



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 5, 2023 – 07:31 AM EDT

PDB ID : 3UFO
Title : Structure of rat nitric oxide synthase heme domain in complex with 6-(((3S,4S)-4-((5-(3-fluorophenyl)pentyl)oxy)pyrrolidin-3-yl)methyl)-4-methylpyridin-2-amine
Authors : Li, H.; Poulos, T.L.
Deposited on : 2011-11-01
Resolution : 2.17 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)
Xtrriage (Phenix) : 1.13
EDS : 2.35
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) : 2.35

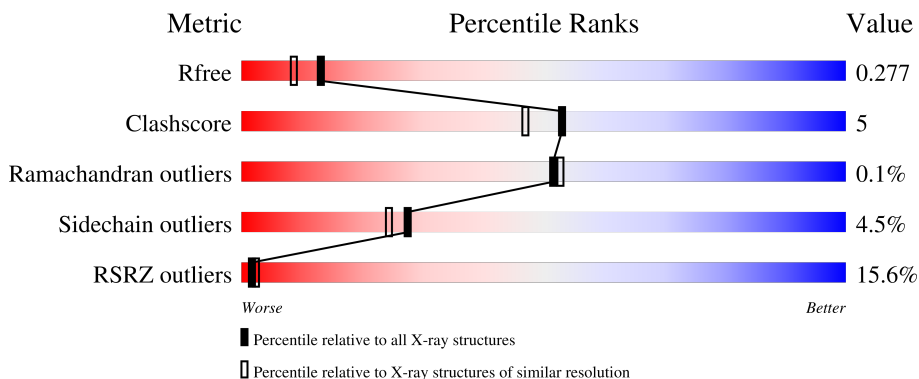
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION



The reported resolution of this entry is 2.17 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	1479 (2.16-2.16)
Clashscore	141614	1585 (2.16-2.16)
Ramachandran outliers	138981	1560 (2.16-2.16)
Sidechain outliers	138945	1559 (2.16-2.16)
RSRZ outliers	127900	1456 (2.16-2.16)

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	

2 Entry composition i

There are 7 unique types of molecules in this entry. The entry contains 7068 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 3322	C 2127	N 567	O 606	S 22	0	2	0
1	B	411	Total 3357	C 2150	N 574	O 611	S 22	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	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 (three-letter code: H4B) (formula: $C_9H_{15}N_5O_3$).



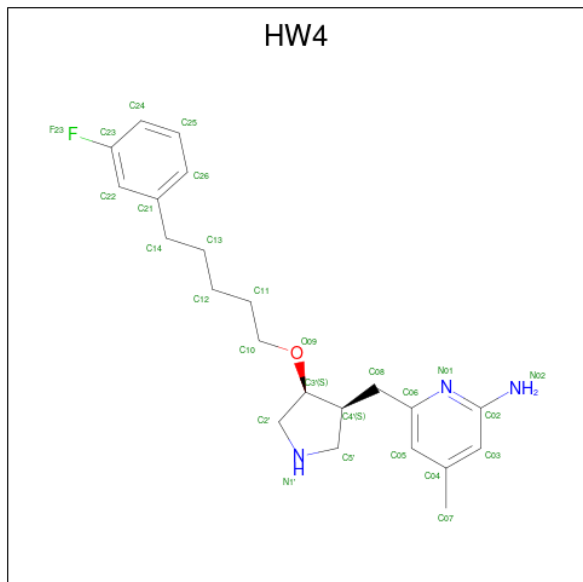
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
3	A	1	17	9	5	3	0	0
3	B	1	17	9	5	3	0	0

- Molecule 4 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



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

- Molecule 5 is 6-{{(3S,4S)-4-{{5-(3-fluorophenyl)pentyl}oxy}pyrrolidin-3-yl}methyl}-4-methylpyridin-2-amine (three-letter code: HW4) (formula: C₂₂H₃₀FN₃O).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	F	N	O		
5	A	1	27	22	1	3	1	0	0
5	B	1	38	32	2	3	1	0	1

- 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

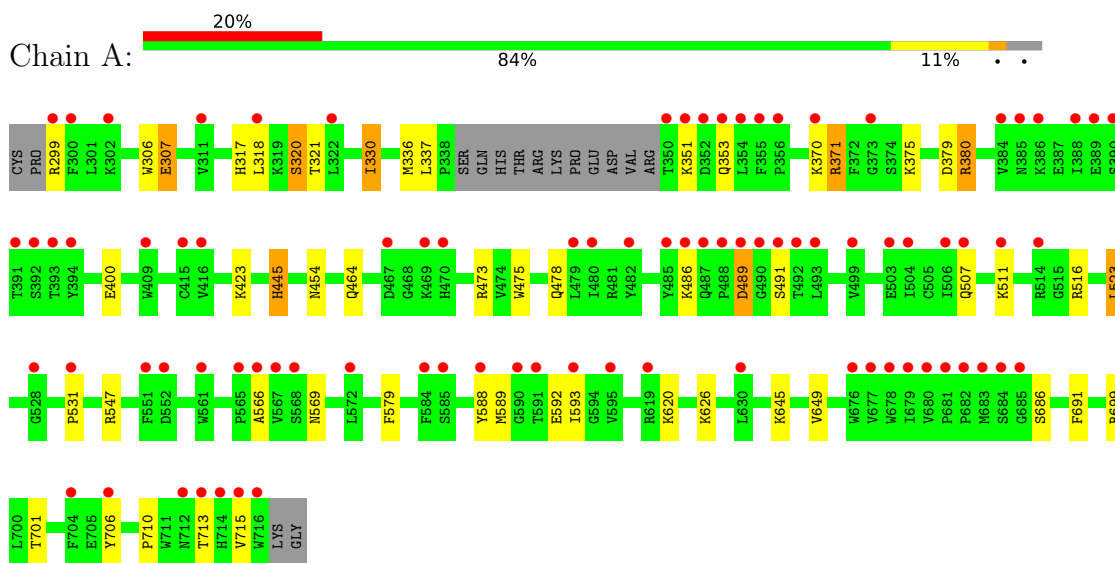
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
7	A	81	81	81	0	0
7	B	114	114	114	0	0

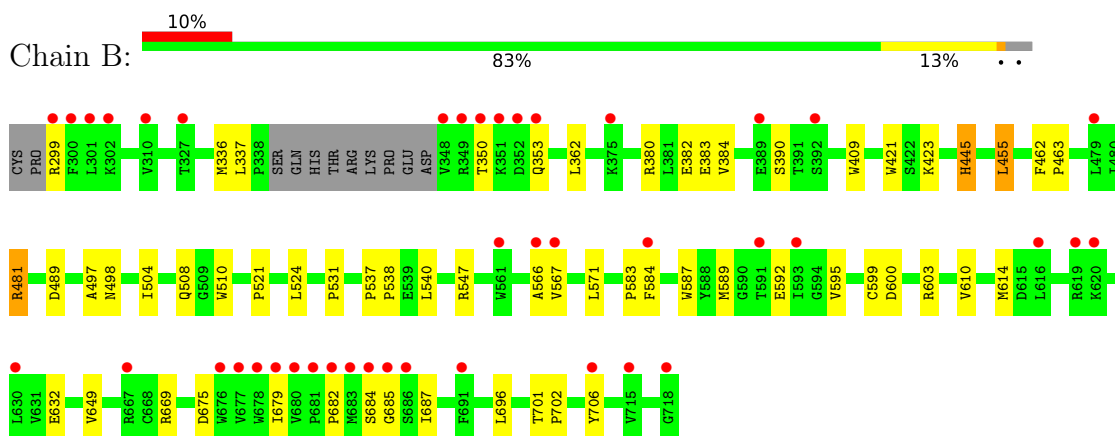
3 Residue-property plots [i](#)

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



- 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	52.19Å 111.58Å 165.01Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.69 – 2.17 38.69 – 2.17	Depositor EDS
% Data completeness (in resolution range)	96.9 (38.69-2.17) 96.9 (38.69-2.17)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.07 (at 2.18Å)	Xtrriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.198 , 0.254 0.237 , 0.277	Depositor DCC
R_{free} test set	2525 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	42.2	Xtrriage
Anisotropy	0.641	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 48.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7068	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.70% 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: ACT, ZN, H4B, HW4, 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	0.72	0/3421	0.72	0/4641
1	B	0.87	1/3457 (0.0%)	0.80	2/4687 (0.0%)
All	All	0.80	1/6878 (0.0%)	0.76	2/9328 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	497	ALA	CA-CB	5.37	1.63	1.52

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	489	ASP	CB-CG-OD2	5.79	123.51	118.30
1	B	669	ARG	NE-CZ-NH2	-5.76	117.42	120.30

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	3322	0	3234	32	0
1	B	3357	0	3273	36	0
2	A	43	0	30	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	43	0	30	8	0
3	A	17	0	15	0	0
3	B	17	0	15	1	0
4	A	4	0	3	0	0
4	B	4	0	3	0	0
5	A	27	0	30	3	0
5	B	38	0	28	4	0
6	A	1	0	0	0	0
7	A	81	0	0	0	0
7	B	114	0	0	1	0
All	All	7068	0	6661	73	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 5.

All (73) 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:371:ARG:HG2	1:A:371:ARG:HH21	1.21	1.00
5:B:800[A]:HW4:H10	5:B:800[A]:HW4:H3	1.46	0.98
1:A:371:ARG:HH21	1:A:371:ARG:CG	1.80	0.94
1:A:371:ARG:HG2	1:A:371:ARG:NH2	1.84	0.88
1:A:523:LEU:HD22	1:A:531:PRO:HB2	1.59	0.85
1:A:330:ILE:HD11	1:B:696:LEU:HD22	1.69	0.75
1:A:686:SER:HB3	1:B:682:PRO:HB2	1.71	0.70
1:A:317:HIS:O	1:A:320:SER:HB3	1.92	0.69
1:A:380:ARG:HD3	1:A:400:GLU:OE1	1.95	0.66
1:B:706[A]:TYR:OH	2:B:750:HEM:O2D	2.17	0.59
1:A:626:LYS:HB3	1:B:687:ILE:HD12	1.85	0.58
1:A:475:TRP:HB2	1:A:523:LEU:HB3	1.85	0.58
1:A:523:LEU:CD2	1:A:531:PRO:HB2	2.33	0.57
1:A:592:GLU:OE1	5:A:800:HW4:H21	2.04	0.57
1:A:706:TYR:OH	2:A:750:HEM:O2D	2.19	0.56
5:B:800[A]:HW4:H3	5:B:800[A]:HW4:C12	2.29	0.56
1:B:504:ILE:O	1:B:508:GLN:HG2	2.06	0.55
5:B:800[A]:HW4:H10	5:B:800[A]:HW4:C26	2.27	0.55
1:B:445:HIS:C	1:B:445:HIS:CD2	2.81	0.54
1:A:473:ARG:NH2	1:A:710:PRO:HD3	2.22	0.54
1:B:409:TRP:CE3	1:B:421:TRP:HA	2.44	0.53
1:B:299:ARG:HB3	1:B:299:ARG:NH1	2.25	0.51
1:B:337:LEU:HD21	1:B:706[B]:TYR:CD2	2.45	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:750:HEM:HBA2	5:A:800:HW4:H23	1.92	0.51
1:B:584:PHE:CD1	2:B:750:HEM:HAC	2.45	0.51
1:A:306:TRP:CD1	1:B:336:MET:HE2	2.46	0.50
1:B:380:ARG:NH1	1:B:383:GLU:OE2	2.43	0.50
1:B:587:TRP:H	2:B:750:HEM:HAB	1.77	0.50
2:A:750:HEM:C1C	5:A:800:HW4:H31	2.47	0.50
1:B:706[A]:TYR:OH	2:B:750:HEM:CGD	2.60	0.49
1:B:675:ASP:O	1:B:679:ILE:HG12	2.13	0.49
1:A:699:ARG:HG3	1:A:699:ARG:NH1	2.28	0.48
1:A:307:GLU:HG2	7:B:32:HOH:O	2.14	0.48
1:B:589:MET:HA	1:B:649:VAL:O	2.14	0.48
1:B:701:THR:HA	1:B:702:PRO:C	2.34	0.48
1:A:489:ASP:HB3	1:A:491:SER:OG	2.13	0.47
1:A:686:SER:HA	1:A:691:PHE:CG	2.50	0.47
1:A:321:THR:HG21	1:A:701:THR:HG22	1.96	0.47
1:B:614:MET:CE	1:B:632:GLU:HG3	2.45	0.47
2:B:750:HEM:O1A	3:B:760:H4B:N2	2.44	0.46
1:A:699:ARG:HG3	1:A:699:ARG:HH11	1.81	0.46
1:A:299:ARG:HE	1:A:318:LEU:HD13	1.81	0.45
1:A:478:GLN:HA	1:A:566:ALA:O	2.17	0.45
1:A:464:GLN:HB3	1:A:579:PHE:CE2	2.52	0.45
1:B:510:TRP:CD1	1:B:521:PRO:HG3	2.52	0.45
1:B:362:LEU:HD11	1:B:384:VAL:HG21	1.99	0.45
1:B:537:PRO:HA	1:B:538:PRO:HD3	1.90	0.45
1:B:571:LEU:C	1:B:571:LEU:HD23	2.37	0.44
1:B:600:ASP:HB2	1:B:603:ARG:HG2	1.98	0.44
1:B:584:PHE:CD1	2:B:750:HEM:CAC	3.01	0.44
1:A:445:HIS:C	1:A:445:HIS:CD2	2.90	0.44
1:A:686:SER:CB	1:B:682:PRO:HB2	2.43	0.43
1:A:423:LYS:HB2	1:A:423:LYS:HE2	1.96	0.43
5:B:800[B]:HW4:H4	5:B:800[B]:HW4:H8	1.76	0.43
1:B:455:LEU:HD23	1:B:587:TRP:HB3	2.00	0.43
1:B:684:SER:O	1:B:685:GLY:C	2.57	0.43
1:A:588:TYR:CD1	1:A:593:ILE:HD11	2.54	0.42
1:B:595:VAL:O	1:B:599:CYS:HB2	2.19	0.42
1:B:610:VAL:O	1:B:614:MET:HG3	2.19	0.42
1:A:371:ARG:CG	1:A:371:ARG:NH2	2.50	0.42
1:A:375:LYS:NZ	1:A:379:ASP:OD1	2.53	0.41
1:B:462:PHE:HB3	1:B:463:PRO:CD	2.50	0.41
1:B:567:VAL:O	1:B:583:PRO:HA	2.20	0.41
2:A:750:HEM:CMC	2:A:750:HEM:HBC2	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:614:MET:HE1	1:B:632:GLU:HG3	2.01	0.41
1:B:566:ALA:HA	1:B:584:PHE:O	2.21	0.41
1:A:706:TYR:OH	2:A:750:HEM:CGD	2.69	0.41
1:B:481:ARG:HD3	1:B:498:ASN:ND2	2.36	0.41
1:A:589:MET:HA	1:A:649:VAL:O	2.21	0.40
1:B:524:LEU:O	1:B:531:PRO:HA	2.20	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	405/422 (96%)	386 (95%)	18 (4%)	1 (0%)	47	46
1	B	409/422 (97%)	388 (95%)	21 (5%)	0	100	100
All	All	814/844 (96%)	774 (95%)	39 (5%)	1 (0%)	51	53

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	351	LYS

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%)	342 (94%)	23 (6%)	18	13
1	B	368/377 (98%)	358 (97%)	10 (3%)	44	46
All	All	733/754 (97%)	700 (96%)	33 (4%)	27	24

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	307	GLU
1	A	320	SER
1	A	330	ILE
1	A	336	MET
1	A	337	LEU
1	A	353	GLN
1	A	370	LYS
1	A	371	ARG
1	A	380	ARG
1	A	445	HIS
1	A	454	ASN
1	A	486	LYS
1	A	489	ASP
1	A	507	GLN
1	A	511	LYS
1	A	516	ARG
1	A	523	LEU
1	A	547	ARG
1	A	569	ASN
1	A	620	LYS
1	A	645	LYS
1	A	713	THR
1	A	715	VAL
1	B	350	THR
1	B	353	GLN
1	B	382	GLU
1	B	390	SER
1	B	423	LYS
1	B	445	HIS
1	B	455	LEU
1	B	481	ARG
1	B	540	LEU
1	B	547	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	454	ASN
1	A	507	GLN
1	A	527	ASN
1	A	569	ASN
1	A	642	GLN
1	A	697	ASN
1	A	712	ASN
1	B	353	GLN
1	B	364	GLN
1	B	385	ASN
1	B	425	GLN
1	B	454	ASN
1	B	508	GLN
1	B	601	ASN
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 10 ligands modelled in this entry, 1 is monoatomic - leaving 9 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
3	H4B	A	760	-	16,18,18	0.86	0	11,26,26	2.85	7 (63%)
5	HW4	B	800[B]	-	28,29,29	0.71	0	32,38,38	1.73	6 (18%)
2	HEM	B	750	1	41,50,50	2.06	10 (24%)	45,82,82	1.56	8 (17%)
5	HW4	A	800	-	28,29,29	0.80	2 (7%)	32,38,38	1.35	5 (15%)
3	H4B	B	760	-	16,18,18	1.26	3 (18%)	11,26,26	3.13	7 (63%)
5	HW4	B	800[A]	-	28,29,29	0.72	0	32,38,38	1.63	6 (18%)
4	ACT	B	860	-	3,3,3	0.75	0	3,3,3	0.84	0
4	ACT	A	860	-	3,3,3	0.93	0	3,3,3	0.33	0
2	HEM	A	750	1	41,50,50	1.97	12 (29%)	45,82,82	2.20	17 (37%)

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
3	H4B	A	760	-	-	0/8/17/17	0/2/2/2
5	HW4	B	800[B]	-	-	4/13/23/23	0/3/3/3
5	HW4	A	800	-	-	5/13/23/23	0/3/3/3
2	HEM	B	750	1	-	2/12/54/54	-
3	H4B	B	760	-	-	0/8/17/17	0/2/2/2
5	HW4	B	800[A]	-	-	2/13/23/23	0/3/3/3
2	HEM	A	750	1	-	6/12/54/54	-

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	750	HEM	C3D-C2D	7.40	1.52	1.36
2	B	750	HEM	C3D-C2D	7.04	1.51	1.36
2	B	750	HEM	C3C-C2C	-5.30	1.33	1.40
2	B	750	HEM	C3C-CAC	4.45	1.56	1.47
2	A	750	HEM	C3C-C2C	-3.12	1.36	1.40
2	B	750	HEM	FE-ND	3.11	2.12	1.96
2	A	750	HEM	CAB-C3B	3.11	1.55	1.47
2	A	750	HEM	CMD-C2D	3.02	1.57	1.50
2	B	750	HEM	CMB-C2B	2.99	1.57	1.50
3	B	760	H4B	C7-N8	2.87	1.49	1.44
2	A	750	HEM	CMB-C2B	2.75	1.56	1.50
2	A	750	HEM	O1D-CGD	2.70	1.31	1.22

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	750	HEM	CAA-C2A	2.66	1.55	1.52
2	B	750	HEM	CMD-C2D	2.64	1.56	1.50
2	A	750	HEM	C3C-CAC	2.38	1.52	1.47
3	B	760	H4B	C4-N3	2.36	1.37	1.33
2	B	750	HEM	C1A-NA	2.35	1.41	1.36
5	A	800	HW4	C22-C23	2.32	1.41	1.37
2	A	750	HEM	CMA-C3A	2.31	1.56	1.51
2	A	750	HEM	CMC-C2C	2.31	1.57	1.51
2	B	750	HEM	CAB-C3B	2.30	1.53	1.47
2	A	750	HEM	FE-NB	2.19	2.07	1.96
2	A	750	HEM	FE-ND	2.18	2.07	1.96
2	B	750	HEM	O1D-CGD	2.16	1.29	1.22
2	B	750	HEM	CAD-C3D	2.14	1.56	1.51
5	A	800	HW4	C24-C23	2.09	1.41	1.37
3	B	760	H4B	C2-N2	2.04	1.38	1.33

All (56) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	760	H4B	C8A-C4A-C4	6.17	120.05	114.57
2	A	750	HEM	C4D-ND-C1D	6.16	111.44	105.07
5	B	800[A]	HW4	C02-N01-C06	5.57	122.32	118.10
5	B	800[B]	HW4	C02-N01-C06	5.57	122.32	118.10
3	A	760	H4B	C8A-C4A-C4	5.37	119.34	114.57
2	B	750	HEM	C4D-ND-C1D	4.47	109.69	105.07
2	A	750	HEM	CBA-CAA-C2A	-4.30	105.28	112.62
2	A	750	HEM	CAD-CBD-CGD	-4.26	104.43	113.60
2	A	750	HEM	CHA-C4D-ND	4.13	129.48	124.38
5	B	800[A]	HW4	C05-C06-N01	-3.94	118.72	122.90
5	B	800[B]	HW4	C05-C06-N01	-3.94	118.72	122.90
3	B	760	H4B	N2-C2-N3	3.89	123.30	117.25
3	A	760	H4B	C4-C4A-N5	3.88	122.37	119.12
3	A	760	H4B	C2-N3-C4	3.86	122.07	115.93
3	B	760	H4B	C4-C4A-N5	3.83	122.33	119.12
2	A	750	HEM	C2C-C3C-C4C	3.67	109.46	106.90
3	B	760	H4B	C2-N3-C4	3.61	121.66	115.93
2	A	750	HEM	CMD-C2D-C1D	3.60	130.52	125.04
2	B	750	HEM	CMD-C2D-C1D	3.54	130.44	125.04
5	A	800	HW4	C5'-N1'-C2'	3.46	113.59	105.42
3	B	760	H4B	N1-C2-N3	-3.14	120.50	125.42
2	B	750	HEM	C4B-C3B-C2B	3.12	109.59	107.11
2	B	750	HEM	CAD-C3D-C4D	3.11	130.09	124.66

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	750	HEM	CAA-CBA-CGA	-3.07	105.17	113.76
5	B	800[B]	HW4	C24-C23-C22	-3.05	119.33	123.29
5	A	800	HW4	C05-C06-N01	-3.04	119.67	122.90
3	A	760	H4B	N1-C2-N3	-2.89	120.88	125.42
2	A	750	HEM	C1B-NB-C4B	2.81	107.98	105.07
3	A	760	H4B	N2-C2-N3	2.75	121.53	117.25
2	A	750	HEM	CMC-C2C-C3C	2.75	129.82	124.68
2	A	750	HEM	C4C-CHD-C1D	2.74	126.18	122.56
3	B	760	H4B	C2-N1-C8A	2.71	120.61	114.54
2	A	750	HEM	C1D-C2D-C3D	-2.67	104.15	106.96
2	B	750	HEM	C4C-CHD-C1D	2.53	125.90	122.56
3	B	760	H4B	C4A-C4-N3	-2.45	117.04	124.01
2	A	750	HEM	C4A-C3A-C2A	2.43	108.68	107.00
5	A	800	HW4	C24-C23-C22	-2.38	120.20	123.29
3	A	760	H4B	C4A-C4-N3	-2.29	117.50	124.01
5	B	800[A]	HW4	C07-C04-C05	-2.25	117.60	120.94
5	B	800[B]	HW4	C07-C04-C05	-2.25	117.60	120.94
5	B	800[A]	HW4	C5'-N1'-C2'	2.25	110.72	105.42
5	B	800[B]	HW4	C5'-N1'-C2'	2.25	110.72	105.42
2	A	750	HEM	C2B-C1B-NB	-2.24	107.18	109.84
2	A	750	HEM	C3B-C2B-C1B	2.23	108.14	106.49
5	B	800[A]	HW4	C5'-C4'-C3'	2.21	106.21	103.40
5	B	800[B]	HW4	C5'-C4'-C3'	2.21	106.21	103.40
2	A	750	HEM	CHC-C4B-C3B	2.20	127.93	124.57
3	A	760	H4B	C2-N1-C8A	2.19	119.45	114.54
5	A	800	HW4	C07-C04-C05	-2.18	117.71	120.94
5	B	800[A]	HW4	C24-C23-C22	-2.17	120.47	123.29
2	A	750	HEM	CAB-C3B-C2B	-2.17	121.46	128.60
5	A	800	HW4	C06-C08-C4'	-2.15	108.46	115.55
2	B	750	HEM	C1B-NB-C4B	2.10	107.25	105.07
2	A	750	HEM	CHD-C1D-C2D	2.08	128.23	124.98
2	A	750	HEM	C3C-C4C-NC	-2.06	107.06	110.94
2	B	750	HEM	C2C-C3C-C4C	2.02	108.31	106.90

There are no chirality outliers.

All (19) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	800[B]	HW4	C11-C10-O09-C3'
2	A	750	HEM	C4D-C3D-CAD-CBD
2	A	750	HEM	C2D-C3D-CAD-CBD
5	A	800	HW4	C12-C13-C14-C21

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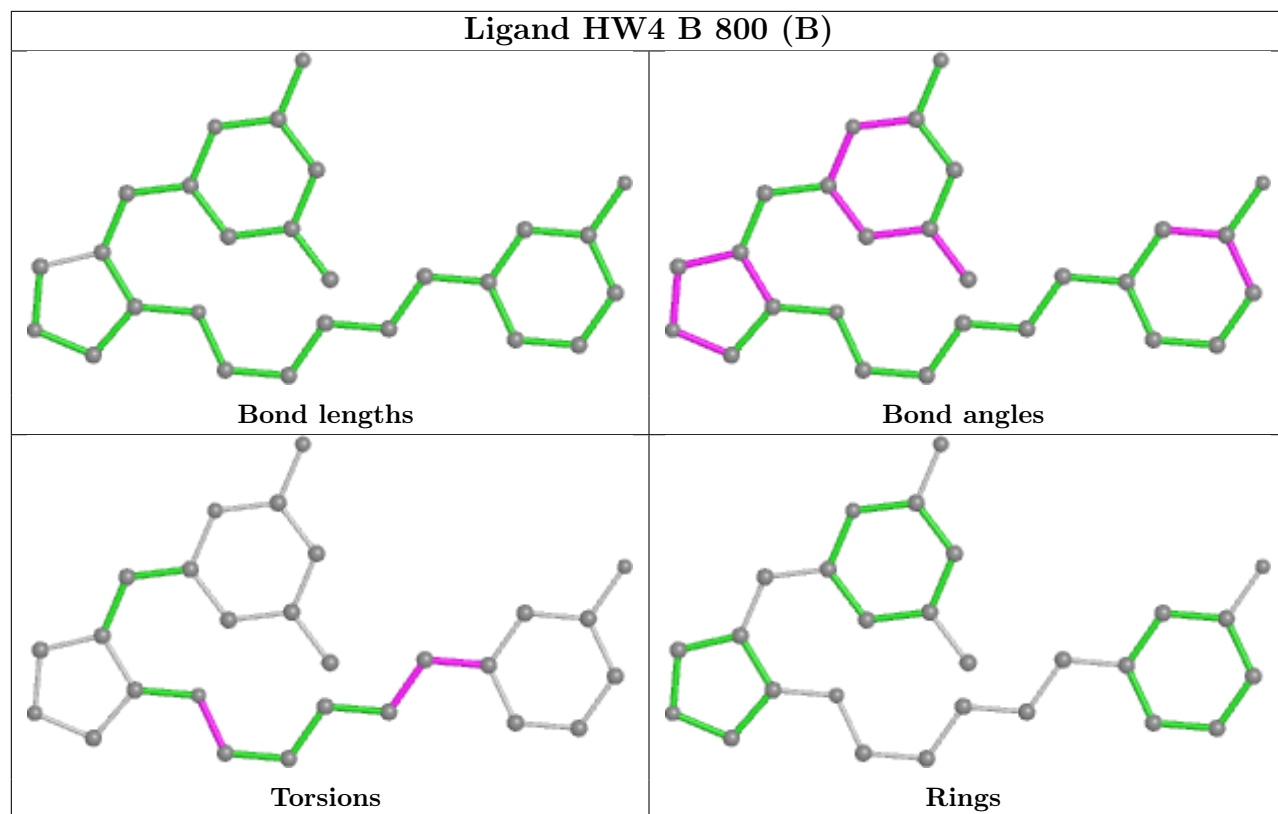
Mol	Chain	Res	Type	Atoms
5	B	800[A]	HW4	C11-C12-C13-C14
5	A	800	HW4	C10-C11-C12-C13
5	A	800	HW4	C11-C12-C13-C14
5	A	800	HW4	C11-C10-O09-C3'
5	B	800[B]	HW4	C13-C14-C21-C22
5	B	800[B]	HW4	C13-C14-C21-C26
5	B	800[B]	HW4	C12-C13-C14-C21
2	A	750	HEM	CAD-CBD-CGD-O1D
2	A	750	HEM	CAD-CBD-CGD-O2D
5	B	800[A]	HW4	C11-C10-O09-C3'
2	A	750	HEM	CAA-CBA-CGA-O2A
5	A	800	HW4	C4'-C3'-O09-C10
2	B	750	HEM	CAA-CBA-CGA-O2A
2	B	750	HEM	CAA-CBA-CGA-O1A
2	A	750	HEM	CAA-CBA-CGA-O1A

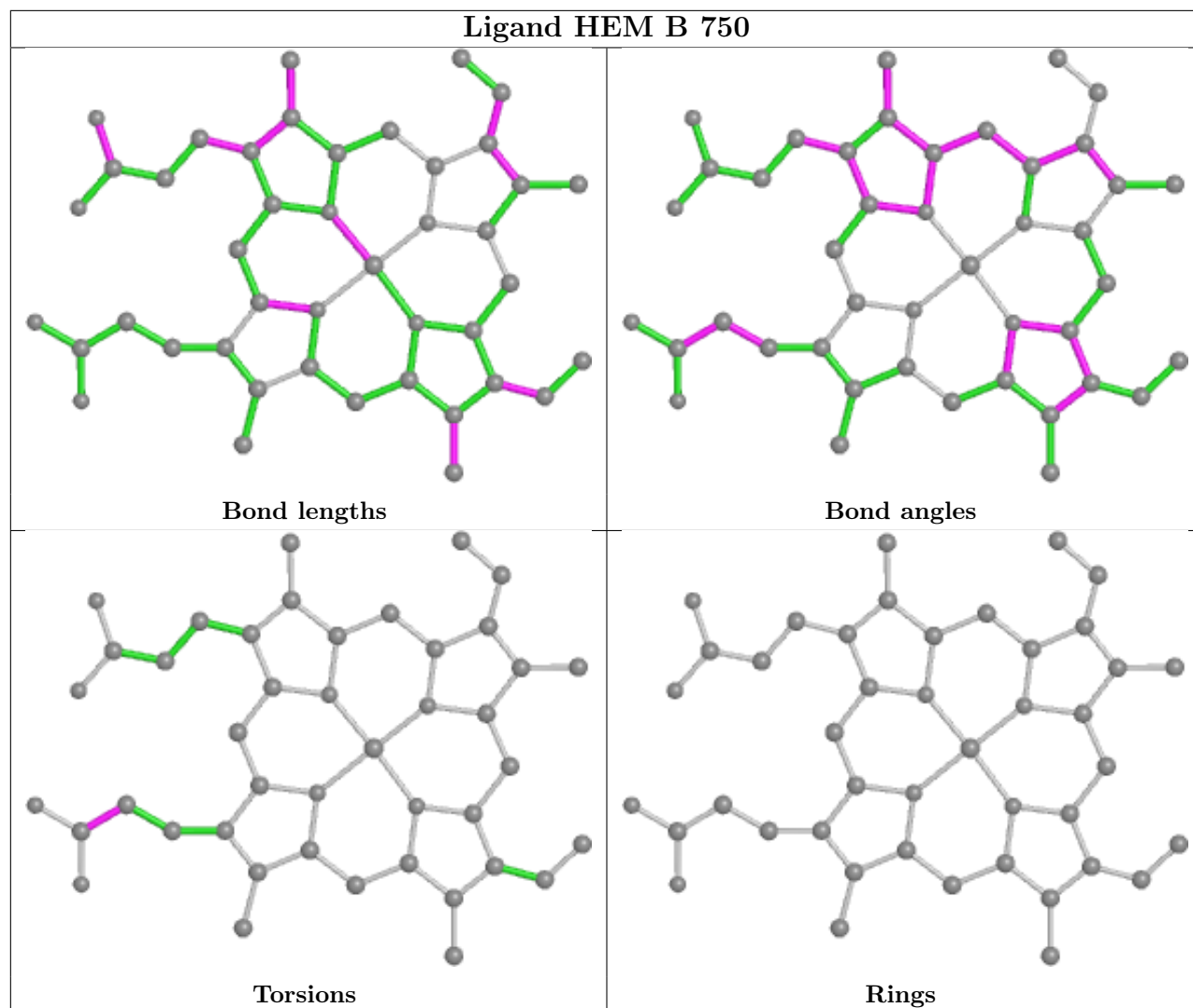
There are no ring outliers.

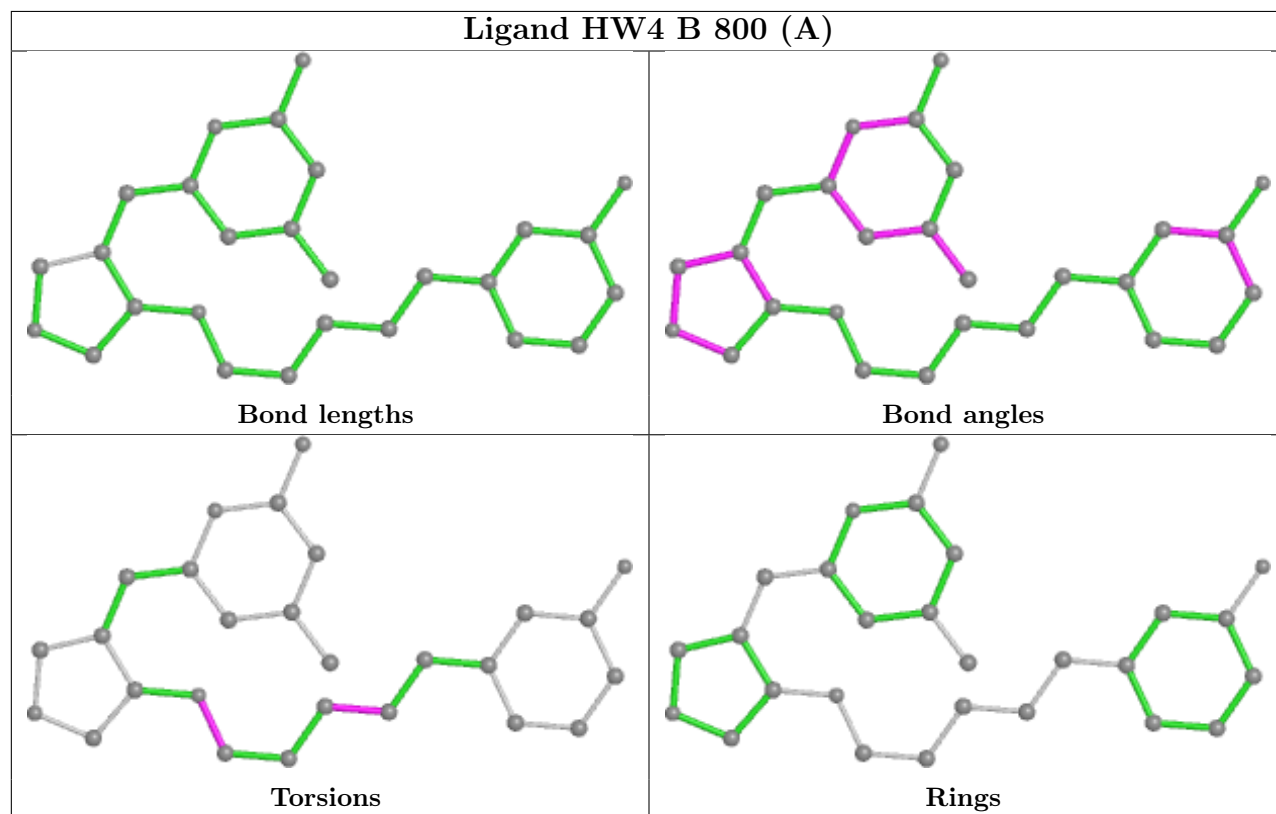
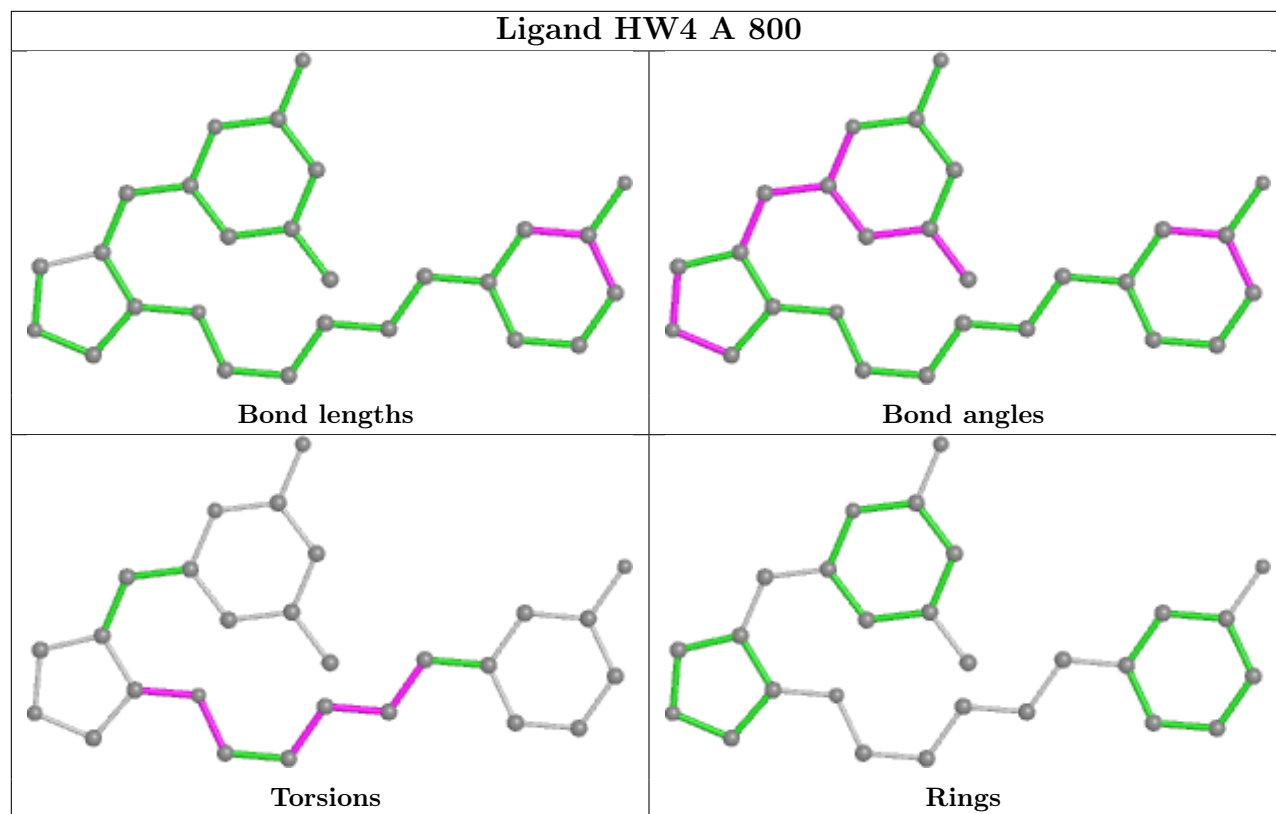
6 monomers are involved in 18 short contacts:

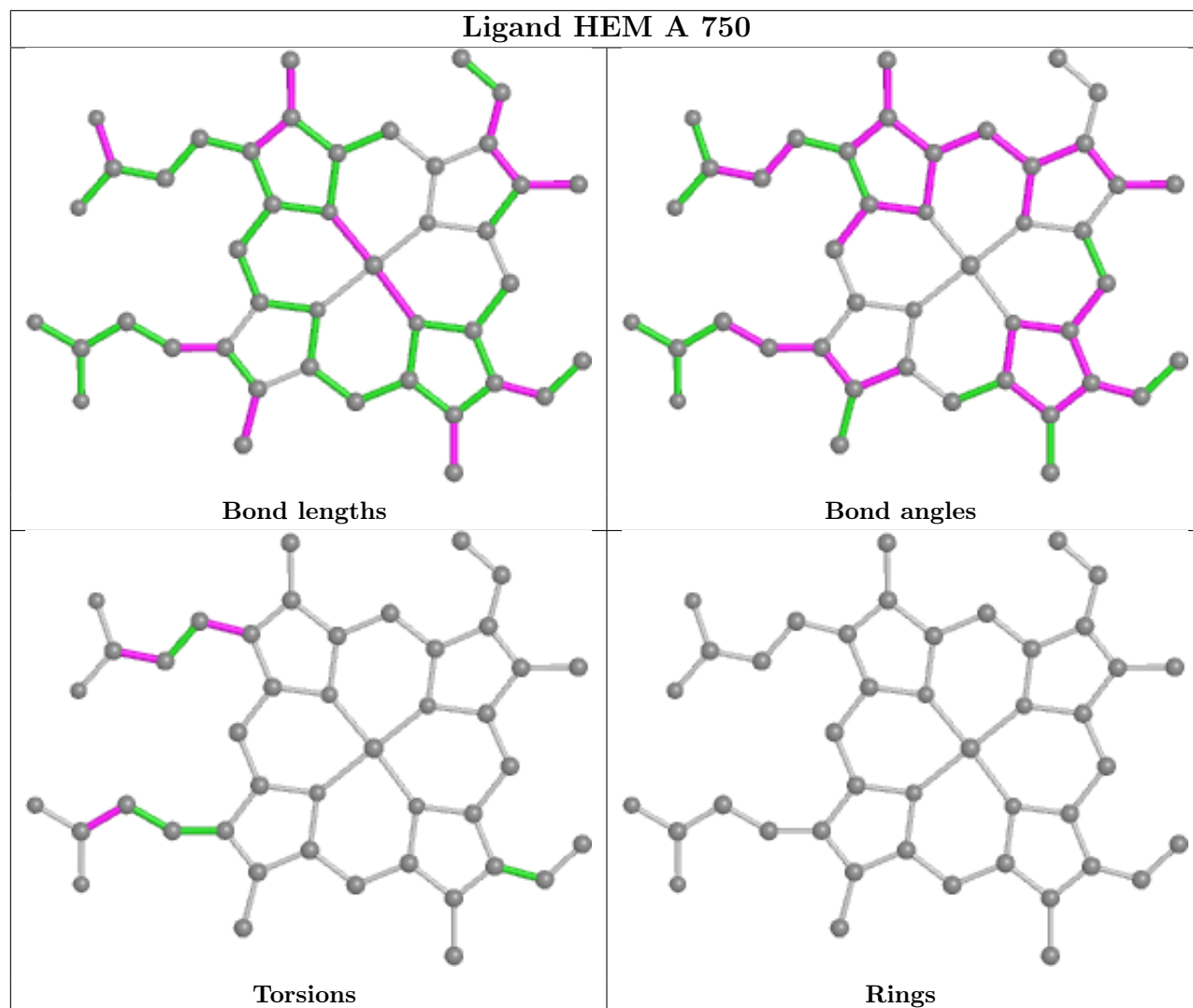
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	800[B]	HW4	1	0
2	B	750	HEM	8	0
5	A	800	HW4	3	0
3	B	760	H4B	1	0
5	B	800[A]	HW4	3	0
2	A	750	HEM	5	0

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 [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%)	1.07	86 (21%) 1 0	31, 61, 103, 131	0
1	B	411/422 (97%)	0.59	42 (10%) 6 10	31, 48, 74, 95	0
All	All	818/844 (96%)	0.83	128 (15%) 2 2	31, 53, 95, 131	0

All (128) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	716	TRP	8.1
1	B	300	PHE	7.5
1	B	348	VAL	6.4
1	A	488	PRO	6.2
1	A	486	LYS	5.5
1	A	352	ASP	5.4
1	A	355	PHE	5.3
1	B	718	GLY	5.3
1	A	351	LYS	5.2
1	B	351	LYS	5.0
1	B	706[A]	TYR	4.8
1	A	390	SER	4.5
1	A	714	HIS	4.5
1	A	300	PHE	4.4
1	A	415	CYS	4.3
1	B	350	THR	4.3
1	A	715	VAL	4.1
1	B	677	VAL	4.1
1	B	681	PRO	4.0
1	A	567	VAL	4.0
1	A	388	ILE	4.0
1	A	350	THR	3.9
1	A	713	THR	3.9
1	A	489	ASP	3.9

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Mol	Chain	Res	Type	RSRZ
1	A	679	ILE	3.9
1	A	392	SER	3.8
1	A	681	PRO	3.7
1	A	370	LYS	3.6
1	A	386	LYS	3.5
1	A	490	GLY	3.5
1	A	593	ILE	3.5
1	A	503	GLU	3.4
1	A	677	VAL	3.4
1	B	619	ARG	3.4
1	B	310	VAL	3.4
1	A	678	TRP	3.4
1	B	667	ARG	3.3
1	B	299	ARG	3.3
1	A	385	ASN	3.3
1	A	299	ARG	3.3
1	A	561	TRP	3.3
1	A	391	THR	3.2
1	A	591	THR	3.1
1	A	506	ILE	3.1
1	A	619	ARG	3.1
1	B	715	VAL	3.1
1	B	479	LEU	3.0
1	A	511	LYS	3.0
1	A	588	TYR	3.0
1	A	469	LYS	3.0
1	B	349	ARG	3.0
1	B	616	LEU	3.0
1	A	682	PRO	3.0
1	A	480	ILE	2.9
1	B	680	VAL	2.9
1	B	302	LYS	2.9
1	A	584	PHE	2.9
1	A	389	GLU	2.9
1	A	479	LEU	2.9
1	A	356	PRO	2.9
1	B	375	LYS	2.9
1	B	353	GLN	2.8
1	A	470	HIS	2.8
1	A	373	GLY	2.8
1	B	561	TRP	2.8
1	A	487	GLN	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	507	GLN	2.8
1	A	630	LEU	2.8
1	B	684	SER	2.7
1	B	352	ASP	2.7
1	B	691	PHE	2.7
1	A	566	ALA	2.7
1	A	393	THR	2.7
1	A	528	GLY	2.7
1	A	706	TYR	2.7
1	A	311	VAL	2.6
1	A	322	LEU	2.6
1	A	482	TYR	2.6
1	A	302	LYS	2.6
1	B	620	LYS	2.6
1	B	685	GLY	2.6
1	A	514	ARG	2.5
1	B	686	SER	2.5
1	A	552	ASP	2.5
1	A	491	SER	2.5
1	B	682	PRO	2.5
1	A	416	VAL	2.5
1	B	566	ALA	2.5
1	B	567	VAL	2.4
1	A	683	MET	2.4
1	B	584	PHE	2.4
1	A	467	ASP	2.4
1	A	704	PHE	2.4
1	A	354	LEU	2.4
1	A	676	TRP	2.4
1	A	493	LEU	2.4
1	A	499	VAL	2.4
1	A	595	VAL	2.4
1	B	679	ILE	2.4
1	A	712	ASN	2.3
1	B	678	TRP	2.3
1	A	680	VAL	2.3
1	A	685	GLY	2.3
1	A	394	TYR	2.3
1	A	572	LEU	2.3
1	A	485	TYR	2.3
1	A	531	PRO	2.3
1	B	389	GLU	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	591	THR	2.2
1	B	392	SER	2.2
1	A	551	PHE	2.2
1	A	585	SER	2.2
1	A	590	GLY	2.2
1	B	676	TRP	2.2
1	A	565	PRO	2.2
1	A	492	THR	2.2
1	A	568	SER	2.2
1	A	384	VAL	2.2
1	A	318	LEU	2.2
1	A	684	SER	2.2
1	A	353	GLN	2.1
1	B	630	LEU	2.1
1	B	683	MET	2.1
1	A	409	TRP	2.1
1	B	301	LEU	2.1
1	B	327	THR	2.0
1	A	504	ILE	2.0
1	B	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.

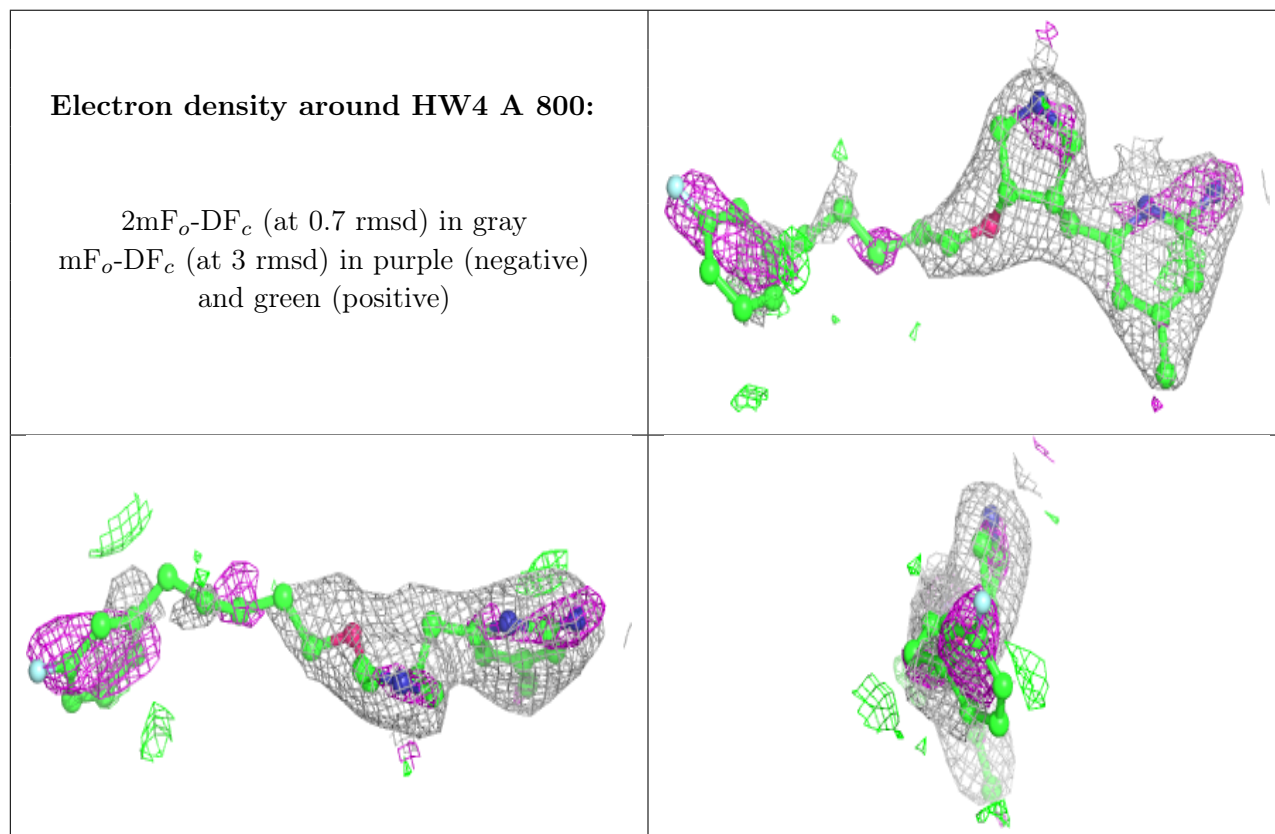
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	HW4	A	800	27/27	0.86	0.21	9,25,65,66	0
5	HW4	B	800[A]	27/27	0.89	0.50	36,55,70,73	11
5	HW4	B	800[B]	27/27	0.89	0.50	36,55,75,75	11

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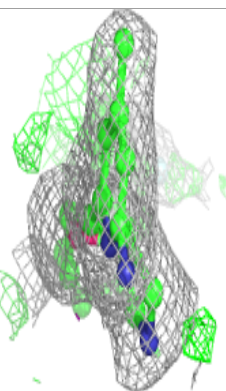
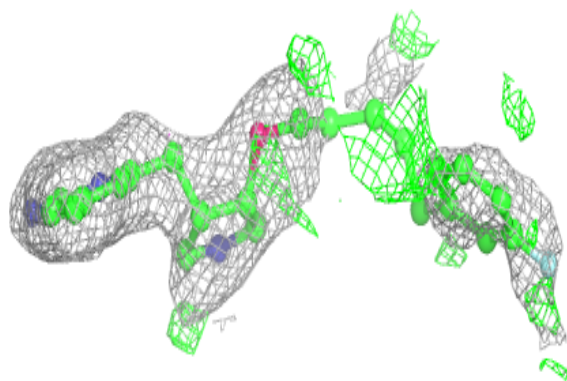
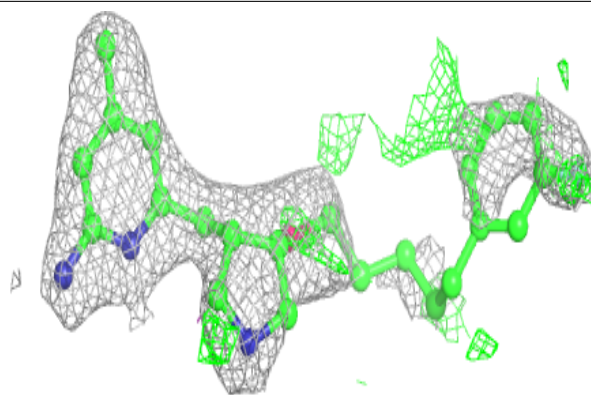
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	H4B	A	760	17/17	0.95	0.18	42,49,53,55	0
3	H4B	B	760	17/17	0.95	0.17	37,41,46,47	0
4	ACT	A	860	4/4	0.95	0.24	59,59,59,61	0
2	HEM	A	750	43/43	0.96	0.26	36,41,58,63	0
4	ACT	B	860	4/4	0.97	0.12	53,54,55,55	0
2	HEM	B	750	43/43	0.97	0.20	33,39,52,55	0
6	ZN	A	900	1/1	0.98	0.09	46,46,46,46	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.

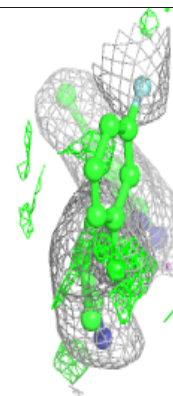
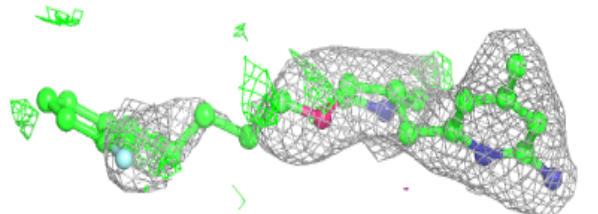
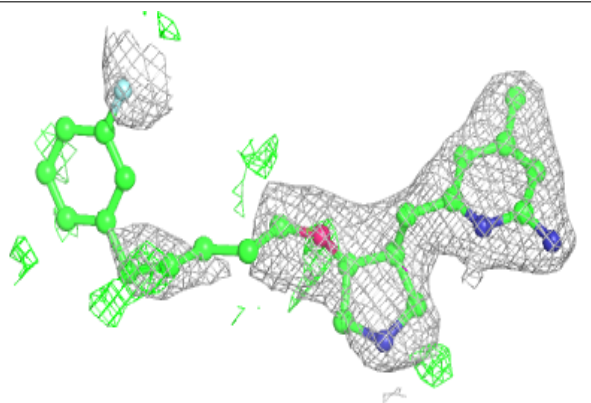


Electron density around HW4 B 800 (A):

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

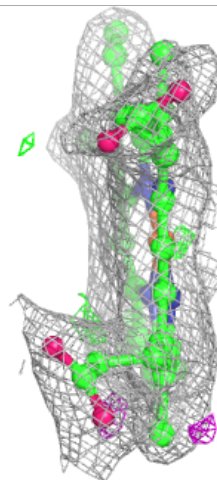
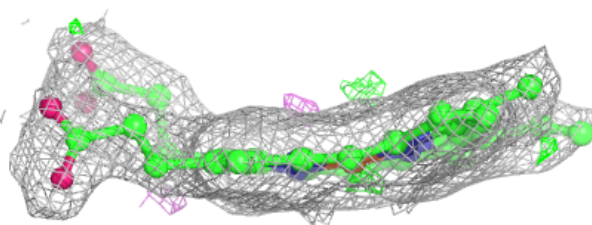
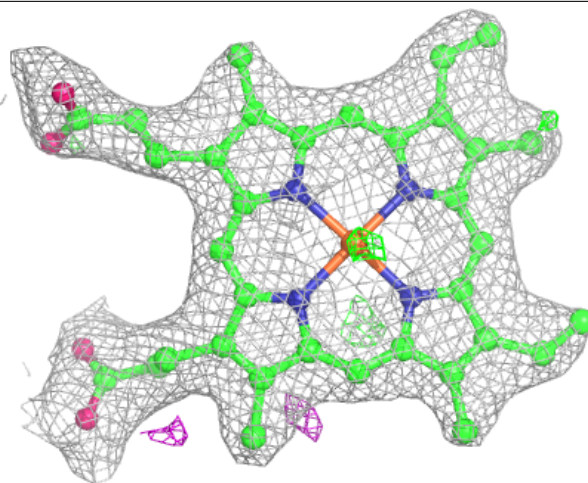
**Electron density around HW4 B 800 (B):**

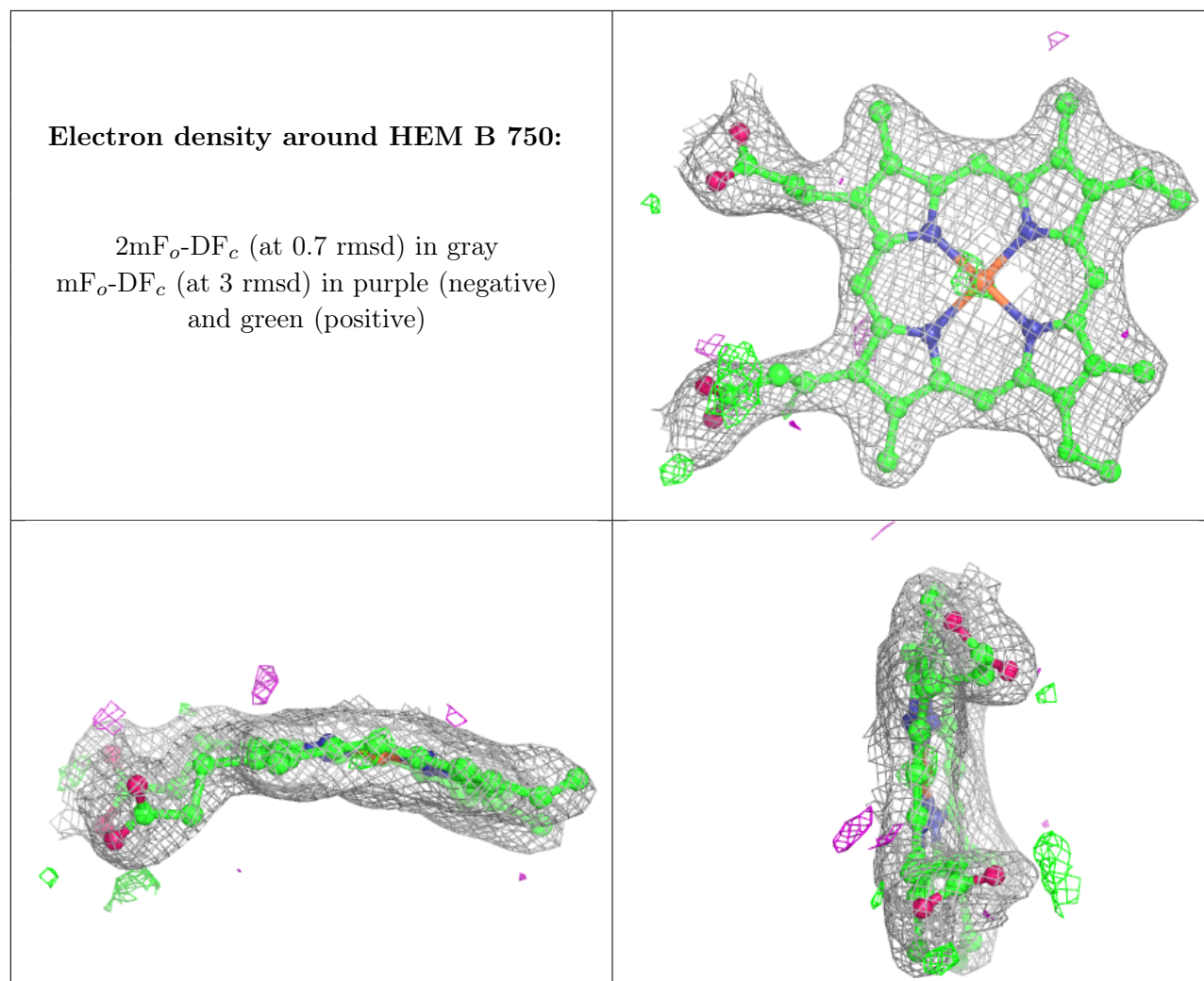
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



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.