



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 5, 2023 – 10:06 AM EST

PDB ID : 5L02
Title : S324T variant of B. pseudomallei KatG
Authors : Loewen, P.C.
Deposited on : 2016-07-26
Resolution : 1.90 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.36
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36

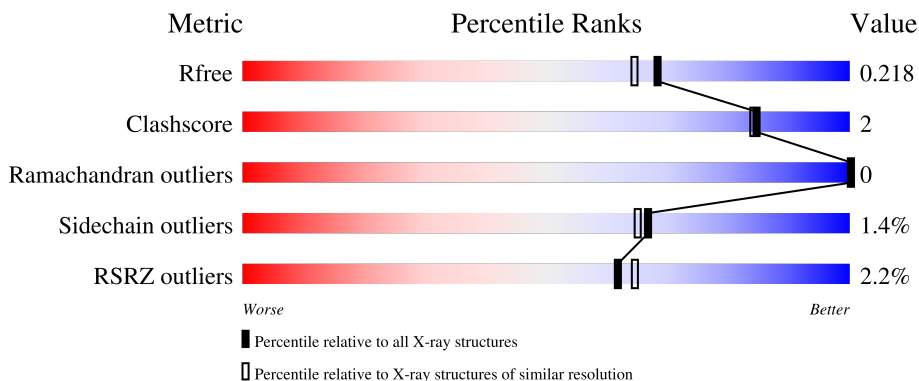
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

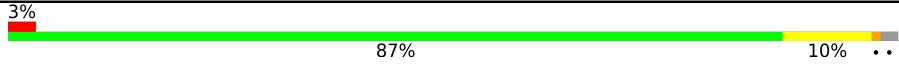
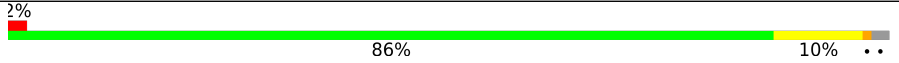
The reported resolution of this entry is 1.90 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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	728	 3% 87% 10% ..
1	B	728	 2% 86% 10% ..

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 12604 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
			Total	C	N	O	S			
1	A	713	Total 5585	C 3521	N 998	O 1052	S 14	0	9	0
1	B	713	Total 5582	C 3519	N 999	O 1050	S 14	0	9	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	324	THR	SER	engineered mutation	UNP Q3JNW6
B	324	THR	SER	engineered mutation	UNP Q3JNW6

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



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

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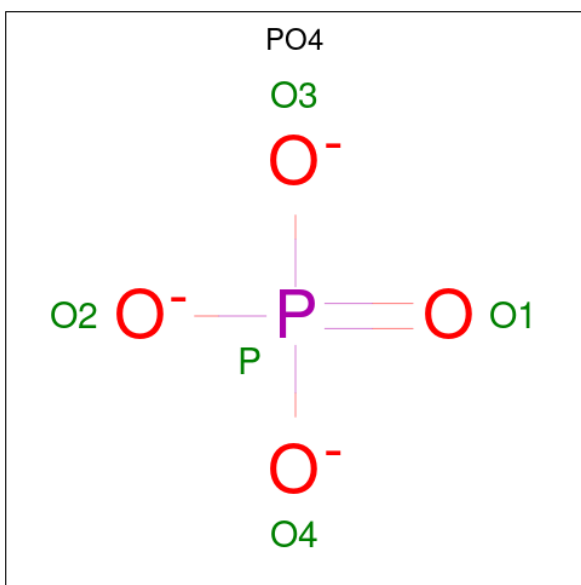
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
			Total	C	Fe	N			O
2	B	1	43	34	1	4	4	0	0

- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Na 1 1	0	0
3	B	1	Total Na 1 1	0	0

- Molecule 4 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



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

- Molecule 5 is (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula: C₆H₁₄O₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total C O 8 6 2	0	0
5	A	1	Total C O 8 6 2	0	0
5	B	1	Total C O 8 6 2	0	0

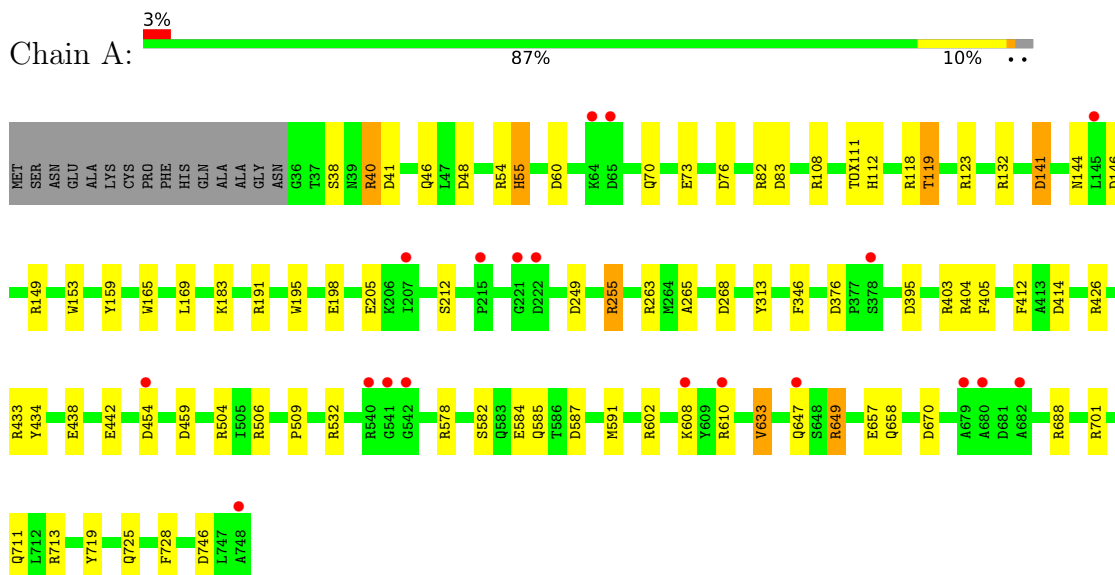
- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	660	Total O 660 660	0	0
6	B	655	Total O 655 655	0	0

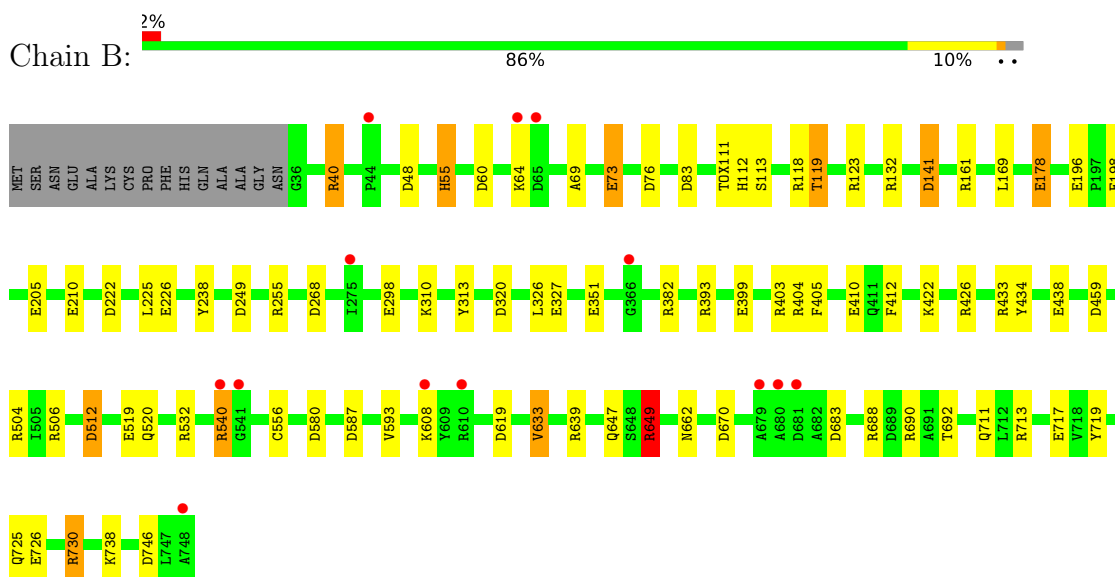
3 Residue-property plots

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



- Molecule 1: Catalase-peroxidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	100.50Å 114.63Å 174.93Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 1.90 29.05 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.1 (30.00-1.90) 99.1 (29.05-1.90)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.92 (at 1.91Å)	Xtrriage
Refinement program	REFMAC 5.8.0155	Depositor
R, R_{free}	0.173 , 0.211 0.184 , 0.218	Depositor DCC
R_{free} test set	7879 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	23.5	Xtrriage
Anisotropy	0.259	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 46.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	12604	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.09% 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: TOX, NA, HEM, MPD, PO4

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	26/5711 (0.5%)	1.26	61/7760 (0.8%)
1	B	1.34	23/5704 (0.4%)	1.24	53/7749 (0.7%)
All	All	1.34	49/11415 (0.4%)	1.25	114/15509 (0.7%)

All (49) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	726	GLU	CG-CD	10.19	1.67	1.51
1	B	512	ASP	CG-OD2	8.11	1.44	1.25
1	B	532	ARG	CD-NE	-8.01	1.32	1.46
1	B	196	GLU	CG-CD	7.84	1.63	1.51
1	B	746	ASP	CB-CG	7.50	1.67	1.51
1	A	584	GLU	CG-CD	7.43	1.63	1.51
1	A	532	ARG	CD-NE	-7.17	1.34	1.46
1	A	728	PHE	CG-CD2	-6.56	1.28	1.38
1	B	438	GLU	CD-OE2	6.48	1.32	1.25
1	B	556	CYS	CB-SG	6.44	1.93	1.82
1	B	399	GLU	CD-OE1	6.33	1.32	1.25
1	A	73	GLU	CG-CD	6.21	1.61	1.51
1	B	327	GLU	CG-CD	6.17	1.61	1.51
1	A	153	TRP	CG-CD1	-6.15	1.28	1.36
1	A	532	ARG	NE-CZ	-6.08	1.25	1.33
1	A	159	TYR	CE1-CZ	6.04	1.46	1.38
1	B	119[A]	THR	CB-CG2	-5.90	1.32	1.52
1	B	119[B]	THR	CB-CG2	-5.90	1.32	1.52
1	A	746	ASP	CB-CG	5.83	1.64	1.51
1	A	313	TYR	CE2-CZ	5.79	1.46	1.38
1	A	195	TRP	CE3-CZ3	5.78	1.48	1.38
1	B	161	ARG	NE-CZ	5.74	1.40	1.33
1	B	351	GLU	CD-OE1	5.69	1.31	1.25
1	B	178	GLU	CG-CD	5.67	1.60	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	442	GLU	CD-OE2	-5.65	1.19	1.25
1	B	210	GLU	CD-OE2	5.62	1.31	1.25
1	B	746	ASP	C-O	-5.51	1.12	1.23
1	A	82	ARG	CZ-NH2	-5.49	1.25	1.33
1	A	119[A]	THR	CB-CG2	-5.47	1.34	1.52
1	A	119[B]	THR	CB-CG2	-5.47	1.34	1.52
1	A	657[A]	GLU	CD-OE2	-5.41	1.19	1.25
1	A	657[B]	GLU	CD-OE2	-5.41	1.19	1.25
1	B	717	GLU	CD-OE1	5.39	1.31	1.25
1	B	506	ARG	CZ-NH2	-5.36	1.26	1.33
1	A	438	GLU	CD-OE2	5.36	1.31	1.25
1	A	602	ARG	CZ-NH2	-5.33	1.26	1.33
1	A	212	SER	CB-OG	-5.31	1.35	1.42
1	B	113	SER	CB-OG	5.31	1.49	1.42
1	B	313	TYR	CG-CD1	5.28	1.46	1.39
1	B	205	GLU	CD-OE1	5.27	1.31	1.25
1	A	38	SER	CA-CB	5.23	1.60	1.52
1	B	327	GLU	CD-OE1	5.18	1.31	1.25
1	A	658	GLN	CD-NE2	-5.17	1.20	1.32
1	A	313	TYR	CG-CD1	5.14	1.45	1.39
1	A	688	ARG	CZ-NH2	-5.07	1.26	1.33
1	A	434	TYR	CE1-CZ	-5.05	1.31	1.38
1	B	226	GLU	CG-CD	5.05	1.59	1.51
1	A	165	TRP	CE2-CZ2	5.03	1.48	1.39
1	A	205	GLU	CG-CD	5.03	1.59	1.51

All (114) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	532	ARG	NE-CZ-NH2	-29.96	105.32	120.30
1	B	532	ARG	NE-CZ-NH2	-23.27	108.67	120.30
1	B	532	ARG	NE-CZ-NH1	19.00	129.80	120.30
1	A	532	ARG	NE-CZ-NH1	17.32	128.96	120.30
1	B	393	ARG	NE-CZ-NH2	-12.63	113.98	120.30
1	B	506	ARG	NE-CZ-NH1	10.47	125.54	120.30
1	B	433	ARG	NE-CZ-NH1	10.34	125.47	120.30
1	B	132	ARG	NE-CZ-NH2	-9.42	115.59	120.30
1	B	512	ASP	CB-CG-OD1	-9.18	110.04	118.30
1	A	40	ARG	NE-CZ-NH1	9.11	124.86	120.30
1	A	376	ASP	CB-CG-OD2	-9.03	110.17	118.30
1	A	633	VAL	CG1-CB-CG2	-9.01	96.49	110.90
1	A	713	ARG	NE-CZ-NH1	8.95	124.77	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	48	ASP	CB-CG-OD1	8.56	126.01	118.30
1	A	433	ARG	NE-CZ-NH1	8.31	124.45	120.30
1	B	268	ASP	CB-CG-OD2	-8.31	110.82	118.30
1	A	454	ASP	CB-CG-OD2	8.26	125.73	118.30
1	B	633	VAL	CG1-CB-CG2	-8.21	97.77	110.90
1	B	688	ARG	NE-CZ-NH1	8.15	124.38	120.30
1	B	619	ASP	CB-CG-OD1	-8.09	111.02	118.30
1	A	191	ARG	NE-CZ-NH2	-8.04	116.28	120.30
1	A	268	ASP	CB-CG-OD2	-7.76	111.31	118.30
1	B	713	ARG	NE-CZ-NH1	7.72	124.16	120.30
1	B	48	ASP	CB-CG-OD1	7.69	125.22	118.30
1	A	713	ARG	NE-CZ-NH2	-7.63	116.48	120.30
1	B	249	ASP	CB-CG-OD1	7.62	125.16	118.30
1	A	191	ARG	NE-CZ-NH1	7.48	124.04	120.30
1	A	434	TYR	CB-CG-CD1	7.45	125.47	121.00
1	B	123	ARG	NE-CZ-NH1	7.41	124.01	120.30
1	B	726	GLU	OE1-CD-OE2	-7.33	114.50	123.30
1	A	132	ARG	NE-CZ-NH1	7.29	123.94	120.30
1	B	268	ASP	CB-CG-OD1	7.28	124.85	118.30
1	A	60	ASP	CB-CG-OD1	7.27	124.85	118.30
1	A	578	ARG	NE-CZ-NH1	7.26	123.93	120.30
1	B	512	ASP	CB-CG-OD2	7.19	124.77	118.30
1	A	587	ASP	CB-CG-OD1	7.18	124.76	118.30
1	A	149	ARG	NE-CZ-NH1	7.13	123.87	120.30
1	A	395	ASP	CB-CG-OD1	7.03	124.63	118.30
1	B	683	ASP	CB-CG-OD2	-6.99	112.01	118.30
1	B	320	ASP	CB-CG-OD1	-6.93	112.07	118.30
1	B	506	ARG	NE-CZ-NH2	-6.85	116.87	120.30
1	A	649	ARG	NE-CZ-NH1	6.84	123.72	120.30
1	B	649	ARG	NE-CZ-NH1	6.82	123.71	120.30
1	A	506	ARG	NE-CZ-NH1	6.81	123.70	120.30
1	B	404	ARG	NE-CZ-NH1	6.76	123.68	120.30
1	B	504	ARG	NE-CZ-NH2	-6.76	116.92	120.30
1	A	183	LYS	CD-CE-NZ	6.75	127.23	111.70
1	A	255	ARG	NE-CZ-NH2	-6.75	116.93	120.30
1	A	532	ARG	CD-NE-CZ	6.73	133.02	123.60
1	B	426	ARG	NE-CZ-NH1	-6.70	116.95	120.30
1	B	60	ASP	CB-CG-OD1	6.65	124.28	118.30
1	B	40	ARG	NE-CZ-NH1	6.61	123.60	120.30
1	A	404	ARG	NE-CZ-NH1	6.59	123.60	120.30
1	A	198[A]	GLU	OE1-CD-OE2	-6.44	115.57	123.30
1	A	198[B]	GLU	OE1-CD-OE2	-6.44	115.57	123.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	73	GLU	OE1-CD-OE2	6.42	131.01	123.30
1	A	249	ASP	CB-CG-OD2	-6.42	112.52	118.30
1	A	198[A]	GLU	CG-CD-OE1	6.38	131.06	118.30
1	A	198[B]	GLU	CG-CD-OE1	6.38	131.06	118.30
1	B	587	ASP	CB-CG-OD1	6.35	124.01	118.30
1	A	670	ASP	CB-CG-OD2	-6.34	112.59	118.30
1	A	83	ASP	CB-CG-OD2	6.31	123.98	118.30
1	A	504	ARG	NE-CZ-NH1	6.28	123.44	120.30
1	A	649	ARG	NE-CZ-NH2	-6.22	117.19	120.30
1	A	688	ARG	NE-CZ-NH1	6.20	123.40	120.30
1	B	83	ASP	CB-CG-OD2	6.19	123.87	118.30
1	A	532	ARG	CG-CD-NE	-6.18	98.83	111.80
1	A	255	ARG	NE-CZ-NH1	6.17	123.38	120.30
1	A	426[A]	ARG	NE-CZ-NH1	-6.14	117.23	120.30
1	A	426[B]	ARG	NE-CZ-NH1	-6.14	117.23	120.30
1	B	169	LEU	CB-CG-CD2	6.14	121.44	111.00
1	A	263	ARG	NE-CZ-NH2	-6.07	117.27	120.30
1	A	123	ARG	NE-CZ-NH1	6.04	123.32	120.30
1	A	433	ARG	NE-CZ-NH2	-5.99	117.30	120.30
1	A	602	ARG	NE-CZ-NH2	-5.95	117.33	120.30
1	A	73	GLU	OE1-CD-OE2	-5.94	116.17	123.30
1	B	670	ASP	CB-CG-OD2	-5.92	112.97	118.30
1	B	532	ARG	CD-NE-CZ	5.89	131.84	123.60
1	B	688	ARG	NE-CZ-NH2	-5.88	117.36	120.30
1	B	434	TYR	CB-CG-CD1	5.86	124.51	121.00
1	A	40	ARG	NE-CZ-NH2	-5.82	117.39	120.30
1	A	602	ARG	NE-CZ-NH1	5.81	123.20	120.30
1	A	249	ASP	CB-CG-OD1	5.80	123.52	118.30
1	B	198	GLU	OE1-CD-OE2	-5.80	116.34	123.30
1	A	54	ARG	NE-CZ-NH1	5.75	123.17	120.30
1	B	198	GLU	CG-CD-OE1	5.72	129.75	118.30
1	B	161	ARG	NE-CZ-NH1	5.65	123.12	120.30
1	A	587	ASP	CB-CG-OD2	-5.57	113.29	118.30
1	B	222	ASP	CB-CG-OD1	-5.51	113.34	118.30
1	A	70	GLN	CA-CB-CG	5.51	125.52	113.40
1	B	639	ARG	NE-CZ-NH1	5.51	123.05	120.30
1	B	580	ASP	CB-CG-OD1	-5.50	113.35	118.30
1	A	41	ASP	CB-CG-OD2	5.39	123.15	118.30
1	A	55	HIS	CB-CA-C	5.36	121.13	110.40
1	B	225	LEU	CB-CG-CD2	-5.36	101.89	111.00
1	B	55	HIS	CB-CA-C	5.36	121.12	110.40
1	B	60	ASP	CB-CG-OD2	-5.34	113.49	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	459	ASP	CB-CG-OD1	5.32	123.09	118.30
1	B	730[A]	ARG	CA-CB-CG	5.29	125.05	113.40
1	B	730[B]	ARG	CA-CB-CG	5.29	125.05	113.40
1	A	404	ARG	NE-CZ-NH2	-5.29	117.66	120.30
1	B	504	ARG	NE-CZ-NH1	5.28	122.94	120.30
1	B	459	ASP	CB-CG-OD1	5.24	123.01	118.30
1	B	532	ARG	CG-CD-NE	-5.22	100.83	111.80
1	A	108	ARG	NE-CZ-NH2	-5.20	117.70	120.30
1	A	532	ARG	NH1-CZ-NH2	5.17	125.09	119.40
1	A	701	ARG	NE-CZ-NH1	5.17	122.88	120.30
1	B	310	LYS	CD-CE-NZ	-5.08	100.02	111.70
1	A	169	LEU	CB-CG-CD2	5.06	119.60	111.00
1	B	619	ASP	CB-CG-OD2	5.02	122.82	118.30
1	A	414	ASP	CB-CG-OD2	5.02	122.82	118.30
1	A	506	ARG	NE-CZ-NH2	-5.01	117.79	120.30
1	B	382	ARG	NE-CZ-NH2	-5.01	117.79	120.30
1	B	326	LEU	CA-CB-CG	-5.01	103.78	115.30

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	5585	0	5388	16	0
1	B	5582	0	5381	29	0
2	A	43	0	30	0	0
2	B	43	0	30	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	5	0	0	0	0
4	B	5	0	0	0	0
5	A	16	0	28	5	0
5	B	8	0	14	2	0
6	A	660	0	0	6	0
6	B	655	0	0	18	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	12604	0	10871	51	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 2.

All (51) 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:520:GLN:HG3	6:B:956:HOH:O	1.62	0.97
1:B:119[B]:THR:HG22	1:B:593:VAL:HG21	1.65	0.79
1:A:711[B]:GLN:OE1	6:A:901:HOH:O	2.00	0.78
1:B:119[B]:THR:HG23	6:B:955:HOH:O	1.83	0.78
1:B:540:ARG:NE	1:B:540:ARG:HA	2.00	0.77
1:A:255:ARG:NH2	6:A:904:HOH:O	2.18	0.76
5:B:804:MPD:HM2	5:B:804:MPD:H52	1.67	0.74
1:B:55:HIS:HD2	6:B:1307:HOH:O	1.69	0.74
1:B:512:ASP:OD1	6:B:901:HOH:O	2.10	0.69
5:A:805:MPD:H52	5:A:805:MPD:H11	1.73	0.69
1:B:519:GLU:OE1	6:B:902:HOH:O	2.11	0.68
5:A:805:MPD:H11	5:A:805:MPD:C5	2.23	0.68
1:A:55:HIS:HD2	6:A:1325:HOH:O	1.75	0.68
1:A:76:ASP:OD1	6:A:903:HOH:O	2.11	0.68
1:B:178:GLU:OE1	6:B:903:HOH:O	2.13	0.65
1:B:76:ASP:OD1	6:B:904:HOH:O	2.15	0.62
5:B:804:MPD:H52	5:B:804:MPD:CM	2.32	0.59
5:A:804:MPD:O4	5:A:804:MPD:CM	2.51	0.59
1:B:119[B]:THR:HG21	6:B:1485:HOH:O	2.04	0.57
1:B:410:GLU:HB2	6:B:1379:HOH:O	2.03	0.56
1:B:69:ALA:O	1:B:73:GLU:HG2	2.06	0.56
1:B:647:GLN:HG2	6:B:1275:HOH:O	2.06	0.56
1:B:540:ARG:NE	1:B:540:ARG:CA	2.69	0.55
1:B:540:ARG:HA	1:B:540:ARG:HE	1.71	0.55
1:A:119[A]:THR:CG2	1:A:265:ALA:HB2	2.37	0.55
1:A:647:GLN:HG2	6:A:1211:HOH:O	2.08	0.54
1:A:633:VAL:CG2	1:A:719:TYR:CZ	2.92	0.53
1:A:46:GLN:NE2	1:B:298:GLU:O	2.43	0.51
1:B:633:VAL:CG2	1:B:719:TYR:CZ	2.94	0.51
1:B:711[A]:GLN:NE2	6:B:913:HOH:O	2.44	0.51
1:B:692:THR:HG21	6:B:905:HOH:O	2.11	0.50
1:B:255:ARG:HD3	6:B:1256:HOH:O	2.11	0.50
1:B:405:PHE:HB3	1:B:412:PHE:HB2	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:804:MPD:O4	5:A:804:MPD:HM1	2.13	0.48
1:A:582:SER:OG	1:A:585:GLN:HG3	2.14	0.48
1:B:403[B]:ARG:NH1	6:B:922:HOH:O	2.47	0.47
1:B:255:ARG:CD	6:B:1256:HOH:O	2.65	0.45
5:A:805:MPD:H52	5:A:805:MPD:C1	2.44	0.45
1:B:112:HIS:CE1	1:B:141:ASP:O	2.71	0.44
1:A:144:ASN:HA	1:A:146:ASP:OD1	2.17	0.44
1:A:405:PHE:HB3	1:A:412:PHE:HB2	1.98	0.43
1:B:662:ASN:H	1:B:725:GLN:HE22	1.65	0.43
1:A:346:PHE:CD2	1:A:403[A]:ARG:CZ	3.01	0.43
1:A:112:HIS:CE1	1:A:141:ASP:O	2.72	0.43
1:B:649:ARG:HG2	6:B:1367:HOH:O	2.18	0.42
1:A:725:GLN:HG3	6:A:1164:HOH:O	2.19	0.42
1:B:738:LYS:NZ	6:B:938:HOH:O	2.52	0.42
1:A:119[A]:THR:HG21	1:A:265:ALA:HB2	2.02	0.41
1:B:422:LYS:NZ	6:B:949:HOH:O	2.54	0.40
1:A:509:PRO:HD2	1:A:591:MET:HG2	2.02	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	719/728 (99%)	709 (99%)	10 (1%)	0	100	100
1	B	718/728 (99%)	706 (98%)	12 (2%)	0	100	100
All	All	1437/1456 (99%)	1415 (98%)	22 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	558/560 (100%)	552 (99%)	6 (1%)	73	73
1	B	557/560 (100%)	547 (98%)	10 (2%)	59	55
All	All	1115/1120 (100%)	1099 (99%)	16 (1%)	67	65

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	40	ARG
1	A	118	ARG
1	A	141	ASP
1	A	608	LYS
1	A	610	ARG
1	A	649	ARG
1	B	40	ARG
1	B	64	LYS
1	B	118	ARG
1	B	141	ASP
1	B	540	ARG
1	B	608	LYS
1	B	649	ARG
1	B	690	ARG
1	B	730[A]	ARG
1	B	730[B]	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	55	HIS
1	A	247	ASN
1	A	647	GLN
1	B	46	GLN
1	B	55	HIS
1	B	544	GLN

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Mol	Chain	Res	Type
1	B	647	GLN
1	B	650	HIS
1	B	725	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](#)

3 non-standard protein/DNA/RNA residues 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
1	TOX	A	111	1,2	10,17,18	2.06	5 (50%)	10,23,25	1.66	3 (30%)
1	TOX	B	111[B]	-	10,17,18	3.41	3 (30%)	10,23,25	1.40	2 (20%)
1	TOX	B	111[A]	2	10,17,18	3.41	3 (30%)	10,23,25	1.40	2 (20%)

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
1	TOX	A	111	1,2	-	2/4/8/10	0/2/2/2
1	TOX	B	111[B]	-	-	2/4/8/10	0/2/2/2
1	TOX	B	111[A]	2	-	2/4/8/10	0/2/2/2

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	111[A]	TOX	CD1-NE1	-8.48	1.31	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	111[B]	TOX	CD1-NE1	-8.48	1.31	1.39
1	B	111[A]	TOX	O-C	5.57	1.42	1.19
1	B	111[B]	TOX	O-C	5.57	1.42	1.19
1	A	111	TOX	O-C	3.78	1.35	1.19
1	B	111[A]	TOX	CA-N	-3.24	1.38	1.48
1	B	111[B]	TOX	CA-N	-3.24	1.38	1.48
1	A	111	TOX	CE3-CD2	-2.84	1.36	1.42
1	A	111	TOX	CZ2-CE2	-2.64	1.35	1.41
1	A	111	TOX	CH2-CZ2	2.24	1.41	1.36
1	A	111	TOX	CH2-CZ3	2.08	1.43	1.38

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	111[A]	TOX	CZ2-CE2-CD2	-2.73	117.22	120.94
1	B	111[B]	TOX	CZ2-CE2-CD2	-2.73	117.22	120.94
1	A	111	TOX	CG-CB-CA	2.64	118.60	114.53
1	A	111	TOX	CB-CA-C	2.61	116.36	111.47
1	A	111	TOX	CZ3-CH2-CZ2	-2.44	117.02	120.44
1	B	111[A]	TOX	CB-CG-CD1	-2.07	125.41	127.97
1	B	111[B]	TOX	CB-CG-CD1	-2.07	125.41	127.97

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	111	TOX	N-CA-CB-CG
1	A	111	TOX	C-CA-CB-CG
1	B	111[A]	TOX	N-CA-CB-CG
1	B	111[A]	TOX	C-CA-CB-CG
1	B	111[B]	TOX	N-CA-CB-CG
1	B	111[B]	TOX	C-CA-CB-CG

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates i

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 2 are monoatomic - leaving 7 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
5	MPD	B	804	-	7,7,7	0.66	0	9,10,10	1.00	0
4	PO4	B	803	-	4,4,4	0.71	0	6,6,6	2.13	1 (16%)
2	HEM	B	801	1	41,50,50	1.60	8 (19%)	45,82,82	1.75	13 (28%)
5	MPD	A	805	-	7,7,7	1.09	0	9,10,10	1.61	3 (33%)
5	MPD	A	804	-	7,7,7	0.83	0	9,10,10	1.67	2 (22%)
4	PO4	A	803	-	4,4,4	0.87	0	6,6,6	1.18	1 (16%)
2	HEM	A	801	1	41,50,50	1.77	11 (26%)	45,82,82	1.90	14 (31%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	HEM	B	801	1	-	2/12/54/54	-
5	MPD	A	805	-	-	2/5/5/5	-
5	MPD	A	804	-	-	0/5/5/5	-
2	HEM	A	801	1	-	3/12/54/54	-
5	MPD	B	804	-	-	2/5/5/5	-

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	801	HEM	C1B-NB	-5.05	1.31	1.40
2	B	801	HEM	C1B-NB	-3.78	1.33	1.40
2	B	801	HEM	CAA-C2A	-3.77	1.46	1.52
2	B	801	HEM	C2A-C3A	-3.49	1.27	1.37
2	A	801	HEM	C4B-NB	-3.29	1.32	1.38
2	A	801	HEM	C4D-ND	-3.24	1.34	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	801	HEM	C3B-C2B	-3.17	1.30	1.37
2	B	801	HEM	CHB-C1B	3.15	1.43	1.35
2	A	801	HEM	FE-NB	3.15	2.12	1.96
2	B	801	HEM	C4B-NB	-3.10	1.32	1.38
2	A	801	HEM	O1D-CGD	3.08	1.32	1.22
2	A	801	HEM	O1A-CGA	2.77	1.31	1.22
2	A	801	HEM	CBA-CGA	2.71	1.56	1.50
2	A	801	HEM	CAA-C2A	-2.45	1.48	1.52
2	B	801	HEM	C1D-ND	-2.20	1.34	1.38
2	B	801	HEM	O1A-CGA	2.19	1.29	1.22
2	A	801	HEM	CHA-C4D	2.16	1.40	1.35
2	A	801	HEM	O2D-CGD	-2.03	1.23	1.30
2	A	801	HEM	CHB-C1B	2.03	1.40	1.35

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	801	HEM	CMA-C3A-C4A	-4.41	121.69	128.46
2	A	801	HEM	C1B-NB-C4B	4.21	109.42	105.07
2	B	801	HEM	CAD-C3D-C4D	4.03	131.71	124.66
4	B	803	PO4	O4-P-O1	-3.79	97.04	110.89
2	A	801	HEM	C4C-CHD-C1D	3.54	127.23	122.56
2	A	801	HEM	C4B-C3B-C2B	-3.51	104.33	107.11
2	B	801	HEM	CMA-C3A-C4A	-3.31	123.38	128.46
2	A	801	HEM	O2A-CGA-CBA	3.25	124.47	114.03
2	A	801	HEM	C4B-CHC-C1C	3.14	126.71	122.56
2	B	801	HEM	C4D-C3D-C2D	-3.07	102.43	106.90
5	A	804	MPD	O4-C4-C3	-2.92	99.58	111.36
2	A	801	HEM	CMA-C3A-C2A	2.84	130.31	124.94
2	B	801	HEM	CMB-C2B-C1B	-2.75	120.85	125.04
5	A	805	MPD	C5-C4-C3	2.60	123.92	111.69
2	A	801	HEM	O1A-CGA-CBA	-2.55	114.88	123.08
2	A	801	HEM	CBD-CAD-C3D	2.49	119.56	112.63
2	A	801	HEM	CAA-CBA-CGA	-2.47	106.83	113.76
2	A	801	HEM	C3C-C4C-NC	-2.47	106.28	110.94
2	B	801	HEM	C4B-CHC-C1C	2.46	125.80	122.56
2	A	801	HEM	CHA-C4D-ND	2.45	127.40	124.38
2	B	801	HEM	CAD-CBD-CGD	2.43	118.83	113.60
2	A	801	HEM	CHA-C4D-C3D	-2.38	120.86	125.33
2	B	801	HEM	CHA-C4D-C3D	-2.32	120.98	125.33
5	A	805	MPD	CM-C2-C1	-2.21	105.97	110.57
5	A	804	MPD	CM-C2-C1	2.15	115.05	110.57

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	801	HEM	CHC-C4B-NB	2.13	126.75	124.43
2	B	801	HEM	CHD-C1D-ND	2.13	126.75	124.43
2	B	801	HEM	C3D-C4D-ND	2.11	112.52	110.17
2	B	801	HEM	C1D-C2D-C3D	2.10	109.16	106.96
2	B	801	HEM	C2B-C1B-NB	-2.08	107.38	109.84
2	B	801	HEM	C1B-NB-C4B	2.04	107.18	105.07
2	A	801	HEM	CMD-C2D-C1D	2.02	128.12	125.04
4	A	803	PO4	O4-P-O2	-2.02	101.48	107.97
5	A	805	MPD	O4-C4-C3	-2.01	103.24	111.36

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	805	MPD	C2-C3-C4-C5
2	A	801	HEM	CAA-CBA-CGA-O2A
2	B	801	HEM	CAA-CBA-CGA-O2A
2	A	801	HEM	CAA-CBA-CGA-O1A
2	B	801	HEM	CAA-CBA-CGA-O1A
5	B	804	MPD	C2-C3-C4-C5
5	A	805	MPD	C2-C3-C4-O4
5	B	804	MPD	C2-C3-C4-O4
2	A	801	HEM	CAD-CBD-CGD-O1D

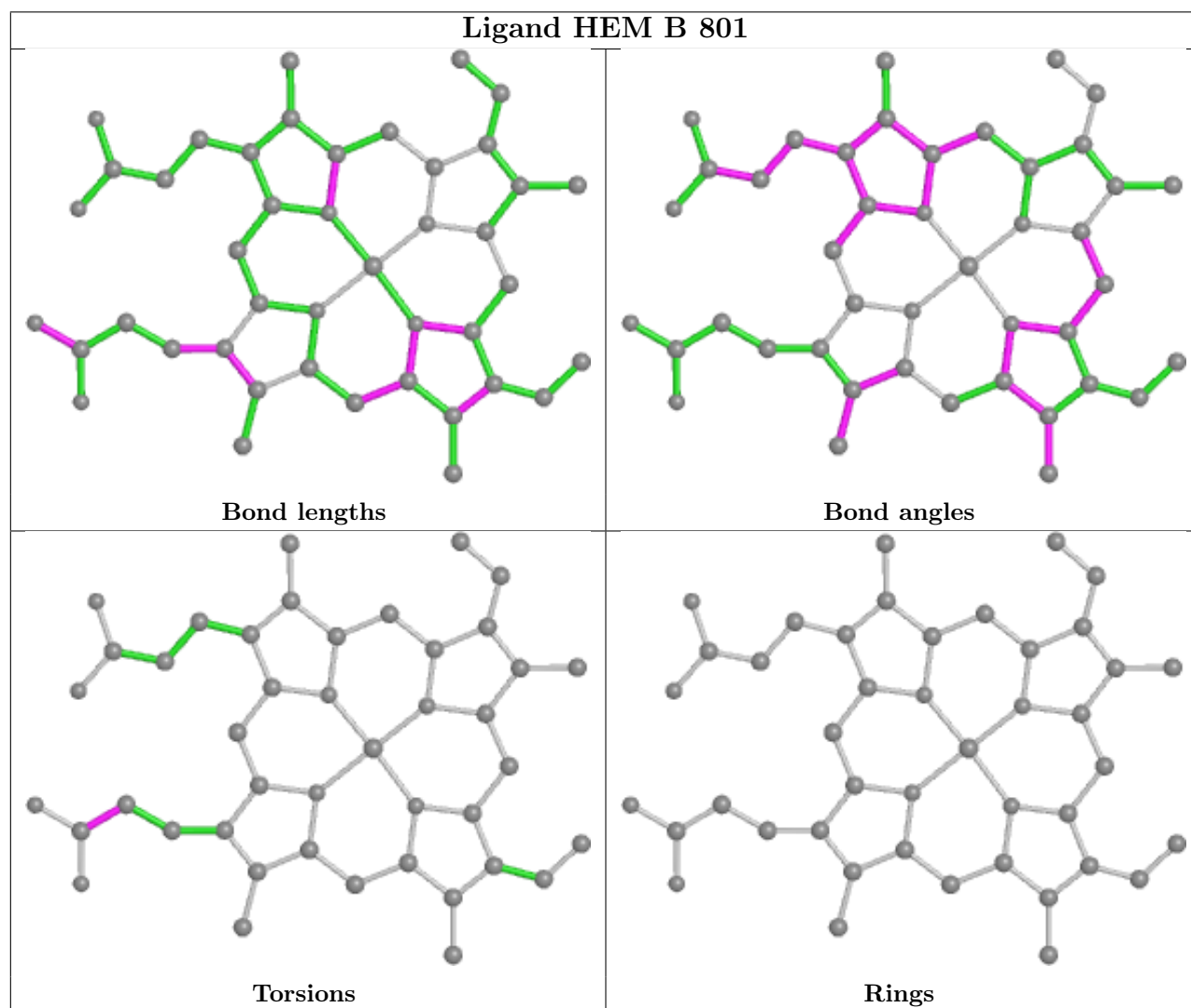
There are no ring outliers.

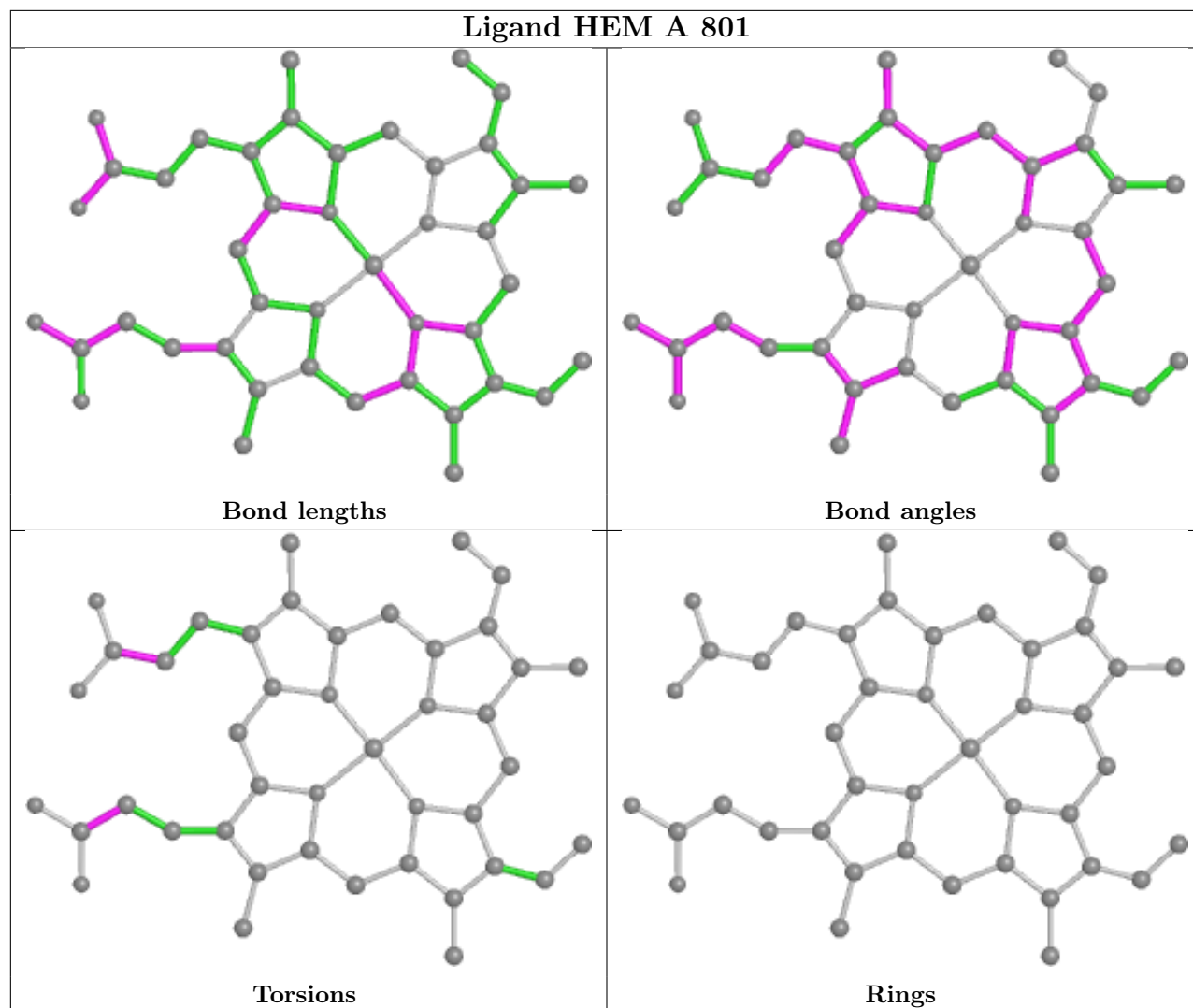
3 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	804	MPD	2	0
5	A	805	MPD	3	0
5	A	804	MPD	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

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	712/728 (97%)	-0.18	19 (2%) 54 57	17, 26, 46, 83	0
1	B	712/728 (97%)	-0.27	13 (1%) 68 71	17, 25, 42, 79	0
All	All	1424/1456 (97%)	-0.22	32 (2%) 62 64	17, 25, 45, 83	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	748	ALA	5.7
1	A	748	ALA	5.4
1	A	540	ARG	4.8
1	A	610	ARG	4.7
1	B	679	ALA	4.7
1	A	541	GLY	4.6
1	B	610	ARG	4.2
1	B	608	LYS	3.9
1	B	540	ARG	3.8
1	A	221	GLY	3.6
1	B	680	ALA	3.6
1	A	608	LYS	3.5
1	A	65	ASP	3.2
1	B	65	ASP	3.2
1	A	64	LYS	3.0
1	A	542	GLY	2.7
1	B	541	GLY	2.6
1	A	222	ASP	2.6
1	A	454	ASP	2.5
1	B	64	LYS	2.5
1	B	275	ILE	2.5
1	A	215	PRO	2.4
1	B	366	GLY	2.4
1	A	378	SER	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	679	ALA	2.3
1	A	207	ILE	2.2
1	A	682	ALA	2.2
1	A	145	LEU	2.2
1	B	681	ASP	2.1
1	A	680	ALA	2.0
1	B	44	PRO	2.0
1	A	647	GLN	2.0

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

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
1	TOX	A	111	16/17	0.97	0.16	17,23,34,37	2
1	TOX	B	111[A]	16/17	0.97	0.13	17,19,28,31	1
1	TOX	B	111[B]	16/17	0.97	0.13	17,19,23,28	1

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

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

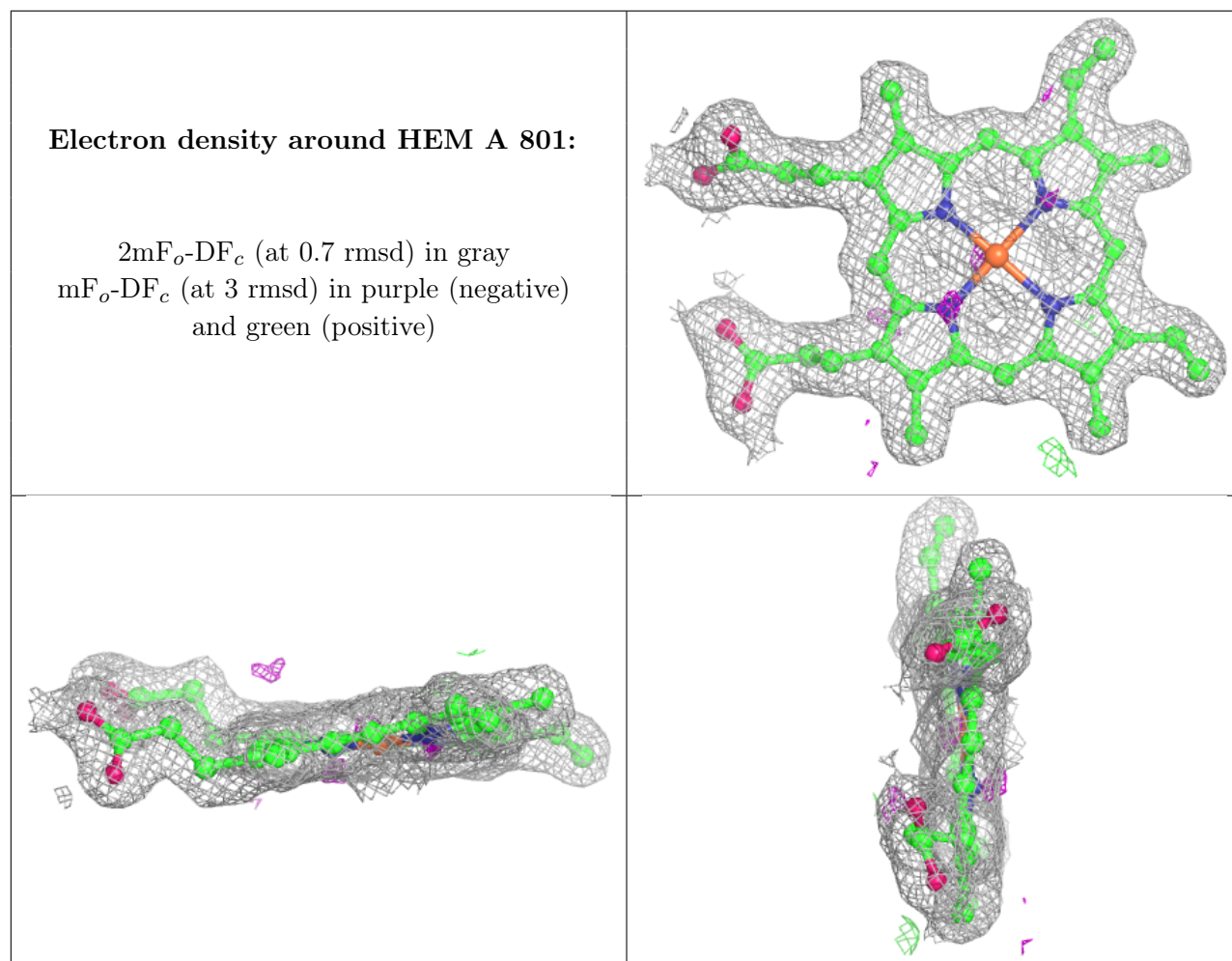
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	MPD	B	804	8/8	0.83	0.20	60,67,77,82	0
5	MPD	A	804	8/8	0.86	0.21	63,69,71,74	0
5	MPD	A	805	8/8	0.91	0.17	49,54,58,60	0
4	PO4	B	803	5/5	0.94	0.16	40,55,56,61	0
4	PO4	A	803	5/5	0.96	0.15	50,60,65,66	0
3	NA	A	802	1/1	0.98	0.04	21,21,21,21	0
2	HEM	A	801	43/43	0.98	0.14	19,22,25,26	0

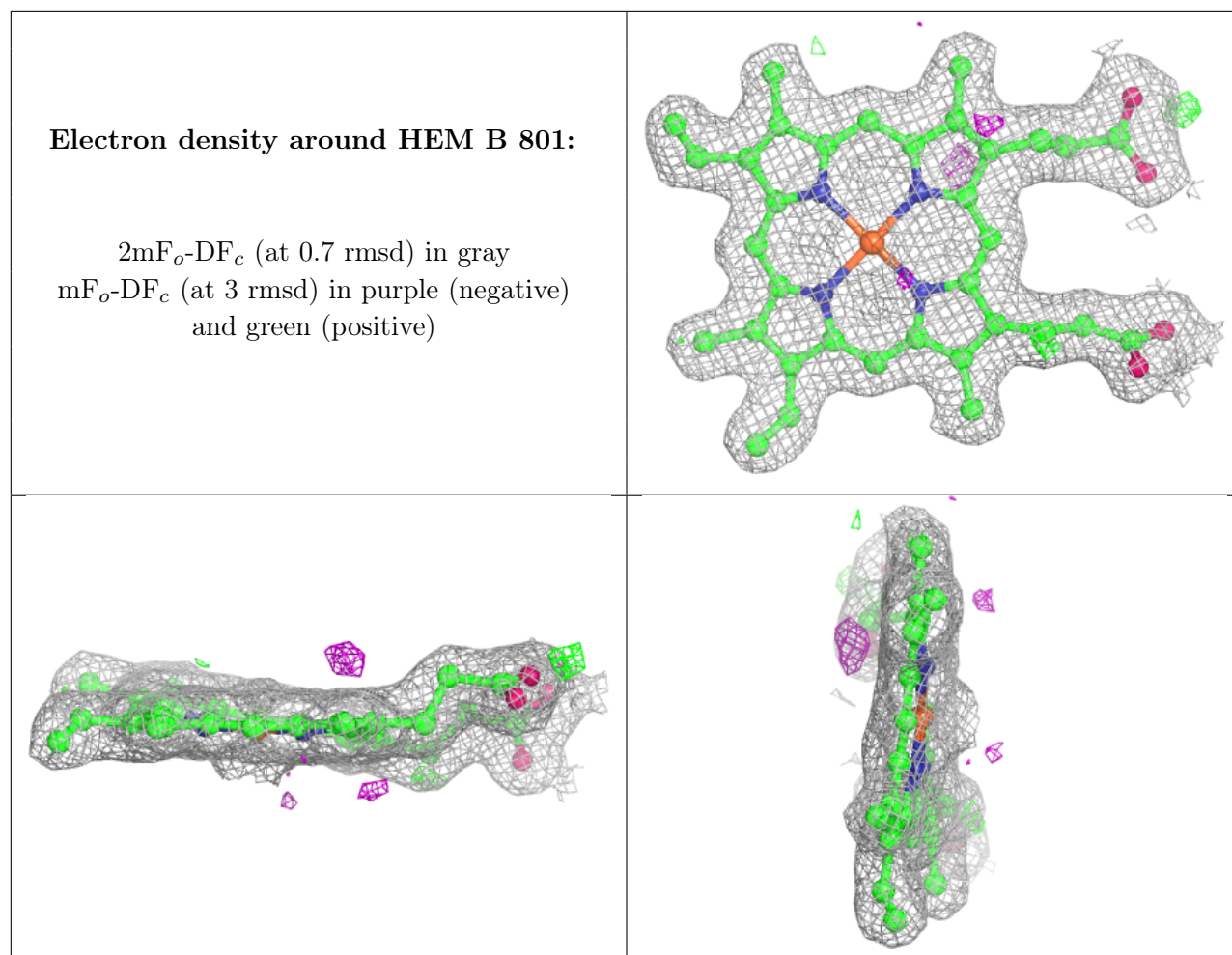
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	HEM	B	801	43/43	0.99	0.13	17,18,20,22	0
3	NA	B	802	1/1	1.00	0.05	22,22,22,22	0

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.