



wwPDB X-ray Structure Validation Summary Report ⓘ

May 28, 2020 – 09:50 pm BST

PDB ID : 2G18
Title : Crystal Structure of Nostoc sp. 7120 phycocyanobilin:ferredoxin oxidoreduc-
tase (PcyA) Apoprotein
Authors : Tu, S.-L.; Lagarias, J.C.; Fisher, A.J.
Deposited on : 2006-02-13
Resolution : 2.50 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

<https://www.wwpdb.org/validation/2017/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
Xtriage (Phenix) : 1.13
EDS : 2.11
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

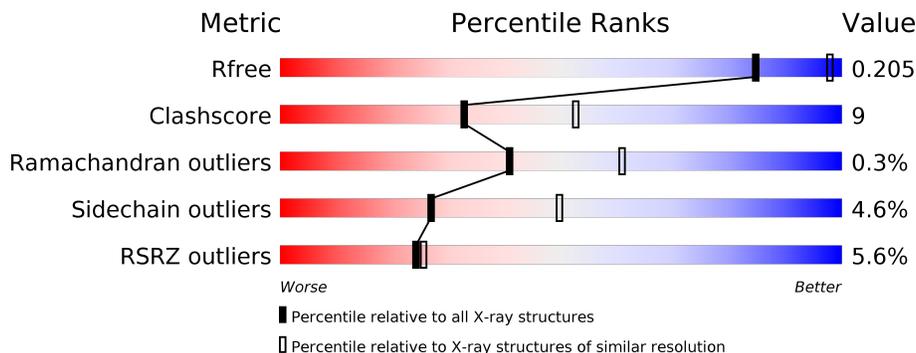
1 Overall quality at a glance

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}	130704	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	253	
1	B	253	
1	C	253	
1	D	253	
1	E	253	
1	F	253	

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Mol	Chain	Length	Quality of chain
1	G	253	<p>4% 79% 18% ••</p>
1	H	253	<p>4% 80% 16% ••</p>
1	I	253	<p>6% 72% 21% • 6%</p>
1	J	253	<p>5% 74% 19% • 6%</p>
1	K	253	<p>8% 72% 21% • 6%</p>
1	L	253	<p>13% 80% 12% •• 6%</p>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 23613 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 Phycocyanobilin:ferredoxin oxidoreductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	245	1964	1242	341	368	13	0	0	0
1	B	245	1964	1242	341	368	13	0	0	0
1	C	245	1964	1242	341	368	13	0	0	0
1	D	248	1991	1261	344	373	13	0	0	0
1	E	245	1964	1242	341	368	13	0	0	0
1	F	245	1964	1242	341	368	13	0	0	0
1	G	248	1991	1261	344	373	13	0	0	0
1	H	245	1964	1242	341	368	13	0	0	0
1	I	238	1914	1209	334	358	13	0	0	0
1	J	238	1914	1209	334	358	13	0	0	0
1	K	238	1914	1209	334	358	13	0	0	0
1	L	238	1914	1209	334	358	13	0	0	0

There are 108 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	GLY	-	cloning artifact	UNP Q93TN0
A	-6	PRO	-	cloning artifact	UNP Q93TN0
A	-5	LEU	-	cloning artifact	UNP Q93TN0
A	-4	GLY	-	cloning artifact	UNP Q93TN0
A	-3	SER	-	cloning artifact	UNP Q93TN0

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	PRO	-	cloning artifact	UNP Q93TN0
A	-1	GLU	-	cloning artifact	UNP Q93TN0
A	0	PHE	-	cloning artifact	UNP Q93TN0
A	1	ILE	-	cloning artifact	UNP Q93TN0
B	-7	GLY	-	cloning artifact	UNP Q93TN0
B	-6	PRO	-	cloning artifact	UNP Q93TN0
B	-5	LEU	-	cloning artifact	UNP Q93TN0
B	-4	GLY	-	cloning artifact	UNP Q93TN0
B	-3	SER	-	cloning artifact	UNP Q93TN0
B	-2	PRO	-	cloning artifact	UNP Q93TN0
B	-1	GLU	-	cloning artifact	UNP Q93TN0
B	0	PHE	-	cloning artifact	UNP Q93TN0
B	1	ILE	-	cloning artifact	UNP Q93TN0
C	-7	GLY	-	cloning artifact	UNP Q93TN0
C	-6	PRO	-	cloning artifact	UNP Q93TN0
C	-5	LEU	-	cloning artifact	UNP Q93TN0
C	-4	GLY	-	cloning artifact	UNP Q93TN0
C	-3	SER	-	cloning artifact	UNP Q93TN0
C	-2	PRO	-	cloning artifact	UNP Q93TN0
C	-1	GLU	-	cloning artifact	UNP Q93TN0
C	0	PHE	-	cloning artifact	UNP Q93TN0
C	1	ILE	-	cloning artifact	UNP Q93TN0
D	-7	GLY	-	cloning artifact	UNP Q93TN0
D	-6	PRO	-	cloning artifact	UNP Q93TN0
D	-5	LEU	-	cloning artifact	UNP Q93TN0
D	-4	GLY	-	cloning artifact	UNP Q93TN0
D	-3	SER	-	cloning artifact	UNP Q93TN0
D	-2	PRO	-	cloning artifact	UNP Q93TN0
D	-1	GLU	-	cloning artifact	UNP Q93TN0
D	0	PHE	-	cloning artifact	UNP Q93TN0
D	1	ILE	-	cloning artifact	UNP Q93TN0
E	-7	GLY	-	cloning artifact	UNP Q93TN0
E	-6	PRO	-	cloning artifact	UNP Q93TN0
E	-5	LEU	-	cloning artifact	UNP Q93TN0
E	-4	GLY	-	cloning artifact	UNP Q93TN0
E	-3	SER	-	cloning artifact	UNP Q93TN0
E	-2	PRO	-	cloning artifact	UNP Q93TN0
E	-1	GLU	-	cloning artifact	UNP Q93TN0
E	0	PHE	-	cloning artifact	UNP Q93TN0
E	1	ILE	-	cloning artifact	UNP Q93TN0
F	-7	GLY	-	cloning artifact	UNP Q93TN0
F	-6	PRO	-	cloning artifact	UNP Q93TN0

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Chain	Residue	Modelled	Actual	Comment	Reference
F	-5	LEU	-	cloning artifact	UNP Q93TN0
F	-4	GLY	-	cloning artifact	UNP Q93TN0
F	-3	SER	-	cloning artifact	UNP Q93TN0
F	-2	PRO	-	cloning artifact	UNP Q93TN0
F	-1	GLU	-	cloning artifact	UNP Q93TN0
F	0	PHE	-	cloning artifact	UNP Q93TN0
F	1	ILE	-	cloning artifact	UNP Q93TN0
G	-7	GLY	-	cloning artifact	UNP Q93TN0
G	-6	PRO	-	cloning artifact	UNP Q93TN0
G	-5	LEU	-	cloning artifact	UNP Q93TN0
G	-4	GLY	-	cloning artifact	UNP Q93TN0
G	-3	SER	-	cloning artifact	UNP Q93TN0
G	-2	PRO	-	cloning artifact	UNP Q93TN0
G	-1	GLU	-	cloning artifact	UNP Q93TN0
G	0	PHE	-	cloning artifact	UNP Q93TN0
G	1	ILE	-	cloning artifact	UNP Q93TN0
H	-7	GLY	-	cloning artifact	UNP Q93TN0
H	-6	PRO	-	cloning artifact	UNP Q93TN0
H	-5	LEU	-	cloning artifact	UNP Q93TN0
H	-4	GLY	-	cloning artifact	UNP Q93TN0
H	-3	SER	-	cloning artifact	UNP Q93TN0
H	-2	PRO	-	cloning artifact	UNP Q93TN0
H	-1	GLU	-	cloning artifact	UNP Q93TN0
H	0	PHE	-	cloning artifact	UNP Q93TN0
H	1	ILE	-	cloning artifact	UNP Q93TN0
I	-7	GLY	-	cloning artifact	UNP Q93TN0
I	-6	PRO	-	cloning artifact	UNP Q93TN0
I	-5	LEU	-	cloning artifact	UNP Q93TN0
I	-4	GLY	-	cloning artifact	UNP Q93TN0
I	-3	SER	-	cloning artifact	UNP Q93TN0
I	-2	PRO	-	cloning artifact	UNP Q93TN0
I	-1	GLU	-	cloning artifact	UNP Q93TN0
I	0	PHE	-	cloning artifact	UNP Q93TN0
I	1	ILE	-	cloning artifact	UNP Q93TN0
J	-7	GLY	-	cloning artifact	UNP Q93TN0
J	-6	PRO	-	cloning artifact	UNP Q93TN0
J	-5	LEU	-	cloning artifact	UNP Q93TN0
J	-4	GLY	-	cloning artifact	UNP Q93TN0
J	-3	SER	-	cloning artifact	UNP Q93TN0
J	-2	PRO	-	cloning artifact	UNP Q93TN0
J	-1	GLU	-	cloning artifact	UNP Q93TN0
J	0	PHE	-	cloning artifact	UNP Q93TN0

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Chain	Residue	Modelled	Actual	Comment	Reference
J	1	ILE	-	cloning artifact	UNP Q93TN0
K	-7	GLY	-	cloning artifact	UNP Q93TN0
K	-6	PRO	-	cloning artifact	UNP Q93TN0
K	-5	LEU	-	cloning artifact	UNP Q93TN0
K	-4	GLY	-	cloning artifact	UNP Q93TN0
K	-3	SER	-	cloning artifact	UNP Q93TN0
K	-2	PRO	-	cloning artifact	UNP Q93TN0
K	-1	GLU	-	cloning artifact	UNP Q93TN0
K	0	PHE	-	cloning artifact	UNP Q93TN0
K	1	ILE	-	cloning artifact	UNP Q93TN0
L	-7	GLY	-	cloning artifact	UNP Q93TN0
L	-6	PRO	-	cloning artifact	UNP Q93TN0
L	-5	LEU	-	cloning artifact	UNP Q93TN0
L	-4	GLY	-	cloning artifact	UNP Q93TN0
L	-3	SER	-	cloning artifact	UNP Q93TN0
L	-2	PRO	-	cloning artifact	UNP Q93TN0
L	-1	GLU	-	cloning artifact	UNP Q93TN0
L	0	PHE	-	cloning artifact	UNP Q93TN0
L	1	ILE	-	cloning artifact	UNP Q93TN0

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	E	1	Total Ca 1 1	0	0

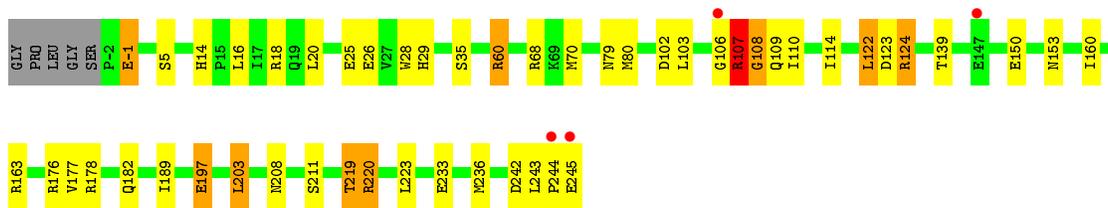
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	19	Total O 19 19	0	0
3	B	7	Total O 7 7	0	0
3	C	16	Total O 16 16	0	0
3	D	31	Total O 31 31	0	0
3	E	18	Total O 18 18	0	0
3	F	17	Total O 17 17	0	0
3	G	18	Total O 18 18	0	0

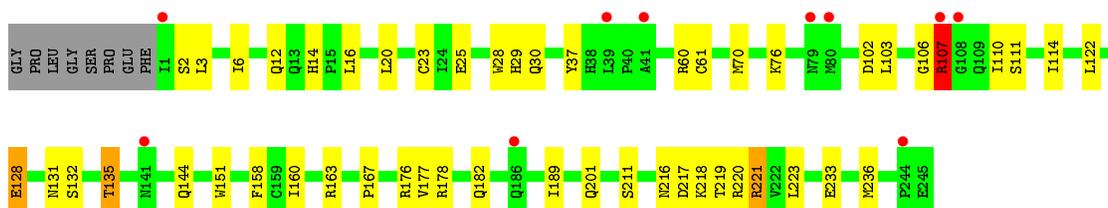
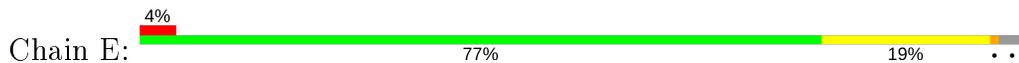
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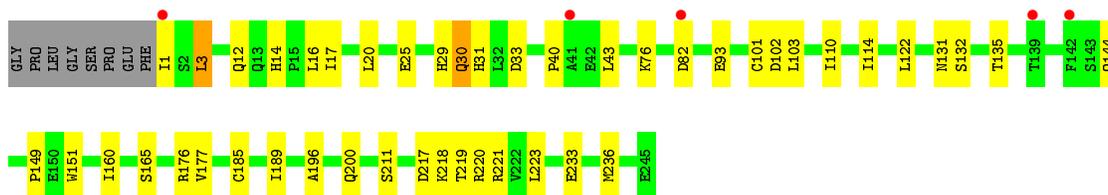
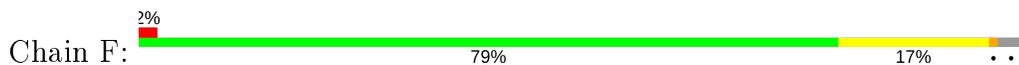
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	H	25	Total 25	O 25	0	0
3	I	14	Total 14	O 14	0	0
3	J	11	Total 11	O 11	0	0
3	K	13	Total 13	O 13	0	0
3	L	1	Total 1	O 1	0	0



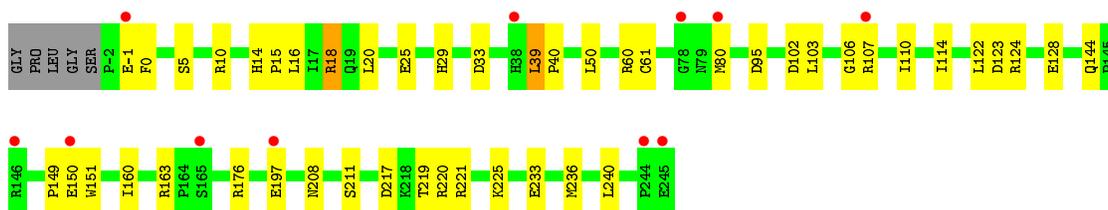
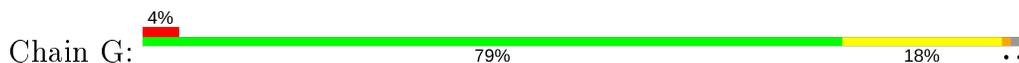
• Molecule 1: Phycocyanobilin:ferredoxin oxidoreductase



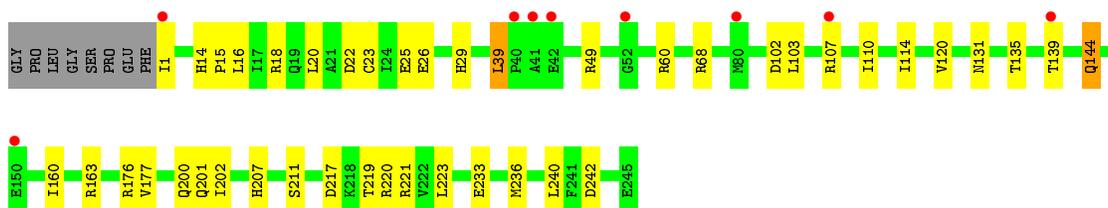
• Molecule 1: Phycocyanobilin:ferredoxin oxidoreductase



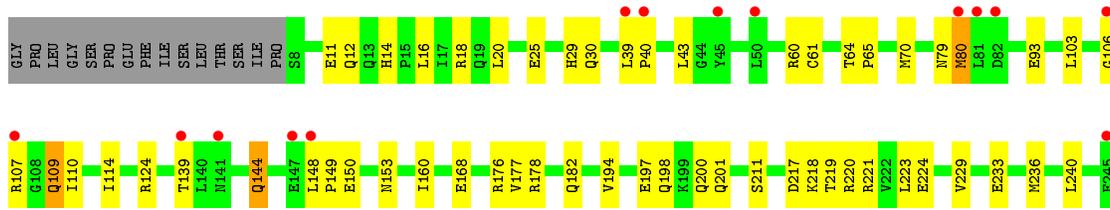
• Molecule 1: Phycocyanobilin:ferredoxin oxidoreductase



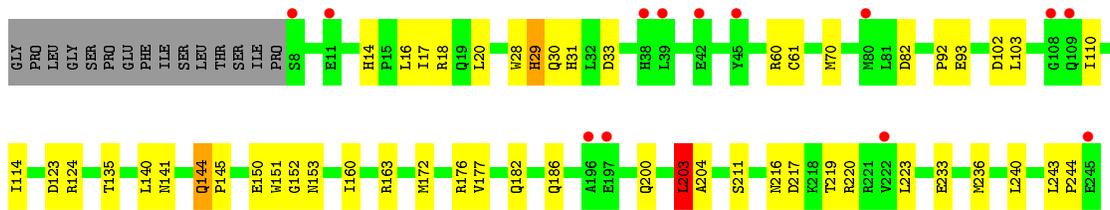
• Molecule 1: Phycocyanobilin:ferredoxin oxidoreductase



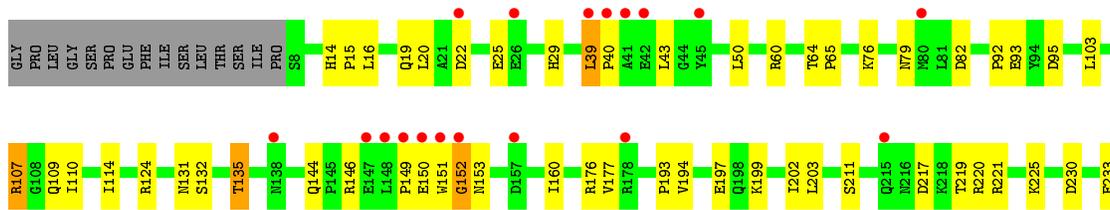
• Molecule 1: Phycocyanobilin:ferredoxin oxidoreductase



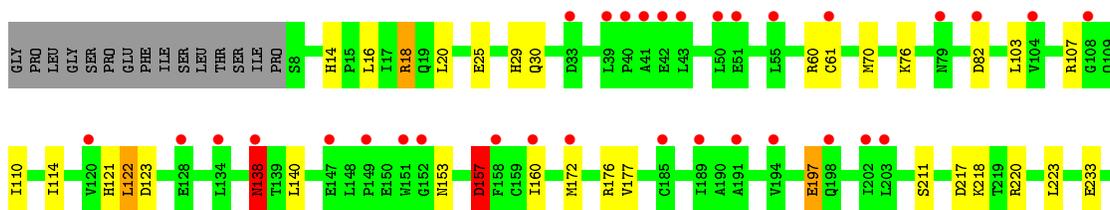
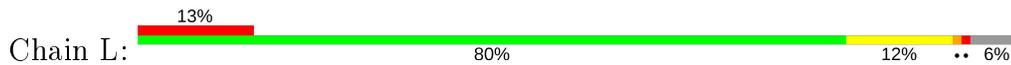
• Molecule 1: Phycocyanobilin:ferredoxin oxidoreductase



• Molecule 1: Phycocyanobilin:ferredoxin oxidoreductase



• Molecule 1: Phycocyanobilin:ferredoxin oxidoreductase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	78.50Å 141.33Å 121.84Å 90.00° 91.29° 90.00°	Depositor
Resolution (Å)	43.94 – 2.50 43.94 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.7 (43.94-2.50) 99.6 (43.94-2.50)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.27 (at 2.51Å)	Xtrriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.211 , 0.257 0.206 , 0.205	Depositor DCC
R_{free} test set	4582 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	44.4	Xtrriage
Anisotropy	0.050	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 55.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	0.018 for h,-k,-l	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	23613	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 27.33 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.2431e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.92	6/2012 (0.3%)	1.00	3/2729 (0.1%)
1	B	0.68	1/2012 (0.0%)	0.71	0/2729
1	C	0.77	1/2012 (0.0%)	0.94	5/2729 (0.2%)
1	D	0.84	1/2041 (0.0%)	0.84	4/2768 (0.1%)
1	E	0.82	0/2012	0.76	1/2729 (0.0%)
1	F	0.79	1/2012 (0.0%)	0.75	1/2729 (0.0%)
1	G	0.72	1/2041 (0.0%)	0.72	0/2768
1	H	0.81	0/2012	0.78	0/2729
1	I	0.73	0/1961	0.74	0/2658
1	J	0.80	2/1961 (0.1%)	0.79	3/2658 (0.1%)
1	K	0.74	0/1961	0.75	0/2658
1	L	1.24	4/1961 (0.2%)	0.70	3/2658 (0.1%)
All	All	0.83	17/23998 (0.1%)	0.79	20/32542 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	3

The worst 5 of 17 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	L	197	GLU	CD-OE1	36.97	1.66	1.25
1	L	197	GLU	CD-OE2	23.40	1.51	1.25
1	L	157	ASP	CB-CG	17.69	1.89	1.51
1	A	178	ARG	NE-CZ	15.40	1.53	1.33
1	A	178	ARG	CZ-NH1	-13.96	1.15	1.33

The worst 5 of 20 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	178	ARG	NE-CZ-NH2	31.77	136.18	120.30
1	C	178	ARG	NE-CZ-NH2	-26.81	106.89	120.30
1	A	178	ARG	NE-CZ-NH1	-13.51	113.54	120.30
1	J	93	GLU	OE1-CD-OE2	-10.95	110.16	123.30
1	L	197	GLU	OE1-CD-OE2	10.62	136.04	123.30

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	39	LEU	Peptide
1	B	40	PRO	Peptide
1	B	41	ALA	Peptide

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1964	0	1916	31	0
1	B	1964	0	1916	38	1
1	C	1964	0	1916	33	0
1	D	1991	0	1938	50	0
1	E	1964	0	1916	39	0
1	F	1964	0	1916	27	2
1	G	1991	0	1938	29	0
1	H	1964	0	1916	26	0
1	I	1914	0	1856	39	1
1	J	1914	0	1856	54	0
1	K	1914	0	1856	48	0
1	L	1914	0	1856	28	0
2	E	1	0	0	0	0
3	A	19	0	0	2	0
3	B	7	0	0	1	0
3	C	16	0	0	4	0
3	D	31	0	0	6	0
3	E	18	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	F	17	0	0	0	0
3	G	18	0	0	1	0
3	H	25	0	0	3	0
3	I	14	0	0	2	0
3	J	11	0	0	0	0
3	K	13	0	0	0	0
3	L	1	0	0	0	0
All	All	23613	0	22796	400	2

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 9.

The worst 5 of 400 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:157:ASP:CB	1:L:157:ASP:CG	1.88	1.39
1:L:197:GLU:CD	1:L:197:GLU:OE1	1.66	1.31
1:K:124:ARG:NH2	1:K:153:ASN:HB2	1.66	1.08
1:D:35:SER:CB	1:J:186:GLN:NE2	2.21	1.04
1:D:35:SER:CB	1:J:186:GLN:HE22	1.69	1.04

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:60:ARG:NH2	1:F:3:LEU:O[2_456]	2.16	0.04
1:F:218:LYS:NZ	1:I:197:GLU:OE1[2_446]	2.19	0.01

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	243/253 (96%)	235 (97%)	8 (3%)	0	100	100
1	B	243/253 (96%)	236 (97%)	4 (2%)	3 (1%)	13	24
1	C	243/253 (96%)	238 (98%)	4 (2%)	1 (0%)	34	54
1	D	246/253 (97%)	238 (97%)	6 (2%)	2 (1%)	19	35
1	E	243/253 (96%)	237 (98%)	5 (2%)	1 (0%)	34	54
1	F	243/253 (96%)	238 (98%)	5 (2%)	0	100	100
1	G	246/253 (97%)	240 (98%)	6 (2%)	0	100	100
1	H	243/253 (96%)	236 (97%)	7 (3%)	0	100	100
1	I	236/253 (93%)	231 (98%)	5 (2%)	0	100	100
1	J	236/253 (93%)	231 (98%)	5 (2%)	0	100	100
1	K	236/253 (93%)	228 (97%)	7 (3%)	1 (0%)	34	54
1	L	236/253 (93%)	228 (97%)	7 (3%)	1 (0%)	34	54
All	All	2894/3036 (95%)	2816 (97%)	69 (2%)	9 (0%)	41	61

5 of 9 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	40	PRO
1	D	108	GLY
1	D	107	ARG
1	K	152	GLY
1	B	41	ALA

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	217/223 (97%)	205 (94%)	12 (6%)	21	41
1	B	217/223 (97%)	206 (95%)	11 (5%)	24	45
1	C	217/223 (97%)	209 (96%)	8 (4%)	34	60
1	D	220/223 (99%)	206 (94%)	14 (6%)	17	33

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	E	217/223 (97%)	206 (95%)	11 (5%)	24	45
1	F	217/223 (97%)	211 (97%)	6 (3%)	43	70
1	G	220/223 (99%)	209 (95%)	11 (5%)	24	46
1	H	217/223 (97%)	207 (95%)	10 (5%)	27	50
1	I	210/223 (94%)	199 (95%)	11 (5%)	23	44
1	J	210/223 (94%)	205 (98%)	5 (2%)	49	74
1	K	210/223 (94%)	198 (94%)	12 (6%)	20	39
1	L	210/223 (94%)	201 (96%)	9 (4%)	29	53
All	All	2582/2676 (96%)	2462 (95%)	120 (5%)	27	50

5 of 120 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	211	SER
1	G	122	LEU
1	K	221	ARG
1	E	221	ARG
1	F	211	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 46 such sidechains are listed below:

Mol	Chain	Res	Type
1	F	31	HIS
1	G	186	GLN
1	K	31	HIS
1	F	144	GLN
1	G	29	HIS

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	A	245/253 (96%)	0.57	15 (6%) 21 22	39, 51, 68, 83	0
1	B	245/253 (96%)	0.56	17 (6%) 16 17	40, 52, 73, 78	0
1	C	245/253 (96%)	0.41	12 (4%) 29 31	40, 52, 69, 79	0
1	D	248/253 (98%)	0.28	4 (1%) 72 74	40, 51, 68, 76	0
1	E	245/253 (96%)	0.36	10 (4%) 37 40	38, 52, 69, 82	0
1	F	245/253 (96%)	0.41	5 (2%) 65 68	40, 52, 66, 73	0
1	G	248/253 (98%)	0.45	11 (4%) 34 37	38, 52, 70, 77	0
1	H	245/253 (96%)	0.42	9 (3%) 41 45	39, 53, 66, 76	0
1	I	238/253 (94%)	0.49	14 (5%) 22 23	39, 52, 68, 77	0
1	J	238/253 (94%)	0.49	13 (5%) 25 26	40, 51, 66, 84	0
1	K	238/253 (94%)	0.65	20 (8%) 11 11	39, 52, 72, 90	0
1	L	238/253 (94%)	1.04	34 (14%) 2 2	39, 53, 67, 74	0
All	All	2918/3036 (96%)	0.51	164 (5%) 24 25	38, 52, 69, 90	0

The worst 5 of 164 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	L	51	GLU	7.5
1	K	151	TRP	6.5
1	F	1	ILE	6.3
1	B	144	GLN	5.5
1	L	42	GLU	5.2

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. 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
2	CA	E	246	1/1	0.69	0.25	23,23,23,23	1

6.5 Other polymers [i](#)

There are no such residues in this entry.