



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 12:11 PM UTC

PDB ID : 3B3P / pdb_00003b3p
Title : Structure of neuronal nos heme domain in complex with a inhibitor (+-)-n1-{
cis-4'-[(6"-amino-4"-methylpyridin-2"-yl)methyl]pyrrolidin-3'-yl}-n2-(4'-chlor
obenzyl)ethane-1,2-diamine
Authors : Igarashi, J.; Li, H.; Poulos, T.L.
Deposited on : 2007-10-22
Resolution : 2.45 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

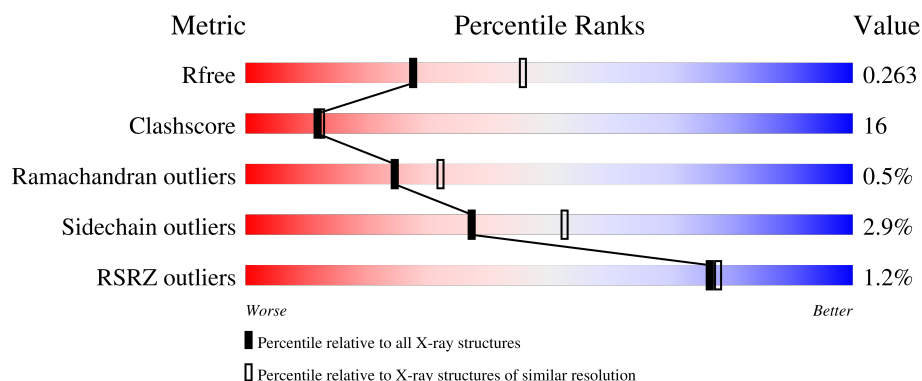
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

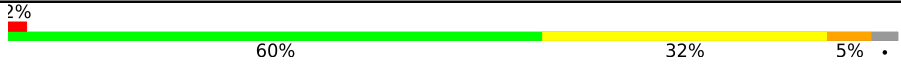
The reported resolution of this entry is 2.45 Å.

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	1190 (2.46-2.46)
Clashscore	190562	1229 (2.46-2.46)
Ramachandran outliers	187476	1218 (2.46-2.46)
Sidechain outliers	187428	1218 (2.46-2.46)
RSRZ outliers	180081	1190 (2.46-2.46)

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	422	
1	B	422	

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
2	ACT	A	860	-	-	X	-

2 Entry composition [i](#)

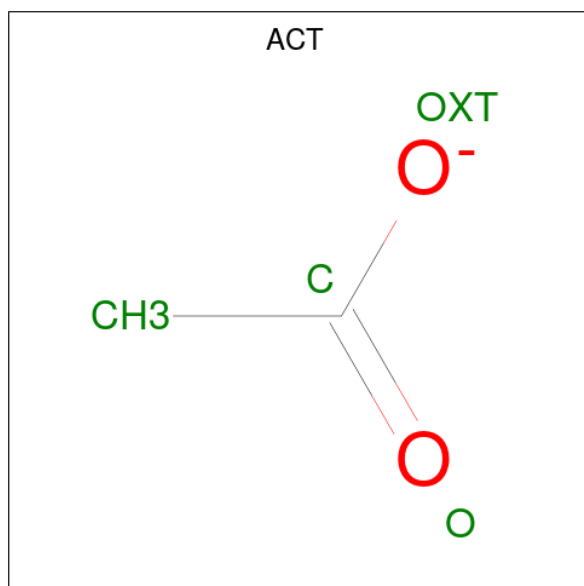
There are 7 unique types of molecules in this entry. The entry contains 7104 atoms, of which 0 are hydrogens and 0 are deuteriums.

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	409	Total	C	N	O	S	0	0	0
			3331	2132	571	607	21			
1	B	411	Total	C	N	O	S	0	0	0
			3345	2140	574	610	21			

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

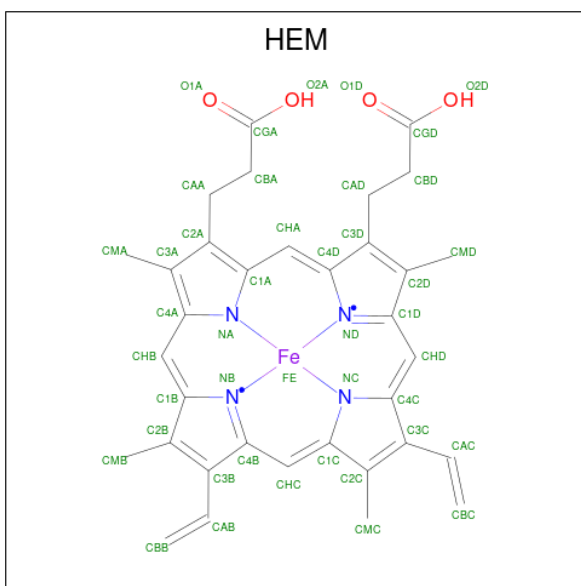


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		

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

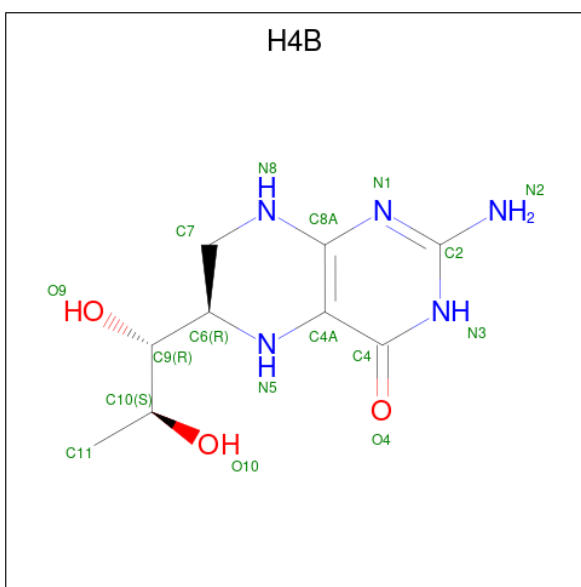
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Zn 1 1	0	0

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



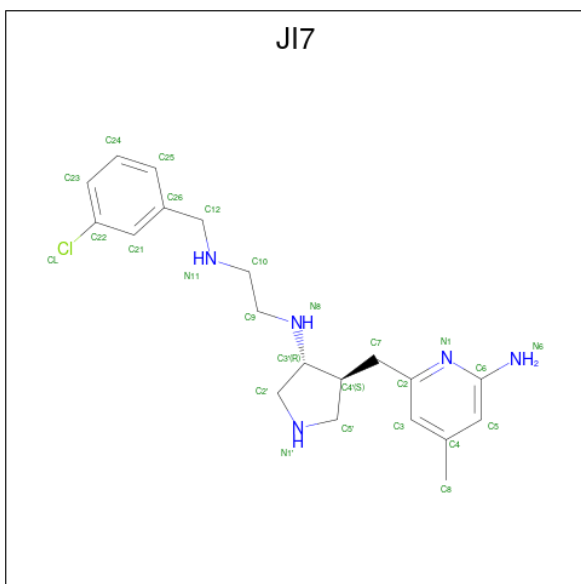
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
4	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- 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 N-[(3R,4S)-4-[(6-amino-4-methylpyridin-2-yl)methyl]pyrrolidin-3-yl]-N'-(3-chlorobenzyl)ethane-1,2-diamine (CCD ID: JI7) (formula: C₂₀H₂₈ClN₅).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	Cl	N	0	0
			26	20	1	5		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	B	1	Total	C	Cl	N	0	0
			26	20	1	5		

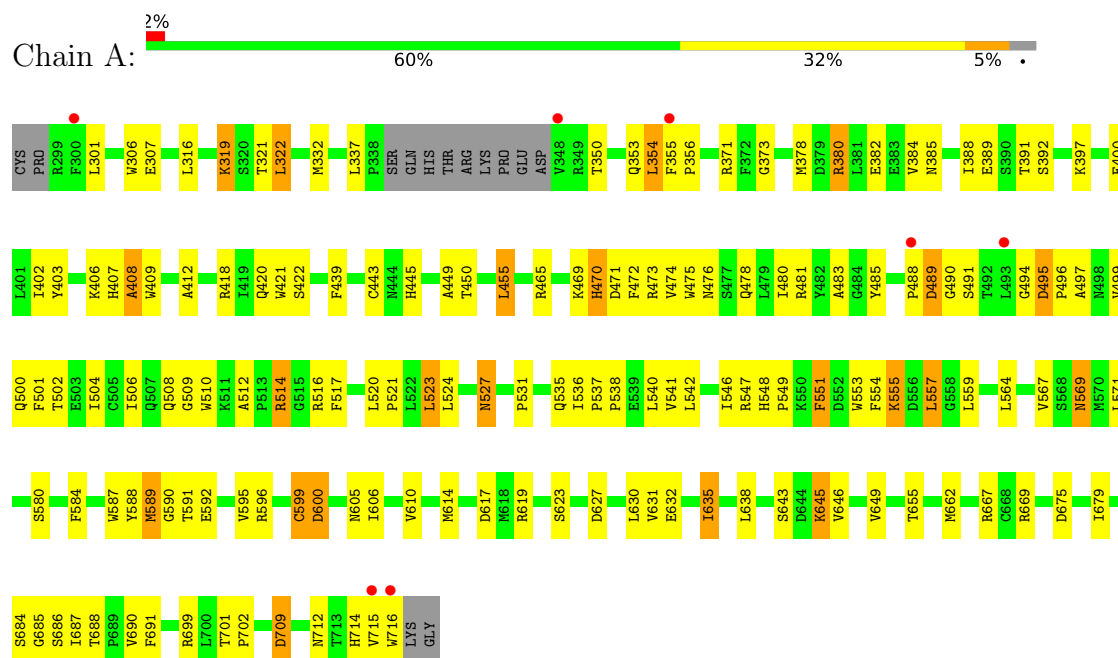
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	113	Total	O	0	0
			113	113		
7	B	134	Total	O	0	0
			134	134		

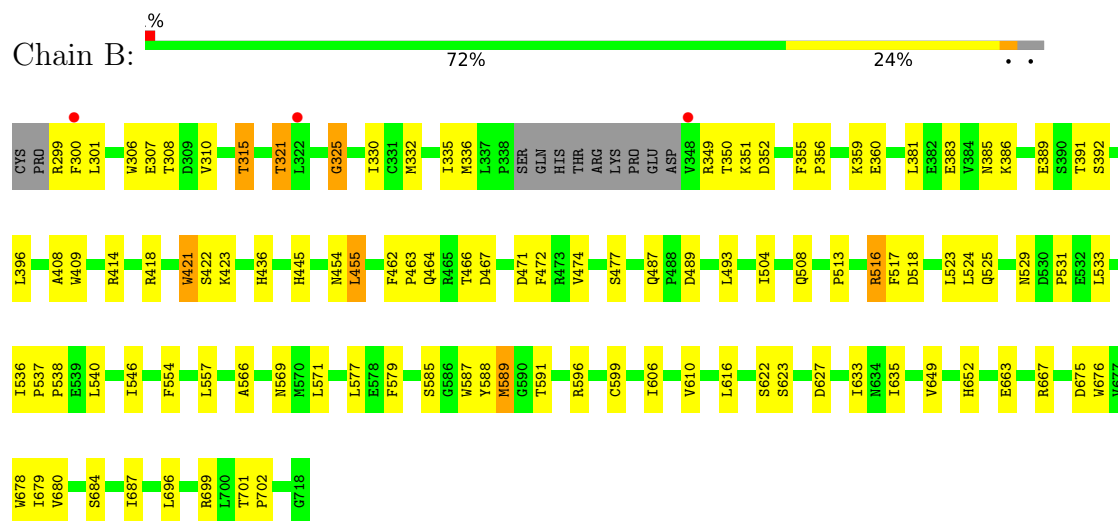
3 Residue-property plots [i](#)

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	51.58Å 110.25Å 164.39Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.86 – 2.45 38.86 – 2.45	Depositor EDS
% Data completeness (in resolution range)	98.8 (38.86-2.45) 98.7 (38.86-2.45)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.12 (at 2.45Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.209 , 0.265 0.204 , 0.263	Depositor DCC
R_{free} test set	1752 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	34.4	Xtriage
Anisotropy	1.160	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 49.6	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	7104	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.78% 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, H4B, ACT, HEM, JI7

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	0.52	0/3424	1.03	21/4645 (0.5%)
1	B	0.57	0/3438	1.03	22/4661 (0.5%)
All	All	0.55	0/6862	1.03	43/9306 (0.5%)

There are no bond length outliers.

All (43) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	354	LEU	N-CA-C	8.36	121.97	111.69
1	B	589	MET	N-CA-C	-8.20	95.88	109.07
1	A	589	MET	N-CA-C	-8.02	96.16	109.07
1	A	474	VAL	N-CA-C	-7.66	96.69	107.80
1	B	599	CYS	N-CA-C	7.41	122.49	113.38
1	B	351	LYS	N-CA-C	-7.28	103.34	111.71
1	A	709	ASP	N-CA-C	-7.10	100.83	109.57
1	B	408	ALA	N-CA-C	-6.88	103.71	111.14
1	A	688	THR	N-CA-C	-6.80	100.93	110.31
1	B	421	TRP	N-CA-C	6.47	119.16	111.33
1	B	588	TYR	N-CA-C	6.47	119.07	110.53
1	A	509	GLY	N-CA-C	6.39	120.42	113.58
1	A	546	ILE	N-CA-C	6.38	117.31	107.99
1	B	623	SER	N-CA-C	-6.35	105.50	113.18
1	B	325	GLY	N-CA-C	-6.29	106.44	114.37
1	A	588	TYR	N-CA-C	6.29	118.83	110.53
1	B	418	ARG	N-CA-C	6.21	120.33	112.87
1	B	349	ARG	N-CA-C	6.05	118.52	110.53
1	B	472	PHE	N-CA-C	-6.00	99.46	109.24
1	A	322	LEU	N-CA-C	5.93	117.99	110.33
1	B	699	ARG	N-CA-C	5.86	118.79	109.24
1	B	315	THR	N-CA-C	-5.86	106.78	114.04

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	474	VAL	N-CA-C	-5.82	98.76	107.37
1	A	455	LEU	N-CA-C	5.82	118.26	110.35
1	B	616	LEU	N-CA-C	-5.77	101.20	110.14
1	A	506	ILE	N-CA-C	-5.72	105.16	110.53
1	A	408	ALA	N-CA-C	-5.67	105.18	111.36
1	A	557	LEU	N-CA-C	-5.62	105.68	112.54
1	A	495	ASP	CA-C-N	5.52	125.87	119.47
1	A	495	ASP	C-N-CA	5.52	125.87	119.47
1	B	596	ARG	N-CA-C	5.46	118.29	111.24
1	B	652	HIS	N-CA-C	-5.44	105.26	111.14
1	B	436	HIS	N-CA-C	-5.44	105.43	111.36
1	B	591	THR	N-CA-C	5.30	118.53	111.75
1	A	319	LYS	N-CA-C	-5.29	106.42	112.92
1	A	635	ILE	N-CA-C	-5.27	105.25	111.00
1	A	564	LEU	N-CA-C	5.24	118.13	108.69
1	A	714	HIS	N-CA-C	5.24	117.77	109.96
1	B	455	LEU	N-CA-C	5.23	117.62	110.24
1	A	623	SER	N-CA-C	-5.20	106.20	112.54
1	B	546	ILE	N-CA-C	5.17	116.14	108.23
1	A	599	CYS	N-CA-C	5.10	119.55	113.23
1	B	471	ASP	N-CA-C	5.01	116.75	110.24

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	3243	140	0
1	B	3345	0	3259	71	0
2	A	4	0	3	2	0
2	B	4	0	3	0	0
3	A	1	0	0	0	0
4	A	43	0	30	4	0
4	B	43	0	30	4	0
5	A	17	0	15	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	B	17	0	15	0	0
6	A	26	0	28	6	0
6	B	26	0	28	4	0
7	A	113	0	0	6	0
7	B	134	0	0	11	0
All	All	7104	0	6654	213	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 16.

All (213) 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	1.36	1.08
1:A:322:LEU:HB3	1:A:699:ARG:HH21	1.40	0.84
1:A:321:THR:HG23	1:A:322:LEU:H	1.46	0.80
1:A:332:MET:HE2	1:B:696:LEU:HD21	1.65	0.77
1:A:470:HIS:HB3	1:A:527:ASN:ND2	1.98	0.77
1:A:384:VAL:O	1:A:388:ILE:HG13	1.85	0.75
1:B:516:ARG:HG2	1:B:517:PHE:CE1	2.22	0.75
1:A:380:ARG:NH1	1:A:397:LYS:HG2	2.03	0.73
1:A:480:ILE:HD13	1:A:541:VAL:HG13	1.70	0.73
1:A:610:VAL:O	1:A:614:MET:HG3	1.89	0.72
1:A:596:ARG:O	1:A:600:ASP:HB2	1.90	0.72
1:A:371:ARG:HH21	1:A:371:ARG:HG3	1.55	0.72
1:A:483:ALA:HB1	1:A:502:THR:CG2	2.22	0.70
1:A:523:LEU:HD22	1:A:531:PRO:CB	2.20	0.69
1:A:555:LYS:NZ	1:A:555:LYS:HB3	2.08	0.69
1:B:525:GLN:HG3	1:B:529:ASN:O	1.94	0.67
1:B:383:GLU:HA	1:B:386:LYS:HE3	1.75	0.67
1:A:655:THR:HB	1:A:690:VAL:HG13	1.76	0.67
1:A:675:ASP:O	1:A:679:ILE:HG12	1.96	0.65
1:B:566:ALA:HB2	1:B:585:SER:HB3	1.78	0.65
1:B:701:THR:HA	1:B:702:PRO:C	2.22	0.65
4:B:750:HEM:HBA2	6:B:800:JI7:H72	1.77	0.65
1:B:359:LYS:HG3	1:B:381:LEU:HD21	1.79	0.65
1:B:321:THR:HB	7:B:36:HOH:O	1.97	0.65
1:A:488:PRO:C	1:A:490:GLY:H	2.04	0.65
1:A:380:ARG:HD3	1:A:400:GLU:OE1	1.96	0.64
1:A:551:PHE:HE1	1:A:614:MET:HE3	1.62	0.64
1:A:630:LEU:HD22	1:B:687:ILE:HD13	1.79	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:537:PRO:HA	7:B:182:HOH:O	1.98	0.63
1:A:373:GLY:HA2	1:A:378:MET:HE3	1.79	0.63
1:A:355:PHE:N	1:A:356:PRO:HD2	2.13	0.63
1:B:391:THR:O	1:B:392:SER:HB2	1.99	0.63
1:A:307:GLU:HG3	7:B:105:HOH:O	1.99	0.62
1:B:325:GLY:O	1:B:332:MET:HE2	1.99	0.62
4:B:750:HEM:HMC2	4:B:750:HEM:HBC2	1.82	0.62
1:A:551:PHE:HB3	1:A:553:TRP:NE1	2.16	0.61
1:B:684:SER:HB3	1:B:687:ILE:HD11	1.82	0.60
1:A:667:ARG:HH11	1:A:667:ARG:HG3	1.67	0.60
1:B:355:PHE:CE1	1:B:385:ASN:HB2	2.36	0.60
1:A:306:TRP:CD1	1:B:336:MET:HE2	2.37	0.59
1:A:699:ARG:HD2	7:A:132:HOH:O	2.03	0.59
1:A:551:PHE:HD1	1:A:553:TRP:CZ2	2.21	0.59
1:B:332:MET:HB3	1:B:335:ILE:HG13	1.84	0.59
1:B:684:SER:HB3	1:B:687:ILE:CG1	2.32	0.58
1:B:386:LYS:O	1:B:389:GLU:HG2	2.03	0.58
1:A:350:THR:OG1	1:A:353:GLN:HG3	2.04	0.58
1:A:569:ASN:ND2	1:A:569:ASN:H	2.02	0.58
4:B:750:HEM:O2A	6:B:800:JI7:H3'	2.04	0.57
1:A:548:HIS:ND1	1:A:549:PRO:HD2	2.19	0.57
1:A:475:TRP:HB2	1:A:523:LEU:HB3	1.86	0.57
1:A:516:ARG:HD3	1:A:517:PHE:CE1	2.39	0.57
1:A:488:PRO:C	1:A:490:GLY:N	2.63	0.57
1:A:569:ASN:H	1:A:569:ASN:HD22	1.52	0.57
1:A:321:THR:HG23	1:A:322:LEU:N	2.20	0.56
1:A:371:ARG:HH21	1:A:371:ARG:CG	2.18	0.56
4:A:750:HEM:HMC2	4:A:750:HEM:HBC2	1.87	0.56
1:B:523:LEU:CD1	1:B:531:PRO:HB2	2.34	0.56
1:A:483:ALA:HB1	1:A:502:THR:HG21	1.87	0.56
1:B:508:GLN:HB2	1:B:533:LEU:HD11	1.86	0.56
1:A:322:LEU:HB3	1:A:699:ARG:NH2	2.17	0.55
1:A:627:ASP:O	1:A:631:VAL:HG23	2.06	0.55
1:A:483:ALA:HB1	1:A:502:THR:HG23	1.89	0.55
1:A:321:THR:HG23	1:A:322:LEU:HG	1.87	0.55
1:A:551:PHE:CD2	1:A:551:PHE:N	2.74	0.55
1:A:569:ASN:HD22	1:A:569:ASN:N	2.03	0.55
1:A:494:GLY:O	1:A:496:PRO:HD3	2.07	0.55
1:A:478:GLN:HB2	1:A:481:ARG:HG3	1.89	0.54
1:A:551:PHE:HD1	1:A:553:TRP:HZ2	1.56	0.54
1:B:566:ALA:CB	1:B:585:SER:HB3	2.38	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:473:ARG:HD2	1:A:580:SER:HB2	1.91	0.53
1:B:383:GLU:HA	1:B:386:LYS:CE	2.38	0.53
1:A:322:LEU:HB2	1:A:699:ARG:HE	1.74	0.53
1:B:356:PRO:O	1:B:360:GLU:HG3	2.10	0.52
1:B:589:MET:HA	1:B:649:VAL:O	2.10	0.52
1:A:551:PHE:N	1:A:551:PHE:HD2	2.08	0.52
1:A:684:SER:HB3	1:A:687:ILE:HG12	1.92	0.52
1:B:504:ILE:O	1:B:508:GLN:HG2	2.10	0.52
1:A:382:GLU:OE1	1:A:385:ASN:HB3	2.09	0.52
1:B:462:PHE:HB3	1:B:463:PRO:CD	2.39	0.52
1:A:606:ILE:O	1:A:610:VAL:HG23	2.11	0.51
1:B:355:PHE:N	1:B:356:PRO:HD2	2.25	0.51
1:B:513:PRO:HG2	1:B:518:ASP:CG	2.36	0.51
1:A:353:GLN:O	1:A:356:PRO:HG2	2.10	0.51
6:B:800:JI7:H2'1	7:B:199:HOH:O	2.10	0.51
1:A:337:LEU:HD11	6:A:800:JI7:CL	2.48	0.51
1:A:589:MET:HE3	1:A:591:THR:OG1	2.11	0.50
1:B:675:ASP:O	1:B:679:ILE:HG12	2.10	0.50
1:A:470:HIS:HB3	1:A:527:ASN:HD22	1.76	0.50
1:A:508:GLN:NE2	1:A:716:TRP:CZ3	2.80	0.50
1:B:538:PRO:HD3	7:B:182:HOH:O	2.10	0.50
1:A:551:PHE:HB3	1:A:553:TRP:CE2	2.47	0.50
1:B:423:LYS:HE3	7:B:142:HOH:O	2.11	0.50
1:A:516:ARG:O	1:A:516:ARG:HG2	2.12	0.50
1:A:465:ARG:HG3	1:A:471:ASP:OD1	2.12	0.49
1:A:501:PHE:CD2	1:A:520:LEU:HD13	2.46	0.49
1:A:523:LEU:CD2	1:A:531:PRO:HB2	2.23	0.49
1:A:508:GLN:HA	1:A:508:GLN:OE1	2.11	0.49
1:A:485:TYR:CE2	1:A:512:ALA:HB1	2.48	0.49
1:A:587:TRP:CZ2	2:A:860:ACT:H3	2.48	0.49
1:A:489:ASP:HB2	7:A:97:HOH:O	2.13	0.48
1:B:299:ARG:HB3	1:B:299:ARG:NH1	2.28	0.48
1:A:445:HIS:C	1:A:445:HIS:CD2	2.91	0.48
1:A:409:TRP:CE3	1:A:421:TRP:HA	2.48	0.48
1:A:402:ILE:O	1:A:403:TYR:C	2.56	0.48
1:A:500:GLN:O	1:A:504:ILE:HG13	2.13	0.48
1:A:686:SER:HA	1:A:691:PHE:CG	2.48	0.48
1:B:487:GLN:OE1	1:B:487:GLN:HA	2.12	0.48
1:A:301:LEU:CD1	1:B:330:ILE:HD13	2.43	0.48
1:B:684:SER:HB3	1:B:687:ILE:CD1	2.42	0.48
1:A:488:PRO:O	1:A:490:GLY:N	2.47	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:750:HEM:O2A	6:A:800:JI7:H3'	2.14	0.48
1:A:510:TRP:CE2	1:A:521:PRO:HD3	2.49	0.47
1:A:406:LYS:HE2	1:A:422:SER:O	2.14	0.47
1:A:555:LYS:HB3	1:A:555:LYS:HZ3	1.79	0.47
1:A:571:LEU:C	1:A:571:LEU:HD23	2.39	0.47
1:A:449:ALA:O	1:A:455:LEU:HA	2.14	0.47
1:B:524:LEU:O	1:B:531:PRO:HA	2.13	0.47
1:A:512:ALA:C	1:A:514:ARG:H	2.21	0.47
1:A:667:ARG:HG3	1:A:667:ARG:NH1	2.29	0.47
1:B:606:ILE:HD11	1:B:633:ILE:HD13	1.96	0.47
1:B:423:LYS:HG3	7:B:142:HOH:O	2.13	0.47
1:A:350:THR:N	1:A:353:GLN:OE1	2.48	0.47
1:B:684:SER:CB	1:B:687:ILE:HD11	2.45	0.47
1:A:557:LEU:HB2	1:A:559:LEU:HG	1.97	0.47
6:B:800:JI7:H21	7:B:14:HOH:O	2.14	0.47
1:A:496:PRO:O	1:A:499:VAL:HG23	2.15	0.47
1:A:306:TRP:CG	1:B:336:MET:HE2	2.50	0.47
1:A:418:ARG:C	1:A:420:GLN:N	2.71	0.46
1:B:300:PHE:N	1:B:300:PHE:CD1	2.83	0.46
1:B:308:THR:OG1	1:B:310:VAL:HG23	2.16	0.46
1:B:610:VAL:HG21	1:B:633:ILE:HD11	1.96	0.46
1:A:355:PHE:N	1:A:356:PRO:CD	2.77	0.46
1:A:508:GLN:NE2	1:A:716:TRP:HZ3	2.12	0.46
1:A:535:GLN:O	1:A:536:ILE:C	2.57	0.46
1:A:555:LYS:HB3	1:A:555:LYS:HZ2	1.79	0.46
1:B:409:TRP:CE3	1:B:421:TRP:HA	2.50	0.46
1:A:684:SER:O	1:A:687:ILE:HG12	2.15	0.46
1:A:450:THR:O	1:A:455:LEU:HD23	2.16	0.46
1:A:470:HIS:ND1	1:A:470:HIS:N	2.64	0.46
1:A:617:ASP:OD2	1:A:619:ARG:HB2	2.15	0.46
1:A:371:ARG:CG	1:A:371:ARG:NH2	2.77	0.46
1:A:455:LEU:HD12	1:A:587:TRP:CB	2.46	0.45
2:A:860:ACT:CH3	4:A:750:HEM:HAB	2.45	0.45
1:B:516:ARG:HD2	7:B:169:HOH:O	2.16	0.45
1:A:501:PHE:HD2	1:A:520:LEU:HD13	1.80	0.45
1:B:477:SER:HB2	1:B:569:ASN:ND2	2.31	0.45
1:A:684:SER:HB3	1:A:687:ILE:CG1	2.46	0.45
1:B:445:HIS:C	1:B:445:HIS:CD2	2.94	0.45
1:A:551:PHE:O	1:A:554:PHE:HB2	2.16	0.45
1:A:569:ASN:ND2	1:A:569:ASN:N	2.61	0.45
1:A:614:MET:CE	1:A:632:GLU:HG3	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:663:GLU:HB3	1:B:667:ARG:NH1	2.32	0.45
6:A:800:JI7:H23	1:B:306:TRP:CZ3	2.52	0.45
1:A:469:LYS:C	1:A:470:HIS:ND1	2.75	0.45
1:A:403:TYR:CE1	1:A:407:HIS:CE1	3.06	0.44
1:A:489:ASP:C	1:A:491:SER:H	2.24	0.44
1:B:299:ARG:HB3	1:B:299:ARG:HH11	1.82	0.44
1:B:462:PHE:HB3	1:B:463:PRO:HD2	1.99	0.44
1:B:493:LEU:HB2	7:B:153:HOH:O	2.18	0.44
1:A:643:SER:C	1:A:645:LYS:H	2.25	0.44
1:A:709:ASP:HB2	1:A:712:ASN:ND2	2.32	0.44
1:B:606:ILE:O	1:B:610:VAL:HG23	2.17	0.44
1:A:391:THR:O	1:A:392:SER:OG	2.24	0.44
1:B:396:LEU:HG	1:B:577:LEU:HD12	2.00	0.44
1:A:455:LEU:HD12	1:A:587:TRP:HB3	1.98	0.44
1:A:547:ARG:HB2	7:A:73:HOH:O	2.17	0.43
1:A:589:MET:HA	1:A:649:VAL:O	2.17	0.43
1:A:715:VAL:O	1:A:715:VAL:HG23	2.18	0.43
1:A:476:ASN:OD1	1:A:476:ASN:N	2.51	0.43
1:A:701:THR:HA	1:A:702:PRO:C	2.44	0.43
1:B:414:ARG:HD3	1:B:678:TRP:CD2	2.54	0.43
1:B:454:ASN:ND2	7:B:234:HOH:O	2.52	0.43
1:A:408:ALA:HA	1:A:702:PRO:O	2.19	0.43
1:A:475:TRP:CE3	1:A:523:LEU:HD13	2.53	0.43
1:A:590:GLY:HA3	1:A:638:LEU:HD21	2.01	0.43
1:A:316:LEU:HD22	1:A:669:ARG:HG2	2.01	0.42
1:A:508:GLN:HE22	1:A:716:TRP:HZ3	1.66	0.42
4:A:750:HEM:HBA2	6:A:800:JI7:H72	2.01	0.42
1:B:489:ASP:OD2	1:B:489:ASP:C	2.62	0.42
1:B:676:TRP:CZ2	1:B:680:VAL:HG21	2.54	0.42
1:A:316:LEU:HD12	1:A:319:LYS:HD2	2.01	0.42
1:A:332:MET:HE1	1:B:301:LEU:HD22	2.00	0.42
1:B:464:GLN:HB3	1:B:579:PHE:CE2	2.54	0.42
1:A:516:ARG:HD3	1:A:517:PHE:HE1	1.84	0.42
1:A:524:LEU:O	1:A:531:PRO:HA	2.20	0.42
1:A:542:LEU:HD21	1:A:646:VAL:HG22	2.01	0.42
1:A:595:VAL:O	1:A:599:CYS:HB2	2.20	0.42
1:A:590:GLY:HA3	7:A:4:HOH:O	2.18	0.42
1:A:592:GLU:OE1	6:A:800:JI7:N1	2.52	0.42
1:A:599:CYS:O	1:A:600:ASP:C	2.61	0.42
1:B:466:THR:OG1	1:B:467:ASP:N	2.52	0.42
7:A:101:HOH:O	1:B:307:GLU:HG3	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:554:PHE:O	1:B:557:LEU:HB2	2.20	0.42
1:A:354:LEU:C	1:A:356:PRO:HD2	2.45	0.42
7:A:143:HOH:O	1:B:635:ILE:HA	2.19	0.42
1:A:472:PHE:C	1:A:473:ARG:HG2	2.45	0.41
1:A:551:PHE:CD1	1:A:553:TRP:HZ2	2.38	0.41
1:A:684:SER:OG	1:B:627:ASP:OD2	2.32	0.41
1:B:571:LEU:C	1:B:571:LEU:HD23	2.45	0.41
1:B:350:THR:HG22	1:B:352:ASP:H	1.85	0.41
1:B:701:THR:CA	1:B:702:PRO:C	2.92	0.41
1:A:337:LEU:HD21	6:A:800:JI7:CL	2.57	0.41
1:A:554:PHE:O	1:A:557:LEU:HG	2.21	0.41
1:A:537:PRO:HB2	1:A:540:LEU:HG	2.02	0.41
1:A:567:VAL:HB	1:A:584:PHE:CE1	2.56	0.41
1:A:472:PHE:O	1:A:473:ARG:HG2	2.21	0.41
1:B:536:ILE:O	1:B:537:PRO:C	2.64	0.41
1:B:684:SER:O	1:B:687:ILE:HG12	2.21	0.41
1:A:495:ASP:OD1	1:A:497:ALA:HB3	2.22	0.40
1:A:380:ARG:NH2	1:A:384:VAL:HG22	2.36	0.40
1:B:587:TRP:O	4:B:750:HEM:HMB3	2.21	0.40
1:A:439:PHE:CZ	1:A:443:CYS:SG	3.14	0.40
1:A:391:THR:O	1:A:392:SER:CB	2.68	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	405/422 (96%)	370 (91%)	31 (8%)	4 (1%)	12	14
1	B	407/422 (96%)	383 (94%)	24 (6%)	0	100	100
All	All	812/844 (96%)	753 (93%)	55 (7%)	4 (0%)	24	32

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	412	ALA
1	A	489	ASP
1	A	605	ASN
1	A	685	GLY

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	365/377 (97%)	351 (96%)	14 (4%)	29	43
1	B	366/377 (97%)	359 (98%)	7 (2%)	50	64
All	All	731/754 (97%)	710 (97%)	21 (3%)	37	52

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	380	ARG
1	A	389	GLU
1	A	470	HIS
1	A	514	ARG
1	A	523	LEU
1	A	527	ASN
1	A	538	PRO
1	A	551	PHE
1	A	555	LYS
1	A	569	ASN
1	A	600	ASP
1	A	635	ILE
1	A	645	LYS
1	A	662	MET
1	B	315	THR
1	B	321	THR
1	B	422	SER
1	B	455	LEU
1	B	516	ARG

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Mol	Chain	Res	Type
1	B	540	LEU
1	B	622	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	425	GLN
1	A	454	ASN
1	A	487	GLN
1	A	508	GLN
1	A	527	ASN
1	A	569	ASN
1	A	605	ASN
1	A	628	GLN
1	A	697	ASN
1	A	712	ASN
1	B	353	GLN
1	B	364	GLN
1	B	385	ASN
1	B	395	GLN
1	B	425	GLN
1	B	454	ASN
1	B	507	GLN
1	B	527	ASN
1	B	535	GLN
1	B	569	ASN
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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	860	-	3,3,3	0.96	0	3,3,3	0.81	0
6	JI7	B	800	-	27,28,28	1.63	9 (33%)	29,37,37	1.21	3 (10%)
2	ACT	B	860	-	3,3,3	0.93	0	3,3,3	0.61	0
5	H4B	B	760	-	17,18,18	1.66	3 (17%)	14,26,26	1.75	3 (21%)
5	H4B	A	760	-	17,18,18	1.65	2 (11%)	14,26,26	1.90	3 (21%)
4	HEM	A	750	1	50,50,50	1.40	8 (16%)	67,82,82	1.35	8 (11%)
6	JI7	A	800	-	27,28,28	1.44	8 (29%)	29,37,37	1.16	2 (6%)
4	HEM	B	750	1	50,50,50	1.26	7 (14%)	67,82,82	1.37	7 (10%)

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
6	JI7	B	800	-	-	4/12/22/22	0/3/3/3
5	H4B	B	760	-	-	0/8/17/17	0/2/2/2
5	H4B	A	760	-	-	0/8/17/17	0/2/2/2
4	HEM	A	750	1	-	5/14/54/54	-
6	JI7	A	800	-	-	3/12/22/22	0/3/3/3
4	HEM	B	750	1	-	2/14/54/54	-

All (37) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	760	H4B	C4A-N5	4.07	1.47	1.37
5	B	760	H4B	C6-N5	3.99	1.54	1.47
5	A	760	H4B	C6-N5	3.99	1.54	1.47
5	B	760	H4B	C4A-N5	3.84	1.46	1.37
4	B	750	HEM	C4D-ND	-3.14	1.34	1.40
6	B	800	JI7	C23-C22	3.12	1.43	1.38
4	A	750	HEM	CHC-C1C	2.96	1.44	1.38
4	B	750	HEM	FE-NB	2.87	2.03	1.94
4	A	750	HEM	CAB-C3B	-2.85	1.39	1.47
6	B	800	JI7	C21-C22	2.79	1.42	1.38
4	A	750	HEM	CAC-C3C	-2.72	1.40	1.47
6	A	800	JI7	C23-C22	2.68	1.43	1.38
4	B	750	HEM	CAC-C3C	-2.63	1.40	1.47
6	B	800	JI7	C3-C4	2.63	1.43	1.39
4	B	750	HEM	CAB-C3B	-2.54	1.40	1.47
6	B	800	JI7	C24-C25	2.53	1.43	1.38
6	B	800	JI7	C5-C6	2.51	1.44	1.40
6	A	800	JI7	C24-C25	2.47	1.43	1.38
4	B	750	HEM	C4C-NC	-2.41	1.35	1.39
6	B	800	JI7	C21-C26	2.40	1.43	1.39
6	B	800	JI7	C25-C26	2.40	1.43	1.38
6	A	800	JI7	C24-C23	2.39	1.43	1.38
6	A	800	JI7	C25-C26	2.38	1.43	1.38
6	A	800	JI7	C5-C6	2.37	1.43	1.40
4	B	750	HEM	C3B-C4B	2.32	1.49	1.44
6	A	800	JI7	C21-C22	2.30	1.42	1.38
6	A	800	JI7	C21-C26	2.29	1.43	1.39
4	A	750	HEM	FE-ND	2.29	2.01	1.94
4	A	750	HEM	C4D-ND	-2.25	1.36	1.40
6	B	800	JI7	C5-C4	2.25	1.42	1.39
4	A	750	HEM	FE-NA	2.22	2.02	1.95
4	A	750	HEM	FE-NB	2.21	2.01	1.94
5	B	760	H4B	C7-N8	2.21	1.48	1.46
4	A	750	HEM	C1C-C2C	2.18	1.49	1.45
4	B	750	HEM	C1D-ND	-2.16	1.34	1.38
6	B	800	JI7	C24-C23	2.10	1.42	1.38
6	A	800	JI7	C5-C4	2.08	1.42	1.39

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	760	H4B	C2-N1-C8A	5.27	122.66	113.36
5	B	760	H4B	C2-N1-C8A	4.73	121.71	113.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	750	HEM	C1D-C2D-C3D	-4.61	102.13	106.98
4	A	750	HEM	CBA-CAA-C2A	-4.30	100.66	112.53
6	B	800	JI7	C6-N1-C2	3.91	120.99	118.07
4	B	750	HEM	C2D-C1D-ND	3.73	114.21	109.90
6	A	800	JI7	N6-C6-N1	3.22	121.76	116.59
6	A	800	JI7	C6-N1-C2	3.20	120.46	118.07
4	B	750	HEM	CBA-CAA-C2A	-3.14	103.84	112.53
4	B	750	HEM	C3B-C4B-NB	3.11	111.70	109.47
4	A	750	HEM	C3C-C2C-C1C	-3.11	104.11	107.05
5	A	760	H4B	N3-C2-N1	-2.88	118.06	123.32
6	B	800	JI7	N6-C6-N1	2.86	121.19	116.59
4	A	750	HEM	CAB-C3B-C2B	-2.77	119.43	128.43
5	B	760	H4B	N3-C2-N1	-2.74	118.31	123.32
5	A	760	H4B	O10-C10-C9	-2.71	105.29	109.77
4	A	750	HEM	CBD-CAD-C3D	-2.45	105.76	112.53
5	B	760	H4B	O10-C10-C9	-2.19	106.15	109.77
6	B	800	JI7	C7-C2-N1	-2.18	113.90	116.57
4	B	750	HEM	CHD-C1D-C2D	-2.13	121.67	125.03
4	B	750	HEM	CHD-C4C-NC	-2.06	122.21	124.45
4	A	750	HEM	CMD-C2D-C1D	2.05	128.24	125.03
4	A	750	HEM	O1A-CGA-CBA	-2.04	116.62	123.09
4	A	750	HEM	C1D-C2D-C3D	-2.04	104.84	106.98
4	A	750	HEM	CAB-C3B-C4B	2.03	133.36	124.39
4	B	750	HEM	O1D-CGD-CBD	-2.02	116.68	123.09

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	800	JI7	C10-C9-N8-C3'
6	B	800	JI7	N11-C10-C9-N8
6	B	800	JI7	C9-C10-N11-C12
6	B	800	JI7	C2'-C3'-N8-C9
6	B	800	JI7	C4'-C3'-N8-C9
4	A	750	HEM	C3D-CAD-CBD-CGD
4	B	750	HEM	CAA-CBA-CGA-O2A
4	A	750	HEM	CAD-CBD-CGD-O2D
4	B	750	HEM	CAA-CBA-CGA-O1A
4	A	750	HEM	CAD-CBD-CGD-O1D
4	A	750	HEM	CAA-CBA-CGA-O2A
6	A	800	JI7	C9-C10-N11-C12
4	A	750	HEM	CAA-CBA-CGA-O1A

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Mol	Chain	Res	Type	Atoms
6	A	800	JI7	N11-C10-C9-N8

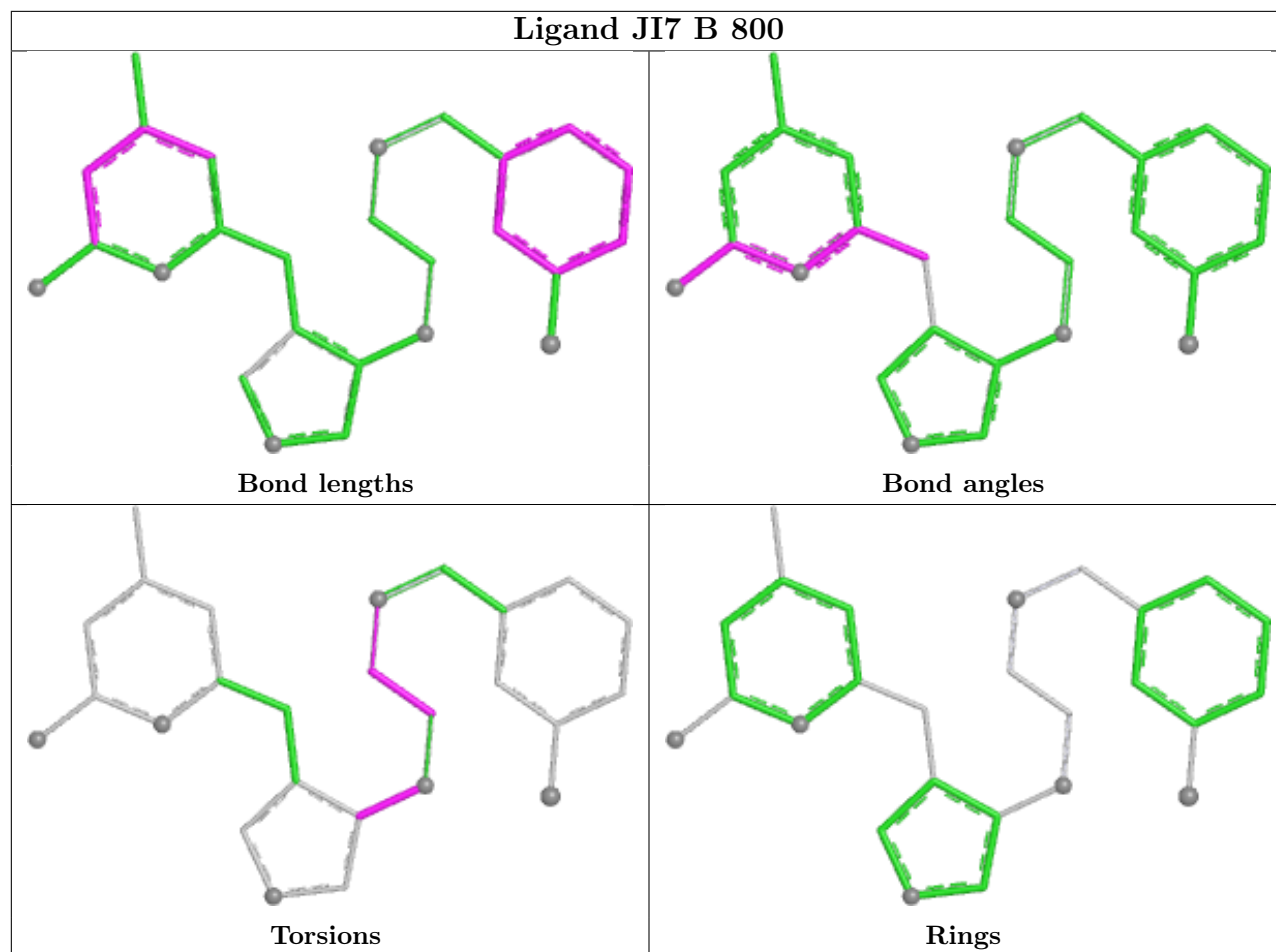
There are no ring outliers.

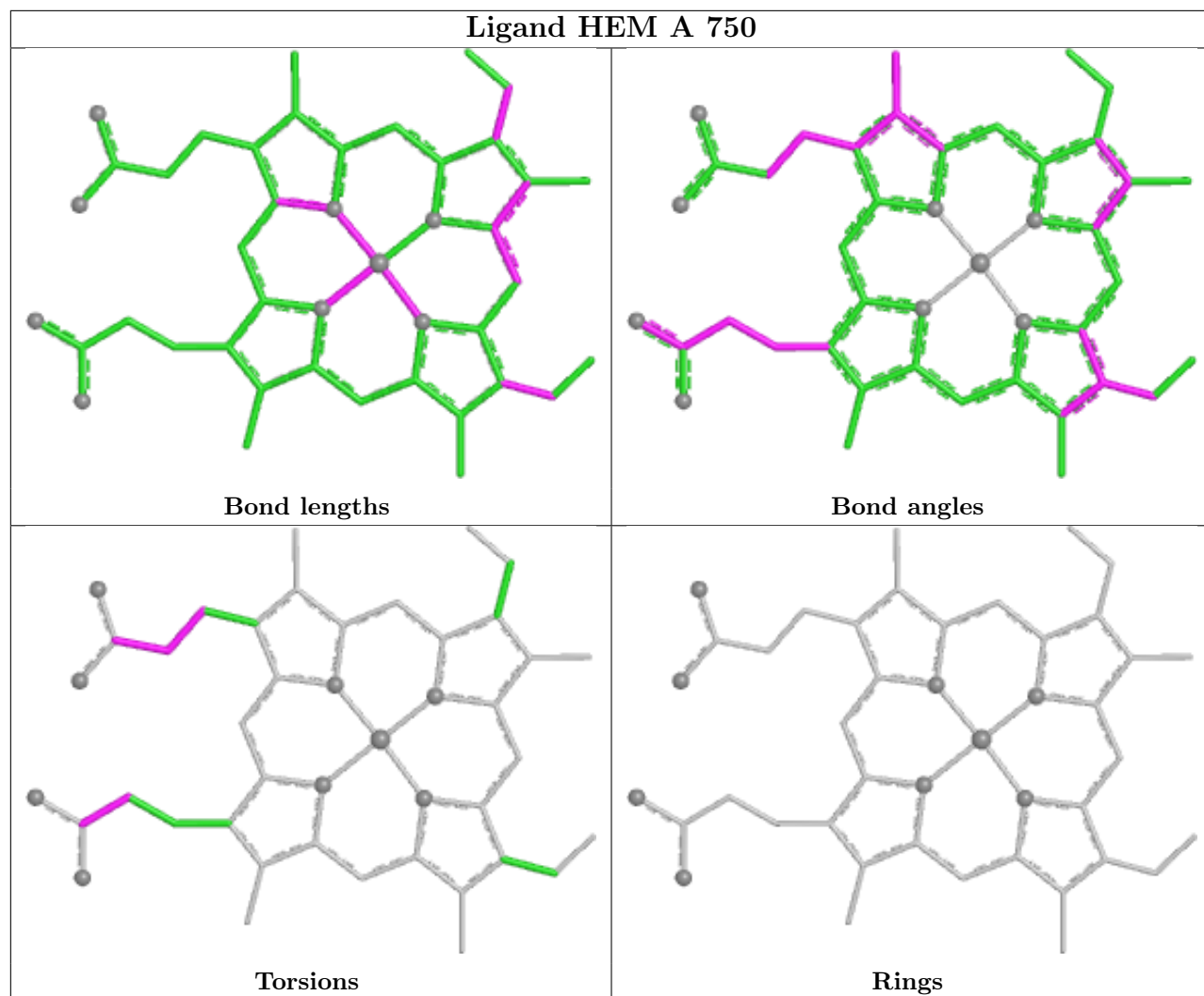
5 monomers are involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	860	ACT	2	0
6	B	800	JI7	4	0
4	A	750	HEM	4	0
6	A	800	JI7	6	0
4	B	750	HEM	4	0

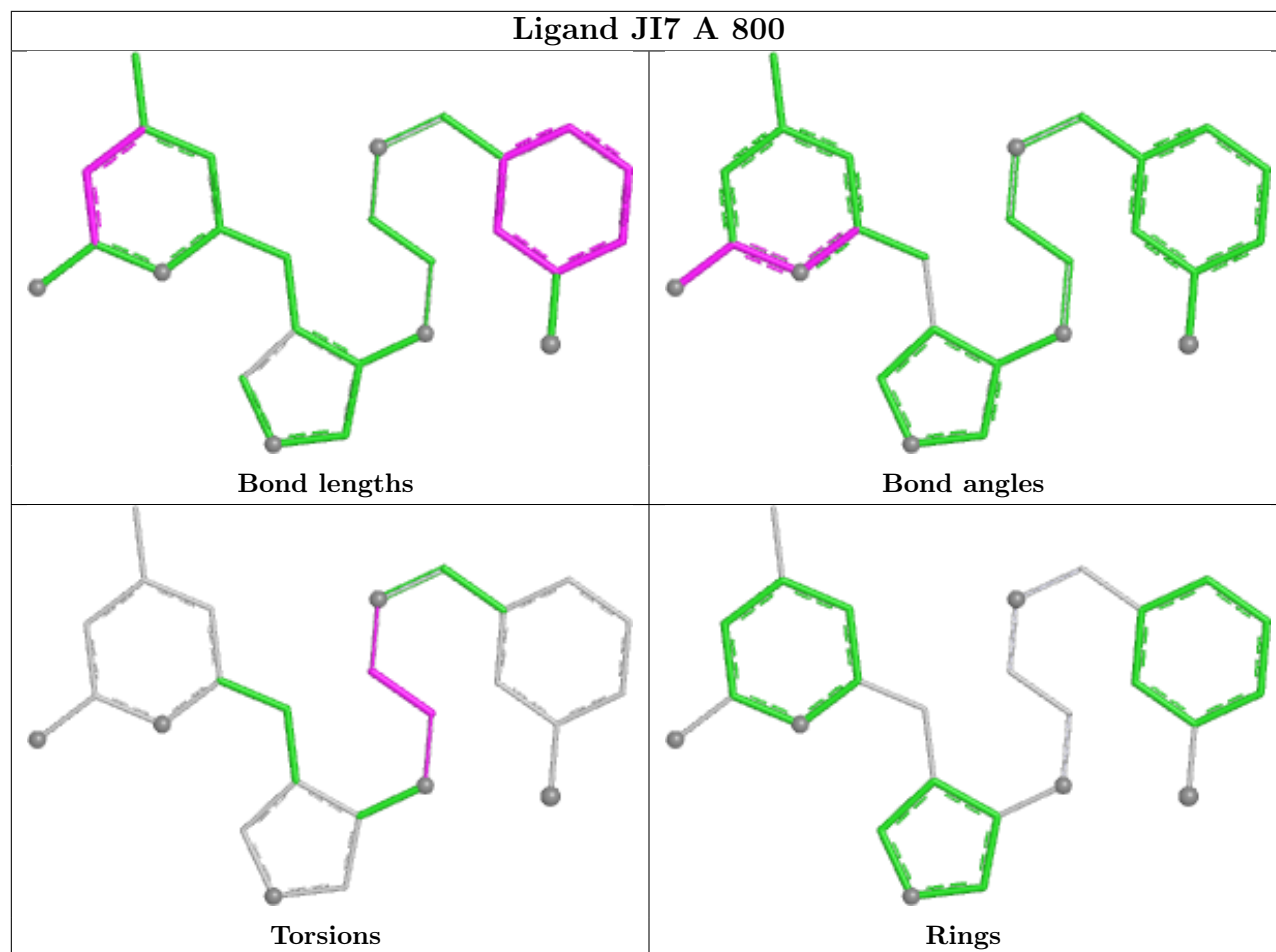
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

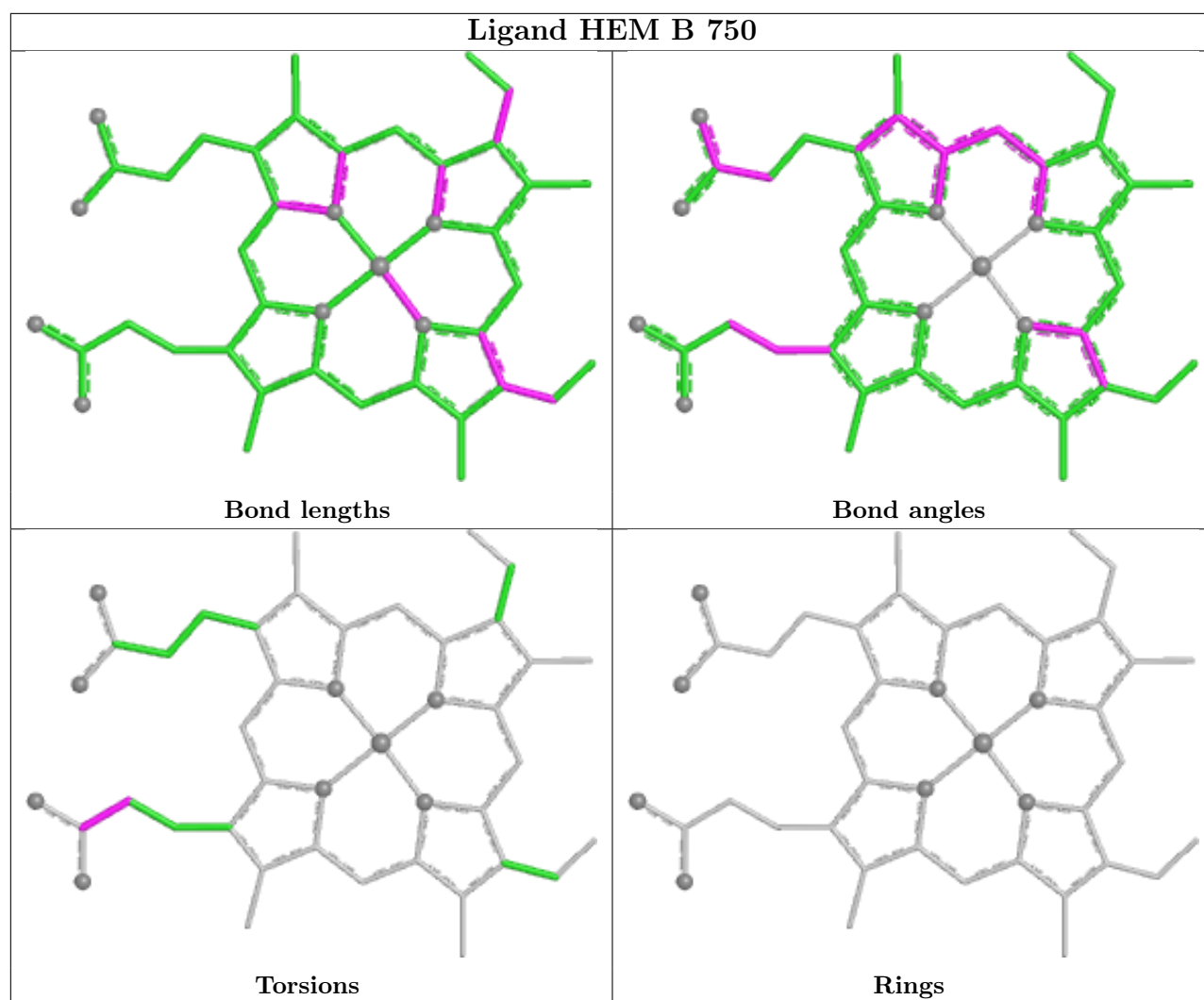
Ligand JI7 B 800





Ligand JI7 A 800





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	409/422 (96%)	0.01	7 (1%) 69 71	27, 47, 73, 90	0
1	B	411/422 (97%)	-0.30	3 (0%) 84 86	21, 40, 64, 80	0
All	All	820/844 (97%)	-0.15	10 (1%) 76 78	21, 43, 69, 90	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	355	PHE	3.2
1	A	348	VAL	2.7
1	B	300	PHE	2.5
1	A	716	TRP	2.5
1	B	348	VAL	2.4
1	A	715	VAL	2.2
1	A	488	PRO	2.2
1	A	493	LEU	2.1
1	B	322	LEU	2.0
1	A	300	PHE	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](#)

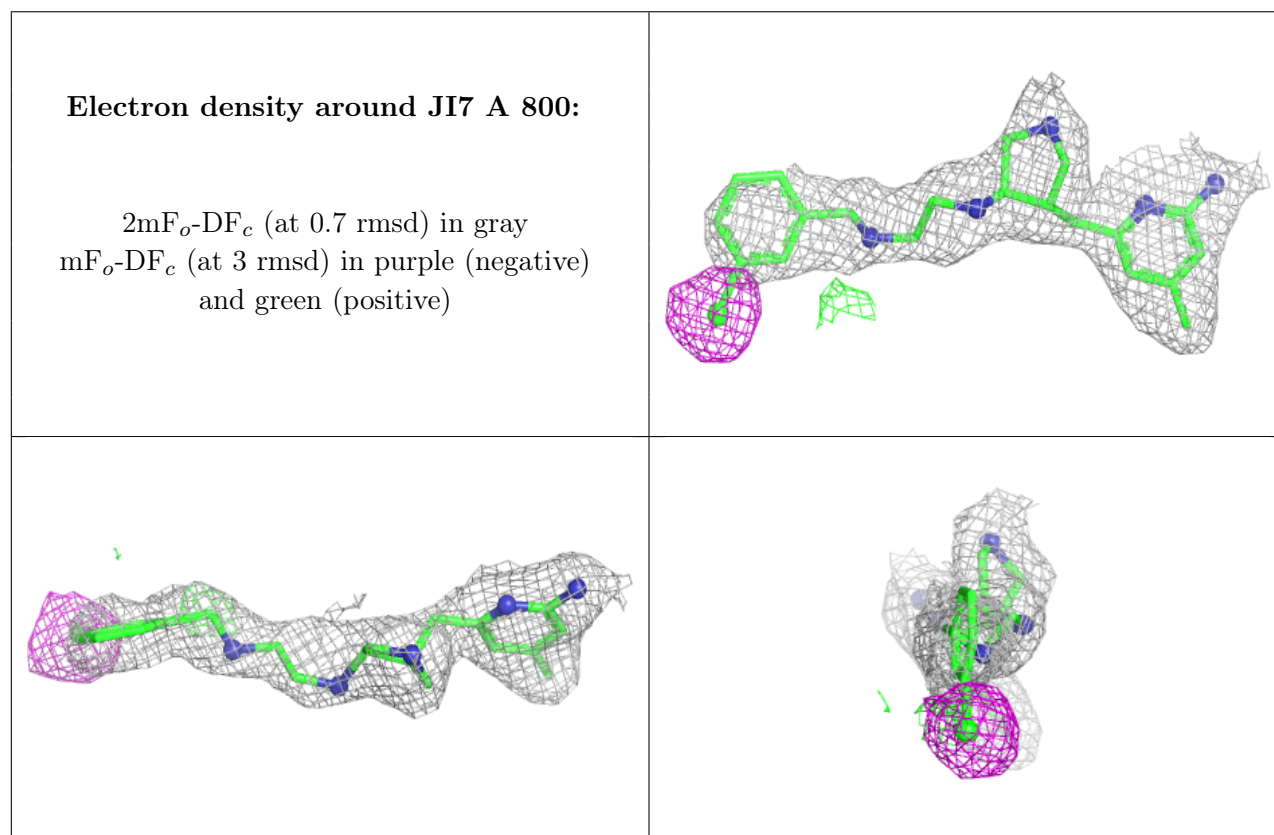
There are no oligosaccharides in this entry.

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.

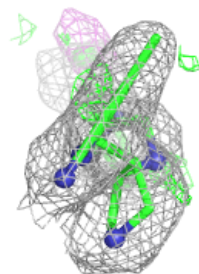
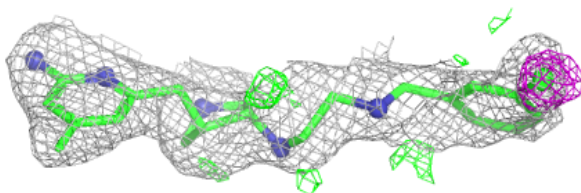
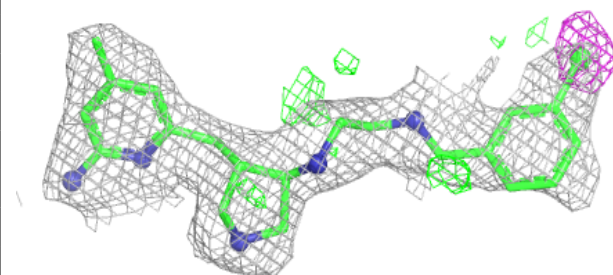
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	JI7	A	800	26/26	0.81	0.15	31,49,78,79	0
6	JI7	B	800	26/26	0.86	0.14	33,48,72,73	0
2	ACT	A	860	4/4	0.90	0.13	62,63,63,64	0
5	H4B	B	760	17/17	0.94	0.07	33,36,39,40	0
5	H4B	A	760	17/17	0.95	0.07	36,39,43,44	0
2	ACT	B	860	4/4	0.95	0.10	36,39,41,43	0
4	HEM	A	750	43/43	0.97	0.07	27,31,39,44	0
4	HEM	B	750	43/43	0.98	0.06	25,31,41,46	0
3	ZN	A	900	1/1	1.00	0.01	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.



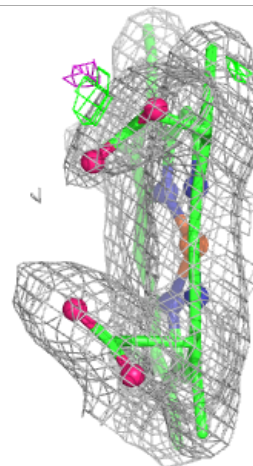
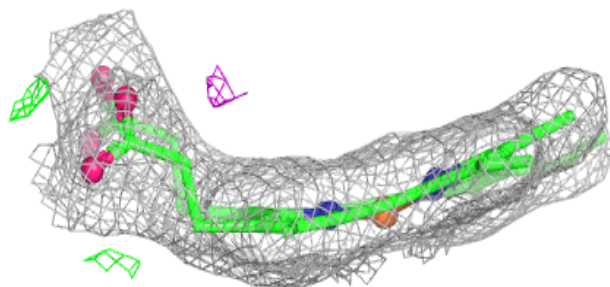
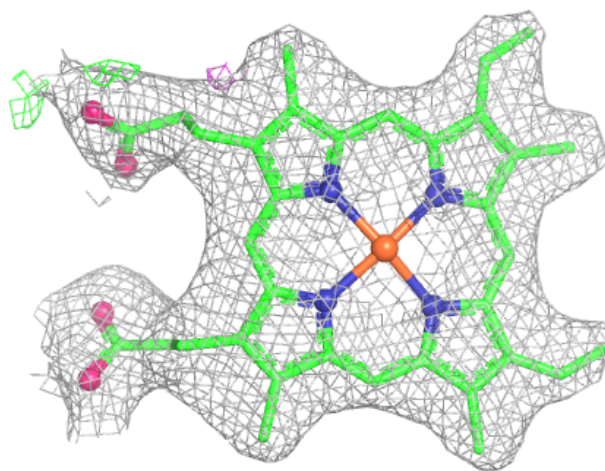
Electron density around JI7 B 800:

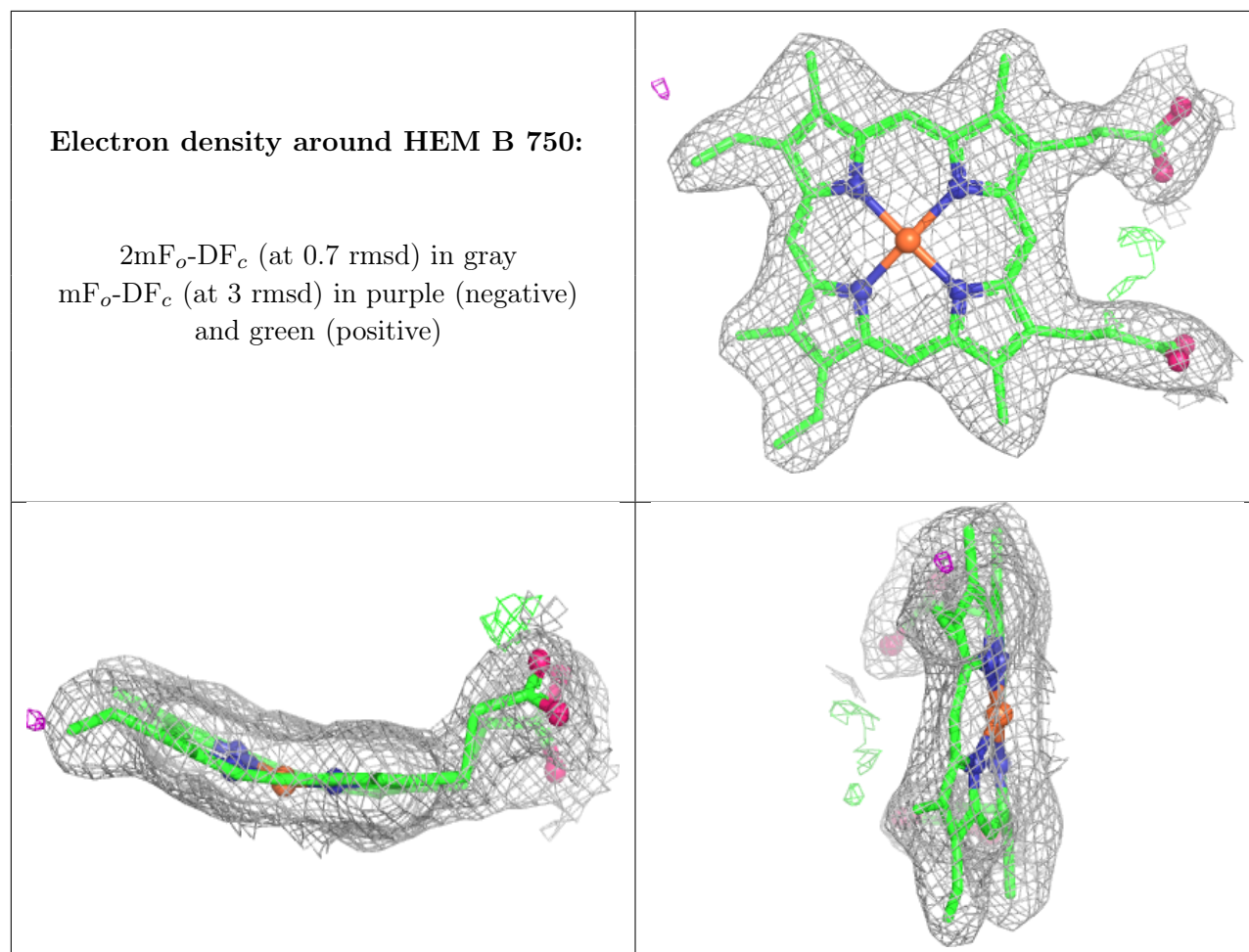
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around HEM A 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.