



# wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 31, 2016 – 08:38 PM GMT

PDB ID : 1L9B  
Title : X-Ray Structure of the Cytochrome-c(2)-Photosynthetic Reaction Center Electron Transfer Complex from Rhodobacter sphaeroides in Type II Co-Crystals  
Authors : Axelrod, H.L.; Abresch, E.C.; Okamura, M.Y.; Yeh, A.P.; Rees, D.C.; Feher, G.  
Deposited on : 2002-03-22  
Resolution : 2.40 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

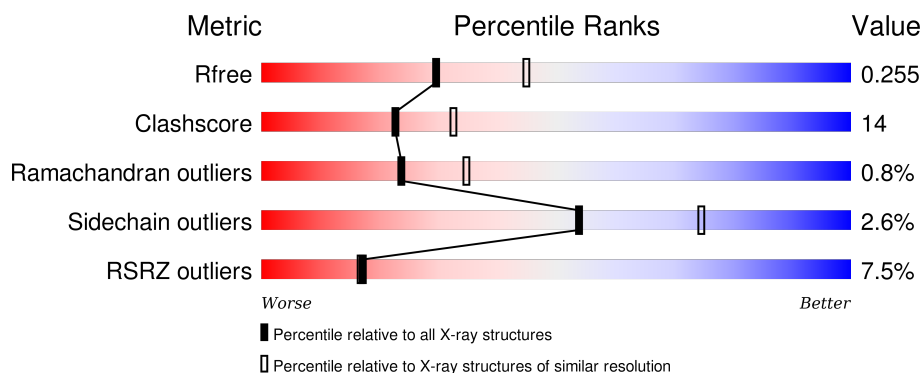
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.40 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))
$R_{free}$	91344	2919 (2.40-2.40)
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (2.40-2.40)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	L	281	<div> <div>6%</div> <div>74%</div> <div>25%</div> <div>.</div> </div>
2	M	307	<div> <div>6%</div> <div>67%</div> <div>19%</div> <div>.</div> <div>13%</div> </div>
3	H	260	<div> <div>11%</div> <div>67%</div> <div>26%</div> <div>.</div> <div>5%</div> </div>
4	C	124	<div> <div>4%</div> <div>72%</div> <div>27%</div> <div>.</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
12	LDA	H	1014	-	-	-	X
12	LDA	L	1017	-	-	-	X
12	LDA	L	1018	-	-	-	X
12	LDA	L	1020	-	-	-	X
12	LDA	M	1012	-	-	-	X
12	LDA	M	1013	-	-	-	X
12	LDA	M	1015	-	-	-	X
13	HTO	C	1022	-	-	X	X
13	HTO	L	1021	-	-	-	X
9	BPH	L	1005	X	-	-	-
9	BPH	M	1006	X	-	-	-

## 2 Entry composition

There are 14 unique types of molecules in this entry. The entry contains 8102 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 REACTION CENTER PROTEIN L CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	281	Total	C	N	O	S	0	0	0
			2232	1507	355	362	8			

- Molecule 2 is a protein called REACTION CENTER PROTEIN M CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	M	267	Total	C	N	O	S	0	0	0
			2150	1450	347	344	9			

- Molecule 3 is a protein called REACTION CENTER PROTEIN H CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	H	246	Total	C	N	O	S	0	0	0
			1871	1197	321	344	9			

- Molecule 4 is a protein called cytochrome c-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	C	124	Total	C	N	O	S	0	0	0
			949	595	166	184	4			

- Molecule 5 is FE (II) ION (three-letter code: FE2) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	M	1	Total	Fe	0	0
			1	1		

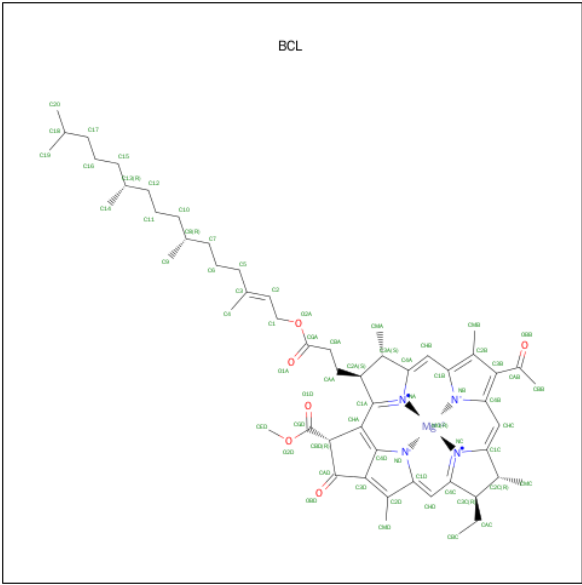
- Molecule 6 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	M	1	Total	Cl	0	0
			1	1		

- Molecule 7 is SODIUM ION (three-letter code: NA) (formula: Na).

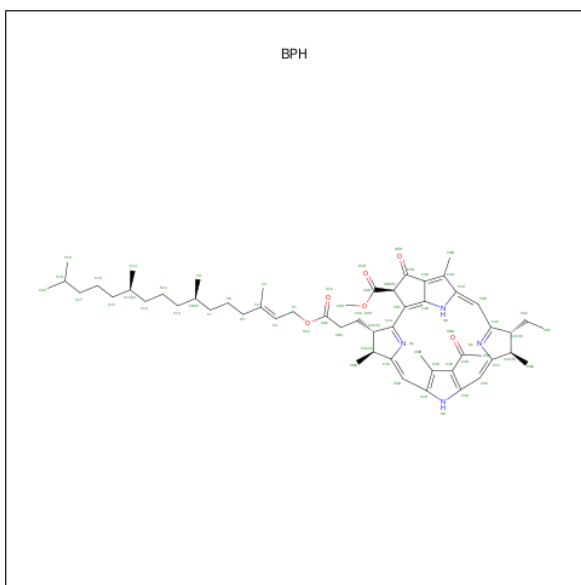
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	L	1	Total	Na	0	0
			1	1		

- Molecule 8 is BACTERIOCHLOROPHYLL A (three-letter code: BCL) (formula: C<sub>55</sub>H<sub>74</sub>MgN<sub>4</sub>O<sub>6</sub>).



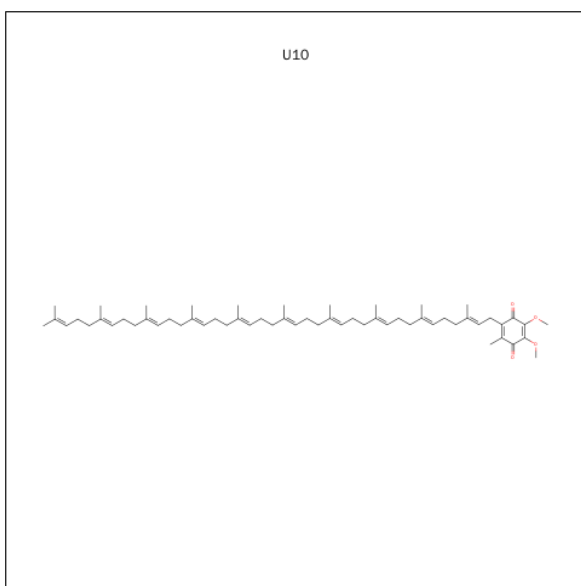
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
8	M	1	Total	C	Mg	N	O	0	0
			50	39	1	4	6		
8	L	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
8	M	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
8	L	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		

- Molecule 9 is BACTERIOPHEOPHYTIN A (three-letter code: BPH) (formula: C<sub>55</sub>H<sub>76</sub>N<sub>4</sub>O<sub>6</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
9	L	1	Total	C	N	O	0	0
			55	45	4	6		
9	M	1	Total	C	N	O	0	0
			65	55	4	6		

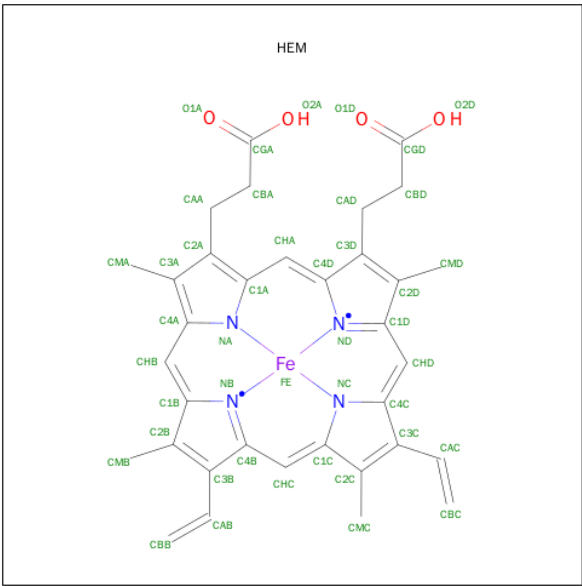
- Molecule 10 is UBIQUINONE-10 (three-letter code: U10) (formula:  $C_{59}H_{90}O_4$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
10	M	1	Total	C	O	0	0
			37	33	4		

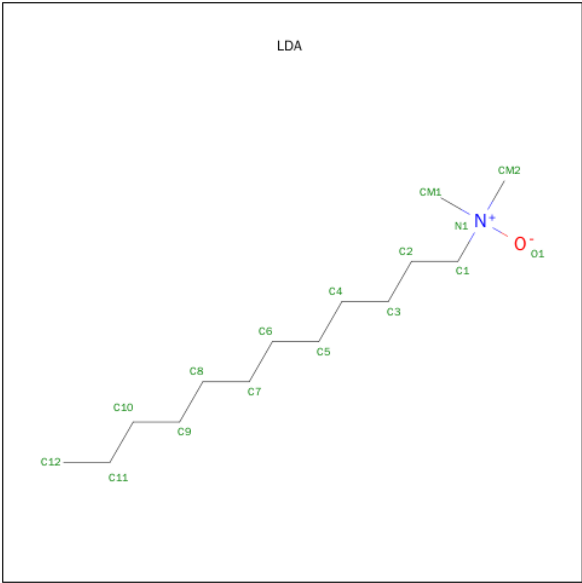
- Molecule 11 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (for-

mula: C<sub>34</sub>H<sub>32</sub>FeN<sub>4</sub>O<sub>4</sub>).



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

- Molecule 12 is LAURYL DIMETHYLAMINE-N-OXIDE (three-letter code: LDA) (formula: C<sub>14</sub>H<sub>31</sub>NO).



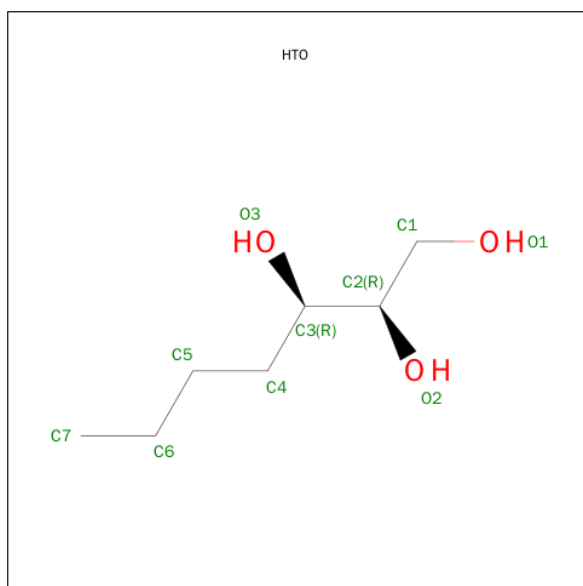
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
12	M	1	Total	C	N	O	0	0
			16	14	1	1		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
12	M	1	Total	C	N	O	0	0
			16	14	1	1		
12	H	1	Total	C	N	O	0	0
			16	14	1	1		
12	M	1	Total	C	N	O	0	0
			16	14	1	1		
12	L	1	Total	C	N	O	0	0
			16	14	1	1		
12	L	1	Total	C	N	O	0	0
			16	14	1	1		
12	L	1	Total	C	N	O	0	0
			16	14	1	1		
12	H	1	Total	C	N	O	0	0
			16	14	1	1		
12	L	1	Total	C	N	O	0	0
			16	14	1	1		

- Molecule 13 is HEPTANE-1,2,3-TRIOL (three-letter code: HTO) (formula: C<sub>7</sub>H<sub>16</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
13	L	1	Total	C	O	0	0
			10	7	3		
13	C	1	Total	C	O	0	0
			10	7	3		

- Molecule 14 is water.

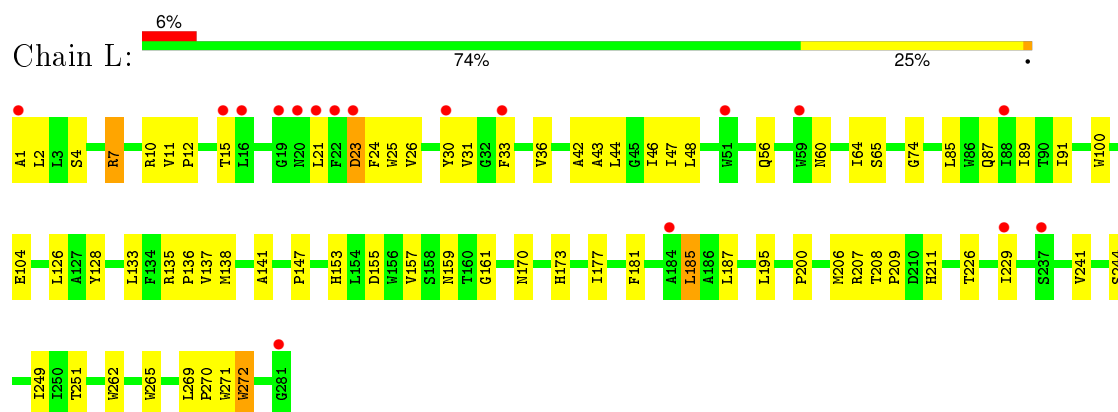


Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
14	C	58	Total 58	O 58	0	0
14	H	66	Total 66	O 66	0	0
14	L	91	Total 91	O 91	0	0
14	M	70	Total 70	O 70	0	0

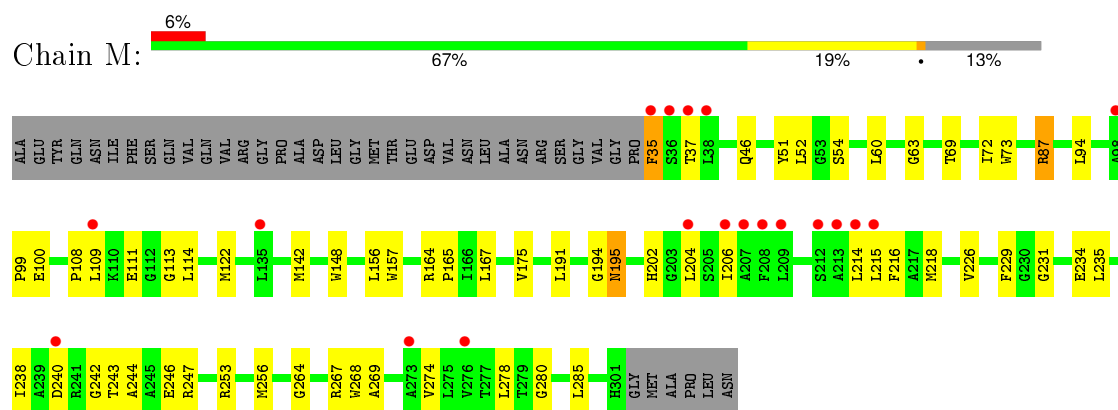
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

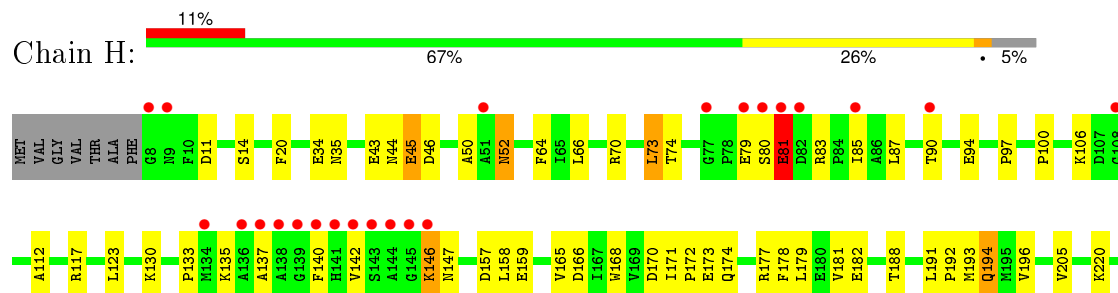
#### • Molecule 1: REACTION CENTER PROTEIN L CHAIN

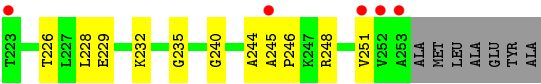


#### • Molecule 2: REACTION CENTER PROTEIN M CHAIN

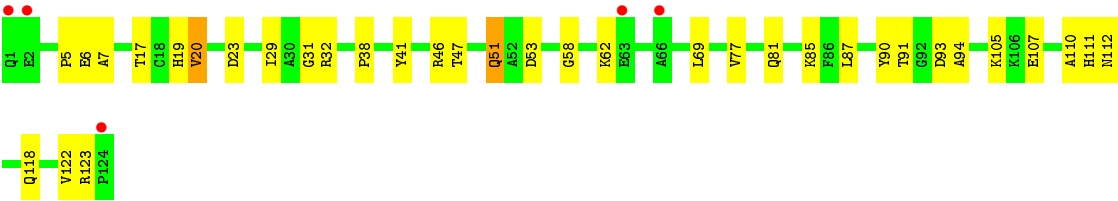


#### • Molecule 3: REACTION CENTER PROTEIN H CHAIN





● Molecule 4: cytochrome c-2



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	78.22Å 115.65Å 79.68Å 90.00° 110.29° 90.00°	Depositor
Resolution (Å)	45.73 – 2.40 45.73 – 2.40	Depositor EDS
% Data completeness (in resolution range)	96.7 (45.73-2.40) 96.7 (45.73-2.40)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.10	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.38 (at 2.39Å)	Xtriage
Refinement program	CNS	Depositor
R, $R_{free}$	0.220 , 0.264 0.213 , 0.255	Depositor DCC
$R_{free}$ test set	2523 reflections (5.28%)	DCC
Wilson B-factor (Å <sup>2</sup> )	49.1	Xtriage
Anisotropy	0.703	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 75.6	EDS
Estimated twinning fraction	0.022 for l,-k,h	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	1 of 50307 reflections (0.002%)	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	8102	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.65% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: BCL, LDA, CL, HTO, BPH, FE2, NA, U10, 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	L	0.37	0/2320	0.56	0/3175
2	M	0.39	0/2238	0.53	0/3057
3	H	0.31	0/1920	0.53	0/2612
4	C	0.35	0/969	0.59	0/1304
All	All	0.36	0/7447	0.55	0/10148

There are no bond length outliers.

There are no bond angle outliers.

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	L	2232	0	2187	58	0
2	M	2150	0	2073	56	0
3	H	1871	0	1877	64	0
4	C	949	0	916	32	0
5	M	1	0	0	0	0
6	M	1	0	0	0	0
7	L	1	0	0	0	0
8	L	132	0	148	10	0
8	M	116	0	115	15	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
9	L	55	0	53	5	0
9	M	65	0	74	3	0
10	M	37	0	47	0	0
11	C	43	0	30	0	0
12	H	32	0	62	1	0
12	L	64	0	124	0	0
12	M	48	0	93	5	0
13	C	10	0	16	6	0
13	L	10	0	16	0	0
14	C	58	0	0	8	0
14	H	66	0	0	3	0
14	L	91	0	0	10	0
14	M	70	0	0	10	0
All	All	8102	0	7831	213	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 14.

The worst 5 of 213 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:M:280:GLY:HA3	8:M:1003:BCL:HED3	1.40	1.02
2:M:280:GLY:HA3	8:M:1003:BCL:CED	1.93	0.98
2:M:157:TRP:HB2	8:M:1003:BCL:H62	1.49	0.93
2:M:218:MET:N	14:M:1085:HOH:O	2.03	0.91
2:M:268:TRP:HE1	3:H:35:ASN:ND2	1.69	0.90

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	L	279/281 (99%)	261 (94%)	15 (5%)	3 (1%)	17	25
2	M	265/307 (86%)	256 (97%)	9 (3%)	0	100	100
3	H	244/260 (94%)	218 (89%)	22 (9%)	4 (2%)	12	16
4	C	122/124 (98%)	112 (92%)	10 (8%)	0	100	100
All	All	910/972 (94%)	847 (93%)	56 (6%)	7 (1%)	24	35

5 of 7 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	H	81	GLU
3	H	220	LYS
3	H	83	ARG
1	L	7	ARG
1	L	23	ASP

### 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	L	220/220 (100%)	216 (98%)	4 (2%)	66	84
2	M	209/240 (87%)	202 (97%)	7 (3%)	45	66
3	H	199/208 (96%)	193 (97%)	6 (3%)	48	70
4	C	93/93 (100%)	91 (98%)	2 (2%)	60	79
All	All	721/761 (95%)	702 (97%)	19 (3%)	54	74

5 of 19 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	M	191	LEU
2	M	216	PHE
3	H	146	LYS
2	M	156	LEU
3	H	194	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 11 such sidechains are listed below:

Mol	Chain	Res	Type
3	H	35	ASN
3	H	52	ASN
4	C	51	GLN
2	M	300	ASN
4	C	16	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

### 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

### 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

### 5.6 Ligand geometry ⓘ

Of 22 ligands modelled in this entry, 3 are monoatomic - leaving 19 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	$\# Z  > 2$
11	HEM	C	1009	4	30,50,50	2.43	12 (40%)	24,82,82	3.12	13 (54%)
13	HTO	C	1022	-	9,9,9	1.18	1 (11%)	8,10,10	0.74	0
12	LDA	H	1014	-	15,15,15	3.91	2 (13%)	16,17,17	2.60	4 (25%)
12	LDA	H	1019	-	15,15,15	3.87	2 (13%)	16,17,17	2.60	4 (25%)
8	BCL	L	1002	1	53,74,74	1.14	3 (5%)	57,115,115	1.84	15 (26%)
8	BCL	L	1004	1	53,74,74	1.13	4 (7%)	57,115,115	1.68	14 (24%)



Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
9	BPH	L	1005	-	54,60,70	1.44	8 (14%)	61,89,101	2.36	17 (27%)
12	LDA	L	1016	-	15,15,15	3.95	2 (13%)	16,17,17	2.54	4 (25%)
12	LDA	L	1017	-	15,15,15	3.97	2 (13%)	16,17,17	2.69	4 (25%)
12	LDA	L	1018	-	15,15,15	3.93	2 (13%)	16,17,17	2.63	4 (25%)
12	LDA	L	1020	-	15,15,15	3.90	2 (13%)	16,17,17	2.60	4 (25%)
13	HTO	L	1021	-	9,9,9	1.34	1 (11%)	8,10,10	0.57	0
8	BCL	M	1001	2	37,58,74	1.39	3 (8%)	39,95,115	2.02	12 (30%)
8	BCL	M	1003	2	53,74,74	1.15	5 (9%)	57,115,115	2.08	18 (31%)
9	BPH	M	1006	-	64,70,70	1.48	10 (15%)	73,101,101	2.24	17 (23%)
10	U10	M	1008	-	37,37,63	2.04	10 (27%)	44,47,79	1.77	8 (18%)
12	LDA	M	1012	-	15,15,15	3.95	2 (13%)	16,17,17	2.55	4 (25%)
12	LDA	M	1013	-	15,15,15	3.78	2 (13%)	16,17,17	2.69	4 (25%)
12	LDA	M	1015	-	15,15,15	3.96	2 (13%)	16,17,17	2.62	4 (25%)

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
11	HEM	C	1009	4	-	0/10/54/54	0/0/8/8
13	HTO	C	1022	-	-	0/10/10/10	0/0/0/0
12	LDA	H	1014	-	-	0/13/13/13	0/0/0/0
12	LDA	H	1019	-	-	0/13/13/13	0/0/0/0
8	BCL	L	1002	1	-	0/37/137/137	0/0/9/9
8	BCL	L	1004	1	-	0/37/137/137	0/0/9/9
9	BPH	L	1005	-	1/1/16/22	0/42/93/105	0/1/6/6
12	LDA	L	1016	-	-	0/13/13/13	0/0/0/0
12	LDA	L	1017	-	-	0/13/13/13	0/0/0/0
12	LDA	L	1018	-	-	0/13/13/13	0/0/0/0
12	LDA	L	1020	-	-	0/13/13/13	0/0/0/0
13	HTO	L	1021	-	-	0/10/10/10	0/0/0/0
8	BCL	M	1001	2	-	0/18/118/137	0/0/9/9
8	BCL	M	1003	2	-	0/37/137/137	0/0/9/9
9	BPH	M	1006	-	2/2/18/22	1/54/105/105	0/1/6/6
10	U10	M	1008	-	-	0/32/56/87	0/1/1/1
12	LDA	M	1012	-	-	0/13/13/13	0/0/0/0
12	LDA	M	1013	-	-	0/13/13/13	0/0/0/0
12	LDA	M	1015	-	-	0/13/13/13	0/0/0/0

The worst 5 of 75 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
12	M	1012	LDA	O1-N1	-14.55	1.25	1.39
12	L	1016	LDA	O1-N1	-14.55	1.25	1.39
12	L	1017	LDA	O1-N1	-14.54	1.25	1.39
12	L	1018	LDA	O1-N1	-14.51	1.25	1.39
12	M	1015	LDA	O1-N1	-14.47	1.25	1.39

The worst 5 of 150 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
12	M	1013	LDA	CM2-N1-CM1	-8.17	99.61	108.83
12	L	1017	LDA	CM2-N1-CM1	-8.03	99.78	108.83
12	L	1018	LDA	CM2-N1-CM1	-7.78	100.05	108.83
12	M	1015	LDA	CM2-N1-CM1	-7.77	100.07	108.83
12	L	1020	LDA	CM2-N1-CM1	-7.61	100.25	108.83

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
9	M	1006	BPH	C8
9	M	1006	BPH	C13
9	L	1005	BPH	C8

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
9	M	1006	BPH	C1-C2-C3-C4

There are no ring outliers.

11 monomers are involved in 43 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
13	C	1022	HTO	6	0
12	H	1019	LDA	1	0
8	L	1002	BCL	8	0
8	L	1004	BCL	5	0
9	L	1005	BPH	5	0
8	M	1001	BCL	7	0
8	M	1003	BCL	11	0
9	M	1006	BPH	3	0
12	M	1012	LDA	3	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
12	M	1013	LDA	1	0
12	M	1015	LDA	1	0

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	L	281/281 (100%)	0.52	17 (6%) 25 25	36, 56, 75, 85	0
2	M	267/307 (86%)	0.58	19 (7%) 19 19	34, 52, 71, 86	0
3	H	246/260 (94%)	0.73	28 (11%) 7 7	50, 73, 113, 139	0
4	C	124/124 (100%)	0.42	5 (4%) 42 43	38, 58, 82, 103	0
All	All	918/972 (94%)	0.58	69 (7%) 17 17	34, 60, 88, 139	0

The worst 5 of 69 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	H	138	ALA	9.6
3	H	253	ALA	9.3
3	H	141	HIS	6.5
3	H	142	VAL	6.3
3	H	252	VAL	6.0

### 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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
12	LDA	L	1018	16/16	0.63	0.52	22.07	94,95,102,102	0
13	HTO	C	1022	10/10	0.62	0.48	12.76	77,81,84,86	0
12	LDA	M	1015	16/16	0.67	0.33	8.82	84,88,90,92	0
12	LDA	M	1013	16/16	0.71	0.47	6.18	68,83,96,97	0
12	LDA	H	1014	16/16	0.61	0.41	6.03	132,133,134,134	0
12	LDA	L	1020	16/16	0.54	0.37	5.14	82,87,93,93	0
13	HTO	L	1021	10/10	0.68	0.36	4.60	97,103,104,104	0
12	LDA	L	1017	16/16	0.78	0.34	4.23	101,105,114,114	0
12	LDA	M	1012	16/16	0.75	0.32	4.01	75,78,87,87	0
9	BPH	M	1006	65/65	0.90	0.26	1.64	42,53,77,81	0
10	U10	M	1008	37/63	0.93	0.26	1.47	41,54,87,87	0
8	BCL	L	1004	66/66	0.94	0.28	1.09	36,48,75,77	0
8	BCL	M	1003	66/66	0.91	0.25	1.03	23,32,78,80	0
8	BCL	L	1002	66/66	0.92	0.23	1.00	17,34,69,74	0
9	BPH	L	1005	55/65	0.90	0.23	0.77	38,46,78,78	0
8	BCL	M	1001	50/66	0.90	0.21	0.70	43,47,71,73	0
11	HEM	C	1009	43/43	0.97	0.16	-0.29	38,42,46,50	0
5	FE2	M	1007	1/1	1.00	0.17	-2.14	45,45,45,45	0
7	NA	L	1011	1/1	0.59	0.21	-	81,81,81,81	0
12	LDA	L	1016	16/16	0.75	0.25	-	120,122,124,125	0
12	LDA	H	1019	16/16	0.69	0.40	-	115,117,120,121	0
6	CL	M	1010	1/1	0.92	0.30	-	72,72,72,72	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.