



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 07:59 AM UTC

PDB ID : 3U1H / pdb_00003u1h
Title : Crystal structure of IPMDH from the last common ancestor of Bacillus
Authors : Haaning, S.; Hobbs, J.K.; Monk, C.R.; Arcus, V.L.
Deposited on : 2011-09-29
Resolution : 2.80 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

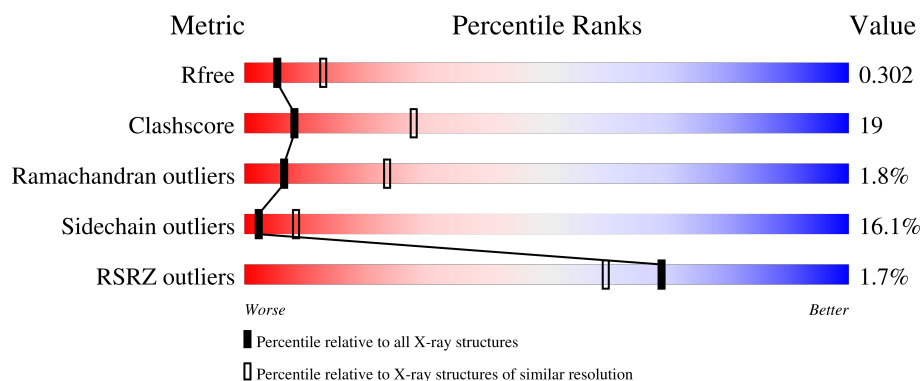
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	390	2% 56% 28% 7% 9%
1	B	390	2% 52% 32% 8% 8%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5228 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 3-isopropylmalate dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	355	Total	C	N	O	S	0	0	0
			2589	1645	428	506	10			
1	B	357	Total	C	N	O	S	0	0	0
			2623	1663	432	518	10			

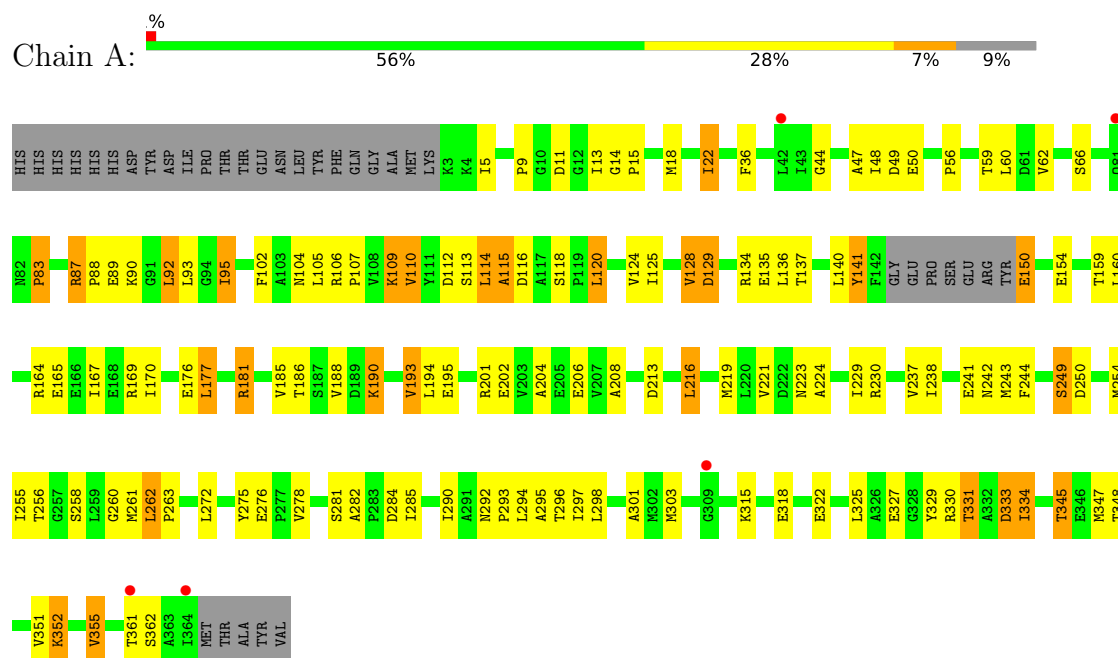
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	7	Total	O	0	0
			7	7		
2	B	9	Total	O	0	0
			9	9		

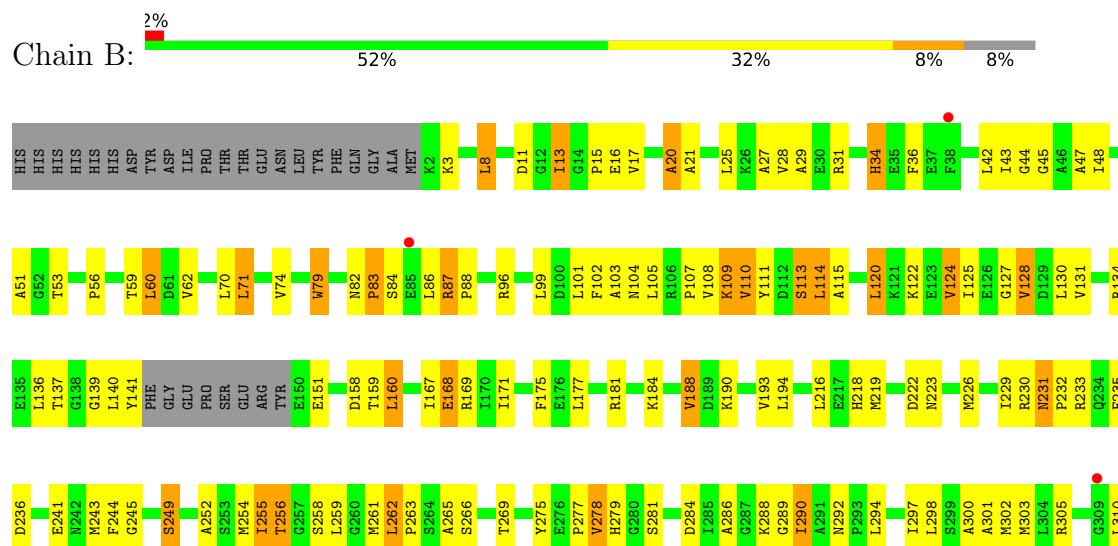
3 Residue-property plots

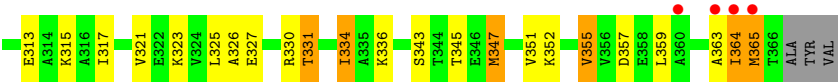
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: 3-isopropylmalate dehydrogenase



• Molecule 1: 3-isopropylmalate dehydrogenase





4 Data and refinement statistics

Property	Value	Source
Space group	P 2 ₁ 2 ₁ 2 ₁	Depositor
Cell constants a, b, c, α , β , γ	71.27Å 75.98Å 171.22Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.43 – 2.80 44.43 – 2.80	Depositor EDS
% Data completeness (in resolution range)	94.5 (44.43-2.80) 94.4 (44.43-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.27 (at 2.81Å)	Xtriage
Refinement program	REFMAC refmac_5.5.0109	Depositor
R, R_{free}	0.233 , 0.309 0.229 , 0.302	Depositor DCC
R_{free} test set	1123 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	65.8	Xtriage
Anisotropy	0.074	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 48.4	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	5228	wwPDB-VP
Average B, all atoms (Å ²)	65.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.75% 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	0.82	0/2628	1.07	6/3572 (0.2%)
1	B	0.83	0/2661	1.07	7/3613 (0.2%)
All	All	0.82	0/5289	1.07	13/7185 (0.2%)

There are no bond length outliers.

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	262	LEU	N-CA-C	6.76	118.31	109.93
1	B	231	ASN	CA-C-N	6.12	126.56	119.47
1	B	231	ASN	C-N-CA	6.12	126.56	119.47
1	B	336	LYS	CA-C-N	-5.96	113.82	120.14
1	B	336	LYS	C-N-CA	-5.96	113.82	120.14
1	B	177	LEU	N-CA-C	5.48	116.93	111.07
1	A	241	GLU	N-CA-C	5.41	117.06	110.41
1	A	93	LEU	N-CA-C	-5.35	102.13	110.10
1	B	188	VAL	N-CA-C	5.32	115.24	107.37
1	B	262	LEU	N-CA-C	5.28	118.74	107.91
1	A	276	GLU	CA-C-N	5.27	125.03	119.76
1	A	276	GLU	C-N-CA	5.27	125.03	119.76
1	A	177	LEU	N-CA-C	5.07	116.80	111.28

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	2589	0	2536	104	0
1	B	2623	0	2591	112	0
2	A	7	0	0	1	0
2	B	9	0	0	1	0
All	All	5228	0	5127	199	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 19.

All (199) 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:139:GLY:HA2	1:B:160:LEU:HD23	1.46	0.98
1:B:256:THR:HG21	1:B:262:LEU:HD21	1.44	0.96
1:A:361:THR:HG23	1:A:362:SER:H	1.29	0.95
1:A:261:MET:CE	1:A:292:ASN:HD21	1.82	0.93
1:B:261:MET:HE1	1:B:294:LEU:HD12	1.53	0.89
1:B:140:LEU:HD12	1:B:160:LEU:HG	1.59	0.85
1:B:261:MET:HE3	1:B:292:ASN:HD21	1.44	0.83
1:A:243:MET:HE1	1:B:190:LYS:HZ2	1.44	0.82
1:B:110:VAL:HG11	1:B:125:ILE:O	1.83	0.79
1:A:361:THR:HG23	1:A:362:SER:N	1.96	0.78
1:A:361:THR:CG2	1:A:362:SER:H	1.97	0.77
1:B:261:MET:CE	1:B:292:ASN:HD21	1.97	0.77
1:B:347:MET:O	1:B:351:VAL:HG23	1.86	0.76
1:A:261:MET:HE3	1:A:292:ASN:HD21	1.48	0.76
1:A:243:MET:HE1	1:B:190:LYS:NZ	2.00	0.76
1:A:110:VAL:CG1	1:A:128:VAL:HG22	2.17	0.75
1:A:230:ARG:HH21	1:B:254:MET:HE1	1.52	0.74
1:B:140:LEU:CD1	1:B:160:LEU:HG	2.17	0.74
1:B:139:GLY:HA2	1:B:160:LEU:CD2	2.17	0.72
1:B:252:ALA:O	1:B:255:ILE:HG12	1.89	0.71
1:A:110:VAL:HG12	1:A:128:VAL:HG22	1.73	0.70
1:B:102:PHE:CE1	1:B:169:ARG:HG2	2.25	0.70
1:B:104:ASN:HB2	1:B:136:LEU:HD11	1.74	0.69
1:A:104:ASN:HB3	1:A:134:ARG:HB3	1.75	0.69
1:A:106:ARG:NH2	1:A:249:SER:HB3	2.08	0.69
1:B:45:GLY:HA3	1:B:79:TRP:CZ2	2.28	0.69
1:B:352:LYS:O	1:B:355:VAL:HG22	1.92	0.68
1:B:108:VAL:HG13	1:B:256:THR:HG22	1.74	0.68
1:B:114:LEU:HD13	1:B:330:ARG:HH21	1.60	0.67
1:A:254:MET:CE	1:B:230:ARG:HB2	2.25	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:87:ARG:N	1:A:88:PRO:HD3	2.10	0.66
1:A:250:ASP:OD2	1:B:222:ASP:HB2	1.96	0.66
1:A:18:MET:O	1:A:22:ILE:HG12	1.96	0.65
1:A:59:THR:O	1:A:62:VAL:HG12	1.97	0.65
1:A:83:PRO:HB2	1:A:88:PRO:HG3	1.78	0.64
1:B:104:ASN:HB3	1:B:134:ARG:HB3	1.78	0.64
1:B:127:GLY:H	1:B:233:ARG:HH11	1.46	0.63
1:A:9:PRO:HB3	1:A:15:PRO:HA	1.78	0.63
1:A:190:LYS:HG2	1:A:221:VAL:HG12	1.82	0.62
1:B:275:TYR:CD1	1:B:303:MET:HG3	2.35	0.61
1:B:313:GLU:N	1:B:313:GLU:OE2	2.32	0.61
1:A:254:MET:HE3	1:B:230:ARG:HB2	1.82	0.61
1:A:56:PRO:HB2	1:A:59:THR:HG23	1.83	0.60
1:B:298:LEU:O	1:B:301:ALA:HB3	2.00	0.60
1:B:29:ALA:HB1	1:B:34:HIS:O	2.01	0.60
1:A:160:LEU:HB3	1:B:158:ASP:HB2	1.84	0.59
1:B:59:THR:O	1:B:62:VAL:HG12	2.01	0.59
1:A:90:LYS:HA	1:A:92:LEU:H	1.66	0.59
1:B:110:VAL:HG12	1:B:128:VAL:HG22	1.85	0.59
1:A:258:SER:C	1:A:260:GLY:H	2.10	0.58
1:B:87:ARG:N	1:B:88:PRO:HD3	2.18	0.58
1:A:102:PHE:CE1	1:A:169:ARG:HG2	2.39	0.58
1:A:331:THR:H	1:A:334:ILE:HD11	1.68	0.58
1:A:190:LYS:HD2	1:A:193:VAL:HG12	1.85	0.57
1:A:329:TYR:O	1:A:330:ARG:HD2	2.04	0.57
1:A:66:SER:O	1:A:272:LEU:HD11	2.05	0.56
1:B:59:THR:O	1:B:62:VAL:N	2.38	0.56
1:B:82:ASN:N	1:B:83:PRO:HD3	2.21	0.56
1:A:87:ARG:H	1:A:88:PRO:HD3	1.71	0.56
1:A:118:SER:HB2	1:A:255:ILE:O	2.06	0.56
1:A:261:MET:HE1	1:A:294:LEU:HD12	1.87	0.55
1:A:275:TYR:CZ	1:A:303:MET:HA	2.42	0.55
1:B:364:ILE:HG12	1:B:364:ILE:O	2.07	0.55
1:A:290:ILE:C	1:A:290:ILE:HD12	2.31	0.54
1:A:258:SER:C	1:A:260:GLY:N	2.64	0.54
1:B:45:GLY:HA2	1:B:48:ILE:HD12	1.89	0.54
1:A:202:GLU:CD	1:B:151:GLU:HG2	2.32	0.54
1:A:281:SER:O	1:A:282:ALA:C	2.51	0.54
1:A:293:PRO:O	1:A:297:ILE:HG13	2.08	0.54
1:B:277:PRO:HB2	1:B:279:HIS:CD2	2.42	0.54
1:B:113:SER:HB2	1:B:114:LEU:HD22	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:130:LEU:CD2	1:B:235:PHE:HB2	2.39	0.53
1:A:150:GLU:HG2	1:A:150:GLU:O	2.09	0.53
1:B:130:LEU:HD12	1:B:130:LEU:O	2.08	0.53
1:B:114:LEU:HD13	1:B:330:ARG:NH2	2.25	0.52
1:A:243:MET:HE3	1:A:243:MET:HA	1.90	0.52
1:B:87:ARG:H	1:B:88:PRO:HD3	1.74	0.52
1:B:13:ILE:HG23	1:B:286:ALA:HA	1.91	0.52
1:A:352:LYS:O	1:A:355:VAL:HG22	2.10	0.51
1:B:103:ALA:O	1:B:266:SER:HA	2.11	0.51
1:A:292:ASN:HB2	1:A:333:ASP:OD1	2.10	0.51
1:B:113:SER:HB3	1:B:330:ARG:HH22	1.75	0.51
1:A:83:PRO:CB	1:A:88:PRO:HG3	2.40	0.51
1:B:45:GLY:HA3	1:B:79:TRP:CH2	2.45	0.51
1:B:114:LEU:HD21	1:B:325:LEU:HD22	1.93	0.51
1:A:44:GLY:O	1:A:47:ALA:N	2.43	0.50
1:B:262:LEU:O	1:B:278:VAL:HG12	2.10	0.50
1:A:164:ARG:NH2	1:A:206:GLU:OE2	2.45	0.50
1:A:194:LEU:HD21	1:B:140:LEU:HD23	1.93	0.50
1:A:140:LEU:HD23	1:B:194:LEU:HD21	1.93	0.50
1:B:130:LEU:HD22	1:B:235:PHE:HB2	1.94	0.50
1:B:261:MET:HA	1:B:261:MET:HE2	1.93	0.49
1:B:223:ASN:O	1:B:226:MET:N	2.45	0.49
1:B:323:LYS:O	1:B:326:ALA:N	2.44	0.49
1:A:188:VAL:HG22	1:A:219:MET:HB3	1.95	0.49
1:B:109:LYS:HB3	2:B:372:HOH:O	2.13	0.49
1:A:177:LEU:O	1:A:181:ARG:HG3	2.13	0.49
1:B:59:THR:O	1:B:60:LEU:C	2.56	0.49
1:B:107:PRO:HA	1:B:131:VAL:HG22	1.95	0.49
1:A:188:VAL:HA	1:A:219:MET:O	2.13	0.48
1:B:130:LEU:HD12	1:B:130:LEU:C	2.39	0.48
1:B:231:ASN:N	1:B:232:PRO:CD	2.77	0.48
1:A:254:MET:HE1	1:B:230:ARG:HE	1.79	0.48
1:B:222:ASP:N	1:B:222:ASP:OD1	2.47	0.48
1:B:262:LEU:HB2	1:B:278:VAL:CG1	2.44	0.47
1:B:140:LEU:HD12	1:B:160:LEU:CG	2.39	0.47
1:A:230:ARG:HB2	1:B:254:MET:HE2	1.96	0.47
1:B:25:LEU:HD13	1:B:300:ALA:HB1	1.96	0.47
1:B:134:ARG:HG3	1:B:245:GLY:HA3	1.96	0.47
1:A:114:LEU:C	1:A:116:ASP:H	2.22	0.47
1:A:89:GLU:HA	2:A:371:HOH:O	2.14	0.47
1:A:193:VAL:HG22	1:B:141:TYR:HB2	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:47:ALA:HB1	1:B:56:PRO:HD3	1.96	0.47
1:A:105:LEU:HD21	1:A:177:LEU:HD21	1.96	0.47
1:B:363:ALA:C	1:B:365:MET:N	2.73	0.47
1:B:83:PRO:O	1:B:84:SER:HB2	2.15	0.47
1:A:109:LYS:HB2	1:A:129:ASP:HB2	1.96	0.47
1:A:221:VAL:HG21	1:A:244:PHE:CG	2.50	0.46
1:B:21:ALA:HA	1:B:297:ILE:HG12	1.97	0.46
1:A:193:VAL:CG2	1:B:141:TYR:HB2	2.45	0.46
1:B:167:ILE:HG22	1:B:171:ILE:HD12	1.96	0.46
1:A:5:ILE:HD12	1:A:36:PHE:CD1	2.50	0.46
1:B:331:THR:H	1:B:334:ILE:HD11	1.80	0.46
1:A:167:ILE:O	1:A:170:ILE:HG22	2.15	0.46
1:B:17:VAL:O	1:B:20:ALA:HB3	2.15	0.46
1:A:345:THR:HA	1:A:348:THR:HB	1.98	0.46
1:A:115:ALA:HA	1:A:125:ILE:HD11	1.96	0.46
1:A:254:MET:HE1	1:B:230:ARG:NE	2.31	0.45
1:A:298:LEU:O	1:A:301:ALA:HB3	2.15	0.45
1:B:29:ALA:HB2	1:B:36:PHE:HE2	1.79	0.45
1:A:154:GLU:O	1:A:154:GLU:HG3	2.16	0.45
1:A:334:ILE:HD13	1:A:334:ILE:H	1.81	0.45
1:A:92:LEU:O	1:A:95:ILE:HB	2.17	0.45
1:B:96:ARG:HA	1:B:101:LEU:HD12	1.98	0.45
1:A:230:ARG:NH2	1:B:254:MET:HE1	2.26	0.45
1:B:330:ARG:HG3	1:B:334:ILE:HG13	1.99	0.44
1:B:184:LYS:HB3	1:B:236:ASP:HB3	1.99	0.44
1:A:154:GLU:O	1:A:154:GLU:CG	2.66	0.44
1:A:330:ARG:HG3	1:A:334:ILE:HG12	1.99	0.44
1:A:262:LEU:HA	1:A:263:PRO:HD3	1.84	0.43
1:B:115:ALA:HB1	1:B:122:LYS:HG3	2.00	0.43
1:A:194:LEU:O	1:A:195:GLU:C	2.61	0.43
1:A:118:SER:OG	1:A:120:LEU:N	2.37	0.43
1:B:15:PRO:O	1:B:16:GLU:C	2.62	0.43
1:A:213:ASP:OD1	1:A:213:ASP:N	2.44	0.43
1:B:31:ARG:CB	1:B:359:LEU:CD2	2.97	0.43
1:A:201:ARG:O	1:A:204:ALA:HB3	2.18	0.43
1:A:243:MET:O	1:A:244:PHE:C	2.61	0.43
1:A:285:ILE:HB	1:A:290:ILE:HD11	1.99	0.43
1:B:245:GLY:O	1:B:249:SER:HB2	2.18	0.43
1:A:104:ASN:HB2	1:A:136:LEU:HD21	1.99	0.43
1:A:290:ILE:C	1:A:290:ILE:CD1	2.92	0.43
1:B:74:VAL:O	1:B:281:SER:HB2	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:265:ALA:HB2	1:B:275:TYR:CD2	2.54	0.43
1:B:42:LEU:O	1:B:59:THR:HG21	2.19	0.42
1:B:124:VAL:HG22	1:B:125:ILE:CG2	2.48	0.42
1:B:323:LYS:O	1:B:325:LEU:N	2.51	0.42
1:A:185:VAL:O	1:A:185:VAL:HG12	2.17	0.42
1:A:351:VAL:O	1:A:355:VAL:HG13	2.19	0.42
1:B:27:ALA:O	1:B:28:VAL:C	2.61	0.42
1:B:262:LEU:HA	1:B:263:PRO:HD3	1.96	0.42
1:A:106:ARG:CZ	1:A:249:SER:HB3	2.48	0.42
1:A:223:ASN:O	1:A:224:ALA:C	2.62	0.42
1:B:8:LEU:HB3	1:B:71:LEU:HD12	2.02	0.42
1:A:135:GLU:O	1:A:242:ASN:HA	2.19	0.42
1:A:361:THR:CG2	1:A:362:SER:N	2.63	0.42
1:B:120:LEU:HB2	1:B:125:ILE:HG21	2.01	0.42
1:B:171:ILE:O	1:B:175:PHE:HD2	2.03	0.42
1:A:48:ILE:HD13	1:A:87:ARG:O	2.20	0.42
1:B:70:LEU:HD23	1:B:275:TYR:HB2	2.01	0.42
1:B:243:MET:O	1:B:244:PHE:C	2.63	0.42
1:A:141:TYR:HD1	1:A:141:TYR:HA	1.80	0.42
1:A:261:MET:HE3	1:A:292:ASN:ND2	2.24	0.42
1:B:216:LEU:HD21	1:B:218:HIS:CE1	2.54	0.42
1:B:310:LEU:HB3	1:B:313:GLU:CD	2.45	0.42
1:A:186:THR:O	1:A:238:ILE:HA	2.20	0.42
1:B:194:LEU:HD23	1:B:194:LEU:HA	1.74	0.42
1:B:43:ILE:HG13	1:B:44:GLY:N	2.34	0.42
1:B:111:TYR:HB2	1:B:114:LEU:HD23	2.02	0.42
1:A:230:ARG:HE	1:B:254:MET:CE	2.33	0.41
1:A:318:GLU:O	1:A:322:GLU:HB2	2.19	0.41
1:A:325:LEU:O	1:A:330:ARG:NH1	2.44	0.41
1:B:188:VAL:HA	1:B:219:MET:O	2.19	0.41
1:A:14:GLY:N	1:A:15:PRO:HD2	2.35	0.41
1:A:261:MET:HE2	1:A:292:ASN:HD21	1.78	0.41
1:A:188:VAL:CG1	1:A:221:VAL:HA	2.51	0.41
1:A:262:LEU:O	1:A:295:ALA:HB1	2.20	0.41
1:A:208:ALA:HB2	1:A:216:LEU:HD22	2.02	0.41
1:A:112:ASP:O	1:A:113:SER:C	2.63	0.41
1:B:84:SER:C	1:B:86:LEU:H	2.28	0.41
1:B:302:MET:O	1:B:305:ARG:HB3	2.20	0.41
1:A:107:PRO:HD2	1:A:263:PRO:O	2.20	0.41
1:A:292:ASN:OD1	1:A:292:ASN:C	2.63	0.41
1:A:329:TYR:C	1:A:330:ARG:HD2	2.44	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:84:SER:C	1:B:86:LEU:N	2.77	0.41
1:B:167:ILE:O	1:B:168:GLU:C	2.64	0.41
1:B:124:VAL:HG22	1:B:125:ILE:HG23	2.02	0.40
1:B:288:LYS:O	1:B:290:ILE:HG12	2.21	0.40
1:A:254:MET:O	1:A:255:ILE:C	2.64	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	351/390 (90%)	297 (85%)	48 (14%)	6 (2%)	7	25
1	B	353/390 (90%)	300 (85%)	46 (13%)	7 (2%)	6	21
All	All	704/780 (90%)	597 (85%)	94 (13%)	13 (2%)	6	23

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	115	ALA
1	B	20	ALA
1	A	83	PRO
1	B	51	ALA
1	B	83	PRO
1	B	365	MET
1	A	49	ASP
1	A	50	GLU
1	A	87	ARG
1	A	193	VAL
1	B	87	ARG
1	B	289	GLY
1	B	364	ILE

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	256/315 (81%)	218 (85%)	38 (15%)	3	10
1	B	265/315 (84%)	219 (83%)	46 (17%)	2	7
All	All	521/630 (83%)	437 (84%)	84 (16%)	2	8

All (84) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	ASP
1	A	13	ILE
1	A	22	ILE
1	A	60	LEU
1	A	92	LEU
1	A	95	ILE
1	A	109	LYS
1	A	110	VAL
1	A	114	LEU
1	A	120	LEU
1	A	124	VAL
1	A	128	VAL
1	A	129	ASP
1	A	137	THR
1	A	141	TYR
1	A	150	GLU
1	A	159	THR
1	A	165	GLU
1	A	176	GLU
1	A	181	ARG
1	A	190	LYS
1	A	216	LEU
1	A	229	ILE
1	A	237	VAL
1	A	249	SER
1	A	256	THR
1	A	278	VAL

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Mol	Chain	Res	Type
1	A	284	ASP
1	A	296	THR
1	A	315	LYS
1	A	327	GLU
1	A	331	THR
1	A	333	ASP
1	A	334	ILE
1	A	345	THR
1	A	347	MET
1	A	352	LYS
1	A	355	VAL
1	B	3	LYS
1	B	8	LEU
1	B	11	ASP
1	B	13	ILE
1	B	34	HIS
1	B	53	THR
1	B	60	LEU
1	B	71	LEU
1	B	79	TRP
1	B	99	LEU
1	B	105	LEU
1	B	109	LYS
1	B	110	VAL
1	B	113	SER
1	B	114	LEU
1	B	120	LEU
1	B	124	VAL
1	B	128	VAL
1	B	137	THR
1	B	159	THR
1	B	160	LEU
1	B	168	GLU
1	B	181	ARG
1	B	193	VAL
1	B	229	ILE
1	B	241	GLU
1	B	249	SER
1	B	255	ILE
1	B	256	THR
1	B	258	SER
1	B	259	LEU

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Mol	Chain	Res	Type
1	B	269	THR
1	B	278	VAL
1	B	284	ASP
1	B	290	ILE
1	B	315	LYS
1	B	317	ILE
1	B	321	VAL
1	B	327	GLU
1	B	331	THR
1	B	334	ILE
1	B	343	SER
1	B	345	THR
1	B	347	MET
1	B	355	VAL
1	B	357	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	34	HIS
1	A	227	GLN
1	A	242	ASN
1	B	218	HIS
1	B	242	ASN
1	B	292	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	355/390 (91%)	0.00	5 (1%) 73 64	39, 63, 94, 105	0
1	B	357/390 (91%)	-0.01	7 (1%) 65 56	38, 64, 95, 104	0
All	All	712/780 (91%)	-0.00	12 (1%) 69 60	38, 64, 94, 105	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	364	ILE	3.6
1	A	364	ILE	3.0
1	B	365	MET	2.9
1	B	360	ALA	2.6
1	B	38	PHE	2.4
1	A	309	GLY	2.2
1	B	309	GLY	2.2
1	B	85	GLU	2.2
1	A	42	LEU	2.2
1	B	363	ALA	2.1
1	A	361	THR	2.0
1	A	81	GLN	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.