



Full wwPDB X-ray Structure Validation Report ⓘ

May 23, 2020 – 01:10 pm BST

PDB ID : 2VHY
Title : Crystal structure of apo L-alanine dehydrogenase from Mycobacterium tuberculosis
Authors : Agren, D.; Schneider, G.
Deposited on : 2007-11-26
Resolution : 2.30 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

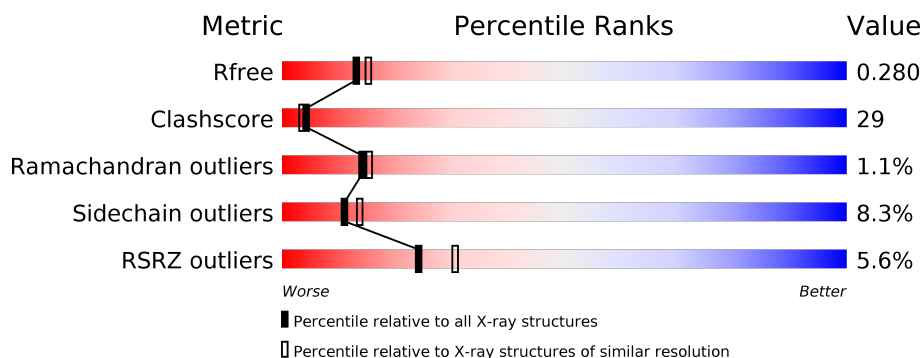
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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}	130704	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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 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	377	<div> <div>4%</div> <div> <div></div> <div>46%</div> <div>36%</div> <div>5%</div> <div>14%</div> </div> </div>
1	B	377	<div> <div>6%</div> <div> <div></div> <div>43%</div> <div>42%</div> <div>•</div> <div>12%</div> </div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4844 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 ALANINE DEHYDROGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	325	Total	C	N	O	S	0	0	1
			2376	1494	419	451	12			
1	B	331	Total	C	N	O	S	0	0	1
			2420	1523	427	458	12			

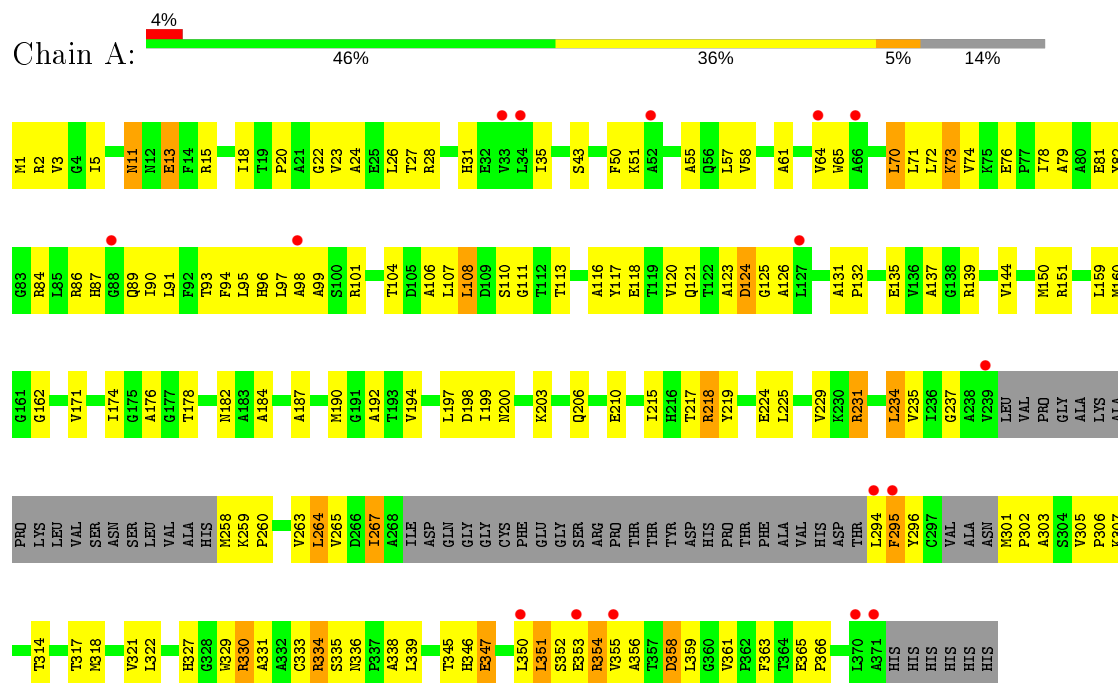
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	29	Total	O	0	0
			29	29		
2	B	19	Total	O	0	0
			19	19		

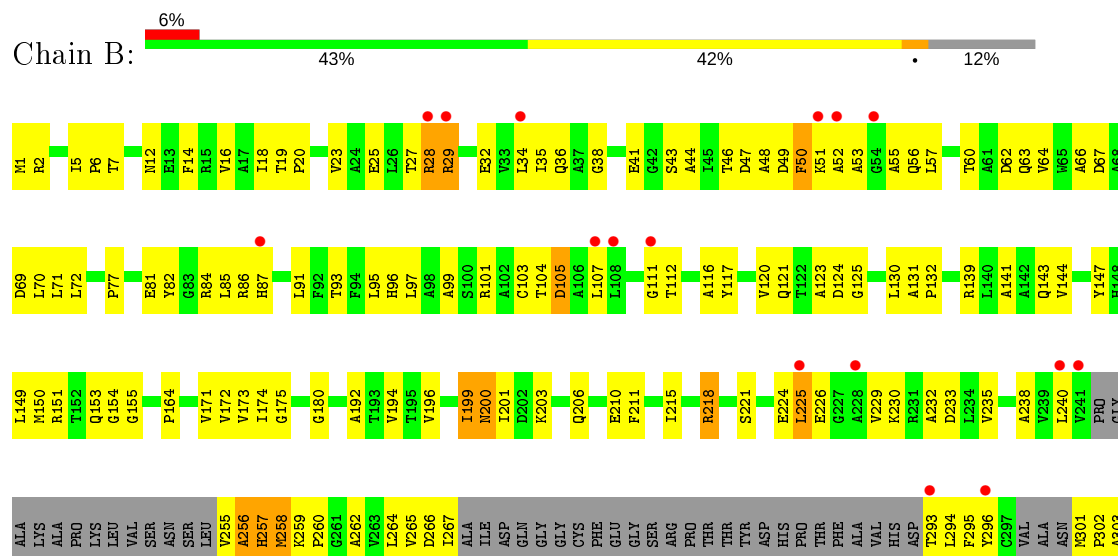
3 Residue-property plots

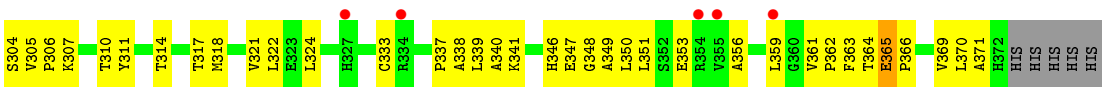
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are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



• Molecule 1: ALANINE DEHYDROGENASE





4 Data and refinement statistics

Property	Value	Source
Space group	P 3 2 1	Depositor
Cell constants a, b, c, α , β , γ	147.11Å 147.11Å 93.66Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	45.00 – 2.30 30.15 – 2.15	Depositor EDS
% Data completeness (in resolution range)	93.2 (45.00-2.30) 71.1 (30.15-2.15)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.51 (at 2.16Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.215 , 0.277 0.227 , 0.280	Depositor DCC
R_{free} test set	2173 reflections (4.80%)	wwPDB-VP
Wilson B-factor (Å ²)	35.1	Xtriage
Anisotropy	0.072	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 48.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.39$, $\langle L^2 \rangle = 0.22$	Xtriage
Estimated twinning fraction	0.419 for -h,-k,l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	4844	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.27% 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	0.43	0/2413	0.66	0/3280
1	B	0.41	0/2458	0.66	0/3343
All	All	0.42	0/4871	0.66	0/6623

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

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	2376	0	2397	147	0
1	B	2420	0	2445	144	0
2	A	29	0	0	3	0
2	B	19	0	0	4	0
All	All	4844	0	4842	282	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 29.

All (282) 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:235:VAL:HG23	1:A:264:LEU:HD23	1.28	1.12

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:98:ALA:HB2	1:A:355:VAL:HG22	1.42	1.01
1:A:235:VAL:CG2	1:A:264:LEU:HD23	1.97	0.94
1:B:199:ILE:H	1:B:199:ILE:HD13	1.32	0.94
1:A:351:LEU:HD12	1:A:366:PRO:HB3	1.50	0.94
1:A:351:LEU:HD23	1:A:351:LEU:H	1.41	0.84
1:B:141:ALA:HB2	1:B:267:ILE:HD11	1.62	0.82
1:A:51:LYS:HD3	1:A:57:LEU:HD23	1.65	0.78
1:B:25:GLU:O	1:B:29:ARG:HB2	1.82	0.77
1:A:160:MET:HB2	1:B:143:GLN:HE22	1.49	0.77
1:B:200:ASN:HB3	1:B:203:LYS:HD2	1.68	0.76
1:B:2:ARG:HG2	1:B:32:GLU:HB3	1.67	0.76
1:B:265:VAL:HG13	1:B:296:TYR:CD2	2.21	0.76
1:B:81:GLU:O	1:B:84:ARG:HB2	1.87	0.75
1:B:346:HIS:CD2	1:B:347:GLU:HG3	2.22	0.74
1:A:113:THR:HA	1:A:345:THR:O	1.86	0.74
1:A:24:ALA:O	1:A:28:ARG:HG2	1.88	0.73
1:B:44:ALA:HB3	1:B:307:LYS:HD3	1.69	0.73
1:B:318:MET:O	1:B:322:LEU:HG	1.89	0.73
1:B:264:LEU:HD23	1:B:295:PHE:HD2	1.53	0.73
1:B:265:VAL:HG13	1:B:296:TYR:HD2	1.51	0.73
1:A:20:PRO:HA	1:A:23:VAL:HG22	1.71	0.72
1:B:206:GLN:O	1:B:210:GLU:HG3	1.89	0.72
1:B:351:LEU:HD13	1:B:369:VAL:HG11	1.71	0.71
1:A:206:GLN:O	1:A:210:GLU:HG3	1.91	0.71
1:B:85:LEU:HD13	1:B:107:LEU:HD13	1.72	0.70
1:A:174:ILE:HD13	1:A:225:LEU:HD11	1.73	0.70
1:A:132:PRO:HA	1:A:135:GLU:HG3	1.72	0.70
1:A:108:LEU:HD11	1:A:350:LEU:HB2	1.73	0.70
1:B:103:CYS:O	1:B:107:LEU:HD23	1.92	0.69
1:B:36:GLN:HB2	1:B:64:VAL:HG21	1.72	0.69
1:B:144:VAL:HG11	1:B:296:TYR:CZ	2.27	0.69
1:B:225:LEU:O	1:B:229:VAL:HG12	1.93	0.69
1:B:20:PRO:HA	1:B:23:VAL:HG22	1.76	0.68
1:B:361:VAL:HG13	1:B:362:PRO:HD2	1.76	0.67
1:A:334:ARG:HG3	1:A:335:SER:N	2.09	0.66
1:A:124:ASP:OD2	1:A:126:ALA:HB2	1.96	0.65
1:A:351:LEU:HD23	1:A:351:LEU:N	2.11	0.65
1:B:337:PRO:O	1:B:341:LYS:HG3	1.97	0.65
1:B:200:ASN:OD1	1:B:203:LYS:HE3	1.97	0.64
1:B:95:LEU:HD12	1:B:116:ALA:HB2	1.77	0.64
1:B:226:GLU:HG2	1:B:230:LYS:HE3	1.78	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:7:THR:H	1:B:36:GLN:NE2	1.96	0.64
1:B:356:ALA:HB1	1:B:361:VAL:O	1.97	0.64
1:B:264:LEU:HD23	1:B:295:PHE:CD2	2.33	0.64
1:B:5:ILE:CD1	1:B:72:LEU:HB2	2.28	0.63
1:B:154:GLY:O	1:B:294:LEU:HD21	1.98	0.63
1:B:47:ASP:HA	1:B:50:PHE:HB2	1.80	0.63
1:A:73:LYS:HB2	1:A:93:THR:HB	1.81	0.63
1:A:73:LYS:HB2	1:A:93:THR:CB	2.29	0.62
1:B:307:LYS:O	1:B:311:TYR:HD2	1.82	0.62
1:A:264:LEU:HD22	1:A:265:VAL:N	2.14	0.61
1:A:264:LEU:O	1:A:264:LEU:HD13	2.00	0.61
1:B:5:ILE:HD13	1:B:72:LEU:HB2	1.83	0.61
1:B:19:THR:O	1:B:23:VAL:HG13	2.00	0.61
1:B:1:MET:HB2	1:B:69:ASP:CB	2.30	0.61
1:B:25:GLU:O	1:B:29:ARG:HD2	2.00	0.60
1:A:346:HIS:NE2	1:A:347:GLU:HG3	2.17	0.59
1:A:20:PRO:HA	1:A:23:VAL:CG2	2.31	0.59
1:A:50:PHE:O	1:A:55:ALA:HB3	2.02	0.59
1:A:197:LEU:HD23	1:A:218:ARG:HB3	1.84	0.59
1:A:79:ALA:HA	1:A:82:TYR:CE2	2.37	0.59
1:A:121:GLN:O	1:A:338:ALA:HA	2.03	0.59
1:A:90:ILE:HD13	1:A:329:TRP:HB3	1.83	0.59
1:A:27:THR:HB	1:A:28:ARG:NH2	2.18	0.58
1:B:1:MET:HB2	1:B:69:ASP:HB2	1.84	0.58
1:A:150:MET:HE2	1:B:304:SER:HB3	1.86	0.58
1:A:123:ALA:C	1:A:125:GLY:H	2.07	0.58
1:A:178:THR:O	1:A:182:ASN:ND2	2.31	0.58
1:B:259:LYS:HB2	1:B:260:PRO:HD2	1.85	0.57
1:B:351:LEU:CD1	1:B:369:VAL:HG11	2.34	0.57
1:B:72:LEU:HD11	1:B:321:VAL:HG11	1.86	0.57
1:B:38:GLY:HA2	1:B:41:GLU:HB2	1.85	0.57
1:A:317:THR:HG22	1:A:321:VAL:HG23	1.86	0.57
1:B:121:GLN:O	1:B:338:ALA:HA	2.04	0.57
1:B:147:TYR:O	1:B:150:MET:HG3	2.05	0.57
1:A:150:MET:CE	1:B:304:SER:HB3	2.33	0.57
1:A:123:ALA:O	1:A:125:GLY:N	2.38	0.57
1:A:330:ARG:O	1:A:334:ARG:HB3	2.05	0.57
1:B:199:ILE:H	1:B:199:ILE:CD1	2.11	0.57
1:B:199:ILE:HD13	1:B:199:ILE:N	2.13	0.57
1:B:333:CYS:O	1:B:340:ALA:HB2	2.05	0.57
1:A:234:LEU:HD21	1:A:265:VAL:HG23	1.87	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:264:LEU:C	1:A:264:LEU:HD13	2.24	0.57
1:A:65:TRP:CD2	1:A:84:ARG:HB3	2.40	0.57
1:B:48:ALA:O	1:B:51:LYS:HG2	2.05	0.56
1:A:217:THR:O	1:A:218:ARG:HD3	2.04	0.56
1:A:365:GLU:OE2	1:A:365:GLU:HA	2.05	0.56
1:A:71:LEU:HD23	1:A:73:LYS:HG2	1.86	0.56
1:B:72:LEU:HD11	1:B:321:VAL:CG1	2.35	0.56
1:A:159:LEU:HD21	1:B:139:ARG:HD3	1.88	0.56
1:B:317:THR:HG21	2:B:2005:HOH:O	2.07	0.55
1:B:62:ASP:CG	1:B:84:ARG:HH21	2.10	0.55
1:A:351:LEU:HD12	1:A:366:PRO:CB	2.32	0.55
1:B:101:ARG:NH1	1:B:359:LEU:HB3	2.22	0.55
1:B:55:ALA:O	1:B:57:LEU:HD22	2.06	0.55
1:B:144:VAL:HG11	1:B:296:TYR:CE2	2.41	0.55
1:B:131:ALA:N	1:B:132:PRO:HD2	2.22	0.55
1:B:240:LEU:HA	2:B:2014:HOH:O	2.06	0.55
1:A:65:TRP:CE3	1:A:84:ARG:HB3	2.42	0.54
1:A:101:ARG:HH22	1:A:359:LEU:HA	1.71	0.54
1:B:256:ALA:O	1:B:257:HIS:ND1	2.40	0.54
1:B:229:VAL:HG22	1:B:258:MET:HB3	1.90	0.54
1:B:25:GLU:HG2	1:B:29:ARG:NH1	2.22	0.54
1:A:101:ARG:NH2	1:A:359:LEU:HA	2.23	0.54
1:A:23:VAL:HG11	1:A:35:ILE:HD11	1.89	0.54
1:B:2:ARG:O	1:B:69:ASP:N	2.38	0.54
1:A:106:ALA:O	1:A:110:SER:HB2	2.08	0.54
1:A:98:ALA:CB	1:A:355:VAL:HG22	2.28	0.54
1:A:43:SER:HB3	1:A:306:PRO:HB2	1.89	0.54
1:B:20:PRO:HA	1:B:23:VAL:CG2	2.38	0.53
1:A:231:ARG:HG2	2:A:2021:HOH:O	2.09	0.53
1:A:317:THR:HG21	2:A:2003:HOH:O	2.08	0.53
1:A:174:ILE:CD1	1:A:225:LEU:HD11	2.37	0.53
1:B:305:VAL:O	1:B:305:VAL:HG23	2.09	0.53
1:A:171:VAL:HB	1:A:194:VAL:HG22	1.90	0.53
1:A:1:MET:H2	1:A:31:HIS:CE1	2.26	0.52
1:B:200:ASN:HD22	1:B:200:ASN:C	2.10	0.52
1:B:139:ARG:O	1:B:143:GLN:HG3	2.10	0.52
1:B:324:LEU:HD21	1:B:339:LEU:HD21	1.91	0.52
1:A:27:THR:HB	1:A:28:ARG:HH22	1.74	0.52
1:B:262:ALA:O	1:B:293:THR:HB	2.09	0.52
1:A:144:VAL:HG11	1:A:296:TYR:CE2	2.44	0.52
1:A:90:ILE:CD1	1:A:329:TRP:HB3	2.39	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:71:LEU:HB3	1:A:91:LEU:HD12	1.93	0.52
1:B:303:ALA:O	1:B:306:PRO:HD3	2.10	0.52
1:A:95:LEU:HB3	1:A:97:LEU:HG	1.92	0.51
1:A:98:ALA:HB2	1:A:355:VAL:CG2	2.29	0.51
1:A:363:PHE:HE2	1:A:365:GLU:CD	2.12	0.51
1:B:28:ARG:CB	1:B:29:ARG:HH21	2.24	0.51
1:B:201:ILE:HG22	1:B:201:ILE:O	2.10	0.51
1:A:70:LEU:HD22	1:A:71:LEU:N	2.26	0.51
1:A:78:ILE:HG12	1:A:81:GLU:HG3	1.91	0.51
1:B:240:LEU:HD21	1:B:267:ILE:C	2.30	0.51
1:A:329:TRP:O	1:A:333:CYS:HB2	2.09	0.51
1:A:70:LEU:HD23	1:A:90:ILE:O	2.11	0.51
1:A:307:LYS:O	1:A:307:LYS:HD3	2.11	0.51
1:B:57:LEU:HD22	1:B:57:LEU:H	1.75	0.51
1:B:123:ALA:C	1:B:125:GLY:H	2.14	0.51
1:B:66:ALA:O	1:B:67:ASP:HB3	2.11	0.51
1:B:120:VAL:HG13	1:B:339:LEU:HA	1.93	0.50
1:B:111:GLY:HA2	1:B:348:GLY:H	1.75	0.50
1:B:259:LYS:O	1:B:293:THR:HG22	2.11	0.50
1:A:3:VAL:HG13	1:A:70:LEU:HD13	1.94	0.50
1:B:258:MET:HE2	1:B:264:LEU:HB2	1.93	0.50
1:B:28:ARG:HB2	1:B:29:ARG:HH21	1.77	0.50
1:A:94:PHE:HB3	1:A:96:HIS:CE1	2.47	0.50
1:B:301:MET:HB3	1:B:302:PRO:HD3	1.94	0.49
1:A:151:ARG:HG2	1:B:14:PHE:CD2	2.48	0.49
1:A:259:LYS:HB2	1:A:260:PRO:HD2	1.94	0.49
1:A:363:PHE:HE2	1:A:365:GLU:OE1	1.96	0.49
1:B:258:MET:CE	1:B:264:LEU:HB2	2.43	0.49
1:A:90:ILE:HA	1:A:113:THR:O	2.12	0.49
1:A:235:VAL:HG22	1:A:258:MET:HE1	1.95	0.49
1:A:137:ALA:HB1	1:A:267:ILE:HD12	1.93	0.49
1:A:18:ILE:HG13	1:A:23:VAL:HG13	1.94	0.49
1:B:150:MET:HB2	1:B:153:GLN:HG2	1.95	0.49
1:A:1:MET:N	1:A:31:HIS:ND1	2.51	0.48
1:B:16:VAL:HG22	1:B:18:ILE:H	1.77	0.48
1:B:255:VAL:O	1:B:256:ALA:HB2	2.12	0.48
1:B:43:SER:HB3	1:B:306:PRO:HB2	1.94	0.48
1:A:18:ILE:HD11	1:A:22:GLY:HA3	1.96	0.48
1:B:353:GLU:HB2	1:B:363:PHE:CE1	2.49	0.48
1:A:74:VAL:O	1:A:94:PHE:HB2	2.14	0.48
1:B:172:VAL:CG2	1:B:232:ALA:HB2	2.44	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:159:LEU:HD21	1:B:139:ARG:HB2	1.96	0.48
1:A:184:ALA:HB1	1:A:215:ILE:HD13	1.95	0.48
1:A:61:ALA:O	1:A:65:TRP:HD1	1.97	0.48
1:A:3:VAL:CG1	1:A:70:LEU:HD13	2.43	0.48
1:A:225:LEU:O	1:A:229:VAL:HG23	2.14	0.48
1:A:1:MET:N	1:A:31:HIS:CE1	2.82	0.47
1:B:41:GLU:OE1	1:B:46:THR:HG22	2.15	0.47
1:A:317:THR:CG2	1:A:321:VAL:HG23	2.44	0.47
1:B:151:ARG:O	1:B:151:ARG:NH1	2.43	0.47
1:A:356:ALA:HA	1:A:361:VAL:CG2	2.44	0.47
1:B:28:ARG:HB2	1:B:29:ARG:NH2	2.29	0.47
1:B:87:HIS:CE1	1:B:347:GLU:HG2	2.50	0.47
1:B:120:VAL:HG13	1:B:338:ALA:O	2.15	0.47
1:A:160:MET:CB	1:B:143:GLN:HE22	2.24	0.46
1:A:131:ALA:N	1:A:132:PRO:HD2	2.29	0.46
1:A:314:THR:O	1:A:318:MET:HB2	2.15	0.46
1:A:231:ARG:HH11	1:A:231:ARG:HG3	1.79	0.46
1:A:234:LEU:HD23	1:A:263:VAL:O	2.15	0.46
1:B:49:ASP:HB2	2:B:2003:HOH:O	2.14	0.46
1:B:63:GLN:O	1:B:66:ALA:O	2.33	0.46
1:B:71:LEU:HD23	1:B:91:LEU:CD1	2.45	0.46
1:B:174:ILE:HD12	1:B:225:LEU:HD21	1.98	0.46
1:B:77:PRO:HA	1:B:81:GLU:OE1	2.16	0.46
1:A:363:PHE:CE2	1:A:365:GLU:CD	2.88	0.46
1:A:13:GLU:HG2	2:A:2002:HOH:O	2.14	0.46
1:A:160:MET:HB2	1:B:143:GLN:NE2	2.25	0.46
1:B:97:LEU:HD13	1:B:104:THR:HG23	1.98	0.46
1:A:264:LEU:HB3	1:A:295:PHE:CE1	2.50	0.46
1:B:180:GLY:N	1:B:238:ALA:HB1	2.31	0.46
1:B:60:THR:O	1:B:64:VAL:HG23	2.15	0.45
1:B:310:THR:O	1:B:314:THR:HG23	2.15	0.45
1:A:97:LEU:C	1:A:99:ALA:H	2.19	0.45
1:A:123:ALA:C	1:A:125:GLY:N	2.68	0.45
1:B:86:ARG:O	1:B:112:THR:HB	2.16	0.45
1:B:200:ASN:OD1	1:B:203:LYS:CE	2.64	0.45
1:B:51:LYS:C	1:B:53:ALA:H	2.19	0.45
1:B:35:ILE:O	1:B:57:LEU:HA	2.17	0.45
1:B:172:VAL:HB	1:B:235:VAL:HG13	1.97	0.45
1:A:354:ARG:O	1:A:358:ASP:OD1	2.35	0.45
1:B:101:ARG:HH12	1:B:359:LEU:HB3	1.82	0.45
1:A:93:THR:HA	1:A:117:TYR:CE1	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:23:VAL:CG1	1:A:35:ILE:HD11	2.47	0.44
1:B:346:HIS:O	1:B:349:ALA:HB3	2.17	0.44
1:A:231:ARG:NH1	1:A:231:ARG:HG3	2.33	0.44
1:B:151:ARG:HA	1:B:155:GLY:O	2.18	0.44
1:A:104:THR:O	1:A:108:LEU:HB2	2.16	0.44
1:A:356:ALA:C	1:A:358:ASP:H	2.21	0.44
1:B:229:VAL:CG2	1:B:258:MET:HB3	2.47	0.44
1:A:162:GLY:O	1:B:139:ARG:NE	2.50	0.44
1:A:351:LEU:CD2	1:A:351:LEU:N	2.81	0.44
1:B:82:TYR:C	1:B:84:ARG:H	2.21	0.44
1:A:187:ALA:O	1:A:192:ALA:HB3	2.17	0.44
1:B:171:VAL:HB	1:B:194:VAL:HG22	1.99	0.43
1:A:20:PRO:CA	1:A:23:VAL:HG22	2.44	0.43
1:B:16:VAL:O	1:B:310:THR:HG23	2.18	0.43
1:A:74:VAL:O	1:A:94:PHE:HD1	2.01	0.43
1:A:327:HIS:O	1:A:331:ALA:HB3	2.19	0.43
1:A:352:SER:O	1:A:356:ALA:HB3	2.19	0.43
1:A:301:MET:HB3	1:A:302:PRO:HD3	1.99	0.43
1:A:5:ILE:HG13	1:A:72:LEU:HB2	1.99	0.43
1:A:135:GLU:O	1:A:139:ARG:HG3	2.19	0.43
1:A:237:GLY:O	1:A:267:ILE:HG23	2.18	0.43
1:A:327:HIS:HB3	1:A:331:ALA:CB	2.49	0.43
1:A:23:VAL:HG11	1:A:35:ILE:CD1	2.49	0.43
1:B:173:VAL:HB	1:B:196:VAL:HG22	2.00	0.43
1:A:197:LEU:CD2	1:A:218:ARG:HB3	2.49	0.43
1:A:305:VAL:O	1:A:305:VAL:HG23	2.19	0.43
1:A:71:LEU:CD2	1:A:73:LYS:HG2	2.49	0.43
1:A:116:ALA:HB1	1:A:118:GLU:OE1	2.18	0.42
1:B:149:LEU:HD21	1:B:233:ASP:O	2.19	0.42
1:B:365:GLU:HA	1:B:366:PRO:HD3	1.93	0.42
1:B:27:THR:C	1:B:29:ARG:H	2.23	0.42
1:A:264:LEU:C	1:A:264:LEU:HD22	2.39	0.42
1:A:317:THR:HG22	1:A:321:VAL:CG2	2.48	0.42
1:A:26:LEU:HD23	1:A:322:LEU:HD23	2.00	0.42
1:A:51:LYS:HA	1:A:57:LEU:HD21	2.00	0.42
1:B:107:LEU:HD12	1:B:112:THR:CG2	2.50	0.42
1:B:211:PHE:CE1	1:B:215:ILE:HD12	2.55	0.42
1:B:333:CYS:HB2	1:B:370:LEU:HD21	2.00	0.42
1:B:364:THR:HG22	1:B:364:THR:O	2.20	0.42
1:A:22:GLY:C	1:A:24:ALA:H	2.22	0.42
1:A:15:ARG:NH1	1:A:303:ALA:HB2	2.33	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:150:MET:O	1:B:153:GLN:N	2.48	0.42
1:B:93:THR:O	1:B:117:TYR:N	2.44	0.42
1:B:32:GLU:OE2	1:B:34:LEU:HD21	2.20	0.42
1:A:58:VAL:HG21	1:A:64:VAL:CG2	2.49	0.41
1:A:346:HIS:CD2	1:A:347:GLU:N	2.88	0.41
1:A:65:TRP:CG	1:A:84:ARG:HD2	2.54	0.41
1:B:6:PRO:HD3	1:B:71:LEU:HD11	2.02	0.41
1:A:73:LYS:HB3	1:A:73:LYS:NZ	2.35	0.41
1:B:28:ARG:C	1:B:29:ARG:HE	2.23	0.41
1:A:176:ALA:HB1	1:A:203:LYS:HB3	2.02	0.41
1:A:294:LEU:HD23	1:A:295:PHE:N	2.36	0.41
1:A:352:SER:O	1:A:356:ALA:CB	2.68	0.41
1:B:337:PRO:HG3	2:B:2019:HOH:O	2.19	0.41
1:A:160:MET:HG3	1:A:190:MET:HE1	2.03	0.41
1:A:87:HIS:HD2	1:A:111:GLY:O	2.03	0.41
1:B:175:GLY:O	1:B:180:GLY:HA3	2.21	0.41
1:B:171:VAL:HG23	1:B:192:ALA:HB1	2.02	0.41
1:B:218:ARG:HG2	1:B:224:GLU:CD	2.40	0.41
1:B:221:SER:O	1:B:225:LEU:HB2	2.21	0.41
1:A:295:PHE:O	1:A:296:TYR:HB2	2.21	0.41
1:B:101:ARG:HG2	1:B:105:ASP:OD1	2.21	0.41
1:B:49:ASP:HA	1:B:52:ALA:HB3	2.02	0.41
1:A:106:ALA:O	1:A:110:SER:CB	2.69	0.41
1:A:108:LEU:HD12	1:A:108:LEU:HA	1.85	0.41
1:A:264:LEU:HD12	1:A:295:PHE:CZ	2.56	0.41
1:B:259:LYS:HG3	1:B:262:ALA:HB2	2.02	0.41
1:A:86:ARG:HB2	1:A:89:GLN:HB2	2.04	0.40
1:A:11:ASN:HD22	1:A:11:ASN:HA	1.66	0.40
1:A:97:LEU:HD21	1:A:107:LEU:HD12	2.03	0.40
1:B:20:PRO:O	1:B:23:VAL:HG22	2.21	0.40
1:A:199:ILE:HA	1:A:219:TYR:CE2	2.56	0.40
1:A:57:LEU:HD22	1:A:57:LEU:N	2.36	0.40
1:B:180:GLY:CA	1:B:238:ALA:CB	2.99	0.40
1:B:96:HIS:O	1:B:99:ALA:HB3	2.22	0.40
1:A:116:ALA:O	1:A:120:VAL:HG23	2.22	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	317/377 (84%)	282 (89%)	33 (10%)	2 (1%)	25	31
1	B	323/377 (86%)	295 (91%)	23 (7%)	5 (2%)	10	10
All	All	640/754 (85%)	577 (90%)	56 (9%)	7 (1%)	14	15

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	371	ALA
1	A	124	ASP
1	B	256	ALA
1	B	28	ARG
1	B	124	ASP
1	B	365	GLU
1	A	347	GLU

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	239/282 (85%)	215 (90%)	24 (10%)	7	9
1	B	244/282 (86%)	228 (93%)	16 (7%)	16	22
All	All	483/564 (86%)	443 (92%)	40 (8%)	11	14

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	ARG
1	A	11	ASN
1	A	13	GLU
1	A	70	LEU
1	A	73	LYS
1	A	76	GLU
1	A	108	LEU
1	A	198	ASP
1	A	200	ASN
1	A	218	ARG
1	A	224	GLU
1	A	231	ARG
1	A	234	LEU
1	A	264	LEU
1	A	267	ILE
1	A	295	PHE
1	A	330	ARG
1	A	334	ARG
1	A	336	ASN
1	A	339	LEU
1	A	351	LEU
1	A	353	GLU
1	A	354	ARG
1	A	358	ASP
1	B	12	ASN
1	B	29	ARG
1	B	50	PHE
1	B	56	GLN
1	B	70	LEU
1	B	105	ASP
1	B	130	LEU
1	B	164	PRO
1	B	199	ILE
1	B	200	ASN
1	B	218	ARG
1	B	225	LEU
1	B	257	HIS
1	B	258	MET
1	B	266	ASP
1	B	350	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	11	ASN
1	A	12	ASN
1	A	96	HIS
1	A	200	ASN
1	A	336	ASN
1	A	346	HIS
1	B	36	GLN
1	B	87	HIS
1	B	89	GLN
1	B	121	GLN
1	B	143	GLN
1	B	148	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	A	325/377 (86%)	0.55	16 (4%) 29 36	21, 47, 70, 85	0
1	B	331/377 (87%)	0.59	21 (6%) 20 25	19, 46, 71, 82	0
All	All	656/754 (87%)	0.57	37 (5%) 24 30	19, 46, 71, 85	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	107	LEU	4.4
1	A	370	LEU	4.3
1	A	350	LEU	4.3
1	A	355	VAL	4.0
1	A	64	VAL	4.0
1	B	296	TYR	3.9
1	B	34	LEU	3.8
1	A	294	LEU	3.4
1	B	334	ARG	3.4
1	B	293	THR	3.2
1	A	33	VAL	2.9
1	B	111	GLY	2.9
1	A	295	PHE	2.9
1	B	51	LYS	2.8
1	A	127	LEU	2.8
1	B	241	VAL	2.8
1	A	34	LEU	2.8
1	A	239	VAL	2.7
1	B	87	HIS	2.7
1	B	54	GLY	2.6
1	A	66	ALA	2.6
1	B	240	LEU	2.5
1	A	353	GLU	2.5
1	B	108	LEU	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	359	LEU	2.4
1	A	371	ALA	2.3
1	B	354	ARG	2.3
1	B	327	HIS	2.3
1	A	52	ALA	2.3
1	B	225	LEU	2.3
1	A	88	GLY	2.2
1	B	355	VAL	2.2
1	A	98	ALA	2.1
1	B	52	ALA	2.1
1	B	28	ARG	2.1
1	B	228	ALA	2.0
1	B	29	ARG	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.