



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 06:11 PM UTC

PDB ID : 1X1O / pdb_00001x1o
Title : Crystal structure of project ID TT0268 from *Thermus thermophilus* HB8
Authors : Shimizu, K.; Kunishima, N.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)
Deposited on : 2005-04-08
Resolution : 1.90 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 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

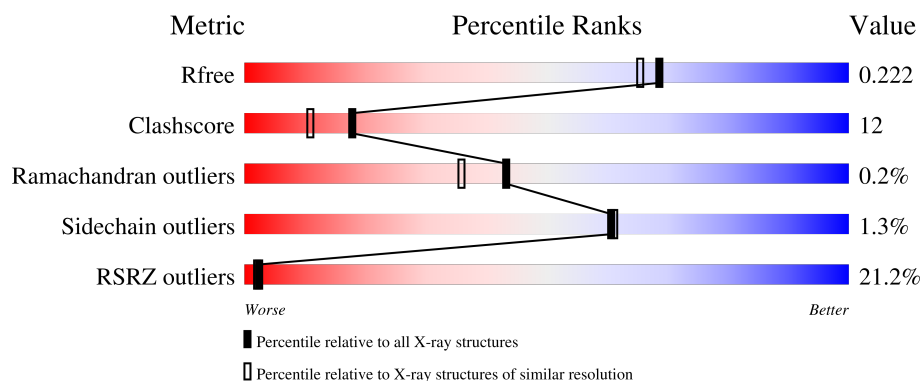
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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}	180053	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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 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	286	3% 78% 18% ..
1	B	286	4% 77% 19% .
1	C	286	55% 63% 31% ..

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7079 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 nicotinate-nucleotide pyrophosphorylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	277	Total	C	N	O	S	0	11	0
			2124	1338	392	393	1			
1	B	276	Total	C	N	O	S	0	4	0
			2089	1314	388	386	1			
1	C	277	Total	C	N	O	S	0	1	0
			2093	1317	388	387	1			

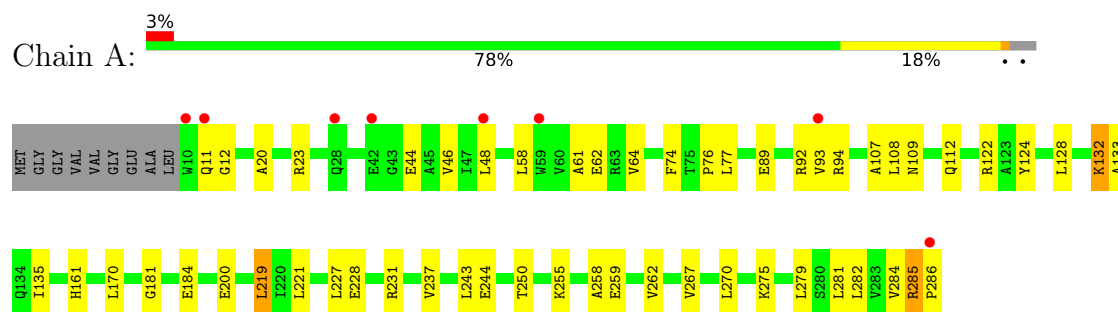
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	307	Total	O	0	0
			307	307		
2	B	357	Total	O	0	0
			357	357		
2	C	109	Total	O	0	0
			109	109		

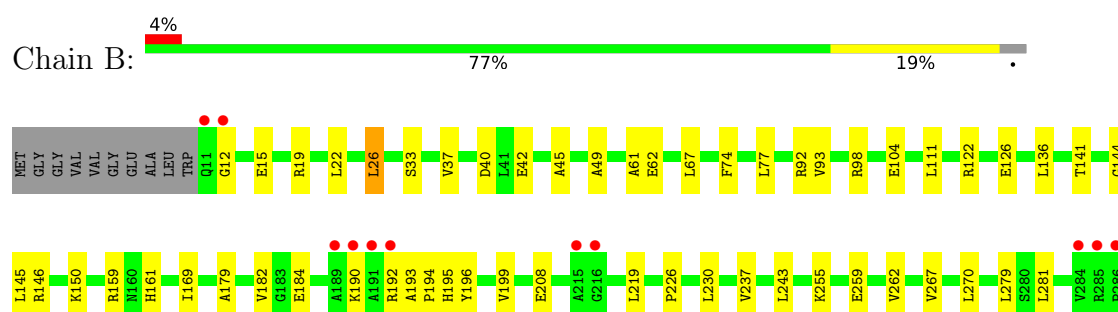
3 Residue-property plots [i](#)

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.

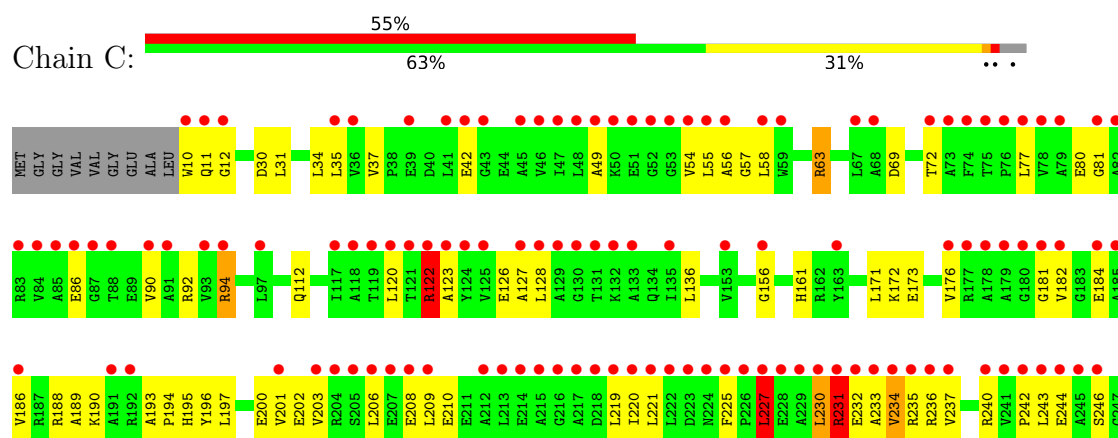
- Molecule 1: nicotinate-nucleotide pyrophosphorylase



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R248	R249	T250	L251	E252	R253	A254	K255	A256	A257	A258	E259	A260	G261	V262	D263	Y264	V265	S266	V267	G268	A269	L270	T271	H272	S273	A274	R275	A276	L277	D278	L279	S280	L281	L282	V283	V284	R285	P286
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4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	112.51Å 73.68Å 108.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.13 – 1.90 39.13 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.7 (39.13-1.90) 99.9 (39.13-1.90)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.14 (at 1.89Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.221 , 0.246 0.225 , 0.222	Depositor DCC
R_{free} test set	3631 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	24.1	Xtriage
Anisotropy	0.334	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 45.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.007 for l,-k,h	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7079	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.87% 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

5.1 Standard geometry

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	0.39	0/2206	0.89	5/2986 (0.2%)
1	B	0.39	0/2134	0.85	4/2888 (0.1%)
1	C	0.50	0/2125	0.99	14/2875 (0.5%)
All	All	0.43	0/6465	0.91	23/8749 (0.3%)

There are no bond length outliers.

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	225	PHE	CA-CB-CG	-10.41	103.39	113.80
1	C	225	PHE	O-C-N	-9.33	112.50	121.18
1	C	227	LEU	O-C-N	-7.74	113.17	122.22
1	A	161	HIS	N-CA-C	-7.43	98.34	109.86
1	A	219[A]	LEU	O-C-N	-7.32	114.47	123.33
1	A	219[B]	LEU	O-C-N	-7.32	114.47	123.33
1	C	234	VAL	CB-CA-C	-7.29	100.86	112.16
1	A	250	THR	N-CA-C	-6.31	99.73	109.14
1	B	104	GLU	N-CA-C	5.84	118.15	111.02
1	C	161	HIS	N-CA-C	-5.83	100.83	109.86
1	B	161	HIS	N-CA-C	-5.81	100.85	109.86
1	B	26	LEU	N-CA-C	5.78	119.51	112.23
1	C	225	PHE	CB-CA-C	5.55	118.97	109.82
1	C	122	ARG	CD-NE-CZ	-5.26	117.04	124.40
1	B	226	PRO	N-CA-C	-5.22	103.58	111.41
1	A	285	ARG	CD-NE-CZ	-5.11	117.24	124.40
1	C	63	ARG	CD-NE-CZ	-5.09	117.28	124.40
1	C	37	VAL	N-CA-C	5.08	113.53	107.73
1	C	94	ARG	CD-NE-CZ	-5.07	117.30	124.40
1	C	231	ARG	CD-NE-CZ	-5.07	117.30	124.40
1	C	230	LEU	CA-C-N	-5.04	112.69	122.06
1	C	230	LEU	C-N-CA	-5.04	112.69	122.06
1	C	225	PHE	CA-C-O	5.01	124.45	119.49

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	2124	0	2216	48	0
1	B	2089	0	2178	36	0
1	C	2093	0	2169	78	0
2	A	307	0	0	1	0
2	B	357	0	0	4	0
2	C	109	0	0	3	0
All	All	7079	0	6563	159	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 12.

All (159) 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:33:SER:O	1:B:37[B]:VAL:HG12	1.61	1.00
1:A:132[A]:LYS:HD2	1:A:133:ALA:N	1.83	0.92
1:A:58[B]:LEU:HD21	1:A:76:PRO:HB3	1.64	0.80
1:C:120:LEU:O	1:C:123:ALA:HB3	1.85	0.75
1:C:49:ALA:HA	1:C:279:LEU:HD23	1.67	0.75
1:C:221:LEU:HD11	1:C:246:SER:HB2	1.69	0.73
1:C:210:GLU:OE1	1:C:236:ARG:HD3	1.89	0.71
1:A:46:VAL:HG12	1:A:92:ARG:HG2	1.72	0.71
1:A:284:VAL:O	1:A:285:ARG:HD3	1.94	0.68
1:B:22:LEU:HG	1:B:26:LEU:HD12	1.77	0.66
1:C:250:THR:HG23	1:C:253:ARG:H	1.61	0.66
1:A:61:ALA:O	1:A:64[A]:VAL:HG22	1.96	0.64
1:A:48:LEU:CD2	1:A:89:GLU:HG2	2.28	0.64
1:C:267:VAL:HB	1:C:270:LEU:HG	1.79	0.64
1:C:31:LEU:O	1:C:35:LEU:HD13	1.97	0.63
1:C:203:VAL:HG22	1:C:220:ILE:HD11	1.81	0.63
1:A:58[B]:LEU:CD2	1:A:76:PRO:HB3	2.28	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:195:HIS:O	1:C:219:LEU:HD11	1.99	0.62
1:C:251:LEU:H	1:C:251:LEU:HD12	1.65	0.61
1:A:227:LEU:HG	1:A:231:ARG:NH1	2.15	0.61
1:A:255:LYS:O	1:A:259:GLU:HG3	2.01	0.60
1:A:200:GLU:CB	1:A:219[B]:LEU:HB3	2.31	0.60
1:C:172:LYS:HA	1:C:202:GLU:HG2	1.83	0.60
1:C:237:VAL:HG21	1:C:243:LEU:HD21	1.83	0.60
1:C:172:LYS:NZ	2:C:376:HOH:O	2.35	0.60
1:C:227:LEU:HD21	1:C:231:ARG:HH11	1.66	0.60
1:B:230:LEU:HD21	1:B:262:VAL:HG13	1.83	0.59
1:C:128:LEU:HB2	1:C:258:ALA:CB	2.33	0.59
1:B:179:ALA:O	1:B:184:GLU:HG2	2.03	0.58
1:C:221:LEU:HD12	1:C:244:GLU:O	2.04	0.58
1:C:171:LEU:HB2	1:C:201:VAL:HG23	1.86	0.58
1:C:58:LEU:HD21	1:C:90:VAL:CG1	2.32	0.58
1:B:37[B]:VAL:CG1	1:B:98:ARG:HG2	2.33	0.58
1:C:58:LEU:HD21	1:C:90:VAL:HG12	1.85	0.57
1:C:227:LEU:HD21	1:C:231:ARG:NH1	2.20	0.57
1:C:232:GLU:HG3	1:C:235:ARG:HH21	1.70	0.57
1:A:200:GLU:CB	1:A:219[A]:LEU:HB2	2.35	0.56
1:A:200:GLU:HB2	1:A:219[A]:LEU:HB2	1.87	0.56
1:A:200:GLU:HA	1:A:219[A]:LEU:HB2	1.88	0.56
1:C:122:ARG:O	1:C:126:GLU:HG3	2.05	0.56
1:C:220:ILE:CG2	1:C:243:LEU:HD23	2.35	0.56
1:C:234:VAL:HG12	1:C:234:VAL:O	2.05	0.55
1:C:128:LEU:HD22	1:C:258:ALA:HB2	1.87	0.55
2:B:580:HOH:O	1:C:197:LEU:HD21	2.06	0.55
1:A:285:ARG:HA	1:A:286:PRO:C	2.32	0.54
1:B:49:ALA:HA	1:B:279:LEU:HD23	1.90	0.53
1:A:64[A]:VAL:HG21	1:A:107:ALA:HB2	1.91	0.53
1:A:228:GLU:H	1:A:228:GLU:CD	2.16	0.53
1:A:275:LYS:HE2	2:A:542:HOH:O	2.08	0.53
1:B:182:VAL:HG21	1:B:208:GLU:HB3	1.91	0.53
1:C:122:ARG:HG3	1:C:156:GLY:O	2.09	0.53
1:C:237:VAL:O	1:C:240:ARG:HG3	2.08	0.53
1:A:200:GLU:CG	1:A:219[B]:LEU:HB3	2.39	0.53
1:C:285:ARG:HA	1:C:286:PRO:C	2.34	0.53
1:B:40:ASP:HA	2:B:434:HOH:O	2.08	0.52
1:C:49:ALA:O	1:C:86:GLU:HA	2.09	0.52
1:C:232:GLU:HG2	2:C:347:HOH:O	2.10	0.52
1:B:37[B]:VAL:HG13	1:B:98:ARG:HG2	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:237:VAL:HG21	1:A:243:LEU:HD21	1.90	0.52
1:A:200:GLU:HB2	1:A:219[A]:LEU:HD12	1.91	0.52
1:B:62:GLU:HG3	1:B:74:PHE:HB3	1.92	0.52
1:C:219:LEU:HD22	1:C:242:PRO:HB2	1.92	0.52
1:A:132[A]:LYS:HE2	1:A:262:VAL:O	2.10	0.51
1:C:112:GLN:HG2	1:C:277:LEU:O	2.10	0.51
1:C:10:TRP:HD1	1:C:12:GLY:H	1.59	0.51
1:C:128:LEU:HA	1:C:255[B]:LYS:HG3	1.93	0.51
1:C:10:TRP:CD1	1:C:11:GLN:HG2	2.46	0.51
1:C:182:VAL:HG21	1:C:208:GLU:HB3	1.92	0.51
1:C:10:TRP:HD1	1:C:12:GLY:N	2.09	0.51
1:C:69:ASP:HB3	1:C:72:THR:HG23	1.94	0.50
1:B:169:ILE:HD11	1:B:194:PRO:HD2	1.93	0.50
1:C:235:ARG:HH11	1:C:235:ARG:HG2	1.77	0.50
1:A:200:GLU:HB2	1:A:219[B]:LEU:HB3	1.93	0.50
1:B:192:ARG:HB2	2:B:555:HOH:O	2.12	0.50
1:C:200:GLU:HB2	1:C:219:LEU:HB2	1.93	0.50
1:B:15:GLU:OE2	1:B:67:LEU:HD21	2.11	0.50
1:B:146:ARG:O	1:B:150:LYS:HG3	2.11	0.50
1:A:200:GLU:CA	1:A:219[A]:LEU:HB2	2.43	0.49
1:C:230:LEU:HD22	1:C:260:ALA:CB	2.42	0.49
1:A:64[A]:VAL:CG2	1:A:107:ALA:HB2	2.43	0.49
1:C:77:LEU:HD11	1:C:92:ARG:HG3	1.94	0.49
1:A:77:LEU:HD11	1:A:92:ARG:HG3	1.93	0.49
1:A:109:ASN:ND2	1:B:141:THR:O	2.46	0.49
1:A:64[A]:VAL:HG21	1:A:107:ALA:HA	1.95	0.49
1:A:267:VAL:HB	1:A:270:LEU:HG	1.95	0.49
1:C:10:TRP:C	1:C:12:GLY:H	2.21	0.48
1:C:203:VAL:CG2	1:C:220:ILE:HD11	2.42	0.48
1:A:48:LEU:HD11	1:A:282:LEU:HD22	1.95	0.48
1:A:64[A]:VAL:HG21	1:A:107:ALA:CB	2.42	0.48
1:B:219:LEU:HD11	1:C:195:HIS:O	2.14	0.47
1:C:282:LEU:C	1:C:282:LEU:HD23	2.39	0.47
1:A:108[B]:LEU:HD11	1:A:281:LEU:HB2	1.96	0.47
1:C:63:ARG:HG2	1:C:63:ARG:NH1	2.30	0.47
1:C:281:LEU:C	1:C:281:LEU:HD23	2.39	0.47
1:B:190:LYS:HG2	1:B:199:VAL:HG21	1.96	0.47
1:C:69:ASP:O	1:C:72:THR:HG23	2.15	0.47
1:C:251:LEU:HD12	1:C:251:LEU:N	2.28	0.47
1:A:200:GLU:HG3	1:A:219[B]:LEU:HB3	1.96	0.46
1:B:267:VAL:HB	1:B:270:LEU:HG	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:176:VAL:HG13	1:C:181:GLY:C	2.41	0.46
1:C:42:GLU:O	1:C:42:GLU:HG3	2.14	0.46
1:C:189:ALA:O	1:C:193:ALA:HB2	2.15	0.46
1:B:15:GLU:O	1:B:19[B]:ARG:HG3	2.15	0.46
1:B:169:ILE:CD1	1:B:193:ALA:HB1	2.45	0.46
1:C:63:ARG:HG2	1:C:63:ARG:HH11	1.79	0.46
1:A:93:VAL:HG13	1:A:93:VAL:O	2.16	0.45
1:C:268:GLY:O	1:C:272:HIS:HB2	2.15	0.45
1:A:112:GLN:HE21	1:A:279:LEU:H	1.64	0.45
1:C:219:LEU:CD2	1:C:242:PRO:HB2	2.47	0.45
1:B:61:ALA:HB2	1:B:111[A]:LEU:HG	1.98	0.45
1:B:77:LEU:HD11	1:B:92:ARG:HD2	1.98	0.45
1:B:37[B]:VAL:HG13	1:B:98:ARG:CG	2.47	0.45
1:A:44:GLU:O	1:A:284:VAL:HG22	2.17	0.45
1:B:122:ARG:O	1:B:126:GLU:HG3	2.17	0.45
1:A:64[A]:VAL:HG21	1:A:107:ALA:CA	2.47	0.44
1:B:255:LYS:O	1:B:259:GLU:HG3	2.17	0.44
1:C:56:ALA:HA	1:C:80:GLU:OE2	2.17	0.44
1:A:124:TYR:HB3	1:A:135:ILE:CD1	2.48	0.44
1:C:55:LEU:CD2	1:C:90:VAL:HG13	2.47	0.44
1:C:194:PRO:HB3	1:C:196:TYR:CE2	2.53	0.44
1:A:20:ALA:HA	1:A:23:ARG:NH1	2.33	0.44
1:C:253:ARG:NH1	2:C:350:HOH:O	2.42	0.43
1:A:92:ARG:HB3	1:A:94:ARG:HH12	1.84	0.43
1:B:136:LEU:HD23	1:B:159:ARG:HB2	2.01	0.43
1:B:144:GLY:C	1:B:145:LEU:HD12	2.43	0.43
1:C:186:VAL:O	1:C:190:LYS:HB2	2.19	0.43
1:C:136:LEU:O	1:C:266:SER:HA	2.19	0.43
1:C:206:LEU:HD23	1:C:209:LEU:HD23	2.00	0.43
1:A:11:GLN:HG3	1:A:12:GLY:H	1.83	0.42
1:B:281:LEU:HD23	1:B:281:LEU:C	2.44	0.42
1:C:173:GLU:N	1:C:202:GLU:OE1	2.39	0.42
1:A:181:GLY:HA3	1:A:184:GLU:OE1	2.20	0.42
1:C:54:VAL:HG13	1:C:81:GLY:O	2.18	0.42
1:B:194:PRO:HB3	1:B:196:TYR:CE2	2.55	0.42
1:C:257:ALA:O	1:C:262:VAL:HG22	2.19	0.42
1:C:184:GLU:OE1	1:C:188:ARG:NH2	2.53	0.42
1:B:15:GLU:HB3	2:B:533:HOH:O	2.19	0.41
1:B:37[B]:VAL:HG11	1:B:98:ARG:HG2	2.00	0.41
1:A:228:GLU:CD	1:A:228:GLU:N	2.78	0.41
1:C:30:ASP:O	1:C:34:LEU:HG	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:45:ALA:HB3	1:B:93:VAL:CG1	2.51	0.41
1:C:220:ILE:HG23	1:C:243:LEU:HD23	2.03	0.41
1:C:57:GLY:N	1:C:80:GLU:OE2	2.47	0.41
1:B:237:VAL:HG21	1:B:243:LEU:HD21	2.02	0.41
1:A:221:LEU:HA	1:A:244:GLU:O	2.20	0.40
1:C:232:GLU:O	1:C:235:ARG:HB3	2.21	0.40
1:C:233:ALA:O	1:C:237:VAL:HG23	2.22	0.40
1:A:128:LEU:HG	1:A:258:ALA:HB2	2.02	0.40
1:C:231:ARG:HH11	1:C:231:ARG:HD3	1.62	0.40
1:B:230:LEU:HD21	1:B:262:VAL:CG1	2.49	0.40
1:C:250:THR:CG2	1:C:253:ARG:H	2.33	0.40
1:A:62:GLU:HB2	1:A:74:PHE:CD2	2.57	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	286/286 (100%)	283 (99%)	3 (1%)	0	100	100
1	B	278/286 (97%)	275 (99%)	2 (1%)	1 (0%)	30	22
1	C	276/286 (96%)	261 (95%)	14 (5%)	1 (0%)	30	22
All	All	840/858 (98%)	819 (98%)	19 (2%)	2 (0%)	43	36

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	12	GLY
1	C	127	ALA

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	216/210 (103%)	212 (98%)	4 (2%)	50	47
1	B	208/210 (99%)	207 (100%)	1 (0%)	81	84
1	C	206/210 (98%)	202 (98%)	4 (2%)	50	47
All	All	630/630 (100%)	621 (99%)	9 (1%)	61	59

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	122	ARG
1	A	132[A]	LYS
1	A	132[B]	LYS
1	A	170	LEU
1	B	42	GLU
1	C	94	ARG
1	C	122	ARG
1	C	227	LEU
1	C	231	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	112	GLN
1	A	195	HIS
1	A	248	ASN
1	B	134	GLN
1	B	224	ASN
1	B	248	ASN
1	C	224	ASN

5.3.3 RNA ⓘ

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	277/286 (96%)	0.13	8 (2%) 53 57	13, 21, 40, 59	11 (3%)
1	B	276/286 (96%)	0.19	11 (3%) 42 45	11, 23, 40, 59	4 (1%)
1	C	277/286 (96%)	2.29	157 (56%) 0 0	25, 49, 67, 75	1 (0%)
All	All	830/858 (96%)	0.87	176 (21%) 2 2	11, 29, 61, 75	16 (1%)

All (176) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	230	LEU	6.8
1	C	234	VAL	6.6
1	C	226	PRO	6.5
1	C	10	TRP	6.1
1	C	225	PHE	6.1
1	A	10	TRP	6.0
1	C	84	VAL	5.7
1	C	233	ALA	5.6
1	C	85	ALA	5.5
1	A	11	GLN	5.1
1	C	241	VAL	5.1
1	C	216	GLY	5.0
1	C	127	ALA	4.9
1	C	48	LEU	4.9
1	C	251	LEU	4.9
1	C	262	VAL	4.8
1	C	49	ALA	4.7
1	C	243	LEU	4.7
1	C	245	ALA	4.6
1	C	237	VAL	4.6
1	C	282	LEU	4.5
1	C	228	GLU	4.5
1	C	213	LEU	4.5

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Mol	Chain	Res	Type	RSRZ
1	C	257	ALA	4.5
1	C	203	VAL	4.3
1	C	229	ALA	4.3
1	C	260	ALA	4.3
1	C	82	ALA	4.3
1	C	265	VAL	4.3
1	C	276	ALA	4.1
1	C	250	THR	4.1
1	C	41	LEU	4.0
1	C	55	LEU	3.9
1	C	47	ILE	3.9
1	C	121	THR	3.9
1	C	261	GLY	3.9
1	C	123	ALA	3.9
1	C	217	ALA	3.9
1	B	286	PRO	3.8
1	C	43	GLY	3.8
1	C	90	VAL	3.8
1	C	220	ILE	3.8
1	C	153	VAL	3.7
1	C	255[A]	LYS	3.7
1	C	128	LEU	3.6
1	C	209	LEU	3.6
1	C	254	ALA	3.6
1	C	125	VAL	3.6
1	C	129	ALA	3.6
1	C	212	ALA	3.6
1	C	249	MET	3.6
1	C	286	PRO	3.5
1	C	120	LEU	3.5
1	C	130	GLY	3.5
1	C	78	VAL	3.5
1	C	182	VAL	3.5
1	C	283	VAL	3.5
1	B	284	VAL	3.5
1	C	206	LEU	3.5
1	C	258	ALA	3.3
1	C	256	ALA	3.3
1	C	93	VAL	3.2
1	C	227	LEU	3.2
1	C	177	ARG	3.2
1	C	231	ARG	3.2

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Mol	Chain	Res	Type	RSRZ
1	C	215	ALA	3.2
1	C	74	PHE	3.2
1	C	87	GLY	3.2
1	C	124	TYR	3.2
1	C	36	VAL	3.2
1	C	11	GLN	3.1
1	C	179	ALA	3.1
1	C	176	VAL	3.1
1	C	284	VAL	3.1
1	C	235	ARG	3.1
1	C	271	THR	3.1
1	B	191	ALA	3.1
1	C	219	LEU	3.0
1	C	272	HIS	3.0
1	C	119	THR	3.0
1	C	221	LEU	3.0
1	C	264	TYR	3.0
1	C	122	ARG	3.0
1	C	184	GLU	3.0
1	C	279	LEU	3.0
1	C	118	ALA	3.0
1	C	91	ALA	2.9
1	C	135	ILE	2.9
1	C	46	VAL	2.9
1	C	131	THR	2.9
1	C	181	GLY	2.9
1	C	268	GLY	2.9
1	C	224	ASN	2.9
1	A	28	GLN	2.9
1	C	205	SER	2.9
1	C	246	SER	2.9
1	C	214	GLU	2.9
1	C	232	GLU	2.9
1	C	277	LEU	2.9
1	C	236	ARG	2.8
1	C	275	LYS	2.8
1	B	215	ALA	2.8
1	C	68	ALA	2.8
1	C	59	TRP	2.8
1	C	274	ALA	2.8
1	C	253	ARG	2.8
1	C	185	ALA	2.7

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Mol	Chain	Res	Type	RSRZ
1	C	252	GLU	2.7
1	C	223	ASP	2.6
1	A	59	TRP	2.6
1	C	222	LEU	2.6
1	B	189	ALA	2.6
1	C	45	ALA	2.6
1	C	204	ARG	2.6
1	A	93	VAL	2.6
1	C	39	GLU	2.6
1	C	186	VAL	2.6
1	C	248	ASN	2.6
1	C	218	ASP	2.6
1	C	51	GLU	2.5
1	C	180	GLY	2.5
1	C	50	LYS	2.5
1	C	81	GLY	2.5
1	A	286	PRO	2.5
1	C	242	PRO	2.5
1	C	281	LEU	2.4
1	C	280	SER	2.4
1	C	79	ALA	2.4
1	C	269	ALA	2.4
1	C	76	PRO	2.4
1	C	259	GLU	2.4
1	C	88	THR	2.4
1	C	133	ALA	2.4
1	C	97	LEU	2.4
1	C	263	ASP	2.4
1	C	192	ARG	2.4
1	C	285	ARG	2.3
1	C	42	GLU	2.3
1	B	285	ARG	2.3
1	C	207	GLU	2.3
1	C	72	THR	2.3
1	C	117	ILE	2.3
1	C	156	GLY	2.3
1	C	266	SER	2.3
1	B	192	ARG	2.3
1	C	163	TYR	2.3
1	C	52	GLY	2.3
1	C	267	VAL	2.3
1	C	270	LEU	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	12	GLY	2.2
1	C	12	GLY	2.2
1	C	53	GLY	2.2
1	C	278	ASP	2.2
1	C	201	VAL	2.2
1	C	35	LEU	2.2
1	C	67	LEU	2.2
1	C	73	ALA	2.2
1	C	240	ARG	2.2
1	C	58	LEU	2.2
1	C	77	LEU	2.2
1	C	94	ARG	2.2
1	B	190	LYS	2.1
1	C	208	GLU	2.1
1	C	54	VAL	2.1
1	C	83	ARG	2.1
1	C	132	LYS	2.1
1	C	178	ALA	2.1
1	C	191	ALA	2.1
1	A	42	GLU	2.1
1	C	244	GLU	2.1
1	B	216	GLY	2.1
1	A	48	LEU	2.1
1	C	75	THR	2.0
1	B	11	GLN	2.0
1	C	56	ALA	2.0
1	C	86	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 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.