



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 12:17 PM UTC

PDB ID : 1K2T / pdb_00001k2t
Title : Structure of rat brain nNOS heme domain complexed with S-ethyl-N-phenyl-isothiourea
Authors : Li, H.; Martasek, P.; Masters, B.S.S.; Poulos, T.L.; Raman, C.S.
Deposited on : 2001-09-28
Resolution : 2.20 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 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

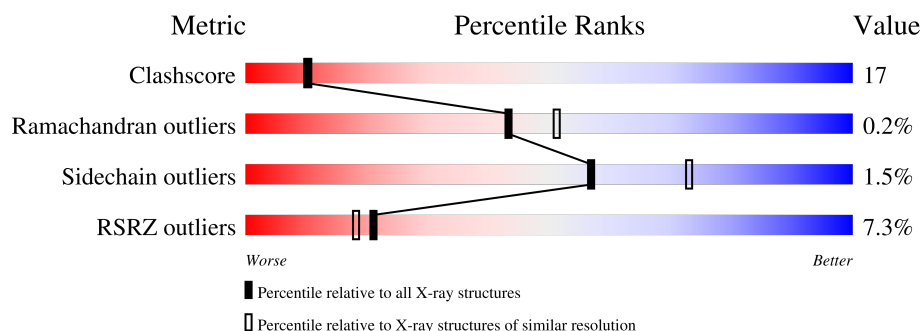
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. 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	419	11% 58% 37% ..
1	B	419	3% 66% 30% ..

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
6	PTU	B	2810	-	X	-	-

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 7129 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	407	Total	C	N	O	S	0	0	0
			3313	2121	566	605	21			
1	B	410	Total	C	N	O	S	0	0	0
			3341	2138	573	609	21			

- Molecule 2 is ACETATE ION (CCD ID: ACT) (formula: $C_2H_3O_2^-$).

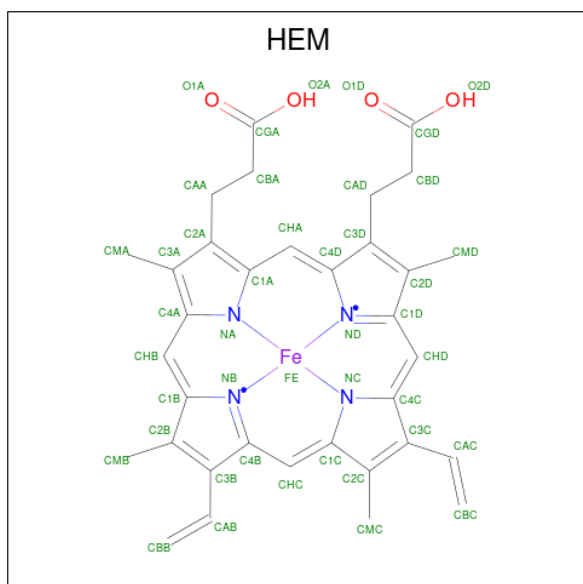


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

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

- Molecule 4 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



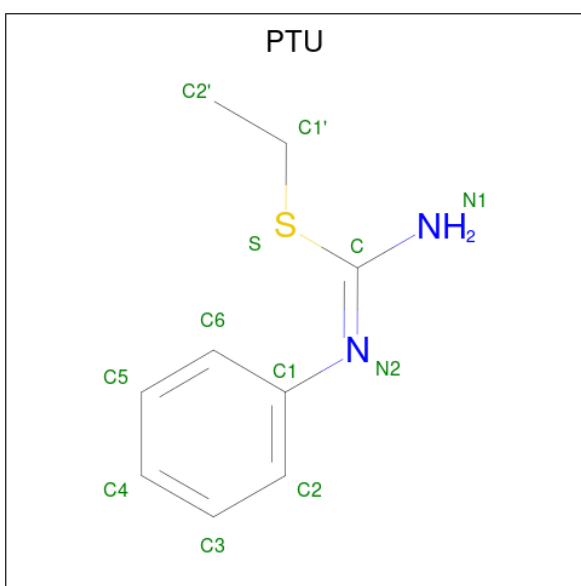
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
4	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 5 is 5,6,7,8-TETRAHYDROBIOPTERIN (CCD ID: H4B) (formula: $\text{C}_9\text{H}_{15}\text{N}_5\text{O}_3$).



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

- Molecule 6 is 2-ETHYL-1-PHENYL-ISOTHIOUREA (CCD ID: PTU) (formula: $C_9H_{12}N_2S$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	N	S	0	0
			12	9	2	1		
6	B	1	Total	C	N	S	0	0
			12	9	2	1		

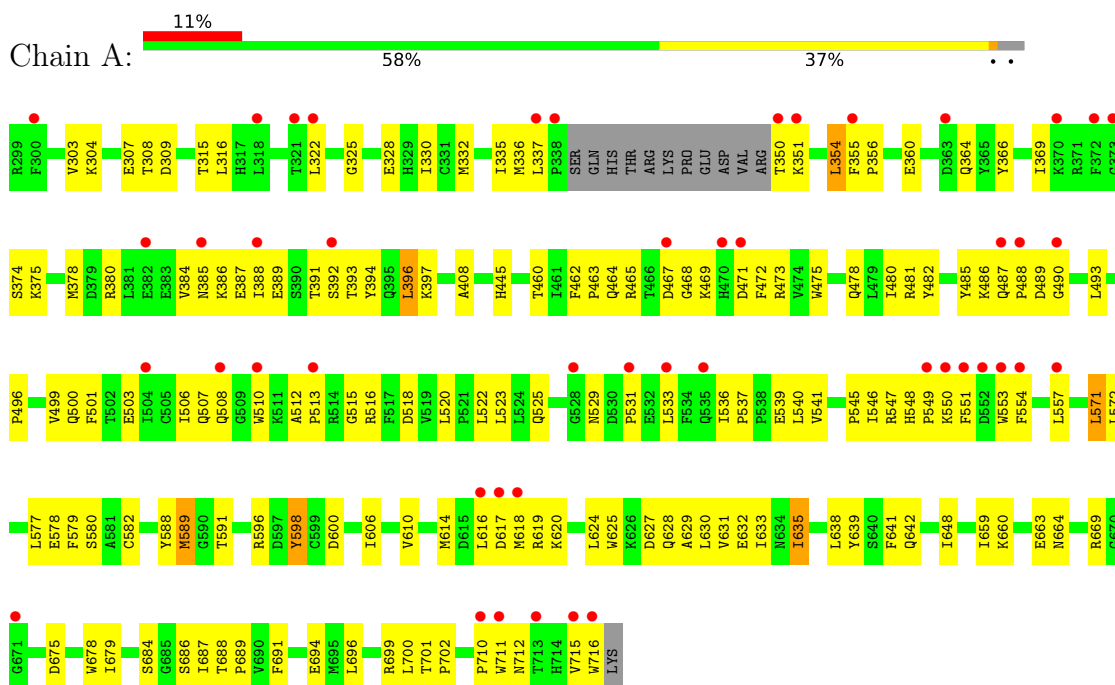
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	130	Total 130	O 130	0	0
7	B	192	Total 192	O 192	0	0

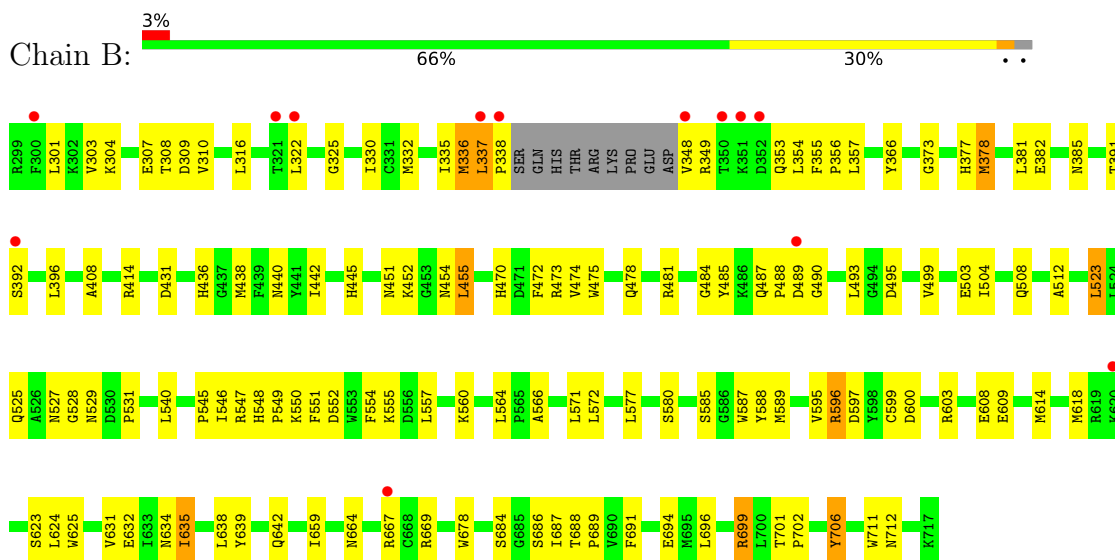
3 Residue-property plots [i](#)

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



• 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	52.00Å 111.24Å 165.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.70 – 2.20 29.70 – 2.20	Depositor EDS
% Data completeness (in resolution range)	92.2 (29.70-2.20) 82.1 (29.70-2.20)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.81 (at 2.20Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.224 , 0.265 0.219 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	30.2	Xtriage
Anisotropy	0.685	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 44.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7129	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.75% 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: PTU, ACT, H4B, ZN, HEM

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	0.39	0/3406	0.95	12/4621 (0.3%)
1	B	0.42	0/3434	0.93	17/4656 (0.4%)
All	All	0.40	0/6840	0.94	29/9277 (0.3%)

There are no bond length outliers.

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	589	MET	N-CA-C	-8.36	95.62	109.24
1	A	589	MET	N-CA-C	-7.88	96.39	109.24
1	A	588	TYR	N-CA-C	7.35	120.86	110.50
1	A	354	LEU	N-CA-C	6.98	118.97	111.36
1	A	408	ALA	N-CA-C	-6.65	104.11	111.36
1	B	354	LEU	N-CA-C	6.41	118.34	111.36
1	B	455	LEU	N-CA-C	6.10	118.84	110.24
1	A	490	GLY	N-CA-C	-6.04	107.07	114.92
1	A	635	ILE	N-CA-C	-6.02	104.64	110.72
1	A	688	THR	N-CA-C	-5.97	102.63	110.39
1	B	408	ALA	N-CA-C	-5.89	104.86	111.28
1	B	474	VAL	N-CA-C	-5.88	99.27	107.80
1	B	588	TYR	N-CA-C	5.60	117.97	110.35
1	A	315	THR	N-CA-C	-5.44	107.65	114.56
1	B	599	CYS	N-CA-C	5.38	120.72	113.72
1	B	635	ILE	N-CA-C	-5.34	105.29	110.42
1	B	472	PHE	N-CA-C	-5.31	100.59	109.24
1	A	375	LYS	N-CA-C	-5.23	105.58	111.28
1	A	546	ILE	N-CA-C	5.20	115.58	107.99
1	A	598	TYR	N-CA-C	5.20	118.09	111.69
1	A	325	GLY	N-CA-C	-5.17	108.54	114.69
1	B	596	ARG	N-CA-C	5.17	117.91	111.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	490	GLY	N-CA-C	-5.16	108.55	114.69
1	B	564	LEU	N-CA-C	5.14	117.94	108.69
1	B	488	PRO	N-CA-C	-5.13	107.17	113.84
1	B	699	ARG	N-CA-C	5.13	117.60	109.24
1	B	436	HIS	N-CA-C	-5.12	105.70	111.28
1	B	706	TYR	N-CA-C	-5.10	103.66	110.55
1	B	688	THR	N-CA-C	-5.08	102.46	110.14

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	3313	0	3221	133	0
1	B	3341	0	3256	106	0
2	A	4	0	3	0	0
2	B	4	0	3	0	0
3	A	1	0	0	0	0
4	A	43	0	30	1	0
4	B	43	0	30	2	0
5	A	17	0	15	1	0
5	B	17	0	15	1	0
6	A	12	0	12	0	0
6	B	12	0	12	0	0
7	A	130	0	0	6	0
7	B	192	0	0	13	0
All	All	7129	0	6597	230	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 17.

All (230) 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:523:LEU:HD22	1:A:531:PRO:HB2	1.40	1.02
1:B:523:LEU:HD22	1:B:531:PRO:HB2	1.49	0.94
1:A:537:PRO:HB2	1:A:540:LEU:HD13	1.49	0.94
1:B:373:GLY:H	1:B:377:HIS:HD2	1.14	0.91
1:A:635:ILE:HD11	1:B:624:LEU:HB2	1.58	0.85
1:B:322:LEU:HB3	1:B:699:ARG:HH21	1.41	0.85
1:A:549:PRO:HG3	1:A:639:TYR:CG	2.16	0.81
1:B:631:VAL:O	1:B:635:ILE:HG12	1.81	0.80
1:B:382:GLU:HG3	7:B:2952:HOH:O	1.85	0.77
1:A:596:ARG:O	1:A:600:ASP:HB2	1.86	0.75
1:B:548:HIS:HD2	1:B:550:LYS:H	1.35	0.74
1:A:385:ASN:O	1:A:389:GLU:HG3	1.87	0.74
1:B:470:HIS:HB3	1:B:527:ASN:ND2	2.03	0.73
1:B:549:PRO:HG3	1:B:639:TYR:CG	2.23	0.73
1:A:635:ILE:CD1	1:B:624:LEU:HB2	2.19	0.72
1:B:348:VAL:HG12	1:B:349:ARG:H	1.54	0.71
1:B:325:GLY:O	1:B:332:MET:HE2	1.90	0.71
1:B:499:VAL:O	1:B:503:GLU:HG3	1.93	0.69
1:B:489:ASP:HB2	7:B:3010:HOH:O	1.93	0.68
1:A:554:PHE:O	1:A:557:LEU:HD13	1.94	0.67
1:A:316:LEU:HD11	1:A:669:ARG:HD3	1.77	0.66
1:A:322:LEU:HB2	1:A:699:ARG:HD3	1.78	0.66
1:B:478:GLN:HB2	1:B:481:ARG:HG3	1.77	0.65
4:A:750:HEM:HBC2	4:A:750:HEM:HMC2	1.79	0.65
1:B:348:VAL:HG12	1:B:349:ARG:N	2.13	0.64
1:A:501:PHE:HD2	1:A:520:LEU:HD13	1.61	0.64
1:B:373:GLY:H	1:B:377:HIS:CD2	2.06	0.64
1:A:553:TRP:CZ3	1:A:557:LEU:HD11	2.33	0.64
1:A:627:ASP:O	1:A:631:VAL:HG23	1.98	0.64
1:A:684:SER:HB3	1:A:687:ILE:HG12	1.79	0.63
1:B:548:HIS:CD2	1:B:550:LYS:H	2.16	0.62
1:B:336:MET:HE2	1:B:678:TRP:HZ2	1.63	0.62
1:B:545:PRO:HG2	1:B:547:ARG:NH1	2.15	0.62
1:A:571:LEU:HD12	1:A:572:LEU:N	2.15	0.61
1:A:467:ASP:OD2	1:A:469:LYS:HB2	2.00	0.61
1:B:701:THR:HA	1:B:702:PRO:C	2.25	0.61
1:A:350:THR:O	1:A:354:LEU:N	2.33	0.61
1:B:504:ILE:O	1:B:508:GLN:HG2	2.00	0.61
1:A:485:TYR:HE2	1:A:512:ALA:HB1	1.65	0.60
1:A:485:TYR:CE2	1:A:512:ALA:HB1	2.36	0.60
1:A:307:GLU:HG3	7:B:2982:HOH:O	2.00	0.60
1:B:608:GLU:HG3	7:B:3028:HOH:O	2.01	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:316:LEU:HD22	1:A:700:LEU:HD11	1.84	0.59
1:A:475:TRP:HB2	1:A:523:LEU:HB3	1.84	0.59
1:A:616:LEU:HD13	1:A:625:TRP:HB2	1.84	0.59
1:A:304:LYS:O	1:A:694:GLU:HG3	2.01	0.59
1:A:501:PHE:CD2	1:A:520:LEU:HD13	2.38	0.59
1:A:638:LEU:O	1:A:642:GLN:HG3	2.02	0.59
1:B:525:GLN:HG3	1:B:529:ASN:O	2.03	0.59
1:A:545:PRO:HG2	1:A:547:ARG:NH1	2.18	0.58
1:B:332:MET:HB3	1:B:335:ILE:HG13	1.84	0.58
1:A:660:LYS:HE2	7:A:1911:HOH:O	2.03	0.57
1:A:496:PRO:HA	1:A:499:VAL:HG23	1.86	0.57
1:A:701:THR:HA	1:A:702:PRO:C	2.30	0.57
1:A:545:PRO:HG2	1:A:547:ARG:HH11	1.69	0.56
1:A:303:VAL:HG13	1:A:694:GLU:HB2	1.87	0.56
1:A:508:GLN:HE22	1:A:716:TRP:HZ3	1.52	0.56
1:B:475:TRP:HB2	1:B:523:LEU:HB3	1.88	0.56
1:A:516:ARG:HG3	1:A:516:ARG:HH11	1.70	0.56
1:B:669:ARG:HD3	7:B:2996:HOH:O	2.06	0.56
1:A:388:ILE:O	1:A:392:SER:HA	2.05	0.56
1:A:571:LEU:HD11	1:A:578:GLU:HB3	1.88	0.56
1:A:316:LEU:HD11	1:A:669:ARG:CD	2.35	0.55
1:A:525:GLN:HG3	1:A:529:ASN:O	2.06	0.55
1:B:684:SER:HB3	1:B:687:ILE:HG12	1.87	0.55
1:A:360:GLU:O	1:A:364:GLN:HG3	2.07	0.55
1:B:664:ASN:O	1:B:667:ARG:HG2	2.07	0.55
1:A:507:GLN:O	1:A:507:GLN:HG2	2.06	0.55
1:B:684:SER:HB3	1:B:687:ILE:HD11	1.87	0.55
1:B:684:SER:HB3	1:B:687:ILE:CG1	2.37	0.55
1:A:696:LEU:HD22	1:B:330:ILE:HD11	1.88	0.54
1:B:478:GLN:HB2	1:B:481:ARG:CG	2.37	0.54
1:A:378:MET:HE2	1:A:378:MET:HA	1.88	0.54
1:B:391:THR:O	1:B:392:SER:HB2	2.08	0.54
1:A:487:GLN:HB3	1:A:488:PRO:HD2	1.89	0.54
1:B:322:LEU:HB2	1:B:699:ARG:HE	1.71	0.54
4:B:750:HEM:HMC1	4:B:750:HEM:HBC2	1.90	0.54
1:A:549:PRO:HG3	1:A:639:TYR:CD1	2.43	0.53
1:B:493:LEU:HD22	7:B:2942:HOH:O	2.07	0.53
1:A:487:GLN:HB3	1:A:488:PRO:CD	2.38	0.53
1:B:445:HIS:C	1:B:445:HIS:CD2	2.87	0.53
1:A:610:VAL:HG21	1:A:633:ILE:HD11	1.91	0.53
1:B:332:MET:HE1	1:B:338:PRO:HB3	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:566:ALA:HB2	1:B:585:SER:HB3	1.91	0.53
4:B:750:HEM:HBC2	4:B:750:HEM:CMC	2.38	0.53
1:A:354:LEU:CD2	1:A:393:THR:HA	2.39	0.52
1:A:554:PHE:HA	1:A:557:LEU:HD13	1.91	0.52
1:A:330:ILE:HD11	1:B:696:LEU:HD22	1.92	0.52
1:A:510:TRP:HB2	1:A:533:LEU:HD13	1.92	0.52
1:B:684:SER:HB3	1:B:687:ILE:CD1	2.40	0.52
1:B:440:ASN:ND2	7:B:3007:HOH:O	2.43	0.51
1:B:473:ARG:HD3	1:B:580:SER:HB2	1.93	0.51
1:B:336:MET:HE2	1:B:678:TRP:CZ2	2.45	0.51
1:A:522:LEU:O	1:A:533:LEU:HA	2.11	0.51
1:B:548:HIS:CD2	1:B:551:PHE:H	2.29	0.51
1:A:335:ILE:HD13	1:B:694:GLU:HB3	1.93	0.50
1:A:624:LEU:HB2	1:B:635:ILE:CD1	2.41	0.50
1:B:355:PHE:CE1	1:B:385:ASN:HB2	2.47	0.50
1:B:470:HIS:HB3	1:B:527:ASN:HD22	1.75	0.50
1:B:325:GLY:C	1:B:332:MET:HE2	2.37	0.50
1:B:336:MET:HE1	5:B:2760:H4B:N5	2.26	0.50
1:A:355:PHE:N	1:A:356:PRO:HD2	2.27	0.50
1:B:308:THR:OG1	1:B:310:VAL:HG23	2.12	0.50
1:B:487:GLN:NE2	7:B:3040:HOH:O	2.45	0.50
1:A:624:LEU:HB2	1:B:635:ILE:HD13	1.94	0.49
1:B:308:THR:O	1:B:309:ASP:HB2	2.12	0.49
1:B:485:TYR:CE2	1:B:512:ALA:HB1	2.46	0.49
1:A:548:HIS:CD2	1:A:549:PRO:HD2	2.47	0.49
1:A:354:LEU:HD23	1:A:388:ILE:HG23	1.95	0.49
1:A:380:ARG:NH1	1:A:397:LYS:HE3	2.27	0.49
1:A:618:MET:HE3	1:A:625:TRP:CH2	2.46	0.49
1:B:332:MET:HB3	1:B:335:ILE:CG1	2.42	0.49
1:A:614:MET:HE3	1:A:632:GLU:HG3	1.95	0.48
1:A:686:SER:HA	1:A:691:PHE:CG	2.48	0.48
1:B:485:TYR:HE2	1:B:512:ALA:HB1	1.78	0.48
1:B:595:VAL:HG23	1:B:634:ASN:HD21	1.78	0.48
1:B:332:MET:CE	1:B:338:PRO:HB3	2.43	0.48
1:A:478:GLN:HB2	1:A:481:ARG:CG	2.44	0.48
1:A:500:GLN:O	1:A:503:GLU:HB2	2.14	0.48
1:B:638:LEU:O	1:B:642:GLN:HG3	2.14	0.48
1:A:486:LYS:HG3	7:A:1889:HOH:O	2.14	0.48
1:A:675:ASP:O	1:A:679:ILE:HG12	2.14	0.48
1:B:597:ASP:OD1	1:B:603:ARG:NH1	2.46	0.48
1:A:473:ARG:NH2	1:A:710:PRO:HD3	2.30	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:641:PHE:CG	1:A:648:ILE:HD12	2.48	0.47
1:B:548:HIS:CD2	1:B:549:PRO:HD2	2.49	0.47
1:A:659:ILE:O	1:A:663:GLU:HG3	2.15	0.47
1:B:596:ARG:O	1:B:600:ASP:HB2	2.14	0.47
1:A:316:LEU:CD2	1:A:700:LEU:HD11	2.43	0.47
1:B:487:GLN:C	1:B:489:ASP:N	2.73	0.47
1:A:617:ASP:OD1	1:A:619:ARG:NE	2.37	0.47
1:B:566:ALA:CB	1:B:585:SER:HB3	2.45	0.47
1:A:512:ALA:HA	1:A:513:PRO:HD3	1.79	0.46
1:A:551:PHE:HB3	1:A:553:TRP:NE1	2.30	0.46
1:A:618:MET:HE3	1:A:625:TRP:CZ3	2.50	0.46
1:B:322:LEU:HB3	1:B:699:ARG:NH2	2.22	0.46
1:A:332:MET:HE1	1:B:301:LEU:HD22	1.97	0.46
1:A:557:LEU:HD12	1:A:557:LEU:N	2.30	0.46
1:B:470:HIS:HA	1:B:528:GLY:HA3	1.96	0.46
1:A:330:ILE:HD11	1:B:696:LEU:HB3	1.96	0.46
1:A:632:GLU:O	1:A:635:ILE:HB	2.15	0.46
1:A:616:LEU:HD12	1:A:616:LEU:O	2.15	0.46
1:A:473:ARG:HD3	1:A:580:SER:HB2	1.98	0.46
1:A:316:LEU:HD23	1:A:316:LEU:O	2.16	0.45
7:A:1896:HOH:O	1:B:337:LEU:HD12	2.15	0.45
1:B:451:ASN:O	1:B:452:LYS:HB2	2.16	0.45
1:B:455:LEU:HD12	1:B:587:TRP:HB3	1.99	0.45
1:B:546:ILE:HG12	1:B:560:LYS:HA	1.98	0.45
1:B:711:TRP:HD1	1:B:712:ASN:ND2	2.15	0.45
1:A:387:GLU:OE2	1:A:394:TYR:HA	2.16	0.45
1:A:354:LEU:HD21	1:A:393:THR:HA	1.98	0.45
1:A:553:TRP:HZ3	1:A:557:LEU:HD11	1.80	0.45
1:A:475:TRP:CE2	1:A:710:PRO:HB2	2.51	0.45
1:A:715:VAL:O	1:A:715:VAL:HG23	2.17	0.45
1:B:348:VAL:CG1	1:B:349:ARG:H	2.27	0.45
1:B:618:MET:HG2	1:B:625:TRP:CD2	2.52	0.45
1:B:686:SER:HA	1:B:691:PHE:CG	2.53	0.44
1:B:355:PHE:N	1:B:356:PRO:HD2	2.31	0.44
1:B:396:LEU:HG	1:B:577:LEU:HD12	1.99	0.44
1:A:468:GLY:HA2	1:A:471:ASP:OD1	2.18	0.44
1:A:493:LEU:HD13	1:A:493:LEU:C	2.43	0.44
1:A:628:GLN:HG2	1:B:631:VAL:HG11	1.99	0.44
1:A:380:ARG:CZ	1:A:397:LYS:HE3	2.48	0.44
1:A:684:SER:HB3	1:A:687:ILE:CG1	2.48	0.44
1:B:304:LYS:O	1:B:694:GLU:HG3	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:445:HIS:C	1:A:445:HIS:CD2	2.96	0.44
1:B:378:MET:HE3	1:B:381:LEU:HD12	1.98	0.44
1:A:350:THR:HG22	1:A:351:LYS:N	2.32	0.44
1:A:464:GLN:HB3	1:A:579:PHE:CE2	2.53	0.44
1:A:589:MET:HE3	1:A:591:THR:OG1	2.18	0.44
1:B:549:PRO:HG3	1:B:639:TYR:CD2	2.53	0.43
1:A:515:GLY:N	1:A:518:ASP:OD2	2.43	0.43
1:B:307:GLU:HG3	7:B:2955:HOH:O	2.17	0.43
1:B:659:ILE:HG13	1:B:689:PRO:HB2	2.00	0.43
1:A:629:ALA:O	1:A:630:LEU:C	2.61	0.43
1:B:451:ASN:HB3	1:B:454:ASN:O	2.18	0.43
1:B:557:LEU:HD23	1:B:609:GLU:OE2	2.18	0.43
1:B:571:LEU:HD12	1:B:572:LEU:N	2.33	0.43
1:A:396:LEU:HD13	1:A:577:LEU:HD12	2.01	0.43
1:A:635:ILE:CD1	1:B:623:SER:O	2.66	0.43
1:A:536:ILE:O	1:A:537:PRO:C	2.60	0.43
1:A:553:TRP:CE3	1:A:557:LEU:HD11	2.52	0.43
1:A:618:MET:HA	1:A:625:TRP:CD1	2.54	0.43
1:A:659:ILE:HG13	1:A:689:PRO:HB2	2.01	0.43
1:B:552:ASP:OD1	1:B:555:LYS:NZ	2.51	0.43
1:A:539:GLU:HG2	1:A:540:LEU:HD12	2.00	0.43
1:A:384:VAL:O	1:A:388:ILE:HG13	2.18	0.43
1:A:322:LEU:HD12	1:A:699:ARG:HD3	2.01	0.43
1:A:478:GLN:HB2	1:A:481:ARG:HG3	2.00	0.43
1:A:548:HIS:HD2	1:A:550:LYS:H	1.66	0.43
1:A:465:ARG:NE	1:A:471:ASP:OD2	2.50	0.42
1:A:523:LEU:HD22	1:A:531:PRO:CB	2.29	0.42
1:B:431:ASP:CG	7:B:2935:HOH:O	2.62	0.42
1:A:463:PRO:HB2	1:A:472:PHE:CE1	2.54	0.42
1:A:598:TYR:O	1:A:606:ILE:HG12	2.20	0.42
1:A:628:GLN:NE2	1:B:632:GLU:OE2	2.52	0.42
1:B:554:PHE:HB3	7:B:2977:HOH:O	2.19	0.42
1:A:462:PHE:HB3	1:A:463:PRO:CD	2.50	0.42
1:A:620:LYS:NZ	7:A:1971:HOH:O	2.52	0.42
1:B:438:MET:O	1:B:442:ILE:HG13	2.19	0.42
1:B:684:SER:O	1:B:687:ILE:HG12	2.20	0.42
1:A:554:PHE:C	1:A:557:LEU:HD13	2.44	0.42
1:A:391:THR:O	1:A:393:THR:HG22	2.19	0.42
1:A:366:TYR:HA	1:A:369:ILE:HG12	2.02	0.41
1:A:391:THR:O	1:A:392:SER:OG	2.36	0.41
1:A:482:TYR:HA	1:A:518:ASP:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:322:LEU:CB	1:B:699:ARG:HE	2.33	0.41
1:A:328:GLU:H	1:A:328:GLU:CD	2.29	0.41
1:A:465:ARG:CZ	7:A:1925:HOH:O	2.68	0.41
1:A:557:LEU:N	1:A:557:LEU:CD1	2.83	0.41
1:A:322:LEU:CB	1:A:699:ARG:HD3	2.48	0.41
1:B:366:TYR:CD2	1:B:377:HIS:HB2	2.56	0.41
1:B:473:ARG:CD	1:B:580:SER:HB2	2.50	0.41
1:B:508:GLN:NE2	7:B:3014:HOH:O	2.50	0.41
1:B:414:ARG:NH1	1:B:706:TYR:OH	2.53	0.41
1:A:460:THR:O	1:A:582:CYS:HA	2.21	0.41
1:B:614:MET:CE	1:B:632:GLU:HG3	2.51	0.41
1:A:386:LYS:HD3	1:A:386:LYS:HA	1.96	0.41
1:A:506:ILE:C	1:A:508:GLN:H	2.27	0.41
1:A:664:ASN:ND2	7:A:1911:HOH:O	2.48	0.41
1:B:632:GLU:O	1:B:635:ILE:HB	2.21	0.41
1:B:706:TYR:HA	7:B:2939:HOH:O	2.20	0.41
1:A:308:THR:O	1:A:309:ASP:HB2	2.21	0.40
1:B:484:GLY:HA3	1:B:495:ASP:O	2.21	0.40
1:A:480:ILE:HD13	1:A:541:VAL:HG13	2.03	0.40
1:B:353:GLN:O	1:B:357:LEU:HG	2.21	0.40
1:A:551:PHE:CD1	1:A:553:TRP:CZ2	3.09	0.40
1:A:614:MET:CE	1:A:632:GLU:HG3	2.52	0.40
1:A:678:TRP:HA	5:A:1760:H4B:N1	2.37	0.40
1:A:711:TRP:CD1	1:A:712:ASN:ND2	2.90	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	403/419 (96%)	370 (92%)	31 (8%)	2 (0%)	24 27

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	406/419 (97%)	385 (95%)	21 (5%)	0	100	100
All	All	809/838 (96%)	755 (93%)	52 (6%)	2 (0%)	43	51

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	374	SER
1	A	489	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	363/375 (97%)	359 (99%)	4 (1%)	65	79
1	B	366/375 (98%)	359 (98%)	7 (2%)	50	66
All	All	729/750 (97%)	718 (98%)	11 (2%)	57	73

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

Mol	Chain	Res	Type
1	A	336	MET
1	A	337	LEU
1	A	396	LEU
1	A	571	LEU
1	B	303	VAL
1	B	316	LEU
1	B	336	MET
1	B	337	LEU
1	B	378	MET
1	B	523	LEU
1	B	540	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	329	HIS
1	A	364	GLN
1	A	425	GLN
1	A	451	ASN
1	A	454	ASN
1	A	548	HIS
1	A	628	GLN
1	A	634	ASN
1	A	697	ASN
1	A	712	ASN
1	A	714	HIS
1	B	377	HIS
1	B	385	ASN
1	B	395	GLN
1	B	425	GLN
1	B	440	ASN
1	B	451	ASN
1	B	454	ASN
1	B	500	GLN
1	B	508	GLN
1	B	527	ASN
1	B	548	HIS
1	B	634	ASN
1	B	664	ASN
1	B	697	ASN
1	B	712	ASN

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 9 ligands modelled in this entry, 1 is monoatomic - leaving 8 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	A	1860	-	3,3,3	0.89	0	3,3,3	0.88	0
4	HEM	B	750	1	50,50,50	1.23	4 (8%)	67,82,82	1.19	3 (4%)
6	PTU	B	2810	-	12,12,12	2.01	6 (50%)	14,14,14	2.61	3 (21%)
5	H4B	B	2760	-	17,18,18	1.65	2 (11%)	14,26,26	1.81	3 (21%)
6	PTU	A	1810	-	12,12,12	1.95	4 (33%)	14,14,14	2.61	3 (21%)
5	H4B	A	1760	-	17,18,18	1.61	2 (11%)	14,26,26	1.96	3 (21%)
2	ACT	B	2860	-	3,3,3	0.89	0	3,3,3	0.88	0
4	HEM	A	750	1	50,50,50	1.17	4 (8%)	67,82,82	1.20	4 (5%)

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
4	HEM	B	750	1	-	1/14/54/54	-
6	PTU	B	2810	-	-	6/6/7/7	0/1/1/1
5	H4B	B	2760	-	-	0/8/17/17	0/2/2/2
6	PTU	A	1810	-	-	3/6/7/7	0/1/1/1
5	H4B	A	1760	-	-	0/8/17/17	0/2/2/2
4	HEM	A	750	1	-	0/14/54/54	-

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	1760	H4B	C6-N5	4.11	1.54	1.47
5	B	2760	H4B	C6-N5	3.90	1.54	1.47
5	A	1760	H4B	C4A-N5	3.88	1.47	1.37
5	B	2760	H4B	C4A-N5	3.68	1.46	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	2810	PTU	C3-C2	3.18	1.44	1.38
6	B	2810	PTU	C6-C1	3.13	1.45	1.39
6	A	1810	PTU	C6-C1	3.09	1.45	1.39
6	A	1810	PTU	C3-C2	3.04	1.44	1.38
4	B	750	HEM	FE-ND	2.64	2.03	1.94
4	A	750	HEM	CAC-C3C	-2.60	1.40	1.47
6	B	2810	PTU	C1-N2	-2.53	1.38	1.42
4	B	750	HEM	CAB-C3B	-2.46	1.40	1.47
4	B	750	HEM	CHB-C1B	2.41	1.43	1.38
4	A	750	HEM	FE-NA	2.36	2.02	1.95
6	A	1810	PTU	C1-N2	-2.34	1.38	1.42
6	A	1810	PTU	C-S	2.19	1.81	1.77
4	A	750	HEM	CMC-C2C	2.18	1.55	1.50
4	A	750	HEM	CAB-C3B	-2.14	1.41	1.47
6	B	2810	PTU	C2-C1	2.13	1.43	1.39
4	B	750	HEM	CAC-C3C	-2.10	1.41	1.47
6	B	2810	PTU	C-S	2.08	1.81	1.77
6	B	2810	PTU	C5-C6	2.04	1.42	1.38

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	1810	PTU	C1'-S-C	8.03	111.50	103.15
6	B	2810	PTU	C1'-S-C	8.00	111.47	103.15
5	A	1760	H4B	C2-N1-C8A	5.38	122.86	113.36
4	B	750	HEM	CMA-C3A-C4A	5.04	133.09	125.42
5	B	2760	H4B	C2-N1-C8A	4.97	122.14	113.36
6	B	2810	PTU	C1-N2-C	4.47	135.27	122.15
6	A	1810	PTU	C1-N2-C	4.32	134.82	122.15
5	A	1760	H4B	O10-C10-C9	-3.02	104.77	109.77
4	A	750	HEM	C3B-C4B-NB	3.01	111.63	109.47
5	A	1760	H4B	N3-C2-N1	-2.84	118.13	123.32
5	B	2760	H4B	N3-C2-N1	-2.74	118.31	123.32
4	A	750	HEM	CMD-C2D-C1D	2.63	129.15	125.03
4	A	750	HEM	CMA-C3A-C4A	2.36	129.01	125.42
6	A	1810	PTU	S-C-N2	2.13	123.47	119.73
5	B	2760	H4B	O9-C9-C6	2.07	114.24	109.28
4	B	750	HEM	C4A-NA-C1A	-2.06	102.46	105.82
4	B	750	HEM	CMA-C3A-C2A	-2.03	121.32	125.62
4	A	750	HEM	C4D-C3D-C2D	2.03	109.84	106.89
6	B	2810	PTU	S-C-N2	2.01	123.26	119.73

There are no chirality outliers.

All (10) torsion outliers are listed below:

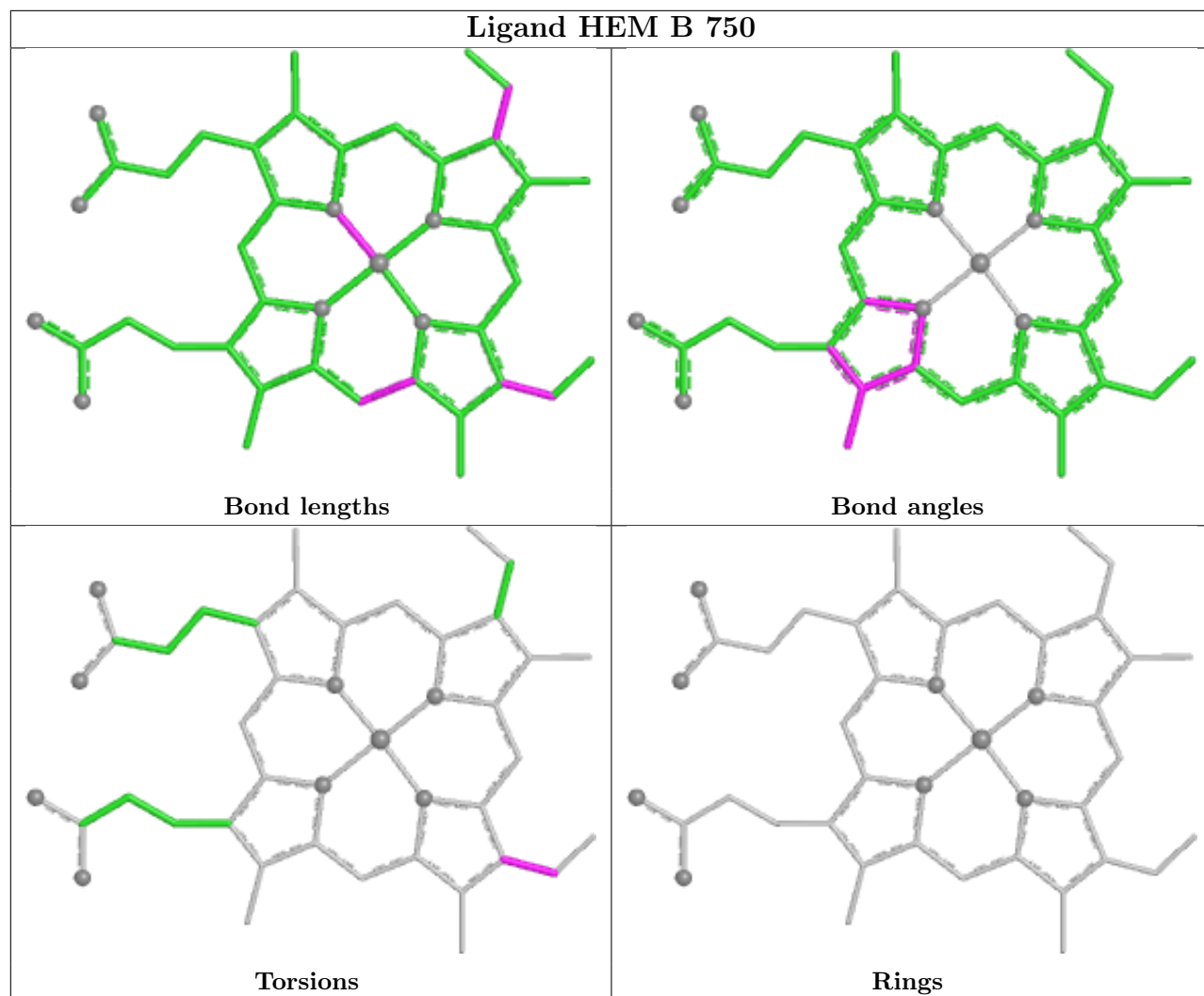
Mol	Chain	Res	Type	Atoms
6	A	1810	PTU	N2-C-S-C1'
6	B	2810	PTU	S-C-N2-C1
6	B	2810	PTU	C2'-C1'-S-C
6	A	1810	PTU	N1-C-S-C1'
6	B	2810	PTU	N2-C-S-C1'
6	A	1810	PTU	C2'-C1'-S-C
6	B	2810	PTU	C6-C1-N2-C
6	B	2810	PTU	N1-C-S-C1'
4	B	750	HEM	C4B-C3B-CAB-CBB
6	B	2810	PTU	C2-C1-N2-C

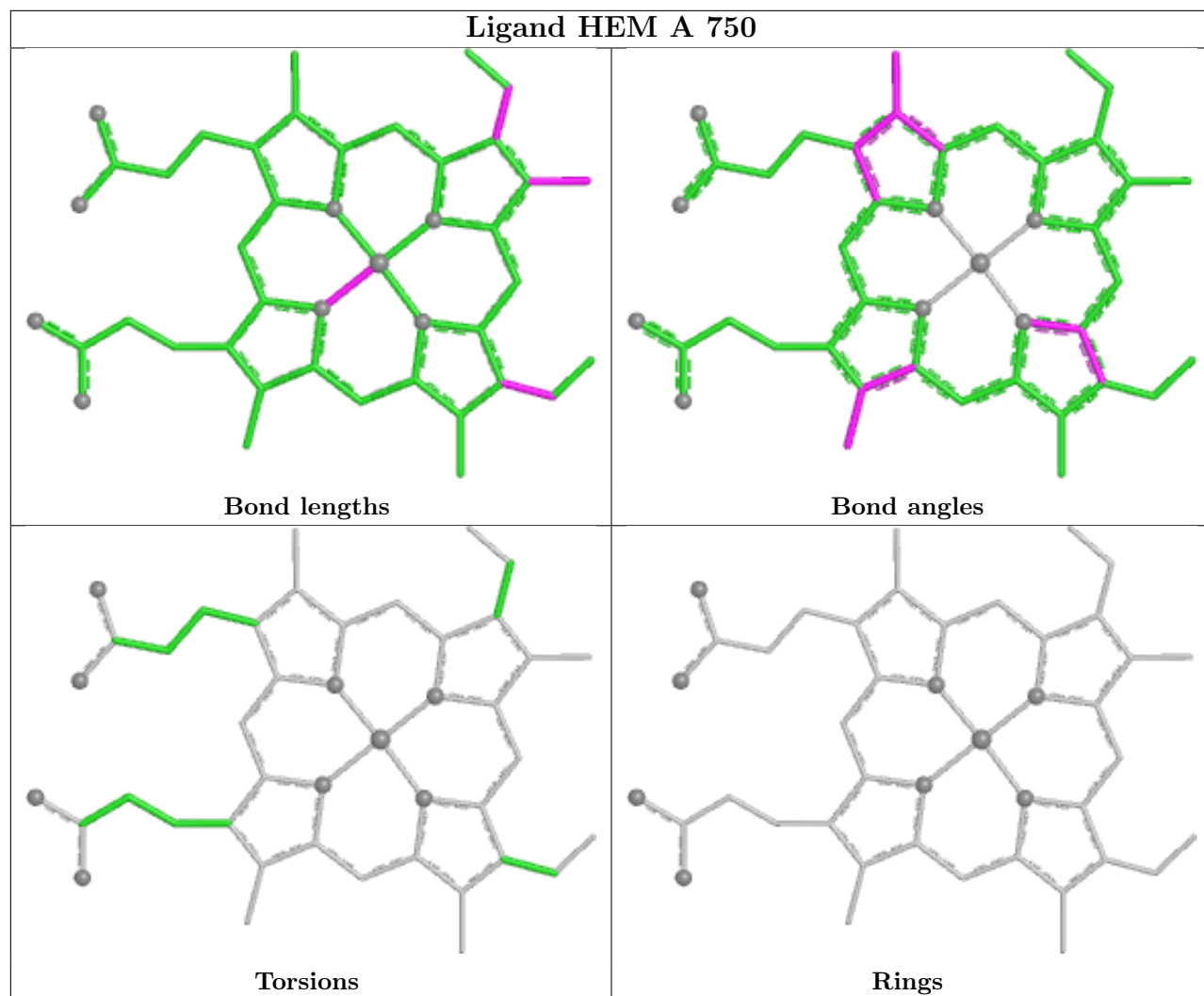
There are no ring outliers.

4 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	750	HEM	2	0
5	B	2760	H4B	1	0
5	A	1760	H4B	1	0
4	A	750	HEM	1	0

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





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	407/419 (97%)	0.88	47 (11%) 9 7	24, 45, 80, 95	0
1	B	410/419 (97%)	0.45	13 (3%) 50 47	21, 39, 66, 84	0
All	All	817/838 (97%)	0.67	60 (7%) 21 18	21, 42, 76, 95	0

All (60) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	322	LEU	4.8
1	A	350	THR	4.6
1	A	715	VAL	4.4
1	B	322	LEU	4.0
1	B	350	THR	4.0
1	A	488	PRO	3.5
1	B	321	THR	3.2
1	A	355	PHE	3.2
1	B	392	SER	3.2
1	A	372	PHE	3.1
1	A	533	LEU	3.1
1	A	528	GLY	3.0
1	B	352	ASP	2.9
1	A	508	GLN	2.8
1	A	300	PHE	2.8
1	A	716	TRP	2.8
1	A	370	LYS	2.7
1	A	557	LEU	2.7
1	B	667	ARG	2.7
1	A	554	PHE	2.7
1	A	617	ASP	2.7
1	A	616	LEU	2.7
1	A	551	PHE	2.6
1	A	338	PRO	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	363	ASP	2.5
1	A	385	ASN	2.5
1	A	490	GLY	2.5
1	A	553	TRP	2.5
1	A	487	GLN	2.5
1	A	550	LYS	2.5
1	A	711	TRP	2.4
1	A	535	GLN	2.4
1	A	318	LEU	2.4
1	B	300	PHE	2.3
1	A	710	PRO	2.3
1	A	337	LEU	2.2
1	A	531	PRO	2.2
1	A	470	HIS	2.2
1	A	618	MET	2.2
1	B	338	PRO	2.2
1	A	382	GLU	2.2
1	A	351	LYS	2.2
1	B	351	LYS	2.2
1	B	337	LEU	2.1
1	A	552	ASP	2.1
1	A	671	GLY	2.1
1	A	392	SER	2.1
1	B	620	LYS	2.1
1	A	373	GLY	2.1
1	A	713	THR	2.1
1	A	388	ILE	2.1
1	A	504	ILE	2.1
1	A	513	PRO	2.1
1	A	549	PRO	2.1
1	A	510	TRP	2.1
1	B	489	ASP	2.0
1	B	348	VAL	2.0
1	A	321	THR	2.0
1	A	467	ASP	2.0
1	A	471	ASP	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](#)

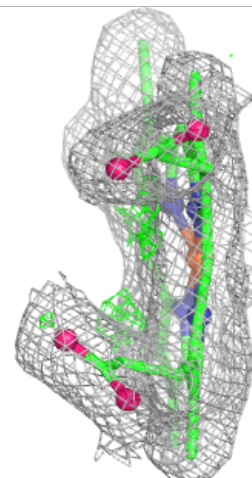
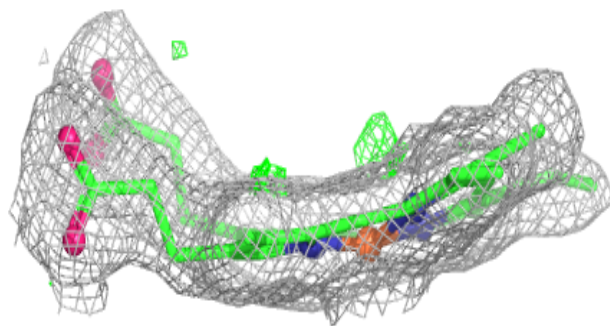
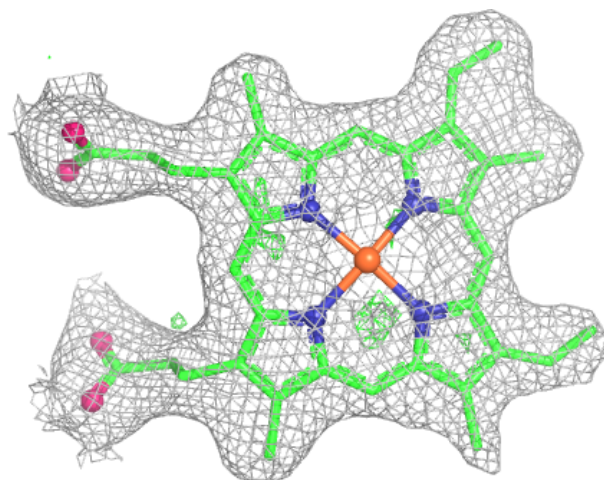
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
5	H4B	A	1760	17/17	0.92	0.09	22,27,33,36	0
2	ACT	B	2860	4/4	0.93	0.09	29,32,33,39	0
5	H4B	B	2760	17/17	0.93	0.08	22,25,31,37	0
2	ACT	A	1860	4/4	0.94	0.12	41,44,44,45	0
6	PTU	B	2810	12/12	0.95	0.09	23,27,29,30	0
6	PTU	A	1810	12/12	0.96	0.08	23,26,29,29	0
4	HEM	A	750	43/43	0.97	0.07	21,29,34,37	0
4	HEM	B	750	43/43	0.97	0.07	16,25,29,33	0
3	ZN	A	900	1/1	0.99	0.03	38,38,38,38	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.

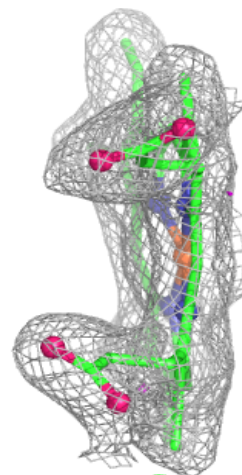
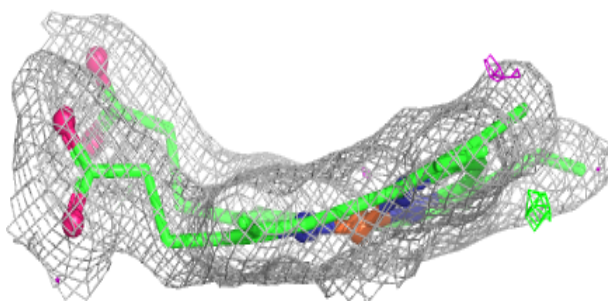
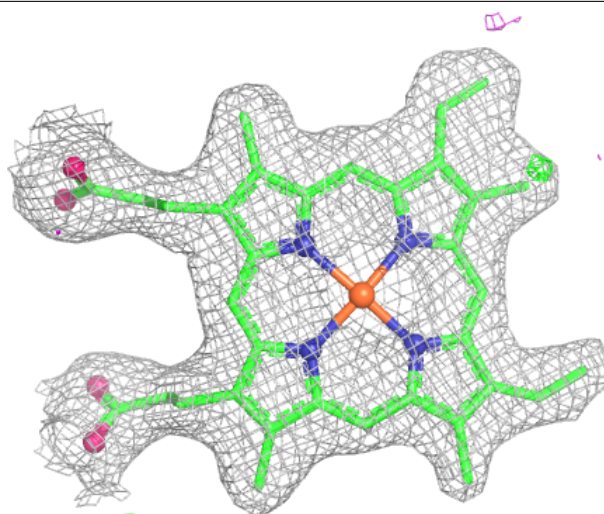
Electron density around HEM A 750:

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



Electron density around HEM B 750:

$2mF_o-DF_c$ (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.