



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:30 AM GMT

PDB ID : 2XAA
Title : ALCOHOL DEHYDROGENASE ADH-'A' FROM RHODOCOCCLUS RUBER DSM 44541 AT PH 8.5 IN COMPLEX WITH NAD AND BUTANE-1,4-DIOL
Authors : Kroutil, W.; Gruber, K.; Grogan, G.
Deposited on : 2010-03-30
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
<http://wwpdb.org/validation/2016/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
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

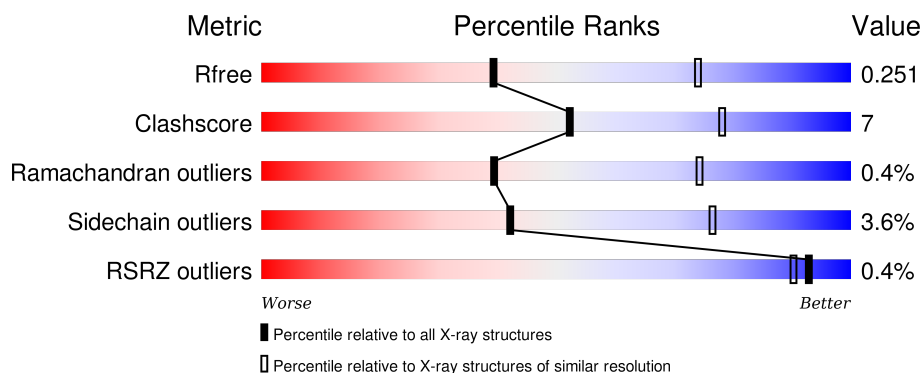
1 Overall quality at a glance

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}	91344	2393 (2.80-2.80)
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (2.80-2.80)
RSRZ outliers	91569	2404 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	345	<div> <div></div> <div>80%15% . .</div> </div>
1	B	345	<div> <div></div> <div>76%22% ..</div> </div>
1	C	345	<div> <div></div> <div>87%12%</div> </div>
1	D	345	<div> <div></div> <div>87%12% .</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	BU1	D	1347	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 10078 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 SECONDARY ALCOHOL DEHYDROGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	333	Total	C	N	O	S	0	1	0
			2370	1491	417	449	13			
1	B	342	Total	C	N	O	S	0	0	0
			2434	1533	429	460	12			
1	C	344	Total	C	N	O	S	0	0	0
			2461	1554	431	464	12			
1	D	345	Total	C	N	O	S	0	0	0
			2469	1558	433	466	12			

There are 36 discrepancies between the modelled and reference sequences:

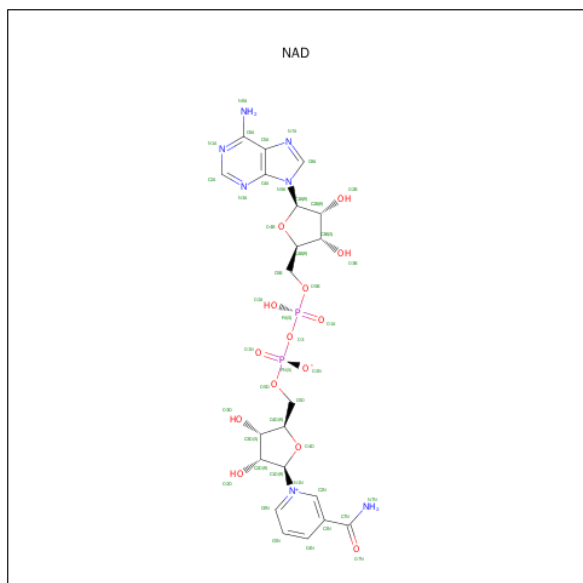
Chain	Residue	Modelled	Actual	Comment	Reference
A	4	VAL	LEU	CONFLICT	UNP Q8KLT9
A	18	ILE	VAL	CONFLICT	UNP Q8KLT9
A	22	THR	ALA	CONFLICT	UNP Q8KLT9
A	73	GLU	ALA	CONFLICT	UNP Q8KLT9
A	80	VAL	THR	CONFLICT	UNP Q8KLT9
A	111	ASP	GLU	CONFLICT	UNP Q8KLT9
A	238	GLN	GLU	CONFLICT	UNP Q8KLT9
A	262	VAL	ILE	CONFLICT	UNP Q8KLT9
A	303	GLU	ASP	CONFLICT	UNP Q8KLT9
B	4	VAL	LEU	CONFLICT	UNP Q8KLT9
B	18	ILE	VAL	CONFLICT	UNP Q8KLT9
B	22	THR	ALA	CONFLICT	UNP Q8KLT9
B	73	GLU	ALA	CONFLICT	UNP Q8KLT9
B	80	VAL	THR	CONFLICT	UNP Q8KLT9
B	111	ASP	GLU	CONFLICT	UNP Q8KLT9
B	238	GLN	GLU	CONFLICT	UNP Q8KLT9
B	262	VAL	ILE	CONFLICT	UNP Q8KLT9
B	303	GLU	ASP	CONFLICT	UNP Q8KLT9
C	4	VAL	LEU	CONFLICT	UNP Q8KLT9
C	18	ILE	VAL	CONFLICT	UNP Q8KLT9
C	22	THR	ALA	CONFLICT	UNP Q8KLT9

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Chain	Residue	Modelled	Actual	Comment	Reference
C	73	GLU	ALA	CONFLICT	UNP Q8KLT9
C	80	VAL	THR	CONFLICT	UNP Q8KLT9
C	111	ASP	GLU	CONFLICT	UNP Q8KLT9
C	238	GLN	GLU	CONFLICT	UNP Q8KLT9
C	262	VAL	ILE	CONFLICT	UNP Q8KLT9
C	303	GLU	ASP	CONFLICT	UNP Q8KLT9
D	4	VAL	LEU	CONFLICT	UNP Q8KLT9
D	18	ILE	VAL	CONFLICT	UNP Q8KLT9
D	22	THR	ALA	CONFLICT	UNP Q8KLT9
D	73	GLU	ALA	CONFLICT	UNP Q8KLT9
D	80	VAL	THR	CONFLICT	UNP Q8KLT9
D	111	ASP	GLU	CONFLICT	UNP Q8KLT9
D	238	GLN	GLU	CONFLICT	UNP Q8KLT9
D	262	VAL	ILE	CONFLICT	UNP Q8KLT9
D	303	GLU	ASP	CONFLICT	UNP Q8KLT9

- Molecule 2 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: C₂₁H₂₇N₇O₁₄P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 44	C 21	N 7	O 14	P 2	0	0
2	B	1	Total 44	C 21	N 7	O 14	P 2	0	0
2	C	1	Total 44	C 21	N 7	O 14	P 2	0	0

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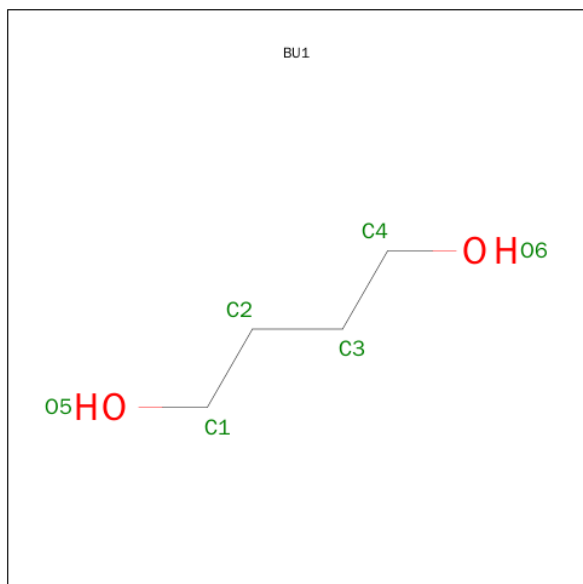
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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	D	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 4 is 1,4-BUTANEDIOL (three-letter code: BU1) (formula: C₄H₁₀O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			6	4	2		
4	C	1	Total	C	O	0	0
			6	4	2		
4	D	1	Total	C	O	0	0
			6	4	2		

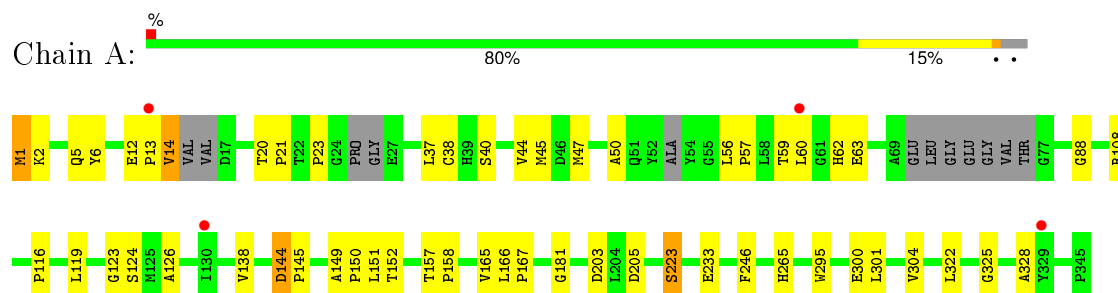
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	28	Total 28	O 28	0	0
5	B	39	Total 39	O 39	0	0
5	C	38	Total 38	O 38	0	0
5	D	37	Total 37	O 37	0	0

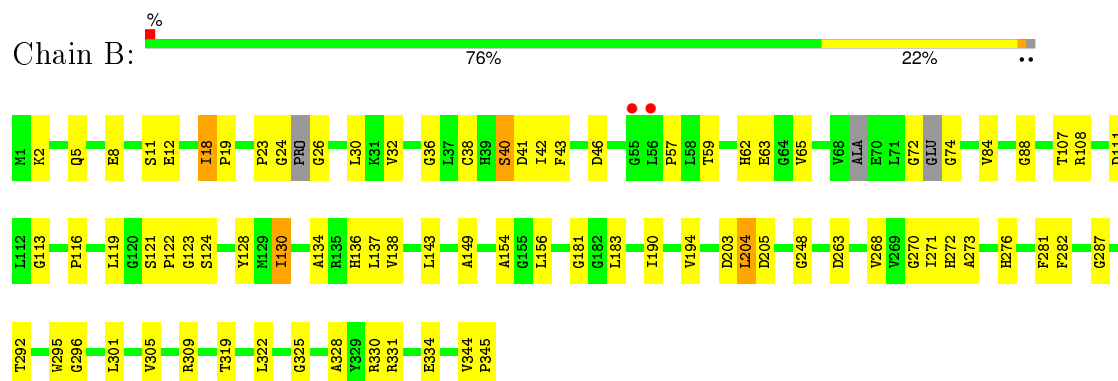
3 Residue-property plots [i](#)

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.

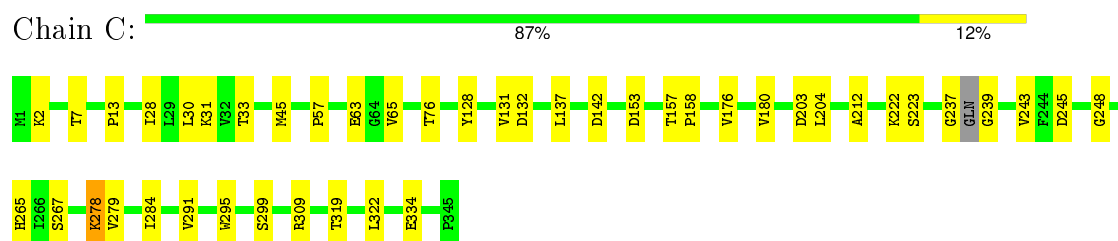
• Molecule 1: SECONDARY ALCOHOL DEHYDROGENASE



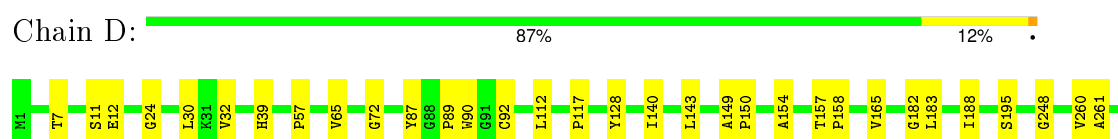
• Molecule 1: SECONDARY ALCOHOL DEHYDROGENASE



• Molecule 1: SECONDARY ALCOHOL DEHYDROGENASE



• Molecule 1: SECONDARY ALCOHOL DEHYDROGENASE



H265	L266	S267	H272	H295	G296	R309	R312	L332	R333	E334	G335	R340	P345
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4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	65.67Å 105.10Å 109.18Å 90.00° 91.27° 90.00°	Depositor
Resolution (Å)	109.15 – 2.80 49.98 – 2.80	Depositor EDS
% Data completeness (in resolution range)	98.7 (109.15-2.80) 98.7 (49.98-2.80)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.48 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.188 , 0.258 0.187 , 0.251	Depositor DCC
R_{free} test set	1807 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	40.7	Xtriage
Anisotropy	0.640	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 30.2	EDS
Estimated twinning fraction	0.006 for -h,l,k 0.018 for -h,-l,-k 0.027 for h,-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 36163 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	10078	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.73% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, BU1, NAD

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	0.78	0/2419	0.79	1/3298 (0.0%)
1	B	0.72	0/2481	0.77	0/3382
1	C	0.64	0/2511	0.74	0/3427
1	D	0.65	0/2520	0.76	0/3441
All	All	0.70	0/9931	0.77	1/13548 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	37	LEU	CA-CB-CG	5.50	127.94	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	272	HIS	Peptide

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	2370	0	2314	34	0
1	B	2434	0	2394	47	0
1	C	2461	0	2452	28	0
1	D	2469	0	2456	23	0
2	A	44	0	26	4	0
2	B	44	0	26	1	0
2	C	44	0	26	2	0
2	D	44	0	26	0	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
3	C	2	0	0	0	0
3	D	2	0	0	0	0
4	B	6	0	10	0	0
4	C	6	0	10	0	0
4	D	6	0	9	0	0
5	A	28	0	0	0	0
5	B	39	0	0	1	0
5	C	38	0	0	4	0
5	D	37	0	0	1	0
All	All	10078	0	9749	131	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 7.

All (131) 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:38[A]:CYS:SG	1:A:62:HIS:NE2	2.47	0.86
1:D:312:ARG:HG2	1:D:312:ARG:HH11	1.44	0.83
1:B:116:PRO:HG2	1:B:119:LEU:HB2	1.65	0.77
1:D:312:ARG:HG2	1:D:312:ARG:NH1	2.00	0.72
1:A:5:GLN:O	1:A:13:PRO:HB3	1.90	0.71
1:C:132:ASP:HB3	5:C:2007:HOH:O	1.94	0.68
1:B:330:ARG:O	1:B:334:GLU:HG2	1.94	0.67
1:D:182:GLY:HA3	1:D:340:ARG:NH1	2.11	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:301:LEU:O	1:B:305:VAL:HG23	1.96	0.66
1:B:88:GLY:HA3	1:B:295:TRP:CZ2	2.30	0.66
1:A:38[A]:CYS:HG	1:A:62:HIS:CD2	2.13	0.65
1:D:157:THR:HB	1:D:158:PRO:HD3	1.79	0.65
1:A:116:PRO:HG2	1:A:119:LEU:HB2	1.78	0.65
1:D:143:LEU:HD13	1:D:309:ARG:HG2	1.79	0.65
1:C:13:PRO:HD3	1:C:45:MET:HE1	1.80	0.64
1:A:149:ALA:HB3	1:A:150:PRO:HD3	1.81	0.63
1:A:59:THR:HB	1:A:123:GLY:H	1.63	0.62
1:A:5:GLN:OE1	1:A:57:PRO:HB2	2.00	0.62
1:A:1:MET:HG3	1:A:2:LYS:O	1.99	0.61
1:D:312:ARG:CG	1:D:312:ARG:HH11	2.14	0.60
1:C:204:LEU:O	1:C:222:LYS:HE2	2.02	0.60
1:A:138:VAL:HG21	1:A:301:LEU:HD23	1.84	0.59
1:B:18:ILE:HB	1:B:19:PRO:HD2	1.84	0.59
1:B:116:PRO:HG2	1:B:119:LEU:CB	2.33	0.58
1:B:203:ASP:OD2	1:B:204:LEU:N	2.36	0.58
1:C:63:GLU:OE2	1:C:153:ASP:HB3	2.04	0.58
1:B:190:ILE:O	1:B:194:VAL:HG22	2.04	0.57
1:B:121:SER:HB3	1:B:122:PRO:CD	2.34	0.57
1:B:41:ASP:C	1:B:43:PHE:H	2.08	0.57
1:B:8:GLU:HB3	1:B:11:SER:HB2	1.86	0.57
1:C:13:PRO:HD3	1:C:45:MET:CE	2.35	0.56
1:B:88:GLY:HA3	1:B:295:TRP:CH2	2.40	0.56
1:C:157:THR:HB	1:C:158:PRO:HD3	1.88	0.55
1:C:237:GLY:O	1:C:239:GLY:N	2.40	0.55
1:A:38[A]:CYS:SG	1:A:63:GLU:HG3	2.47	0.54
1:A:88:GLY:HA3	1:A:295:TRP:CZ2	2.41	0.54
1:B:328:ALA:O	1:B:331:ARG:HB2	2.07	0.54
1:D:32:VAL:HG12	1:D:345:PRO:HG2	1.89	0.54
1:B:38:CYS:HB3	1:B:62:HIS:CE1	2.43	0.54
1:D:39:HIS:HA	1:D:332:LEU:HD21	1.89	0.53
1:A:300:GLU:O	1:A:304:VAL:HG23	2.09	0.53
1:B:156:LEU:HG	1:B:296:GLY:HA3	1.89	0.53
1:D:335:GLY:HA2	5:D:2036:HOH:O	2.08	0.52
1:B:36:GLY:HA3	1:B:63:GLU:OE1	2.08	0.52
1:D:87:TYR:CZ	1:D:89:PRO:HG2	2.44	0.52
1:A:1:MET:SD	1:A:126:ALA:HB1	2.50	0.52
1:A:181:GLY:HA3	2:A:503:NAD:O5B	2.09	0.52
1:A:325:GLY:O	1:A:328:ALA:HB3	2.09	0.52
1:B:5:GLN:OE1	1:B:57:PRO:HB2	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:299:SER:HB2	5:C:2029:HOH:O	2.10	0.51
1:B:38:CYS:SG	1:B:40:SER:HB2	2.51	0.51
1:B:263:ASP:HA	1:B:287:GLY:O	2.10	0.51
1:C:180:VAL:O	1:C:180:VAL:HG12	2.10	0.51
1:D:157:THR:HB	1:D:158:PRO:CD	2.40	0.50
1:A:203:ASP:HA	2:A:503:NAD:H2A	1.92	0.50
1:B:30:LEU:O	1:B:128:TYR:HA	2.11	0.50
1:B:32:VAL:O	1:B:345:PRO:HG2	2.12	0.50
1:D:154:ALA:HA	1:D:183:LEU:HG	1.93	0.50
1:B:268:VAL:HB	1:B:292:THR:HG22	1.93	0.50
1:A:40:SER:O	1:A:44:VAL:HG23	2.12	0.49
1:C:7:THR:HG22	1:C:57:PRO:HB3	1.94	0.49
1:B:325:GLY:O	1:B:328:ALA:HB3	2.12	0.49
1:C:204:LEU:HD23	1:C:223:SER:HB3	1.94	0.49
1:B:121:SER:HB3	1:B:122:PRO:HD2	1.94	0.49
1:B:59:THR:HB	1:B:123:GLY:H	1.77	0.49
1:C:279:VAL:HG13	1:C:284:ILE:HG21	1.95	0.49
1:D:65:VAL:HG13	1:D:149:ALA:HA	1.95	0.48
1:C:319:THR:HG22	5:C:2009:HOH:O	2.13	0.48
1:A:203:ASP:C	1:A:223:SER:HB2	2.34	0.48
1:A:157:THR:HB	1:A:158:PRO:HD3	1.94	0.48
1:C:2:LYS:HB3	1:C:322:LEU:HD13	1.94	0.48
1:B:84:VAL:HG21	1:B:137:LEU:HD22	1.96	0.47
1:A:44:VAL:O	1:A:47:MET:HB2	2.15	0.47
1:A:203:ASP:HA	2:A:503:NAD:C2A	2.45	0.47
1:B:65:VAL:HG23	1:B:149:ALA:HA	1.96	0.47
1:C:131:VAL:HG21	1:C:137:LEU:HD21	1.96	0.46
1:C:203:ASP:HA	2:C:503:NAD:C2A	2.45	0.46
1:A:60:LEU:O	1:A:124:SER:N	2.48	0.46
1:B:134:ALA:O	1:B:136:HIS:N	2.48	0.46
1:A:165:VAL:HG11	1:A:265:HIS:CG	2.51	0.46
1:B:41:ASP:C	1:B:43:PHE:N	2.69	0.46
1:A:38[A]:CYS:SG	1:A:62:HIS:CD2	3.06	0.46
1:B:319:THR:HG23	1:B:344:VAL:HG23	1.97	0.46
1:B:111:ASP:C	1:B:113:GLY:H	2.20	0.45
1:C:176:VAL:HB	1:C:243:VAL:HG22	1.98	0.45
1:D:90:TRP:CZ2	1:D:117:PRO:HD3	2.51	0.45
1:B:331:ARG:O	1:B:334:GLU:HG3	2.16	0.45
1:D:165:VAL:HG11	1:D:265:HIS:CG	2.52	0.45
1:B:72:GLY:O	1:B:74:GLY:N	2.49	0.45
1:B:276:HIS:CD2	1:C:278:LYS:HG2	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:281:PHE:O	1:B:282:PHE:HB2	2.17	0.45
1:C:203:ASP:OD2	2:C:503:NAD:O2B	2.31	0.45
1:C:31:LYS:HD2	1:C:128:TYR:CZ	2.51	0.45
1:A:14:VAL:O	1:A:14:VAL:HG23	2.17	0.45
1:D:7:THR:HG22	1:D:57:PRO:HB3	1.97	0.44
1:A:20:THR:HA	1:A:21:PRO:HD3	1.73	0.44
1:A:6:TYR:CE1	1:A:45:MET:HG2	2.53	0.44
1:B:12:GLU:OE1	1:B:330:ARG:NH2	2.51	0.44
1:C:33:THR:OG1	1:C:65:VAL:HG12	2.17	0.44
1:B:248:GLY:HA3	1:B:270:GLY:O	2.18	0.44
1:D:149:ALA:HB3	1:D:150:PRO:HD3	2.00	0.44
1:C:245:ASP:CG	1:C:248:GLY:H	2.20	0.43
1:B:143:LEU:N	5:B:2014:HOH:O	2.51	0.43
1:A:144:ASP:HA	1:A:145:PRO:HD3	1.81	0.43
1:B:40:SER:O	1:B:43:PHE:HB3	2.19	0.43
1:B:130:ILE:HD13	1:B:130:ILE:HA	1.76	0.43
1:C:180:VAL:HG21	1:C:212:ALA:HB2	2.01	0.43
1:A:246:PHE:O	2:A:503:NAD:H4D	2.19	0.42
1:C:319:THR:CG2	5:C:2009:HOH:O	2.67	0.42
1:C:142:ASP:H	1:C:309:ARG:HH12	1.65	0.42
1:B:23:PRO:HB2	1:B:72:GLY:HA2	2.02	0.42
1:B:181:GLY:HA3	2:B:503:NAD:O5B	2.20	0.42
1:A:166:LEU:HB3	1:A:167:PRO:HD3	2.00	0.42
1:A:12:GLU:HA	1:A:13:PRO:HD3	1.74	0.42
1:D:30:LEU:O	1:D:128:TYR:HA	2.20	0.42
1:D:140:ILE:O	1:D:143:LEU:HB2	2.19	0.42
1:B:322:LEU:HD11	1:B:345:PRO:HB3	2.02	0.42
1:B:154:ALA:HA	1:B:183:LEU:HG	2.01	0.42
1:A:150:PRO:O	1:A:151:LEU:C	2.58	0.41
1:C:265:HIS:CE1	1:C:291:VAL:HG12	2.54	0.41
1:D:260:VAL:HG22	1:D:261:ALA:N	2.34	0.41
1:B:32:VAL:HG11	1:B:124:SER:O	2.20	0.41
1:D:295:TRP:HB3	1:D:296:GLY:H	1.60	0.41
1:B:138:VAL:HG21	1:B:301:LEU:HD23	2.03	0.41
1:A:138:VAL:CG2	1:A:301:LEU:HD23	2.48	0.41
1:C:28:ILE:N	1:C:28:ILE:HD12	2.36	0.41
1:D:24:GLY:O	1:D:72:GLY:HA3	2.21	0.41
1:B:24:GLY:O	1:B:26:GLY:N	2.53	0.41
1:C:30:LEU:O	1:C:128:TYR:HA	2.21	0.40
1:D:248:GLY:O	1:D:272:HIS:HB2	2.22	0.40
1:A:63:GLU:HA	1:A:152:THR:OG1	2.21	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	324/345 (94%)	291 (90%)	31 (10%)	2 (1%)	30	65
1	B	334/345 (97%)	310 (93%)	22 (7%)	2 (1%)	30	65
1	C	340/345 (99%)	325 (96%)	15 (4%)	0	100	100
1	D	343/345 (99%)	323 (94%)	19 (6%)	1 (0%)	46	79
All	All	1341/1380 (97%)	1249 (93%)	87 (6%)	5 (0%)	39	74

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	23	PRO
1	A	50	ALA
1	B	273	ALA
1	D	188	ILE
1	B	42	ILE

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	232/248 (94%)	223 (96%)	9 (4%)	39	74
1	B	239/248 (96%)	228 (95%)	11 (5%)	33	67
1	C	246/248 (99%)	241 (98%)	5 (2%)	63	90

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	246/248 (99%)	236 (96%)	10 (4%)	37	72
All	All	963/992 (97%)	928 (96%)	35 (4%)	42	76

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	14	VAL
1	A	56	LEU
1	A	108	ARG
1	A	144	ASP
1	A	205	ASP
1	A	223	SER
1	A	233	GLU
1	A	322	LEU
1	B	2	LYS
1	B	18	ILE
1	B	40	SER
1	B	46	ASP
1	B	107	THR
1	B	108	ARG
1	B	130	ILE
1	B	204	LEU
1	B	205	ASP
1	B	271	ILE
1	B	309	ARG
1	C	76	THR
1	C	267	SER
1	C	278	LYS
1	C	295	TRP
1	C	334	GLU
1	D	11	SER
1	D	12	GLU
1	D	92	CYS
1	D	112	LEU
1	D	195	SER
1	D	267	SER
1	D	295	TRP
1	D	312	ARG
1	D	334	GLU
1	D	340	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	258	GLN
1	C	276	HIS
1	D	39	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](#)

Of 15 ligands modelled in this entry, 8 are monoatomic - leaving 7 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
2	NAD	A	503	-	38,48,48	1.66	2 (5%)	47,73,73	2.60	7 (14%)
4	BU1	B	1347	-	5,5,5	0.62	0	4,4,4	0.21	0
2	NAD	B	503	-	38,48,48	1.62	2 (5%)	47,73,73	2.18	5 (10%)
4	BU1	C	1348	-	5,5,5	0.51	0	4,4,4	0.36	0
2	NAD	C	503	-	38,48,48	1.59	3 (7%)	47,73,73	2.42	9 (19%)
4	BU1	D	1347	3	5,5,5	0.69	0	4,4,4	0.43	0
2	NAD	D	503	-	38,48,48	1.69	3 (7%)	47,73,73	2.16	8 (17%)

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
2	NAD	A	503	-	-	0/22/62/62	0/5/5/5
4	BU1	B	1347	-	-	0/3/3/3	0/0/0/0
2	NAD	B	503	-	-	0/22/62/62	0/5/5/5
4	BU1	C	1348	-	-	0/3/3/3	0/0/0/0
2	NAD	C	503	-	-	0/22/62/62	0/5/5/5
4	BU1	D	1347	3	-	0/3/3/3	0/0/0/0
2	NAD	D	503	-	-	0/22/62/62	0/5/5/5

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	503	NAD	C2A-N1A	2.03	1.37	1.33
2	D	503	NAD	C2A-N1A	2.55	1.38	1.33
2	B	503	NAD	C2A-N3A	3.27	1.38	1.32
2	A	503	NAD	C2A-N3A	3.36	1.38	1.32
2	C	503	NAD	C2A-N3A	3.42	1.38	1.32
2	D	503	NAD	C2A-N3A	3.52	1.38	1.32
2	C	503	NAD	O7N-C7N	7.11	1.39	1.24
2	A	503	NAD	O7N-C7N	8.01	1.41	1.24
2	B	503	NAD	O7N-C7N	8.23	1.41	1.24
2	D	503	NAD	O7N-C7N	8.33	1.41	1.24

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	503	NAD	N3A-C2A-N1A	-12.85	119.05	128.89
2	B	503	NAD	N3A-C2A-N1A	-12.29	119.48	128.89
2	A	503	NAD	N3A-C2A-N1A	-12.22	119.53	128.89
2	D	503	NAD	N3A-C2A-N1A	-10.90	120.55	128.89
2	B	503	NAD	PN-O3-PA	-5.02	118.63	132.73
2	A	503	NAD	C1B-N9A-C4A	-4.53	120.11	126.94
2	C	503	NAD	C1B-N9A-C4A	-3.90	121.06	126.94
2	D	503	NAD	C1B-N9A-C4A	-3.88	121.09	126.94
2	D	503	NAD	C2B-C1B-N9A	-3.63	108.75	114.29
2	D	503	NAD	PN-O3-PA	-3.54	122.79	132.73
2	A	503	NAD	PN-O3-PA	-3.44	123.08	132.73
2	C	503	NAD	O3-PN-O5D	-3.13	94.63	102.94
2	C	503	NAD	C2B-C1B-N9A	-3.07	109.60	114.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	503	NAD	C4A-C5A-N7A	-3.05	106.67	109.48
2	C	503	NAD	PN-O3-PA	-2.83	124.77	132.73
2	B	503	NAD	C1B-N9A-C4A	-2.63	122.97	126.94
2	D	503	NAD	C4B-O4B-C1B	-2.57	106.89	109.72
2	C	503	NAD	O7N-C7N-N7N	-2.53	119.04	122.59
2	C	503	NAD	O3D-C3D-C4D	-2.23	104.35	111.05
2	A	503	NAD	C4A-C5A-N7A	-2.21	107.44	109.48
2	D	503	NAD	C5B-C4B-C3B	-2.21	106.45	115.21
2	A	503	NAD	C5N-C4N-C3N	-2.11	117.68	120.33
2	B	503	NAD	O4B-C1B-N9A	2.12	112.53	108.10
2	A	503	NAD	C3N-C2N-N1N	2.22	122.92	120.36
2	D	503	NAD	O4D-C1D-N1N	2.32	110.68	108.13
2	C	503	NAD	O4B-C1B-N9A	2.47	113.27	108.10
2	D	503	NAD	O4B-C1B-N9A	2.75	113.86	108.10
2	B	503	NAD	O4D-C1D-N1N	2.95	111.37	108.13
2	A	503	NAD	O4D-C1D-N1N	9.89	119.00	108.13

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	503	NAD	4	0
2	B	503	NAD	1	0
2	C	503	NAD	2	0

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](#)

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	333/345 (96%)	-0.23	4 (1%) 81 73	23, 38, 57, 61	0
1	B	342/345 (99%)	-0.31	2 (0%) 90 86	23, 39, 58, 71	0
1	C	344/345 (99%)	-0.44	0 100 100	23, 38, 54, 60	0
1	D	345/345 (100%)	-0.56	0 100 100	23, 38, 55, 60	0
All	All	1364/1380 (98%)	-0.39	6 (0%) 93 90	23, 38, 57, 71	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	55	GLY	2.4
1	A	60	LEU	2.3
1	B	56	LEU	2.2
1	A	13	PRO	2.2
1	A	130	ILE	2.1
1	A	329	TYR	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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron

density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
4	BU1	D	1347	6/6	0.91	0.23	5.08	25,33,33,35	0
4	BU1	B	1347	6/6	0.91	0.23	1.62	49,52,53,53	0
4	BU1	C	1348	6/6	0.92	0.19	0.13	38,41,44,45	0
2	NAD	D	503	44/44	0.98	0.13	-0.33	27,32,36,36	0
2	NAD	B	503	44/44	0.97	0.13	-0.38	34,38,42,43	0
2	NAD	A	503	44/44	0.97	0.12	-0.81	28,37,39,40	0
3	ZN	D	1348	1/1	1.00	0.11	-0.85	34,34,34,34	0
2	NAD	C	503	44/44	0.98	0.12	-1.03	22,29,31,32	0
3	ZN	A	1347	1/1	0.99	0.09	-1.53	33,33,33,33	0
3	ZN	C	1347	1/1	1.00	0.08	-1.71	31,31,31,31	0
3	ZN	B	1348	1/1	1.00	0.06	-2.05	32,32,32,32	0
3	ZN	A	1346	1/1	0.99	0.05	-4.10	51,51,51,51	0
3	ZN	B	1346	1/1	0.99	0.07	-	42,42,42,42	0
3	ZN	D	1346	1/1	0.99	0.12	-	35,35,35,35	0
3	ZN	C	1346	1/1	1.00	0.07	-	37,37,37,37	0

6.5 Other polymers [i](#)

There are no such residues in this entry.