



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:57 PM UTC

PDB ID : 7BFE / pdb_00007bfe
Title : Circular permutant of ribosomal protein S6, P54-55 truncated, L21A mutant.
Authors : Wang, H.; Logan, D.T.; Oliveberg, M.
Deposited on : 2021-01-02
Resolution : 1.95 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4-5-2 with Phenix2.0
Xtrriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

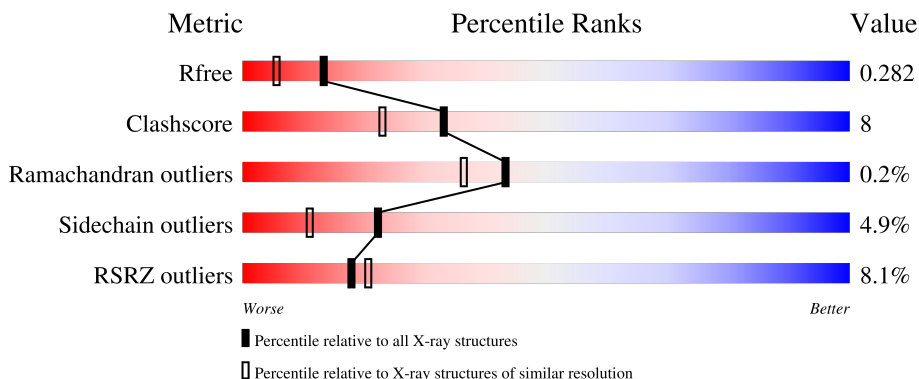
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)
RSRZ outliers	180081	3495 (1.96-1.96)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	86	<div style="display: flex; align-items: center;"> <div style="width: 3%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 80%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 14%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 3%; height: 10px; background-color: grey;"></div> </div> <p style="margin-left: 10px;">3% 80% 14% . .</p>
1	B	86	<div style="display: flex; align-items: center;"> <div style="width: 5%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 67%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 24%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 8%; height: 10px; background-color: orange;"></div> </div> <p style="margin-left: 10px;">5% 67% 24% 8%</p>
1	C	86	<div style="display: flex; align-items: center;"> <div style="width: 2%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 72%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 26%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: grey;"></div> </div> <p style="margin-left: 10px;">2% 72% 26% ..</p>
1	D	86	<div style="display: flex; align-items: center;"> <div style="width: 15%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 66%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 28%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 3%; height: 10px; background-color: orange;"></div> <div style="width: 2%; height: 10px; background-color: grey;"></div> </div> <p style="margin-left: 10px;">15% 66% 28% . . .</p>
1	E	86	<div style="display: flex; align-items: center;"> <div style="width: 15%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 80%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 13%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: orange;"></div> <div style="width: 2%; height: 10px; background-color: grey;"></div> </div> <p style="margin-left: 10px;">15% 80% 13% . . .</p>

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Mol	Chain	Length	Quality of chain
1	F	86	

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
2	IOD	B	101	-	-	X	-

2 Entry composition i

There are 3 unique types of molecules in this entry. The entry contains 4235 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 30S ribosomal protein S6,30S ribosomal protein S6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	83	Total 679	C 425	N 123	O 129	S 2	0	0	0
1	B	86	Total 708	C 441	N 130	O 134	S 3	0	0	0
1	C	85	Total 700	C 436	N 129	O 133	S 2	0	0	0
1	D	85	Total 696	C 435	N 126	O 132	S 3	0	0	0
1	E	83	Total 679	C 425	N 123	O 129	S 2	0	0	0
1	F	85	Total 700	C 436	N 129	O 133	S 2	0	0	0

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	initiating methionine	UNP Q5SLP8
A	36	ALA	-	linker	UNP Q5SLP8
A	37	SER	-	linker	UNP Q5SLP8
A	38	THR	-	linker	UNP Q5SLP8
A	39	THR	-	linker	UNP Q5SLP8
A	40	PRO	-	linker	UNP Q5SLP8
A	41	GLY	-	linker	UNP Q5SLP8
A	60	ALA	LEU	engineered mutation	UNP Q5SLP8
B	0	MET	-	initiating methionine	UNP Q5SLP8
B	36	ALA	-	linker	UNP Q5SLP8
B	37	SER	-	linker	UNP Q5SLP8
B	38	THR	-	linker	UNP Q5SLP8
B	39	THR	-	linker	UNP Q5SLP8
B	40	PRO	-	linker	UNP Q5SLP8
B	41	GLY	-	linker	UNP Q5SLP8
B	60	ALA	LEU	engineered mutation	UNP Q5SLP8
C	0	MET	-	initiating methionine	UNP Q5SLP8

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Chain	Residue	Modelled	Actual	Comment	Reference
C	36	ALA	-	linker	UNP Q5SLP8
C	37	SER	-	linker	UNP Q5SLP8
C	38	THR	-	linker	UNP Q5SLP8
C	39	THR	-	linker	UNP Q5SLP8
C	40	PRO	-	linker	UNP Q5SLP8
C	41	GLY	-	linker	UNP Q5SLP8
C	60	ALA	LEU	engineered mutation	UNP Q5SLP8
D	0	MET	-	initiating methionine	UNP Q5SLP8
D	36	ALA	-	linker	UNP Q5SLP8
D	37	SER	-	linker	UNP Q5SLP8
D	38	THR	-	linker	UNP Q5SLP8
D	39	THR	-	linker	UNP Q5SLP8
D	40	PRO	-	linker	UNP Q5SLP8
D	41	GLY	-	linker	UNP Q5SLP8
D	60	ALA	LEU	engineered mutation	UNP Q5SLP8
E	0	MET	-	initiating methionine	UNP Q5SLP8
E	36	ALA	-	linker	UNP Q5SLP8
E	37	SER	-	linker	UNP Q5SLP8
E	38	THR	-	linker	UNP Q5SLP8
E	39	THR	-	linker	UNP Q5SLP8
E	40	PRO	-	linker	UNP Q5SLP8
E	41	GLY	-	linker	UNP Q5SLP8
E	60	ALA	LEU	engineered mutation	UNP Q5SLP8
F	0	MET	-	initiating methionine	UNP Q5SLP8
F	36	ALA	-	linker	UNP Q5SLP8
F	37	SER	-	linker	UNP Q5SLP8
F	38	THR	-	linker	UNP Q5SLP8
F	39	THR	-	linker	UNP Q5SLP8
F	40	PRO	-	linker	UNP Q5SLP8
F	41	GLY	-	linker	UNP Q5SLP8
F	60	ALA	LEU	engineered mutation	UNP Q5SLP8

- Molecule 2 is IODIDE ION (CCD ID: IOD) (formula: I).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total I 1 1	0	0
2	B	1	Total I 1 1	0	0
2	C	1	Total I 1 1	0	0
2	D	1	Total I 1 1	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	E	1	Total I 1 1	0	0
2	F	1	Total I 1 1	0	0

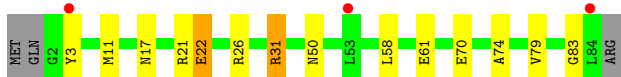
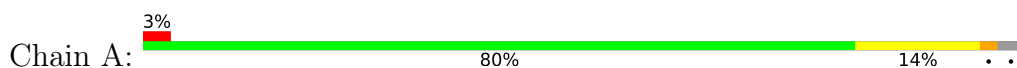
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	19	Total O 19 19	0	0
3	B	16	Total O 16 16	0	0
3	C	11	Total O 11 11	0	0
3	D	3	Total O 3 3	0	0
3	E	4	Total O 4 4	0	0
3	F	14	Total O 14 14	0	0

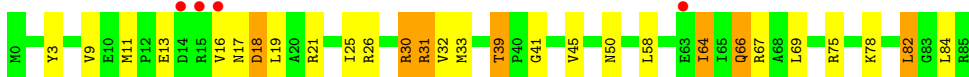
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

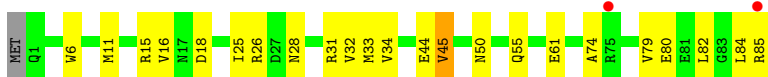
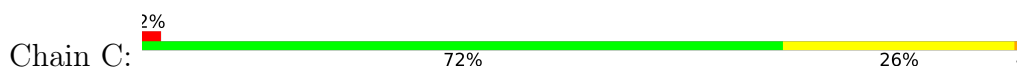
- Molecule 1: 30S ribosomal protein S6,30S ribosomal protein S6



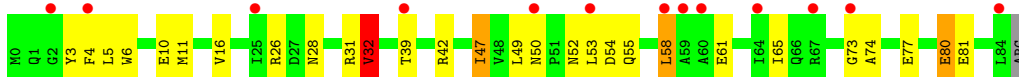
- Molecule 1: 30S ribosomal protein S6,30S ribosomal protein S6



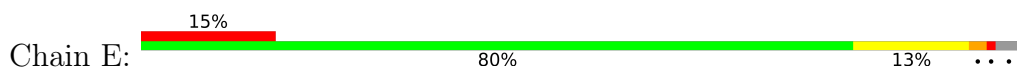
- Molecule 1: 30S ribosomal protein S6,30S ribosomal protein S6




- Molecule 1: 30S ribosomal protein S6,30S ribosomal protein S6

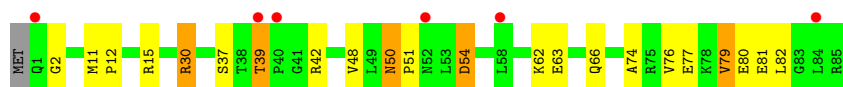


- Molecule 1: 30S ribosomal protein S6,30S ribosomal protein S6



- Molecule 1: 30S ribosomal protein S6,30S ribosomal protein S6

Chain F: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	39.00Å 67.49Å 221.95Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	64.65 – 1.95 64.65 – 1.95	Depositor EDS
% Data completeness (in resolution range)	97.5 (64.65-1.95) 97.5 (64.65-1.95)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.41 (at 1.95Å)	Xtrriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.233 , 0.280 0.238 , 0.282	Depositor DCC
R_{free} test set	2156 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å ²)	28.9	Xtrriage
Anisotropy	0.914	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 34.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4235	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 12.52% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality i

5.1 Standard geometry i

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

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	1.40	8/688 (1.2%)	1.67	10/931 (1.1%)
1	B	1.23	2/717 (0.3%)	1.55	5/967 (0.5%)
1	C	1.27	3/709 (0.4%)	1.42	2/957 (0.2%)
1	D	1.27	6/705 (0.9%)	1.49	1/953 (0.1%)
1	E	1.62	5/688 (0.7%)	1.50	4/931 (0.4%)
1	F	1.28	2/709 (0.3%)	1.43	6/957 (0.6%)
All	All	1.35	26/4216 (0.6%)	1.51	28/5696 (0.5%)

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	14	ASP	CG-OD1	26.72	1.76	1.25
1	A	79	VAL	C-O	7.91	1.32	1.23
1	C	45	VAL	C-O	7.31	1.32	1.24
1	A	31	ARG	C-O	7.28	1.33	1.23
1	D	32	VAL	C-O	7.10	1.31	1.24
1	A	22	GLU	CD-OE1	6.99	1.38	1.25
1	D	4	PHE	C-O	6.66	1.32	1.24
1	C	79	VAL	C-O	6.51	1.30	1.24
1	F	79	VAL	C-O	6.28	1.30	1.24
1	D	47	ILE	C-O	6.19	1.30	1.24
1	A	22	GLU	CG-CD	6.06	1.67	1.52
1	A	22	GLU	CB-CG	5.97	1.70	1.52
1	A	22	GLU	N-CA	5.91	1.53	1.46
1	E	45	VAL	C-O	5.75	1.30	1.24
1	B	31	ARG	C-O	5.74	1.30	1.23
1	A	22	GLU	CA-CB	5.74	1.62	1.53
1	E	14	ASP	CG-OD2	5.68	1.36	1.25
1	B	45	VAL	C-O	5.60	1.30	1.24
1	D	80	GLU	C-O	5.59	1.30	1.23
1	C	44	GLU	C-O	5.34	1.30	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	48	VAL	C-O	5.30	1.29	1.24
1	E	6	TRP	C-O	5.20	1.29	1.24
1	E	48	VAL	C-O	5.17	1.29	1.24
1	D	81	GLU	C-O	5.10	1.30	1.23
1	A	22	GLU	C-O	-5.09	1.18	1.24
1	D	6	TRP	C-O	5.03	1.30	1.24

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	22	GLU	CB-CG-CD	10.91	131.15	112.60
1	A	21	ARG	CA-C-O	-10.84	109.73	120.90
1	A	22	GLU	N-CA-C	-10.27	98.90	111.33
1	B	39	THR	CB-CA-C	9.78	120.99	109.85
1	E	10	GLU	CB-CA-C	8.42	124.14	110.16
1	B	17	ASN	CB-CA-C	-8.16	96.97	110.85
1	B	66	GLN	CB-CG-CD	7.54	125.42	112.60
1	F	50	ASN	CB-CA-C	7.49	121.18	109.42
1	A	22	GLU	CA-CB-CG	7.37	128.84	114.10
1	D	54	ASP	CA-CB-CG	7.06	119.66	112.60
1	F	39	THR	CB-CA-C	-6.75	99.31	108.68
1	A	22	GLU	N-CA-CB	6.52	119.83	110.06
1	A	50	ASN	CA-CB-CG	6.51	119.11	112.60
1	C	50	ASN	CB-CA-C	6.06	119.10	109.52
1	C	18	ASP	N-CA-C	-6.05	104.68	111.28
1	A	22	GLU	CG-CD-OE1	6.01	132.22	118.40
1	E	14	ASP	OD1-CG-OD2	5.84	136.92	122.90
1	B	18	ASP	N-CA-C	-5.59	105.27	111.36
1	A	21	ARG	CA-C-N	5.56	128.00	120.38
1	A	21	ARG	C-N-CA	5.56	128.00	120.38
1	F	51	PRO	N-CA-C	5.33	120.77	113.84
1	A	83	GLY	CA-C-O	-5.32	115.96	122.82
1	F	54	ASP	CA-C-N	5.31	127.66	120.38
1	F	54	ASP	C-N-CA	5.31	127.66	120.38
1	B	30	ARG	N-CA-C	-5.15	107.55	113.88
1	E	13	GLU	CA-C-N	5.11	128.86	120.63
1	E	13	GLU	C-N-CA	5.11	128.86	120.63
1	F	30	ARG	CB-CA-C	5.03	117.83	109.38

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	A	679	0	679	7	0
1	B	708	0	712	20	0
1	C	700	0	703	18	0
1	D	696	0	699	18	1
1	E	679	0	679	5	1
1	F	700	0	703	12	0
2	A	1	0	0	1	0
2	B	1	0	0	3	0
2	C	1	0	0	1	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
3	A	19	0	0	0	0
3	B	16	0	0	1	0
3	C	11	0	0	1	0
3	D	3	0	0	1	0
3	E	4	0	0	0	0
3	F	14	0	0	0	0
All	All	4235	0	4175	65	1

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 8.

All (65) 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:E:14:ASP:OD1	1:E:14:ASP:CG	1.76	1.29
1:C:31:ARG:NH1	1:F:81:GLU:OE1	1.92	1.03
1:F:37:SER:HB3	1:F:42:ARG:HB3	1.60	0.83
1:C:84:LEU:HD21	1:D:16:VAL:HG11	1.63	0.80
1:D:50:ASN:HD21	1:D:52:ASN:HB2	1.60	0.67
1:C:15:ARG:HD2	3:C:201:HOH:O	1.95	0.65
1:B:26:ARG:NE	1:B:64:ILE:HD11	2.13	0.63
1:B:13:GLU:O	1:B:16:VAL:HG12	2.02	0.60
1:B:11:MET:HE1	1:B:19:LEU:HD22	1.86	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:31:ARG:NH2	2:C:101:IOD:I	3.05	0.58
1:B:3:TYR:CG	1:B:58:LEU:HD11	2.38	0.57
1:D:3:TYR:CG	1:D:58:LEU:HD11	2.40	0.56
1:C:28:ASN:HD22	1:C:61:GLU:HG2	1.69	0.56
1:B:18:ASP:OD1	1:B:21:ARG:NH1	2.35	0.56
1:B:32:VAL:O	1:E:80:GLU:HA	2.07	0.55
1:F:54:ASP:OD1	1:F:54:ASP:C	2.49	0.55
1:D:26:ARG:HD3	1:D:61:GLU:OE1	2.07	0.55
1:C:16:VAL:CG1	1:C:34:VAL:HG11	2.38	0.53
1:C:11:MET:HB3	1:C:74:ALA:HA	1.91	0.53
1:B:75:ARG:NH2	3:B:201:HOH:O	2.41	0.52
1:D:28:ASN:OD1	1:D:53:LEU:HD21	2.10	0.52
1:B:31:ARG:HD3	2:B:101:IOD:I	2.81	0.50
1:C:84:LEU:HD21	1:D:16:VAL:CG1	2.39	0.50
1:A:3:TYR:CE2	1:A:58:LEU:HD22	2.47	0.50
1:D:80:GLU:HA	1:E:32:VAL:O	2.11	0.50
1:B:11:MET:O	1:B:41:GLY:HA2	2.12	0.49
1:B:82:LEU:HD22	1:B:84:LEU:HG	1.93	0.49
1:D:11:MET:HB3	1:D:74:ALA:HA	1.94	0.49
1:A:17:ASN:HD22	1:B:82:LEU:HD11	1.78	0.49
1:D:5:LEU:HD11	1:D:49:LEU:HD11	1.93	0.49
1:D:47:ILE:HD13	1:D:65:ILE:HD13	1.94	0.49
1:A:17:ASN:ND2	1:B:82:LEU:CD1	2.76	0.49
1:B:30:ARG:HG3	1:B:50:ASN:OD1	2.13	0.48
1:A:17:ASN:ND2	1:B:82:LEU:HD11	2.28	0.48
1:B:31:ARG:CD	2:B:101:IOD:I	3.32	0.47
1:C:33:MET:HE3	1:F:81:GLU:OE2	2.14	0.47
1:C:26:ARG:NH1	1:C:61:GLU:OE2	2.37	0.47
1:B:9:VAL:HA	1:B:75:ARG:O	2.14	0.47
1:C:28:ASN:HD22	1:C:61:GLU:CG	2.27	0.47
1:D:55:GLN:HA	1:D:55:GLN:OE1	2.15	0.46
1:D:31:ARG:NE	3:D:201:HOH:O	2.48	0.46
1:B:66:GLN:HA	1:B:69:LEU:HD12	1.99	0.45
1:C:85:ARG:NH1	1:D:42:ARG:O	2.50	0.45
1:D:3:TYR:CZ	1:D:58:LEU:HD21	2.52	0.45
1:D:10:GLU:OE1	1:D:42:ARG:HD3	2.18	0.44
1:A:11:MET:HB3	1:A:74:ALA:HA	1.99	0.44
1:F:30:ARG:HB3	1:F:50:ASN:OD1	2.18	0.44
1:F:42:ARG:NE	1:F:77:GLU:OE2	2.49	0.44
1:C:33:MET:HG3	1:F:81:GLU:HB2	1.99	0.44
1:C:80:GLU:HA	1:D:32:VAL:O	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:31:ARG:NH2	2:A:101:IOD:I	3.20	0.43
1:C:32:VAL:O	1:F:80:GLU:HA	2.19	0.43
1:B:33:MET:HE2	2:B:101:IOD:I	2.88	0.43
1:C:25:ILE:HD12	1:C:26:ARG:N	2.34	0.43
1:A:26:ARG:O	1:B:78:LYS:NZ	2.47	0.42
1:C:33:MET:CE	1:F:62:LYS:NZ	2.83	0.42
1:E:49:LEU:HD22	1:E:61:GLU:HG2	2.02	0.42
1:B:30:ARG:CG	1:B:50:ASN:OD1	2.68	0.41
1:F:12:PRO:HB2	1:F:15:ARG:HG3	2.03	0.41
1:F:11:MET:HB3	1:F:74:ALA:HA	2.02	0.41
1:E:26:ARG:CZ	1:E:26:ARG:HB3	2.50	0.41
1:D:3:TYR:CE2	1:D:58:LEU:HD21	2.55	0.41
1:C:6:TRP:HA	1:C:45:VAL:O	2.21	0.40
1:F:76:VAL:HG11	1:F:79:VAL:CG2	2.51	0.40
1:D:42:ARG:HD2	1:D:77:GLU:OE1	2.21	0.40

All (1) 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:D:73:GLY:O	1:E:75:ARG:NH2[3_445]	2.13	0.07

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	81/86 (94%)	80 (99%)	1 (1%)	0	100	100
1	B	84/86 (98%)	83 (99%)	1 (1%)	0	100	100
1	C	83/86 (96%)	82 (99%)	1 (1%)	0	100	100
1	D	83/86 (96%)	78 (94%)	5 (6%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	E	81/86 (94%)	79 (98%)	2 (2%)	0	100	100
1	F	83/86 (96%)	79 (95%)	3 (4%)	1 (1%)	10	4
All	All	495/516 (96%)	481 (97%)	13 (3%)	1 (0%)	43	36

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	2	GLY

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	73/76 (96%)	70 (96%)	3 (4%)	27	17
1	B	76/76 (100%)	71 (93%)	5 (7%)	15	5
1	C	75/76 (99%)	73 (97%)	2 (3%)	39	32
1	D	75/76 (99%)	72 (96%)	3 (4%)	28	17
1	E	73/76 (96%)	68 (93%)	5 (7%)	14	5
1	F	75/76 (99%)	71 (95%)	4 (5%)	20	9
All	All	447/456 (98%)	425 (95%)	22 (5%)	22	11

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	GLU
1	A	61	GLU
1	A	70	GLU
1	B	25	ILE
1	B	39	THR
1	B	64	ILE
1	B	67	ARG
1	B	82	LEU
1	C	55	GLN

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Mol	Chain	Res	Type
1	C	82	LEU
1	D	32	VAL
1	D	39	THR
1	D	58	LEU
1	E	13	GLU
1	E	14	ASP
1	E	26	ARG
1	E	46	ASN
1	E	63	GLU
1	F	39	THR
1	F	63	GLU
1	F	66	GLN
1	F	82	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	17	ASN
1	B	17	ASN
1	B	52	ASN
1	B	55	GLN
1	C	28	ASN
1	C	46	ASN
1	D	50	ASN
1	E	46	ASN
1	E	71	ASN
1	F	66	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 6 ligands modelled in this entry, 6 are 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

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	83/86 (96%)	0.35	3 (3%) 46 53	25, 34, 51, 66	0
1	B	86/86 (100%)	0.58	4 (4%) 36 42	24, 41, 64, 72	0
1	C	85/86 (98%)	0.33	2 (2%) 59 66	28, 36, 56, 78	0
1	D	85/86 (98%)	1.34	13 (15%) 5 6	32, 53, 81, 95	0
1	E	83/86 (96%)	1.13	13 (15%) 5 5	27, 51, 76, 86	0
1	F	85/86 (98%)	0.57	6 (7%) 22 25	27, 39, 59, 73	0
All	All	507/516 (98%)	0.72	41 (8%) 18 21	24, 41, 72, 95	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	25	ILE	5.4
1	F	40	PRO	4.1
1	E	84	LEU	3.9
1	D	60	ALA	3.7
1	D	2	GLY	3.6
1	D	84	LEU	3.5
1	B	16	VAL	3.5
1	F	39	THR	3.4
1	A	84	LEU	3.4
1	D	53	LEU	3.4
1	E	25	ILE	3.4
1	E	83	GLY	2.8
1	D	4	PHE	2.8
1	E	72	TYR	2.7
1	E	35	VAL	2.7
1	B	15	ARG	2.7
1	D	73	GLY	2.6
1	E	17	ASN	2.6
1	D	59	ALA	2.6

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Mol	Chain	Res	Type	RSRZ
1	D	58	LEU	2.6
1	F	52	ASN	2.6
1	A	53	LEU	2.5
1	F	58	LEU	2.5
1	E	23	LEU	2.4
1	D	64	ILE	2.4
1	A	3	TYR	2.3
1	D	50	ASN	2.3
1	E	21	ARG	2.3
1	E	69	LEU	2.3
1	C	75	ARG	2.2
1	E	53	LEU	2.2
1	E	67	ARG	2.2
1	E	74	ALA	2.2
1	D	39	THR	2.1
1	C	85	ARG	2.1
1	D	67	ARG	2.1
1	B	14	ASP	2.1
1	E	71	ASN	2.1
1	F	1	GLN	2.1
1	B	63	GLU	2.0
1	F	84	LEU	2.0

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 oligosaccharides in this entry.

6.4 Ligands [\(i\)](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	IOD	E	101	1/1	0.95	0.09	64,64,64,64	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	IOD	C	101	1/1	0.97	0.07	54,54,54,54	1
2	IOD	D	101	1/1	0.98	0.06	62,62,62,62	1
2	IOD	B	101	1/1	0.99	0.09	55,55,55,55	1
2	IOD	A	101	1/1	0.99	0.06	42,42,42,42	1
2	IOD	F	101	1/1	0.99	0.03	44,44,44,44	1

6.5 Other polymers [i](#)

There are no such residues in this entry.