Full wwPDB X-ray Structure Validation Report (i)

Apr 27, 2024 – 11:37 pm BST

PDB ID	:	4V3W
Title	:	Structure of rat neuronal nitric oxide synthase heme domain in complex with 2-(2-(1H-imidazol-1-yl)pyrimidin-4-yl)-N-(3- fluorophenethyl)ethan-1-amine
Authors	:	Li, H.; Poulos, T.L.
Deposited on	:	2014-10-20
Resolution	:	2.13 Å(reported)

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

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

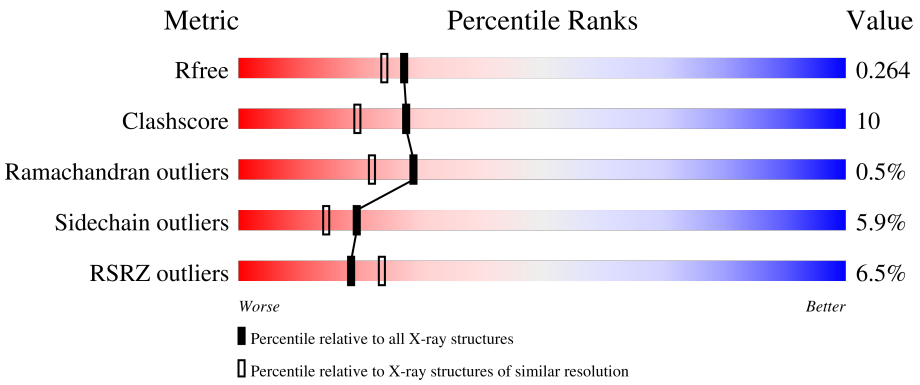
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36.2
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.13 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}	130704	2523 (2.16-2.12)
Clashscore	141614	2653 (2.16-2.12)
Ramachandran outliers	138981	2618 (2.16-2.12)
Sidechain outliers	138945	2617 (2.16-2.12)
RSRZ outliers	127900	2485 (2.16-2.12)

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

Mol	Chain	Length	Quality of chain
1	A	422	9% 71% 24%
1	B	422	3% 73% 22%

2 Entry composition (i)

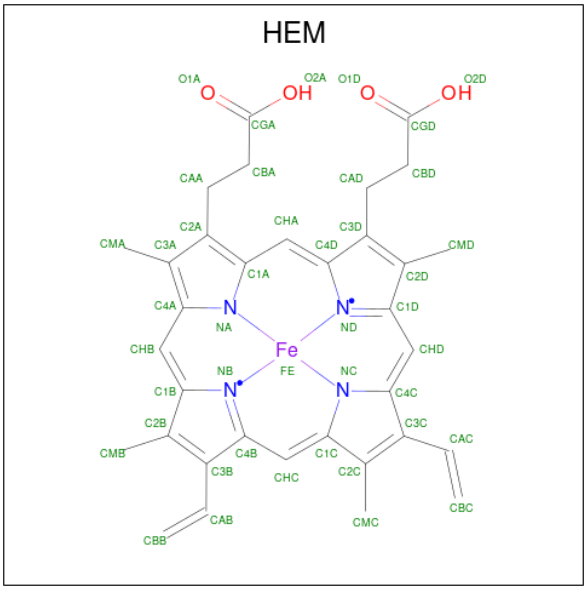
There are 7 unique types of molecules in this entry. The entry contains 7004 atoms, of which 0 are hydrogens and 0 are deuteriums.

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.

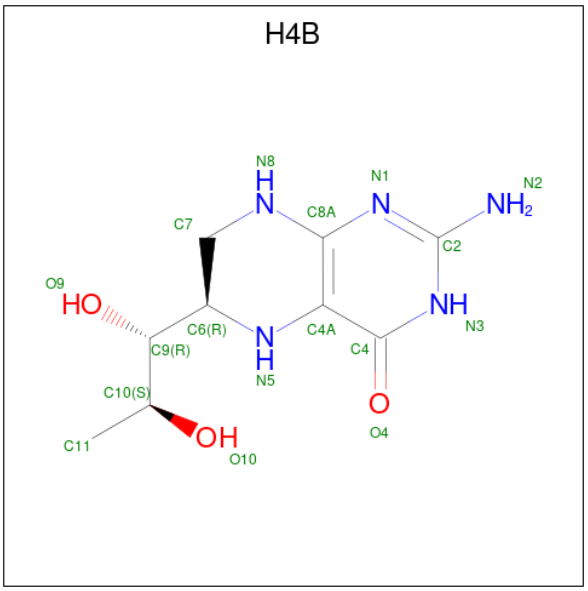
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	408	3323	2127	567	607	22	0	2	1
1	B	411	3354	2146	574	613	21	0	2	0

- Molecule 2 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



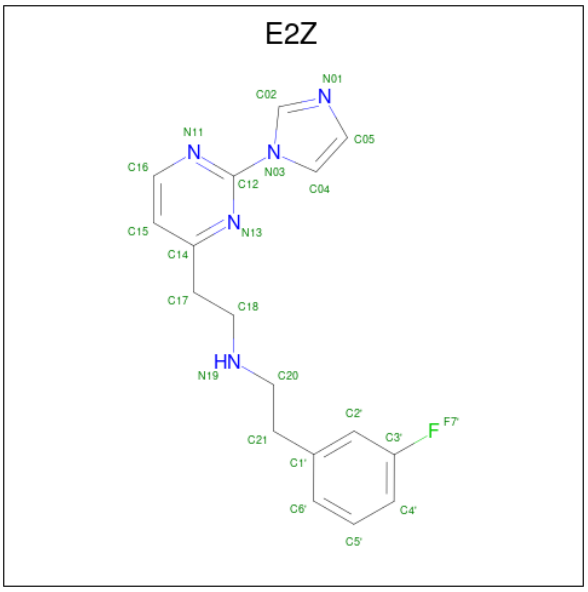
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	Fe	N	O		
2	A	1	43	34	1	4	4	0	0
2	B	1	43	34	1	4	4	0	0

- Molecule 3 is 5,6,7,8-TETRAHYDROBIOPTERIN (three-letter code: H4B) (formula: $C_9H_{15}N_5O_3$).



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

- Molecule 4 is 2-(3-fluorophenyl)-N-{2-[2-(1H-imidazol-1-yl)pyrimidin-4-yl]ethyl}ethanamine (three-letter code: E2Z) (formula: $C_{17}H_{18}FN_5$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	F	N		
4	A	1	23	17	1	5	0	0

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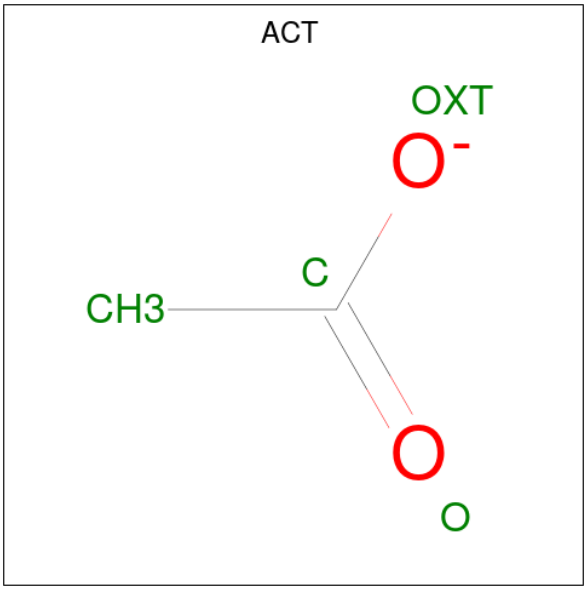
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	F	N		
4	B	1	23	17	1	5	0	0

- Molecule 5 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



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

- Molecule 6 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 7 is water.

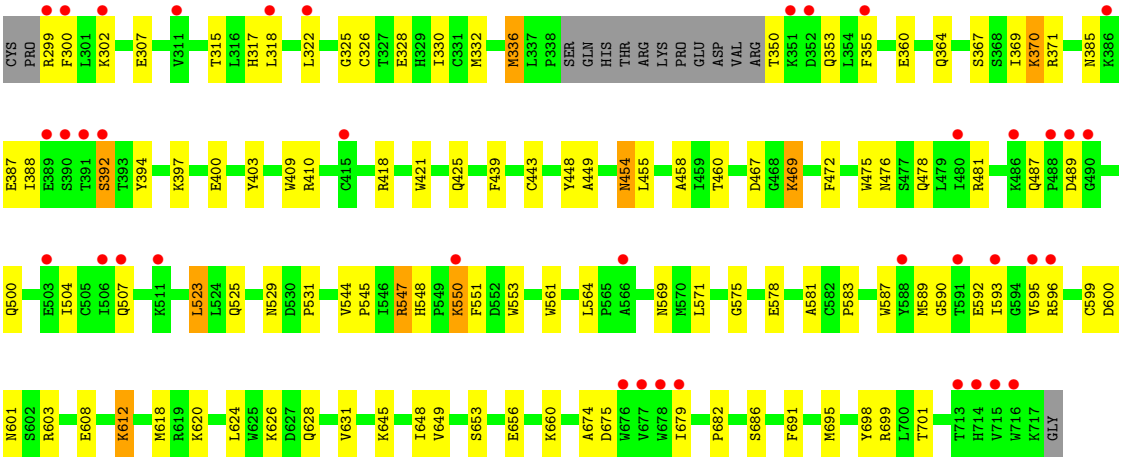
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
7	A	72	72	72	0	0
7	B	80	80	80	0	0

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.

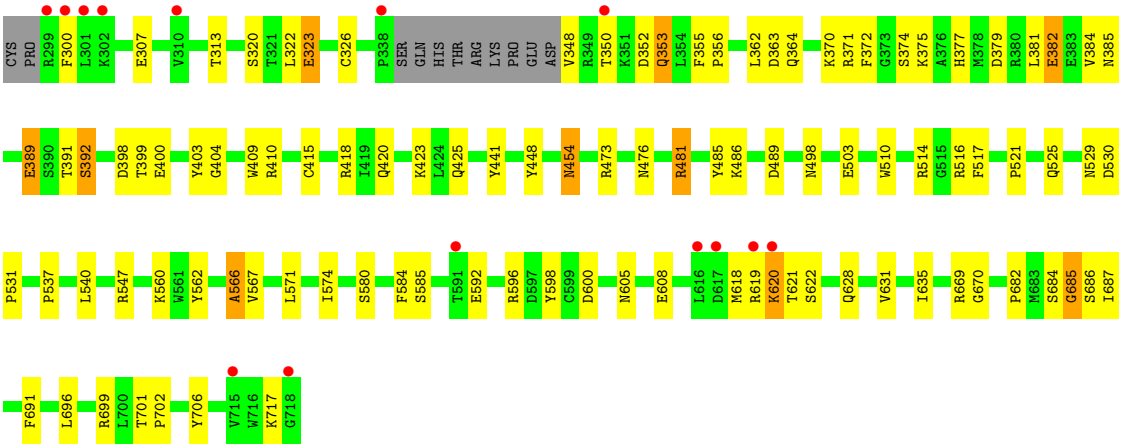
• Molecule 1: NITRIC OXIDE SYNTHASE, BRAIN

Chain A:



• Molecule 1: NITRIC OXIDE SYNTHASE, BRAIN

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	51.84Å 110.25Å 164.32Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	91.72 – 2.13 38.85 – 2.13	Depositor EDS
% Data completeness (in resolution range)	98.6 (91.72-2.13) 98.7 (38.85-2.13)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.62 (at 2.14Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.206 , 0.265 0.206 , 0.264	Depositor DCC
R_{free} test set	2640 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å^2)	40.2	Xtriage
Anisotropy	0.627	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 44.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	7004	wwPDB-VP
Average B, all atoms (Å^2)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.16% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: E2Z, ACT, HEM, ZN, H4B

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	0.89	3/3422 (0.1%)	0.98	7/4643 (0.2%)
1	B	1.05	8/3453 (0.2%)	1.04	9/4681 (0.2%)
All	All	0.97	11/6875 (0.2%)	1.01	16/9324 (0.2%)

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	592	GLU	CG-CD	9.10	1.65	1.51
1	A	596	ARG	CZ-NH2	7.51	1.42	1.33
1	B	400	GLU	CD-OE1	6.23	1.32	1.25
1	B	403	TYR	CG-CD2	6.00	1.47	1.39
1	B	562	TYR	CE2-CZ	5.87	1.46	1.38
1	B	566	ALA	C-O	5.76	1.34	1.23
1	B	592	GLU	CG-CD	5.43	1.60	1.51
1	A	600	ASP	CB-CG	5.41	1.63	1.51
1	B	371	ARG	NE-CZ	5.39	1.40	1.33
1	B	441	TYR	CG-CD1	5.24	1.46	1.39
1	B	670	GLY	N-CA	5.20	1.53	1.46

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	596	ARG	NE-CZ-NH1	-12.49	114.06	120.30
1	A	600	ASP	CB-CG-OD2	8.79	126.21	118.30
1	A	418	ARG	NE-CZ-NH1	-6.90	116.85	120.30
1	B	489	ASP	CB-CG-OD2	-6.81	112.17	118.30
1	A	592	GLU	OE1-CD-OE2	-6.46	115.55	123.30
1	A	410	ARG	NE-CZ-NH2	-6.13	117.23	120.30
1	B	489	ASP	CB-CG-OD1	6.11	123.80	118.30
1	B	481	ARG	NE-CZ-NH1	-6.04	117.28	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	596	ARG	NE-CZ-NH2	5.74	123.17	120.30
1	B	530	ASP	C-N-CD	5.64	140.25	128.40
1	B	592	GLU	OE1-CD-OE2	-5.56	116.63	123.30
1	B	371	ARG	NE-CZ-NH2	-5.39	117.60	120.30
1	A	410	ARG	NE-CZ-NH1	5.32	122.96	120.30
1	B	410	ARG	NE-CZ-NH1	-5.32	117.64	120.30
1	B	669	ARG	NE-CZ-NH2	-5.04	117.78	120.30
1	B	560	LYS	CD-CE-NZ	-5.03	100.12	111.70

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3323	0	3231	69	0
1	B	3354	0	3270	66	0
2	A	43	0	30	4	0
2	B	43	0	30	11	0
3	A	17	0	15	2	0
3	B	17	0	15	0	0
4	A	23	0	18	1	0
4	B	23	0	18	1	0
5	A	4	0	3	1	0
5	B	4	0	3	0	0
6	A	1	0	0	0	0
7	A	72	0	0	0	0
7	B	80	0	0	5	0
All	All	7004	0	6633	135	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 10.

All (135) 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)
2:B:750:HEM:HMC2	2:B:750:HEM:HBC2	1.34	1.08
1:B:596:ARG:NH2	1:B:600:ASP:OD2	1.98	0.95
1:B:537:PRO:HG2	1:B:540:LEU:HD22	1.46	0.95
1:A:336:MET:HE2	3:A:760:H4B:O4	1.75	0.84
1:A:336:MET:CE	3:A:760:H4B:O4	2.28	0.81
1:A:350:THR:N	1:A:353:GLN:HE21	1.83	0.76
1:A:322:LEU:HB2	1:A:699:ARG:HB2	1.73	0.70
1:B:364:GLN:NE2	7:B:2003:HOH:O	2.24	0.69
1:B:391:THR:O	1:B:392:SER:HB2	1.92	0.69
1:A:571:LEU:HD21	1:A:578:GLU:HB3	1.74	0.69
1:A:487:GLN:HB2	1:A:489:ASP:HB3	1.78	0.66
1:A:328:GLU:HB3	1:B:323:GLU:HG2	1.78	0.65
1:B:398:ASP:HB2	7:B:2008:HOH:O	1.95	0.65
2:A:750:HEM:HBC2	2:A:750:HEM:CMC	2.28	0.63
2:A:750:HEM:HBC2	2:A:750:HEM:HMC1	1.81	0.63
1:B:415:CYS:HB2	2:B:750:HEM:C4D	2.34	0.63
1:B:415:CYS:HB2	2:B:750:HEM:ND	2.12	0.63
2:B:750:HEM:HBC2	2:B:750:HEM:CMC	2.17	0.62
1:A:355:PHE:CE1	1:A:385:ASN:HB2	2.34	0.61
1:A:523:LEU:HD22	1:A:531:PRO:HB2	1.82	0.61
1:A:589:MET:HA	1:A:649:VAL:O	2.00	0.61
1:A:475:TRP:HB2	1:A:523:LEU:HB3	1.83	0.61
1:B:510:TRP:CE2	1:B:521:PRO:HD3	2.37	0.60
1:B:355:PHE:CE2	1:B:385:ASN:HB2	2.37	0.59
1:B:353:GLN:O	1:B:356:PRO:HG2	2.02	0.59
1:A:454:ASN:C	1:A:454:ASN:HD22	2.07	0.59
1:B:404:GLY:HA3	1:B:574:ILE:HD13	1.84	0.58
1:B:620:LYS:HE3	1:B:620:LYS:HA	1.86	0.58
1:B:391:THR:O	1:B:392:SER:CB	2.51	0.58
1:B:375:LYS:NZ	1:B:379:ASP:OD2	2.37	0.56
1:B:350:THR:O	1:B:353:GLN:HG2	2.05	0.56
1:A:631:VAL:HG11	1:B:628:GLN:HG3	1.87	0.56
1:A:350:THR:N	1:A:353:GLN:NE2	2.53	0.56
1:A:409:TRP:CE3	1:A:421:TRP:HA	2.42	0.55
1:B:619:ARG:HH11	1:B:619:ARG:HB2	1.71	0.55
1:A:487:GLN:C	1:A:489:ASP:H	2.11	0.54
1:A:325:GLY:O	1:A:332:MET:HG3	2.07	0.53
1:B:701:THR:HB	1:B:702:PRO:HA	1.90	0.53
1:B:619:ARG:HB2	1:B:619:ARG:NH1	2.24	0.53
1:A:388:ILE:O	1:A:392:SER:HA	2.07	0.53
1:A:626:LYS:HB3	1:B:687:ILE:HD12	1.91	0.53
1:B:476:ASN:HB3	7:B:2039:HOH:O	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:608:GLU:HG3	1:B:618:MET:HE1	1.91	0.53
1:B:409:TRP:CE2	2:B:750:HEM:C2C	2.97	0.52
1:A:628:GLN:HG3	1:B:631:VAL:HG11	1.90	0.52
2:B:750:HEM:HHC	2:B:750:HEM:HBB2	1.91	0.52
1:A:388:ILE:O	1:A:392:SER:N	2.43	0.52
1:B:525:GLN:HG3	1:B:529:ASN:O	2.10	0.51
1:A:369:ILE:O	1:A:370:LYS:HE3	2.10	0.51
1:A:603:ARG:HH11	1:A:603:ARG:HG3	1.76	0.51
1:B:374:SER:O	1:B:377:HIS:HB3	2.11	0.50
1:A:525:GLN:HG3	1:A:529:ASN:O	2.12	0.50
1:A:500:GLN:O	1:A:504:ILE:HG13	2.12	0.49
1:A:460:THR:O	1:A:583:PRO:HD2	2.12	0.49
1:B:481:ARG:NE	1:B:498:ASN:HD21	2.11	0.49
1:A:317:HIS:CE1	1:A:318:LEU:HG	2.48	0.49
1:A:350:THR:OG1	1:A:353:GLN:NE2	2.45	0.49
1:A:550:LYS:HA	1:A:550:LYS:NZ	2.28	0.48
1:A:487:GLN:C	1:A:489:ASP:N	2.66	0.48
1:A:439:PHE:CE1	1:A:443:CYS:SG	3.07	0.48
1:B:382:GLU:HA	1:B:382:GLU:OE1	2.12	0.48
1:A:322:LEU:HB3	1:A:699:ARG:HH21	1.78	0.48
2:A:750:HEM:O1D	4:A:800:E2Z:N19	2.46	0.48
1:B:701:THR:HA	1:B:702:PRO:C	2.34	0.48
1:B:385:ASN:O	1:B:389:GLU:HG2	2.14	0.47
1:B:353:GLN:HG2	1:B:353:GLN:H	1.40	0.47
1:A:364:GLN:HG3	1:A:403:TYR:OH	2.14	0.47
1:A:686:SER:HA	1:A:691:PHE:CG	2.49	0.47
1:B:307[B]:GLU:HA	1:B:307[B]:GLU:OE1	2.15	0.47
1:B:485:TYR:CZ	1:B:514:ARG:HA	2.50	0.47
1:A:601:ASN:HB2	1:B:307[B]:GLU:CD	2.35	0.46
1:A:675:ASP:O	1:A:679:ILE:HG12	2.15	0.46
1:A:561:TRP:CD1	1:A:593:ILE:HD13	2.51	0.46
1:A:318:LEU:HD23	1:A:318:LEU:HA	1.74	0.46
1:A:478:GLN:HB2	1:A:481:ARG:HG3	1.98	0.46
1:B:454:ASN:ND2	7:B:2026:HOH:O	2.49	0.46
1:A:387:GLU:OE2	1:A:394:TYR:HA	2.16	0.45
1:A:590:GLY:HA2	1:A:648:ILE:HD11	1.99	0.45
2:B:750:HEM:C1A	4:B:800:E2Z:H02	2.52	0.45
1:A:425:GLN:HG2	1:A:448:TYR:CE2	2.52	0.45
1:B:686:SER:HA	1:B:691:PHE:CG	2.51	0.45
1:A:300:PHE:CD1	1:A:300:PHE:N	2.85	0.44
1:B:596:ARG:HH21	1:B:600:ASP:CG	2.08	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:455:LEU:HD13	1:A:564:LEU:HD12	1.98	0.44
1:A:551:PHE:HB3	1:A:553:TRP:CE2	2.52	0.44
2:A:750:HEM:HMC1	2:A:750:HEM:CBC	2.47	0.44
1:B:706:TYR:OH	2:B:750:HEM:O2D	2.19	0.44
1:B:567:VAL:HB	1:B:584:PHE:CZ	2.52	0.44
1:B:363:ASP:HB3	1:B:372:PHE:HE1	1.82	0.44
1:B:415:CYS:HB2	2:B:750:HEM:C1D	2.52	0.44
2:B:750:HEM:HMC2	2:B:750:HEM:CBC	2.22	0.44
1:B:516:ARG:HG2	1:B:517:PHE:CE2	2.53	0.44
1:A:469:LYS:CE	1:A:469:LYS:HA	2.47	0.44
1:A:472:PHE:O	1:A:581:ALA:HB2	2.18	0.43
1:A:603:ARG:HG3	1:A:603:ARG:NH1	2.33	0.43
1:B:571:LEU:C	1:B:571:LEU:HD23	2.38	0.43
1:B:322:LEU:HB2	1:B:699:ARG:HB2	2.00	0.43
1:B:619:ARG:HH11	1:B:619:ARG:CB	2.30	0.43
1:B:567:VAL:HB	1:B:584:PHE:CE1	2.53	0.43
2:B:750:HEM:HBD1	2:B:750:HEM:HHA	1.99	0.43
1:B:418:ARG:C	1:B:420:GLN:H	2.22	0.43
1:A:674:ALA:HB3	1:A:695:MET:HB3	2.01	0.43
1:A:322:LEU:CB	1:A:699:ARG:HE	2.31	0.42
1:A:449:ALA:O	1:A:455:LEU:HA	2.19	0.42
1:A:682:PRO:HB2	1:B:686:SER:HB3	2.01	0.42
1:B:448:TYR:CD1	1:B:448:TYR:C	2.92	0.42
1:A:612:LYS:HG3	1:A:618:MET:SD	2.59	0.42
1:B:566:ALA:HB2	1:B:585:SER:HB3	2.00	0.42
1:A:624:LEU:HD12	1:B:635:ILE:HG13	2.02	0.42
1:B:473:ARG:HD2	1:B:580:SER:O	2.20	0.42
1:A:326:CYS:HB3	1:B:326:CYS:HB3	2.01	0.42
1:A:317:HIS:HA	1:A:698:TYR:CE2	2.55	0.42
1:A:587:TRP:CZ2	5:A:860:ACT:H1	2.54	0.42
1:B:370:LYS:HZ3	1:B:370:LYS:HG3	1.64	0.42
1:A:322:LEU:HB2	1:A:699:ARG:CB	2.45	0.41
1:A:425:GLN:O	1:A:458:ALA:HA	2.20	0.41
1:A:467:ASP:OD1	1:A:469:LYS:HB2	2.20	0.41
1:A:544:VAL:HA	1:A:545:PRO:HD2	1.82	0.41
1:B:353:GLN:HE21	1:B:353:GLN:HB3	1.68	0.41
1:B:684:SER:O	1:B:685:GLY:C	2.59	0.41
1:B:399:THR:HG23	7:B:2008:HOH:O	2.21	0.41
1:A:548:HIS:CE1	1:A:550:LYS:HB3	2.56	0.41
1:A:561:TRP:CD1	1:A:593:ILE:CD1	3.03	0.41
1:A:397:LYS:HB2	1:A:400:GLU:HG3	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:300:PHE:HB3	1:B:313:THR:CG2	2.50	0.41
1:A:595:VAL:O	1:A:599:CYS:HB2	2.21	0.41
1:B:322:LEU:HD13	1:B:699:ARG:NH2	2.36	0.41
1:B:362:LEU:HD11	1:B:384:VAL:HG21	2.03	0.41
1:B:425:GLN:HG2	1:B:448:TYR:CZ	2.55	0.40
1:A:547:ARG:HD3	1:A:547:ARG:H	1.86	0.40
1:B:598:TYR:O	1:B:605:ASN:N	2.55	0.40
1:A:330:ILE:HD11	1:B:696:LEU:HB3	2.03	0.40
1:B:485:TYR:CE2	1:B:514:ARG:HA	2.57	0.40
1:A:388:ILE:O	1:A:392:SER:CA	2.69	0.40
1:A:686:SER:OG	1:B:682:PRO:HB2	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	406/422 (96%)	382 (94%)	22 (5%)	2 (0%)	29	22
1	B	409/422 (97%)	386 (94%)	21 (5%)	2 (0%)	29	22
All	All	815/844 (97%)	768 (94%)	43 (5%)	4 (0%)	29	22

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	352	ASP
1	A	575	GLY
1	A	392	SER
1	B	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	365/377 (97%)	339 (93%)	26 (7%)	14 9
1	B	368/377 (98%)	350 (95%)	18 (5%)	25 20
All	All	733/754 (97%)	689 (94%)	44 (6%)	19 14

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	299	ARG
1	A	302	LYS
1	A	307[A]	GLU
1	A	307[B]	GLU
1	A	315	THR
1	A	336	MET
1	A	360	GLU
1	A	367	SER
1	A	370	LYS
1	A	371	ARG
1	A	454	ASN
1	A	469	LYS
1	A	476	ASN
1	A	507	GLN
1	A	523	LEU
1	A	547	ARG
1	A	550	LYS
1	A	569	ASN
1	A	608	GLU
1	A	612	LYS
1	A	620	LYS
1	A	645	LYS
1	A	653	SER
1	A	656	GLU
1	A	660	LYS
1	A	701	THR
1	B	320	SER

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Mol	Chain	Res	Type
1	B	323	GLU
1	B	348	VAL
1	B	353	GLN
1	B	381	LEU
1	B	382	GLU
1	B	389	GLU
1	B	392	SER
1	B	423	LYS
1	B	454	ASN
1	B	486	LYS
1	B	503	GLU
1	B	531	PRO
1	B	547	ARG
1	B	620	LYS
1	B	621	THR
1	B	622	SER
1	B	717	LYS

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

Mol	Chain	Res	Type
1	A	317	HIS
1	A	353	GLN
1	A	436	HIS
1	A	454	ASN
1	A	487	GLN
1	A	507	GLN
1	A	569	ASN
1	A	605	ASN
1	A	642	GLN
1	A	697	ASN
1	B	364	GLN
1	B	385	ASN
1	B	454	ASN
1	B	507	GLN
1	B	508	GLN
1	B	605	ASN
1	B	642	GLN
1	B	697	ASN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 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	HEM	A	750	4,1	41,50,50	1.45	6 (14%)	45,82,82	2.05	12 (26%)
3	H4B	A	760	-	16,18,18	1.01	1 (6%)	11,26,26	3.02	6 (54%)
2	HEM	B	750	4,1	41,50,50	1.36	5 (12%)	45,82,82	2.12	13 (28%)
5	ACT	A	860	-	3,3,3	0.78	0	3,3,3	0.84	0
4	E2Z	B	800	2	24,25,25	1.46	3 (12%)	28,32,32	3.09	13 (46%)
3	H4B	B	760	-	16,18,18	1.24	1 (6%)	11,26,26	2.41	4 (36%)
4	E2Z	A	800	2	24,25,25	1.74	8 (33%)	28,32,32	3.00	12 (42%)
5	ACT	B	860	-	3,3,3	0.85	0	3,3,3	0.41	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	HEM	A	750	4,1	-	1/12/54/54	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	H4B	A	760	-	-	0/8/17/17	0/2/2/2
2	HEM	B	750	4,1	-	2/12/54/54	-
4	E2Z	B	800	2	-	1/8/12/12	0/3/3/3
3	H4B	B	760	-	-	2/8/17/17	0/2/2/2
4	E2Z	A	800	2	-	0/8/12/12	0/3/3/3

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	800	E2Z	C12-N13	4.02	1.37	1.32
2	A	750	HEM	C1B-NB	-3.81	1.33	1.40
2	A	750	HEM	C4D-ND	-3.70	1.33	1.40
4	A	800	E2Z	C4'-C3'	3.67	1.44	1.37
4	B	800	E2Z	C04-N03	-3.60	1.34	1.39
4	B	800	E2Z	C16-N11	3.15	1.41	1.34
2	B	750	HEM	C1B-NB	-2.89	1.35	1.40
2	B	750	HEM	C4D-ND	-2.86	1.35	1.40
3	B	760	H4B	C4-N3	2.76	1.37	1.33
2	A	750	HEM	CHB-C1B	2.65	1.41	1.35
3	A	760	H4B	C4-N3	2.48	1.37	1.33
4	B	800	E2Z	C12-N11	2.40	1.35	1.31
4	A	800	E2Z	C2'-C3'	2.40	1.41	1.37
2	A	750	HEM	C3D-C2D	-2.37	1.31	1.36
4	A	800	E2Z	C04-N03	-2.37	1.35	1.39
2	B	750	HEM	C1D-ND	-2.34	1.34	1.38
2	B	750	HEM	CHB-C1B	2.31	1.40	1.35
2	A	750	HEM	FE-NB	2.21	2.07	1.96
2	B	750	HEM	O1D-CGD	2.14	1.29	1.22
2	A	750	HEM	C1D-ND	-2.09	1.34	1.38
4	A	800	E2Z	C02-N03	-2.08	1.34	1.36
4	A	800	E2Z	C6'-C1'	2.07	1.43	1.38
4	A	800	E2Z	C12-N11	2.06	1.34	1.31
4	A	800	E2Z	C5'-C6'	2.04	1.43	1.38

All (60) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	B	800	E2Z	N11-C12-N13	-7.76	117.28	126.08
4	A	800	E2Z	N13-C12-N03	7.31	124.24	114.78
4	B	800	E2Z	C16-N11-C12	7.03	123.36	114.04
4	A	800	E2Z	N11-C12-N13	-6.91	118.24	126.08
3	A	760	H4B	C8A-C4A-C4	6.36	120.22	114.57

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	750	HEM	C1B-NB-C4B	5.61	110.87	105.07
3	B	760	H4B	C8A-C4A-C4	5.60	119.54	114.57
2	B	750	HEM	C1B-NB-C4B	5.59	110.85	105.07
4	B	800	E2Z	C02-N03-C04	5.37	118.24	108.50
4	B	800	E2Z	C04-N03-C12	-5.13	119.71	125.50
4	B	800	E2Z	N11-C12-N03	5.04	122.62	114.81
4	A	800	E2Z	C16-N11-C12	4.94	120.59	114.04
2	A	750	HEM	CHC-C4B-NB	4.93	129.79	124.43
2	B	750	HEM	CHC-C4B-NB	4.92	129.78	124.43
4	A	800	E2Z	C04-N03-C12	-4.80	120.08	125.50
2	A	750	HEM	CHA-C4D-ND	4.38	129.79	124.38
2	B	750	HEM	CHA-C4D-ND	4.34	129.74	124.38
4	A	800	E2Z	C21-C1'-C2'	-4.24	113.58	120.54
4	B	800	E2Z	N13-C12-N03	4.10	120.08	114.78
2	B	750	HEM	CHA-C4D-C3D	-4.01	117.80	125.33
3	A	760	H4B	C4-C4A-N5	3.93	122.42	119.12
4	A	800	E2Z	C02-N03-C04	3.91	115.60	108.50
4	A	800	E2Z	F7'-C3'-C4'	3.73	124.87	118.54
2	B	750	HEM	CHB-C1B-NB	3.69	128.94	124.38
2	B	750	HEM	CHD-C1D-ND	3.52	128.26	124.43
4	B	800	E2Z	C05-C04-N03	-3.36	100.56	106.50
4	B	800	E2Z	C15-C16-N11	-3.34	119.80	123.96
3	A	760	H4B	N1-C2-N3	-3.31	120.23	125.42
2	B	750	HEM	C4C-CHD-C1D	-3.22	118.31	122.56
2	B	750	HEM	CHD-C1D-C2D	-3.21	119.96	124.98
2	A	750	HEM	CMD-C2D-C1D	3.21	129.93	125.04
2	A	750	HEM	CHA-C4D-C3D	-3.19	119.35	125.33
3	A	760	H4B	C2-N1-C8A	3.15	121.61	114.54
2	B	750	HEM	CBA-CAA-C2A	-3.14	107.26	112.62
2	A	750	HEM	CHD-C1D-ND	3.09	127.78	124.43
4	A	800	E2Z	C17-C14-N13	3.02	120.45	115.95
3	A	760	H4B	N2-C2-N3	2.96	121.86	117.25
2	B	750	HEM	O2A-CGA-CBA	2.91	123.39	114.03
2	A	750	HEM	CHD-C1D-C2D	-2.89	120.46	124.98
3	A	760	H4B	C2-N3-C4	2.83	120.43	115.93
2	A	750	HEM	CBD-CAD-C3D	-2.82	104.78	112.63
4	A	800	E2Z	C4'-C3'-C2'	-2.76	119.71	123.29
4	A	800	E2Z	C18-C17-C14	-2.70	102.74	112.55
2	B	750	HEM	O2A-CGA-O1A	-2.64	116.73	123.30
4	B	800	E2Z	C05-N01-C02	2.58	113.94	106.20
2	A	750	HEM	CHB-C1B-NB	2.57	127.56	124.38
2	A	750	HEM	O2D-CGD-CBD	2.49	122.03	114.03

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	B	760	H4B	C2-N3-C4	2.49	119.88	115.93
4	B	800	E2Z	C6'-C1'-C2'	2.48	122.02	118.54
2	B	750	HEM	O2D-CGD-O1D	-2.47	117.15	123.30
3	B	760	H4B	N1-C2-N3	-2.43	121.61	125.42
3	B	760	H4B	C4-C4A-N5	2.41	121.14	119.12
2	A	750	HEM	O2A-CGA-CBA	2.38	121.67	114.03
4	B	800	E2Z	C16-C15-C14	-2.27	114.49	116.62
4	A	800	E2Z	C6'-C1'-C2'	2.25	121.70	118.54
4	A	800	E2Z	C05-C04-N03	-2.14	102.73	106.50
2	B	750	HEM	C2C-C3C-C4C	2.09	108.36	106.90
2	A	750	HEM	C4A-C3A-C2A	2.09	108.45	107.00
4	B	800	E2Z	C02-N03-C12	-2.03	122.04	126.02
4	B	800	E2Z	C21-C1'-C2'	-2.03	117.21	120.54

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	760	H4B	O10-C10-C9-O9
3	B	760	H4B	C11-C10-C9-O9
4	B	800	E2Z	C17-C18-N19-C20
2	B	750	HEM	CAD-CBD-CGD-O2D
2	B	750	HEM	CAD-CBD-CGD-O1D
2	A	750	HEM	CAD-CBD-CGD-O2D

There are no ring outliers.

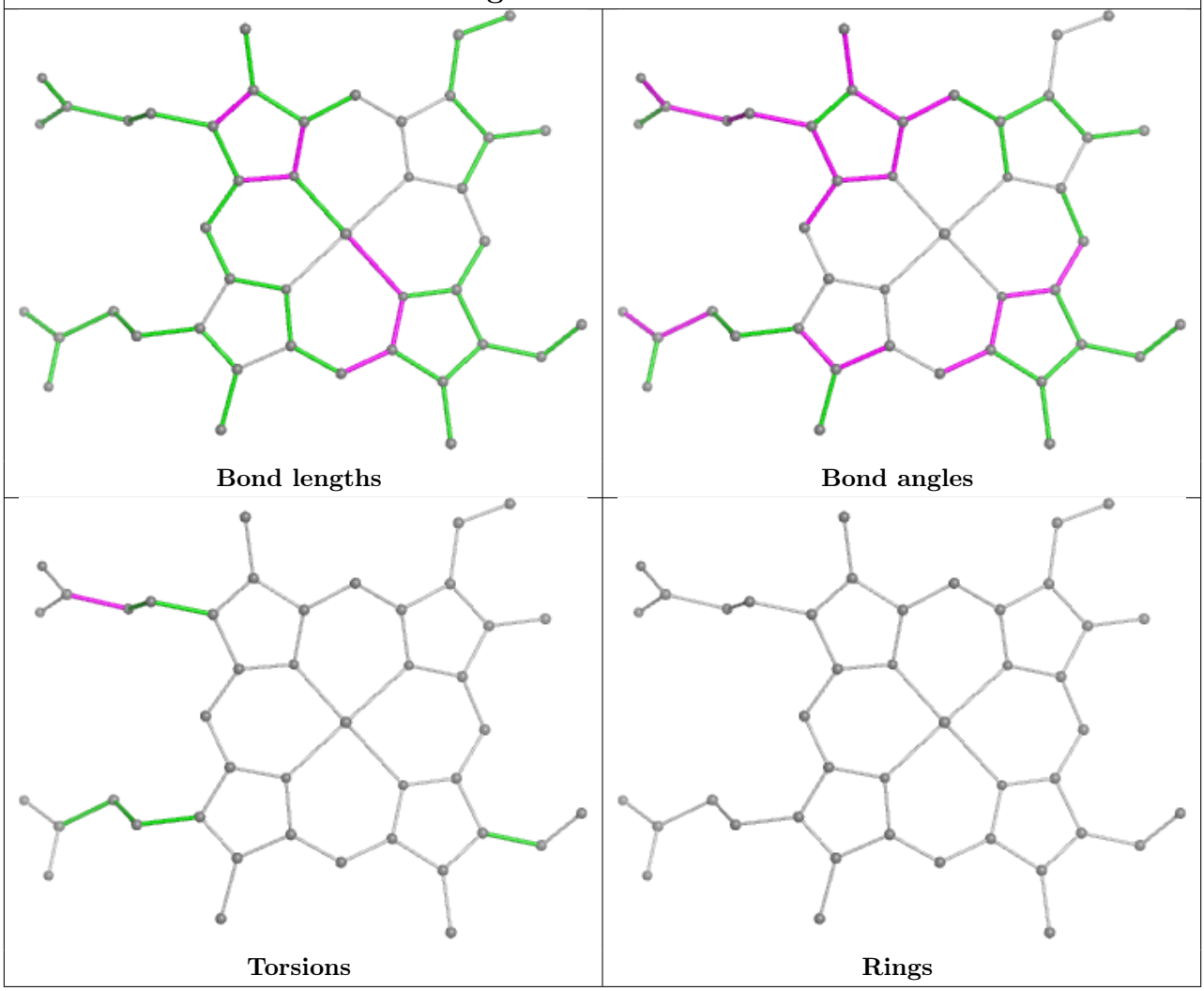
6 monomers are involved in 18 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	750	HEM	4	0
3	A	760	H4B	2	0
2	B	750	HEM	11	0
5	A	860	ACT	1	0
4	B	800	E2Z	1	0
4	A	800	E2Z	1	0

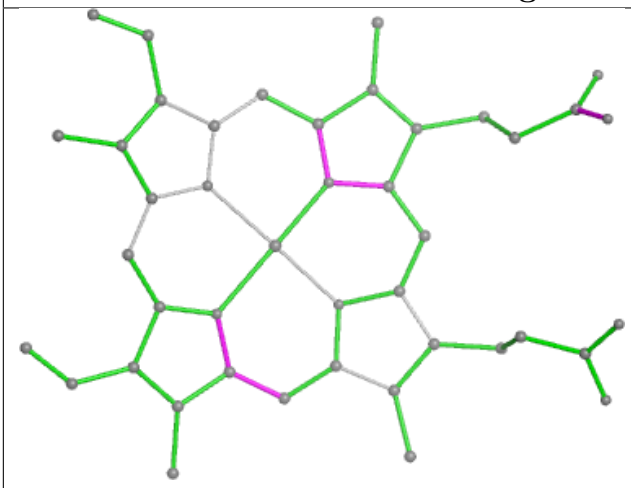
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be

highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

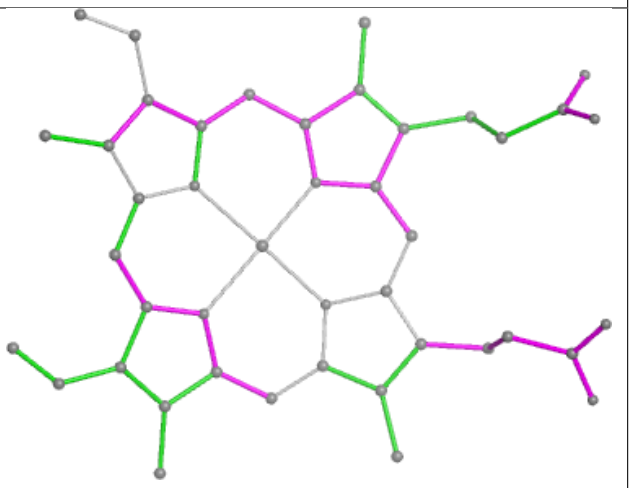
Ligand HEM A 750



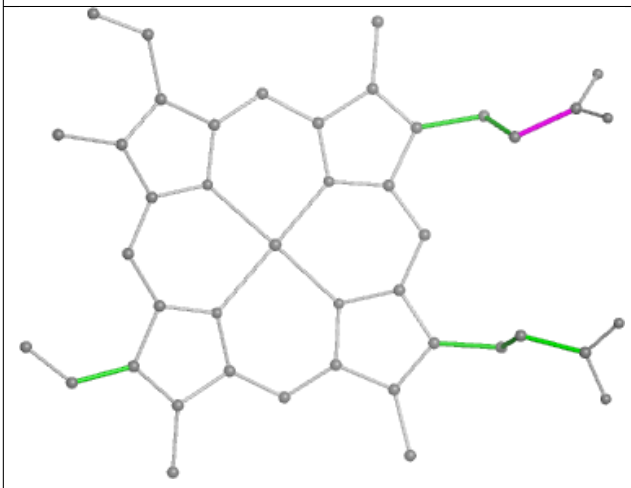
Ligand HEM B 750



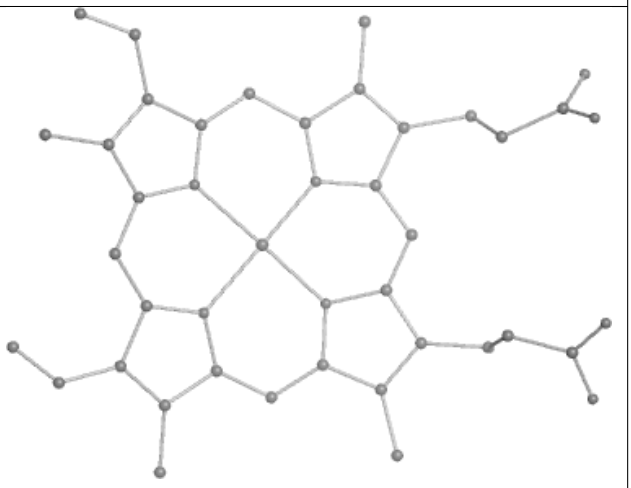
Bond lengths



Bond angles

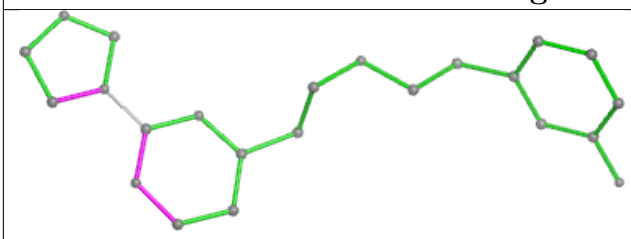


Torsions

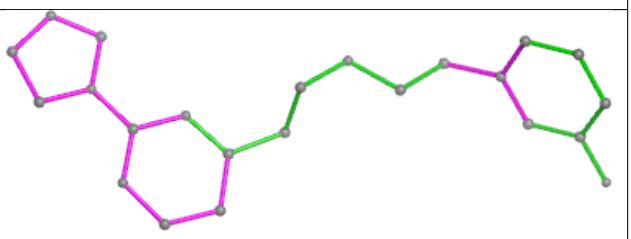


Rings

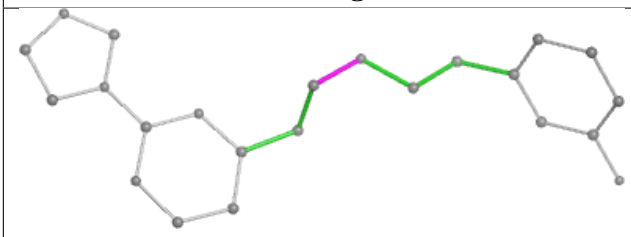
Ligand E2Z B 800



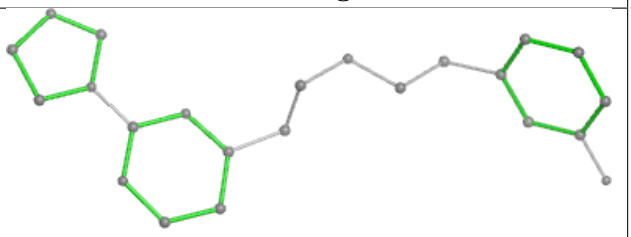
Bond lengths



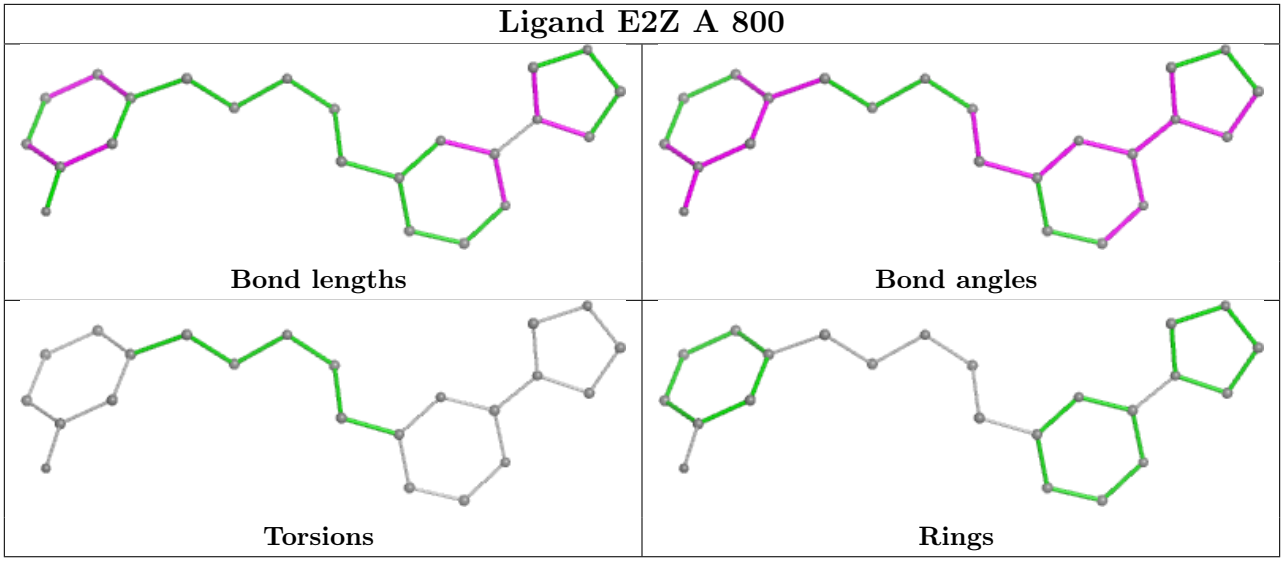
Bond angles



Torsions



Rings



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	408/422 (96%)	0.49	39 (9%)	8	10	32, 53, 82, 105	0
1	B	411/422 (97%)	0.17	14 (3%)	45	52	31, 45, 70, 93	1 (0%)
All	All	819/844 (97%)	0.33	53 (6%)	18	23	31, 49, 79, 105	1 (0%)

All (53) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	300	PHE	7.0
1	B	718	GLY	5.8
1	A	488	PRO	5.4
1	A	299	ARG	4.8
1	A	352	ASP	4.6
1	A	715	VAL	4.5
1	A	716	TRP	4.4
1	B	619	ARG	4.0
1	A	355	PHE	4.0
1	A	322	LEU	3.9
1	A	351	LYS	3.9
1	A	300	PHE	3.5
1	A	713	THR	3.5
1	B	715	VAL	3.5
1	A	489	ASP	3.5
1	B	616	LEU	3.1
1	B	301	LEU	3.1
1	A	311	VAL	2.8
1	A	506	ILE	2.8
1	A	588	TYR	2.7
1	A	392	SER	2.7
1	B	620	LYS	2.7
1	A	511	LYS	2.7
1	A	389	GLU	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	415	CYS	2.6
1	A	566	ALA	2.6
1	A	390	SER	2.6
1	B	350	THR	2.6
1	B	302	LYS	2.6
1	A	318	LEU	2.6
1	A	677	VAL	2.5
1	B	591	THR	2.5
1	A	302	LYS	2.5
1	B	338	PRO	2.5
1	A	595	VAL	2.5
1	A	714	HIS	2.5
1	A	503	GLU	2.4
1	A	386	LYS	2.4
1	A	486	LYS	2.4
1	A	391	THR	2.4
1	B	617	ASP	2.3
1	A	679	ILE	2.3
1	A	591	THR	2.3
1	A	490	GLY	2.3
1	A	678	TRP	2.2
1	B	310	VAL	2.2
1	A	596	ARG	2.1
1	A	507	GLN	2.1
1	B	299	ARG	2.1
1	A	550	LYS	2.0
1	A	480	ILE	2.0
1	A	676	TRP	2.0
1	A	593	ILE	2.0

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

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

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

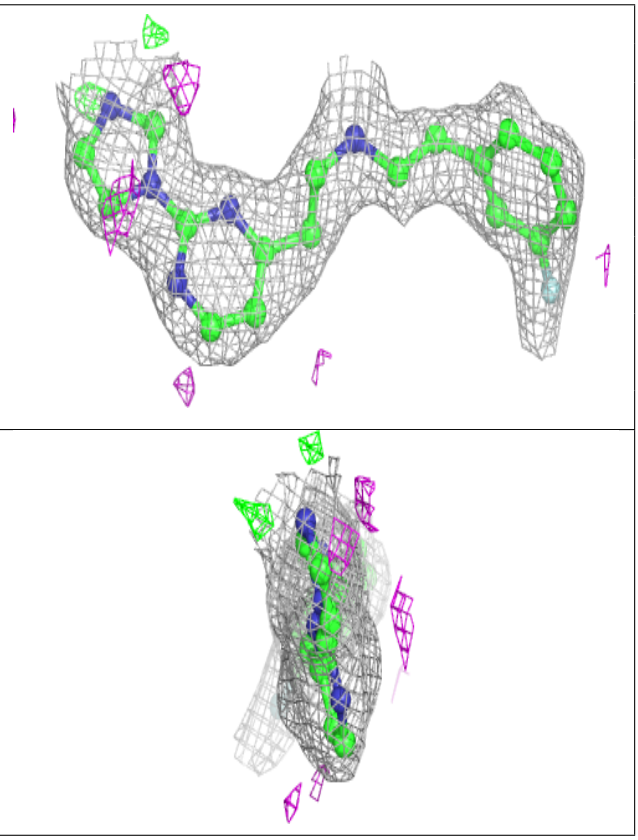
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	H4B	A	760	17/17	0.81	0.29	58,70,87,96	0
5	ACT	B	860	4/4	0.86	0.16	60,67,68,69	0
3	H4B	B	760	17/17	0.90	0.19	50,54,55,55	0
5	ACT	A	860	4/4	0.92	0.30	90,90,91,95	0
4	E2Z	A	800	23/23	0.93	0.22	40,42,65,69	0
4	E2Z	B	800	23/23	0.94	0.20	36,39,63,63	0
2	HEM	A	750	43/43	0.97	0.20	38,40,43,46	0
2	HEM	B	750	43/43	0.98	0.15	32,34,42,45	0
6	ZN	A	900	1/1	0.99	0.11	55,55,55,55	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.

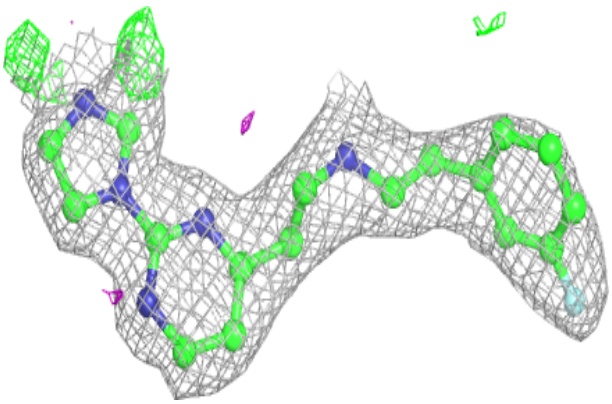
Electron density around E2Z A 800:

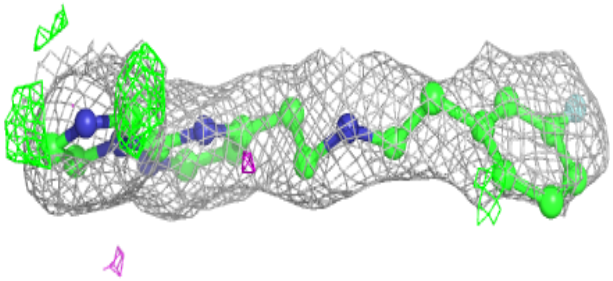
2mF$_o$-DF$_c$ (at 0.7 rmsd) in gray
mF$_o$-DF$_c$ (at 3 rmsd) in purple (negative) and green (positive)

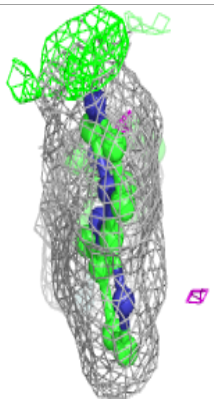


Electron density around E2Z B 800:

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

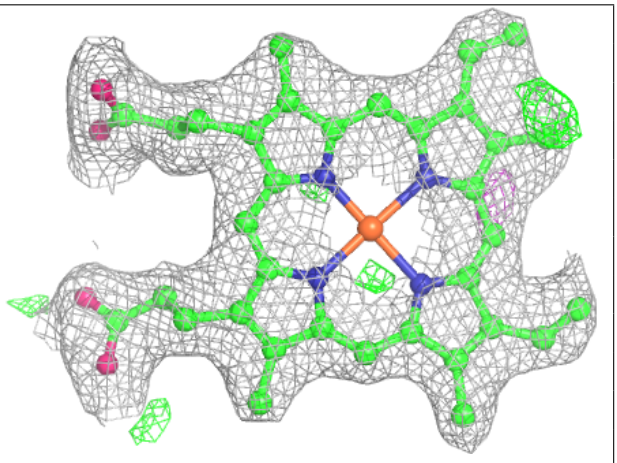


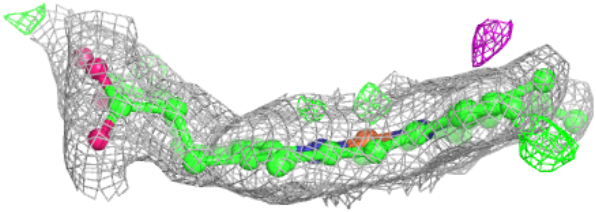


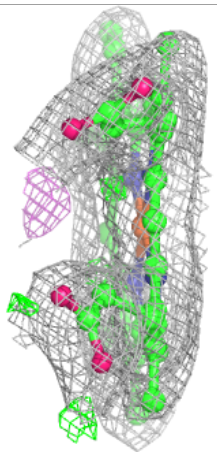


Electron density around HEM A 750:

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

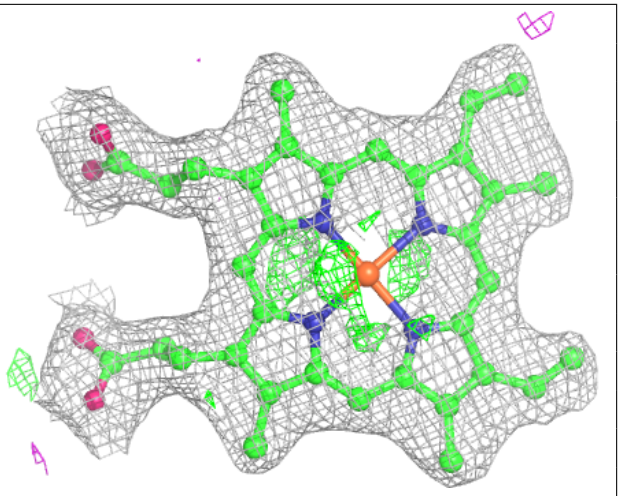


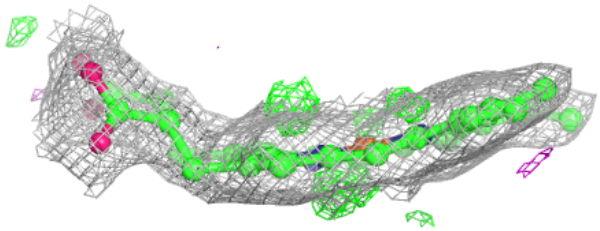


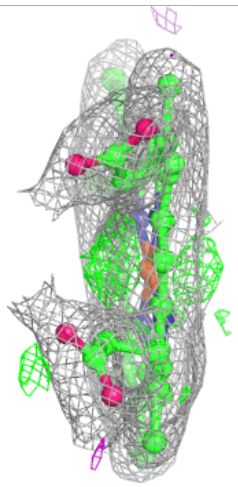


Electron density around HEM B 750:

$2mF_o$-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.