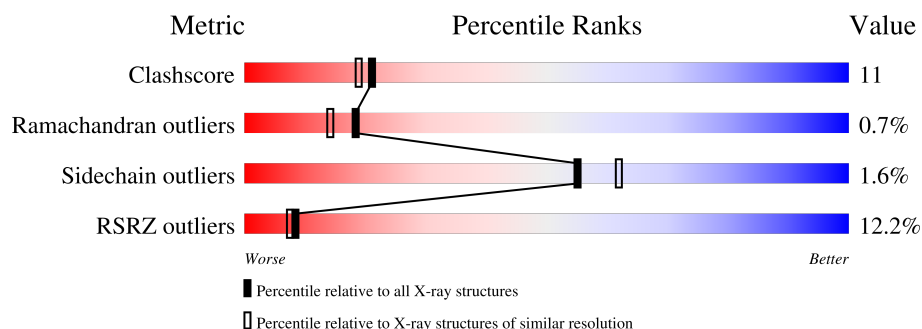
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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

Mol	Chain	Length	Quality of chain
1	A	444	 10% 72% 19% • 6%
1	B	444	 13% 73% 19% • 7%

2 Entry composition [i](#)

There are 9 unique types of molecules in this entry. The entry contains 7148 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.

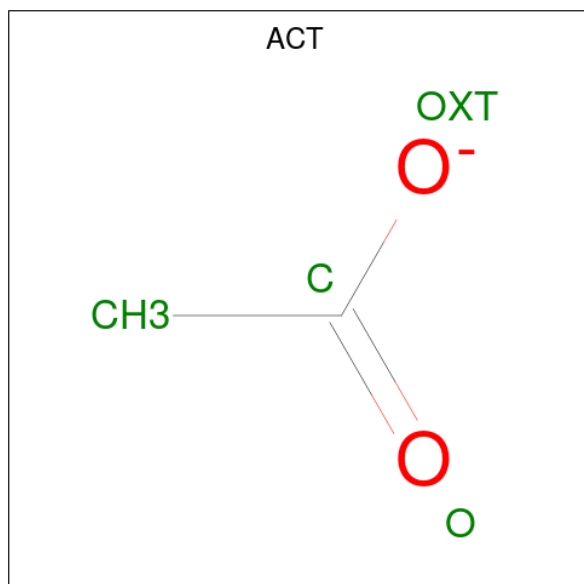
- Molecule 1 is a protein called NITRIC OXIDE SYNTHASE.

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

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 (CCD ID: ACT) (formula: C₂H₃O₂).



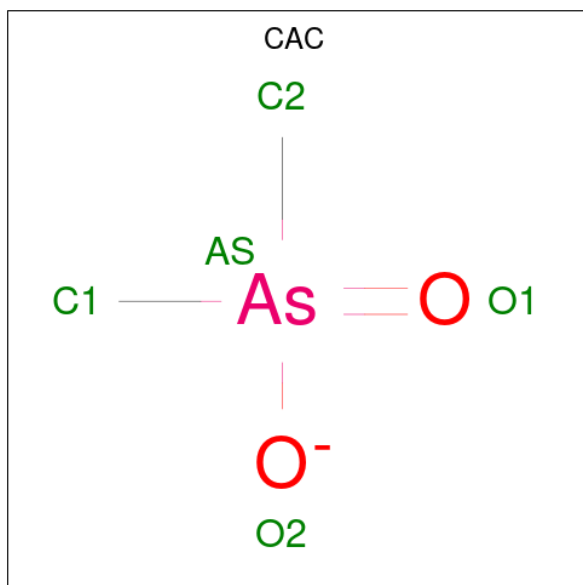
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			4	2	2		

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

- Molecule 3 is CACODYLATE ION (CCD ID: CAC) (formula: $C_2H_6AsO_2$).

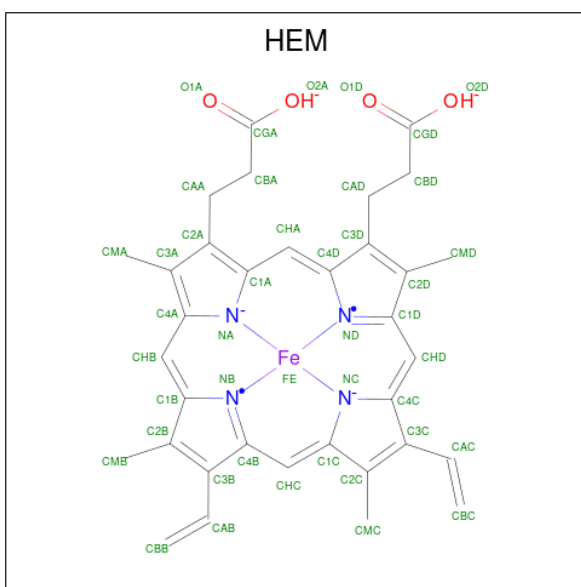


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	As	C	0	0
			3	1	2		
3	B	1	Total	As	C	0	0
			3	1	2		

- Molecule 4 is ZINC ION (CCD ID: ZN) (formula: Zn).

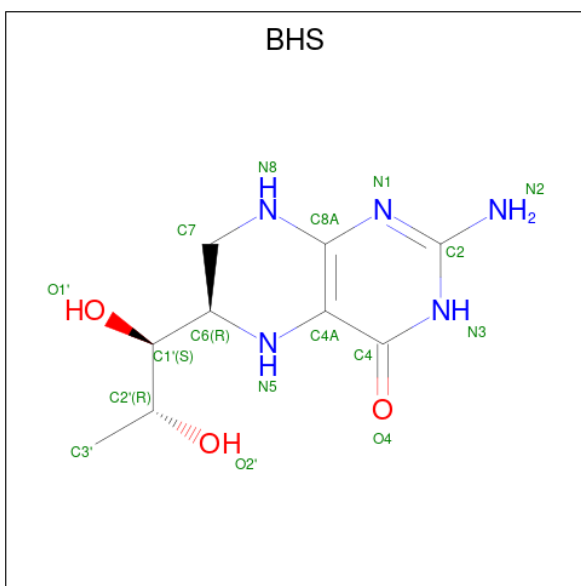
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Zn	0	0
			1	1		

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



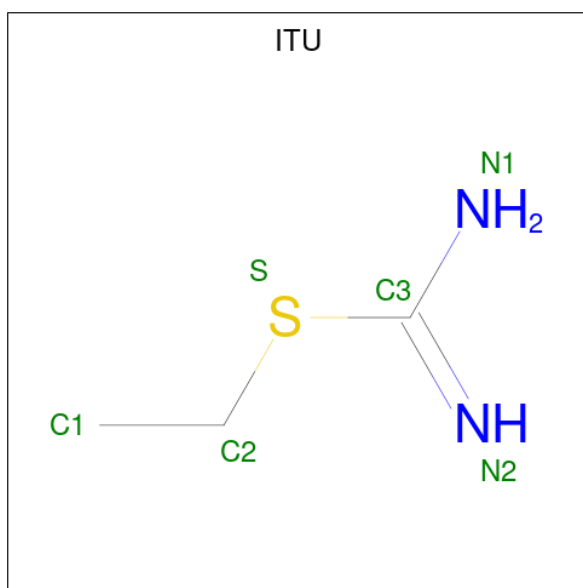
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
5	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 6 is 6S-5,6,7,8-TETRAHYDROBIOPTERIN (CCD ID: BHS) (formula: $C_9H_{15}N_5O_3$).



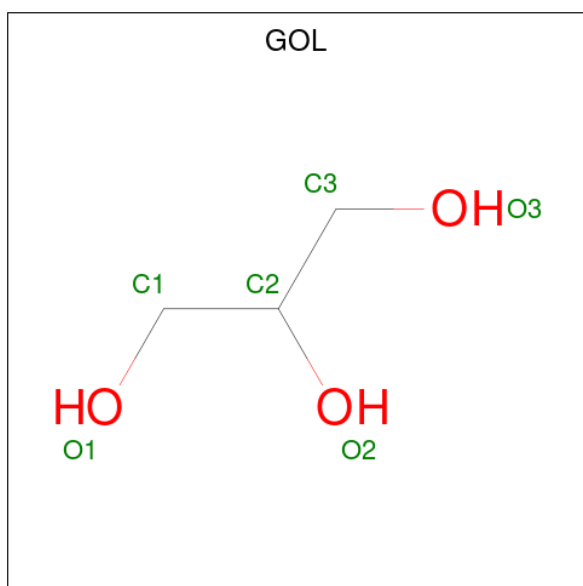
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total 17	C 9	N 5	O 3	0	0
6	B	1	Total 17	C 9	N 5	O 3	0	0

- Molecule 7 is ETHYLISOTHIOUREA (CCD ID: ITU) (formula: $C_3H_8N_2S$).



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

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



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

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

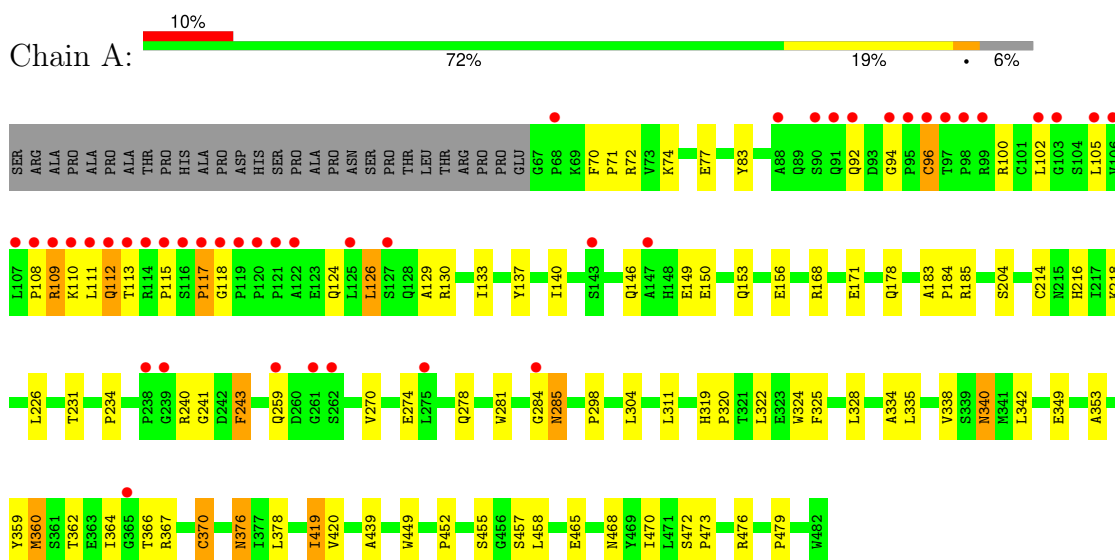
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	199	Total	O	0	0
			199	199		
9	B	189	Total	O	0	0
			189	189		

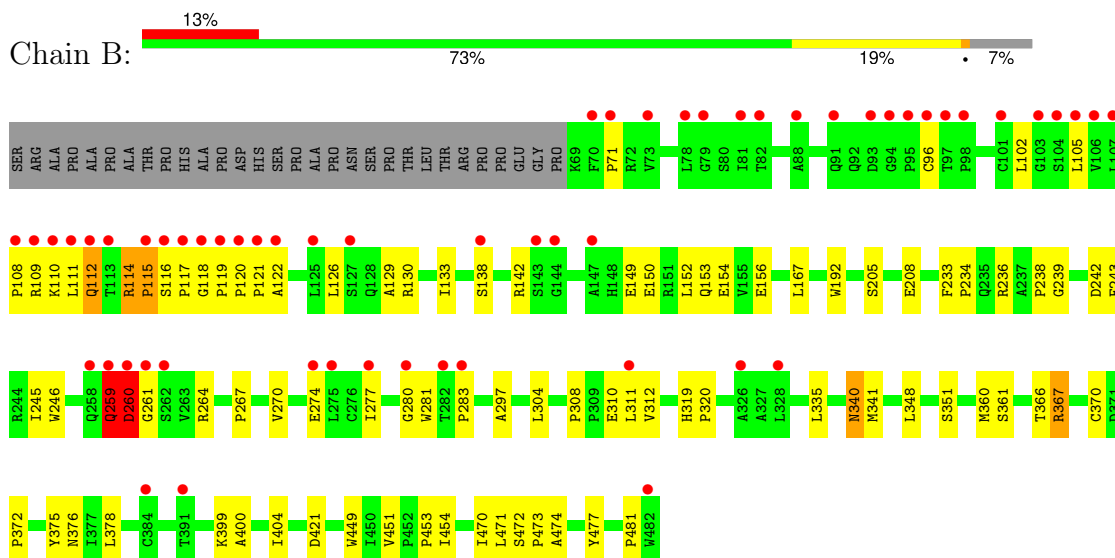
3 Residue-property plots

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

• Molecule 1: NITRIC OXIDE SYNTHASE



• Molecule 1: NITRIC OXIDE SYNTHASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	58.63Å 106.56Å 156.79Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	32.59 – 2.00 32.59 – 2.00	Depositor EDS
% Data completeness (in resolution range)	87.2 (32.59-2.00) 80.1 (32.59-2.00)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.51 (at 2.00Å)	Xtriage
Refinement program	CNS 0.9	Depositor
R, R_{free}	0.229 , 0.264 0.225 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	35.5	Xtriage
Anisotropy	0.477	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 48.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7148	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.47% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: HEM, CAC, ITU, BHS, ACT, ZN, 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	0.46	0/3397	0.95	15/4631 (0.3%)
1	B	0.44	1/3385 (0.0%)	0.96	14/4614 (0.3%)
All	All	0.45	1/6782 (0.0%)	0.95	29/9245 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	122	ALA	CA-CB	-6.79	1.42	1.53

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	360	MET	N-CA-C	-7.86	96.42	109.07
1	B	360	MET	N-CA-C	-7.46	97.08	109.24
1	A	243	PHE	N-CA-C	-6.86	98.07	109.46
1	B	243	PHE	N-CA-C	-6.86	98.86	109.76
1	B	451	VAL	CA-C-N	6.79	124.62	119.66
1	B	451	VAL	C-N-CA	6.79	124.62	119.66
1	B	245	ILE	N-CA-C	-6.63	98.19	107.80
1	B	335	LEU	N-CA-C	6.52	120.43	108.69
1	A	109	ARG	N-CA-C	6.34	119.48	109.39
1	B	470	ILE	N-CA-C	6.11	116.91	107.99
1	A	183	ALA	CA-C-N	6.03	125.91	119.28
1	A	183	ALA	C-N-CA	6.03	125.91	119.28
1	A	94	GLY	CA-C-N	5.62	125.15	119.19
1	A	94	GLY	C-N-CA	5.62	125.15	119.19
1	B	367	ARG	N-CA-C	5.59	117.23	111.03
1	B	138	SER	N-CA-C	-5.58	104.83	111.69
1	B	280	GLY	N-CA-C	5.55	122.13	114.92
1	B	192	TRP	N-CA-C	5.54	118.85	111.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	335	LEU	N-CA-C	5.40	118.41	108.69
1	B	261	GLY	N-CA-C	-5.39	108.28	114.69
1	A	439	ALA	N-CA-C	5.37	117.21	111.36
1	A	96	CYS	N-CA-C	5.25	117.45	110.53
1	A	226	LEU	N-CA-C	5.22	117.60	110.24
1	A	284	GLY	N-CA-C	-5.16	105.22	112.60
1	B	312	VAL	N-CA-C	5.16	114.03	106.55
1	B	297	ALA	N-CA-C	-5.12	103.50	110.36
1	A	370	CYS	N-CA-C	5.12	120.37	113.72
1	A	452	PRO	N-CA-C	5.04	115.31	110.47
1	A	334	ALA	N-CA-C	5.02	117.65	111.82

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	3302	0	3215	77	0
1	B	3291	0	3205	77	0
2	A	8	0	6	0	0
2	B	8	0	6	0	0
3	A	3	0	0	0	0
3	B	3	0	0	0	0
4	A	1	0	0	0	0
5	A	43	0	30	0	0
5	B	43	0	30	2	0
6	A	17	0	15	1	0
6	B	17	0	15	2	0
7	A	6	0	7	2	0
7	B	6	0	7	0	0
8	A	6	0	8	0	0
8	B	6	0	8	0	0
9	A	199	0	0	4	0
9	B	189	0	0	3	0
All	All	7148	0	6552	147	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 (147) 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	1.28	1.15
1:A:126:LEU:HD11	1:A:156:GLU:HG2	1.37	1.02
1:B:126:LEU:HD11	1:B:156:GLU:HG2	1.53	0.89
1:B:108:PRO:HB3	1:B:111:LEU:HD13	1.57	0.84
1:B:114:ARG:HA	1:B:114:ARG:HH11	1.41	0.84
1:B:116:SER:HB2	1:B:236:ARG:NH2	1.98	0.78
1:B:270:VAL:O	1:B:274:GLU:HG2	1.82	0.77
1:B:281:TRP:HB2	1:B:304:LEU:HD21	1.66	0.77
1:A:92:GLN:NE2	1:A:111:LEU:HD23	1.99	0.76
1:A:204:SER:OG	9:A:2062:HOH:O	2.05	0.73
1:B:114:ARG:HA	1:B:114:ARG:NH1	2.07	0.70
1:A:338:VAL:HG23	7:A:1800:ITU:H11	1.73	0.69
1:A:77:GLU:HG3	1:B:372:PRO:HG2	1.77	0.66
1:A:111:LEU:HD22	1:A:476:ARG:NH1	2.11	0.66
1:B:378:LEU:HB2	9:B:3033:HOH:O	1.96	0.66
1:A:178:GLN:HG2	9:A:2058:HOH:O	1.95	0.65
1:B:236:ARG:HD3	1:B:351:SER:HB3	1.79	0.65
1:A:146:GLN:O	1:A:150:GLU:HG3	1.97	0.64
1:A:340:ASN:H	1:A:340:ASN:HD22	1.46	0.64
1:B:114:ARG:HH11	1:B:115:PRO:HD3	1.64	0.63
1:B:259:GLN:HG2	1:B:260:ASP:H	1.64	0.62
1:B:149:GLU:O	1:B:153:GLN:HG3	1.99	0.62
1:A:378:LEU:HB2	9:A:2070:HOH:O	1.98	0.62
1:A:240:ARG:HD3	1:A:298:PRO:HB3	1.80	0.62
1:A:96:CYS:HB3	1:B:96:CYS:HB3	1.82	0.61
1:B:126:LEU:O	1:B:130:ARG:HG3	2.00	0.60
1:A:111:LEU:HD21	1:A:470:ILE:HG21	1.82	0.60
1:A:168:ARG:HB2	1:A:171:GLU:HG3	1.83	0.59
1:A:115:PRO:C	1:A:117:PRO:HD3	2.27	0.59
1:A:270:VAL:O	1:A:274:GLU:HG3	2.03	0.59
1:B:236:ARG:HD2	1:B:242:ASP:CG	2.27	0.59
1:B:370:CYS:SG	1:B:378:LEU:HD13	2.41	0.59
1:A:115:PRO:O	1:A:117:PRO:HD3	2.04	0.58
1:A:214:CYS:O	1:A:218:LYS:HG3	2.04	0.57
1:A:322:LEU:HD13	1:A:324:TRP:CZ2	2.39	0.57
1:B:310:GLU:HG2	1:B:311:LEU:HD12	1.87	0.56
1:B:119:PRO:HA	1:B:238:PRO:CG	2.36	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:218:LYS:HG2	1:A:311:LEU:HD22	1.88	0.55
1:A:113:THR:HG21	1:A:342:LEU:HD22	1.88	0.55
1:A:149:GLU:O	1:A:153:GLN:HG3	2.07	0.55
1:A:105:LEU:CB	1:A:108:PRO:HG3	2.37	0.55
1:A:285:ASN:HD22	1:A:285:ASN:C	2.13	0.55
1:B:150:GLU:O	1:B:154:GLU:HG3	2.07	0.55
1:B:449:TRP:HA	6:B:2610:BHS:N1	2.22	0.55
1:A:111:LEU:HD11	1:A:470:ILE:CD1	2.38	0.54
1:B:102:LEU:HB3	1:B:105:LEU:HD22	1.90	0.54
1:A:111:LEU:HD11	1:A:470:ILE:HD12	1.90	0.54
1:A:240:ARG:HD3	1:A:298:PRO:CB	2.37	0.54
1:B:108:PRO:CB	1:B:111:LEU:HB2	2.38	0.54
1:B:472:SER:HA	1:B:473:PRO:C	2.33	0.53
1:A:360:MET:HE3	1:A:362:THR:OG1	2.08	0.53
1:A:455:SER:HB3	1:A:458:LEU:HD12	1.91	0.53
1:A:342:LEU:HD11	1:A:349:GLU:HB3	1.90	0.52
1:A:105:LEU:HB2	1:A:108:PRO:HG3	1.90	0.52
1:B:400:ALA:O	1:B:404:ILE:HG13	2.10	0.52
1:A:74:LYS:O	1:A:465:GLU:HG3	2.09	0.52
1:A:111:LEU:HD21	1:A:470:ILE:HG13	1.92	0.52
1:B:111:LEU:O	1:B:112:GLN:O	2.26	0.52
1:B:152:LEU:O	1:B:156:GLU:HG3	2.09	0.52
1:B:167:LEU:HG	1:B:348:LEU:HD12	1.93	0.51
1:B:119:PRO:N	1:B:238:PRO:HB3	2.26	0.51
1:B:114:ARG:NH1	1:B:115:PRO:HD3	2.27	0.50
1:A:109:ARG:O	1:A:110:LYS:HD2	2.12	0.50
1:A:472:SER:HA	1:A:473:PRO:C	2.36	0.50
1:B:119:PRO:HA	1:B:238:PRO:HG3	1.94	0.50
1:A:274:GLU:O	1:A:278:GLN:HG3	2.11	0.50
1:B:236:ARG:HD2	1:B:242:ASP:OD1	2.12	0.50
1:B:319:HIS:CG	1:B:320:PRO:HD2	2.47	0.49
1:B:130:ARG:HB3	1:B:130:ARG:NH1	2.27	0.49
1:A:340:ASN:H	1:A:340:ASN:ND2	2.10	0.49
1:A:457:SER:OG	1:B:453:PRO:HB2	2.12	0.49
1:B:471:LEU:O	1:B:474:ALA:HB2	2.12	0.49
1:A:340:ASN:HD22	1:A:340:ASN:N	2.07	0.49
1:A:111:LEU:O	1:A:112:GLN:C	2.56	0.49
1:A:419:ILE:HG13	1:A:420:VAL:N	2.28	0.49
1:B:116:SER:HB2	1:B:236:ARG:HH22	1.76	0.49
1:A:231:THR:O	1:A:353:ALA:HA	2.13	0.48
1:A:285:ASN:C	1:A:285:ASN:ND2	2.72	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:100:ARG:HD3	1:A:100:ARG:H	1.79	0.48
1:A:115:PRO:HD3	1:A:479:PRO:CG	2.20	0.47
1:A:102:LEU:HD21	1:B:71:PRO:HB3	1.96	0.47
1:A:338:VAL:HG23	7:A:1800:ITU:C1	2.43	0.47
1:A:126:LEU:O	1:A:130:ARG:HG3	2.15	0.47
1:B:117:PRO:O	1:B:239:GLY:N	2.48	0.47
1:B:118:GLY:C	1:B:238:PRO:HB3	2.40	0.47
1:B:112:GLN:NE2	1:B:477:TYR:C	2.73	0.47
1:A:240:ARG:HD2	1:A:241:GLY:O	2.14	0.47
1:B:308:PRO:HB2	1:B:311:LEU:HD13	1.97	0.46
1:B:116:SER:HB2	1:B:236:ARG:CZ	2.45	0.46
1:B:281:TRP:CB	1:B:304:LEU:HD21	2.42	0.46
1:B:120:PRO:HA	1:B:121:PRO:HD3	1.88	0.46
1:B:399:LYS:NZ	9:B:3039:HOH:O	2.48	0.46
1:A:234:PRO:HB2	1:A:243:PHE:CE1	2.51	0.46
1:B:129:ALA:O	1:B:133:ILE:HG12	2.16	0.46
1:B:205:SER:OG	1:B:208:GLU:HG3	2.16	0.46
1:B:236:ARG:HD2	1:B:242:ASP:OD2	2.16	0.45
1:A:92:GLN:HE22	1:A:111:LEU:HD23	1.77	0.45
1:A:72:ARG:HB2	1:B:109:ARG:NH2	2.32	0.45
1:A:83:TYR:HE2	1:B:109:ARG:HH22	1.65	0.45
1:B:277:ILE:HD13	1:B:283:PRO:HG3	1.99	0.45
1:A:117:PRO:O	1:A:118:GLY:C	2.58	0.45
1:A:325:PHE:O	1:A:328:LEU:HB2	2.17	0.45
1:A:77:GLU:HG3	1:B:372:PRO:CG	2.46	0.45
1:A:319:HIS:CG	1:A:320:PRO:HD2	2.51	0.45
1:B:340:ASN:HD22	1:B:340:ASN:C	2.25	0.45
1:B:366:THR:HG21	1:B:454:ILE:HG23	1.99	0.45
1:A:281:TRP:HB2	1:A:304:LEU:HD21	1.99	0.44
1:A:111:LEU:HD22	1:A:476:ARG:HH12	1.83	0.44
1:B:117:PRO:C	1:B:239:GLY:H	2.25	0.44
1:A:216:HIS:CD2	1:A:216:HIS:C	2.95	0.44
1:B:108:PRO:HB3	1:B:111:LEU:HB2	2.00	0.43
1:B:340:ASN:HD22	1:B:341:MET:N	2.16	0.43
1:B:116:SER:N	1:B:117:PRO:CD	2.81	0.43
1:B:366:THR:O	1:B:370:CYS:HB2	2.18	0.43
1:B:246:TRP:CD1	1:B:481:PRO:HG3	2.54	0.43
1:A:113:THR:HG21	1:A:476:ARG:HD2	2.01	0.43
1:A:70:PHE:O	1:B:109:ARG:NH2	2.52	0.42
1:B:110:LYS:C	1:B:112:GLN:N	2.76	0.42
1:A:109:ARG:HD3	1:A:109:ARG:HA	1.86	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:109:ARG:HG3	1:B:110:LYS:HG3	2.01	0.42
1:B:267:PRO:O	1:B:270:VAL:HG23	2.19	0.42
1:A:359:TYR:CD2	1:A:364:ILE:HD11	2.54	0.42
1:B:117:PRO:O	1:B:239:GLY:CA	2.67	0.42
1:B:233:PHE:HB3	1:B:234:PRO:CD	2.50	0.42
1:A:71:PRO:HB3	1:B:102:LEU:HD21	2.02	0.42
1:A:185:ARG:HD3	1:A:449:TRP:CD2	2.54	0.42
1:B:310:GLU:HG2	1:B:311:LEU:CD1	2.48	0.42
1:A:129:ALA:O	1:A:133:ILE:HG12	2.21	0.41
1:A:184:PRO:HB3	1:A:468:ASN:ND2	2.36	0.41
1:A:366:THR:O	1:A:370:CYS:HB2	2.20	0.41
5:B:2500:HEM:CMC	5:B:2500:HEM:HBC2	2.50	0.41
1:A:113:THR:CG2	1:A:476:ARG:CD	2.99	0.41
1:B:375:TYR:O	1:B:376:ASN:C	2.63	0.41
1:B:264:ARG:HG3	1:B:264:ARG:HH11	1.85	0.41
1:A:111:LEU:CD2	1:A:476:ARG:NH1	2.83	0.41
1:B:114:ARG:HD3	1:B:115:PRO:HD3	2.02	0.41
5:B:2500:HEM:HBC2	5:B:2500:HEM:HMC1	2.03	0.40
1:A:137:TYR:HA	1:A:140:ILE:HG12	2.02	0.40
1:B:233:PHE:HB3	1:B:234:PRO:HD2	2.02	0.40
1:B:259:GLN:HG2	1:B:260:ASP:N	2.34	0.40
1:B:361:SER:OG	1:B:421:ASP:HA	2.21	0.40
1:A:110:LYS:HE3	1:A:110:LYS:HA	2.04	0.40
1:A:376:ASN:ND2	9:A:2070:HOH:O	2.41	0.40
1:A:367:ARG:HH12	6:A:1610:BHS:C4	2.34	0.40
1:B:142:ARG:HG2	1:B:142:ARG:NH1	2.35	0.40
1:B:367:ARG:HH12	6:B:2610:BHS:C4	2.34	0.40
1:A:96:CYS:HB2	9:B:2987:HOH:O	2.22	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%)	392 (95%)	20 (5%)	2 (0%)	24	21
1	B	412/444 (93%)	387 (94%)	21 (5%)	4 (1%)	12	8
All	All	826/888 (93%)	779 (94%)	41 (5%)	6 (1%)	18	14

All (6) 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
1	B	115	PRO
1	B	260	ASP

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	354/377 (94%)	347 (98%)	7 (2%)	48	54
1	B	353/377 (94%)	349 (99%)	4 (1%)	65	73
All	All	707/754 (94%)	696 (98%)	11 (2%)	55	62

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	124	GLN
1	A	126	LEU
1	A	259	GLN
1	A	285	ASN
1	A	340	ASN
1	A	376	ASN
1	A	419	ILE
1	B	114	ARG
1	B	259	GLN
1	B	260	ASP

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Mol	Chain	Res	Type
1	B	340	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	112	GLN
1	A	124	GLN
1	A	153	GLN
1	A	191	GLN
1	A	259	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	191	GLN
1	B	196	GLN
1	B	222	ASN
1	B	225	ASN
1	B	259	GLN
1	B	340	ASN
1	B	376	ASN
1	B	405	ASN
1	B	410	HIS

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 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
2	ACT	B	2860	-	3,3,3	0.86	0	3,3,3	0.90	0
2	ACT	B	2850	-	3,3,3	1.06	0	3,3,3	0.69	0
5	HEM	A	1500	1	50,50,50	1.18	4 (8%)	67,82,82	1.18	4 (5%)
8	GOL	A	1880	-	5,5,5	0.47	0	5,5,5	1.66	1 (20%)
7	ITU	B	2800	-	4,5,5	2.81	1 (25%)	3,5,5	2.83	1 (33%)
7	ITU	A	1800	-	4,5,5	1.69	1 (25%)	3,5,5	2.57	1 (33%)
6	BHS	A	1610	-	17,18,18	1.59	3 (17%)	14,26,26	1.85	3 (21%)
2	ACT	A	1850	-	3,3,3	0.93	0	3,3,3	0.82	0
5	HEM	B	2500	1	50,50,50	1.12	4 (8%)	67,82,82	1.06	2 (2%)
6	BHS	B	2610	-	17,18,18	1.73	3 (17%)	14,26,26	1.80	3 (21%)
8	GOL	B	2880	-	5,5,5	0.28	0	5,5,5	0.30	0
3	CAC	A	950	1	0,2,4	-	-	0,1,6	-	-
2	ACT	A	1860	-	3,3,3	0.90	0	3,3,3	0.73	0
3	CAC	B	950	1	0,2,4	-	-	0,1,6	-	-

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	A	1500	1	-	2/14/54/54	-
8	GOL	A	1880	-	-	4/4/4/4	-
7	ITU	B	2800	-	-	0/3/3/3	-
7	ITU	A	1800	-	-	0/3/3/3	-
6	BHS	A	1610	-	-	4/8/17/17	0/2/2/2
5	HEM	B	2500	1	-	1/14/54/54	-
6	BHS	B	2610	-	-	4/8/17/17	0/2/2/2
8	GOL	B	2880	-	-	0/4/4/4	-

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	B	2800	ITU	C3-S	5.56	1.81	1.75
6	B	2610	BHS	C7-N8	-4.86	1.41	1.46
6	A	1610	BHS	C7-N8	-4.24	1.42	1.46
6	B	2610	BHS	C4-N3	3.31	1.45	1.38
7	A	1800	ITU	C3-S	3.28	1.79	1.75
6	A	1610	BHS	C4-N3	3.19	1.44	1.38
5	B	2500	HEM	CHB-C1B	2.62	1.43	1.38
5	A	1500	HEM	CAB-C3B	-2.59	1.40	1.47
5	A	1500	HEM	C3B-C4B	2.52	1.49	1.44
5	A	1500	HEM	CMA-C3A	2.50	1.55	1.50
6	B	2610	BHS	C8A-N1	2.49	1.40	1.36
6	A	1610	BHS	C8A-N1	2.30	1.39	1.36
5	B	2500	HEM	CAB-C3B	-2.28	1.41	1.47
5	B	2500	HEM	CAC-C3C	-2.22	1.41	1.47
5	B	2500	HEM	FE-NA	2.20	2.02	1.95
5	A	1500	HEM	CAC-C3C	-2.10	1.41	1.47

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	1610	BHS	C2-N1-C8A	5.40	122.88	113.36
6	B	2610	BHS	C2-N1-C8A	5.25	122.63	113.36
7	B	2800	ITU	C2-S-C3	4.89	108.23	103.15
7	A	1800	ITU	C2-S-C3	4.45	107.77	103.15
5	A	1500	HEM	C3B-C4B-NB	3.28	111.82	109.47
5	B	2500	HEM	C3B-C4B-NB	3.08	111.68	109.47
6	A	1610	BHS	N3-C2-N1	-2.73	118.32	123.32
6	B	2610	BHS	N3-C2-N1	-2.72	118.34	123.32
5	A	1500	HEM	C2B-C1B-NB	2.71	112.96	109.84
5	B	2500	HEM	C3B-C2B-C1B	2.56	108.33	106.41
5	A	1500	HEM	CMA-C3A-C4A	2.44	129.13	125.42
8	A	1880	GOL	O3-C3-C2	2.29	120.70	110.38
6	B	2610	BHS	N2-C2-N3	2.18	121.37	116.76
6	A	1610	BHS	N2-C2-N3	2.06	121.12	116.76
5	A	1500	HEM	C1B-NB-C4B	-2.05	102.78	105.21

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	1610	BHS	O1'-C1'-C6-N5

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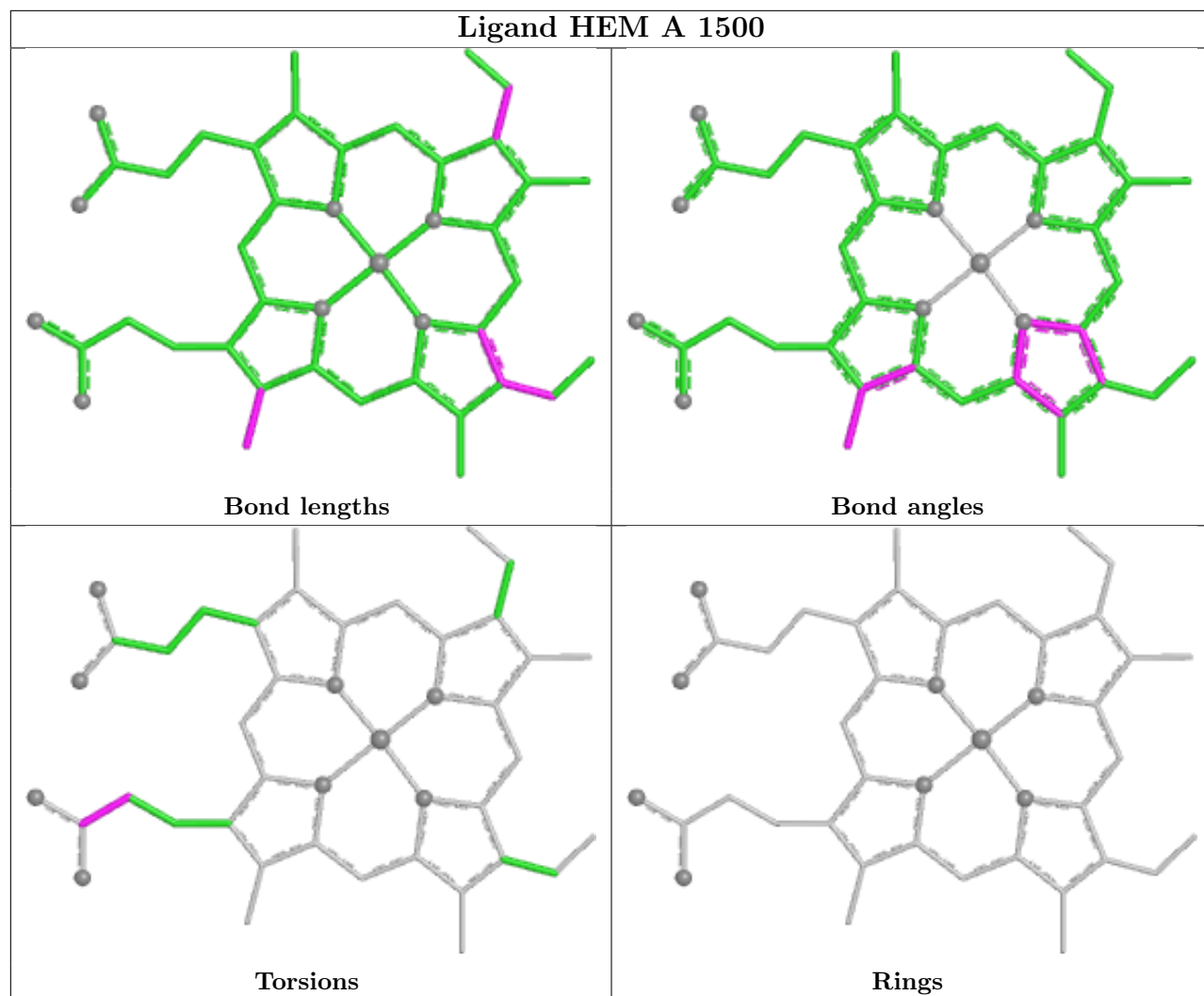
Mol	Chain	Res	Type	Atoms
6	A	1610	BHS	C2'-C1'-C6-N5
6	A	1610	BHS	O1'-C1'-C6-C7
6	B	2610	BHS	O1'-C1'-C6-N5
6	B	2610	BHS	C2'-C1'-C6-N5
6	B	2610	BHS	O1'-C1'-C6-C7
6	B	2610	BHS	C2'-C1'-C6-C7
8	A	1880	GOL	C1-C2-C3-O3
8	A	1880	GOL	O1-C1-C2-C3
8	A	1880	GOL	O1-C1-C2-O2
6	A	1610	BHS	C2'-C1'-C6-C7
8	A	1880	GOL	O2-C2-C3-O3
5	A	1500	HEM	CAA-CBA-CGA-O2A
5	A	1500	HEM	CAA-CBA-CGA-O1A
5	B	2500	HEM	CAA-CBA-CGA-O2A

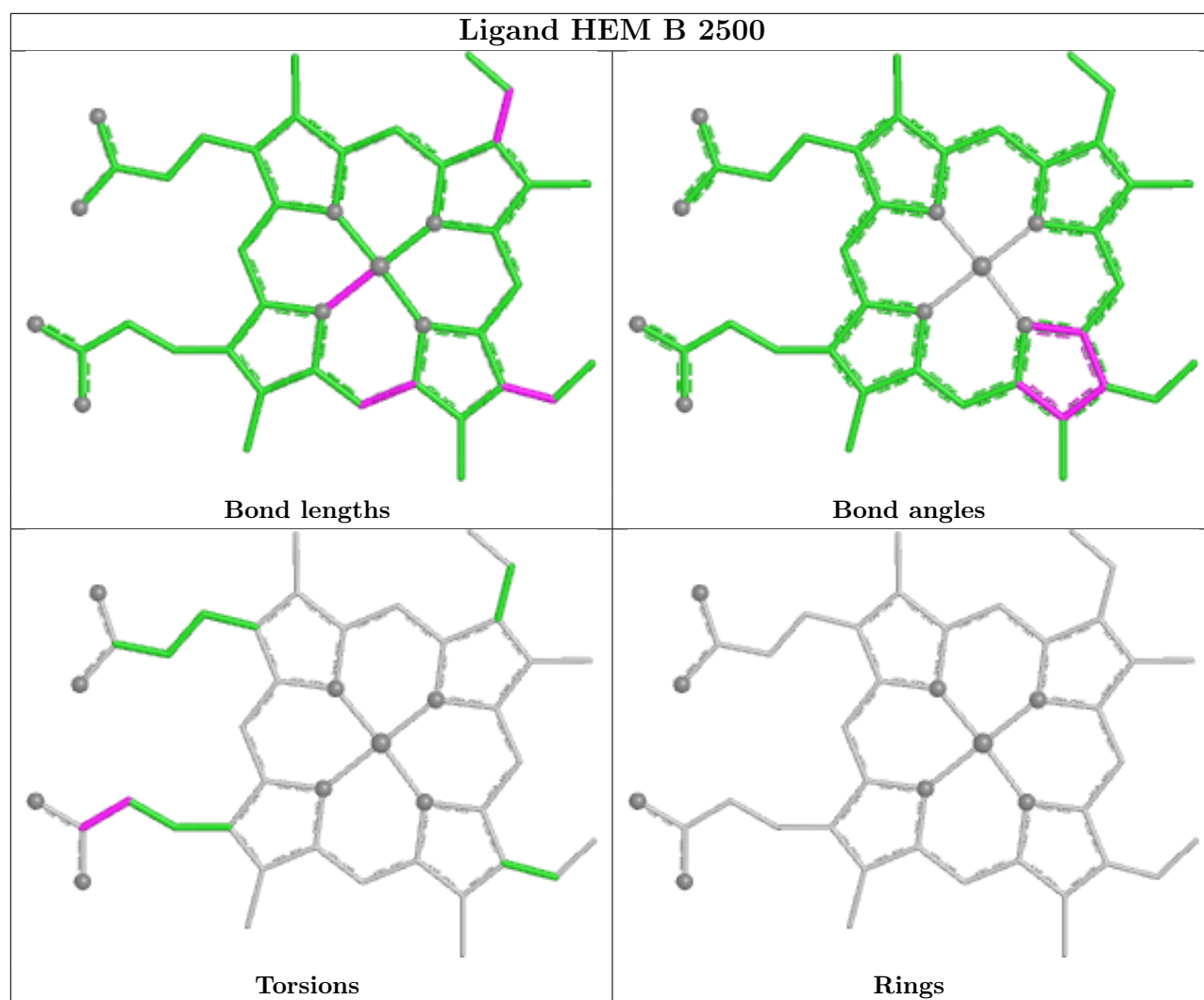
There are no ring outliers.

4 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	A	1800	ITU	2	0
6	A	1610	BHS	1	0
5	B	2500	HEM	2	0
6	B	2610	BHS	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.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	416/444 (93%)	0.69	43 (10%) 12 11	32, 47, 89, 99	0
1	B	414/444 (93%)	0.95	58 (14%) 6 5	35, 52, 88, 99	0
All	All	830/888 (93%)	0.82	101 (12%) 8 7	32, 49, 89, 99	0

All (101) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	113	THR	8.3
1	B	118	GLY	7.5
1	B	117	PRO	7.5
1	B	113	THR	7.4
1	A	117	PRO	6.5
1	A	115	PRO	6.1
1	B	120	PRO	5.5
1	B	119	PRO	5.1
1	B	98	PRO	4.7
1	B	111	LEU	4.7
1	B	261	GLY	4.6
1	A	111	LEU	4.6
1	A	118	GLY	4.4
1	A	119	PRO	4.1
1	A	108	PRO	4.1
1	B	115	PRO	4.1
1	A	107	LEU	4.0
1	B	116	SER	4.0
1	B	103	GLY	3.9
1	A	106	VAL	3.8
1	B	258	GLN	3.8
1	B	112	GLN	3.6
1	B	121	PRO	3.6
1	B	105	LEU	3.5

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Mol	Chain	Res	Type	RSRZ
1	B	73	VAL	3.5
1	A	109	ARG	3.4
1	A	121	PRO	3.3
1	A	97	THR	3.2
1	A	239	GLY	3.2
1	B	79	GLY	3.2
1	A	120	PRO	3.1
1	B	282	THR	3.1
1	B	108	PRO	3.1
1	A	261	GLY	3.0
1	A	116	SER	2.9
1	B	384	CYS	2.9
1	B	275	LEU	2.8
1	A	95	PRO	2.8
1	A	110	LYS	2.8
1	A	94	GLY	2.8
1	B	78	LEU	2.8
1	A	68	PRO	2.8
1	A	127	SER	2.7
1	B	70	PHE	2.7
1	B	97	THR	2.7
1	A	262	SER	2.7
1	B	93	ASP	2.7
1	B	311	LEU	2.7
1	A	98	PRO	2.7
1	A	284	GLY	2.7
1	B	91	GLN	2.7
1	B	107	LEU	2.7
1	B	260	ASP	2.7
1	B	110	LYS	2.7
1	B	94	GLY	2.7
1	A	99	ARG	2.6
1	A	105	LEU	2.6
1	B	283	PRO	2.6
1	B	259	GLN	2.6
1	A	96	CYS	2.6
1	A	114	ARG	2.6
1	B	326	ALA	2.5
1	A	147	ALA	2.5
1	A	238	PRO	2.5
1	B	95	PRO	2.5
1	B	277	ILE	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	143	SER	2.5
1	B	71	PRO	2.5
1	B	109	ARG	2.5
1	B	280	GLY	2.4
1	B	106	VAL	2.4
1	B	482	TRP	2.4
1	A	90	SER	2.4
1	B	82	THR	2.4
1	B	125	LEU	2.4
1	B	96	CYS	2.4
1	B	391	THR	2.4
1	B	88	ALA	2.3
1	A	103	GLY	2.3
1	A	143	SER	2.3
1	A	91	GLN	2.3
1	B	147	ALA	2.3
1	A	275	LEU	2.3
1	A	88	ALA	2.3
1	A	122	ALA	2.3
1	B	328	LEU	2.3
1	B	274	GLU	2.2
1	A	112	GLN	2.2
1	B	81	ILE	2.2
1	B	101	CYS	2.2
1	B	144	GLY	2.2
1	B	262	SER	2.2
1	A	259	GLN	2.2
1	A	125	LEU	2.1
1	A	365	GLY	2.1
1	A	92	GLN	2.1
1	B	138	SER	2.1
1	B	122	ALA	2.1
1	B	104	SER	2.0
1	B	127	SER	2.0
1	A	102	LEU	2.0

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](#)

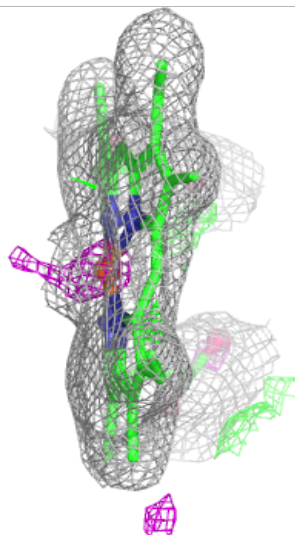
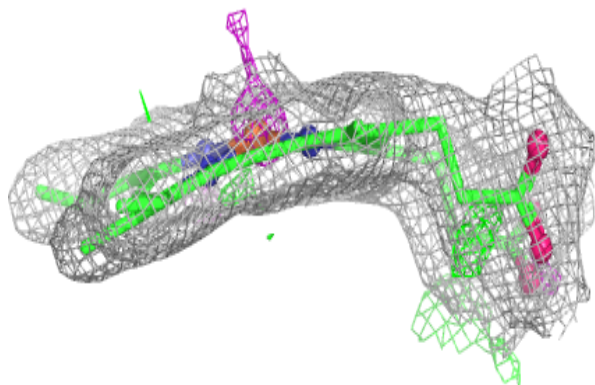
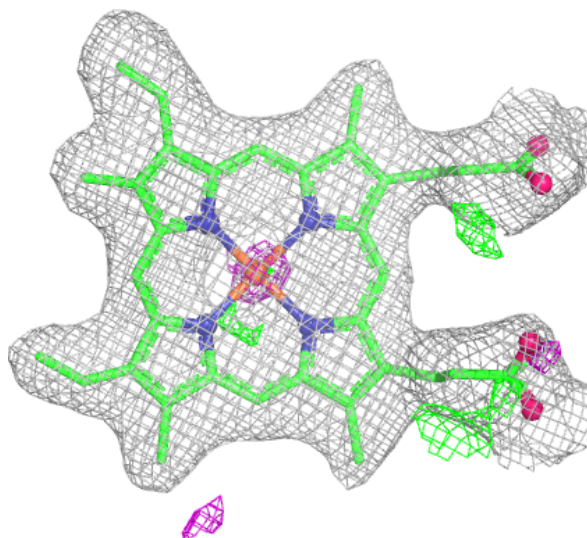
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
8	GOL	A	1880	6/6	0.85	0.13	63,64,65,65	0
2	ACT	B	2850	4/4	0.88	0.11	49,50,50,50	0
8	GOL	B	2880	6/6	0.89	0.13	61,65,66,68	0
3	CAC	B	950	3/5	0.90	0.15	85,85,85,89	0
6	BHS	A	1610	17/17	0.90	0.12	63,65,66,67	0
2	ACT	A	1850	4/4	0.91	0.11	50,50,51,52	0
6	BHS	B	2610	17/17	0.92	0.10	66,66,67,67	0
2	ACT	B	2860	4/4	0.93	0.08	55,55,55,56	0
2	ACT	A	1860	4/4	0.93	0.08	52,52,53,54	0
3	CAC	A	950	3/5	0.94	0.19	69,69,69,75	0
7	ITU	B	2800	6/6	0.97	0.09	44,45,46,47	0
5	HEM	A	1500	43/43	0.97	0.08	29,34,45,48	0
5	HEM	B	2500	43/43	0.97	0.07	33,35,46,51	0
7	ITU	A	1800	6/6	0.98	0.07	37,39,40,40	0
4	ZN	A	900	1/1	0.98	0.12	69,69,69,69	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.

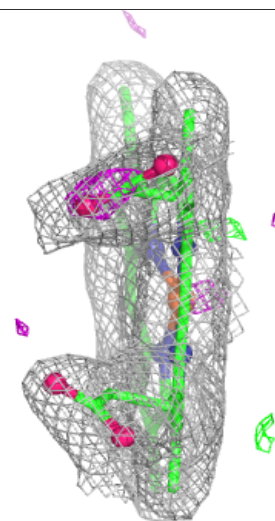
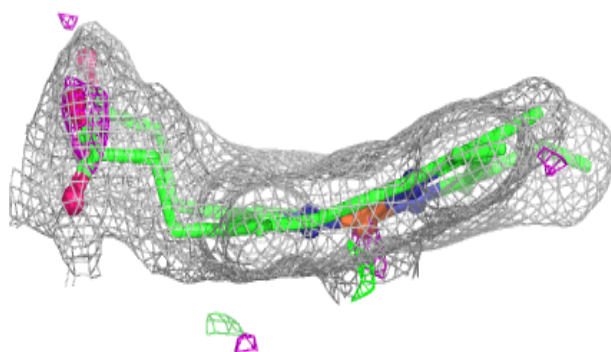
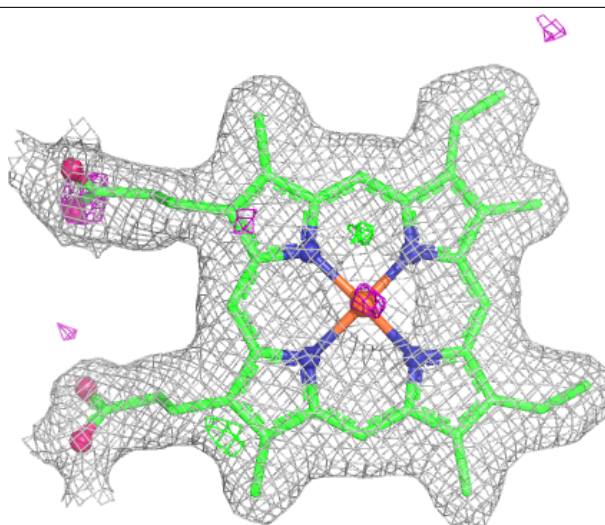
Electron density around HEM A 1500:

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



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)



6.5 Other polymers [i](#)

There are no such residues in this entry.