



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2014 – 09:34 AM GMT

PDB ID : 1ADB
Title : CRYSTALLOGRAPHIC STUDIES OF ISOSTERIC NAD ANALOGUES
BOUND TO ALCOHOL DEHYDROGENASE: SPECIFICITY AND SUB-
STRATE BINDING IN TWO TERNARY COMPLEXES
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stein, B.M.
Deposited on : 1993-12-13
Resolution : 2.40 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

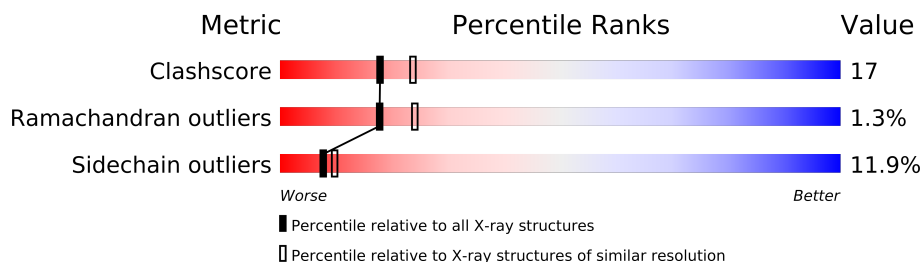
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 21963
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

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	79885	2789 (2.40-2.40)
Ramachandran outliers	78287	2736 (2.40-2.40)
Sidechain outliers	78261	2737 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	374	
1	B	374	

2 Entry composition i

There are 5 unique types of molecules in this entry. The entry contains 7165 atoms, of which 1410 are hydrogens and 0 are deuterium.

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.

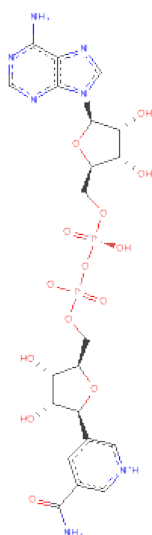
- Molecule 1 is a protein called ALCOHOL DEHYDROGENASE.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	374	Total	C	H	N	O	S	0	0	0
			3391	1769	607	472	520	23			
1	B	374	Total	C	H	N	O	S	0	0	0
			3392	1769	607	472	521	23			

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

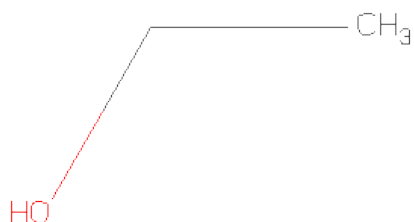
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	2	Total	Zn	0	0
			2	2		
2	A	2	Total	Zn	0	0
			2	2		

- Molecule 3 is 5-BETA-D-RIBOFURANOSYLNICOTINAMIDEADENINE DINUCLEOTIDE (three-letter code: CND) (formula: C₂₁H₂₇N₇O₁₄P₂).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	H	N	O	P	0	0
			52	21	8	7	14	2		
3	B	1	Total	C	H	N	O	P	0	0
			52	21	8	7	14	2		

- Molecule 4 is ETHANOL (three-letter code: EOH) (formula: C₂H₆O).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	H	O	0	0
			5	2	2	1		
4	B	1	Total	C	H	O	0	0
			5	2	2	1		

- Molecule 5 is water.

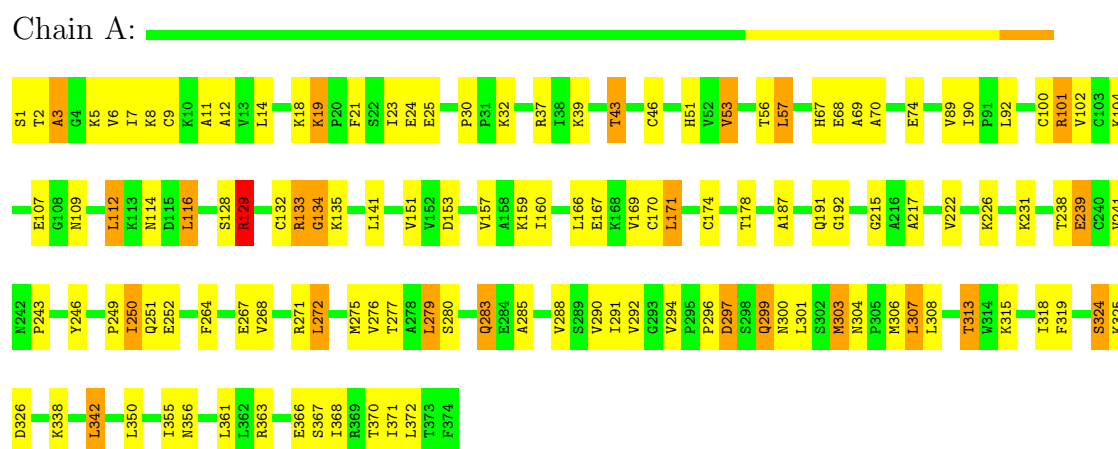
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	50	Total	H	O	0	0
			150	100	50		
5	B	38	Total	H	O	0	0
			114	76	38		

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 errors 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.

Note EDS was not executed.

• Molecule 1: ALCOHOL DEHYDROGENASE



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	52.10Å 44.70Å 93.50Å 103.30° 87.90° 70.40°	Depositor
Resolution (Å)	(Not available) – 2.40	Depositor
% Data completeness (in resolution range)	(Not available) ((Not available)-2.40)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.180 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7165	wwPDB-VP
Average B, all atoms (Å ²)	16.0	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, EOH, CND

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	0.53	0/2836	0.78	2/3834 (0.1%)
1	B	0.51	0/2837	0.80	1/3834 (0.0%)
All	All	0.52	0/5673	0.79	3/7668 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	57	LEU	CA-CB-CG	5.18	127.21	115.30
1	A	368	ILE	N-CA-C	-5.07	97.31	111.00
1	B	123	MET	N-CA-C	-5.05	97.36	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2784	607	2241	92	0
1	B	2785	607	2241	109	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	44	8	17	1	0
3	B	44	8	16	1	0
4	A	3	2	4	1	0
4	B	3	2	4	2	0
5	A	50	100	0	7	0
5	B	38	76	0	2	0
All	All	5755	1410	4523	196	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 17.

All (196) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:76:ILE:HD11	1:B:83:VAL:HG23	1.48	0.95
1:A:43:THR:HG23	1:A:69:ALA:HB2	1.49	0.93
1:B:129:ARG:HG2	1:B:151:VAL:HG11	1.55	0.89
1:A:251:GLN:HG3	1:A:277:THR:HG23	1.61	0.82
1:A:178:THR:HG22	1:A:319:PHE:HA	1.63	0.80
1:B:153:ASP:HB3	1:B:155:ILE:HG22	1.63	0.80
1:B:178:THR:HG22	1:B:319:PHE:HA	1.64	0.79
1:A:301:LEU:HD11	1:B:303:MET:HE2	1.68	0.76
1:B:313:THR:HG22	5:B:412:HOH:O	1.84	0.75
1:B:160:ILE:HD11	1:B:328:VAL:HG23	1.70	0.74
1:A:92:LEU:HD13	1:A:324:SER:HB3	1.69	0.74
1:B:251:GLN:HG3	1:B:277:THR:HG23	1.73	0.71
1:A:272:LEU:HD21	1:A:299:GLN:HB3	1.75	0.69
1:B:160:ILE:HG12	1:B:332:VAL:HG21	1.75	0.68
1:B:203:VAL:HG23	1:B:268:VAL:HG23	1.75	0.68
1:A:56:THR:HG23	1:A:296:PRO:HA	1.74	0.67
1:B:100:CYS:O	1:B:104:LYS:HG2	1.94	0.67
1:A:291:ILE:HG22	1:A:315:LYS:O	1.94	0.67
1:A:275:MET:SD	1:A:291:ILE:HD12	2.35	0.66
1:B:348:HIS:HB2	1:B:370:THR:HB	1.78	0.66
1:A:174:CYS:O	1:A:178:THR:HG23	1.96	0.65
1:B:120:ARG:HH21	5:B:397:HOH:H2	1.42	0.65
1:B:89:VAL:HG21	1:B:157:VAL:HG22	1.78	0.65
1:A:90:ILE:HD11	1:A:169:VAL:HG13	1.79	0.65
1:B:294:VAL:HG21	4:B:378:EOH:H23	1.79	0.64
1:A:169:VAL:HA	5:A:384:HOH:O	1.96	0.64
1:A:249:PRO:HB2	1:A:252:GLU:HG3	1.79	0.64
1:B:89:VAL:CG2	1:B:157:VAL:HG22	2.27	0.63
1:B:47:ARG:HD3	3:B:377:CND:O1A	1.99	0.63

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:7:ILE:HG13	1:B:37:ARG:NH2	2.13	0.63
1:A:361:LEU:HB3	1:A:367:SER:OG	1.98	0.63
1:B:129:ARG:HG2	1:B:151:VAL:CG1	2.31	0.60
1:A:68:GLU:HB3	1:A:174:CYS:HB3	1.84	0.60
1:B:203:VAL:HG23	1:B:268:VAL:CG2	2.32	0.59
1:A:338:LYS:HA	5:A:420:HOH:O	2.02	0.59
1:B:97:CYS:HB3	1:B:113:LYS:HG3	1.84	0.58
1:A:157:VAL:O	1:A:325:LYS:HE3	2.03	0.58
1:A:275:MET:SD	1:A:291:ILE:CD1	2.93	0.57
1:B:105:HIS:HD2	1:B:107:GLU:H	1.52	0.57
1:A:5:LYS:HA	1:A:30:PRO:HG3	1.86	0.57
1:A:21:PHE:H	1:A:356:ASN:HD21	1.52	0.57
1:A:114:ASN:ND2	1:A:116:LEU:H	2.02	0.57
1:B:160:ILE:CG1	1:B:332:VAL:HG21	2.35	0.56
1:A:187:ALA:HB2	1:A:290:VAL:HG21	1.87	0.56
1:A:153:ASP:HB3	5:A:403:HOH:O	2.06	0.56
1:A:2:THR:O	1:A:3:ALA:HB2	2.06	0.56
1:B:253:VAL:O	1:B:257:MET:HG3	2.05	0.56
1:A:222:VAL:HG22	1:A:241:VAL:CG1	2.35	0.56
1:A:102:VAL:HG23	1:A:112:LEU:HD22	1.88	0.55
1:A:251:GLN:CG	1:A:277:THR:HG23	2.35	0.55
1:A:178:THR:HG22	1:A:319:PHE:HD1	1.71	0.55
1:A:51:HIS:HB3	1:A:57:LEU:HB2	1.88	0.55
1:A:12:ALA:HB1	1:A:21:PHE:HB3	1.88	0.55
1:A:114:ASN:HD22	1:A:116:LEU:H	1.54	0.55
1:A:271:ARG:O	1:A:275:MET:HG3	2.07	0.54
1:A:170:CYS:SG	1:A:371:ILE:HD12	2.48	0.54
1:B:9:CYS:HB2	1:B:148:GLN:OE1	2.07	0.54
1:B:39:LYS:HG3	1:B:149:TYR:CE1	2.42	0.54
1:A:37:ARG:HG3	1:A:151:VAL:HG22	1.90	0.54
1:A:7:ILE:HG13	1:A:37:ARG:NH2	2.22	0.54
1:B:64:ILE:HG13	1:B:137:ILE:HG21	1.89	0.54
1:A:191:GLN:HG2	1:A:215:GLY:HA3	1.89	0.54
1:A:292:VAL:O	3:A:377:CND:H2N	2.09	0.53
1:B:11:ALA:O	1:B:23:ILE:HA	2.08	0.53
1:B:18:LYS:NZ	1:B:18:LYS:HB3	2.24	0.53
1:A:160:ILE:HD12	1:A:169:VAL:HG11	1.90	0.53
1:A:299:GLN:O	1:B:304:ASN:ND2	2.42	0.52
1:B:303:MET:HE1	1:B:305:PRO:HA	1.91	0.52
1:A:283:GLN:NE2	1:A:285:ALA:H	2.07	0.52
1:A:8:LYS:NZ	1:A:25:GLU:HG2	2.24	0.52
1:A:8:LYS:HZ2	1:A:25:GLU:HG2	1.74	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:31:PRO:HG3	1:B:37:ARG:HB2	1.91	0.52
1:B:165:PRO:HD2	1:B:336:MET:HE1	1.91	0.52
1:B:200:LEU:HD13	1:B:232:ALA:HB2	1.92	0.51
1:B:105:HIS:CD2	1:B:107:GLU:H	2.28	0.51
1:A:32:LYS:HG3	1:A:129:ARG:HH12	1.76	0.51
1:B:63:VAL:HG22	1:B:64:ILE:N	2.25	0.51
1:B:2:THR:O	1:B:3:ALA:HB2	2.11	0.51
1:A:178:THR:HG22	1:A:319:PHE:CD1	2.46	0.51
1:B:198:PHE:O	1:B:268:VAL:HG13	2.10	0.51
1:A:285:ALA:HB1	5:A:390:HOH:O	2.10	0.51
1:A:272:LEU:O	1:A:276:VAL:HG23	2.10	0.50
1:A:107:GLU:O	1:A:107:GLU:HG2	2.10	0.50
1:B:343:ASP:HB2	1:B:344:PRO:HD3	1.92	0.50
1:B:357:GLU:O	1:B:361:LEU:HD22	2.11	0.50
1:B:97:CYS:HA	1:B:155:ILE:HG13	1.93	0.50
1:B:308:LEU:HD12	1:B:312:ARG:O	2.12	0.50
1:A:297:ASP:HA	5:A:380:HOH:H1	1.74	0.50
1:B:33:ALA:HA	1:B:78:GLU:O	2.12	0.50
1:B:203:VAL:CG2	1:B:268:VAL:HG23	2.41	0.50
1:B:225:ASN:HD21	1:B:227:ASP:HB2	1.76	0.50
1:A:313:THR:HB	1:B:315:LYS:HG2	1.94	0.49
1:B:225:ASN:ND2	1:B:227:ASP:HB2	2.27	0.49
1:B:249:PRO:HB3	1:B:251:GLN:HE21	1.77	0.49
1:B:9:CYS:O	1:B:25:GLU:HA	2.11	0.49
1:B:164:SER:HA	1:B:336:MET:HE3	1.93	0.49
1:B:359:PHE:HB3	1:B:363:ARG:NH2	2.28	0.49
1:A:23:ILE:HD11	1:A:355:ILE:CG2	2.42	0.49
1:B:10:LYS:HA	1:B:24:GLU:O	2.13	0.49
1:B:88:LYS:HD2	1:B:162:ALA:HA	1.95	0.49
1:A:192:GLY:HA2	1:A:217:ALA:HB2	1.94	0.49
1:A:32:LYS:HG3	1:A:129:ARG:NH1	2.28	0.48
1:A:39:LYS:HE2	5:A:396:HOH:O	2.13	0.48
1:B:31:PRO:CG	1:B:37:ARG:HB2	2.43	0.48
1:B:43:THR:HG23	1:B:69:ALA:HB2	1.94	0.48
1:B:5:LYS:HG3	1:B:6:VAL:H	1.78	0.48
1:B:196:ALA:HB2	1:B:262:VAL:HG21	1.94	0.48
1:A:2:THR:HG22	1:A:3:ALA:H	1.79	0.48
1:A:350:LEU:HD13	1:A:370:THR:HG23	1.95	0.48
1:B:272:LEU:HG	1:B:301:LEU:HB3	1.96	0.47
1:B:10:LYS:O	1:B:148:GLN:HG3	2.15	0.47
1:A:267:GLU:HG3	1:A:275:MET:HA	1.96	0.47
1:B:39:LYS:HG3	1:B:149:TYR:CZ	2.49	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:238:THR:HG22	1:A:239:GLU:OE1	2.14	0.47
1:B:279:LEU:HD23	1:B:312:ARG:HD2	1.97	0.47
1:B:92:LEU:HD13	1:B:324:SER:HB2	1.97	0.46
1:A:46:CYS:HB3	1:A:67:HIS:CE1	2.50	0.46
1:A:283:GLN:HE21	1:A:285:ALA:HB3	1.79	0.46
1:B:73:VAL:HG11	1:B:76:ILE:HD13	1.98	0.46
1:A:222:VAL:HG22	1:A:241:VAL:HG13	1.98	0.46
1:A:116:LEU:HG	1:A:141:LEU:CD2	2.46	0.46
1:B:5:LYS:HA	1:B:5:LYS:HE3	1.99	0.45
1:B:48:SER:HA	1:B:51:HIS:CD2	2.51	0.45
1:B:174:CYS:SG	1:B:175:GLY:N	2.89	0.45
1:A:9:CYS:O	1:A:25:GLU:HA	2.16	0.45
1:A:303:MET:HG2	1:A:304:ASN:N	2.31	0.45
1:A:315:LYS:HG2	1:B:313:THR:HB	1.98	0.45
1:B:324:SER:O	1:B:328:VAL:HG13	2.15	0.45
1:B:187:ALA:HB2	1:B:290:VAL:HG21	1.98	0.45
1:B:36:VAL:O	1:B:151:VAL:HA	2.15	0.45
1:A:308:LEU:O	1:B:316:GLY:HA3	2.16	0.45
1:B:105:HIS:CD2	1:B:108:GLY:H	2.34	0.45
1:A:68:GLU:HB3	1:A:174:CYS:CB	2.47	0.45
1:A:100:CYS:O	1:A:104:LYS:HG2	2.17	0.45
1:B:204:GLY:HA2	1:B:268:VAL:HG21	1.99	0.45
1:B:335:PHE:CD1	1:B:342:LEU:HD22	2.52	0.45
1:B:294:VAL:CG2	4:B:378:EOH:H23	2.47	0.44
1:B:268:VAL:O	1:B:268:VAL:CG2	2.66	0.44
1:A:241:VAL:HG23	1:A:246:TYR:CE1	2.53	0.44
1:B:45:ILE:HG23	1:B:359:PHE:CZ	2.52	0.44
1:B:267:GLU:HG3	1:B:275:MET:HA	1.99	0.44
1:B:89:VAL:HG22	1:B:90:ILE:N	2.33	0.44
1:B:231:LYS:HG3	1:B:368:ILE:CD1	2.47	0.44
1:A:361:LEU:HA	1:A:361:LEU:HD23	1.90	0.44
1:B:105:HIS:HD2	1:B:107:GLU:N	2.14	0.44
1:A:5:LYS:HD2	1:A:30:PRO:HG3	2.00	0.44
1:B:165:PRO:HD2	1:B:336:MET:CE	2.47	0.44
1:B:229:PHE:CD1	1:B:229:PHE:N	2.85	0.44
1:A:57:LEU:HD23	1:A:294:VAL:HG11	2.00	0.43
1:B:307:LEU:HD23	1:B:312:ARG:NH2	2.32	0.43
1:A:128:SER:OG	1:A:129:ARG:N	2.51	0.43
1:B:135:LYS:HA	1:B:136:PRO:HD3	1.78	0.43
1:B:40:MET:CE	1:B:40:MET:HA	2.49	0.43
1:A:116:LEU:HD11	4:A:378:EOH:H21	1.99	0.43
1:B:61:LEU:HD12	1:B:63:VAL:HG12	2.00	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:101:ARG:HD2	1:A:101:ARG:HA	1.79	0.43
1:A:6:VAL:N	1:A:30:PRO:HD3	2.34	0.43
1:A:24:GLU:OE1	1:A:132:CYS:SG	2.77	0.43
1:A:226:LYS:HE2	1:A:226:LYS:HB3	1.90	0.43
1:B:294:VAL:HA	1:B:295:PRO:HD2	1.95	0.42
1:B:48:SER:HB2	1:B:67:HIS:HE1	1.84	0.42
1:B:231:LYS:O	1:B:235:VAL:HG22	2.18	0.42
1:B:331:LEU:HD23	1:B:331:LEU:HA	1.83	0.42
1:B:192:GLY:HA2	1:B:217:ALA:HB2	2.01	0.42
1:A:342:LEU:CD2	5:A:384:HOH:O	2.67	0.42
1:B:37:ARG:HG2	1:B:74:GLU:OE1	2.20	0.42
1:B:272:LEU:HD21	1:B:299:GLN:HB3	2.01	0.42
1:A:18:LYS:O	1:A:19:LYS:HD2	2.19	0.42
1:B:175:GLY:HA2	1:B:203:VAL:HG12	2.00	0.42
1:A:264:PHE:HA	1:A:288:VAL:O	2.19	0.42
1:A:53:VAL:HG21	1:A:363:ARG:NH2	2.34	0.42
1:A:70:ALA:HB1	1:A:166:LEU:HD22	2.01	0.42
1:A:89:VAL:HG12	1:A:159:LYS:HA	2.01	0.42
1:B:345:LEU:O	1:B:369:ARG:HB2	2.19	0.42
1:A:132:CYS:O	1:A:134:GLY:N	2.54	0.41
1:B:165:PRO:HG2	1:B:168:LYS:HB2	2.02	0.41
1:A:283:GLN:HE21	1:A:285:ALA:H	1.67	0.41
1:A:279:LEU:HD13	1:A:307:LEU:HD23	2.02	0.41
1:B:242:ASN:HD22	1:B:243:PRO:N	2.17	0.41
1:B:242:ASN:HD22	1:B:242:ASN:C	2.24	0.41
1:A:114:ASN:HD22	1:A:116:LEU:N	2.17	0.41
1:B:48:SER:O	1:B:52:VAL:HG23	2.21	0.41
1:B:269:ILE:HG21	1:B:271:ARG:HD2	2.02	0.41
1:B:5:LYS:HG3	1:B:6:VAL:N	2.35	0.41
1:A:243:PRO:HG3	1:A:250:ILE:HG13	2.03	0.41
1:B:204:GLY:CA	1:B:268:VAL:HG21	2.51	0.41
1:A:171:LEU:HD12	1:A:171:LEU:HA	1.91	0.41
1:B:195:CYS:HA	1:B:264:PHE:O	2.21	0.40
1:B:268:VAL:O	1:B:268:VAL:HG22	2.21	0.40
1:A:32:LYS:CG	1:A:129:ARG:HH12	2.35	0.40
1:B:42:ALA:HA	1:B:372:LEU:O	2.22	0.40
1:B:328:VAL:N	1:B:329:PRO:HD2	2.36	0.40
1:A:37:ARG:HD3	1:A:74:GLU:OE1	2.22	0.40
1:B:63:VAL:CG2	1:B:64:ILE:N	2.85	0.40
1:A:11:ALA:O	1:A:23:ILE:HA	2.22	0.40
1:B:366:GLU:H	1:B:366:GLU:HG2	1.74	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	372/374 (100%)	331 (89%)	36 (10%)	5 (1%)	18	24
1	B	372/374 (100%)	325 (87%)	42 (11%)	5 (1%)	18	24
All	All	744/748 (100%)	656 (88%)	78 (10%)	10 (1%)	18	24

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	297	ASP
1	B	174	CYS
1	B	297	ASP
1	B	2	THR
1	B	175	GLY
1	A	3	ALA
1	A	133	ARG
1	A	129	ARG
1	B	4	GLY
1	A	134	GLY

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	308/308 (100%)	274 (89%)	34 (11%)	9	12
1	B	308/308 (100%)	269 (87%)	39 (13%)	6	7
All	All	616/616 (100%)	543 (88%)	73 (12%)	8	10

All (73) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	SER
1	A	14	LEU
1	A	19	LYS
1	A	43	THR
1	A	53	VAL
1	A	101	ARG
1	A	109	ASN
1	A	112	LEU
1	A	116	LEU
1	A	129	ARG
1	A	133	ARG
1	A	135	LYS
1	A	167	GLU
1	A	171	LEU
1	A	231	LYS
1	A	239	GLU
1	A	250	ILE
1	A	268	VAL
1	A	272	LEU
1	A	279	LEU
1	A	280	SER
1	A	283	GLN
1	A	299	GLN
1	A	300	ASN
1	A	303	MET
1	A	306	MET
1	A	307	LEU
1	A	313	THR
1	A	318	ILE
1	A	324	SER
1	A	326	ASP
1	A	342	LEU
1	A	366	GLU
1	A	372	LEU
1	B	5	LYS
1	B	14	LEU
1	B	18	LYS
1	B	37	ARG
1	B	40	MET
1	B	84	ARG
1	B	85	PRO
1	B	112	LEU
1	B	122	THR

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Mol	Chain	Res	Type
1	B	129	ARG
1	B	133	ARG
1	B	141	LEU
1	B	151	VAL
1	B	153	ASP
1	B	159	LYS
1	B	168	LYS
1	B	171	LEU
1	B	186	VAL
1	B	190	THR
1	B	212	LYS
1	B	225	ASN
1	B	226	LYS
1	B	242	ASN
1	B	268	VAL
1	B	271	ARG
1	B	272	LEU
1	B	274	THR
1	B	303	MET
1	B	306	MET
1	B	308	LEU
1	B	313	THR
1	B	330	LYS
1	B	342	LEU
1	B	345	LEU
1	B	361	LEU
1	B	362	LEU
1	B	366	GLU
1	B	370	THR
1	B	372	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	114	ASN
1	A	124	GLN
1	A	283	GLN
1	A	356	ASN
1	B	105	HIS
1	B	138	HIS
1	B	225	ASN
1	B	242	ASN

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Mol	Chain	Res	Type
1	B	251	GLN
1	B	304	ASN

5.3.3 RNA ⓘ

There are no RNA chains 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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	CND	A	377	2	48,48,48	1.27	6 (12%)	73,73,73	2.15	18 (24%)
4	EOH	A	378	-	2,2,2	0.54	0	1,1,1	0.37	0
3	CND	B	377	2	48,48,48	1.45	8 (16%)	73,73,73	2.05	17 (23%)
4	EOH	B	378	-	2,2,2	0.61	0	1,1,1	0.28	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	CND	A	377	2	-	0/30/62/62	0/3/5/5
4	EOH	A	378	-	-	0/0/0/0	0/0/0/0
3	CND	B	377	2	-	0/30/62/62	0/3/5/5
4	EOH	B	378	-	-	0/0/0/0	0/0/0/0

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	377	CND	C6N-N5N	3.85	1.42	1.34
3	A	377	CND	C4A-N9A	-3.84	1.32	1.37
3	B	377	CND	C4A-N9A	-3.48	1.32	1.37
3	B	377	CND	C3N-C7N	3.25	1.56	1.50
3	B	377	CND	C4N-C3N	3.13	1.42	1.38
3	A	377	CND	C4N-N5N	2.87	1.40	1.34
3	A	377	CND	C2D-C1D	-2.76	1.51	1.53
3	A	377	CND	C6N-N5N	2.54	1.39	1.34
3	B	377	CND	PA-O3	2.50	1.64	1.59
3	B	377	CND	C6N-C1N	2.48	1.40	1.38
3	A	377	CND	C1N-C1D	-2.43	1.46	1.51
3	B	377	CND	C4N-N5N	2.32	1.39	1.34
3	B	377	CND	O4B-C1B	2.30	1.44	1.41
3	A	377	CND	PN-O3	-2.09	1.56	1.60

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	377	CND	N3A-C2A-N1A	-7.95	122.06	128.71
3	B	377	CND	N3A-C4A-N9A	7.01	138.08	125.43
3	A	377	CND	N3A-C4A-N9A	6.56	137.28	125.43
3	A	377	CND	N3A-C2A-N1A	-5.44	124.16	128.71
3	B	377	CND	C4A-C5A-N7A	5.42	114.17	109.52
3	A	377	CND	C3D-C2D-C1D	-5.10	95.87	101.85
3	A	377	CND	O4D-C1D-C1N	-5.10	102.86	110.26
3	B	377	CND	O4D-C1D-C1N	4.43	116.68	110.26
3	A	377	CND	O4B-C1B-N9A	4.24	112.38	108.44
3	A	377	CND	C6N-C1N-C2N	4.03	123.43	117.16
3	A	377	CND	C4A-C5A-N7A	4.02	112.97	109.52
3	A	377	CND	C6A-C5A-C4A	3.84	124.29	117.25
3	A	377	CND	C5A-C4A-N3A	-3.69	117.67	125.70
3	B	377	CND	O4B-C1B-N9A	3.64	111.82	108.44
3	B	377	CND	C5A-C4A-N3A	-3.43	118.24	125.70
3	A	377	CND	O2D-C2D-C1D	3.30	119.47	111.93
3	A	377	CND	C4B-O4B-C1B	-3.14	106.34	109.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	377	CND	C6N-N5N-C4N	-3.08	117.02	122.65
3	A	377	CND	C3N-C2N-C1N	-2.97	118.41	121.14
3	B	377	CND	PN-O3-PA	2.96	145.67	132.95
3	A	377	CND	C3N-C7N-N7N	-2.81	114.57	117.77
3	B	377	CND	C6A-C5A-C4A	2.70	122.20	117.25
3	B	377	CND	N6A-C6A-N1A	2.52	124.31	119.36
3	A	377	CND	C1N-C1D-C2D	2.47	119.74	115.63
3	B	377	CND	C5A-C4A-N9A	-2.41	103.68	107.16
3	B	377	CND	C8A-N9A-C4A	2.39	108.72	106.90
3	A	377	CND	C4D-O4D-C1D	-2.28	106.74	109.45
3	A	377	CND	C6A-C5A-N7A	-2.25	122.74	131.34
3	B	377	CND	C3D-C2D-C1D	-2.18	99.29	101.85
3	A	377	CND	N6A-C6A-N1A	2.10	123.48	119.36
3	B	377	CND	C1B-N9A-C4A	-2.08	123.03	126.64
3	B	377	CND	C5B-C4B-C3B	-2.05	107.02	115.21
3	B	377	CND	O2A-PA-O1A	2.04	123.63	112.21
3	A	377	CND	C6N-C1N-C1D	-2.04	117.53	121.43
3	B	377	CND	C6A-C5A-N7A	-2.02	123.62	131.34

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.