



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 07:38 PM UTC

PDB ID : 1QW5 / pdb_00001qw5
Title : Murine inducible nitric oxide synthase oxygenase domain in complex with W1400 inhibitor.
Authors : Fedorov, R.; Hartmann, E.; Ghosh, D.K.; Schlichting, I.
Deposited on : 2003-08-31
Resolution : 2.70 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

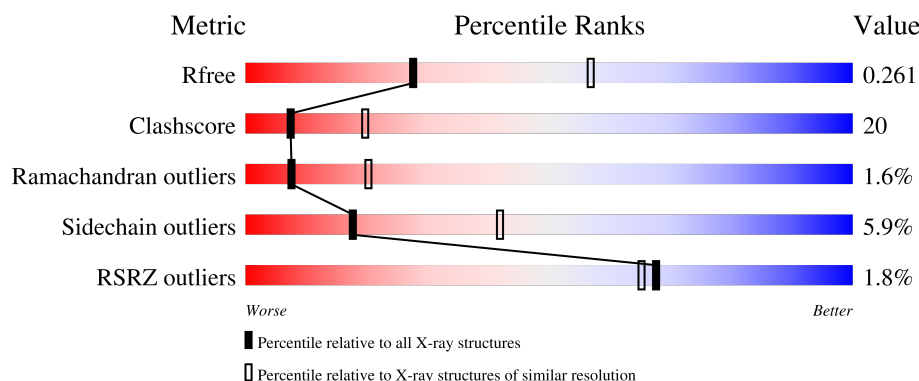
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	419	
1	B	419	

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 7241 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.

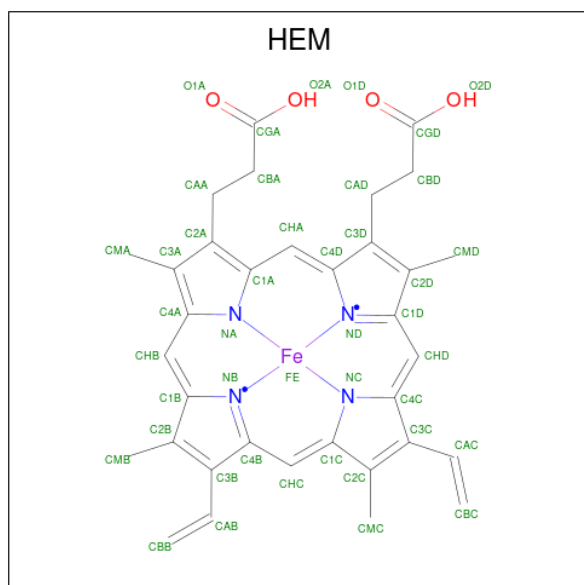
- Molecule 1 is a protein called Nitric oxide synthase, inducible.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	419	Total	C	N	O	S	0	0	0
			3415	2189	588	617	21			
1	B	419	Total	C	N	O	S	0	0	0
			3415	2189	588	617	21			

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

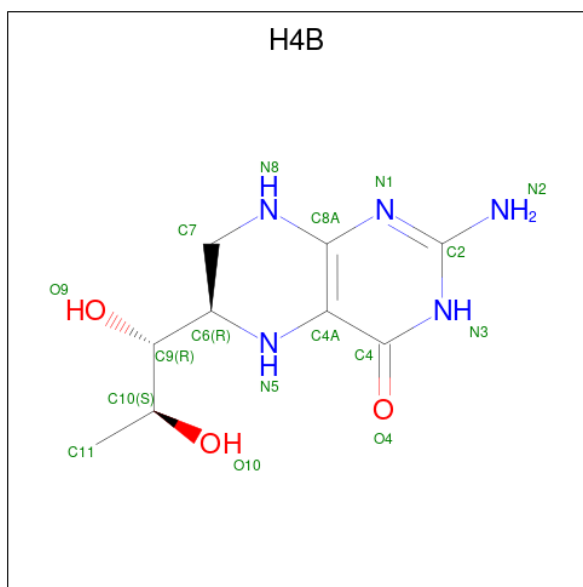
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		
2	B	1	Total	Zn	0	0
			1	1		

- Molecule 3 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



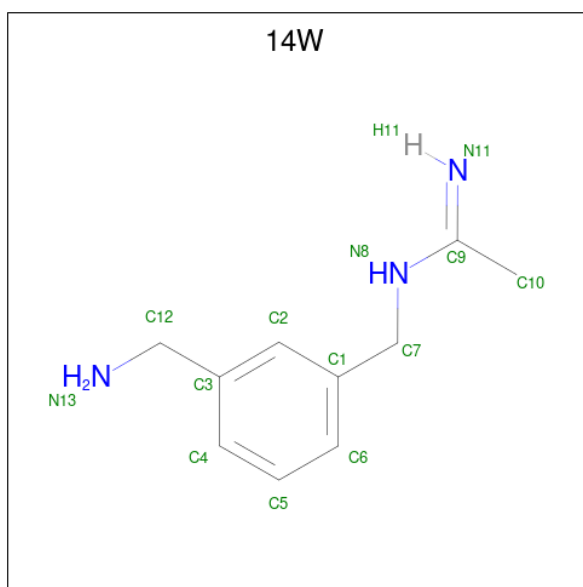
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
3	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 4 is 5,6,7,8-TETRAHYDROBIOPTERIN (CCD ID: H4B) (formula: $C_9H_{15}N_5O_3$).



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

- Molecule 5 is N-(3-(AMINOMETHYL)BENZYL)ACETAMIDINE (CCD ID: 14W) (formula: $C_{10}H_{15}N_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	N	0	0
			13	10	3		
5	B	1	Total	C	N	0	0
			13	10	3		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	135	Total	O	0	0
			135	135		
6	B	128	Total	O	0	0
			128	128		

4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	214.56Å 214.56Å 117.39Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	8.00 – 2.70 8.00 – 2.70	Depositor EDS
% Data completeness (in resolution range)	(Not available) (8.00-2.70) 95.6 (8.00-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.11 (at 2.71Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.211 , 0.264 0.209 , 0.261	Depositor DCC
R_{free} test set	2107 reflections (4.81%)	wwPDB-VP
Wilson B-factor (Å ²)	51.6	Xtriage
Anisotropy	0.579	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.42 , 80.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7241	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.62% 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 ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: HEM, 14W, H4B, 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	0.47	0/3515	0.95	9/4778 (0.2%)
1	B	0.48	1/3515 (0.0%)	0.96	14/4778 (0.3%)
All	All	0.47	1/7030 (0.0%)	0.95	23/9556 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	428	MET	SD-CE	-5.92	1.64	1.79

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	149	LYS	N-CA-C	-7.72	96.65	109.46
1	A	386	LEU	N-CA-C	7.29	119.31	111.36
1	B	122	GLY	N-CA-C	6.99	120.69	112.23
1	B	122	GLY	CA-C-N	6.92	126.90	119.78
1	B	122	GLY	C-N-CA	6.92	126.90	119.78
1	A	348	ASN	N-CA-C	6.68	120.74	112.59
1	B	367	TYR	N-CA-C	6.31	118.86	110.53
1	B	125	ASP	N-CA-C	-6.14	106.39	114.31
1	B	348	ASN	N-CA-C	5.92	120.10	112.41
1	B	151	ALA	N-CA-C	5.81	117.79	110.24
1	A	188	TRP	N-CA-C	-5.75	104.93	111.14
1	B	253	LEU	N-CA-C	-5.72	99.20	108.52
1	B	473	GLU	N-CA-C	-5.65	102.53	110.50
1	B	386	LEU	N-CA-C	5.60	117.46	111.36
1	A	375	ARG	N-CA-C	5.58	117.23	111.03
1	B	325	MET	N-CA-C	5.54	117.72	109.25
1	B	368	MET	N-CA-C	-5.38	99.64	108.41
1	A	316	PRO	CA-C-N	5.35	125.67	119.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	316	PRO	C-N-CA	5.35	125.67	119.47
1	B	176	THR	N-CA-C	-5.16	103.59	110.55
1	A	176	THR	N-CA-C	-5.07	103.36	110.50
1	B	234	ILE	N-CA-C	5.04	116.17	109.37
1	A	485	TYR	N-CA-C	-5.01	102.59	110.36

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	3415	0	3313	147	0
1	B	3415	0	3313	128	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	43	0	30	3	0
3	B	43	0	30	2	0
4	A	17	0	15	0	0
4	B	17	0	15	2	0
5	A	13	0	14	1	0
5	B	13	0	14	0	0
6	A	135	0	0	3	0
6	B	128	0	0	6	0
All	All	7241	0	6744	280	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 20.

All (280) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:153:ILE:HD12	1:B:153:ILE:H	1.13	1.14
1:A:153:ILE:HD12	1:A:153:ILE:H	1.28	0.96
1:B:285:ILE:HD11	1:B:291:PRO:HB3	1.53	0.91

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:494:ILE:H	1:B:494:ILE:HD12	1.36	0.91
1:A:494:ILE:HD12	1:A:494:ILE:H	1.34	0.91
1:A:80:ARG:HE	1:A:89:ILE:HG21	1.38	0.89
1:A:77:GLN:HG2	1:A:78:TYR:H	1.37	0.88
1:A:81:ILE:HD11	1:A:475:LEU:HD13	1.61	0.83
1:B:83:ASN:C	1:B:83:ASN:HD22	1.87	0.82
1:A:134:LEU:O	1:A:138:ILE:HG12	1.80	0.81
1:B:153:ILE:H	1:B:153:ILE:CD1	1.90	0.80
1:A:103:THR:HG22	1:A:110:LEU:HB2	1.63	0.79
1:B:83:ASN:ND2	1:B:85:GLY:H	1.81	0.79
1:B:301:LEU:HD13	1:B:315:ILE:HD11	1.66	0.78
1:B:83:ASN:HD22	1:B:85:GLY:H	1.34	0.76
1:B:103:THR:HG21	1:B:116:PRO:HB2	1.66	0.76
1:B:387:GLU:OE1	1:B:397:THR:HG21	1.86	0.75
1:B:387:GLU:OE2	1:B:391:ARG:HD2	1.85	0.75
1:A:149:LYS:O	1:A:150:GLU:HG2	1.86	0.75
1:A:77:GLN:HG2	1:A:78:TYR:N	2.01	0.75
1:B:107:LYS:NZ	1:B:107:LYS:HB3	2.03	0.74
1:A:130:LEU:HD13	1:A:167:ILE:HG22	1.69	0.74
1:A:154:GLU:CD	1:A:154:GLU:H	1.96	0.73
1:A:221:ILE:HG21	1:A:301:LEU:HD21	1.68	0.73
1:B:494:ILE:HD12	1:B:494:ILE:N	2.03	0.73
1:A:129:PRO:HB2	1:A:132:GLU:HG3	1.72	0.71
1:B:110:LEU:HD23	1:B:110:LEU:N	2.05	0.70
1:A:104:CYS:HA	1:A:110:LEU:HD12	1.74	0.70
1:A:134:LEU:HB3	1:A:135:PRO:HD3	1.74	0.70
1:B:103:THR:O	1:B:110:LEU:HG	1.92	0.69
1:A:438:MET:HE3	1:A:469:VAL:HG12	1.72	0.69
1:B:165:LYS:O	1:B:169:THR:HG23	1.93	0.69
1:A:159:ARG:O	1:A:163:VAL:HG23	1.93	0.69
1:A:266:MET:HE2	1:A:272:ARG:CD	2.23	0.68
1:A:104:CYS:HA	1:A:110:LEU:CD1	2.24	0.67
1:B:438:MET:HE2	1:B:438:MET:HA	1.75	0.67
1:B:88:GLU:C	1:B:89:ILE:HD12	2.21	0.66
1:A:188:TRP:CE3	1:A:200:TRP:HA	2.31	0.65
1:A:176:THR:OG1	1:A:179:GLU:HG3	1.96	0.65
1:A:266:MET:HE2	1:A:272:ARG:HD3	1.77	0.65
1:B:266:MET:HE2	1:B:272:ARG:HD3	1.78	0.65
1:B:375:ARG:O	1:B:379:ASP:HB2	1.96	0.65
1:A:138:ILE:HG22	1:A:142:ASN:HD21	1.62	0.65
1:A:262:ALA:HB2	1:A:299:LEU:CD2	2.27	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:153:ILE:HD12	1:B:153:ILE:N	1.98	0.64
1:A:387:GLU:O	1:A:391:ARG:HG3	1.98	0.63
1:A:138:ILE:HG23	1:A:160:LEU:HD22	1.81	0.63
1:A:215:GLN:HG3	1:A:219:GLN:NE2	2.14	0.62
1:B:110:LEU:HD23	1:B:110:LEU:H	1.62	0.62
1:A:252:ARG:HD3	1:A:359:PRO:HB2	1.80	0.62
1:A:330:TYR:HB3	1:A:332:TRP:CE2	2.34	0.62
1:B:260:ARG:HG3	6:B:986:HOH:O	1.99	0.62
1:A:271:ILE:HD12	1:A:271:ILE:N	2.16	0.61
1:B:266:MET:HE2	1:B:272:ARG:CD	2.30	0.61
1:A:375:ARG:O	1:A:379:ASP:HB2	2.00	0.61
1:B:171:GLY:O	1:B:172:THR:HB	2.01	0.61
1:B:285:ILE:HD11	1:B:291:PRO:CB	2.28	0.61
1:A:258:LEU:HB2	1:A:345:ALA:HB3	1.83	0.60
3:B:900:HEM:HBC2	3:B:900:HEM:HMC1	1.83	0.60
1:B:149:LYS:HD2	1:B:149:LYS:O	2.01	0.60
1:B:81:ILE:HD11	1:B:475:LEU:CD1	2.32	0.59
1:A:132:GLU:O	1:A:135:PRO:HD2	2.02	0.59
1:A:116:PRO:HG2	1:A:119:LEU:HB2	1.85	0.59
1:B:264:TYR:CE2	1:B:293:TYR:HA	2.38	0.59
1:B:116:PRO:CG	1:B:119:LEU:HD22	2.33	0.59
1:A:114:MET:HE2	1:A:457:TRP:HZ2	1.68	0.59
1:B:107:LYS:HB3	1:B:107:LYS:HZ3	1.66	0.59
1:A:217:MET:HB3	1:A:303:LEU:HD13	1.83	0.58
3:A:900:HEM:HBA2	6:A:963:HOH:O	2.02	0.58
1:A:266:MET:CE	1:A:272:ARG:HH11	2.15	0.58
1:A:251:PHE:O	1:A:360:ALA:HB2	2.02	0.58
1:A:103:THR:HG23	1:A:116:PRO:HB3	1.84	0.58
1:A:494:ILE:HD12	1:A:494:ILE:N	2.13	0.58
1:A:103:THR:HG22	1:A:110:LEU:CB	2.34	0.57
1:A:266:MET:HE2	1:A:272:ARG:HH11	1.69	0.57
1:B:407:ARG:HD3	6:B:983:HOH:O	2.05	0.57
1:B:188:TRP:CE3	1:B:200:TRP:HA	2.39	0.57
1:B:166:GLU:OE2	1:B:173:TYR:HA	2.05	0.57
1:B:143:GLN:HE21	1:B:186:MET:HE2	1.68	0.56
1:A:417:LEU:O	1:A:421:GLN:HG3	2.05	0.56
1:B:78:TYR:CD1	1:B:78:TYR:C	2.83	0.56
1:A:154:GLU:HG2	1:A:155:GLU:H	1.71	0.56
1:A:271:ILE:HD12	1:A:271:ILE:H	1.71	0.56
1:A:223:ARG:HH11	1:A:223:ARG:HG2	1.70	0.56
1:A:254:TRP:HE3	1:A:283:LEU:HD21	1.70	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:163:VAL:HG13	1:A:173:TYR:CD2	2.41	0.56
1:A:230:ASN:HB3	1:A:233:ASN:O	2.06	0.56
1:B:304:GLN:HG3	1:B:308:GLN:O	2.06	0.56
1:B:103:THR:CG2	1:B:116:PRO:HB2	2.35	0.55
1:A:138:ILE:HG23	1:A:160:LEU:CD2	2.36	0.55
1:A:303:LEU:O	1:A:310:PRO:HA	2.07	0.55
1:B:254:TRP:HE3	1:B:283:LEU:HD21	1.72	0.55
1:A:124:ARG:HH22	1:A:128:THR:HG22	1.72	0.55
1:B:137:ALA:O	1:B:141:ILE:HG12	2.07	0.55
1:A:153:ILE:O	1:A:157:LEU:HG	2.07	0.55
1:B:153:ILE:O	1:B:156:HIS:HB3	2.06	0.55
1:B:176:THR:OG1	1:B:179:GLU:HG3	2.06	0.55
1:A:133:LEU:HD22	1:A:167:ILE:HD13	1.88	0.55
1:A:105:LYS:HG3	1:A:110:LEU:HD11	1.88	0.54
1:A:131:GLU:O	1:A:135:PRO:HG2	2.07	0.54
1:B:301:LEU:CD1	1:B:315:ILE:HD11	2.36	0.54
1:A:344:PRO:HB2	5:A:902:14W:H6	1.90	0.54
1:B:103:THR:HG23	1:B:118:SER:HB3	1.90	0.54
1:B:283:LEU:O	1:B:287:LEU:HG	2.08	0.54
1:A:459:VAL:HG22	1:A:469:VAL:HG23	1.90	0.53
1:B:116:PRO:HG2	1:B:119:LEU:HD22	1.90	0.53
1:B:221:ILE:HG21	1:B:301:LEU:HD21	1.89	0.53
1:B:241:PHE:HB3	1:B:242:PRO:CD	2.38	0.53
1:B:489:PRO:C	1:B:491:LYS:H	2.16	0.53
1:A:107:LYS:HB2	1:A:107:LYS:NZ	2.22	0.53
1:B:134:LEU:O	1:B:138:ILE:HG13	2.08	0.53
1:B:154:GLU:N	1:B:154:GLU:CD	2.66	0.53
1:A:132:GLU:C	1:A:135:PRO:HD2	2.33	0.52
1:A:246:ASP:OD2	1:A:248:LYS:HB2	2.09	0.52
1:B:438:MET:HE3	1:B:469:VAL:HG12	1.91	0.52
1:A:342:ALA:HB1	1:A:425:VAL:HG11	1.92	0.52
1:B:244:ARG:HA	6:B:1002:HOH:O	2.10	0.51
1:B:330:TYR:HD2	1:B:332:TRP:NE1	2.08	0.51
1:B:124:ARG:HB2	1:B:247:GLY:HA3	1.92	0.51
1:B:251:PHE:O	1:B:360:ALA:HB2	2.10	0.51
1:A:104:CYS:CB	1:A:109:CYS:HA	2.40	0.51
1:A:138:ILE:O	1:A:142:ASN:ND2	2.43	0.51
1:B:198:ILE:HB	1:B:440:HIS:HB2	1.91	0.51
1:A:124:ARG:NH2	1:A:128:THR:HG22	2.25	0.51
1:B:417:LEU:O	1:B:421:GLN:HG3	2.11	0.51
1:B:81:ILE:HD11	1:B:475:LEU:HD12	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:148:PHE:N	1:A:148:PHE:CD2	2.79	0.51
1:A:266:MET:HE2	1:A:272:ARG:HD2	1.91	0.51
1:A:153:ILE:HG22	1:A:157:LEU:HD11	1.92	0.50
1:B:154:GLU:CD	1:B:155:GLU:H	2.18	0.50
1:A:330:TYR:HE1	1:A:393:MET:HE3	1.75	0.50
1:A:330:TYR:CD2	1:A:330:TYR:N	2.76	0.50
1:B:132:GLU:O	1:B:135:PRO:HD2	2.11	0.50
1:B:266:MET:HB3	1:B:267:PRO:HD2	1.93	0.50
1:A:410:THR:O	1:A:414:VAL:HG23	2.12	0.50
1:B:268:ASP:O	1:B:270:THR:N	2.37	0.50
1:A:241:PHE:HB3	1:A:242:PRO:HD2	1.93	0.50
1:B:138:ILE:HG22	1:B:142:ASN:HD21	1.77	0.50
1:B:327:HIS:ND1	1:B:328:PRO:HD2	2.27	0.50
1:B:330:TYR:HD2	1:B:332:TRP:HE1	1.58	0.50
1:B:83:ASN:C	1:B:83:ASN:ND2	2.60	0.50
1:B:215:GLN:HG3	1:B:219:GLN:NE2	2.27	0.50
1:B:410:THR:O	1:B:414:VAL:HG23	2.11	0.49
1:A:104:CYS:HA	1:A:110:LEU:H	1.76	0.49
1:A:241:PHE:HB3	1:A:242:PRO:CD	2.42	0.49
1:A:262:ALA:HB2	1:A:299:LEU:HD21	1.91	0.49
1:B:387:GLU:CD	1:B:391:ARG:HD2	2.36	0.49
1:B:246:ASP:OD2	1:B:248:LYS:HB2	2.12	0.49
1:A:180:LEU:O	1:A:184:THR:HG23	2.12	0.49
1:A:130:LEU:CD1	1:A:167:ILE:HG22	2.42	0.49
1:A:78:TYR:C	1:A:78:TYR:CD1	2.91	0.49
1:B:350:LEU:C	1:B:350:LEU:HD23	2.38	0.48
1:B:453:ALA:HB1	1:B:458:LEU:CD1	2.43	0.48
1:A:264:TYR:CD2	1:A:293:TYR:HA	2.47	0.48
1:A:305:ALA:O	1:A:308:GLN:HB2	2.14	0.48
1:B:110:LEU:N	1:B:110:LEU:CD2	2.75	0.48
1:B:215:GLN:HG3	1:B:219:GLN:HE22	1.79	0.48
1:A:153:ILE:HD12	1:A:153:ILE:N	2.11	0.48
1:A:379:ASP:HB3	1:A:382:ARG:HG2	1.94	0.48
1:B:173:TYR:CE1	1:B:356:LEU:HD21	2.48	0.48
1:A:154:GLU:CD	1:A:154:GLU:N	2.69	0.48
1:A:304:GLN:HG3	1:A:308:GLN:O	2.13	0.48
1:A:301:LEU:HD13	1:A:315:ILE:HD11	1.96	0.48
1:A:494:ILE:H	1:A:494:ILE:CD1	2.11	0.48
1:B:195:ILE:N	6:B:1021:HOH:O	2.44	0.48
1:B:271:ILE:HD13	1:B:278:LEU:HD11	1.96	0.48
1:B:480:SER:HA	1:B:481:PRO:C	2.39	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:124:ARG:HG3	1:B:244:ARG:HD3	1.96	0.47
1:B:267:PRO:HG2	1:B:268:ASP:OD1	2.14	0.47
1:A:138:ILE:HG22	1:A:142:ASN:ND2	2.26	0.47
1:A:252:ARG:CD	1:A:359:PRO:HB2	2.44	0.47
1:A:104:CYS:HB2	1:A:109:CYS:HA	1.96	0.47
1:A:379:ASP:CB	1:A:382:ARG:HG2	2.44	0.47
1:B:239:THR:O	1:B:361:CYS:HA	2.15	0.47
1:A:81:ILE:HD12	1:A:475:LEU:HD22	1.97	0.47
1:A:438:MET:HA	1:A:438:MET:HE2	1.96	0.47
1:A:457:TRP:HZ3	6:A:963:HOH:O	1.97	0.47
1:B:98:ALA:O	1:B:99:THR:O	2.32	0.47
1:B:241:PHE:HB3	1:B:242:PRO:HD2	1.96	0.47
1:B:487:ILE:N	1:B:487:ILE:HD12	2.30	0.46
1:A:161:GLU:OE2	1:A:165:LYS:NZ	2.47	0.46
1:A:254:TRP:CE3	1:A:283:LEU:HD21	2.50	0.46
1:A:349:MET:HE3	1:A:485:TYR:CD1	2.51	0.46
3:A:900:HEM:HBD1	6:A:963:HOH:O	2.15	0.46
1:B:230:ASN:HB3	1:B:233:ASN:O	2.16	0.46
1:A:348:ASN:H	1:A:348:ASN:ND2	2.14	0.46
1:B:80:ARG:NH2	6:B:1042:HOH:O	2.44	0.46
1:B:83:ASN:HB2	1:B:472:GLN:OE1	2.15	0.46
1:B:89:ILE:HD12	1:B:89:ILE:N	2.31	0.46
3:A:900:HEM:HHC	3:A:900:HEM:HBB2	1.97	0.45
1:A:264:TYR:CE2	1:A:293:TYR:HA	2.52	0.45
1:A:103:THR:HG23	1:A:116:PRO:CB	2.47	0.45
1:A:249:HIS:HA	1:A:307:GLY:HA3	1.99	0.45
1:B:466:ILE:O	1:B:466:ILE:HG22	2.16	0.45
1:A:480:SER:HA	1:A:481:PRO:C	2.41	0.45
1:A:302:VAL:C	1:A:303:LEU:HD23	2.42	0.45
1:B:79:VAL:HG11	1:B:475:LEU:HD11	1.98	0.45
1:A:330:TYR:HB3	1:A:332:TRP:NE1	2.32	0.45
1:B:83:ASN:ND2	1:B:85:GLY:N	2.59	0.44
1:B:258:LEU:HD22	1:B:345:ALA:HB1	1.98	0.44
1:B:264:TYR:CD2	1:B:293:TYR:HA	2.52	0.44
3:B:900:HEM:HBC2	3:B:900:HEM:CMC	2.46	0.44
1:B:154:GLU:CD	1:B:154:GLU:H	2.24	0.44
1:B:489:PRO:C	1:B:491:LYS:N	2.75	0.44
1:A:104:CYS:HA	1:A:110:LEU:HD13	1.99	0.44
1:A:137:ALA:O	1:A:141:ILE:HG12	2.17	0.44
1:A:134:LEU:HD11	1:A:138:ILE:HD11	1.99	0.44
1:A:350:LEU:C	1:A:350:LEU:HD23	2.43	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:116:PRO:CG	1:A:119:LEU:HD22	2.48	0.44
1:B:181:ILE:O	1:B:185:LYS:HG3	2.17	0.44
1:A:130:LEU:HD13	1:A:167:ILE:CG2	2.44	0.44
1:A:131:GLU:O	1:A:135:PRO:CD	2.65	0.44
1:A:103:THR:CG2	1:A:116:PRO:HB3	2.48	0.43
1:A:346:VAL:HB	1:A:363:PHE:CZ	2.53	0.43
1:A:244:ARG:HG3	1:A:250:ASP:OD1	2.18	0.43
1:A:262:ALA:CB	1:A:299:LEU:HD21	2.48	0.43
1:A:373:GLY:HA3	1:A:413:ASN:ND2	2.34	0.43
1:B:227:TYR:CD1	1:B:227:TYR:C	2.95	0.43
1:B:258:LEU:HB2	1:B:345:ALA:HB3	1.99	0.43
1:B:327:HIS:ND1	1:B:328:PRO:CD	2.80	0.43
1:A:227:TYR:CD1	1:A:227:TYR:C	2.95	0.43
1:B:383:TYR:O	1:B:384:ASN:C	2.60	0.43
1:B:494:ILE:H	1:B:494:ILE:CD1	2.17	0.43
1:A:131:GLU:O	1:A:135:PRO:CG	2.67	0.43
1:B:327:HIS:CG	1:B:328:PRO:HD2	2.53	0.43
1:A:80:ARG:HG3	1:A:89:ILE:CG2	2.49	0.43
1:B:138:ILE:HG23	1:B:160:LEU:HD22	2.01	0.43
1:A:148:PHE:N	1:A:148:PHE:HD2	2.17	0.42
1:A:202:ASN:N	1:A:202:ASN:HD22	2.16	0.42
1:A:154:GLU:HG2	1:A:155:GLU:N	2.34	0.42
1:A:368:MET:O	1:A:369:GLY:C	2.62	0.42
1:B:159:ARG:O	1:B:163:VAL:HG23	2.19	0.42
1:A:123:PRO:HD2	1:A:487:ILE:HD12	2.01	0.42
1:A:466:ILE:HG22	1:A:466:ILE:O	2.18	0.42
1:A:260:ARG:HD2	1:A:277:THR:HG21	2.02	0.42
1:B:367:TYR:CD2	1:B:372:ILE:HD11	2.54	0.42
1:A:114:MET:CE	1:A:457:TRP:HZ2	2.32	0.42
1:A:207:ASP:OD1	1:A:207:ASP:C	2.63	0.42
1:A:266:MET:HB3	1:A:267:PRO:HD2	2.02	0.42
1:A:344:PRO:O	1:A:344:PRO:HG2	2.19	0.42
1:B:327:HIS:CE1	1:B:330:TYR:HD1	2.37	0.42
1:A:103:THR:HG21	1:A:110:LEU:O	2.20	0.41
1:B:256:SER:HA	1:B:348:ASN:ND2	2.35	0.41
1:B:459:VAL:HG22	1:B:469:VAL:HG23	2.01	0.41
1:B:121:ARG:HD3	1:B:121:ARG:HA	1.90	0.41
1:B:375:ARG:HH12	4:B:903:H4B:C4	2.33	0.41
1:A:153:ILE:O	1:A:156:HIS:HB3	2.21	0.41
1:B:166:GLU:O	1:B:170:THR:OG1	2.31	0.41
1:A:193:ARG:HD3	1:A:457:TRP:CD2	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:330:TYR:HD1	1:A:332:TRP:CZ2	2.38	0.41
1:A:413:ASN:HD22	1:A:413:ASN:HA	1.56	0.41
1:B:266:MET:HE2	1:B:272:ARG:HD2	2.00	0.41
1:A:248:LYS:HE3	1:A:248:LYS:HB3	1.81	0.41
1:B:422:LYS:NZ	1:B:422:LYS:HB3	2.36	0.41
1:B:77:GLN:O	1:B:96:HIS:HE1	2.02	0.41
1:B:83:ASN:O	1:B:87:GLY:N	2.53	0.41
1:A:214:ALA:O	1:A:217:MET:HB2	2.20	0.41
1:B:103:THR:HG23	1:B:118:SER:CB	2.50	0.41
1:B:132:GLU:C	1:B:135:PRO:HD2	2.46	0.41
1:B:195:ILE:HG12	1:B:195:ILE:O	2.20	0.41
1:B:211:CYS:SG	1:B:216:GLU:HB3	2.60	0.41
1:A:102:PHE:CG	1:A:103:THR:N	2.82	0.41
1:A:110:LEU:HD12	1:A:110:LEU:N	2.35	0.41
1:A:176:THR:HG23	1:A:179:GLU:CD	2.46	0.41
1:A:330:TYR:N	1:A:330:TYR:HD2	2.17	0.41
1:A:330:TYR:CD1	1:A:332:TRP:CZ2	3.09	0.41
1:A:418:HIS:HA	1:A:421:GLN:OE1	2.21	0.41
1:B:457:TRP:HA	4:B:903:H4B:N1	2.36	0.41
1:A:494:ILE:N	1:A:494:ILE:CD1	2.79	0.40
1:B:116:PRO:HG2	1:B:119:LEU:HB2	2.02	0.40
1:B:149:LYS:HD2	1:B:149:LYS:C	2.46	0.40
1:B:282:GLN:O	1:B:283:LEU:C	2.64	0.40
1:A:78:TYR:HB3	1:A:93:THR:HG22	2.04	0.40
1:B:243:GLN:HB3	1:B:358:PHE:CE2	2.56	0.40
1:B:369:GLY:HA3	6:B:1004:HOH:O	2.21	0.40
1:A:134:LEU:CD1	1:A:138:ILE:HD11	2.50	0.40
1:A:244:ARG:HG3	1:A:247:GLY:HA2	2.04	0.40
1:A:374:VAL:O	1:A:378:CYS:HB2	2.21	0.40
1:B:103:THR:HG21	1:B:116:PRO:CB	2.46	0.40
1:B:138:ILE:O	1:B:142:ASN:ND2	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	417/419 (100%)	374 (90%)	37 (9%)	6 (1%)	9	23
1	B	417/419 (100%)	376 (90%)	34 (8%)	7 (2%)	7	19
All	All	834/838 (100%)	750 (90%)	71 (8%)	13 (2%)	7	20

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	100	SER
1	A	105	LYS
1	B	99	THR
1	B	172	THR
1	B	269	GLY
1	A	150	GLU
1	A	369	GLY
1	A	397	THR
1	B	100	SER
1	A	268	ASP
1	B	464	GLY
1	B	307	GLY
1	B	369	GLY

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	367/367 (100%)	343 (94%)	24 (6%)	15	37
1	B	367/367 (100%)	348 (95%)	19 (5%)	21	47
All	All	734/734 (100%)	691 (94%)	43 (6%)	18	42

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	81	ILE
1	A	99	THR
1	A	107	LYS
1	A	119	LEU
1	A	124	ARG
1	A	128	THR
1	A	130	LEU
1	A	149	LYS
1	A	153	ILE
1	A	154	GLU
1	A	178	ASP
1	A	201	SER
1	A	256	SER
1	A	258	LEU
1	A	260	ARG
1	A	301	LEU
1	A	330	TYR
1	A	348	ASN
1	A	380	THR
1	A	413	ASN
1	A	435	GLU
1	A	475	LEU
1	A	488	GLU
1	A	494	ILE
1	B	83	ASN
1	B	103	THR
1	B	107	LYS
1	B	110	LEU
1	B	119	LEU
1	B	128	THR
1	B	153	ILE
1	B	154	GLU
1	B	178	ASP
1	B	219	GLN
1	B	258	LEU
1	B	268	ASP
1	B	301	LEU
1	B	324	THR
1	B	413	ASN
1	B	422	LYS
1	B	474	MET
1	B	475	LEU
1	B	494	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	96	HIS
1	A	142	ASN
1	A	143	GLN
1	A	174	GLN
1	A	202	ASN
1	A	215	GLN
1	A	219	GLN
1	A	231	ASN
1	A	308	GLN
1	A	334	GLN
1	A	348	ASN
1	A	398	HIS
1	A	413	ASN
1	A	486	GLN
1	B	83	ASN
1	B	96	HIS
1	B	115	ASN
1	B	142	ASN
1	B	143	GLN
1	B	174	GLN
1	B	202	ASN
1	B	215	GLN
1	B	219	GLN
1	B	220	HIS
1	B	231	ASN
1	B	249	HIS
1	B	476	ASN
1	B	486	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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 8 ligands modelled in this entry, 2 are monoatomic - leaving 6 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	H4B	A	901	-	17,18,18	1.47	3 (17%)	14,26,26	1.96	3 (21%)
5	14W	B	904	-	13,13,13	1.54	3 (23%)	13,16,16	1.08	1 (7%)
4	H4B	B	903	-	17,18,18	1.63	4 (23%)	14,26,26	1.87	3 (21%)
5	14W	A	902	-	13,13,13	1.53	5 (38%)	13,16,16	1.24	2 (15%)
3	HEM	A	900	1	50,50,50	1.31	6 (12%)	67,82,82	1.38	9 (13%)
3	HEM	B	900	1	50,50,50	1.39	4 (8%)	67,82,82	1.28	4 (5%)

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	H4B	A	901	-	-	0/8/17/17	0/2/2/2
5	14W	B	904	-	-	1/6/7/7	0/1/1/1
4	H4B	B	903	-	-	0/8/17/17	0/2/2/2
5	14W	A	902	-	-	1/6/7/7	0/1/1/1
3	HEM	A	900	1	-	3/14/54/54	-
3	HEM	B	900	1	-	3/14/54/54	-

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	900	HEM	FE-NB	5.74	2.12	1.94
3	A	900	HEM	FE-NB	4.85	2.09	1.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	903	H4B	C4A-N5	3.74	1.46	1.37
4	A	901	H4B	C4A-N5	3.29	1.45	1.37
3	B	900	HEM	CAC-C3C	-2.89	1.39	1.47
3	A	900	HEM	CAC-C3C	-2.67	1.40	1.47
4	B	903	H4B	C4-N3	2.66	1.43	1.38
5	B	904	14W	C5-C4	2.63	1.43	1.38
4	B	903	H4B	C6-N5	2.58	1.52	1.47
5	A	902	14W	C2-C3	2.49	1.43	1.39
4	A	901	H4B	C6-N5	2.49	1.51	1.47
5	B	904	14W	C6-C1	2.45	1.43	1.38
3	B	900	HEM	FE-ND	-2.39	1.87	1.94
4	A	901	H4B	C8A-N1	2.35	1.39	1.36
3	A	900	HEM	CAB-C3B	-2.33	1.41	1.47
5	A	902	14W	C5-C6	2.22	1.42	1.38
5	A	902	14W	C5-C4	2.20	1.42	1.38
5	A	902	14W	C6-C1	2.17	1.43	1.38
3	A	900	HEM	C3B-C4B	2.13	1.49	1.44
3	A	900	HEM	C3D-C2D	-2.12	1.32	1.36
3	B	900	HEM	CMC-C2C	2.11	1.55	1.50
5	B	904	14W	C4-C3	2.11	1.43	1.38
4	B	903	H4B	C8A-N1	2.10	1.39	1.36
3	A	900	HEM	C4A-C3A	2.04	1.47	1.43
5	A	902	14W	C4-C3	2.03	1.42	1.38

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	901	H4B	C2-N1-C8A	5.17	122.49	113.36
4	B	903	H4B	C2-N1-C8A	5.16	122.46	113.36
3	B	900	HEM	C3B-C4B-NB	4.10	112.41	109.47
3	A	900	HEM	CMD-C2D-C1D	3.77	130.93	125.03
5	A	902	14W	C10-C9-N8	3.71	120.08	114.81
3	A	900	HEM	C3B-C4B-NB	3.60	112.05	109.47
3	A	900	HEM	C4B-C3B-C2B	-3.53	104.03	107.28
3	B	900	HEM	C4B-C3B-C2B	-3.51	104.06	107.28
4	A	901	H4B	N3-C2-N1	-3.39	117.11	123.32
3	B	900	HEM	CHC-C4B-NB	-3.27	120.91	124.42
3	A	900	HEM	CAD-C3D-C2D	-3.24	121.81	127.87
4	B	903	H4B	N3-C2-N1	-3.14	117.57	123.32
5	B	904	14W	C10-C9-N8	3.10	119.22	114.81
3	A	900	HEM	CHC-C4B-NB	-2.71	121.51	124.42
3	A	900	HEM	CAD-C3D-C4D	2.58	129.19	124.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	901	H4B	O10-C10-C9	-2.56	105.53	109.77
3	A	900	HEM	C3B-C2B-C1B	2.46	108.26	106.41
5	A	902	14W	C1-C7-N8	-2.29	108.25	113.07
4	B	903	H4B	O10-C10-C9	-2.29	105.99	109.77
3	B	900	HEM	C2D-C1D-ND	2.25	112.50	109.90
3	A	900	HEM	CHD-C4C-NC	-2.20	122.06	124.45
3	A	900	HEM	CAA-C2A-C1A	2.09	129.01	124.94

There are no chirality outliers.

All (8) torsion outliers are listed below:

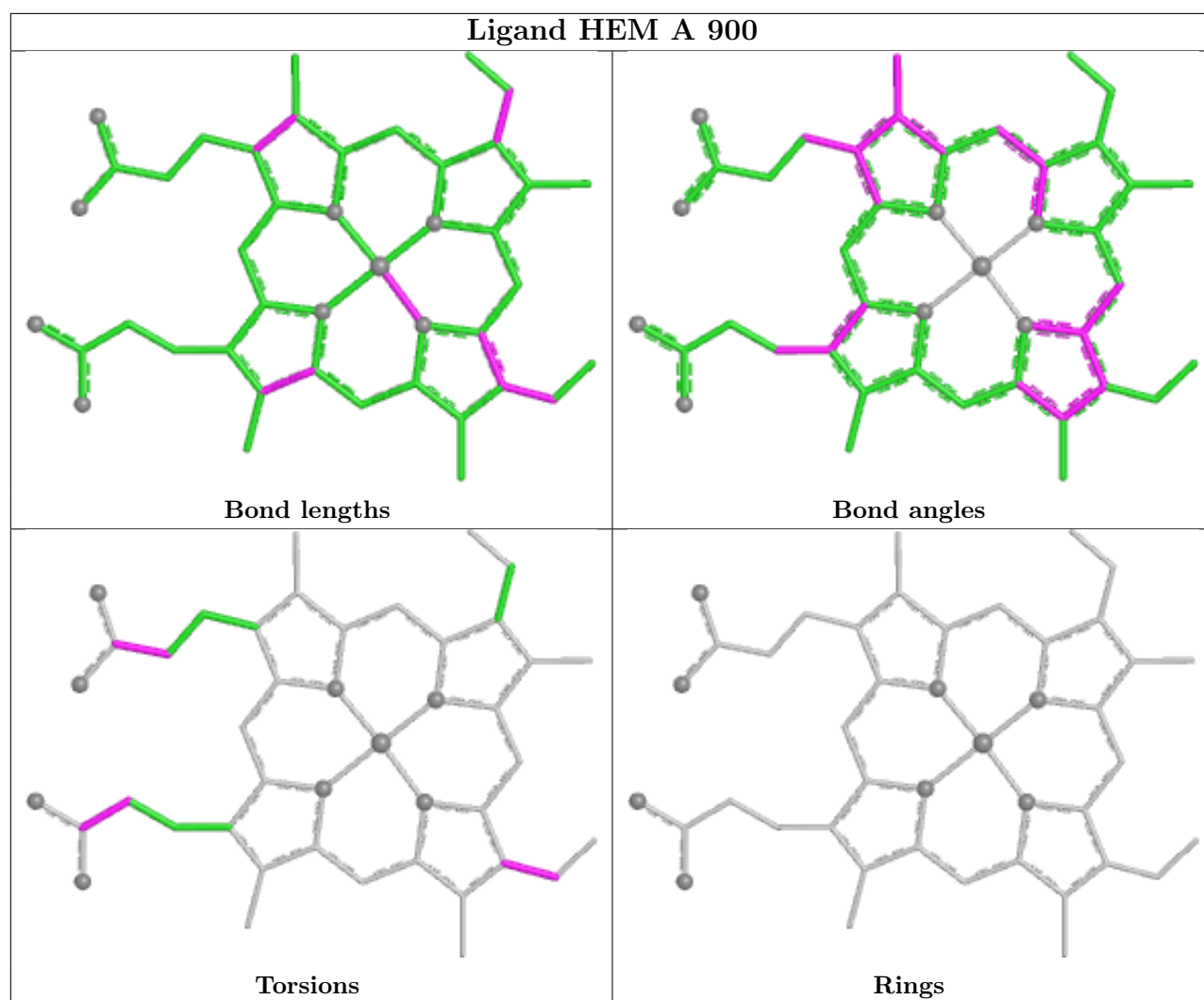
Mol	Chain	Res	Type	Atoms
3	A	900	HEM	C4B-C3B-CAB-CBB
3	B	900	HEM	C4B-C3B-CAB-CBB
5	A	902	14W	C1-C7-N8-C9
5	B	904	14W	C1-C7-N8-C9
3	B	900	HEM	CAD-CBD-CGD-O2D
3	B	900	HEM	CAD-CBD-CGD-O1D
3	A	900	HEM	CAD-CBD-CGD-O2D
3	A	900	HEM	CAA-CBA-CGA-O2A

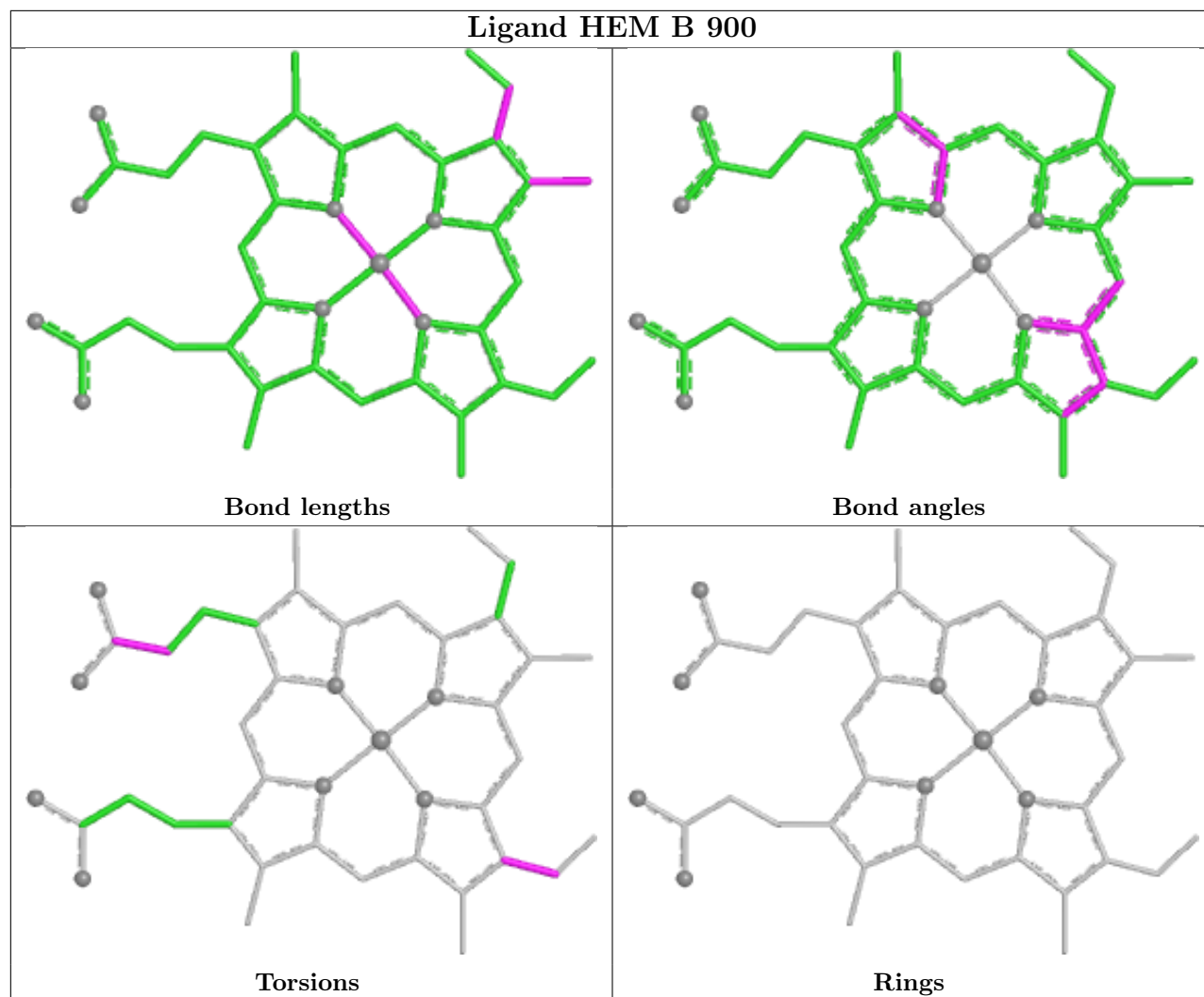
There are no ring outliers.

4 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	903	H4B	2	0
5	A	902	14W	1	0
3	A	900	HEM	3	0
3	B	900	HEM	2	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	419/419 (100%)	-0.35	5 (1%) 76 75	32, 49, 72, 89	0
1	B	419/419 (100%)	-0.37	10 (2%) 59 56	32, 49, 69, 93	0
All	All	838/838 (100%)	-0.36	15 (1%) 67 65	32, 49, 71, 93	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	104	CYS	5.0
1	A	106	SER	4.6
1	B	103	THR	4.4
1	B	106	SER	4.0
1	A	103	THR	3.4
1	B	104	CYS	2.9
1	B	101	ASP	2.8
1	B	269	GLY	2.6
1	B	154	GLU	2.5
1	A	102	PHE	2.5
1	A	100	SER	2.5
1	B	107	LYS	2.3
1	B	268	ASP	2.3
1	B	102	PHE	2.3
1	B	109	CYS	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

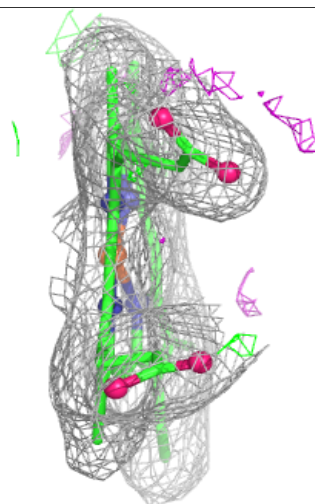
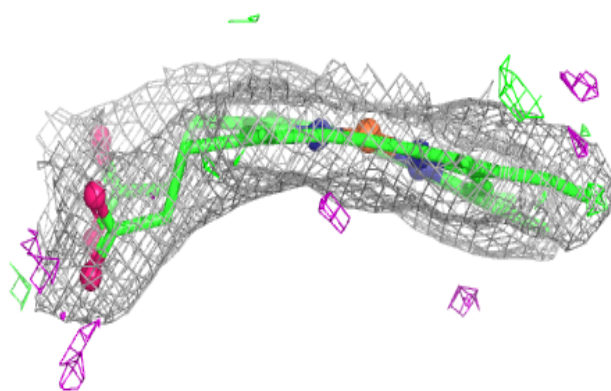
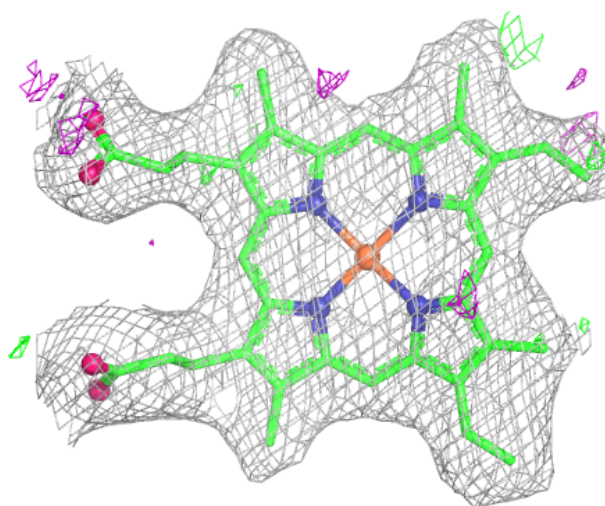
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

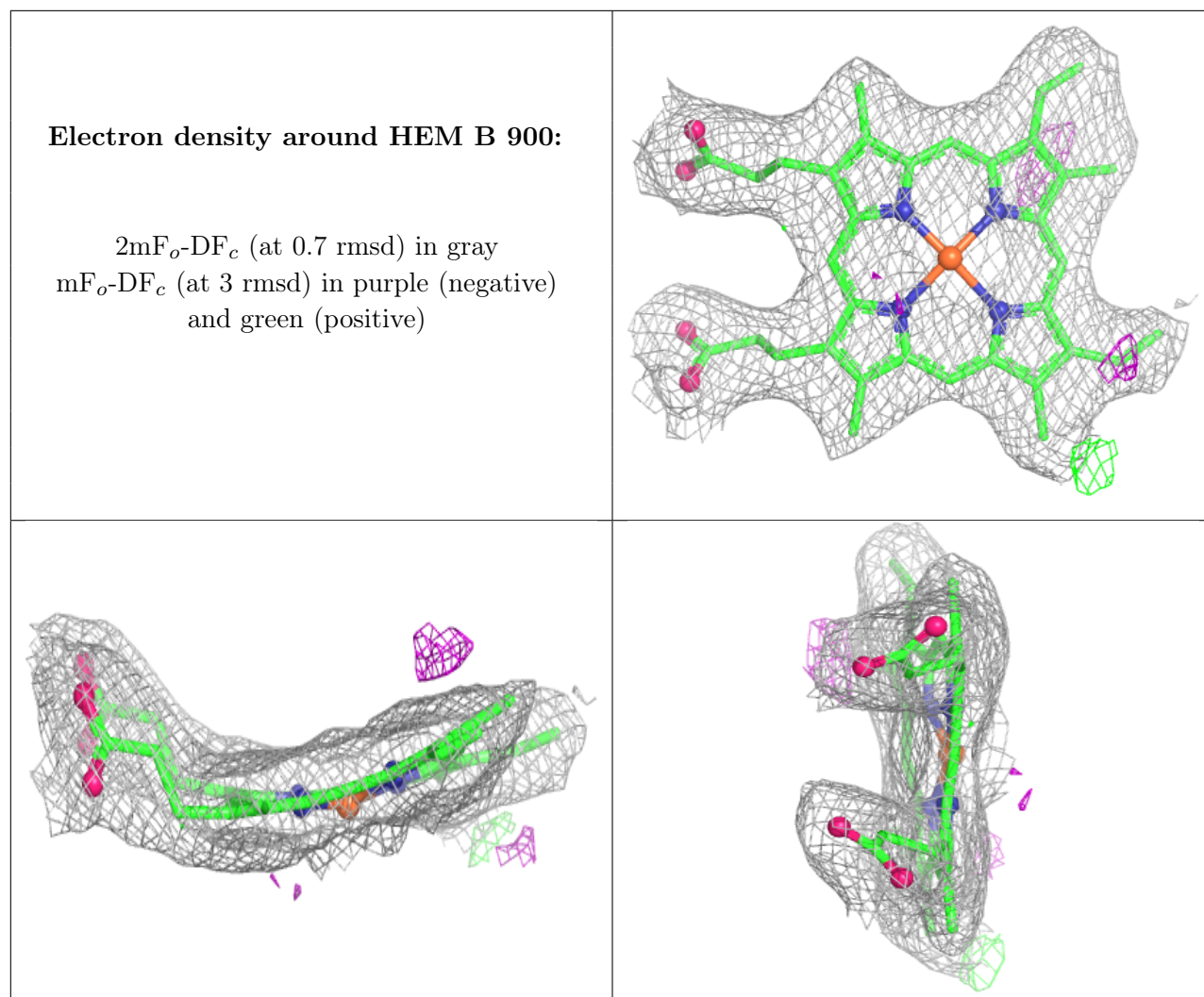
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	ZN	B	950	1/1	0.93	0.36	45,45,45,45	1
5	14W	A	902	13/13	0.94	0.07	37,39,40,41	0
5	14W	B	904	13/13	0.94	0.08	38,40,43,43	0
4	H4B	B	903	17/17	0.95	0.06	22,35,40,42	0
2	ZN	A	950	1/1	0.96	0.36	49,49,49,49	1
4	H4B	A	901	17/17	0.97	0.05	18,35,47,48	0
3	HEM	A	900	43/43	0.97	0.06	26,43,57,64	0
3	HEM	B	900	43/43	0.98	0.05	30,42,48,53	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 HEM A 900:

$2mF_o-DF_c$ (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.