



wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 31, 2016 – 09:53 PM GMT

PDB ID : 1R2C
Title : PHOTOSYNTHETIC REACTION CENTER BLASTOCHLORIS VIRIDIS (ATCC)
Authors : Baxter, R.H.; Ponomarenko, N.; Pahl, R.; Srajer, V.; Moffat, K.; Norris, J.R.
Deposited on : 2003-09-26
Resolution : 2.86 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
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

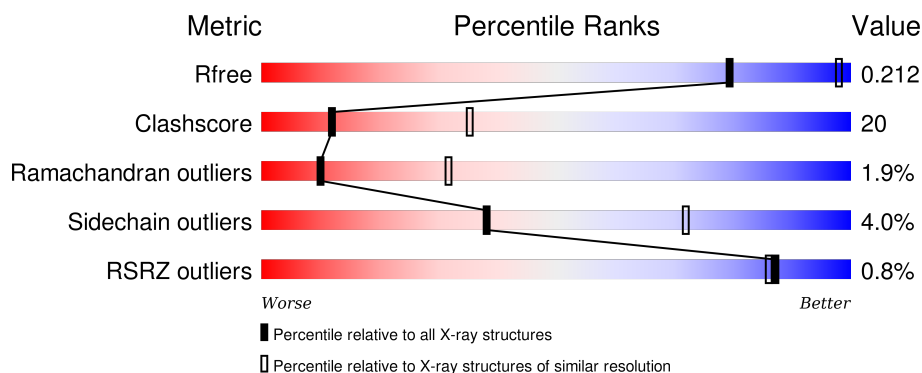
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.86 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	2228 (2.90-2.82)
Clashscore	102246	2499 (2.90-2.82)
Ramachandran outliers	100387	2439 (2.90-2.82)
Sidechain outliers	100360	2442 (2.90-2.82)
RSRZ outliers	91569	2236 (2.90-2.82)

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 ≥ 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	C	336	<div> <div>2%</div> <div>67%</div> <div>31%</div> <div>•</div> </div>
2	L	273	<div> <div>67%</div> <div>31%</div> <div>•</div> </div>
3	M	323	<div> <div>65%</div> <div>33%</div> <div>•</div> </div>
4	H	258	<div> <div>2%</div> <div>53%</div> <div>42%</div> <div>5%</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
10	MQ7	M	501	-	-	-	X
11	UQ2	L	502	-	-	-	X
12	NS5	M	600	-	-	-	X
13	LDA	H	703	-	-	-	X
13	LDA	L	705	-	-	-	X
13	LDA	M	702	-	-	-	X
13	LDA	M	704	-	-	-	X
13	LDA	M	706	-	-	-	X
8	BCB	L	304	X	-	-	-
8	BCB	M	1303	X	-	-	-
9	BPB	M	401	X	-	X	X

2 Entry composition

There are 13 unique types of molecules in this entry. The entry contains 10140 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 Photosynthetic reaction center cytochrome C subunit precursor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	332	Total	C	N	O	S	41	0	0
			2602	1640	466	478	18			

- Molecule 2 is a protein called Reaction center protein L chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	273	Total	C	N	O	S	10	0	0
			2171	1459	350	355	7			

- Molecule 3 is a protein called Reaction center protein M chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	M	323	Total	C	N	O	S	15	0	0
			2555	1702	419	423	11			

- Molecule 4 is a protein called Reaction center protein H chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	H	258	Total	C	N	O	S	98	0	0
			2018	1292	344	380	2			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
H	1	FME	MET	MODIFIED RESIDUE	UNP P06008

- 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 SULFATE ION (three-letter code: SO4) (formula: O₄S).



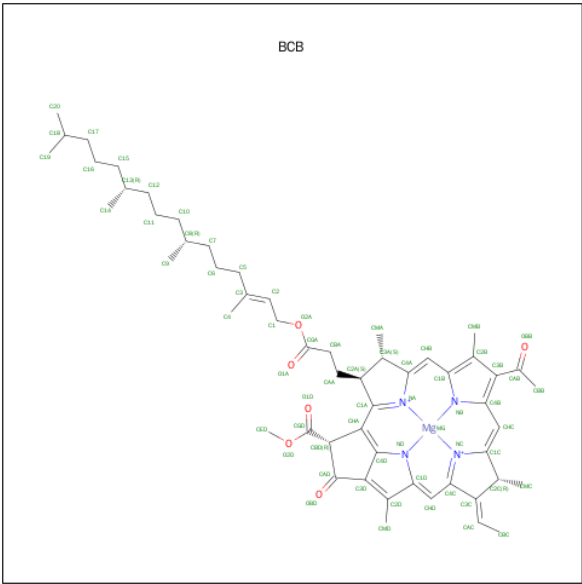
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	H	1	Total	O	S	0	0
			5	4	1		
6	M	1	Total	O	S	0	0
			5	4	1		
6	M	1	Total	O	S	0	0
			5	4	1		
6	M	1	Total	O	S	0	0
			5	4	1		

- Molecule 7 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: C₃₄H₃₂FeN₄O₄).



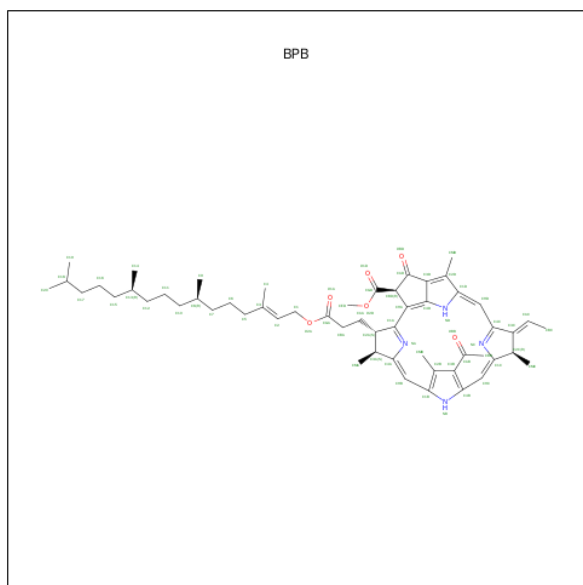
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	C	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
7	C	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
7	C	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
7	C	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 8 is BACTERIOCHLOROPHYLL B (three-letter code: BCB) (formula: $C_{55}H_{72}MgN_4O_6$).



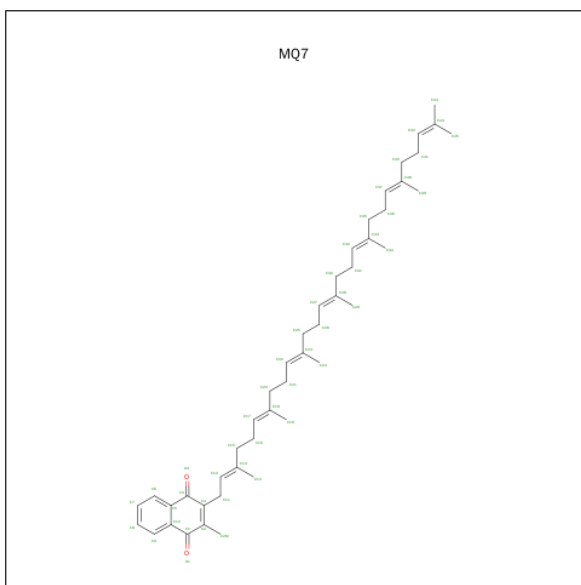
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	M	1	Total	C	Mg	N	O	
			66	55	1	4	6	
8	L	1	Total	C	Mg	N	O	
			66	55	1	4	6	
8	M	1	Total	C	Mg	N	O	
			66	55	1	4	6	
8	L	1	Total	C	Mg	N	O	
			66	55	1	4	6	

- Molecule 9 is BACTERIOPHEOPHYTIN B (three-letter code: BPB) (formula: $C_{55}H_{74}N_4O_6$).



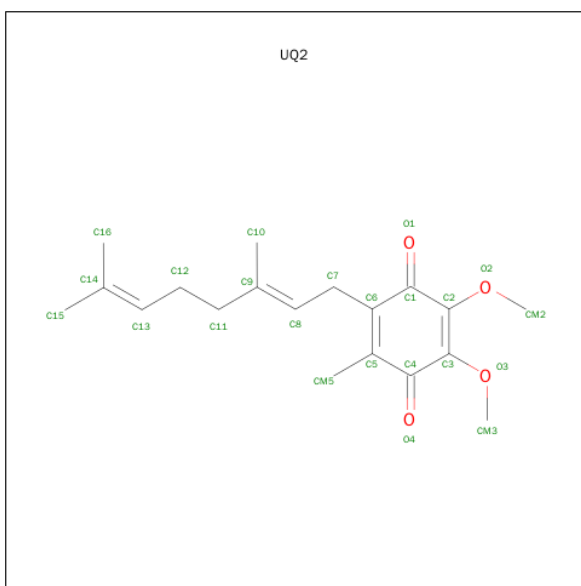
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
9	M	1	Total	C	N	O		
			65	55	4	6	0	0
9	L	1	Total	C	N	O		
			65	55	4	6	0	0

- Molecule 10 is MENAQUINONE-7 (three-letter code: MQ7) (formula: $C_{46}H_{64}O_2$).



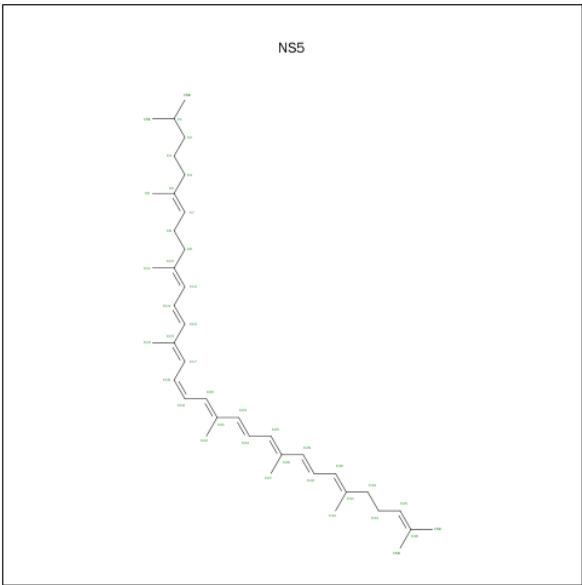
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
10	M	1	Total	C	O	0	0
			48	46	2		

- Molecule 11 is UBIQUINONE-2 (three-letter code: UQ2) (formula: $C_{19}H_{26}O_4$).



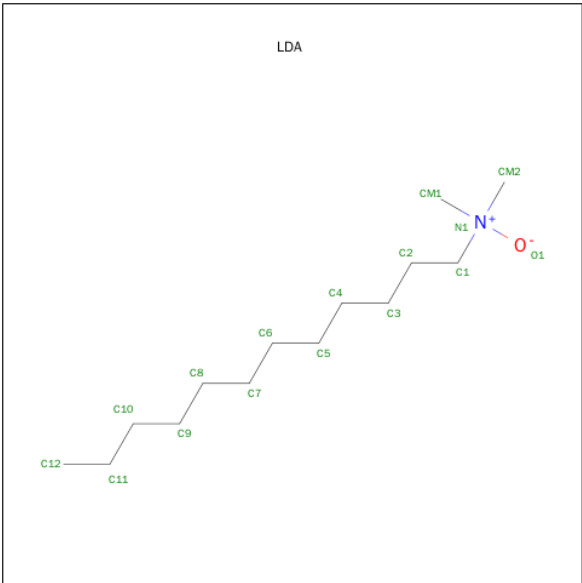
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
11	L	1	Total	C	O	0	0
			23	19	4		

- Molecule 12 is 15-CIS-1,2-DIHYDRONEUROSPORENE (three-letter code: NS5) (formula: $C_{40}H_{60}$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
12	M	1	Total	C			11	0
			40	40				

- Molecule 13 is LAURYL DIMETHYLAMINE-N-OXIDE (three-letter code: LDA) (formula: C₁₄H₃₁NO).



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

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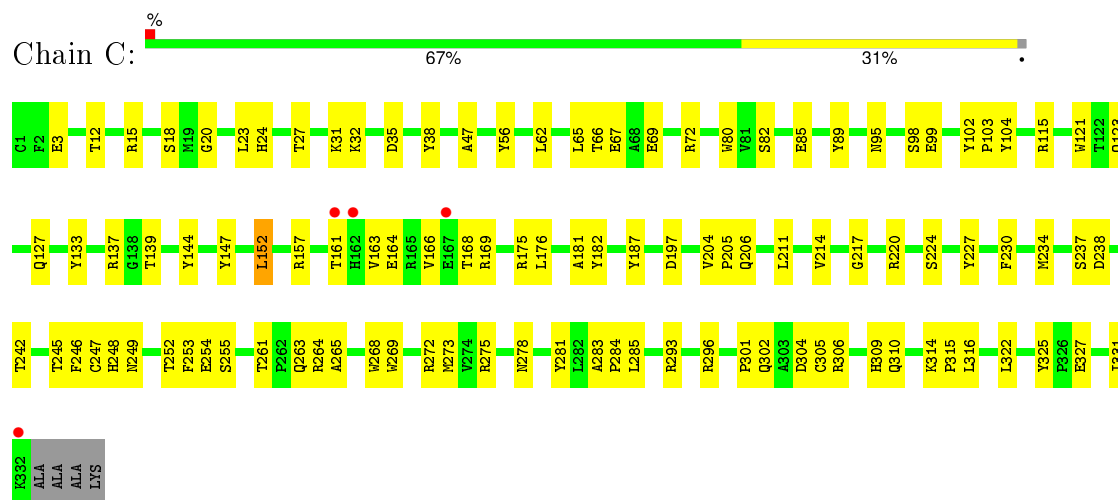
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
13	H	1	Total	C	N	O	0	0
			16	14	1	1		
13	M	1	Total	C	N	O	3	0
			16	14	1	1		
13	L	1	Total	C	N	O	3	0
			16	14	1	1		
13	M	1	Total	C	N	O	0	0
			16	14	1	1		

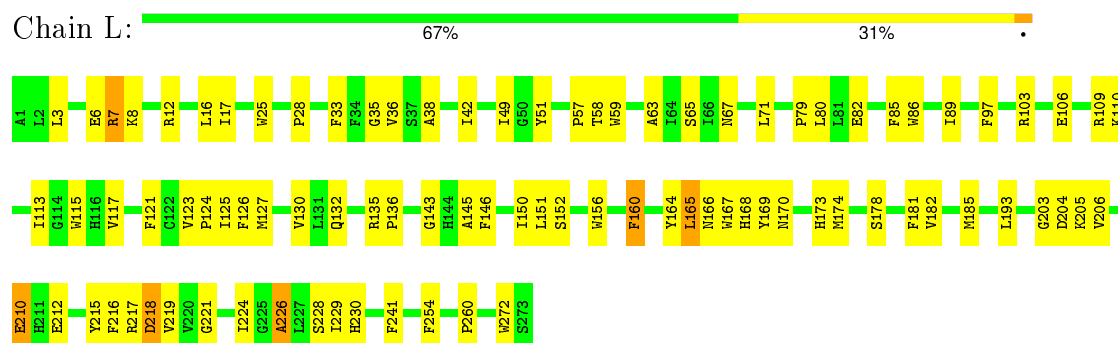
3 Residue-property plots

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 ($\text{RSRZ} > 2$). 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.

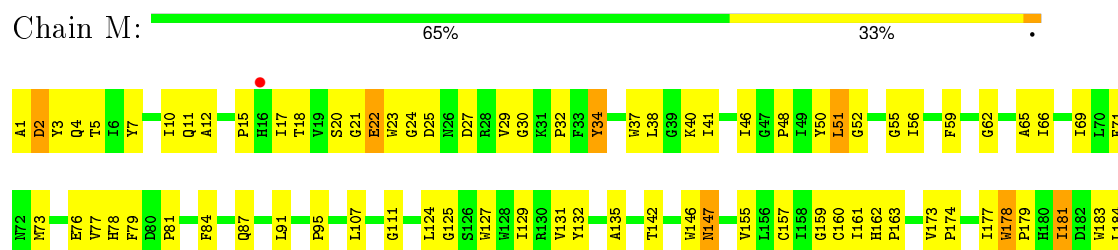
- Molecule 1: Photosynthetic reaction center cytochrome C subunit precursor

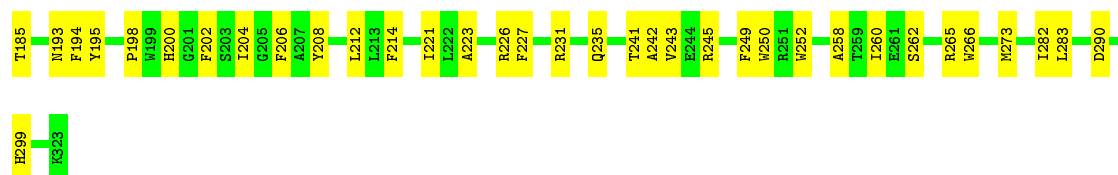


- Molecule 2: Reaction center protein L chain

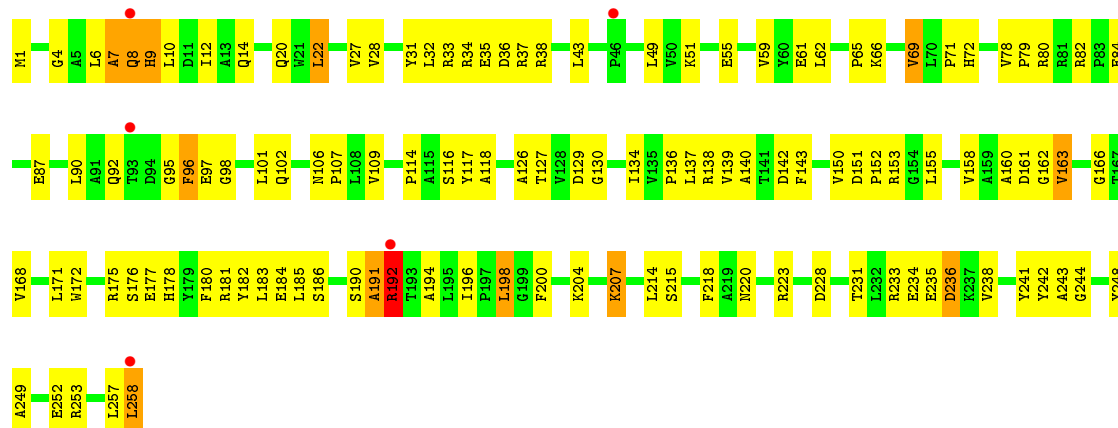


- Molecule 3: Reaction center protein M chain





• Molecule 4: Reaction center protein H chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	223.50 Å 223.50 Å 112.50 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.83 – 2.86 48.83 – 2.86	Depositor EDS
% Data completeness (in resolution range)	97.3 (48.83-2.86) 97.4 (48.83-2.86)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.90 (at 2.86 Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.202 , 0.228 0.190 , 0.212	Depositor DCC
R_{free} test set	6311 reflections (10.89%)	DCC
Wilson B-factor (Å ²)	59.3	Xtriage
Anisotropy	0.312	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 56.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 64248 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	10140	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.10% 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, 0.375 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: LDA, BPB, BCB, FE2, SO4, MQ7, HEM, FME, UQ2, NS5

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	C	0.35	0/2669	0.62	1/3637 (0.0%)
2	L	0.41	0/2259	0.60	0/3084
3	M	0.42	0/2659	0.58	0/3637
4	H	0.35	0/2055	0.62	0/2807
All	All	0.38	0/9642	0.61	1/13165 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	249	ASN	N-CA-C	-6.50	93.44	111.00

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	C	2602	0	2579	74	0
2	L	2171	0	2098	90	0
3	M	2555	0	2452	112	0
4	H	2018	0	2020	113	0
5	M	1	0	0	0	0
6	H	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	M	15	0	0	0	0
7	C	172	0	120	7	0
8	L	132	0	144	22	0
8	M	132	0	143	29	0
9	L	65	0	74	12	0
9	M	65	0	74	22	0
10	M	48	0	64	0	0
11	L	23	0	26	4	0
12	M	40	0	60	5	0
13	H	16	0	31	4	0
13	L	16	0	31	2	0
13	M	64	0	124	11	0
All	All	10140	0	10040	393	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 20.

The worst 5 of 393 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)
4:H:152:PRO:HA	4:H:155:LEU:HD12	1.31	1.05
9:M:401:BPB:H14	9:M:401:BPB:H17	1.39	1.03
3:M:32:PRO:HG3	3:M:48:PRO:HD3	1.39	1.02
8:L:302:BCB:HBA1	8:L:302:BCB:HBD	1.46	0.94
9:L:402:BPB:HBBA	3:M:208:TYR:HB3	1.47	0.94

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	C	330/336 (98%)	299 (91%)	30 (9%)	1 (0%)	46	76
2	L	271/273 (99%)	251 (93%)	17 (6%)	3 (1%)	17	47
3	M	321/323 (99%)	299 (93%)	19 (6%)	3 (1%)	21	52
4	H	256/258 (99%)	216 (84%)	25 (10%)	15 (6%)	2	5
All	All	1178/1190 (99%)	1065 (90%)	91 (8%)	22 (2%)	10	32

5 of 22 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	L	57	PRO
4	H	7	ALA
4	H	8	GLN
4	H	9	HIS
4	H	96	PHE

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	C	281/282 (100%)	276 (98%)	5 (2%)	66	89
2	L	218/218 (100%)	209 (96%)	9 (4%)	37	71
3	M	249/249 (100%)	239 (96%)	10 (4%)	38	72
4	H	212/212 (100%)	198 (93%)	14 (7%)	21	48
All	All	960/961 (100%)	922 (96%)	38 (4%)	38	72

5 of 38 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	M	51	LEU
3	M	194	PHE
4	H	236	ASP
3	M	178	TRP
3	M	214	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 16 such sidechains are listed below:

Mol	Chain	Res	Type
2	L	239	ASN
3	M	16	HIS
4	H	72	HIS
2	L	213	ASN
4	H	102	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is 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	FME	H	1	4	8,9,10	0.62	0	6,9,11	4.40	3 (50%)

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	FME	H	1	4	-	0/6/9/11	0/0/0/0

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	H	1	FME	CA-N-CN	-7.21	111.74	122.82

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	H	1	FME	O1-CN-N	-6.95	114.75	124.76
4	H	1	FME	CE-SD-CG	-3.31	89.07	100.37

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	H	1	FME	1	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 24 ligands modelled in this entry, 1 is monoatomic - leaving 23 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
7	HEM	C	1201	1	30,50,50	2.59	9 (30%)	24,82,82	2.78	8 (33%)
7	HEM	C	1202	1	30,50,50	2.58	8 (26%)	24,82,82	3.38	8 (33%)
7	HEM	C	1203	1	30,50,50	2.64	9 (30%)	24,82,82	2.98	9 (37%)
7	HEM	C	1204	1	30,50,50	2.65	9 (30%)	24,82,82	3.56	12 (50%)
13	LDA	H	703	-	15,15,15	4.88	3 (20%)	16,17,17	0.55	0
6	SO4	H	801	-	4,4,4	0.78	0	6,6,6	0.36	0
8	BCB	L	302	2	56,74,74	1.57	7 (12%)	57,115,115	2.34	14 (24%)
8	BCB	L	304	2	56,74,74	1.55	6 (10%)	57,115,115	2.15	10 (17%)
9	BPB	L	402	-	63,70,70	1.28	6 (9%)	63,101,101	2.54	13 (20%)
11	UQ2	L	502	-	23,23,23	1.78	5 (21%)	28,31,31	1.26	3 (10%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
13	LDA	L	705	-	15,15,15	4.15	2 (13%)	16,17,17	0.71	0
8	BCB	M	1301	3	56,74,74	1.51	6 (10%)	57,115,115	1.83	7 (12%)
8	BCB	M	1303	3	56,74,74	1.60	7 (12%)	57,115,115	1.92	11 (19%)
9	BPB	M	401	-	63,70,70	1.43	7 (11%)	63,101,101	2.08	12 (19%)
10	MQ7	M	501	-	49,49,49	1.78	13 (26%)	62,63,63	1.58	10 (16%)
12	NS5	M	600	-	39,39,39	0.77	0	44,46,46	1.04	4 (9%)
13	LDA	M	701	-	15,15,15	4.72	3 (20%)	16,17,17	0.81	0
13	LDA	M	702	-	15,15,15	4.27	1 (6%)	16,17,17	0.57	0
13	LDA	M	704	-	15,15,15	4.82	4 (26%)	16,17,17	0.62	0
13	LDA	M	706	-	15,15,15	4.62	2 (13%)	16,17,17	0.57	0
6	SO4	M	802	-	4,4,4	1.11	0	6,6,6	0.44	0
6	SO4	M	803	-	4,4,4	1.02	0	6,6,6	0.07	0
6	SO4	M	804	-	4,4,4	1.09	0	6,6,6	0.27	0

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
7	HEM	C	1201	1	-	0/10/54/54	0/0/8/8
7	HEM	C	1202	1	-	0/10/54/54	0/0/8/8
7	HEM	C	1203	1	-	0/10/54/54	0/0/8/8
7	HEM	C	1204	1	-	0/10/54/54	0/0/8/8
13	LDA	H	703	-	-	0/13/13/13	0/0/0/0
6	SO4	H	801	-	-	0/0/0/0	0/0/0/0
8	BCB	L	302	2	-	0/37/137/137	0/0/9/9
8	BCB	L	304	2	1/1/21/26	0/37/137/137	0/0/9/9
9	BPB	L	402	-	-	0/46/105/105	0/1/6/6
11	UQ2	L	502	-	-	0/15/39/39	0/1/1/1
13	LDA	L	705	-	-	0/13/13/13	0/0/0/0
8	BCB	M	1301	3	-	0/37/137/137	0/0/9/9
8	BCB	M	1303	3	1/1/21/26	0/37/137/137	0/0/9/9
9	BPB	M	401	-	1/1/18/23	0/46/105/105	0/1/6/6
10	MQ7	M	501	-	-	0/41/61/61	0/2/2/2
12	NS5	M	600	-	-	0/43/43/43	0/0/0/0
13	LDA	M	701	-	-	0/13/13/13	0/0/0/0
13	LDA	M	702	-	-	0/13/13/13	0/0/0/0
13	LDA	M	704	-	-	0/13/13/13	0/0/0/0
13	LDA	M	706	-	-	0/13/13/13	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	SO4	M	802	-	-	0/0/0/0	0/0/0/0
6	SO4	M	803	-	-	0/0/0/0	0/0/0/0
6	SO4	M	804	-	-	0/0/0/0	0/0/0/0

The worst 5 of 107 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
13	H	703	LDA	O1-N1	-18.39	1.22	1.39
13	M	704	LDA	O1-N1	-18.09	1.22	1.39
13	M	701	LDA	O1-N1	-17.74	1.22	1.39
13	M	706	LDA	O1-N1	-17.50	1.22	1.39
13	M	702	LDA	O1-N1	-16.32	1.24	1.39

The worst 5 of 121 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	C	1204	HEM	C3C-CAC-CBC	-10.52	108.33	124.46
7	C	1202	HEM	C3C-CAC-CBC	-9.83	109.37	124.46
7	C	1203	HEM	C3B-CAB-CBB	-8.99	110.67	124.46
7	C	1202	HEM	C3B-CAB-CBB	-8.95	110.73	124.46
7	C	1204	HEM	C3B-CAB-CBB	-8.33	111.68	124.46

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
9	M	401	BPB	C13
8	M	1303	BCB	C13
8	L	304	BCB	C8

There are no torsion outliers.

There are no ring outliers.

17 monomers are involved in 113 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	C	1201	HEM	2	0
7	C	1202	HEM	1	0
7	C	1203	HEM	1	0
7	C	1204	HEM	3	0
13	H	703	LDA	4	0
8	L	302	BCB	15	0
8	L	304	BCB	8	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	L	402	BPB	12	0
11	L	502	UQ2	4	0
13	L	705	LDA	2	0
8	M	1301	BCB	12	0
8	M	1303	BCB	19	0
9	M	401	BPB	22	0
12	M	600	NS5	5	0
13	M	701	LDA	3	0
13	M	702	LDA	7	0
13	M	706	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, 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	C	332/336 (98%)	-0.35	4 (1%) 81 78	23, 45, 67, 78	15 (4%)
2	L	273/273 (100%)	-0.52	0 100 100	25, 43, 65, 76	6 (2%)
3	M	323/323 (100%)	-0.54	1 (0%) 94 93	25, 42, 68, 77	8 (2%)
4	H	251/258 (97%)	-0.18	5 (1%) 68 64	30, 56, 84, 93	19 (7%)
All	All	1179/1190 (99%)	-0.40	10 (0%) 87 86	23, 46, 71, 93	48 (4%)

The worst 5 of 10 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	H	8	GLN	5.3
4	H	46	PRO	3.4
4	H	258	LEU	2.9
4	H	93	THR	2.7
1	C	162	HIS	2.5

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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, 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	LLDF	B-factors(Å ²)	Q<0.9
4	FME	H	1	10/11	0.96	0.32	-	53,55,63,65	0

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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, 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	LLDF	B-factors(\AA^2)	Q<0.9
13	LDA	M	704	16/16	0.82	0.43	7.80	79,82,83,84	3
13	LDA	M	702	16/16	0.49	0.49	6.86	73,87,100,101	0
11	UQ2	L	502	23/23	0.93	0.28	6.69	48,49,50,50	23
13	LDA	M	706	16/16	0.73	0.39	5.15	99,105,114,114	0
12	NS5	M	600	40/40	0.90	0.23	3.85	39,58,67,68	11
13	LDA	L	705	16/16	0.74	0.36	3.19	71,78,94,95	3
13	LDA	H	703	16/16	0.90	0.31	2.81	74,79,85,85	0
10	MQ7	M	501	48/48	0.93	0.22	2.74	36,45,67,70	0
9	BPB	M	401	65/65	0.94	0.21	2.71	37,45,94,95	0
9	BPB	L	402	65/65	0.96	0.20	2.00	31,41,45,47	0
13	LDA	M	701	16/16	0.95	0.20	0.83	48,50,52,52	0
8	BCB	L	304	66/66	0.97	0.18	0.52	23,29,54,57	0
8	BCB	M	1301	66/66	0.97	0.18	0.35	33,43,59,63	15
7	HEM	C	1203	43/43	0.99	0.16	0.34	26,33,41,44	0
8	BCB	M	1303	66/66	0.97	0.16	0.07	17,25,46,48	0
8	BCB	L	302	66/66	0.98	0.16	0.05	10,20,42,43	0
7	HEM	C	1201	43/43	0.98	0.14	-0.45	47,51,61,65	0
7	HEM	C	1202	43/43	0.99	0.14	-0.48	39,43,55,58	0
7	HEM	C	1204	43/43	0.99	0.13	-0.55	32,36,48,53	0
6	SO4	M	804	5/5	0.96	0.12	-0.83	81,82,82,83	0
6	SO4	H	801	5/5	0.97	0.10	-2.22	73,76,76,77	0
5	FE2	M	500	1/1	1.00	0.06	-3.07	33,33,33,33	0
6	SO4	M	803	5/5	0.95	0.13	-	102,102,103,103	0
6	SO4	M	802	5/5	0.98	0.11	-	57,59,61,61	0

6.5 Other polymers ⓘ

There are no such residues in this entry.