wwPDB X-ray Structure Validation Summary Report (i)

May 22, 2020 – 09:20 am BST

PDB ID	:	6SCF
Title	:	A viral anti-CRISPR subverts type III CRISPR immunity by rapid degradation of cyclic oligoadenylate
Authors	:	McMahon, S.A.; Athukoralage, J.S.; Graham, S.; White, M.F.; Gloster, T.M.
Deposited on	:	2019-07-24
Resolution	:	1.55 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

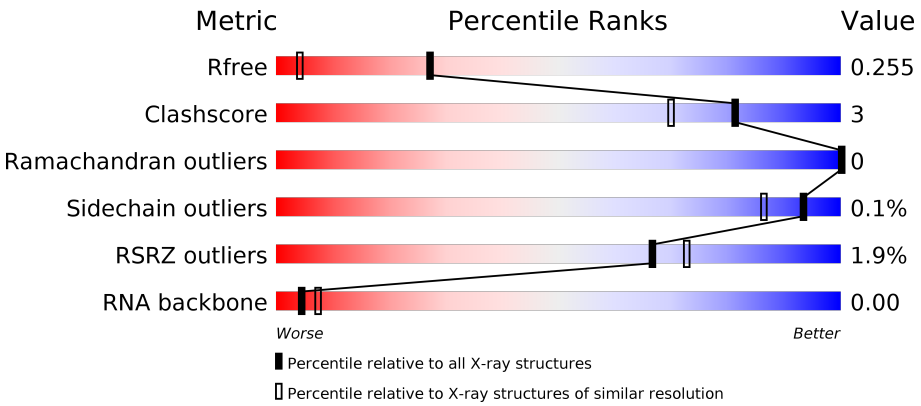
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 1.55 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	1483 (1.56-1.56)
Clashscore	141614	1529 (1.56-1.56)
Ramachandran outliers	138981	1498 (1.56-1.56)
Sidechain outliers	138945	1495 (1.56-1.56)
RSRZ outliers	127900	1465 (1.56-1.56)
RNA backbone	3102	1015 (2.36-0.86)

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 <=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	138	75% 7% 18%
1	B	138	% 76% 6% 18%
1	C	138	3% 78% • 19%
1	D	138	% 80% • 19%

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Mol	Chain	Length	Quality of chain
1	E	138	3% 76% 19%
1	F	138	74% 9% 17%
1	G	138	2% 76% 5% 19%
1	H	138	3% 75% 21%
2	I	4	50% 25% 25%
2	K	4	25% 50% 25%
2	L	4	50% 25% 25%
2	M	4	50% 50%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 8312 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 Uncharacterized protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	113	916	587	146	180	3	0	0	0
1	B	113	950	610	151	186	3	0	4	0
1	C	112	928	598	146	181	3	0	3	0
1	D	112	920	599	145	173	3	0	4	0
1	E	112	902	581	142	176	3	0	1	0
1	F	114	950	609	150	187	4	0	4	0
1	G	112	917	591	147	176	3	0	2	0
1	H	109	882	570	141	168	3	0	0	0

There are 200 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-23	MET	-	initiating methionine	UNP Q8QL27
A	-22	SER	-	expression tag	UNP Q8QL27
A	-21	TYR	-	expression tag	UNP Q8QL27
A	-20	TYR	-	expression tag	UNP Q8QL27
A	-19	HIS	-	expression tag	UNP Q8QL27
A	-18	HIS	-	expression tag	UNP Q8QL27
A	-17	HIS	-	expression tag	UNP Q8QL27
A	-16	HIS	-	expression tag	UNP Q8QL27
A	-15	HIS	-	expression tag	UNP Q8QL27
A	-14	HIS	-	expression tag	UNP Q8QL27
A	-13	ASP	-	expression tag	UNP Q8QL27
A	-12	TYR	-	expression tag	UNP Q8QL27
A	-11	ASP	-	expression tag	UNP Q8QL27

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-10	ILE	-	expression tag	UNP Q8QL27
A	-9	PRO	-	expression tag	UNP Q8QL27
A	-8	THR	-	expression tag	UNP Q8QL27
A	-7	THR	-	expression tag	UNP Q8QL27
A	-6	GLU	-	expression tag	UNP Q8QL27
A	-5	ASN	-	expression tag	UNP Q8QL27
A	-4	LEU	-	expression tag	UNP Q8QL27
A	-3	TYR	-	expression tag	UNP Q8QL27
A	-2	PHE	-	expression tag	UNP Q8QL27
A	-1	GLN	-	expression tag	UNP Q8QL27
A	0	GLY	-	expression tag	UNP Q8QL27
A	47	ALA	HIS	conflict	UNP Q8QL27
B	-23	MET	-	initiating methionine	UNP Q8QL27
B	-22	SER	-	expression tag	UNP Q8QL27
B	-21	TYR	-	expression tag	UNP Q8QL27
B	-20	TYR	-	expression tag	UNP Q8QL27
B	-19	HIS	-	expression tag	UNP Q8QL27
B	-18	HIS	-	expression tag	UNP Q8QL27
B	-17	HIS	-	expression tag	UNP Q8QL27
B	-16	HIS	-	expression tag	UNP Q8QL27
B	-15	HIS	-	expression tag	UNP Q8QL27
B	-14	HIS	-	expression tag	UNP Q8QL27
B	-13	ASP	-	expression tag	UNP Q8QL27
B	-12	TYR	-	expression tag	UNP Q8QL27
B	-11	ASP	-	expression tag	UNP Q8QL27
B	-10	ILE	-	expression tag	UNP Q8QL27
B	-9	PRO	-	expression tag	UNP Q8QL27
B	-8	THR	-	expression tag	UNP Q8QL27
B	-7	THR	-	expression tag	UNP Q8QL27
B	-6	GLU	-	expression tag	UNP Q8QL27
B	-5	ASN	-	expression tag	UNP Q8QL27
B	-4	LEU	-	expression tag	UNP Q8QL27
B	-3	TYR	-	expression tag	UNP Q8QL27
B	-2	PHE	-	expression tag	UNP Q8QL27
B	-1	GLN	-	expression tag	UNP Q8QL27
B	0	GLY	-	expression tag	UNP Q8QL27
B	47	ALA	HIS	conflict	UNP Q8QL27
C	-23	MET	-	initiating methionine	UNP Q8QL27
C	-22	SER	-	expression tag	UNP Q8QL27
C	-21	TYR	-	expression tag	UNP Q8QL27
C	-20	TYR	-	expression tag	UNP Q8QL27
C	-19	HIS	-	expression tag	UNP Q8QL27

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-18	HIS	-	expression tag	UNP Q8QL27
C	-17	HIS	-	expression tag	UNP Q8QL27
C	-16	HIS	-	expression tag	UNP Q8QL27
C	-15	HIS	-	expression tag	UNP Q8QL27
C	-14	HIS	-	expression tag	UNP Q8QL27
C	-13	ASP	-	expression tag	UNP Q8QL27
C	-12	TYR	-	expression tag	UNP Q8QL27
C	-11	ASP	-	expression tag	UNP Q8QL27
C	-10	ILE	-	expression tag	UNP Q8QL27
C	-9	PRO	-	expression tag	UNP Q8QL27
C	-8	THR	-	expression tag	UNP Q8QL27
C	-7	THR	-	expression tag	UNP Q8QL27
C	-6	GLU	-	expression tag	UNP Q8QL27
C	-5	ASN	-	expression tag	UNP Q8QL27
C	-4	LEU	-	expression tag	UNP Q8QL27
C	-3	TYR	-	expression tag	UNP Q8QL27
C	-2	PHE	-	expression tag	UNP Q8QL27
C	-1	GLN	-	expression tag	UNP Q8QL27
C	0	GLY	-	expression tag	UNP Q8QL27
C	47	ALA	HIS	conflict	UNP Q8QL27
D	-23	MET	-	initiating methionine	UNP Q8QL27
D	-22	SER	-	expression tag	UNP Q8QL27
D	-21	TYR	-	expression tag	UNP Q8QL27
D	-20	TYR	-	expression tag	UNP Q8QL27
D	-19	HIS	-	expression tag	UNP Q8QL27
D	-18	HIS	-	expression tag	UNP Q8QL27
D	-17	HIS	-	expression tag	UNP Q8QL27
D	-16	HIS	-	expression tag	UNP Q8QL27
D	-15	HIS	-	expression tag	UNP Q8QL27
D	-14	HIS	-	expression tag	UNP Q8QL27
D	-13	ASP	-	expression tag	UNP Q8QL27
D	-12	TYR	-	expression tag	UNP Q8QL27
D	-11	ASP	-	expression tag	UNP Q8QL27
D	-10	ILE	-	expression tag	UNP Q8QL27
D	-9	PRO	-	expression tag	UNP Q8QL27
D	-8	THR	-	expression tag	UNP Q8QL27
D	-7	THR	-	expression tag	UNP Q8QL27
D	-6	GLU	-	expression tag	UNP Q8QL27
D	-5	ASN	-	expression tag	UNP Q8QL27
D	-4	LEU	-	expression tag	UNP Q8QL27
D	-3	TYR	-	expression tag	UNP Q8QL27
D	-2	PHE	-	expression tag	UNP Q8QL27

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-1	GLN	-	expression tag	UNP Q8QL27
D	0	GLY	-	expression tag	UNP Q8QL27
D	47	ALA	HIS	conflict	UNP Q8QL27
E	-23	MET	-	initiating methionine	UNP Q8QL27
E	-22	SER	-	expression tag	UNP Q8QL27
E	-21	TYR	-	expression tag	UNP Q8QL27
E	-20	TYR	-	expression tag	UNP Q8QL27
E	-19	HIS	-	expression tag	UNP Q8QL27
E	-18	HIS	-	expression tag	UNP Q8QL27
E	-17	HIS	-	expression tag	UNP Q8QL27
E	-16	HIS	-	expression tag	UNP Q8QL27
E	-15	HIS	-	expression tag	UNP Q8QL27
E	-14	HIS	-	expression tag	UNP Q8QL27
E	-13	ASP	-	expression tag	UNP Q8QL27
E	-12	TYR	-	expression tag	UNP Q8QL27
E	-11	ASP	-	expression tag	UNP Q8QL27
E	-10	ILE	-	expression tag	UNP Q8QL27
E	-9	PRO	-	expression tag	UNP Q8QL27
E	-8	THR	-	expression tag	UNP Q8QL27
E	-7	THR	-	expression tag	UNP Q8QL27
E	-6	GLU	-	expression tag	UNP Q8QL27
E	-5	ASN	-	expression tag	UNP Q8QL27
E	-4	LEU	-	expression tag	UNP Q8QL27
E	-3	TYR	-	expression tag	UNP Q8QL27
E	-2	PHE	-	expression tag	UNP Q8QL27
E	-1	GLN	-	expression tag	UNP Q8QL27
E	0	GLY	-	expression tag	UNP Q8QL27
E	47	ALA	HIS	conflict	UNP Q8QL27
F	-23	MET	-	initiating methionine	UNP Q8QL27
F	-22	SER	-	expression tag	UNP Q8QL27
F	-21	TYR	-	expression tag	UNP Q8QL27
F	-20	TYR	-	expression tag	UNP Q8QL27
F	-19	HIS	-	expression tag	UNP Q8QL27
F	-18	HIS	-	expression tag	UNP Q8QL27
F	-17	HIS	-	expression tag	UNP Q8QL27
F	-16	HIS	-	expression tag	UNP Q8QL27
F	-15	HIS	-	expression tag	UNP Q8QL27
F	-14	HIS	-	expression tag	UNP Q8QL27
F	-13	ASP	-	expression tag	UNP Q8QL27
F	-12	TYR	-	expression tag	UNP Q8QL27
F	-11	ASP	-	expression tag	UNP Q8QL27
F	-10	ILE	-	expression tag	UNP Q8QL27

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Chain	Residue	Modelled	Actual	Comment	Reference
F	-9	PRO	-	expression tag	UNP Q8QL27
F	-8	THR	-	expression tag	UNP Q8QL27
F	-7	THR	-	expression tag	UNP Q8QL27
F	-6	GLU	-	expression tag	UNP Q8QL27
F	-5	ASN	-	expression tag	UNP Q8QL27
F	-4	LEU	-	expression tag	UNP Q8QL27
F	-3	TYR	-	expression tag	UNP Q8QL27
F	-2	PHE	-	expression tag	UNP Q8QL27
F	-1	GLN	-	expression tag	UNP Q8QL27
F	0	GLY	-	expression tag	UNP Q8QL27
F	47	ALA	HIS	conflict	UNP Q8QL27
G	-23	MET	-	initiating methionine	UNP Q8QL27
G	-22	SER	-	expression tag	UNP Q8QL27
G	-21	TYR	-	expression tag	UNP Q8QL27
G	-20	TYR	-	expression tag	UNP Q8QL27
G	-19	HIS	-	expression tag	UNP Q8QL27
G	-18	HIS	-	expression tag	UNP Q8QL27
G	-17	HIS	-	expression tag	UNP Q8QL27
G	-16	HIS	-	expression tag	UNP Q8QL27
G	-15	HIS	-	expression tag	UNP Q8QL27
G	-14	HIS	-	expression tag	UNP Q8QL27
G	-13	ASP	-	expression tag	UNP Q8QL27
G	-12	TYR	-	expression tag	UNP Q8QL27
G	-11	ASP	-	expression tag	UNP Q8QL27
G	-10	ILE	-	expression tag	UNP Q8QL27
G	-9	PRO	-	expression tag	UNP Q8QL27
G	-8	THR	-	expression tag	UNP Q8QL27
G	-7	THR	-	expression tag	UNP Q8QL27
G	-6	GLU	-	expression tag	UNP Q8QL27
G	-5	ASN	-	expression tag	UNP Q8QL27
G	-4	LEU	-	expression tag	UNP Q8QL27
G	-3	TYR	-	expression tag	UNP Q8QL27
G	-2	PHE	-	expression tag	UNP Q8QL27
G	-1	GLN	-	expression tag	UNP Q8QL27
G	0	GLY	-	expression tag	UNP Q8QL27
G	47	ALA	HIS	conflict	UNP Q8QL27
H	-23	MET	-	initiating methionine	UNP Q8QL27
H	-22	SER	-	expression tag	UNP Q8QL27
H	-21	TYR	-	expression tag	UNP Q8QL27
H	-20	TYR	-	expression tag	UNP Q8QL27
H	-19	HIS	-	expression tag	UNP Q8QL27
H	-18	HIS	-	expression tag	UNP Q8QL27

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Chain	Residue	Modelled	Actual	Comment	Reference
H	-17	HIS	-	expression tag	UNP Q8QL27
H	-16	HIS	-	expression tag	UNP Q8QL27
H	-15	HIS	-	expression tag	UNP Q8QL27
H	-14	HIS	-	expression tag	UNP Q8QL27
H	-13	ASP	-	expression tag	UNP Q8QL27
H	-12	TYR	-	expression tag	UNP Q8QL27
H	-11	ASP	-	expression tag	UNP Q8QL27
H	-10	ILE	-	expression tag	UNP Q8QL27
H	-9	PRO	-	expression tag	UNP Q8QL27
H	-8	THR	-	expression tag	UNP Q8QL27
H	-7	THR	-	expression tag	UNP Q8QL27
H	-6	GLU	-	expression tag	UNP Q8QL27
H	-5	ASN	-	expression tag	UNP Q8QL27
H	-4	LEU	-	expression tag	UNP Q8QL27
H	-3	TYR	-	expression tag	UNP Q8QL27
H	-2	PHE	-	expression tag	UNP Q8QL27
H	-1	GLN	-	expression tag	UNP Q8QL27
H	0	GLY	-	expression tag	UNP Q8QL27
H	47	ALA	HIS	conflict	UNP Q8QL27

- Molecule 2 is a RNA chain called cyclic oligoadenylate.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	K	4	88	40	20	24	4	0	0	0
			Total	C	N	O	P			
2	I	4	88	40	20	24	4	0	0	0
			Total	C	N	O	P			
2	L	4	88	40	20	24	4	0	0	0
			Total	C	N	O	P			
2	M	4	88	40	20	24	4	0	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	80	80	80	0	0
			Total	O		
3	B	77	78	78	0	1
			Total	O		
3	C	66	66	66	0	0
			Total	O		
3	D	51	51	51	0	0

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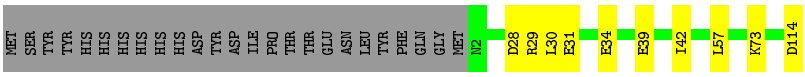
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	E	56	Total O 56 56	0	0
3	F	110	Total O 111 111	0	1
3	G	75	Total O 75 75	0	0
3	H	41	Total O 41 41	0	0
3	K	10	Total O 10 10	0	0
3	I	8	Total O 8 8	0	0
3	L	9	Total O 10 10	0	1
3	M	9	Total O 9 9	0	0

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 the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

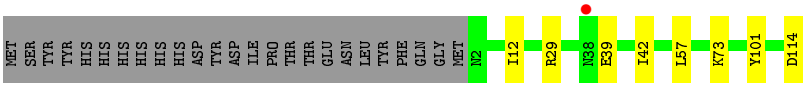
- Molecule 1: Uncharacterized protein

Chain A:



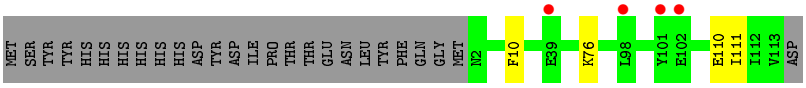
- Molecule 1: Uncharacterized protein

Chain B:



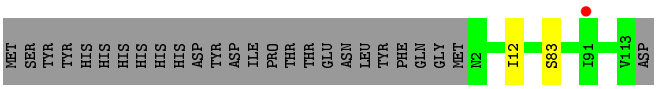
- Molecule 1: Uncharacterized protein

Chain C:



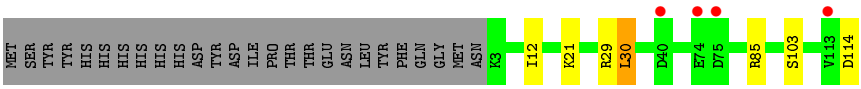
- Molecule 1: Uncharacterized protein

Chain D:



- Molecule 1: Uncharacterized protein

Chain E:



- Molecule 1: Uncharacterized protein

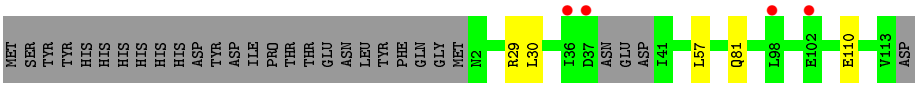
Chain F:

- Molecule 1: Uncharacterized protein

Chain G:

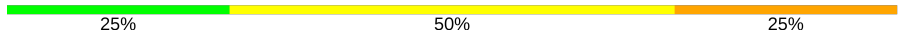
- Molecule 1: Uncharacterized protein

Chain H:



- Molecule 2: cyclic oligoadenylate

Chain K:



- Molecule 2: cyclic oligoadenylate

Chain I:

- Molecule 2: cyclic oligoadenylate

Chain L:

- Molecule 2: cyclic oligoadenylate

Chain M:

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	49.83Å 51.73Å 85.61Å 80.22° 89.68° 83.38°	Depositor
Resolution (Å)	50.64 – 1.55 50.63 – 1.55	Depositor EDS
% Data completeness (in resolution range)	98.6 (50.64-1.55) 98.6 (50.63-1.55)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.11 (at 1.55Å)	Xtriage
Refinement program	REFMAC 5.8.0218	Depositor
R, R_{free}	0.202 , 0.249 0.210 , 0.255	Depositor DCC
R_{free} test set	5883 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å^2)	13.7	Xtriage
Anisotropy	0.250	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 45.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	8312	wwPDB-VP
Average B, all atoms (Å^2)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 82.38 % 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 2.8279e-07. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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)

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	0.65	0/925	0.78	0/1243
1	B	0.65	0/959	0.79	0/1288
1	C	0.67	0/937	0.76	0/1262
1	D	0.60	0/932	0.77	0/1256
1	E	0.69	0/911	0.81	1/1226 (0.1%)
1	F	0.70	0/959	0.84	1/1291 (0.1%)
1	G	0.59	0/926	0.77	0/1246
1	H	0.64	1/890 (0.1%)	0.72	0/1195
2	I	0.82	0/99	1.51	1/152 (0.7%)
2	K	0.72	0/99	1.53	1/152 (0.7%)
2	L	0.89	0/99	1.48	2/152 (1.3%)
2	M	0.95	0/99	1.50	1/152 (0.7%)
All	All	0.66	1/7835 (0.0%)	0.84	7/10615 (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	C	0	1
2	L	0	1
All	All	0	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	110	GLU	CD-OE1	-5.71	1.19	1.25

The worst 5 of 7 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	M	2	A	O4'-C1'-N9	-6.73	102.81	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	K	2	A	O4'-C1'-N9	-6.48	103.02	108.20
2	L	4	A	O4'-C1'-N9	-6.19	103.25	108.20
2	L	1	A	O5'-P-OP2	-6.04	100.26	105.70
1	F	70	LYS	CD-CE-NZ	5.49	124.33	111.70

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	10	PHE	Mainchain
2	L	4	A	Sidechain

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	916	0	942	14	0
1	B	950	0	979	7	0
1	C	928	0	949	2	0
1	D	920	0	956	2	0
1	E	902	0	919	7	0
1	F	950	0	972	12	0
1	G	917	0	947	5	0
1	H	882	0	921	4	0
2	I	88	0	44	0	0
2	K	88	0	44	0	0
2	L	88	0	44	2	0
2	M	88	0	44	2	0
3	A	80	0	0	0	0
3	B	78	0	0	0	0
3	C	66	0	0	0	0
3	D	51	0	0	0	0
3	E	56	0	0	1	0
3	F	111	0	0	4	0
3	G	75	0	0	2	0
3	H	41	0	0	1	0
3	I	8	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	K	10	0	0	0	0
3	L	10	0	0	0	0
3	M	9	0	0	0	0
All	All	8312	0	7761	41	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 3.

The worst 5 of 41 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:A:34:GLU:OE2	1:F:112:ILE:HD13	1.98	0.64
1:B:12:ILE:HD11	1:G:111:ILE:HD11	1.78	0.64
1:H:29:ARG:HE	1:H:57:LEU:HD22	1.68	0.58
1:A:31:GLU:CG	3:F:218:HOH:O	2.53	0.56
1:D:12:ILE:HD11	1:F:111:ILE:HD11	1.87	0.56

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	A	111/138 (80%)	110 (99%)	1 (1%)	0	100	100
1	B	115/138 (83%)	114 (99%)	1 (1%)	0	100	100
1	C	113/138 (82%)	112 (99%)	1 (1%)	0	100	100
1	D	114/138 (83%)	113 (99%)	1 (1%)	0	100	100
1	E	111/138 (80%)	110 (99%)	1 (1%)	0	100	100
1	F	116/138 (84%)	112 (97%)	4 (3%)	0	100	100
1	G	112/138 (81%)	110 (98%)	2 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	H	105/138 (76%)	103 (98%)	2 (2%)	0	100	100
All	All	897/1104 (81%)	884 (99%)	13 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	106/130 (82%)	105 (99%)	1 (1%)	78	61
1	B	110/130 (85%)	110 (100%)	0	100	100
1	C	106/130 (82%)	106 (100%)	0	100	100
1	D	104/130 (80%)	104 (100%)	0	100	100
1	E	102/130 (78%)	102 (100%)	0	100	100
1	F	110/130 (85%)	110 (100%)	0	100	100
1	G	105/130 (81%)	105 (100%)	0	100	100
1	H	102/130 (78%)	102 (100%)	0	100	100
All	All	845/1040 (81%)	844 (100%)	1 (0%)	93	86

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	13	ASN

5.3.3 RNA (i)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	I	3/4 (75%)	2 (66%)	0
2	K	3/4 (75%)	3 (100%)	0
2	L	3/4 (75%)	3 (100%)	0
2	M	3/4 (75%)	2 (66%)	0
All	All	12/16 (75%)	10 (83%)	0

5 of 10 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	K	2	A
2	K	3	A
2	K	4	A
2	I	2	A
2	I	3	A

There are no RNA pucker outliers to report.

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)

There are no ligands in this entry.

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	113/138 (81%)	-0.01	0	100	100	8, 16, 28, 38	0
1	B	113/138 (81%)	0.07	1 (0%)	84	87	10, 16, 29, 34	0
1	C	112/138 (81%)	0.32	4 (3%)	42	50	11, 21, 33, 43	0
1	D	112/138 (81%)	0.33	1 (0%)	84	87	11, 22, 36, 45	0
1	E	112/138 (81%)	0.26	4 (3%)	42	50	9, 19, 45, 56	0
1	F	114/138 (82%)	0.11	0	100	100	10, 16, 30, 44	0
1	G	112/138 (81%)	0.14	3 (2%)	54	62	10, 19, 32, 39	0
1	H	109/138 (78%)	0.28	4 (3%)	41	48	12, 24, 37, 45	0
2	I	4/4 (100%)	-0.52	0	100	100	11, 12, 13, 13	0
2	K	4/4 (100%)	-0.43	0	100	100	13, 13, 15, 16	0
2	L	4/4 (100%)	-0.40	0	100	100	11, 14, 14, 16	0
2	M	4/4 (100%)	-0.51	0	100	100	12, 12, 13, 15	0
All	All	913/1120 (81%)	0.18	17 (1%)	66	73	8, 18, 33, 56	0

The worst 5 of 17 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	102	GLU	3.8
1	E	74	GLU	3.8
1	C	101	TYR	3.4
1	G	37	ASP	3.2
1	E	75	ASP	3.2

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)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.