



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:33 PM GMT

PDB ID : 3ZQO
Title : Crystal structure of the small terminase oligomerization core domain from a SPP1-like bacteriophage (crystal form 3)
Authors : Buttner, C.R.; Chechik, M.; Ortiz-Lombardia, M.; Smits, C.; Chechik, V.; Jeschke, G.; Dykeman, E.; Benini, S.; Alonso, J.C.; Antson, A.A.
Deposited on : 2011-06-10
Resolution : 1.68 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
<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

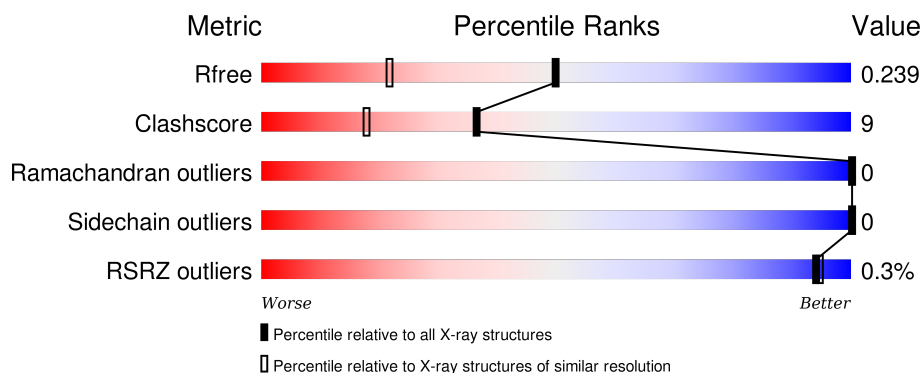
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.68 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	4802 (1.70-1.66)
Clashscore	102246	5317 (1.70-1.66)
Ramachandran outliers	100387	5225 (1.70-1.66)
Sidechain outliers	100360	5224 (1.70-1.66)
RSRZ outliers	91569	4813 (1.70-1.66)

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	A	72	
1	B	72	
1	C	72	
1	D	72	
1	E	72	

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Mol	Chain	Length	Quality of chain
1	F	72	
1	G	72	
1	H	72	
1	I	72	
1	J	72	
1	K	72	
1	L	72	
1	M	72	
1	N	72	
1	O	72	
1	P	72	
1	Q	72	
1	R	72	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 8478 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 TERMINASE SMALL SUBUNIT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	58	Total	C	N	O	Se	0	0	0
			447	279	81	86	1			
1	B	56	Total	C	N	O	Se	0	0	0
			430	270	78	81	1			
1	C	56	Total	C	N	O	Se	0	0	0
			430	270	78	81	1			
1	D	58	Total	C	N	O	Se	0	0	0
			447	279	81	86	1			
1	E	56	Total	C	N	O	Se	0	0	0
			430	270	78	81	1			
1	F	56	Total	C	N	O	Se	0	0	0
			430	270	78	81	1			
1	G	57	Total	C	N	O	Se	0	0	0
			439	275	79	84	1			
1	H	57	Total	C	N	O	Se	0	0	0
			439	275	79	84	1			
1	I	57	Total	C	N	O	Se	0	0	0
			439	275	79	84	1			
1	J	58	Total	C	N	O	Se	0	0	0
			447	279	81	86	1			
1	K	56	Total	C	N	O	Se	0	0	0
			430	270	78	81	1			
1	L	56	Total	C	N	O	Se	0	0	0
			430	270	78	81	1			
1	M	57	Total	C	N	O	Se	0	0	0
			439	275	79	84	1			
1	N	56	Total	C	N	O	Se	0	0	0
			430	270	78	81	1			
1	O	57	Total	C	N	O	Se	0	0	0
			439	275	79	84	1			
1	P	58	Total	C	N	O	Se	0	0	0
			447	279	81	86	1			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	Q	58	Total	C	N	O	Se	0	0	0
			447	279	81	86	1			
1	R	56	Total	C	N	O	Se	0	0	0
			430	270	78	81	1			

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	49	GLY	-	EXPRESSION TAG	UNP Q1EJR8
A	50	SER	-	EXPRESSION TAG	UNP Q1EJR8
A	51	HIS	-	EXPRESSION TAG	UNP Q1EJR8
A	52	MSE	-	EXPRESSION TAG	UNP Q1EJR8
B	49	GLY	-	EXPRESSION TAG	UNP Q1EJR8
B	50	SER	-	EXPRESSION TAG	UNP Q1EJR8
B	51	HIS	-	EXPRESSION TAG	UNP Q1EJR8
B	52	MSE	-	EXPRESSION TAG	UNP Q1EJR8
C	49	GLY	-	EXPRESSION TAG	UNP Q1EJR8
C	50	SER	-	EXPRESSION TAG	UNP Q1EJR8
C	51	HIS	-	EXPRESSION TAG	UNP Q1EJR8
C	52	MSE	-	EXPRESSION TAG	UNP Q1EJR8
D	49	GLY	-	EXPRESSION TAG	UNP Q1EJR8
D	50	SER	-	EXPRESSION TAG	UNP Q1EJR8
D	51	HIS	-	EXPRESSION TAG	UNP Q1EJR8
D	52	MSE	-	EXPRESSION TAG	UNP Q1EJR8
E	49	GLY	-	EXPRESSION TAG	UNP Q1EJR8
E	50	SER	-	EXPRESSION TAG	UNP Q1EJR8
E	51	HIS	-	EXPRESSION TAG	UNP Q1EJR8
E	52	MSE	-	EXPRESSION TAG	UNP Q1EJR8
F	49	GLY	-	EXPRESSION TAG	UNP Q1EJR8
F	50	SER	-	EXPRESSION TAG	UNP Q1EJR8
F	51	HIS	-	EXPRESSION TAG	UNP Q1EJR8
F	52	MSE	-	EXPRESSION TAG	UNP Q1EJR8
G	49	GLY	-	EXPRESSION TAG	UNP Q1EJR8
G	50	SER	-	EXPRESSION TAG	UNP Q1EJR8
G	51	HIS	-	EXPRESSION TAG	UNP Q1EJR8
G	52	MSE	-	EXPRESSION TAG	UNP Q1EJR8
H	49	GLY	-	EXPRESSION TAG	UNP Q1EJR8
H	50	SER	-	EXPRESSION TAG	UNP Q1EJR8
H	51	HIS	-	EXPRESSION TAG	UNP Q1EJR8
H	52	MSE	-	EXPRESSION TAG	UNP Q1EJR8
I	49	GLY	-	EXPRESSION TAG	UNP Q1EJR8
I	50	SER	-	EXPRESSION TAG	UNP Q1EJR8

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Chain	Residue	Modelled	Actual	Comment	Reference
I	51	HIS	-	EXPRESSION TAG	UNP Q1EJR8
I	52	MSE	-	EXPRESSION TAG	UNP Q1EJR8
J	49	GLY	-	EXPRESSION TAG	UNP Q1EJR8
J	50	SER	-	EXPRESSION TAG	UNP Q1EJR8
J	51	HIS	-	EXPRESSION TAG	UNP Q1EJR8
J	52	MSE	-	EXPRESSION TAG	UNP Q1EJR8
K	49	GLY	-	EXPRESSION TAG	UNP Q1EJR8
K	50	SER	-	EXPRESSION TAG	UNP Q1EJR8
K	51	HIS	-	EXPRESSION TAG	UNP Q1EJR8
K	52	MSE	-	EXPRESSION TAG	UNP Q1EJR8
L	49	GLY	-	EXPRESSION TAG	UNP Q1EJR8
L	50	SER	-	EXPRESSION TAG	UNP Q1EJR8
L	51	HIS	-	EXPRESSION TAG	UNP Q1EJR8
L	52	MSE	-	EXPRESSION TAG	UNP Q1EJR8
M	49	GLY	-	EXPRESSION TAG	UNP Q1EJR8
M	50	SER	-	EXPRESSION TAG	UNP Q1EJR8
M	51	HIS	-	EXPRESSION TAG	UNP Q1EJR8
M	52	MSE	-	EXPRESSION TAG	UNP Q1EJR8
N	49	GLY	-	EXPRESSION TAG	UNP Q1EJR8
N	50	SER	-	EXPRESSION TAG	UNP Q1EJR8
N	51	HIS	-	EXPRESSION TAG	UNP Q1EJR8
N	52	MSE	-	EXPRESSION TAG	UNP Q1EJR8
O	49	GLY	-	EXPRESSION TAG	UNP Q1EJR8
O	50	SER	-	EXPRESSION TAG	UNP Q1EJR8
O	51	HIS	-	EXPRESSION TAG	UNP Q1EJR8
O	52	MSE	-	EXPRESSION TAG	UNP Q1EJR8
P	49	GLY	-	EXPRESSION TAG	UNP Q1EJR8
P	50	SER	-	EXPRESSION TAG	UNP Q1EJR8
P	51	HIS	-	EXPRESSION TAG	UNP Q1EJR8
P	52	MSE	-	EXPRESSION TAG	UNP Q1EJR8
Q	49	GLY	-	EXPRESSION TAG	UNP Q1EJR8
Q	50	SER	-	EXPRESSION TAG	UNP Q1EJR8
Q	51	HIS	-	EXPRESSION TAG	UNP Q1EJR8
Q	52	MSE	-	EXPRESSION TAG	UNP Q1EJR8
R	49	GLY	-	EXPRESSION TAG	UNP Q1EJR8
R	50	SER	-	EXPRESSION TAG	UNP Q1EJR8
R	51	HIS	-	EXPRESSION TAG	UNP Q1EJR8
R	52	MSE	-	EXPRESSION TAG	UNP Q1EJR8

- Molecule 2 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	R	1	Total K 1 1	0	0
2	K	1	Total K 1 1	0	0

- Molecule 3 is water.

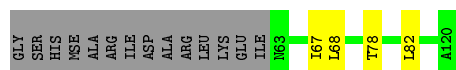
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	52	Total O 52 52	0	0
3	B	33	Total O 33 33	0	0
3	C	32	Total O 32 32	0	0
3	D	39	Total O 39 39	0	0
3	E	32	Total O 32 32	0	0
3	F	32	Total O 32 32	0	0
3	G	39	Total O 39 39	0	0
3	H	35	Total O 36 36	0	1
3	I	27	Total O 27 27	0	0
3	J	46	Total O 47 47	0	1
3	K	21	Total O 21 21	0	0
3	L	37	Total O 37 37	0	0
3	M	33	Total O 33 33	0	0
3	N	20	Total O 20 20	0	0
3	O	33	Total O 33 33	0	0
3	P	49	Total O 49 49	0	0
3	Q	25	Total O 25 25	0	0
3	R	19	Total O 19 19	0	0

3 Residue-property plots [i](#)

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

• Molecule 1: TERMINASE SMALL SUBUNIT

Chain A: 



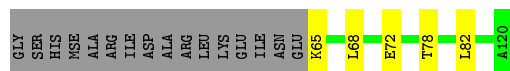
• Molecule 1: TERMINASE SMALL SUBUNIT

Chain B: 



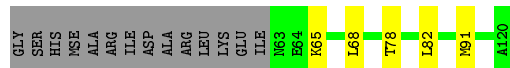
• Molecule 1: TERMINASE SMALL SUBUNIT

Chain C: 



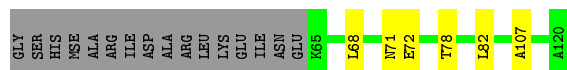
• Molecule 1: TERMINASE SMALL SUBUNIT

Chain D: 



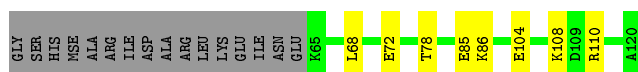
• Molecule 1: TERMINASE SMALL SUBUNIT

Chain E: 

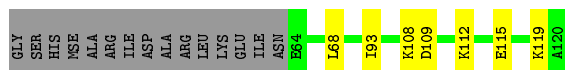


• Molecule 1: TERMINASE SMALL SUBUNIT

Chain F: 



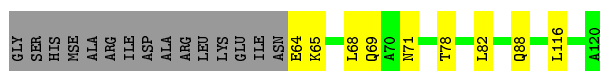
• Molecule 1: TERMINASE SMALL SUBUNIT



• Molecule 1: TERMINASE SMALL SUBUNIT



• Molecule 1: TERMINASE SMALL SUBUNIT



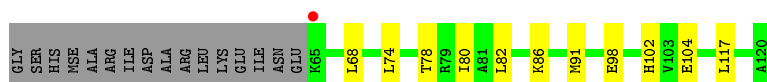
• Molecule 1: TERMINASE SMALL SUBUNIT



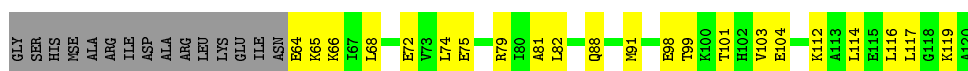
• Molecule 1: TERMINASE SMALL SUBUNIT



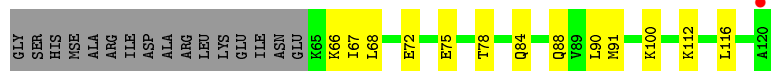
• Molecule 1: TERMINASE SMALL SUBUNIT



• Molecule 1: TERMINASE SMALL SUBUNIT



- Molecule 1: TERMINASE SMALL SUBUNIT



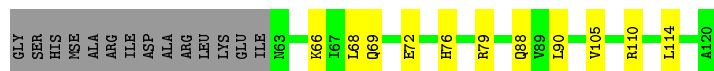
- Molecule 1: TERMINASE SMALL SUBUNIT



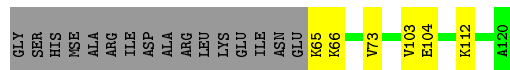
- Molecule 1: TERMINASE SMALL SUBUNIT



- Molecule 1: TERMINASE SMALL SUBUNIT



- Molecule 1: TERMINASE SMALL SUBUNIT



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	59.20 Å 198.20 Å 59.20 Å 90.00° 90.81° 90.00°	Depositor
Resolution (Å)	22.10 – 1.68 22.10 – 1.68	Depositor EDS
% Data completeness (in resolution range)	83.0 (22.10-1.68) 83.1 (22.10-1.68)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.18 (at 1.69 Å)	Xtriage
Refinement program	REFMAC 5.6.0116	Depositor
R, R_{free}	0.183 , 0.234 0.187 , 0.239	Depositor DCC
R_{free} test set	1279 reflections (1.01%)	DCC
Wilson B-factor (Å ²)	16.8	Xtriage
Anisotropy	0.429	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 34.0	EDS
Estimated twinning fraction	0.534 for H, K, L 0.466 for L, -K, H 0.053 for l,k,-h 0.055 for h,-k,-l 0.439 for l,-k,h	Xtriage
Reported twinning fraction	0.534 for H, K, L 0.466 for L, -K, H	Depositor
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.26$	Xtriage
Outliers	0 of 127549 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	8478	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.27% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

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

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	0.73	0/447	0.76	0/592
1	B	0.73	0/430	0.71	0/569
1	C	0.71	0/430	0.76	0/569
1	D	0.68	0/447	0.77	0/592
1	E	0.69	0/430	0.69	0/569
1	F	0.69	0/430	0.73	0/569
1	G	0.70	0/439	0.74	0/581
1	H	0.69	0/439	0.70	0/581
1	I	0.68	0/439	0.74	0/581
1	J	0.68	0/447	0.73	0/592
1	K	0.68	0/430	0.76	0/569
1	L	0.67	0/430	0.69	0/569
1	M	0.71	0/439	0.67	0/581
1	N	0.66	0/430	0.70	0/569
1	O	0.68	0/439	0.79	0/581
1	P	0.67	0/447	0.76	0/592
1	Q	0.67	0/447	0.71	0/592
1	R	0.68	0/430	0.74	0/569
All	All	0.69	0/7870	0.73	0/10417

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	447	0	479	5	0
1	B	430	0	467	4	0
1	C	430	0	467	5	0
1	D	447	0	479	9	0
1	E	430	0	467	14	0
1	F	430	0	467	16	0
1	G	439	0	473	10	0
1	H	439	0	473	8	0
1	I	439	0	473	12	0
1	J	447	0	479	12	0
1	K	430	0	467	10	0
1	L	430	0	467	15	0
1	M	439	0	473	27	0
1	N	430	0	467	16	0
1	O	439	0	473	15	0
1	P	447	0	479	15	0
1	Q	447	0	479	18	0
1	R	430	0	467	10	0
2	K	1	0	0	0	0
2	R	1	0	0	0	0
3	A	52	0	0	0	0
3	B	33	0	0	0	0
3	C	32	0	0	1	0
3	D	39	0	0	0	0
3	E	32	0	0	1	0
3	F	32	0	0	2	0
3	G	39	0	0	5	0
3	H	36	0	0	0	0
3	I	27	0	0	0	0
3	J	47	0	0	8	0
3	K	21	0	0	0	0
3	L	37	0	0	0	0
3	M	33	0	0	4	0
3	N	20	0	0	0	0
3	O	33	0	0	3	0
3	P	49	0	0	4	0
3	Q	25	0	0	3	0
3	R	19	0	0	1	0
All	All	8478	0	8496	155	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 9.

All (155) 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:M:101:THR:HG23	3:M:2022:HOH:O	1.44	1.14
1:D:82:LEU:HD21	1:E:68:LEU:HD21	1.39	1.04
1:M:74:LEU:HB3	1:N:67:ILE:HD13	1.42	0.97
1:A:67:ILE:HD11	1:I:71:ASN:OD1	1.67	0.94
3:M:2022:HOH:O	1:N:100:LYS:HE2	1.75	0.86
1:L:91:MSE:HE3	1:M:98:GLU:O	1.77	0.85
1:J:79:ARG:HE	1:K:65:LYS:HE3	1.46	0.80
1:M:103:VAL:HG13	3:M:2022:HOH:O	1.82	0.78
1:Q:79:ARG:HG3	1:R:65:LYS:HE2	1.65	0.77
1:J:88:GLN:HE21	1:R:104:GLU:H	1.34	0.75
1:N:75:GLU:HG3	1:O:67:ILE:HD11	1.69	0.75
1:P:66:LYS:HD3	3:P:2004:HOH:O	1.88	0.73
1:F:68:LEU:HD22	1:F:72:GLU:HG2	1.71	0.73
1:D:78:THR:HG23	1:E:68:LEU:HG	1.71	0.72
1:R:65:LYS:HG2	1:R:66:LYS:N	2.05	0.71
1:N:91:MSE:HG3	1:O:90:LEU:HD22	1.73	0.71
1:Q:68:LEU:HD22	1:Q:72:GLU:HG3	1.72	0.70
1:J:71:ASN:OD1	1:J:72:GLU:N	2.25	0.70
1:A:78:THR:O	1:A:82:LEU:HD13	1.91	0.69
1:N:68:LEU:HD22	1:N:72:GLU:HB3	1.77	0.65
1:Q:105:VAL:O	1:Q:110:ARG:NH1	2.29	0.65
1:K:65:LYS:HG2	1:K:66:LYS:N	2.11	0.65
1:I:69:GLN:NE2	1:I:71:ASN:HB3	2.13	0.64
1:P:63:ASN:O	1:P:64:GLU:HG3	1.98	0.64
1:I:78:THR:O	1:I:82:LEU:HD13	1.99	0.63
1:Q:79:ARG:CG	1:R:65:LYS:HE2	2.28	0.63
1:H:75:GLU:OE2	1:I:64:GLU:HA	1.98	0.63
1:G:93:ILE:C	3:G:2021:HOH:O	2.37	0.63
1:E:82:LEU:HD11	1:F:68:LEU:HD21	1.80	0.63
1:O:86:LYS:HE3	1:O:102:HIS:HB3	1.81	0.62
1:J:88:GLN:NE2	1:R:104:GLU:H	1.96	0.62
1:F:108:LYS:HG3	3:F:2026:HOH:O	1.99	0.62
1:O:79:ARG:NH2	1:P:65:LYS:HE2	2.15	0.62
1:E:82:LEU:CD1	1:F:68:LEU:HD11	2.29	0.62
1:O:69:GLN:NE2	3:O:2002:HOH:O	2.13	0.62
1:P:66:LYS:CD	3:P:2004:HOH:O	2.44	0.61
1:G:115:GLU:O	1:G:119:LYS:HG2	2.00	0.61
1:H:69:GLN:HB2	1:H:72:GLU:OE1	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:P:91:MSE:SE	3:Q:2018:HOH:O	2.71	0.59
1:K:65:LYS:CG	1:K:66:LYS:H	2.15	0.58
1:A:68:LEU:HG	1:I:78:THR:HG23	1.85	0.58
1:D:82:LEU:HD21	1:E:68:LEU:CD2	2.26	0.58
1:K:65:LYS:HG2	1:K:66:LYS:H	1.68	0.58
1:O:64:GLU:O	1:O:65:LYS:HB2	2.04	0.58
1:L:117:LEU:HD13	1:M:68:LEU:HB2	1.87	0.57
1:M:116:LEU:HD23	1:M:119:LYS:HE3	1.87	0.57
1:F:108:LYS:HG2	1:G:109:ASP:OD1	2.05	0.57
1:Q:68:LEU:HD22	1:Q:72:GLU:CG	2.35	0.56
1:E:107:ALA:HB1	3:E:2026:HOH:O	2.04	0.56
1:M:75:GLU:HB3	1:M:79:ARG:HH12	1.71	0.56
1:L:78:THR:HG23	1:M:68:LEU:HG	1.87	0.56
1:J:69:GLN:O	1:J:73:VAL:HG23	2.05	0.56
1:J:68:LEU:HD22	1:J:72:GLU:HB3	1.87	0.56
1:Q:90:LEU:HD12	1:Q:90:LEU:N	2.22	0.55
1:E:82:LEU:HD11	1:F:68:LEU:HD11	1.89	0.55
1:G:93:ILE:HA	3:G:2021:HOH:O	2.07	0.55
1:K:65:LYS:CG	1:K:66:LYS:N	2.69	0.55
1:P:104:GLU:H	1:Q:88:GLN:HE21	1.54	0.54
1:E:78:THR:HG23	1:F:68:LEU:HG	1.89	0.53
1:L:91:MSE:CE	1:M:98:GLU:O	2.54	0.53
1:L:74:LEU:HD22	1:L:117:LEU:HD23	1.90	0.53
1:E:82:LEU:HD11	1:F:68:LEU:CG	2.38	0.53
1:H:104:GLU:H	1:I:88:GLN:NE2	2.07	0.53
1:P:63:ASN:O	1:P:64:GLU:CG	2.57	0.53
1:N:84:GLN:OE1	1:O:64:GLU:OE2	2.28	0.52
1:N:75:GLU:CG	1:O:67:ILE:HD11	2.38	0.52
1:B:77:LEU:HD22	1:B:113:ALA:HA	1.91	0.52
1:M:104:GLU:H	1:N:88:GLN:HE21	1.57	0.51
1:P:63:ASN:C	1:P:64:GLU:HG3	2.31	0.51
3:J:2020:HOH:O	1:R:103:VAL:HG11	2.11	0.51
1:R:65:LYS:HG2	1:R:66:LYS:H	1.76	0.51
1:M:68:LEU:HD22	1:M:72:GLU:HB3	1.91	0.51
1:A:67:ILE:CD1	1:I:71:ASN:OD1	2.51	0.50
1:J:63:ASN:N	3:J:2001:HOH:O	2.43	0.50
1:G:112:LYS:HG3	3:G:2035:HOH:O	2.10	0.50
1:C:78:THR:HG23	1:D:68:LEU:HG	1.93	0.50
1:F:68:LEU:HD22	1:F:72:GLU:CG	2.40	0.50
1:M:112:LYS:NZ	1:M:116:LEU:HD11	2.27	0.49
1:I:69:GLN:HE21	1:I:71:ASN:HB3	1.77	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:90:LEU:HD23	3:J:2020:HOH:O	2.12	0.49
1:E:82:LEU:HD11	1:F:68:LEU:CD2	2.43	0.49
1:G:112:LYS:NZ	1:H:112:LYS:NZ	2.61	0.48
1:P:104:GLU:H	1:Q:88:GLN:NE2	2.11	0.48
1:M:74:LEU:HD22	1:M:117:LEU:HD23	1.95	0.48
1:Q:76:HIS:HA	1:Q:79:ARG:HH11	1.79	0.48
1:F:86:LYS:HD3	1:F:104:GLU:HA	1.96	0.47
1:L:91:MSE:HE3	1:M:98:GLU:C	2.32	0.47
1:C:82:LEU:HD21	1:D:68:LEU:HD21	1.96	0.47
1:L:104:GLU:H	1:M:88:GLN:HE21	1.63	0.47
1:L:86:LYS:HE2	1:L:102:HIS:HB3	1.97	0.47
1:M:91:MSE:HG3	1:N:90:LEU:HD22	1.97	0.46
3:J:2020:HOH:O	1:R:103:VAL:CG1	2.63	0.46
1:M:91:MSE:SE	1:N:90:LEU:HD22	2.65	0.46
1:M:74:LEU:HB3	1:N:67:ILE:CD1	2.30	0.46
1:M:101:THR:CG2	3:M:2022:HOH:O	2.24	0.45
1:O:86:LYS:NZ	3:O:2016:HOH:O	2.44	0.45
1:O:93:ILE:HG21	1:O:99:THR:HG22	1.97	0.45
1:Q:90:LEU:HG	3:Q:2018:HOH:O	2.16	0.45
1:H:75:GLU:OE2	1:I:65:LYS:N	2.47	0.45
1:L:91:MSE:HE3	1:M:99:THR:HA	1.99	0.45
1:N:112:LYS:NZ	1:N:116:LEU:HD21	2.32	0.45
1:P:78:THR:HG23	1:Q:68:LEU:HG	2.00	0.45
1:O:93:ILE:CG2	1:O:99:THR:HG22	2.46	0.45
1:P:82:LEU:HD23	1:Q:76:HIS:NE2	2.32	0.44
1:N:78:THR:HG21	1:O:66:LYS:HG3	1.98	0.44
1:R:112:LYS:HD3	3:R:2018:HOH:O	2.17	0.44
1:M:81:ALA:HB1	1:M:114:LEU:HD21	2.00	0.44
1:N:84:GLN:NE2	1:O:64:GLU:OE2	2.50	0.44
1:J:81:ALA:HB1	1:J:114:LEU:CD2	2.48	0.44
1:F:78:THR:HG23	1:G:68:LEU:HG	2.00	0.44
1:K:111:ILE:HD11	1:L:80:ILE:HD13	1.99	0.44
1:A:68:LEU:HG	1:I:82:LEU:HD11	2.00	0.44
1:G:93:ILE:CA	3:G:2021:HOH:O	2.66	0.44
1:L:82:LEU:HD11	1:M:66:LYS:HE2	2.00	0.44
1:H:115:GLU:HG3	1:I:116:LEU:HD21	1.98	0.44
1:J:69:GLN:HG3	3:J:2009:HOH:O	2.17	0.44
1:H:78:THR:HG23	1:I:68:LEU:HG	2.00	0.43
1:B:79:ARG:CG	1:C:65:LYS:HE3	2.47	0.43
1:E:82:LEU:HD12	1:F:68:LEU:HD11	1.98	0.43
3:J:2045:HOH:O	1:K:112:LYS:HE2	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:78:THR:HG22	1:L:82:LEU:HD12	2.01	0.43
1:M:82:LEU:HD11	1:N:66:LYS:HE2	2.00	0.43
1:L:104:GLU:H	1:M:88:GLN:NE2	2.16	0.43
3:C:2006:HOH:O	1:D:65:LYS:HB2	2.18	0.43
1:O:65:LYS:HD3	1:O:65:LYS:HA	1.60	0.43
1:P:82:LEU:HD11	1:Q:66:LYS:HZ3	1.84	0.43
1:F:85:GLU:HG3	1:F:86:LYS:N	2.34	0.43
1:C:68:LEU:HD22	1:C:72:GLU:CB	2.49	0.43
1:Q:69:GLN:O	1:Q:72:GLU:HG2	2.20	0.42
1:M:81:ALA:HB1	1:M:114:LEU:CD2	2.50	0.42
1:B:69:GLN:OE1	1:B:72:GLU:OE1	2.36	0.42
1:K:117:LEU:HD13	1:L:68:LEU:HB2	2.00	0.42
1:D:91:MSE:HB3	1:D:91:MSE:HE2	1.92	0.42
1:Q:114:LEU:HB3	1:R:73:VAL:HG13	2.01	0.42
1:M:64:GLU:HG2	1:M:65:LYS:H	1.84	0.42
1:E:82:LEU:HD11	1:F:68:LEU:CD1	2.49	0.42
1:K:91:MSE:HG3	1:L:98:GLU:HB2	2.01	0.42
1:Q:90:LEU:CD1	1:Q:90:LEU:N	2.83	0.41
1:J:87:GLU:HA	3:J:2017:HOH:O	2.19	0.41
1:O:67:ILE:O	3:O:2002:HOH:O	2.21	0.41
1:G:112:LYS:HZ3	1:H:112:LYS:NZ	2.18	0.41
1:J:120:ALA:HA	3:J:2047:HOH:O	2.19	0.41
1:K:68:LEU:HG	1:K:72:GLU:HB3	2.02	0.41
1:D:78:THR:O	1:D:82:LEU:HG	2.21	0.41
1:P:66:LYS:NZ	3:P:2004:HOH:O	2.29	0.41
1:C:68:LEU:HA	1:C:68:LEU:HD23	1.83	0.41
1:P:108:LYS:HE2	3:P:2044:HOH:O	2.21	0.41
1:Q:110:ARG:HD2	3:Q:2005:HOH:O	2.19	0.41
1:G:108:LYS:HG3	3:G:2032:HOH:O	2.21	0.41
1:B:74:LEU:HD22	1:B:117:LEU:HD23	2.03	0.41
1:D:82:LEU:CD2	1:E:68:LEU:HD21	2.28	0.41
1:M:104:GLU:H	1:N:88:GLN:NE2	2.18	0.40
1:F:110:ARG:NH1	3:F:2025:HOH:O	2.54	0.40
1:P:104:GLU:HG3	1:Q:88:GLN:NE2	2.36	0.40
1:E:71:ASN:OD1	1:E:72:GLU:N	2.55	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	56/72 (78%)	55 (98%)	1 (2%)	0	100	100
1	B	54/72 (75%)	54 (100%)	0	0	100	100
1	C	54/72 (75%)	53 (98%)	1 (2%)	0	100	100
1	D	56/72 (78%)	56 (100%)	0	0	100	100
1	E	54/72 (75%)	54 (100%)	0	0	100	100
1	F	54/72 (75%)	53 (98%)	1 (2%)	0	100	100
1	G	55/72 (76%)	55 (100%)	0	0	100	100
1	H	55/72 (76%)	55 (100%)	0	0	100	100
1	I	55/72 (76%)	54 (98%)	1 (2%)	0	100	100
1	J	56/72 (78%)	56 (100%)	0	0	100	100
1	K	54/72 (75%)	54 (100%)	0	0	100	100
1	L	54/72 (75%)	52 (96%)	2 (4%)	0	100	100
1	M	55/72 (76%)	55 (100%)	0	0	100	100
1	N	54/72 (75%)	53 (98%)	1 (2%)	0	100	100
1	O	55/72 (76%)	55 (100%)	0	0	100	100
1	P	56/72 (78%)	56 (100%)	0	0	100	100
1	Q	56/72 (78%)	56 (100%)	0	0	100	100
1	R	54/72 (75%)	53 (98%)	1 (2%)	0	100	100
All	All	987/1296 (76%)	979 (99%)	8 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	A	47/56 (84%)	47 (100%)	0	100	100
1	B	45/56 (80%)	45 (100%)	0	100	100
1	C	45/56 (80%)	45 (100%)	0	100	100
1	D	47/56 (84%)	47 (100%)	0	100	100
1	E	45/56 (80%)	45 (100%)	0	100	100
1	F	45/56 (80%)	45 (100%)	0	100	100
1	G	46/56 (82%)	46 (100%)	0	100	100
1	H	46/56 (82%)	46 (100%)	0	100	100
1	I	46/56 (82%)	46 (100%)	0	100	100
1	J	47/56 (84%)	47 (100%)	0	100	100
1	K	45/56 (80%)	45 (100%)	0	100	100
1	L	45/56 (80%)	45 (100%)	0	100	100
1	M	46/56 (82%)	46 (100%)	0	100	100
1	N	45/56 (80%)	45 (100%)	0	100	100
1	O	46/56 (82%)	46 (100%)	0	100	100
1	P	47/56 (84%)	47 (100%)	0	100	100
1	Q	47/56 (84%)	47 (100%)	0	100	100
1	R	45/56 (80%)	45 (100%)	0	100	100
All	All	825/1008 (82%)	825 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	76	HIS
1	C	84	GLN
1	I	69	GLN
1	I	88	GLN
1	J	88	GLN
1	M	69	GLN
1	M	71	ASN
1	M	88	GLN
1	N	88	GLN

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Mol	Chain	Res	Type
1	P	63	ASN
1	Q	88	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 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 [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 ⓘ

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	57/72 (79%)	-0.51	0	100	100	12, 16, 28, 41	0
1	B	55/72 (76%)	-0.56	0	100	100	12, 16, 27, 43	0
1	C	55/72 (76%)	-0.51	0	100	100	12, 16, 29, 35	0
1	D	57/72 (79%)	-0.53	0	100	100	12, 16, 29, 44	0
1	E	55/72 (76%)	-0.52	0	100	100	10, 16, 28, 32	0
1	F	55/72 (76%)	-0.46	0	100	100	10, 16, 31, 44	0
1	G	56/72 (77%)	-0.58	0	100	100	11, 16, 26, 43	0
1	H	56/72 (77%)	-0.52	0	100	100	11, 17, 28, 39	0
1	I	56/72 (77%)	-0.47	0	100	100	11, 17, 28, 42	0
1	J	57/72 (79%)	-0.52	0	100	100	14, 19, 29, 30	0
1	K	55/72 (76%)	-0.49	0	100	100	13, 18, 30, 33	0
1	L	55/72 (76%)	-0.49	1 (1%)	71	75	14, 18, 29, 36	0
1	M	56/72 (77%)	-0.45	0	100	100	13, 18, 29, 36	0
1	N	55/72 (76%)	-0.44	1 (1%)	71	75	14, 18, 32, 34	0
1	O	56/72 (77%)	-0.45	1 (1%)	71	75	13, 19, 30, 40	0
1	P	57/72 (79%)	-0.51	0	100	100	13, 18, 31, 33	0
1	Q	57/72 (79%)	-0.51	0	100	100	13, 18, 29, 43	0
1	R	55/72 (76%)	-0.48	0	100	100	14, 18, 30, 34	0
All	All	1005/1296 (77%)	-0.50	3 (0%)	94	95	10, 17, 30, 44	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	N	120	ALA	2.4
1	O	120	ALA	2.2
1	L	65	LYS	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 carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 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
2	K	K	1121	1/1	0.99	0.14	-	51,51,51,51	0
2	K	R	1121	1/1	0.99	0.09	-	26,26,26,26	0

6.5 Other polymers [i](#)

There are no such residues in this entry.