



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 12, 2024 – 01:19 PM EST

PDB ID : 3KS0
Title : Crystal structure of the heme domain of flavocytochrome b2 in complex with Fab B2B4
Authors : Golinelli-Pimpaneau, B.; Lederer, F.; Le, K.H.D.
Deposited on : 2009-11-20
Resolution : 2.70 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

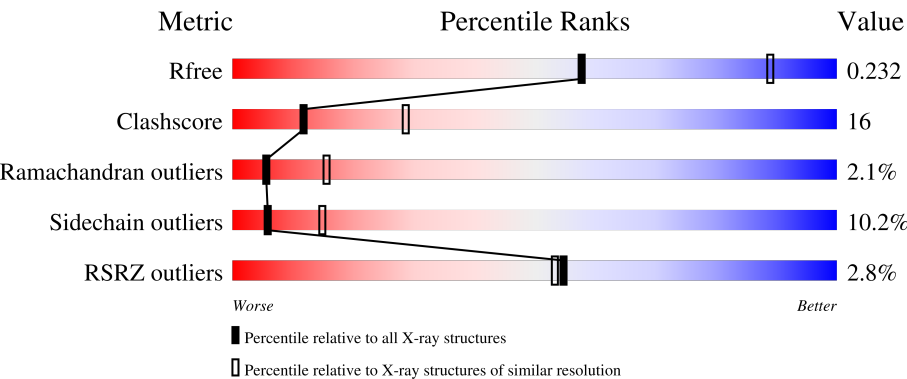
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	3.0
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	164625	3333 (2.70-2.70)
Clashscore	180529	3684 (2.70-2.70)
Ramachandran outliers	177936	3633 (2.70-2.70)
Sidechain outliers	177891	3633 (2.70-2.70)
RSRZ outliers	164620	3333 (2.70-2.70)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	J	214	<div><div></div><div><div></div><div></div><div></div><div></div></div><div>71%27%..</div></div>
1	L	214	<div><div></div><div><div></div><div></div><div></div><div></div></div><div>3%73%24%..</div></div>
2	H	225	<div><div></div><div><div></div><div></div><div></div><div></div><div></div></div><div>2%58%29%7%6%</div></div>
2	K	225	<div><div></div><div><div></div><div></div><div></div><div></div><div></div></div><div>4%60%27%7%5%</div></div>
3	A	95	<div><div></div><div><div></div><div></div><div></div><div></div></div><div>3%60%36%..</div></div>

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Mol	Chain	Length	Quality of chain
3	B	95	<div><div></div><div>3%</div><div>61%</div><div>31%</div><div>5%</div><div></div></div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8021 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 heme domain of flavocytochrome b2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	212	Total	C	N	O	S	0	0	0
			1579	988	263	323	5			
1	J	212	Total	C	N	O	S	0	0	0
			1574	985	260	323	6			

- Molecule 2 is a protein called Fragment Antigen Binding B2B4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	211	Total	C	N	O	S	0	0	0
			1613	1032	257	318	6			
2	K	214	Total	C	N	O	S	0	0	0
			1617	1034	259	318	6			

- Molecule 3 is a protein called Cytochrome b2, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	92	Total	C	N	O	S	0	0	0
			710	461	116	129	4			
3	B	92	Total	C	N	O	S	0	0	0
			710	461	116	129	4			

- Molecule 4 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



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

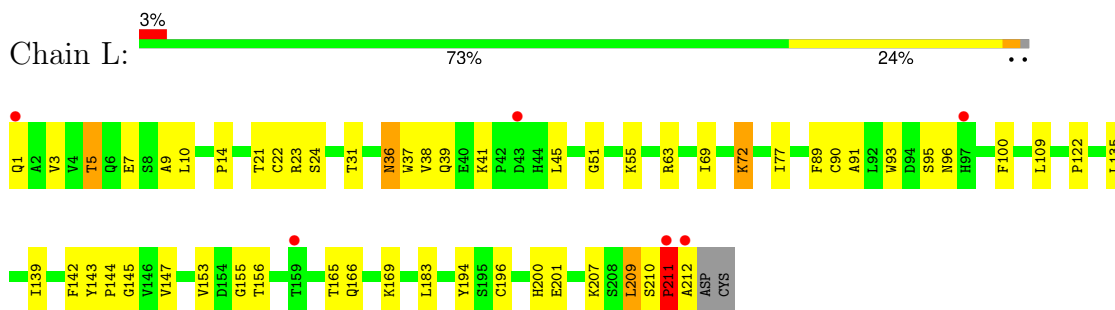
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	L	31	Total	O	0	0
			31	31		
5	H	21	Total	O	0	0
			21	21		
5	A	18	Total	O	0	0
			18	18		
5	J	31	Total	O	0	0
			31	31		
5	K	18	Total	O	0	0
			18	18		
5	B	13	Total	O	0	0
			13	13		

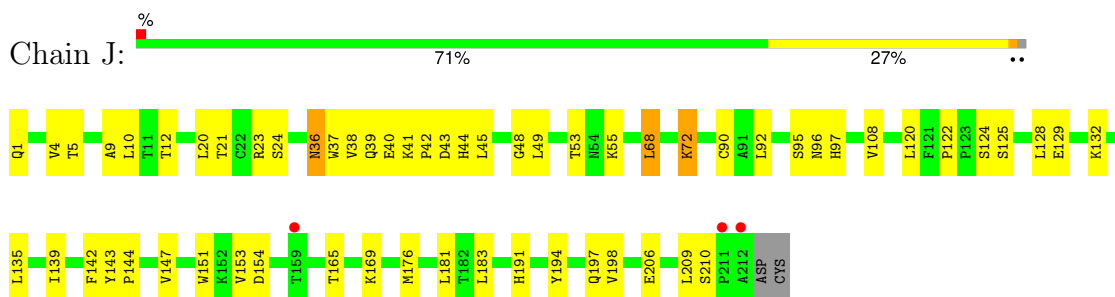
3 Residue-property plots

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.

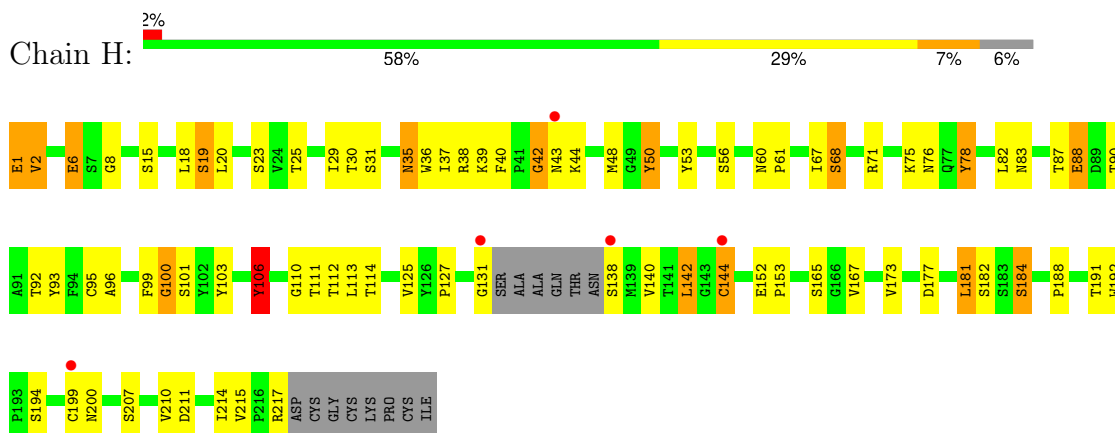
- Molecule 1: heme domain of flavocytochrome b2



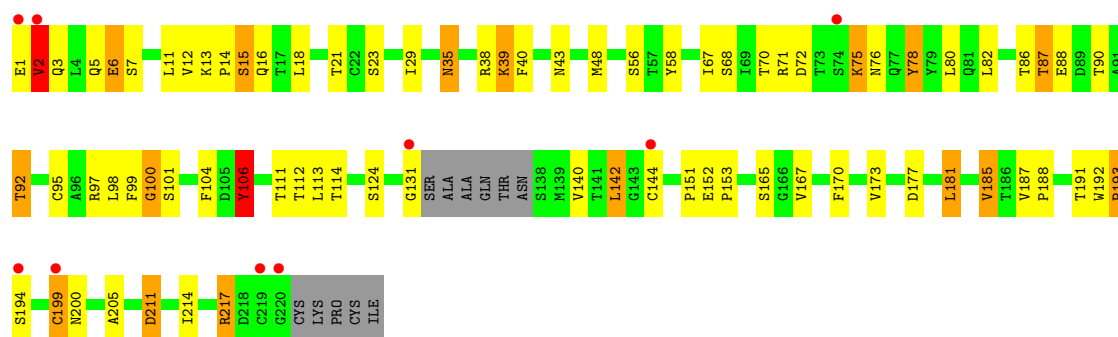
- Molecule 1: heme domain of flavocytochrome b2



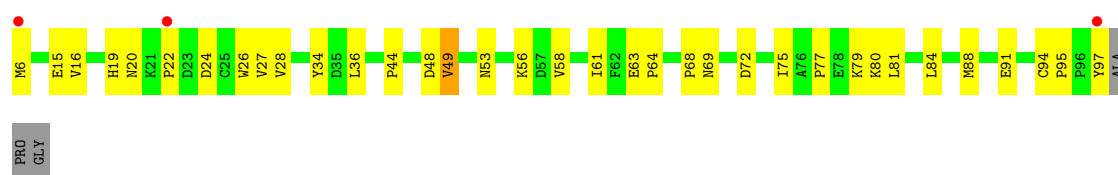
- Molecule 2: Fragment Antigen Binding B2B4



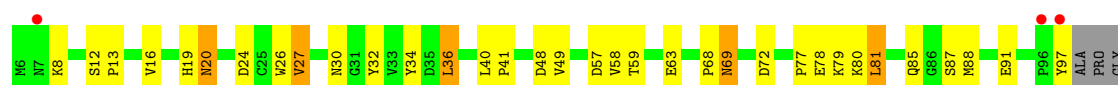
- Molecule 2: Fragment Antigen Binding B2B4



• Molecule 3: Cytochrome b2, mitochondrial



• Molecule 3: Cytochrome b2, mitochondrial



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	72.58Å 83.68Å 92.17Å 90.00° 96.80° 90.00°	Depositor
Resolution (Å)	24.75 – 2.70 24.75 – 2.70	Depositor EDS
% Data completeness (in resolution range)	96.6 (24.75-2.70) 96.6 (24.75-2.70)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	0.17	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.82 (at 2.72Å)	Xtriage
Refinement program	REFMAC 5.5.0102, CNS	Depositor
R, R_{free}	0.214 , 0.289 0.221 , 0.232	Depositor DCC
R_{free} test set	2914 reflections (9.96%)	wwPDB-VP
Wilson B-factor (Å ²)	31.5	Xtriage
Anisotropy	0.192	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 50.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	8021	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 17.32% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: HEM

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	J	0.70	1/1609 (0.1%)	0.79	0/2201
1	L	0.69	1/1615 (0.1%)	0.78	1/2210 (0.0%)
2	H	0.71	1/1659 (0.1%)	0.82	2/2275 (0.1%)
2	K	0.71	1/1663 (0.1%)	0.84	2/2281 (0.1%)
3	A	0.76	0/733	0.78	0/1003
3	B	0.78	0/733	0.95	1/1003 (0.1%)
All	All	0.72	4/8012 (0.0%)	0.82	6/10973 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	199	CYS	CB-SG	-6.19	1.71	1.82
2	K	199	CYS	CB-SG	-5.88	1.72	1.81
1	L	90	CYS	CB-SG	-5.86	1.72	1.81
1	J	90	CYS	CB-SG	-5.29	1.73	1.81

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	97	TYR	CA-C-O	-18.34	81.58	120.10
2	K	181	LEU	CA-CB-CG	6.74	130.81	115.30
1	L	209	LEU	CA-CB-CG	5.85	128.75	115.30
2	H	181	LEU	CA-CB-CG	5.50	127.95	115.30
2	K	106	TYR	CA-CB-CG	5.49	123.83	113.40
2	H	106	TYR	CA-CB-CG	5.33	123.53	113.40

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	J	1574	0	1515	39	0
1	L	1579	0	1525	44	0
2	H	1613	0	1548	56	0
2	K	1617	0	1538	68	0
3	A	710	0	697	28	0
3	B	710	0	697	30	0
4	A	43	0	30	6	0
4	B	43	0	30	4	0
5	A	18	0	0	0	0
5	B	13	0	0	0	0
5	H	21	0	0	1	0
5	J	31	0	0	2	0
5	K	18	0	0	1	0
5	L	31	0	0	2	0
All	All	8021	0	7580	241	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 16.

All (241) 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)
2:H:18:LEU:HD22	2:H:113:LEU:HD22	1.38	1.00
2:K:18:LEU:HD22	2:K:113:LEU:HD22	1.53	0.90
2:H:38:ARG:HG3	2:H:48:MET:SD	2.14	0.87
3:A:75:ILE:CD1	4:A:101:HEM:HAB	2.05	0.87
2:K:90:THR:HG23	2:K:114:THR:HA	1.59	0.82
1:J:5:THR:HG23	1:J:23:ARG:HB3	1.62	0.80
3:B:58:VAL:HG21	4:B:101:HEM:HMC2	1.63	0.79
3:A:75:ILE:HD11	4:A:101:HEM:HAB	1.64	0.79
3:B:77:PRO:HA	3:B:80:LYS:HE3	1.66	0.77
3:B:69:ASN:HD22	3:B:69:ASN:H	1.29	0.77
2:H:90:THR:HG23	2:H:114:THR:HA	1.66	0.76
2:H:2:VAL:HG21	2:H:106:TYR:CE1	2.21	0.76
2:H:29:ILE:H	2:H:76:ASN:HD21	1.32	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:16:VAL:HB	3:B:88:MET:HE1	1.69	0.75
2:K:188:PRO:O	2:K:191:THR:HG22	1.86	0.75
3:A:16:VAL:HB	3:A:88:MET:HE1	1.68	0.74
1:J:124:SER:HA	2:K:217:ARG:HH12	1.51	0.74
1:L:23:ARG:NH1	2:K:5:GLN:HB3	2.03	0.74
1:L:210:SER:O	1:L:211:PRO:C	2.26	0.74
2:H:200:ASN:HD22	2:H:211:ASP:HB3	1.54	0.73
1:J:122:PRO:HA	1:J:135:LEU:HD23	1.72	0.72
2:K:38:ARG:HG3	2:K:48:MET:SD	2.30	0.71
1:L:122:PRO:HA	1:L:135:LEU:HD23	1.73	0.70
2:K:39:LYS:HE2	2:K:43:ASN:HA	1.71	0.70
2:H:142:LEU:HG	2:H:214:ILE:HG21	1.73	0.69
3:B:16:VAL:HB	3:B:88:MET:CE	2.23	0.69
2:K:187:VAL:HB	2:K:191:THR:HG21	1.72	0.69
3:A:75:ILE:HD12	4:A:101:HEM:HAB	1.73	0.68
3:A:72:ASP:OD1	3:A:80:LYS:NZ	2.26	0.67
2:K:167:VAL:HG22	2:K:185:VAL:HB	1.76	0.66
2:H:18:LEU:CD2	2:H:113:LEU:HD22	2.21	0.66
2:H:152:GLU:HB3	2:H:153:PRO:HA	1.78	0.66
3:B:13:PRO:HA	3:B:88:MET:HE2	1.77	0.66
2:H:18:LEU:HD22	2:H:113:LEU:CD2	2.21	0.65
1:L:155:GLY:O	5:L:238:HOH:O	2.13	0.65
1:L:139:ILE:HG22	1:L:142:PHE:CE1	2.32	0.65
2:H:131:GLY:HA2	2:H:217:ARG:HB2	1.78	0.65
2:K:39:LYS:CE	2:K:43:ASN:HA	2.27	0.65
2:K:187:VAL:HB	2:K:191:THR:CG2	2.28	0.64
1:J:21:THR:HB	1:J:72:LYS:HD3	1.79	0.64
1:J:125:SER:HA	1:J:128:LEU:HD12	1.79	0.63
1:L:210:SER:O	1:L:212:ALA:N	2.32	0.63
2:H:60:ASN:OD1	2:H:61:PRO:HD2	1.98	0.63
2:K:152:GLU:HB3	2:K:153:PRO:HA	1.80	0.63
2:K:140:VAL:HG13	2:K:187:VAL:HG23	1.81	0.62
3:A:63:GLU:O	2:K:101:SER:HB2	1.99	0.62
2:K:140:VAL:HG21	2:K:192:TRP:CE3	2.35	0.61
2:H:188:PRO:HD2	2:H:191:THR:HG21	1.83	0.61
3:A:53:ASN:HA	3:A:56:LYS:HE3	1.83	0.61
2:K:86:THR:HG22	2:K:88:GLU:HG2	1.80	0.61
2:H:2:VAL:HG11	2:H:106:TYR:CZ	2.36	0.61
3:A:84:LEU:HD11	3:A:88:MET:HE2	1.82	0.61
1:J:197:GLN:HG2	1:J:206:GLU:HG3	1.83	0.59
1:L:153:VAL:HG22	1:L:194:TYR:CD2	2.37	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:188:PRO:HD2	2:H:191:THR:CG2	2.31	0.59
1:J:165:THR:HG22	2:K:173:VAL:HG13	1.84	0.59
2:K:13:LYS:O	2:K:16:GLN:HG3	2.04	0.58
2:H:92:THR:HG22	2:H:112:THR:HA	1.85	0.58
1:J:53:THR:CG2	1:J:68:LEU:HD13	2.33	0.58
1:L:166:GLN:HE21	2:H:42:GLY:HA3	1.69	0.58
3:A:68:PRO:HD3	2:K:101:SER:HA	1.85	0.58
2:H:125:VAL:HG21	2:H:210:VAL:CG2	2.34	0.58
2:K:29:ILE:H	2:K:76:ASN:HD21	1.52	0.57
1:L:93:TRP:CE2	3:B:68:PRO:HG2	2.39	0.57
1:J:20:LEU:HD23	5:J:279:HOH:O	2.03	0.57
1:L:9:ALA:O	1:L:10:LEU:HD23	2.05	0.57
3:A:44:PRO:HG2	4:A:101:HEM:HMA2	1.88	0.56
1:L:5:THR:HG23	1:L:23:ARG:HB3	1.87	0.56
1:J:139:ILE:HG22	1:J:142:PHE:CD1	2.41	0.56
2:K:199:CYS:O	2:K:211:ASP:HA	2.07	0.55
1:L:165:THR:HG22	2:H:173:VAL:HG13	1.89	0.55
3:A:91:GLU:CD	3:A:91:GLU:H	2.10	0.55
1:L:7:GLU:HG2	1:L:10:LEU:HD21	1.89	0.55
3:B:19:HIS:HD2	3:B:24:ASP:OD1	1.89	0.55
1:L:23:ARG:HH11	2:K:5:GLN:HB3	1.70	0.55
2:H:92:THR:HG22	2:H:112:THR:HG23	1.89	0.54
2:H:211:ASP:OD2	2:H:211:ASP:N	2.39	0.54
3:B:13:PRO:HA	3:B:88:MET:CE	2.37	0.54
2:H:2:VAL:HG21	2:H:106:TYR:CD1	2.42	0.54
1:J:4:VAL:HG22	1:J:92:LEU:HD12	1.89	0.54
1:L:122:PRO:HA	1:L:135:LEU:CD2	2.36	0.54
2:H:71:ARG:HA	2:H:78:TYR:HA	1.88	0.54
2:K:211:ASP:OD2	2:K:211:ASP:N	2.40	0.54
2:H:100:GLY:O	2:H:101:SER:HB2	2.07	0.54
1:L:22:CYS:O	1:L:72:LYS:HB3	2.07	0.54
1:L:36:ASN:C	1:L:36:ASN:HD22	2.11	0.54
2:H:78:TYR:OH	2:H:95:CYS:HB2	2.06	0.54
3:A:15:GLU:OE1	3:A:19:HIS:HE1	1.90	0.54
2:K:142:LEU:HG	2:K:214:ILE:HG21	1.89	0.54
2:K:39:LYS:HE2	2:K:43:ASN:CA	2.38	0.54
3:A:44:PRO:HG2	4:A:101:HEM:CMA	2.38	0.53
3:B:40:LEU:N	3:B:41:PRO:HD2	2.23	0.53
2:H:101:SER:HA	3:B:68:PRO:HD3	1.89	0.53
1:J:53:THR:HG21	1:J:68:LEU:HD13	1.89	0.53
3:B:27:VAL:HG22	3:B:58:VAL:HG22	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:36:TRP:HB3	2:H:48:MET:HG3	1.90	0.53
2:K:86:THR:CG2	2:K:88:GLU:HG2	2.38	0.53
2:H:38:ARG:CG	2:H:48:MET:SD	2.94	0.52
3:B:69:ASN:HD22	3:B:69:ASN:N	1.99	0.52
1:L:3:VAL:HG11	2:K:3:GLN:CD	2.30	0.52
1:J:139:ILE:HG22	1:J:142:PHE:CE1	2.45	0.52
1:J:154:ASP:OD2	1:J:191:HIS:HD2	1.93	0.51
2:H:144:CYS:O	2:H:182:SER:HA	2.09	0.51
2:K:86:THR:HG22	2:K:88:GLU:H	1.76	0.51
2:H:1:GLU:CG	2:H:1:GLU:O	2.57	0.51
1:L:22:CYS:O	1:L:72:LYS:CB	2.59	0.51
1:L:93:TRP:CD2	3:B:68:PRO:HG2	2.45	0.51
1:L:209:LEU:CD1	1:L:212:ALA:H	2.24	0.51
2:H:35:ASN:ND2	2:H:96:ALA:O	2.43	0.51
3:A:79:LYS:O	3:A:81:LEU:HD13	2.11	0.51
1:L:36:ASN:HD22	1:L:37:TRP:N	2.08	0.51
1:L:3:VAL:HG11	2:K:3:GLN:NE2	2.26	0.51
3:A:16:VAL:HB	3:A:88:MET:CE	2.40	0.50
1:J:48:GLY:HA3	2:K:104:PHE:O	2.11	0.50
2:K:140:VAL:HG22	2:K:142:LEU:HD13	1.94	0.50
2:K:71:ARG:HA	2:K:78:TYR:HA	1.94	0.50
1:J:139:ILE:HG12	1:J:198:VAL:HG21	1.94	0.49
2:H:127:PRO:CB	2:H:214:ILE:HD13	2.42	0.49
3:A:19:HIS:HD2	3:A:24:ASP:OD1	1.96	0.49
2:H:39:LYS:HE2	2:H:44:LYS:H	1.77	0.49
1:J:40:GLU:HA	1:J:45:LEU:O	2.11	0.49
2:K:131:GLY:HA2	2:K:217:ARG:HB2	1.93	0.49
3:B:69:ASN:H	3:B:69:ASN:ND2	2.04	0.49
3:A:28:VAL:HG11	3:A:84:LEU:HD22	1.95	0.49
2:H:1:GLU:O	2:H:1:GLU:HG3	2.13	0.49
2:K:7:SER:HB2	2:K:21:THR:HB	1.95	0.49
2:K:100:GLY:O	2:K:101:SER:HB2	2.12	0.49
4:B:101:HEM:HHC	4:B:101:HEM:HBB2	1.95	0.48
1:J:39:GLN:OE1	1:J:41:LYS:HG2	2.13	0.48
2:K:35:ASN:N	2:K:35:ASN:HD22	2.11	0.48
2:K:14:PRO:O	2:K:15:SER:CB	2.61	0.48
1:J:36:ASN:HD22	1:J:37:TRP:N	2.12	0.48
2:K:1:GLU:O	2:K:2:VAL:CG2	2.61	0.48
2:K:1:GLU:O	2:K:2:VAL:HG23	2.14	0.48
2:K:11:LEU:HD12	2:K:114:THR:HB	1.96	0.48
3:B:58:VAL:HG21	4:B:101:HEM:CMC	2.39	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:K:35:ASN:N	2:K:35:ASN:ND2	2.61	0.48
2:H:40:PHE:CB	2:H:43:ASN:HB2	2.43	0.48
2:H:19:SER:O	2:H:20:LEU:HD23	2.13	0.47
2:H:93:TYR:O	2:H:110:GLY:HA2	2.14	0.47
1:J:44:HIS:NE2	2:K:92:THR:HG21	2.29	0.47
2:H:40:PHE:HE2	2:H:88:GLU:O	1.98	0.47
3:A:26:TRP:HA	3:A:34:TYR:O	2.15	0.47
3:A:61:ILE:O	3:A:64:PRO:HD2	2.15	0.47
2:K:2:VAL:HG21	2:K:106:TYR:CE1	2.50	0.47
1:L:211:PRO:O	1:L:212:ALA:CB	2.63	0.47
3:A:84:LEU:HD21	3:A:88:MET:HG2	1.97	0.47
1:J:9:ALA:O	1:J:10:LEU:HD23	2.15	0.47
3:A:77:PRO:HA	3:A:80:LYS:HE3	1.97	0.47
1:L:14:PRO:HD3	1:L:109:LEU:O	2.14	0.47
2:K:78:TYR:OH	2:K:95:CYS:HB2	2.14	0.47
2:H:30:THR:C	2:H:53:TYR:HB2	2.36	0.47
3:A:58:VAL:HG21	4:A:101:HEM:HMC2	1.97	0.47
2:H:127:PRO:HB2	2:H:214:ILE:HD13	1.97	0.46
1:J:97:HIS:HA	2:K:58:TYR:OH	2.15	0.46
2:K:6:GLU:OE1	2:K:6:GLU:N	2.39	0.46
1:J:165:THR:HG22	2:K:173:VAL:CG1	2.46	0.46
1:L:183:LEU:HD22	3:A:97:TYR:OH	2.16	0.46
1:J:122:PRO:HA	1:J:135:LEU:CD2	2.45	0.46
1:J:44:HIS:HD2	5:K:299:HOH:O	1.99	0.46
2:K:72:ASP:OD2	2:K:75:LYS:HG3	2.16	0.46
2:H:50:TYR:CD1	2:H:50:TYR:C	2.90	0.45
4:B:101:HEM:HHC	4:B:101:HEM:CBB	2.47	0.45
2:H:6:GLU:OE1	2:H:6:GLU:N	2.47	0.45
3:B:57:ASP:OD2	3:B:59:THR:OG1	2.17	0.45
2:H:40:PHE:HB2	2:H:43:ASN:HB2	1.97	0.45
2:H:35:ASN:ND2	2:H:35:ASN:N	2.63	0.45
2:H:35:ASN:N	2:H:35:ASN:HD22	2.14	0.45
1:L:63:ARG:HG2	1:L:63:ARG:HH11	1.81	0.45
1:L:200:HIS:O	1:L:201:GLU:C	2.55	0.45
2:H:92:THR:CG2	2:H:112:THR:HG23	2.47	0.44
1:J:124:SER:HA	2:K:217:ARG:NH1	2.27	0.44
2:K:38:ARG:CG	2:K:48:MET:SD	3.02	0.44
1:L:209:LEU:HD12	1:L:212:ALA:H	1.82	0.44
2:K:97:ARG:HG2	2:K:98:LEU:N	2.32	0.44
3:B:20:ASN:H	3:B:20:ASN:HD22	1.66	0.44
1:J:5:THR:CG2	1:J:23:ARG:HB3	2.42	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:85:GLN:OE1	3:B:85:GLN:HA	2.18	0.44
2:H:101:SER:HB2	3:B:63:GLU:O	2.18	0.44
2:K:192:TRP:HA	2:K:193:PRO:C	2.38	0.43
3:B:34:TYR:HB3	3:B:36:LEU:HD13	2.00	0.43
1:L:139:ILE:HG22	1:L:142:PHE:CD1	2.52	0.43
1:L:155:GLY:C	3:A:49:VAL:HG13	2.39	0.43
1:J:151:TRP:CE3	1:J:181:LEU:HD22	2.53	0.43
2:K:67:ILE:HG12	2:K:68:SER:N	2.33	0.43
3:B:72:ASP:OD1	3:B:80:LYS:NZ	2.48	0.43
2:H:140:VAL:HG11	2:H:192:TRP:HB3	1.99	0.43
3:B:79:LYS:O	3:B:81:LEU:HD13	2.18	0.43
2:K:140:VAL:CG1	2:K:187:VAL:HG23	2.46	0.43
2:H:25:THR:HG23	5:H:251:HOH:O	2.17	0.43
1:J:120:LEU:HA	1:J:120:LEU:HD12	1.75	0.43
1:L:91:ALA:HB2	1:L:100:PHE:CD2	2.53	0.43
1:L:63:ARG:O	1:L:77:ILE:HA	2.18	0.43
3:A:53:ASN:ND2	3:A:94:CYS:SG	2.91	0.43
3:A:69:ASN:HB2	1:J:96:ASN:HB3	2.01	0.42
1:J:12:THR:O	1:J:108:VAL:HA	2.18	0.42
2:K:200:ASN:HD22	2:K:211:ASP:HB3	1.84	0.42
2:K:86:THR:O	2:K:88:GLU:N	2.52	0.42
1:L:96:ASN:HB3	3:B:69:ASN:HB2	2.00	0.42
1:J:153:VAL:HG22	1:J:194:TYR:CD2	2.54	0.42
2:H:152:GLU:CB	2:H:153:PRO:HA	2.46	0.42
1:L:39:GLN:OE1	1:L:41:LYS:HG2	2.19	0.42
2:K:12:VAL:HG13	2:K:16:GLN:HB2	2.01	0.42
2:K:39:LYS:HE2	2:K:43:ASN:C	2.40	0.42
1:L:31:THR:HG23	3:B:30:ASN:HD21	1.85	0.42
1:L:38:VAL:HG22	1:L:89:PHE:HB2	2.01	0.42
3:A:94:CYS:HA	3:A:95:PRO:HD3	1.86	0.42
2:K:70:THR:O	2:K:78:TYR:HB2	2.19	0.42
3:B:26:TRP:HA	3:B:34:TYR:O	2.20	0.42
1:J:143:TYR:CD1	1:J:144:PRO:HA	2.54	0.41
3:B:20:ASN:HD22	3:B:20:ASN:N	2.18	0.41
3:B:69:ASN:N	3:B:69:ASN:ND2	2.68	0.41
3:B:91:GLU:CD	3:B:91:GLU:H	2.23	0.41
1:L:36:ASN:ND2	1:L:37:TRP:N	2.68	0.41
1:J:1:GLN:OE1	1:J:97:HIS:ND1	2.53	0.41
1:L:143:TYR:HA	1:L:144:PRO:C	2.41	0.41
3:A:34:TYR:HE1	3:A:80:LYS:HG2	1.85	0.41
1:J:43:ASP:O	1:J:45:LEU:HG	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:K:86:THR:O	2:K:87:THR:C	2.59	0.41
3:B:8:LYS:HE3	3:B:8:LYS:HB2	1.74	0.41
1:L:63:ARG:HD3	5:L:236:HOH:O	2.20	0.41
1:L:69:ILE:O	1:L:72:LYS:HD2	2.21	0.41
2:K:40:PHE:HE2	2:K:88:GLU:O	2.03	0.41
2:K:131:GLY:HA2	2:K:217:ARG:CG	2.50	0.41
2:K:151:PRO:HD2	2:K:205:ALA:CB	2.51	0.41
2:H:8:GLY:HA3	2:H:20:LEU:CD2	2.51	0.41
2:H:188:PRO:HD2	2:H:191:THR:HG22	2.03	0.41
2:K:2:VAL:HG11	2:K:106:TYR:CZ	2.56	0.41
2:K:92:THR:HG22	2:K:112:THR:HA	2.02	0.41
1:J:42:PRO:O	1:J:45:LEU:HD12	2.22	0.40
1:J:176:MET:HB2	2:K:170:PHE:HE1	1.85	0.40
2:K:67:ILE:HG13	2:K:80:LEU:HD11	2.03	0.40
2:K:152:GLU:CB	2:K:153:PRO:HA	2.47	0.40
1:J:169:LYS:HE3	5:J:292:HOH:O	2.22	0.40
1:L:21:THR:HB	1:L:72:LYS:HD3	2.02	0.40
1:L:63:ARG:HG2	1:L:63:ARG:NH1	2.36	0.40
2:H:167:VAL:HA	2:H:184:SER:O	2.21	0.40
1:J:68:LEU:HD12	1:J:68:LEU:HA	1.86	0.40
2:H:39:LYS:HE3	2:H:44:LYS:HA	2.03	0.40
2:H:67:ILE:HG12	2:H:68:SER:N	2.36	0.40
1:L:51:GLY:HA3	2:H:103:TYR:N	2.37	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	J	210/214 (98%)	193 (92%)	16 (8%)	1 (0%)	25	49
1	L	210/214 (98%)	190 (90%)	18 (9%)	2 (1%)	13	33

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	H	207/225 (92%)	187 (90%)	12 (6%)	8 (4%)	2	5
2	K	210/225 (93%)	186 (89%)	15 (7%)	9 (4%)	2	4
3	A	90/95 (95%)	82 (91%)	8 (9%)	0	100	100
3	B	90/95 (95%)	78 (87%)	11 (12%)	1 (1%)	12	30
All	All	1017/1068 (95%)	916 (90%)	80 (8%)	21 (2%)	5	15

All (21) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	L	211	PRO
2	K	194	SER
2	H	2	VAL
2	H	42	GLY
2	H	99	PHE
2	H	100	GLY
2	K	15	SER
2	K	87	THR
2	K	100	GLY
3	B	78	GLU
2	H	87	THR
2	H	194	SER
2	K	2	VAL
2	K	177	ASP
2	H	106	TYR
2	K	217	ARG
2	H	15	SER
1	J	129	GLU
2	K	99	PHE
1	L	145	GLY
2	K	193	PRO

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	J	176/181 (97%)	163 (93%)	13 (7%)	11	28
1	L	176/181 (97%)	162 (92%)	14 (8%)	10	24
2	H	183/200 (92%)	157 (86%)	26 (14%)	2	7
2	K	180/200 (90%)	161 (89%)	19 (11%)	5	13
3	A	78/81 (96%)	71 (91%)	7 (9%)	8	19
3	B	78/81 (96%)	68 (87%)	10 (13%)	3	9
All	All	871/924 (94%)	782 (90%)	89 (10%)	6	15

All (89) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	1	GLN
1	L	5	THR
1	L	24	SER
1	L	36	ASN
1	L	45	LEU
1	L	55	LYS
1	L	72	LYS
1	L	95	SER
1	L	147	VAL
1	L	156	THR
1	L	169	LYS
1	L	196	CYS
1	L	207	LYS
1	L	211	PRO
2	H	1	GLU
2	H	6	GLU
2	H	19	SER
2	H	23	SER
2	H	31	SER
2	H	35	ASN
2	H	37	ILE
2	H	50	TYR
2	H	56	SER
2	H	68	SER
2	H	75	LYS
2	H	78	TYR
2	H	82	LEU
2	H	83	ASN
2	H	88	GLU
2	H	106	TYR

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Mol	Chain	Res	Type
2	H	111	THR
2	H	138	SER
2	H	142	LEU
2	H	144	CYS
2	H	165	SER
2	H	177	ASP
2	H	181	LEU
2	H	184	SER
2	H	207	SER
2	H	215	VAL
3	A	6	MET
3	A	20	ASN
3	A	22	PRO
3	A	27	VAL
3	A	36	LEU
3	A	48	ASP
3	A	49	VAL
1	J	24	SER
1	J	36	ASN
1	J	38	VAL
1	J	49	LEU
1	J	55	LYS
1	J	68	LEU
1	J	72	LYS
1	J	95	SER
1	J	132	LYS
1	J	147	VAL
1	J	183	LEU
1	J	209	LEU
1	J	210	SER
2	K	2	VAL
2	K	6	GLU
2	K	23	SER
2	K	35	ASN
2	K	39	LYS
2	K	56	SER
2	K	75	LYS
2	K	78	TYR
2	K	82	LEU
2	K	92	THR
2	K	106	TYR
2	K	111	THR

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Mol	Chain	Res	Type
2	K	124	SER
2	K	142	LEU
2	K	144	CYS
2	K	165	SER
2	K	181	LEU
2	K	185	VAL
2	K	211	ASP
3	B	12	SER
3	B	20	ASN
3	B	27	VAL
3	B	32	TYR
3	B	36	LEU
3	B	48	ASP
3	B	49	VAL
3	B	69	ASN
3	B	81	LEU
3	B	87	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	1	GLN
1	L	36	ASN
1	L	44	HIS
1	L	166	GLN
1	L	173	ASN
1	L	191	HIS
2	H	5	GLN
2	H	35	ASN
2	H	76	ASN
2	H	81	GLN
2	H	83	ASN
3	A	19	HIS
3	A	20	ASN
3	A	30	ASN
3	A	53	ASN
1	J	1	GLN
1	J	36	ASN
1	J	44	HIS
1	J	191	HIS
2	K	76	ASN
2	K	175	GLN

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Mol	Chain	Res	Type
3	B	19	HIS
3	B	20	ASN
3	B	30	ASN
3	B	53	ASN
3	B	69	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	$\# Z > 2$
4	HEM	B	101	3	42,50,50	1.86	6 (14%)	46,82,82	1.63	9 (19%)
4	HEM	A	101	3	42,50,50	1.88	6 (14%)	46,82,82	1.74	11 (23%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	HEM	B	101	3	-	5/12/54/54	-
4	HEM	A	101	3	-	2/12/54/54	-

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	101	HEM	C3D-C2D	7.98	1.54	1.36
4	B	101	HEM	C3D-C2D	7.81	1.53	1.36
4	A	101	HEM	C3C-C2C	-3.92	1.35	1.40
4	B	101	HEM	C3C-C2C	-3.17	1.36	1.40
4	B	101	HEM	CAB-C3B	2.98	1.55	1.47
4	A	101	HEM	CAB-C3B	2.95	1.55	1.47
4	B	101	HEM	FE-NB	2.65	2.12	1.98
4	B	101	HEM	C3C-CAC	2.53	1.53	1.47
4	A	101	HEM	FE-NB	2.31	2.10	1.98
4	A	101	HEM	C3C-CAC	2.25	1.52	1.47
4	A	101	HEM	C3B-C2B	-2.23	1.32	1.37
4	B	101	HEM	C4A-NA	2.07	1.40	1.36

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	101	HEM	C4D-ND-C1D	4.85	110.94	105.21
4	B	101	HEM	C4D-ND-C1D	4.65	110.72	105.21
4	A	101	HEM	CBA-CAA-C2A	-4.08	105.68	112.54
4	A	101	HEM	CAD-CBD-CGD	-3.41	104.62	113.67
4	B	101	HEM	CAD-CBD-CGD	-3.07	105.52	113.67
4	B	101	HEM	C4B-CHC-C1C	3.07	126.61	122.56
4	A	101	HEM	C4B-CHC-C1C	3.04	126.58	122.56
4	A	101	HEM	C3B-C4B-NB	-2.54	107.64	109.47
4	B	101	HEM	C1B-NB-C4B	2.53	108.20	105.21
4	B	101	HEM	CBA-CAA-C2A	-2.41	108.49	112.54
4	B	101	HEM	C3B-C2B-C1B	2.40	108.21	106.41
4	A	101	HEM	CHD-C1D-ND	2.37	126.98	124.44
4	B	101	HEM	C1D-C2D-C3D	-2.36	104.50	106.98
4	B	101	HEM	O1D-CGD-CBD	-2.34	115.68	123.09
4	A	101	HEM	O1A-CGA-CBA	-2.24	115.98	123.09
4	A	101	HEM	C2C-C3C-C4C	2.24	108.46	106.90
4	A	101	HEM	C1D-C2D-C3D	-2.15	104.72	106.98
4	A	101	HEM	CMA-C3A-C4A	-2.14	125.32	128.46
4	A	101	HEM	O2A-CGA-CBA	2.08	120.58	114.00
4	B	101	HEM	CHD-C1D-ND	2.01	126.59	124.44

There are no chirality outliers.

All (7) torsion outliers are listed below:

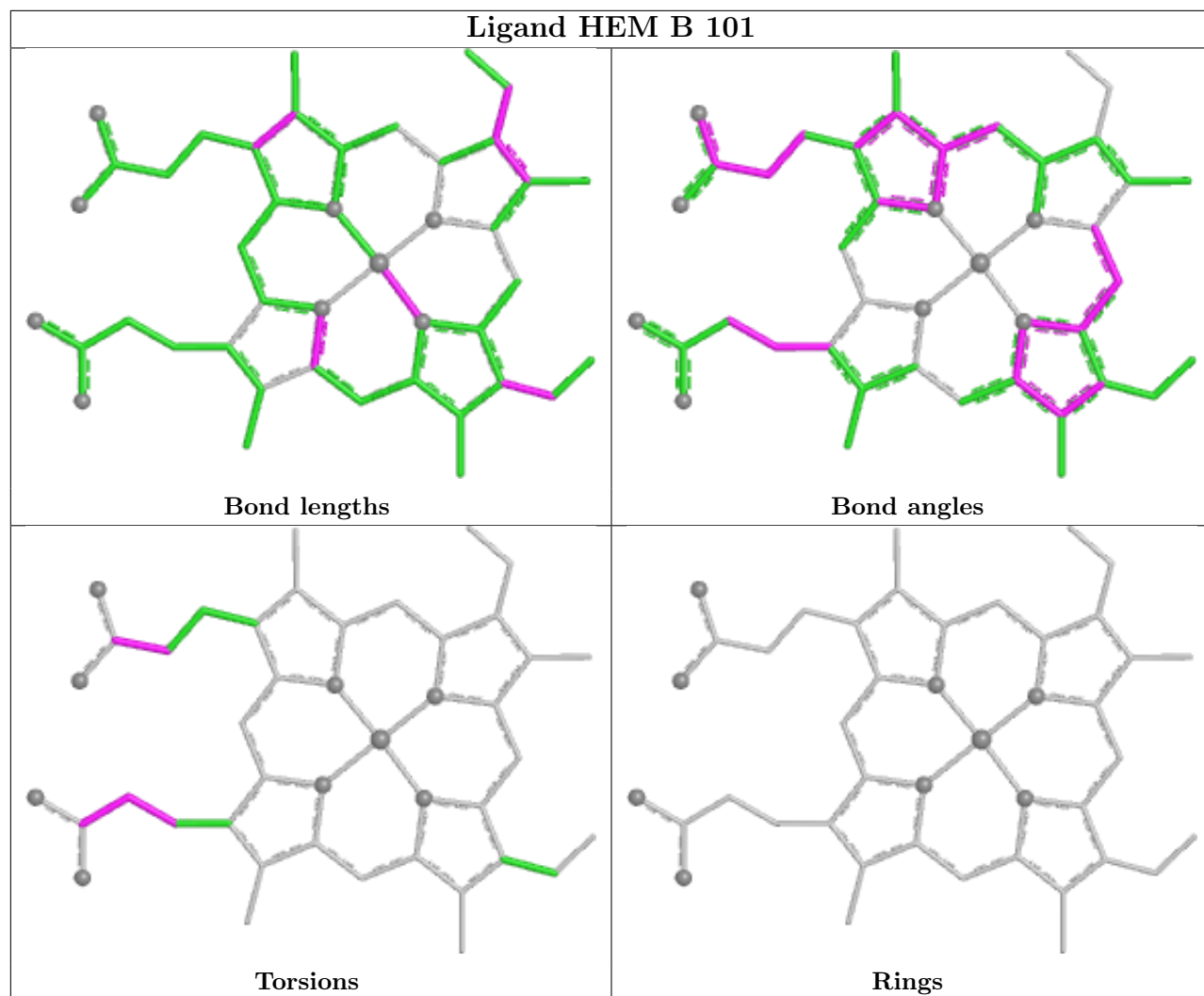
Mol	Chain	Res	Type	Atoms
4	B	101	HEM	C2A-CAA-CBA-CGA
4	A	101	HEM	CAD-CBD-CGD-O1D
4	B	101	HEM	CAA-CBA-CGA-O1A
4	A	101	HEM	CAD-CBD-CGD-O2D
4	B	101	HEM	CAA-CBA-CGA-O2A
4	B	101	HEM	CAD-CBD-CGD-O2D
4	B	101	HEM	CAD-CBD-CGD-O1D

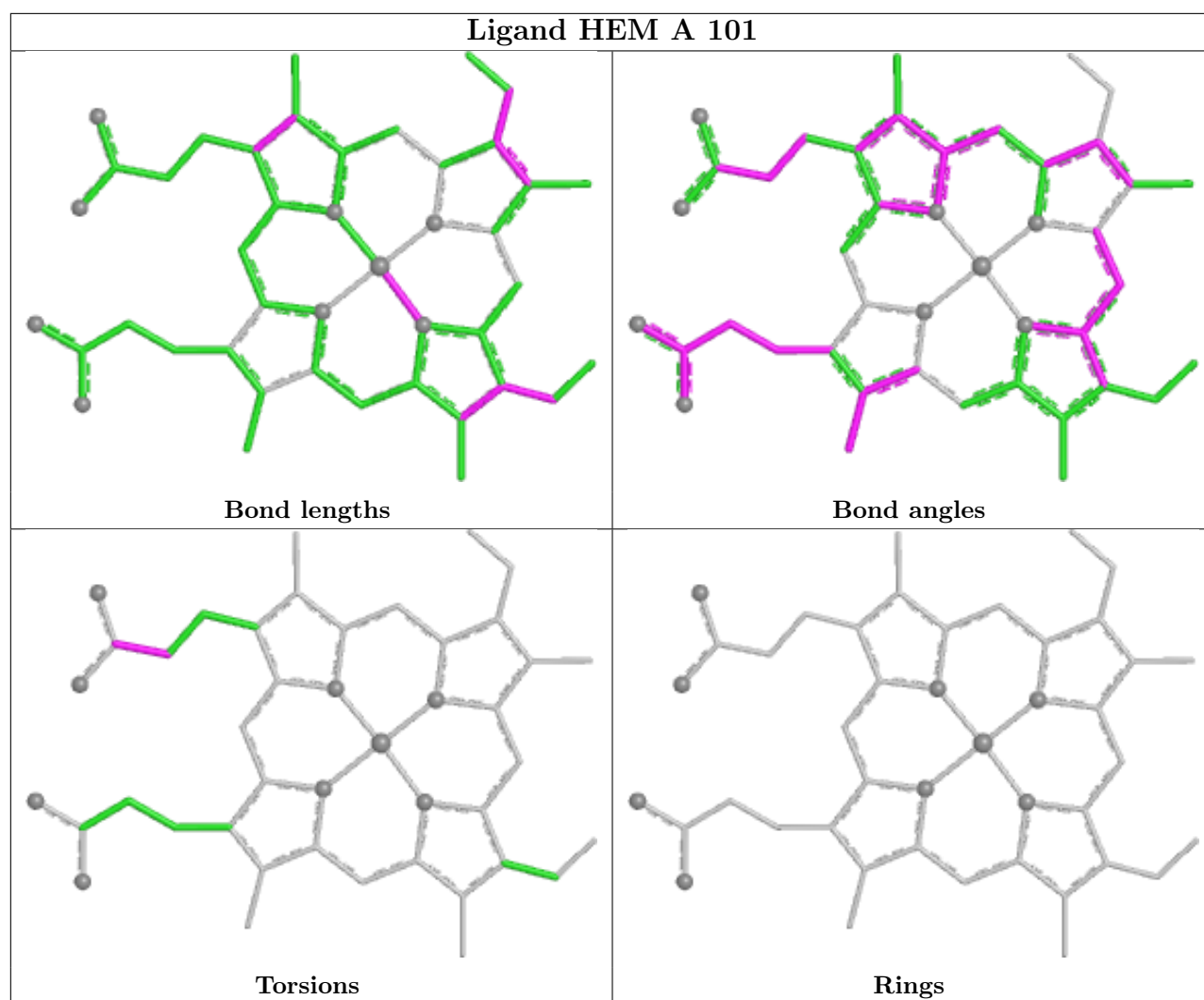
There are no ring outliers.

2 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	101	HEM	4	0
4	A	101	HEM	6	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	J	212/214 (99%)	0.11	3 (1%) 73 73	31, 42, 55, 63	0
1	L	212/214 (99%)	0.19	6 (2%) 55 53	31, 43, 55, 66	0
2	H	211/225 (93%)	0.28	5 (2%) 59 58	18, 44, 60, 68	3 (1%)
2	K	214/225 (95%)	0.29	9 (4%) 41 39	18, 44, 62, 81	3 (1%)
3	A	92/95 (96%)	0.21	3 (3%) 49 47	31, 44, 56, 69	0
3	B	92/95 (96%)	0.32	3 (3%) 49 47	33, 43, 58, 72	0
All	All	1033/1068 (96%)	0.23	29 (2%) 55 53	18, 43, 58, 81	6 (0%)

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	199	CYS	6.0
2	K	220	GLY	5.3
2	K	199	CYS	4.8
1	L	212	ALA	4.6
2	K	219	CYS	3.9
2	H	131	GLY	3.7
1	L	211	PRO	3.5
2	H	144	CYS	3.5
2	K	1	GLU	3.3
3	B	97	TYR	3.2
3	A	97	TYR	3.2
2	K	144	CYS	3.0
3	B	7	ASN	2.8
1	L	97	HIS	2.6
1	L	43	ASP	2.6
3	B	96	PRO	2.3
1	L	1	GLN	2.3
2	H	138	SER	2.3
3	A	6	MET	2.1

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Mol	Chain	Res	Type	RSRZ
1	J	212	ALA	2.1
1	L	159	THR	2.1
2	K	194	SER	2.1
2	K	2	VAL	2.1
1	J	211	PRO	2.1
1	J	159	THR	2.0
2	H	43	ASN	2.0
2	K	131	GLY	2.0
3	A	22	PRO	2.0
2	K	74	SER	2.0

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 monosaccharides in this entry.

6.4 Ligands [i](#)

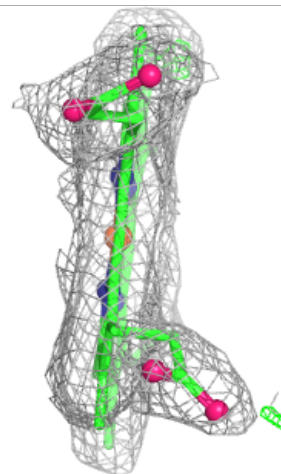
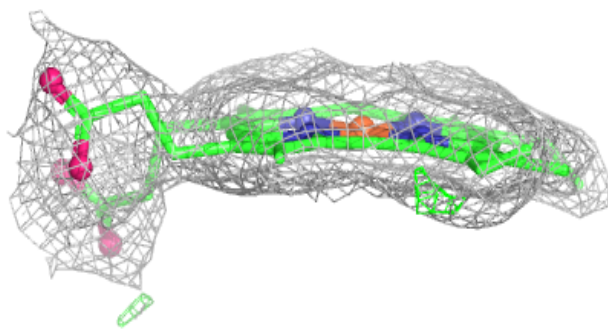
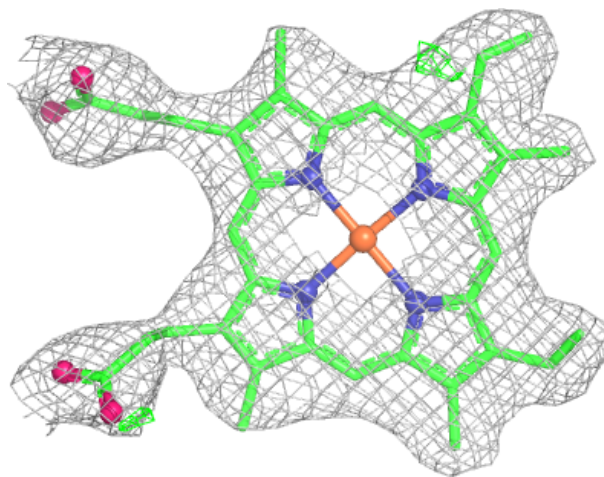
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

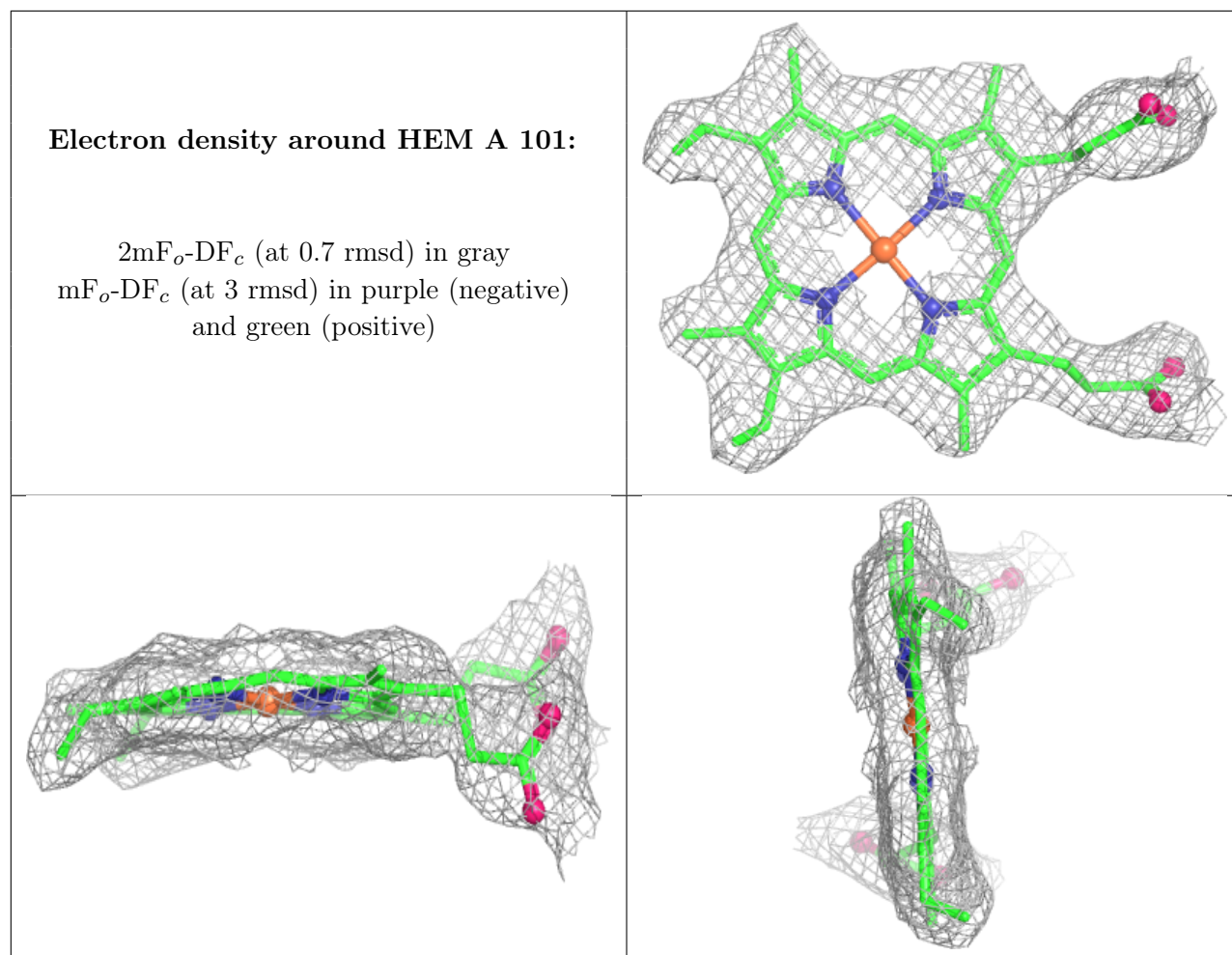
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	HEM	B	101	43/43	0.96	0.10	30,39,51,54	0
4	HEM	A	101	43/43	0.97	0.08	29,38,49,55	0

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 B 101:

$2mF_o-DF_c$ (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.