



Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 02:04 pm BST

PDB ID : 2DLA
Title : Primase large subunit amino terminal domain from *Pyrococcus horikoshii*
Authors : Ito, N.
Deposited on : 2006-04-17
Resolution : 2.90 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Xtrriage (Phenix) : 1.13
EDS : 2.11
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) : 2.11

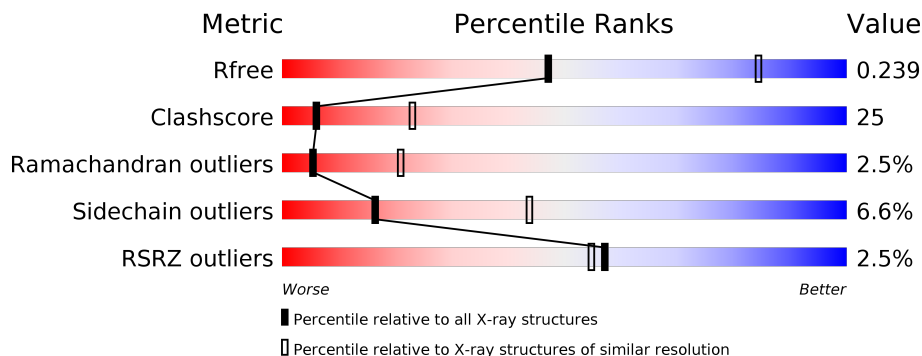
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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 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	222	
1	B	222	
1	C	222	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5289 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 397aa long hypothetical protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	222	Total 1842	C 1193	N 297	O 345	S 7	0	0	0
1	B	214	Total 1776	C 1152	N 284	O 333	S 7	0	0	0
1	C	200	Total 1654	C 1072	N 266	O 309	S 7	0	0	0

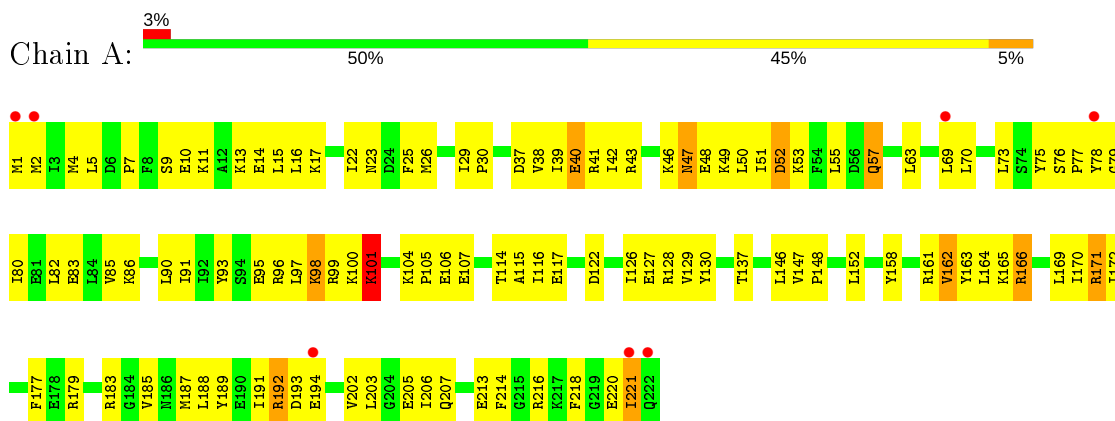
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	12	Total 12	O 12	0	0
2	B	2	Total 2	O 2	0	0
2	C	3	Total 3	O 3	0	0

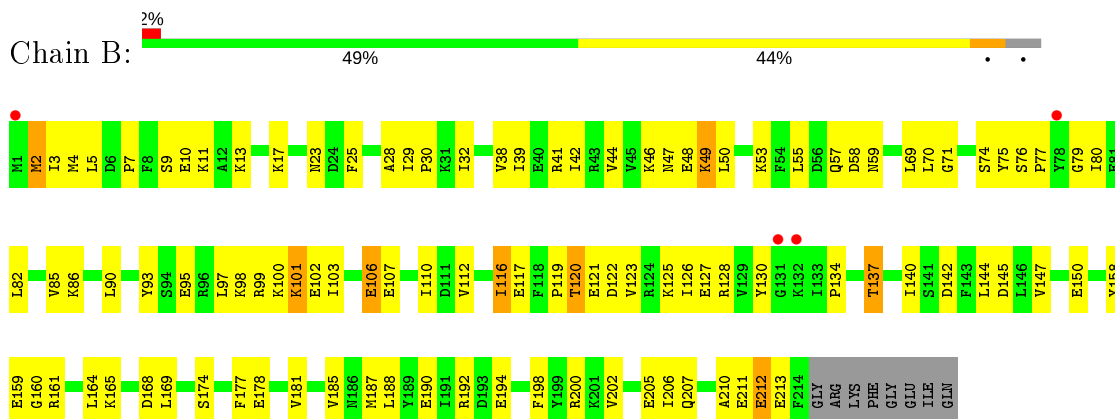
3 Residue-property plots

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.

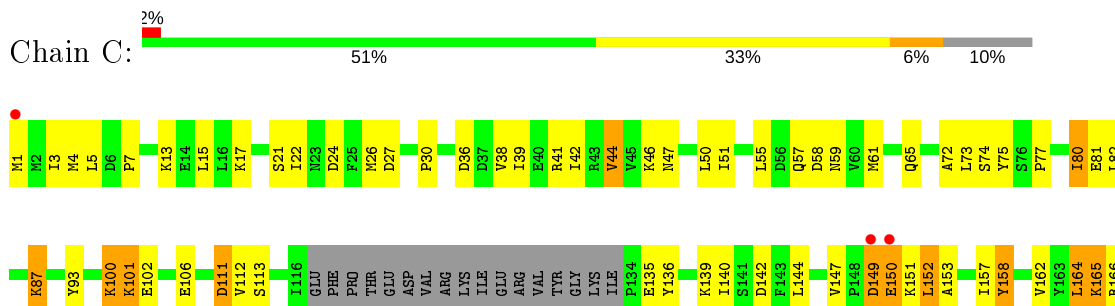
- Molecule 1: 397aa long hypothetical protein



- Molecule 1: 397aa long hypothetical protein



- Molecule 1: 397aa long hypothetical protein





4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	89.25Å 89.25Å 252.67Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.90 19.96 – 2.72	Depositor EDS
% Data completeness (in resolution range)	(Not available) (20.00-2.90) 99.0 (19.96-2.72)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.15 (at 2.71Å)	Xtrriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.248 , 0.309 0.240 , 0.239	Depositor DCC
R_{free} test set	2009 reflections (7.17%)	wwPDB-VP
Wilson B-factor (Å ²)	65.1	Xtrriage
Anisotropy	0.191	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 51.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	5289	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.78% 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	0.44	0/1875	0.64	0/2517
1	B	0.42	0/1808	0.65	0/2431
1	C	0.42	0/1682	0.60	0/2257
All	All	0.43	0/5365	0.63	0/7205

There are no bond length outliers.

There are no bond angle outliers.

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	1842	0	1877	99	0
1	B	1776	0	1811	88	0
1	C	1654	0	1691	80	1
2	A	12	0	0	0	0
2	B	2	0	0	0	0
2	C	3	0	0	0	0
All	All	5289	0	5379	263	1

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 25.

All (263) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:101:LYS:HD2	1:C:101:LYS:H	1.29	0.94
1:C:7:PRO:HG3	1:C:93:TYR:CZ	2.11	0.86
1:B:39:ILE:HG12	1:B:205:GLU:HB3	1.57	0.85
1:A:166:ARG:HH11	1:A:166:ARG:HG2	1.42	0.85
1:B:55:LEU:HD21	1:B:95:GLU:HG3	1.60	0.84
1:A:216:ARG:O	1:A:220:GLU:HB2	1.79	0.82
1:B:101:LYS:CD	1:B:101:LYS:H	1.93	0.82
1:A:221:ILE:HD11	1:B:17:LYS:HG2	1.65	0.78
1:B:112:VAL:HG21	1:B:169:LEU:HD11	1.66	0.77
1:A:220:GLU:HG2	1:A:221:ILE:H	1.50	0.76
1:B:5:LEU:HD12	1:B:107:GLU:HB3	1.66	0.76
1:A:39:ILE:HG12	1:A:205:GLU:HG3	1.68	0.74
1:C:165:LYS:HE2	1:C:166:ARG:H	1.53	0.74
1:A:171:ARG:HB3	1:A:171:ARG:HH11	1.52	0.72
1:B:120:THR:HA	1:B:123:VAL:HG23	1.72	0.71
1:B:80:ILE:HD12	1:B:80:ILE:N	2.05	0.71
1:C:21:SER:HB3	1:C:24:ASP:OD2	1.93	0.69
1:B:55:LEU:HD21	1:B:95:GLU:CG	2.22	0.69
1:A:104:LYS:HB2	1:A:107:GLU:HG3	1.75	0.68
1:A:80:ILE:HA	1:A:83:GLU:HB2	1.74	0.68
1:B:101:LYS:HD3	1:B:101:LYS:N	2.08	0.68
1:B:147:VAL:HB	1:B:150:GLU:HG2	1.76	0.68
1:C:164:LEU:N	1:C:164:LEU:HD23	2.08	0.68
1:B:101:LYS:H	1:B:101:LYS:HD3	1.60	0.67
1:B:101:LYS:CD	1:B:101:LYS:N	2.57	0.66
1:A:221:ILE:CD1	1:B:17:LYS:HG2	2.26	0.66
1:C:147:VAL:HG13	1:C:149:ASP:HB3	1.78	0.66
1:B:174:SER:O	1:B:178:GLU:HG3	1.96	0.66
1:A:39:ILE:HD11	1:A:205:GLU:HB3	1.78	0.66
1:B:9:SER:HB2	1:B:106:GLU:O	1.95	0.65
1:C:75:TYR:C	1:C:77:PRO:HD3	2.16	0.65
1:A:57:GLN:NE2	1:A:96:ARG:HH22	1.95	0.64
1:C:39:ILE:HG13	1:C:205:GLU:HB3	1.79	0.64
1:A:179:ARG:HH11	1:A:183:ARG:NH2	1.95	0.64
1:C:15:LEU:CD2	1:C:59:ASN:HD22	2.10	0.64
1:C:144:LEU:HD21	1:C:152:LEU:HD22	1.77	0.64
1:B:4:MET:CE	1:B:11:LYS:HE3	2.28	0.64
1:B:2:MET:HB3	1:B:11:LYS:NZ	2.13	0.63
1:A:166:ARG:NH1	1:A:166:ARG:HG2	2.07	0.62
1:A:47:ASN:ND2	1:A:49:LYS:H	1.96	0.62
1:A:7:PRO:HD3	1:A:93:TYR:CZ	2.34	0.62
1:A:2:MET:HG3	1:A:11:LYS:NZ	2.15	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:46:LYS:HZ1	1:A:213:GLU:HG3	1.64	0.61
1:B:130:TYR:HD2	1:B:165:LYS:HG3	1.67	0.60
1:C:106:GLU:CD	1:C:106:GLU:H	2.05	0.60
1:C:151:LYS:C	1:C:153:ALA:H	2.05	0.60
1:C:15:LEU:HD21	1:C:59:ASN:HD22	1.66	0.60
1:B:181:VAL:O	1:B:185:VAL:HG23	2.01	0.60
1:C:100:LYS:HD3	1:C:101:LYS:HD2	1.84	0.60
1:B:120:THR:CA	1:B:123:VAL:HG23	2.31	0.59
1:B:74:SER:OG	1:B:192:ARG:HB3	2.02	0.59
1:B:7:PRO:HG3	1:B:93:TYR:CZ	2.38	0.59
1:A:26:MET:HE2	1:A:188:LEU:HD21	1.83	0.59
1:A:76:SER:N	1:A:77:PRO:HD3	2.17	0.59
1:B:4:MET:HE2	1:B:11:LYS:HE3	1.84	0.59
1:A:4:MET:HE3	1:A:63:LEU:HD11	1.85	0.58
1:A:55:LEU:HD13	1:A:99:ARG:NH1	2.17	0.58
1:A:130:TYR:HB3	1:A:165:LYS:NZ	2.18	0.58
1:C:77:PRO:HG2	1:C:189:TYR:CD1	2.38	0.58
1:A:26:MET:CE	1:A:188:LEU:HD21	2.34	0.58
1:A:86:LYS:O	1:A:90:LEU:HB2	2.04	0.58
1:B:122:ASP:O	1:B:126:ILE:HG13	2.04	0.58
1:B:76:SER:N	1:B:77:PRO:HD3	2.19	0.58
1:B:101:LYS:H	1:B:101:LYS:HD2	1.69	0.58
1:B:10:GLU:HG2	1:B:106:GLU:HB2	1.85	0.58
1:B:70:LEU:HD11	1:B:181:VAL:HG13	1.86	0.57
1:C:164:LEU:HD23	1:C:164:LEU:H	1.69	0.57
1:C:47:ASN:O	1:C:50:LEU:HB2	2.03	0.57
1:B:116:ILE:H	1:B:116:ILE:HD12	1.68	0.57
1:C:202:VAL:O	1:C:206:ILE:HG13	2.04	0.57
1:A:99:ARG:HG3	1:A:99:ARG:O	2.05	0.57
1:A:179:ARG:HH11	1:A:183:ARG:HH21	1.52	0.57
1:C:164:LEU:CD1	1:C:172:ILE:HD12	2.34	0.57
1:A:37:ASP:O	1:A:40:GLU:HB2	2.04	0.57
1:C:22:ILE:O	1:C:26:MET:HG2	2.05	0.56
1:C:75:TYR:O	1:C:77:PRO:HD3	2.05	0.56
1:A:137:THR:OG1	1:A:161:ARG:NH1	2.39	0.56
1:A:98:LYS:HB3	1:A:170:ILE:HD11	1.87	0.56
1:A:7:PRO:HD3	1:A:93:TYR:CE2	2.40	0.56
1:C:73:LEU:HD13	1:C:82:LEU:HD12	1.87	0.56
1:C:3:ILE:HD13	1:C:55:LEU:O	2.07	0.55
1:C:213:GLU:O	1:C:216:ARG:HG2	2.07	0.55
1:A:221:ILE:O	1:A:221:ILE:HG23	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:191:ILE:O	1:C:191:ILE:HG22	2.07	0.55
1:A:25:PHE:CE2	1:A:29:ILE:HD11	2.42	0.55
1:B:75:TYR:HA	1:B:192:ARG:HH11	1.71	0.55
1:A:183:ARG:O	1:A:187:MET:HG3	2.08	0.54
1:B:210:ALA:O	1:B:212:GLU:N	2.40	0.54
1:B:116:ILE:N	1:B:116:ILE:HD12	2.21	0.54
1:C:152:LEU:HD23	1:C:152:LEU:H	1.72	0.54
1:B:47:ASN:HB3	1:B:49:LYS:CD	2.37	0.54
1:A:2:MET:HG3	1:A:11:LYS:HZ3	1.73	0.54
1:B:80:ILE:CD1	1:B:80:ILE:N	2.70	0.54
1:B:95:GLU:HG3	1:B:99:ARG:HE	1.73	0.54
1:A:70:LEU:HD22	1:A:185:VAL:HG23	1.89	0.54
1:A:98:LYS:HB2	1:A:166:ARG:NH1	2.23	0.54
1:B:39:ILE:CG1	1:B:205:GLU:HB3	2.33	0.53
1:C:112:VAL:HG12	1:C:113:SER:H	1.74	0.53
1:C:5:LEU:HD11	1:C:100:LYS:HG3	1.90	0.53
1:A:52:ASP:HA	1:A:55:LEU:HG	1.91	0.53
1:C:149:ASP:O	1:C:151:LYS:N	2.42	0.53
1:C:112:VAL:HG12	1:C:113:SER:N	2.23	0.53
1:A:90:LEU:HD13	1:A:177:PHE:CD2	2.44	0.52
1:C:101:LYS:N	1:C:101:LYS:HD2	2.12	0.52
1:A:127:GLU:C	1:A:129:VAL:H	2.12	0.52
1:C:164:LEU:N	1:C:164:LEU:CD2	2.72	0.52
1:A:41:ARG:HH12	1:A:57:GLN:HE22	1.55	0.52
1:A:46:LYS:HE3	1:A:216:ARG:HH21	1.75	0.52
1:B:198:PHE:O	1:B:202:VAL:HG23	2.09	0.52
1:B:47:ASN:HB3	1:B:49:LYS:HD2	1.92	0.52
1:A:90:LEU:HD13	1:A:177:PHE:HD2	1.76	0.51
1:A:105:PRO:HG2	1:A:106:GLU:OE2	2.10	0.51
1:A:77:PRO:O	1:A:82:LEU:HD22	2.10	0.51
1:A:22:ILE:O	1:A:26:MET:HB2	2.10	0.51
1:B:4:MET:HG2	1:B:59:ASN:OD1	2.11	0.51
1:C:46:LYS:HG2	1:C:47:ASN:ND2	2.24	0.51
1:A:46:LYS:HZ1	1:A:213:GLU:CG	2.24	0.51
1:C:151:LYS:O	1:C:153:ALA:N	2.44	0.51
1:A:189:TYR:O	1:A:192:ARG:HG3	2.11	0.51
1:A:218:PHE:HZ	1:B:187:MET:HB3	1.76	0.51
1:B:119:PRO:C	1:B:121:GLU:H	2.14	0.51
1:A:203:LEU:HG	1:A:207:GLN:NE2	2.25	0.51
1:A:146:LEU:C	1:A:148:PRO:HD3	2.32	0.51
1:A:39:ILE:O	1:A:43:ARG:HG3	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:28:ALA:O	1:B:32:ILE:HG13	2.11	0.50
1:C:192:ARG:O	1:C:195:LEU:HB2	2.10	0.50
1:B:202:VAL:O	1:B:206:ILE:HG13	2.12	0.50
1:B:41:ARG:O	1:B:44:VAL:HG22	2.11	0.50
1:A:29:ILE:N	1:A:30:PRO:HD2	2.26	0.50
1:A:46:LYS:NZ	1:A:213:GLU:HG3	2.27	0.50
1:C:165:LYS:HE2	1:C:166:ARG:N	2.24	0.50
1:B:119:PRO:HB2	1:B:121:GLU:OE2	2.12	0.49
1:B:142:ASP:O	1:B:145:ASP:HB3	2.13	0.49
1:B:75:TYR:CD2	1:B:192:ARG:HD2	2.48	0.49
1:C:44:VAL:HG22	1:C:51:ILE:HG12	1.94	0.49
1:C:7:PRO:HG3	1:C:93:TYR:OH	2.13	0.49
1:C:140:ILE:HG12	1:C:162:VAL:CG2	2.43	0.48
1:C:152:LEU:HD23	1:C:152:LEU:N	2.28	0.48
1:A:46:LYS:HZ1	1:A:216:ARG:HH21	1.62	0.48
1:C:101:LYS:CD	1:C:101:LYS:H	2.03	0.48
1:A:116:ILE:HG22	1:A:117:GLU:HG3	1.95	0.48
1:B:119:PRO:O	1:B:121:GLU:N	2.47	0.48
1:B:125:LYS:HD3	1:B:125:LYS:N	2.29	0.48
1:C:147:VAL:C	1:C:149:ASP:N	2.67	0.47
1:C:140:ILE:HD11	1:C:157:ILE:CG2	2.44	0.47
1:C:183:ARG:O	1:C:187:MET:HE2	2.14	0.47
1:A:51:ILE:HG21	1:A:95:GLU:HG2	1.97	0.47
1:B:137:THR:OG1	1:B:161:ARG:HD3	2.13	0.47
1:C:190:GLU:C	1:C:191:ILE:HG13	2.34	0.47
1:A:14:GLU:HA	1:A:17:LYS:HB3	1.95	0.47
1:A:158:TYR:HB3	1:A:163:TYR:CE1	2.50	0.47
1:B:101:LYS:HG2	1:B:102:GLU:H	1.78	0.47
1:C:61:MET:O	1:C:65:GLN:HG3	2.14	0.47
1:A:158:TYR:HD2	1:A:163:TYR:HH	1.62	0.47
1:C:13:LYS:O	1:C:17:LYS:HD3	2.14	0.47
1:C:151:LYS:C	1:C:153:ALA:N	2.68	0.47
1:A:163:TYR:O	1:A:164:LEU:HD23	2.14	0.47
1:A:5:LEU:HD22	1:A:97:LEU:HA	1.96	0.47
1:A:5:LEU:HD22	1:A:97:LEU:HD12	1.97	0.47
1:A:9:SER:O	1:A:13:LYS:HG2	2.15	0.47
1:B:177:PHE:O	1:B:181:VAL:HG23	2.14	0.47
1:B:47:ASN:C	1:B:49:LYS:H	2.18	0.47
1:C:74:SER:O	1:C:192:ARG:HD3	2.15	0.47
1:A:115:ALA:HB2	1:A:137:THR:HG23	1.97	0.47
1:B:207:GLN:O	1:B:210:ALA:HB3	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:47:ASN:HB2	1:C:50:LEU:HD12	1.97	0.47
1:B:25:PHE:CE2	1:B:29:ILE:HD11	2.50	0.47
1:B:30:PRO:HG3	1:B:198:PHE:CD2	2.50	0.47
1:B:49:LYS:H	1:B:49:LYS:HD2	1.79	0.47
1:B:5:LEU:HD21	1:B:100:LYS:HG3	1.96	0.47
1:B:116:ILE:HG22	1:B:116:ILE:O	2.14	0.46
1:C:147:VAL:CG1	1:C:149:ASP:HB3	2.44	0.46
1:B:98:LYS:HB2	1:B:98:LYS:HE3	1.67	0.46
1:C:135:GLU:HG3	1:C:136:TYR:CD1	2.51	0.46
1:A:39:ILE:HG12	1:A:205:GLU:CG	2.43	0.46
1:C:80:ILE:HD12	1:C:80:ILE:HA	1.71	0.46
1:A:202:VAL:O	1:A:206:ILE:HG13	2.15	0.46
1:C:157:ILE:HG22	1:C:158:TYR:N	2.31	0.46
1:A:97:LEU:C	1:A:99:ARG:H	2.19	0.46
1:C:149:ASP:O	1:C:150:GLU:C	2.54	0.46
1:C:27:ASP:O	1:C:30:PRO:HD2	2.16	0.46
1:C:81:GLU:CD	1:C:81:GLU:H	2.19	0.46
1:A:75:TYR:CE2	1:A:192:ARG:HB2	2.51	0.46
1:B:38:VAL:HG13	1:B:69:LEU:HG	1.96	0.46
1:A:47:ASN:C	1:A:47:ASN:HD22	2.20	0.46
1:B:116:ILE:HG21	1:B:134:PRO:O	2.16	0.46
1:B:10:GLU:HA	1:B:13:LYS:HB2	1.98	0.46
1:C:177:PHE:O	1:C:181:VAL:HG23	2.16	0.45
1:C:165:LYS:O	1:C:168:ASP:HB2	2.16	0.45
1:C:44:VAL:HG22	1:C:51:ILE:CG1	2.46	0.45
1:A:214:PHE:CZ	1:B:190:GLU:HG3	2.51	0.45
1:A:42:ILE:HD11	1:A:69:LEU:HD22	1.99	0.45
1:A:47:ASN:HD22	1:A:48:GLU:N	2.14	0.45
1:C:38:VAL:O	1:C:41:ARG:HB3	2.16	0.45
1:A:146:LEU:HD23	1:A:179:ARG:NH2	2.32	0.45
1:B:164:LEU:HD22	1:B:168:ASP:HB3	1.98	0.45
1:A:128:ARG:O	1:A:128:ARG:HG2	2.17	0.45
1:A:39:ILE:CD1	1:A:205:GLU:HB3	2.45	0.45
1:B:90:LEU:HD11	1:B:174:SER:HB3	1.99	0.45
1:C:100:LYS:O	1:C:166:ARG:NH1	2.48	0.45
1:C:186:ASN:O	1:C:189:TYR:HB3	2.16	0.45
1:B:100:LYS:HD3	1:B:101:LYS:HD3	2.00	0.44
1:C:190:GLU:O	1:C:191:ILE:HG13	2.17	0.44
1:C:57:GLN:HG2	1:C:58:ASP:N	2.33	0.44
1:A:80:ILE:HA	1:A:83:GLU:CB	2.46	0.44
1:B:2:MET:HB3	1:B:11:LYS:HZ1	1.80	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:57:GLN:HE21	1:A:96:ARG:HH22	1.62	0.44
1:B:158:TYR:CD2	1:B:159:GLU:HG3	2.52	0.44
1:B:57:GLN:HG2	1:B:58:ASP:O	2.18	0.44
1:A:46:LYS:CE	1:A:216:ARG:HH21	2.31	0.44
1:C:111:ASP:C	1:C:111:ASP:OD1	2.56	0.44
1:B:39:ILE:HD11	1:B:205:GLU:HB2	2.01	0.43
1:C:72:ALA:O	1:C:203:LEU:HD13	2.18	0.43
1:C:15:LEU:HD22	1:C:59:ASN:HD22	1.80	0.43
1:B:50:LEU:O	1:B:53:LYS:HG2	2.17	0.43
1:C:136:TYR:HE1	1:C:166:ARG:HB2	1.83	0.43
1:A:47:ASN:ND2	1:A:47:ASN:C	2.72	0.43
1:B:46:LYS:HD2	1:B:213:GLU:OE1	2.19	0.43
1:A:164:LEU:HD13	1:A:172:ILE:HD12	2.01	0.43
1:C:139:LYS:HB2	1:C:142:ASP:OD2	2.19	0.43
1:A:38:VAL:HG12	1:A:69:LEU:HD21	2.00	0.43
1:C:87:LYS:HA	1:C:87:LYS:CE	2.49	0.43
1:A:101:LYS:N	1:A:101:LYS:HD2	2.34	0.43
1:B:5:LEU:HD12	1:B:107:GLU:CB	2.45	0.43
1:A:5:LEU:HD11	1:A:100:LYS:HE2	2.00	0.42
1:A:97:LEU:O	1:A:99:ARG:N	2.52	0.42
1:B:71:GLY:HA2	1:B:188:LEU:HD13	2.01	0.42
1:B:101:LYS:HG2	1:B:102:GLU:N	2.34	0.42
1:B:123:VAL:O	1:B:127:GLU:HG3	2.19	0.42
1:B:140:ILE:HG13	1:B:160:GLY:O	2.19	0.42
1:C:140:ILE:HD11	1:C:157:ILE:HG23	2.00	0.42
1:A:4:MET:HE3	1:A:63:LEU:HD21	2.00	0.42
1:A:91:ILE:O	1:A:95:GLU:HB2	2.20	0.42
1:C:44:VAL:O	1:C:44:VAL:HG13	2.19	0.42
1:B:86:LYS:O	1:B:90:LEU:HB2	2.19	0.42
1:B:4:MET:HE1	1:B:11:LYS:HE3	1.99	0.42
1:C:3:ILE:HG22	1:C:4:MET:N	2.34	0.42
1:A:137:THR:HA	1:A:162:VAL:O	2.20	0.42
1:A:70:LEU:HD22	1:A:185:VAL:CG2	2.50	0.42
1:C:189:TYR:O	1:C:192:ARG:HG3	2.20	0.41
1:A:46:LYS:NZ	1:A:216:ARG:HH21	2.18	0.41
1:A:50:LEU:HD23	1:A:53:LYS:NZ	2.35	0.41
1:B:125:LYS:HD3	1:B:128:ARG:HH12	1.85	0.41
1:C:140:ILE:HG12	1:C:162:VAL:HG22	2.01	0.41
1:A:152:LEU:HA	1:A:152:LEU:HD23	1.91	0.41
1:A:15:LEU:HD23	1:A:63:LEU:CD1	2.51	0.41
1:C:195:LEU:HA	1:C:196:PRO:HD3	1.94	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:211:GLU:HG3	1:C:212:GLU:N	2.36	0.41
1:B:110:ILE:HG13	1:B:112:VAL:HG23	2.03	0.41
1:A:191:ILE:O	1:A:193:ASP:N	2.54	0.41
1:B:25:PHE:O	1:B:29:ILE:HG13	2.20	0.41
1:C:42:ILE:C	1:C:44:VAL:H	2.24	0.41
1:A:130:TYR:HB3	1:A:165:LYS:HZ1	1.85	0.41
1:A:46:LYS:HE3	1:A:216:ARG:NH2	2.36	0.41
1:B:97:LEU:HD22	1:B:103:ILE:HD13	2.03	0.41
1:B:82:LEU:HG	1:B:86:LYS:NZ	2.35	0.41
1:C:100:LYS:HD2	1:C:102:GLU:O	2.21	0.41
1:A:122:ASP:O	1:A:126:ILE:HG12	2.21	0.40
1:A:75:TYR:HB2	1:A:203:LEU:HD22	2.03	0.40
1:B:42:ILE:HD11	1:B:85:VAL:HG22	2.03	0.40
1:A:147:VAL:N	1:A:148:PRO:HD3	2.35	0.40
1:A:73:LEU:HD11	1:A:85:VAL:HG21	2.03	0.40
1:B:144:LEU:HA	1:B:144:LEU:HD23	1.83	0.40

All (1) 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)
1:C:58:ASP:OD2	1:C:58:ASP:OD2[8_664]	1.95	0.25

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	220/222 (99%)	194 (88%)	22 (10%)	4 (2%)	8	29
1	B	212/222 (96%)	192 (91%)	13 (6%)	7 (3%)	4	15
1	C	196/222 (88%)	171 (87%)	20 (10%)	5 (3%)	5	20
All	All	628/666 (94%)	557 (89%)	55 (9%)	16 (2%)	5	21

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	221	ILE
1	C	149	ASP
1	C	150	GLU
1	A	79	GLY
1	A	98	LYS
1	B	120	THR
1	B	212	GLU
1	C	152	LEU
1	C	191	ILE
1	A	101	LYS
1	B	48	GLU
1	B	79	GLY
1	C	158	TYR
1	B	3	ILE
1	B	211	GLU
1	B	116	ILE

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	201/201 (100%)	184 (92%)	17 (8%)	10	31
1	B	195/201 (97%)	186 (95%)	9 (5%)	27	60
1	C	181/201 (90%)	169 (93%)	12 (7%)	16	44
All	All	577/603 (96%)	539 (93%)	38 (7%)	16	44

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	10	GLU
1	A	16	LEU
1	A	23	ASN
1	A	40	GLU

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Mol	Chain	Res	Type
1	A	47	ASN
1	A	52	ASP
1	A	57	GLN
1	A	78	TYR
1	A	101	LYS
1	A	114	THR
1	A	162	VAL
1	A	166	ARG
1	A	169	LEU
1	A	171	ARG
1	A	192	ARG
1	A	194	GLU
1	B	2	MET
1	B	23	ASN
1	B	49	LYS
1	B	101	LYS
1	B	106	GLU
1	B	117	GLU
1	B	137	THR
1	B	194	GLU
1	B	200	ARG
1	C	1	MET
1	C	36	ASP
1	C	44	VAL
1	C	80	ILE
1	C	87	LYS
1	C	100	LYS
1	C	101	LYS
1	C	111	ASP
1	C	164	LEU
1	C	165	LYS
1	C	182	GLU
1	C	195	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	23	ASN
1	A	47	ASN
1	A	57	GLN
1	A	59	ASN
1	A	65	GLN

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Mol	Chain	Res	Type
1	A	180	ASN
1	B	207	GLN
1	C	47	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 carbohydrates 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 [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	222/222 (100%)	-0.14	7 (3%) 47 43	27, 58, 83, 101	0
1	B	214/222 (96%)	-0.20	4 (1%) 66 65	28, 53, 89, 99	0
1	C	200/222 (90%)	-0.07	5 (2%) 57 55	34, 58, 87, 109	0
All	All	636/666 (95%)	-0.14	16 (2%) 57 55	27, 56, 87, 109	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	222	GLN	9.0
1	C	149	ASP	4.6
1	C	216	ARG	4.1
1	C	150	GLU	3.5
1	C	217	LYS	3.2
1	B	131	GLY	3.1
1	B	132	LYS	3.0
1	A	78	TYR	2.7
1	C	1	MET	2.7
1	A	194	GLU	2.5
1	A	1	MET	2.5
1	B	1	MET	2.4
1	A	2	MET	2.3
1	A	221	ILE	2.2
1	A	69	LEU	2.1
1	B	78	TYR	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

There are no carbohydrates in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.