



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:26 AM UTC

PDB ID : 6G6N / pdb_00006g6n
Title : Crystal structure of the computationally designed Tako8 protein in C2
Authors : Noguchi, H.; Addy, C.; Simoncini, D.; Van Meervelt, L.; Schiex, T.; Zhang, K.Y.J.; Tame, J.R.H.; Voet, A.R.D.
Deposited on : 2018-04-01
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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

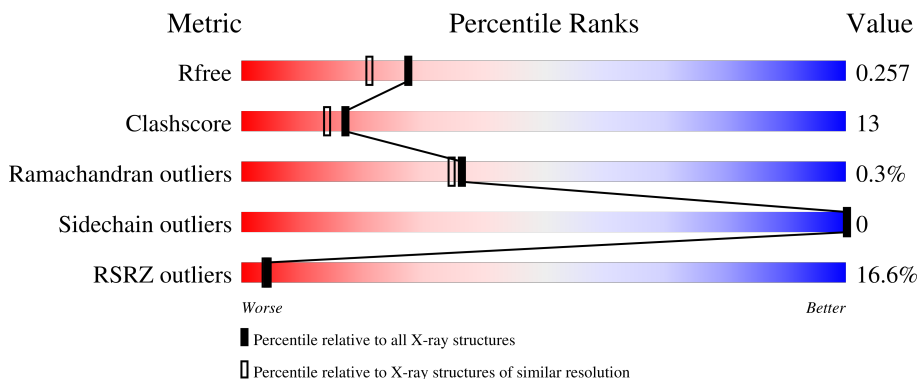
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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}	180053	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (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 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	325	18% 76% 22% .
1	B	325	14% 75% 23% .
1	C	325	16% 71% 26% .

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7475 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 Tako8.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	A	319	Total	C	N	O	0	0	0
			2425	1484	431	510			
1	B	318	Total	C	N	O	0	0	0
			2416	1479	429	508			
1	C	318	Total	C	N	O	0	0	0
			2416	1479	429	508			

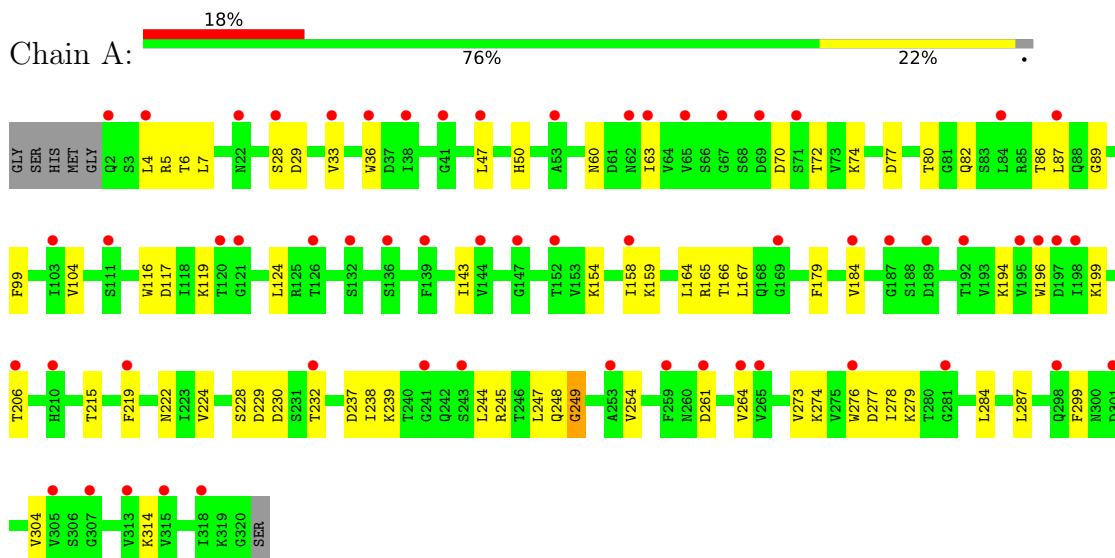
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	70	Total	O	0	0
			70	70		
2	B	76	Total	O	0	0
			76	76		
2	C	72	Total	O	0	0
			72	72		

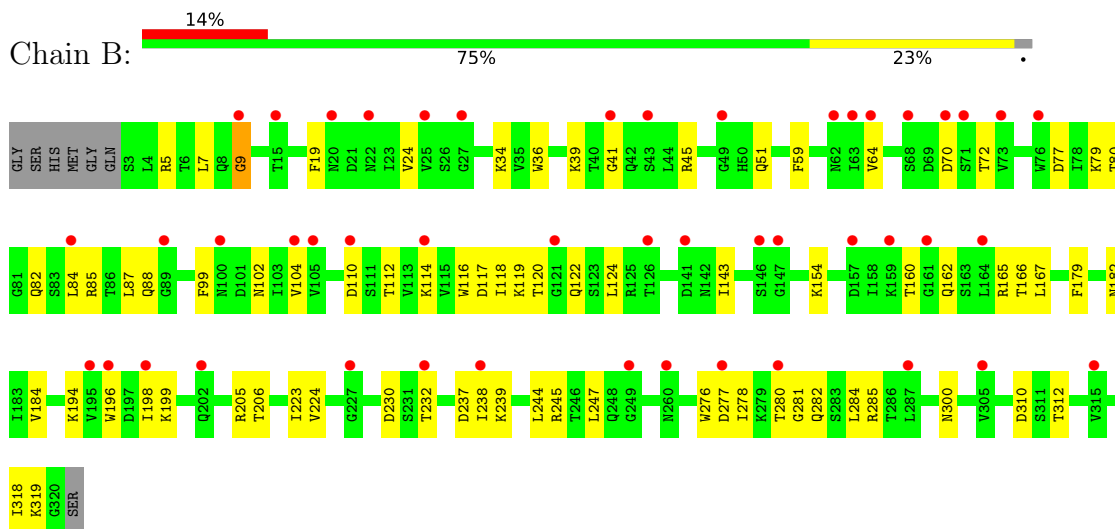
3 Residue-property plots

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: Tako8

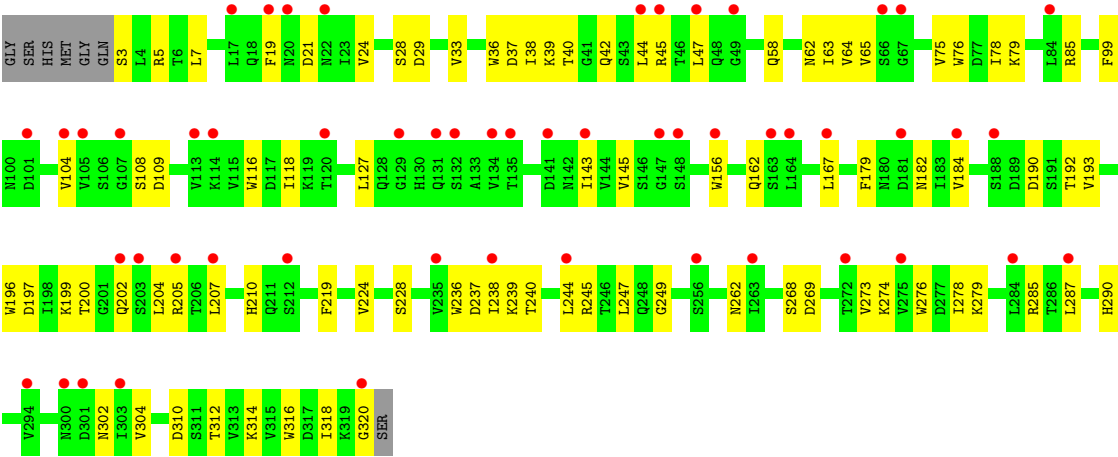


• Molecule 1: Tako8



• Molecule 1: Tako8





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	93.55Å 132.30Å 81.02Å 90.00° 125.26° 90.00°	Depositor
Resolution (Å)	46.77 – 2.00 46.77 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.9 (46.77-2.00) 99.9 (46.77-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	111.59 (at 2.00Å)	Xtriage
Refinement program	PHENIX (1.13_2998: ???)	Depositor
R, R_{free}	0.243 , 0.266 0.225 , 0.257	Depositor DCC
R_{free} test set	2485 reflections (4.58%)	wwPDB-VP
Wilson B-factor (Å ²)	45.4	Xtriage
Anisotropy	0.027	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 237.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.469 for -h,-h-2*k,1/2*h-1/2*k 0.468 for -h,h+2*k,1/2*h+1/2*k 0.468 for -1/2*h+1/2*k-l,3/2*h+1/2*k+l,1/2*h-1/2*k 0.469 for 1/2*h-1/2*k+l,-1/2*h+1/2*k+l,-h-1/2*h+1/2*k+l,1/2*h+1/2*k-l,-h-1/2*h-1/2*k-l,-3/2*h+1/2*k-l,1/2*h+1/2*k 0.469 for 1/2*h+1/2*k+l,3/2*h-1/2*k+l,-l 0.469 for 1/2*h-1/2*k+l,-3/2*h-1/2*k-l,-l 0.470 for -1/2*h+1/2*k-l,1/2*h-1/2*k-l,-1/2*h-1/2*k 0.469 for -1/2*h-1/2*k-l,-1/2*h-1/2*k+l,-1/2*h+1/2*k 0.469 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7475	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

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.

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.80% of the height of the origin peak. No significant pseudotranslation is detected.*

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	0.13	0/2456	0.38	0/3339
1	B	0.13	0/2447	0.39	0/3327
1	C	0.16	0/2447	0.40	0/3327
All	All	0.14	0/7350	0.39	0/9993

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	2425	0	2368	54	0
1	B	2416	0	2360	58	0
1	C	2416	0	2360	78	0
2	A	70	0	0	0	0
2	B	76	0	0	3	0
2	C	72	0	0	3	0
All	All	7475	0	7088	180	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 13.

All (180) 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:B:119:LYS:HE3	1:C:62:ASN:HB3	1.56	0.86
1:C:62:ASN:HA	1:C:79:LYS:HZ1	1.43	0.83
1:A:224:VAL:HG23	1:A:238:ILE:HG22	1.60	0.82
1:B:88:GLN:O	1:B:116:TRP:HH2	1.64	0.80
1:B:277:ASP:HB2	1:B:284:LEU:HD21	1.65	0.77
1:C:78:ILE:HG22	1:C:79:LYS:HE3	1.64	0.77
1:C:238:ILE:HG22	1:C:239:LYS:HD2	1.71	0.73
1:B:179:PHE:HB2	1:B:184:VAL:HG12	1.72	0.71
1:B:277:ASP:HB3	1:B:280:THR:HB	1.74	0.68
1:C:37:ASP:HB2	1:C:44:LEU:HD21	1.78	0.66
1:C:237:ASP:HB3	1:C:240:THR:HB	1.78	0.66
1:C:99:PHE:HB3	1:C:104:VAL:HG23	1.77	0.65
1:A:60:ASN:HD21	1:A:63:ILE:HD12	1.62	0.65
1:A:165:ARG:NH2	1:A:199:LYS:O	2.30	0.65
1:A:222:ASN:HB3	1:C:39:LYS:HD3	1.79	0.65
1:C:143:ILE:HD12	1:C:143:ILE:C	2.22	0.65
1:A:77:ASP:HB3	1:A:80:THR:HG22	1.78	0.64
1:C:62:ASN:HA	1:C:79:LYS:NZ	2.12	0.64
1:C:304:VAL:HB	1:C:316:TRP:HB2	1.78	0.63
1:A:117:ASP:HB2	1:A:124:LEU:HD11	1.79	0.63
1:C:219:PHE:HB2	1:C:238:ILE:HD11	1.80	0.62
1:B:143:ILE:HD12	1:B:198:ILE:HD11	1.83	0.61
1:A:80:THR:HG23	1:A:82:GLN:H	1.66	0.60
1:A:167:LEU:HD21	1:A:184:VAL:HG11	1.83	0.60
1:B:85:ARG:HH22	1:C:85:ARG:NH2	2.00	0.59
1:C:167:LEU:HD23	1:C:196:TRP:CG	2.37	0.59
1:C:58:GLN:NE2	2:C:401:HOH:O	2.33	0.59
1:B:99:PHE:HB2	1:B:104:VAL:HG12	1.84	0.59
1:A:4:LEU:HB3	1:A:5:ARG:HH11	1.67	0.59
1:A:165:ARG:NH2	1:B:165:ARG:HH22	2.01	0.59
1:A:74:LYS:NZ	1:A:86:THR:OG1	2.36	0.58
1:B:114:LYS:HB3	1:B:116:TRP:HE1	1.69	0.58
1:C:33:VAL:HB	1:C:47:LEU:HB2	1.86	0.58
1:A:247:LEU:HD23	1:A:276:TRP:CG	2.40	0.57
1:B:19:PHE:HB2	1:B:24:VAL:HG12	1.86	0.57
1:C:200:THR:HG23	1:C:202:GLN:H	1.70	0.57
1:C:62:ASN:CA	1:C:79:LYS:HZ1	2.17	0.57
1:A:99:PHE:HB2	1:A:104:VAL:HG12	1.86	0.57
1:B:194:LYS:NZ	1:B:206:THR:OG1	2.38	0.56
1:B:154:LYS:NZ	1:B:166:THR:OG1	2.26	0.56
1:C:104:VAL:HG12	1:C:116:TRP:HB2	1.88	0.56
1:A:274:LYS:HB3	1:A:276:TRP:NE1	2.21	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:3:SER:N	2:C:403:HOH:O	2.39	0.56
1:C:37:ASP:HB3	1:C:40:THR:HG1	1.72	0.56
1:C:247:LEU:HD23	1:C:276:TRP:CG	2.41	0.55
1:C:285:ARG:HD3	1:C:320:GLY:C	2.31	0.55
1:A:219:PHE:HB3	1:A:224:VAL:HG22	1.88	0.55
1:B:102:ASN:OD1	1:C:62:ASN:ND2	2.33	0.55
1:A:299:PHE:HB3	1:A:304:VAL:HG13	1.88	0.55
1:C:78:ILE:HG22	1:C:79:LYS:CE	2.35	0.55
1:A:238:ILE:HD12	1:A:239:LYS:HD3	1.89	0.55
1:B:285:ARG:CZ	1:B:319:LYS:HA	2.37	0.55
1:B:119:LYS:HE3	1:C:62:ASN:CB	2.33	0.54
1:B:77:ASP:HB2	1:B:84:LEU:HD21	1.90	0.54
1:B:280:THR:HG22	1:B:282:GLN:HG2	1.89	0.54
1:C:75:VAL:HG22	1:C:85:ARG:H	1.73	0.54
1:C:47:LEU:HD13	1:C:76:TRP:CG	2.44	0.53
1:C:205:ARG:HE	1:C:240:THR:C	2.17	0.53
1:A:87:LEU:HD23	1:A:116:TRP:CG	2.44	0.53
1:B:5:ARG:NH1	1:B:39:LYS:O	2.41	0.53
1:C:290:HIS:CE1	1:C:314:LYS:HG3	2.42	0.53
1:B:230:ASP:OD1	1:B:232:THR:OG1	2.21	0.53
1:C:247:LEU:HD23	1:C:276:TRP:CD2	2.43	0.53
1:A:274:LYS:HB3	1:A:276:TRP:HE1	1.72	0.53
1:B:114:LYS:HB3	1:B:116:TRP:NE1	2.24	0.53
1:C:224:VAL:HG22	1:C:238:ILE:HD13	1.90	0.53
1:A:7:LEU:HD23	1:A:36:TRP:CG	2.44	0.53
1:C:37:ASP:HB3	1:C:40:THR:OG1	2.09	0.52
1:B:122:GLN:NE2	2:B:405:HOH:O	2.42	0.52
1:A:249:GLY:C	1:A:274:LYS:HE3	2.34	0.52
1:C:143:ILE:HD13	1:C:145:VAL:HG23	1.91	0.52
1:A:154:LYS:NZ	1:A:166:THR:OG1	2.39	0.52
1:C:310:ASP:OD1	1:C:312:THR:OG1	2.22	0.51
1:A:248:GLN:C	1:A:276:TRP:HH2	2.18	0.51
1:C:19:PHE:HB2	1:C:24:VAL:HG22	1.93	0.51
1:A:179:PHE:HB3	1:A:184:VAL:HG23	1.92	0.51
1:C:63:ILE:HD12	1:C:118:ILE:HD11	1.93	0.51
1:B:167:LEU:HD23	1:B:196:TRP:CG	2.45	0.51
1:B:165:ARG:NH2	1:B:199:LYS:O	2.30	0.51
1:B:223:ILE:HD12	1:B:278:ILE:HD11	1.91	0.51
1:C:197:ASP:HB3	1:C:200:THR:HG22	1.92	0.51
1:A:194:LYS:NZ	1:A:206:THR:OG1	2.43	0.50
1:A:273:VAL:HB	1:A:287:LEU:HB2	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:184:VAL:HG13	1:B:198:ILE:HB	1.92	0.50
1:C:273:VAL:HB	1:C:287:LEU:HB2	1.92	0.50
1:C:190:ASP:OD1	1:C:192:THR:HG22	2.12	0.50
1:B:224:VAL:HG23	1:B:238:ILE:HG12	1.93	0.49
1:C:40:THR:OG1	1:C:42:GLN:HG2	2.12	0.49
1:C:143:ILE:CD1	1:C:145:VAL:HG23	2.42	0.49
1:C:5:ARG:NH1	1:C:38:ILE:O	2.45	0.49
1:A:7:LEU:HD23	1:A:36:TRP:CD2	2.47	0.49
1:A:230:ASP:OD1	1:A:232:THR:OG1	2.22	0.49
1:B:87:LEU:HD23	1:B:116:TRP:CG	2.49	0.48
1:A:6:THR:OG1	1:A:314:LYS:NZ	2.32	0.48
1:A:165:ARG:CZ	1:B:165:ARG:HH22	2.26	0.48
1:A:237:ASP:HB2	1:A:244:LEU:HD21	1.94	0.48
1:A:184:VAL:HG12	1:A:196:TRP:HB2	1.96	0.47
1:C:193:VAL:HB	1:C:207:LEU:HB2	1.95	0.47
1:C:21:ASP:OD1	1:C:21:ASP:N	2.39	0.47
1:A:159:LYS:HD3	1:A:159:LYS:HA	1.62	0.47
1:B:160:THR:HG22	1:B:162:GLN:HG2	1.95	0.47
1:C:79:LYS:HD2	1:C:79:LYS:H	1.79	0.47
1:A:117:ASP:OD1	1:A:119:LYS:HG2	2.14	0.47
1:A:167:LEU:CD2	1:A:184:VAL:HG11	2.45	0.47
1:C:65:VAL:HG11	1:C:104:VAL:HG21	1.97	0.47
1:B:117:ASP:HB3	1:B:120:THR:OG1	2.14	0.47
1:A:215:THR:OG1	1:A:254:VAL:O	2.26	0.47
1:B:104:VAL:HG13	1:B:118:ILE:HB	1.97	0.47
1:B:82:GLN:NE2	2:B:408:HOH:O	2.48	0.46
1:C:179:PHE:HB2	1:C:184:VAL:HG22	1.97	0.46
1:C:47:LEU:HB3	1:C:76:TRP:CE3	2.51	0.46
1:A:143:ILE:HD11	1:A:179:PHE:CE1	2.51	0.46
1:A:165:ARG:HH21	1:B:165:ARG:HH12	1.64	0.46
1:B:9:GLY:C	1:B:34:LYS:HE3	2.40	0.46
1:B:310:ASP:OD1	1:B:312:THR:OG1	2.27	0.46
1:C:45:ARG:NH1	1:C:79:LYS:O	2.48	0.46
1:A:158:ILE:HG23	1:A:159:LYS:HE3	1.98	0.45
1:B:179:PHE:CB	1:B:184:VAL:HG12	2.42	0.45
1:B:88:GLN:O	1:B:116:TRP:CH2	2.55	0.45
1:C:162:GLN:NE2	2:C:406:HOH:O	2.46	0.45
1:C:184:VAL:N	1:C:196:TRP:O	2.49	0.45
1:A:277:ASP:HB2	1:A:284:LEU:HD11	1.99	0.45
1:C:78:ILE:CG2	1:C:79:LYS:HE3	2.40	0.45
1:A:50:HIS:CE1	1:A:74:LYS:HD2	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:63:ILE:HD11	1:C:99:PHE:CE2	2.51	0.45
1:A:164:LEU:HB3	1:A:165:ARG:HD2	1.99	0.45
1:C:99:PHE:CB	1:C:104:VAL:HG23	2.44	0.45
1:C:249:GLY:C	1:C:274:LYS:HE2	2.42	0.45
1:A:228:SER:OG	1:A:229:ASP:N	2.50	0.45
1:C:197:ASP:H	1:C:204:LEU:HD13	1.82	0.45
1:C:205:ARG:HG3	1:C:205:ARG:HH11	1.81	0.45
1:A:165:ARG:NH2	1:B:165:ARG:HH12	2.14	0.45
1:A:247:LEU:HD23	1:A:276:TRP:CD2	2.51	0.45
1:C:64:VAL:CG2	1:C:78:ILE:HG13	2.47	0.45
1:B:318:ILE:HG23	1:B:319:LYS:HG3	1.99	0.44
1:A:104:VAL:HG22	1:A:116:TRP:HB2	1.99	0.44
1:A:165:ARG:CZ	1:A:199:LYS:O	2.65	0.44
1:A:28:SER:OG	1:A:29:ASP:N	2.51	0.43
1:C:7:LEU:HD13	1:C:36:TRP:CG	2.53	0.43
1:C:7:LEU:HD13	1:C:36:TRP:CD2	2.53	0.43
1:C:182:ASN:HA	1:C:199:LYS:NZ	2.34	0.43
1:C:302:ASN:HA	1:C:318:ILE:HG22	2.01	0.43
1:B:300:ASN:ND2	2:B:404:HOH:O	2.41	0.43
1:C:207:LEU:HB3	1:C:236:TRP:CE3	2.54	0.43
1:C:179:PHE:CB	1:C:184:VAL:HG22	2.49	0.42
1:A:245:ARG:NE	1:A:279:LYS:O	2.51	0.42
1:B:77:ASP:HB3	1:B:80:THR:OG1	2.19	0.42
1:B:247:LEU:HD13	1:B:276:TRP:CG	2.54	0.42
1:C:108:SER:OG	1:C:109:ASP:N	2.51	0.42
1:C:127:LEU:HB3	1:C:156:TRP:CE3	2.54	0.42
1:B:87:LEU:HD23	1:B:116:TRP:CD2	2.54	0.42
1:C:287:LEU:HB3	1:C:316:TRP:CE3	2.55	0.42
1:A:33:VAL:HB	1:A:47:LEU:HB2	2.01	0.42
1:C:274:LYS:HD3	1:C:276:TRP:CZ2	2.55	0.42
1:B:117:ASP:HB2	1:B:124:LEU:HD11	2.02	0.42
1:B:237:ASP:HB2	1:B:244:LEU:HD11	2.02	0.42
1:B:245:ARG:NH1	1:B:281:GLY:HA3	2.34	0.42
1:C:224:VAL:HG22	1:C:238:ILE:CD1	2.50	0.42
1:B:7:LEU:HD22	1:B:41:GLY:HA2	2.02	0.42
1:C:268:SER:OG	1:C:269:ASP:N	2.53	0.42
1:B:70:ASP:OD1	1:B:72:THR:OG1	2.26	0.41
1:C:210:HIS:CE1	1:C:228:SER:HB3	2.55	0.41
1:A:70:ASP:OD1	1:A:72:THR:OG1	2.33	0.41
1:A:261:ASP:O	1:A:279:LYS:NZ	2.53	0.41
1:B:7:LEU:HB3	1:B:36:TRP:CE3	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:182:ASN:O	1:B:198:ILE:HG22	2.20	0.41
1:A:279:LYS:HE3	1:C:302:ASN:HB2	2.01	0.41
1:B:59:PHE:HB3	1:B:64:VAL:HG22	2.02	0.41
1:C:28:SER:OG	1:C:29:ASP:N	2.54	0.41
1:A:264:VAL:HG23	1:A:278:ILE:HB	2.02	0.41
1:B:99:PHE:CB	1:B:104:VAL:HG12	2.50	0.41
1:C:262:ASN:O	1:C:278:ILE:HG22	2.21	0.41
1:B:110:ASP:OD1	1:B:112:THR:OG1	2.37	0.41
1:B:205:ARG:NE	1:B:239:LYS:O	2.54	0.41
1:C:197:ASP:HB3	1:C:200:THR:CG2	2.51	0.41
1:C:245:ARG:NH1	1:C:279:LYS:O	2.54	0.41
1:B:45:ARG:CZ	1:B:79:LYS:HD2	2.51	0.40
1:C:237:ASP:HB2	1:C:244:LEU:HD11	2.03	0.40
1:B:51:GLN:H	1:B:51:GLN:HG2	1.59	0.40
1:B:45:ARG:NH2	1:B:79:LYS:HD2	2.37	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	317/325 (98%)	295 (93%)	20 (6%)	2 (1%)	21	17
1	B	316/325 (97%)	292 (92%)	23 (7%)	1 (0%)	36	35
1	C	316/325 (97%)	295 (93%)	21 (7%)	0	100	100
All	All	949/975 (97%)	882 (93%)	64 (7%)	3 (0%)	36	35

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	89	GLY
1	A	249	GLY

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Mol	Chain	Res	Type
1	B	9	GLY

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	287/291 (99%)	287 (100%)	0	100	100
1	B	286/291 (98%)	286 (100%)	0	100	100
1	C	286/291 (98%)	286 (100%)	0	100	100
All	All	859/873 (98%)	859 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	128	GLN
1	A	142	ASN
1	A	182	ASN
1	A	202	GLN
1	A	302	ASN
1	B	91	GLN
1	B	162	GLN
1	B	182	ASN
1	C	128	GLN
1	C	298	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	319/325 (98%)	1.22	59 (18%) 3 3	31, 55, 85, 119	0
1	B	318/325 (97%)	1.13	47 (14%) 5 5	36, 52, 77, 96	0
1	C	318/325 (97%)	1.20	53 (16%) 4 4	35, 57, 81, 96	0
All	All	955/975 (97%)	1.18	159 (16%) 4 4	31, 54, 81, 119	0

All (159) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	147	GLY	5.1
1	C	163	SER	5.1
1	C	84	LEU	5.0
1	A	120	THR	4.9
1	A	313	VAL	4.8
1	C	143	ILE	4.5
1	A	192	THR	4.5
1	A	206	THR	4.5
1	A	253	ALA	4.5
1	A	126	THR	4.2
1	B	315	VAL	4.1
1	A	132	SER	3.8
1	A	63	ILE	3.8
1	B	63	ILE	3.7
1	A	71	SER	3.7
1	C	164	LEU	3.7
1	B	146	SER	3.7
1	B	110	ASP	3.6
1	A	84	LEU	3.5
1	B	164	LEU	3.5
1	C	134	VAL	3.5
1	C	131	GLN	3.5
1	B	89	GLY	3.4

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Mol	Chain	Res	Type	RSRZ
1	C	47	LEU	3.4
1	A	111	SER	3.4
1	C	300	ASN	3.4
1	C	114	LYS	3.4
1	C	104	VAL	3.3
1	A	67	GLY	3.3
1	B	280	THR	3.2
1	C	184	VAL	3.2
1	A	158	ILE	3.2
1	B	104	VAL	3.2
1	A	152	THR	3.1
1	B	249	GLY	3.1
1	B	196	TRP	3.1
1	A	241	GLY	3.1
1	B	121	GLY	3.1
1	B	25	VAL	3.1
1	B	159	LYS	3.1
1	C	49	GLY	3.1
1	B	277	ASP	3.1
1	C	235	VAL	3.1
1	A	121	GLY	3.0
1	A	187	GLY	3.0
1	A	219	PHE	3.0
1	A	4	LEU	3.0
1	C	120	THR	3.0
1	B	64	VAL	3.0
1	C	20	ASN	3.0
1	C	303	ILE	3.0
1	C	135	THR	2.9
1	A	136	SER	2.9
1	C	17	LEU	2.9
1	A	189	ASP	2.9
1	B	73	VAL	2.9
1	B	198	ILE	2.9
1	C	67	GLY	2.9
1	C	44	LEU	2.8
1	C	113	VAL	2.8
1	A	232	THR	2.8
1	C	207	LEU	2.8
1	A	210	HIS	2.7
1	B	147	GLY	2.7
1	C	129	GLY	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	103	ILE	2.7
1	A	195	VAL	2.7
1	C	188	SER	2.7
1	C	294	VAL	2.7
1	C	156	TRP	2.7
1	B	20	ASN	2.6
1	B	15	THR	2.6
1	C	287	LEU	2.6
1	C	205	ARG	2.6
1	A	169	GLY	2.6
1	B	100	ASN	2.6
1	B	76	TRP	2.6
1	B	84	LEU	2.6
1	A	305	VAL	2.5
1	C	238	ILE	2.5
1	B	141	ASP	2.5
1	A	265	VAL	2.5
1	A	22	ASN	2.5
1	C	301	ASP	2.5
1	A	243	SER	2.5
1	B	27	GLY	2.4
1	B	232	THR	2.4
1	C	147	GLY	2.4
1	C	22	ASN	2.4
1	C	203	SER	2.4
1	C	284	LEU	2.4
1	B	9	GLY	2.4
1	A	65	VAL	2.4
1	B	238	ILE	2.4
1	C	275	VAL	2.4
1	B	287	LEU	2.3
1	A	281	GLY	2.3
1	B	114	LYS	2.3
1	A	36	TRP	2.3
1	B	22	ASN	2.3
1	A	198	ILE	2.3
1	A	53	ALA	2.3
1	A	261	ASP	2.3
1	B	49	GLY	2.3
1	B	227	GLY	2.3
1	C	132	SER	2.3
1	A	318	ILE	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	184	VAL	2.2
1	A	315	VAL	2.2
1	B	105	VAL	2.2
1	B	126	THR	2.2
1	B	195	VAL	2.2
1	B	260	ASN	2.2
1	A	87	LEU	2.2
1	A	298	GLN	2.2
1	C	167	LEU	2.2
1	A	2	GLN	2.2
1	C	202	GLN	2.2
1	B	43	SER	2.2
1	A	276	TRP	2.2
1	A	62	ASN	2.2
1	A	197	ASP	2.2
1	A	33	VAL	2.2
1	A	144	VAL	2.2
1	A	41	GLY	2.2
1	A	28	SER	2.2
1	C	45	ARG	2.2
1	B	62	ASN	2.2
1	B	157	ASP	2.2
1	C	141	ASP	2.2
1	C	272	THR	2.2
1	A	38	ILE	2.1
1	B	202	GLN	2.1
1	B	71	SER	2.1
1	A	196	TRP	2.1
1	A	307	GLY	2.1
1	B	161	GLY	2.1
1	C	256	SER	2.1
1	A	69	ASP	2.1
1	A	301	ASP	2.1
1	B	70	ASP	2.1
1	C	181	ASP	2.1
1	C	263	ILE	2.1
1	C	320	GLY	2.1
1	B	305	VAL	2.1
1	C	148	SER	2.1
1	C	244	LEU	2.1
1	A	259	PHE	2.1
1	A	264	VAL	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	68	SER	2.1
1	C	66	SER	2.1
1	A	47	LEU	2.0
1	A	139	PHE	2.0
1	C	19	PHE	2.0
1	C	105	VAL	2.0
1	C	212	SER	2.0
1	C	101	ASP	2.0
1	B	41	GLY	2.0
1	C	107	GLY	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.