



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 11:45 PM UTC

PDB ID : 3N6A / pdb_00003n6a
Title : Structure of endothelial nitric oxide synthase N368D/V106M double mutant heme domain complexed with 4-(2-(5-(2-(6-amino-4-methylpyridin-2-yl)ethyl)pyridin-3-yl)ethyl)-6-methylpyridin-2-amine
Authors : Delker, S.L.; Li, H.; Poulos, T.L.
Deposited on : 2010-05-25
Resolution : 2.49 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

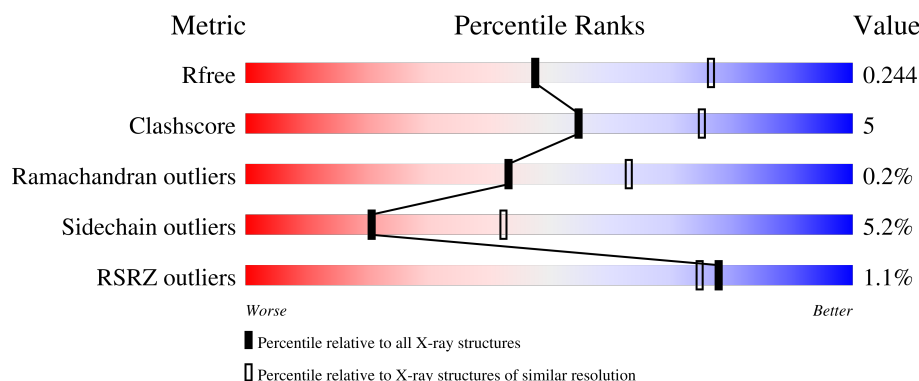
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.49 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	444	78% 11% • 9%
1	B	444	79% 11% • 9%

2 Entry composition [i](#)

There are 9 unique types of molecules in this entry. The entry contains 6785 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.

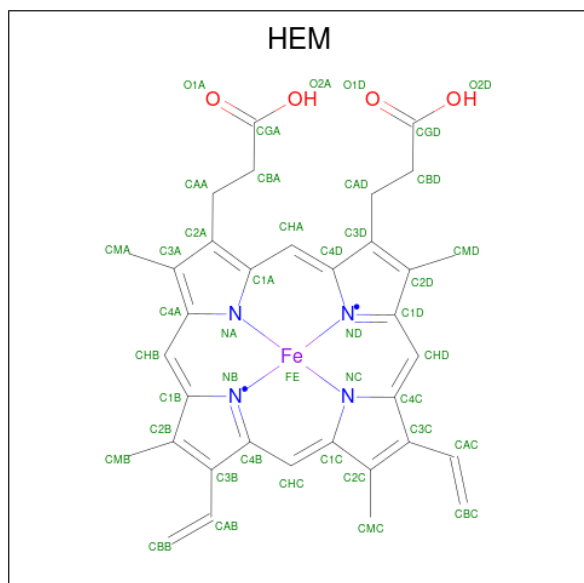
- Molecule 1 is a protein called Nitric oxide synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	406	Total	C	N	O	S	0	0	0
			3228	2052	568	591	17			
1	B	403	Total	C	N	O	S	0	0	0
			3210	2040	565	588	17			

There are 6 discrepancies between the modelled and reference sequences:

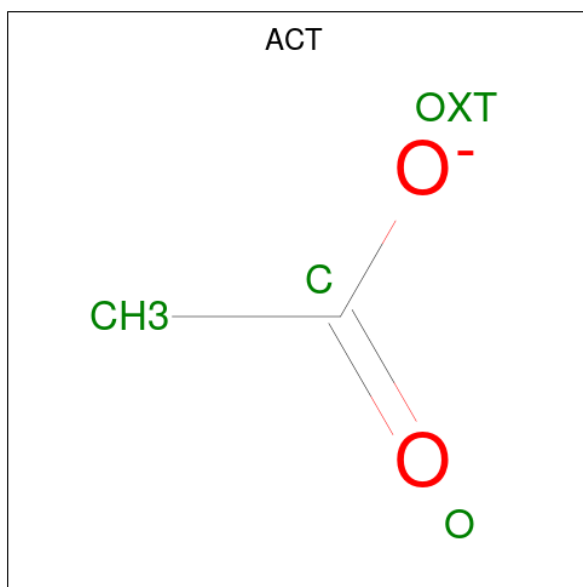
Chain	Residue	Modelled	Actual	Comment	Reference
A	100	ARG	CYS	SEE REMARK 999	UNP P29476
A	106	MET	VAL	engineered mutation	UNP P29476
A	368	ASP	ASN	engineered mutation	UNP P29476
B	100	ARG	CYS	SEE REMARK 999	UNP P29476
B	106	MET	VAL	engineered mutation	UNP P29476
B	368	ASP	ASN	engineered mutation	UNP P29476

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



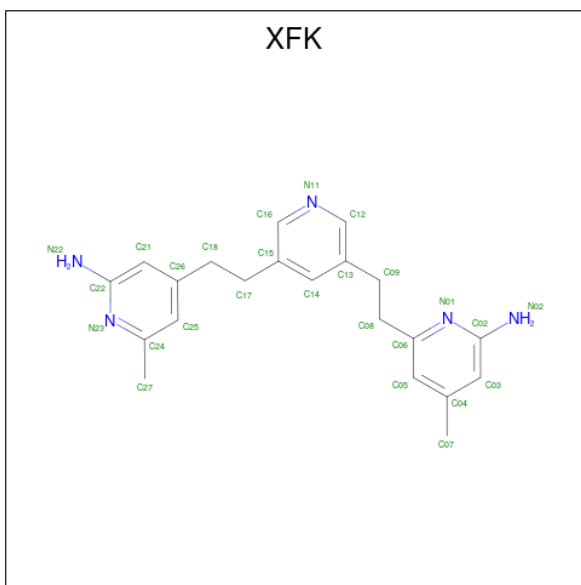
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 3 is ACETATE ION (CCD ID: ACT) (formula: $C_2H_3O_2^-$).



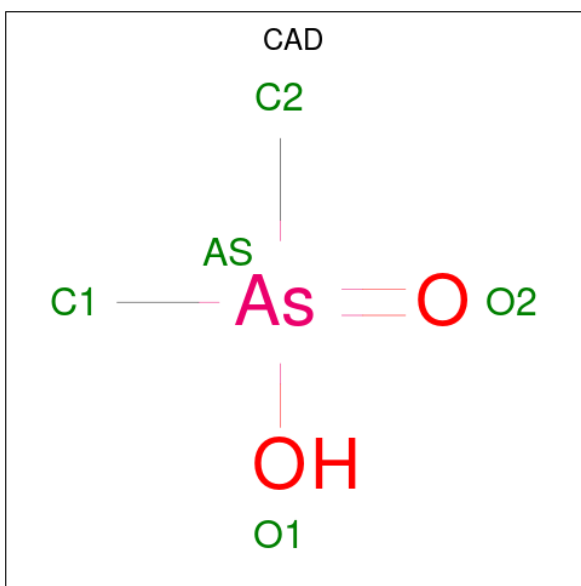
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 4 is 6-(2-{5-[2-(2-amino-6-methylpyridin-4-yl)ethyl]pyridin-3-yl}ethyl)-4-methylpyridin-2-amine (CCD ID: XFK) (formula: $C_{21}H_{25}N_5$).



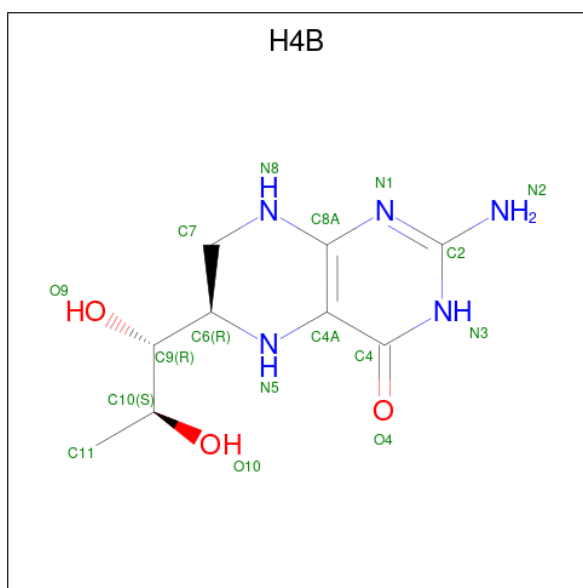
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	N	0	0
			26	21	5		
4	B	1	Total	C	N	0	0
			26	21	5		

- Molecule 5 is CACODYLIC ACID (CCD ID: CAD) (formula: $C_2H_7AsO_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	As	C	0	0
			3	1	2		
5	B	1	Total	As	C	0	0
			3	1	2		

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



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

- Molecule 7 is ZINC ION (CCD ID: ZN) (formula: Zn).

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

- Molecule 8 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	B	1	Total	C	O	0	0
			6	3	3		

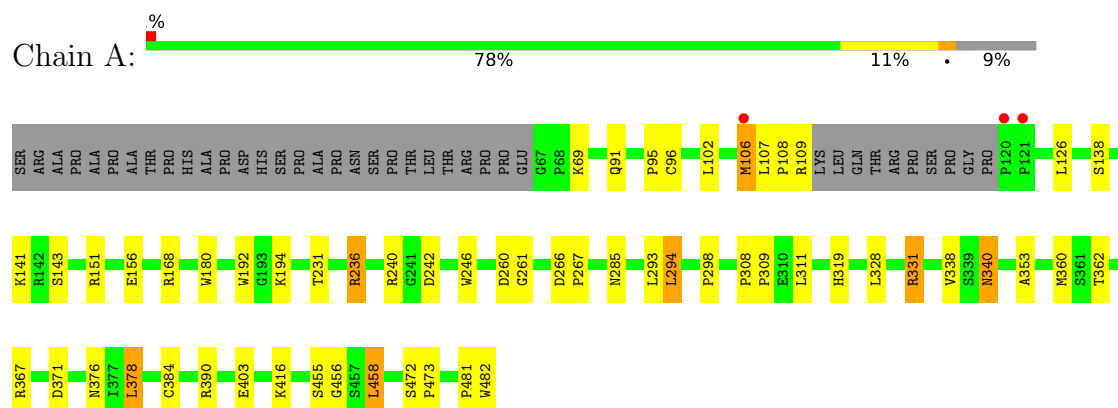
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	78	Total	O	0	0
			78	78		
9	B	76	Total	O	0	0
			76	76		

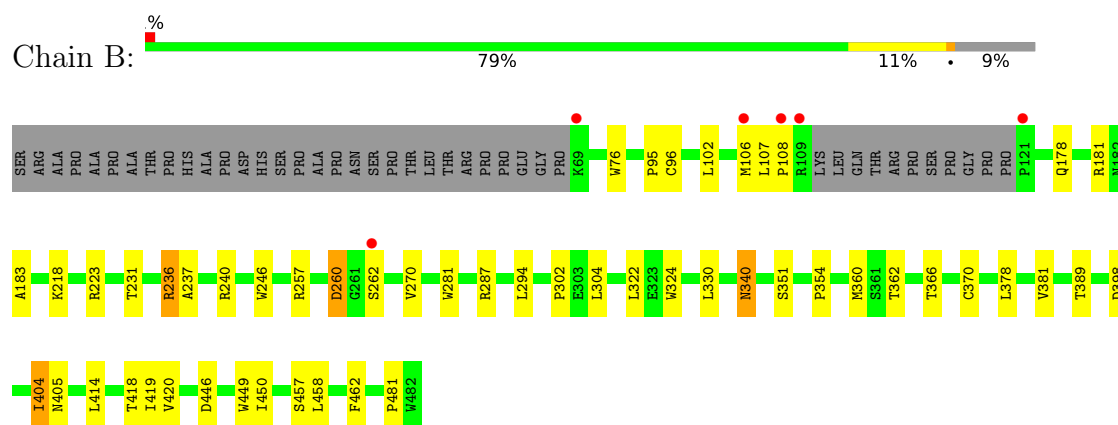
3 Residue-property plots [i](#)

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



• Molecule 1: 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	58.07Å 106.61Å 156.89Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.76 – 2.49 42.76 – 2.49	Depositor EDS
% Data completeness (in resolution range)	98.8 (42.76-2.49) 98.8 (42.76-2.49)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.57 (at 2.48Å)	Xtriage
Refinement program	REFMAC 5.5.0089, CNS	Depositor
R, R_{free}	0.184 , 0.243 0.191 , 0.244	Depositor DCC
R_{free} test set	1702 reflections (4.89%)	wwPDB-VP
Wilson B-factor (Å ²)	42.2	Xtriage
Anisotropy	0.111	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 42.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6785	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.68% 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: GOL, XFK, HEM, CAD, H4B, ZN, ACT

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	0.75	0/3319	0.89	0/4520
1	B	0.74	0/3299	0.95	8/4491 (0.2%)
All	All	0.75	0/6618	0.92	8/9011 (0.1%)

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	107	LEU	CA-C-N	-7.55	112.14	120.14
1	B	107	LEU	C-N-CA	-7.55	112.14	120.14
1	B	183	ALA	CA-C-N	-5.52	113.98	119.56
1	B	183	ALA	C-N-CA	-5.52	113.98	119.56
1	B	237	ALA	CA-C-N	5.52	125.87	119.47
1	B	237	ALA	C-N-CA	5.52	125.87	119.47
1	B	404	ILE	N-CA-C	-5.19	105.66	110.53
1	B	240	ARG	N-CA-C	5.07	116.14	108.07

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	3228	0	3132	30	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	3210	0	3115	28	0
2	A	43	0	30	2	0
2	B	43	0	30	4	0
3	A	4	0	3	0	0
3	B	4	0	3	0	0
4	A	26	0	25	3	0
4	B	26	0	25	0	0
5	A	3	0	0	2	0
5	B	3	0	0	0	0
6	A	17	0	15	0	0
6	B	17	0	15	0	0
7	A	1	0	0	0	0
8	B	6	0	8	0	0
9	A	78	0	0	4	0
9	B	76	0	0	1	0
All	All	6785	0	6401	63	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 (63) 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:384:CYS:SG	5:A:950:CAD:AS	2.47	1.33
1:B:106:MET:HE1	1:B:449:TRP:CZ2	1.82	1.14
1:B:106:MET:HE1	1:B:449:TRP:HZ2	1.02	1.09
1:B:106:MET:CE	1:B:449:TRP:HZ2	1.81	0.94
1:A:285:ASN:HB3	9:A:1007:HOH:O	1.67	0.93
2:B:500:HEM:HBB2	2:B:500:HEM:HHC	1.56	0.87
1:A:360:MET:HE3	1:A:362:THR:OG1	1.81	0.81
1:A:106:MET:HG3	1:B:76:TRP:NE1	2.03	0.73
1:B:366:THR:O	1:B:370:CYS:HB2	1.95	0.67
1:B:281:TRP:HB2	1:B:304:LEU:HD21	1.80	0.63
1:B:95:PRO:HB3	1:B:108:PRO:HG2	1.81	0.63
1:A:472:SER:HA	1:A:473:PRO:C	2.29	0.58
2:B:500:HEM:HHC	2:B:500:HEM:CBB	2.33	0.57
1:B:178:GLN:HE22	1:B:181:ARG:HH11	1.53	0.56
1:A:331:ARG:HD2	9:A:1020:HOH:O	2.06	0.56
1:A:367:ARG:NH2	1:A:371:ASP:OD2	2.38	0.55
1:B:322:LEU:HD13	1:B:324:TRP:CZ2	2.42	0.55
1:B:231:THR:O	1:B:354:PRO:HD2	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:370:CYS:SG	1:B:378:LEU:HD13	2.50	0.52
1:A:481:PRO:HD2	1:A:482:TRP:CE3	2.44	0.52
2:B:500:HEM:HBB2	2:B:500:HEM:CHC	2.31	0.51
1:A:236:ARG:HG3	1:A:242:ASP:OD1	2.11	0.50
1:B:360:MET:HE3	1:B:362:THR:OG1	2.12	0.50
1:A:240:ARG:HD3	1:A:298:PRO:HB3	1.93	0.50
1:B:446:ASP:O	1:B:450:ILE:HG12	2.12	0.49
2:A:500:HEM:HBA1	4:A:800:XFK:H08	1.94	0.49
1:B:95:PRO:CB	1:B:108:PRO:HG2	2.42	0.49
1:A:340:ASN:H	1:A:340:ASN:HD22	1.61	0.49
1:A:96:CYS:HB3	1:B:96:CYS:HB3	1.96	0.48
1:B:294:LEU:HD22	1:B:302:PRO:HB2	1.96	0.48
1:B:457:SER:HA	1:B:462:PHE:CG	2.49	0.48
1:B:236:ARG:HD3	1:B:351:SER:HB3	1.96	0.47
1:B:260:ASP:C	1:B:262:SER:H	2.22	0.47
1:B:362:THR:HA	1:B:405:ASN:HD21	1.80	0.47
1:A:231:THR:O	1:A:353:ALA:HA	2.15	0.47
1:A:378:LEU:HB2	9:A:1049:HOH:O	2.15	0.46
1:A:180:TRP:CE3	1:A:192:TRP:HA	2.51	0.45
1:A:246:TRP:HB2	1:A:294:LEU:HB3	1.98	0.45
1:B:381:VAL:HG21	1:B:404:ILE:HD11	1.99	0.45
1:B:106:MET:CA	1:B:106:MET:HE3	2.44	0.44
1:B:340:ASN:HD22	1:B:340:ASN:C	2.25	0.44
4:A:800:XFK:H12	4:A:800:XFK:H08A	1.78	0.44
1:A:106:MET:HB3	1:A:106:MET:HE3	1.17	0.43
1:B:246:TRP:CD1	1:B:481:PRO:HG2	2.53	0.43
1:B:106:MET:HE3	1:B:106:MET:HB3	1.69	0.43
1:A:376:ASN:ND2	9:A:1049:HOH:O	2.44	0.43
1:A:151:ARG:HD3	1:A:168:ARG:NH2	2.34	0.43
1:B:362:THR:HA	1:B:405:ASN:ND2	2.33	0.43
1:B:418:THR:HA	9:B:1151:HOH:O	2.19	0.43
2:B:500:HEM:HBC2	2:B:500:HEM:CMC	2.49	0.43
1:A:126:LEU:HD12	1:A:126:LEU:HA	1.80	0.42
1:A:156:GLU:H	1:A:156:GLU:HG2	1.76	0.41
1:B:360:MET:HA	1:B:420:VAL:O	2.20	0.41
1:A:126:LEU:HD11	1:A:156:GLU:HA	2.02	0.41
1:A:319:HIS:NE2	1:A:403:GLU:OE1	2.50	0.41
2:A:500:HEM:HBB2	2:A:500:HEM:HHC	2.02	0.41
1:A:266:ASP:HA	1:A:267:PRO:HD3	1.90	0.41
1:A:384:CYS:CB	5:A:950:CAD:AS	3.27	0.41
1:A:455:SER:O	1:A:458:LEU:HB2	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:95:PRO:HB3	1:A:108:PRO:CB	2.51	0.41
1:A:141:LYS:HA	1:A:141:LYS:HD3	1.94	0.40
1:A:338:VAL:CG2	4:A:800:XFK:H09	2.52	0.40
1:A:308:PRO:HA	1:A:309:PRO:HD3	1.98	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	402/444 (90%)	386 (96%)	14 (4%)	2 (0%)	24	43
1	B	399/444 (90%)	384 (96%)	15 (4%)	0	100	100
All	All	801/888 (90%)	770 (96%)	29 (4%)	2 (0%)	43	63

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	456	GLY
1	A	261	GLY

5.3.2 Protein sidechains [i](#)

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

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

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	345/377 (92%)	324 (94%)	21 (6%)	17	35
1	B	343/377 (91%)	328 (96%)	15 (4%)	25	50
All	All	688/754 (91%)	652 (95%)	36 (5%)	21	42

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

Mol	Chain	Res	Type
1	A	69	LYS
1	A	91	GLN
1	A	102	LEU
1	A	106	MET
1	A	107	LEU
1	A	109	ARG
1	A	138	SER
1	A	143	SER
1	A	194	LYS
1	A	236	ARG
1	A	260	ASP
1	A	293	LEU
1	A	294	LEU
1	A	311	LEU
1	A	328	LEU
1	A	331	ARG
1	A	340	ASN
1	A	378	LEU
1	A	390	ARG
1	A	416	LYS
1	A	458	LEU
1	B	102	LEU
1	B	218	LYS
1	B	223	ARG
1	B	236	ARG
1	B	257	ARG
1	B	260	ASP
1	B	270	VAL
1	B	287	ARG
1	B	330	LEU
1	B	340	ASN
1	B	389	THR

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Mol	Chain	Res	Type
1	B	398	ASP
1	B	414	LEU
1	B	419	ILE
1	B	458	LEU

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	166	HIS
1	A	207	GLN
1	A	340	ASN
1	A	376	ASN
1	A	413	GLN
1	A	468	ASN
1	B	124	GLN
1	B	128	GLN
1	B	178	GLN
1	B	222	ASN
1	B	225	ASN
1	B	279	HIS
1	B	340	ASN
1	B	376	ASN
1	B	405	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 12 ligands modelled in this entry, 1 is monoatomic - leaving 11 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
4	XFK	A	800	-	28,28,28	0.63	0	37,38,38	1.79	10 (27%)
3	ACT	A	860	-	3,3,3	0.86	0	3,3,3	0.72	0
5	CAD	B	950	-	0,2,4	-	-	0,1,6	-	-
6	H4B	B	600	-	17,18,18	0.88	1 (5%)	14,26,26	1.87	5 (35%)
6	H4B	A	600	-	17,18,18	0.83	0	14,26,26	1.84	5 (35%)
5	CAD	A	950	-	0,2,4	-	-	0,1,6	-	-
2	HEM	A	500	1	50,50,50	1.81	9 (18%)	67,82,82	1.25	6 (8%)
2	HEM	B	500	1	50,50,50	1.87	9 (18%)	67,82,82	1.48	9 (13%)
4	XFK	B	800	-	28,28,28	0.39	0	37,38,38	1.86	10 (27%)
8	GOL	B	880	-	5,5,5	0.41	0	5,5,5	0.66	0
3	ACT	B	860	-	3,3,3	0.76	0	3,3,3	0.83	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
4	XFK	A	800	-	-	2/10/10/10	0/3/3/3
6	H4B	B	600	-	-	2/8/17/17	0/2/2/2
6	H4B	A	600	-	-	1/8/17/17	0/2/2/2
2	HEM	A	500	1	-	1/14/54/54	-
2	HEM	B	500	1	-	0/14/54/54	-
4	XFK	B	800	-	-	3/10/10/10	0/3/3/3
8	GOL	B	880	-	-	2/4/4/4	-

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	500	HEM	C3D-C2D	7.66	1.53	1.36
2	A	500	HEM	C3D-C2D	6.90	1.51	1.36
2	A	500	HEM	FE-ND	5.68	2.12	1.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	500	HEM	FE-NB	4.12	2.07	1.94
2	B	500	HEM	FE-NC	3.92	2.08	1.95
2	B	500	HEM	FE-NA	3.79	2.07	1.95
2	A	500	HEM	FE-NA	3.28	2.06	1.95
2	B	500	HEM	CAB-C3B	3.00	1.55	1.47
2	B	500	HEM	CAC-C3C	2.98	1.55	1.47
2	A	500	HEM	CAB-C3B	2.81	1.54	1.47
2	A	500	HEM	CAC-C3C	2.65	1.54	1.47
2	A	500	HEM	C3B-C2B	-2.62	1.31	1.37
2	A	500	HEM	FE-NB	2.61	2.02	1.94
2	A	500	HEM	FE-NC	2.57	2.03	1.95
2	B	500	HEM	FE-ND	2.47	2.02	1.94
2	B	500	HEM	CMD-C2D	2.27	1.55	1.50
6	B	600	H4B	C4-N3	-2.17	1.34	1.38
2	A	500	HEM	CMC-C2C	2.12	1.55	1.50
2	B	500	HEM	C2A-C3A	-2.06	1.33	1.38

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	500	HEM	CBA-CAA-C2A	-5.67	96.85	112.53
4	A	800	XFK	C02-N01-C06	5.12	121.90	118.07
2	A	500	HEM	C4D-ND-C1D	4.85	110.95	105.21
4	B	800	XFK	C02-N01-C06	4.49	121.43	118.07
6	A	600	H4B	C2-N1-C8A	4.36	121.05	113.36
2	B	500	HEM	C4D-ND-C1D	4.31	110.31	105.21
6	B	600	H4B	C2-N1-C8A	3.84	120.14	113.36
4	B	800	XFK	C14-C15-C16	3.66	120.29	116.90
4	B	800	XFK	C26-C25-C24	-3.62	117.64	120.32
4	A	800	XFK	C14-C15-C16	3.56	120.19	116.90
4	B	800	XFK	C14-C13-C12	3.42	120.07	116.90
2	B	500	HEM	C1D-C2D-C3D	-3.23	103.58	106.98
4	B	800	XFK	C15-C14-C13	-3.09	117.36	121.20
2	A	500	HEM	CAD-C3D-C4D	2.95	129.84	124.70
2	B	500	HEM	CBD-CAD-C3D	-2.90	104.51	112.53
6	B	600	H4B	C4A-C4-N3	2.85	119.95	112.13
4	A	800	XFK	N22-C22-N23	2.74	120.99	116.59
4	A	800	XFK	C08-C06-C05	2.73	125.61	121.65
2	A	500	HEM	CMD-C2D-C1D	2.61	129.10	125.03
6	A	600	H4B	C4-C4A-N5	2.58	123.30	116.27
2	A	500	HEM	CBA-CAA-C2A	-2.54	105.52	112.53
6	B	600	H4B	O4-C4-C4A	-2.49	121.26	127.26

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	500	HEM	CHD-C4C-NC	2.49	127.16	124.45
4	B	800	XFK	C18-C17-C15	-2.45	104.88	113.23
2	B	500	HEM	C2A-C1A-NA	-2.45	107.44	110.15
6	B	600	H4B	C2-N3-C4	-2.41	120.73	125.11
4	B	800	XFK	N22-C22-N23	2.38	120.42	116.59
4	A	800	XFK	C14-C13-C12	2.38	119.11	116.90
4	A	800	XFK	C22-N23-C24	2.28	122.43	117.73
2	A	500	HEM	CBD-CAD-C3D	-2.27	106.26	112.53
4	A	800	XFK	C18-C26-C21	-2.23	116.80	120.54
6	B	600	H4B	C4-C4A-N5	2.20	122.28	116.27
6	A	600	H4B	N3-C2-N1	-2.16	119.37	123.32
4	A	800	XFK	C26-C25-C24	-2.13	118.75	120.32
6	A	600	H4B	C4A-C4-N3	2.13	117.97	112.13
4	A	800	XFK	C15-C14-C13	-2.12	118.56	121.20
4	A	800	XFK	C16-N11-C12	2.12	120.39	117.51
2	B	500	HEM	CMD-C2D-C1D	2.11	128.34	125.03
2	B	500	HEM	C3B-C2B-C1B	2.11	108.00	106.41
6	A	600	H4B	N2-C2-N3	2.06	121.11	116.76
4	B	800	XFK	C26-C21-C22	2.03	119.37	117.86
2	B	500	HEM	CHA-C4D-ND	2.03	126.88	124.37
4	B	800	XFK	C05-C06-N01	-2.02	120.42	122.73
2	B	500	HEM	CHC-C4B-NB	2.02	126.59	124.42
4	B	800	XFK	C09-C13-C12	-2.01	118.16	121.73

There are no chirality outliers.

All (11) torsion outliers are listed below:

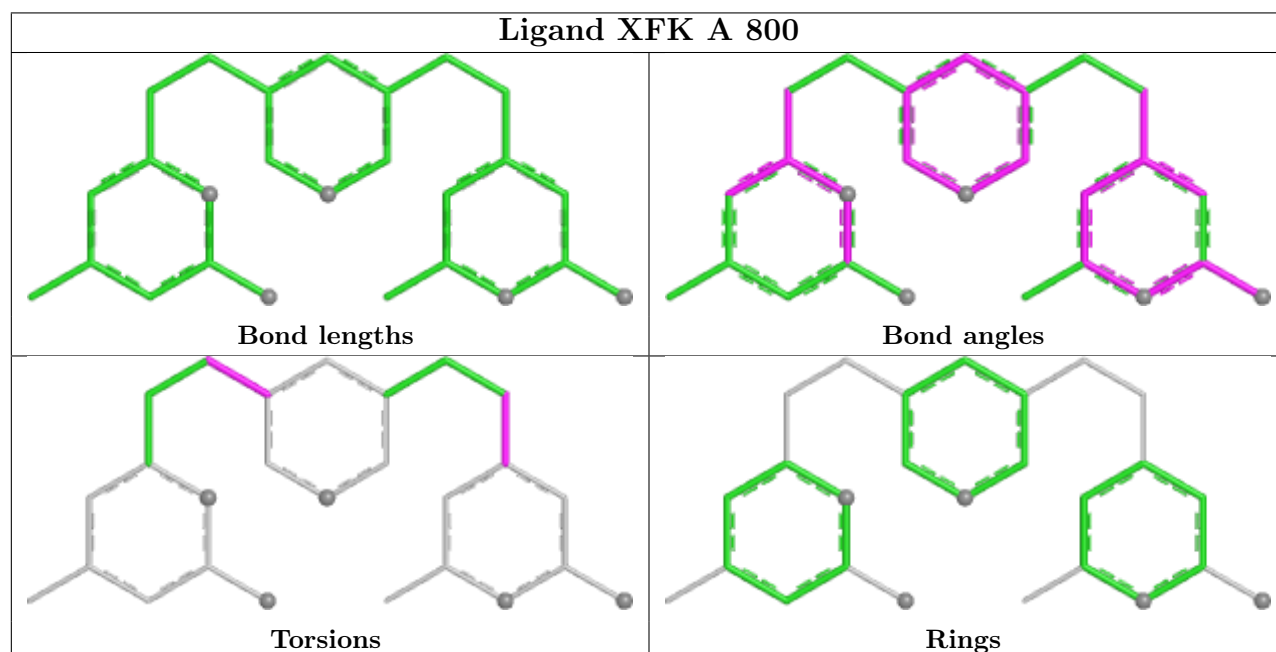
Mol	Chain	Res	Type	Atoms
8	B	880	GOL	O1-C1-C2-C3
8	B	880	GOL	O1-C1-C2-O2
6	B	600	H4B	O10-C10-C9-O9
4	A	800	XFK	C08-C09-C13-C14
2	A	500	HEM	C2A-CAA-CBA-CGA
4	B	800	XFK	C14-C15-C17-C18
4	B	800	XFK	C05-C06-C08-C09
4	B	800	XFK	N01-C06-C08-C09
6	A	600	H4B	N5-C6-C9-O9
6	B	600	H4B	N5-C6-C9-O9
4	A	800	XFK	C17-C18-C26-C21

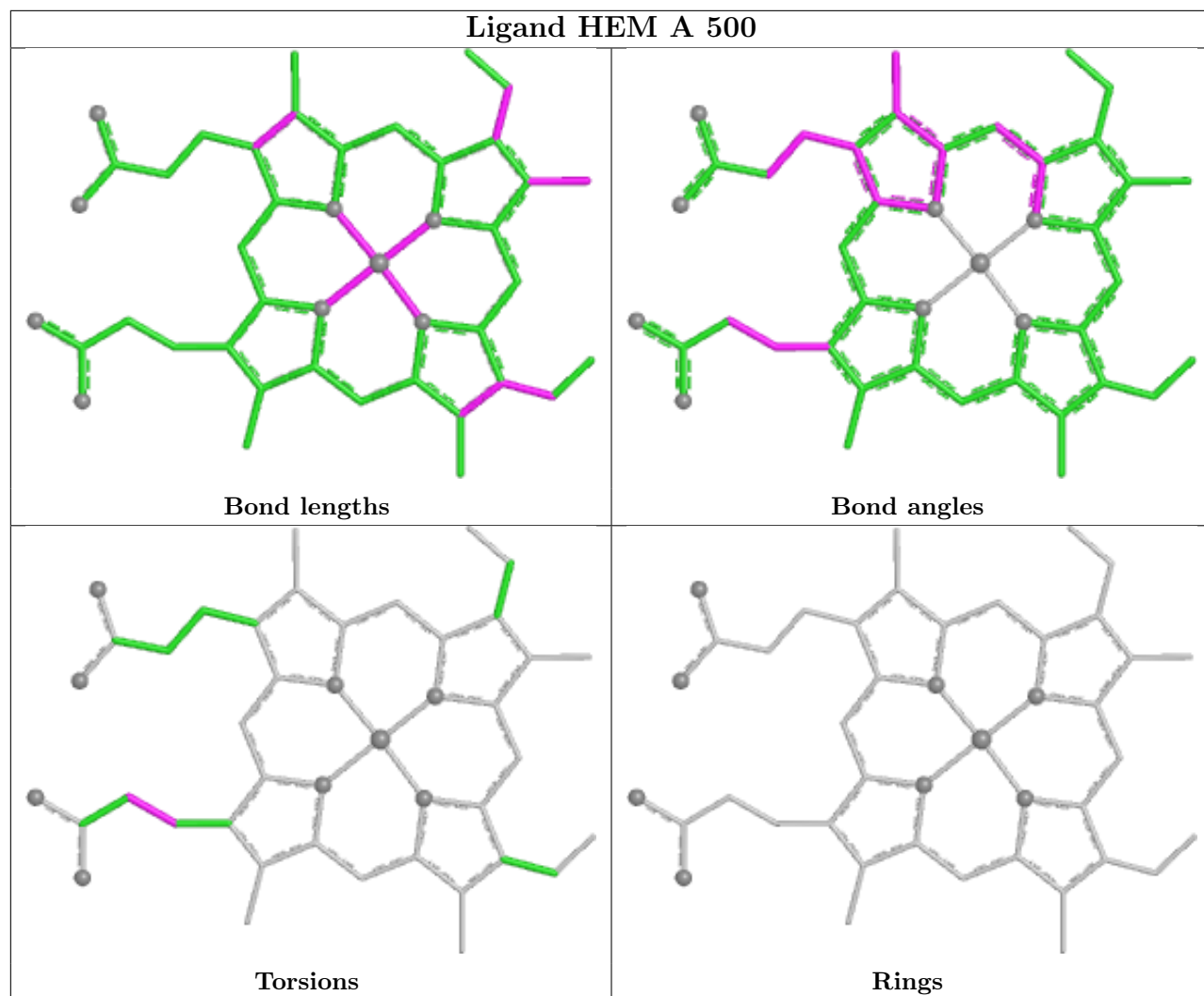
There are no ring outliers.

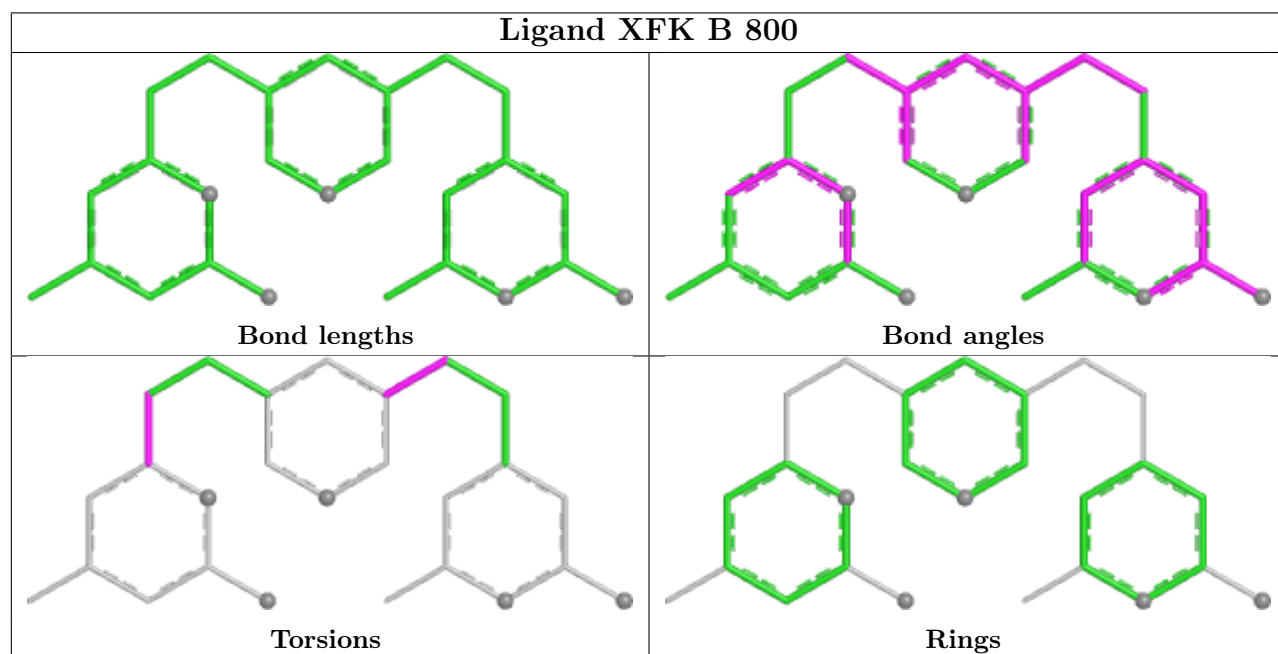
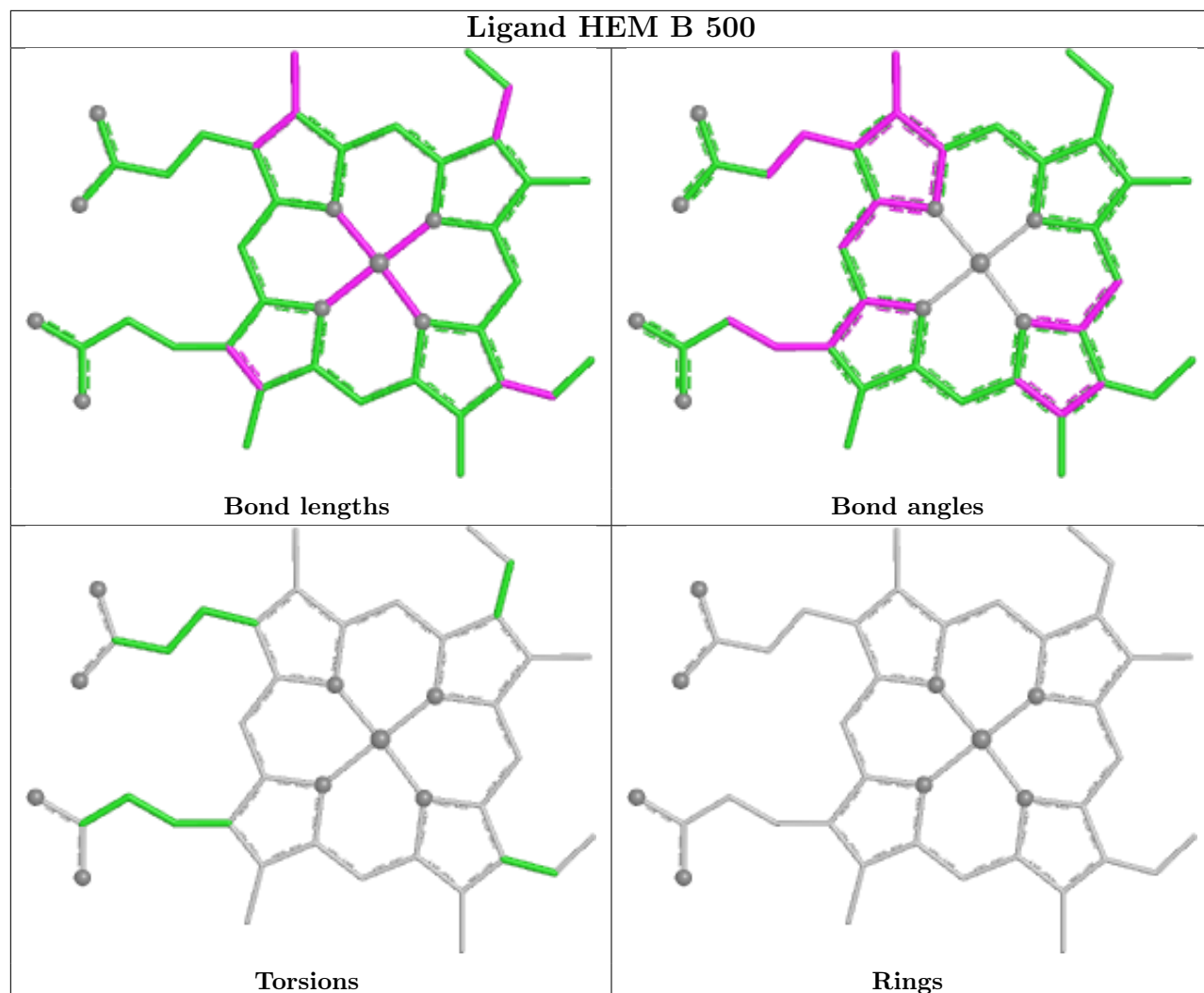
4 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	800	XFK	3	0
5	A	950	CAD	2	0
2	A	500	HEM	2	0
2	B	500	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 [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	406/444 (91%)	-0.28	3 (0%) 84 81	26, 39, 65, 86	0
1	B	403/444 (90%)	-0.19	6 (1%) 72 68	26, 42, 67, 96	0
All	All	809/888 (91%)	-0.23	9 (1%) 78 75	26, 40, 66, 96	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	120	PRO	4.8
1	B	121	PRO	3.6
1	A	121	PRO	2.9
1	B	108	PRO	2.9
1	B	106	MET	2.7
1	A	106	MET	2.6
1	B	69	LYS	2.4
1	B	109	ARG	2.3
1	B	262	SER	2.1

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.

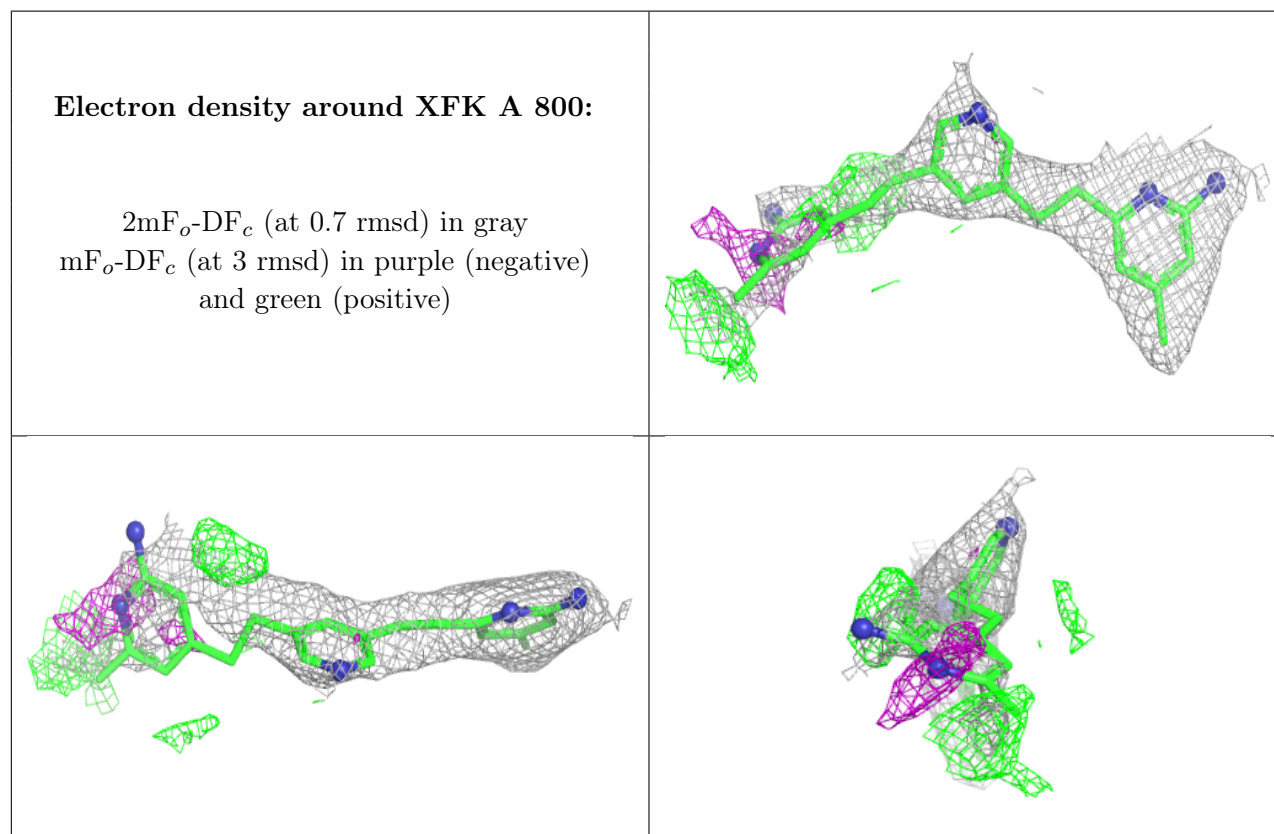
6.4 Ligands [i](#)

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.

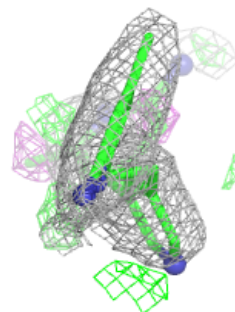
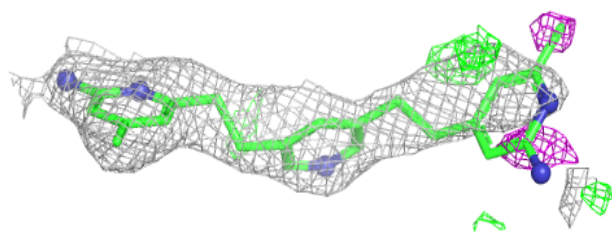
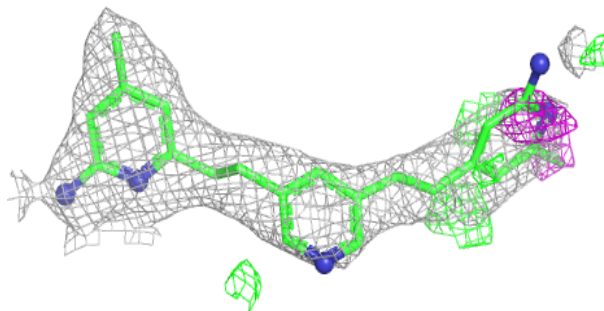
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	XFK	A	800	26/26	0.80	0.19	33,65,88,89	0
3	ACT	A	860	4/4	0.89	0.11	35,38,40,40	0
4	XFK	B	800	26/26	0.89	0.17	44,67,89,90	0
8	GOL	B	880	6/6	0.91	0.11	52,56,56,57	0
3	ACT	B	860	4/4	0.92	0.14	46,48,48,50	0
6	H4B	B	600	17/17	0.97	0.06	38,40,42,43	0
6	H4B	A	600	17/17	0.97	0.06	35,38,41,43	0
2	HEM	B	500	43/43	0.98	0.06	21,29,41,46	0
2	HEM	A	500	43/43	0.98	0.06	23,28,42,47	0
5	CAD	B	950	3/5	0.99	0.08	55,55,56,56	0
5	CAD	A	950	3/5	0.99	0.07	51,51,52,53	0
7	ZN	A	900	1/1	1.00	0.01	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.



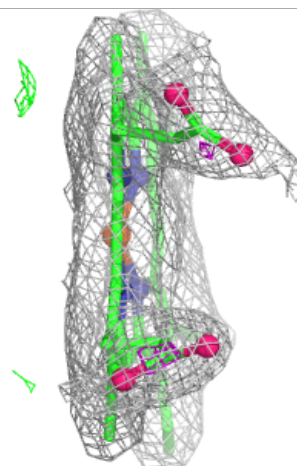
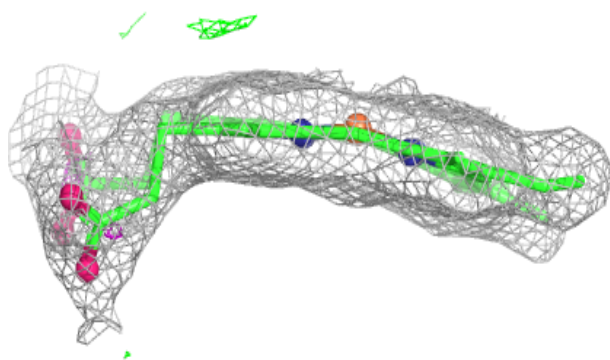
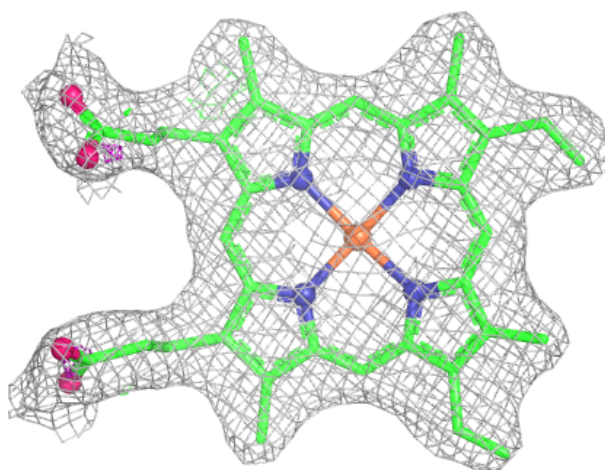
Electron density around XFK B 800:

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



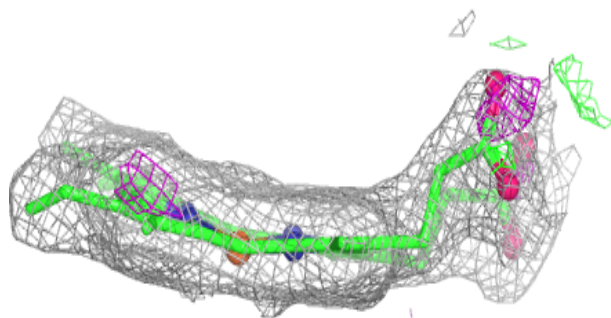
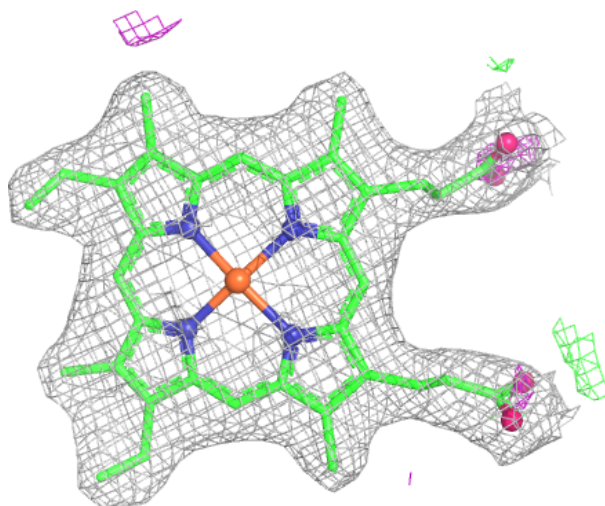
Electron density around HEM B 500:

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



Electron density around HEM A 500:

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



6.5 Other polymers ⓘ

There are no such residues in this entry.