



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:22 PM UTC

PDB ID : 6PK5 / pdb_00006pk5
Title : Adenylate kinase from Methanococcus igneus - apo form
Authors : Moon, S.; Kim, J.; Bae, E.; Phillips Jr., G.N.
Deposited on : 2019-06-28
Resolution : 2.30 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
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

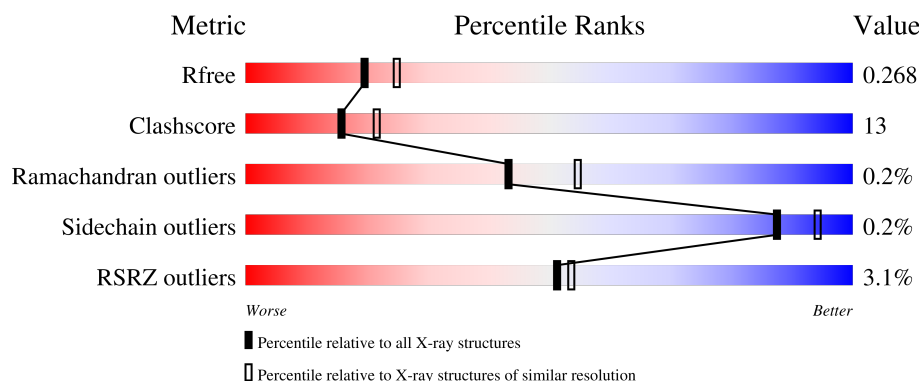
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	200	5% 72% 22% 6%
1	B	200	3% 70% 24% 6%
1	C	200	3% 74% 22% 5%
1	D	200	2% 74% 20% • •
1	E	200	2% 76% 18% 6%

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Mol	Chain	Length	Quality of chain
1	F	200	

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	CL	F	303	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 18234 atoms, of which 9236 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 Adenylate kinase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	188	Total	C	H	N	O	S	0	0	0
			2988	920	1528	249	283	8			
1	B	189	Total	C	H	N	O	S	0	0	0
			3002	924	1534	251	285	8			
1	C	190	Total	C	H	N	O	S	0	0	0
			3019	928	1540	255	288	8			
1	D	191	Total	C	H	N	O	S	0	0	0
			3038	934	1551	256	289	8			
1	E	189	Total	C	H	N	O	S	0	0	0
			3007	924	1536	254	285	8			
1	F	190	Total	C	H	N	O	S	0	0	0
			3026	930	1547	255	286	8			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	193	LEU	-	expression tag	UNP P43408
A	194	GLU	-	expression tag	UNP P43408
A	195	HIS	-	expression tag	UNP P43408
A	196	HIS	-	expression tag	UNP P43408
A	197	HIS	-	expression tag	UNP P43408
A	198	HIS	-	expression tag	UNP P43408
A	199	HIS	-	expression tag	UNP P43408
A	200	HIS	-	expression tag	UNP P43408
B	193	LEU	-	expression tag	UNP P43408
B	194	GLU	-	expression tag	UNP P43408
B	195	HIS	-	expression tag	UNP P43408
B	196	HIS	-	expression tag	UNP P43408
B	197	HIS	-	expression tag	UNP P43408
B	198	HIS	-	expression tag	UNP P43408
B	199	HIS	-	expression tag	UNP P43408
B	200	HIS	-	expression tag	UNP P43408
C	193	LEU	-	expression tag	UNP P43408

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Chain	Residue	Modelled	Actual	Comment	Reference
C	194	GLU	-	expression tag	UNP P43408
C	195	HIS	-	expression tag	UNP P43408
C	196	HIS	-	expression tag	UNP P43408
C	197	HIS	-	expression tag	UNP P43408
C	198	HIS	-	expression tag	UNP P43408
C	199	HIS	-	expression tag	UNP P43408
C	200	HIS	-	expression tag	UNP P43408
D	193	LEU	-	expression tag	UNP P43408
D	194	GLU	-	expression tag	UNP P43408
D	195	HIS	-	expression tag	UNP P43408
D	196	HIS	-	expression tag	UNP P43408
D	197	HIS	-	expression tag	UNP P43408
D	198	HIS	-	expression tag	UNP P43408
D	199	HIS	-	expression tag	UNP P43408
D	200	HIS	-	expression tag	UNP P43408
E	193	LEU	-	expression tag	UNP P43408
E	194	GLU	-	expression tag	UNP P43408
E	195	HIS	-	expression tag	UNP P43408
E	196	HIS	-	expression tag	UNP P43408
E	197	HIS	-	expression tag	UNP P43408
E	198	HIS	-	expression tag	UNP P43408
E	199	HIS	-	expression tag	UNP P43408
E	200	HIS	-	expression tag	UNP P43408
F	193	LEU	-	expression tag	UNP P43408
F	194	GLU	-	expression tag	UNP P43408
F	195	HIS	-	expression tag	UNP P43408
F	196	HIS	-	expression tag	UNP P43408
F	197	HIS	-	expression tag	UNP P43408
F	198	HIS	-	expression tag	UNP P43408
F	199	HIS	-	expression tag	UNP P43408
F	200	HIS	-	expression tag	UNP P43408

- Molecule 2 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	2	Total Cl 2 2	0	0
2	B	2	Total Cl 2 2	0	0
2	C	3	Total Cl 3 3	0	0
2	D	2	Total Cl 2 2	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	E	3	Total 3	Cl 3	0	0
2	F	3	Total 3	Cl 3	0	0

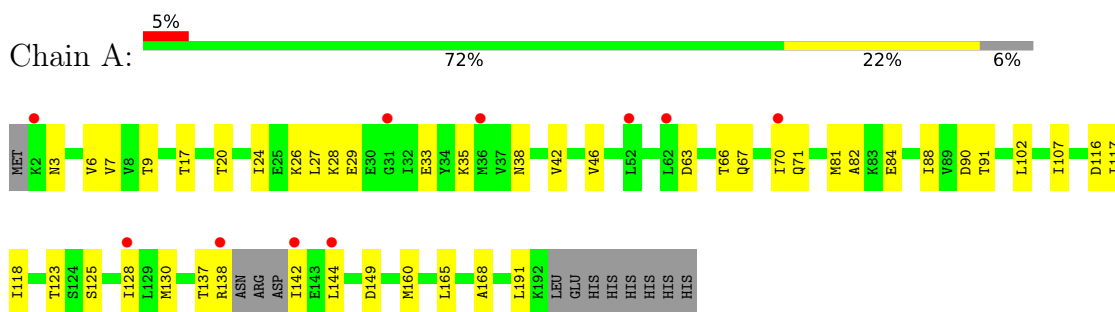
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	18	Total 18	O 18	0	0
3	B	11	Total 11	O 11	0	0
3	C	17	Total 17	O 17	0	0
3	D	33	Total 33	O 33	0	0
3	E	37	Total 37	O 37	0	0
3	F	23	Total 23	O 23	0	0

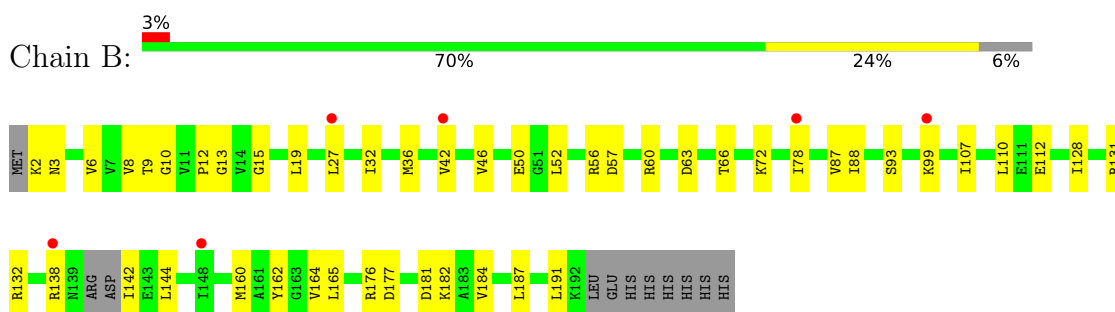
3 Residue-property plots [i](#)

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.

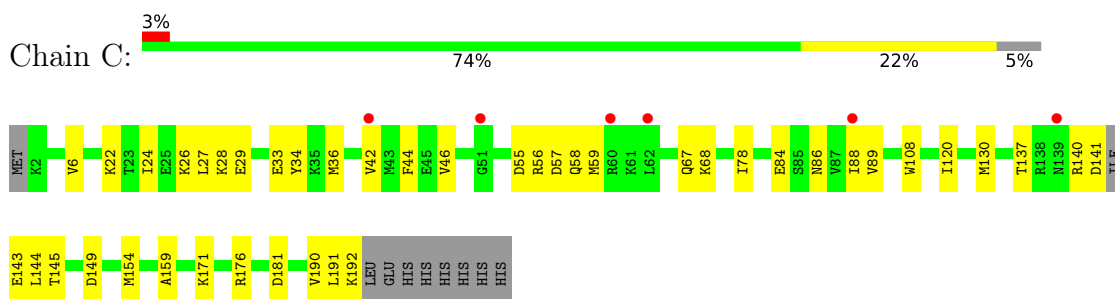
- Molecule 1: Adenylate kinase



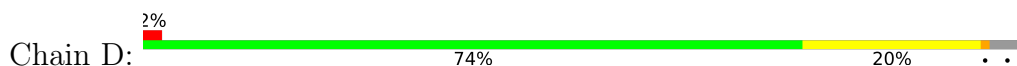
- Molecule 1: Adenylate kinase

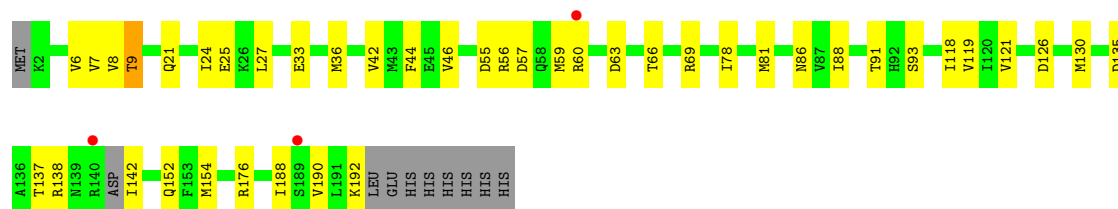


- Molecule 1: Adenylate kinase



- Molecule 1: Adenylate kinase





4 Data and refinement statistics

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, α , β , γ	68.71Å 68.71Å 257.98Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	31.33 – 2.30 31.33 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.9 (31.33-2.30) 99.9 (31.33-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.98 (at 2.29Å)	Xtriage
Refinement program	PHENIX (1.15.2_3472: ???)	Depositor
R, R_{free}	0.229 , 0.267 0.231 , 0.268	Depositor DCC
R_{free} test set	1903 reflections (3.15%)	wwPDB-VP
Wilson B-factor (Å ²)	55.5	Xtriage
Anisotropy	0.092	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 49.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.029 for -h,-k,l 0.066 for h,-h-k,-l 0.046 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	18234	wwPDB-VP
Average B, all atoms (Å ²)	63.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 13.33% 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: CL

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	0.60	0/1473	0.83	0/1982
1	B	0.59	0/1481	0.86	1/1993 (0.1%)
1	C	0.62	0/1492	0.85	0/2007
1	D	0.64	0/1501	0.86	0/2021
1	E	0.70	0/1484	0.86	0/1996
1	F	0.70	0/1492	0.92	0/2007
All	All	0.64	0/8923	0.86	1/12006 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	60	ARG	CB-CG-CD	5.42	123.76	111.30

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	1460	1528	1526	43	0
1	B	1468	1534	1532	48	0
1	C	1479	1540	1538	45	0
1	D	1487	1551	1550	48	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	1471	1536	1534	38	0
1	F	1479	1547	1545	42	0
2	A	2	0	0	0	0
2	B	2	0	0	1	0
2	C	3	0	0	0	0
2	D	2	0	0	1	0
2	E	3	0	0	0	0
2	F	3	0	0	3	0
3	A	18	0	0	7	0
3	B	11	0	0	4	0
3	C	17	0	0	14	0
3	D	33	0	0	6	0
3	E	37	0	0	5	0
3	F	23	0	0	7	0
All	All	8998	9236	9225	241	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 13.

All (241) 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:138:ARG:O	3:A:401:HOH:O	1.75	1.03
1:D:154:MET:HE3	1:E:160:MET:HE2	1.42	0.98
1:F:142:ILE:O	3:F:401:HOH:O	1.83	0.95
1:D:160:MET:CE	1:F:154:MET:HE3	2.00	0.91
1:E:132:ARG:C	1:E:138:ARG:HH21	1.79	0.90
1:C:6:VAL:HB	1:C:88:ILE:HD12	1.51	0.89
1:A:160:MET:CE	1:C:154:MET:HE3	2.06	0.85
1:D:154:MET:HE3	1:E:160:MET:CE	2.06	0.84
1:E:27:LEU:HD11	1:E:191:LEU:HD12	1.57	0.84
1:C:33:GLU:OE2	3:C:401:HOH:O	1.97	0.82
1:E:132:ARG:HB3	1:E:138:ARG:CZ	2.09	0.81
1:B:12:PRO:HB3	1:B:138:ARG:HH12	1.46	0.81
1:D:160:MET:HE3	1:F:154:MET:HE3	1.61	0.80
1:B:142:ILE:N	3:B:402:HOH:O	2.15	0.80
1:D:154:MET:HB3	1:E:160:MET:HE3	1.65	0.79
1:B:176:ARG:NH2	3:B:403:HOH:O	2.16	0.79
1:C:130:MET:HE2	1:E:144:LEU:HD21	1.66	0.78
1:B:15:GLY:HA2	2:B:301:CL:CL	2.22	0.77
1:B:6:VAL:HB	1:B:88:ILE:HD12	1.68	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:63:ASP:OD1	1:F:66:THR:HG23	1.87	0.75
1:A:160:MET:HE3	1:C:154:MET:HE3	1.68	0.73
1:D:6:VAL:HB	1:D:88:ILE:HD12	1.68	0.73
1:A:149:ASP:OD1	3:A:402:HOH:O	2.07	0.73
1:C:144:LEU:HD23	1:E:130:MET:HE3	1.70	0.73
1:A:24:ILE:HD11	1:A:88:ILE:HG21	1.69	0.72
1:B:32:ILE:O	3:B:401:HOH:O	2.09	0.71
1:D:137:THR:O	1:D:137:THR:HG22	1.88	0.71
1:A:144:LEU:HD11	1:D:130:MET:HB2	1.72	0.71
1:D:42:VAL:O	1:D:46:VAL:HG23	1.90	0.70
1:D:63:ASP:OD1	1:D:66:THR:HG23	1.91	0.70
1:D:192:LYS:O	3:D:401:HOH:O	2.09	0.70
1:B:9:THR:HG21	1:B:93:SER:HB2	1.74	0.69
1:E:192:LYS:O	3:E:401:HOH:O	2.09	0.69
1:F:8:VAL:HG23	1:F:88:ILE:HD11	1.74	0.69
1:A:137:THR:HG22	1:A:137:THR:O	1.92	0.68
1:C:58:GLN:OE1	3:C:402:HOH:O	2.12	0.68
2:F:303:CL:CL	3:F:415:HOH:O	2.47	0.68
1:C:137:THR:HG22	1:C:137:THR:O	1.95	0.67
1:A:142:ILE:HG23	3:A:403:HOH:O	1.94	0.66
1:A:160:MET:HE2	1:C:154:MET:HE3	1.77	0.66
2:D:301:CL:CL	3:D:425:HOH:O	2.52	0.65
1:A:6:VAL:HB	1:A:88:ILE:HD12	1.79	0.65
1:E:50:GLU:OE2	3:E:402:HOH:O	2.14	0.65
1:B:36:MET:HA	1:B:88:ILE:HG23	1.78	0.64
1:E:132:ARG:O	1:E:138:ARG:NH2	2.26	0.64
1:F:24:ILE:HD11	1:F:88:ILE:HG21	1.79	0.64
1:F:44:PHE:HA	1:F:59:MET:SD	2.38	0.64
1:D:160:MET:HE2	1:F:154:MET:HE3	1.81	0.63
1:A:107:ILE:HD13	1:A:165:LEU:HD13	1.81	0.62
1:B:9:THR:HG22	1:B:10:GLY:N	2.14	0.62
1:C:84:GLU:O	3:C:401:HOH:O	2.15	0.62
1:A:26:LYS:HD3	1:A:29:GLU:OE2	2.00	0.62
1:F:135:ASP:HB3	1:F:138:ARG:HG2	1.80	0.62
1:C:6:VAL:CB	1:C:88:ILE:HD12	2.28	0.62
1:C:24:ILE:HD11	1:C:88:ILE:HG21	1.81	0.62
1:E:3:ASN:HB2	1:E:82:ALA:O	2.00	0.61
1:B:50:GLU:HG3	1:B:50:GLU:O	1.98	0.61
1:F:118:ILE:N	1:F:118:ILE:HD12	2.15	0.61
1:E:137:THR:HG22	1:E:137:THR:O	1.99	0.61
1:E:143:GLU:N	3:E:405:HOH:O	2.33	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:19:LEU:HD22	1:B:184:VAL:HG22	1.82	0.61
1:D:19:LEU:CD1	1:D:187:LEU:HD22	2.30	0.61
1:F:188:ILE:O	1:F:192:LYS:HG3	2.01	0.60
1:D:20:THR:O	1:D:24:ILE:HG13	2.02	0.59
1:D:55:ASP:OD2	1:D:56:ARG:N	2.36	0.59
1:C:26:LYS:HD3	1:C:29:GLU:OE2	2.03	0.59
1:B:42:VAL:O	1:B:46:VAL:HG23	2.02	0.59
1:C:145:THR:OG1	1:E:126:ASP:OD1	2.19	0.59
1:E:26:LYS:HB3	1:E:188:ILE:HD11	1.85	0.58
1:C:55:ASP:HB3	1:C:58:GLN:HG3	1.86	0.58
1:A:125:SER:OG	1:A:149:ASP:OD1	2.22	0.58
1:D:26:LYS:HD3	1:D:29:GLU:OE2	2.03	0.58
1:F:9:THR:CG2	3:F:411:HOH:O	2.51	0.58
1:B:99:LYS:HE3	1:C:190:VAL:O	2.05	0.57
1:C:192:LYS:O	3:C:403:HOH:O	2.17	0.57
1:A:142:ILE:N	3:A:403:HOH:O	2.39	0.56
1:B:9:THR:HG23	1:B:93:SER:H	1.70	0.56
1:B:99:LYS:O	1:C:171:LYS:HB2	2.06	0.56
1:E:132:ARG:C	1:E:138:ARG:NH2	2.58	0.56
1:B:8:VAL:HG23	1:B:88:ILE:HD11	1.89	0.55
1:B:9:THR:HG21	1:B:93:SER:CB	2.35	0.55
1:B:50:GLU:HG3	1:B:52:LEU:HG	1.88	0.55
1:B:56:ARG:HD3	1:B:57:ASP:N	2.21	0.55
1:C:42:VAL:O	1:C:46:VAL:HG23	2.06	0.55
1:A:81:MET:HE1	3:A:417:HOH:O	2.05	0.54
1:C:140:ARG:O	1:C:141:ASP:OD1	2.24	0.54
1:C:22:LYS:NZ	3:C:409:HOH:O	2.40	0.54
1:C:27:LEU:HD11	1:C:191:LEU:HD12	1.90	0.54
1:F:60:ARG:HB2	3:F:402:HOH:O	2.08	0.54
1:F:137:THR:HG22	1:F:137:THR:O	2.08	0.53
1:A:3:ASN:HB2	1:A:82:ALA:O	2.09	0.53
1:B:9:THR:CG2	1:B:93:SER:H	2.21	0.53
1:B:144:LEU:HD23	1:F:130:MET:HE3	1.89	0.53
1:E:24:ILE:HD11	1:E:88:ILE:HD12	1.90	0.53
1:E:66:THR:HG22	1:E:69:ARG:HH12	1.73	0.53
1:A:42:VAL:O	1:A:46:VAL:HG23	2.09	0.53
1:C:144:LEU:CD2	1:E:130:MET:HE3	2.39	0.53
1:B:99:LYS:NZ	3:C:403:HOH:O	2.41	0.53
1:E:132:ARG:HB3	1:E:138:ARG:NE	2.24	0.53
1:D:181:ASP:N	1:D:181:ASP:OD1	2.41	0.52
1:D:24:ILE:HD11	1:D:88:ILE:HG21	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:107:ILE:HA	1:D:110:LEU:HD12	1.92	0.52
1:F:66:THR:HG22	1:F:69:ARG:HH22	1.75	0.52
1:B:182:LYS:HD2	1:B:182:LYS:N	2.24	0.52
1:B:8:VAL:CG2	1:B:88:ILE:HD11	2.40	0.52
1:F:9:THR:HB	1:F:119:VAL:O	2.11	0.51
1:B:181:ASP:N	1:B:181:ASP:OD1	2.44	0.51
1:E:132:ARG:HB3	1:E:138:ARG:NH2	2.24	0.51
1:B:36:MET:HA	1:B:88:ILE:CG2	2.41	0.51
1:C:143:GLU:N	3:C:406:HOH:O	2.42	0.51
1:A:24:ILE:HG22	1:A:28:LYS:HE3	1.92	0.51
1:B:13:GLY:N	1:B:138:ARG:NH2	2.59	0.51
1:F:7:VAL:HG13	1:F:91:THR:HG22	1.93	0.51
1:D:56:ARG:H	1:D:56:ARG:HD3	1.76	0.51
1:A:142:ILE:HD12	1:A:142:ILE:C	2.36	0.51
1:A:38:ASN:O	1:A:42:VAL:HG23	2.12	0.50
1:E:52:LEU:HD12	1:E:70:ILE:HD11	1.94	0.50
1:B:131:ARG:NH2	1:B:177:ASP:OD1	2.40	0.50
1:B:72:LYS:HE2	3:B:409:HOH:O	2.11	0.50
1:D:180:LEU:HB3	3:D:420:HOH:O	2.10	0.50
1:C:24:ILE:HG23	1:C:28:LYS:HE3	1.93	0.49
1:F:56:ARG:HD3	1:F:57:ASP:N	2.27	0.49
1:B:63:ASP:OD1	1:B:66:THR:HG23	2.13	0.49
1:D:60:ARG:HB2	3:D:409:HOH:O	2.11	0.49
1:D:96:LYS:HE2	1:D:101:TYR:CE1	2.48	0.49
1:D:137:THR:O	1:D:137:THR:CG2	2.59	0.49
1:B:2:LYS:O	1:B:3:ASN:C	2.54	0.49
1:A:123:THR:HG23	1:A:128:ILE:HG13	1.95	0.48
1:C:58:GLN:NE2	3:C:402:HOH:O	2.46	0.48
1:F:21:GLN:O	1:F:25:GLU:HG3	2.13	0.48
1:C:24:ILE:CD1	1:C:36:MET:HB2	2.43	0.48
1:B:13:GLY:N	1:B:138:ARG:HH22	2.11	0.48
1:D:19:LEU:HD12	1:D:187:LEU:HD22	1.96	0.48
1:D:59:MET:C	1:D:61:LYS:H	2.20	0.48
1:A:46:VAL:HG12	1:A:70:ILE:HD12	1.94	0.48
1:D:19:LEU:HD23	3:D:420:HOH:O	2.13	0.48
1:A:28:LYS:HE2	1:A:33:GLU:OE1	2.14	0.48
1:F:60:ARG:O	3:F:402:HOH:O	2.20	0.48
1:C:176:ARG:HH22	1:F:176:ARG:HE	1.62	0.48
1:B:27:LEU:HD11	1:B:191:LEU:HD12	1.96	0.47
1:D:26:LYS:CB	1:D:188:ILE:HD11	2.45	0.47
1:F:135:ASP:HB3	1:F:138:ARG:CG	2.44	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:99:LYS:HE3	1:F:190:VAL:O	2.14	0.47
1:B:9:THR:CG2	1:B:10:GLY:N	2.77	0.47
1:D:63:ASP:OD1	1:D:66:THR:CG2	2.61	0.47
1:A:130:MET:HE1	1:D:142:ILE:O	2.14	0.47
1:F:63:ASP:OD1	1:F:66:THR:CG2	2.61	0.47
1:D:135:ASP:C	1:D:137:THR:H	2.22	0.47
1:A:7:VAL:HG13	1:A:91:THR:CG2	2.45	0.47
1:B:99:LYS:HE3	3:C:403:HOH:O	2.15	0.47
1:B:110:LEU:HD11	1:B:162:TYR:CD1	2.50	0.47
1:C:141:ASP:O	1:C:143:GLU:N	2.48	0.47
1:C:120:ILE:HD11	1:C:159:ALA:HB3	1.97	0.47
1:C:137:THR:O	1:C:137:THR:CG2	2.63	0.47
1:E:34:TYR:CD1	1:E:86:ASN:HB3	2.49	0.47
1:D:36:MET:HA	1:D:88:ILE:HG22	1.97	0.46
1:C:44:PHE:HA	1:C:59:MET:SD	2.54	0.46
1:F:9:THR:HG21	1:F:93:SER:HB2	1.97	0.46
1:A:17:THR:OG1	1:A:90:ASP:OD2	2.20	0.46
1:E:34:TYR:CD1	1:E:88:ILE:HG13	2.50	0.46
1:D:188:ILE:O	1:D:192:LYS:HG3	2.16	0.46
1:A:20:THR:O	1:A:24:ILE:HG13	2.16	0.46
1:A:118:ILE:N	1:A:118:ILE:HD12	2.30	0.46
1:B:12:PRO:HB3	1:B:138:ARG:NH1	2.23	0.46
1:B:19:LEU:CD1	1:B:187:LEU:HD22	2.46	0.46
1:D:26:LYS:HB2	1:D:188:ILE:HD11	1.97	0.46
1:D:42:VAL:HG13	1:D:77:LYS:HD3	1.97	0.46
1:A:9:THR:HA	1:A:91:THR:O	2.16	0.46
1:B:99:LYS:CE	3:C:403:HOH:O	2.63	0.45
1:E:68:LYS:NZ	3:E:403:HOH:O	2.18	0.45
1:D:131:ARG:NH2	1:D:177:ASP:OD1	2.31	0.45
1:A:116:ASP:C	1:A:117:ILE:HG13	2.41	0.45
1:F:152:GLN:NE2	2:F:303:CL:CL	2.78	0.45
1:D:60:ARG:C	3:D:409:HOH:O	2.60	0.45
1:E:181:ASP:OD1	1:E:181:ASP:N	2.46	0.45
1:F:142:ILE:O	1:F:142:ILE:HD12	2.16	0.45
1:C:58:GLN:CD	3:C:402:HOH:O	2.56	0.45
1:A:46:VAL:HG12	1:A:70:ILE:CD1	2.47	0.45
1:D:59:MET:HE3	1:D:62:LEU:HD11	1.98	0.45
1:A:7:VAL:HG13	1:A:91:THR:HG22	1.98	0.45
1:C:34:TYR:CE1	1:C:86:ASN:HB3	2.52	0.45
1:E:19:LEU:HD22	1:E:184:VAL:HG22	1.99	0.45
1:C:68:LYS:HE2	1:C:108:TRP:CG	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:68:LYS:HA	1:D:71:GLN:OE1	2.17	0.45
1:F:24:ILE:HD11	1:F:88:ILE:CG2	2.47	0.45
1:A:35:LYS:NZ	3:A:406:HOH:O	2.50	0.44
1:A:160:MET:HE3	1:C:154:MET:HB3	1.98	0.44
1:C:55:ASP:OD1	1:C:56:ARG:N	2.51	0.44
1:B:50:GLU:HG2	1:B:52:LEU:CD1	2.47	0.44
1:D:36:MET:HA	1:D:88:ILE:CG2	2.48	0.44
1:F:6:VAL:HB	1:F:88:ILE:HD12	1.99	0.44
1:F:60:ARG:C	3:F:402:HOH:O	2.59	0.44
1:B:78:ILE:HG23	1:B:87:VAL:HG11	1.99	0.44
1:D:138:ARG:HB3	1:D:141:ASP:HB2	1.99	0.44
1:A:142:ILE:HD13	1:D:130:MET:HE1	1.98	0.44
1:C:181:ASP:OD1	1:C:181:ASP:N	2.50	0.44
1:C:192:LYS:C	3:C:403:HOH:O	2.61	0.43
1:D:101:TYR:HB2	1:E:160:MET:HE1	2.00	0.43
1:F:42:VAL:O	1:F:46:VAL:HG23	2.18	0.43
1:F:121:VAL:O	2:F:303:CL:CL	2.73	0.43
1:F:27:LEU:HD21	1:F:188:ILE:HG23	2.01	0.43
1:F:9:THR:HG22	3:F:411:HOH:O	2.18	0.43
1:F:33:GLU:O	1:F:86:ASN:HB2	2.18	0.43
1:F:7:VAL:HG13	1:F:91:THR:CG2	2.47	0.43
1:B:160:MET:O	1:B:164:VAL:HG23	2.19	0.43
1:E:19:LEU:CD1	1:E:187:LEU:HD22	2.49	0.43
1:E:24:ILE:HG23	1:E:28:LYS:HE3	2.00	0.43
1:A:102:LEU:HD12	3:A:413:HOH:O	2.19	0.42
1:A:137:THR:O	1:A:137:THR:CG2	2.62	0.42
1:C:55:ASP:OD1	1:C:56:ARG:HD3	2.19	0.42
1:D:59:MET:O	1:D:62:LEU:HG	2.19	0.42
1:B:50:GLU:CG	1:B:52:LEU:HG	2.49	0.42
1:C:56:ARG:HD3	1:C:57:ASP:N	2.35	0.42
1:A:84:GLU:O	1:A:84:GLU:HG3	2.18	0.42
1:B:107:ILE:HD13	1:B:165:LEU:HD13	2.00	0.42
1:E:17:THR:O	1:E:21:GLN:HG3	2.19	0.42
1:B:63:ASP:OD1	1:B:63:ASP:C	2.62	0.42
1:E:137:THR:OG1	3:E:404:HOH:O	2.21	0.42
1:A:63:ASP:OD1	1:A:66:THR:HG23	2.19	0.42
1:C:28:LYS:NZ	1:C:34:TYR:O	2.45	0.42
1:E:128:ILE:O	1:E:132:ARG:HG3	2.19	0.42
1:D:66:THR:O	1:D:70:ILE:HG12	2.19	0.42
1:B:72:LYS:CG	1:B:112:GLU:HG3	2.50	0.41
1:E:66:THR:O	1:E:70:ILE:HG12	2.19	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:78:ILE:HD13	1:C:89:VAL:CG2	2.50	0.41
1:B:128:ILE:O	1:B:132:ARG:HG3	2.20	0.41
1:D:23:THR:HG23	1:D:188:ILE:HG13	2.03	0.41
1:F:55:ASP:OD2	1:F:56:ARG:N	2.53	0.41
1:C:67:GLN:NE2	3:C:410:HOH:O	2.53	0.41
1:C:149:ASP:HA	3:C:405:HOH:O	2.21	0.41
1:A:67:GLN:O	1:A:71:GLN:HG3	2.21	0.41
1:A:118:ILE:CD1	1:A:168:ALA:HB3	2.50	0.41
1:D:154:MET:CE	1:E:160:MET:HE2	2.31	0.41
1:F:24:ILE:CD1	1:F:36:MET:HB2	2.49	0.41
1:B:144:LEU:HB3	1:F:126:ASP:OD2	2.21	0.41
1:F:8:VAL:CG2	1:F:88:ILE:HD11	2.47	0.41
1:A:27:LEU:HD11	1:A:191:LEU:HD12	2.03	0.40
1:D:55:ASP:OD2	1:D:56:ARG:HD3	2.21	0.40
1:A:118:ILE:CD1	1:A:168:ALA:CB	2.99	0.40
1:E:76:ARG:NH2	1:E:112:GLU:OE2	2.52	0.40
1:F:78:ILE:HA	1:F:81:MET:HG2	2.03	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	184/200 (92%)	180 (98%)	4 (2%)	0	100	100
1	B	185/200 (92%)	181 (98%)	4 (2%)	0	100	100
1	C	186/200 (93%)	181 (97%)	5 (3%)	0	100	100
1	D	189/200 (94%)	178 (94%)	9 (5%)	2 (1%)	11	13
1	E	185/200 (92%)	180 (97%)	5 (3%)	0	100	100
1	F	186/200 (93%)	181 (97%)	5 (3%)	0	100	100
All	All	1115/1200 (93%)	1081 (97%)	32 (3%)	2 (0%)	43	55

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	191	LEU
1	D	141	ASP

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	163/175 (93%)	163 (100%)	0	100	100
1	B	164/175 (94%)	164 (100%)	0	100	100
1	C	165/175 (94%)	165 (100%)	0	100	100
1	D	166/175 (95%)	165 (99%)	1 (1%)	78	89
1	E	164/175 (94%)	164 (100%)	0	100	100
1	F	165/175 (94%)	164 (99%)	1 (1%)	78	89
All	All	987/1050 (94%)	985 (100%)	2 (0%)	87	94

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	88	ILE
1	F	9	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	21	GLN
1	A	152	GLN
1	B	152	GLN
1	E	152	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](#)

Of 15 ligands modelled in this entry, 15 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	188/200 (94%)	0.66	10 (5%) 32 34	57, 65, 79, 91	0
1	B	189/200 (94%)	0.72	6 (3%) 50 52	56, 66, 82, 93	0
1	C	190/200 (95%)	0.63	6 (3%) 50 52	55, 63, 76, 89	0
1	D	191/200 (95%)	0.60	5 (2%) 57 59	54, 64, 75, 82	0
1	E	189/200 (94%)	0.41	5 (2%) 57 59	53, 61, 72, 78	0
1	F	190/200 (95%)	0.45	3 (1%) 70 72	52, 59, 69, 77	0
All	All	1137/1200 (94%)	0.58	35 (3%) 51 53	52, 63, 76, 93	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	42	VAL	3.4
1	B	138	ARG	3.4
1	F	140	ARG	3.4
1	D	142	ILE	3.3
1	C	60	ARG	3.2
1	C	62	LEU	3.0
1	C	51	GLY	2.9
1	B	42	VAL	2.9
1	F	60	ARG	2.9
1	A	142	ILE	2.8
1	E	138	ARG	2.8
1	F	189	SER	2.8
1	A	52	LEU	2.7
1	B	99	LYS	2.7
1	A	31	GLY	2.6
1	B	78	ILE	2.5
1	A	138	ARG	2.5
1	C	88	ILE	2.4
1	E	189	SER	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	87	VAL	2.3
1	D	129	LEU	2.3
1	D	140	ARG	2.3
1	A	2	LYS	2.3
1	B	27	LEU	2.3
1	E	60	ARG	2.2
1	A	70	ILE	2.2
1	A	128	ILE	2.2
1	D	60	ARG	2.1
1	E	140	ARG	2.1
1	C	139	ASN	2.1
1	A	144	LEU	2.1
1	A	36	MET	2.0
1	A	62	LEU	2.0
1	B	148	ILE	2.0
1	E	64	PRO	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	CL	D	301	1/1	0.89	0.09	59,59,59,59	0
2	CL	C	302	1/1	0.90	0.10	60,60,60,60	0
2	CL	B	301	1/1	0.90	0.08	62,62,62,62	0
2	CL	E	302	1/1	0.91	0.10	57,57,57,57	0
2	CL	C	301	1/1	0.93	0.32	57,57,57,57	0
2	CL	B	302	1/1	0.93	0.09	62,62,62,62	0
2	CL	F	303	1/1	0.93	0.24	57,57,57,57	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	CL	E	303	1/1	0.94	0.20	57,57,57,57	0
2	CL	F	301	1/1	0.95	0.08	57,57,57,57	0
2	CL	E	301	1/1	0.96	0.13	59,59,59,59	0
2	CL	D	302	1/1	0.97	0.11	57,57,57,57	0
2	CL	A	302	1/1	0.97	0.10	58,58,58,58	0
2	CL	C	303	1/1	0.98	0.11	60,60,60,60	0
2	CL	F	302	1/1	0.98	0.07	57,57,57,57	0
2	CL	A	301	1/1	0.98	0.16	60,60,60,60	0

6.5 Other polymers [i](#)

There are no such residues in this entry.