



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 12:03 PM UTC

PDB ID : 1P74 / pdb_00001p74
Title : CRYSTAL STRUCTURE OF SHIKIMATE DEHYDROGENASE (AROE)
FROM HAEMOPHILUS INFLUENZAE
Authors : Ye, S.; von Delft, F.; Brooun, A.; Knuth, M.W.; Swanson, R.V.; McRee, D.E.
Deposited on : 2003-04-30
Resolution : 2.40 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) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

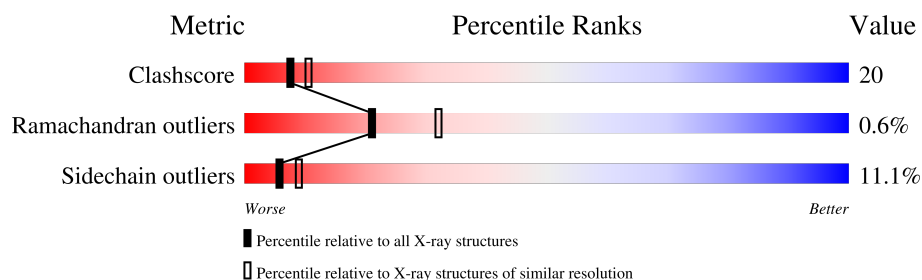
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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))
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	272	 65% 25% 7% ..
1	B	272	 62% 28% 7% ..

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4257 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 Shikimate 5-dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	267	Total	C	N	O	S	0	0	0
			2064	1317	347	390	10			
1	B	266	Total	C	N	O	S	0	0	0
			2060	1315	346	389	10			

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	67	Total	O	0	0
			67	67		
2	B	66	Total	O	0	0
			66	66		

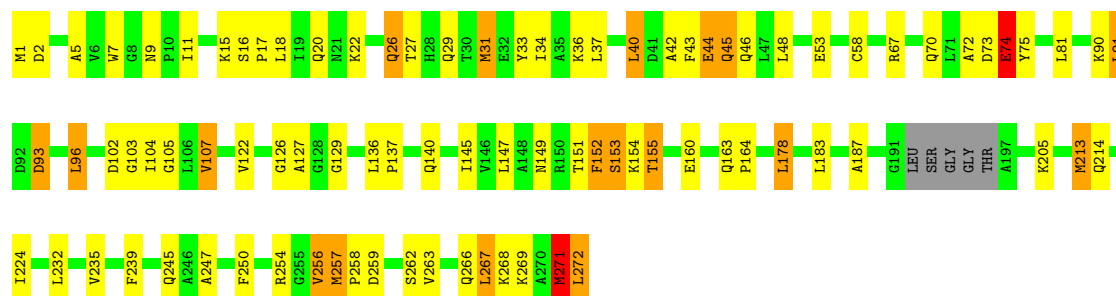
3 Residue-property plots

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. 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. 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.

Note EDS was not executed.

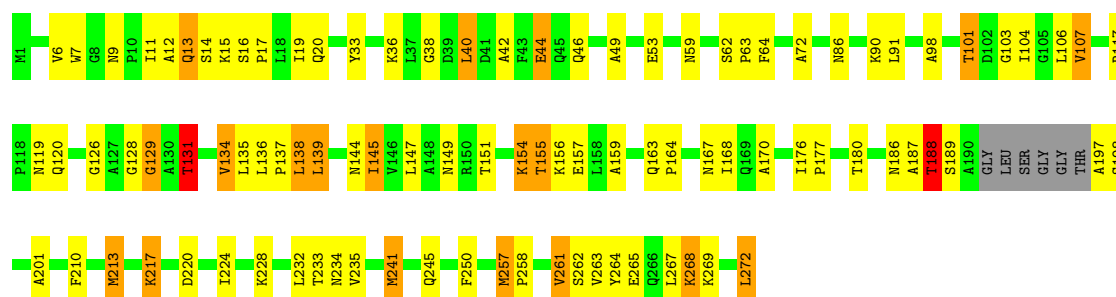
• Molecule 1: Shikimate 5-dehydrogenase

Chain A: 



• Molecule 1: Shikimate 5-dehydrogenase

Chain B: 



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	78.22Å 86.58Å 92.92Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.40	Depositor
% Data completeness (in resolution range)	99.7 (20.00-2.40)	Depositor
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
Refinement program	REFMAC 5.0	Depositor
R, R_{free}	0.228 , 0.276	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4257	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

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	0.71	3/2102 (0.1%)	1.22	20/2845 (0.7%)
1	B	0.67	3/2098 (0.1%)	1.03	7/2840 (0.2%)
All	All	0.69	6/4200 (0.1%)	1.13	27/5685 (0.5%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	241	MET	SD-CE	-12.13	1.49	1.79
1	A	31	MET	SD-CE	-11.12	1.51	1.79
1	B	213	MET	SD-CE	-10.05	1.54	1.79
1	A	257	MET	SD-CE	-8.47	1.58	1.79
1	B	257	MET	SD-CE	-6.93	1.62	1.79
1	A	213	MET	SD-CE	-6.86	1.62	1.79

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	74	GLU	O-C-N	-19.53	99.45	123.27
1	A	74	GLU	CA-C-O	16.68	139.84	120.66
1	A	43	PHE	CA-C-N	11.78	135.75	120.44
1	A	43	PHE	C-N-CA	11.78	135.75	120.44
1	A	45	GLN	O-C-N	-10.78	109.02	122.27
1	A	45	GLN	CA-C-O	9.11	130.45	119.97
1	B	261	VAL	N-CA-C	7.88	118.43	110.23
1	A	145	ILE	N-CA-C	7.25	118.26	108.11
1	A	45	GLN	CB-CG-CD	7.23	124.89	112.60
1	A	43	PHE	CA-CB-CG	7.05	120.85	113.80
1	B	201	ALA	N-CA-C	6.72	118.61	111.28
1	A	73	ASP	N-CA-C	-6.67	104.78	113.12
1	A	45	GLN	CA-CB-CG	6.47	127.04	114.10
1	A	74	GLU	CB-CG-CD	6.31	123.32	112.60
1	B	145	ILE	N-CA-C	6.22	116.87	108.17
1	A	45	GLN	OE1-CD-NE2	-6.19	116.41	122.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	74	GLU	CA-CB-CG	6.15	126.40	114.10
1	B	189	SER	N-CA-C	6.08	119.47	110.46
1	A	44	GLU	CB-CG-CD	6.05	122.89	112.60
1	B	14	SER	N-CA-C	5.62	118.60	110.28
1	A	105	GLY	N-CA-C	-5.58	105.62	112.77
1	A	72	ALA	CA-C-O	-5.49	115.83	121.87
1	B	188	THR	N-CA-C	5.41	117.28	110.24
1	B	131	THR	N-CA-C	5.36	118.03	111.82
1	A	45	GLN	CG-CD-NE2	5.20	124.19	116.40
1	A	72	ALA	CA-C-N	5.12	132.36	122.53
1	A	72	ALA	C-N-CA	5.12	132.36	122.53

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	2064	0	2069	94	0
1	B	2060	0	2066	75	0
2	A	67	0	0	5	0
2	B	66	0	0	5	0
All	All	4257	0	4135	167	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 20.

All (167) 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:44:GLU:OE2	1:A:67:ARG:NH2	1.77	1.18
1:A:178:LEU:HD13	1:B:167:ASN:HB3	1.38	1.06
1:B:147:LEU:HG	1:B:155:THR:HG22	1.33	1.05
1:B:220:ASP:OD2	1:B:228:LYS:NZ	1.89	1.05
1:A:31:MET:HE1	1:A:254:ARG:CD	1.88	1.03

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:257:MET:HE3	1:A:258:PRO:HD2	1.38	1.02
1:B:106:LEU:HD13	1:B:241:MET:HE2	1.40	1.02
1:A:187:ALA:HA	1:A:213:MET:HE3	1.43	1.00
1:A:151:THR:HG22	1:A:153:SER:H	1.23	0.99
1:B:106:LEU:HD13	1:B:241:MET:CE	1.97	0.94
1:A:151:THR:CG2	1:A:154:LYS:H	1.85	0.90
1:A:31:MET:HE1	1:A:254:ARG:HD2	1.53	0.87
1:A:151:THR:HG21	1:A:154:LYS:H	1.39	0.87
1:A:152:PHE:O	1:A:152:PHE:CD2	2.30	0.85
1:A:151:THR:HG22	1:A:153:SER:N	1.92	0.83
1:A:31:MET:CE	1:A:254:ARG:HD2	2.10	0.81
1:B:147:LEU:HG	1:B:155:THR:CG2	2.11	0.79
1:A:271:MET:O	1:A:272:LEU:HD13	1.82	0.79
1:A:29:GLN:HB3	1:A:31:MET:HE3	1.64	0.78
1:B:213:MET:HG2	1:B:241:MET:HE1	1.66	0.77
1:A:126:GLY:HA2	1:A:149:ASN:HD22	1.51	0.74
1:A:81:LEU:HD12	1:A:136:LEU:HD22	1.68	0.74
1:B:106:LEU:CD1	1:B:241:MET:HE2	2.18	0.73
1:B:163:GLN:O	2:B:330:HOH:O	2.05	0.73
1:A:31:MET:CE	1:A:254:ARG:CD	2.63	0.73
1:B:126:GLY:O	1:B:131:THR:HG21	1.87	0.73
1:A:31:MET:CE	1:A:254:ARG:NE	2.52	0.72
1:A:31:MET:HE1	1:A:254:ARG:NE	2.05	0.72
1:A:239:PHE:HZ	1:A:267:LEU:HD22	1.56	0.70
1:A:271:MET:HE2	1:A:272:LEU:CD2	2.23	0.69
1:A:149:ASN:HB3	1:A:155:THR:HG21	1.75	0.69
1:B:163:GLN:NE2	1:B:168:ILE:O	2.26	0.67
1:A:178:LEU:HD13	1:B:167:ASN:CB	2.21	0.67
1:A:152:PHE:O	1:A:152:PHE:CG	2.49	0.66
1:B:147:LEU:CG	1:B:155:THR:HG22	2.21	0.65
1:B:188:THR:HG1	1:B:197:ALA:N	1.94	0.64
1:A:44:GLU:CD	1:A:67:ARG:HH21	1.95	0.64
1:A:29:GLN:CB	1:A:31:MET:HE3	2.28	0.64
1:B:117:ARG:O	1:B:120:GLN:HG3	1.97	0.64
1:A:90:LYS:HE3	2:A:318:HOH:O	1.97	0.63
1:B:104:ILE:HA	1:B:107:VAL:HG13	1.80	0.63
1:B:263:VAL:HG12	1:B:267:LEU:HD12	1.80	0.63
1:A:187:ALA:HA	1:A:213:MET:CE	2.26	0.61
1:A:81:LEU:CD1	1:A:136:LEU:HD22	2.30	0.61
1:A:271:MET:HE2	1:A:272:LEU:HD23	1.84	0.60
1:A:136:LEU:HB3	1:A:137:PRO:HD3	1.84	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:205:LYS:HD2	1:A:232:LEU:HD21	1.83	0.60
1:B:128:GLY:O	1:B:129:GLY:C	2.45	0.59
1:B:139:LEU:HD13	1:B:145:ILE:HD12	1.82	0.59
1:B:136:LEU:HB3	1:B:137:PRO:HD3	1.85	0.59
1:A:16:SER:HB2	1:A:17:PRO:HD3	1.85	0.58
1:A:31:MET:HE1	1:A:254:ARG:CG	2.32	0.58
1:A:151:THR:HG21	1:A:154:LYS:HB2	1.84	0.58
1:A:187:ALA:CA	1:A:213:MET:HE3	2.27	0.57
1:A:151:THR:HG21	1:A:154:LYS:N	2.16	0.56
1:A:250:PHE:CD1	1:A:258:PRO:HG3	2.40	0.56
1:A:151:THR:CG2	1:A:153:SER:HB3	2.36	0.56
1:A:151:THR:HG23	1:A:153:SER:HB3	1.87	0.56
1:B:131:THR:HG22	1:B:187:ALA:CB	2.36	0.56
1:A:70:GLN:NE2	2:A:315:HOH:O	2.39	0.55
1:B:6:VAL:HG13	1:B:11:ILE:HD11	1.88	0.55
1:A:93:ASP:OD2	1:A:93:ASP:N	2.31	0.55
1:B:90:LYS:NZ	2:B:315:HOH:O	2.39	0.54
1:A:239:PHE:CZ	1:A:267:LEU:HD22	2.39	0.54
1:B:159:ALA:O	1:B:163:GLN:HG2	2.08	0.54
1:B:163:GLN:N	1:B:164:PRO:CD	2.70	0.54
1:A:271:MET:C	1:A:272:LEU:HD22	2.34	0.53
1:A:224:ILE:HD12	1:A:235:VAL:HB	1.90	0.53
1:A:5:ALA:O	1:A:58:CYS:HB2	2.08	0.53
1:B:103:GLY:O	1:B:107:VAL:HG12	2.08	0.53
1:A:263:VAL:O	1:A:267:LEU:HB2	2.09	0.52
1:A:151:THR:CG2	1:A:154:LYS:N	2.65	0.52
1:A:259:ASP:OD2	1:A:262:SER:CB	2.57	0.52
1:A:187:ALA:CB	1:A:213:MET:CE	2.88	0.52
1:B:154:LYS:O	1:B:157:GLU:HG2	2.10	0.51
1:B:250:PHE:CD1	1:B:258:PRO:HG3	2.46	0.51
1:A:122:VAL:HG22	1:A:183:LEU:HB3	1.92	0.51
1:B:217:LYS:O	1:B:217:LYS:HG3	2.09	0.51
1:A:257:MET:HE3	1:A:258:PRO:CD	2.26	0.51
1:A:31:MET:HE2	1:A:254:ARG:NE	2.25	0.50
1:A:205:LYS:CD	1:A:232:LEU:HD21	2.41	0.50
1:A:259:ASP:OD2	1:A:262:SER:HB3	2.10	0.50
1:A:11:ILE:HG22	1:A:37:LEU:HD12	1.93	0.50
1:A:102:ASP:HA	1:A:245:GLN:HE21	1.76	0.50
1:A:187:ALA:HB2	1:A:213:MET:CE	2.42	0.50
1:A:152:PHE:CD2	1:A:152:PHE:C	2.90	0.50
1:B:269:LYS:O	1:B:272:LEU:HD23	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:31:MET:HE2	1:A:254:ARG:CZ	2.42	0.49
1:B:261:VAL:HG23	2:B:282:HOH:O	2.11	0.49
1:B:15:LYS:O	1:B:19:ILE:HG12	2.12	0.49
1:B:154:LYS:HA	1:B:157:GLU:OE1	2.12	0.49
1:A:247:ALA:HB1	1:A:257:MET:HE1	1.95	0.49
1:B:20:GLN:HG3	1:B:33:TYR:CE1	2.48	0.49
1:B:261:VAL:O	1:B:264:TYR:HB3	2.13	0.49
1:A:22:LYS:NZ	1:A:266:GLN:HE22	2.10	0.48
1:A:187:ALA:CB	1:A:213:MET:HE1	2.43	0.48
1:A:163:GLN:N	1:A:164:PRO:CD	2.77	0.48
1:B:233:THR:O	1:B:235:VAL:HG22	2.13	0.48
1:A:7:TRP:CZ2	1:A:36:LYS:HD3	2.49	0.48
1:A:259:ASP:OD2	1:A:262:SER:N	2.40	0.48
1:A:26:GLN:HG3	1:A:27:THR:HG23	1.96	0.48
1:A:126:GLY:HA2	1:A:149:ASN:ND2	2.26	0.48
1:B:261:VAL:HG23	1:B:262:SER:N	2.28	0.48
1:B:49:ALA:O	1:B:53:GLU:HG3	2.14	0.47
1:A:90:LYS:NZ	2:A:313:HOH:O	2.47	0.47
1:B:6:VAL:HB	1:B:33:TYR:HH	1.79	0.47
1:B:261:VAL:CG2	2:B:282:HOH:O	2.62	0.47
1:B:13:GLN:H	1:B:13:GLN:HG2	1.47	0.47
1:A:267:LEU:HD23	1:A:271:MET:HG3	1.96	0.47
1:A:103:GLY:HA3	1:A:137:PRO:HG2	1.96	0.47
1:A:136:LEU:O	1:A:140:GLN:HG3	2.14	0.47
1:B:7:TRP:CZ2	1:B:36:LYS:HD3	2.49	0.47
1:B:101:THR:HG21	1:B:245:GLN:O	2.14	0.47
1:B:269:LYS:HA	1:B:272:LEU:CD2	2.45	0.47
1:A:7:TRP:CH2	1:A:36:LYS:HD3	2.49	0.47
1:B:42:ALA:O	1:B:46:GLN:HG3	2.14	0.47
1:B:131:THR:HG22	1:B:187:ALA:HB2	1.96	0.47
1:A:91:LEU:HB3	1:A:93:ASP:OD2	2.15	0.47
1:B:131:THR:CG2	1:B:187:ALA:CB	2.93	0.47
1:A:44:GLU:OE1	1:A:67:ARG:NE	2.27	0.46
1:B:155:THR:HB	1:B:170:ALA:HB1	1.98	0.46
1:A:149:ASN:CB	1:A:155:THR:HG21	2.45	0.46
1:A:187:ALA:HB2	1:A:213:MET:HE1	1.97	0.46
1:B:11:ILE:O	1:B:12:ALA:C	2.59	0.46
1:B:72:ALA:HB2	1:B:98:ALA:HB2	1.99	0.45
1:A:104:ILE:HA	1:A:107:VAL:HG13	1.98	0.45
1:B:62:SER:HA	1:B:63:PRO:HA	1.67	0.45
1:B:131:THR:O	1:B:135:LEU:HG	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:149:ASN:HB2	1:B:155:THR:HG23	1.98	0.45
1:A:102:ASP:HA	1:A:245:GLN:NE2	2.31	0.45
1:B:40:LEU:HA	1:B:40:LEU:HD12	1.72	0.45
1:A:96:LEU:HD12	1:A:96:LEU:HA	1.80	0.45
1:A:44:GLU:O	1:A:48:LEU:HD12	2.16	0.45
1:A:127:ALA:HB1	2:A:289:HOH:O	2.17	0.45
1:A:247:ALA:HB1	1:A:257:MET:CE	2.47	0.44
1:A:20:GLN:HG3	1:A:33:TYR:CE1	2.53	0.44
1:A:151:THR:HB	1:A:155:THR:HG23	1.99	0.44
1:B:186:ASN:ND2	1:B:188:THR:OG1	2.50	0.44
1:B:63:PRO:HD2	1:B:64:PHE:CE2	2.53	0.44
1:B:176:ILE:HA	1:B:177:PRO:HD3	1.89	0.44
1:A:74:GLU:O	1:A:75:TYR:HB3	2.18	0.44
1:A:187:ALA:CA	1:A:213:MET:CE	2.92	0.43
1:B:144:ASN:OD1	1:B:167:ASN:ND2	2.51	0.43
1:B:44:GLU:H	1:B:44:GLU:HG2	1.31	0.43
1:B:59:ASN:HA	1:B:86:ASN:O	2.18	0.43
1:B:257:MET:HA	1:B:258:PRO:HD3	1.87	0.43
1:B:33:TYR:OH	1:B:59:ASN:HB2	2.19	0.43
1:B:210:PHE:CZ	1:B:232:LEU:HD13	2.55	0.42
1:B:103:GLY:HA2	1:B:134:VAL:HG22	2.01	0.42
1:B:138:LEU:HB3	1:B:145:ILE:HD11	2.01	0.42
1:B:9:ASN:ND2	1:B:38:GLY:O	2.38	0.42
1:B:268:LYS:O	1:B:272:LEU:HD22	2.20	0.42
1:A:2:ASP:OD1	1:A:254:ARG:NH1	2.47	0.42
1:A:9:ASN:ND2	1:A:40:LEU:HD13	2.34	0.42
1:A:256:VAL:O	1:A:258:PRO:HD3	2.19	0.42
1:A:129:GLY:N	2:A:307:HOH:O	2.53	0.42
1:B:187:ALA:HA	1:B:213:MET:HE2	2.02	0.41
1:B:134:VAL:HG13	1:B:138:LEU:HD22	2.01	0.41
1:B:134:VAL:CG1	1:B:138:LEU:HD22	2.49	0.41
1:B:198:SER:O	1:B:198:SER:OG	2.27	0.41
1:B:117:ARG:HG3	1:B:120:GLN:HG3	2.02	0.41
1:B:86:ASN:HD21	1:B:101:THR:HG22	1.85	0.41
1:B:16:SER:HB2	1:B:17:PRO:HD3	2.01	0.40
1:A:74:GLU:O	1:A:75:TYR:CB	2.62	0.40
1:A:102:ASP:CB	1:A:245:GLN:NE2	2.84	0.40
1:A:42:ALA:O	1:A:46:GLN:HG3	2.21	0.40
1:B:262:SER:HB2	2:B:282:HOH:O	2.20	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	263/272 (97%)	253 (96%)	9 (3%)	1 (0%)	30	43
1	B	262/272 (96%)	250 (95%)	10 (4%)	2 (1%)	16	25
All	All	525/544 (96%)	503 (96%)	19 (4%)	3 (1%)	21	32

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	271	MET
1	B	129	GLY
1	B	234	ASN

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	217/220 (99%)	191 (88%)	26 (12%)	5	7
1	B	217/220 (99%)	195 (90%)	22 (10%)	7	11
All	All	434/440 (99%)	386 (89%)	48 (11%)	6	9

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	15	LYS
1	A	18	LEU

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Mol	Chain	Res	Type
1	A	26	GLN
1	A	34	ILE
1	A	40	LEU
1	A	45	GLN
1	A	53	GLU
1	A	74	GLU
1	A	91	LEU
1	A	93	ASP
1	A	96	LEU
1	A	107	VAL
1	A	147	LEU
1	A	152	PHE
1	A	153	SER
1	A	155	THR
1	A	160	GLU
1	A	178	LEU
1	A	214	GLN
1	A	256	VAL
1	A	267	LEU
1	A	268	LYS
1	A	269	LYS
1	A	271	MET
1	A	272	LEU
1	B	13	GLN
1	B	40	LEU
1	B	44	GLU
1	B	91	LEU
1	B	101	THR
1	B	107	VAL
1	B	119	ASN
1	B	131	THR
1	B	134	VAL
1	B	138	LEU
1	B	139	LEU
1	B	151	THR
1	B	154	LYS
1	B	155	THR
1	B	156	LYS
1	B	180	THR
1	B	188	THR
1	B	217	LYS
1	B	224	ILE

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Mol	Chain	Res	Type
1	B	265	GLU
1	B	268	LYS
1	B	272	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	77	GLN
1	A	120	GLN
1	A	149	ASN
1	A	214	GLN
1	A	245	GLN
1	A	266	GLN
1	B	20	GLN
1	B	59	ASN
1	B	119	ASN
1	B	120	GLN
1	B	140	GLN
1	B	149	ASN
1	B	245	GLN
1	B	266	GLN

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates [i](#)

EDS was not executed - this section is therefore empty.

6.4 Ligands [i](#)

EDS was not executed - this section is therefore empty.

6.5 Other polymers [i](#)

EDS was not executed - this section is therefore empty.