



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 24, 2024 – 11:01 PM EDT

PDB ID : 6F5A
Title : Crystal structure of H. pylori purine nucleoside phosphorylase
Authors : Stefanic, Z.
Deposited on : 2017-12-01
Resolution : 2.20 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	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

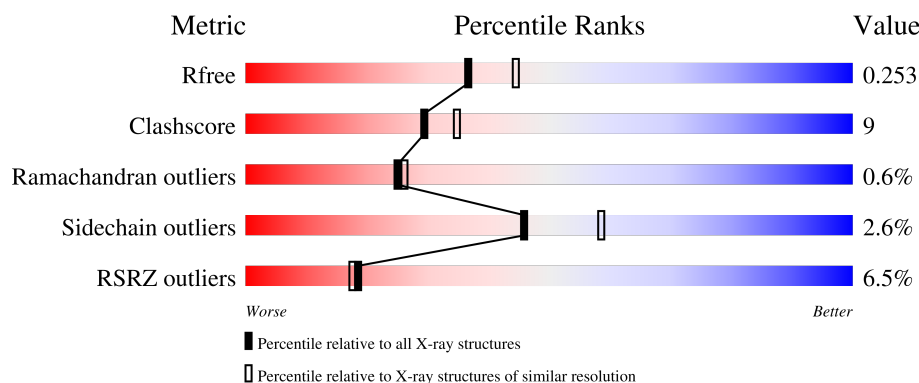
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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	233	<div> <div>10%</div> <div>75%</div> <div>23%</div> <div>.</div> </div>
1	B	233	<div> <div>12%</div> <div>83%</div> <div>17%</div> </div>
1	C	233	<div> <div>90%</div> <div>10%</div> </div>
1	D	233	<div> <div>16%</div> <div>62%</div> <div>33%</div> <div>..</div> </div>
1	E	233	<div> <div>91%</div> <div>9%</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	233	<div><div>%</div><div><div></div><div>86%</div><div>13%</div><div></div></div><div></div></div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 11356 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 Purine nucleoside phosphorylase DeoD-type.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	233	Total	C	N	O	S	0	0	0
			1802	1149	302	334	17			
1	B	233	Total	C	N	O	S	0	0	0
			1802	1149	302	334	17			
1	C	233	Total	C	N	O	S	0	0	0
			1802	1149	302	334	17			
1	D	233	Total	C	N	O	S	0	0	0
			1802	1149	302	334	17			
1	E	233	Total	C	N	O	S	0	0	0
			1802	1149	302	334	17			
1	F	233	Total	C	N	O	S	0	0	0
			1802	1149	302	334	17			

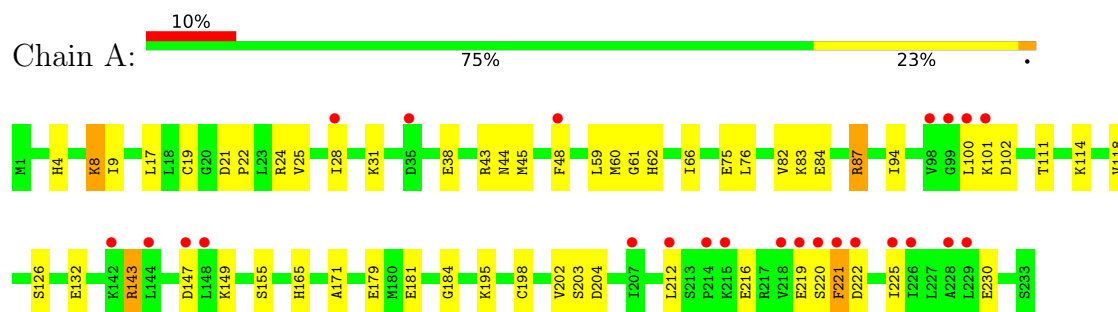
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	65	Total	O	0	0
			65	65		
2	B	53	Total	O	0	0
			53	53		
2	C	144	Total	O	0	0
			144	144		
2	D	58	Total	O	0	0
			58	58		
2	E	133	Total	O	0	0
			133	133		
2	F	91	Total	O	0	0
			91	91		

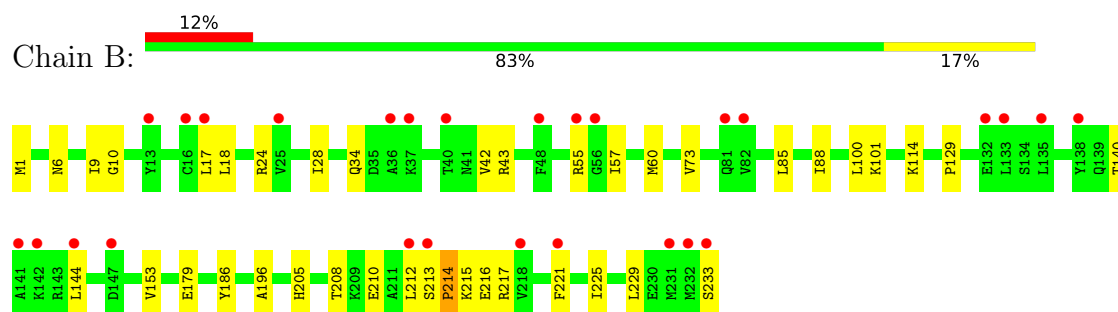
3 Residue-property plots

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.

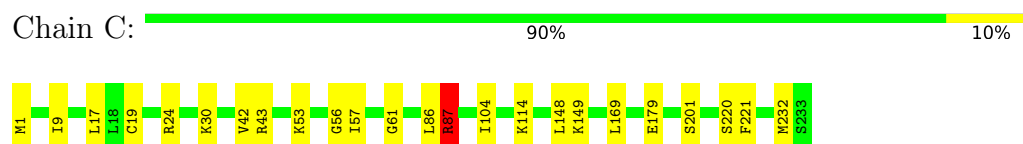
• Molecule 1: Purine nucleoside phosphorylase DeoD-type



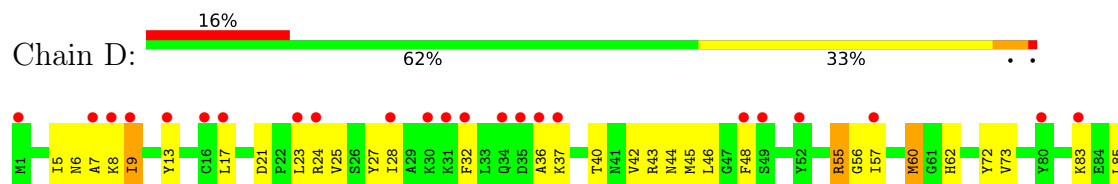
• Molecule 1: Purine nucleoside phosphorylase DeoD-type

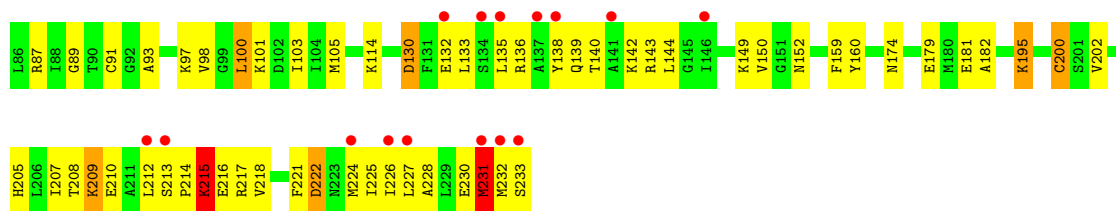


• Molecule 1: Purine nucleoside phosphorylase DeoD-type



• Molecule 1: Purine nucleoside phosphorylase DeoD-type





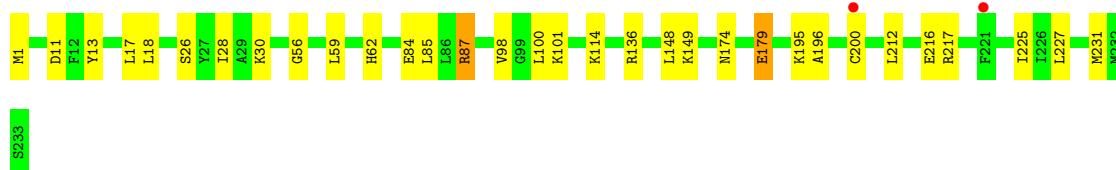
- Molecule 1: Purine nucleoside phosphorylase DeoD-type

Chain E: 91% 9%



- Molecule 1: Purine nucleoside phosphorylase DeoD-type

Chain F: 86% 13%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	83.75Å 65.09Å 140.19Å 90.00° 99.62° 90.00°	Depositor
Resolution (Å)	43.30 – 2.20 47.38 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.2 (43.30-2.20) 99.3 (47.38-2.20)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.57 (at 2.20Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.202 , 0.252 0.202 , 0.253	Depositor DCC
R_{free} test set	3658 reflections (4.86%)	wwPDB-VP
Wilson B-factor (Å ²)	32.3	Xtriage
Anisotropy	0.578	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 43.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	11356	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.82% 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](#)

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	0.38	0/1833	0.58	0/2467
1	B	0.41	0/1833	0.56	0/2467
1	C	0.44	0/1833	0.61	1/2467 (0.0%)
1	D	0.45	0/1833	0.66	1/2467 (0.0%)
1	E	0.45	0/1833	0.59	0/2467
1	F	0.43	1/1833 (0.1%)	0.62	1/2467 (0.0%)
All	All	0.43	1/10998 (0.0%)	0.60	3/14802 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	200	CYS	CB-SG	-5.57	1.72	1.81

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	200	CYS	CA-CB-SG	8.03	128.46	114.00
1	D	231	MET	CB-CG-SD	-5.37	96.29	112.40
1	C	87	ARG	NE-CZ-NH1	-5.07	117.77	120.30

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	1802	0	1838	52	0
1	B	1802	0	1838	30	0
1	C	1802	0	1838	16	0
1	D	1802	0	1837	88	0
1	E	1802	0	1838	15	0
1	F	1802	0	1836	18	0
2	A	65	0	0	6	0
2	B	53	0	0	1	0
2	C	144	0	0	5	0
2	D	58	0	0	5	0
2	E	133	0	0	5	0
2	F	91	0	0	3	0
All	All	11356	0	11025	207	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 (207) 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:D:55:ARG:NH2	1:D:232:MET:CE	1.71	1.49
1:D:55:ARG:CZ	1:D:232:MET:CE	2.12	1.27
1:D:55:ARG:NH2	1:D:232:MET:SD	2.10	1.23
1:D:55:ARG:CZ	1:D:232:MET:HE1	1.69	1.20
1:D:55:ARG:NE	1:D:232:MET:HE3	1.58	1.17
1:D:55:ARG:NH2	1:D:232:MET:HE1	1.54	1.09
1:D:55:ARG:HE	1:D:232:MET:HE3	1.10	1.06
1:D:55:ARG:NE	1:D:232:MET:CE	2.19	1.00
1:D:55:ARG:NH2	1:D:232:MET:HE3	1.73	0.99
1:D:55:ARG:CZ	1:D:232:MET:HE3	1.84	0.96
1:A:45:MET:HE3	1:A:61:GLY:HA3	1.56	0.87
1:B:212:LEU:HD13	1:B:216:GLU:HB2	1.54	0.87
1:D:215:LYS:CD	1:D:216:GLU:H	1.87	0.86
1:D:215:LYS:NZ	1:D:216:GLU:OE1	2.09	0.86
1:A:143:ARG:NH2	1:A:230:GLU:OE1	2.09	0.85
1:D:222:ASP:HA	1:D:225:ILE:HD12	1.59	0.84
1:A:147:ASP:O	2:A:301:HOH:O	1.94	0.84
1:F:1:MET:N	2:F:302:HOH:O	2.10	0.83
1:A:44:ASN:HB3	1:D:44:ASN:OD1	1.78	0.83
1:A:165:HIS:ND1	2:A:302:HOH:O	2.12	0.82
1:D:215:LYS:HD2	1:D:216:GLU:H	1.44	0.82
1:C:201:SER:OG	2:C:401:HOH:O	1.97	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:6:ASN:OD1	2:D:301:HOH:O	1.97	0.82
1:A:44:ASN:OD1	1:D:23:LEU:CD1	2.30	0.80
1:D:55:ARG:HE	1:D:232:MET:CE	1.88	0.80
1:B:208:THR:HG23	1:B:210:GLU:H	1.48	0.79
1:A:83:LYS:HD3	1:A:84:GLU:HG3	1.65	0.78
1:A:44:ASN:OD1	1:D:23:LEU:HD12	1.84	0.78
1:D:215:LYS:HD3	1:D:216:GLU:HG3	1.64	0.77
1:A:212:LEU:HD12	1:A:216:GLU:HB3	1.65	0.77
1:D:213:SER:OG	1:D:215:LYS:HE3	1.85	0.76
1:C:1:MET:N	2:C:402:HOH:O	2.20	0.75
1:D:97:LYS:O	1:D:149:LYS:NZ	2.21	0.74
1:E:43:ARG:HD3	2:E:305:HOH:O	1.87	0.74
1:F:11:ASP:OD2	2:F:301:HOH:O	2.06	0.72
1:D:100:LEU:H	1:D:100:LEU:HD23	1.55	0.71
1:D:87:ARG:NH2	1:D:181:GLU:OE2	2.25	0.69
1:D:27:TYR:OH	1:D:222:ASP:OD1	2.10	0.69
1:F:212:LEU:O	1:F:217:ARG:NH2	2.27	0.67
1:E:210:GLU:OE1	2:E:301:HOH:O	2.12	0.66
1:A:100:LEU:HD13	1:A:212:LEU:HD22	1.78	0.65
1:D:100:LEU:HD12	1:D:212:LEU:HD21	1.77	0.65
1:B:205:HIS:HB3	1:B:208:THR:HG22	1.79	0.65
1:F:84:GLU:OE2	1:F:195:LYS:NZ	2.30	0.65
1:A:181:GLU:OE1	2:A:303:HOH:O	2.15	0.64
1:A:126:SER:HB2	2:F:311:HOH:O	1.99	0.63
1:B:24:ARG:HD3	1:B:221:PHE:CZ	2.33	0.63
1:A:87:ARG:NH2	1:A:181:GLU:OE1	2.32	0.63
1:B:100:LEU:O	1:B:101:LYS:HB3	1.97	0.63
1:C:114:LYS:HB2	1:F:114:LYS:HB2	1.80	0.62
1:D:132:GLU:HA	1:D:135:LEU:HD22	1.83	0.61
1:D:132:GLU:HA	1:D:135:LEU:CD2	2.31	0.60
1:E:28:ILE:HG12	1:E:225:ILE:HD13	1.83	0.60
1:D:40:THR:HG21	1:D:60:MET:CE	2.33	0.59
1:D:215:LYS:HD2	1:D:216:GLU:N	2.17	0.59
1:D:17:LEU:HD11	1:D:228:ALA:HB1	1.84	0.58
1:B:24:ARG:HD3	1:B:221:PHE:CE2	2.38	0.58
1:D:8:LYS:HD2	1:D:9:ILE:N	2.18	0.58
1:E:62:HIS:NE2	1:E:87:ARG:HD2	2.18	0.57
1:A:45:MET:CE	1:A:61:GLY:HA3	2.31	0.57
1:A:28:ILE:HG12	1:A:225:ILE:HD12	1.86	0.57
1:D:215:LYS:HD3	1:D:216:GLU:H	1.69	0.57
1:A:45:MET:HE2	1:A:62:HIS:N	2.20	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:42:VAL:HG12	1:C:43:ARG:HG3	1.86	0.56
1:D:208:THR:O	1:D:210:GLU:N	2.39	0.56
1:D:213:SER:HB2	1:D:214:PRO:HD2	1.88	0.55
1:F:85:LEU:O	1:F:196:ALA:HA	2.07	0.55
1:D:159:PHE:N	2:D:303:HOH:O	2.40	0.55
1:A:87:ARG:HD2	1:A:198:CYS:SG	2.45	0.55
1:A:25:VAL:HG13	1:A:59:LEU:HB3	1.88	0.55
1:D:230:GLU:O	1:D:233:SER:N	2.40	0.54
1:E:69:CYS:O	1:E:73:VAL:HG13	2.08	0.54
1:B:213:SER:O	1:B:215:LYS:N	2.39	0.54
1:F:28:ILE:HG12	1:F:225:ILE:HD13	1.89	0.54
1:D:217:ARG:NH1	2:D:305:HOH:O	2.25	0.53
1:A:100:LEU:O	1:A:101:LYS:HB3	2.09	0.53
1:B:212:LEU:HD11	1:B:217:ARG:HG2	1.92	0.53
1:F:136:ARG:NH1	1:F:231:MET:O	2.34	0.52
1:A:31:LYS:NZ	1:A:222:ASP:OD2	2.34	0.52
1:D:208:THR:C	1:D:210:GLU:H	2.12	0.52
1:D:215:LYS:HZ3	1:D:216:GLU:CD	2.08	0.52
1:F:87:ARG:NH1	1:F:179:GLU:OE1	2.43	0.52
1:A:114:LYS:HB2	1:D:114:LYS:HB2	1.92	0.52
1:B:18:LEU:HD23	1:B:60:MET:HB3	1.92	0.52
1:C:24:ARG:HD2	1:C:221:PHE:CE1	2.46	0.51
1:D:205:HIS:CE1	1:D:207:ILE:HB	2.44	0.51
1:D:135:LEU:HD23	1:D:136:ARG:H	1.76	0.51
1:B:114:LYS:HB2	1:E:114:LYS:HB2	1.93	0.51
1:D:215:LYS:HZ2	1:D:216:GLU:HB2	1.76	0.51
1:A:62:HIS:CE1	1:A:181:GLU:HG2	2.46	0.50
1:A:4:HIS:NE2	2:A:310:HOH:O	2.34	0.50
1:A:82:VAL:O	2:A:305:HOH:O	2.18	0.50
1:E:122:ASN:ND2	2:E:306:HOH:O	2.45	0.50
1:B:73:VAL:HG13	1:B:85:LEU:HD22	1.94	0.50
1:D:93:ALA:HB2	1:D:202:VAL:HG13	1.93	0.50
1:B:100:LEU:HD11	1:B:210:GLU:HB3	1.94	0.49
1:D:89:GLY:O	1:D:200:CYS:HA	2.12	0.49
1:D:130:ASP:OD2	1:D:133:LEU:HD13	2.12	0.49
1:D:24:ARG:HA	1:D:27:TYR:HB3	1.95	0.49
1:C:221:PHE:HB3	2:C:420:HOH:O	2.13	0.49
1:A:203:SER:HA	1:A:212:LEU:HD23	1.94	0.49
1:C:19:CYS:O	1:C:61:GLY:HA2	2.13	0.49
2:A:334:HOH:O	1:F:174:ASN:HB2	2.13	0.48
1:A:60:MET:HG2	1:A:76:LEU:HD21	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:87:ARG:NH2	1:A:181:GLU:HB2	2.29	0.48
1:B:6:ASN:HB2	1:B:42:VAL:HG23	1.95	0.48
1:C:56:GLY:O	1:C:57:ILE:HD12	2.14	0.48
1:D:21:ASP:OD1	1:D:23:LEU:HB2	2.14	0.48
1:B:212:LEU:CD1	1:B:217:ARG:HG2	2.44	0.47
1:D:205:HIS:HE1	1:D:207:ILE:HB	1.80	0.47
1:A:21:ASP:OD1	1:D:43:ARG:HA	2.15	0.47
1:A:24:ARG:CG	1:A:221:PHE:HE1	2.27	0.47
1:D:105:MET:HE3	1:D:150:VAL:HG13	1.96	0.47
1:B:140:THR:O	1:B:144:LEU:HG	2.15	0.47
1:B:85:LEU:O	1:B:196:ALA:HA	2.15	0.46
1:D:62:HIS:HA	1:D:72:TYR:CD2	2.50	0.46
1:A:75:GLU:OE1	1:D:160:TYR:OH	2.30	0.46
1:A:100:LEU:HD23	1:A:202:VAL:HG12	1.96	0.46
1:D:13:TYR:CZ	1:D:56:GLY:HA3	2.50	0.46
1:B:1:MET:N	2:B:306:HOH:O	2.47	0.46
1:B:214:PRO:HA	1:B:217:ARG:HG3	1.97	0.46
1:B:28:ILE:HG12	1:B:225:ILE:HD13	1.97	0.46
1:F:13:TYR:CE1	1:F:56:GLY:HA3	2.50	0.46
1:D:28:ILE:HG12	1:D:225:ILE:HG21	1.97	0.46
1:D:25:VAL:HG11	1:D:46:LEU:O	2.16	0.46
1:A:24:ARG:HG2	1:A:221:PHE:HE1	1.81	0.46
1:D:212:LEU:HD23	1:D:212:LEU:HA	1.64	0.46
1:B:55:ARG:HD3	1:B:233:SER:HA	1.98	0.46
1:A:101:LYS:HD2	1:A:220:SER:OG	2.16	0.45
1:C:17:LEU:HD23	1:C:86:LEU:O	2.16	0.45
1:A:22:PRO:HD3	1:A:45:MET:HE1	1.98	0.45
1:E:91:CYS:SG	1:E:200:CYS:HB3	2.57	0.45
1:E:221:PHE:O	1:E:225:ILE:HG13	2.16	0.45
1:A:83:LYS:HD3	1:A:84:GLU:CG	2.43	0.45
1:C:53:LYS:NZ	2:C:414:HOH:O	2.48	0.45
1:A:83:LYS:HE2	1:A:195:LYS:HD3	1.97	0.45
1:B:9:ILE:HD12	1:B:10:GLY:H	1.81	0.45
1:C:169:LEU:HD23	1:E:121:LEU:HD21	1.99	0.45
1:C:17:LEU:HG	1:C:232:MET:HE1	1.98	0.45
1:C:104:ILE:HA	1:C:149:LYS:O	2.17	0.45
1:A:17:LEU:HB2	1:A:59:LEU:HD23	1.99	0.45
1:B:114:LYS:HE2	1:E:112:ASP:O	2.17	0.44
1:D:83:LYS:O	1:D:195:LYS:HD2	2.16	0.44
1:D:179:GLU:CD	1:D:181:GLU:H	2.21	0.44
1:D:36:ALA:C	1:D:37:LYS:HD2	2.38	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:40:THR:HG21	1:D:60:MET:HE2	1.99	0.44
1:D:98:VAL:HA	1:D:149:LYS:NZ	2.32	0.44
1:D:179:GLU:OE2	1:D:182:ALA:N	2.24	0.44
1:D:221:PHE:CD2	1:D:225:ILE:HD11	2.53	0.44
1:F:212:LEU:HD22	1:F:216:GLU:HB3	2.00	0.44
1:D:98:VAL:HA	1:D:149:LYS:HZ3	1.83	0.44
1:E:147:ASP:HB3	2:E:319:HOH:O	2.16	0.44
1:D:140:THR:O	1:D:144:LEU:HD13	2.18	0.44
1:F:17:LEU:HB2	1:F:59:LEU:HD23	2.00	0.44
1:A:8:LYS:HD2	1:A:9:ILE:O	2.16	0.43
1:B:43:ARG:HA	1:E:21:ASP:OD1	2.18	0.43
1:C:87:ARG:HD3	1:C:87:ARG:C	2.38	0.43
1:A:25:VAL:HG12	1:A:48:PHE:HD1	1.82	0.43
1:A:203:SER:O	1:A:212:LEU:HD23	2.18	0.43
1:B:213:SER:C	1:B:215:LYS:H	2.19	0.43
1:B:57:ILE:HD13	1:B:229:LEU:HD23	2.00	0.43
1:C:1:MET:HG2	1:C:9:ILE:CD1	2.49	0.43
1:C:30:LYS:O	2:C:403:HOH:O	2.20	0.43
1:D:5:ILE:HG22	1:D:7:ALA:H	1.83	0.43
1:A:44:ASN:CB	1:D:44:ASN:OD1	2.58	0.43
1:B:17:LEU:HD11	1:B:88:ILE:HG23	2.00	0.43
1:E:62:HIS:CE1	1:E:181:GLU:HG2	2.54	0.43
1:D:45:MET:O	2:D:302:HOH:O	2.20	0.43
1:D:73:VAL:HG13	1:D:85:LEU:HD22	2.00	0.43
1:D:231:MET:HE2	1:D:231:MET:HB3	1.99	0.43
1:F:18:LEU:HB3	1:F:62:HIS:HD2	1.84	0.43
1:A:66:ILE:HG23	1:A:184:GLY:HA3	2.00	0.43
1:D:7:ALA:HB2	1:D:40:THR:HB	2.00	0.42
1:F:98:VAL:HG12	1:F:149:LYS:HG3	2.01	0.42
1:A:38:GLU:HB2	1:A:48:PHE:CE2	2.55	0.42
1:A:102:ASP:OD2	1:A:149:LYS:HE2	2.19	0.42
1:D:40:THR:O	1:D:46:LEU:HA	2.20	0.42
1:D:42:VAL:HG12	1:D:43:ARG:HG3	2.02	0.42
1:E:221:PHE:HB3	2:E:313:HOH:O	2.19	0.42
1:D:221:PHE:HD2	1:D:225:ILE:HD11	1.84	0.42
1:D:224:MET:HA	1:D:227:LEU:HB3	2.00	0.42
1:A:8:LYS:HB3	1:A:8:LYS:HE3	1.83	0.42
1:D:28:ILE:O	1:D:32:PHE:HB2	2.20	0.42
1:F:26:SER:O	1:F:30:LYS:HG2	2.20	0.42
1:B:129:PRO:HB3	1:B:186:TYR:CZ	2.55	0.42
1:D:135:LEU:O	1:D:139:GLN:HG3	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:91:CYS:SG	1:D:200:CYS:HB2	2.60	0.42
1:F:227:LEU:O	1:F:231:MET:HG2	2.20	0.41
1:A:19:CYS:O	1:A:61:GLY:HA2	2.20	0.41
1:A:111:THR:HA	1:A:155:SER:O	2.21	0.41
1:D:56:GLY:O	1:D:57:ILE:HD13	2.20	0.41
1:F:100:LEU:O	1:F:101:LYS:HB2	2.20	0.41
1:B:221:PHE:O	1:B:225:ILE:HG13	2.20	0.41
1:A:94:ILE:HD12	1:A:171:ALA:HB2	2.01	0.41
1:D:25:VAL:HG12	1:D:48:PHE:CE1	2.55	0.41
1:D:103:ILE:HD13	1:D:227:LEU:HD22	2.02	0.41
1:D:152:ASN:HB2	1:D:174:ASN:O	2.21	0.41
1:A:43:ARG:HD3	2:D:318:HOH:O	2.20	0.41
1:B:215:LYS:O	1:B:215:LYS:HD3	2.20	0.41
1:D:224:MET:HB2	1:D:227:LEU:HD23	2.01	0.41
1:B:205:HIS:HB3	1:B:208:THR:CG2	2.47	0.41
1:D:226:ILE:O	1:D:230:GLU:N	2.41	0.41
1:A:114:LYS:O	1:A:118:VAL:HG23	2.20	0.41
1:D:93:ALA:HB1	1:D:98:VAL:HG23	2.03	0.41
1:A:45:MET:HE2	1:A:61:GLY:C	2.42	0.40
1:D:98:VAL:HG12	1:D:149:LYS:CE	2.51	0.40
1:A:132:GLU:HB3	1:A:195:LYS:NZ	2.36	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	231/233 (99%)	215 (93%)	15 (6%)	1 (0%)	34	37
1	B	231/233 (99%)	220 (95%)	10 (4%)	1 (0%)	34	37
1	C	231/233 (99%)	225 (97%)	6 (3%)	0	100	100
1	D	231/233 (99%)	209 (90%)	16 (7%)	6 (3%)	5	3

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	E	231/233 (99%)	227 (98%)	4 (2%)	0	100	100
1	F	231/233 (99%)	224 (97%)	7 (3%)	0	100	100
All	All	1386/1398 (99%)	1320 (95%)	58 (4%)	8 (1%)	25	26

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	219	GLU
1	D	209	LYS
1	D	215	LYS
1	D	231	MET
1	B	214	PRO
1	D	101	LYS
1	D	9	ILE
1	D	218	VAL

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	196/196 (100%)	190 (97%)	6 (3%)	40	51
1	B	196/196 (100%)	193 (98%)	3 (2%)	65	78
1	C	196/196 (100%)	192 (98%)	4 (2%)	55	69
1	D	196/196 (100%)	184 (94%)	12 (6%)	18	21
1	E	196/196 (100%)	193 (98%)	3 (2%)	65	78
1	F	196/196 (100%)	193 (98%)	3 (2%)	65	78
All	All	1176/1176 (100%)	1145 (97%)	31 (3%)	46	58

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	LYS
1	A	87	ARG

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Mol	Chain	Res	Type
1	A	143	ARG
1	A	179	GLU
1	A	204	ASP
1	A	221	PHE
1	B	34	GLN
1	B	153	VAL
1	B	179	GLU
1	C	87	ARG
1	C	148	LEU
1	C	179	GLU
1	C	220	SER
1	D	55	ARG
1	D	60	MET
1	D	100	LEU
1	D	130	ASP
1	D	138	TYR
1	D	142	LYS
1	D	143	ARG
1	D	195	LYS
1	D	200	CYS
1	D	209	LYS
1	D	215	LYS
1	D	222	ASP
1	E	30	LYS
1	E	55	ARG
1	E	179	GLU
1	F	87	ARG
1	F	148	LEU
1	F	179	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	122	ASN
1	E	122	ASN

5.3.3 RNA ⓘ

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

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	233/233 (100%)	0.72	24 (10%) 6 5	27, 49, 72, 98	0
1	B	233/233 (100%)	0.67	27 (11%) 4 4	26, 49, 68, 90	0
1	C	233/233 (100%)	0.12	0 100 100	22, 31, 43, 57	0
1	D	233/233 (100%)	1.06	38 (16%) 1 1	29, 55, 77, 96	0
1	E	233/233 (100%)	0.12	0 100 100	21, 31, 44, 55	0
1	F	233/233 (100%)	0.17	2 (0%) 84 83	25, 39, 59, 78	0
All	All	1398/1398 (100%)	0.48	91 (6%) 18 17	21, 40, 69, 98	0

All (91) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	215	LYS	7.0
1	A	220	SER	6.8
1	D	57	ILE	6.4
1	D	212	LEU	5.4
1	D	233	SER	4.7
1	A	214	PRO	4.7
1	D	7	ALA	4.6
1	D	231	MET	4.5
1	D	48	PHE	4.4
1	A	207	ILE	3.8
1	B	213	SER	3.6
1	D	227	LEU	3.6
1	B	233	SER	3.5
1	B	13	TYR	3.5
1	D	135	LEU	3.4
1	D	49	SER	3.4
1	D	23	LEU	3.4
1	D	134	SER	3.4
1	A	212	LEU	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	221	PHE	3.4
1	D	28	ILE	3.3
1	B	232	MET	3.2
1	D	35	ASP	3.2
1	D	226	ILE	3.1
1	A	219	GLU	3.1
1	F	200	CYS	3.1
1	D	232	MET	3.0
1	B	56	GLY	3.0
1	D	146	ILE	3.0
1	A	218	VAL	3.0
1	D	36	ALA	3.0
1	B	25	VAL	2.9
1	A	28	ILE	2.9
1	D	13	TYR	2.9
1	D	30	LYS	2.9
1	B	48	PHE	2.9
1	D	132	GLU	2.9
1	D	9	ILE	2.8
1	D	32	PHE	2.8
1	B	17	LEU	2.8
1	B	212	LEU	2.7
1	A	35	ASP	2.7
1	A	100	LEU	2.7
1	A	226	ILE	2.7
1	B	132	GLU	2.7
1	B	55	ARG	2.6
1	A	144	LEU	2.6
1	B	133	LEU	2.6
1	B	221	PHE	2.6
1	B	36	ALA	2.6
1	F	221	PHE	2.6
1	A	101	LYS	2.6
1	D	138	TYR	2.5
1	D	37	LYS	2.5
1	D	137	ALA	2.5
1	A	48	PHE	2.5
1	D	8	LYS	2.5
1	A	99	GLY	2.4
1	B	138	TYR	2.4
1	D	16	CYS	2.4
1	D	1	MET	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	82	VAL	2.4
1	A	147	ASP	2.4
1	B	37	LYS	2.3
1	B	231	MET	2.3
1	A	222	ASP	2.3
1	D	141	ALA	2.3
1	D	34	GLN	2.3
1	B	144	LEU	2.2
1	D	17	LEU	2.2
1	D	224	MET	2.2
1	B	40	THR	2.2
1	D	24	ARG	2.2
1	A	98	VAL	2.2
1	B	142	LYS	2.2
1	D	52	TYR	2.2
1	D	80	TYR	2.2
1	D	213	SER	2.2
1	A	228	ALA	2.2
1	B	135	LEU	2.2
1	B	147	ASP	2.2
1	A	225	ILE	2.1
1	A	148	LEU	2.1
1	A	229	LEU	2.1
1	B	81	GLN	2.1
1	B	16	CYS	2.1
1	A	142	LYS	2.1
1	D	83	LYS	2.0
1	B	218	VAL	2.0
1	D	31	LYS	2.0
1	B	141	ALA	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.