



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 03:05 am GMT

PDB ID : 2IHR
Title : RF2 of Thermus thermophilus
Authors : Dobbek, H.; Voertler, C.S.; Sprinzl, M.
Deposited on : 2006-09-27
Resolution : 2.50 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 : trunk28620
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) : recalc28949

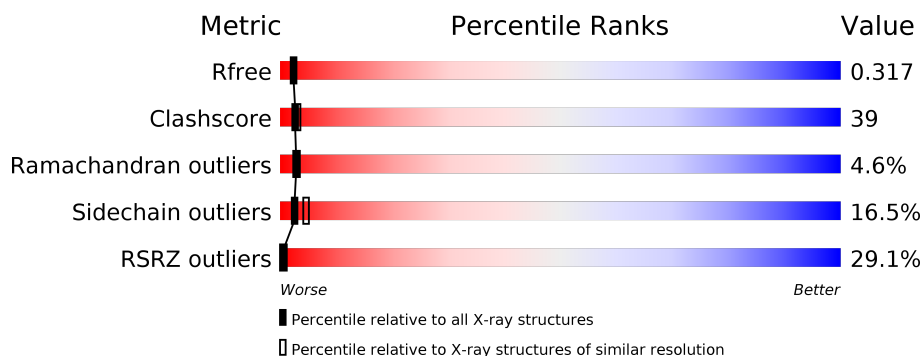
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	3846 (2.50-2.50)
Clashscore	112137	4554 (2.50-2.50)
Ramachandran outliers	110173	4463 (2.50-2.50)
Sidechain outliers	110143	4465 (2.50-2.50)
RSRZ outliers	101464	3876 (2.50-2.50)

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	1	365	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2823 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 Peptide chain release factor 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	1	351	Total	C	N	O	S	0	0	0
			2792	1748	503	533	8			

There are 13 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
1	-4	MET	-	CLONING ARTIFACT	UNP Q5SM01
1	-3	ASP	-	CLONING ARTIFACT	UNP Q5SM01
1	-2	LEU	-	CLONING ARTIFACT	UNP Q5SM01
1	-1	GLU	-	CLONING ARTIFACT	UNP Q5SM01
1	0	ARG	-	CLONING ARTIFACT	UNP Q5SM01
1	1	LEU	-	CLONING ARTIFACT	UNP Q5SM01
1	2	ALA	-	CLONING ARTIFACT	UNP Q5SM01
1	3	GLN	-	CLONING ARTIFACT	UNP Q5SM01
1	4	ARG	-	CLONING ARTIFACT	UNP Q5SM01
1	5	LEU	-	CLONING ARTIFACT	UNP Q5SM01
1	6	GLU	-	CLONING ARTIFACT	UNP Q5SM01
1	7	GLY	-	CLONING ARTIFACT	UNP Q5SM01
1	8	LEU	-	CLONING ARTIFACT	UNP Q5SM01

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	1	31	Total	O	0	0
			31	31		

4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	144.65Å 144.65Å 54.47Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.50 24.81 – 2.37	Depositor EDS
% Data completeness (in resolution range)	99.0 (20.00-2.50) 98.8 (24.81-2.37)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.99 (at 2.36Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.269 , 0.317 0.268 , 0.317	Depositor DCC
R_{free} test set	1018 reflections (5.00%)	DCC
Wilson B-factor (Å ²)	55.6	Xtriage
Anisotropy	0.418	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 73.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	2823	wwPDB-VP
Average B, all atoms (Å ²)	120.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.56% 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	1	0.60	0/2840	0.94	5/3837 (0.1%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	1	104	PHE	C-N-CD	-16.95	83.31	120.60
1	1	104	PHE	C-N-CA	10.22	164.91	122.00
1	1	333	ASP	CB-CG-OD2	-5.55	113.30	118.30
1	1	277	ARG	NE-CZ-NH2	-5.46	117.57	120.30
1	1	314	ASN	N-CA-C	5.01	124.52	111.00

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	1	2792	0	2809	221	0
2	1	31	0	0	3	0
All	All	2823	0	2809	221	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 39.

All (221) 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)
1:1:49:ARG:HA	1:1:52:ARG:HD2	1.43	1.00
1:1:21:LEU:HD13	1:1:43:VAL:HG23	1.52	0.90
1:1:17:LYS:HE3	1:1:21:LEU:HD21	1.50	0.89
1:1:297:VAL:HG12	1:1:298:ARG:H	1.40	0.87
1:1:96:LEU:O	1:1:100:THR:HG22	1.77	0.84
1:1:235:GLN:HA	1:1:238:ASN:HB2	1.56	0.84
1:1:88:GLU:O	1:1:92:LYS:HB2	1.81	0.81
1:1:183:ARG:HB2	1:1:305:GLN:HG2	1.65	0.78
1:1:348:ALA:O	1:1:350:ARG:N	2.17	0.77
1:1:253:ILE:HD11	1:1:281:LEU:HD12	1.66	0.76
1:1:18:GLU:OE1	1:1:19:THR:HG23	1.87	0.74
1:1:245:ARG:HH11	1:1:245:ARG:HB3	1.53	0.74
1:1:210:ASP:OD1	1:1:210:ASP:N	2.16	0.73
1:1:64:LEU:HD12	1:1:64:LEU:O	1.87	0.73
1:1:17:LYS:HA	1:1:20:ARG:HD2	1.69	0.72
1:1:316:VAL:HG23	1:1:328:PRO:HB3	1.71	0.72
1:1:5:LEU:HD12	1:1:5:LEU:O	1.89	0.72
1:1:88:GLU:O	1:1:92:LYS:HD3	1.90	0.72
1:1:218:LYS:HB2	1:1:221:GLU:CG	2.21	0.71
1:1:257:CYS:SG	1:1:269:LEU:HD23	2.29	0.71
1:1:54:VAL:HA	1:1:57:PHE:CD2	2.26	0.70
1:1:18:GLU:HB3	1:1:47:ALA:HA	1.73	0.70
1:1:70:LEU:O	1:1:73:GLU:O	2.08	0.70
1:1:274:LEU:O	1:1:274:LEU:HD22	1.92	0.69
1:1:61:GLU:HG2	1:1:64:LEU:HD23	1.75	0.69
1:1:228:ARG:HH21	1:1:240:THR:HA	1.57	0.69
1:1:5:LEU:C	1:1:5:LEU:HD12	2.13	0.69
1:1:76:ALA:HB1	1:1:79:ARG:HB2	1.75	0.69
1:1:227:MET:SD	1:1:245:ARG:NH1	2.66	0.68
1:1:14:ILE:N	1:1:15:PRO:HD2	2.09	0.68
1:1:291:LYS:NZ	1:1:299:PRO:HG3	2.08	0.68
1:1:321:THR:CG2	1:1:323:LEU:H	2.07	0.68
1:1:107:ALA:O	1:1:170:ASN:HA	1.94	0.68
1:1:284:LYS:O	1:1:288:GLU:HG2	1.95	0.67
1:1:54:VAL:HG13	1:1:57:PHE:CZ	2.30	0.67
1:1:28:LEU:HD12	1:1:28:LEU:N	2.11	0.66
1:1:208:GLU:HG2	1:1:208:GLU:O	1.96	0.66
1:1:61:GLU:HG2	1:1:64:LEU:CD2	2.27	0.65
1:1:218:LYS:HB2	1:1:221:GLU:HG2	1.78	0.65
1:1:105:PRO:O	1:1:107:ALA:N	2.25	0.64
1:1:61:GLU:O	1:1:64:LEU:HB3	1.97	0.64
1:1:6:GLU:O	1:1:11:ILE:HD12	1.97	0.64

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:5:LEU:HD21	1:1:96:LEU:CD2	2.28	0.63
1:1:49:ARG:HG2	1:1:49:ARG:HH11	1.63	0.63
1:1:129:GLU:OE1	1:1:158:GLY:HA2	1.99	0.63
1:1:21:LEU:HD22	1:1:43:VAL:HG21	1.81	0.62
1:1:84:PRO:O	1:1:88:GLU:HG2	2.00	0.62
1:1:280:GLU:OE1	1:1:283:ARG:NH2	2.32	0.61
1:1:245:ARG:HH11	1:1:245:ARG:CB	2.13	0.61
1:1:91:LYS:O	1:1:95:GLU:HG2	2.00	0.61
1:1:22:LYS:O	1:1:26:ARG:HG3	2.01	0.61
1:1:30:ASP:OD1	1:1:37:PRO:HB3	2.01	0.61
1:1:348:ALA:C	1:1:350:ARG:H	2.04	0.61
1:1:11:ILE:HD13	1:1:11:ILE:H	1.66	0.60
1:1:54:VAL:HG13	1:1:57:PHE:CE2	2.36	0.60
1:1:83:LYS:N	1:1:84:PRO:HD2	2.17	0.59
1:1:46:GLU:CG	1:1:49:ARG:HH21	2.16	0.59
1:1:204:GLU:OE2	1:1:205:VAL:N	2.36	0.59
1:1:18:GLU:CD	1:1:19:THR:HG23	2.21	0.59
1:1:21:LEU:HD22	1:1:43:VAL:CG2	2.33	0.59
1:1:300:ILE:HD13	1:1:300:ILE:H	1.68	0.59
1:1:321:THR:CG2	1:1:323:LEU:HB2	2.33	0.59
1:1:18:GLU:CB	1:1:47:ALA:HA	2.33	0.59
1:1:57:PHE:C	1:1:57:PHE:CD1	2.76	0.58
1:1:11:ILE:CD1	1:1:11:ILE:H	2.17	0.58
1:1:57:PHE:HA	1:1:60:LEU:HB2	1.86	0.57
1:1:44:SER:O	1:1:47:ALA:HB3	2.03	0.57
1:1:183:ARG:CB	1:1:305:GLN:HG2	2.33	0.57
1:1:228:ARG:HB3	1:1:241:ASP:OD1	2.04	0.57
1:1:59:SER:O	1:1:62:SER:HB2	2.05	0.57
1:1:321:THR:HG23	1:1:323:LEU:H	1.68	0.56
1:1:17:LYS:HE3	1:1:21:LEU:CD2	2.30	0.56
1:1:154:GLY:HA2	1:1:160:ASP:OD2	2.04	0.56
1:1:49:ARG:O	1:1:52:ARG:HB2	2.05	0.56
1:1:51:ARG:O	1:1:55:ASP:HB2	2.05	0.56
1:1:70:LEU:HD11	1:1:74:LEU:O	2.06	0.56
1:1:210:ASP:O	1:1:212:GLU:N	2.36	0.55
1:1:103:ASN:O	1:1:103:ASN:CG	2.45	0.55
1:1:73:GLU:OE1	1:1:75:PRO:O	2.24	0.55
1:1:155:PRO:CD	1:1:160:ASP:OD2	2.55	0.55
1:1:14:ILE:N	1:1:15:PRO:CD	2.69	0.54
1:1:211:GLU:HA	1:1:214:GLU:OE1	2.07	0.54
1:1:25:GLU:OE2	1:1:40:ALA:HB3	2.07	0.54

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:249:LEU:HB2	1:1:250:PRO:HD3	1.90	0.54
1:1:252:GLY:O	1:1:254:THR:HG22	2.08	0.54
1:1:70:LEU:HA	1:1:73:GLU:O	2.09	0.53
1:1:57:PHE:O	1:1:61:GLU:HB2	2.09	0.53
1:1:228:ARG:O	1:1:228:ARG:HG2	2.09	0.53
1:1:299:PRO:HB2	1:1:300:ILE:HD13	1.91	0.53
1:1:300:ILE:O	1:1:300:ILE:HG12	2.09	0.52
1:1:46:GLU:HG3	1:1:49:ARG:HH21	1.73	0.52
1:1:62:SER:O	1:1:66:GLY:N	2.43	0.52
1:1:111:ALA:HB2	1:1:172:TYR:HA	1.90	0.52
1:1:63:ASP:C	1:1:65:GLN:H	2.12	0.52
1:1:113:LEU:HA	1:1:204:GLU:O	2.10	0.52
1:1:262:SER:HB2	1:1:264:ILE:CD1	2.40	0.52
1:1:291:LYS:HZ3	1:1:299:PRO:HG3	1.73	0.52
1:1:110:ASN:HD22	1:1:110:ASN:N	2.08	0.52
1:1:5:LEU:C	1:1:7:GLY:N	2.63	0.52
1:1:5:LEU:C	1:1:7:GLY:H	2.12	0.52
1:1:31:PRO:HA	1:1:37:PRO:HG2	1.91	0.52
1:1:6:GLU:OE1	1:1:53:THR:CG2	2.58	0.52
1:1:76:ALA:HB1	1:1:79:ARG:CB	2.39	0.52
1:1:155:PRO:HD3	1:1:160:ASP:OD2	2.10	0.52
1:1:169:GLU:HG2	1:1:170:ASN:ND2	2.25	0.51
1:1:297:VAL:HG12	1:1:298:ARG:N	2.19	0.51
1:1:34:TRP:O	1:1:38:GLU:HG2	2.10	0.51
1:1:73:GLU:CD	1:1:75:PRO:O	2.48	0.51
1:1:213:VAL:HG12	1:1:213:VAL:O	2.10	0.51
1:1:95:GLU:O	1:1:99:GLN:N	2.39	0.51
1:1:335:ASP:HA	2:1:1029:HOH:O	2.10	0.51
1:1:318:ASP:HB3	1:1:321:THR:HG22	1.92	0.51
1:1:321:THR:HG22	1:1:323:LEU:H	1.75	0.51
1:1:348:ALA:C	1:1:350:ARG:N	2.63	0.51
1:1:228:ARG:O	1:1:229:ALA:C	2.49	0.51
1:1:12:PHE:C	1:1:15:PRO:HD2	2.31	0.50
1:1:208:GLU:CG	1:1:208:GLU:O	2.59	0.50
1:1:174:LEU:O	1:1:177:PRO:HD2	2.10	0.50
1:1:253:ILE:HD12	1:1:278:LEU:HD23	1.93	0.50
1:1:51:ARG:HG2	1:1:51:ARG:HH11	1.77	0.50
1:1:210:ASP:C	1:1:212:GLU:H	2.15	0.50
1:1:50:LEU:O	1:1:53:THR:HB	2.11	0.50
1:1:282:GLU:HA	1:1:282:GLU:OE1	2.12	0.49
1:1:28:LEU:CD1	1:1:28:LEU:N	2.75	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:263:GLN:NE2	1:1:267:LYS:HE3	2.27	0.49
1:1:323:LEU:HD13	1:1:339:LEU:HD23	1.94	0.49
1:1:40:ALA:HA	1:1:43:VAL:HG22	1.95	0.49
1:1:73:GLU:OE2	1:1:76:ALA:HB2	2.12	0.49
1:1:127:TRP:CE2	1:1:184:LEU:HB2	2.48	0.49
1:1:211:GLU:HA	1:1:214:GLU:HB2	1.94	0.49
1:1:48:ALA:HB1	1:1:52:ARG:NH2	2.28	0.49
1:1:5:LEU:HD21	1:1:96:LEU:HD23	1.95	0.48
1:1:218:LYS:O	1:1:222:LEU:HD13	2.13	0.48
1:1:27:ARG:HG2	1:1:29:GLU:H	1.79	0.48
1:1:300:ILE:HD13	1:1:300:ILE:N	2.27	0.48
1:1:6:GLU:OE2	1:1:11:ILE:HG21	2.14	0.48
1:1:97:TYR:O	1:1:100:THR:HG23	2.14	0.48
1:1:70:LEU:CD1	1:1:74:LEU:O	2.62	0.48
1:1:63:ASP:C	1:1:65:GLN:N	2.66	0.48
1:1:83:LYS:N	1:1:84:PRO:CD	2.76	0.48
1:1:76:ALA:HB1	1:1:79:ARG:CG	2.44	0.47
1:1:34:TRP:O	1:1:37:PRO:HD2	2.14	0.47
1:1:274:LEU:HD22	1:1:274:LEU:C	2.36	0.47
1:1:28:LEU:H	1:1:28:LEU:HD12	1.80	0.46
1:1:36:ASP:N	1:1:37:PRO:CD	2.78	0.46
1:1:33:LEU:HD12	1:1:33:LEU:N	2.30	0.46
1:1:95:GLU:HA	1:1:95:GLU:OE1	2.13	0.46
1:1:306:ILE:HG13	1:1:307:ARG:N	2.30	0.46
1:1:101:LEU:HD21	1:1:345:GLU:HG2	1.98	0.46
1:1:54:VAL:HA	1:1:57:PHE:CE2	2.51	0.46
1:1:13:ASP:N	1:1:15:PRO:HD2	2.31	0.46
1:1:123:GLU:HB3	1:1:199:SER:OG	2.16	0.46
1:1:222:LEU:CD1	1:1:222:LEU:N	2.79	0.46
1:1:218:LYS:HB2	1:1:221:GLU:HG3	1.98	0.46
1:1:40:ALA:HA	1:1:43:VAL:CG2	2.46	0.46
1:1:69:GLU:O	1:1:73:GLU:HG3	2.16	0.46
1:1:105:PRO:C	1:1:107:ALA:H	2.17	0.45
1:1:193:SER:C	1:1:195:ARG:H	2.19	0.45
1:1:21:LEU:HD13	1:1:43:VAL:C	2.36	0.45
1:1:53:THR:HA	1:1:56:THR:HG22	1.98	0.45
1:1:46:GLU:HG2	1:1:49:ARG:HH21	1.79	0.45
1:1:185:VAL:HB	1:1:310:VAL:HB	1.99	0.45
1:1:56:THR:O	1:1:60:LEU:HB2	2.16	0.45
1:1:264:ILE:H	1:1:264:ILE:HD13	1.81	0.45
1:1:137:ARG:O	1:1:141:ARG:HG2	2.16	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:66:GLY:C	1:1:68:LEU:H	2.19	0.45
1:1:51:ARG:HH22	1:1:52:ARG:HE	1.64	0.45
1:1:1:LEU:C	1:1:3:GLN:N	2.70	0.45
1:1:327:ASP:OD1	1:1:327:ASP:N	2.48	0.45
1:1:289:GLU:O	1:1:292:ALA:N	2.41	0.44
1:1:57:PHE:C	1:1:57:PHE:HD1	2.18	0.44
1:1:221:GLU:C	1:1:222:LEU:HD12	2.38	0.44
1:1:36:ASP:H	1:1:37:PRO:CD	2.30	0.44
1:1:53:THR:CA	1:1:56:THR:HG22	2.47	0.44
1:1:218:LYS:O	1:1:221:GLU:HB2	2.18	0.44
1:1:289:GLU:C	1:1:291:LYS:N	2.70	0.44
1:1:191:ASP:C	1:1:191:ASP:OD1	2.56	0.43
1:1:287:GLU:C	1:1:289:GLU:N	2.72	0.43
1:1:21:LEU:CD1	1:1:43:VAL:HG23	2.36	0.43
1:1:11:ILE:CD1	1:1:11:ILE:N	2.79	0.43
1:1:172:TYR:CG	1:1:207:PRO:HB3	2.54	0.43
1:1:289:GLU:C	1:1:291:LYS:H	2.22	0.43
1:1:220:GLU:HG3	2:1:1023:HOH:O	2.19	0.43
1:1:221:GLU:OE1	1:1:221:GLU:HA	2.18	0.43
1:1:49:ARG:HG2	1:1:49:ARG:NH1	2.33	0.43
1:1:218:LYS:O	1:1:221:GLU:N	2.36	0.43
1:1:103:ASN:O	1:1:103:ASN:ND2	2.52	0.43
1:1:241:ASP:O	1:1:242:SER:HB2	2.18	0.43
1:1:341:TRP:HA	1:1:341:TRP:CE3	2.54	0.43
1:1:291:LYS:HZ1	1:1:299:PRO:HG3	1.82	0.43
1:1:127:TRP:HA	1:1:130:MET:HE3	2.01	0.43
1:1:18:GLU:HB2	1:1:47:ALA:HB1	2.00	0.42
1:1:110:ASN:N	1:1:110:ASN:ND2	2.67	0.42
1:1:34:TRP:C	1:1:37:PRO:HD2	2.39	0.42
1:1:244:VAL:HG21	1:1:267:LYS:HG2	2.01	0.42
1:1:28:LEU:CD1	1:1:28:LEU:H	2.31	0.42
1:1:93:LEU:O	1:1:94:ASP:C	2.57	0.42
1:1:84:PRO:O	1:1:88:GLU:CG	2.65	0.42
1:1:175:LEU:HD23	1:1:175:LEU:HA	1.81	0.42
1:1:245:ARG:HH11	1:1:245:ARG:CG	2.33	0.41
1:1:262:SER:HB2	1:1:264:ILE:HD11	2.01	0.41
1:1:126:ASP:O	1:1:130:MET:HG3	2.19	0.41
1:1:95:GLU:O	1:1:96:LEU:C	2.58	0.41
1:1:248:HIS:CE1	1:1:251:THR:HG23	2.55	0.41
1:1:210:ASP:O	1:1:211:GLU:HG2	2.20	0.41
1:1:68:LEU:O	1:1:72:GLU:HB2	2.20	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:69:GLU:O	1:1:73:GLU:CG	2.69	0.41
1:1:94:ASP:O	1:1:95:GLU:C	2.58	0.41
1:1:289:GLU:O	1:1:291:LYS:N	2.53	0.41
1:1:195:ARG:NH2	2:1:1015:HOH:O	2.54	0.41
1:1:310:VAL:CG1	1:1:315:TYR:CZ	3.03	0.41
1:1:73:GLU:HG3	1:1:73:GLU:O	2.21	0.41
1:1:13:ASP:C	1:1:15:PRO:HD2	2.41	0.41
1:1:70:LEU:HD12	1:1:74:LEU:HA	2.02	0.41
1:1:281:LEU:O	1:1:281:LEU:HD23	2.20	0.41
1:1:287:GLU:C	1:1:289:GLU:H	2.24	0.41
1:1:53:THR:C	1:1:56:THR:HG22	2.42	0.40
1:1:321:THR:HG23	1:1:323:LEU:HB2	2.02	0.40
1:1:172:TYR:O	1:1:176:SER:HB3	2.21	0.40
1:1:193:SER:O	1:1:195:ARG:N	2.48	0.40
1:1:228:ARG:NH2	1:1:240:THR:HA	2.29	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	1	349/365 (96%)	290 (83%)	43 (12%)	16 (5%)	 

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	1	65	GLN
1	1	105	PRO
1	1	106	HIS
1	1	211	GLU
1	1	229	ALA
1	1	10	GLY

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	1	34	TRP
1	1	230	SER
1	1	299	PRO
1	1	349	GLY
1	1	104	PHE
1	1	232	PRO
1	1	242	SER
1	1	209	VAL
1	1	295	GLY
1	1	215	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	1	297/309 (96%)	248 (84%)	49 (16%)	2 4

All (49) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	1	5	LEU
1	1	11	ILE
1	1	23	GLU
1	1	24	LEU
1	1	45	GLN
1	1	50	LEU
1	1	57	PHE
1	1	61	GLU
1	1	62	SER
1	1	64	LEU
1	1	65	GLN
1	1	69	GLU
1	1	70	LEU
1	1	73	GLU
1	1	78	GLU
1	1	94	ASP

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	1	100	THR
1	1	103	ASN
1	1	104	PHE
1	1	105	PRO
1	1	108	GLU
1	1	132	LEU
1	1	142	GLN
1	1	147	GLU
1	1	160	ASP
1	1	169	GLU
1	1	184	LEU
1	1	195	ARG
1	1	204	GLU
1	1	210	ASP
1	1	215	VAL
1	1	216	VAL
1	1	222	LEU
1	1	228	ARG
1	1	230	SER
1	1	244	VAL
1	1	245	ARG
1	1	254	THR
1	1	258	GLN
1	1	260	THR
1	1	264	ILE
1	1	274	LEU
1	1	300	ILE
1	1	305	GLN
1	1	310	VAL
1	1	321	THR
1	1	323	LEU
1	1	324	MET
1	1	350	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	1	65	GLN
1	1	103	ASN
1	1	110	ASN
1	1	142	GLN
1	1	170	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	1	258	GLN
1	1	263	GLN
1	1	314	ASN
1	1	326	HIS

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 ⓘ

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	1	351/365 (96%)	2.25	102 (29%) 1 0	41, 90, 239, 276	0

All (102) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	1	236	GLY	23.2
1	1	82	LEU	22.2
1	1	43	VAL	19.7
1	1	26	ARG	16.1
1	1	85	GLU	15.8
1	1	68	LEU	14.0
1	1	78	GLU	13.8
1	1	35	ASN	13.6
1	1	29	GLU	13.5
1	1	229	ALA	13.4
1	1	23	GLU	13.0
1	1	76	ALA	13.0
1	1	75	PRO	12.9
1	1	34	TRP	12.9
1	1	71	MET	12.4
1	1	42	LYS	12.3
1	1	28	LEU	12.1
1	1	239	THR	11.4
1	1	67	LEU	11.3
1	1	84	PRO	11.1
1	1	30	ASP	11.0
1	1	44	SER	10.6
1	1	240	THR	10.4
1	1	14	ILE	10.3
1	1	233	GLY	10.2
1	1	237	VAL	9.8
1	1	31	PRO	9.6

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	1	232	PRO	9.6
1	1	74	LEU	9.5
1	1	39	ALA	9.3
1	1	19	THR	9.1
1	1	235	GLN	8.8
1	1	20	ARG	8.6
1	1	33	LEU	8.5
1	1	297	VAL	8.4
1	1	72	GLU	7.9
1	1	17	LYS	7.8
1	1	12	PHE	7.6
1	1	27	ARG	7.5
1	1	36	ASP	7.0
1	1	293	LEU	7.0
1	1	38	GLU	6.8
1	1	81	ALA	6.7
1	1	230	SER	6.6
1	1	45	GLN	6.4
1	1	234	GLY	6.4
1	1	238	ASN	6.2
1	1	296	GLU	6.2
1	1	22	LYS	6.2
1	1	80	GLU	6.1
1	1	79	ARG	6.1
1	1	294	ARG	5.7
1	1	86	LEU	5.7
1	1	24	LEU	5.6
1	1	40	ALA	5.5
1	1	298	ARG	5.5
1	1	46	GLU	5.4
1	1	32	SER	5.2
1	1	11	ILE	5.2
1	1	66	GLY	5.1
1	1	77	GLU	4.9
1	1	41	ARG	4.8
1	1	1	LEU	4.6
1	1	301	GLU	4.6
1	1	192	ALA	4.2
1	1	231	GLY	4.2
1	1	55	ASP	4.1
1	1	295	GLY	4.0
1	1	37	PRO	3.9

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	1	299	PRO	3.9
1	1	13	ASP	3.8
1	1	241	ASP	3.8
1	1	111	ALA	3.7
1	1	83	LYS	3.6
1	1	18	GLU	3.5
1	1	291	LYS	3.3
1	1	112	ILE	3.2
1	1	3	GLN	3.2
1	1	224	ILE	3.1
1	1	89	ALA	3.1
1	1	135	TYR	3.0
1	1	15	PRO	2.9
1	1	300	ILE	2.9
1	1	6	GLU	2.8
1	1	257	CYS	2.7
1	1	65	GLN	2.6
1	1	164	ILE	2.6
1	1	287	GLU	2.5
1	1	134	MET	2.4
1	1	70	LEU	2.4
1	1	163	GLN	2.4
1	1	246	VAL	2.4
1	1	274	LEU	2.4
1	1	206	ILE	2.3
1	1	58	ARG	2.2
1	1	52	ARG	2.2
1	1	349	GLY	2.2
1	1	292	ALA	2.1
1	1	113	LEU	2.1
1	1	131	LEU	2.1
1	1	256	THR	2.0
1	1	25	GLU	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.