



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2018 – 12:07 PM EST

PDB ID : 6F52
Title : Crystal structure of H. pylori purine nucleoside phosphorylase in complex with PO4 and formycin A
Authors : Stefanic, Z.
Deposited on : 2017-11-30
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

<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
Xtriage (Phenix) : 1.9-1692
EDS : rb-20030736
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : rb-20030736

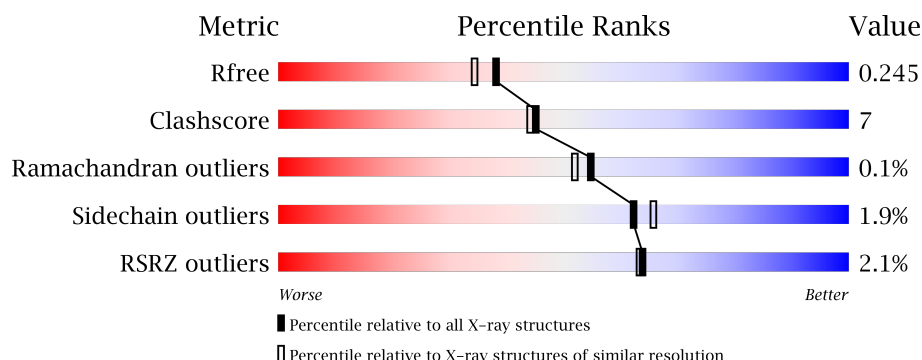
1 Overall quality at a glance

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}	100719	6609 (2.00-2.00)
Clashscore	112137	7775 (2.00-2.00)
Ramachandran outliers	110173	7679 (2.00-2.00)
Sidechain outliers	110143	7678 (2.00-2.00)
RSRZ outliers	101464	6696 (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 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	233	<div> <div>6%</div> <div> <div></div> <div>80%</div> <div>19%</div> <div>.</div> </div> </div>
1	B	233	<div> <div>%</div> <div> <div></div> <div>84%</div> <div>15%</div> <div>.</div> </div> </div>
1	C	233	<div> <div></div> <div> <div></div> <div>91%</div> <div>8%</div> <div>.</div> </div> </div>
1	D	233	<div> <div>5%</div> <div> <div></div> <div>71%</div> <div>28%</div> </div> </div>
1	E	233	<div> <div></div> <div> <div></div> <div>87%</div> <div>12%</div> <div>.</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	F	233	 89% 10% •

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 11548 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			

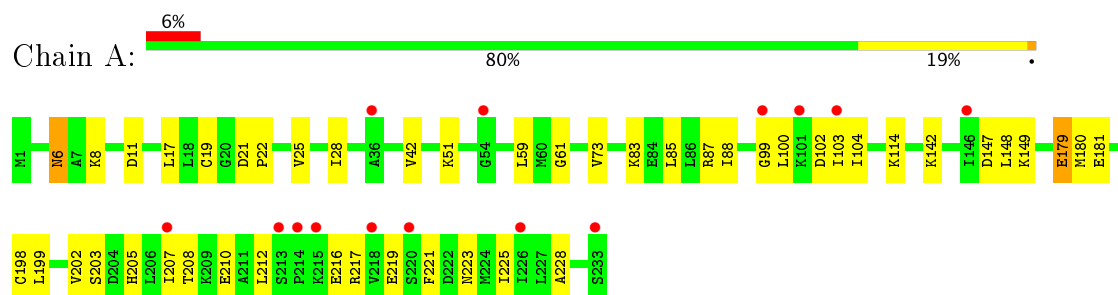
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	81	Total	O	0	0
			81	81		
2	B	82	Total	O	0	0
			82	82		
2	C	184	Total	O	0	0
			184	184		
2	D	83	Total	O	0	0
			83	83		
2	E	193	Total	O	0	0
			193	193		
2	F	113	Total	O	0	0
			113	113		

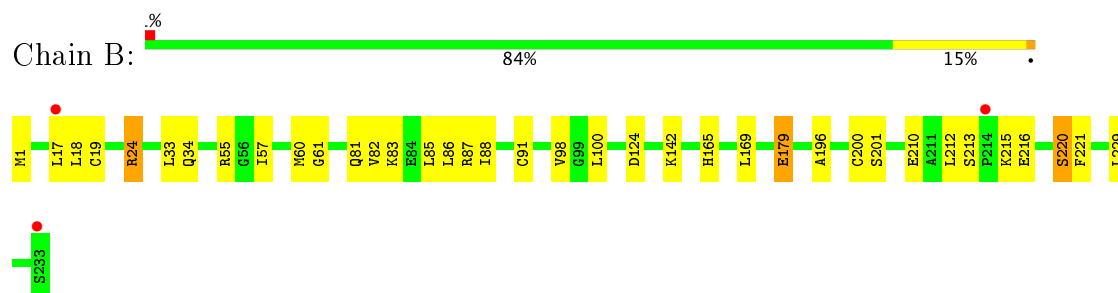
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

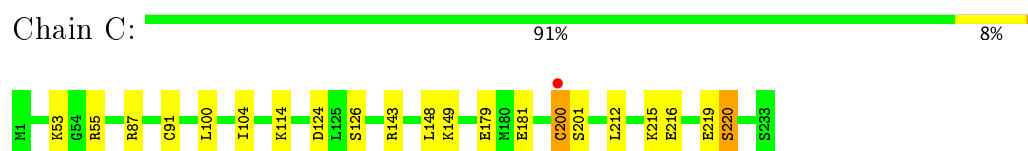
- Molecule 1: Purine nucleoside phosphorylase DeoD-type



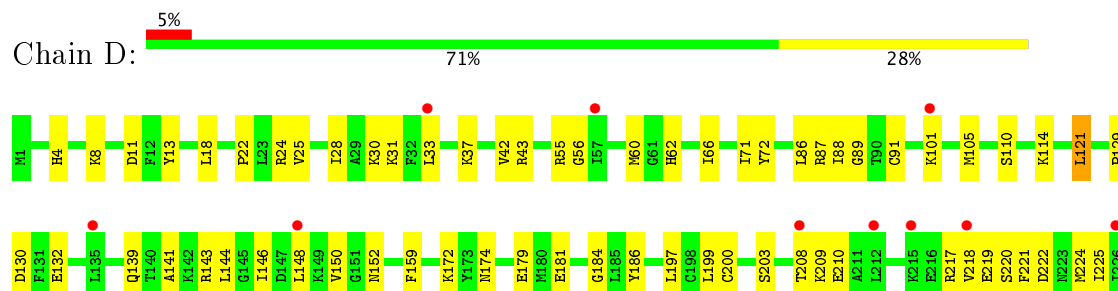
- Molecule 1: Purine nucleoside phosphorylase DeoD-type

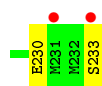


- Molecule 1: Purine nucleoside phosphorylase DeoD-type



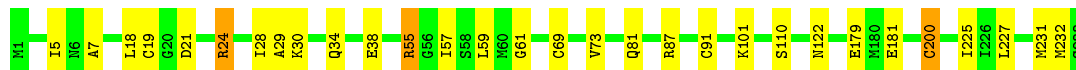
- Molecule 1: Purine nucleoside phosphorylase DeoD-type





- Molecule 1: Purine nucleoside phosphorylase DeoD-type

Chain E: 87% 12%



- Molecule 1: Purine nucleoside phosphorylase DeoD-type

Chain F: 89% 10%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	83.62Å 64.79Å 140.73Å 90.00° 99.37° 90.00°	Depositor
Resolution (Å)	47.37 – 2.00 47.37 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.1 (47.37-2.00) 99.1 (47.37-2.00)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.63 (at 2.00Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.205 , 0.246 0.204 , 0.245	Depositor DCC
R_{free} test set	5021 reflections (5.03%)	DCC
Wilson B-factor (Å ²)	28.3	Xtriage
Anisotropy	0.429	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 45.3	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	11548	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.34% 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	0.55	0/2467
1	B	0.39	0/1833	0.54	0/2467
1	C	0.45	0/1833	0.63	1/2467 (0.0%)
1	D	0.38	0/1833	0.58	0/2467
1	E	0.47	1/1833 (0.1%)	0.59	0/2467
1	F	0.48	3/1833 (0.2%)	0.58	0/2467
All	All	0.42	4/10998 (0.0%)	0.58	1/14802 (0.0%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	24	ARG	NE-CZ	-7.34	1.23	1.33
1	E	200	CYS	CB-SG	-6.98	1.70	1.82
1	F	24	ARG	CZ-NH1	-6.67	1.24	1.33
1	F	200	CYS	CB-SG	-5.76	1.72	1.81

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	200	CYS	CA-CB-SG	-8.67	98.39	114.00

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	41	0
1	B	1802	0	1838	31	0
1	C	1802	0	1836	12	0
1	D	1802	0	1838	50	0
1	E	1802	0	1838	19	0
1	F	1802	0	1836	16	0
2	A	81	0	0	4	1
2	B	82	0	0	8	0
2	C	184	0	0	2	1
2	D	83	0	0	9	1
2	E	193	0	0	7	1
2	F	113	0	0	2	0
All	All	11548	0	11024	162	2

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 7.

All (162) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:91:CYS:SG	1:E:200:CYS:HB3	2.06	0.94
1:A:208:THR:HG23	1:A:210:GLU:OE2	1.69	0.90
1:D:33:LEU:O	2:D:301:HOH:O	1.92	0.87
1:D:139:GLN:O	2:D:302:HOH:O	1.95	0.85
1:E:122:ASN:ND2	2:E:302:HOH:O	2.08	0.85
1:D:208:THR:OG1	1:D:210:GLU:OE2	1.96	0.83
1:D:210:GLU:OE2	1:D:210:GLU:N	2.10	0.83
1:E:21:ASP:HB3	1:E:24:ARG:HG3	1.60	0.83
1:A:203:SER:HB2	1:A:217:ARG:HH11	1.45	0.80
1:B:1:MET:N	2:B:304:HOH:O	2.14	0.79
1:A:208:THR:HG23	1:A:210:GLU:CD	2.03	0.78
1:B:33:LEU:HD21	1:B:57:ILE:HD11	1.68	0.75
1:E:38:GLU:OE1	2:E:301:HOH:O	2.04	0.75
1:B:142:LYS:O	2:B:301:HOH:O	2.04	0.75
1:A:51:LYS:NZ	2:A:302:HOH:O	2.19	0.75
1:B:82:VAL:O	2:B:302:HOH:O	2.05	0.72
1:E:81:GLN:NE2	2:E:303:HOH:O	2.22	0.72
1:D:105:MET:HB2	1:D:148:LEU:HD21	1.71	0.72
1:D:101:LYS:HE3	1:D:220:SER:HB3	1.72	0.71
1:A:147:ASP:OD2	1:A:149:LYS:NZ	2.22	0.70
1:D:87:ARG:HH12	1:D:181:GLU:HB2	1.57	0.70
1:E:55:ARG:NH2	2:E:304:HOH:O	2.24	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:210:GLU:N	1:A:210:GLU:OE2	2.26	0.68
1:B:34:GLN:NE2	2:B:307:HOH:O	2.23	0.67
1:B:98:VAL:O	2:B:303:HOH:O	2.13	0.67
1:F:104:ILE:HB	1:F:200:CYS:HB3	1.77	0.66
1:A:205:HIS:CE1	1:A:207:ILE:HG12	2.30	0.66
1:D:144:LEU:HD21	1:D:230:GLU:HG3	1.79	0.65
1:A:87:ARG:HH12	1:A:181:GLU:HB2	1.62	0.65
1:D:88:ILE:HB	1:D:224:MET:HE3	1.80	0.64
1:F:130:ASP:OD2	1:F:195:LYS:NZ	2.30	0.64
1:A:142:LYS:O	2:A:301:HOH:O	2.15	0.64
2:A:331:HOH:O	1:F:126:SER:HB2	1.99	0.63
1:B:34:GLN:OE1	2:B:306:HOH:O	2.16	0.62
1:A:114:LYS:HB2	1:D:114:LYS:HB2	1.81	0.62
1:D:91:CYS:SG	1:D:200:CYS:HB3	2.40	0.61
1:C:215:LYS:O	1:C:219:GLU:HG2	2.01	0.60
1:B:165:HIS:ND1	2:B:308:HOH:O	2.31	0.60
1:A:100:LEU:HD11	1:A:210:GLU:HB3	1.81	0.60
1:A:205:HIS:HE1	1:A:207:ILE:HG12	1.65	0.60
1:A:17:LEU:HD21	1:A:228:ALA:HB1	1.83	0.59
1:C:126:SER:HB2	2:E:367:HOH:O	2.00	0.59
1:D:203:SER:OG	2:D:303:HOH:O	2.17	0.59
1:B:212:LEU:HB3	1:B:216:GLU:HG3	1.85	0.58
1:A:203:SER:HB2	1:A:217:ARG:NH1	2.18	0.57
1:A:8:LYS:HG2	1:A:11:ASP:OD1	2.05	0.56
1:A:21:ASP:OD1	1:D:43:ARG:HA	2.06	0.56
1:A:205:HIS:CG	1:A:208:THR:HG22	2.40	0.56
1:A:219:GLU:O	1:A:223:ASN:ND2	2.38	0.56
1:E:227:LEU:O	1:E:231:MET:HG2	2.06	0.55
1:C:114:LYS:HB2	1:F:114:LYS:HB2	1.87	0.55
1:B:57:ILE:HD12	1:B:229:LEU:HD22	1.87	0.55
1:D:141:ALA:HB1	1:D:146:ILE:HB	1.89	0.55
1:F:24:ARG:NH2	2:F:306:HOH:O	2.41	0.54
1:D:13:TYR:CE1	1:D:56:GLY:HA3	2.44	0.53
1:D:55:ARG:HD3	1:D:233:SER:HA	1.91	0.53
1:C:200:CYS:SG	1:C:201:SER:N	2.80	0.53
1:E:57:ILE:HD11	1:E:232:MET:HE1	1.92	0.52
1:D:222:ASP:HA	1:D:225:ILE:HD12	1.92	0.52
1:E:87:ARG:HH12	1:E:181:GLU:HB2	1.75	0.52
1:B:33:LEU:CD2	1:B:57:ILE:HD11	2.40	0.51
1:B:216:GLU:O	1:B:220:SER:OG	2.27	0.51
1:D:105:MET:CB	1:D:148:LEU:HD21	2.39	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:143:ARG:N	2:D:302:HOH:O	2.36	0.51
1:C:55:ARG:NH2	2:C:304:HOH:O	2.44	0.50
1:E:57:ILE:HD11	1:E:232:MET:CE	2.41	0.50
1:A:208:THR:CG2	1:A:210:GLU:OE1	2.60	0.50
1:E:69:CYS:O	1:E:73:VAL:HG13	2.11	0.50
1:B:213:SER:H	1:B:216:GLU:HG3	1.76	0.50
1:F:13:TYR:CE1	1:F:56:GLY:HA3	2.47	0.49
1:A:19:CYS:O	1:A:61:GLY:HA2	2.13	0.49
1:F:98:VAL:HG12	1:F:149:LYS:HG3	1.93	0.49
1:A:87:ARG:NH2	1:A:198:CYS:SG	2.85	0.49
1:B:213:SER:OG	1:B:216:GLU:HG2	2.12	0.48
1:A:6:ASN:HB2	1:A:42:VAL:HG23	1.94	0.48
1:E:34:GLN:HG3	2:E:318:HOH:O	2.13	0.48
1:B:213:SER:H	1:B:216:GLU:CG	2.26	0.48
1:D:208:THR:O	1:D:209:LYS:HB2	2.13	0.48
1:F:227:LEU:O	1:F:231:MET:HG2	2.13	0.48
1:A:99:GLY:O	1:A:202:VAL:HG21	2.13	0.48
1:B:18:LEU:HD23	1:B:60:MET:HB3	1.96	0.48
1:E:28:ILE:HG12	1:E:225:ILE:HD13	1.95	0.48
1:D:4:HIS:NE2	2:D:306:HOH:O	2.35	0.48
1:B:17:LEU:HD22	1:B:86:LEU:O	2.14	0.48
1:D:42:VAL:HG12	1:D:43:ARG:HG3	1.96	0.47
1:D:159:PHE:N	2:D:305:HOH:O	2.47	0.47
1:B:216:GLU:HG2	1:B:216:GLU:H	1.54	0.47
1:D:18:LEU:HA	1:D:60:MET:O	2.15	0.47
1:A:102:ASP:O	1:A:202:VAL:HG22	2.14	0.46
1:D:8:LYS:O	1:D:11:ASP:HB2	2.15	0.46
1:C:87:ARG:HH12	1:C:181:GLU:HB2	1.80	0.46
1:B:200:CYS:HB3	2:B:337:HOH:O	2.14	0.46
1:A:22:PRO:HA	1:A:25:VAL:HG22	1.97	0.46
1:F:85:LEU:O	1:F:196:ALA:HA	2.15	0.46
1:D:130:ASP:OD2	1:D:132:GLU:HG3	2.16	0.46
1:C:100:LEU:HB3	1:C:212:LEU:HD21	1.97	0.46
1:A:100:LEU:HD11	1:A:210:GLU:CB	2.46	0.46
1:D:152:ASN:HB2	1:D:174:ASN:O	2.16	0.46
1:A:28:ILE:HG12	1:A:225:ILE:HD13	1.97	0.46
1:B:91:CYS:SG	1:B:200:CYS:SG	3.13	0.46
1:A:179:GLU:HB2	2:A:327:HOH:O	2.15	0.45
1:C:91:CYS:N	1:C:200:CYS:SG	2.89	0.45
1:B:24:ARG:HD2	1:B:221:PHE:CE1	2.51	0.45
1:D:221:PHE:HD1	1:D:225:ILE:HD11	1.81	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:24:ARG:HD2	1:B:221:PHE:CZ	2.52	0.45
1:D:66:ILE:HG23	1:D:184:GLY:HA3	1.99	0.45
1:A:17:LEU:HB2	1:A:59:LEU:HD23	1.99	0.45
1:C:216:GLU:O	1:C:220:SER:HB2	2.17	0.44
1:D:208:THR:OG1	1:D:210:GLU:CD	2.53	0.44
1:D:13:TYR:CZ	1:D:56:GLY:HA3	2.52	0.44
1:C:124:ASP:O	1:E:110:SER:HB3	2.18	0.44
1:F:163:GLU:OE2	2:F:301:HOH:O	2.21	0.44
1:C:143:ARG:HD2	2:C:302:HOH:O	2.18	0.44
1:E:18:LEU:HB2	1:E:87:ARG:HG2	1.99	0.44
1:F:57:ILE:HG13	1:F:58:SER:N	2.31	0.44
1:B:87:ARG:HD2	1:B:88:ILE:N	2.33	0.44
1:A:208:THR:C	1:A:210:GLU:OE2	2.56	0.43
1:A:221:PHE:HD1	1:A:225:ILE:HD11	1.84	0.43
1:D:172:LYS:HE3	2:D:378:HOH:O	2.17	0.43
1:E:5:ILE:HG22	1:E:7:ALA:H	1.83	0.43
1:D:208:THR:HG1	1:D:210:GLU:CD	2.12	0.43
1:D:62:HIS:HA	1:D:72:TYR:CD2	2.54	0.43
1:D:143:ARG:HB3	2:D:302:HOH:O	2.17	0.43
1:D:89:GLY:H	1:D:224:MET:HE3	1.83	0.43
1:D:86:LEU:HD23	1:D:197:LEU:HB3	2.01	0.42
1:B:215:LYS:HB3	1:B:215:LYS:HE3	1.80	0.42
1:B:81:GLN:O	1:B:83:LYS:HD2	2.19	0.42
1:F:130:ASP:CG	1:F:195:LYS:HG2	2.39	0.42
1:B:124:ASP:O	1:D:110:SER:HB3	2.20	0.42
1:B:100:LEU:HD11	1:B:210:GLU:HB3	2.02	0.42
1:D:30:LYS:HG3	1:D:31:LYS:N	2.35	0.42
1:F:5:ILE:HG22	1:F:7:ALA:H	1.85	0.42
1:A:180:MET:CE	1:D:71:ILE:HD11	2.50	0.42
1:A:73:VAL:HG13	1:A:85:LEU:HD22	2.02	0.42
1:D:217:ARG:C	1:D:219:GLU:H	2.22	0.42
1:A:103:ILE:HG23	1:A:148:LEU:HD12	2.00	0.42
1:F:8:LYS:HB2	1:F:8:LYS:HE2	1.76	0.42
1:D:129:PRO:HB3	1:D:186:TYR:CZ	2.54	0.41
1:A:212:LEU:HB3	1:A:216:GLU:HB3	2.02	0.41
1:A:19:CYS:HB2	1:A:25:VAL:HG13	2.02	0.41
1:C:104:ILE:HA	1:C:149:LYS:O	2.20	0.41
1:D:87:ARG:NH1	1:D:181:GLU:HB2	2.30	0.41
1:D:208:THR:CB	1:D:210:GLU:OE2	2.68	0.41
1:A:88:ILE:HA	1:A:199:LEU:O	2.20	0.41
1:B:85:LEU:O	1:B:196:ALA:HA	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:148:LEU:HD22	1:D:150:VAL:HG22	2.01	0.41
1:F:140:THR:HA	1:F:143:ARG:HG2	2.01	0.41
1:B:87:ARG:NH2	1:B:179:GLU:OE2	2.53	0.41
1:F:213:SER:O	1:F:217:ARG:HG3	2.20	0.41
1:A:104:ILE:HA	1:A:149:LYS:O	2.19	0.41
1:D:88:ILE:HG22	1:D:199:LEU:HB2	2.02	0.41
1:E:101:LYS:HD3	2:E:454:HOH:O	2.19	0.41
1:A:205:HIS:ND1	1:A:208:THR:HG22	2.36	0.41
1:B:169:LEU:HD23	1:D:121:LEU:CD1	2.50	0.41
1:D:24:ARG:O	1:D:28:ILE:HG13	2.21	0.41
1:D:22:PRO:O	1:D:25:VAL:HB	2.20	0.41
1:E:19:CYS:O	1:E:61:GLY:HA2	2.20	0.41
1:A:208:THR:HG23	1:A:210:GLU:OE1	2.19	0.40
1:A:205:HIS:HB3	1:A:208:THR:HG22	2.03	0.40
1:D:132:GLU:HG2	2:D:355:HOH:O	2.20	0.40
1:B:19:CYS:O	1:B:61:GLY:HA2	2.22	0.40
1:E:29:ALA:HB2	1:E:59:LEU:HD12	2.03	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:460:HOH:O	2:E:447:HOH:O[2_546]	1.99	0.21
2:A:375:HOH:O	2:D:322:HOH:O[2_645]	2.14	0.06

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%)	220 (95%)	11 (5%)	0	100	100
1	B	231/233 (99%)	222 (96%)	9 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	231/233 (99%)	226 (98%)	5 (2%)	0	100	100
1	D	231/233 (99%)	220 (95%)	10 (4%)	1 (0%)	38	33
1	E	231/233 (99%)	226 (98%)	5 (2%)	0	100	100
1	F	231/233 (99%)	224 (97%)	7 (3%)	0	100	100
All	All	1386/1398 (99%)	1338 (96%)	47 (3%)	1 (0%)	55	52

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
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%)	193 (98%)	3 (2%)	70	74
1	B	196/196 (100%)	191 (97%)	5 (3%)	51	52
1	C	196/196 (100%)	192 (98%)	4 (2%)	60	64
1	D	196/196 (100%)	193 (98%)	3 (2%)	70	74
1	E	196/196 (100%)	192 (98%)	4 (2%)	60	64
1	F	196/196 (100%)	193 (98%)	3 (2%)	70	74
All	All	1176/1176 (100%)	1154 (98%)	22 (2%)	62	66

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

Mol	Chain	Res	Type
1	A	6	ASN
1	A	83	LYS
1	A	179	GLU
1	B	24	ARG
1	B	55	ARG
1	B	179	GLU

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Mol	Chain	Res	Type
1	B	201	SER
1	B	220	SER
1	C	53	LYS
1	C	148	LEU
1	C	179	GLU
1	C	220	SER
1	D	37	LYS
1	D	121	LEU
1	D	179	GLU
1	E	24	ARG
1	E	30	LYS
1	E	55	ARG
1	E	179	GLU
1	F	57	ILE
1	F	179	GLU
1	F	200	CYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	15	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](#)

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å ²)	Q<0.9
1	A	233/233 (100%)	0.39	14 (6%)	23 23	24, 43, 61, 76	0
1	B	233/233 (100%)	0.17	3 (1%)	77 77	23, 41, 57, 70	0
1	C	233/233 (100%)	-0.23	1 (0%)	92 92	18, 27, 39, 48	0
1	D	233/233 (100%)	0.45	12 (5%)	28 28	26, 44, 66, 75	0
1	E	233/233 (100%)	-0.26	0	100 100	18, 27, 37, 48	0
1	F	233/233 (100%)	-0.19	0	100 100	21, 35, 53, 62	0
All	All	1398/1398 (100%)	0.06	30 (2%)	64 63	18, 35, 57, 76	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	218	VAL	4.4
1	A	99	GLY	4.1
1	D	212	LEU	3.5
1	A	103	ILE	3.4
1	B	17	LEU	3.2
1	A	214	PRO	3.1
1	D	226	ILE	3.1
1	A	207	ILE	3.0
1	B	233	SER	3.0
1	D	208	THR	2.9
1	D	135	LEU	2.8
1	A	215	LYS	2.8
1	D	101	LYS	2.6
1	B	214	PRO	2.5
1	A	226	ILE	2.4
1	A	220	SER	2.4
1	A	213	SER	2.3
1	A	36	ALA	2.3
1	D	148	LEU	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	146	ILE	2.3
1	D	215	LYS	2.2
1	C	200	CYS	2.2
1	D	33	LEU	2.1
1	D	231	MET	2.1
1	A	233	SER	2.1
1	A	218	VAL	2.1
1	A	54	GLY	2.1
1	D	233	SER	2.1
1	A	101	LYS	2.0
1	D	57	ILE	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.