



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 03:14 AM UTC

PDB ID : 1WY5 / pdb_00001wy5
Title : Crystal structure of isoleucyl-tRNA lysidine synthetase
Authors : Nakanishi, K.; Fukai, S.; Ikeuchi, Y.; Soma, A.; Sekine, Y.; Suzuki, T.; Nureki, O.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)
Deposited on : 2005-02-06
Resolution : 2.42 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

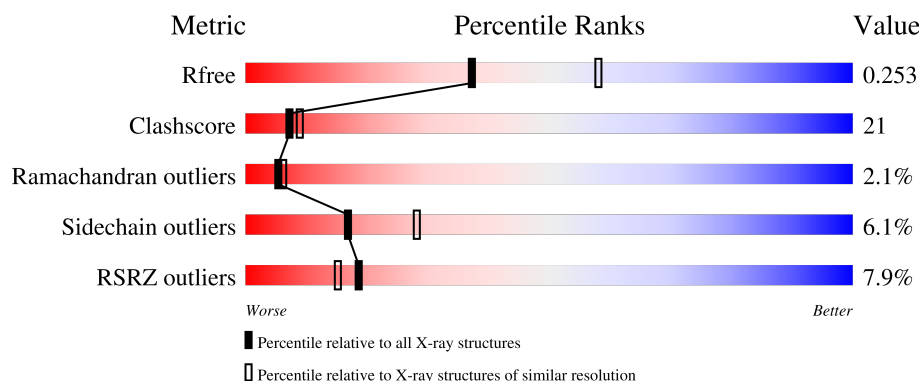
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.42 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	6062 (2.44-2.40)
Clashscore	190562	6562 (2.44-2.40)
Ramachandran outliers	187476	6481 (2.44-2.40)
Sidechain outliers	187428	6482 (2.44-2.40)
RSRZ outliers	180081	6066 (2.44-2.40)

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	317	
1	B	317	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5243 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 Hypothetical UPF0072 protein AQ_1887.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	311	Total	C	N	O	S	0	0	0
			2583	1652	461	462	8			
1	B	311	Total	C	N	O	S	0	0	0
			2583	1652	461	462	8			

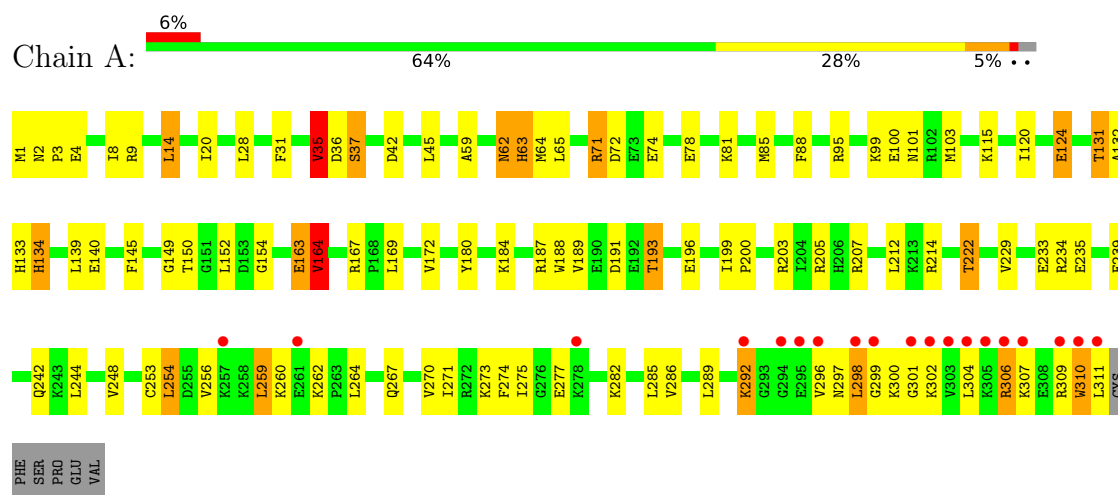
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	46	Total	O	0	0
			46	46		
2	B	31	Total	O	0	0
			31	31		

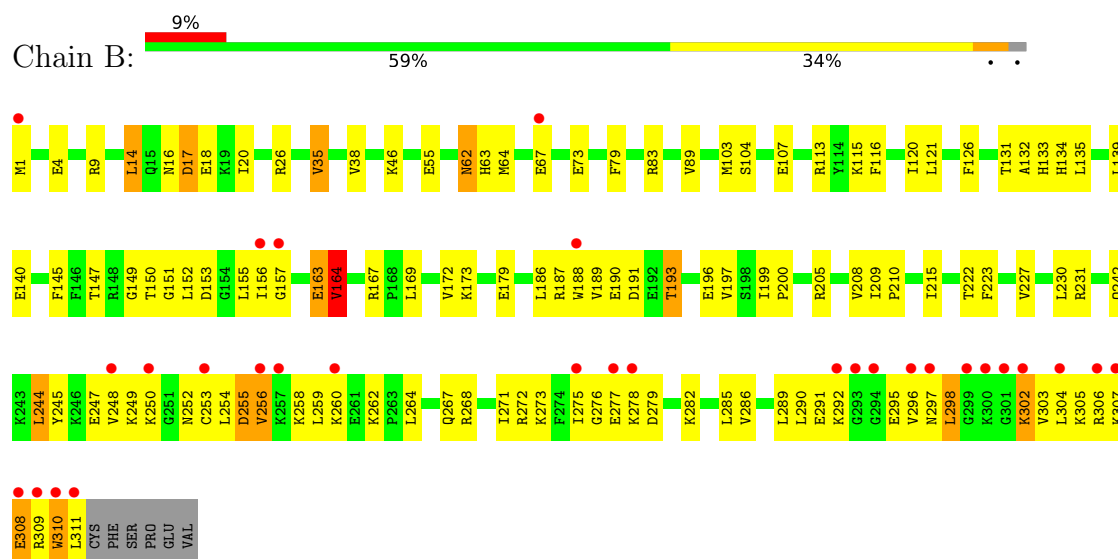
3 Residue-property plots

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: Hypothetical UPF0072 protein AQ_1887



• Molecule 1: Hypothetical UPF0072 protein AQ_1887



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	81.96Å 81.96Å 302.58Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.56 – 2.42 39.56 – 2.42	Depositor EDS
% Data completeness (in resolution range)	98.7 (39.56-2.42) 98.8 (39.56-2.42)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.59 (at 2.42Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.218 , 0.253 0.218 , 0.253	Depositor DCC
R_{free} test set	2829 reflections (7.07%)	wwPDB-VP
Wilson B-factor (Å ²)	43.4	Xtriage
Anisotropy	0.388	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 36.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5243	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.40% 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	0.43	0/2623	0.91	8/3503 (0.2%)
1	B	0.41	0/2623	0.87	6/3503 (0.2%)
All	All	0.42	0/5246	0.89	14/7006 (0.2%)

There are no bond length outliers.

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	167	ARG	CA-C-N	6.92	126.90	119.28
1	A	167	ARG	C-N-CA	6.92	126.90	119.28
1	A	115	LYS	N-CA-C	-6.01	104.64	111.07
1	B	215	ILE	N-CA-C	-5.79	106.58	111.56
1	A	35	VAL	CB-CA-C	-5.71	104.58	111.88
1	B	208	VAL	N-CA-C	5.50	115.59	110.42
1	B	256	VAL	N-CA-C	5.48	116.12	110.36
1	B	115	LYS	N-CA-C	-5.38	105.33	111.14
1	A	310	TRP	N-CA-C	5.30	118.64	110.42
1	B	167	ARG	CA-C-N	5.30	125.61	119.47
1	B	167	ARG	C-N-CA	5.30	125.61	119.47
1	A	298	LEU	N-CA-C	-5.07	105.87	113.89
1	A	134	HIS	N-CA-C	5.02	116.30	109.18
1	A	214	ARG	N-CA-C	-5.01	106.68	112.89

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	2583	0	2683	106	0
1	B	2583	0	2683	118	0
2	A	46	0	0	0	0
2	B	31	0	0	0	0
All	All	5243	0	5366	218	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 21.

All (218) 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:A:191:ASP:OD1	1:A:193:THR:HB	1.59	1.02
1:B:199:ILE:H	1:B:199:ILE:HD12	1.37	0.87
1:A:289:LEU:HD12	1:A:309:ARG:HH21	1.40	0.86
1:A:302:LYS:HG3	1:A:311:LEU:HD13	1.57	0.85
1:A:289:LEU:HD12	1:A:309:ARG:NH2	1.92	0.85
1:B:131:THR:HG22	1:B:133:HIS:H	1.42	0.84
1:B:309:ARG:HG2	1:B:310:TRP:H	1.46	0.81
1:A:289:LEU:HD11	1:A:296:VAL:HG23	1.63	0.80
1:A:286:VAL:HG13	1:A:309:ARG:NH2	1.98	0.79
1:A:292:LYS:HE3	1:A:292:LYS:HA	1.62	0.78
1:B:260:LYS:HE3	1:B:291:GLU:HG2	1.64	0.77
1:B:62:ASN:ND2	1:B:64:MET:H	1.84	0.76
1:B:302:LYS:NZ	1:B:302:LYS:HB3	2.01	0.76
1:B:79:PHE:O	1:B:83:ARG:HD2	1.89	0.73
1:B:62:ASN:C	1:B:62:ASN:HD22	1.96	0.72
1:A:163:GLU:O	1:A:164:VAL:HB	1.88	0.72
1:A:140:GLU:OE2	1:A:205:ARG:HD3	1.90	0.72
1:A:298:LEU:HD11	1:A:302:LYS:HZ2	1.55	0.71
1:A:297:ASN:ND2	1:A:299:GLY:H	1.88	0.70
1:B:121:LEU:HD12	1:B:126:PHE:HB2	1.74	0.70
1:B:191:ASP:OD1	1:B:193:THR:HB	1.91	0.70
1:B:135:LEU:HD23	1:B:173:LYS:HG2	1.74	0.69
1:A:62:ASN:C	1:A:62:ASN:HD22	2.00	0.69
1:B:295:GLU:HG2	1:B:303:VAL:HG13	1.75	0.69
1:A:62:ASN:ND2	1:A:64:MET:H	1.92	0.68
1:B:16:ASN:O	1:B:17:ASP:HB2	1.92	0.67
1:A:235:GLU:O	1:A:239:GLU:HG3	1.94	0.67
1:A:302:LYS:HG3	1:A:311:LEU:CD1	2.25	0.67
1:A:310:TRP:CG	1:A:311:LEU:N	2.64	0.66
1:A:62:ASN:HD22	1:A:64:MET:H	1.45	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:131:THR:HG22	1:A:133:HIS:H	1.61	0.64
1:B:156:ILE:HD11	1:B:231:ARG:HD2	1.78	0.64
1:A:286:VAL:HG21	1:A:302:LYS:NZ	2.14	0.63
1:A:100:GLU:O	1:A:100:GLU:HG2	1.97	0.63
1:A:244:LEU:HD22	1:A:270:VAL:HG21	1.79	0.63
1:B:305:LYS:C	1:B:307:LYS:H	2.04	0.63
1:A:311:LEU:O	1:A:311:LEU:HD23	1.99	0.62
1:A:256:VAL:HG21	1:A:304:LEU:HD21	1.79	0.62
1:A:253:CYS:HA	1:A:309:ARG:O	1.99	0.62
1:B:305:LYS:HZ1	1:B:309:ARG:HD3	1.65	0.61
1:A:271:ILE:O	1:A:275:ILE:HG12	2.01	0.61
1:A:286:VAL:HG21	1:A:302:LYS:HZ3	1.66	0.61
1:B:244:LEU:HD11	1:B:262:LYS:HG3	1.81	0.61
1:B:199:ILE:HD12	1:B:199:ILE:N	2.13	0.60
1:A:88:PHE:CZ	1:A:124:GLU:HG2	2.36	0.60
1:A:302:LYS:CG	1:A:311:LEU:HD13	2.30	0.60
1:B:254:LEU:H	1:B:308:GLU:HG2	1.66	0.60
1:B:247:GLU:O	1:B:250:LYS:HE3	2.01	0.59
1:B:249:LYS:NZ	1:B:252:ASN:HA	2.18	0.58
1:A:306:ARG:O	1:A:307:LYS:HB3	2.03	0.58
1:A:187:ARG:HG2	1:A:187:ARG:HH11	1.69	0.58
1:A:35:VAL:HG13	1:A:188:TRP:CD2	2.39	0.58
1:B:73:GLU:OE2	1:B:89:VAL:HG11	2.04	0.57
1:B:199:ILE:H	1:B:199:ILE:CD1	2.11	0.57
1:B:121:LEU:HD12	1:B:126:PHE:CB	2.33	0.57
1:B:156:ILE:HD11	1:B:231:ARG:CD	2.34	0.57
1:A:163:GLU:O	1:A:164:VAL:CB	2.53	0.57
1:B:62:ASN:HD22	1:B:64:MET:H	1.49	0.57
1:B:253:CYS:HB2	1:B:308:GLU:HB3	1.87	0.57
1:A:42:ASP:OD1	1:A:184:LYS:HE2	2.04	0.57
1:A:222:THR:HG23	1:B:149:GLY:CA	2.35	0.57
1:A:297:ASN:HD22	1:A:298:LEU:N	2.02	0.57
1:B:259:LEU:HD23	1:B:259:LEU:O	2.05	0.57
1:A:300:LYS:HD2	1:A:311:LEU:HD21	1.86	0.56
1:B:62:ASN:HD22	1:B:63:HIS:N	2.03	0.56
1:A:300:LYS:HB2	1:A:311:LEU:HD11	1.86	0.56
1:B:254:LEU:H	1:B:308:GLU:HB3	1.70	0.56
1:B:272:ARG:HD3	1:B:278:LYS:HD3	1.88	0.56
1:B:302:LYS:HB3	1:B:302:LYS:HZ3	1.69	0.55
1:A:310:TRP:CD1	1:A:311:LEU:H	2.24	0.55
1:A:302:LYS:HE3	1:A:311:LEU:HD13	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:254:LEU:H	1:B:308:GLU:CB	2.20	0.55
1:A:298:LEU:HD11	1:A:302:LYS:NZ	2.21	0.55
1:A:28:LEU:HD11	1:A:59:ALA:HB2	1.89	0.55
1:B:145:PHE:HB3	1:B:150:THR:HG21	1.87	0.55
1:B:156:ILE:CD1	1:B:231:ARG:HD2	2.36	0.55
1:B:249:LYS:HZ2	1:B:252:ASN:HA	1.71	0.55
1:A:297:ASN:HD22	1:A:299:GLY:H	1.54	0.55
1:A:1:MET:HE1	1:A:9:ARG:CZ	2.37	0.54
1:B:116:PHE:CE1	1:B:120:ILE:HD11	2.42	0.54
1:B:254:LEU:H	1:B:308:GLU:CG	2.20	0.54
1:A:36:ASP:HB3	1:A:132:ALA:HB1	1.88	0.54
1:B:156:ILE:HG23	1:B:157:GLY:N	2.21	0.54
1:A:222:THR:HG23	1:B:149:GLY:HA2	1.90	0.54
1:B:305:LYS:NZ	1:B:309:ARG:HD3	2.22	0.54
1:A:256:VAL:HG21	1:A:304:LEU:CD2	2.38	0.53
1:B:259:LEU:CD2	1:B:267:GLN:HG2	2.38	0.53
1:A:244:LEU:O	1:A:244:LEU:HD23	2.08	0.53
1:A:212:LEU:HD22	1:B:147:THR:OG1	2.08	0.53
1:A:310:TRP:O	1:A:311:LEU:HB2	2.08	0.53
1:A:101:ASN:HB3	1:A:103:MET:HE3	1.90	0.53
1:A:134:HIS:HB2	1:A:172:VAL:O	2.09	0.53
1:A:31:PHE:HA	1:A:37:SER:OG	2.09	0.53
1:B:163:GLU:HG3	1:B:163:GLU:O	2.09	0.53
1:B:289:LEU:HD21	1:B:296:VAL:HG23	1.91	0.53
1:A:45:LEU:HD21	1:A:85:MET:HE3	1.91	0.53
1:B:4:GLU:OE2	1:B:46:LYS:HE2	2.09	0.52
1:B:163:GLU:O	1:B:164:VAL:HB	2.10	0.52
1:B:116:PHE:CZ	1:B:120:ILE:HD11	2.44	0.52
1:B:259:LEU:HD22	1:B:290:LEU:HD11	1.92	0.52
1:A:65:LEU:HD11	1:A:196:GLU:HG2	1.90	0.52
1:A:277:GLU:HG2	1:A:282:LYS:HD2	1.91	0.51
1:A:131:THR:CG2	1:A:133:HIS:H	2.23	0.51
1:A:285:LEU:HD12	1:A:296:VAL:CG1	2.39	0.51
1:A:306:ARG:O	1:A:306:ARG:HG3	2.09	0.51
1:B:255:ASP:CG	1:B:258:LYS:HD3	2.35	0.51
1:A:300:LYS:HG3	1:A:311:LEU:HD11	1.91	0.51
1:A:302:LYS:CE	1:A:311:LEU:HD13	2.41	0.51
1:B:197:VAL:HG21	1:B:205:ARG:HH22	1.76	0.51
1:B:256:VAL:HG21	1:B:304:LEU:CD2	2.41	0.51
1:B:259:LEU:HD23	1:B:267:GLN:HG2	1.91	0.51
1:A:163:GLU:O	1:A:163:GLU:HG3	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:103:MET:HB2	1:B:107:GLU:HG3	1.93	0.51
1:B:305:LYS:HE3	1:B:309:ARG:HB2	1.93	0.51
1:A:301:GLY:C	1:A:311:LEU:HD12	2.37	0.50
1:B:309:ARG:CG	1:B:310:TRP:H	2.21	0.50
1:A:199:ILE:HB	1:A:200:PRO:HD3	1.93	0.50
1:B:63:HIS:HE1	1:B:196:GLU:OE2	1.95	0.50
1:B:150:THR:HG23	1:B:151:GLY:N	2.27	0.50
1:B:197:VAL:HG21	1:B:205:ARG:NH2	2.26	0.50
1:A:149:GLY:HA2	1:B:222:THR:HG23	1.94	0.49
1:B:256:VAL:HG22	1:B:290:LEU:HD23	1.94	0.49
1:B:307:LYS:O	1:B:308:GLU:HB2	2.12	0.49
1:A:101:ASN:CB	1:A:103:MET:HE3	2.41	0.49
1:B:302:LYS:HB3	1:B:302:LYS:HZ2	1.77	0.49
1:A:2:ASN:HB2	1:A:3:PRO:CD	2.41	0.49
1:A:2:ASN:HB2	1:A:3:PRO:HD2	1.93	0.49
1:B:140:GLU:OE2	1:B:205:ARG:HD3	2.12	0.49
1:B:309:ARG:NH2	1:B:311:LEU:HB3	2.28	0.49
1:A:95:ARG:HG2	1:A:99:LYS:HE3	1.95	0.49
1:A:254:LEU:HD22	1:A:274:PHE:CD2	2.48	0.49
1:B:309:ARG:HG2	1:B:310:TRP:N	2.23	0.49
1:A:14:LEU:HD22	1:A:20:ILE:HD11	1.93	0.48
1:A:286:VAL:O	1:A:309:ARG:NH1	2.44	0.48
1:A:300:LYS:CG	1:A:311:LEU:HD11	2.43	0.48
1:B:121:LEU:HD21	1:B:164:VAL:HG22	1.96	0.48
1:B:189:VAL:HG22	1:B:190:GLU:H	1.78	0.48
1:B:199:ILE:HB	1:B:200:PRO:CD	2.44	0.48
1:B:153:ASP:O	1:B:156:ILE:HG22	2.14	0.48
1:B:275:ILE:O	1:B:277:GLU:N	2.43	0.48
1:A:180:TYR:CE2	1:A:184:LYS:HD3	2.49	0.47
1:B:242:GLN:HG2	1:B:273:LYS:NZ	2.30	0.47
1:A:244:LEU:HD11	1:A:262:LYS:HG3	1.95	0.47
1:B:152:LEU:HD11	1:B:230:LEU:HD22	1.96	0.47
1:B:305:LYS:C	1:B:307:LYS:N	2.71	0.47
1:A:242:GLN:OE1	1:A:273:LYS:NZ	2.43	0.47
1:B:18:GLU:CD	1:B:231:ARG:HH22	2.23	0.47
1:B:259:LEU:HD23	1:B:259:LEU:C	2.40	0.47
1:A:289:LEU:HB2	1:A:309:ARG:CZ	2.45	0.46
1:A:300:LYS:CB	1:A:311:LEU:HD11	2.45	0.46
1:B:227:VAL:O	1:B:231:ARG:HB2	2.15	0.46
1:B:295:GLU:HG2	1:B:303:VAL:HG22	1.96	0.46
1:A:150:THR:OG1	1:A:154:GLY:HA3	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:156:ILE:CG2	1:B:157:GLY:N	2.78	0.46
1:B:191:ASP:CG	1:B:193:THR:HB	2.40	0.46
1:A:62:ASN:C	1:A:62:ASN:ND2	2.72	0.46
1:B:296:VAL:HG12	1:B:297:ASN:N	2.30	0.46
1:A:286:VAL:HA	1:A:309:ARG:HH22	1.80	0.46
1:A:149:GLY:HA2	1:B:222:THR:CG2	2.47	0.45
1:A:285:LEU:HD12	1:A:296:VAL:HG13	1.97	0.45
1:A:286:VAL:HG13	1:A:309:ARG:CZ	2.45	0.45
1:B:1:MET:SD	1:B:9:ARG:CZ	3.05	0.45
1:B:277:GLU:OE1	1:B:282:LYS:HE3	2.17	0.45
1:A:78:GLU:HA	1:A:81:LYS:HE3	1.98	0.45
1:A:297:ASN:ND2	1:A:299:GLY:N	2.63	0.44
1:B:134:HIS:HB2	1:B:172:VAL:O	2.17	0.44
1:A:120:ILE:O	1:A:124:GLU:HB2	2.16	0.44
1:A:63:HIS:N	1:A:63:HIS:CD2	2.85	0.44
1:A:74:GLU:O	1:A:78:GLU:HG3	2.17	0.44
1:A:248:VAL:HG11	1:A:259:LEU:HG	2.00	0.44
1:A:285:LEU:HD13	1:A:285:LEU:O	2.18	0.44
1:A:302:LYS:HA	1:A:310:TRP:O	2.17	0.44
1:A:260:LYS:HA	1:A:267:GLN:NE2	2.32	0.43
1:A:310:TRP:CD1	1:A:311:LEU:N	2.87	0.43
1:B:14:LEU:HD13	1:B:20:ILE:HG12	1.99	0.43
1:B:282:LYS:HD2	1:B:298:LEU:HD22	2.00	0.43
1:B:285:LEU:HG	1:B:296:VAL:HG11	1.99	0.43
1:B:62:ASN:ND2	1:B:64:MET:N	2.61	0.43
1:B:131:THR:CG2	1:B:132:ALA:N	2.81	0.43
1:B:62:ASN:ND2	1:B:62:ASN:C	2.67	0.43
1:A:229:VAL:O	1:A:233:GLU:HG3	2.19	0.43
1:B:104:SER:OG	1:B:107:GLU:HG2	2.18	0.43
1:B:1:MET:SD	1:B:9:ARG:NH2	2.92	0.43
1:B:35:VAL:HG13	1:B:188:TRP:CD2	2.53	0.43
1:B:295:GLU:CG	1:B:303:VAL:HG13	2.46	0.43
1:A:149:GLY:CA	1:B:222:THR:HG23	2.49	0.43
1:B:151:GLY:O	1:B:155:LEU:HG	2.19	0.43
1:B:152:LEU:CD1	1:B:230:LEU:HD22	2.49	0.43
1:B:223:PHE:O	1:B:227:VAL:HG23	2.19	0.43
1:B:259:LEU:HD22	1:B:290:LEU:CD1	2.49	0.43
1:A:297:ASN:HD22	1:A:298:LEU:H	1.66	0.42
1:B:103:MET:HB2	1:B:107:GLU:CG	2.49	0.42
1:B:248:VAL:HG13	1:B:255:ASP:N	2.34	0.42
1:B:286:VAL:HG22	1:B:302:LYS:HD2	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:209:ILE:HB	1:B:210:PRO:HD3	2.01	0.42
1:B:4:GLU:CD	1:B:46:LYS:HE2	2.45	0.42
1:B:268:ARG:HB3	1:B:272:ARG:HH12	1.84	0.42
1:B:38:VAL:HG11	1:B:186:LEU:HB3	2.02	0.42
1:B:253:CYS:HB2	1:B:308:GLU:C	2.44	0.41
1:B:295:GLU:OE1	1:B:303:VAL:HG22	2.19	0.41
1:A:286:VAL:CA	1:A:309:ARG:HH22	2.33	0.41
1:A:311:LEU:HD23	1:A:311:LEU:C	2.45	0.41
1:A:310:TRP:O	1:A:311:LEU:CB	2.67	0.41
1:B:277:GLU:HG2	1:B:279:ASP:OD2	2.21	0.41
1:A:4:GLU:HA	1:A:180:TYR:CD1	2.55	0.41
1:A:35:VAL:HG13	1:A:188:TRP:CE2	2.56	0.41
1:A:71:ARG:NH1	1:A:72:ASP:OD1	2.50	0.41
1:B:245:TYR:O	1:B:249:LYS:HB3	2.21	0.41
1:A:248:VAL:O	1:A:248:VAL:HG12	2.20	0.41
1:B:16:ASN:O	1:B:17:ASP:CB	2.64	0.41
1:B:305:LYS:O	1:B:306:ARG:HB2	2.20	0.41
1:A:145:PHE:HB3	1:A:150:THR:HB	2.03	0.40
1:A:203:ARG:O	1:A:207:ARG:HB2	2.22	0.40
1:B:256:VAL:HG23	1:B:307:LYS:O	2.22	0.40
1:B:271:ILE:O	1:B:275:ILE:HG12	2.21	0.40
1:A:285:LEU:HD13	1:A:285:LEU:C	2.47	0.40
1:B:26:ARG:NH1	1:B:55:GLU:OE1	2.55	0.40
1:B:279:ASP:HB2	1:B:282:LYS:HG2	2.04	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	309/317 (98%)	297 (96%)	9 (3%)	3 (1%)	12 18

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	309/317 (98%)	281 (91%)	18 (6%)	10 (3%)	3	2
All	All	618/634 (98%)	578 (94%)	27 (4%)	13 (2%)	5	6

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	164	VAL
1	A	306	ARG
1	B	17	ASP
1	B	163	GLU
1	A	163	GLU
1	B	164	VAL
1	B	310	TRP
1	B	292	LYS
1	B	67	GLU
1	B	276	GLY
1	B	298	LEU
1	B	308	GLU
1	B	255	ASP

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	279/285 (98%)	258 (92%)	21 (8%)	12	20
1	B	279/285 (98%)	266 (95%)	13 (5%)	23	39
All	All	558/570 (98%)	524 (94%)	34 (6%)	17	28

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	ILE
1	A	14	LEU
1	A	35	VAL

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Mol	Chain	Res	Type
1	A	37	SER
1	A	62	ASN
1	A	63	HIS
1	A	71	ARG
1	A	124	GLU
1	A	131	THR
1	A	139	LEU
1	A	152	LEU
1	A	164	VAL
1	A	169	LEU
1	A	189	VAL
1	A	193	THR
1	A	222	THR
1	A	234	ARG
1	A	254	LEU
1	A	259	LEU
1	A	264	LEU
1	A	292	LYS
1	B	14	LEU
1	B	35	VAL
1	B	62	ASN
1	B	113	ARG
1	B	139	LEU
1	B	164	VAL
1	B	169	LEU
1	B	179	GLU
1	B	187	ARG
1	B	193	THR
1	B	244	LEU
1	B	264	LEU
1	B	302	LYS

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

Mol	Chain	Res	Type
1	A	62	ASN
1	A	134	HIS
1	A	206	HIS
1	A	267	GLN
1	A	297	ASN
1	B	62	ASN
1	B	63	HIS

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Mol	Chain	Res	Type
1	B	267	GLN
1	B	297	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 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	311/317 (98%)	0.21	19 (6%) 27 23	27, 44, 91, 114	0
1	B	311/317 (98%)	0.42	30 (9%) 13 10	28, 48, 101, 127	0
All	All	622/634 (98%)	0.32	49 (7%) 18 15	27, 46, 97, 127	0

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	311	LEU	7.6
1	A	309	ARG	7.1
1	A	310	TRP	6.8
1	A	311	LEU	6.6
1	B	304	LEU	5.3
1	B	310	TRP	5.2
1	A	303	VAL	4.9
1	A	299	GLY	4.7
1	B	156	ILE	4.5
1	B	307	LYS	4.4
1	A	292	LYS	4.3
1	B	299	GLY	4.0
1	A	304	LEU	3.8
1	B	309	ARG	3.8
1	B	302	LYS	3.8
1	A	302	LYS	3.7
1	B	308	GLU	3.4
1	B	296	VAL	3.4
1	B	250	LYS	3.3
1	B	248	VAL	3.1
1	B	292	LYS	3.0
1	A	294	GLY	3.0
1	A	306	ARG	2.9
1	A	261	GLU	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	298	LEU	2.9
1	B	157	GLY	2.8
1	A	278	LYS	2.8
1	B	300	LYS	2.8
1	B	294	GLY	2.7
1	B	278	LYS	2.7
1	B	67	GLU	2.7
1	B	293	GLY	2.7
1	B	257	LYS	2.6
1	B	297	ASN	2.5
1	B	253	CYS	2.5
1	B	301	GLY	2.5
1	A	301	GLY	2.4
1	B	1	MET	2.4
1	B	275	ILE	2.3
1	A	295	GLU	2.3
1	A	307	LYS	2.3
1	B	256	VAL	2.2
1	B	277	GLU	2.2
1	A	305	LYS	2.2
1	B	306	ARG	2.2
1	B	188	TRP	2.2
1	B	260	LYS	2.1
1	A	296	VAL	2.1
1	A	257	LYS	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 ⓘ

There are no such residues in this entry.