



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 10:10 AM UTC

PDB ID : 2G6J / pdb_00002g6j
Title : Structure of rat nNOS (L337N) heme domain (4-aminobiopterin bound) complexed with NO
Authors : Li, H.; Igarashi, J.; Jamal, J.; Yang, W.; Poulos, T.L.
Deposited on : 2006-02-24
Resolution : 2.30 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

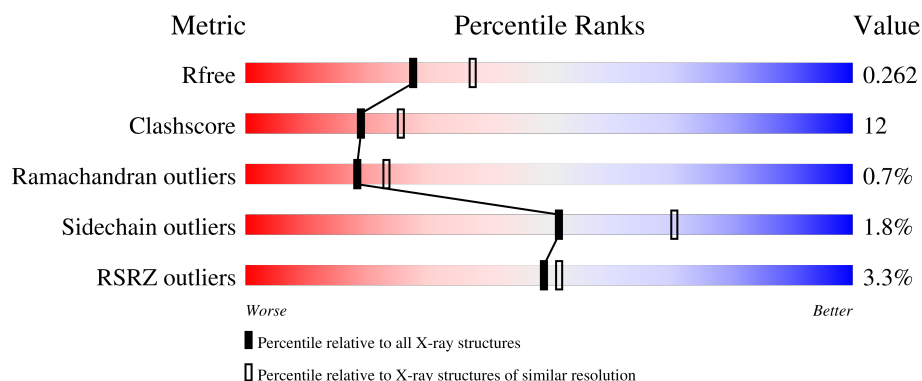
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

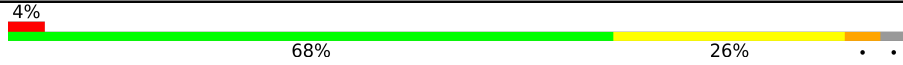
The reported resolution of this entry is 2.30 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	420	
1	B	420	

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 7175 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.

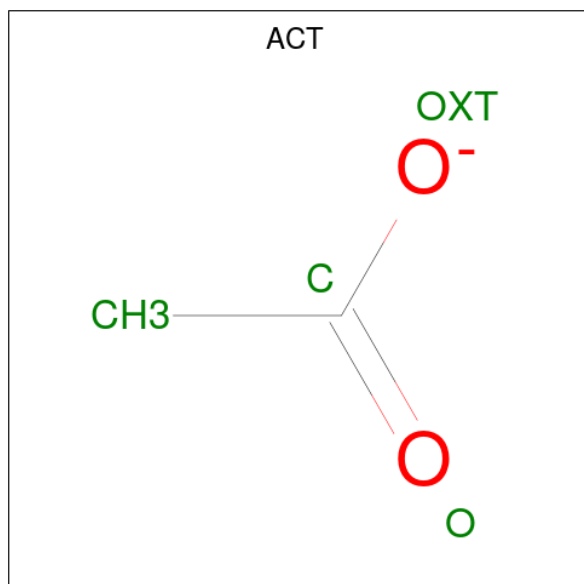
- Molecule 1 is a protein called Nitric-oxide synthase, brain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	409	Total	C	N	O	S	0	0	0
			3331	2130	572	608	21			
1	B	411	Total	C	N	O	S	0	0	0
			3345	2138	575	611	21			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	337	ASN	LEU	engineered mutation	UNP P29476
B	337	ASN	LEU	engineered mutation	UNP P29476

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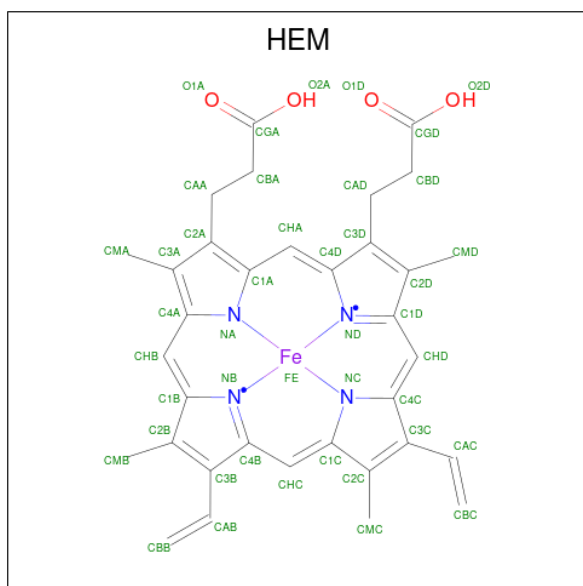
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	C	O	0	0
			4	2	2		

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

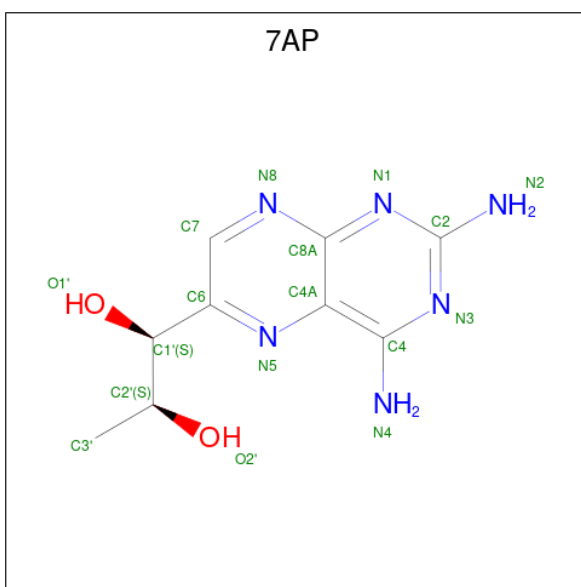
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Zn	0	0
			1	1		

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



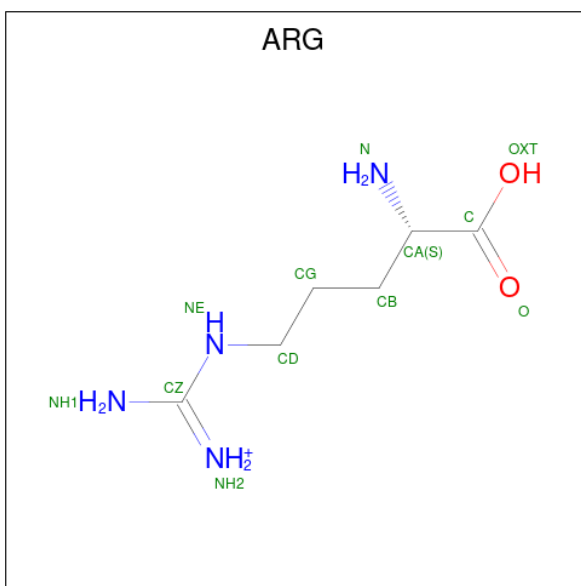
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
4	B	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- Molecule 5 is (1S,2S)-1-(2,4-DIAMINOPTERIDIN-6-YL)PROPANE-1,2-DIOL (CCD ID: 7AP) (formula: $C_9H_{12}N_6O_2$).



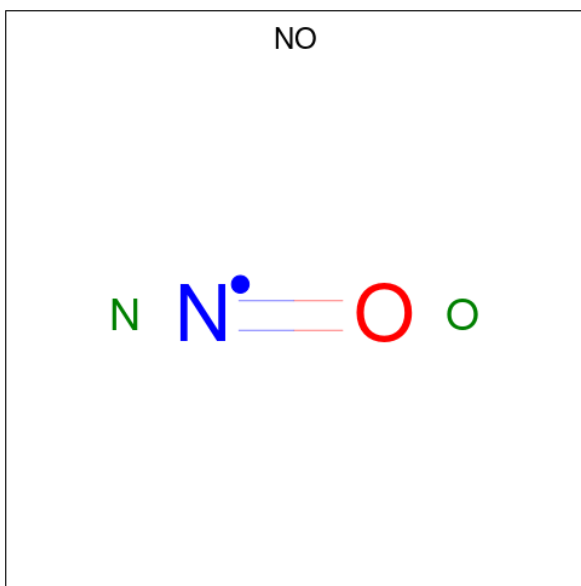
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			17	9	6	2		
5	B	1	Total	C	N	O	0	0
			17	9	6	2		

- Molecule 6 is ARGinine (CCD ID: ARG) (formula: $C_6H_{15}N_4O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	N	O	0	0
			12	6	4	2		
6	B	1	Total	C	N	O	0	0
			12	6	4	2		

- Molecule 7 is NITRIC OXIDE (CCD ID: NO) (formula: NO).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	N	O	0	0
			2	1	1		
7	B	1	Total	N	O	0	0
			2	1	1		

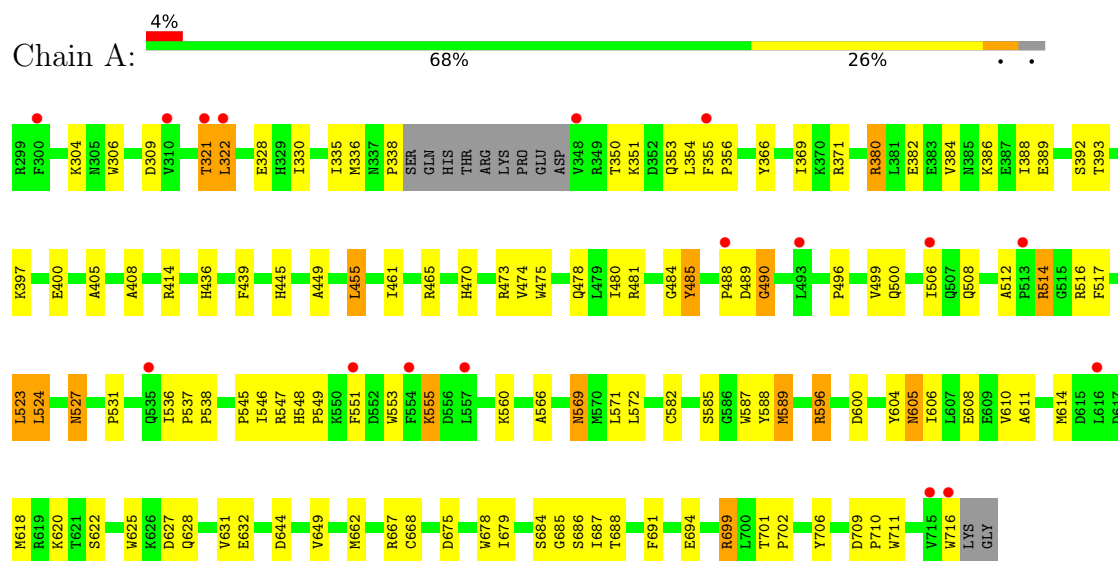
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	153	Total	O	0	0
			153	153		
8	B	189	Total	O	0	0
			189	189		

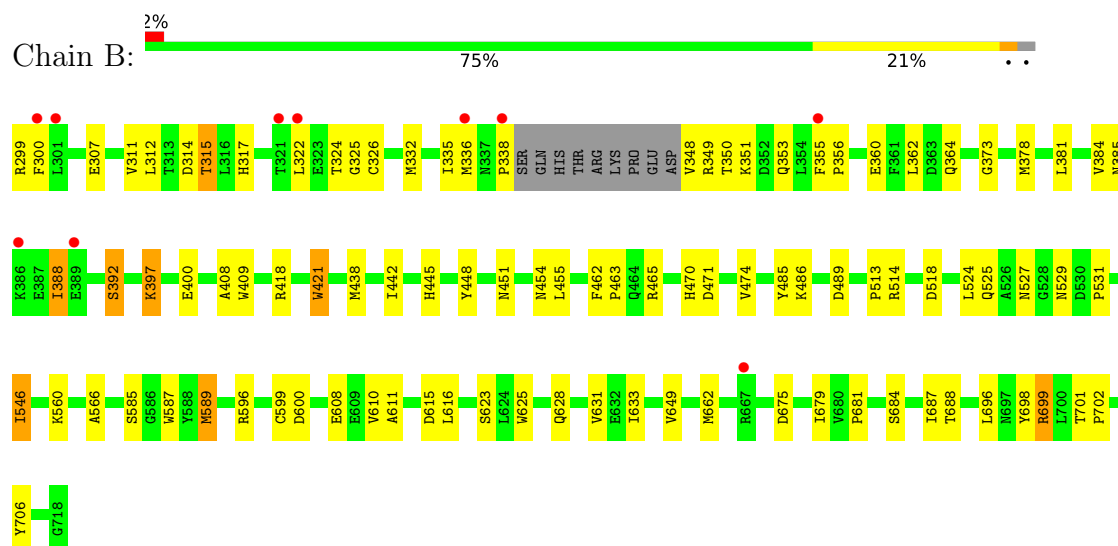
3 Residue-property plots [i](#)

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, brain



- Molecule 1: Nitric-oxide synthase, brain



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	51.52Å 110.49Å 164.91Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.22 – 2.30 49.22 – 2.30	Depositor EDS
% Data completeness (in resolution range)	98.7 (49.22-2.30) 98.8 (49.22-2.30)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.67 (at 2.29Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.213 , 0.269 0.209 , 0.262	Depositor DCC
R_{free} test set	2103 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	36.4	Xtriage
Anisotropy	0.951	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 49.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7175	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.82% 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: ZN, 7AP, HEM, NO, ACT

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	0.49	0/3424	1.01	15/4645 (0.3%)
1	B	0.56	0/3438	1.01	20/4661 (0.4%)
All	All	0.52	0/6862	1.01	35/9306 (0.4%)

There are no bond length outliers.

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	589	MET	N-CA-C	-8.77	94.94	109.24
1	B	589	MET	N-CA-C	-8.72	95.03	109.07
1	B	326	CYS	CA-CB-SG	7.46	131.56	114.40
1	A	688	THR	N-CA-C	-7.05	100.32	110.40
1	A	546	ILE	N-CA-C	6.79	117.90	107.99
1	B	471	ASP	N-CA-C	6.67	118.91	110.24
1	B	408	ALA	N-CA-C	-6.62	104.07	111.28
1	B	688	THR	N-CA-C	-6.61	100.16	110.14
1	A	596	ARG	N-CA-C	6.57	119.41	111.40
1	A	588	TYR	N-CA-C	6.40	119.52	110.50
1	B	421	TRP	N-CA-C	6.26	118.90	111.33
1	B	455	LEU	N-CA-C	6.17	119.17	110.23
1	A	706	TYR	N-CA-C	-6.13	102.28	110.55
1	A	524	LEU	N-CA-C	6.00	118.73	109.07
1	A	474	VAL	N-CA-C	-5.96	97.81	107.28
1	B	623	SER	N-CA-C	-5.91	106.07	113.28
1	B	315	THR	N-CA-C	-5.91	106.71	114.04
1	B	474	VAL	N-CA-C	-5.89	98.66	107.37
1	A	408	ALA	N-CA-C	-5.86	104.97	111.36
1	B	600	ASP	N-CA-C	-5.78	102.49	110.35
1	B	681	PRO	CA-C-N	5.63	126.88	119.84
1	B	681	PRO	C-N-CA	5.63	126.88	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	418	ARG	N-CA-C	5.59	119.58	112.87
1	B	599	CYS	N-CA-C	5.57	120.95	113.72
1	A	699	ARG	N-CA-C	5.55	117.68	108.20
1	B	706	TYR	N-CA-C	-5.45	103.19	110.55
1	B	546	ILE	N-CA-C	5.44	116.22	107.73
1	A	354	LEU	N-CA-C	5.39	117.23	111.36
1	A	455	LEU	N-CA-C	5.34	117.78	110.24
1	A	709	ASP	N-CA-C	-5.32	101.83	109.48
1	B	397	LYS	N-CA-C	-5.30	102.67	110.52
1	B	699	ARG	N-CA-C	5.30	117.60	109.07
1	A	461	ILE	N-CA-C	5.14	114.98	107.37
1	A	490	GLY	N-CA-C	-5.13	108.97	115.08
1	B	349	ARG	N-CA-C	5.04	117.21	110.35

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	3331	0	3238	101	0
1	B	3345	0	3254	63	0
2	A	4	0	3	1	0
2	B	4	0	3	0	0
3	A	1	0	0	0	0
4	A	43	0	30	3	0
4	B	43	0	30	1	0
5	A	17	0	12	1	0
5	B	17	0	12	1	0
6	A	12	0	12	0	0
6	B	12	0	12	0	0
7	A	2	0	0	0	0
7	B	2	0	0	0	0
8	A	153	0	0	8	0
8	B	189	0	0	10	0
All	All	7175	0	6606	159	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 12.

All (159) 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	1.37	1.03
1:A:405:ALA:HA	8:A:1034:HOH:O	1.61	1.01
1:A:569:ASN:HD22	1:A:569:ASN:H	1.30	0.80
1:B:322:LEU:HB3	1:B:699:ARG:HH21	1.49	0.76
1:A:470:HIS:HB3	1:A:527:ASN:ND2	2.02	0.74
1:B:465:ARG:HB2	8:B:992:HOH:O	1.88	0.72
4:A:750:HEM:HMC2	4:A:750:HEM:HBC2	1.73	0.70
1:A:572:LEU:HD11	8:A:1034:HOH:O	1.91	0.70
1:B:325:GLY:O	1:B:332:MET:HE2	1.92	0.69
1:A:569:ASN:HD22	1:A:569:ASN:N	1.90	0.68
1:A:523:LEU:CD2	1:A:531:PRO:HB2	2.18	0.68
1:A:478:GLN:HB2	1:A:481:ARG:HG3	1.76	0.68
1:A:569:ASN:H	1:A:569:ASN:ND2	1.91	0.67
1:A:473:ARG:NH2	1:A:710:PRO:HD3	2.11	0.65
1:B:587:TRP:H	4:B:750:HEM:HAB	1.61	0.65
1:A:631:VAL:HG11	1:B:628:GLN:HG3	1.78	0.64
1:A:701:THR:HB	8:A:993:HOH:O	1.98	0.64
1:A:350:THR:OG1	1:A:353:GLN:HG3	1.99	0.63
1:B:311:VAL:O	1:B:312:LEU:HD23	1.97	0.63
1:B:356:PRO:O	1:B:360:GLU:HG3	2.00	0.62
1:B:610:VAL:HG21	1:B:633:ILE:HD11	1.82	0.61
1:B:364:GLN:NE2	8:B:1096:HOH:O	2.32	0.61
1:A:514:ARG:HH21	1:A:514:ARG:HG3	1.66	0.61
1:A:620:LYS:HE3	1:A:622:SER:OG	2.02	0.60
1:B:438:MET:O	1:B:442:ILE:HG13	2.00	0.60
1:B:332:MET:CE	1:B:338:PRO:HB3	2.32	0.59
1:B:513:PRO:HG2	1:B:518:ASP:CG	2.27	0.59
1:A:485:TYR:HE2	1:A:512:ALA:HB1	1.67	0.59
1:A:436:HIS:ND1	8:A:1058:HOH:O	2.31	0.59
1:A:675:ASP:O	1:A:679:ILE:HG12	2.03	0.59
1:A:321:THR:HG23	1:A:322:LEU:H	1.67	0.59
1:A:355:PHE:HB2	1:A:356:PRO:HD3	1.84	0.59
1:A:322:LEU:HD12	1:A:699:ARG:HE	1.67	0.58
1:A:606:ILE:O	1:A:610:VAL:HG23	2.03	0.58
1:A:380:ARG:HD3	1:A:400:GLU:OE1	2.03	0.58
1:A:455:LEU:HD12	1:A:587:TRP:HB3	1.86	0.57
1:A:551:PHE:HB3	1:A:553:TRP:CE2	2.39	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:351:LYS:O	1:A:355:PHE:HD1	1.87	0.56
1:A:475:TRP:HB2	1:A:523:LEU:HB3	1.86	0.56
1:A:632:GLU:OE2	1:B:628:GLN:NE2	2.38	0.56
1:A:485:TYR:CE2	1:A:512:ALA:HB1	2.41	0.56
1:B:373:GLY:HA2	1:B:378:MET:HE3	1.86	0.56
1:B:388:ILE:O	1:B:392:SER:N	2.34	0.56
1:A:514:ARG:HH21	1:A:514:ARG:CG	2.19	0.56
1:B:470:HIS:O	1:B:527:ASN:HA	2.06	0.55
1:B:486:LYS:HE2	8:B:985:HOH:O	2.06	0.55
1:A:335:ILE:HB	1:A:338:PRO:HG3	1.88	0.55
1:B:355:PHE:HB2	1:B:356:PRO:HD3	1.88	0.55
1:B:470:HIS:HB3	1:B:527:ASN:ND2	2.22	0.54
1:B:351:LYS:HE3	1:B:392:SER:OG	2.07	0.54
1:A:322:LEU:HB2	1:A:699:ARG:HB2	1.90	0.53
1:A:322:LEU:HD12	1:A:699:ARG:HB3	1.90	0.53
1:A:572:LEU:CD1	8:A:1034:HOH:O	2.50	0.53
1:A:330:ILE:HD11	1:B:696:LEU:HB3	1.89	0.53
1:A:694:GLU:HB3	1:B:335:ILE:HD13	1.90	0.53
1:A:555:LYS:NZ	1:A:555:LYS:HB3	2.25	0.52
1:B:589:MET:HA	1:B:649:VAL:O	2.10	0.52
1:B:701:THR:HA	1:B:702:PRO:C	2.34	0.52
1:B:325:GLY:C	1:B:332:MET:HE2	2.35	0.52
1:A:628:GLN:HG2	1:B:631:VAL:HG11	1.92	0.51
1:A:547:ARG:NH1	1:A:644:ASP:OD1	2.39	0.51
1:A:610:VAL:O	1:A:614:MET:HG3	2.10	0.51
1:A:667:ARG:NH1	1:A:668:CYS:SG	2.84	0.51
1:B:362:LEU:HD11	1:B:384:VAL:HG21	1.93	0.51
1:B:675:ASP:O	1:B:679:ILE:HG12	2.10	0.50
1:A:489:ASP:OD2	1:A:489:ASP:O	2.28	0.50
1:A:551:PHE:HB3	1:A:553:TRP:NE1	2.26	0.50
1:B:566:ALA:HB2	1:B:585:SER:HB3	1.93	0.50
1:B:299:ARG:HB3	1:B:299:ARG:NH1	2.27	0.50
1:B:445:HIS:C	1:B:445:HIS:CD2	2.90	0.49
1:A:524:LEU:O	1:A:531:PRO:HA	2.11	0.49
1:A:516:ARG:HD3	1:A:517:PHE:CE1	2.46	0.49
1:A:614:MET:CE	1:A:632:GLU:HG3	2.43	0.49
1:B:451:ASN:HB3	1:B:454:ASN:O	2.12	0.49
1:B:596:ARG:HH12	5:B:761:7AP:C4	2.25	0.49
1:A:304:LYS:HE3	1:A:309:ASP:OD1	2.13	0.49
1:B:314:ASP:OD1	1:B:317:HIS:ND1	2.41	0.49
1:B:397:LYS:HB2	1:B:400:GLU:HG3	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:351:LYS:HE2	1:A:392:SER:HB3	1.94	0.49
1:B:465:ARG:NE	8:B:992:HOH:O	2.45	0.49
1:A:384:VAL:O	1:A:388:ILE:HG13	2.13	0.48
1:A:716:TRP:H	1:A:716:TRP:CD1	2.30	0.48
2:A:860:ACT:H2	4:A:750:HEM:HAB	1.95	0.48
1:A:627:ASP:O	1:A:631:VAL:HG23	2.13	0.48
1:A:455:LEU:HD12	1:A:587:TRP:CB	2.44	0.48
1:A:523:LEU:HD22	1:A:531:PRO:CB	2.26	0.47
4:A:750:HEM:HBC2	4:A:750:HEM:CMC	2.44	0.47
1:A:380:ARG:NH1	1:A:397:LYS:HG2	2.30	0.47
1:A:536:ILE:O	1:A:537:PRO:C	2.55	0.47
1:B:616:LEU:HD22	8:B:1089:HOH:O	2.15	0.47
1:B:485:TYR:CZ	1:B:514:ARG:HA	2.49	0.47
1:B:409:TRP:CE3	1:B:421:TRP:HA	2.50	0.46
1:A:351:LYS:O	1:A:355:PHE:CD1	2.68	0.46
1:A:366:TYR:HA	1:A:369:ILE:HG12	1.98	0.46
1:A:445:HIS:C	1:A:445:HIS:CD2	2.93	0.46
1:A:596:ARG:O	1:A:600:ASP:HB2	2.15	0.46
1:A:684:SER:HB3	1:A:687:ILE:HD11	1.97	0.46
1:A:449:ALA:O	1:A:455:LEU:HA	2.16	0.46
1:A:545:PRO:HG2	1:A:547:ARG:NH2	2.31	0.46
1:A:484:GLY:O	1:A:499:VAL:HA	2.16	0.46
1:A:571:LEU:C	1:A:571:LEU:HD23	2.41	0.46
1:B:546:ILE:HG12	1:B:560:LYS:HA	1.97	0.46
1:B:462:PHE:HB3	1:B:463:PRO:HD2	1.99	0.45
1:A:618:MET:HE3	1:A:625:TRP:CZ3	2.52	0.45
1:B:489:ASP:OD2	1:B:489:ASP:C	2.60	0.45
1:A:480:ILE:HA	8:A:950:HOH:O	2.15	0.45
1:A:321:THR:HG21	1:A:701:THR:HG22	1.98	0.45
1:A:614:MET:HE1	1:A:632:GLU:HG3	1.98	0.45
1:A:517:PHE:HB2	1:A:560:LYS:CE	2.47	0.45
1:A:382:GLU:HG3	1:A:386:LYS:HE3	1.98	0.44
1:B:350:THR:HG22	1:B:353:GLN:CD	2.43	0.44
1:B:300:PHE:N	1:B:300:PHE:CD1	2.85	0.44
1:B:317:HIS:HB3	1:B:698:TYR:CE2	2.52	0.44
1:A:306:TRP:CZ2	1:B:336:MET:HG2	2.52	0.44
1:A:701:THR:HA	1:A:702:PRO:C	2.43	0.44
1:A:353:GLN:O	1:A:356:PRO:HD2	2.18	0.44
1:A:488:PRO:C	1:A:490:GLY:N	2.76	0.44
1:B:699:ARG:HA	8:B:994:HOH:O	2.17	0.44
1:A:514:ARG:CG	1:A:514:ARG:NH2	2.78	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:548:HIS:ND1	1:A:549:PRO:HD2	2.33	0.43
1:B:662:MET:HE3	8:B:922:HOH:O	2.17	0.43
1:A:439:PHE:HZ	1:A:537:PRO:HD3	1.82	0.43
1:A:500:GLN:NE2	1:A:711:TRP:NE1	2.66	0.43
1:B:524:LEU:O	1:B:531:PRO:HA	2.18	0.43
1:A:551:PHE:HE1	1:A:614:MET:HE3	1.84	0.43
1:A:589:MET:HA	1:A:649:VAL:O	2.18	0.43
1:B:307:GLU:HG3	8:B:964:HOH:O	2.16	0.43
1:A:465:ARG:NE	8:A:1015:HOH:O	2.51	0.43
1:B:348:VAL:N	8:B:1063:HOH:O	2.51	0.43
1:A:508:GLN:OE1	1:A:716:TRP:CZ3	2.70	0.43
1:A:322:LEU:CD1	1:A:699:ARG:HE	2.31	0.43
1:A:414:ARG:HD3	1:A:678:TRP:CD2	2.54	0.43
1:B:332:MET:HE3	1:B:338:PRO:HB3	2.00	0.42
1:B:381:LEU:O	1:B:385:ASN:ND2	2.52	0.42
1:A:701:THR:OG1	1:A:702:PRO:HA	2.20	0.42
1:B:325:GLY:CA	1:B:332:MET:HE2	2.50	0.42
1:A:328:GLU:O	1:B:324:THR:HG23	2.19	0.42
1:A:608:GLU:O	1:A:611:ALA:HB3	2.20	0.42
1:A:566:ALA:HB2	1:A:585:SER:HB3	2.01	0.42
1:B:525:GLN:HG3	1:B:529:ASN:O	2.19	0.42
1:A:517:PHE:HB2	1:A:560:LYS:HE2	2.01	0.41
1:A:684:SER:HB3	1:A:687:ILE:CG1	2.50	0.41
1:B:485:TYR:CE1	1:B:514:ARG:HA	2.55	0.41
1:A:604:TYR:O	1:A:605:ASN:C	2.63	0.41
1:A:618:MET:HE3	1:A:625:TRP:CH2	2.54	0.41
1:B:625:TRP:HA	8:B:1089:HOH:O	2.18	0.41
1:A:582:CYS:O	1:A:582:CYS:SG	2.78	0.41
1:A:631:VAL:HG11	1:B:628:GLN:CG	2.50	0.41
1:B:566:ALA:CB	1:B:585:SER:HB3	2.50	0.41
1:B:448:TYR:CD2	1:B:448:TYR:C	2.98	0.41
1:A:371:ARG:NH2	1:A:371:ARG:HG2	2.34	0.41
1:A:596:ARG:HE	1:A:600:ASP:CG	2.29	0.41
1:A:336:MET:SD	5:A:760:7AP:H3'2	2.60	0.41
1:A:380:ARG:NH2	8:A:1005:HOH:O	2.48	0.41
1:B:608:GLU:O	1:B:611:ALA:HB3	2.21	0.41
1:A:686:SER:HA	1:A:691:PHE:CG	2.57	0.40
1:A:496:PRO:HA	1:A:499:VAL:HG23	2.02	0.40
1:A:506:ILE:C	1:A:508:GLN:H	2.29	0.40
1:B:684:SER:HB3	1:B:687:ILE:HD11	2.03	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	405/420 (96%)	379 (94%)	22 (5%)	4 (1%)	12	15
1	B	407/420 (97%)	381 (94%)	24 (6%)	2 (0%)	24	31
All	All	812/840 (97%)	760 (94%)	46 (6%)	6 (1%)	18	23

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	322	LEU
1	A	605	ASN
1	A	393	THR
1	B	388	ILE
1	B	615	ASP
1	A	685	GLY

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	365/375 (97%)	354 (97%)	11 (3%)	36	53
1	B	366/375 (98%)	364 (100%)	2 (0%)	81	90
All	All	731/750 (98%)	718 (98%)	13 (2%)	51	70

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	321	THR
1	A	380	ARG
1	A	389	GLU
1	A	485	TYR
1	A	514	ARG
1	A	523	LEU
1	A	527	ASN
1	A	538	PRO
1	A	555	LYS
1	A	569	ASN
1	A	662	MET
1	B	315	THR
1	B	392	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	385	ASN
1	A	420	GLN
1	A	454	ASN
1	A	500	GLN
1	A	507	GLN
1	A	508	GLN
1	A	527	ASN
1	A	569	ASN
1	A	605	ASN
1	A	642	GLN
1	A	697	ASN
1	B	337	ASN
1	B	364	GLN
1	B	407	HIS
1	B	425	GLN
1	B	454	ASN
1	B	507	GLN
1	B	527	ASN
1	B	535	GLN
1	B	601	ASN
1	B	605	ASN
1	B	642	GLN
1	B	697	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 11 ligands modelled in this entry, 1 is monoatomic - leaving 10 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
5	7AP	B	761	-	17,18,18	3.27	5 (29%)	16,26,26	3.34	8 (50%)
7	NO	A	910	4	0,1,1	-	-	-		
7	NO	B	910	4	0,1,1	-	-	-		
4	HEM	B	750	1,7	50,50,50	1.17	4 (8%)	67,82,82	1.13	5 (7%)
6	ARG	A	770	-	10,11,11	0.90	0	9,13,13	0.67	0
2	ACT	B	861	-	3,3,3	0.96	0	3,3,3	0.79	0
6	ARG	B	771	-	10,11,11	0.81	0	9,13,13	0.63	0
4	HEM	A	750	1,7	50,50,50	1.40	11 (22%)	67,82,82	1.15	3 (4%)
2	ACT	A	860	-	3,3,3	0.91	0	3,3,3	0.82	0
5	7AP	A	760	-	17,18,18	3.13	5 (29%)	16,26,26	3.40	9 (56%)

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	7AP	B	761	-	-	1/8/8/8	0/2/2/2
4	HEM	B	750	1,7	-	4/14/54/54	-
6	ARG	A	770	-	-	0/11/11/11	-
6	ARG	B	771	-	-	0/11/11/11	-
4	HEM	A	750	1,7	-	5/14/54/54	-
5	7AP	A	760	-	-	3/8/8/8	0/2/2/2

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	761	7AP	C6-N5	10.47	1.41	1.32
5	A	760	7AP	C6-N5	9.72	1.41	1.32
5	B	761	7AP	C7-N8	5.13	1.40	1.31
5	A	760	7AP	C7-N8	5.09	1.40	1.31
5	A	760	7AP	C2'-C1'	3.66	1.57	1.53
5	B	761	7AP	C4-N3	3.62	1.40	1.33
4	A	750	HEM	FE-NB	3.30	2.05	1.94
5	B	761	7AP	C2'-C1'	3.28	1.57	1.53
5	A	760	7AP	C4-N3	3.19	1.39	1.33
4	A	750	HEM	CHC-C1C	2.88	1.44	1.38
5	B	761	7AP	C2-N3	2.68	1.40	1.35
4	B	750	HEM	CAC-C3C	-2.63	1.40	1.47
5	A	760	7AP	C2-N3	2.60	1.39	1.35
4	A	750	HEM	CMC-C2C	2.57	1.56	1.50
4	B	750	HEM	CAB-C3B	-2.55	1.40	1.47
4	A	750	HEM	CAC-C3C	-2.47	1.40	1.47
4	A	750	HEM	CHD-C4C	2.45	1.43	1.38
4	A	750	HEM	CAB-C3B	-2.29	1.41	1.47
4	B	750	HEM	C2A-C3A	-2.26	1.33	1.38
4	A	750	HEM	CMB-C2B	2.24	1.55	1.50
4	B	750	HEM	CMD-C2D	2.20	1.55	1.50
4	A	750	HEM	C3C-C4C	-2.20	1.42	1.46
4	A	750	HEM	FE-ND	2.03	2.01	1.94
4	A	750	HEM	CHC-C4B	2.03	1.43	1.39
4	A	750	HEM	C3C-C2C	2.03	1.41	1.37

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	760	7AP	N1-C2-N3	-8.16	116.83	127.21
5	B	761	7AP	N1-C2-N3	-8.05	116.97	127.21
5	A	760	7AP	C2-N1-C8A	6.95	122.97	115.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	761	7AP	C2-N1-C8A	6.87	122.89	115.48
5	A	760	7AP	N8-C8A-N1	4.92	121.13	115.77
5	B	761	7AP	N8-C8A-N1	4.65	120.84	115.77
4	A	750	HEM	CBA-CAA-C2A	-4.44	100.25	112.53
4	B	750	HEM	CBA-CAA-C2A	-4.16	101.03	112.53
5	A	760	7AP	N2-C2-N1	3.21	122.81	117.79
5	B	761	7AP	N2-C2-N1	3.10	122.64	117.79
5	B	761	7AP	C7-N8-C8A	2.96	121.33	117.20
5	A	760	7AP	C7-N8-C8A	2.95	121.31	117.20
4	A	750	HEM	CAB-C3B-C2B	-2.69	119.67	128.43
4	A	750	HEM	CHC-C4B-NB	-2.65	121.57	124.42
4	B	750	HEM	CBD-CAD-C3D	-2.42	105.83	112.53
5	B	761	7AP	C4A-C4-N3	-2.42	117.98	120.84
4	B	750	HEM	C2D-C1D-ND	2.40	112.67	109.90
5	A	760	7AP	C2-N3-C4	2.38	123.51	116.72
5	A	760	7AP	C4A-C4-N3	-2.35	118.06	120.84
5	B	761	7AP	C2-N3-C4	2.33	123.36	116.72
4	B	750	HEM	C3B-C4B-NB	2.23	111.07	109.47
5	B	761	7AP	N2-C2-N3	2.11	120.39	117.22
4	B	750	HEM	CHC-C4B-NB	-2.10	122.16	124.42
5	A	760	7AP	N2-C2-N3	2.09	120.36	117.22
5	A	760	7AP	C4A-C8A-N1	-2.07	118.54	121.74

There are no chirality outliers.

All (13) torsion outliers are listed below:

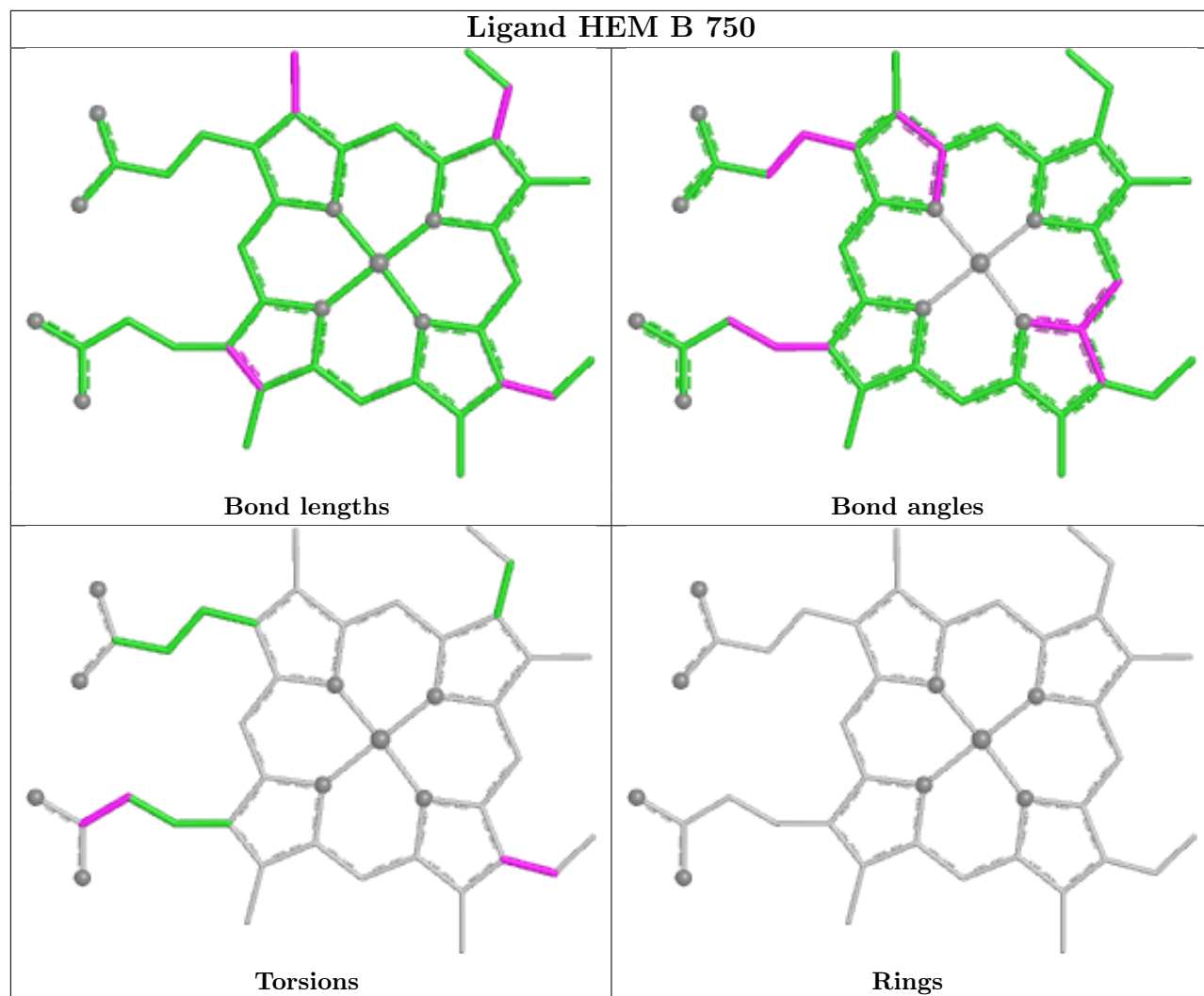
Mol	Chain	Res	Type	Atoms
4	B	750	HEM	C2B-C3B-CAB-CBB
5	A	760	7AP	O1'-C1'-C6-N5
5	A	760	7AP	O1'-C1'-C6-C7
5	B	761	7AP	O1'-C1'-C6-C7
4	B	750	HEM	C4B-C3B-CAB-CBB
4	B	750	HEM	CAA-CBA-CGA-O2A
4	A	750	HEM	CAA-CBA-CGA-O2A
4	A	750	HEM	C4B-C3B-CAB-CBB
5	A	760	7AP	O1'-C1'-C2'-O2'
4	B	750	HEM	CAA-CBA-CGA-O1A
4	A	750	HEM	CAA-CBA-CGA-O1A
4	A	750	HEM	CAD-CBD-CGD-O2D
4	A	750	HEM	CAD-CBD-CGD-O1D

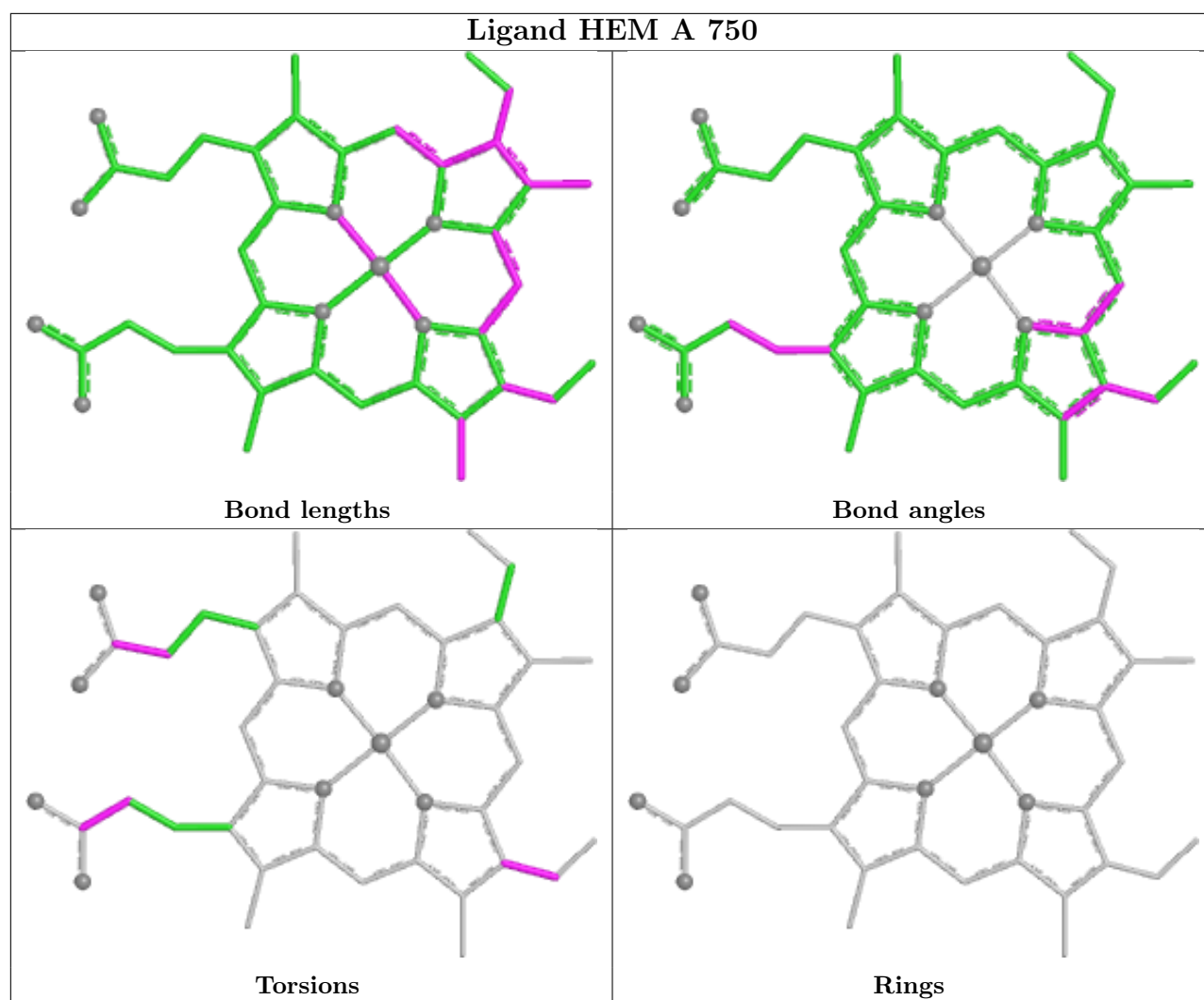
There are no ring outliers.

5 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	761	7AP	1	0
4	B	750	HEM	1	0
4	A	750	HEM	3	0
2	A	860	ACT	1	0
5	A	760	7AP	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	409/420 (97%)	0.44	17 (4%) 40 42	26, 49, 72, 87	0
1	B	411/420 (97%)	0.08	10 (2%) 59 62	24, 41, 66, 76	0
All	All	820/840 (97%)	0.26	27 (3%) 49 51	24, 44, 70, 87	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	300	PHE	3.9
1	A	322	LEU	3.3
1	A	554	PHE	3.1
1	A	715	VAL	3.0
1	A	716	TRP	3.0
1	B	322	LEU	3.0
1	A	488	PRO	3.0
1	A	557	LEU	3.0
1	A	321	THR	2.7
1	B	355	PHE	2.5
1	B	300	PHE	2.5
1	A	348	VAL	2.5
1	A	535	GLN	2.5
1	A	513	PRO	2.5
1	A	506	ILE	2.3
1	B	336	MET	2.3
1	A	310	VAL	2.3
1	A	551	PHE	2.2
1	B	321	THR	2.2
1	A	616	LEU	2.2
1	A	355	PHE	2.2
1	B	389	GLU	2.2
1	B	338	PRO	2.1
1	B	301	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	667	ARG	2.1
1	A	493	LEU	2.0
1	B	386	LYS	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

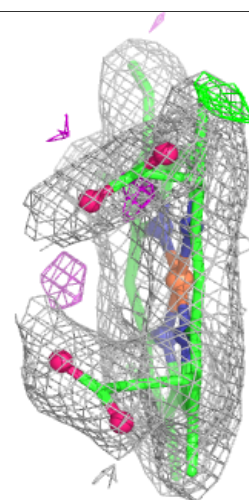
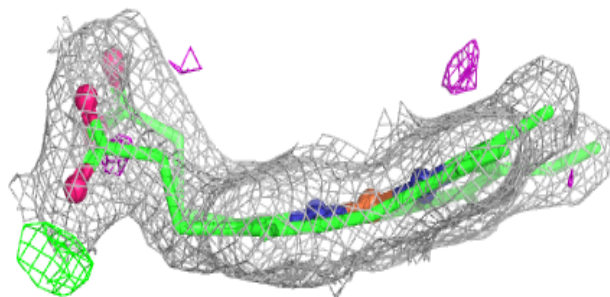
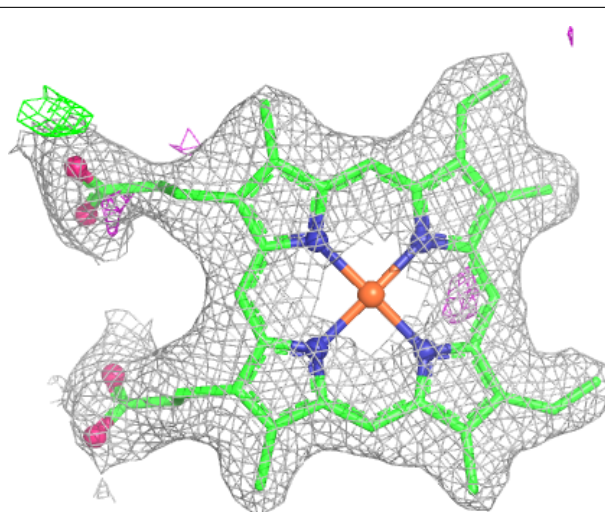
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

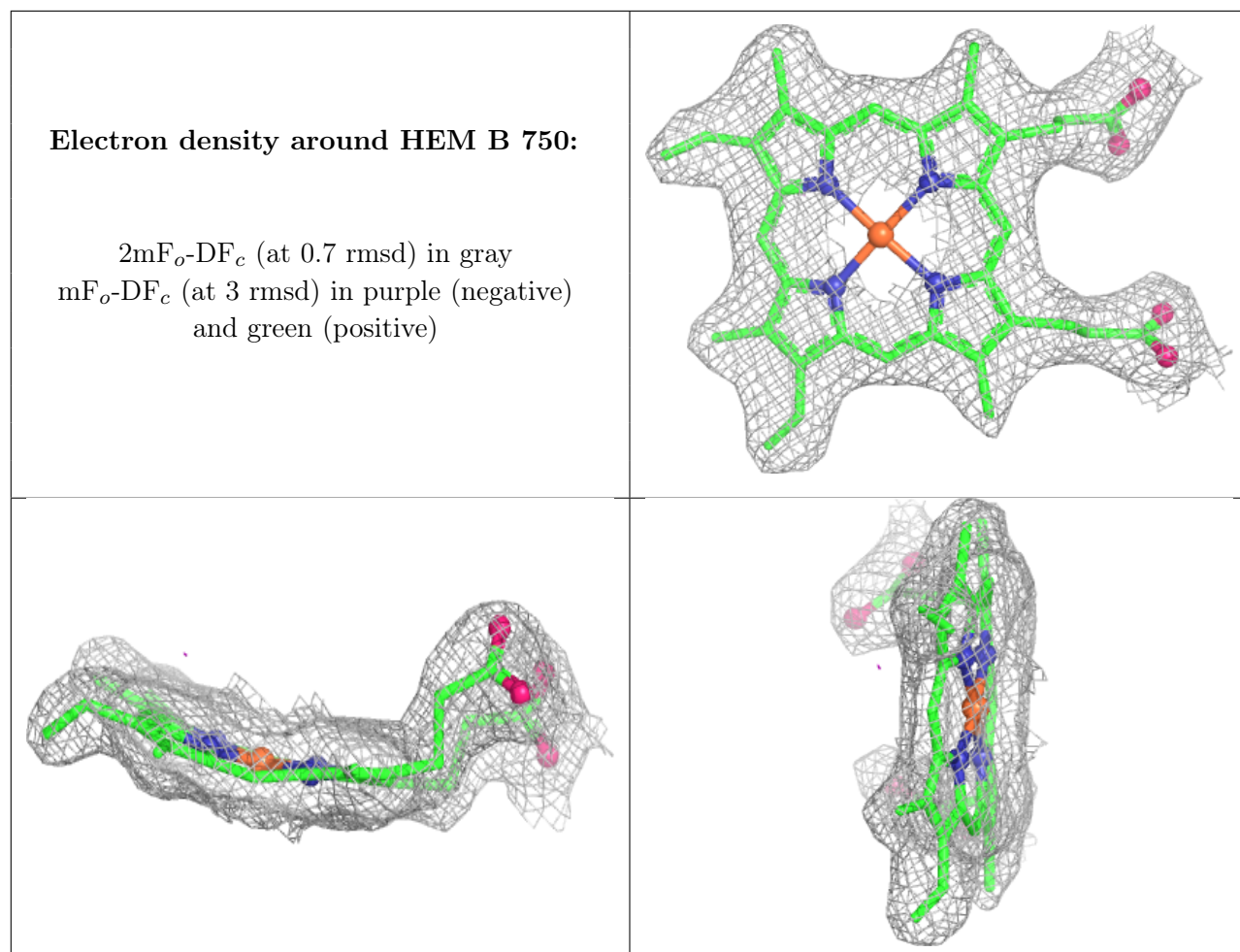
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	7AP	A	760	17/17	0.82	0.15	63,63,64,65	0
5	7AP	B	761	17/17	0.82	0.14	49,54,60,60	0
2	ACT	B	861	4/4	0.90	0.13	58,59,59,61	0
2	ACT	A	860	4/4	0.91	0.14	60,61,62,62	0
6	ARG	B	771	12/12	0.91	0.08	32,38,42,44	0
6	ARG	A	770	12/12	0.93	0.08	42,44,45,46	0
7	NO	B	910	2/2	0.96	0.06	35,35,35,37	0
7	NO	A	910	2/2	0.97	0.06	38,38,38,38	0
4	HEM	A	750	43/43	0.97	0.07	30,34,43,46	0
4	HEM	B	750	43/43	0.98	0.06	25,32,44,47	0
3	ZN	A	900	1/1	1.00	0.03	53,53,53,53	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)





6.5 Other polymers [i](#)

There are no such residues in this entry.