



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 17, 2023 – 04:41 AM EDT

PDB ID : 2CQZ
Title : Crystal Structure of PH0347 protein from *Pyrococcus horikoshii* OT3
Authors : Lokanath, N.K.; Terao, Y.; Kunishima, N.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)
Deposited on : 2005-05-20
Resolution : 2.60 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 : 2.36
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) : 2.36

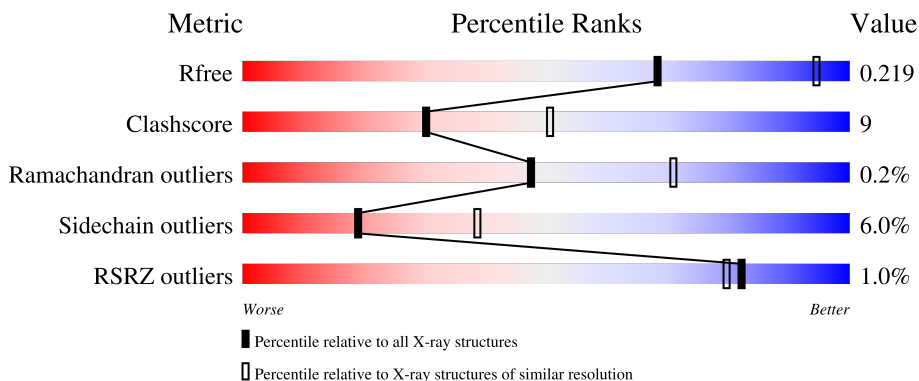
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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.



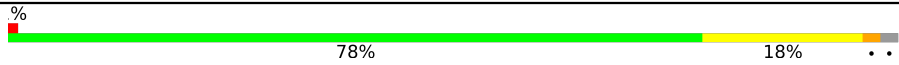
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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	177	 76% 20% ..
1	B	177	 75% 20% ..
1	C	177	 72% 25% ..
1	D	177	 72% 23% ..
1	E	177	 71% 23% ...

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Mol	Chain	Length	Quality of chain
1	F	177	 <p>A horizontal bar chart representing the quality of chain. The bar is divided into segments: a small red segment at the start, followed by a large green segment labeled '78%', then a yellow segment labeled '18%', and finally a small grey segment at the end. A '%' symbol is positioned above the bar on the left, and two dots are positioned below the bar on the right.</p>

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 8625 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 177aa long hypothetical protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	173	1408	912	229	262	5	0	0	0
1	B	173	1408	912	229	262	5	0	0	0
1	C	173	1408	912	229	262	5	0	0	0
1	D	173	1408	912	229	262	5	0	0	0
1	E	173	1408	912	229	262	5	0	0	0
1	F	173	1408	912	229	262	5	0	0	0

- Molecule 2 is NICKEL (II) ION (three-letter code: NI) (formula: Ni).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total 1	Ni 1	0	0
2	B	1	Total 1	Ni 1	0	0
2	C	1	Total 1	Ni 1	0	0
2	D	1	Total 1	Ni 1	0	0
2	E	1	Total 1	Ni 1	0	0
2	F	1	Total 1	Ni 1	0	0

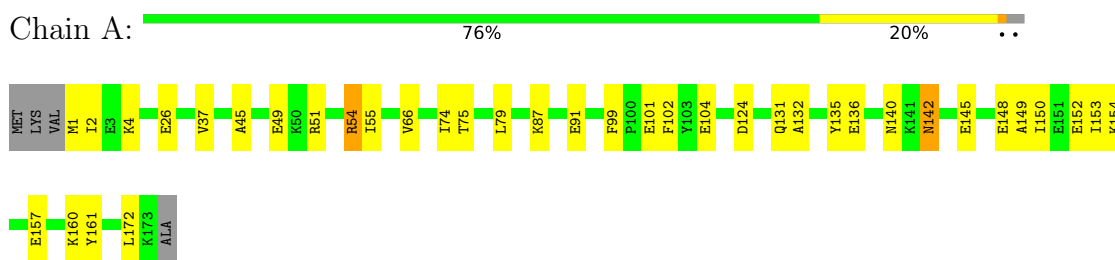
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	29	Total O 29 29	0	0
3	B	26	Total O 26 26	0	0
3	C	28	Total O 28 28	0	0
3	D	37	Total O 37 37	0	0
3	E	24	Total O 24 24	0	0
3	F	27	Total O 27 27	0	0

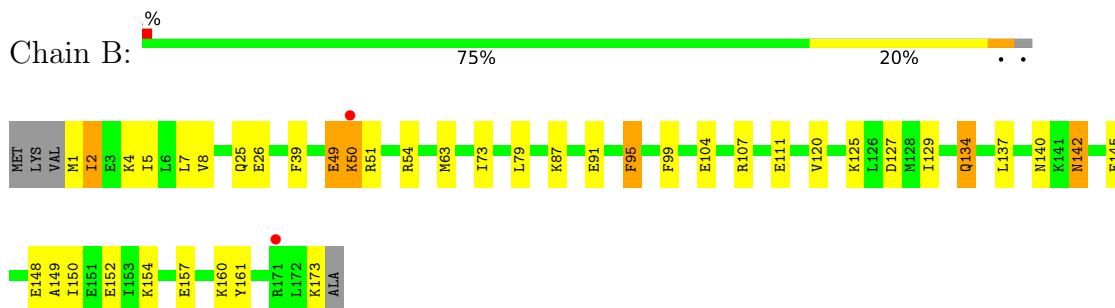
3 Residue-property plots

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.

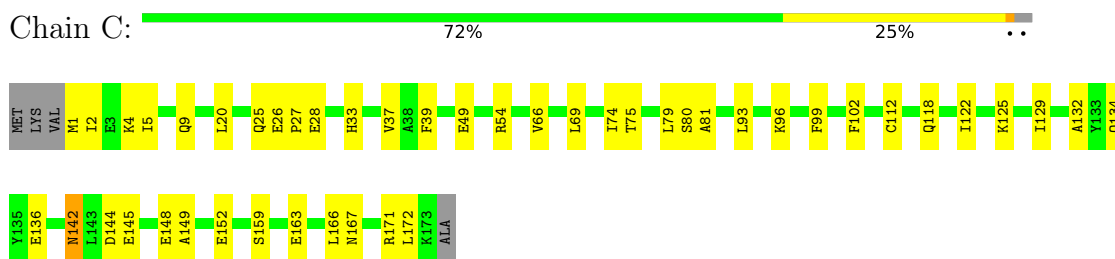
- Molecule 1: 177aa long hypothetical protein



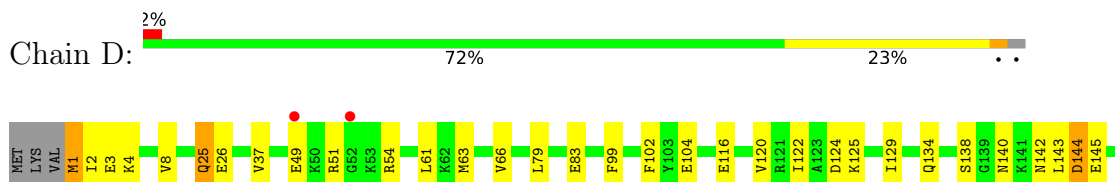
- Molecule 1: 177aa long hypothetical protein

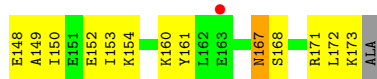


- Molecule 1: 177aa long hypothetical protein



- Molecule 1: 177aa long hypothetical protein

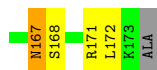
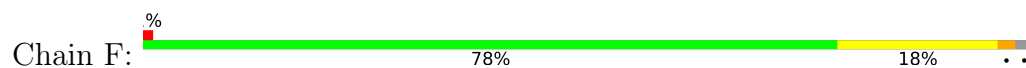




- Molecule 1: 177aa long hypothetical protein



- Molecule 1: 177aa long hypothetical protein



4 Data and refinement statistics i

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	78.51Å 101.17Å 78.65Å 90.00° 119.78° 90.00°	Depositor
Resolution (Å)	40.00 – 2.60 40.64 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.7 (40.00-2.60) 99.4 (40.64-2.60)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.31 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.1.27	Depositor
R, R_{free}	0.195 , 0.225 0.194 , 0.219	Depositor DCC
R_{free} test set	1613 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	37.3	Xtriage
Anisotropy	0.612	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 45.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.063 for -h-l,k,h 0.063 for l,k,-h-l 0.067 for h,-k,-h-l 0.126 for -h-l,-k,l 0.079 for l,-k,h	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	8625	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.53% 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](#)

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

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	0.43	0/1431	0.68	0/1927
1	B	0.43	0/1431	0.71	1/1927 (0.1%)
1	C	0.45	0/1431	0.70	0/1927
1	D	0.41	0/1431	0.68	0/1927
1	E	0.49	0/1431	0.76	1/1927 (0.1%)
1	F	0.43	0/1431	0.67	0/1927
All	All	0.44	0/8586	0.70	2/11562 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	25	GLN	O-C-N	-6.01	113.08	122.70
1	B	95	PHE	O-C-N	-5.56	113.80	122.70

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	1408	0	1454	26	0
1	B	1408	0	1454	29	0
1	C	1408	0	1454	30	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	1408	0	1454	36	0
1	E	1408	0	1454	40	0
1	F	1408	0	1454	29	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
3	A	29	0	0	0	0
3	B	26	0	0	4	0
3	C	28	0	0	1	0
3	D	37	0	0	2	0
3	E	24	0	0	1	0
3	F	27	0	0	0	0
All	All	8625	0	8724	160	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 (160) 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:163:GLU:OE2	1:E:163:GLU:HA	1.47	1.08
1:E:163:GLU:OE2	1:E:163:GLU:CA	2.21	0.87
1:B:145:GLU:O	1:B:148:GLU:HG2	1.75	0.87
1:F:145:GLU:O	1:F:148:GLU:HG2	1.77	0.85
1:A:54:ARG:HD2	1:A:55:ILE:N	1.93	0.84
1:D:145:GLU:O	1:D:148:GLU:HG2	1.81	0.81
1:E:146:PHE:HZ	1:E:170:GLY:HA2	1.48	0.79
1:D:140:ASN:HD22	1:D:143:LEU:HG	1.47	0.79
1:C:145:GLU:O	1:C:148:GLU:HG2	1.82	0.79
1:A:145:GLU:O	1:A:148:GLU:HG2	1.83	0.78
1:B:25:GLN:HB2	1:C:25:GLN:HG2	1.65	0.77
1:D:25:GLN:HE21	1:F:25:GLN:HG3	1.56	0.71
1:E:25:GLN:HG3	3:E:616:HOH:O	1.92	0.68
1:B:134:GLN:NE2	3:B:627:HOH:O	2.29	0.65
1:A:101:GLU:OE1	1:A:101:GLU:N	2.26	0.63
1:F:4:LYS:HD2	1:F:99:PHE:CE1	2.34	0.63
1:D:149:ALA:HA	1:D:152:GLU:HG3	1.81	0.62
1:E:146:PHE:CZ	1:E:170:GLY:HA2	2.32	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:142:ASN:HD22	1:F:137:LEU:HD13	1.66	0.61
1:E:122:ILE:HG12	1:E:153:ILE:HG23	1.82	0.61
1:A:45:ALA:O	1:A:49:GLU:HG3	2.00	0.61
1:D:144:ASP:HB2	3:D:635:HOH:O	2.00	0.61
1:A:150:ILE:O	1:A:154:LYS:HG2	1.99	0.61
1:D:142:ASN:ND2	1:F:137:LEU:HB3	2.15	0.60
1:B:125:LYS:O	1:B:129:ILE:HG13	2.01	0.60
1:B:39:PHE:CZ	1:E:8:VAL:HG21	2.37	0.59
3:B:621:HOH:O	1:C:25:GLN:HG3	2.02	0.58
1:F:101:GLU:OE1	1:F:101:GLU:N	2.34	0.58
1:A:172:LEU:HD11	1:F:2:ILE:HG21	1.84	0.58
1:F:4:LYS:HD2	1:F:99:PHE:HE1	1.67	0.58
1:C:2:ILE:HG21	1:D:172:LEU:HD11	1.85	0.58
1:A:37:VAL:HG22	1:A:124:ASP:HA	1.86	0.58
1:A:1:MET:HG3	1:A:4:LYS:H	1.69	0.57
1:B:4:LYS:HD2	1:B:99:PHE:HE1	1.69	0.57
1:D:1:MET:SD	1:D:3:GLU:HB3	2.43	0.57
1:B:8:VAL:HG21	1:E:39:PHE:CZ	2.39	0.57
1:B:149:ALA:HA	1:B:152:GLU:HG3	1.87	0.56
1:B:4:LYS:HD2	1:B:99:PHE:CE1	2.40	0.56
1:D:173:LYS:NZ	3:D:638:HOH:O	2.38	0.56
1:C:81:ALA:CB	1:E:74:ILE:HA	2.36	0.56
1:F:1:MET:HG3	1:F:4:LYS:H	1.70	0.56
1:E:150:ILE:HG22	1:E:151:GLU:N	2.20	0.56
1:F:149:ALA:HA	1:F:152:GLU:HG3	1.88	0.55
1:D:25:GLN:NE2	1:F:25:GLN:HG3	2.21	0.55
1:E:126:LEU:O	1:E:130:LEU:HB2	2.06	0.55
1:A:66:VAL:HG11	1:A:102:PHE:HB3	1.89	0.55
1:F:5:ILE:O	1:F:9:GLN:HG3	2.07	0.54
1:C:81:ALA:HB1	1:E:74:ILE:HA	1.90	0.54
1:D:4:LYS:HD2	1:D:99:PHE:HE1	1.73	0.54
1:C:74:ILE:O	1:C:75:THR:OG1	2.24	0.54
1:E:4:LYS:HD2	1:E:99:PHE:CE1	2.43	0.54
1:C:132:ALA:O	1:C:136:GLU:HG3	2.08	0.54
1:D:25:GLN:HG2	1:F:25:GLN:HG3	1.89	0.53
1:A:54:ARG:CD	1:A:55:ILE:N	2.70	0.53
1:D:140:ASN:ND2	1:D:143:LEU:HG	2.20	0.53
1:E:4:LYS:HD2	1:E:99:PHE:HE1	1.73	0.53
1:C:39:PHE:CZ	1:D:8:VAL:HG21	2.43	0.53
1:D:4:LYS:HD2	1:D:99:PHE:CE1	2.43	0.53
1:C:81:ALA:HB2	1:E:73:ILE:O	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:66:VAL:HG11	1:C:102:PHE:HB3	1.90	0.52
1:C:1:MET:HG3	1:C:4:LYS:H	1.75	0.52
1:E:163:GLU:O	1:E:167:ASN:ND2	2.43	0.52
1:A:87:LYS:O	1:A:91:GLU:HG3	2.11	0.51
1:C:172:LEU:HD11	1:D:2:ILE:HG21	1.92	0.51
1:C:118:GLN:O	1:C:122:ILE:HG13	2.10	0.51
1:C:5:ILE:O	1:C:9:GLN:HG3	2.10	0.51
1:E:146:PHE:CD2	1:E:146:PHE:C	2.84	0.51
1:E:66:VAL:HG11	1:E:102:PHE:HB3	1.93	0.51
1:E:144:ASP:OD1	1:E:145:GLU:N	2.43	0.51
1:B:1:MET:HG3	1:B:4:LYS:H	1.76	0.50
1:E:83:GLU:O	1:E:83:GLU:HG2	2.10	0.50
1:E:129:ILE:HG23	1:E:146:PHE:HD1	1.76	0.50
1:E:125:LYS:O	1:E:129:ILE:HG13	2.11	0.50
1:B:173:LYS:HE3	3:B:616:HOH:O	2.12	0.50
1:A:140:ASN:OD1	1:A:142:ASN:ND2	2.44	0.50
1:D:25:GLN:OE1	1:D:25:GLN:N	2.42	0.50
1:B:51:ARG:NH1	1:B:161:TYR:O	2.46	0.49
1:E:25:GLN:H	1:E:25:GLN:CD	2.15	0.49
1:A:26:GLU:HG2	1:C:26:GLU:OE1	2.12	0.49
1:F:125:LYS:HZ1	1:F:152:GLU:HB3	1.77	0.49
1:A:54:ARG:HD2	1:A:54:ARG:C	2.33	0.48
1:D:66:VAL:HG11	1:D:102:PHE:HB3	1.93	0.48
1:D:140:ASN:ND2	1:D:142:ASN:OD1	2.40	0.48
1:B:95:PHE:HA	1:B:99:PHE:HB2	1.95	0.48
1:E:1:MET:HG3	1:E:4:LYS:H	1.79	0.48
1:A:2:ILE:HG21	1:F:172:LEU:HD11	1.95	0.47
1:C:125:LYS:O	1:C:129:ILE:HG13	2.14	0.47
1:A:74:ILE:O	1:A:75:THR:OG1	2.26	0.47
1:B:157:GLU:O	1:B:160:LYS:HB3	2.15	0.47
1:D:26:GLU:OE1	1:E:26:GLU:HG2	2.14	0.47
1:B:49:GLU:HG3	1:B:50:LYS:N	2.28	0.47
1:C:167:ASN:O	1:C:171:ARG:HG3	2.14	0.47
1:A:157:GLU:O	1:A:160:LYS:HB3	2.16	0.46
1:F:167:ASN:HD22	1:F:167:ASN:N	2.14	0.46
1:B:150:ILE:O	1:B:154:LYS:HG2	2.15	0.46
1:F:25:GLN:H	1:F:25:GLN:HG2	1.19	0.46
1:B:2:ILE:HG21	1:E:172:LEU:HD11	1.98	0.46
1:B:5:ILE:HD13	1:E:40:ILE:HG12	1.99	0.45
1:B:63:MET:HB3	1:B:120:VAL:HG21	1.98	0.45
1:A:149:ALA:HA	1:A:152:GLU:HG3	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:26:GLU:OE1	1:B:26:GLU:HG2	2.17	0.45
1:A:4:LYS:HD2	1:A:99:PHE:CE1	2.52	0.45
1:C:33:HIS:O	1:C:37:VAL:HG23	2.16	0.45
1:C:80:SER:OG	3:C:617:HOH:O	2.21	0.45
1:E:37:VAL:HG22	1:E:124:ASP:HA	1.98	0.45
1:F:63:MET:O	1:F:67:HIS:HB3	2.17	0.45
1:F:125:LYS:NZ	1:F:152:GLU:HB3	2.31	0.44
1:E:137:LEU:HB3	1:F:142:ASN:OD1	2.17	0.44
1:D:63:MET:HB3	1:D:120:VAL:HG21	1.98	0.44
1:C:93:LEU:HB3	1:E:84:PHE:HD1	1.82	0.44
1:F:168:SER:O	1:F:171:ARG:HB3	2.18	0.44
1:D:25:GLN:CG	1:F:25:GLN:HG3	2.48	0.44
1:E:144:ASP:OD1	1:E:145:GLU:HG2	2.18	0.44
1:D:2:ILE:N	1:D:2:ILE:CD1	2.80	0.43
1:C:149:ALA:HA	1:C:152:GLU:HG3	1.99	0.43
1:B:73:ILE:O	1:F:81:ALA:HB2	2.17	0.43
1:F:132:ALA:O	1:F:136:GLU:HG3	2.18	0.43
1:D:138:SER:HB3	1:E:25:GLN:NE2	2.33	0.43
1:A:51:ARG:NH1	1:A:161:TYR:O	2.49	0.43
1:D:1:MET:HG3	1:D:4:LYS:H	1.83	0.43
1:A:2:ILE:HD12	1:A:2:ILE:N	2.34	0.43
1:E:25:GLN:OE1	1:F:25:GLN:NE2	2.52	0.43
1:F:125:LYS:NZ	1:F:152:GLU:CB	2.81	0.43
1:A:131:GLN:NE2	1:A:135:TYR:CZ	2.87	0.43
1:B:26:GLU:OE1	1:C:26:GLU:HG2	2.18	0.43
1:D:122:ILE:HG12	1:D:153:ILE:HG23	2.00	0.43
1:E:170:GLY:O	1:E:173:LYS:HB3	2.18	0.43
1:F:22:LYS:NZ	1:F:145:GLU:OE2	2.45	0.43
1:B:107:ARG:HD2	1:B:111:GLU:OE2	2.20	0.42
1:D:125:LYS:O	1:D:129:ILE:HG13	2.19	0.42
1:C:4:LYS:HD2	1:C:99:PHE:CE1	2.55	0.42
1:B:140:ASN:OD1	1:B:142:ASN:ND2	2.53	0.42
1:B:2:ILE:N	1:B:2:ILE:CD1	2.83	0.42
1:D:150:ILE:O	1:D:154:LYS:HG2	2.20	0.42
1:D:37:VAL:HG22	1:D:124:ASP:HA	2.02	0.42
1:D:51:ARG:NH1	1:D:161:TYR:O	2.50	0.42
1:F:83:GLU:O	1:F:83:GLU:HG2	2.19	0.42
1:C:74:ILE:C	1:C:75:THR:HG23	2.40	0.42
1:E:20:LEU:HG	1:E:27:PRO:HG3	2.01	0.42
1:C:159:SER:HB2	1:C:166:LEU:HD11	2.02	0.41
1:D:167:ASN:N	1:D:167:ASN:HD22	2.18	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:137:LEU:HB3	1:C:142:ASN:OD1	2.21	0.41
1:B:4:LYS:HE2	1:B:4:LYS:HB2	1.95	0.41
1:D:83:GLU:HG2	1:D:83:GLU:O	2.20	0.41
1:E:51:ARG:NH1	1:E:161:TYR:O	2.50	0.41
1:E:61:LEU:O	1:E:65:ILE:HG13	2.19	0.41
1:F:150:ILE:O	1:F:154:LYS:HG2	2.19	0.41
1:B:87:LYS:O	1:B:91:GLU:HG3	2.21	0.41
1:A:132:ALA:O	1:A:136:GLU:HG3	2.20	0.41
3:B:621:HOH:O	1:C:25:GLN:CG	2.67	0.41
1:C:20:LEU:HG	1:C:27:PRO:HG3	2.02	0.41
1:D:116:GLU:H	1:D:116:GLU:CD	2.23	0.41
1:C:66:VAL:O	1:C:69:LEU:HB2	2.20	0.41
1:D:61:LEU:HD23	1:D:61:LEU:HA	1.90	0.41
1:D:168:SER:O	1:D:171:ARG:HB3	2.21	0.41
1:E:4:LYS:O	1:E:7:LEU:HB3	2.21	0.41
1:A:149:ALA:O	1:A:153:ILE:HG13	2.20	0.40
1:E:134:GLN:HE21	1:E:134:GLN:HB3	1.67	0.40
1:A:1:MET:CE	1:A:4:LYS:HE2	2.51	0.40
1:B:127:ASP:OD1	1:E:9:GLN:NE2	2.54	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	171/177 (97%)	167 (98%)	4 (2%)	0	100	100
1	B	171/177 (97%)	167 (98%)	4 (2%)	0	100	100
1	C	171/177 (97%)	168 (98%)	3 (2%)	0	100	100
1	D	171/177 (97%)	166 (97%)	4 (2%)	1 (1%)	25	47
1	E	171/177 (97%)	162 (95%)	8 (5%)	1 (1%)	25	47

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	F	171/177 (97%)	167 (98%)	4 (2%)	0	100	100
All	All	1026/1062 (97%)	997 (97%)	27 (3%)	2 (0%)	47	71

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	25	GLN
1	E	25	GLN

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	153/156 (98%)	149 (97%)	4 (3%)	46	72
1	B	153/156 (98%)	144 (94%)	9 (6%)	19	39
1	C	153/156 (98%)	143 (94%)	10 (6%)	17	34
1	D	153/156 (98%)	144 (94%)	9 (6%)	19	39
1	E	153/156 (98%)	140 (92%)	13 (8%)	10	21
1	F	153/156 (98%)	143 (94%)	10 (6%)	17	34
All	All	918/936 (98%)	863 (94%)	55 (6%)	19	39

All (55) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	54	ARG
1	A	79	LEU
1	A	104	GLU
1	A	142	ASN
1	B	2	ILE
1	B	7	LEU
1	B	49	GLU
1	B	50	LYS
1	B	54	ARG

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Mol	Chain	Res	Type
1	B	79	LEU
1	B	104	GLU
1	B	134	GLN
1	B	142	ASN
1	C	28	GLU
1	C	49	GLU
1	C	54	ARG
1	C	79	LEU
1	C	96	LYS
1	C	112	CYS
1	C	134	GLN
1	C	142	ASN
1	C	144	ASP
1	C	163	GLU
1	D	1	MET
1	D	49	GLU
1	D	54	ARG
1	D	79	LEU
1	D	104	GLU
1	D	134	GLN
1	D	144	ASP
1	D	160	LYS
1	D	167	ASN
1	E	49	GLU
1	E	79	LEU
1	E	104	GLU
1	E	130	LEU
1	E	134	GLN
1	E	136	GLU
1	E	142	ASN
1	E	145	GLU
1	E	146	PHE
1	E	148	GLU
1	E	150	ILE
1	E	163	GLU
1	E	173	LYS
1	F	25	GLN
1	F	49	GLU
1	F	54	ARG
1	F	79	LEU
1	F	107	ARG
1	F	113	SER

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Mol	Chain	Res	Type
1	F	134	GLN
1	F	142	ASN
1	F	144	ASP
1	F	167	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	25	GLN
1	A	140	ASN
1	A	142	ASN
1	B	82	GLN
1	B	134	GLN
1	B	140	ASN
1	B	142	ASN
1	B	167	ASN
1	C	82	GLN
1	C	134	GLN
1	D	82	GLN
1	D	134	GLN
1	D	140	ASN
1	D	142	ASN
1	D	167	ASN
1	E	25	GLN
1	E	110	GLN
1	E	134	GLN
1	E	167	ASN
1	F	25	GLN
1	F	134	GLN
1	F	167	ASN

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 6 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 [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	173/177 (97%)	-0.36	0 100 100	23, 36, 53, 68	0
1	B	173/177 (97%)	-0.35	2 (1%) 79 76	26, 37, 54, 67	0
1	C	173/177 (97%)	-0.38	0 100 100	24, 35, 52, 66	0
1	D	173/177 (97%)	-0.25	3 (1%) 70 66	25, 38, 54, 72	0
1	E	173/177 (97%)	-0.17	4 (2%) 60 54	25, 36, 57, 86	0
1	F	173/177 (97%)	-0.33	1 (0%) 89 88	25, 36, 52, 63	0
All	All	1038/1062 (97%)	-0.31	10 (0%) 82 80	23, 36, 54, 86	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	163	GLU	4.1
1	E	170	GLY	4.0
1	B	50	LYS	3.1
1	F	163	GLU	2.6
1	D	163	GLU	2.3
1	E	26	GLU	2.2
1	B	171	ARG	2.1
1	D	52	GLY	2.1
1	E	172	LEU	2.1
1	D	49	GLU	2.1

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 monosaccharides 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	NI	E	605	1/1	0.86	0.10	85,85,85,85	0
2	NI	B	602	1/1	0.87	0.12	76,76,76,76	0
2	NI	D	604	1/1	0.88	0.08	76,76,76,76	0
2	NI	A	601	1/1	0.93	0.11	83,83,83,83	0
2	NI	F	606	1/1	0.93	0.12	77,77,77,77	0
2	NI	C	603	1/1	0.95	0.07	76,76,76,76	0

6.5 Other polymers [i](#)

There are no such residues in this entry.