



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 12:53 PM UTC

PDB ID : 6G94 / pdb_00006g94
Title : Structure of E. coli hydrogenase-1 C19G variant in complex with cytochrome b
Authors : Volbeda, A.; Fontecilla-Camps, J.C.
Deposited on : 2018-04-10
Resolution : 2.50 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

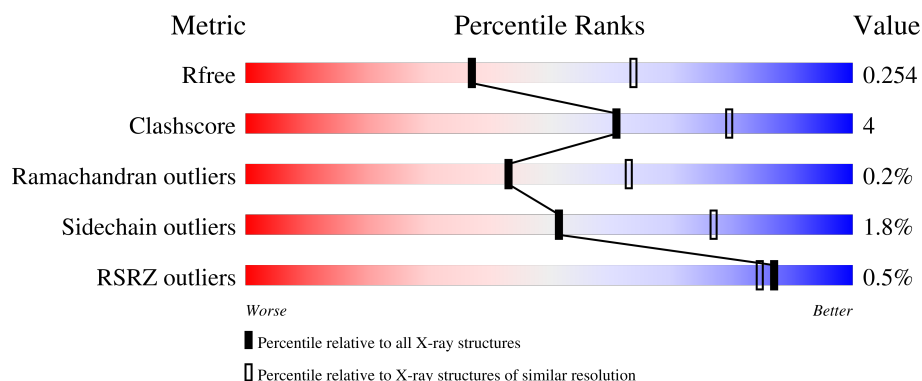
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	Q	335	 80% 8% 12%
1	R	335	 76% 10% • 13%
1	S	335	 79% 9% • 12%
1	T	335	 75% 13% • 12%
2	J	582	 88% 11%

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Mol	Chain	Length	Quality of chain
2	K	582	89% 10%
2	L	582	90% 9%
2	M	582	87% 12%
3	A	235	4% 60% 13% 27%
3	B	235	4% 57% 10% 32%

2 Entry composition

There are 12 unique types of molecules in this entry. The entry contains 30463 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 Hydrogenase-1 small chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	S	296	Total	C	N	O	S	0	0	0
			2242	1420	385	417	20			
1	T	296	Total	C	N	O	S	0	0	0
			2220	1408	379	413	20			
1	Q	295	Total	C	N	O	S	0	0	0
			2234	1416	383	415	20			
1	R	291	Total	C	N	O	S	0	0	0
			2203	1394	378	411	20			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
S	19	GLY	CYS	conflict	UNP P69739
S	328	ARG	-	expression tag	UNP P69739
S	329	SER	-	expression tag	UNP P69739
S	330	HIS	-	expression tag	UNP P69739
S	331	HIS	-	expression tag	UNP P69739
S	332	HIS	-	expression tag	UNP P69739
S	333	HIS	-	expression tag	UNP P69739
S	334	HIS	-	expression tag	UNP P69739
S	335	HIS	-	expression tag	UNP P69739
T	19	GLY	CYS	conflict	UNP P69739
T	328	ARG	-	expression tag	UNP P69739
T	329	SER	-	expression tag	UNP P69739
T	330	HIS	-	expression tag	UNP P69739
T	331	HIS	-	expression tag	UNP P69739
T	332	HIS	-	expression tag	UNP P69739
T	333	HIS	-	expression tag	UNP P69739
T	334	HIS	-	expression tag	UNP P69739
T	335	HIS	-	expression tag	UNP P69739
Q	19	GLY	CYS	conflict	UNP P69739
Q	328	ARG	-	expression tag	UNP P69739
Q	329	SER	-	expression tag	UNP P69739

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Chain	Residue	Modelled	Actual	Comment	Reference
Q	330	HIS	-	expression tag	UNP P69739
Q	331	HIS	-	expression tag	UNP P69739
Q	332	HIS	-	expression tag	UNP P69739
Q	333	HIS	-	expression tag	UNP P69739
Q	334	HIS	-	expression tag	UNP P69739
Q	335	HIS	-	expression tag	UNP P69739
R	19	GLY	CYS	conflict	UNP P69739
R	328	ARG	-	expression tag	UNP P69739
R	329	SER	-	expression tag	UNP P69739
R	330	HIS	-	expression tag	UNP P69739
R	331	HIS	-	expression tag	UNP P69739
R	332	HIS	-	expression tag	UNP P69739
R	333	HIS	-	expression tag	UNP P69739
R	334	HIS	-	expression tag	UNP P69739
R	335	HIS	-	expression tag	UNP P69739

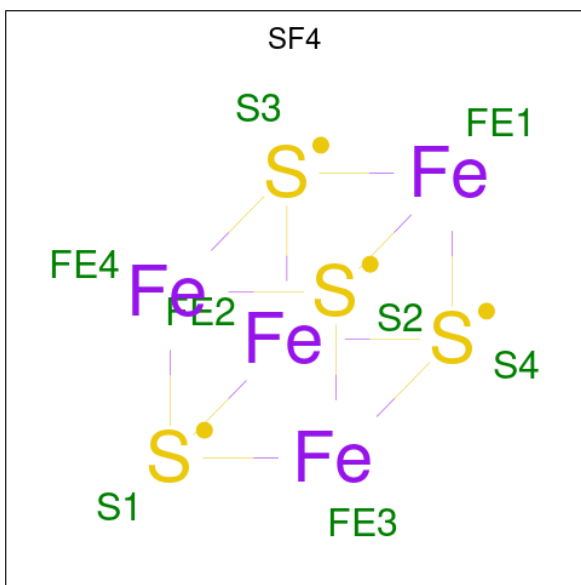
- Molecule 2 is a protein called Hydrogenase-1 large chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	581	Total	C	N	O	S	0	1	0
			4526	2874	787	838	27			
2	M	581	Total	C	N	O	S	0	0	0
			4512	2869	787	829	27			
2	J	581	Total	C	N	O	S	0	0	0
			4528	2876	792	833	27			
2	K	581	Total	C	N	O	S	0	0	0
			4526	2873	791	835	27			

- Molecule 3 is a protein called Probable Ni/Fe-hydrogenase 1 B-type cytochrome subunit.

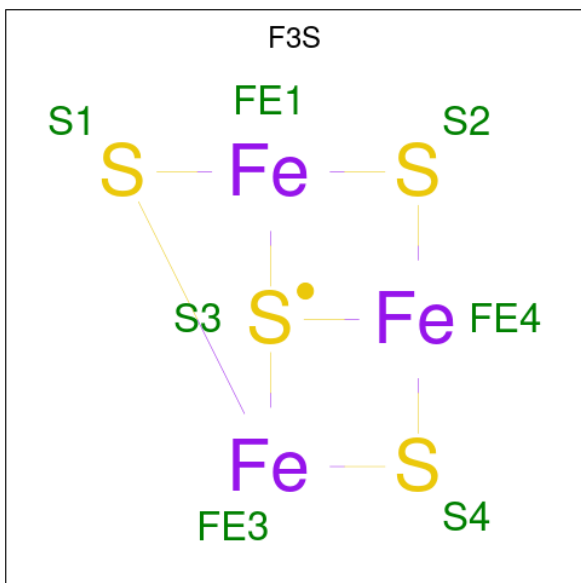
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	172	Total	C	N	O	S	0	0	0
			1310	888	204	207	11			
3	B	160	Total	C	N	O	S	0	0	0
			1242	842	196	194	10			

- Molecule 4 is IRON/SULFUR CLUSTER (CCD ID: SF4) (formula: Fe₄S₄).



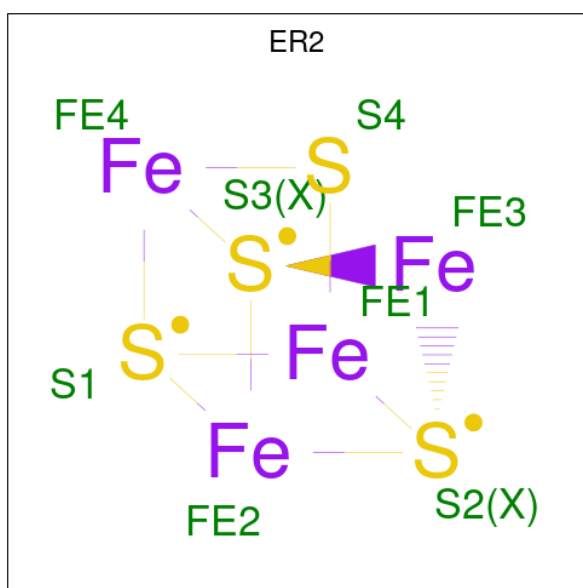
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	S	1	Total	Fe	S	0	0
			8	4	4		
4	T	1	Total	Fe	S	0	0
			8	4	4		
4	Q	1	Total	Fe	S	0	0
			8	4	4		
4	R	1	Total	Fe	S	0	0
			8	4	4		

- Molecule 5 is FE3-S4 CLUSTER (CCD ID: F3S) (formula: Fe_3S_4).



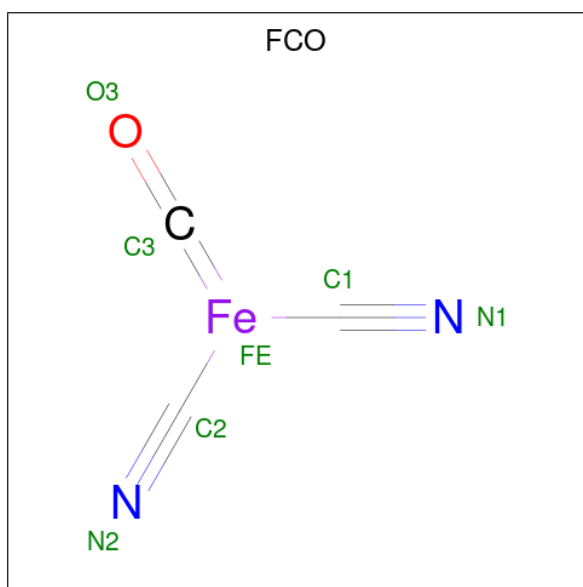
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	S	1	Total	Fe	S	0	0
			7	3	4		
5	T	1	Total	Fe	S	0	0
			7	3	4		
5	Q	1	Total	Fe	S	0	0
			7	3	4		
5	R	1	Total	Fe	S	0	0
			7	3	4		

- Molecule 6 is Fe₄S₄ (CCD ID: ER2) (formula: Fe₄S₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	S	1	Total	Fe	S	0	0
			8	4	4		
6	T	1	Total	Fe	S	0	0
			8	4	4		
6	Q	1	Total	Fe	S	0	0
			8	4	4		
6	R	1	Total	Fe	S	0	0
			8	4	4		

- Molecule 7 is CARBONMONOXIDE-(DICYANO) IRON (CCD ID: FCO) (formula: C₃FeN₂O).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	L	1	Total	C	Fe	N	O	0	0
			7	3	1	2	1		
7	M	1	Total	C	Fe	N	O	0	0
			7	3	1	2	1		
7	J	1	Total	C	Fe	N	O	0	0
			7	3	1	2	1		
7	K	1	Total	C	Fe	N	O	0	0
			7	3	1	2	1		

- Molecule 8 is NICKEL (II) ION (CCD ID: NI) (formula: Ni).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	L	1	Total	Ni	0	0
			1	1		
8	M	1	Total	Ni	0	0
			1	1		
8	J	1	Total	Ni	0	0
			1	1		
8	K	1	Total	Ni	0	0
			1	1		

- Molecule 9 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	L	1	Total	Mg	0	0
			1	1		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	M	1	Total Mg 1 1	0	0
9	J	1	Total Mg 1 1	0	0
9	K	1	Total Mg 1 1	0	0

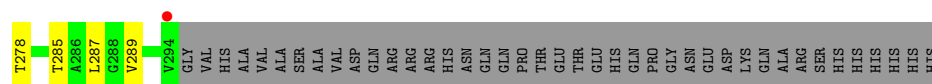
- | Mol | Chain | Residues | Atoms | ZeroOcc | AltConf |
|-----|-------|----------|-----------------|---------|---------|
| 10 | L | 1 | Total Cl
1 1 | 0 | 0 |
| 10 | M | 1 | Total Cl
1 1 | 0 | 0 |
| 10 | J | 1 | Total Cl
1 1 | 0 | 0 |
| 10 | K | 1 | Total Cl
1 1 | 0 | 0 |

- # HEM

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

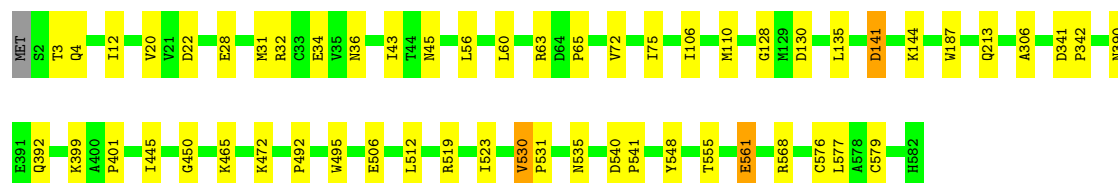
- Molecule 12 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
12	S	62	Total 62	O 62	0	0
12	L	111	Total 111	O 111	0	0
12	T	59	Total 59	O 59	0	0
12	M	113	Total 113	O 113	0	0
12	Q	62	Total 62	O 62	0	0
12	J	112	Total 112	O 112	0	0
12	R	63	Total 63	O 63	0	0
12	K	108	Total 108	O 108	0	0
12	A	6	Total 6	O 6	0	0
12	B	6	Total 6	O 6	0	0



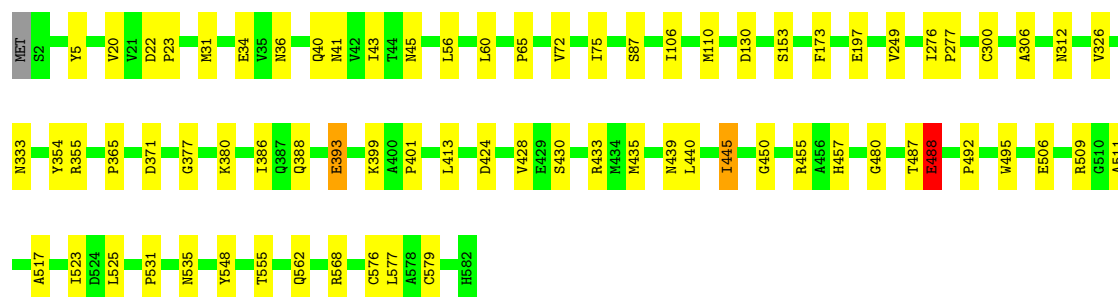
• Molecule 2: Hydrogenase-1 large chain

Chain L: 90% 9%



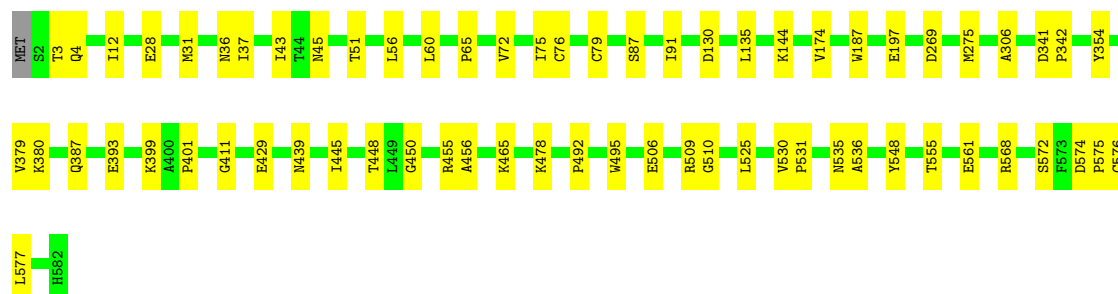
• Molecule 2: Hydrogenase-1 large chain

Chain M: 87% 12%



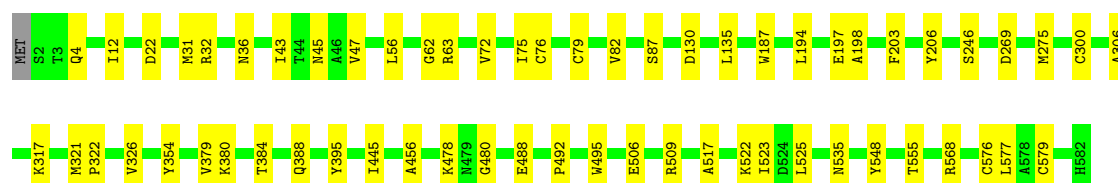
• Molecule 2: Hydrogenase-1 large chain

Chain J: 88% 11%

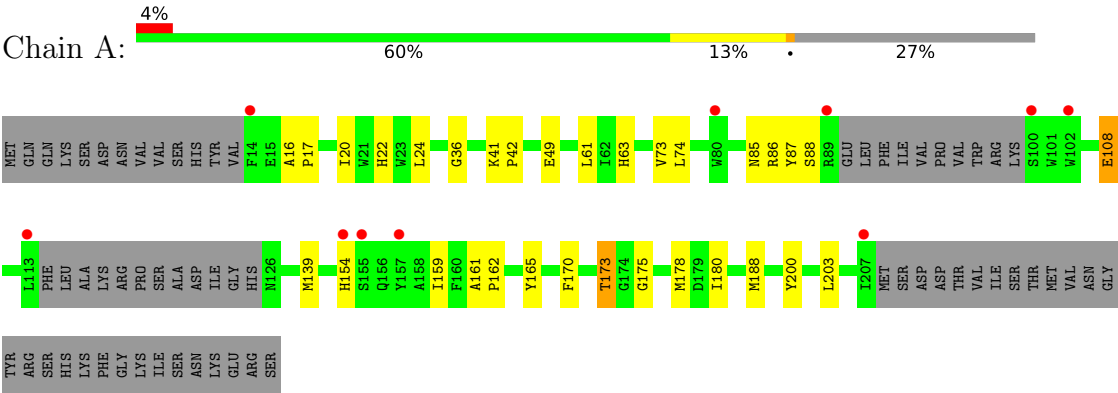


• Molecule 2: Hydrogenase-1 large chain

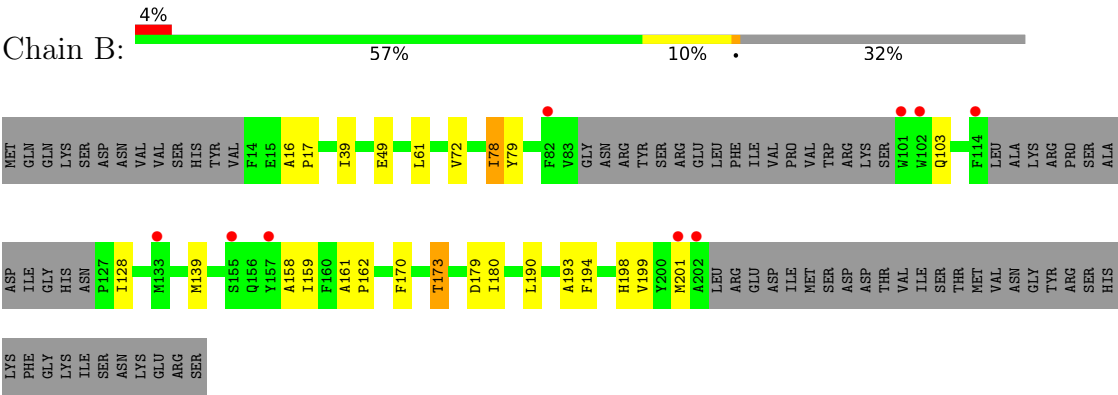
Chain K: 89% 10%



● Molecule 3: Probable Ni/Fe-hydrogenase 1 B-type cytochrome subunit



● Molecule 3: Probable Ni/Fe-hydrogenase 1 B-type cytochrome subunit



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	124.41Å 165.03Å 206.22Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.00 – 2.50 25.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	97.6 (25.00-2.50) 97.6 (25.00-2.50)	Depositor EDS
R_{merge}	0.30	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.50 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.8.0189	Depositor
R, R_{free}	0.222 , 0.253 0.223 , 0.254	Depositor DCC
R_{free} test set	7076 reflections (4.82%)	wwPDB-VP
Wilson B-factor (Å ²)	42.3	Xtriage
Anisotropy	0.490	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 33.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	30463	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.81% 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: F3S, FCO, ER2, NI, SF4, CL, HEM, MG

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	Q	0.67	0/2292	0.87	0/3120
1	R	0.68	0/2260	0.89	3/3077 (0.1%)
1	S	0.68	0/2300	0.89	1/3131 (0.0%)
1	T	0.67	0/2278	0.92	4/3104 (0.1%)
2	J	0.66	0/4647	0.88	4/6328 (0.1%)
2	K	0.63	0/4645	0.85	2/6327 (0.0%)
2	L	0.67	0/4648	0.86	3/6332 (0.0%)
2	M	0.65	0/4631	0.88	8/6308 (0.1%)
3	A	0.79	0/1356	1.03	5/1848 (0.3%)
3	B	0.74	0/1288	1.02	0/1753
All	All	0.67	0/30345	0.89	30/41328 (0.1%)

There are no bond length outliers.

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	165	TYR	N-CA-C	-7.64	104.48	113.88
3	A	22	HIS	N-CA-C	-6.24	105.57	113.43
3	A	88	SER	N-CA-C	6.05	119.38	110.30
2	K	384	THR	N-CA-C	-5.87	105.47	114.16
2	L	128	GLY	N-CA-C	5.86	119.73	112.64
3	A	108	GLU	N-CA-C	5.85	117.46	111.14
2	M	5	TYR	N-CA-C	5.83	117.94	109.07
2	M	153	SER	N-CA-C	5.77	118.05	108.99
2	J	510	GLY	N-CA-C	5.76	118.21	110.43
2	J	379	VAL	N-CA-C	-5.71	100.08	108.99
2	M	393	GLU	CB-CA-C	-5.70	108.86	117.07
1	T	270	ILE	N-CA-C	-5.55	102.05	107.55
2	L	141	ASP	CA-C-N	5.55	125.38	119.28
2	L	141	ASP	C-N-CA	5.55	125.38	119.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	K	379	VAL	CB-CA-C	-5.50	105.54	111.59
2	M	393	GLU	CA-C-O	5.38	121.06	117.94
1	T	121	VAL	N-CA-C	5.38	116.11	110.62
3	A	85	ASN	N-CA-C	5.36	122.20	110.80
2	M	488	GLU	N-CA-C	5.27	122.03	110.80
1	T	296	VAL	N-CA-C	-5.26	105.72	110.82
1	R	152	ILE	CA-C-N	5.25	124.69	119.24
1	R	152	ILE	C-N-CA	5.25	124.69	119.24
1	T	268	VAL	N-CA-C	5.21	116.37	108.96
2	J	411	GLY	CA-C-N	5.21	125.26	119.32
2	J	411	GLY	C-N-CA	5.21	125.26	119.32
1	R	268	VAL	N-CA-C	5.20	116.34	108.96
1	S	38	SER	N-CA-C	5.05	120.30	114.04
2	M	386	ILE	CB-CA-C	-5.04	104.58	111.33
2	M	562	GLN	CA-C-N	5.03	126.13	119.84
2	M	562	GLN	C-N-CA	5.03	126.13	119.84

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	Q	2234	0	2164	21	0
1	R	2203	0	2121	23	0
1	S	2242	0	2170	21	0
1	T	2220	0	2129	30	0
2	J	4528	0	4390	35	0
2	K	4526	0	4379	34	0
2	L	4526	0	4376	33	0
2	M	4512	0	4371	39	0
3	A	1310	0	1188	20	0
3	B	1242	0	1141	15	0
4	Q	8	0	0	0	0
4	R	8	0	0	0	0
4	S	8	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	T	8	0	0	0	0
5	Q	7	0	0	0	0
5	R	7	0	0	0	0
5	S	7	0	0	0	0
5	T	7	0	0	0	0
6	Q	8	0	0	0	0
6	R	8	0	0	0	0
6	S	8	0	0	0	0
6	T	8	0	0	0	0
7	J	7	0	0	0	0
7	K	7	0	0	0	0
7	L	7	0	0	0	0
7	M	7	0	0	0	0
8	J	1	0	0	0	0
8	K	1	0	0	0	0
8	L	1	0	0	0	0
8	M	1	0	0	0	0
9	J	1	0	0	0	0
9	K	1	0	0	0	0
9	L	1	0	0	0	0
9	M	1	0	0	0	0
10	J	1	0	0	0	0
10	K	1	0	0	0	0
10	L	1	0	0	0	0
10	M	1	0	0	0	0
11	A	43	0	30	5	0
11	B	43	0	30	0	0
12	A	6	0	0	1	0
12	B	6	0	0	0	0
12	J	112	0	0	6	0
12	K	108	0	0	3	0
12	L	111	0	0	1	0
12	M	113	0	0	7	0
12	Q	62	0	0	1	0
12	R	63	0	0	0	0
12	S	62	0	0	1	0
12	T	59	0	0	0	0
All	All	30463	0	28489	242	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 4.

All (242) 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)
3:B:78:ILE:HG13	3:B:79:TYR:H	1.21	1.02
3:B:78:ILE:HG13	3:B:79:TYR:N	1.77	0.95
2:J:3:THR:HB	12:J:717:HOH:O	1.70	0.91
3:B:139:MET:HE2	3:B:139:MET:HA	1.59	0.85
11:A:301:HEM:HBB2	11:A:301:HEM:HMB2	1.61	0.80
3:A:16:ALA:HB1	3:A:17:PRO:HD2	1.65	0.79
1:T:260:ARG:HH21	3:A:175:GLY:HA2	1.54	0.72
3:B:16:ALA:HB1	3:B:17:PRO:CD	2.25	0.67
1:S:186:ILE:HD11	1:S:228:ASN:HB3	1.74	0.67
3:A:139:MET:HE2	3:A:139:MET:HA	1.76	0.66
2:L:519:ARG:HG2	2:L:519:ARG:HH11	1.62	0.65
1:R:186:ILE:HD11	1:R:228:ASN:HB3	1.80	0.64
2:J:43:ILE:HD12	2:J:555:THR:HB	1.79	0.64
1:T:186:ILE:HD11	1:T:228:ASN:HB3	1.81	0.62
1:T:260:ARG:HH21	3:A:175:GLY:CA	2.13	0.61
3:B:16:ALA:HB1	3:B:17:PRO:HD2	1.83	0.60
2:M:36:ASN:HB2	2:M:45:ASN:HB3	1.83	0.59
1:Q:186:ILE:HD11	1:Q:228:ASN:HB3	1.84	0.59
3:A:41:LYS:HG3	12:A:405:HOH:O	2.03	0.59
3:B:78:ILE:CG1	3:B:79:TYR:N	2.61	0.58
2:J:275:MET:HE2	2:J:456:ALA:HA	1.85	0.57
2:K:56:LEU:HD22	2:K:75:ILE:HG13	1.85	0.57
2:L:530:VAL:HG12	2:L:531:PRO:HD2	1.86	0.57
1:T:133:PRO:HG2	1:T:136:LYS:HG2	1.87	0.57
1:S:187:HIS:CD2	1:S:193:ARG:NH1	2.73	0.56
2:L:36:ASN:HB2	2:L:45:ASN:HB3	1.88	0.55
2:M:517:ALA:HB3	2:M:525:LEU:HB3	1.89	0.55
2:K:354:TYR:HB3	12:K:774:HOH:O	2.07	0.55
1:R:268:VAL:C	1:R:278:THR:HG22	2.32	0.55
2:L:141:ASP:HB3	2:L:144:LYS:HD2	1.89	0.54
2:K:275:MET:HE2	2:K:456:ALA:HA	1.89	0.54
1:S:275:THR:HA	3:A:61:LEU:HD12	1.89	0.54
2:J:4:GLN:HA	2:J:12:ILE:O	2.07	0.54
2:L:135:LEU:HD13	2:L:187:TRP:CD1	2.43	0.54
2:J:56:LEU:HD22	2:J:75:ILE:HG13	1.89	0.54
2:M:424:ASP:O	2:M:428:VAL:HG23	2.08	0.53
2:K:36:ASN:HB2	2:K:45:ASN:HB3	1.90	0.53
2:K:130:ASP:HB3	2:K:568:ARG:HG2	1.89	0.53
1:S:193:ARG:HG3	1:T:194:ALA:HB2	1.90	0.53
1:Q:194:ALA:HB2	1:R:193:ARG:HG3	1.91	0.53
1:Q:165:THR:CG2	2:K:478:LYS:HE2	2.39	0.52
2:L:530:VAL:CG1	2:L:531:PRO:HD2	2.39	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:R:261:GLY:HA3	1:R:268:VAL:HB	1.92	0.52
1:T:159:ILE:HG22	1:T:163:MET:HE2	1.91	0.52
2:L:576:CYS:CB	12:L:808:HOH:O	2.57	0.52
2:M:56:LEU:HD22	2:M:75:ILE:HG13	1.92	0.52
2:K:535:ASN:HB3	2:K:548:TYR:CE1	2.44	0.52
2:L:130:ASP:HB3	2:L:568:ARG:HG2	1.92	0.52
1:T:280:ASP:O	1:T:284:LEU:HB2	2.10	0.52
2:K:62:GLY:O	2:K:522:LYS:HD2	2.09	0.52
1:T:262:SER:HB3	1:T:265:SER:OG	2.10	0.51
1:R:35:VAL:HA	1:R:39:LEU:HB2	1.91	0.51
1:S:171:ASP:OD1	2:K:480:GLY:HA3	2.11	0.51
1:T:298:ALA:HB2	3:A:200:TYR:CE1	2.46	0.51
2:M:43:ILE:HD12	2:M:555:THR:HB	1.92	0.50
2:M:31:MET:HB2	2:M:577:LEU:HG	1.93	0.50
2:J:36:ASN:HB2	2:J:45:ASN:HB3	1.92	0.50
2:L:60:LEU:HD11	2:L:72:VAL:HG13	1.94	0.50
2:L:56:LEU:HD22	2:L:75:ILE:HG13	1.94	0.50
1:T:260:ARG:NH2	3:A:175:GLY:HA2	2.24	0.50
2:M:430:SER:OG	2:M:457:HIS:CE1	2.65	0.50
2:J:56:LEU:HD22	2:J:75:ILE:CG1	2.41	0.50
2:L:31:MET:HB2	2:L:577:LEU:HG	1.93	0.50
1:S:232:SER:OG	2:M:249:VAL:HB	2.12	0.49
2:L:399:LYS:O	2:L:401:PRO:HD3	2.11	0.49
1:Q:183:GLY:HA3	2:K:246:SER:HB2	1.94	0.49
2:L:22:ASP:HB2	2:L:32:ARG:HG3	1.95	0.49
2:M:197:GLU:CD	2:M:197:GLU:H	2.20	0.49
2:M:435:MET:HE3	2:M:440:LEU:HB2	1.94	0.49
2:M:576:CYS:CB	12:M:812:HOH:O	2.60	0.48
1:Q:193:ARG:HG3	1:R:194:ALA:HB2	1.95	0.48
2:M:535:ASN:HB3	2:M:548:TYR:CE1	2.48	0.48
2:L:43:ILE:HD12	2:L:555:THR:HB	1.95	0.48
1:T:272:GLN:NE2	3:A:173:THR:HG23	2.28	0.48
2:M:365:PRO:HA	12:M:751:HOH:O	2.12	0.48
1:S:294:VAL:HG12	1:S:296:VAL:HG23	1.96	0.48
1:S:22:GLU:O	1:S:26:ARG:HG2	2.14	0.48
2:K:76:CYS:HB3	2:K:79:CYS:SG	2.53	0.48
2:L:306:ALA:HB2	2:L:506:GLU:H	1.79	0.48
1:R:226:THR:HG23	1:R:254:GLU:HG2	1.94	0.48
1:R:285:THR:O	1:R:289:VAL:HG23	2.14	0.47
2:J:306:ALA:HB2	2:J:506:GLU:H	1.80	0.47
3:B:170:PHE:CD2	3:B:180:ILE:HG12	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:519:ARG:HG2	2:L:519:ARG:NH1	2.25	0.47
2:J:393:GLU:HB3	12:J:784:HOH:O	2.14	0.47
2:L:65:PRO:HA	2:L:523:ILE:HD11	1.97	0.47
2:L:535:ASN:HB3	2:L:548:TYR:CE1	2.50	0.47
1:T:273:MET:HB2	1:T:273:MET:HE3	1.66	0.47
2:J:354:TYR:HB3	12:J:729:HOH:O	2.15	0.47
1:Q:218:LYS:HE2	1:Q:266:ARG:HH22	1.80	0.47
2:K:72:VAL:HG21	2:K:87:SER:HB2	1.96	0.47
2:M:72:VAL:HG21	2:M:87:SER:HB2	1.97	0.46
1:R:215:CYS:SG	1:R:217:TYR:HB2	2.55	0.46
2:K:63:ARG:O	2:K:523:ILE:HG13	2.16	0.46
2:K:509:ARG:HD2	12:K:805:HOH:O	2.13	0.46
1:T:275:THR:C	1:T:276:HIS:HD1	2.23	0.46
2:M:355:ARG:NH1	2:M:371:ASP:OD2	2.44	0.46
3:B:194:PHE:O	3:B:198:HIS:N	2.42	0.46
1:T:14:GLY:O	1:T:16:GLU:HG2	2.16	0.46
2:J:135:LEU:HD13	2:J:187:TRP:CD1	2.50	0.46
2:K:43:ILE:HD12	2:K:555:THR:HB	1.97	0.46
3:A:24:LEU:HD23	3:A:73:VAL:HG21	1.97	0.46
1:S:159:ILE:HG22	1:S:163:MET:HE2	1.97	0.46
1:T:7:ILE:HD12	1:T:8:PRO:HD2	1.98	0.46
2:M:480:GLY:HA3	1:Q:171:ASP:OD1	2.16	0.46
2:J:72:VAL:HG21	2:J:87:SER:HB2	1.97	0.46
2:L:492:PRO:HA	2:L:495:TRP:CD2	2.50	0.46
1:Q:7:ILE:HD12	1:Q:8:PRO:HD2	1.97	0.46
2:J:130:ASP:HB3	2:J:568:ARG:HG2	1.98	0.46
2:M:377:GLY:HA3	2:M:393:GLU:OE2	2.16	0.45
1:R:273:MET:HB3	1:R:273:MET:HE2	1.65	0.45
3:A:36:GLY:HA3	11:A:301:HEM:C4A	2.50	0.45
11:A:301:HEM:HBB2	11:A:301:HEM:CMB	2.37	0.45
1:Q:159:ILE:HG22	1:Q:163:MET:HE2	1.96	0.45
2:M:399:LYS:O	2:M:401:PRO:HD3	2.16	0.45
1:S:7:ILE:HD12	1:S:8:PRO:HD2	1.99	0.45
3:B:190:LEU:O	3:B:193:ALA:HB3	2.17	0.45
1:S:174:ARG:HH11	2:K:488:GLU:CD	2.25	0.45
1:T:297:HIS:C	1:T:297:HIS:CD2	2.94	0.45
2:M:455:ARG:HD2	12:M:768:HOH:O	2.16	0.45
2:J:380:LYS:HB2	2:J:387:GLN:HB2	1.99	0.45
1:S:3:ASN:CG	1:S:4:LYS:H	2.24	0.45
2:M:354:TYR:HB3	12:M:753:HOH:O	2.16	0.45
2:J:399:LYS:O	2:J:401:PRO:HD3	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:T:272:GLN:HE22	3:A:173:THR:HG23	1.81	0.45
1:Q:125:ARG:HA	1:Q:126:PRO:HA	1.78	0.45
2:J:509:ARG:HD2	12:J:806:HOH:O	2.16	0.45
3:B:139:MET:HA	3:B:139:MET:CE	2.37	0.45
2:M:40:GLN:O	2:M:41:ASN:HB2	2.17	0.45
2:J:197:GLU:CD	2:J:197:GLU:H	2.24	0.45
2:K:47:VAL:HG13	2:K:395:TYR:CE2	2.52	0.45
1:T:260:ARG:NH1	1:T:271:PRO:HG2	2.32	0.45
2:M:492:PRO:HA	2:M:495:TRP:CD2	2.52	0.44
3:A:16:ALA:HB1	3:A:17:PRO:CD	2.37	0.44
2:M:60:LEU:HD11	2:M:72:VAL:HG13	2.00	0.44
2:M:511:ALA:O	2:M:531:PRO:HD3	2.18	0.44
1:T:79:PRO:HD2	1:T:132:THR:O	2.16	0.44
2:J:445:ILE:O	2:J:450:GLY:HA3	2.18	0.44
2:J:492:PRO:HA	2:J:495:TRP:CD2	2.52	0.44
1:S:194:ALA:HB2	1:T:193:ARG:HG3	1.99	0.44
2:J:530:VAL:CG1	2:J:531:PRO:HD2	2.47	0.44
2:K:492:PRO:HA	2:K:495:TRP:CD2	2.52	0.44
2:L:561:GLU:CD	2:L:561:GLU:H	2.24	0.44
2:K:300:CYS:HA	2:K:326:VAL:O	2.17	0.44
2:K:517:ALA:HB3	2:K:525:LEU:HB3	1.99	0.44
2:J:576:CYS:CB	12:J:806:HOH:O	2.66	0.44
2:K:22:ASP:HB2	2:K:32:ARG:HG3	2.00	0.44
1:T:273:MET:O	1:T:277:SER:HB2	2.18	0.44
1:Q:274:GLY:O	3:B:61:LEU:HD12	2.17	0.44
2:K:31:MET:HB2	2:K:577:LEU:HG	2.00	0.44
2:M:488:GLU:OE1	1:Q:174:ARG:NH1	2.51	0.43
2:K:4:GLN:HA	2:K:12:ILE:O	2.17	0.43
2:K:380:LYS:HB2	2:K:388:GLN:HB2	1.99	0.43
3:B:161:ALA:N	3:B:162:PRO:HD2	2.34	0.43
2:M:20:VAL:HG22	2:M:34:GLU:HG2	2.00	0.43
1:S:174:ARG:NH1	2:K:488:GLU:CD	2.76	0.43
2:L:445:ILE:O	2:L:450:GLY:HA3	2.19	0.43
1:Q:189:LYS:O	1:Q:231:SER:HB2	2.19	0.43
1:R:217:TYR:OH	3:B:39:ILE:O	2.25	0.43
2:M:106:ILE:O	2:M:110:MET:HG2	2.19	0.43
1:R:272:GLN:NE2	3:B:173:THR:HG23	2.33	0.43
2:K:79:CYS:O	2:K:82:VAL:HG12	2.19	0.43
2:K:203:PHE:O	2:K:206:TYR:HB3	2.19	0.43
1:S:118:TRP:CD2	1:S:255:ASN:HB2	2.54	0.43
1:T:225:THR:HG21	3:A:178:MET:HA	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:M:312:ASN:HA	12:M:743:HOH:O	2.19	0.43
1:Q:13:HIS:HD2	12:Q:510:HOH:O	2.02	0.42
2:J:455:ARG:HD2	12:J:773:HOH:O	2.18	0.42
1:S:197:ASP:OD2	1:T:197:ASP:OD2	2.37	0.42
2:K:306:ALA:HB2	2:K:506:GLU:H	1.84	0.42
1:T:193:ARG:HD2	1:T:193:ARG:HA	1.82	0.42
2:J:60:LEU:HD11	2:J:72:VAL:CG1	2.49	0.42
2:K:197:GLU:H	2:K:197:GLU:CD	2.27	0.42
1:S:297:HIS:O	1:S:298:ALA:HB3	2.19	0.42
2:L:512:LEU:HD13	2:L:530:VAL:HG13	2.01	0.42
2:M:445:ILE:O	2:M:450:GLY:HA3	2.20	0.42
1:R:77:GLY:HA2	1:R:115:CYS:HB3	2.02	0.42
1:R:260:ARG:NH2	3:B:179:ASP:OD2	2.48	0.42
1:Q:165:THR:HG23	2:K:478:LYS:HE2	2.01	0.42
1:Q:174:ARG:HG2	1:Q:174:ARG:HH21	1.85	0.42
1:T:37:LEU:HD22	2:M:173:PHE:CD1	2.55	0.42
2:M:276:ILE:HB	2:M:277:PRO:HD3	2.01	0.42
2:M:509:ARG:HD2	12:M:812:HOH:O	2.19	0.42
1:Q:29:HIS:HA	1:Q:30:PRO:HA	1.91	0.42
1:Q:79:PRO:HD2	1:Q:132:THR:O	2.20	0.42
3:A:170:PHE:CD2	3:A:180:ILE:HG12	2.55	0.42
1:S:297:HIS:HD2	12:S:523:HOH:O	2.01	0.42
1:R:78:ASN:HB2	1:R:132:THR:O	2.20	0.42
1:R:185:ARG:HD3	1:R:227:TYR:CZ	2.55	0.42
3:A:24:LEU:HD12	3:A:24:LEU:HA	1.83	0.42
2:L:390:ASN:OD1	2:L:392:GLN:HB2	2.19	0.41
3:A:161:ALA:N	3:A:162:PRO:HD2	2.35	0.41
1:T:25:ILE:HA	1:T:32:ALA:HB2	2.02	0.41
2:M:300:CYS:HA	2:M:326:VAL:O	2.20	0.41
2:J:28:GLU:HB3	2:J:576:CYS:HA	2.02	0.41
1:R:22:GLU:O	1:R:26:ARG:HG2	2.20	0.41
2:L:56:LEU:HD22	2:L:75:ILE:CG1	2.50	0.41
2:J:65:PRO:HB3	2:J:91:ILE:HD12	2.01	0.41
2:J:448:THR:HG23	2:J:572:SER:HB3	2.03	0.41
1:T:218:LYS:HE3	3:A:42:PRO:HB2	2.02	0.41
2:M:22:ASP:HA	2:M:23:PRO:HA	1.94	0.41
2:M:130:ASP:HB3	2:M:568:ARG:HG2	2.02	0.41
2:M:306:ALA:HB2	2:M:506:GLU:H	1.85	0.41
3:A:63:HIS:CE1	11:A:301:HEM:NA	2.88	0.41
2:L:341:ASP:HA	2:L:342:PRO:HD3	1.94	0.41
1:T:125:ARG:HA	1:T:126:PRO:HA	1.66	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:K:194:LEU:HB3	2:K:198:ALA:HB3	2.01	0.41
2:J:535:ASN:HB3	2:J:548:TYR:CE1	2.55	0.41
1:R:133:PRO:HG2	1:R:136:LYS:HG2	2.02	0.41
1:R:79:PRO:HD2	1:R:132:THR:O	2.21	0.41
2:K:576:CYS:CB	12:K:805:HOH:O	2.69	0.41
3:A:36:GLY:HA3	11:A:301:HEM:C3A	2.55	0.41
2:L:28:GLU:HB3	2:L:576:CYS:HA	2.03	0.41
2:L:106:ILE:O	2:L:110:MET:HG2	2.21	0.41
2:L:540:ASP:HB2	2:L:541:PRO:CD	2.51	0.41
2:J:536:ALA:HB2	2:J:548:TYR:CE2	2.56	0.41
2:L:4:GLN:HA	2:L:12:ILE:O	2.20	0.41
2:K:135:LEU:HD13	2:K:187:TRP:CD1	2.56	0.41
1:S:25:ILE:O	2:L:213:GLN:NE2	2.54	0.40
2:M:65:PRO:HA	2:M:523:ILE:HD11	2.03	0.40
2:J:465:LYS:HD3	2:J:465:LYS:HA	1.95	0.40
2:K:321:MET:HA	2:K:322:PRO:HD3	1.94	0.40
1:T:22:GLU:O	1:T:26:ARG:HG2	2.21	0.40
1:T:226:THR:HG23	1:T:254:GLU:HG2	2.04	0.40
2:M:393:GLU:HB3	12:M:783:HOH:O	2.21	0.40
1:Q:189:LYS:HD2	1:R:191:TYR:CD2	2.56	0.40
2:J:144:LYS:HB3	2:J:197:GLU:HG2	2.03	0.40
2:J:478:LYS:HE2	1:R:165:THR:CG2	2.51	0.40
2:J:574:ASP:N	2:J:575:PRO:HD3	2.36	0.40
2:M:413:LEU:CD1	2:M:445:ILE:HG23	2.51	0.40
1:Q:193:ARG:HD2	1:Q:193:ARG:HA	1.92	0.40
2:J:31:MET:HB2	2:J:577:LEU:HG	2.02	0.40
1:R:187:HIS:CB	1:R:224:PRO:HA	2.51	0.40
1:R:193:ARG:HA	1:R:193:ARG:HD2	1.92	0.40
1:S:125:ARG:HA	1:S:126:PRO:HA	1.71	0.40
2:L:20:VAL:HG22	2:L:34:GLU:HG2	2.02	0.40
2:L:60:LEU:HA	2:L:63:ARG:HG3	2.03	0.40
1:Q:7:ILE:HA	1:Q:8:PRO:HD3	1.93	0.40
2:J:76:CYS:HB3	2:J:79:CYS:SG	2.62	0.40
2:J:341:ASP:HA	2:J:342:PRO:HD3	1.93	0.40
1:S:233:THR:O	1:S:234:ARG:HB2	2.20	0.40
2:L:465:LYS:HD3	2:L:465:LYS:HA	1.95	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	Q	293/335 (88%)	280 (96%)	13 (4%)	0	100	100
1	R	289/335 (86%)	275 (95%)	13 (4%)	1 (0%)	36	55
1	S	294/335 (88%)	280 (95%)	14 (5%)	0	100	100
1	T	294/335 (88%)	279 (95%)	14 (5%)	1 (0%)	36	55
2	J	579/582 (100%)	558 (96%)	21 (4%)	0	100	100
2	K	579/582 (100%)	562 (97%)	17 (3%)	0	100	100
2	L	580/582 (100%)	564 (97%)	16 (3%)	0	100	100
2	M	579/582 (100%)	560 (97%)	18 (3%)	1 (0%)	43	63
3	A	166/235 (71%)	140 (84%)	24 (14%)	2 (1%)	10	20
3	B	154/235 (66%)	135 (88%)	15 (10%)	4 (3%)	4	7
All	All	3807/4138 (92%)	3633 (95%)	165 (4%)	9 (0%)	43	63

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	86	ARG
2	M	488	GLU
3	A	154	HIS
3	B	201	MET
1	R	272	GLN
3	B	158	ALA
1	T	298	ALA
3	B	128	ILE
3	B	199	VAL

5.3.2 Protein sidechains

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

resolution.

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	Q	235/273 (86%)	230 (98%)	5 (2%)	47	74
1	R	231/273 (85%)	228 (99%)	3 (1%)	61	82
1	S	236/273 (86%)	232 (98%)	4 (2%)	53	78
1	T	230/273 (84%)	226 (98%)	4 (2%)	53	78
2	J	474/481 (98%)	466 (98%)	8 (2%)	53	78
2	K	474/481 (98%)	470 (99%)	4 (1%)	73	88
2	L	475/481 (99%)	470 (99%)	5 (1%)	65	84
2	M	471/481 (98%)	463 (98%)	8 (2%)	53	78
3	A	114/203 (56%)	105 (92%)	9 (8%)	11	24
3	B	111/203 (55%)	105 (95%)	6 (5%)	20	41
All	All	3051/3422 (89%)	2995 (98%)	56 (2%)	51	77

All (56) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	S	16	GLU
1	S	37	LEU
1	S	193	ARG
1	S	296	VAL
2	L	3	THR
2	L	472	LYS
2	L	530	VAL
2	L	561	GLU
2	L	579	CYS
1	T	232	SER
1	T	284	LEU
1	T	289	VAL
1	T	294	VAL
2	M	333	ASN
2	M	380	LYS
2	M	388	GLN
2	M	433	ARG
2	M	439	ASN
2	M	445	ILE
2	M	487	THR

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Mol	Chain	Res	Type
2	M	579	CYS
1	Q	16	GLU
1	Q	232	SER
1	Q	266	ARG
1	Q	275	THR
1	Q	296	VAL
2	J	37	ILE
2	J	51	THR
2	J	174	VAL
2	J	269	ASP
2	J	429	GLU
2	J	439	ASN
2	J	525	LEU
2	J	561	GLU
1	R	232	SER
1	R	277	SER
1	R	287	LEU
2	K	269	ASP
2	K	317	LYS
2	K	445	ILE
2	K	579	CYS
3	A	20	ILE
3	A	49	GLU
3	A	74	LEU
3	A	87	TYR
3	A	108	GLU
3	A	159	ILE
3	A	173	THR
3	A	188	MET
3	A	203	LEU
3	B	49	GLU
3	B	72	VAL
3	B	78	ILE
3	B	103	GLN
3	B	159	ILE
3	B	173	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	S	68	ASN
2	L	125	GLN

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Mol	Chain	Res	Type
2	L	261	GLN
2	L	467	GLN
1	T	68	ASN
1	T	203	GLN
1	T	297	HIS
2	M	125	GLN
2	M	387	GLN
2	M	388	GLN
2	M	446	GLN
1	Q	68	ASN
2	J	150	GLN
2	J	261	GLN
2	J	385	ASN
2	J	457	HIS
2	J	467	GLN
2	J	544	GLN
2	J	562	GLN
1	R	68	ASN
1	R	276	HIS
2	K	125	GLN
2	K	150	GLN
2	K	213	GLN
2	K	460	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 30 ligands modelled in this entry, 12 are monoatomic - leaving 18 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	$\# Z > 2$
6	ER2	S	403	1	0,11,11	-	-	-		
11	HEM	A	301	3	50,50,50	1.72	14 (28%)	67,82,82	1.43	10 (14%)
5	F3S	T	402	1	0,9,9	-	-	-		
4	SF4	R	401	1	0,12,12	-	-	-		
7	FCO	M	601	12,2	0,6,6	-	-	-		
11	HEM	B	301	3	50,50,50	1.63	9 (18%)	67,82,82	1.82	14 (20%)
7	FCO	K	601	12,2	0,6,6	-	-	-		
7	FCO	J	601	12,2	0,6,6	-	-	-		
5	F3S	R	402	1	0,9,9	-	-	-		
6	ER2	Q	403	1	0,11,11	-	-	-		
7	FCO	L	601	12,2	0,6,6	-	-	-		
4	SF4	S	401	1	0,12,12	-	-	-		
5	F3S	Q	402	1	0,9,9	-	-	-		
4	SF4	T	401	1	0,12,12	-	-	-		
6	ER2	T	403	1	0,11,11	-	-	-		
5	F3S	S	402	1	0,9,9	-	-	-		
6	ER2	R	403	1	0,11,11	-	-	-		
4	SF4	Q	401	1	0,12,12	-	-	-		

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
6	ER2	S	403	1	-	-	0/4/4/4
11	HEM	A	301	3	-	8/14/54/54	-
5	F3S	T	402	1	-	-	0/3/3/3
4	SF4	R	401	1	-	-	0/6/5/5
11	HEM	B	301	3	-	9/14/54/54	-
5	F3S	R	402	1	-	-	0/3/3/3
6	ER2	Q	403	1	-	-	0/4/4/4
4	SF4	S	401	1	-	-	0/6/5/5
5	F3S	Q	402	1	-	-	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	SF4	T	401	1	-	-	0/6/5/5
6	ER2	T	403	1	-	-	0/4/4/4
5	F3S	S	402	1	-	-	0/3/3/3
6	ER2	R	403	1	-	-	0/4/4/4
4	SF4	Q	401	1	-	-	0/6/5/5

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
11	A	301	HEM	FE-NB	5.15	2.10	1.94
11	B	301	HEM	FE-NB	4.50	2.08	1.94
11	B	301	HEM	C1B-NB	-4.41	1.32	1.40
11	B	301	HEM	FE-NC	3.96	2.08	1.95
11	A	301	HEM	C1B-NB	-3.90	1.33	1.40
11	A	301	HEM	C4B-NB	-3.27	1.32	1.38
11	A	301	HEM	C4D-ND	-3.25	1.34	1.40
11	A	301	HEM	FE-NC	3.25	2.05	1.95
11	B	301	HEM	FE-ND	-3.14	1.85	1.94
11	A	301	HEM	FE-ND	-3.12	1.85	1.94
11	B	301	HEM	C4D-ND	-3.08	1.34	1.40
11	B	301	HEM	C1D-ND	-2.84	1.33	1.38
11	A	301	HEM	C1C-C2C	-2.64	1.40	1.45
11	B	301	HEM	C1C-C2C	-2.63	1.40	1.45
11	A	301	HEM	C3C-C4C	-2.60	1.41	1.46
11	A	301	HEM	C4C-NC	-2.46	1.35	1.39
11	A	301	HEM	C1A-C2A	-2.24	1.40	1.44
11	B	301	HEM	C1B-C2B	-2.16	1.40	1.44
11	A	301	HEM	C1D-ND	-2.15	1.34	1.38
11	A	301	HEM	O2D-CGD	-2.14	1.23	1.30
11	A	301	HEM	O2A-CGA	-2.14	1.23	1.30
11	A	301	HEM	C1C-NC	-2.04	1.35	1.39
11	B	301	HEM	O2D-CGD	-2.03	1.24	1.30

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
11	B	301	HEM	CHC-C4B-NB	6.32	131.22	124.42
11	A	301	HEM	CHC-C4B-NB	4.55	129.32	124.42
11	B	301	HEM	CHD-C1D-ND	3.72	128.43	124.42
11	A	301	HEM	CHD-C1D-ND	3.59	128.28	124.42
11	B	301	HEM	CHA-C4D-ND	3.32	128.47	124.37
11	B	301	HEM	CHD-C1D-C2D	-3.27	119.87	125.03

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
11	B	301	HEM	C1B-NB-C4B	3.16	108.95	105.21
11	A	301	HEM	CHD-C1D-C2D	-3.03	120.25	125.03
11	B	301	HEM	O2D-CGD-O1D	-2.92	115.83	123.33
11	B	301	HEM	CHB-C1B-NB	2.84	127.88	124.37
11	A	301	HEM	CHA-C4D-ND	2.78	127.80	124.37
11	B	301	HEM	CHA-C4D-C3D	-2.74	120.17	125.23
11	B	301	HEM	O2D-CGD-CBD	2.72	122.58	114.00
11	A	301	HEM	C1B-NB-C4B	2.69	108.39	105.21
11	B	301	HEM	CMC-C2C-C1C	2.67	129.43	124.73
11	B	301	HEM	C3B-C4B-NB	-2.52	107.66	109.47
11	B	301	HEM	CAD-C3D-C2D	2.50	132.54	127.87
11	B	301	HEM	CHD-C4C-NC	2.50	127.17	124.45
11	A	301	HEM	CHD-C4C-NC	2.45	127.12	124.45
11	B	301	HEM	C1A-CHA-C4D	-2.30	120.83	126.25
11	A	301	HEM	O2A-CGA-O1A	-2.21	117.65	123.33
11	A	301	HEM	CHA-C4D-C3D	-2.18	121.21	125.23
11	A	301	HEM	C1A-CHA-C4D	-2.09	121.33	126.25
11	A	301	HEM	O2D-CGD-O1D	-2.06	118.04	123.33

There are no chirality outliers.

All (17) torsion outliers are listed below:

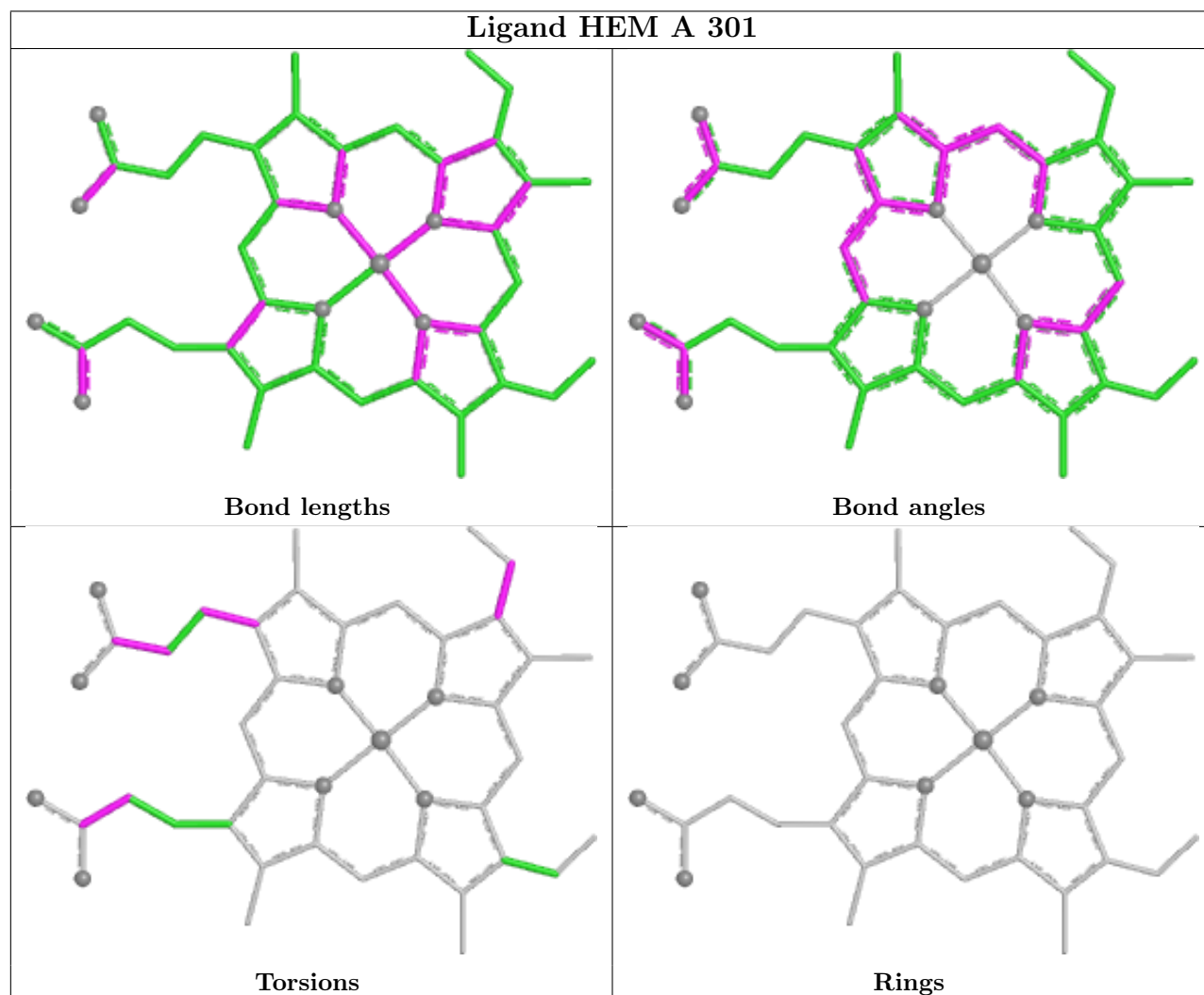
Mol	Chain	Res	Type	Atoms
11	B	301	HEM	C2D-C3D-CAD-CBD
11	B	301	HEM	C4D-C3D-CAD-CBD
11	A	301	HEM	C2C-C3C-CAC-CBC
11	B	301	HEM	C2C-C3C-CAC-CBC
11	A	301	HEM	C4D-C3D-CAD-CBD
11	A	301	HEM	C2D-C3D-CAD-CBD
11	B	301	HEM	C3D-CAD-CBD-CGD
11	A	301	HEM	C4C-C3C-CAC-CBC
11	B	301	HEM	C4C-C3C-CAC-CBC
11	A	301	HEM	CAD-CBD-CGD-O1D
11	B	301	HEM	CAD-CBD-CGD-O1D
11	B	301	HEM	CAD-CBD-CGD-O2D
11	A	301	HEM	CAD-CBD-CGD-O2D
11	A	301	HEM	CAA-CBA-CGA-O2A
11	A	301	HEM	CAA-CBA-CGA-O1A
11	B	301	HEM	CAA-CBA-CGA-O2A
11	B	301	HEM	CAA-CBA-CGA-O1A

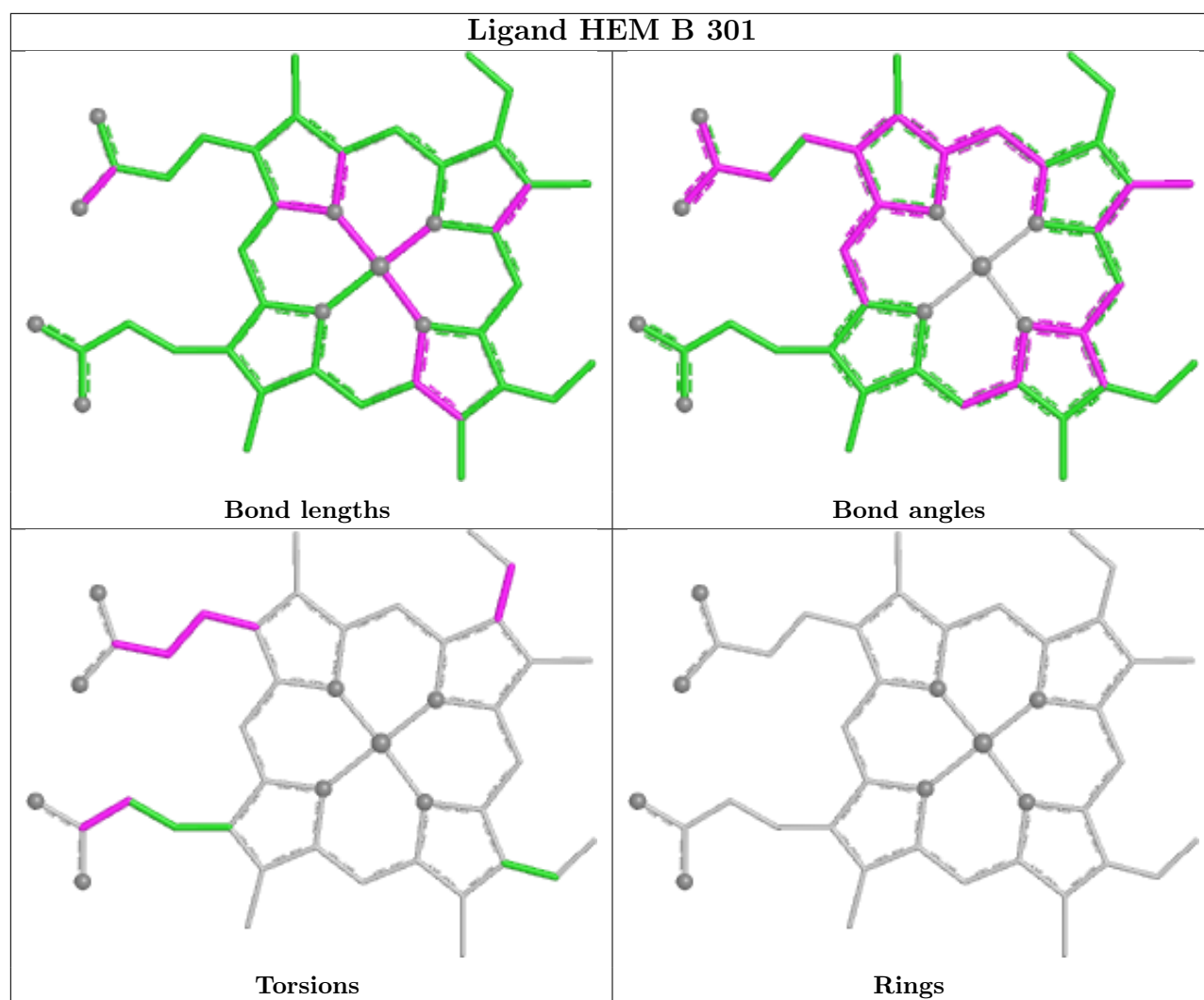
There are no ring outliers.

1 monomer is involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
11	A	301	HEM	5	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	Q	295/335 (88%)	-0.28	0 100 100	38, 48, 64, 87	0
1	R	291/335 (86%)	-0.13	1 (0%) 90 87	38, 52, 73, 101	0
1	S	296/335 (88%)	-0.35	0 100 100	33, 43, 62, 87	0
1	T	296/335 (88%)	-0.12	1 (0%) 90 87	35, 51, 76, 117	0
2	J	581/582 (99%)	-0.22	0 100 100	40, 52, 66, 78	0
2	K	581/582 (99%)	-0.30	0 100 100	37, 49, 65, 85	0
2	L	581/582 (99%)	-0.36	0 100 100	32, 46, 59, 76	1 (0%)
2	M	581/582 (99%)	-0.25	0 100 100	36, 52, 70, 85	0
3	A	172/235 (73%)	0.38	10 (5%) 29 25	40, 61, 102, 112	0
3	B	160/235 (68%)	0.42	9 (5%) 30 26	45, 62, 106, 132	0
All	All	3834/4138 (92%)	-0.20	21 (0%) 87 85	32, 50, 73, 132	1 (0%)

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	B	155	SER	4.5
3	A	154	HIS	4.0
3	B	114	PHE	3.7
3	A	155	SER	3.6
3	B	201	MET	3.4
3	A	14	PHE	3.3
3	B	202	ALA	3.0
3	A	102	TRP	2.9
3	A	113	LEU	2.9
3	A	100	SER	2.9
3	B	101	TRP	2.8
3	A	207	ILE	2.6
1	T	299	VAL	2.6

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Mol	Chain	Res	Type	RSRZ
3	B	157	TYR	2.6
3	B	102	TRP	2.5
3	B	82	PHE	2.4
3	A	157	TYR	2.4
1	R	294	VAL	2.2
3	B	133	MET	2.2
3	A	80	TRP	2.1
3	A	89	ARG	2.1

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [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
10	CL	K	604	1/1	0.75	0.11	43,43,43,43	1
10	CL	J	604	1/1	0.78	0.13	43,43,43,43	1
10	CL	L	604	1/1	0.90	0.09	41,41,41,41	1
10	CL	M	604	1/1	0.91	0.09	35,35,35,35	1
9	MG	L	603	1/1	0.96	0.05	42,42,42,42	0
11	HEM	A	301	43/43	0.97	0.08	47,50,54,59	0
6	ER2	R	403	8/8	0.98	0.03	44,47,51,52	0
5	F3S	S	402	7/7	0.98	0.04	38,40,40,45	0
9	MG	M	603	1/1	0.98	0.04	50,50,50,50	0
5	F3S	T	402	7/7	0.98	0.04	40,41,44,47	0
11	HEM	B	301	43/43	0.98	0.07	43,48,54,57	0
7	FCO	M	601	7/7	0.99	0.05	42,45,49,49	0
7	FCO	J	601	7/7	0.99	0.06	41,41,49,53	0
7	FCO	K	601	7/7	0.99	0.07	39,41,45,46	0
8	NI	J	602	1/1	0.99	0.04	43,43,43,43	0

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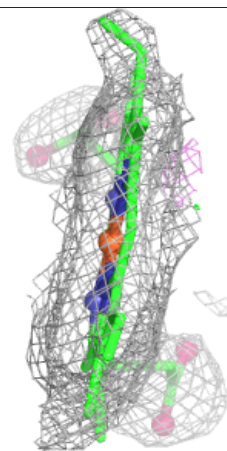
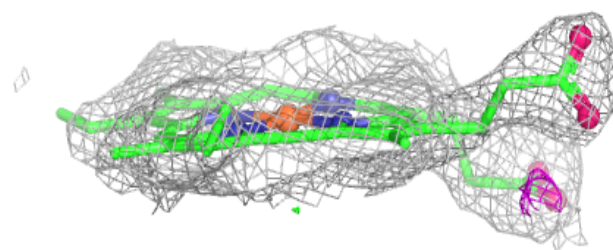
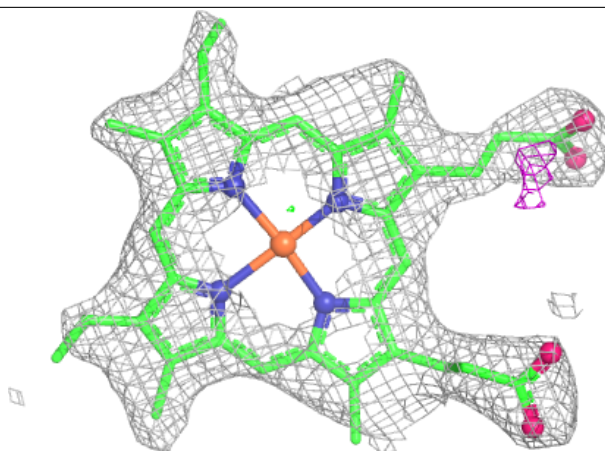
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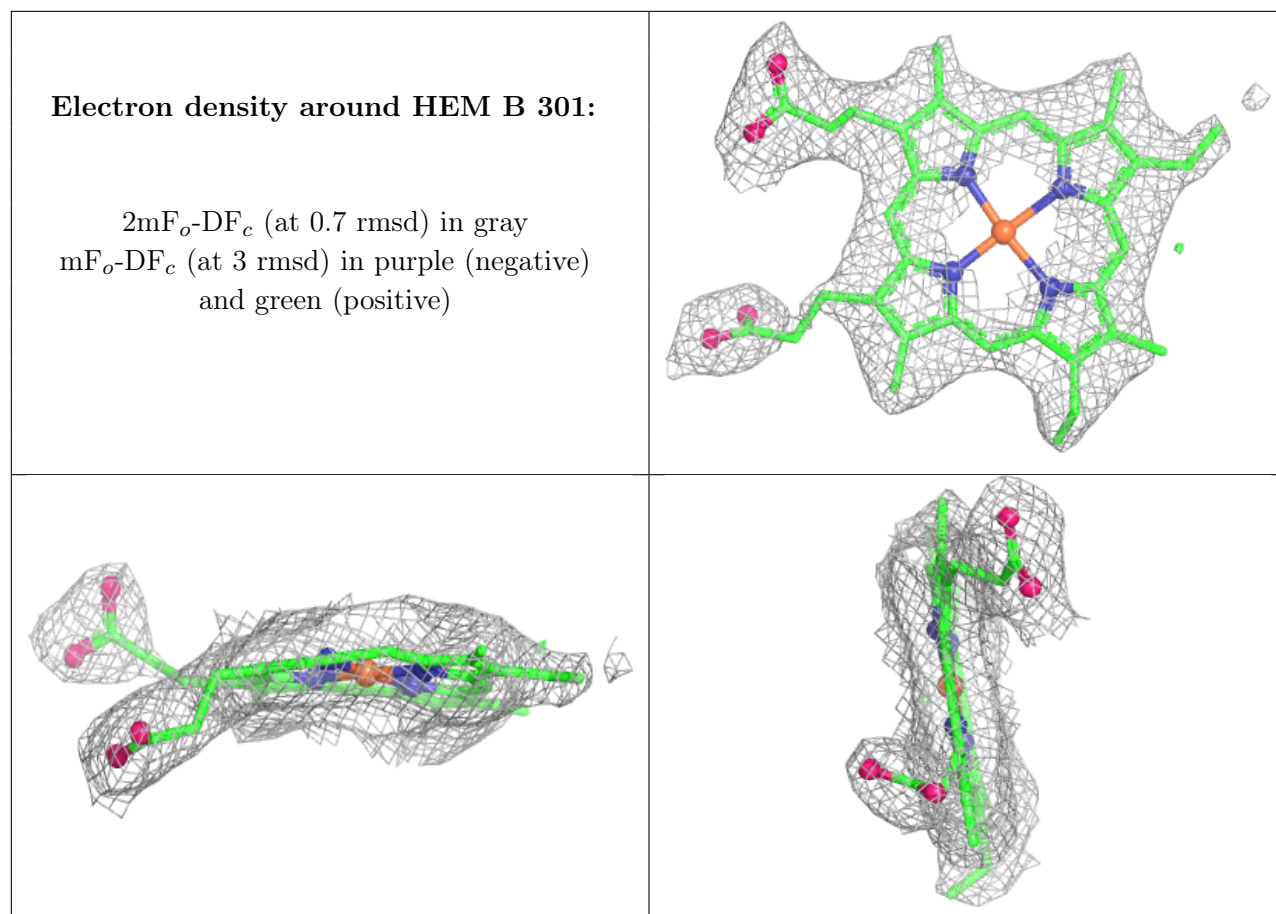
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	SF4	R	401	8/8	0.99	0.02	40,42,44,45	0
4	SF4	S	401	8/8	0.99	0.02	39,40,41,41	0
9	MG	J	603	1/1	0.99	0.05	44,44,44,44	0
9	MG	K	603	1/1	0.99	0.02	45,45,45,45	0
4	SF4	T	401	8/8	0.99	0.03	39,40,41,45	0
5	F3S	Q	402	7/7	0.99	0.04	39,40,41,42	0
6	ER2	S	403	8/8	0.99	0.03	39,40,42,45	0
6	ER2	T	403	8/8	0.99	0.03	42,43,47,49	0
6	ER2	Q	403	8/8	0.99	0.03	41,44,49,50	0
4	SF4	Q	401	8/8	0.99	0.03	41,44,46,48	0
5	F3S	R	402	7/7	1.00	0.02	39,41,42,42	0
8	NI	L	602	1/1	1.00	0.02	39,39,39,39	0
8	NI	M	602	1/1	1.00	0.04	44,44,44,44	0
7	FCO	L	601	7/7	1.00	0.04	39,39,46,50	0
8	NI	K	602	1/1	1.00	0.03	43,43,43,43	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around HEM A 301:

$2mF_o-DF_c$ (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.