



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 04:02 PM UTC

PDB ID : 3VLM / pdb_00003vlm
Title : Crystal Structure Analysis of the Met244Ala Variant of KatG from *Haloarcula marismortui*
Authors : Sato, T.; Ten-i, T.; Higuchi, W.; Yoshimatsu, K.; Fujiwara, T.
Deposited on : 2011-12-01
Resolution : 2.33 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

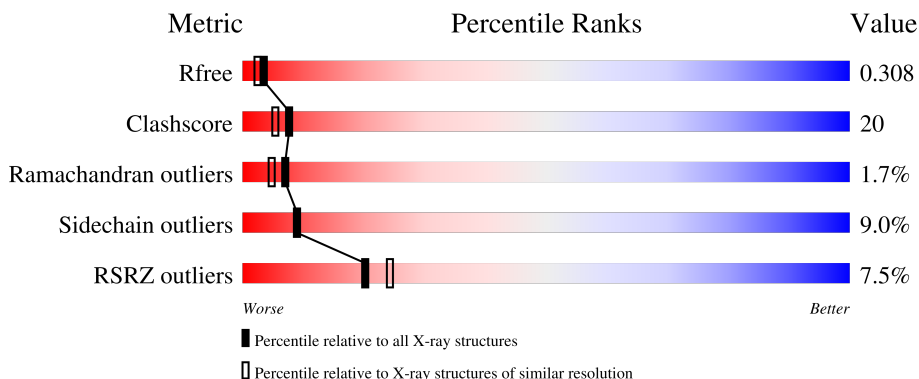
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.33 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	3031 (2.36-2.32)
Clashscore	190562	3127 (2.36-2.32)
Ramachandran outliers	187476	3095 (2.36-2.32)
Sidechain outliers	187428	3095 (2.36-2.32)
RSRZ outliers	180081	3033 (2.36-2.32)

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	737	5% 51% 34% 5% 10%
1	B	737	9% 54% 31% 10%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10623 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.

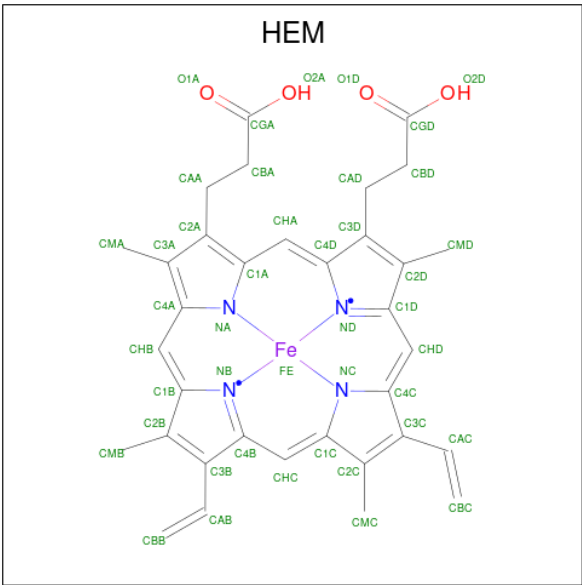
- Molecule 1 is a protein called Catalase-peroxidase 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	663	Total	C	N	O	S	0	0	0
			5201	3257	872	1054	18			
1	B	660	Total	C	N	O	S	0	0	0
			5185	3247	872	1049	17			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	244	ALA	MET	engineered mutation	UNP O59651
A	732	HIS	-	expression tag	UNP O59651
A	733	HIS	-	expression tag	UNP O59651
A	734	HIS	-	expression tag	UNP O59651
A	735	HIS	-	expression tag	UNP O59651
A	736	HIS	-	expression tag	UNP O59651
A	737	HIS	-	expression tag	UNP O59651
B	244	ALA	MET	engineered mutation	UNP O59651
B	732	HIS	-	expression tag	UNP O59651
B	733	HIS	-	expression tag	UNP O59651
B	734	HIS	-	expression tag	UNP O59651
B	735	HIS	-	expression tag	UNP O59651
B	736	HIS	-	expression tag	UNP O59651
B	737	HIS	-	expression tag	UNP O59651

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



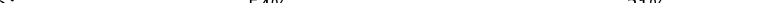
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

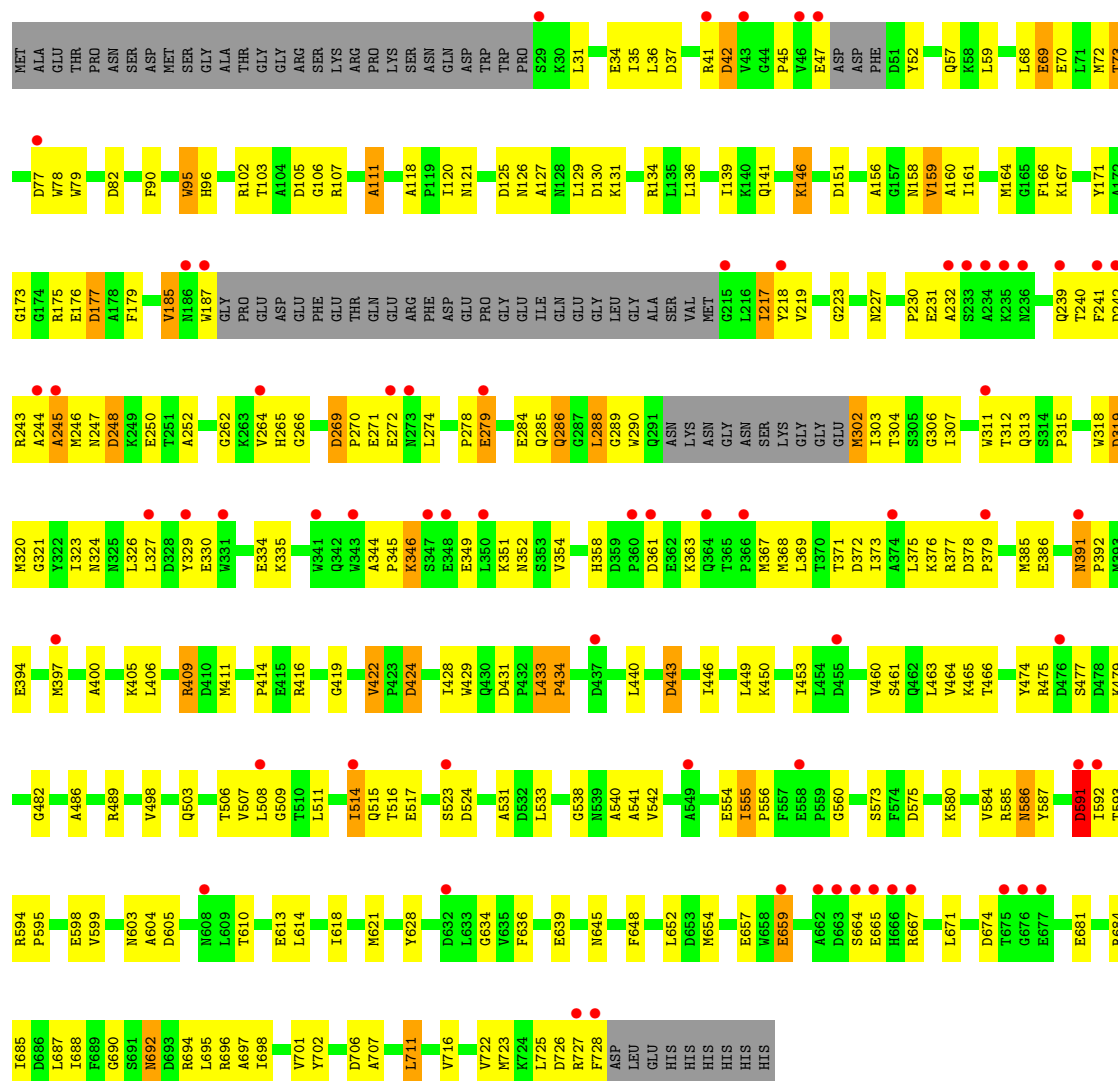
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	100	Total	O	0	0
			100	100		
3	B	51	Total	O	0	0
			51	51		

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.

- [illegible]

- Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	315.83Å 81.05Å 74.84Å 90.00° 100.04° 90.00°	Depositor
Resolution (Å)	24.94 – 2.33 24.94 – 2.33	Depositor EDS
% Data completeness (in resolution range)	94.8 (24.94-2.33) 94.8 (24.94-2.33)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.15 (at 2.01Å)	Xtriage
Refinement program	CNS, REFMAC 5.5.0109	Depositor
R, R_{free}	0.283 , 0.325 0.267 , 0.308	Depositor DCC
R_{free} test set	5841 reflections (4.64%)	wwPDB-VP
Wilson B-factor (Å ²)	28.8	Xtriage
Anisotropy	0.174	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 45.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.53$, $\langle L^2 \rangle = 0.37$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	10623	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.60% 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: 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	1.16	8/5322 (0.2%)	1.27	29/7230 (0.4%)
1	B	1.03	4/5307 (0.1%)	1.20	28/7211 (0.4%)
All	All	1.10	12/10629 (0.1%)	1.24	57/14441 (0.4%)

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	431	ASP	CA-C	6.28	1.60	1.52
1	B	139	ILE	CA-CB	-6.21	1.47	1.54
1	A	618	ILE	CA-C	-6.14	1.45	1.52
1	B	697	ALA	CA-CB	-6.04	1.43	1.53
1	A	253	ALA	CA-CB	5.95	1.62	1.53
1	A	446	ILE	CA-CB	5.76	1.60	1.54
1	A	697	ALA	CA-CB	-5.61	1.44	1.53
1	A	261	PHE	CA-C	5.59	1.59	1.52
1	B	422	VAL	CA-C	-5.56	1.48	1.53
1	B	159	VAL	CA-CB	-5.36	1.48	1.54
1	A	471	ALA	CA-C	5.29	1.59	1.52
1	A	360	PRO	N-CA	5.06	1.53	1.47

All (57) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	359	ASP	N-CA-C	11.25	122.62	108.11
1	A	500	GLU	CA-C-N	8.53	128.11	119.24
1	A	500	GLU	C-N-CA	8.53	128.11	119.24
1	A	571	ALA	CA-C-N	-8.20	111.20	119.56
1	A	571	ALA	C-N-CA	-8.20	111.20	119.56
1	A	136	LEU	N-CA-C	-7.86	102.77	113.30
1	A	86	TYR	N-CA-C	7.83	120.71	111.71

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	620	GLY	N-CA-C	-7.80	103.37	112.73
1	A	419	GLY	CA-C-N	7.71	127.43	119.56
1	A	419	GLY	C-N-CA	7.71	127.43	119.56
1	B	422	VAL	CB-CA-C	-7.47	103.99	110.94
1	B	79	TRP	CA-C-N	7.03	127.00	119.76
1	B	79	TRP	C-N-CA	7.03	127.00	119.76
1	B	269	ASP	CA-C-N	7.02	126.51	118.85
1	B	269	ASP	C-N-CA	7.02	126.51	118.85
1	A	256	ALA	N-CA-C	6.64	119.34	111.71
1	A	237	ILE	N-CA-C	-6.60	104.42	110.82
1	B	618	ILE	CB-CA-C	-6.44	103.64	111.88
1	B	422	VAL	N-CA-C	-6.39	103.13	108.63
1	A	313	GLN	CA-C-N	-6.36	115.77	122.89
1	A	313	GLN	C-N-CA	-6.36	115.77	122.89
1	B	391	ASN	N-CA-C	6.27	120.47	108.12
1	B	95	TRP	CA-C-O	-6.20	114.27	120.90
1	A	36	LEU	N-CA-C	-6.18	104.80	112.90
1	A	229	ASP	CA-C-N	6.09	125.57	119.24
1	A	229	ASP	C-N-CA	6.09	125.57	119.24
1	A	143	TYR	N-CA-C	-5.99	106.95	114.56
1	B	431	ASP	CA-C-N	5.81	125.78	120.03
1	B	431	ASP	C-N-CA	5.81	125.78	120.03
1	B	614	LEU	N-CA-C	-5.79	104.16	111.11
1	B	303	ILE	N-CA-C	5.74	117.02	108.46
1	B	433	LEU	CA-C-N	5.74	127.02	119.84
1	B	433	LEU	C-N-CA	5.74	127.02	119.84
1	A	370	THR	N-CA-C	-5.71	105.13	111.36
1	B	400	ALA	N-CA-C	-5.70	104.97	111.07
1	A	468	TRP	N-CA-C	-5.68	105.01	111.14
1	B	346	LYS	N-CA-C	-5.66	105.64	112.54
1	A	313	GLN	N-CA-C	5.64	119.27	112.38
1	A	527	GLN	N-CA-C	5.60	117.30	108.96
1	B	111	ALA	N-CA-C	5.46	117.94	111.33
1	B	706	ASP	N-CA-C	5.46	119.64	112.92
1	A	555	ILE	N-CA-CB	5.42	115.12	110.08
1	B	279	GLU	N-CA-C	5.39	118.95	112.38
1	A	414	PRO	N-CA-C	5.37	120.82	113.84
1	B	141	GLN	N-CA-C	5.35	117.80	111.33
1	A	311	TRP	CA-C-N	-5.33	114.22	122.59
1	A	311	TRP	C-N-CA	-5.33	114.22	122.59
1	A	166	PHE	N-CA-C	5.30	117.72	108.76
1	B	555	ILE	CA-C-N	-5.23	114.34	119.99

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	555	ILE	C-N-CA	-5.23	114.34	119.99
1	B	659	GLU	CA-C-N	5.16	125.07	119.85
1	B	659	GLU	C-N-CA	5.16	125.07	119.85
1	B	232	ALA	N-CA-C	5.16	116.98	111.36
1	A	273	ASN	N-CA-C	-5.14	106.94	114.39
1	B	156	ALA	N-CA-C	-5.07	105.84	111.36
1	B	240	THR	N-CA-C	-5.03	105.68	111.07
1	A	422	VAL	CB-CA-C	-5.03	105.46	110.89

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5201	0	4891	230	0
1	B	5185	0	4880	183	0
2	A	43	0	30	4	0
2	B	43	0	30	5	0
3	A	100	0	0	12	0
3	B	51	0	0	4	0
All	All	10623	0	9831	408	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 (408) 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:95:TRP:CH2	1:B:218:TYR:HE1	1.12	1.62
1:B:95:TRP:HH2	1:B:218:TYR:CE1	1.19	1.59
1:B:95:TRP:CH2	1:B:218:TYR:CE1	1.94	1.52
1:B:95:TRP:CZ2	1:B:218:TYR:HE1	1.58	1.22
1:B:95:TRP:HH2	1:B:218:TYR:CZ	1.69	1.08
1:B:584:VAL:HG13	1:B:621:MET:HG2	1.38	1.03

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:105:ASP:OD2	1:A:107:ARG:HD2	1.63	0.98
1:B:95:TRP:CZ2	1:B:218:TYR:CE1	2.43	0.94
1:A:474:TYR:CD1	1:A:725:LEU:HB3	2.08	0.89
1:A:474:TYR:CE1	1:A:725:LEU:HB3	2.10	0.87
1:B:555:ILE:HD12	1:B:716:VAL:HG13	1.57	0.86
1:A:245:ALA:HB3	1:A:409:ARG:NH1	1.94	0.82
1:A:41:ARG:HD2	1:B:41:ARG:HD2	1.62	0.81
1:A:101:TYR:CD2	1:A:409:ARG:HD3	2.19	0.78
1:A:101:TYR:CE2	1:A:409:ARG:HD3	2.20	0.77
1:A:470:SER:OG	1:A:483:ALA:HB1	1.82	0.77
1:B:486:ALA:HB2	1:B:531:ALA:HB2	1.68	0.75
1:B:367:MET:HE1	1:B:369:LEU:HD23	1.68	0.75
1:B:246:MET:HE1	1:B:405:LYS:HD3	1.66	0.75
1:A:581:VAL:HG12	1:A:581:VAL:O	1.85	0.75
1:A:320:MET:HE2	1:A:389:GLN:HA	1.67	0.74
1:A:405:LYS:HE3	1:A:429:TRP:CD1	2.22	0.74
1:A:245:ALA:HB3	1:A:409:ARG:HH12	1.52	0.74
1:B:692:ASN:ND2	1:B:695:LEU:H	1.85	0.74
1:A:34:GLU:OE2	1:A:183:LYS:HE2	1.88	0.73
1:A:95:TRP:HH2	1:A:218:TYR:CD1	2.07	0.72
1:B:449:LEU:O	1:B:453:ILE:HG13	1.90	0.72
1:B:318:TRP:O	1:B:319:ASP:HB3	1.88	0.72
1:A:677:GLU:O	1:A:677:GLU:HG2	1.89	0.72
1:A:308:GLU:H	1:A:342:GLN:HE22	1.34	0.72
1:A:479:LYS:HD2	1:A:728:PHE:CZ	2.24	0.71
1:A:565:GLY:H	1:A:568:HIS:HD2	1.38	0.71
1:A:412:GLY:HA3	1:A:479:LYS:HB2	1.73	0.71
1:B:34:GLU:HB2	3:B:753:HOH:O	1.90	0.70
1:B:95:TRP:HE3	1:B:406:LEU:HD21	1.56	0.69
1:B:319:ASP:OD1	1:B:319:ASP:C	2.35	0.69
1:A:486:ALA:HB3	1:A:489:ARG:NH1	2.09	0.68
1:A:597:GLU:O	1:A:600:LEU:HB3	1.93	0.68
1:A:239:GLN:NE2	1:A:243:ARG:HH21	1.92	0.68
1:B:160:ALA:O	1:B:164:MET:HE2	1.94	0.67
1:A:613:GLU:HA	3:A:760:HOH:O	1.94	0.67
1:B:610:THR:OG1	1:B:613:GLU:HG3	1.94	0.67
1:B:440:LEU:HD21	1:B:560:GLY:HA2	1.77	0.67
1:B:131:LYS:NZ	1:B:289:GLY:O	2.28	0.66
1:B:376:LYS:HG2	1:B:376:LYS:O	1.96	0.66
1:B:90:PHE:CE1	1:B:136:LEU:HD21	2.31	0.66
1:B:368:MET:HG3	1:B:372:ASP:HB2	1.78	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:645:ASN:HA	1:A:711:LEU:HD23	1.78	0.65
1:B:95:TRP:CH2	1:B:218:TYR:CZ	2.58	0.65
1:B:685:ILE:O	1:B:688:ILE:HG22	1.97	0.65
1:A:30:LYS:HG2	3:A:802:HOH:O	1.95	0.65
1:A:95:TRP:HH2	1:A:218:TYR:HD1	1.43	0.64
1:A:471:ALA:N	1:A:483:ALA:HB2	2.13	0.64
1:B:391:ASN:OD1	1:B:394:GLU:HB2	1.97	0.64
1:B:701:VAL:O	1:B:707:ALA:HB2	1.97	0.64
1:A:122:SER:HB3	1:A:277:GLU:HG3	1.79	0.64
1:A:310:PRO:HG3	1:A:354:VAL:HG11	1.80	0.64
1:A:102:ARG:HH21	1:A:105:ASP:CG	2.06	0.64
1:A:308:GLU:H	1:A:342:GLN:NE2	1.96	0.64
1:A:38:GLN:NE2	1:A:183:LYS:HD2	2.13	0.63
1:A:474:TYR:CD2	1:A:475:ARG:N	2.66	0.63
1:A:388:PHE:CG	1:A:395:PHE:HB2	2.33	0.63
1:A:264:VAL:HG23	1:A:307:ILE:O	1.99	0.63
1:A:259:HIS:HB3	1:A:311:TRP:CD2	2.34	0.63
1:A:470:SER:OG	1:A:483:ALA:CB	2.47	0.63
1:A:470:SER:C	1:A:483:ALA:HB2	2.24	0.63
1:B:591:ASP:OD2	1:B:591:ASP:N	2.31	0.63
1:A:259:HIS:HB3	1:A:311:TRP:CE2	2.33	0.62
1:B:584:VAL:HG13	1:B:621:MET:CG	2.23	0.62
1:B:371:THR:HG21	2:B:800:HEM:HMB3	1.80	0.62
1:B:465:LYS:HD3	3:B:749:HOH:O	1.98	0.62
1:B:95:TRP:CH2	1:B:218:TYR:OH	2.52	0.62
1:A:333:PRO:HD3	1:A:343:TRP:CZ3	2.35	0.62
2:B:800:HEM:HMB1	2:B:800:HEM:HBB2	1.82	0.62
1:A:486:ALA:HB3	1:A:489:ARG:HH11	1.63	0.61
1:A:87:GLY:HA2	1:A:160:ALA:HB1	1.82	0.61
1:A:386:GLU:HG3	3:A:761:HOH:O	2.00	0.61
1:A:128:ASN:HA	1:A:130:ASP:OD2	2.00	0.60
1:A:242:ASP:O	1:A:244:ALA:N	2.32	0.60
1:A:409:ARG:HH21	1:A:409:ARG:HG3	1.66	0.60
1:B:450:LYS:HG3	1:B:540:ALA:HB2	1.82	0.60
1:A:65:LYS:HE3	1:A:159:VAL:HG22	1.83	0.60
1:A:594:ARG:H	1:A:594:ARG:HE	1.50	0.60
1:A:474:TYR:HD1	1:A:725:LEU:CB	2.14	0.60
1:A:140:LYS:HG2	1:A:149:TRP:CE2	2.37	0.59
1:A:460:VAL:HG23	3:A:741:HOH:O	2.01	0.59
1:B:120:ILE:HD11	1:B:187:TRP:CZ2	2.37	0.59
1:A:162:GLU:HA	1:A:166:PHE:O	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:474:TYR:HD1	1:A:725:LEU:HB3	1.62	0.59
1:B:223:GLY:HA3	1:B:227:ASN:O	2.01	0.59
1:B:274:LEU:HD22	1:B:290:TRP:HB3	1.85	0.59
1:B:368:MET:HG3	1:B:372:ASP:CB	2.33	0.59
1:B:244:ALA:O	1:B:245:ALA:HB3	2.02	0.59
1:B:35:ILE:HG13	1:B:36:LEU:N	2.17	0.59
1:A:320:MET:HE3	1:A:323:ILE:HB	1.83	0.59
1:A:35:ILE:HD11	1:A:605:ASP:HB2	1.83	0.59
1:B:134:ARG:CZ	1:B:288:LEU:HD21	2.33	0.59
1:B:479:LYS:HB3	1:B:728:PHE:CZ	2.38	0.58
1:A:86:TYR:O	1:A:87:GLY:C	2.47	0.58
1:A:245:ALA:CB	1:A:409:ARG:HH12	2.15	0.58
1:A:361:ASP:O	1:A:362:GLU:HB2	2.03	0.58
1:B:711:LEU:O	1:B:711:LEU:HG	2.03	0.58
1:B:68:LEU:HB3	1:B:159:VAL:HG11	1.86	0.58
1:A:359:ASP:OD1	1:A:360:PRO:HD2	2.03	0.58
1:A:474:TYR:CD1	1:A:725:LEU:CB	2.84	0.58
1:A:320:MET:CE	1:A:323:ILE:HB	2.34	0.58
1:B:313:GLN:HA	1:B:354:VAL:HG22	1.86	0.58
1:A:41:ARG:H	1:A:41:ARG:HD3	1.69	0.58
1:A:335:LYS:HE2	1:A:339:GLY:O	2.04	0.58
1:A:344:ALA:HB2	3:A:778:HOH:O	2.04	0.58
1:A:692:ASN:ND2	1:A:694:ARG:H	2.03	0.57
1:A:467:ALA:HA	1:A:534:ILE:HG22	1.86	0.57
1:A:115:GLN:O	1:A:121:ASN:HB3	2.05	0.57
1:B:270:PRO:HB3	1:B:274:LEU:HD12	1.84	0.57
1:A:354:VAL:HG23	1:A:366:PRO:HD3	1.86	0.57
1:A:474:TYR:HE1	1:A:725:LEU:HB3	1.67	0.56
1:A:354:VAL:O	1:A:363:LYS:HA	2.05	0.56
1:A:648:PHE:HE1	1:A:711:LEU:CD2	2.18	0.56
1:B:330:GLU:HB3	1:B:346:LYS:HE3	1.87	0.56
2:B:800:HEM:HBB2	2:B:800:HEM:CMB	2.36	0.56
1:B:244:ALA:O	1:B:245:ALA:CB	2.53	0.56
1:B:449:LEU:HD22	1:B:533:LEU:HD21	1.87	0.56
1:A:452:GLU:HG2	1:A:514:ILE:HG23	1.87	0.56
1:B:166:PHE:HD1	1:B:318:TRP:CD2	2.24	0.55
1:A:412:GLY:CA	1:A:479:LYS:HB2	2.35	0.55
1:B:134:ARG:CZ	1:B:288:LEU:CD2	2.84	0.55
1:A:245:ALA:HB3	1:A:409:ARG:CZ	2.36	0.55
1:A:405:LYS:HA	1:A:429:TRP:CZ2	2.41	0.55
1:B:344:ALA:HB1	1:B:345:PRO:HD2	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:159:VAL:O	1:B:160:ALA:C	2.49	0.55
1:A:498:VAL:HG23	1:A:587:TYR:CG	2.43	0.54
1:A:587:TYR:O	1:A:588:ILE:HG13	2.06	0.54
1:B:270:PRO:C	1:B:272:GLU:H	2.15	0.54
1:B:460:VAL:O	1:B:461:SER:C	2.49	0.54
1:B:262:GLY:HA3	2:B:800:HEM:O2D	2.07	0.54
1:B:95:TRP:CE3	1:B:406:LEU:HD21	2.42	0.54
1:B:511:LEU:HA	1:B:514:ILE:HG13	1.89	0.54
1:A:256:ALA:HB2	1:A:385:MET:HE1	1.89	0.54
1:A:463:LEU:HB2	1:A:541:ALA:HB2	1.90	0.54
1:A:582:ASP:C	1:A:582:ASP:OD1	2.50	0.54
1:B:443:ASP:HA	1:B:446:ILE:HD12	1.88	0.54
1:B:323:ILE:HG23	1:B:327:LEU:HD12	1.89	0.54
1:A:87:GLY:CA	1:A:160:ALA:HB1	2.38	0.53
1:A:125:ASP:HB2	1:A:216:LEU:HD23	1.91	0.53
1:B:118:ALA:HB2	1:B:279:GLU:HG3	1.91	0.53
1:B:482:GLY:HA3	1:B:726:ASP:OD2	2.09	0.53
1:A:95:TRP:NE1	1:A:96:HIS:HD2	2.06	0.53
1:A:82:ASP:C	1:A:84:GLY:H	2.16	0.53
1:A:405:LYS:HA	1:A:429:TRP:CH2	2.43	0.53
1:B:161:ILE:HG22	1:B:166:PHE:HB3	1.90	0.53
1:B:692:ASN:HD22	1:B:695:LEU:H	1.54	0.53
1:A:95:TRP:NE1	1:A:96:HIS:CD2	2.76	0.53
1:B:367:MET:HE1	1:B:369:LEU:CD2	2.37	0.53
1:A:95:TRP:CH2	1:A:218:TYR:CD1	2.94	0.53
1:A:87:GLY:HA2	1:A:160:ALA:CB	2.39	0.53
1:A:416:ARG:HD3	1:A:479:LYS:HG3	1.91	0.52
1:A:438:TYR:CD1	1:A:527:GLN:HB2	2.44	0.52
1:A:458:LEU:HA	1:A:462:GLN:OE1	2.09	0.52
2:A:800:HEM:HBC2	2:A:800:HEM:CMC	2.39	0.52
1:B:440:LEU:CD2	1:B:560:GLY:HA2	2.40	0.52
1:B:95:TRP:CD1	1:B:96:HIS:HD2	2.27	0.52
1:B:177:ASP:HB2	3:B:755:HOH:O	2.08	0.52
1:A:241:PHE:O	1:A:244:ALA:HB3	2.09	0.52
1:A:386:GLU:CD	3:A:761:HOH:O	2.53	0.52
1:A:720:SER:OG	1:A:724:LYS:NZ	2.42	0.52
1:B:474:TYR:CD1	1:B:474:TYR:C	2.88	0.52
1:A:134:ARG:HG2	1:B:654:MET:HE2	1.91	0.52
1:A:66:SER:O	1:A:69:GLU:HB2	2.10	0.52
1:A:645:ASN:HA	1:A:711:LEU:CD2	2.38	0.52
1:B:466:THR:OG1	1:B:507:VAL:HG11	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:524:ASP:OD2	1:B:524:ASP:C	2.53	0.51
1:B:587:TYR:HD1	1:B:628:TYR:CD1	2.28	0.51
1:B:555:ILE:HD12	1:B:716:VAL:CG1	2.35	0.51
1:A:249:LYS:HG2	1:A:398:ASN:OD1	2.11	0.51
1:A:35:ILE:CD1	1:A:605:ASP:HB2	2.40	0.51
1:A:610:THR:OG1	1:A:613:GLU:HG3	2.10	0.51
1:B:242:ASP:C	1:B:244:ALA:H	2.19	0.51
1:B:326:LEU:O	1:B:376:LYS:NZ	2.36	0.51
1:A:82:ASP:C	1:A:84:GLY:N	2.67	0.51
1:A:95:TRP:CD1	1:A:96:HIS:HD2	2.29	0.51
1:A:149:TRP:O	1:A:153:MET:HG3	2.11	0.51
1:A:486:ALA:O	1:A:489:ARG:HG2	2.10	0.51
1:A:82:ASP:OD1	1:A:265:HIS:NE2	2.43	0.50
1:A:213:VAL:HB	1:A:216:LEU:HD12	1.93	0.50
1:B:722:VAL:HA	1:B:725:LEU:HG	1.92	0.50
1:A:692:ASN:HD22	1:A:695:LEU:H	1.58	0.50
1:A:45:PRO:HG3	1:B:698:ILE:HG23	1.92	0.50
1:A:502:GLU:CD	1:A:502:GLU:H	2.18	0.50
1:B:692:ASN:ND2	1:B:692:ASN:C	2.69	0.50
1:A:648:PHE:HE1	1:A:711:LEU:HD21	1.77	0.50
1:B:185:VAL:HG11	1:B:187:TRP:CZ3	2.47	0.50
1:B:270:PRO:CB	1:B:274:LEU:HD12	2.41	0.50
1:A:320:MET:CE	1:A:389:GLN:HA	2.38	0.50
1:A:56:PHE:C	1:A:58:LYS:H	2.19	0.50
1:A:171:TYR:HD2	3:A:757:HOH:O	1.94	0.50
1:A:174:GLY:O	3:A:746:HOH:O	2.20	0.50
1:B:95:TRP:CD1	1:B:96:HIS:CD2	3.00	0.50
1:A:314:SER:O	1:A:316:THR:N	2.44	0.49
1:B:405:LYS:HE3	1:B:429:TRP:CD1	2.46	0.49
1:B:474:TYR:CD2	1:B:726:ASP:HB3	2.47	0.49
1:A:148:SER:HB2	1:A:177:ASP:CG	2.38	0.49
1:B:684:ARG:HG3	3:B:757:HOH:O	2.13	0.49
1:A:593:THR:H	1:A:594:ARG:HH21	1.59	0.49
1:B:311:TRP:NE1	1:B:372:ASP:OD2	2.39	0.49
1:A:666:HIS:H	1:A:667:ARG:HD3	1.77	0.49
1:A:701:VAL:O	1:A:704:SER:HB2	2.12	0.49
1:B:538:GLY:O	1:B:542:VAL:HG23	2.12	0.49
1:A:264:VAL:HG23	1:A:307:ILE:C	2.38	0.49
1:B:125:ASP:OD1	1:B:217:ILE:HG12	2.13	0.49
1:B:486:ALA:HB2	1:B:531:ALA:CB	2.41	0.49
1:B:603:ASN:C	1:B:605:ASP:N	2.71	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:125:ASP:OD2	1:A:125:ASP:O	2.30	0.49
1:A:264:VAL:HG23	1:A:264:VAL:O	2.12	0.49
1:A:644:THR:OG1	1:A:646:ASP:HB2	2.13	0.49
1:A:665:GLU:HB3	1:A:666:HIS:CE1	2.48	0.49
1:A:254:LEU:HG	1:A:402:ALA:HB1	1.95	0.48
1:B:158:ASN:ND2	1:B:171:TYR:HB2	2.28	0.48
1:A:694:ARG:O	1:A:698:ILE:HG13	2.14	0.48
1:B:265:HIS:HD2	1:B:302:MET:HG2	1.77	0.48
1:B:409:ARG:NH2	1:B:409:ARG:HG3	2.28	0.48
1:B:508:LEU:O	1:B:509:GLY:C	2.55	0.48
1:A:473:THR:O	1:A:481:GLY:HA3	2.13	0.48
1:B:68:LEU:O	1:B:72:MET:HG2	2.13	0.48
1:B:694:ARG:O	1:B:698:ILE:HG13	2.13	0.48
1:A:519:ASN:HD21	1:A:528:VAL:H	1.62	0.48
1:B:603:ASN:O	1:B:604:ALA:C	2.57	0.48
1:B:648:PHE:O	1:B:652:LEU:HG	2.13	0.48
1:A:140:LYS:HD3	1:A:149:TRP:CD2	2.49	0.48
1:A:474:TYR:HD2	1:A:475:ARG:H	1.62	0.48
1:A:484:ASN:HB3	1:A:726:ASP:OD2	2.14	0.48
1:A:259:HIS:CE1	2:A:800:HEM:NB	2.80	0.48
1:A:449:LEU:O	1:A:453:ILE:HG13	2.13	0.48
1:A:715:PHE:C	1:A:715:PHE:CD2	2.91	0.48
1:B:241:PHE:O	1:B:244:ALA:HB3	2.14	0.48
1:B:329:TYR:O	1:B:376:LYS:NZ	2.46	0.48
1:B:252:ALA:HB1	1:B:385:MET:SD	2.54	0.47
1:A:30:LYS:O	1:A:31:LEU:C	2.56	0.47
1:B:373:ILE:O	1:B:376:LYS:N	2.47	0.47
1:A:352:ASN:HB3	1:A:363:LYS:HB2	1.96	0.47
1:B:424:ASP:N	1:B:424:ASP:OD1	2.47	0.47
1:A:688:ILE:C	1:A:690:GLY:H	2.23	0.47
1:A:56:PHE:C	1:A:58:LYS:N	2.72	0.47
1:B:230:PRO:O	1:B:231:GLU:C	2.58	0.47
1:A:48:ASP:OD2	1:A:48:ASP:N	2.47	0.47
1:A:409:ARG:HG3	1:A:409:ARG:NH2	2.30	0.47
1:B:320:MET:HB2	1:B:324:ASN:ND2	2.30	0.47
1:B:688:ILE:C	1:B:690:GLY:H	2.23	0.47
1:A:218:TYR:HD2	1:A:218:TYR:HA	1.53	0.47
1:A:486:ALA:CB	1:A:489:ARG:HH11	2.27	0.47
1:A:673:ARG:HD3	1:B:78:TRP:HB3	1.97	0.47
1:A:444:GLU:HG3	1:A:445:GLU:H	1.79	0.46
1:B:486:ALA:CB	1:B:531:ALA:HB2	2.44	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:175:ARG:NH2	1:A:407:THR:O	2.41	0.46
1:B:111:ALA:O	1:B:175:ARG:HB3	2.14	0.46
1:B:134:ARG:NH1	1:B:288:LEU:HD21	2.29	0.46
1:B:664:SER:OG	1:B:667:ARG:HB2	2.15	0.46
1:B:313:GLN:C	1:B:315:PRO:HD3	2.40	0.46
1:B:634:GLY:HA2	1:B:636:PHE:CE2	2.51	0.46
1:B:727:ARG:HD2	1:B:727:ARG:HA	1.61	0.46
1:A:140:LYS:HG2	1:A:149:TRP:CZ2	2.51	0.46
1:A:361:ASP:O	1:A:362:GLU:CB	2.62	0.46
1:A:664:SER:OG	1:A:667:ARG:HB2	2.15	0.46
1:B:95:TRP:CZ3	1:B:218:TYR:OH	2.67	0.46
1:B:127:ALA:HB3	1:B:304:THR:HG23	1.96	0.46
1:B:285:GLN:O	1:B:286:GLN:C	2.56	0.46
1:A:137:LEU:HB3	1:A:138:PRO:HD3	1.97	0.46
1:A:367:MET:HE2	1:A:367:MET:HB2	1.53	0.46
1:A:395:PHE:O	1:A:395:PHE:CD1	2.68	0.46
1:A:42:ASP:OD1	1:A:42:ASP:N	2.49	0.46
1:A:386:GLU:CG	3:A:761:HOH:O	2.61	0.46
1:A:544:GLN:O	1:A:547:ALA:N	2.49	0.46
1:A:471:ALA:O	1:A:472:SER:C	2.59	0.46
1:B:474:TYR:HD2	1:B:726:ASP:HB3	1.81	0.46
1:A:443:ASP:OD1	1:A:443:ASP:N	2.41	0.45
1:A:718:THR:O	1:A:719:TRP:C	2.55	0.45
1:A:471:ALA:CA	1:A:483:ALA:HB2	2.47	0.45
1:A:509:GLY:O	1:A:513:ASN:OD1	2.34	0.45
1:A:587:TYR:C	1:A:588:ILE:HG13	2.41	0.45
1:B:34:GLU:OE1	1:B:34:GLU:N	2.50	0.45
1:A:269:ASP:OD1	1:A:271:GLU:HB2	2.17	0.45
1:A:440:LEU:HD23	1:A:440:LEU:HA	1.81	0.45
1:B:312:THR:OG1	1:B:315:PRO:HA	2.16	0.45
1:A:258:GLY:O	1:A:260:THR:N	2.50	0.45
1:A:379:PRO:HD2	3:A:832:HOH:O	2.17	0.45
1:B:42:ASP:OD1	1:B:42:ASP:N	2.50	0.45
1:B:105:ASP:OD2	1:B:107:ARG:HD2	2.15	0.45
1:B:173:GLY:O	1:B:411:MET:HE1	2.17	0.45
1:A:216:LEU:O	1:A:217:ILE:C	2.60	0.45
1:B:166:PHE:HE1	1:B:318:TRP:CZ2	2.35	0.45
1:B:264:VAL:HG12	2:B:800:HEM:HAA1	1.98	0.45
1:B:106:GLY:C	1:B:477:SER:OG	2.60	0.44
1:B:118:ALA:HB2	1:B:279:GLU:CG	2.46	0.44
1:A:248:ASP:HA	3:A:739:HOH:O	2.16	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:58:LYS:HB3	1:A:146:LYS:HZ2	1.82	0.44
1:A:67:ASP:HB3	1:A:139:ILE:HD13	2.00	0.44
1:A:213:VAL:HG12	1:A:214:MET:N	2.33	0.44
1:B:82:ASP:OD2	1:B:358:HIS:CE1	2.70	0.44
1:B:555:ILE:CG2	1:B:556:PRO:HD2	2.48	0.44
1:A:67:ASP:HB3	1:A:139:ILE:CD1	2.47	0.44
1:A:47:GLU:H	1:A:47:GLU:HG2	1.55	0.44
1:A:230:PRO:HG2	1:A:231:GLU:OE2	2.18	0.44
1:A:283:ILE:HD11	1:B:688:ILE:N	2.33	0.44
1:B:69:GLU:O	1:B:70:GLU:C	2.60	0.44
1:B:134:ARG:NH2	1:B:288:LEU:HD23	2.33	0.44
1:B:645:ASN:OD1	1:B:645:ASN:C	2.60	0.44
1:B:659:GLU:OE1	1:B:671:LEU:HD11	2.18	0.44
1:A:147:ILE:O	1:A:147:ILE:HG23	2.18	0.44
1:A:408:HIS:HB3	1:A:417:PHE:CE1	2.52	0.44
1:B:405:LYS:O	1:B:409:ARG:HB2	2.17	0.44
1:A:134:ARG:CZ	1:B:696:ARG:NH2	2.81	0.44
1:A:72:MET:HE1	1:A:90:PHE:CD2	2.53	0.43
1:B:45:PRO:HD2	1:B:179:PHE:CZ	2.52	0.43
1:B:52:TYR:OH	1:B:151:ASP:OD2	2.29	0.43
1:B:166:PHE:CD1	1:B:318:TRP:CD2	3.05	0.43
1:A:578:LYS:HA	1:A:579:PRO:HD3	1.88	0.43
1:A:658:TRP:CD1	1:A:658:TRP:N	2.85	0.43
1:B:242:ASP:C	1:B:242:ASP:OD2	2.61	0.43
1:B:515:GLN:O	1:B:516:THR:C	2.60	0.43
1:A:412:GLY:HA3	1:A:479:LYS:CB	2.46	0.43
1:A:656:THR:HA	1:A:671:LEU:O	2.18	0.43
1:B:166:PHE:HD2	1:B:167:LYS:H	1.59	0.43
1:B:585:ARG:O	1:B:586:ASN:C	2.62	0.43
1:A:92:ARG:HA	1:A:95:TRP:HD1	1.84	0.43
1:A:351:LYS:HA	1:A:365:THR:OG1	2.19	0.43
1:A:78:TRP:CD2	1:B:654:MET:HE3	2.54	0.43
1:B:59:LEU:HB2	1:B:146:LYS:HB3	1.99	0.43
1:A:489:ARG:NH2	1:A:564:ALA:O	2.49	0.43
1:B:103:THR:O	1:B:475:ARG:NH1	2.51	0.43
1:A:264:VAL:O	1:A:308:GLU:HA	2.19	0.43
1:A:479:LYS:HD2	1:A:728:PHE:CE2	2.53	0.43
1:A:285:GLN:NE2	1:A:291:GLN:HE22	2.17	0.43
1:B:103:THR:O	1:B:475:ARG:NH2	2.51	0.43
1:A:41:ARG:HD2	1:B:41:ARG:CD	2.43	0.43
1:A:265:HIS:ND1	1:A:265:HIS:N	2.67	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:302:MET:N	3:A:811:HOH:O	2.51	0.42
1:A:443:ASP:HA	1:A:446:ILE:HD12	2.01	0.42
2:A:800:HEM:HBC2	2:A:800:HEM:HMC1	2.01	0.42
1:B:166:PHE:CD2	1:B:167:LYS:N	2.77	0.42
1:B:329:TYR:CD1	1:B:329:TYR:N	2.87	0.42
1:B:373:ILE:HG22	1:B:377:ARG:HG3	2.00	0.42
1:B:419:GLY:O	1:B:422:VAL:HG23	2.19	0.42
1:B:595:PRO:O	1:B:598:GLU:HB2	2.19	0.42
1:A:405:LYS:HB2	1:A:429:TRP:CZ2	2.55	0.42
1:B:70:GLU:O	1:B:72:MET:N	2.51	0.42
1:B:246:MET:HA	1:B:250:GLU:OE1	2.20	0.42
1:A:259:HIS:CG	1:A:311:TRP:CZ2	3.07	0.42
1:A:665:GLU:HB3	1:A:666:HIS:ND1	2.34	0.42
1:B:503:GLN:O	1:B:506:THR:HB	2.20	0.42
1:B:702:TYR:O	1:B:707:ALA:HB3	2.19	0.42
1:B:247:ASN:OD1	1:B:247:ASN:C	2.62	0.42
1:A:126:ASN:HA	1:A:129:LEU:HD12	2.01	0.42
1:B:307:ILE:HG22	1:B:367:MET:HE3	2.02	0.42
1:B:463:LEU:O	1:B:464:VAL:C	2.61	0.42
1:A:65:LYS:HD2	1:A:155:LEU:HD11	2.02	0.42
1:A:185:VAL:HG11	1:A:187:TRP:CE2	2.54	0.42
1:A:354:VAL:CG2	1:A:366:PRO:HD3	2.50	0.42
1:B:95:TRP:HZ3	1:B:406:LEU:HD11	1.84	0.42
1:B:414:PRO:C	1:B:416:ARG:H	2.28	0.42
1:B:463:LEU:HB2	1:B:541:ALA:HB2	2.01	0.42
1:A:283:ILE:HD11	1:B:688:ILE:HA	2.02	0.41
1:B:266:GLY:HA3	1:B:306:GLY:C	2.44	0.41
1:A:127:ALA:HA	1:A:290:TRP:CH2	2.54	0.41
1:A:38:GLN:NE2	1:A:180:GLU:CD	2.78	0.41
1:A:93:MET:HB3	1:A:93:MET:HE2	1.77	0.41
1:A:107:ARG:CZ	1:A:606:LEU:HD21	2.51	0.41
1:B:73:THR:O	1:B:73:THR:OG1	2.38	0.41
1:B:247:ASN:O	1:B:248:ASP:C	2.64	0.41
1:A:220:ASN:OD1	1:A:222:GLU:HB2	2.21	0.41
1:A:536:LEU:HD12	1:A:536:LEU:HA	1.88	0.41
1:B:376:LYS:O	1:B:376:LYS:CG	2.66	0.41
1:A:264:VAL:HG12	2:A:800:HEM:HAA1	2.01	0.41
1:A:391:ASN:O	1:A:392:PRO:C	2.63	0.41
1:A:405:LYS:HE3	1:A:429:TRP:CG	2.53	0.41
1:A:461:SER:O	1:A:465:LYS:HB3	2.21	0.41
1:A:622:ARG:HB3	1:A:634:GLY:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:217:ILE:N	1:B:217:ILE:HD13	2.36	0.41
1:A:360:PRO:HB2	1:A:361:ASP:H	1.28	0.41
1:A:428:ILE:HD12	1:A:428:ILE:HA	1.80	0.41
1:A:464:VAL:HG23	1:A:541:ALA:HB1	2.02	0.41
1:A:617:LEU:HA	1:A:715:PHE:CD1	2.55	0.41
1:B:247:ASN:N	1:B:250:GLU:OE1	2.51	0.41
1:A:217:ILE:O	1:A:218:TYR:CD2	2.74	0.41
1:A:268:ASP:H	1:A:303:ILE:HG12	1.85	0.41
1:A:277:GLU:O	1:A:278:PRO:C	2.63	0.41
1:B:555:ILE:HG22	1:B:556:PRO:HD2	2.01	0.41
1:A:147:ILE:HD13	1:A:152:LEU:HD13	2.03	0.41
1:A:375:LEU:HD22	1:A:385:MET:HE2	2.02	0.41
1:A:621:MET:HE2	1:A:715:PHE:CE1	2.56	0.41
1:B:95:TRP:HZ3	1:B:218:TYR:HH	1.58	0.41
1:B:126:ASN:O	1:B:129:LEU:HB2	2.21	0.41
1:B:352:ASN:HB3	1:B:363:LYS:HB3	2.03	0.41
1:B:375:LEU:HD23	1:B:375:LEU:HA	1.90	0.41
1:B:433:LEU:HA	1:B:434:PRO:HD2	1.68	0.41
1:A:441:ILE:HA	1:A:445:GLU:OE1	2.21	0.41
1:A:512:GLU:HG3	1:A:530:LEU:HD22	2.02	0.41
1:B:34:GLU:HA	1:B:37:ASP:OD2	2.21	0.41
1:A:95:TRP:CD1	1:A:96:HIS:CD2	3.09	0.40
1:A:510:THR:O	1:A:514:ILE:HG12	2.22	0.40
1:B:378:ASP:O	1:B:379:PRO:C	2.62	0.40
1:B:409:ARG:CG	1:B:409:ARG:HH21	2.34	0.40
1:A:536:LEU:O	1:A:537:GLY:C	2.62	0.40
1:B:31:LEU:HA	1:B:31:LEU:HD23	1.65	0.40
1:A:132:ALA:O	1:A:135:LEU:HB2	2.21	0.40
1:A:116:ARG:HD3	1:A:149:TRP:HB3	2.03	0.40
1:A:274:LEU:HD13	1:A:290:TRP:HB3	2.03	0.40
1:A:474:TYR:CD2	1:A:474:TYR:C	2.99	0.40
1:B:320:MET:SD	1:B:392:PRO:HB3	2.62	0.40
1:B:391:ASN:HA	1:B:392:PRO:HD3	1.94	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	653/737 (89%)	573 (88%)	68 (10%)	12 (2%)	6	4
1	B	652/737 (88%)	558 (86%)	84 (13%)	10 (2%)	8	6
All	All	1305/1474 (88%)	1131 (87%)	152 (12%)	22 (2%)	7	5

All (22) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	360	PRO
1	A	361	ASP
1	B	321	GLY
1	B	592	ILE
1	A	87	GLY
1	A	409	ARG
1	A	630	ASP
1	B	245	ALA
1	B	271	GLU
1	A	243	ARG
1	A	244	ALA
1	B	434	PRO
1	A	259	HIS
1	A	581	VAL
1	A	592	ILE
1	B	243	ARG
1	B	319	ASP
1	B	674	ASP
1	A	217	ILE
1	B	286	GLN
1	B	591	ASP
1	A	678	VAL

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	545/607 (90%)	502 (92%)	43 (8%)	11	12
1	B	543/607 (90%)	488 (90%)	55 (10%)	7	6
All	All	1088/1214 (90%)	990 (91%)	98 (9%)	9	9

All (98) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	31	LEU
1	A	35	ILE
1	A	42	ASP
1	A	47	GLU
1	A	48	ASP
1	A	51	ASP
1	A	57	GLN
1	A	69	GLU
1	A	95	TRP
1	A	103	THR
1	A	139	ILE
1	A	176	GLU
1	A	233	SER
1	A	240	THR
1	A	269	ASP
1	A	291	GLN
1	A	302	MET
1	A	342	GLN
1	A	361	ASP
1	A	367	MET
1	A	368	MET
1	A	378	ASP
1	A	409	ARG
1	A	440	LEU
1	A	443	ASP
1	A	445	GLU
1	A	448	GLU

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Mol	Chain	Res	Type
1	A	455	ASP
1	A	494	LYS
1	A	500	GLU
1	A	517	GLU
1	A	529	SER
1	A	578	LYS
1	A	586	ASN
1	A	594	ARG
1	A	638	ASP
1	A	640	PRO
1	A	651	LEU
1	A	659	GLU
1	A	664	SER
1	A	669	LYS
1	A	677	GLU
1	A	679	LYS
1	B	42	ASP
1	B	47	GLU
1	B	57	GLN
1	B	69	GLU
1	B	73	THR
1	B	77	ASP
1	B	102	ARG
1	B	121	ASN
1	B	130	ASP
1	B	146	LYS
1	B	176	GLU
1	B	177	ASP
1	B	185	VAL
1	B	217	ILE
1	B	219	VAL
1	B	239	GLN
1	B	248	ASP
1	B	269	ASP
1	B	278	PRO
1	B	284	GLU
1	B	288	LEU
1	B	302	MET
1	B	334	GLU
1	B	335	LYS
1	B	349	GLU
1	B	351	LYS

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Mol	Chain	Res	Type
1	B	361	ASP
1	B	386	GLU
1	B	397	MET
1	B	409	ARG
1	B	424	ASP
1	B	428	ILE
1	B	443	ASP
1	B	489	ARG
1	B	498	VAL
1	B	514	ILE
1	B	517	GLU
1	B	523	SER
1	B	554	GLU
1	B	573	SER
1	B	575	ASP
1	B	580	LYS
1	B	586	ASN
1	B	591	ASP
1	B	593	THR
1	B	594	ARG
1	B	599	VAL
1	B	639	GLU
1	B	657	GLU
1	B	665	GLU
1	B	681	GLU
1	B	687	LEU
1	B	692	ASN
1	B	711	LEU
1	B	723	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (33) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	32	ASN
1	A	38	GLN
1	A	141	GLN
1	A	239	GLN
1	A	285	GLN
1	A	286	GLN
1	A	291	GLN
1	A	342	GLN
1	A	358	HIS

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Mol	Chain	Res	Type
1	A	515	GLN
1	A	519	ASN
1	A	568	HIS
1	A	586	ASN
1	A	666	HIS
1	A	692	ASN
1	B	32	ASN
1	B	38	GLN
1	B	57	GLN
1	B	96	HIS
1	B	126	ASN
1	B	141	GLN
1	B	286	GLN
1	B	291	GLN
1	B	325	ASN
1	B	513	ASN
1	B	515	GLN
1	B	519	ASN
1	B	544	GLN
1	B	586	ASN
1	B	589	GLN
1	B	608	ASN
1	B	666	HIS
1	B	692	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](#)

2 ligands are modelled in this entry.

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	A	800	1	50,50,50	2.13	10 (20%)	67,82,82	1.64	15 (22%)
2	HEM	B	800	1	50,50,50	1.78	9 (18%)	67,82,82	1.40	8 (11%)

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	A	800	1	-	6/14/54/54	-
2	HEM	B	800	1	-	8/14/54/54	-

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	800	HEM	C3D-C2D	7.61	1.53	1.36
2	B	800	HEM	C3D-C2D	7.08	1.52	1.36
2	A	800	HEM	FE-NC	5.70	2.14	1.95
2	A	800	HEM	FE-NA	4.47	2.09	1.95
2	A	800	HEM	C4A-NA	-3.55	1.32	1.39
2	B	800	HEM	FE-ND	3.34	2.05	1.94
2	B	800	HEM	FE-NA	3.30	2.06	1.95
2	A	800	HEM	C4C-NC	-3.25	1.33	1.39
2	A	800	HEM	CAC-C3C	3.20	1.55	1.47
2	A	800	HEM	C2A-C3A	-3.03	1.31	1.38
2	B	800	HEM	FE-NC	2.90	2.04	1.95
2	A	800	HEM	CMB-C2B	2.59	1.56	1.50
2	A	800	HEM	CAB-C3B	2.57	1.54	1.47
2	B	800	HEM	CMA-C3A	2.53	1.56	1.50
2	B	800	HEM	CAC-C3C	2.51	1.54	1.47
2	B	800	HEM	CMC-C2C	2.41	1.55	1.50
2	B	800	HEM	CAB-C3B	2.37	1.53	1.47
2	A	800	HEM	C1D-ND	2.15	1.43	1.38
2	B	800	HEM	CHA-C1A	-2.06	1.34	1.39

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	800	HEM	CAA-CBA-CGA	-5.21	99.86	113.67
2	B	800	HEM	C4D-ND-C1D	4.35	110.36	105.21
2	A	800	HEM	C4D-ND-C1D	4.21	110.19	105.21
2	A	800	HEM	CMD-C2D-C1D	4.17	131.56	125.03
2	A	800	HEM	CAD-CBD-CGD	-3.82	103.53	113.67
2	A	800	HEM	CAA-CBA-CGA	-3.59	104.14	113.67
2	A	800	HEM	C4C-CHD-C1D	3.11	132.62	126.02
2	B	800	HEM	CHB-C1B-NB	3.03	128.12	124.37
2	A	800	HEM	CMC-C2C-C1C	-2.92	119.59	124.73
2	A	800	HEM	C4A-C3A-C2A	2.86	110.10	106.82
2	B	800	HEM	CBA-CAA-C2A	2.73	120.09	112.53
2	B	800	HEM	CHA-C4D-ND	2.60	127.58	124.37
2	A	800	HEM	CHB-C1B-NB	2.53	127.50	124.37
2	A	800	HEM	C1C-CHC-C4B	-2.49	120.73	126.02
2	A	800	HEM	CBC-CAC-C3C	-2.40	115.51	127.53
2	B	800	HEM	C3B-C2B-C1B	2.40	108.21	106.41
2	A	800	HEM	C1D-C2D-C3D	-2.36	104.50	106.98
2	A	800	HEM	CMC-C2C-C3C	2.32	134.03	128.43
2	A	800	HEM	C3A-C4A-NA	-2.30	106.46	110.14
2	B	800	HEM	CAD-CBD-CGD	-2.28	107.61	113.67
2	A	800	HEM	CAA-C2A-C1A	2.27	129.38	124.94
2	B	800	HEM	CMD-C2D-C1D	2.25	128.54	125.03
2	A	800	HEM	C3B-C4B-NB	2.10	110.98	109.47

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	800	HEM	C1A-C2A-CAA-CBA
2	B	800	HEM	C2C-C3C-CAC-CBC
2	B	800	HEM	C3A-C2A-CAA-CBA
2	A	800	HEM	CAA-CBA-CGA-O2A
2	A	800	HEM	C4B-C3B-CAB-CBB
2	B	800	HEM	C4C-C3C-CAC-CBC
2	A	800	HEM	CAA-CBA-CGA-O1A
2	B	800	HEM	CAA-CBA-CGA-O2A
2	B	800	HEM	CAD-CBD-CGD-O1D
2	B	800	HEM	CAD-CBD-CGD-O2D
2	B	800	HEM	CAA-CBA-CGA-O1A
2	A	800	HEM	C2D-C3D-CAD-CBD
2	A	800	HEM	CAD-CBD-CGD-O2D

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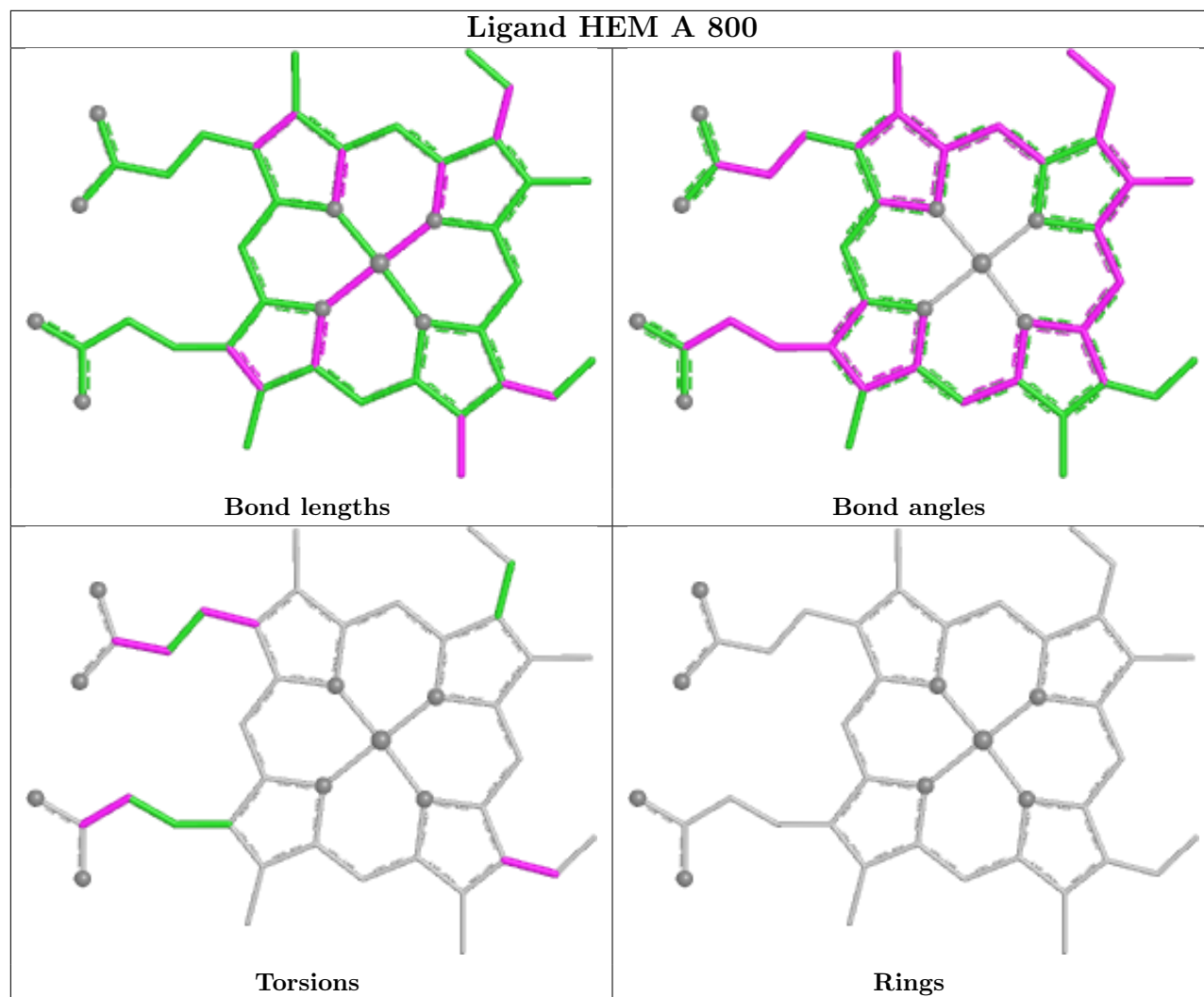
Mol	Chain	Res	Type	Atoms
2	A	800	HEM	C2B-C3B-CAB-CBB

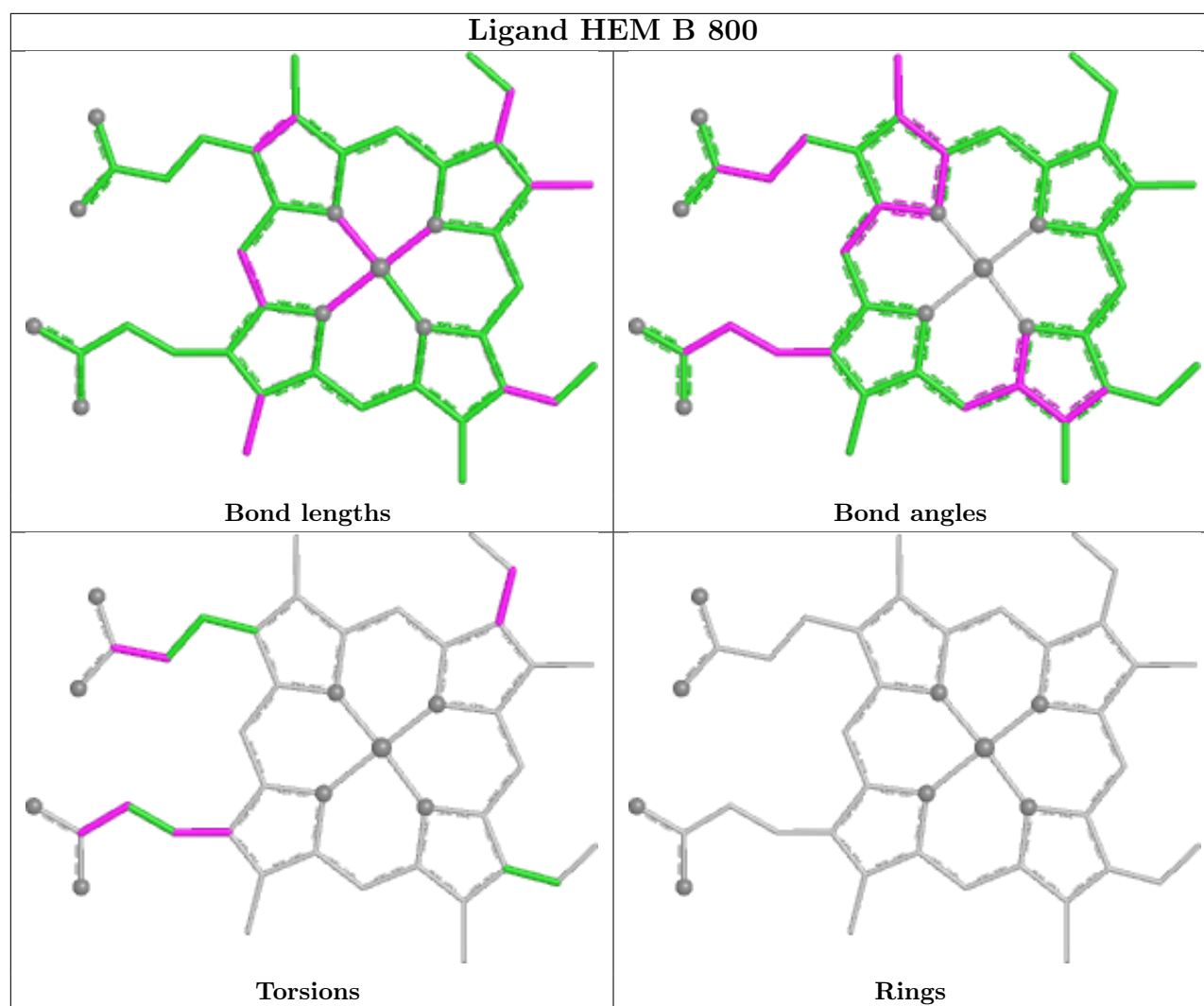
There are no ring outliers.

2 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	800	HEM	4	0
2	B	800	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

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	663/737 (89%)	0.62	34 (5%) 33 39	14, 35, 55, 68	0
1	B	660/737 (89%)	0.86	65 (9%) 13 16	24, 43, 60, 71	0
All	All	1323/1474 (89%)	0.74	99 (7%) 20 24	14, 39, 58, 71	0

All (99) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	360	PRO	8.7
1	A	361	ASP	7.3
1	A	245	ALA	5.0
1	B	215	GLY	4.4
1	B	29	SER	4.3
1	B	728	PHE	4.0
1	B	397	MET	3.8
1	A	218	TYR	3.7
1	B	476	ASP	3.5
1	A	728	PHE	3.5
1	A	176	GLU	3.4
1	B	663	ASP	3.4
1	A	212	SER	3.4
1	B	361	ASP	3.4
1	A	216	LEU	3.2
1	B	523	SER	3.2
1	A	215	GLY	3.2
1	A	41	ARG	3.2
1	B	77	ASP	3.1
1	B	311	TRP	3.1
1	B	665	GLU	3.0
1	A	213	VAL	2.9
1	B	727	ARG	2.9
1	A	362	GLU	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	186	ASN	2.9
1	B	662	ALA	2.9
1	A	706	ASP	2.9
1	B	233	SER	2.8
1	B	592	ILE	2.8
1	A	48	ASP	2.8
1	A	525	GLY	2.8
1	B	664	SER	2.8
1	B	187	TRP	2.8
1	B	234	ALA	2.7
1	B	239	GLN	2.7
1	A	186	ASN	2.7
1	B	341	TRP	2.7
1	B	667	ARG	2.7
1	B	360	PRO	2.7
1	B	379	PRO	2.6
1	A	187	TRP	2.6
1	B	331	TRP	2.6
1	A	38	GLN	2.5
1	B	366	PRO	2.5
1	A	214	MET	2.5
1	B	244	ALA	2.5
1	B	264	VAL	2.5
1	A	437	ASP	2.5
1	A	632	ASP	2.5
1	B	348	GLU	2.4
1	B	236	ASN	2.4
1	B	232	ALA	2.4
1	A	217	ILE	2.4
1	B	242	ASP	2.4
1	B	675	THR	2.4
1	B	591	ASP	2.4
1	B	632	ASP	2.4
1	B	47	GLU	2.4
1	B	245	ALA	2.3
1	B	558	GLU	2.3
1	B	437	ASP	2.3
1	A	503	GLN	2.3
1	B	272	GLU	2.3
1	B	364	GLN	2.3
1	A	30	LYS	2.3
1	A	524	ASP	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	527	GLN	2.3
1	B	350	LEU	2.2
1	B	455	ASP	2.2
1	B	374	ALA	2.2
1	A	521	SER	2.2
1	B	273	ASN	2.2
1	A	556	PRO	2.2
1	A	667	ARG	2.2
1	B	241	PHE	2.2
1	A	359	ASP	2.2
1	B	608	ASN	2.2
1	A	34	GLU	2.2
1	B	41	ARG	2.2
1	B	659	GLU	2.2
1	B	549	ALA	2.2
1	B	343	TRP	2.1
1	B	43	VAL	2.1
1	A	271	GLU	2.1
1	B	666	HIS	2.1
1	B	218	TYR	2.1
1	B	508	LEU	2.1
1	B	329	TYR	2.1
1	A	123	TRP	2.1
1	B	677	GLU	2.1
1	B	327	LEU	2.1
1	A	591	ASP	2.1
1	B	347	SER	2.0
1	B	279	GLU	2.0
1	B	514	ILE	2.0
1	B	46	VAL	2.0
1	B	676	GLY	2.0
1	B	391	ASN	2.0
1	B	235	LYS	2.0

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

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

6.3 Carbohydrates ⓘ

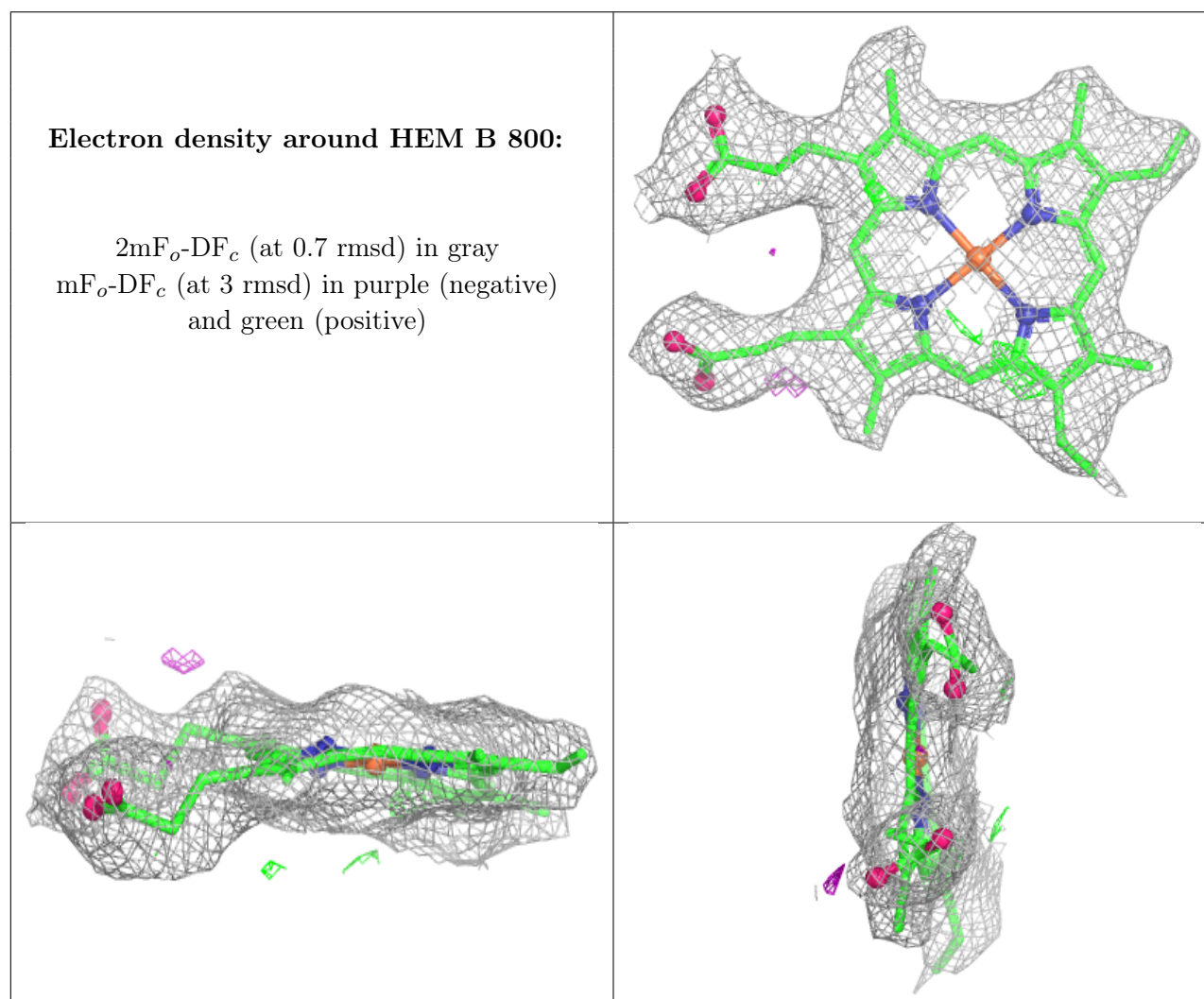
There are no oligosaccharides in this entry.

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.

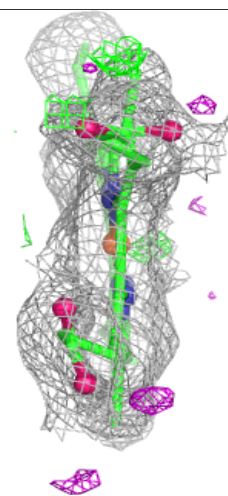
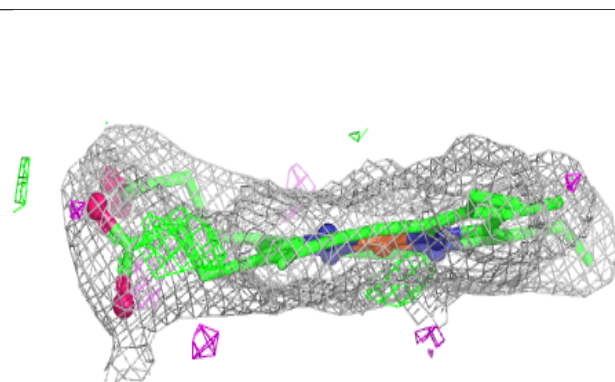
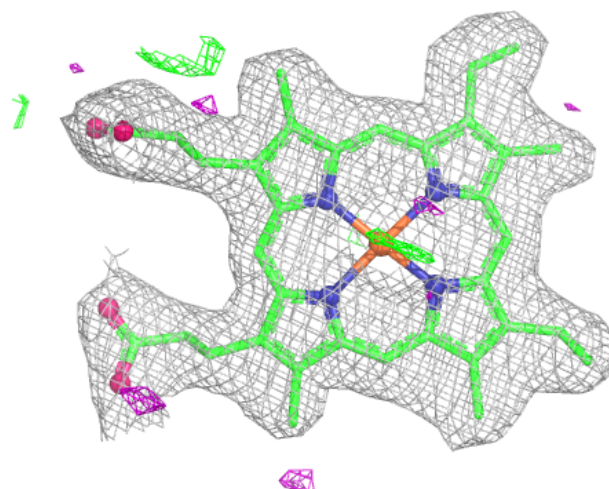
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	HEM	B	800	43/43	0.95	0.11	39,42,46,48	0
2	HEM	A	800	43/43	0.96	0.08	17,22,25,29	0

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 800:

$2mF_o-DF_c$ (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.