



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 3, 2024 – 04:42 pm GMT

PDB ID : 4UGZ
Title : Structure of rat neuronal nitric oxide synthase heme domain in complex with N1-(3-(2-(6-Amino-4-methylpyridin-2-yl)ethyl)phenyl)-N1, N2-dimethylethane-1,2-diamine
Authors : Li, H.; Poulos, T.L.
Deposited on : 2015-03-23
Resolution : 2.08 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 : 1.8.4, 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

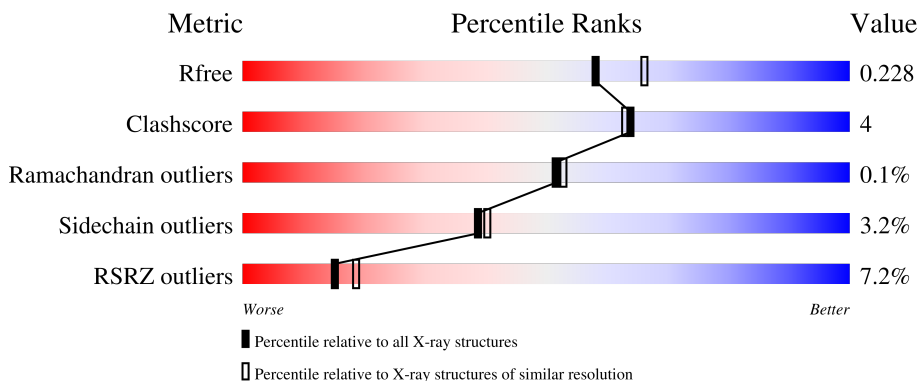
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION


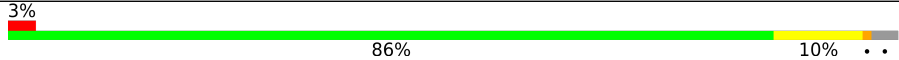
The reported resolution of this entry is 2.08 Å.

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	6189 (2.10-2.06)
Clashscore	141614	6738 (2.10-2.06)
Ramachandran outliers	138981	6663 (2.10-2.06)
Sidechain outliers	138945	6664 (2.10-2.06)
RSRZ outliers	127900	6057 (2.10-2.06)

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	

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	N	0	0
			22	18	4		

- Molecule 5 is ACETATE ION (three-letter code: ACT) (formula: C₂H₃O₂).



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

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

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

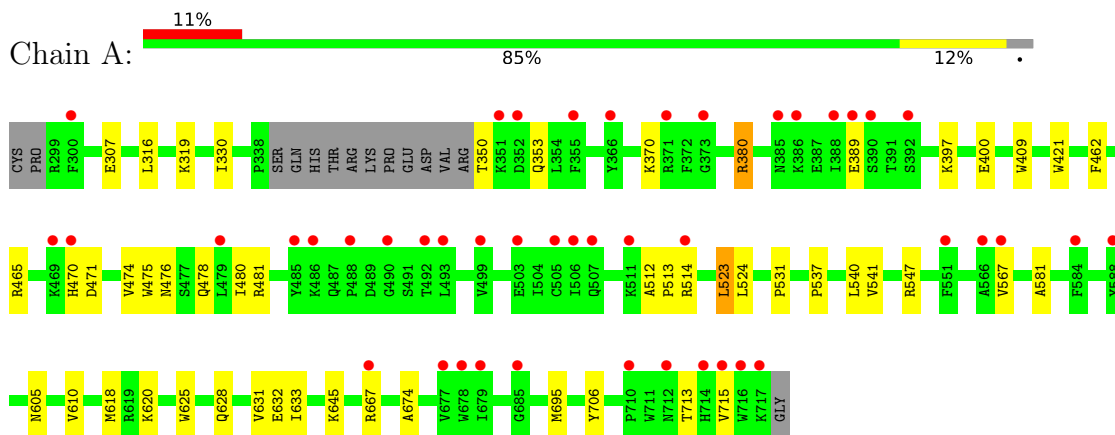
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	107	Total	O	0	0
			107	107		
7	B	149	Total	O	0	0
			149	149		

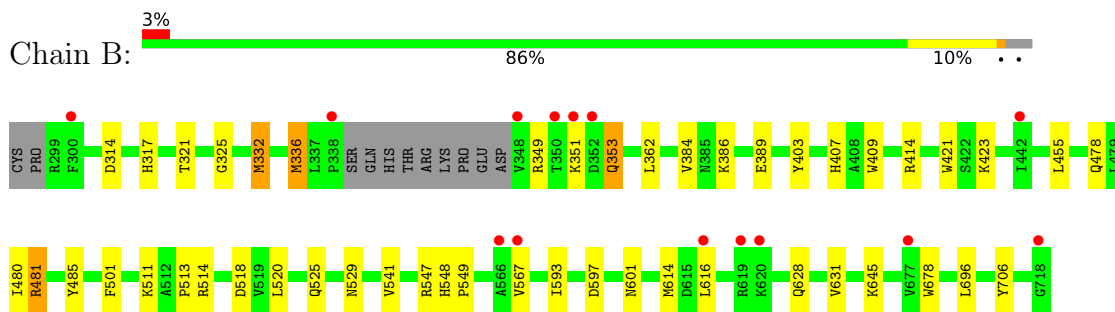
3 Residue-property plots [i](#)

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: NEURONAL NITRIC OXIDE SYNTHASE



• Molecule 1: NEURONAL NITRIC OXIDE SYNTHASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	51.87Å 111.23Å 164.88Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.84 – 2.08 40.84 – 2.08	Depositor EDS
% Data completeness (in resolution range)	97.7 (40.84-2.08) 97.9 (40.84-2.08)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.28 (at 2.08Å)	Xtrriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.188 , 0.227 0.188 , 0.228	Depositor DCC
R_{free} test set	2835 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	39.8	Xtrriage
Anisotropy	1.052	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 51.6	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	0.97	EDS
Total number of atoms	7102	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.11% 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, 4V4, H4B, HEM, ZN

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	0.38	0/3430	0.51	0/4654
1	B	0.40	0/3438	0.51	0/4661
All	All	0.39	0/6868	0.51	0/9315

There are no bond length outliers.

There are no bond angle outliers.

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	3328	0	3238	27	0
1	B	3345	0	3259	26	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	1	0
4	A	22	0	26	3	0
4	B	22	0	26	3	0
5	A	4	0	3	0	0
5	B	4	0	3	1	0
6	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	A	107	0	0	1	0
7	B	149	0	0	0	0
All	All	7102	0	6645	56	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 4.

All (56) 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:330:ILE:HD11	1:B:696:LEU:HD22	1.56	0.86
1:A:523:LEU:HD22	1:A:531:PRO:HB2	1.61	0.80
1:B:706:TYR:OH	2:B:750:HEM:O1D	2.03	0.74
2:B:750:HEM:HHC	2:B:750:HEM:HBB2	1.73	0.70
1:A:706:TYR:OH	2:A:750:HEM:O1D	2.10	0.69
1:B:478:GLN:HB2	1:B:481:ARG:HG3	1.75	0.68
2:A:750:HEM:HBB2	2:A:750:HEM:HHC	1.77	0.66
2:A:750:HEM:HMC2	2:A:750:HEM:HBC2	1.78	0.66
1:B:336:MET:HE2	3:B:760:H4B:H9	1.82	0.60
1:A:478:GLN:HB2	1:A:481:ARG:HG3	1.84	0.59
1:B:567:VAL:HG23	4:B:800:4V4:H05	1.85	0.57
1:A:316:LEU:HD12	1:A:319:LYS:HD2	1.86	0.57
1:A:307[A]:GLU:OE2	1:B:601:ASN:ND2	2.37	0.55
1:A:380:ARG:HD3	1:A:400:GLU:OE2	2.06	0.55
2:B:750:HEM:HBD2	4:B:800:4V4:C14	2.37	0.54
1:B:593:ILE:HA	1:B:597:ASP:HB2	1.92	0.52
1:A:480:ILE:HD13	1:A:541:VAL:HG13	1.93	0.51
1:B:414:ARG:HD3	1:B:678:TRP:CD2	2.45	0.51
1:A:475:TRP:HB2	1:A:523:LEU:HB3	1.92	0.51
1:A:350:THR:N	1:A:353:GLN:OE1	2.44	0.50
1:A:632:GLU:OE2	1:B:628:GLN:NE2	2.46	0.48
1:A:465:ARG:NH2	1:A:471:ASP:OD2	2.46	0.48
1:A:537:PRO:HB2	1:A:540:LEU:HG	1.96	0.47
1:B:353:GLN:HE21	1:B:353:GLN:HB3	1.53	0.47
1:A:628:GLN:HG3	1:B:631:VAL:HG11	1.95	0.47
1:A:567:VAL:HG23	4:A:800:4V4:H05	1.97	0.46
1:A:380:ARG:NH1	1:A:397:LYS:HG2	2.31	0.46
1:A:605:ASN:ND2	7:A:2076:HOH:O	2.48	0.46
1:A:674:ALA:HB3	1:A:695:MET:HB3	1.98	0.46
1:B:525:GLN:HG3	1:B:529:ASN:O	2.16	0.46
1:A:370:LYS:HE3	1:A:370:LYS:HB2	1.73	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:462:PHE:HB2	1:A:581:ALA:HB3	1.96	0.45
2:B:750:HEM:HBA1	4:B:800:4V4:H09	1.97	0.45
1:B:336:MET:HE2	1:B:336:MET:HB2	1.86	0.45
4:A:800:4V4:H16	4:A:800:4V4:H18	1.80	0.44
1:B:480:ILE:HD13	1:B:541:VAL:HG13	1.99	0.44
1:B:614:MET:HE2	1:B:616:LEU:HD21	1.99	0.44
1:B:314:ASP:OD1	1:B:317:HIS:ND1	2.39	0.44
1:B:513:PRO:HG2	1:B:518:ASP:CG	2.38	0.44
1:A:409:TRP:CE3	1:A:421:TRP:HA	2.52	0.43
1:A:512:ALA:HA	1:A:513:PRO:HD3	1.91	0.43
1:A:474:VAL:HG22	1:A:524:LEU:HD22	2.00	0.43
1:B:325:GLY:O	1:B:332:MET:HG3	2.19	0.43
1:B:501:PHE:HD2	1:B:520:LEU:HD13	1.83	0.42
1:A:524:LEU:O	1:A:531:PRO:HA	2.19	0.42
2:A:750:HEM:HBD1	4:A:800:4V4:C14	2.49	0.42
1:A:610:VAL:HG21	1:A:633:ILE:HD11	2.01	0.42
1:A:618:MET:HG2	1:A:625:TRP:CD2	2.54	0.42
1:A:631:VAL:HG11	1:B:628:GLN:HG3	2.02	0.41
1:B:409:TRP:CE3	1:B:421:TRP:HA	2.55	0.41
1:B:548:HIS:CG	1:B:549:PRO:HD2	2.56	0.41
2:B:750:HEM:HMB2	5:B:860:ACT:H2	2.01	0.41
1:B:386:LYS:HA	1:B:389:GLU:HG2	2.02	0.41
1:B:403:TYR:CE1	1:B:407:HIS:CE1	3.09	0.41
1:B:362:LEU:HD11	1:B:384:VAL:HG21	2.04	0.40
1:B:485:TYR:CZ	1:B:514:ARG:HA	2.56	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	407/422 (96%)	392 (96%)	14 (3%)	1 (0%)	47 47

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	407/422 (96%)	394 (97%)	13 (3%)	0	100	100
All	All	814/844 (96%)	786 (97%)	27 (3%)	1 (0%)	51	53

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	514	ARG

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	366/377 (97%)	355 (97%)	11 (3%)	41	43
1	B	366/377 (97%)	354 (97%)	12 (3%)	38	39
All	All	732/754 (97%)	709 (97%)	23 (3%)	39	41

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	380	ARG
1	A	389	GLU
1	A	470	HIS
1	A	476	ASN
1	A	523	LEU
1	A	547	ARG
1	A	620	LYS
1	A	645	LYS
1	A	667	ARG
1	A	713	THR
1	A	715	VAL
1	B	321	THR
1	B	332	MET
1	B	336	MET
1	B	349	ARG

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Mol	Chain	Res	Type
1	B	351	LYS
1	B	353	GLN
1	B	423	LYS
1	B	455	LEU
1	B	481	ARG
1	B	511	LYS
1	B	547	ARG
1	B	645	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	425	GLN
1	A	605	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 1 is monoatomic - leaving 8 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	HEM	B	750	1	41,50,50	1.98	9 (21%)	45,82,82	1.94	12 (26%)
4	4V4	B	800	-	23,23,23	0.54	0	30,30,30	1.31	5 (16%)
5	ACT	B	860	-	3,3,3	0.83	0	3,3,3	0.74	0
5	ACT	A	860	-	3,3,3	0.87	0	3,3,3	1.11	0
4	4V4	A	800	-	23,23,23	0.47	0	30,30,30	1.32	3 (10%)
3	H4B	A	760	-	16,18,18	0.79	0	11,26,26	2.66	5 (45%)
3	H4B	B	760	-	16,18,18	1.17	2 (12%)	11,26,26	2.39	4 (36%)
2	HEM	A	750	1	41,50,50	1.98	5 (12%)	45,82,82	2.01	11 (24%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	HEM	B	750	1	-	2/12/54/54	-
4	4V4	B	800	-	-	1/13/13/13	0/2/2/2
4	4V4	A	800	-	-	2/13/13/13	0/2/2/2
3	H4B	A	760	-	-	0/8/17/17	0/2/2/2
3	H4B	B	760	-	-	0/8/17/17	0/2/2/2
2	HEM	A	750	1	-	4/12/54/54	-

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	750	HEM	C3D-C2D	7.97	1.53	1.36
2	B	750	HEM	C3D-C2D	7.76	1.53	1.36
2	B	750	HEM	C3C-CAC	4.08	1.56	1.47
2	A	750	HEM	C3C-CAC	3.87	1.55	1.47
2	A	750	HEM	FE-NB	3.80	2.15	1.96
2	B	750	HEM	C3C-C2C	-3.63	1.35	1.40
2	A	750	HEM	C3C-C2C	-3.40	1.35	1.40
2	B	750	HEM	FE-ND	3.27	2.13	1.96
2	A	750	HEM	CAB-C3B	3.15	1.56	1.47
2	B	750	HEM	CAB-C3B	2.89	1.55	1.47
3	B	760	H4B	C7-C6	2.51	1.54	1.52
3	B	760	H4B	C4A-C4	-2.37	1.38	1.41
2	B	750	HEM	FE-NB	2.26	2.08	1.96
2	B	750	HEM	CMB-C2B	2.19	1.55	1.50
2	B	750	HEM	CMD-C2D	2.16	1.55	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	750	HEM	CAA-C2A	2.13	1.55	1.52

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	750	HEM	C4D-ND-C1D	6.70	111.99	105.07
2	B	750	HEM	C4D-ND-C1D	6.69	111.98	105.07
3	B	760	H4B	C8A-C4A-C4	5.54	119.49	114.57
2	A	750	HEM	C4C-CHD-C1D	5.03	129.19	122.56
3	A	760	H4B	C4-C4A-N5	4.53	122.92	119.12
3	A	760	H4B	C8A-C4A-C4	4.37	118.45	114.57
4	A	800	4V4	C02-N01-C06	3.99	121.12	118.10
4	B	800	4V4	C02-N01-C06	3.88	121.04	118.10
2	B	750	HEM	CAD-C3D-C4D	3.67	131.07	124.66
2	A	750	HEM	CAD-CBD-CGD	-3.62	105.82	113.60
3	A	760	H4B	N1-C2-N3	-3.32	120.20	125.42
2	B	750	HEM	CAD-CBD-CGD	-3.28	106.55	113.60
2	B	750	HEM	C4C-CHD-C1D	3.27	126.87	122.56
2	B	750	HEM	C4B-CHC-C1C	3.26	126.85	122.56
2	A	750	HEM	CBA-CAA-C2A	-3.01	107.49	112.62
3	A	760	H4B	C2-N1-C8A	2.98	121.21	114.54
3	B	760	H4B	N1-C2-N3	-2.96	120.77	125.42
2	B	750	HEM	C1B-NB-C4B	2.95	108.12	105.07
2	A	750	HEM	C3B-C2B-C1B	2.91	108.65	106.49
3	A	760	H4B	C2-N3-C4	2.91	120.55	115.93
2	B	750	HEM	CAD-C3D-C2D	-2.81	122.64	127.88
2	A	750	HEM	CMD-C2D-C1D	2.74	129.21	125.04
2	A	750	HEM	CHA-C4D-ND	2.69	127.71	124.38
4	B	800	4V4	C21-N20-C19	2.60	119.46	111.95
3	B	760	H4B	C2-N1-C8A	2.59	120.35	114.54
3	B	760	H4B	C2-N3-C4	2.59	120.04	115.93
2	A	750	HEM	CHB-C1B-NB	2.54	127.51	124.38
4	A	800	4V4	C21-N20-C19	2.52	119.23	111.95
4	B	800	4V4	C19-C18-N17	-2.52	103.19	112.31
2	B	750	HEM	C3D-C4D-ND	-2.51	107.38	110.17
4	B	800	4V4	C05-C06-N01	-2.35	120.41	122.90
2	B	750	HEM	CHC-C4B-C3B	2.32	128.13	124.57
2	A	750	HEM	C1B-NB-C4B	2.32	107.47	105.07
4	A	800	4V4	C05-C06-N01	-2.23	120.54	122.90
4	B	800	4V4	C08-C06-N01	2.22	119.26	115.95
2	A	750	HEM	C4B-CHC-C1C	2.18	125.44	122.56
2	B	750	HEM	C1D-C2D-C3D	-2.08	104.77	106.96

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	750	HEM	CHC-C4B-C3B	2.08	127.76	124.57
2	B	750	HEM	CMD-C2D-C1D	2.08	128.20	125.04
2	B	750	HEM	CBA-CAA-C2A	-2.01	109.19	112.62

There are no chirality outliers.

All (9) torsion outliers are listed below:

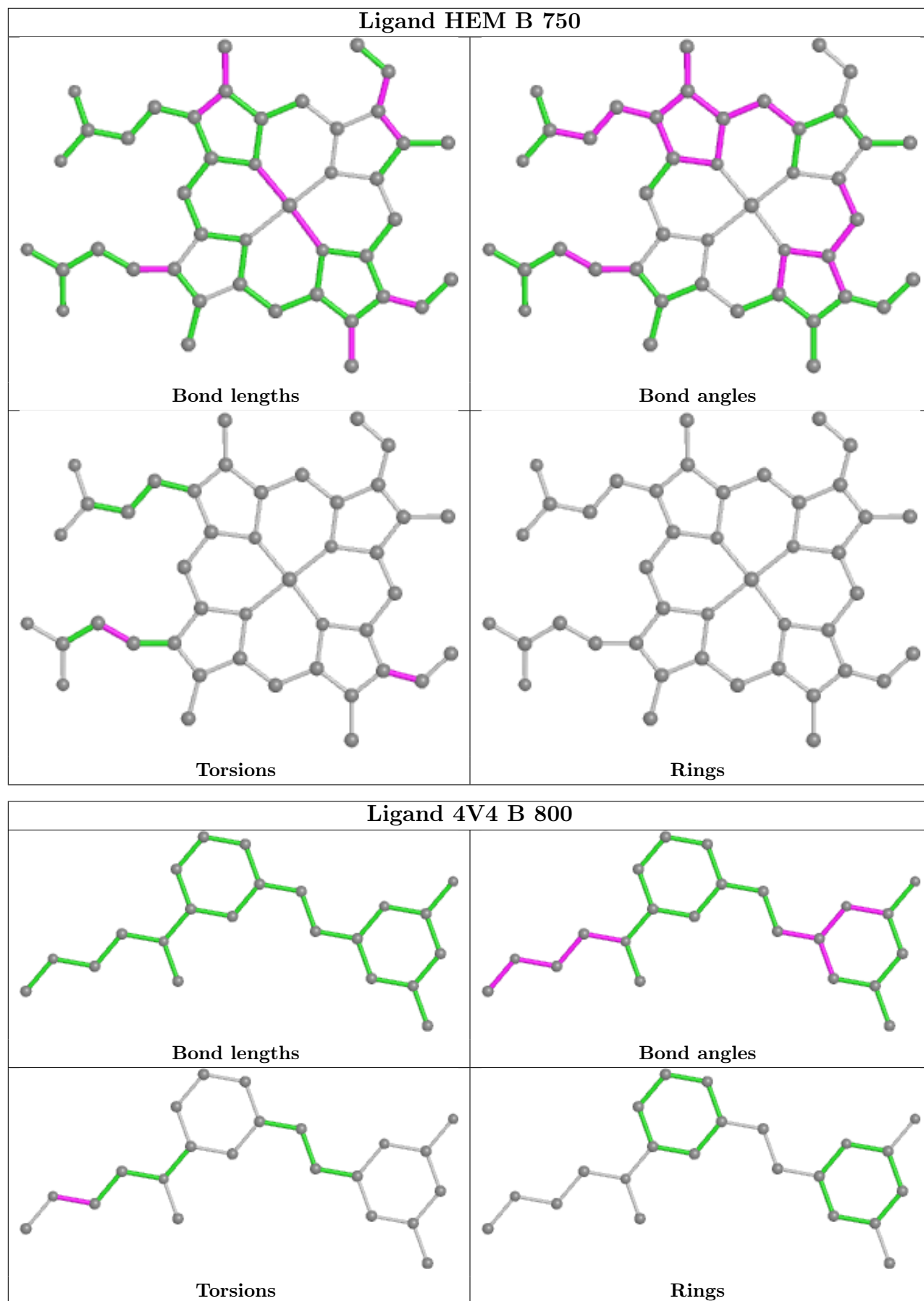
Mol	Chain	Res	Type	Atoms
2	A	750	HEM	C2A-CAA-CBA-CGA
4	A	800	4V4	N17-C18-C19-N20
2	B	750	HEM	C2A-CAA-CBA-CGA
4	B	800	4V4	C18-C19-N20-C21
2	A	750	HEM	C4B-C3B-CAB-CBB
2	B	750	HEM	C4B-C3B-CAB-CBB
4	A	800	4V4	C18-C19-N20-C21
2	A	750	HEM	CAD-CBD-CGD-O2D
2	A	750	HEM	CAD-CBD-CGD-O1D

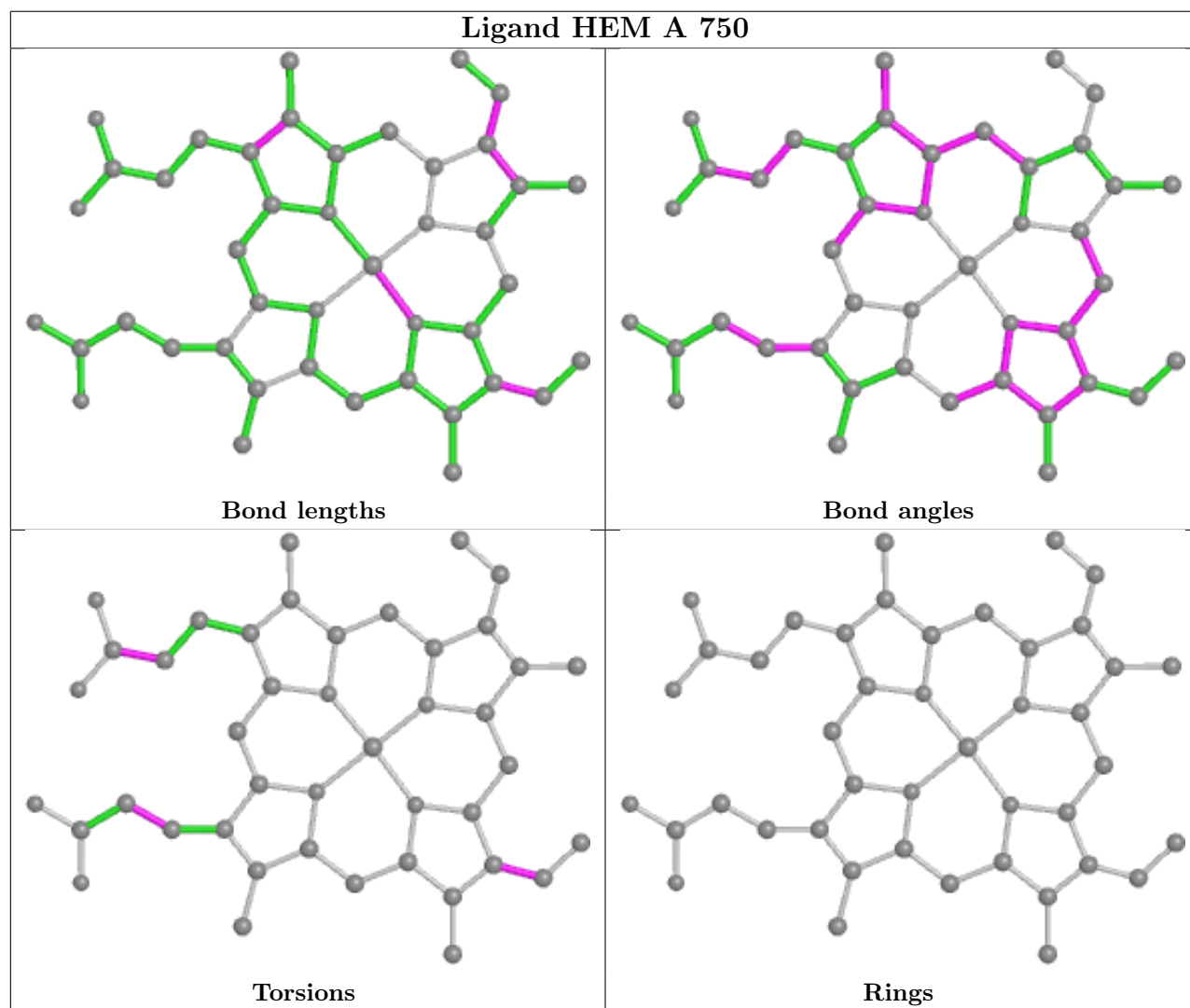
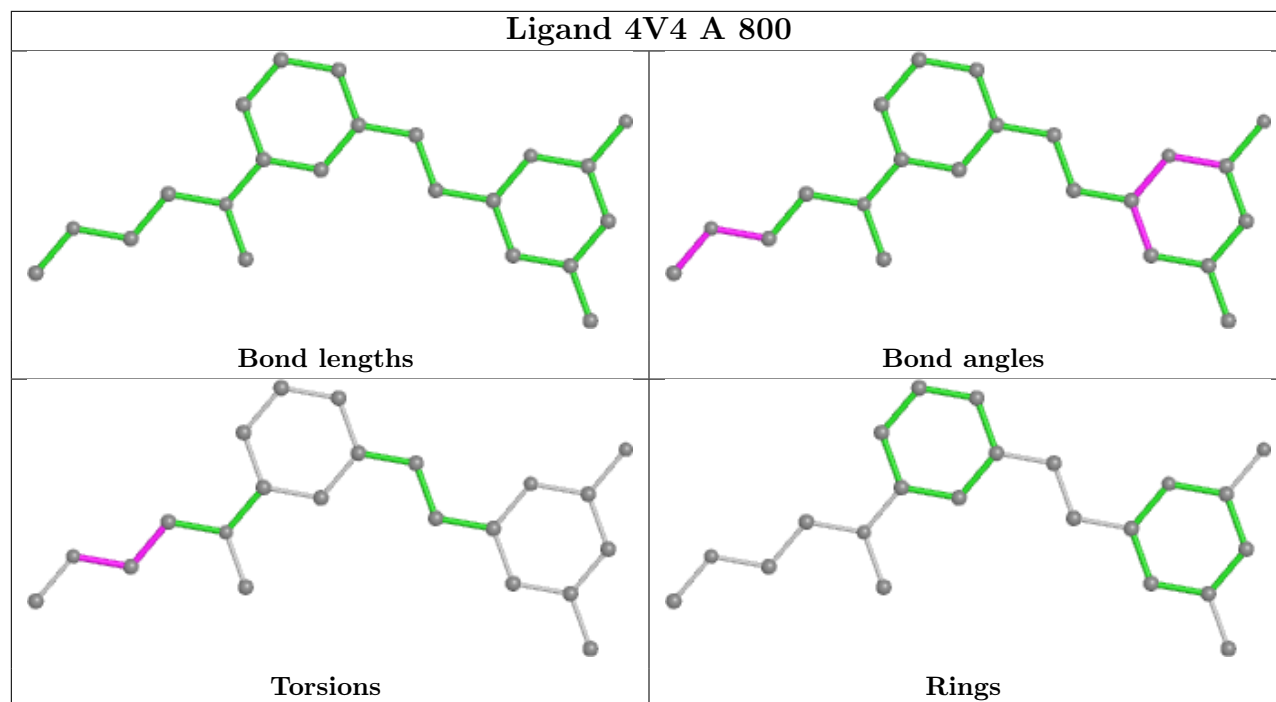
There are no ring outliers.

6 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	750	HEM	5	0
4	B	800	4V4	3	0
5	B	860	ACT	1	0
4	A	800	4V4	3	0
3	B	760	H4B	1	0
2	A	750	HEM	4	0

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





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	408/422 (96%)	0.53	45 (11%) 5 7	19, 49, 94, 125	0
1	B	411/422 (97%)	0.07	14 (3%) 45 50	20, 39, 70, 95	0
All	All	819/844 (97%)	0.30	59 (7%) 15 19	19, 43, 88, 125	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	300	PHE	7.0
1	A	355	PHE	5.4
1	A	715	VAL	5.2
1	A	717	LYS	5.1
1	A	488	PRO	4.8
1	A	716	TRP	4.8
1	B	350	THR	3.8
1	A	493	LEU	3.7
1	B	619	ARG	3.6
1	A	503	GLU	3.4
1	A	392	SER	3.4
1	A	551	PHE	3.4
1	A	714	HIS	3.3
1	A	712	ASN	3.2
1	A	678	TRP	3.2
1	A	386	LYS	3.1
1	A	506	ILE	3.0
1	A	492	THR	2.9
1	A	352	ASP	2.9
1	A	567	VAL	2.9
1	A	390	SER	2.8
1	A	490	GLY	2.8
1	A	679	ILE	2.8
1	A	486	LYS	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	351	LYS	2.8
1	A	389	GLU	2.8
1	B	348	VAL	2.7
1	A	507	GLN	2.7
1	A	388	ILE	2.6
1	A	373	GLY	2.6
1	A	366	TYR	2.6
1	A	677	VAL	2.6
1	A	469	LYS	2.5
1	B	351	LYS	2.4
1	A	385	ASN	2.4
1	A	566	ALA	2.3
1	B	338	PRO	2.3
1	B	677	VAL	2.3
1	B	718	GLY	2.3
1	B	616	LEU	2.3
1	B	352	ASP	2.3
1	A	371	ARG	2.2
1	A	479	LEU	2.2
1	B	567	VAL	2.2
1	A	667	ARG	2.2
1	A	584	PHE	2.1
1	A	485	TYR	2.1
1	A	514	ARG	2.1
1	A	588	TYR	2.1
1	A	710	PRO	2.1
1	A	505	CYS	2.1
1	A	685	GLY	2.1
1	A	470	HIS	2.1
1	B	442	ILE	2.0
1	A	499	VAL	2.0
1	A	300	PHE	2.0
1	B	566	ALA	2.0
1	A	511	LYS	2.0
1	B	620	LYS	2.0

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

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

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

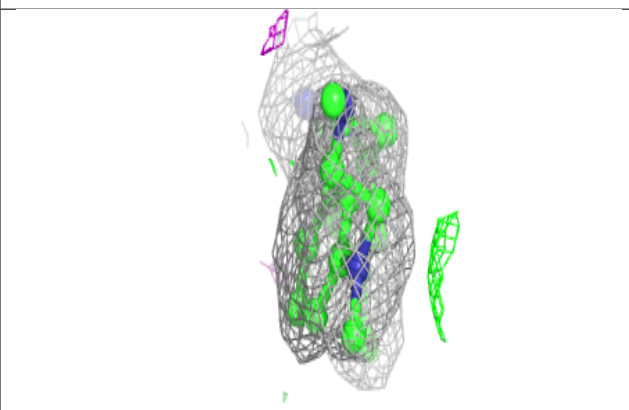
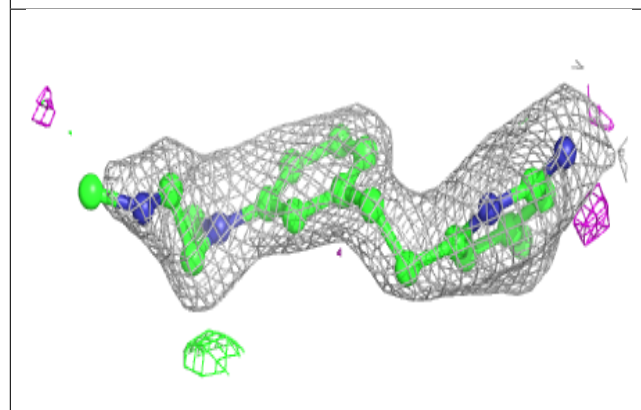
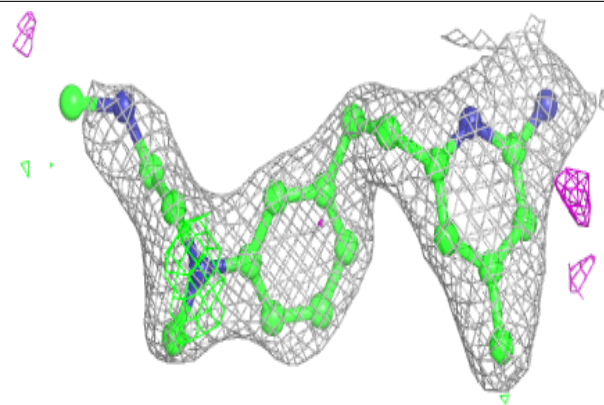
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	H4B	A	760	17/17	0.93	0.16	31,36,42,42	0
4	4V4	B	800	22/22	0.93	0.20	23,34,52,57	0
5	ACT	B	860	4/4	0.93	0.23	28,32,36,43	0
3	H4B	B	760	17/17	0.94	0.17	18,31,37,40	0
4	4V4	A	800	22/22	0.95	0.26	27,40,61,66	0
2	HEM	B	750	43/43	0.96	0.20	18,31,50,52	0
5	ACT	A	860	4/4	0.96	0.25	17,20,36,39	0
2	HEM	A	750	43/43	0.96	0.22	21,30,46,52	0
6	ZN	A	900	1/1	1.00	0.07	33,33,33,33	0

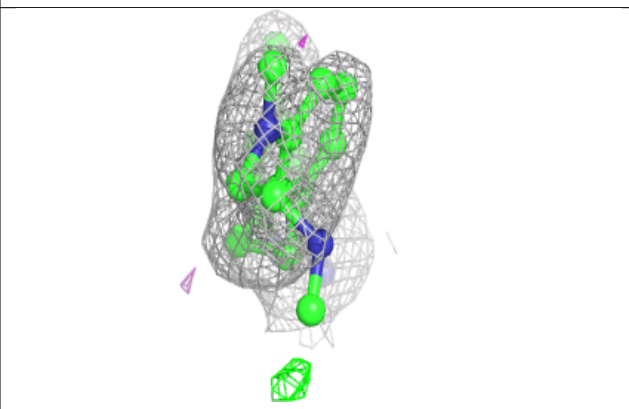
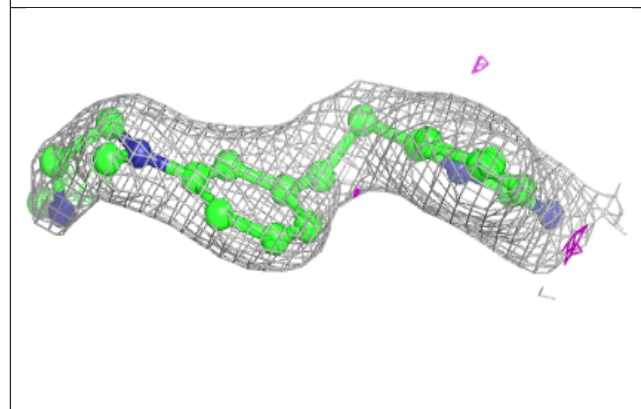
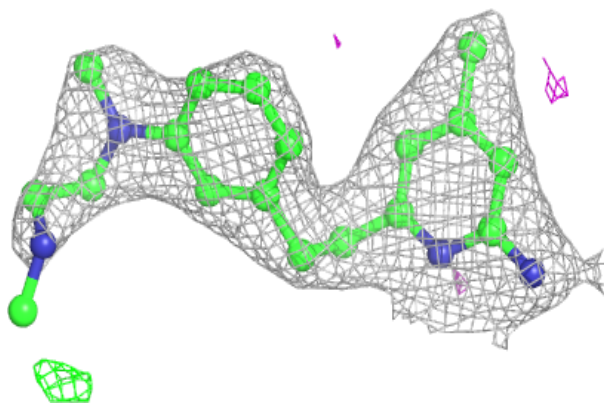
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around 4V4 B 800:

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

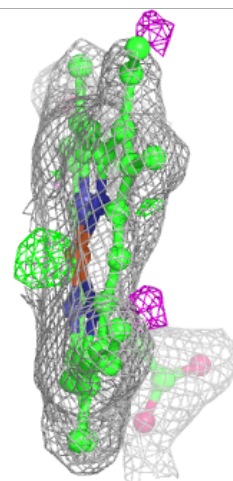
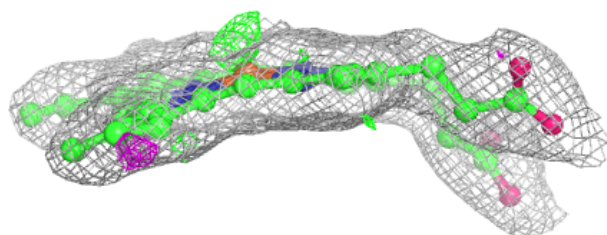
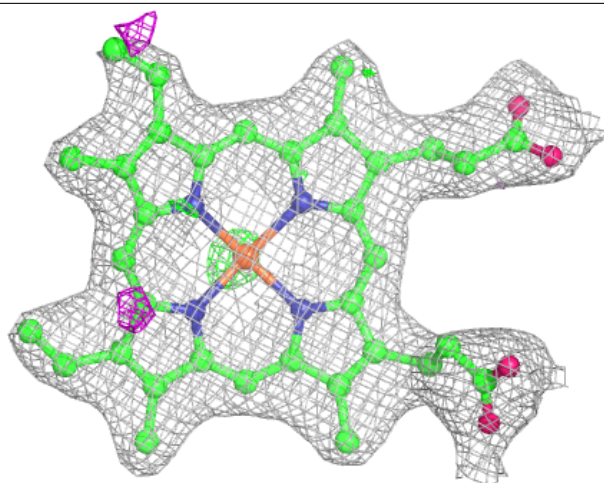
**Electron density around 4V4 A 800:**

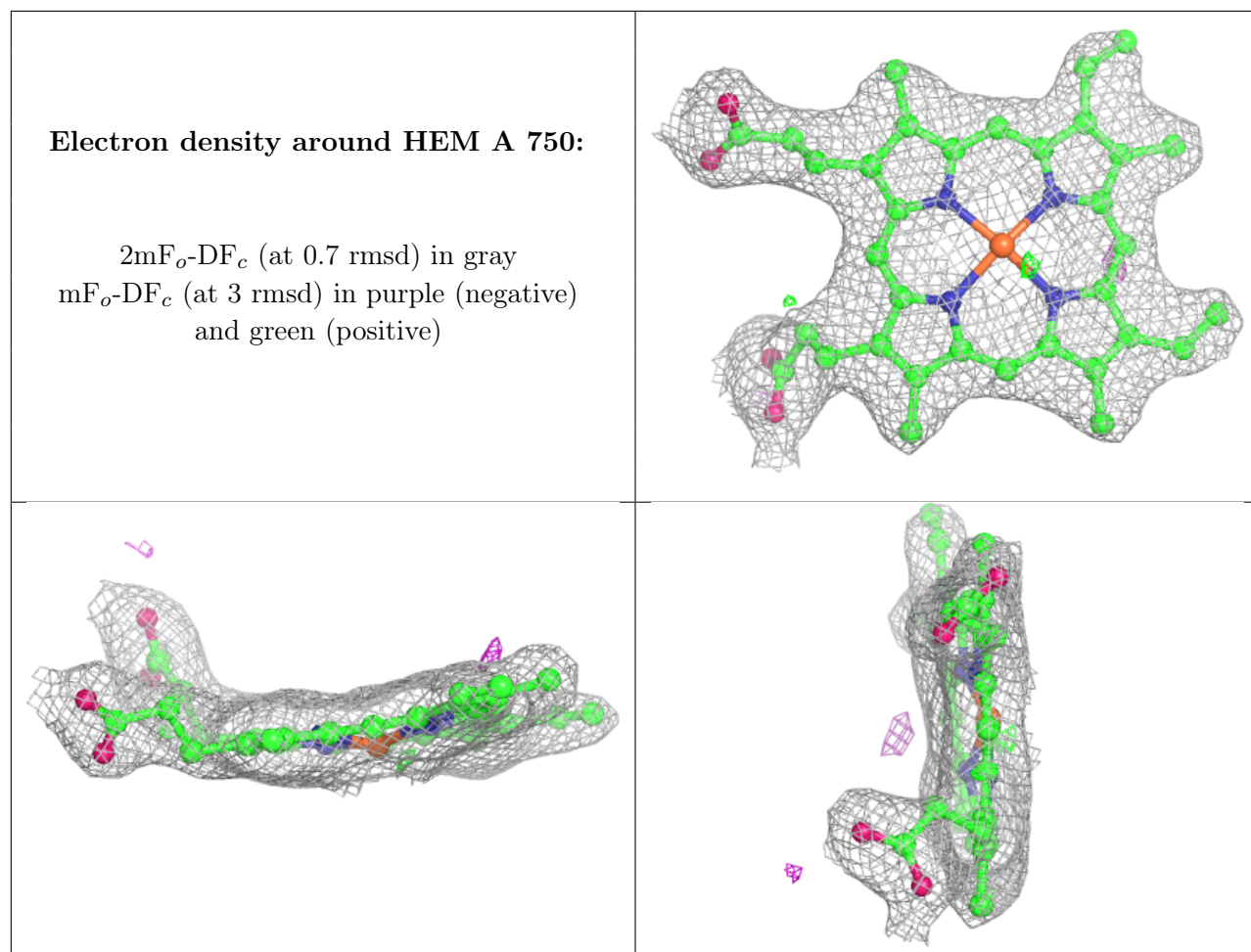
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around HEM B 750:

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





6.5 Other polymers [i](#)

There are no such residues in this entry.