



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 3, 2024 – 11:30 PM EST

PDB ID : 1P6H
Title : Rat neuronal NOS heme domain with L-N(omega)-nitroarginine-2,4-L-diaminobutyric amide bound
Authors : Flinspach, M.L.; Li, H.; Jamal, J.; Yang, W.; Huang, H.; Hah, J.-M.; Gomez-Vidal, J.A.; Litzinger, E.A.; Silverman, R.B.; Poulos, T.L.
Deposited on : 2003-04-29
Resolution : 1.98 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.36
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

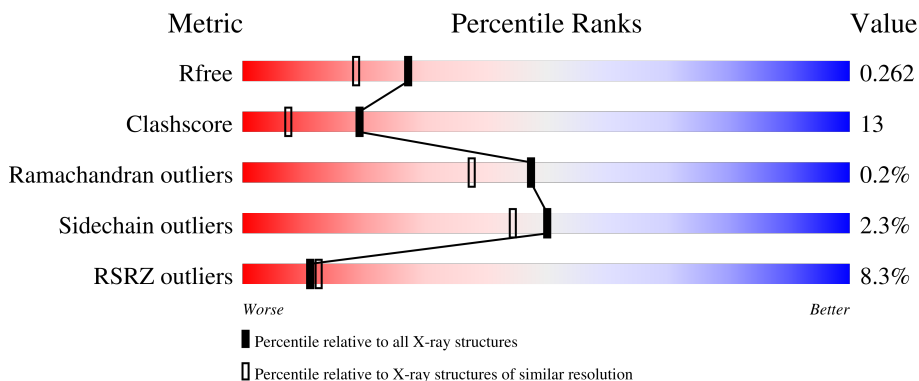
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.98 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	11647 (2.00-1.96)
Clashscore	141614	1014 (1.98-1.98)
Ramachandran outliers	138981	1006 (1.98-1.98)
Sidechain outliers	138945	1006 (1.98-1.98)
RSRZ outliers	127900	11410 (2.00-1.96)

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	421	
1	B	421	

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 7314 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	407	Total 3313	C 2121	N 566	O 605	S 21	0	0	0
1	B	410	Total 3341	C 2138	N 573	O 609	S 21	0	0	0

- Molecule 2 is ACETATE ION (three-letter code: ACT) (formula: C₂H₃O₂).

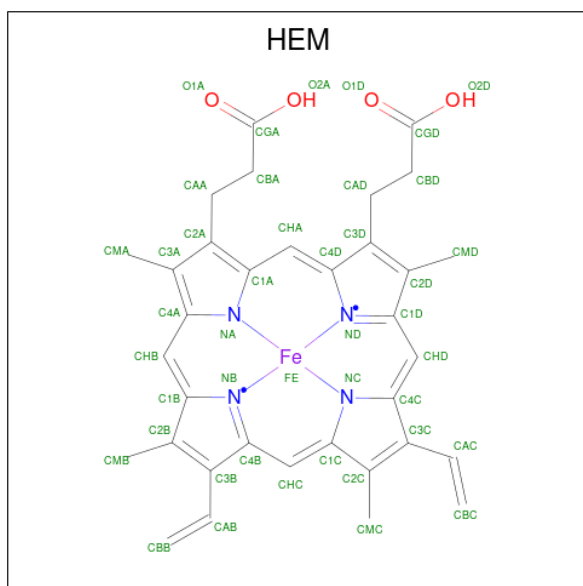


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	A	1	Total 4	C 2	O 2	0	0
2	B	1	Total 4	C 2	O 2	0	0

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

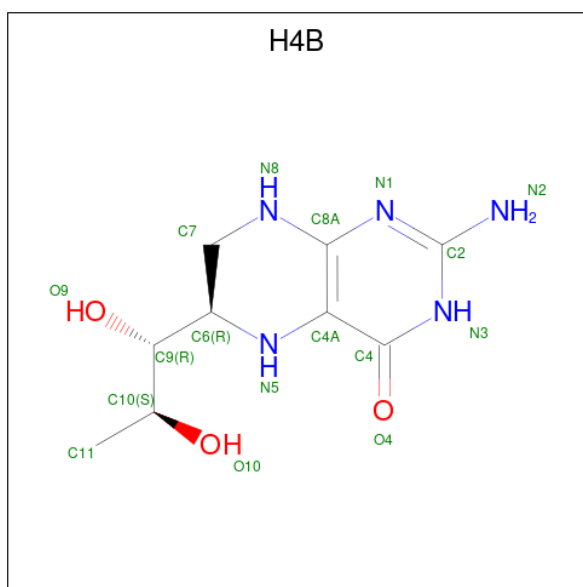
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Zn 1 1	0	0

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



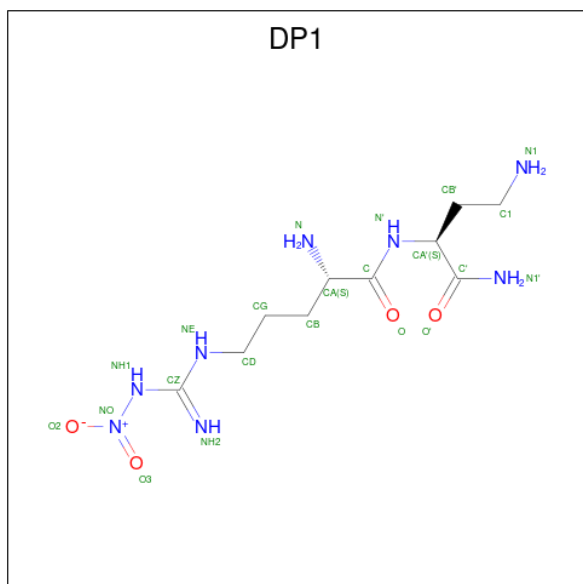
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C Fe N O 43 34 1 4 4	0	0
4	B	1	Total C Fe N O 43 34 1 4 4	0	0

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



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 L-N(OMEGA)-NITROARGININE-2,4-L-DIAMINO BUTYRIC AMIDE (three-letter code: DP1) (formula: $C_{10}H_{22}N_8O_4$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	N	O	0	0
			22	10	8	4		

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
6	B	1	22	10	8	4	0	0

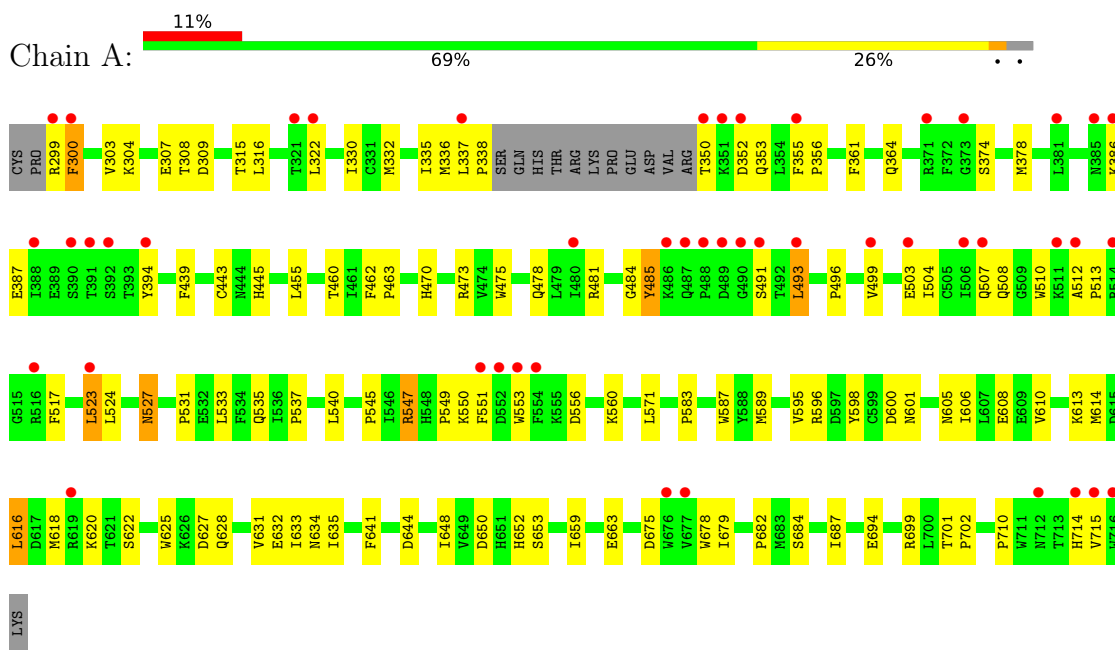
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	214	Total	O	0	0
			214	214		
7	B	273	Total	O	0	0
			273	273		

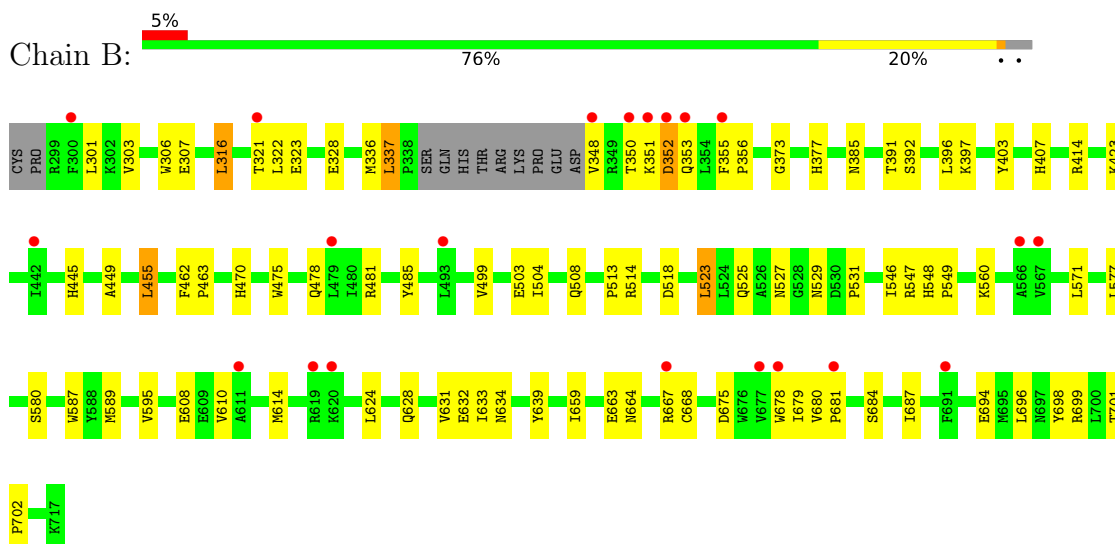
3 Residue-property plots [i](#)

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	51.94Å 110.22Å 164.77Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.94 – 1.98 43.94 – 1.95	Depositor EDS
% Data completeness (in resolution range)	94.9 (43.94-1.98) 95.0 (43.94-1.95)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.25 (at 1.95Å)	Xtrriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.231 , 0.273 0.223 , 0.262	Depositor DCC
R_{free} test set	3322 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	28.1	Xtrriage
Anisotropy	0.723	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 48.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7314	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.55% 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 i

5.1 Standard geometry i

Bond lengths and bond angles in the following residue types are not validated in this section: HEM, ZN, ACT, DP1, 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	0.41	0/3406	0.65	1/4621 (0.0%)
1	B	0.45	0/3434	0.67	1/4656 (0.0%)
All	All	0.43	0/6840	0.66	2/9277 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	589	MET	N-CA-C	-5.87	95.15	111.00
1	A	589	MET	N-CA-C	-5.57	95.96	111.00

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	3313	0	3221	102	0
1	B	3341	0	3256	74	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	3	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	17	0	15	1	0
5	B	17	0	15	0	0
6	A	22	0	21	1	0
6	B	22	0	21	1	0
7	A	214	0	0	13	0
7	B	273	0	0	6	0
All	All	7314	0	6615	173	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 13.

All (173) 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:B:373:GLY:H	1:B:377:HIS:HD2	1.09	0.96
1:B:523:LEU:HD22	1:B:531:PRO:HB2	1.50	0.91
1:A:350:THR:HB	1:A:353:GLN:HG3	1.58	0.85
1:A:523:LEU:HD22	1:A:531:PRO:HB2	1.60	0.83
1:B:373:GLY:H	1:B:377:HIS:CD2	1.98	0.81
1:A:613:LYS:HE2	7:A:977:HOH:O	1.82	0.78
1:B:668:CYS:HB3	7:B:1012:HOH:O	1.87	0.73
1:A:473:ARG:NH2	1:A:710:PRO:HD3	2.05	0.72
1:A:616:LEU:HD13	1:A:625:TRP:HB2	1.72	0.71
1:A:353:GLN:O	1:A:356:PRO:HD2	1.94	0.68
1:B:350:THR:HB	1:B:353:GLN:OE1	1.95	0.66
1:A:350:THR:HG22	1:A:352:ASP:H	1.60	0.66
1:A:307:GLU:HG3	7:B:904:HOH:O	1.95	0.65
1:A:455:LEU:HD12	1:A:587:TRP:HB3	1.79	0.65
1:A:545:PRO:HD3	7:A:1050:HOH:O	1.96	0.64
1:B:337:LEU:N	1:B:337:LEU:HD23	2.13	0.64
1:B:684:SER:HB3	1:B:687:ILE:HD11	1.79	0.63
1:A:304:LYS:O	1:A:694:GLU:HG3	1.99	0.62
4:B:750:HEM:HMC1	4:B:750:HEM:HBC2	1.81	0.62
1:A:508:GLN:HG3	7:A:1025:HOH:O	2.00	0.61
1:B:355:PHE:CE1	1:B:385:ASN:HB2	2.36	0.61
1:A:332:MET:HE1	1:B:301:LEU:HD22	1.83	0.60
1:A:517:PHE:HB2	1:A:560:LYS:HE2	1.84	0.60
1:B:322:LEU:HD13	1:B:699:ARG:HH21	1.67	0.60
1:B:571:LEU:HD12	1:B:580:SER:HA	1.83	0.59
1:A:507:GLN:O	1:A:507:GLN:HG2	2.03	0.59
1:A:598:TYR:O	1:A:606:ILE:HG12	2.02	0.58

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:350:THR:HG22	1:B:352:ASP:H	1.68	0.58
1:A:682:PRO:HG3	7:A:903:HOH:O	2.02	0.58
1:A:303:VAL:HG13	1:A:694:GLU:HB2	1.85	0.58
1:A:610:VAL:O	1:A:614:MET:HG3	2.03	0.58
1:A:595:VAL:HG23	1:A:634:ASN:HD21	1.68	0.57
1:B:391:THR:O	1:B:392:SER:HB2	2.05	0.57
1:A:595:VAL:HG11	1:A:682:PRO:HB2	1.85	0.57
7:A:1001:HOH:O	1:B:307:GLU:HG3	2.05	0.56
1:B:455:LEU:HD23	1:B:455:LEU:N	2.19	0.56
1:A:684:SER:HB3	1:A:687:ILE:HD11	1.86	0.56
1:A:491:SER:HB2	7:A:1102:HOH:O	2.05	0.56
1:A:300:PHE:CD2	1:A:315:THR:HG22	2.41	0.56
1:A:493:LEU:HA	7:A:950:HOH:O	2.05	0.55
1:A:684:SER:HB3	1:A:687:ILE:CG1	2.37	0.55
1:A:556:ASP:HB2	7:A:977:HOH:O	2.05	0.55
4:B:750:HEM:HBC2	4:B:750:HEM:CMC	2.37	0.55
1:B:352:ASP:HB3	1:B:353:GLN:OE1	2.07	0.55
4:A:750:HEM:HMC2	4:A:750:HEM:HBC2	1.87	0.55
1:B:684:SER:HB3	1:B:687:ILE:CG1	2.38	0.54
1:B:355:PHE:N	1:B:356:PRO:HD2	2.22	0.54
1:A:455:LEU:HD12	1:A:587:TRP:CB	2.38	0.54
1:B:316:LEU:HB3	1:B:698:TYR:OH	2.06	0.54
1:A:610:VAL:HG21	1:A:633:ILE:HD11	1.90	0.53
1:A:620:LYS:HE3	1:A:622:SER:OG	2.07	0.53
1:A:499:VAL:O	1:A:503:GLU:HG2	2.09	0.53
1:A:524:LEU:O	1:A:531:PRO:HA	2.08	0.53
1:B:350:THR:HB	1:B:352:ASP:HB3	1.91	0.53
1:B:513:PRO:HG2	1:B:518:ASP:CG	2.29	0.53
1:A:350:THR:HB	1:A:353:GLN:CG	2.36	0.52
1:A:485:TYR:CE2	1:A:512:ALA:HB1	2.45	0.52
1:A:571:LEU:C	1:A:571:LEU:HD23	2.30	0.52
1:B:481:ARG:HG2	7:B:984:HOH:O	2.09	0.52
1:A:478:GLN:HG3	6:A:790:DP1:O'	2.09	0.52
1:B:470:HIS:HB3	1:B:527:ASN:ND2	2.24	0.52
1:A:475:TRP:HB2	1:A:523:LEU:HB3	1.92	0.52
1:B:478:GLN:HB2	1:B:481:ARG:HG3	1.92	0.52
1:A:361:PHE:O	1:A:364:GLN:HG2	2.10	0.52
1:A:545:PRO:HG2	1:A:547:ARG:NH1	2.25	0.52
1:B:549:PRO:HG3	1:B:639:TYR:CG	2.45	0.51
1:B:475:TRP:HB2	1:B:523:LEU:HB3	1.92	0.51
1:B:323:GLU:O	1:B:699:ARG:HD3	2.11	0.51

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:373:GLY:N	1:B:377:HIS:HD2	1.93	0.51
1:A:631:VAL:HG11	1:B:628:GLN:HG2	1.94	0.50
1:A:714:HIS:ND1	7:A:1114:HOH:O	2.34	0.50
1:B:608:GLU:HB2	7:B:1062:HOH:O	2.12	0.50
1:B:546:ILE:HG12	1:B:560:LYS:HA	1.94	0.49
1:A:535:GLN:NE2	7:A:956:HOH:O	2.46	0.49
2:A:860:ACT:H3	4:A:750:HEM:HMB3	1.95	0.49
1:B:321:THR:HG23	1:B:322:LEU:N	2.27	0.49
1:B:659:ILE:O	1:B:663:GLU:HG3	2.12	0.49
1:B:414:ARG:HD3	1:B:678:TRP:CD2	2.47	0.49
1:A:659:ILE:O	1:A:663:GLU:HG3	2.13	0.49
1:A:715:VAL:O	1:A:715:VAL:HG23	2.12	0.49
1:A:300:PHE:CD1	1:A:300:PHE:N	2.80	0.49
1:A:387:GLU:OE2	1:A:394:TYR:HA	2.13	0.48
1:B:403:TYR:O	1:B:407:HIS:HD2	1.96	0.48
1:A:523:LEU:CD2	1:A:531:PRO:HB2	2.39	0.48
1:B:462:PHE:HB3	1:B:463:PRO:CD	2.43	0.48
1:B:595:VAL:HG23	1:B:634:ASN:HD21	1.79	0.48
1:B:684:SER:HB3	1:B:687:ILE:CD1	2.42	0.48
1:B:664:ASN:O	1:B:667:ARG:HG2	2.14	0.48
1:A:620:LYS:HB3	7:A:1020:HOH:O	2.14	0.47
1:A:601:ASN:HB3	7:A:1023:HOH:O	2.13	0.47
1:B:351:LYS:HE2	1:B:392:SER:OG	2.15	0.47
1:B:485:TYR:CZ	1:B:514:ARG:HA	2.50	0.47
1:A:308:THR:O	1:A:309:ASP:HB2	2.15	0.47
1:A:653:SER:HB2	7:B:1073:HOH:O	2.14	0.47
1:A:551:PHE:HE1	1:A:614:MET:HE3	1.79	0.47
1:A:439:PHE:CZ	1:A:443:CYS:SG	3.08	0.47
1:A:330:ILE:HD11	1:B:696:LEU:HB3	1.96	0.46
1:A:545:PRO:HD2	1:A:644:ASP:OD2	2.15	0.46
1:A:299:ARG:CG	1:A:300:PHE:N	2.78	0.46
1:B:614:MET:CE	1:B:632:GLU:HG3	2.45	0.46
1:A:478:GLN:HB2	1:A:481:ARG:HG3	1.96	0.46
1:B:667:ARG:HG3	7:B:1012:HOH:O	2.15	0.46
1:B:322:LEU:HB2	1:B:699:ARG:HB2	1.97	0.46
1:A:684:SER:HB3	1:A:687:ILE:CD1	2.46	0.45
1:B:610:VAL:HG21	1:B:633:ILE:HD11	1.97	0.45
1:A:523:LEU:HG	1:A:533:LEU:HD23	1.97	0.45
1:A:460:THR:O	1:A:583:PRO:HD2	2.16	0.45
1:B:675:ASP:O	1:B:679:ILE:HG12	2.16	0.45
1:A:338:PRO:HB3	7:A:1051:HOH:O	2.17	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:860:ACT:H1	4:A:750:HEM:HAB	1.98	0.45
1:B:610:VAL:O	1:B:614:MET:HG3	2.17	0.45
1:A:335:ILE:HD13	1:B:694:GLU:HB3	1.99	0.44
1:A:504:ILE:O	1:A:508:GLN:HB2	2.17	0.44
1:A:374:SER:O	1:A:378:MET:HG2	2.17	0.44
1:A:496:PRO:HA	1:A:499:VAL:HG23	1.97	0.44
1:B:504:ILE:O	1:B:508:GLN:HG2	2.18	0.44
1:A:332:MET:HE2	1:B:696:LEU:HD21	1.97	0.44
1:A:614:MET:HE3	1:A:632:GLU:HG3	1.98	0.44
1:A:675:ASP:O	1:A:679:ILE:HG12	2.17	0.44
1:A:336:MET:HE2	1:B:306:TRP:CG	2.53	0.44
1:B:337:LEU:N	1:B:337:LEU:CD2	2.80	0.44
1:B:396:LEU:HG	1:B:577:LEU:HD12	2.00	0.44
1:A:537:PRO:HB2	1:A:540:LEU:HG	2.00	0.44
1:B:336:MET:HB3	1:B:337:LEU:HD23	1.99	0.44
1:B:548:HIS:ND1	1:B:549:PRO:HD2	2.32	0.44
1:A:330:ILE:HD11	1:B:696:LEU:HD22	1.99	0.44
4:A:750:HEM:HBC2	4:A:750:HEM:CMC	2.47	0.44
1:B:336:MET:HB3	1:B:337:LEU:CD2	2.47	0.44
1:B:445:HIS:CD2	1:B:445:HIS:C	2.91	0.44
1:B:462:PHE:HB3	1:B:463:PRO:HD2	1.99	0.44
1:A:303:VAL:CG1	1:A:694:GLU:O	2.66	0.44
1:B:587:TRP:H	4:B:750:HEM:HAB	1.83	0.44
1:A:484:GLY:O	1:A:499:VAL:HA	2.18	0.43
1:B:350:THR:C	1:B:352:ASP:N	2.70	0.43
1:A:322:LEU:HB2	1:A:699:ARG:HB2	2.00	0.43
1:A:551:PHE:HB3	1:A:553:TRP:NE1	2.34	0.43
1:A:549:PRO:O	1:A:550:LYS:HG3	2.18	0.43
1:A:618:MET:HA	1:A:625:TRP:CD1	2.53	0.43
1:A:627:ASP:O	1:A:631:VAL:HG23	2.19	0.43
1:B:321:THR:HG23	1:B:322:LEU:H	1.83	0.43
1:A:678:TRP:HA	5:A:760:H4B:N1	2.33	0.43
1:A:470:HIS:HB3	1:A:527:ASN:OD1	2.19	0.43
1:A:596:ARG:O	1:A:600:ASP:HB2	2.19	0.43
1:B:499:VAL:O	1:B:503:GLU:HG3	2.19	0.43
1:A:510:TRP:HB2	1:A:533:LEU:HD13	2.01	0.43
1:B:348:VAL:HG22	1:B:348:VAL:O	2.19	0.42
1:B:701:THR:HA	1:B:702:PRO:C	2.39	0.42
1:A:496:PRO:O	1:A:499:VAL:HG23	2.19	0.42
1:B:525:GLN:HG3	1:B:529:ASN:O	2.19	0.42
1:A:635:ILE:HG13	1:B:624:LEU:HD12	2.01	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:478:GLN:HG3	6:B:791:DP1:O'	2.20	0.42
1:A:701:THR:HA	1:A:702:PRO:C	2.40	0.42
1:A:355:PHE:N	1:A:356:PRO:CD	2.83	0.42
1:A:462:PHE:HB3	1:A:463:PRO:CD	2.50	0.42
1:A:512:ALA:HA	1:A:513:PRO:HD3	1.85	0.41
1:A:551:PHE:CD2	1:A:551:PHE:N	2.86	0.41
1:A:485:TYR:HE2	1:A:512:ALA:HB1	1.84	0.41
1:A:595:VAL:HG23	1:A:634:ASN:ND2	2.34	0.41
1:B:449:ALA:O	1:B:455:LEU:HA	2.21	0.41
1:A:386:LYS:HD3	1:A:386:LYS:HA	1.75	0.41
1:A:299:ARG:HG2	1:A:300:PHE:N	2.35	0.41
1:A:523:LEU:HG	1:A:533:LEU:CD2	2.50	0.41
1:A:608:GLU:HG2	1:A:618:MET:CE	2.50	0.41
1:A:628:GLN:CG	1:B:631:VAL:HG11	2.51	0.41
1:A:614:MET:CE	1:A:632:GLU:HG3	2.50	0.41
1:B:328:GLU:H	1:B:328:GLU:CD	2.25	0.41
1:A:641:PHE:CG	1:A:648:ILE:HD12	2.55	0.40
1:A:445:HIS:CD2	1:A:445:HIS:C	2.95	0.40
1:A:608:GLU:HG2	1:A:618:MET:HE3	2.04	0.40
1:B:548:HIS:HA	1:B:549:PRO:HD3	1.93	0.40
1:B:680:VAL:HA	1:B:681:PRO:HD3	1.94	0.40
1:A:650:ASP:OD2	1:A:652:HIS:HB2	2.21	0.40
1:A:684:SER:HB3	1:A:687:ILE:HG12	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	403/421 (96%)	371 (92%)	31 (8%)	1 (0%)	47 38
1	B	406/421 (96%)	393 (97%)	12 (3%)	1 (0%)	47 38

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	809/842 (96%)	764 (94%)	43 (5%)	2 (0%)	47 38

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	605	ASN
1	B	352	ASP

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	363/377 (96%)	354 (98%)	9 (2%)	47 39
1	B	366/377 (97%)	358 (98%)	8 (2%)	52 46
All	All	729/754 (97%)	712 (98%)	17 (2%)	50 44

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	300	PHE
1	A	316	LEU
1	A	337	LEU
1	A	485	TYR
1	A	493	LEU
1	A	523	LEU
1	A	527	ASN
1	A	547	ARG
1	A	616	LEU
1	B	303	VAL
1	B	316	LEU
1	B	337	LEU
1	B	397	LYS
1	B	423	LYS
1	B	455	LEU
1	B	523	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	547	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	353	GLN
1	A	451	ASN
1	A	454	ASN
1	A	500	GLN
1	A	535	GLN
1	A	634	ASN
1	A	697	ASN
1	B	364	GLN
1	B	377	HIS
1	B	407	HIS
1	B	425	GLN
1	B	451	ASN
1	B	454	ASN
1	B	507	GLN
1	B	527	ASN
1	B	535	GLN
1	B	634	ASN
1	B	664	ASN
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

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	0.87	0	3,3,3	0.82	0
2	ACT	B	861	-	3,3,3	0.91	0	3,3,3	0.70	0
4	HEM	A	750	1	41,50,50	1.07	2 (4%)	45,82,82	1.29	5 (11%)
5	H4B	A	760	-	16,18,18	2.26	3 (18%)	11,26,26	4.15	8 (72%)
4	HEM	B	750	1	41,50,50	1.45	6 (14%)	45,82,82	1.26	2 (4%)
6	DP1	B	791	-	17,21,21	0.99	1 (5%)	20,26,26	0.96	0
6	DP1	A	790	-	17,21,21	1.03	1 (5%)	20,26,26	0.86	0
5	H4B	B	761	-	16,18,18	2.41	5 (31%)	11,26,26	4.33	9 (81%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	HEM	A	750	1	-	0/12/54/54	-
5	H4B	A	760	-	-	0/8/17/17	0/2/2/2
4	HEM	B	750	1	-	2/12/54/54	-
6	DP1	B	791	-	-	6/23/26/26	-
6	DP1	A	790	-	-	7/23/26/26	-
5	H4B	B	761	-	-	0/8/17/17	0/2/2/2

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	760	H4B	C4-N3	5.88	1.43	1.33
5	B	761	H4B	C4-N3	5.46	1.42	1.33
5	B	761	H4B	C6-N5	4.85	1.55	1.45
5	B	761	H4B	C4A-N5	4.46	1.47	1.38

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	760	H4B	C6-N5	4.41	1.54	1.45
5	A	760	H4B	C4A-N5	4.34	1.46	1.38
4	B	750	HEM	C4A-NA	3.40	1.43	1.36
4	B	750	HEM	CHB-C1B	3.24	1.43	1.35
4	B	750	HEM	C3C-CAC	-3.18	1.41	1.47
4	B	750	HEM	CAB-C3B	-3.00	1.39	1.47
6	A	790	DP1	CZ-NH1	-2.80	1.27	1.35
4	A	750	HEM	C3C-CAC	-2.72	1.42	1.47
6	B	791	DP1	CZ-NH1	-2.69	1.28	1.35
5	B	761	H4B	C8A-N1	2.69	1.39	1.34
4	B	750	HEM	CMD-C2D	2.53	1.56	1.50
5	B	761	H4B	C7-N8	2.51	1.49	1.44
4	A	750	HEM	C1A-NA	2.33	1.41	1.36
4	B	750	HEM	C3C-C2C	-2.04	1.37	1.40

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	761	H4B	C8A-C4A-C4	9.15	122.70	114.57
5	A	760	H4B	C8A-C4A-C4	8.48	122.10	114.57
5	B	761	H4B	C2-N3-C4	5.87	125.25	115.93
5	A	760	H4B	C2-N3-C4	5.50	124.66	115.93
5	A	760	H4B	N1-C2-N3	-5.10	117.42	125.42
5	B	761	H4B	N1-C2-N3	-4.69	118.06	125.42
5	A	760	H4B	C2-N1-C8A	3.97	123.44	114.54
5	B	761	H4B	C4A-C4-N3	-3.68	113.54	124.01
5	A	760	H4B	C4A-C4-N3	-3.51	114.05	124.01
5	A	760	H4B	N2-C2-N1	3.43	122.58	117.25
5	B	761	H4B	C4-C4A-N5	-3.42	116.25	119.12
5	A	760	H4B	C4A-N5-C6	-3.27	112.26	121.16
5	B	761	H4B	C2-N1-C8A	3.27	121.86	114.54
5	B	761	H4B	N2-C2-N1	3.24	122.30	117.25
4	B	750	HEM	CBA-CAA-C2A	-3.18	107.19	112.62
5	B	761	H4B	C4A-N5-C6	-3.18	112.50	121.16
4	A	750	HEM	C4B-C3B-C2B	-2.94	104.78	107.11
4	B	750	HEM	C4A-C3A-C2A	-2.76	105.07	107.00
5	B	761	H4B	O9-C9-C6	2.67	115.37	108.98
4	A	750	HEM	CMC-C2C-C3C	2.49	129.34	124.68
4	A	750	HEM	C3B-C2B-C1B	2.31	108.20	106.49
4	A	750	HEM	CAB-C3B-C2B	-2.25	121.19	128.60
5	A	760	H4B	O9-C9-C6	2.21	114.25	108.98
4	A	750	HEM	CHC-C4B-NB	-2.06	122.20	124.43

There are no chirality outliers.

All (15) torsion outliers are listed below:

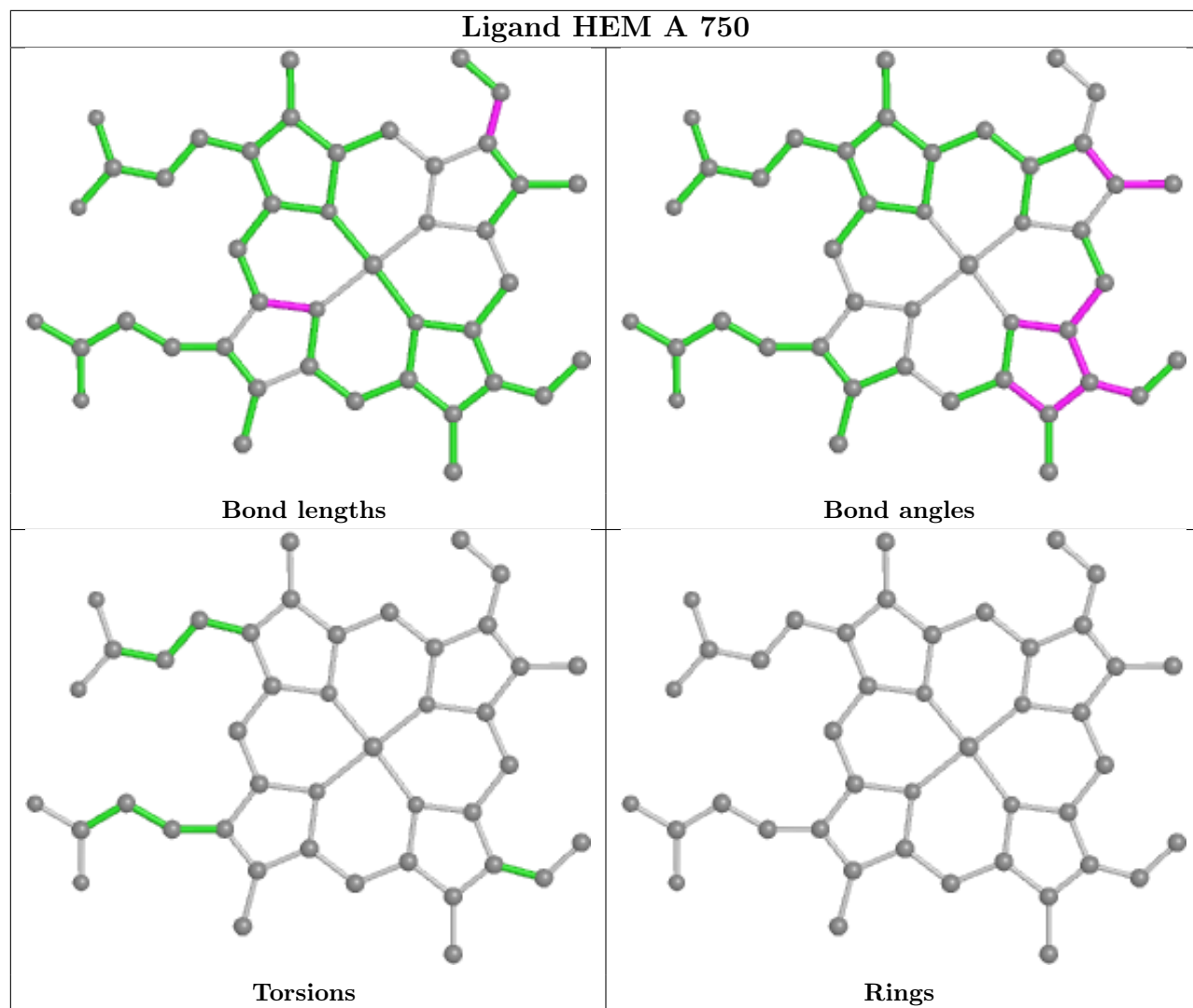
Mol	Chain	Res	Type	Atoms
4	B	750	HEM	C4B-C3B-CAB-CBB
6	A	790	DP1	O'-C'-CA'-CB'
6	A	790	DP1	N1'-C'-CA'-CB'
6	B	791	DP1	N-CA-CB-CG
6	B	791	DP1	N1'-C'-CA'-CB'
6	B	791	DP1	C-CA-CB-CG
6	B	791	DP1	O'-C'-CA'-CB'
4	B	750	HEM	C2B-C3B-CAB-CBB
6	A	790	DP1	O'-C'-CA'-N'
6	A	790	DP1	N1'-C'-CA'-N'
6	B	791	DP1	O'-C'-CA'-N'
6	B	791	DP1	N1'-C'-CA'-N'
6	A	790	DP1	C-CA-CB-CG
6	A	790	DP1	N-CA-CB-CG
6	A	790	DP1	N1-C1-CB'-CA'

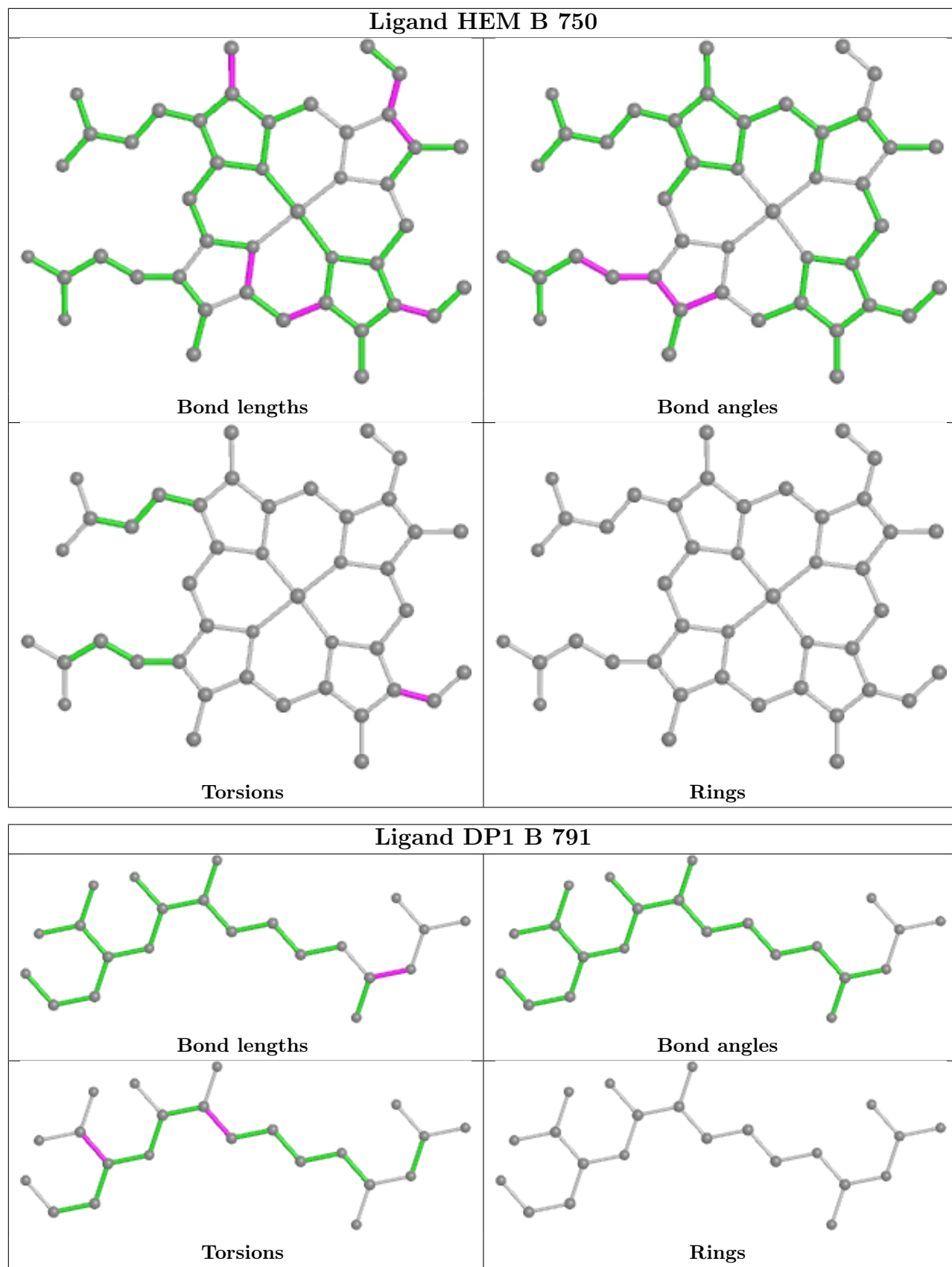
There are no ring outliers.

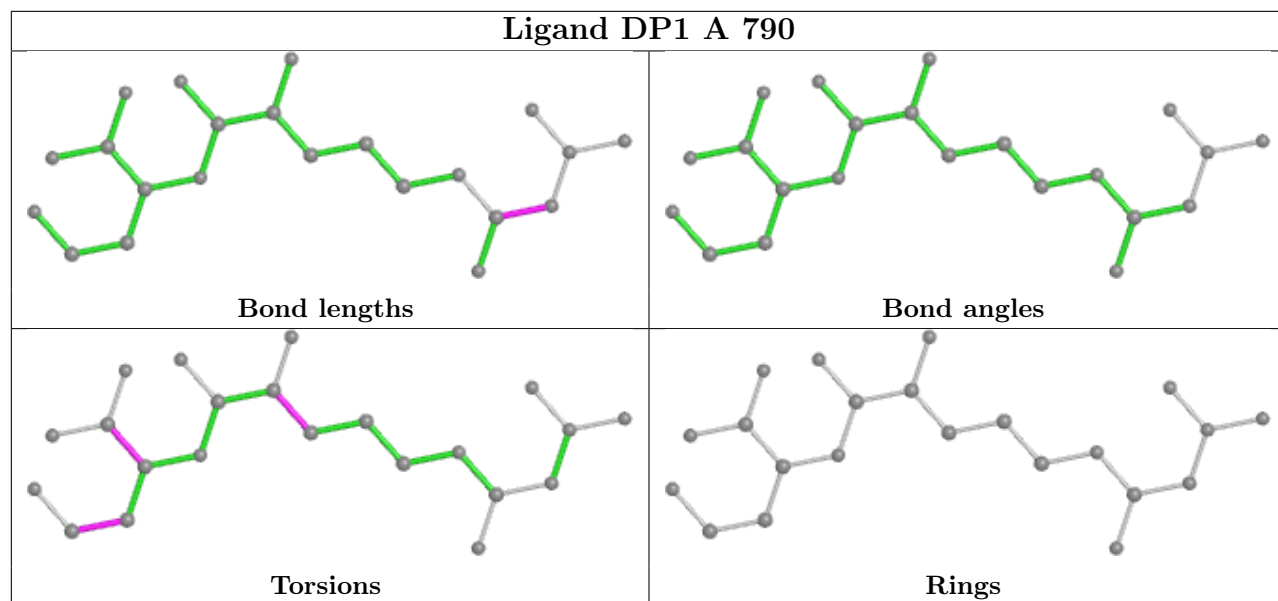
6 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	860	ACT	2	0
4	A	750	HEM	4	0
5	A	760	H4B	1	0
4	B	750	HEM	3	0
6	B	791	DP1	1	0
6	A	790	DP1	1	0

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







5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	407/421 (96%)	0.82	47 (11%) 4 5	22, 41, 66, 78	0
1	B	410/421 (97%)	0.43	21 (5%) 28 30	21, 33, 56, 71	0
All	All	817/842 (97%)	0.63	68 (8%) 11 12	21, 37, 63, 78	0

All (68) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	355	PHE	5.6
1	A	486	LYS	4.8
1	B	300	PHE	4.8
1	B	348	VAL	4.7
1	A	715	VAL	4.6
1	A	716	TRP	4.5
1	A	488	PRO	4.5
1	B	351	LYS	4.5
1	A	300	PHE	4.1
1	B	350	THR	3.9
1	A	489	ASP	3.8
1	B	619	ARG	3.8
1	A	381	LEU	3.7
1	A	490	GLY	3.7
1	A	299	ARG	3.7
1	A	493	LEU	3.6
1	A	388	ILE	3.5
1	A	514	ARG	3.4
1	B	620	LYS	3.3
1	A	512	ALA	3.3
1	B	352	ASP	3.3
1	B	355	PHE	3.3
1	B	321	THR	3.2
1	B	611	ALA	3.2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	499	VAL	3.1
1	A	507	GLN	3.1
1	A	350	THR	3.0
1	A	551	PHE	3.0
1	A	321	THR	3.0
1	A	491	SER	2.9
1	A	506	ILE	2.9
1	B	667	ARG	2.8
1	A	511	LYS	2.6
1	A	677	VAL	2.6
1	A	390	SER	2.6
1	A	480	ILE	2.6
1	A	351	LYS	2.6
1	B	677	VAL	2.6
1	B	566	ALA	2.6
1	A	392	SER	2.6
1	A	523	LEU	2.5
1	B	567	VAL	2.4
1	A	712	ASN	2.4
1	A	322	LEU	2.4
1	A	554	PHE	2.4
1	B	353	GLN	2.4
1	A	352	ASP	2.4
1	B	479	LEU	2.4
1	A	553	TRP	2.4
1	A	373	GLY	2.4
1	B	681	PRO	2.3
1	A	487	GLN	2.3
1	A	516	ARG	2.3
1	A	552	ASP	2.3
1	A	503	GLU	2.3
1	A	676	TRP	2.3
1	A	386	LYS	2.3
1	B	442	ILE	2.2
1	A	337	LEU	2.1
1	A	391	THR	2.1
1	B	493	LEU	2.1
1	A	385	ASN	2.1
1	A	371	ARG	2.1
1	A	394	TYR	2.1
1	A	619	ARG	2.1
1	B	678	TRP	2.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	691	PHE	2.0
1	A	714	HIS	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](#)

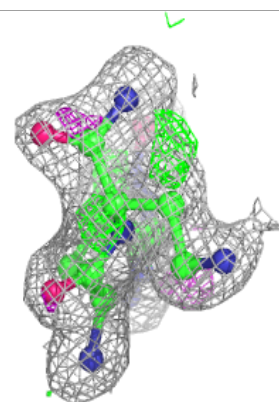
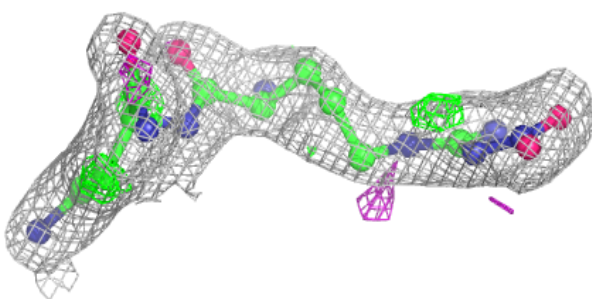
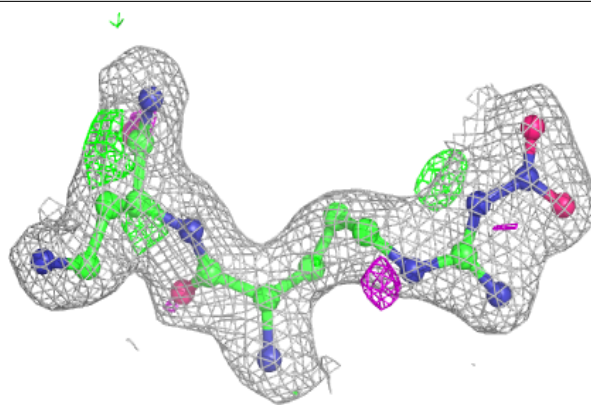
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q < 0.9
6	DP1	B	791	22/22	0.89	0.19	28,34,37,38	0
6	DP1	A	790	22/22	0.90	0.19	31,37,43,44	0
2	ACT	A	860	4/4	0.90	0.17	64,64,64,64	0
2	ACT	B	861	4/4	0.91	0.12	42,43,43,44	0
5	H4B	B	761	17/17	0.96	0.18	22,24,29,30	0
4	HEM	A	750	43/43	0.96	0.14	24,27,28,28	0
5	H4B	A	760	17/17	0.96	0.18	24,25,30,31	0
4	HEM	B	750	43/43	0.97	0.14	22,24,27,29	0
3	ZN	A	900	1/1	0.99	0.08	33,33,33,33	0

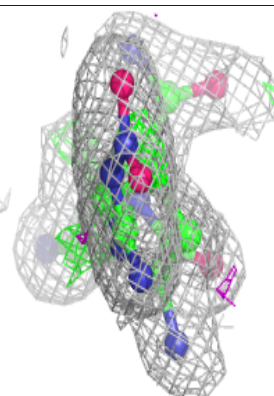
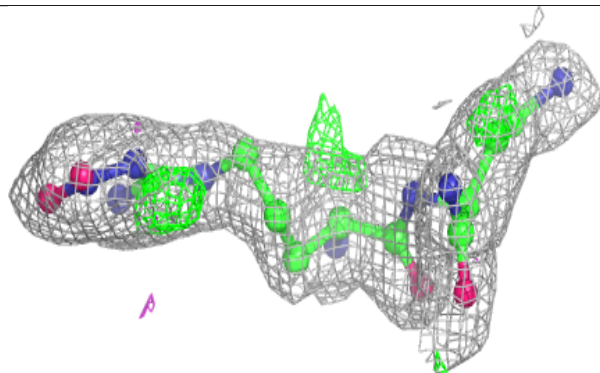
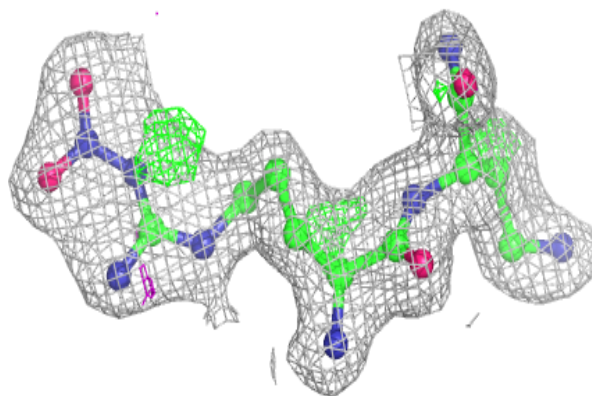
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around DP1 B 791:

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

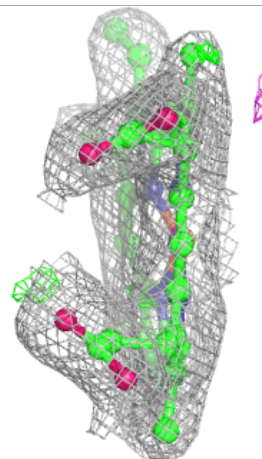
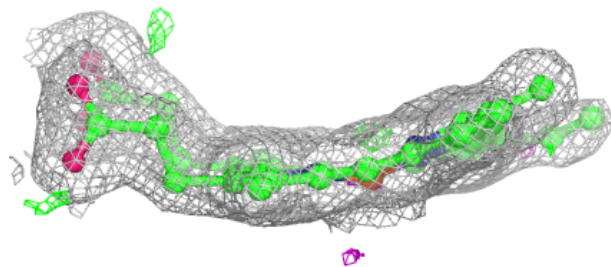
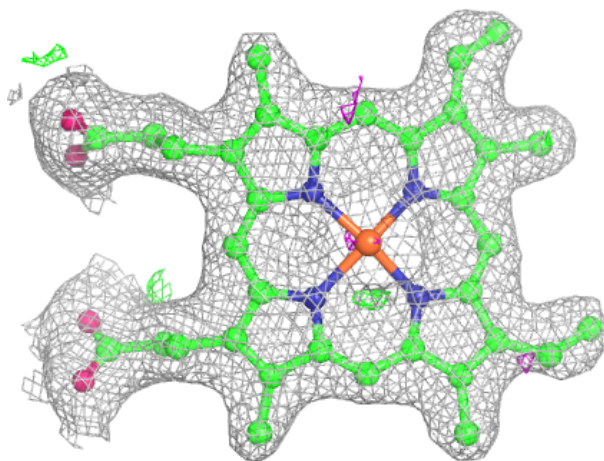
**Electron density around DP1 A 790:**

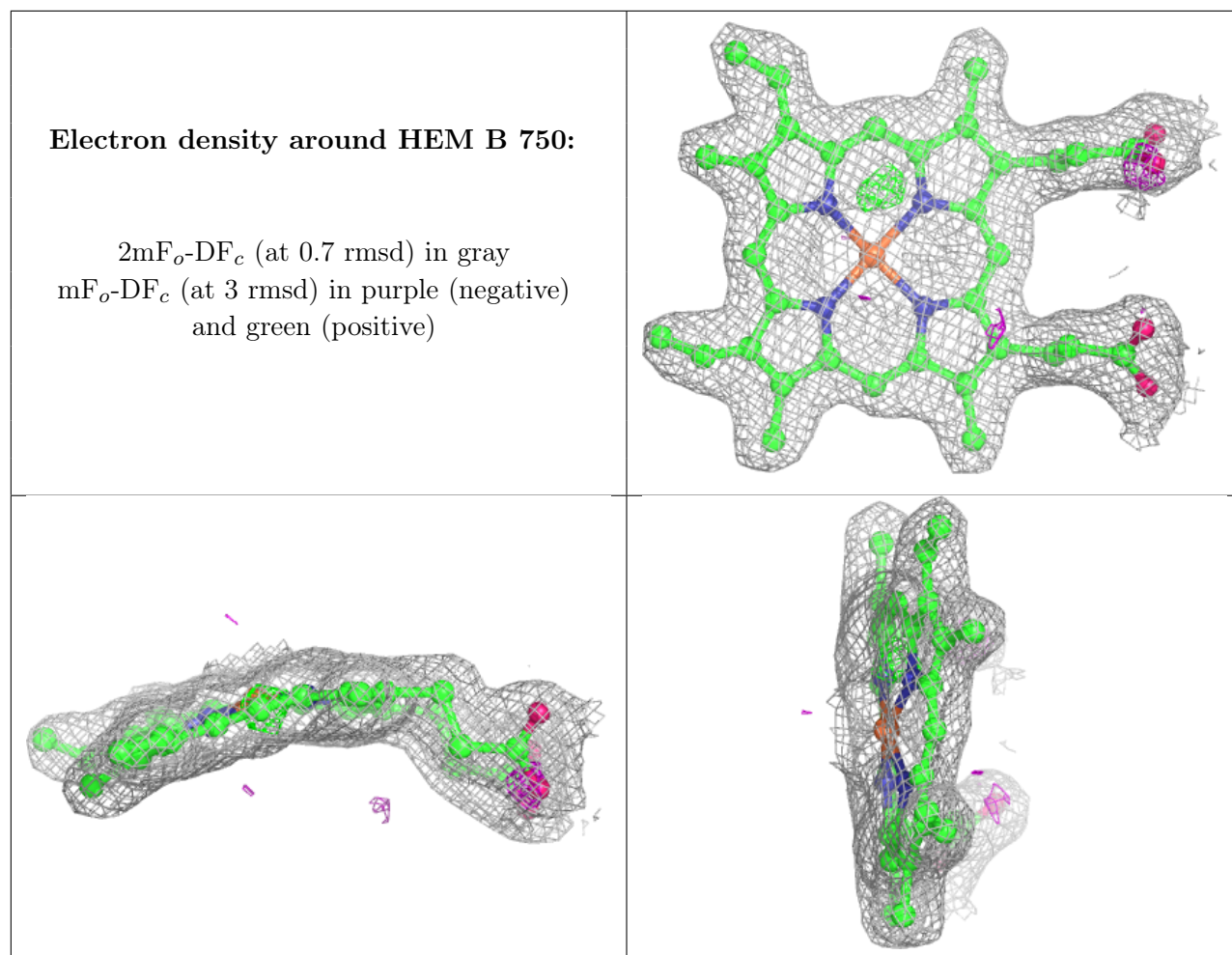
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



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.