



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 11:57 AM UTC

PDB ID : 4K5F / pdb_00004k5f
Title : Structure of neuronal nitric oxide synthase heme domain in complex with (S)-1,3-bis((2-amino-4-methylpyridin-6-yl)-methoxy)-butan-4-amine
Authors : Li, H.; Poulos, T.L.
Deposited on : 2013-04-14
Resolution : 2.20 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

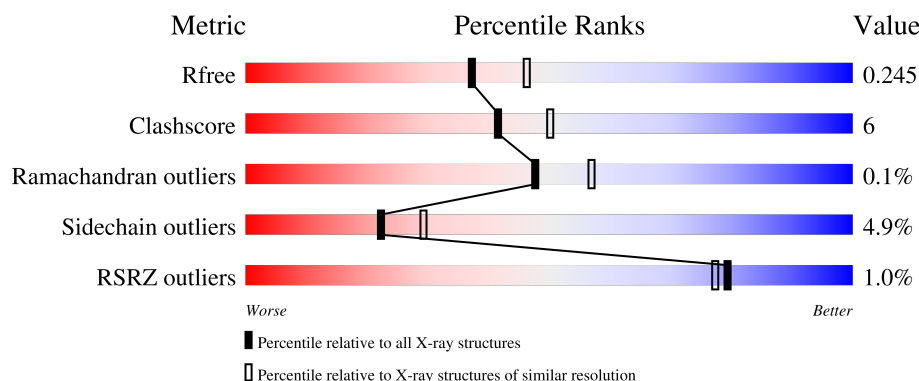
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

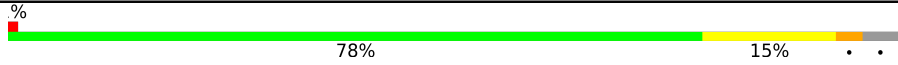
The reported resolution of this entry is 2.20 Å.

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



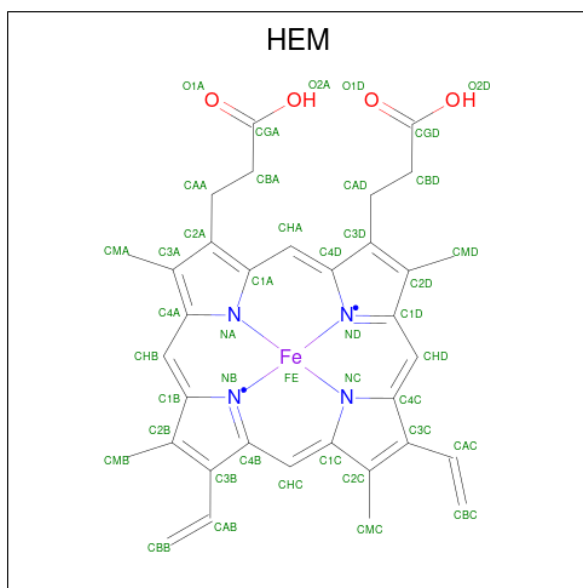
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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

Mol	Chain	Length	Quality of chain
1	A	422	
1	B	422	

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



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

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



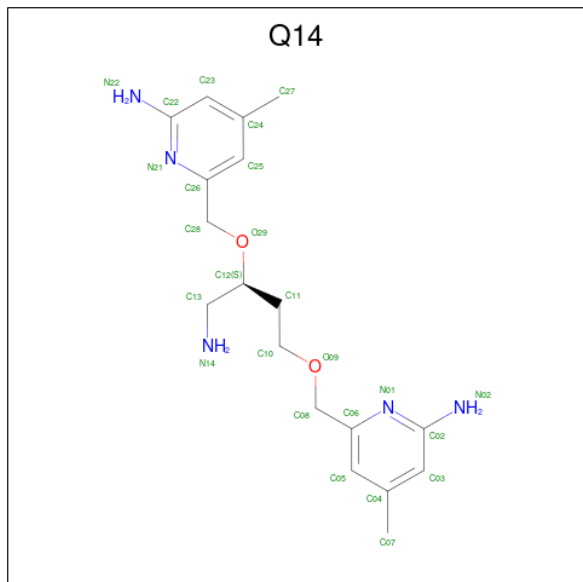
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			17	9	5	3		
3	B	1	Total	C	N	O	0	0
			17	9	5	3		

- Molecule 4 is ACETATE ION (CCD ID: ACT) (formula: $C_2H_3O_2$).



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

- Molecule 5 is 6-[(2S)-1-amino-4-[(6-amino-4-methylpyridin-2-yl)methoxy]butan-2-yl]oxy methyl]-4-methylpyridin-2-amine (CCD ID: Q14) (formula: C₁₈H₂₇N₅O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			25	18	5	2		
5	B	1	Total	C	N	O	0	0
			25	18	5	2		

- Molecule 6 is ZINC ION (CCD ID: ZN) (formula: Zn).

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

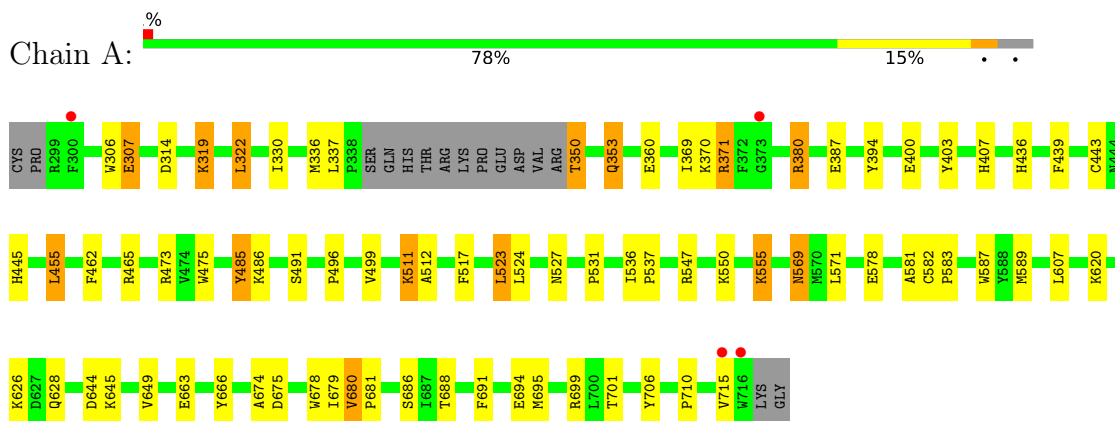
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	60	Total	O	0	0
			60	60		
7	B	88	Total	O	0	0
			88	88		

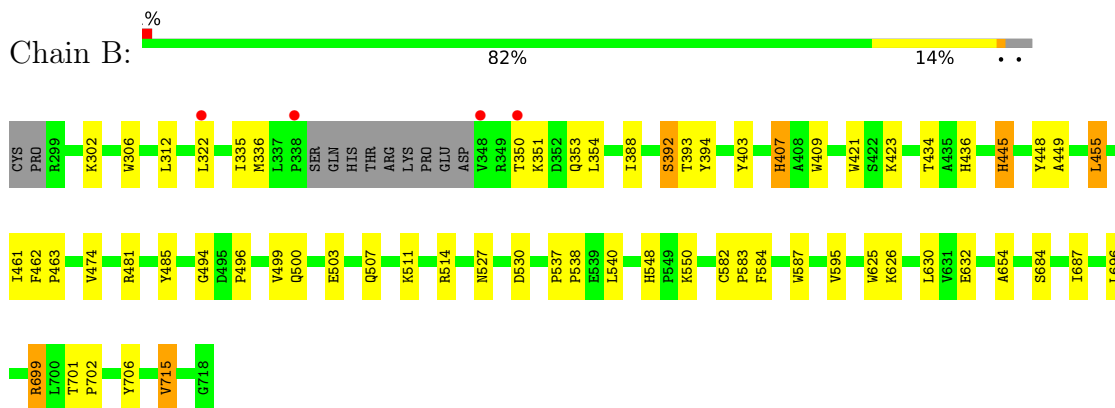
3 Residue-property plots [i](#)

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	51.59Å 110.74Å 164.35Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	92.19 – 2.20 91.84 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.1 (92.19-2.20) 99.2 (91.84-2.20)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.75 (at 2.20Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.192 , 0.248 0.190 , 0.245	Depositor DCC
R_{free} test set	2430 reflections (2.40%)	wwPDB-VP
Wilson B-factor (Å ²)	42.3	Xtriage
Anisotropy	0.807	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 36.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6985	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.46% 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: H4B, ZN, Q14, ACT, HEM

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.92	2/3406 (0.1%)	1.05	7/4621 (0.2%)
1	B	1.00	3/3438 (0.1%)	1.07	7/4661 (0.2%)
All	All	0.96	5/6844 (0.1%)	1.06	14/9282 (0.2%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	436	HIS	CG-CD2	5.34	1.41	1.35
1	A	436	HIS	ND1-CE1	5.29	1.37	1.32
1	A	436	HIS	CG-CD2	5.21	1.41	1.35
1	B	407	HIS	ND1-CE1	5.21	1.37	1.32
1	B	445	HIS	CG-CD2	5.12	1.41	1.35

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	688	THR	CA-C-N	-6.13	113.42	120.04
1	A	688	THR	C-N-CA	-6.13	113.42	120.04
1	A	701	THR	CB-CA-C	6.02	117.93	108.84
1	B	354	LEU	N-CA-C	5.94	117.56	111.14
1	B	474	VAL	N-CA-C	-5.74	98.87	107.37
1	B	530	ASP	CA-C-N	-5.70	114.72	120.31
1	B	530	ASP	C-N-CA	-5.70	114.72	120.31
1	A	455	LEU	N-CA-C	5.63	118.18	110.24
1	B	434	THR	CA-C-N	5.57	127.68	120.44
1	B	434	THR	C-N-CA	5.57	127.68	120.44
1	A	680	VAL	CA-C-N	5.31	123.54	119.66
1	A	680	VAL	C-N-CA	5.31	123.54	119.66
1	B	461	ILE	N-CA-C	5.30	115.22	107.37
1	A	307	GLU	N-CA-C	-5.20	107.00	113.55

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3313	0	3221	53	0
1	B	3345	0	3259	32	0
2	A	43	0	30	4	0
2	B	43	0	30	5	0
3	A	17	0	15	0	0
3	B	17	0	15	0	0
4	A	4	0	3	0	0
4	B	4	0	3	1	0
5	A	25	0	27	1	0
5	B	25	0	27	1	0
6	A	1	0	0	0	0
7	A	60	0	0	1	0
7	B	88	0	0	2	0
All	All	6985	0	6630	86	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 6.

All (86) 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:336:MET:CE	1:A:678:TRP:HZ2	1.88	0.86
1:B:587:TRP:H	2:B:801:HEM:HAB	1.44	0.83
1:B:499:VAL:O	1:B:503:GLU:HG3	1.80	0.81
1:B:706:TYR:OH	2:B:801:HEM:O1D	2.00	0.80
2:A:801:HEM:HMC2	2:A:801:HEM:HBC2	1.62	0.79
1:A:336:MET:CE	1:A:678:TRP:CZ2	2.66	0.78
1:A:350:THR:N	1:A:353:GLN:NE2	2.33	0.77
2:A:801:HEM:HBA1	5:A:804:Q14:H9	1.67	0.75
2:B:801:HEM:HBA1	5:B:803:Q14:H9	1.76	0.66
2:A:801:HEM:HBC2	2:A:801:HEM:CMC	2.26	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:511:LYS:O	1:A:511:LYS:HG2	1.96	0.64
1:A:336:MET:HE2	1:A:678:TRP:HZ2	1.62	0.64
1:B:699:ARG:HB2	1:B:699:ARG:HH11	1.62	0.63
1:A:523:LEU:HD22	1:A:531:PRO:HB2	1.83	0.60
1:A:380:ARG:HD3	1:A:400:GLU:OE1	2.02	0.60
1:A:336:MET:HG3	1:B:306:TRP:NE1	2.18	0.57
1:A:350:THR:N	1:A:353:GLN:HE21	2.02	0.56
1:A:675:ASP:O	1:A:679:ILE:HG12	2.06	0.55
1:A:706:TYR:OH	2:A:801:HEM:O1D	2.18	0.55
1:B:322:LEU:HD12	1:B:699:ARG:HG2	1.89	0.55
1:B:548:HIS:NE2	1:B:632:GLU:OE1	2.40	0.54
1:A:319:LYS:HD2	1:A:319:LYS:N	2.22	0.53
1:A:455:LEU:HD12	1:A:587:TRP:HB3	1.90	0.53
1:A:322:LEU:HD12	1:A:699:ARG:HB3	1.91	0.53
1:A:306:TRP:CD1	1:B:336:MET:HE2	2.43	0.53
1:A:555:LYS:NZ	1:A:555:LYS:HB3	2.24	0.52
1:B:455:LEU:HD23	1:B:587:TRP:HB3	1.91	0.52
1:A:473:ARG:NH2	1:A:710:PRO:HD3	2.25	0.52
1:B:715:VAL:HG23	7:B:952:HOH:O	2.09	0.52
1:A:475:TRP:HB2	1:A:523:LEU:HB3	1.92	0.51
1:B:584:PHE:CD1	2:B:801:HEM:CAC	2.95	0.50
1:A:462:PHE:HB2	1:A:581:ALA:HB3	1.94	0.49
1:A:307:GLU:HG3	7:B:914:HOH:O	2.12	0.49
1:B:449:ALA:O	1:B:455:LEU:HA	2.13	0.49
2:B:801:HEM:HBC2	2:B:801:HEM:CMC	2.44	0.48
1:B:302:LYS:HA	1:B:312:LEU:O	2.14	0.48
1:A:369:ILE:HG13	1:A:371:ARG:HB2	1.95	0.48
1:A:547:ARG:HH12	1:A:644:ASP:CG	2.21	0.48
1:B:403:TYR:CE1	1:B:407:HIS:CE1	3.02	0.47
1:A:314:ASP:HB2	1:A:666:TYR:HE2	1.79	0.47
1:A:330:ILE:HD11	1:B:696:LEU:HB3	1.96	0.47
1:B:684:SER:HB3	1:B:687:ILE:HD11	1.96	0.47
1:A:455:LEU:HD12	1:A:587:TRP:CB	2.44	0.47
1:A:686:SER:HA	1:A:691:PHE:CG	2.50	0.47
1:B:445:HIS:C	1:B:445:HIS:CD2	2.93	0.47
1:A:336:MET:HE3	1:A:678:TRP:CZ2	2.47	0.47
1:B:595:VAL:HG22	1:B:630:LEU:HD11	1.97	0.47
1:A:536:ILE:O	1:A:537:PRO:C	2.56	0.47
1:A:336:MET:HE3	1:A:678:TRP:HZ2	1.76	0.46
1:A:694:GLU:HB3	1:B:335:ILE:HD13	1.98	0.46
1:B:448:TYR:CD2	1:B:448:TYR:C	2.93	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:322:LEU:HB2	1:A:699:ARG:HB2	1.97	0.46
1:B:485:TYR:CZ	1:B:514:ARG:HA	2.51	0.46
1:A:445:HIS:C	1:A:445:HIS:CD2	2.94	0.46
1:B:654:ALA:HA	4:B:804:ACT:H2	1.98	0.46
1:B:582:CYS:O	1:B:583:PRO:C	2.59	0.45
1:B:494:GLY:O	1:B:496:PRO:HD3	2.16	0.45
1:A:571:LEU:C	1:A:571:LEU:HD23	2.42	0.45
1:A:569:ASN:HD22	1:A:569:ASN:H	1.64	0.45
1:A:517:PHE:O	7:A:947:HOH:O	2.20	0.44
1:A:680:VAL:HA	1:A:681:PRO:HD3	1.86	0.44
1:A:569:ASN:H	1:A:569:ASN:ND2	2.16	0.44
1:A:524:LEU:O	1:A:531:PRO:HA	2.16	0.44
1:A:589:MET:HA	1:A:649:VAL:O	2.17	0.44
1:A:319:LYS:HD2	1:A:319:LYS:H	1.83	0.44
1:A:322:LEU:HB2	1:A:699:ARG:HE	1.84	0.43
1:A:674:ALA:HB3	1:A:695:MET:HB3	2.01	0.43
1:B:302:LYS:HE3	1:B:302:LYS:HB2	1.39	0.43
1:B:625:TRP:CZ3	1:B:626:LYS:HG2	2.55	0.42
1:B:462:PHE:HB3	1:B:463:PRO:CD	2.50	0.42
1:A:496:PRO:O	1:A:499:VAL:HG23	2.20	0.42
1:B:701:THR:HA	1:B:702:PRO:C	2.45	0.41
1:A:607:LEU:HD13	1:A:626:LYS:HG2	2.02	0.41
1:B:409:TRP:CE3	1:B:421:TRP:HA	2.55	0.41
1:A:582:CYS:O	1:A:583:PRO:C	2.63	0.41
1:B:393:THR:OG1	1:B:394:TYR:N	2.52	0.41
1:A:353:GLN:H	1:A:353:GLN:HG3	1.67	0.41
1:A:439:PHE:CZ	1:A:443:CYS:SG	3.14	0.41
1:A:485:TYR:CE2	1:A:512:ALA:HB1	2.56	0.41
1:B:388:ILE:O	1:B:392:SER:HA	2.20	0.41
1:B:537:PRO:HA	1:B:538:PRO:HD3	1.83	0.41
1:A:403:TYR:CE1	1:A:407:HIS:CE1	3.09	0.40
1:A:569:ASN:HD22	1:A:569:ASN:N	2.19	0.40
1:A:387:GLU:OE1	1:A:394:TYR:HA	2.21	0.40
1:A:550:LYS:HE3	1:A:550:LYS:HB2	1.58	0.40
1:A:465:ARG:NH1	1:A:578:GLU:OE1	2.55	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	403/422 (96%)	387 (96%)	15 (4%)	1 (0%)	43	51
1	B	407/422 (96%)	395 (97%)	12 (3%)	0	100	100
All	All	810/844 (96%)	782 (96%)	27 (3%)	1 (0%)	48	57

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	491	SER

5.3.2 Protein sidechains

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	363/377 (96%)	342 (94%)	21 (6%)	18	22
1	B	366/377 (97%)	351 (96%)	15 (4%)	27	37
All	All	729/754 (97%)	693 (95%)	36 (5%)	22	29

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	319	LYS
1	A	322	LEU
1	A	337	LEU
1	A	350	THR
1	A	353	GLN

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Mol	Chain	Res	Type
1	A	360	GLU
1	A	370	LYS
1	A	371	ARG
1	A	380	ARG
1	A	485	TYR
1	A	486	LYS
1	A	511	LYS
1	A	523	LEU
1	A	527	ASN
1	A	555	LYS
1	A	569	ASN
1	A	620	LYS
1	A	628	GLN
1	A	645	LYS
1	A	663	GLU
1	A	715	VAL
1	B	350	THR
1	B	351	LYS
1	B	353	GLN
1	B	392	SER
1	B	423	LYS
1	B	455	LEU
1	B	481	ARG
1	B	500	GLN
1	B	507	GLN
1	B	511	LYS
1	B	527	ASN
1	B	540	LEU
1	B	550	LYS
1	B	699	ARG
1	B	715	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	353	GLN
1	A	436	HIS
1	A	454	ASN
1	A	527	ASN
1	A	569	ASN
1	A	697	ASN
1	A	712	ASN

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Mol	Chain	Res	Type
1	B	385	ASN
1	B	407	HIS
1	B	454	ASN
1	B	507	GLN
1	B	535	GLN
1	B	601	ASN
1	B	605	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 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	$\# Z > 2$
5	Q14	B	803	-	25,26,26	0.73	0	28,34,34	2.15	8 (28%)
3	H4B	B	802	-	17,18,18	1.25	2 (11%)	14,26,26	2.13	5 (35%)
4	ACT	A	803	-	3,3,3	0.90	0	3,3,3	0.52	0
5	Q14	A	804	-	25,26,26	0.65	0	28,34,34	1.77	6 (21%)
3	H4B	A	802	-	17,18,18	0.90	1 (5%)	14,26,26	2.08	5 (35%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	HEM	B	801	1	50,50,50	2.52	24 (48%)	67,82,82	3.08	26 (38%)
2	HEM	A	801	1	50,50,50	2.50	19 (38%)	67,82,82	2.97	26 (38%)
4	ACT	B	804	-	3,3,3	0.89	0	3,3,3	0.73	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
5	Q14	B	803	-	-	2/14/14/14	0/2/2/2
3	H4B	B	802	-	-	0/8/17/17	0/2/2/2
5	Q14	A	804	-	-	1/14/14/14	0/2/2/2
3	H4B	A	802	-	-	0/8/17/17	0/2/2/2
2	HEM	B	801	1	-	3/14/54/54	-
2	HEM	A	801	1	-	3/14/54/54	-

All (46) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	801	HEM	C3C-C2C	5.62	1.48	1.37
2	B	801	HEM	FE-NB	5.19	2.10	1.94
2	B	801	HEM	FE-ND	5.16	2.10	1.94
2	A	801	HEM	FE-NB	5.10	2.10	1.94
2	B	801	HEM	C3C-C2C	4.96	1.47	1.37
2	A	801	HEM	FE-ND	4.92	2.10	1.94
2	B	801	HEM	CHC-C1C	4.51	1.47	1.38
2	A	801	HEM	CHD-C4C	4.50	1.47	1.38
2	B	801	HEM	C3B-C2B	4.48	1.46	1.37
2	A	801	HEM	C3D-C2D	4.47	1.46	1.36
2	A	801	HEM	CHC-C1C	4.40	1.47	1.38
2	B	801	HEM	CHD-C4C	4.31	1.46	1.38
2	B	801	HEM	C4A-NA	4.19	1.47	1.39
2	A	801	HEM	C3B-C2B	4.13	1.45	1.37
2	A	801	HEM	C4D-ND	-3.88	1.33	1.40
2	B	801	HEM	C3D-C2D	3.80	1.45	1.36
2	A	801	HEM	CHC-C4B	3.68	1.47	1.39
2	A	801	HEM	C1C-NC	3.65	1.46	1.39
2	B	801	HEM	CHC-C4B	3.50	1.47	1.39
2	B	801	HEM	CHA-C4D	3.42	1.45	1.38
3	B	802	H4B	C7-N8	3.41	1.49	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	801	HEM	CHB-C4A	3.38	1.47	1.39
2	B	801	HEM	C2A-C3A	3.31	1.47	1.38
2	A	801	HEM	CHD-C1D	3.23	1.46	1.39
2	B	801	HEM	C4D-ND	-3.17	1.34	1.40
2	B	801	HEM	C1A-NA	3.10	1.45	1.39
2	A	801	HEM	CHA-C4D	3.08	1.44	1.38
2	B	801	HEM	C4A-C3A	2.98	1.50	1.43
2	A	801	HEM	C4A-NA	2.93	1.45	1.39
2	A	801	HEM	C1A-NA	2.92	1.45	1.39
2	A	801	HEM	C2A-C3A	2.90	1.46	1.38
2	B	801	HEM	CHD-C1D	2.83	1.45	1.39
2	B	801	HEM	CHB-C4A	2.81	1.45	1.39
2	B	801	HEM	CHA-C1A	2.66	1.45	1.39
2	B	801	HEM	C1D-ND	-2.59	1.33	1.38
2	B	801	HEM	CHB-C1B	2.56	1.43	1.38
2	A	801	HEM	CHB-C1B	2.39	1.43	1.38
2	A	801	HEM	C4A-C3A	2.37	1.48	1.43
3	B	802	H4B	O4-C4	2.34	1.28	1.23
2	B	801	HEM	C1C-NC	2.33	1.44	1.39
2	B	801	HEM	O2A-CGA	-2.28	1.23	1.30
2	B	801	HEM	C1D-C2D	2.24	1.49	1.44
2	B	801	HEM	C1A-C2A	2.21	1.49	1.44
2	B	801	HEM	C3B-C4B	2.18	1.49	1.44
2	A	801	HEM	C1B-NB	-2.12	1.36	1.40
3	A	802	H4B	C4-N3	-2.00	1.35	1.38

All (76) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	801	HEM	C3B-C2B-C1B	-9.95	98.94	106.41
2	B	801	HEM	C2B-C1B-NB	8.67	119.81	109.84
2	A	801	HEM	C3B-C2B-C1B	-8.66	99.91	106.41
2	A	801	HEM	C2B-C1B-NB	7.99	119.03	109.84
2	B	801	HEM	C3D-C4D-ND	7.59	118.49	110.17
2	B	801	HEM	C2D-C1D-ND	7.26	118.30	109.90
2	A	801	HEM	C2D-C1D-ND	7.06	118.07	109.90
2	B	801	HEM	C3B-C4B-NB	7.06	114.54	109.47
2	A	801	HEM	C3B-C4B-NB	6.99	114.49	109.47
5	B	803	Q14	C02-N01-C06	6.73	123.10	118.07
2	A	801	HEM	C3D-C4D-ND	6.16	116.93	110.17
2	A	801	HEM	C3C-C2C-C1C	-6.01	101.36	107.05
2	A	801	HEM	C1D-C2D-C3D	-5.66	101.03	106.98

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	803	Q14	C22-N21-C26	5.63	122.28	118.07
2	B	801	HEM	C1D-C2D-C3D	-5.20	101.51	106.98
2	A	801	HEM	C2A-C1A-NA	4.83	115.52	110.15
3	A	802	H4B	C2-N1-C8A	4.82	121.87	113.36
2	B	801	HEM	C4D-ND-C1D	-4.64	99.71	105.21
5	A	804	Q14	C02-N01-C06	4.62	121.53	118.07
3	B	802	H4B	C2-N1-C8A	4.60	121.48	113.36
5	A	804	Q14	C22-N21-C26	4.57	121.49	118.07
2	B	801	HEM	C2A-C1A-NA	4.44	115.07	110.15
2	B	801	HEM	C1B-NB-C4B	-3.99	100.49	105.21
2	B	801	HEM	CMD-C2D-C1D	3.93	131.18	125.03
2	B	801	HEM	C3C-C2C-C1C	-3.83	103.42	107.05
2	A	801	HEM	CMB-C2B-C1B	3.78	130.93	125.03
2	B	801	HEM	CHA-C4D-C3D	-3.76	118.29	125.23
2	A	801	HEM	CAD-CBD-CGD	-3.70	103.86	113.67
2	B	801	HEM	C4A-NA-C1A	-3.68	99.82	105.82
2	B	801	HEM	CMA-C3A-C4A	3.65	130.98	125.42
2	B	801	HEM	CHB-C1B-C2B	-3.62	116.66	126.95
2	B	801	HEM	CBA-CAA-C2A	-3.60	102.57	112.53
2	A	801	HEM	C1B-NB-C4B	-3.58	100.97	105.21
2	B	801	HEM	CHD-C1D-ND	-3.56	120.60	124.42
2	A	801	HEM	C4A-NA-C1A	-3.51	100.09	105.82
2	A	801	HEM	C4D-ND-C1D	-3.49	101.07	105.21
3	B	802	H4B	O4-C4-C4A	-3.46	118.91	127.26
2	B	801	HEM	C4D-C3D-C2D	-3.41	101.92	106.89
2	A	801	HEM	CHD-C1D-C2D	-3.25	119.89	125.03
5	B	803	Q14	C05-C06-N01	-3.19	119.10	122.73
2	A	801	HEM	CHA-C4D-C3D	-3.16	119.40	125.23
2	B	801	HEM	C3A-C4A-NA	3.12	115.14	110.14
2	A	801	HEM	CMD-C2D-C1D	2.96	129.66	125.03
5	A	804	Q14	N22-C22-N21	2.94	121.31	116.59
3	A	802	H4B	C4-C4A-N5	2.87	124.08	116.27
3	B	802	H4B	N2-C2-N3	2.84	122.76	116.76
2	A	801	HEM	CHA-C1A-C2A	-2.81	119.14	125.30
2	A	801	HEM	C4C-C3C-C2C	2.80	109.24	106.81
5	B	803	Q14	C25-C26-N21	-2.79	119.55	122.73
2	A	801	HEM	C4D-C3D-C2D	-2.74	102.90	106.89
2	A	801	HEM	C1A-CHA-C4D	-2.68	119.94	126.25
2	B	801	HEM	C4A-C3A-C2A	-2.60	103.84	106.82
3	A	802	H4B	C4A-C4-N3	2.60	119.26	112.13
5	A	804	Q14	C05-C06-N01	-2.60	119.77	122.73
2	B	801	HEM	CBD-CAD-C3D	-2.59	105.36	112.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	801	HEM	CHB-C1B-C2B	-2.59	119.58	126.95
3	A	802	H4B	N3-C2-N1	-2.59	118.58	123.32
2	B	801	HEM	CHD-C1D-C2D	-2.56	120.99	125.03
2	B	801	HEM	CMB-C2B-C1B	2.55	129.02	125.03
5	A	804	Q14	C25-C26-N21	-2.48	119.90	122.73
2	A	801	HEM	C2C-C1C-NC	2.47	114.20	109.64
2	A	801	HEM	CBA-CAA-C2A	-2.46	105.72	112.53
2	A	801	HEM	C3A-C4A-NA	2.42	114.02	110.14
3	B	802	H4B	C4A-C4-N3	2.41	118.75	112.13
3	B	802	H4B	N3-C2-N1	-2.41	118.91	123.32
2	A	801	HEM	CHB-C1B-NB	-2.41	121.39	124.37
3	A	802	H4B	N2-C2-N3	2.39	121.80	116.76
5	B	803	Q14	C28-C26-N21	2.38	120.50	115.69
2	B	801	HEM	CHA-C1A-C2A	-2.26	120.35	125.30
2	A	801	HEM	CHD-C1D-ND	-2.24	122.02	124.42
5	B	803	Q14	N22-C22-N21	2.20	120.13	116.59
2	B	801	HEM	CMC-C2C-C1C	2.18	128.57	124.73
5	A	804	Q14	C28-C26-N21	2.15	120.02	115.69
2	B	801	HEM	CBB-CAB-C3B	-2.10	117.03	127.53
5	B	803	Q14	C08-C06-N01	2.04	119.81	115.69
5	B	803	Q14	N02-C02-N01	2.02	119.84	116.59

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	801	HEM	C2B-C3B-CAB-CBB
2	A	801	HEM	C2A-CAA-CBA-CGA
2	B	801	HEM	C2A-CAA-CBA-CGA
2	A	801	HEM	C2B-C3B-CAB-CBB
2	A	801	HEM	C4B-C3B-CAB-CBB
2	B	801	HEM	C4B-C3B-CAB-CBB
5	B	803	Q14	N01-C06-C08-O09
5	B	803	Q14	C05-C06-C08-O09
5	A	804	Q14	C11-C10-O09-C08

There are no ring outliers.

5 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	803	Q14	1	0

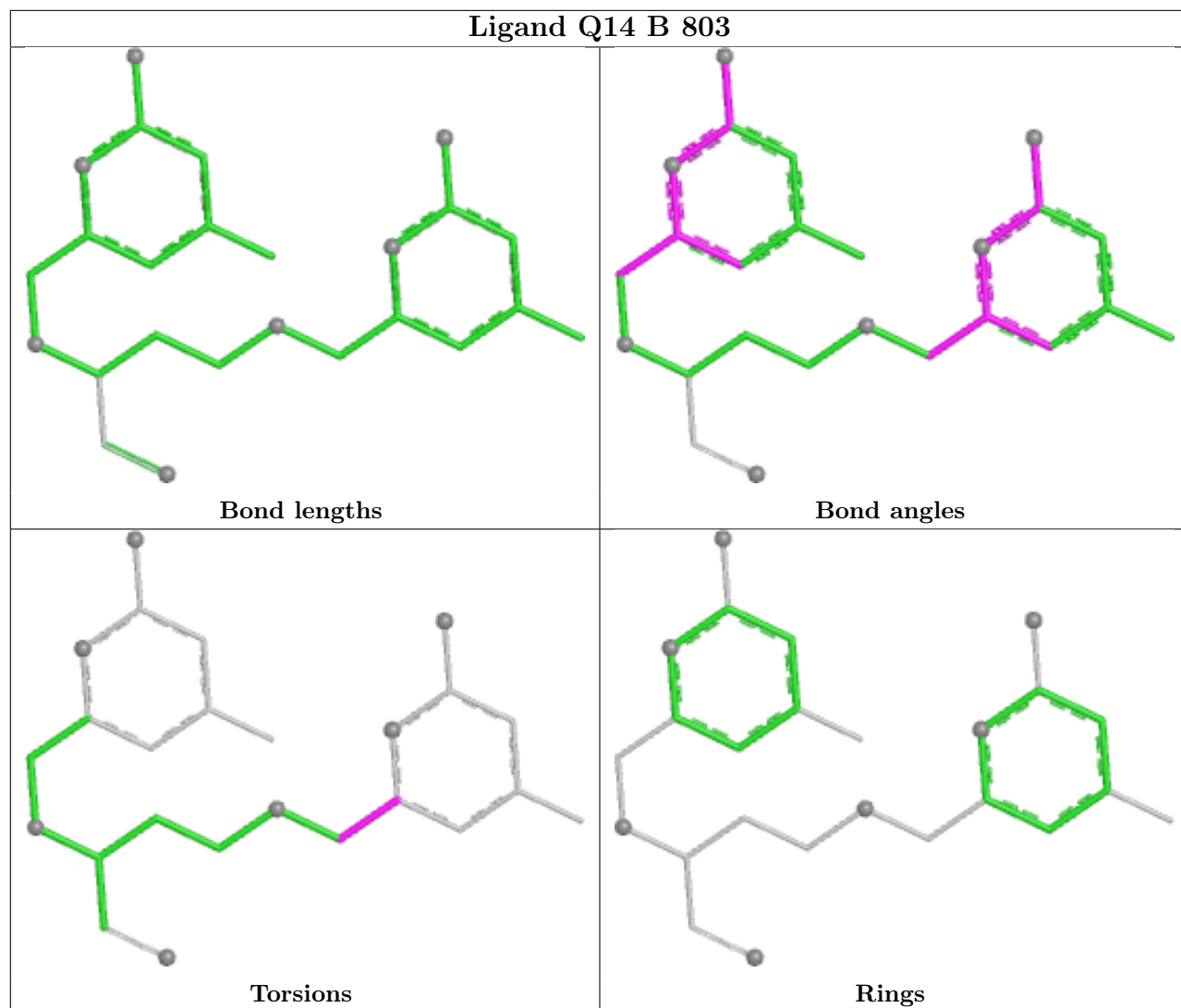
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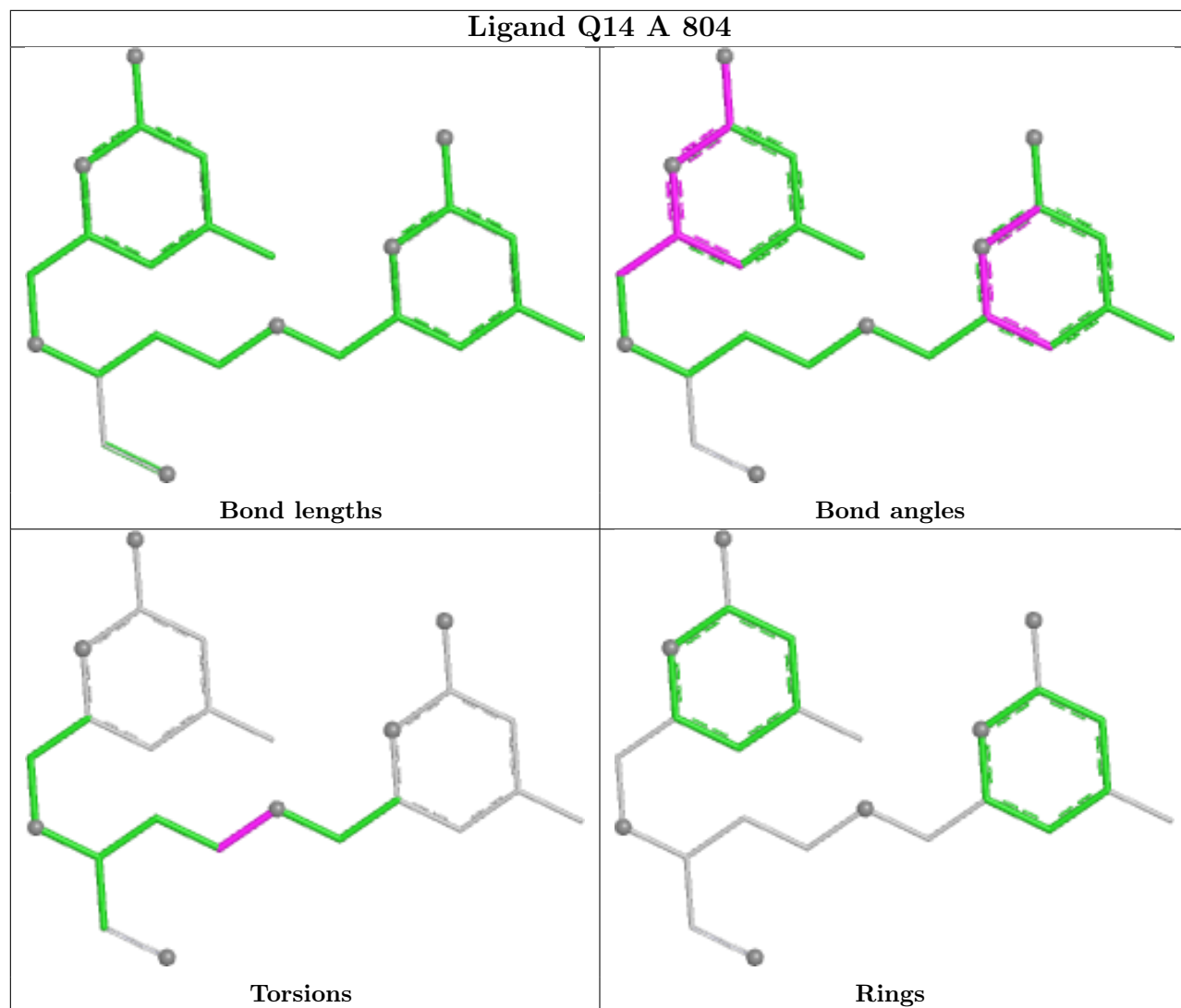
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	804	Q14	1	0
2	B	801	HEM	5	0
2	A	801	HEM	4	0
4	B	804	ACT	1	0

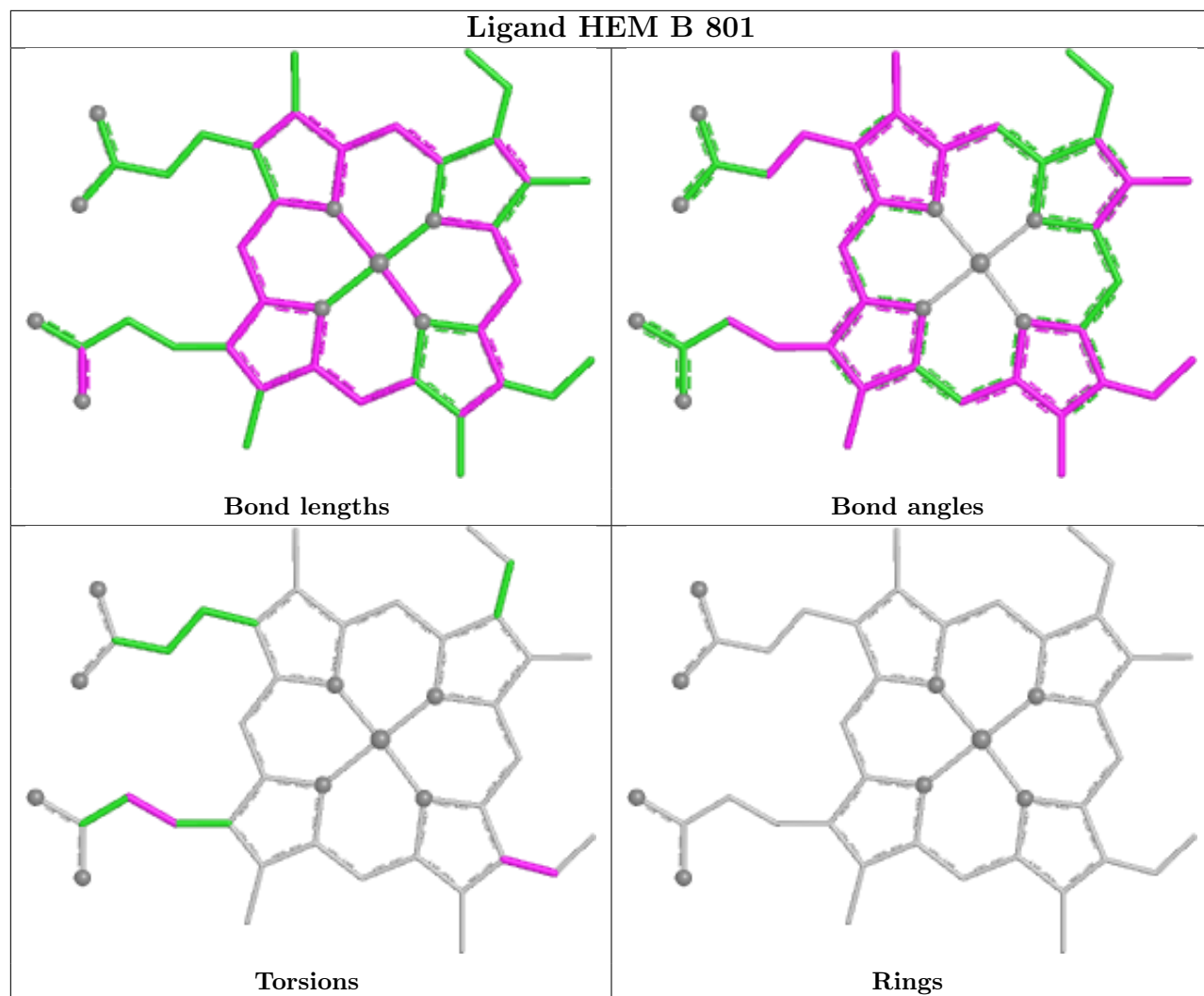
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

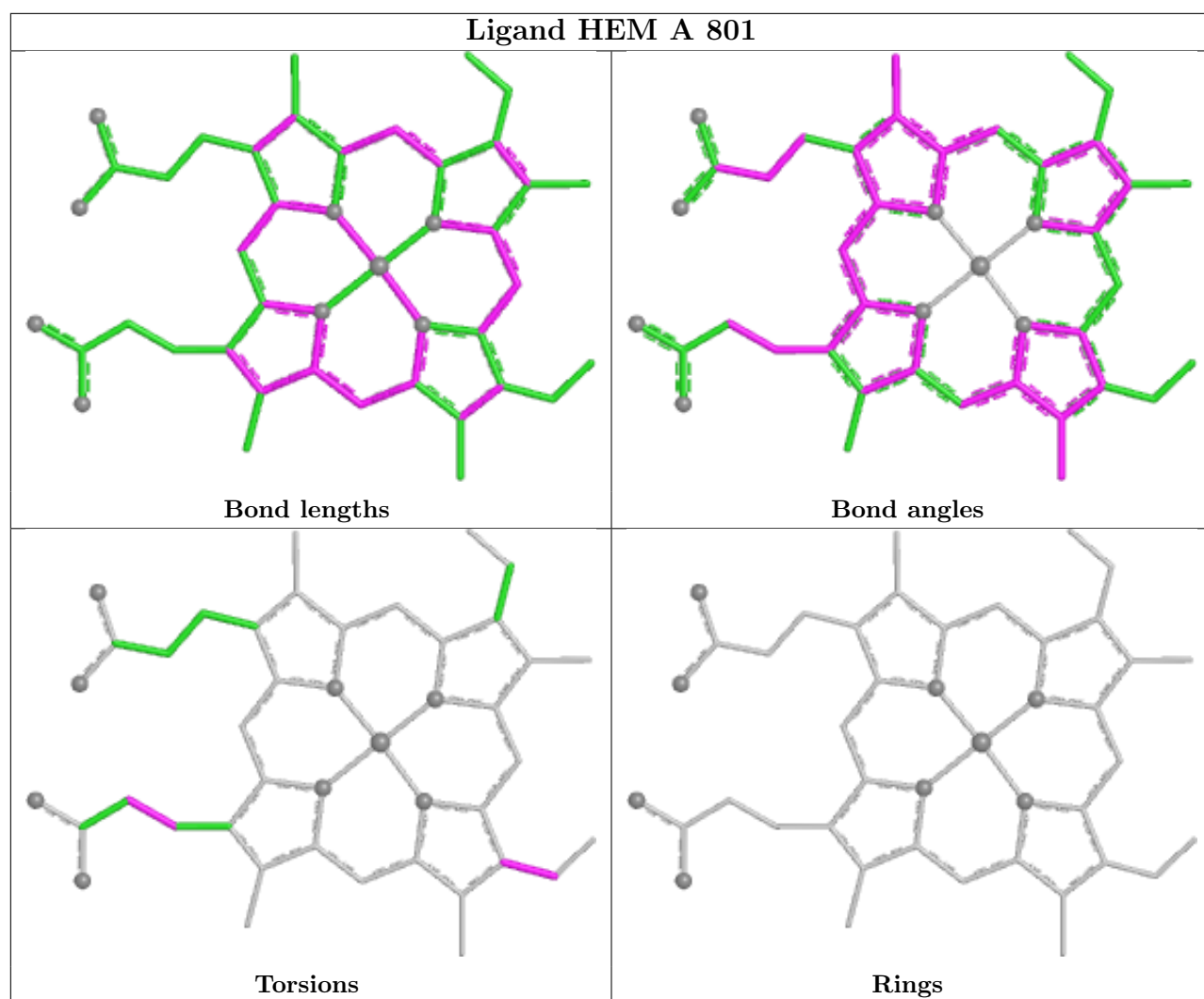
Ligand Q14 B 803



Ligand Q14 A 804







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	407/422 (96%)	-0.02	4 (0%) 79 77	36, 61, 100, 129	0
1	B	411/422 (97%)	-0.25	4 (0%) 79 77	33, 51, 81, 101	0
All	All	818/844 (96%)	-0.13	8 (0%) 79 77	33, 55, 94, 129	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	300	PHE	4.0
1	B	348	VAL	4.0
1	B	338	PRO	3.8
1	A	716	TRP	2.2
1	B	350	THR	2.2
1	A	715	VAL	2.1
1	B	322	LEU	2.0
1	A	373	GLY	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

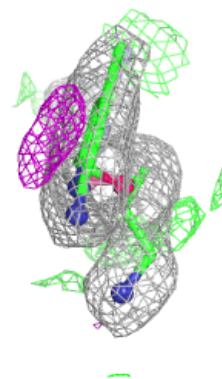
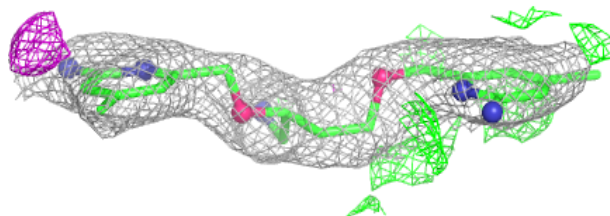
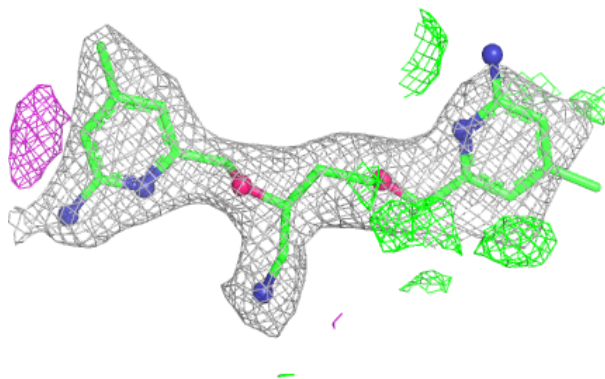
median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	Q14	B	803	25/25	0.84	0.14	44,54,85,92	0
4	ACT	B	804	4/4	0.88	0.15	65,65,67,73	0
5	Q14	A	804	25/25	0.90	0.11	33,48,99,102	0
3	H4B	B	802	17/17	0.92	0.10	44,49,55,60	0
4	ACT	A	803	4/4	0.93	0.19	67,68,73,75	0
3	H4B	A	802	17/17	0.93	0.07	43,51,56,57	0
2	HEM	A	801	43/43	0.98	0.07	35,44,63,72	0
2	HEM	B	801	43/43	0.98	0.08	36,45,60,70	0
6	ZN	A	805	1/1	1.00	0.02	53,53,53,53	0

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

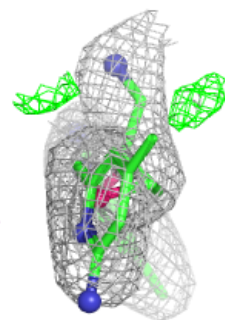
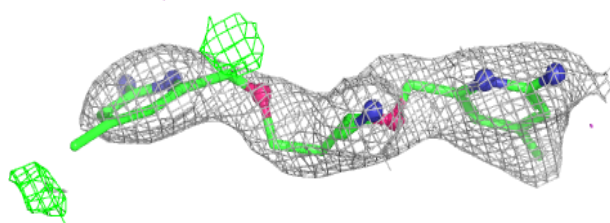
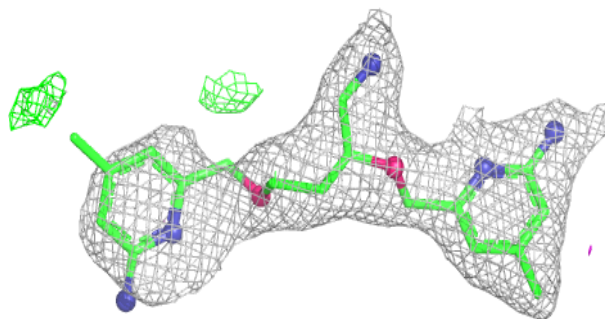
Electron density around Q14 B 803:

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



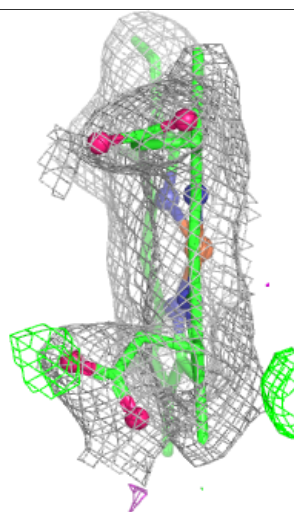
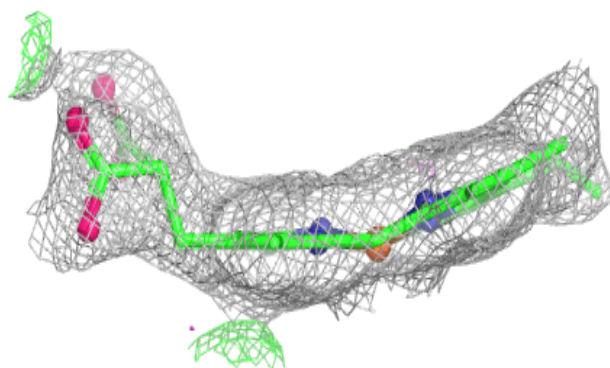
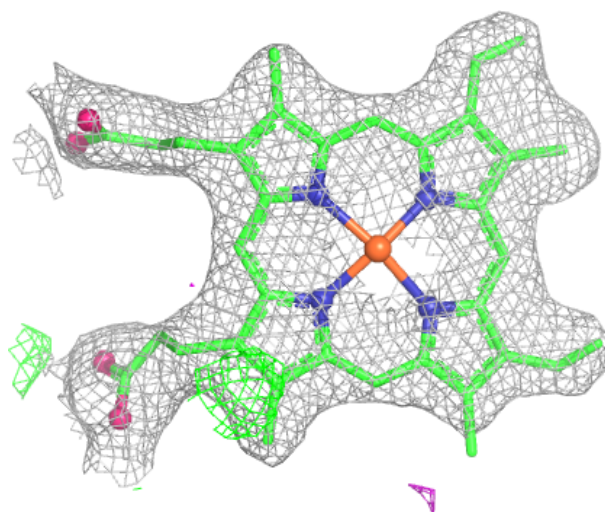
Electron density around Q14 A 804:

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



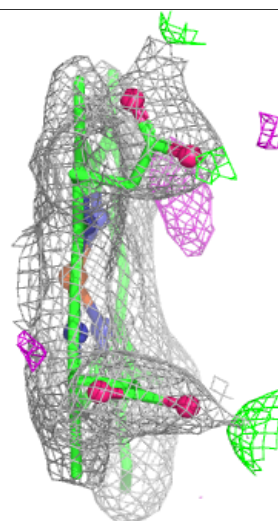
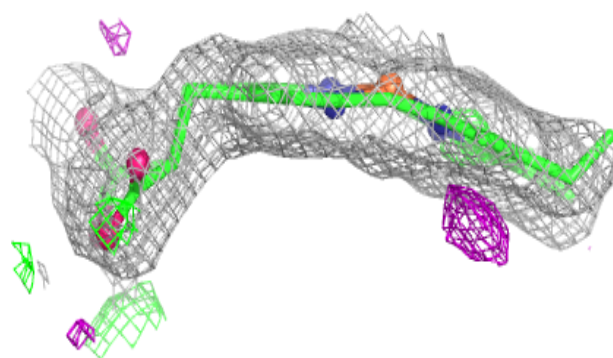
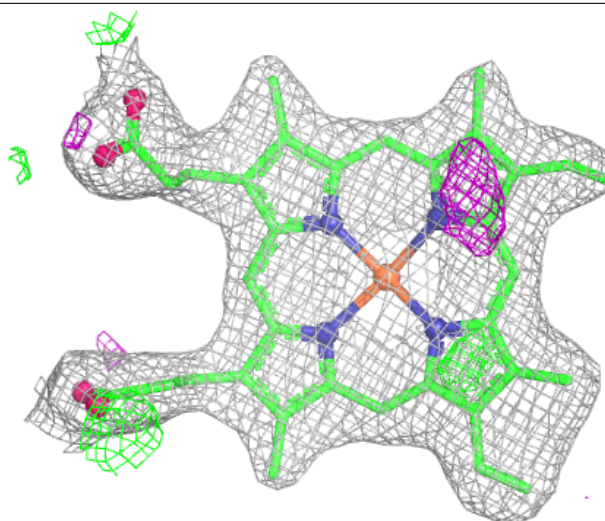
Electron density around HEM A 801:

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



Electron density around HEM B 801:

$2mF_o-DF_c$ (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.