Full wwPDB X-ray Structure Validation Report

Jan 31, 2016 – 07:52 PM GMT

PDB ID	:	1HLD
Title	:	STRUCTURES OF HORSE LIVER ALCOHOL DEHYDROGENASE COMPLEXED WITH NAD+ AND SUBSTITUTED BENZYL ALCOHOLS
Authors	:	Ramaswamy, S.; Eklund, H.; Plapp, B.V.
Deposited on	:	1993-11-23
Resolution	:	2.10 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:

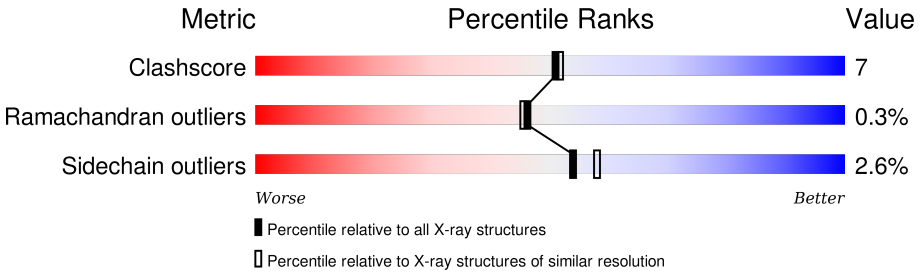
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	102246	4460 (2.10-2.10)
Ramachandran outliers	100387	4413 (2.10-2.10)
Sidechain outliers	100360	4414 (2.10-2.10)

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 <=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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	374	82% 17%
1	B	374	84% 16%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6100 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.

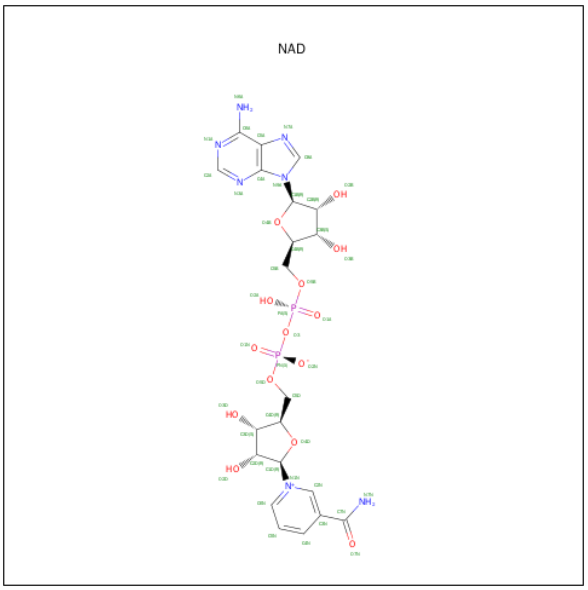
- Molecule 1 is a protein called ALCOHOL DEHYDROGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	374	Total	C	N	O	S	0	0	0
			2785	1769	472	521	23			
1	B	374	Total	C	N	O	S	0	0	0
			2785	1769	472	521	23			

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

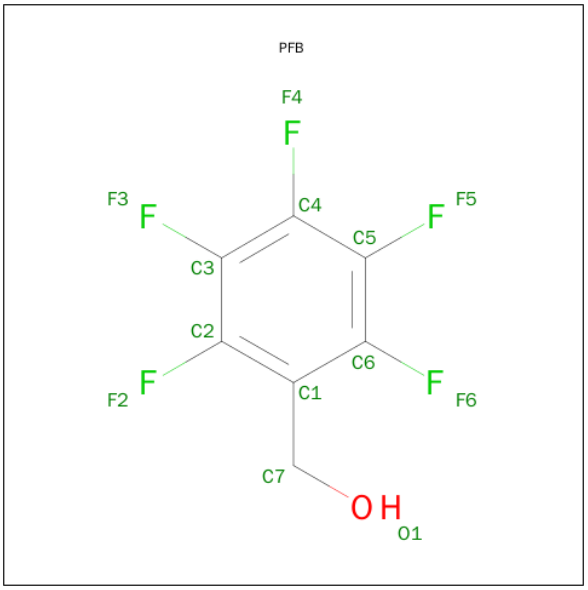
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	2	Total	Zn	0	0
			2	2		
2	A	2	Total	Zn	0	0
			2	2		

- Molecule 3 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: $C_{21}H_{27}N_7O_{14}P_2$).



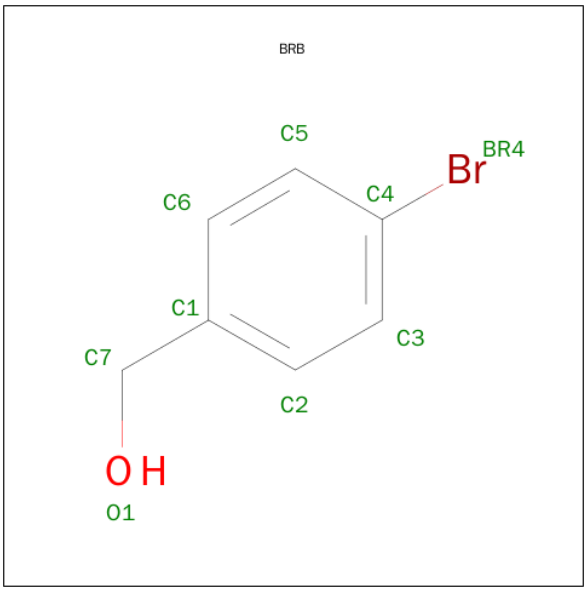
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
3	B	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

- Molecule 4 is 2,3,4,5,6-PENTAFLUOROBENZYL ALCOHOL (three-letter code: PFB) (formula: $C_7H_3F_5O$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	F	O	0	0
			13	7	5	1		
4	B	1	Total	C	F	O	0	0
			13	7	5	1		

- Molecule 5 is PARA-BROMOBENZYL ALCOHOL (three-letter code: BRB) (formula: C_7H_7BrO).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	Br	C	O	0	0
			9	1	7	1		
5	B	1	Total	Br	C	O	0	0
			9	1	7	1		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	218	Total	O	0	0
			218	218		
6	B	176	Total	O	0	0
			176	176		

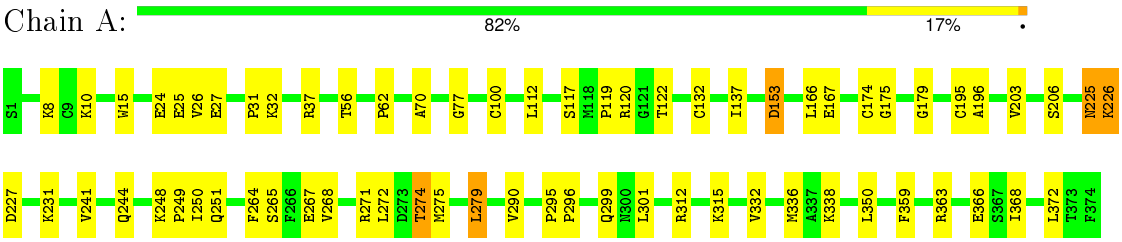
3 Residue-property plots (i)

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.

Note EDS was not executed.

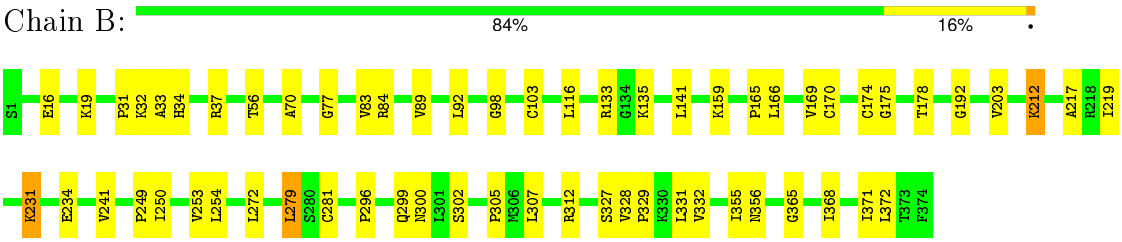
- Molecule 1: ALCOHOL DEHYDROGENASE

Chain A:



- Molecule 1: ALCOHOL DEHYDROGENASE

Chain B:



4 Data and refinement statistics (i)

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

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	51.10Å 180.70Å 44.30Å 90.00° 108.00° 90.00°	Depositor
Resolution (Å)	8.00 – 2.10	Depositor
% Data completeness (in resolution range)	(Not available) (8.00-2.10)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.183 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6100	wwPDB-VP
Average B, all atoms (Å^2)	26.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, PFB, BRB, 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	0.51	0/2837	0.71	0/3834
1	B	0.48	0/2837	0.71	0/3834
All	All	0.49	0/5674	0.71	0/7668

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	2785	0	2848	42	0
1	B	2785	0	2848	40	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	44	0	26	0	0
3	B	44	0	26	1	0
4	A	13	0	1	0	0
4	B	13	0	1	0	0
5	A	9	0	0	0	0
5	B	9	0	1	0	0
6	A	218	0	0	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	B	176	0	0	2	0
All	All	6100	0	5751	79	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 (79) 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:B:272:LEU:HD11	1:B:299:GLN:HB3	1.59	0.84
1:B:241:VAL:HG11	1:B:250:ILE:HD11	1.59	0.84
1:B:170:CYS:SG	1:B:371:ILE:HD12	2.21	0.80
1:B:212:LYS:HA	1:B:212:LYS:HE3	1.65	0.76
1:B:296:PRO:HB2	1:B:299:GLN:HG3	1.70	0.73
1:A:132:CYS:HB3	1:A:137:ILE:HD11	1.77	0.65
1:A:153:ASP:HB3	6:A:2193:HOH:O	1.96	0.64
1:B:241:VAL:CG1	1:B:250:ILE:HD11	2.27	0.62
1:B:92:LEU:HA	6:B:2022:HOH:O	1.99	0.62
1:B:327:SER:O	1:B:331:LEU:HG	2.02	0.59
1:A:10:LYS:HA	1:A:24:GLU:O	2.03	0.58
1:A:8:LYS:HE2	1:A:27:GLU:HG2	1.84	0.58
1:B:133:ARG:HH11	1:B:135:LYS:HZ2	1.53	0.57
1:B:56:THR:HG23	1:B:296:PRO:HA	1.87	0.56
1:A:295:PRO:HB2	6:A:2142:HOH:O	2.05	0.56
1:A:301:LEU:HD23	1:B:305:PRO:HG3	1.87	0.56
1:A:167:GLU:HB2	6:A:2290:HOH:O	2.06	0.56
1:B:70:ALA:HB1	1:B:166:LEU:HD22	1.88	0.55
1:A:267:GLU:HG3	1:A:275:MET:HA	1.88	0.55
1:B:192:GLY:HA2	1:B:217:ALA:HB2	1.88	0.54
1:A:272:LEU:HD11	1:A:299:GLN:HB3	1.89	0.54
1:A:226:LYS:HG3	6:A:2386:HOH:O	2.08	0.54
1:A:8:LYS:HB3	1:A:8:LYS:HZ2	1.74	0.53
1:A:8:LYS:NZ	1:A:25:GLU:HG2	2.24	0.53
1:A:8:LYS:CE	1:A:27:GLU:HG2	2.38	0.52
1:A:175:GLY:HA2	1:A:203:VAL:HG22	1.90	0.52
1:B:328:VAL:O	1:B:332:VAL:HG23	2.10	0.52
1:A:301:LEU:CD2	1:B:305:PRO:HG3	2.40	0.51
1:B:365:GLY:HA2	6:B:2029:HOH:O	2.11	0.51
1:B:133:ARG:HH11	1:B:135:LYS:NZ	2.09	0.50
1:A:359:PHE:HB3	1:A:363:ARG:CZ	2.42	0.50
1:A:167:GLU:H	1:A:167:GLU:CD	2.15	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:241:VAL:HG11	1:A:250:ILE:HD11	1.94	0.49
1:A:100:CYS:HB2	1:A:112:LEU:HD12	1.95	0.48
1:B:231:LYS:O	1:B:234:GLU:HB3	2.11	0.48
1:B:83:VAL:HG12	1:B:159:LYS:HB2	1.96	0.48
1:A:231:LYS:HG2	1:A:368:ILE:CD1	2.44	0.47
1:A:70:ALA:HB1	1:A:166:LEU:HD22	1.95	0.47
1:A:290:VAL:HA	1:A:315:LYS:O	2.15	0.47
1:B:165:PRO:O	1:B:169:VAL:HG22	2.15	0.47
1:A:301:LEU:O	1:B:302:SER:HA	2.15	0.47
1:A:332:VAL:O	1:A:336:MET:HG2	2.15	0.46
1:B:175:GLY:HA2	1:B:203:VAL:HG22	1.96	0.46
1:B:178:THR:HG21	3:B:901:NAD:C4N	2.46	0.46
1:B:89:VAL:HG12	1:B:159:LYS:HA	1.96	0.46
1:A:32:LYS:O	1:A:77:GLY:HA3	2.16	0.46
1:B:328:VAL:N	1:B:329:PRO:HD2	2.31	0.46
1:B:16:GLU:HB2	1:B:19:LYS:HG3	1.99	0.45
1:B:254:LEU:HD12	1:B:281:CYS:SG	2.57	0.44
1:B:249:PRO:O	1:B:253:VAL:HG23	2.17	0.44
1:A:271:ARG:HB2	1:A:274:THR:OG1	2.18	0.44
1:A:179:GLY:HA3	1:A:206:SER:HB2	2.00	0.44
1:B:116:LEU:HG	1:B:141:LEU:HD22	1.98	0.44
1:B:92:LEU:CD2	1:B:328:VAL:HG21	2.47	0.44
1:A:120:ARG:NH1	1:A:122:THR:OG1	2.51	0.44
1:A:279:LEU:HD22	1:A:312:ARG:HD3	1.99	0.43
1:A:56:THR:HG23	1:A:296:PRO:HA	2.01	0.43
1:A:31:PRO:HD3	1:A:37:ARG:HB2	2.01	0.43
1:A:350:LEU:O	1:A:372:LEU:HA	2.19	0.42
1:B:250:ILE:HA	1:B:250:ILE:HD12	1.87	0.42
1:A:26:VAL:HG12	1:A:132:CYS:HB2	2.01	0.42
1:B:32:LYS:O	1:B:77:GLY:HA3	2.19	0.42
1:B:231:LYS:HB2	1:B:231:LYS:NZ	2.33	0.42
1:B:212:LYS:HD2	1:B:219:ILE:HD12	2.02	0.42
1:B:279:LEU:HD22	1:B:312:ARG:HD3	2.02	0.42
1:A:338:LYS:HB3	1:A:338:LYS:HE2	1.84	0.42
1:B:31:PRO:HD3	1:B:37:ARG:HB2	2.02	0.42
1:A:195:CYS:HA	1:A:264:PHE:O	2.20	0.41
1:B:98:GLY:HA2	1:B:103:CYS:HB3	2.02	0.41
1:A:231:LYS:HG2	1:A:368:ILE:HD11	2.03	0.41
1:A:15:TRP:HA	1:A:62:PRO:HB3	2.02	0.41
1:A:225:ASN:ND2	1:A:227:ASP:H	2.19	0.41
1:A:248:LYS:HG3	1:A:249:PRO:HD2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:231:LYS:HG2	1:B:368:ILE:CD1	2.51	0.41
1:A:117:SER:C	1:A:119:PRO:HD3	2.41	0.40
1:B:33:ALA:O	1:B:34:HIS:HB2	2.22	0.40
1:A:117:SER:O	1:A:119:PRO:HD3	2.21	0.40
1:B:355:ILE:HG23	1:B:356:ASN:N	2.36	0.40
1:A:196:ALA:O	1:A:265:SER:HA	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	372/374 (100%)	354 (95%)	17 (5%)	1 (0%)	46	45
1	B	372/374 (100%)	352 (95%)	19 (5%)	1 (0%)	46	45
All	All	744/748 (100%)	706 (95%)	36 (5%)	2 (0%)	46	45

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	174	CYS
1	B	174	CYS

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	308/308 (100%)	299 (97%)	9 (3%)	50 53
1	B	308/308 (100%)	301 (98%)	7 (2%)	58 62
All	All	616/616 (100%)	600 (97%)	16 (3%)	54 58

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	153	ASP
1	A	225	ASN
1	A	226	LYS
1	A	244	GLN
1	A	251	GLN
1	A	268	VAL
1	A	274	THR
1	A	279	LEU
1	A	366	GLU
1	B	84	ARG
1	B	212	LYS
1	B	231	LYS
1	B	279	LEU
1	B	300	ASN
1	B	307	LEU
1	B	372	LEU

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

Mol	Chain	Res	Type
1	A	225	ASN
1	A	300	ASN
1	B	300	ASN

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 10 ligands modelled in this entry, 4 are monoatomic - leaving 6 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	$\#\|Z\|>2$
4	PFB	A	1001	2,5	13,13,13	2.15	5 (38%)	19,19,19	1.60	5 (26%)
5	BRB	A	1003	2,4	9,9,9	2.60	2 (22%)	11,11,11	1.16	1 (9%)
3	NAD	A	801	-	38,48,48	0.99	2 (5%)	47,73,73	2.35	10 (21%)
4	PFB	B	1002	2,5	13,13,13	2.46	6 (46%)	19,19,19	2.18	4 (21%)
5	BRB	B	1004	2,4	9,9,9	3.17	2 (22%)	11,11,11	1.23	1 (9%)
3	NAD	B	901	-	38,48,48	1.11	3 (7%)	47,73,73	1.93	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
4	PFB	A	1001	2,5	-	0/2/2/2	0/1/1/1
5	BRB	A	1003	2,4	-	0/2/2/2	0/1/1/1
3	NAD	A	801	-	-	0/22/62/62	0/5/5/5
4	PFB	B	1002	2,5	-	0/2/2/2	0/1/1/1
5	BRB	B	1004	2,4	-	0/2/2/2	0/1/1/1
3	NAD	B	901	-	-	0/22/62/62	0/5/5/5

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	1004	BRB	BR4-C4	-8.57	1.71	1.90
5	A	1003	BRB	BR4-C4	-6.93	1.75	1.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	1002	PFB	C7-C1	-6.37	1.42	1.51
4	A	1001	PFB	C7-C1	-5.12	1.44	1.51
4	A	1001	PFB	F5-C5	-3.09	1.30	1.35
4	B	1002	PFB	F2-C2	-2.75	1.30	1.35
4	A	1001	PFB	F3-C3	-2.74	1.30	1.35
5	B	1004	BRB	C7-C1	-2.66	1.40	1.51
4	B	1002	PFB	F5-C5	-2.58	1.30	1.35
4	A	1001	PFB	F6-C6	-2.58	1.30	1.35
4	B	1002	PFB	F3-C3	-2.56	1.30	1.35
4	A	1001	PFB	F2-C2	-2.29	1.31	1.35
4	B	1002	PFB	F4-C4	-2.24	1.31	1.35
3	A	801	NAD	PA-O2A	-2.04	1.46	1.54
3	A	801	NAD	O4D-C1D	2.10	1.43	1.41
3	B	901	NAD	C6N-N1N	2.25	1.41	1.35
4	B	1002	PFB	C5-C4	2.40	1.42	1.37
5	A	1003	BRB	C3-C2	2.53	1.43	1.38
3	B	901	NAD	O4B-C1B	2.55	1.44	1.41
3	B	901	NAD	O4D-C1D	3.59	1.45	1.41

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	801	NAD	N3A-C2A-N1A	-10.21	121.08	128.89
3	B	901	NAD	N3A-C2A-N1A	-9.63	121.52	128.89
3	A	801	NAD	C4B-O4B-C1B	-7.23	101.77	109.72
4	B	1002	PFB	C7-C1-C2	-5.73	116.15	121.81
4	A	1001	PFB	C7-C1-C6	-4.16	117.70	121.81
4	B	1002	PFB	C1-C6-C5	-4.16	117.23	122.33
3	A	801	NAD	O5B-PA-O1A	-3.07	97.69	109.62
3	A	801	NAD	C1B-N9A-C4A	-2.86	122.63	126.94
3	B	901	NAD	O3-PN-O5D	-2.31	96.81	102.94
3	A	801	NAD	C4D-O4D-C1D	-2.26	107.23	109.72
5	A	1003	BRB	C5-C4-C3	-2.22	117.32	121.41
3	B	901	NAD	O5D-PN-O1N	-2.18	101.14	109.62
4	A	1001	PFB	C1-C6-C5	-2.12	119.73	122.33
3	A	801	NAD	O3-PN-O5D	-2.11	97.35	102.94
4	A	1001	PFB	C1-C2-C3	-2.10	119.76	122.33
3	B	901	NAD	O3D-C3D-C2D	-2.02	105.24	111.83
3	B	901	NAD	C3N-C7N-N7N	-2.02	115.61	117.82
3	B	901	NAD	O4B-C1B-N9A	2.14	112.57	108.10
4	A	1001	PFB	C6-C5-C4	2.18	122.11	119.53
3	B	901	NAD	O7N-C7N-N7N	2.35	125.90	122.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	B	1002	PFB	F6-C6-C1	2.55	122.53	119.23
5	B	1004	BRB	BR4-C4-C5	2.59	123.37	119.28
3	A	801	NAD	O2A-PA-O5B	2.65	121.84	108.46
3	A	801	NAD	C2A-N1A-C6A	2.91	123.96	118.77
3	A	801	NAD	O4B-C1B-N9A	2.95	114.27	108.10
4	A	1001	PFB	C6-C1-C2	3.21	119.79	115.84
3	B	901	NAD	C4A-C5A-N7A	3.29	112.51	109.48
3	A	801	NAD	C4A-C5A-N7A	4.05	113.20	109.48
4	B	1002	PFB	C6-C1-C2	4.60	121.51	115.84

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	901	NAD	1	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

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.