



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 13, 2024 – 11:45 AM EDT

PDB ID : 1PED
Title : BACTERIAL SECONDARY ALCOHOL DEHYDROGENASE (APO-FORM)
Authors : Korkhin, Y.; Frolow, F.
Deposited on : 1995-12-28
Resolution : 2.15 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 : 4.02b-467
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36.2

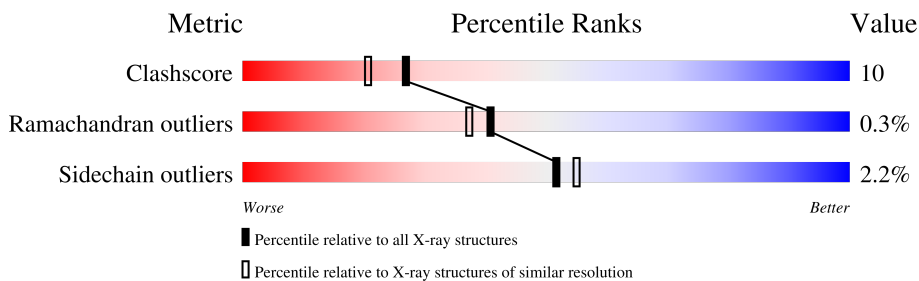
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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	141614	1585 (2.16-2.16)
Ramachandran outliers	138981	1560 (2.16-2.16)
Sidechain outliers	138945	1559 (2.16-2.16)

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	351	
1	B	351	
1	C	351	
1	D	351	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 11122 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 NADP-DEPENDENT ALCOHOL DEHYDROGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	351	Total	C	N	O	S	0	0	0
			2640	1675	460	482	23			
1	B	351	Total	C	N	O	S	0	0	0
			2640	1675	460	482	23			
1	C	351	Total	C	N	O	S	0	0	0
			2640	1675	460	482	23			
1	D	351	Total	C	N	O	S	0	0	0
			2640	1675	460	482	23			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	154	THR	SER	CONFLICT	UNP P25984
A	234	LYS	GLU	CONFLICT	UNP P25984
B	154	THR	SER	CONFLICT	UNP P25984
B	234	LYS	GLU	CONFLICT	UNP P25984
C	154	THR	SER	CONFLICT	UNP P25984
C	234	LYS	GLU	CONFLICT	UNP P25984
D	154	THR	SER	CONFLICT	UNP P25984
D	234	LYS	GLU	CONFLICT	UNP P25984

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		
2	B	1	Total	Zn	0	0
			1	1		
2	C	1	Total	Zn	0	0
			1	1		
2	D	1	Total	Zn	0	0
			1	1		

- Molecule 3 is water.

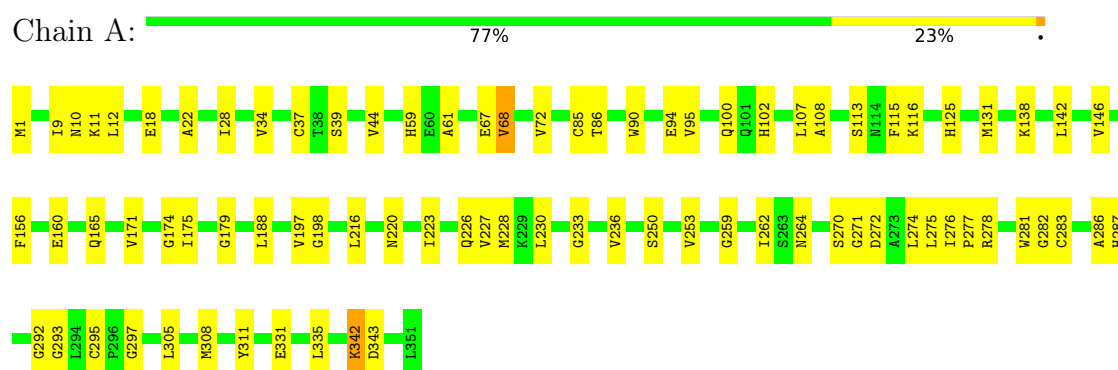
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	145	Total 145	O 145	0	0
3	B	141	Total 141	O 141	0	0
3	C	145	Total 145	O 145	0	0
3	D	127	Total 127	O 127	0	0

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.

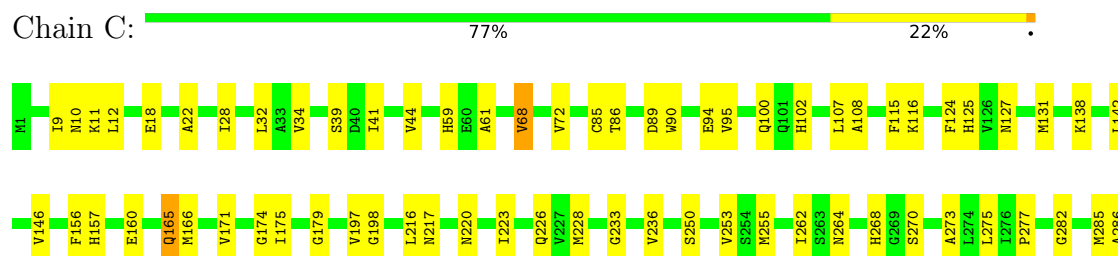
• Molecule 1: NADP-DEPENDENT ALCOHOL DEHYDROGENASE



• Molecule 1: NADP-DEPENDENT ALCOHOL DEHYDROGENASE

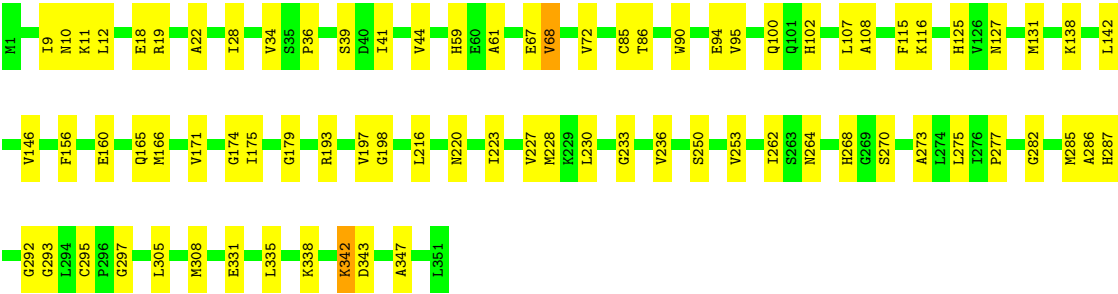
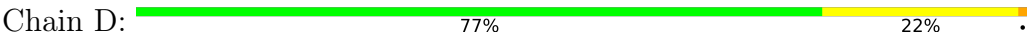


• Molecule 1: NADP-DEPENDENT ALCOHOL DEHYDROGENASE





● Molecule 1: NADP-DEPENDENT ALCOHOL DEHYDROGENASE



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	80.40Å 102.26Å 193.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.15	Depositor
% Data completeness (in resolution range)	79.0 (50.00-2.15)	Depositor
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.215 , 0.258	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	11122	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ZN

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	0.32	0/2689	0.58	0/3626
1	B	0.31	0/2689	0.57	0/3626
1	C	0.30	0/2689	0.58	0/3626
1	D	0.30	0/2689	0.58	0/3626
All	All	0.31	0/10756	0.58	0/14504

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 [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	2640	0	2681	61	0
1	B	2640	0	2681	64	0
1	C	2640	0	2681	64	0
1	D	2640	0	2681	60	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	145	0	0	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	141	0	0	8	0
3	C	145	0	0	8	0
3	D	127	0	0	4	0
All	All	11122	0	10724	215	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 10.

All (215) 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:197:VAL:HG21	1:A:223:ILE:HD11	1.53	0.90
1:C:275:LEU:HD23	1:D:273:ALA:HB1	1.52	0.89
1:C:273:ALA:HB1	1:D:275:LEU:HD23	1.55	0.89
1:C:197:VAL:HG21	1:C:223:ILE:HD11	1.53	0.88
1:D:197:VAL:HG21	1:D:223:ILE:HD11	1.55	0.88
1:B:197:VAL:HG21	1:B:223:ILE:HD11	1.58	0.84
1:B:91:ARG:HD3	3:B:482:HOH:O	1.83	0.77
1:B:68:VAL:HG13	1:B:72:VAL:HB	1.68	0.76
1:C:275:LEU:HG	1:D:275:LEU:HG	1.67	0.75
1:A:68:VAL:HG13	1:A:72:VAL:HB	1.69	0.74
1:C:68:VAL:HG13	1:C:72:VAL:HB	1.69	0.74
1:D:67:GLU:HB2	3:D:406:HOH:O	1.87	0.73
1:C:102:HIS:HD2	1:C:107:LEU:H	1.36	0.73
1:A:275:LEU:HG	1:B:275:LEU:HG	1.71	0.72
1:A:102:HIS:HD2	1:A:107:LEU:H	1.37	0.72
1:A:276:ILE:HD11	1:B:276:ILE:HD11	1.72	0.71
1:D:68:VAL:HG13	1:D:72:VAL:HB	1.71	0.71
1:A:228:MET:HG3	1:A:233:GLY:HA2	1.73	0.71
1:D:102:HIS:HD2	1:D:107:LEU:H	1.37	0.71
1:B:102:HIS:HD2	1:B:107:LEU:H	1.38	0.69
1:A:67:GLU:HB2	3:A:448:HOH:O	1.91	0.69
1:D:9:ILE:HG22	1:D:10:ASN:HD22	1.63	0.63
1:A:281:TRP:HZ3	1:B:274:LEU:HD22	1.63	0.63
1:C:228:MET:HG3	1:C:233:GLY:HA2	1.82	0.62
1:D:228:MET:HG3	1:D:233:GLY:HA2	1.81	0.62
1:A:278:ARG:HD3	1:B:271:GLY:O	1.98	0.62
1:B:9:ILE:HG22	1:B:10:ASN:HD22	1.64	0.62
1:A:278:ARG:HE	1:B:272:ASP:HA	1.64	0.62
1:C:9:ILE:HG22	1:C:10:ASN:HD22	1.63	0.61
1:C:217:ASN:HB2	3:C:444:HOH:O	2.00	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:285:MET:HA	1:D:293:GLY:HA2	1.81	0.61
1:B:228:MET:HG3	1:B:233:GLY:HA2	1.83	0.60
1:A:9:ILE:HG22	1:A:10:ASN:HD22	1.67	0.60
1:A:287:HIS:HA	1:B:292:GLY:O	2.02	0.60
1:A:305:LEU:HD23	1:A:308:MET:CE	2.33	0.59
1:C:285:MET:HB2	1:D:107:LEU:HD21	1.85	0.59
1:B:305:LEU:HD23	1:B:308:MET:CE	2.32	0.58
1:A:171:VAL:HG23	1:A:236:VAL:HG11	1.86	0.58
1:C:285:MET:HA	1:D:293:GLY:CA	2.34	0.57
1:A:275:LEU:HB3	3:A:463:HOH:O	2.04	0.57
1:C:85:CYS:SG	1:C:295:CYS:SG	3.03	0.57
1:D:85:CYS:SG	1:D:295:CYS:SG	3.02	0.56
1:C:197:VAL:HG11	1:C:223:ILE:HD12	1.87	0.56
1:A:175:ILE:HG12	1:A:198:GLY:HA3	1.87	0.56
1:D:197:VAL:HG11	1:D:223:ILE:HD12	1.88	0.56
1:C:107:LEU:HD21	1:D:285:MET:HB2	1.88	0.56
1:C:305:LEU:HD23	1:C:308:MET:CE	2.35	0.56
1:C:175:ILE:HG12	1:C:198:GLY:HA3	1.88	0.55
1:B:34:VAL:HG12	1:B:61:ALA:HB2	1.87	0.55
1:C:86:THR:HB	1:C:297:GLY:HA3	1.89	0.55
1:D:171:VAL:HG23	1:D:236:VAL:HG11	1.87	0.55
1:A:197:VAL:HG11	1:A:223:ILE:HD12	1.88	0.54
1:D:34:VAL:HG12	1:D:61:ALA:HB2	1.90	0.54
1:A:271:GLY:O	1:B:278:ARG:HD3	2.07	0.54
1:B:171:VAL:HG23	1:B:236:VAL:HG11	1.89	0.54
1:C:102:HIS:HE1	1:D:287:HIS:HB2	1.73	0.54
1:C:171:VAL:HG23	1:C:236:VAL:HG11	1.88	0.54
1:D:305:LEU:HD23	1:D:308:MET:CE	2.38	0.53
1:A:281:TRP:CZ3	1:B:274:LEU:HD22	2.43	0.53
1:B:175:ILE:HG12	1:B:198:GLY:HA3	1.90	0.53
1:B:127:ASN:HB3	3:B:372:HOH:O	2.08	0.53
1:A:287:HIS:CE1	1:B:293:GLY:HA3	2.43	0.53
1:D:175:ILE:HG12	1:D:198:GLY:HA3	1.90	0.53
1:A:272:ASP:HA	1:B:278:ARG:HE	1.74	0.53
1:C:287:HIS:HB2	1:D:102:HIS:HE1	1.74	0.53
1:A:34:VAL:HG12	1:A:61:ALA:HB2	1.91	0.52
1:C:34:VAL:HG12	1:C:61:ALA:HB2	1.91	0.52
1:B:22:ALA:HB2	1:B:28:ILE:HG12	1.91	0.52
1:B:329:HIS:HE1	3:B:487:HOH:O	1.91	0.52
1:A:85:CYS:SG	1:A:295:CYS:SG	3.08	0.52
1:A:86:THR:HB	1:A:297:GLY:HA3	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:86:THR:HB	1:D:297:GLY:HA3	1.93	0.51
1:A:39:SER:OG	1:A:59:HIS:HE1	1.94	0.51
1:C:116:LYS:NZ	1:C:125:HIS:HD2	2.09	0.51
1:B:197:VAL:HG11	1:B:223:ILE:HD12	1.93	0.51
1:B:308:MET:HE1	1:C:166:MET:HB3	1.92	0.51
1:C:90:TRP:HA	1:C:95:VAL:HG21	1.93	0.51
1:C:22:ALA:HB2	1:C:28:ILE:HG12	1.93	0.50
1:C:39:SER:OG	1:C:59:HIS:HE1	1.94	0.50
1:A:1:MET:HA	3:A:476:HOH:O	2.11	0.50
1:D:127:ASN:HB3	3:D:388:HOH:O	2.12	0.50
1:C:89:ASP:HB3	3:C:460:HOH:O	2.11	0.50
1:B:86:THR:HB	1:B:297:GLY:HA3	1.93	0.50
1:D:22:ALA:HB2	1:D:28:ILE:HG12	1.93	0.50
1:B:116:LYS:NZ	1:B:125:HIS:HD2	2.10	0.50
1:C:127:ASN:HB3	3:C:416:HOH:O	2.11	0.50
1:A:116:LYS:NZ	1:A:125:HIS:HD2	2.10	0.50
1:B:108:ALA:HA	1:B:115:PHE:HE2	1.77	0.49
1:A:278:ARG:NE	1:B:272:ASP:HA	2.26	0.49
1:A:22:ALA:HB2	1:A:28:ILE:HG12	1.92	0.49
1:D:39:SER:OG	1:D:59:HIS:HE1	1.96	0.49
1:D:116:LYS:NZ	1:D:125:HIS:HD2	2.10	0.49
1:A:90:TRP:CD1	1:A:131:MET:HB3	2.48	0.49
1:A:250:SER:HA	1:A:277:PRO:HD3	1.95	0.48
1:B:39:SER:OG	1:B:59:HIS:HE1	1.96	0.48
1:A:274:LEU:HD22	1:B:281:TRP:HZ3	1.79	0.48
1:A:216:LEU:HD21	1:A:230:LEU:HD11	1.96	0.48
1:C:293:GLY:HA2	1:D:285:MET:HA	1.96	0.48
1:C:108:ALA:HA	1:C:115:PHE:HE2	1.78	0.47
1:B:90:TRP:HA	1:B:95:VAL:HG21	1.96	0.47
1:D:156:PHE:O	1:D:160:GLU:HG3	2.14	0.47
1:A:292:GLY:O	1:B:287:HIS:HA	2.14	0.47
1:A:253:VAL:HA	1:A:262:ILE:HD11	1.97	0.47
1:A:108:ALA:HA	1:A:115:PHE:HE2	1.80	0.47
3:B:440:HOH:O	1:C:165:GLN:HG3	2.15	0.47
1:D:90:TRP:CD1	1:D:131:MET:HB3	2.50	0.47
1:B:218:TYR:HB2	3:B:485:HOH:O	2.14	0.46
1:D:216:LEU:HD21	1:D:230:LEU:HD11	1.97	0.46
1:A:90:TRP:HA	1:A:95:VAL:HG21	1.97	0.46
1:A:308:MET:HE1	1:D:166:MET:HB3	1.97	0.46
1:C:90:TRP:CD1	1:C:131:MET:HB3	2.51	0.46
1:C:124:PHE:HA	3:C:417:HOH:O	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:253:VAL:HA	1:C:262:ILE:HD11	1.97	0.46
1:A:331:GLU:O	1:A:335:LEU:HG	2.16	0.46
1:B:108:ALA:HA	1:B:115:PHE:CE2	2.51	0.46
1:A:188:LEU:HD13	3:A:475:HOH:O	2.15	0.46
1:B:90:TRP:CD1	1:B:131:MET:HB3	2.51	0.46
1:A:12:LEU:HD21	1:A:44:VAL:HG21	1.98	0.46
1:A:174:GLY:O	1:A:179:GLY:HA3	2.16	0.46
1:B:12:LEU:HD21	1:B:44:VAL:HG21	1.98	0.46
1:D:11:LYS:NZ	1:D:11:LYS:HB3	2.31	0.46
1:C:138:LYS:NZ	1:C:138:LYS:HB3	2.31	0.46
1:B:11:LYS:HB3	1:B:11:LYS:NZ	2.30	0.45
1:C:285:MET:CA	1:D:293:GLY:HA2	2.45	0.45
1:D:12:LEU:HD21	1:D:44:VAL:HG21	1.99	0.45
1:B:174:GLY:O	1:B:179:GLY:HA3	2.17	0.45
1:C:108:ALA:HA	1:C:115:PHE:CE2	2.51	0.45
1:C:264:ASN:OD1	1:C:268:HIS:HE1	2.00	0.45
1:D:90:TRP:HA	1:D:95:VAL:HG21	1.97	0.45
1:A:108:ALA:HA	1:A:115:PHE:CE2	2.52	0.45
1:B:253:VAL:HA	1:B:262:ILE:HD11	1.99	0.45
1:B:86:THR:O	1:B:297:GLY:HA3	2.16	0.45
1:D:108:ALA:HA	1:D:115:PHE:HE2	1.81	0.45
1:B:211:GLY:HA2	3:C:360:HOH:O	2.16	0.45
1:C:142:LEU:O	1:C:146:VAL:HG23	2.16	0.45
1:C:216:LEU:HD22	1:C:226:GLN:NE2	2.31	0.44
1:D:174:GLY:O	1:D:179:GLY:HA3	2.17	0.44
1:A:156:PHE:O	1:A:160:GLU:HG3	2.17	0.44
1:B:250:SER:HA	1:B:277:PRO:HD3	1.98	0.44
1:C:11:LYS:NZ	1:C:11:LYS:HB3	2.33	0.44
1:C:12:LEU:HD21	1:C:44:VAL:HG21	2.00	0.44
1:C:86:THR:O	1:C:297:GLY:HA3	2.17	0.44
1:C:102:HIS:CE1	1:D:287:HIS:HB2	2.51	0.44
1:C:327:PHE:HB3	3:C:419:HOH:O	2.16	0.44
1:D:142:LEU:O	1:D:146:VAL:HG23	2.17	0.44
1:A:86:THR:O	1:A:297:GLY:HA3	2.17	0.44
1:A:138:LYS:NZ	1:A:138:LYS:HB3	2.33	0.44
1:D:253:VAL:HA	1:D:262:ILE:HD11	1.99	0.44
1:A:342:LYS:HD3	1:A:343:ASP:N	2.33	0.44
1:A:113:SER:HB3	3:A:439:HOH:O	2.17	0.44
1:A:216:LEU:HD22	1:A:226:GLN:NE2	2.33	0.44
1:B:331:GLU:O	1:B:335:LEU:HG	2.17	0.44
1:C:156:PHE:O	1:C:160:GLU:HG3	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:250:SER:HA	1:D:277:PRO:HD3	1.99	0.44
1:A:11:LYS:NZ	1:A:11:LYS:HB3	2.32	0.44
1:D:108:ALA:HA	1:D:115:PHE:CE2	2.53	0.44
1:D:331:GLU:O	1:D:335:LEU:HG	2.18	0.44
1:B:138:LYS:NZ	1:B:138:LYS:HB3	2.33	0.43
1:C:174:GLY:O	1:C:179:GLY:HA3	2.19	0.43
1:D:86:THR:O	1:D:297:GLY:HA3	2.19	0.43
1:D:305:LEU:HD23	1:D:308:MET:HE3	2.00	0.43
1:D:264:ASN:OD1	1:D:268:HIS:HE1	2.02	0.43
1:C:250:SER:HA	1:C:277:PRO:HD3	2.00	0.43
1:C:287:HIS:HB2	1:D:102:HIS:CE1	2.53	0.43
1:C:264:ASN:O	1:C:292:GLY:HA2	2.18	0.43
1:C:285:MET:C	1:D:293:GLY:HA2	2.37	0.43
1:C:331:GLU:O	1:C:335:LEU:HG	2.19	0.43
1:D:138:LYS:NZ	1:D:138:LYS:HB3	2.34	0.43
1:A:282:GLY:HA3	1:A:286:ALA:HB2	2.01	0.43
1:B:113:SER:HB3	3:B:395:HOH:O	2.18	0.42
1:B:171:VAL:HG11	1:B:255:MET:SD	2.58	0.42
1:B:282:GLY:HA3	1:B:286:ALA:HB2	2.00	0.42
1:D:36:PRO:HD2	1:D:347:ALA:O	2.20	0.42
1:B:216:LEU:HD21	1:B:230:LEU:HD11	2.00	0.42
1:B:223:ILE:O	1:B:227:VAL:HG23	2.19	0.42
1:A:37:CYS:HA	3:A:364:HOH:O	2.19	0.42
1:D:223:ILE:O	1:D:227:VAL:HG23	2.19	0.42
1:D:264:ASN:O	1:D:292:GLY:HA2	2.19	0.42
1:A:223:ILE:O	1:A:227:VAL:HG23	2.19	0.42
1:B:89:ASP:HB3	3:B:436:HOH:O	2.19	0.42
1:B:2:LYS:HE3	3:B:486:HOH:O	2.19	0.42
1:B:41:ILE:HD12	1:B:338:LYS:HB2	2.02	0.42
1:C:94:GLU:HG2	1:C:102:HIS:O	2.20	0.42
1:C:282:GLY:HA3	1:C:286:ALA:HB2	2.02	0.42
1:C:292:GLY:O	1:D:287:HIS:HA	2.20	0.42
1:B:156:PHE:O	1:B:160:GLU:HG3	2.20	0.42
1:D:41:ILE:HD12	1:D:338:LYS:HB2	2.02	0.42
1:C:157:HIS:HA	1:C:160:GLU:OE1	2.20	0.41
1:C:305:LEU:HA	1:C:308:MET:HE3	2.00	0.41
1:C:342:LYS:HD3	1:C:343:ASP:N	2.35	0.41
1:B:264:ASN:OD1	1:B:268:HIS:HE1	2.03	0.41
1:B:264:ASN:O	1:B:292:GLY:HA2	2.19	0.41
1:D:342:LYS:HD3	1:D:343:ASP:N	2.34	0.41
1:B:305:LEU:HA	1:B:308:MET:HE3	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:171:VAL:HG11	1:C:255:MET:SD	2.61	0.41
1:D:19:ARG:NH2	3:D:406:HOH:O	2.53	0.41
1:A:264:ASN:O	1:A:292:GLY:HA2	2.21	0.41
1:D:94:GLU:HG2	1:D:102:HIS:O	2.21	0.41
1:A:94:GLU:HG2	1:A:102:HIS:O	2.21	0.41
1:C:138:LYS:HA	3:C:476:HOH:O	2.21	0.41
1:A:283:CYS:HB3	1:B:268:HIS:O	2.20	0.41
1:A:287:HIS:HB2	1:B:102:HIS:HE1	1.86	0.41
1:C:32:LEU:HD21	3:C:477:HOH:O	2.21	0.41
1:D:282:GLY:HA3	1:D:286:ALA:HB2	2.02	0.41
1:B:151:MET:SD	1:B:346:LYS:HD2	2.61	0.41
1:B:305:LEU:HD23	1:B:308:MET:HE3	1.99	0.41
1:A:142:LEU:O	1:A:146:VAL:HG23	2.21	0.41
1:A:259:GLY:HA2	1:A:287:HIS:HB3	2.03	0.41
1:D:286:ALA:N	3:D:421:HOH:O	2.54	0.41
1:A:311:TYR:CE2	1:D:193:ARG:HB2	2.56	0.40
1:B:216:LEU:HD22	1:B:226:GLN:NE2	2.35	0.40
1:A:293:GLY:HA3	1:B:287:HIS:CE1	2.56	0.40
1:B:9:ILE:O	1:B:10:ASN:HB2	2.22	0.40
1:C:41:ILE:HD12	1:C:338:LYS:HB2	2.01	0.40
1:C:305:LEU:HD23	1:C:308:MET:HE3	2.02	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	349/351 (99%)	328 (94%)	20 (6%)	1 (0%)	41	37
1	B	349/351 (99%)	328 (94%)	20 (6%)	1 (0%)	41	37
1	C	349/351 (99%)	328 (94%)	20 (6%)	1 (0%)	41	37
1	D	349/351 (99%)	327 (94%)	21 (6%)	1 (0%)	41	37

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	1396/1404 (99%)	1311 (94%)	81 (6%)	4 (0%)	41	37

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	270	SER
1	B	270	SER
1	C	270	SER
1	D	270	SER

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	279/279 (100%)	273 (98%)	6 (2%)	52	55
1	B	279/279 (100%)	273 (98%)	6 (2%)	52	55
1	C	279/279 (100%)	273 (98%)	6 (2%)	52	55
1	D	279/279 (100%)	273 (98%)	6 (2%)	52	55
All	All	1116/1116 (100%)	1092 (98%)	24 (2%)	52	55

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	GLU
1	A	68	VAL
1	A	100	GLN
1	A	165	GLN
1	A	220	ASN
1	A	342	LYS
1	B	18	GLU
1	B	68	VAL
1	B	100	GLN
1	B	165	GLN
1	B	220	ASN

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Mol	Chain	Res	Type
1	B	342	LYS
1	C	18	GLU
1	C	68	VAL
1	C	100	GLN
1	C	165	GLN
1	C	220	ASN
1	C	342	LYS
1	D	18	GLU
1	D	68	VAL
1	D	100	GLN
1	D	165	GLN
1	D	220	ASN
1	D	342	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (39) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	10	ASN
1	A	102	HIS
1	A	125	HIS
1	A	144	ASN
1	A	157	HIS
1	A	220	ASN
1	A	222	HIS
1	A	226	GLN
1	A	251	GLN
1	A	268	HIS
1	B	10	ASN
1	B	102	HIS
1	B	125	HIS
1	B	144	ASN
1	B	157	HIS
1	B	220	ASN
1	B	222	HIS
1	B	226	GLN
1	B	232	ASN
1	B	268	HIS
1	C	10	ASN
1	C	54	ASN
1	C	102	HIS
1	C	125	HIS
1	C	144	ASN

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Mol	Chain	Res	Type
1	C	157	HIS
1	C	220	ASN
1	C	222	HIS
1	C	226	GLN
1	C	268	HIS
1	D	10	ASN
1	D	102	HIS
1	D	125	HIS
1	D	144	ASN
1	D	157	HIS
1	D	220	ASN
1	D	226	GLN
1	D	232	ASN
1	D	268	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.