



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 06:33 AM UTC

PDB ID : 4R0N / pdb_00004r0n
Title : Hexagonal form of phosphopantetheine adenylyltransferase from *Mycobacterium tuberculosis*
Authors : Timofeev, V.I.; Smirnova, E.A.; Chupova, L.A.; Esipov, R.S.; Kuranova, I.P.
Deposited on : 2014-08-01
Resolution : 2.00 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
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (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

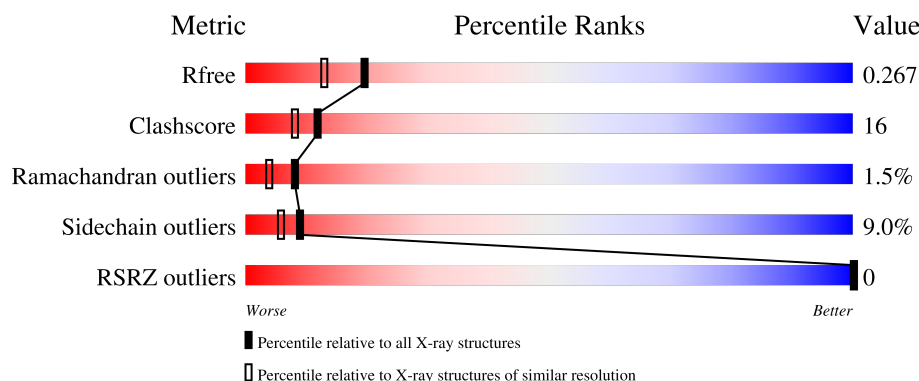
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	156	 80% 16% . .
1	C	156	 75% 19% . .
1	E	156	 78% 15% 5% .
1	G	156	 80% 15% . .
1	I	156	 81% 13% 6%

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Mol	Chain	Length	Quality of chain
1	K	156	76% 17% 6% •

2 Entry composition [i](#)

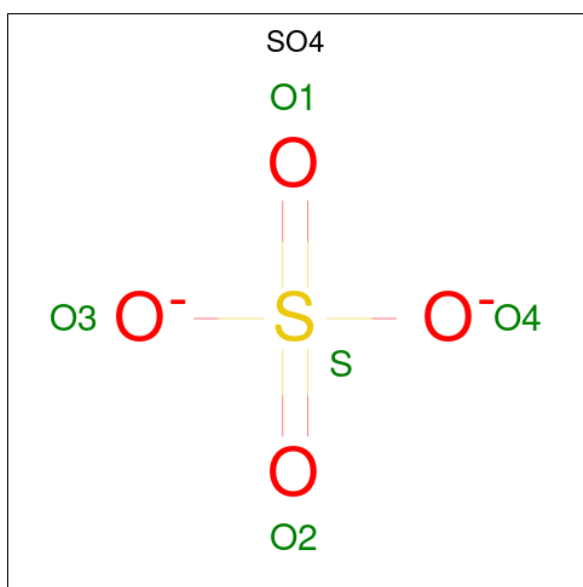
There are 3 unique types of molecules in this entry. The entry contains 7482 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Phosphopantetheine adenylyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	156	Total	C	N	O	S	0	1	0
			1200	758	209	225	8			
1	C	156	Total	C	N	O	S	0	0	0
			1193	753	208	224	8			
1	E	156	Total	C	N	O	S	0	0	0
			1193	753	208	224	8			
1	G	156	Total	C	N	O	S	0	0	0
			1193	753	208	224	8			
1	I	156	Total	C	N	O	S	0	1	0
			1200	758	209	225	8			
1	K	156	Total	C	N	O	S	0	1	0
			1200	757	209	226	8			

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total O S 5 4 1	0	0
2	C	1	Total O S 5 4 1	0	0
2	C	1	Total O S 5 4 1	0	0
2	E	1	Total O S 5 4 1	0	0
2	G	1	Total O S 5 4 1	0	0
2	G	1	Total O S 5 4 1	0	0
2	I	1	Total O S 5 4 1	0	0
2	K	1	Total O S 5 4 1	0	0

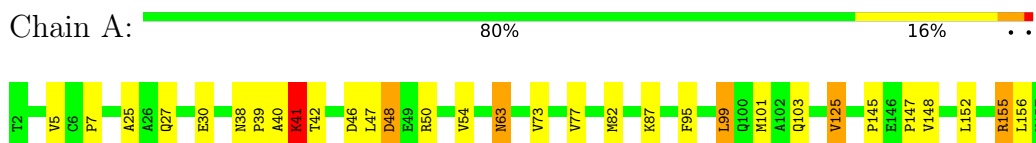
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	47	Total O 47 47	0	0
3	C	53	Total O 53 53	0	0
3	E	43	Total O 43 43	0	0
3	G	37	Total O 37 37	0	0
3	I	44	Total O 44 44	0	0
3	K	39	Total O 39 39	0	0

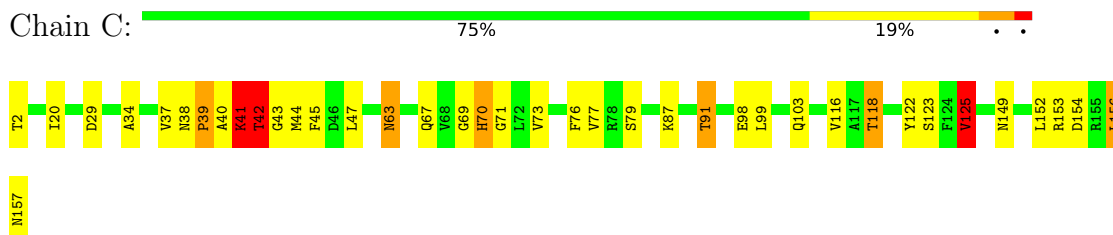
3 Residue-property plots

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.

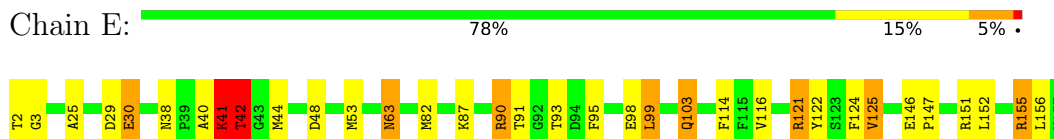
- Molecule 1: Phosphopantetheine adenylyltransferase



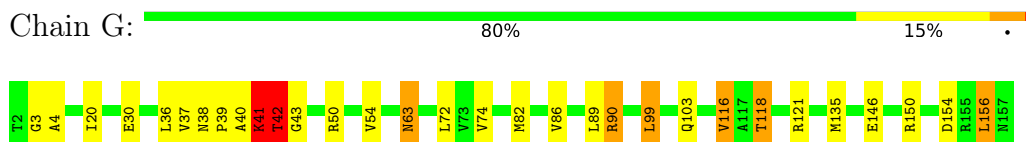
- Molecule 1: Phosphopantetheine adenylyltransferase



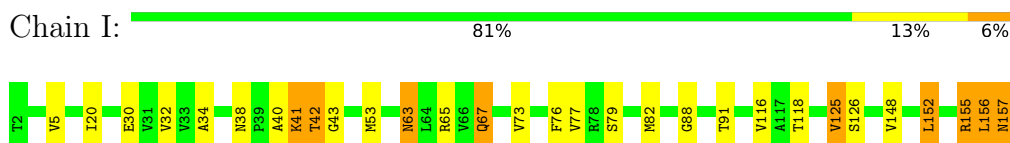
- Molecule 1: Phosphopantetheine adenylyltransferase



- Molecule 1: Phosphopantetheine adenylyltransferase



- Molecule 1: Phosphopantetheine adenylyltransferase



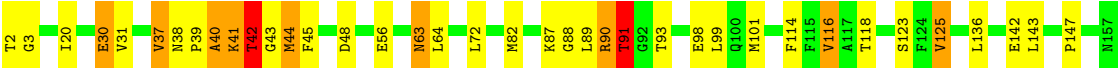
- Molecule 1: Phosphopantetheine adenylyltransferase

Chain K:

76%

17%

6% •



4 Data and refinement statistics

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, α , β , γ	106.47Å 106.47Å 71.32Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	19.64 – 2.00 19.64 – 2.00	Depositor EDS
% Data completeness (in resolution range)	94.3 (19.64-2.00) 94.4 (19.64-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.39 (at 2.01Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.210 , 0.266 0.211 , 0.267	Depositor DCC
R_{free} test set	2916 reflections (4.78%)	wwPDB-VP
Wilson B-factor (Å ²)	22.5	Xtriage
Anisotropy	0.247	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 23.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.031 for -h,-k,l 0.480 for h,-h-k,-l 0.031 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	7482	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.15% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SO4

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	0.76	0/1221	0.91	1/1654 (0.1%)
1	C	0.75	0/1214	0.90	1/1644 (0.1%)
1	E	0.78	0/1214	0.92	1/1644 (0.1%)
1	G	0.73	0/1214	0.89	1/1644 (0.1%)
1	I	0.75	0/1221	0.87	2/1654 (0.1%)
1	K	0.73	0/1221	0.93	2/1654 (0.1%)
All	All	0.75	0/7305	0.91	8/9894 (0.1%)

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	I	125	VAL	CB-CA-C	-6.62	102.35	111.49
1	E	125	VAL	CB-CA-C	-5.92	103.46	111.63
1	G	116	VAL	CB-CA-C	-5.90	101.71	110.33
1	A	125	VAL	CB-CA-C	-5.73	103.37	111.28
1	K	116	VAL	CB-CA-C	-5.31	102.58	110.33
1	I	126	SER	N-CA-C	-5.13	101.27	109.07
1	K	91	THR	CB-CA-C	5.09	118.04	109.89
1	C	125	VAL	CB-CA-C	-5.07	104.63	111.63

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	1200	0	1203	34	0
1	C	1193	0	1195	50	0
1	E	1193	0	1195	38	0
1	G	1193	0	1195	37	0
1	I	1200	0	1203	30	0
1	K	1200	0	1201	43	0
2	A	5	0	0	0	0
2	C	10	0	0	1	0
2	E	5	0	0	0	0
2	G	10	0	0	2	0
2	I	5	0	0	0	0
2	K	5	0	0	0	0
3	A	47	0	0	2	0
3	C	53	0	0	3	0
3	E	43	0	0	1	0
3	G	37	0	0	4	0
3	I	44	0	0	0	0
3	K	39	0	0	6	0
All	All	7482	0	7192	228	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 16.

All (228) 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:C:41:LYS:CD	1:C:43:GLY:H	1.55	1.17
1:A:42:THR:HA	3:A:331:HOH:O	1.45	1.15
1:C:41:LYS:CG	1:C:42:THR:H	1.60	1.13
1:G:42:THR:HB	1:G:43:GLY:CA	1.79	1.13
1:I:42:THR:HG22	1:I:43:GLY:HA2	1.14	1.12
1:G:43:GLY:H	1:G:50:ARG:NH1	1.48	1.10
1:I:42:THR:CG2	1:I:43:GLY:HA2	1.81	1.10
1:C:41:LYS:HG3	1:C:42:THR:N	1.58	1.08
1:C:41:LYS:O	1:C:42:THR:HG23	1.51	1.07
1:I:5[B]:VAL:HG11	1:I:77:VAL:HG22	1.32	1.05
1:A:5[B]:VAL:HG11	1:A:77:VAL:HG22	1.40	1.02
1:G:42:THR:CB	1:G:43:GLY:HA2	1.86	1.01
1:C:41:LYS:HD2	1:C:43:GLY:H	1.24	0.99

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:42:THR:HB	1:G:43:GLY:C	1.88	0.98
1:G:118:THR:HG21	3:G:306:HOH:O	1.65	0.97
1:A:38:ASN:O	1:A:40:ALA:HA	1.65	0.95
1:E:124:PHE:HB3	1:G:103:GLN:HE21	1.29	0.94
1:C:41:LYS:HD2	1:C:43:GLY:N	1.83	0.94
1:C:41:LYS:HG3	1:C:42:THR:H	0.79	0.93
1:G:40:ALA:O	1:G:41:LYS:HB2	1.66	0.91
1:A:38:ASN:HD21	1:A:41:LYS:HB2	1.38	0.88
1:I:155:ARG:HH11	1:I:155:ARG:CG	1.87	0.88
1:K:37:VAL:HG23	1:K:41:LYS:CB	2.05	0.87
1:I:42:THR:HG22	1:I:43:GLY:CA	2.02	0.86
1:G:63:ASN:HD22	1:G:63:ASN:H	1.21	0.86
1:C:41:LYS:CD	1:C:43:GLY:N	2.37	0.86
1:E:38:ASN:CG	1:E:41:LYS:NZ	2.34	0.85
1:E:124:PHE:HB3	1:G:103:GLN:NE2	1.90	0.85
1:K:44:MET:O	1:K:44:MET:HG3	1.76	0.85
1:I:63:ASN:HD22	1:I:63:ASN:H	1.27	0.83
1:I:42:THR:CB	1:I:43:GLY:HA2	2.05	0.82
1:E:155:ARG:CG	1:E:155:ARG:HH11	1.91	0.82
1:E:38:ASN:ND2	1:E:41:LYS:HZ1	1.77	0.81
1:K:89:LEU:O	1:K:118[B]:THR:HG23	1.81	0.81
1:C:41:LYS:HD3	1:C:43:GLY:H	1.44	0.81
1:K:37:VAL:HG23	1:K:41:LYS:HB2	1.62	0.80
1:C:63:ASN:H	1:C:63:ASN:HD22	1.30	0.80
1:K:63:ASN:HD22	1:K:63:ASN:H	1.27	0.80
1:C:41:LYS:O	1:C:42:THR:CG2	2.30	0.79
1:A:155:ARG:HG2	1:A:155:ARG:HH11	1.49	0.77
3:E:337:HOH:O	1:K:2:THR:HB	1.84	0.77
1:G:43:GLY:N	1:G:50:ARG:NH1	2.30	0.76
1:A:155:ARG:HH11	1:A:155:ARG:CG	1.98	0.76
1:K:37:VAL:HG23	1:K:41:LYS:HG3	1.67	0.76
1:G:43:GLY:H	1:G:50:ARG:HH11	1.28	0.76
1:I:155:ARG:HH11	1:I:155:ARG:HG3	1.51	0.76
1:G:40:ALA:O	1:G:41:LYS:CB	2.33	0.76
1:K:37:VAL:HG23	1:K:41:LYS:CG	2.15	0.75
1:C:41:LYS:HD2	1:C:43:GLY:CA	2.16	0.75
1:G:43:GLY:H	1:G:50:ARG:HH12	1.30	0.74
1:A:145:PRO:HB2	1:A:147:PRO:HD2	1.70	0.73
1:E:155:ARG:HH11	1:E:155:ARG:HG2	1.52	0.73
1:G:42:THR:HB	1:G:43:GLY:O	1.89	0.73
1:I:40:ALA:C	1:I:42:THR:H	1.95	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:38:ASN:ND2	1:A:41:LYS:HB2	2.02	0.73
1:A:101:MET:HE2	3:A:336:HOH:O	1.90	0.72
1:K:42:THR:HG22	1:K:42:THR:O	1.90	0.72
1:E:90:ARG:HG3	1:E:93:THR:CG2	2.21	0.71
1:E:38:ASN:CG	1:E:41:LYS:HZ3	1.96	0.71
1:C:87:LYS:HE2	1:C:98:GLU:OE2	1.91	0.70
1:E:38:ASN:ND2	1:E:41:LYS:NZ	2.40	0.69
1:E:41:LYS:O	1:E:41:LYS:HG3	1.92	0.69
1:E:40:ALA:O	1:E:41:LYS:C	2.35	0.68
1:K:41:LYS:O	1:K:41:LYS:HD3	1.93	0.67
1:G:43:GLY:N	1:G:50:ARG:HH12	1.92	0.67
1:C:41:LYS:HD2	1:C:43:GLY:HA2	1.76	0.66
1:C:37:VAL:HG22	1:C:38:ASN:N	2.10	0.66
1:K:38:ASN:HD21	1:K:40:ALA:HB3	1.59	0.66
1:C:38:ASN:OD1	1:C:40:ALA:N	2.30	0.65
1:C:38:ASN:OD1	1:C:38:ASN:C	2.40	0.64
1:I:40:ALA:O	1:I:42:THR:N	2.31	0.64
1:K:37:VAL:CG2	1:K:41:LYS:HG3	2.28	0.64
1:C:125:VAL:HG22	3:C:332:HOH:O	1.99	0.63
1:C:38:ASN:C	1:C:40:ALA:H	2.07	0.63
1:E:103:GLN:OE1	1:E:103:GLN:HA	1.98	0.63
1:C:44:MET:HE1	1:C:152:LEU:CD2	2.29	0.63
1:I:155:ARG:HH11	1:I:155:ARG:HG2	1.63	0.63
1:I:42:THR:CB	1:I:43:GLY:CA	2.76	0.62
1:K:44:MET:HG2	1:K:45:PHE:CE1	2.34	0.62
1:G:103:GLN:HG2	3:G:325:HOH:O	1.99	0.61
1:K:44:MET:HG2	1:K:45:PHE:CD1	2.34	0.61
1:K:91:THR:HG23	1:K:123:SER:O	2.00	0.61
1:E:38:ASN:CG	1:E:41:LYS:HZ1	2.01	0.61
1:E:90:ARG:H	1:E:93:THR:CG2	2.13	0.61
1:I:157:ASN:OD1	1:I:157:ASN:N	2.33	0.61
1:I:76:PHE:O	1:I:79:SER:HB2	2.01	0.60
1:A:5[B]:VAL:CG1	1:A:77:VAL:HG22	2.25	0.60
1:A:63:ASN:H	1:A:63:ASN:HD22	1.50	0.60
1:A:38:ASN:HD21	1:A:41:LYS:CB	2.13	0.59
1:K:101:MET:CE	3:K:319:HOH:O	2.49	0.59
1:G:63:ASN:HD22	1:G:63:ASN:N	1.94	0.59
1:E:63:ASN:H	1:E:63:ASN:HD22	1.50	0.59
1:G:42:THR:OG1	1:G:43:GLY:HA2	2.02	0.59
1:C:40:ALA:O	1:C:41:LYS:C	2.47	0.58
1:K:37:VAL:CG2	1:K:41:LYS:CG	2.81	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:38:ASN:OD1	1:G:39:PRO:HD2	2.04	0.57
1:I:155:ARG:CG	1:I:155:ARG:NH1	2.56	0.57
1:A:5[B]:VAL:HG11	1:A:77:VAL:CG2	2.26	0.57
1:G:89:LEU:O	1:G:118:THR:HG23	2.05	0.57
1:E:41:LYS:O	1:E:41:LYS:CG	2.53	0.57
1:C:20:ILE:HD11	1:C:118:THR:HG22	1.86	0.56
1:E:44:MET:O	1:E:155:ARG:NH2	2.39	0.56
1:E:3:GLY:HA3	1:E:82:MET:HE2	1.88	0.56
1:E:155:ARG:CG	1:E:155:ARG:NH1	2.59	0.56
1:K:37:VAL:CG2	1:K:41:LYS:CB	2.82	0.55
1:I:73:VAL:O	1:I:77:VAL:HG23	2.05	0.55
1:C:41:LYS:CG	1:C:42:THR:N	2.32	0.55
1:K:125:VAL:HG21	1:K:143:LEU:HB3	1.88	0.55
1:I:30:GLU:OE2	1:I:65:ARG:NH1	2.41	0.54
1:C:37:VAL:HG22	1:C:38:ASN:H	1.71	0.54
1:E:25:ALA:HB1	1:E:63:ASN:HD21	1.72	0.54
1:E:90:ARG:O	1:E:93:THR:HG23	2.08	0.54
2:C:202:SO4:O1	2:G:202:SO4:O2	2.25	0.54
1:E:30:GLU:HG2	1:E:82:MET:CE	2.38	0.54
1:C:149:ASN:HB3	1:C:153:ARG:HH11	1.71	0.53
1:A:155:ARG:CG	1:A:155:ARG:NH1	2.66	0.53
1:K:20:ILE:HD11	1:K:118[B]:THR:HG22	1.91	0.53
1:G:41:LYS:C	1:G:42:THR:OG1	2.51	0.53
1:C:44:MET:HE1	1:C:152:LEU:HD21	1.89	0.53
1:C:76:PHE:O	1:C:79:SER:HB2	2.09	0.53
1:I:65:ARG:HH21	1:I:67:GLN:HE22	1.58	0.52
1:E:90:ARG:HG3	1:E:93:THR:HG22	1.91	0.52
1:C:38:ASN:OD1	1:C:39:PRO:N	2.43	0.52
1:A:38:ASN:O	1:A:40:ALA:CA	2.49	0.51
1:K:39:PRO:C	1:K:41:LYS:H	2.18	0.51
1:C:44:MET:HE2	1:C:45:PHE:CE1	2.46	0.51
1:C:103:GLN:HA	1:C:103:GLN:OE1	2.10	0.51
1:E:41:LYS:O	1:E:42:THR:C	2.54	0.51
1:K:44:MET:O	1:K:45:PHE:CD2	2.63	0.51
1:A:48:ASP:OD1	1:A:48:ASP:N	2.38	0.51
1:I:63:ASN:HD22	1:I:63:ASN:N	2.01	0.51
1:K:101:MET:HE2	3:K:319:HOH:O	2.09	0.51
1:C:2:THR:HG22	1:C:29:ASP:H	1.75	0.50
1:I:155:ARG:HG2	1:I:155:ARG:NH1	2.24	0.50
1:E:155:ARG:HH11	1:E:155:ARG:HG3	1.74	0.50
1:C:41:LYS:O	1:C:42:THR:CB	2.58	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:40:ALA:O	1:C:42:THR:N	2.45	0.50
1:I:65:ARG:HE	1:I:67:GLN:NE2	2.09	0.50
1:A:38:ASN:CG	1:A:41:LYS:HB2	2.37	0.49
1:G:121:ARG:NH1	3:G:335:HOH:O	2.43	0.49
1:I:40:ALA:C	1:I:42:THR:N	2.61	0.49
1:A:50:ARG:O	1:A:54:VAL:HG23	2.13	0.49
1:A:95:PHE:CZ	1:A:99:LEU:HD23	2.48	0.49
1:E:90:ARG:H	1:E:93:THR:HG23	1.76	0.49
1:I:5[B]:VAL:CG1	1:I:77:VAL:HG22	2.23	0.49
1:C:37:VAL:CG2	1:C:38:ASN:N	2.75	0.49
1:I:32:VAL:HG23	1:I:82:MET:HE1	1.95	0.49
1:A:103:GLN:HA	1:A:103:GLN:OE1	2.14	0.48
1:E:95:PHE:CZ	1:E:99:LEU:HD23	2.48	0.48
1:E:90:ARG:HG3	1:E:93:THR:HG23	1.93	0.48
1:C:73:VAL:O	1:C:77:VAL:HG23	2.14	0.47
1:A:30:GLU:HG3	1:A:82:MET:CE	2.44	0.47
1:C:63:ASN:HD22	1:C:63:ASN:N	2.05	0.47
1:A:41:LYS:O	1:A:42:THR:C	2.55	0.47
1:K:44:MET:CG	1:K:45:PHE:CE1	2.97	0.47
1:E:146:GLU:N	1:E:147:PRO:CD	2.78	0.47
1:A:39:PRO:HA	1:A:40:ALA:HB2	1.97	0.47
1:G:63:ASN:H	1:G:63:ASN:ND2	2.01	0.47
1:E:53:MET:SD	1:E:151:ARG:HD2	2.55	0.46
1:A:38:ASN:OD1	1:A:41:LYS:HB2	2.15	0.46
1:C:38:ASN:O	1:C:40:ALA:N	2.47	0.46
1:K:37:VAL:O	1:K:37:VAL:HG22	2.16	0.46
1:K:30:GLU:O	1:K:31:VAL:HG23	2.16	0.46
1:C:20:ILE:HD12	1:C:116:VAL:HG13	1.98	0.45
1:G:20:ILE:HD11	1:G:118:THR:HG22	1.98	0.45
1:K:56:GLU:OE1	1:K:147:PRO:HB3	2.16	0.45
1:A:38:ASN:O	1:A:38:ASN:CG	2.60	0.45
1:K:118[B]:THR:HG21	3:K:303:HOH:O	2.16	0.45
1:G:90:ARG:HD3	2:G:201:SO4:O2	2.16	0.45
1:A:25:ALA:HB1	1:A:63:ASN:HD21	1.80	0.45
1:E:2:THR:O	1:E:29:ASP:HB2	2.17	0.45
1:I:20:ILE:HG13	1:I:88:GLY:HA3	1.98	0.45
1:K:41:LYS:O	1:K:42:THR:OG1	2.33	0.45
1:C:2:THR:HB	1:C:29:ASP:OD1	2.17	0.45
1:C:69:GLY:C	1:C:70:HIS:CD2	2.95	0.45
1:K:125:VAL:HG22	3:K:332:HOH:O	2.17	0.45
1:G:154:ASP:C	1:G:156:LEU:H	2.25	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:121:ARG:HD3	1:E:122:TYR:CE1	2.52	0.44
1:E:90:ARG:H	1:E:93:THR:HG21	1.80	0.44
1:E:114:PHE:HB3	1:K:114:PHE:HB3	1.99	0.44
1:C:157:ASN:OD1	1:C:157:ASN:C	2.61	0.44
1:K:101:MET:HE3	3:K:319:HOH:O	2.15	0.44
1:G:4:ALA:HB1	1:G:86:VAL:HG23	1.99	0.44
1:G:135:MET:HE2	1:G:135:MET:HA	2.00	0.44
1:I:53:MET:SD	1:I:148:VAL:HG22	2.58	0.44
1:C:71:GLY:HA2	1:K:136:LEU:HD22	2.00	0.44
1:G:3:GLY:HA3	1:G:82:MET:HE2	2.00	0.44
1:G:41:LYS:C	1:G:42:THR:HG1	2.25	0.44
1:G:74:VAL:HG22	3:G:310:HOH:O	2.19	0.43
1:G:50:ARG:O	1:G:54:VAL:HG23	2.18	0.43
1:K:90:ARG:HG3	1:K:93:THR:OG1	2.17	0.43
1:K:44:MET:O	1:K:45:PHE:CG	2.71	0.43
1:C:91:THR:HG21	3:C:341:HOH:O	2.19	0.43
1:K:3:GLY:HA3	1:K:82:MET:HE2	1.99	0.43
1:C:154:ASP:O	1:C:157:ASN:HB3	2.17	0.43
3:C:311:HOH:O	1:K:142:GLU:HB2	2.19	0.43
1:A:63:ASN:HD22	1:A:63:ASN:N	2.10	0.43
1:C:37:VAL:CG2	1:C:38:ASN:H	2.31	0.43
1:E:124:PHE:CB	1:G:103:GLN:HE21	2.15	0.43
1:G:41:LYS:O	1:G:42:THR:OG1	2.30	0.43
1:K:42:THR:HA	1:K:43:GLY:HA3	1.58	0.42
1:C:34:ALA:HB2	1:C:76:PHE:CE2	2.54	0.42
1:A:38:ASN:O	1:A:38:ASN:OD1	2.38	0.42
1:A:73:VAL:O	1:A:77:VAL:HG23	2.18	0.42
1:G:146:GLU:O	1:G:150:ARG:HB2	2.19	0.42
1:I:34:ALA:HB2	1:I:76:PHE:CE2	2.54	0.42
1:K:87:LYS:HE3	1:K:98:GLU:OE2	2.19	0.42
1:C:122:TYR:O	1:C:125:VAL:HG23	2.20	0.42
1:K:63:ASN:H	1:K:63:ASN:ND2	2.06	0.42
1:C:41:LYS:CG	1:C:43:GLY:H	2.26	0.42
1:E:41:LYS:NZ	1:E:41:LYS:HB3	2.34	0.42
1:I:42:THR:HB	1:I:43:GLY:CA	2.50	0.42
1:K:20:ILE:HG13	1:K:88:GLY:HA3	2.02	0.42
1:E:155:ARG:NH1	1:E:155:ARG:HG3	2.30	0.41
1:A:46:ASP:OD1	1:A:46:ASP:C	2.63	0.41
1:A:39:PRO:HA	1:A:40:ALA:CB	2.51	0.41
1:A:47:LEU:HD12	1:A:47:LEU:HA	1.89	0.41
1:C:91:THR:HG23	1:C:123:SER:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:87:LYS:HD3	1:E:98:GLU:OE2	2.21	0.41
1:C:156:LEU:HD23	1:C:156:LEU:HA	1.83	0.41
1:G:42:THR:HB	1:G:43:GLY:HA2	1.52	0.41
1:K:31:VAL:O	1:K:64:LEU:HA	2.20	0.41
1:I:38:ASN:N	1:I:38:ASN:HD22	2.19	0.40
1:I:152:LEU:HD12	1:I:152:LEU:HA	1.97	0.40
1:A:7:PRO:HG2	1:A:87:LYS:HE2	2.02	0.40
1:G:99:LEU:HD22	1:G:99:LEU:HA	1.95	0.40
1:K:125:VAL:CG2	3:K:332:HOH:O	2.69	0.40
1:A:145:PRO:HG2	1:A:148:VAL:HG23	2.02	0.40

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	155/156 (99%)	147 (95%)	6 (4%)	2 (1%)	9	5
1	C	154/156 (99%)	146 (95%)	5 (3%)	3 (2%)	6	3
1	E	154/156 (99%)	147 (96%)	4 (3%)	3 (2%)	6	3
1	G	154/156 (99%)	150 (97%)	2 (1%)	2 (1%)	9	5
1	I	155/156 (99%)	148 (96%)	5 (3%)	2 (1%)	9	5
1	K	155/156 (99%)	149 (96%)	4 (3%)	2 (1%)	9	5
All	All	927/936 (99%)	887 (96%)	26 (3%)	14 (2%)	8	4

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	41	LYS
1	C	42	THR

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Mol	Chain	Res	Type
1	G	41	LYS
1	G	42	THR
1	I	41	LYS
1	K	42	THR
1	E	41	LYS
1	E	42	THR
1	E	156	LEU
1	I	156	LEU
1	A	41	LYS
1	A	156	LEU
1	C	39	PRO
1	K	40	ALA

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	131/130 (101%)	123 (94%)	8 (6%)	17	14
1	C	130/130 (100%)	119 (92%)	11 (8%)	10	7
1	E	130/130 (100%)	116 (89%)	14 (11%)	6	3
1	G	130/130 (100%)	118 (91%)	12 (9%)	8	5
1	I	131/130 (101%)	119 (91%)	12 (9%)	8	5
1	K	131/130 (101%)	118 (90%)	13 (10%)	7	4
All	All	783/780 (100%)	713 (91%)	70 (9%)	9	6

All (70) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	GLN
1	A	41	LYS
1	A	48	ASP
1	A	63	ASN
1	A	99	LEU
1	A	125	VAL

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Mol	Chain	Res	Type
1	A	152	LEU
1	A	155	ARG
1	C	41	LYS
1	C	42	THR
1	C	47	LEU
1	C	63	ASN
1	C	67	GLN
1	C	70	HIS
1	C	91	THR
1	C	99	LEU
1	C	118	THR
1	C	125	VAL
1	C	156	LEU
1	E	30	GLU
1	E	41	LYS
1	E	42	THR
1	E	48	ASP
1	E	63	ASN
1	E	90	ARG
1	E	91	THR
1	E	99	LEU
1	E	103	GLN
1	E	116	VAL
1	E	121	ARG
1	E	125	VAL
1	E	152	LEU
1	E	155	ARG
1	G	30	GLU
1	G	36	LEU
1	G	37	VAL
1	G	41	LYS
1	G	42	THR
1	G	63	ASN
1	G	72	LEU
1	G	90	ARG
1	G	99	LEU
1	G	116	VAL
1	G	118	THR
1	G	156	LEU
1	I	41	LYS
1	I	42	THR
1	I	63	ASN

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Mol	Chain	Res	Type
1	I	67	GLN
1	I	91	THR
1	I	116	VAL
1	I	118	THR
1	I	125	VAL
1	I	152	LEU
1	I	155	ARG
1	I	156	LEU
1	I	157	ASN
1	K	30	GLU
1	K	37	VAL
1	K	41	LYS
1	K	42	THR
1	K	44	MET
1	K	48	ASP
1	K	63	ASN
1	K	72	LEU
1	K	90	ARG
1	K	91	THR
1	K	99	LEU
1	K	116	VAL
1	K	125	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	63	ASN
1	A	70	HIS
1	A	157	ASN
1	C	63	ASN
1	C	67	GLN
1	C	70	HIS
1	C	100	GLN
1	E	38	ASN
1	E	63	ASN
1	E	70	HIS
1	E	100	GLN
1	G	63	ASN
1	G	100	GLN
1	G	103	GLN
1	G	157	ASN
1	I	38	ASN

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Mol	Chain	Res	Type
1	I	63	ASN
1	I	67	GLN
1	K	38	ASN
1	K	63	ASN
1	K	67	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 [i](#)

8 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SO4	K	201	-	4,4,4	0.49	0	6,6,6	0.28	0
2	SO4	C	201	-	4,4,4	0.44	0	6,6,6	0.28	0
2	SO4	C	202	-	4,4,4	0.54	0	6,6,6	0.43	0
2	SO4	A	201	-	4,4,4	0.51	0	6,6,6	0.28	0
2	SO4	E	201	-	4,4,4	0.36	0	6,6,6	0.30	0
2	SO4	G	201	-	4,4,4	0.30	0	6,6,6	0.35	0
2	SO4	I	201	-	4,4,4	0.38	0	6,6,6	0.18	0
2	SO4	G	202	-	4,4,4	0.48	0	6,6,6	0.36	0

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.

3 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	202	SO4	1	0
2	G	201	SO4	1	0
2	G	202	SO4	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	156/156 (100%)	-1.37	0 100 100	11, 28, 58, 92	1 (0%)
1	C	156/156 (100%)	-1.42	0 100 100	20, 29, 47, 73	0
1	E	156/156 (100%)	-1.39	0 100 100	19, 28, 56, 74	0
1	G	156/156 (100%)	-1.38	0 100 100	19, 28, 61, 79	0
1	I	156/156 (100%)	-1.38	0 100 100	12, 29, 52, 71	1 (0%)
1	K	156/156 (100%)	-1.38	0 100 100	11, 29, 69, 83	1 (0%)
All	All	936/936 (100%)	-1.39	0 100 100	11, 29, 56, 92	3 (0%)

There are no RSRZ outliers to report.

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	SO4	C	201	5/5	0.99	0.02	28,29,33,34	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	SO4	C	202	5/5	0.99	0.05	40,43,46,48	0
2	SO4	G	202	5/5	0.99	0.07	46,46,53,54	0
2	SO4	I	201	5/5	0.99	0.03	28,28,29,31	0
2	SO4	G	201	5/5	1.00	0.02	28,28,29,29	0
2	SO4	A	201	5/5	1.00	0.02	28,29,29,29	0
2	SO4	E	201	5/5	1.00	0.04	30,31,32,32	0
2	SO4	K	201	5/5	1.00	0.02	26,26,27,30	0

6.5 Other polymers [i](#)

There are no such residues in this entry.