



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 4, 2026 – 08:26 PM EDT

PDB ID : 3WNU / pdb_00003wnu
Title : The crystal structure of catalase-peroxidase, KatG, from *Synechococcus* PCC7942
Authors : Tada, T.; Wada, K.; Kamachi, S.
Deposited on : 2013-12-17
Resolution : 2.20 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	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

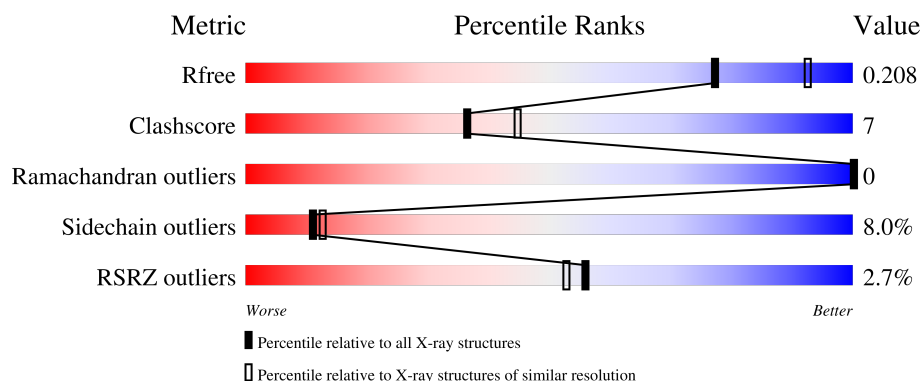
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	720	

2 Entry composition [i](#)

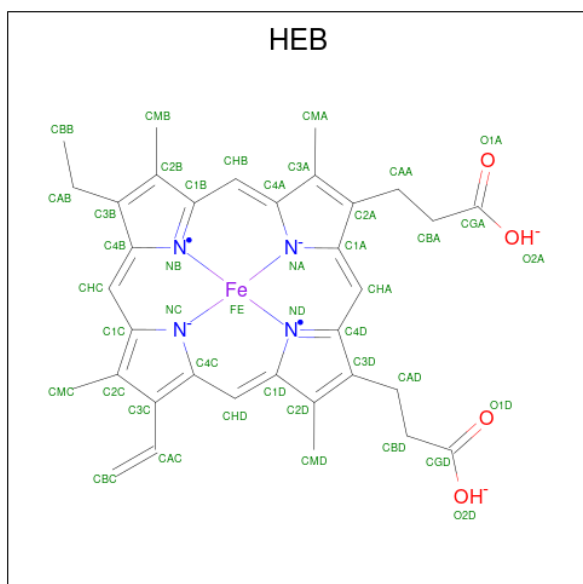
There are 4 unique types of molecules in this entry. The entry contains 6149 atoms, of which 0 are hydrogens and 0 are deuteriums.

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.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	710	Total	C	N	O	S	0	0	0
			5582	3523	980	1060	19			

- Molecule 2 is HEME B/C (CCD ID: HEB) (formula: $C_{34}H_{34}FeN_4O_4$).



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

- Molecule 3 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Na	0	0
			2	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	522	Total 522	O 522	0	0

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Catalase-peroxidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	108.74Å 108.74Å 202.89Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	34.39 – 2.20 34.39 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.8 (34.39-2.20) 99.9 (34.39-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.48 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.168 , 0.206 0.173 , 0.208	Depositor DCC
R_{free} test set	3159 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	34.3	Xtriage
Anisotropy	0.016	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 35.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6149	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.00% 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: HEB, NA

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	1.33	18/5729 (0.3%)	1.23	34/7803 (0.4%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	327	ARG	NE-CZ	11.43	1.45	1.33
1	A	350	ASN	C-O	-9.48	1.19	1.23
1	A	694	ASP	C-O	-7.13	1.15	1.24
1	A	327	ARG	CD-NE	6.99	1.56	1.46
1	A	161	GLU	CA-C	-6.65	1.44	1.52
1	A	179	HIS	N-CA	-6.30	1.40	1.46
1	A	350	ASN	CA-C	-6.12	1.49	1.53
1	A	130	LYS	N-CA	5.99	1.53	1.46
1	A	327	ARG	CZ-NH2	5.84	1.41	1.33
1	A	362	ASP	C-O	-5.66	1.20	1.25
1	A	176	ASP	C-O	-5.49	1.16	1.23
1	A	16	THR	CA-C	5.48	1.60	1.52
1	A	576	ARG	CD-NE	-5.34	1.38	1.46
1	A	251	ASN	CA-C	5.24	1.60	1.53
1	A	321	TRP	C-O	-5.22	1.17	1.23
1	A	685	ARG	C-O	-5.19	1.18	1.24
1	A	21	SER	CA-C	5.16	1.58	1.52
1	A	37	GLN	CD-NE2	-5.08	1.22	1.33

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	327	ARG	NE-CZ-NH2	11.21	129.29	119.20
1	A	327	ARG	CG-CD-NE	9.77	133.50	112.00
1	A	643	THR	N-CA-CB	-8.98	97.20	110.49
1	A	99	THR	N-CA-CB	-8.59	96.97	110.46
1	A	576	ARG	NE-CZ-NH2	-8.27	111.76	119.20
1	A	370	VAL	CB-CA-C	-8.21	95.71	110.48
1	A	327	ARG	NH1-CZ-NH2	-8.16	108.69	119.30
1	A	628	THR	CB-CA-C	-8.14	92.59	109.94
1	A	353	GLU	CB-CA-C	-8.12	97.82	110.81
1	A	327	ARG	N-CA-C	-7.35	102.29	111.11
1	A	242	ARG	NE-CZ-NH1	6.74	128.24	121.50
1	A	242	ARG	NE-CZ-NH2	-6.56	113.30	119.20
1	A	21	SER	N-CA-C	6.45	118.01	108.60
1	A	643	THR	OG1-CB-CG2	6.08	121.46	109.30
1	A	211	ASN	CA-C-N	-6.07	112.76	119.19
1	A	211	ASN	C-N-CA	-6.07	112.76	119.19
1	A	414	MET	CG-SD-CE	5.97	114.04	100.90
1	A	99	THR	CB-CA-C	5.77	120.80	109.72
1	A	161	GLU	N-CA-C	-5.69	104.98	111.07
1	A	21	SER	CA-C-N	5.61	128.35	120.28
1	A	21	SER	C-N-CA	5.61	128.35	120.28
1	A	320	GLN	N-CA-C	5.53	118.46	109.06
1	A	16	THR	N-CA-CB	-5.43	102.44	110.53
1	A	567	VAL	CG1-CB-CG2	5.42	122.73	110.80
1	A	467	TRP	N-CA-C	-5.33	105.36	111.07
1	A	242	ARG	CD-NE-CZ	5.33	131.86	124.40
1	A	643	THR	CB-CA-C	5.24	118.84	110.09
1	A	455	SER	N-CA-C	-5.20	104.02	110.41
1	A	637	ASP	N-CA-C	5.19	117.61	111.33
1	A	294	ILE	CB-CA-C	-5.15	104.03	111.34
1	A	576	ARG	CD-NE-CZ	5.15	131.60	124.40
1	A	327	ARG	CB-CG-CD	5.08	122.99	111.30
1	A	622	THR	N-CA-CB	-5.05	101.95	110.49
1	A	99	THR	OG1-CB-CG2	5.02	119.34	109.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	210	GLU	Peptide

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	5582	0	5365	78	0
2	A	43	0	32	1	0
3	A	2	0	0	0	0
4	A	522	0	0	11	1
All	All	6149	0	5397	79	1

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 7.

All (79) 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:94:TRP:CH2	1:A:222:TYR:HE2	1.00	1.65
1:A:94:TRP:HH2	1:A:222:TYR:CE2	0.93	1.58
1:A:94:TRP:CH2	1:A:222:TYR:CE2	1.83	1.31
1:A:94:TRP:HH2	1:A:222:TYR:CD2	1.83	0.95
1:A:99:THR:HG22	1:A:108:GLY:H	1.36	0.88
1:A:94:TRP:CZ3	1:A:222:TYR:HE2	1.88	0.88
1:A:591:LEU:O	1:A:595:THR:HG23	1.77	0.84
1:A:94:TRP:CH2	1:A:222:TYR:CD2	2.64	0.81
1:A:106:ARG:HH21	1:A:596:GLN:HE21	1.32	0.78
1:A:114:GLN:HE21	1:A:132:ARG:HH11	1.28	0.76
1:A:94:TRP:CZ3	1:A:222:TYR:CE2	2.70	0.73
1:A:99:THR:CG2	1:A:108:GLY:H	2.02	0.71
1:A:251:ASN:HD22	1:A:251:ASN:C	2.00	0.69
1:A:38:HIS:H	1:A:179:HIS:HE1	1.42	0.67
1:A:179:HIS:HD2	1:A:180:PRO:O	1.78	0.66
1:A:198:ASN:OD1	1:A:199:PRO:O	2.14	0.66
1:A:171:ALA:H	1:A:411:HIS:HE1	1.42	0.65
1:A:628:THR:HB	1:A:630:ARG:H	1.66	0.60
1:A:492:GLN:HG2	1:A:495:TRP:CZ3	2.36	0.60
1:A:251:ASN:ND2	1:A:254:GLU:H	2.00	0.59
1:A:639:PHE:O	1:A:643:THR:HB	2.02	0.59
1:A:114:GLN:HE21	1:A:132:ARG:NH1	2.00	0.58
1:A:224:ASN:C	1:A:224:ASN:HD22	2.12	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:242:ARG:HD3	1:A:252:ASP:OD1	2.03	0.57
1:A:106:ARG:HH21	1:A:596:GLN:NE2	2.00	0.56
1:A:648:LEU:C	1:A:648:LEU:HD12	2.31	0.56
1:A:502:ARG:HD3	4:A:1327:HOH:O	2.06	0.55
1:A:628:THR:HG21	4:A:1254:HOH:O	2.06	0.55
1:A:99:THR:HG21	4:A:1053:HOH:O	2.05	0.55
1:A:285:ASP:CB	4:A:1344:HOH:O	2.54	0.54
2:A:801:HEB:HHC	2:A:801:HEB:HBB3	1.90	0.54
1:A:595:THR:HG21	1:A:605:MET:SD	2.47	0.54
1:A:94:TRP:HZ3	1:A:222:TYR:HH	1.53	0.54
1:A:440:ASN:ND2	1:A:442:ASN:H	2.06	0.53
1:A:38:HIS:H	1:A:179:HIS:CE1	2.25	0.53
1:A:445:VAL:HG13	1:A:527:LEU:HD21	1.91	0.53
1:A:591:LEU:O	1:A:595:THR:CG2	2.55	0.53
1:A:171:ALA:H	1:A:411:HIS:CE1	2.26	0.52
1:A:366:ARG:NH1	4:A:1199:HOH:O	2.43	0.51
1:A:310:LEU:HD22	1:A:343:ALA:HB1	1.93	0.50
1:A:561:ASP:C	1:A:561:ASP:OD1	2.54	0.50
1:A:224:ASN:HD22	1:A:226:GLU:H	1.58	0.49
1:A:628:THR:CG2	1:A:637:ASP:OD2	2.60	0.49
1:A:618:ASN:OD1	1:A:622:THR:HG22	2.12	0.49
1:A:113:ASN:HD22	1:A:114:GLN:N	2.11	0.48
1:A:120:ASN:HD22	1:A:121:SER:N	2.12	0.47
1:A:385:ARG:NH2	1:A:389:GLU:OE1	2.38	0.47
1:A:507:LEU:O	1:A:511:GLU:HB2	2.15	0.47
1:A:312:GLY:C	1:A:370:VAL:HG22	2.39	0.46
1:A:224:ASN:ND2	1:A:226:GLU:H	2.14	0.46
1:A:618:ASN:ND2	4:A:1046:HOH:O	2.48	0.46
1:A:655:ASN:ND2	1:A:673:ARG:H	2.14	0.46
1:A:217:THR:HG22	1:A:220:LEU:HD12	1.97	0.45
1:A:611:GLY:HA3	1:A:704:PHE:CE1	2.51	0.45
1:A:514:ALA:HB1	1:A:519:ALA:O	2.17	0.45
1:A:68:GLN:HE21	1:A:68:GLN:HB2	1.62	0.45
1:A:48:ASP:N	4:A:1421:HOH:O	2.16	0.44
1:A:179:HIS:HB2	1:A:180:PRO:HD2	1.99	0.44
1:A:238:ALA:HB2	1:A:377:ALA:HB1	2.00	0.44
1:A:492:GLN:HG2	1:A:495:TRP:CH2	2.52	0.44
1:A:50:ASN:OD1	1:A:50:ASN:C	2.61	0.44
1:A:537:ALA:O	1:A:538:ARG:C	2.61	0.43
1:A:127:ASN:HA	1:A:129:ASP:OD1	2.18	0.43
1:A:94:TRP:CZ2	1:A:222:TYR:CD2	3.07	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:25:TRP:HB2	1:A:26:TRP:CE3	2.53	0.43
1:A:317:HIS:HE1	4:A:998:HOH:O	2.01	0.43
1:A:285:ASP:HB3	4:A:1344:HOH:O	2.16	0.42
1:A:412:ARG:NH1	4:A:1387:HOH:O	2.44	0.42
1:A:92:LEU:HD23	1:A:114:GLN:HE22	1.84	0.42
1:A:622:THR:HG23	1:A:624:HIS:H	1.84	0.42
1:A:628:THR:HG22	1:A:637:ASP:OD2	2.20	0.41
1:A:602:ALA:HB3	1:A:603:PRO:HD3	2.01	0.41
1:A:439:GLY:HA3	1:A:552:ARG:O	2.20	0.41
1:A:502:ARG:CD	4:A:1327:HOH:O	2.67	0.41
1:A:488:ARG:O	1:A:493:LYS:HD3	2.21	0.41
1:A:628:THR:HG23	1:A:637:ASP:OD2	2.21	0.41
1:A:37:GLN:NE2	1:A:182:LYS:H	2.18	0.41
1:A:251:ASN:HD21	1:A:254:GLU:H	1.68	0.40
1:A:224:ASN:HD22	1:A:225:PRO:N	2.19	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:1332:HOH:O	4:A:1334:HOH:O[7_554]	2.17	0.03

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	708/720 (98%)	687 (97%)	21 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	573/581 (99%)	527 (92%)	46 (8%)	11	13

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	MET
1	A	16	THR
1	A	22	THR
1	A	24	GLU
1	A	28	LYS
1	A	32	LEU
1	A	53	GLU
1	A	65	GLN
1	A	67	LEU
1	A	68	GLN
1	A	92	LEU
1	A	99	THR
1	A	113	ASN
1	A	120	ASN
1	A	124	ASP
1	A	165	LEU
1	A	204	THR
1	A	206	ASP
1	A	207	ARG
1	A	224	ASN
1	A	251	ASN
1	A	298	GLN
1	A	310	LEU
1	A	327	ARG
1	A	370	VAL
1	A	373	ASP
1	A	427	GLN
1	A	430	LEU
1	A	431	ILE
1	A	442	ASN

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Mol	Chain	Res	Type
1	A	511	GLU
1	A	527	LEU
1	A	544	ILE
1	A	562	THR
1	A	563	GLU
1	A	567	VAL
1	A	582	ASP
1	A	595	THR
1	A	608	LEU
1	A	622	THR
1	A	628	THR
1	A	643	THR
1	A	656	LEU
1	A	663	LYS
1	A	719	LEU
1	A	720	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	31	ASN
1	A	37	GLN
1	A	68	GLN
1	A	113	ASN
1	A	114	GLN
1	A	120	ASN
1	A	125	ASN
1	A	179	HIS
1	A	198	ASN
1	A	224	ASN
1	A	251	ASN
1	A	304	ASN
1	A	317	HIS
1	A	350	ASN
1	A	411	HIS
1	A	440	ASN
1	A	572	HIS
1	A	596	GLN
1	A	619	HIS
1	A	655	ASN
1	A	666	GLN
1	A	693	GLN

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 3 ligands modelled in this entry, 2 are monoatomic - leaving 1 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	HEB	A	801	1	50,50,50	1.72	13 (26%)	66,82,82	2.17	23 (34%)

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	HEB	A	801	1	-	4/14/54/54	-

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	801	HEB	C4B-NB	-5.00	1.31	1.40
2	A	801	HEB	CBB-CAB	-4.43	1.32	1.51
2	A	801	HEB	C1D-ND	-3.44	1.34	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	801	HEB	C1D-C2D	3.11	1.50	1.44
2	A	801	HEB	FE-NC	2.74	2.04	1.95
2	A	801	HEB	CAB-C3B	-2.56	1.44	1.51
2	A	801	HEB	C4C-NC	-2.35	1.35	1.39
2	A	801	HEB	CMA-C3A	-2.29	1.46	1.50
2	A	801	HEB	FE-NB	2.27	2.01	1.94
2	A	801	HEB	CMB-C2B	2.18	1.55	1.50
2	A	801	HEB	C1B-NB	-2.09	1.34	1.38
2	A	801	HEB	C3C-C2C	-2.08	1.34	1.41
2	A	801	HEB	C1C-NC	-2.02	1.35	1.39

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	801	HEB	CAB-C3B-C4B	5.66	132.15	124.79
2	A	801	HEB	C3D-C4D-ND	5.41	115.58	110.35
2	A	801	HEB	CHA-C4D-C3D	-4.95	117.56	124.77
2	A	801	HEB	CAB-C3B-C2B	-4.21	119.81	127.56
2	A	801	HEB	CHC-C4B-NB	4.10	129.44	124.37
2	A	801	HEB	CHB-C1B-NB	3.93	128.65	124.42
2	A	801	HEB	CMB-C2B-C1B	3.91	131.14	125.03
2	A	801	HEB	CBD-CAD-C3D	3.61	122.53	112.53
2	A	801	HEB	CBB-CAB-C3B	3.53	121.98	112.42
2	A	801	HEB	C4B-NB-C1B	3.47	109.32	105.21
2	A	801	HEB	C1A-CHA-C4D	-3.44	118.72	126.02
2	A	801	HEB	O1D-CGD-CBD	-2.95	113.73	123.09
2	A	801	HEB	C1D-ND-C4D	-2.87	101.81	105.21
2	A	801	HEB	C3B-C4B-NB	-2.85	107.05	110.17
2	A	801	HEB	C1C-CHC-C4B	-2.63	120.05	126.25
2	A	801	HEB	C4D-C3D-C2D	-2.51	103.23	106.89
2	A	801	HEB	O2D-CGD-CBD	2.47	121.82	114.00
2	A	801	HEB	CHB-C1B-C2B	-2.43	121.19	125.03
2	A	801	HEB	CAD-C3D-C4D	2.41	128.91	124.70
2	A	801	HEB	CAA-CBA-CGA	-2.37	107.37	113.67
2	A	801	HEB	CMC-C2C-C1C	-2.20	122.07	125.42
2	A	801	HEB	CMD-C2D-C1D	2.19	128.45	125.03
2	A	801	HEB	O2A-CGA-CBA	2.08	120.56	114.00

There are no chirality outliers.

All (4) torsion outliers are listed below:

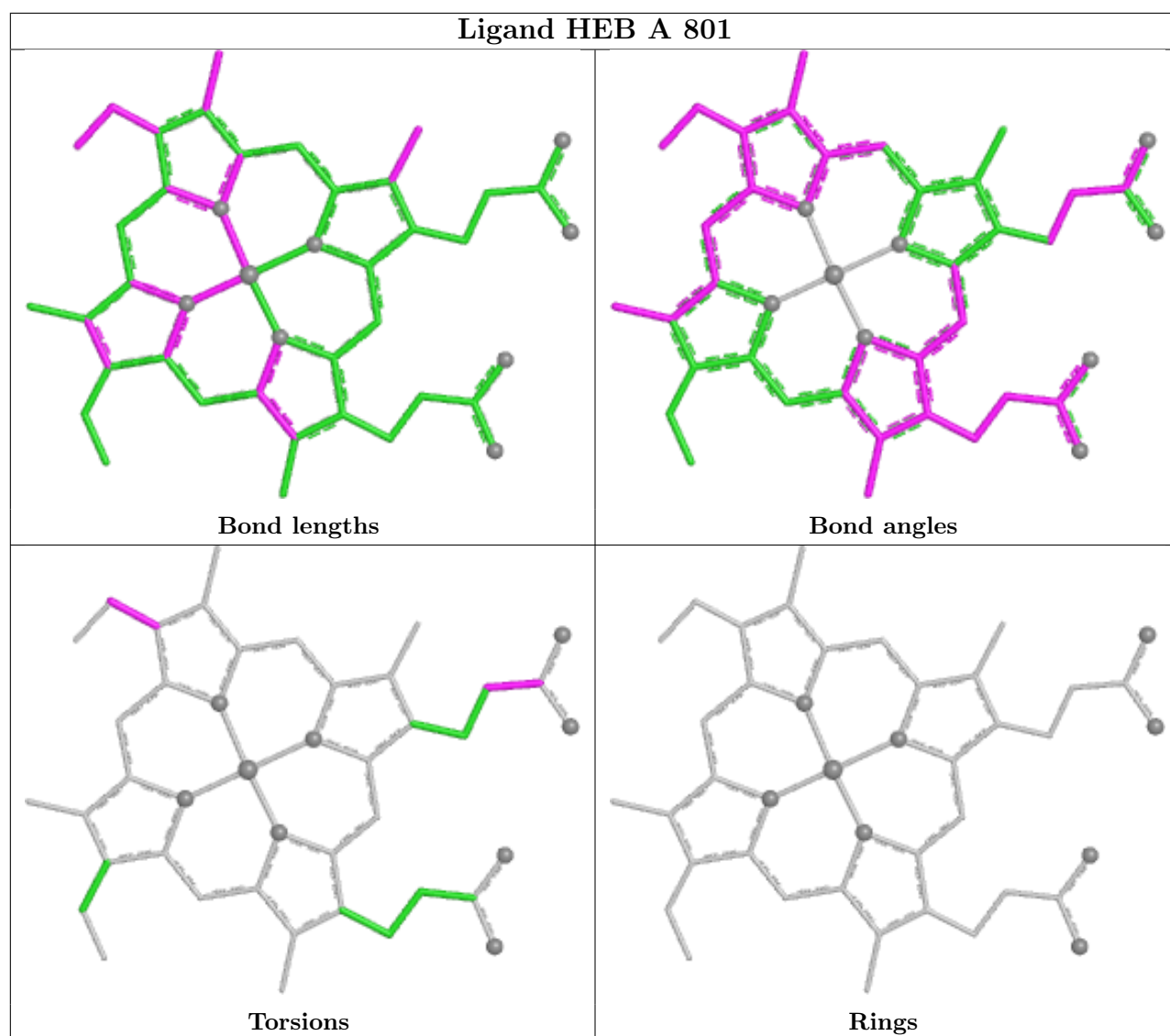
Mol	Chain	Res	Type	Atoms
2	A	801	HEB	C4B-C3B-CAB-CBB
2	A	801	HEB	C2B-C3B-CAB-CBB
2	A	801	HEB	CAA-CBA-CGA-O2A
2	A	801	HEB	CAA-CBA-CGA-O1A

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	801	HEB	1	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	710/720 (98%)	-0.26	19 (2%) 56 53	23, 33, 57, 97	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	11	MET	6.4
1	A	20	ILE	4.1
1	A	199	PRO	3.6
1	A	22	THR	3.5
1	A	720	ASP	3.5
1	A	211	ASN	3.5
1	A	327	ARG	3.1
1	A	12	HIS	3.0
1	A	23	ALA	3.0
1	A	18	VAL	3.0
1	A	582	ASP	3.0
1	A	197	THR	2.8
1	A	542	VAL	2.8
1	A	663	LYS	2.7
1	A	24	GLU	2.6
1	A	14	GLY	2.6
1	A	544	ILE	2.6
1	A	19	ASN	2.6
1	A	15	ALA	2.2

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](#)

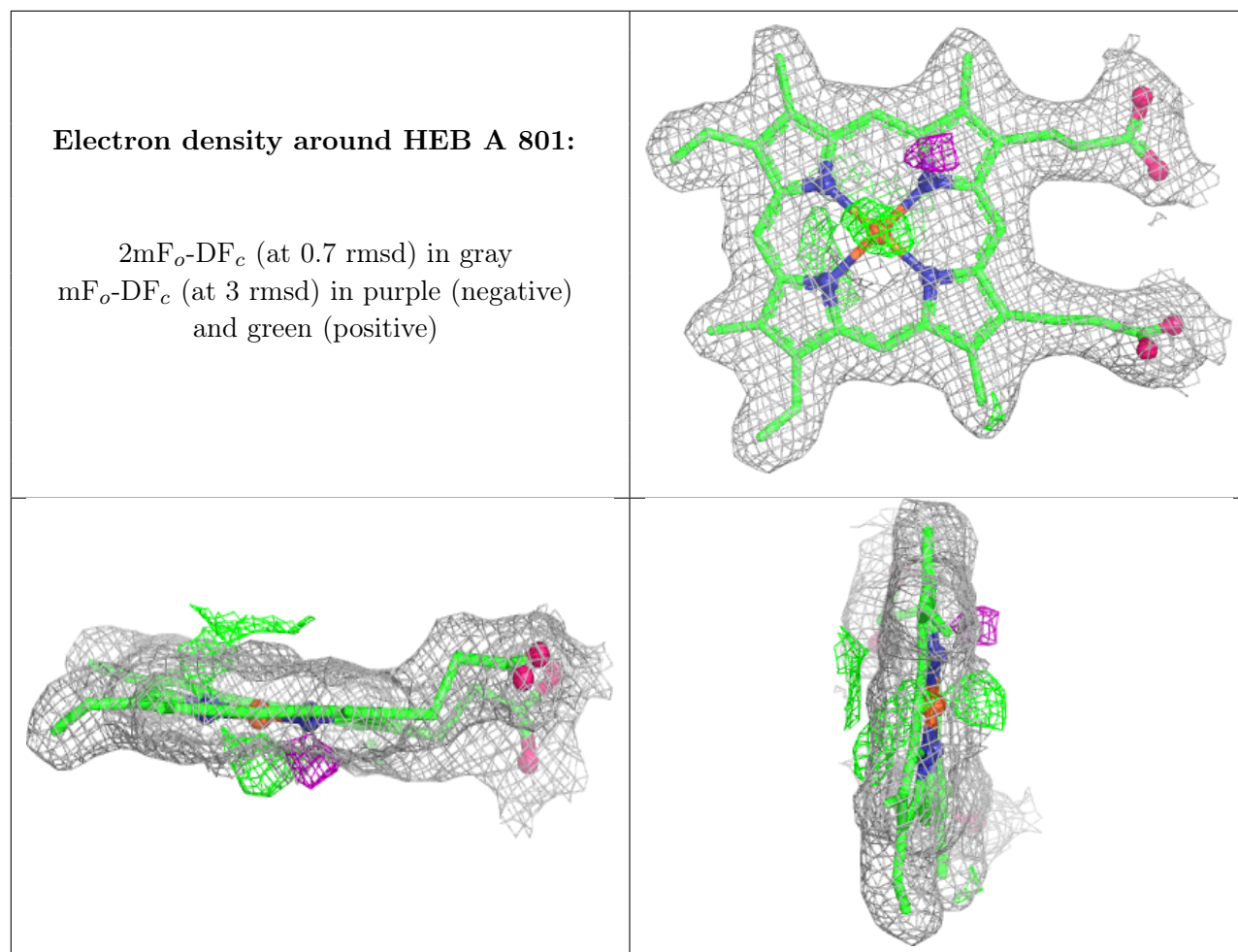
There are no oligosaccharides in this entry.

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($\text{\AA}^2$)	Q<0.9
2	HEB	A	801	43/43	0.98	0.06	22,26,29,33	0
3	NA	A	802	1/1	0.99	0.04	27,27,27,27	0
3	NA	A	803	1/1	0.99	0.06	27,27,27,27	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.



6.5 Other polymers [i](#)

There are no such residues in this entry.