



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:06 PM GMT

PDB ID : 1E3S
Title : RAT BRAIN 3-HYDROXYACYL-COA DEHYDROGENASE BINARY
COMPLEX WITH NADH
Authors : Powell, A.J.; Read, J.A.; Banfield, M.J.; Brady, R.L.
Deposited on : 2000-06-22
Resolution : 2.00 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

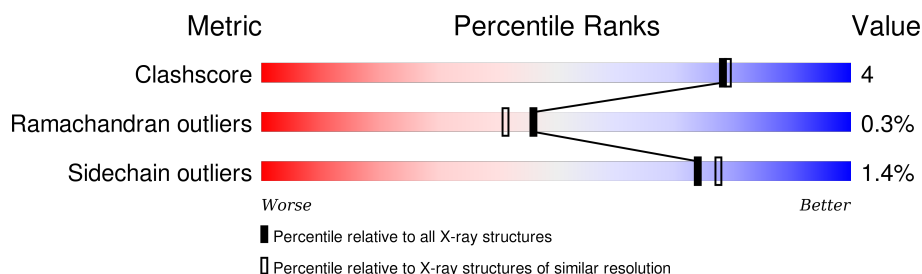
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)

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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	261	 82% 9% • 8%
1	B	261	 82% 8% • 8%
1	C	261	 83% 8% • 8%
1	D	261	 82% 9% • 8%

2 Entry composition [i](#)

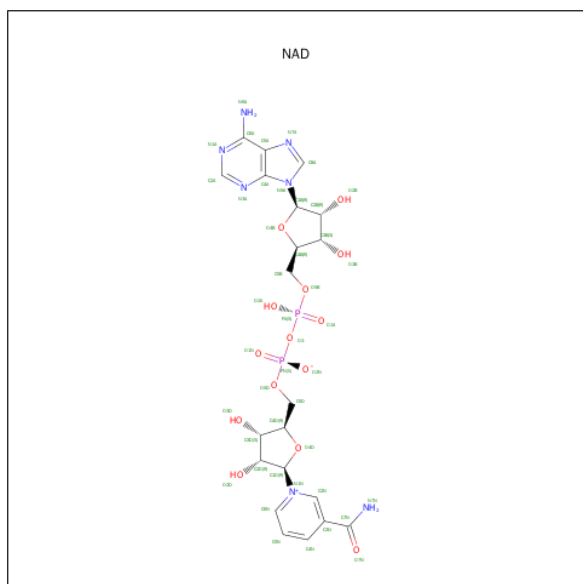
There are 3 unique types of molecules in this entry. The entry contains 7605 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 SHORT CHAIN 3-HYDROXYACYL-COA DEHYDROGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	241	Total	C	N	O	S	3	0	0
			1730	1093	307	324	6			
1	B	241	Total	C	N	O	S	0	0	0
			1733	1094	307	326	6			
1	C	241	Total	C	N	O	S	0	0	0
			1735	1096	307	326	6			
1	D	241	Total	C	N	O	S	3	0	0
			1739	1098	308	327	6			

- Molecule 2 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: $C_{21}H_{27}N_7O_{14}P_2$).



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

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

- Molecule 3 is water.

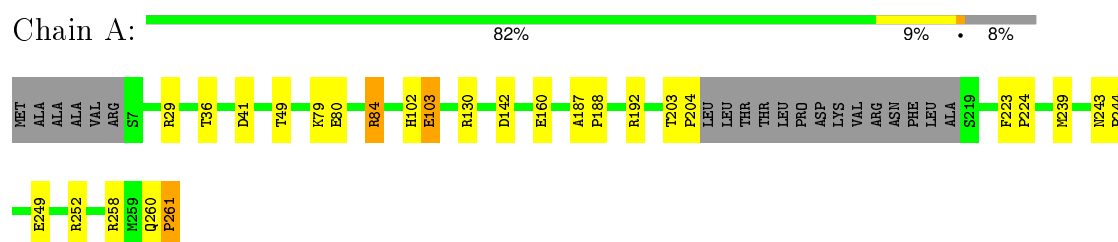
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	109	Total	O	0	0
			109	109		
3	B	130	Total	O	0	0
			130	130		
3	C	124	Total	O	0	0
			124	124		
3	D	129	Total	O	0	0
			129	129		

3 Residue-property plots [i](#)

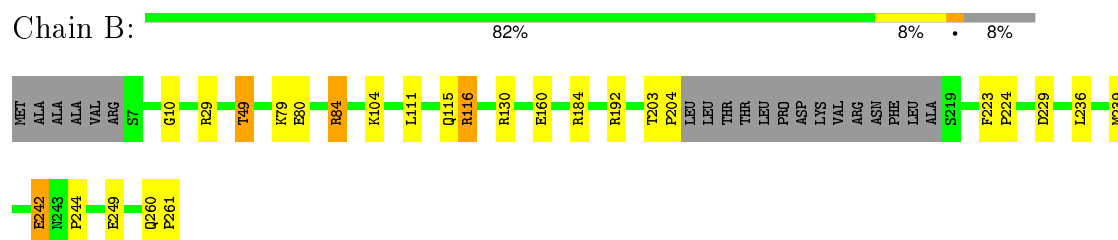
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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.

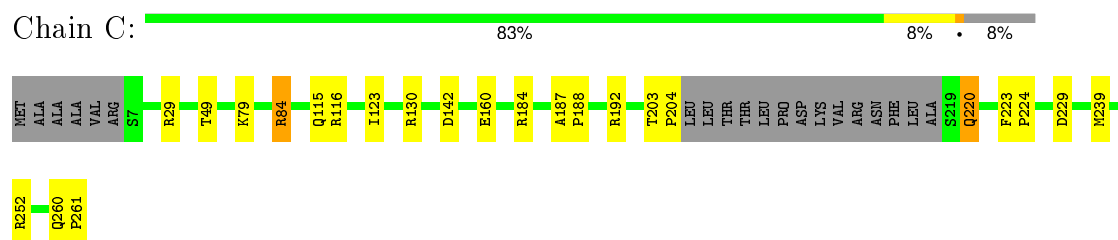
• Molecule 1: SHORT CHAIN 3-HYDROXYACYL-COA DEHYDROGENASE



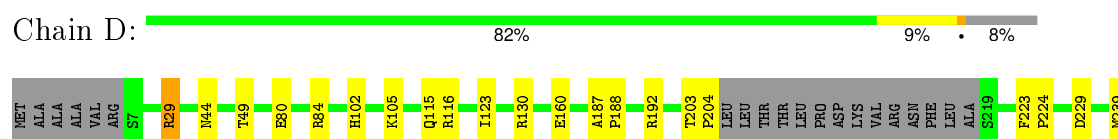
• Molecule 1: SHORT CHAIN 3-HYDROXYACYL-COA DEHYDROGENASE



• Molecule 1: SHORT CHAIN 3-HYDROXYACYL-COA DEHYDROGENASE



• Molecule 1: SHORT CHAIN 3-HYDROXYACYL-COA DEHYDROGENASE



E249	D254	R258	Q259	Q260	P261
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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	59.28 Å 68.96 Å 69.55 Å 64.84° 72.51° 74.98°	Depositor
Resolution (Å)	40.00 – 2.00	Depositor
% Data completeness (in resolution range)	94.9 (40.00-2.00)	Depositor
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
Refinement program	REFMAC	Depositor
R, R_{free}	0.189 , 0.226	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7605	wwPDB-VP
Average B, all atoms (Å ²)	18.0	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: 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	0.71	2/1754 (0.1%)	1.43	14/2380 (0.6%)
1	B	0.68	1/1757 (0.1%)	1.50	13/2383 (0.5%)
1	C	0.67	1/1759 (0.1%)	1.41	10/2387 (0.4%)
1	D	0.66	1/1763 (0.1%)	1.21	12/2393 (0.5%)
All	All	0.68	5/7033 (0.1%)	1.39	49/9543 (0.5%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	79	LYS	CG-CD	-9.19	1.21	1.52
1	B	261	PRO	N-CD	6.74	1.57	1.47
1	D	261	PRO	N-CD	6.56	1.57	1.47
1	A	261	PRO	N-CD	6.17	1.56	1.47
1	C	261	PRO	N-CD	5.24	1.55	1.47

All (49) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	84	ARG	CD-NE-CZ	35.41	173.17	123.60
1	B	84	ARG	CD-NE-CZ	32.14	168.60	123.60
1	A	84	ARG	CD-NE-CZ	31.75	168.04	123.60
1	B	160	GLU	OE1-CD-OE2	-23.83	94.71	123.30
1	A	79	LYS	CB-CG-CD	12.95	145.26	111.60
1	B	84	ARG	NE-CZ-NH2	11.64	126.12	120.30
1	A	130	ARG	NE-CZ-NH2	-11.42	114.59	120.30
1	A	160	GLU	OE1-CD-OE2	-9.17	112.30	123.30
1	B	192	ARG	NE-CZ-NH1	8.95	124.78	120.30
1	D	160	GLU	OE1-CD-OE2	-7.87	113.86	123.30
1	B	130	ARG	NE-CZ-NH2	-7.79	116.41	120.30
1	C	160	GLU	OE1-CD-OE2	-7.64	114.13	123.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	258	ARG	NE-CZ-NH1	7.52	124.06	120.30
1	B	160	GLU	CG-CD-OE1	7.43	133.16	118.30
1	C	130	ARG	NE-CZ-NH2	-7.21	116.69	120.30
1	A	84	ARG	NE-CZ-NH1	7.21	123.90	120.30
1	D	84	ARG	NE-CZ-NH1	-7.18	116.71	120.30
1	D	84	ARG	NE-CZ-NH2	7.09	123.85	120.30
1	C	29	ARG	NE-CZ-NH1	6.99	123.80	120.30
1	D	116	ARG	NE-CZ-NH2	-6.95	116.83	120.30
1	B	160	GLU	CG-CD-OE2	6.92	132.13	118.30
1	C	229	ASP	CB-CG-OD1	6.41	124.07	118.30
1	C	142	ASP	CB-CG-OD1	6.36	124.02	118.30
1	A	142	ASP	CB-CG-OD1	6.22	123.90	118.30
1	A	84	ARG	NE-CZ-NH2	6.07	123.33	120.30
1	D	130	ARG	NE-CZ-NH2	-6.06	117.27	120.30
1	A	84	ARG	NH1-CZ-NH2	-6.05	112.75	119.40
1	A	249	GLU	OE1-CD-OE2	-6.00	116.10	123.30
1	B	249	GLU	OE1-CD-OE2	-5.97	116.13	123.30
1	D	192	ARG	NE-CZ-NH1	5.87	123.23	120.30
1	B	192	ARG	CD-NE-CZ	5.79	131.70	123.60
1	A	29	ARG	NE-CZ-NH1	5.71	123.15	120.30
1	B	29	ARG	NE-CZ-NH1	5.66	123.13	120.30
1	D	44	ASN	CB-CA-C	5.63	121.67	110.40
1	A	41	ASP	CB-CG-OD1	5.60	123.34	118.30
1	D	229	ASP	CB-CG-OD1	5.54	123.29	118.30
1	C	252	ARG	NE-CZ-NH2	5.48	123.04	120.30
1	A	192	ARG	CD-NE-CZ	5.40	131.16	123.60
1	D	29	ARG	NE-CZ-NH1	5.34	122.97	120.30
1	B	229	ASP	CB-CG-OD1	5.32	123.09	118.30
1	D	258	ARG	NE-CZ-NH1	5.30	122.95	120.30
1	D	249	GLU	OE1-CD-OE2	-5.29	116.96	123.30
1	A	252	ARG	NE-CZ-NH1	5.21	122.90	120.30
1	B	116	ARG	NE-CZ-NH2	-5.18	117.71	120.30
1	C	116	ARG	NE-CZ-NH2	-5.18	117.71	120.30
1	D	254	ASP	CB-CG-OD1	5.17	122.95	118.30
1	C	192	ARG	NE-CZ-NH2	5.14	122.87	120.30
1	C	220	GLN	N-CA-C	5.13	124.86	111.00
1	B	242	GLU	OE1-CD-OE2	-5.00	117.30	123.30

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	1730	0	1762	18	0
1	B	1733	0	1765	21	0
1	C	1735	0	1767	15	0
1	D	1739	0	1772	22	0
2	A	44	0	26	0	0
2	B	44	0	26	0	0
2	C	44	0	26	0	0
2	D	44	0	26	0	0
3	A	109	0	0	1	0
3	B	130	0	0	3	0
3	C	124	0	0	0	0
3	D	129	0	0	1	0
All	All	7605	0	7170	51	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 4.

All (51) 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:236:LEU:HD12	1:D:239:MET:HE1	1.23	1.11
1:B:239:MET:HB2	1:D:239:MET:HE3	1.43	1.01
1:B:236:LEU:HD12	1:D:239:MET:CE	1.93	0.99
1:A:260:GLN:HE21	1:D:260:GLN:HE21	1.20	0.90
1:B:260:GLN:HE21	1:C:260:GLN:HE21	1.20	0.89
1:B:239:MET:HB2	1:D:239:MET:CE	2.06	0.85
1:A:239:MET:HG3	1:C:239:MET:HG3	1.64	0.79
1:B:236:LEU:CD1	1:D:239:MET:HE1	2.11	0.79
1:A:260:GLN:HG3	1:D:260:GLN:NE2	1.99	0.76
1:B:260:GLN:NE2	1:C:260:GLN:HG3	2.06	0.70
1:A:260:GLN:HE21	1:D:260:GLN:NE2	1.91	0.69
1:A:260:GLN:NE2	1:D:260:GLN:HG3	2.08	0.68
1:A:260:GLN:NE2	1:D:260:GLN:HE21	1.90	0.68
1:B:239:MET:CB	1:D:239:MET:HE3	2.25	0.64
1:A:102:HIS:HD2	1:A:103:GLU:N	1.95	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:260:GLN:NE2	1:C:260:GLN:HE21	1.96	0.63
1:B:49:THR:HG23	3:B:2028:HOH:O	1.99	0.62
1:B:260:GLN:HE21	1:C:260:GLN:NE2	1.95	0.62
1:A:260:GLN:HG3	1:D:260:GLN:HE21	1.64	0.62
1:B:260:GLN:HG3	1:C:260:GLN:NE2	2.17	0.59
1:B:260:GLN:HE21	1:C:260:GLN:HG3	1.67	0.58
1:D:102:HIS:CD2	1:D:105:LYS:H	2.23	0.57
1:D:102:HIS:HD2	1:D:105:LYS:H	1.53	0.56
1:A:102:HIS:C	1:A:102:HIS:CD2	2.81	0.52
1:A:239:MET:HG3	1:C:239:MET:CG	2.39	0.52
1:A:36:THR:HG23	3:A:2012:HOH:O	2.09	0.51
1:B:116:ARG:HD3	3:B:2075:HOH:O	2.09	0.51
1:A:260:GLN:HE21	1:D:260:GLN:HG3	1.75	0.49
1:A:223:PHE:CD1	1:A:224:PRO:HA	2.48	0.48
1:A:261:PRO:HG3	1:B:184:ARG:NE	2.28	0.47
1:C:115:GLN:OE1	1:D:123:ILE:HD13	2.15	0.47
1:C:223:PHE:CD1	1:C:224:PRO:HA	2.49	0.47
1:A:187:ALA:HB3	1:A:188:PRO:HD3	1.96	0.47
1:C:187:ALA:HB3	1:C:188:PRO:HD3	1.97	0.47
1:A:203:THR:HB	1:A:204:PRO:CD	2.47	0.45
1:B:10:GLY:O	1:B:84:ARG:NH1	2.50	0.44
1:C:223:PHE:HA	1:C:224:PRO:C	2.38	0.44
1:D:223:PHE:CD1	1:D:224:PRO:HA	2.53	0.44
1:B:111:LEU:O	1:B:115:GLN:HG3	2.17	0.43
1:B:79:LYS:HE3	3:B:2081:HOH:O	2.19	0.43
1:A:102:HIS:C	1:A:102:HIS:HD2	2.21	0.42
1:C:203:THR:HB	1:C:204:PRO:CD	2.49	0.42
1:B:203:THR:HB	1:B:204:PRO:CD	2.49	0.42
1:D:203:THR:HB	1:D:204:PRO:CD	2.49	0.42
1:B:223:PHE:CD1	1:B:224:PRO:HA	2.55	0.41
1:A:243:ASN:HA	1:A:244:PRO:HD3	1.85	0.41
1:D:187:ALA:HB3	1:D:188:PRO:HD3	2.03	0.41
1:C:123:ILE:HD13	1:D:115:GLN:OE1	2.21	0.40
1:C:184:ARG:NE	1:D:261:PRO:HG3	2.37	0.40
1:D:29:ARG:HD2	3:D:2014:HOH:O	2.21	0.40
1:B:242:GLU:O	1:B:244:PRO:HD3	2.21	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	237/261 (91%)	230 (97%)	6 (2%)	1 (0%)	39	33
1	B	237/261 (91%)	227 (96%)	9 (4%)	1 (0%)	39	33
1	C	237/261 (91%)	229 (97%)	7 (3%)	1 (0%)	39	33
1	D	237/261 (91%)	229 (97%)	8 (3%)	0	100	100
All	All	948/1044 (91%)	915 (96%)	30 (3%)	3 (0%)	46	41

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	103	GLU
1	C	220	GLN
1	B	104	LYS

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	176/200 (88%)	173 (98%)	3 (2%)	68	71
1	B	177/200 (88%)	175 (99%)	2 (1%)	80	83
1	C	177/200 (88%)	174 (98%)	3 (2%)	68	71
1	D	178/200 (89%)	176 (99%)	2 (1%)	80	83
All	All	708/800 (88%)	698 (99%)	10 (1%)	74	77

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	49	THR
1	A	80	GLU
1	A	84	ARG
1	B	49	THR
1	B	80	GLU
1	C	49	THR
1	C	79	LYS
1	C	84	ARG
1	D	49	THR
1	D	80	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	102	HIS
1	A	260	GLN
1	B	260	GLN
1	C	44	ASN
1	D	102	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](#)

4 ligands are modelled in this entry.

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	301	-	38,48,48	1.03	2 (5%)	47,73,73	1.78	8 (17%)
2	NAD	B	301	-	38,48,48	1.08	3 (7%)	47,73,73	2.14	10 (21%)
2	NAD	C	301	-	38,48,48	1.15	3 (7%)	47,73,73	1.85	9 (19%)
2	NAD	D	301	-	38,48,48	1.13	2 (5%)	47,73,73	2.38	12 (25%)

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	301	-	-	0/22/62/62	0/5/5/5
2	NAD	B	301	-	-	0/22/62/62	0/5/5/5
2	NAD	C	301	-	-	0/22/62/62	0/5/5/5
2	NAD	D	301	-	-	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	301	NAD	PN-O2N	-2.10	1.46	1.54
2	B	301	NAD	C2A-N1A	2.12	1.37	1.33
2	A	301	NAD	C6N-N1N	2.35	1.41	1.35
2	B	301	NAD	C6N-N1N	2.37	1.41	1.35
2	D	301	NAD	C6N-N1N	2.84	1.43	1.35
2	C	301	NAD	C6N-N1N	3.06	1.43	1.35
2	A	301	NAD	C3N-C7N	3.40	1.55	1.50
2	C	301	NAD	C3N-C7N	3.97	1.56	1.50
2	B	301	NAD	C3N-C7N	4.01	1.56	1.50
2	D	301	NAD	C3N-C7N	4.12	1.57	1.50

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	301	NAD	C5N-C4N-C3N	-7.02	111.50	120.33
2	B	301	NAD	C5N-C4N-C3N	-6.80	111.78	120.33
2	C	301	NAD	C5N-C4N-C3N	-5.97	112.83	120.33
2	A	301	NAD	C5N-C4N-C3N	-5.76	113.09	120.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	301	NAD	C4N-C3N-C7N	-5.09	107.65	121.09
2	D	301	NAD	O7N-C7N-C3N	-4.84	114.31	119.59
2	B	301	NAD	C4N-C3N-C7N	-4.30	109.74	121.09
2	A	301	NAD	C4N-C3N-C7N	-4.28	109.78	121.09
2	D	301	NAD	C5N-C6N-N1N	-3.99	113.57	120.47
2	B	301	NAD	C3N-C7N-N7N	-3.96	113.48	117.82
2	C	301	NAD	C4N-C3N-C7N	-3.90	110.78	121.09
2	B	301	NAD	C2B-C1B-N9A	-3.85	108.40	114.29
2	C	301	NAD	C5N-C6N-N1N	-3.78	113.94	120.47
2	A	301	NAD	C5N-C6N-N1N	-3.47	114.47	120.47
2	D	301	NAD	C4B-O4B-C1B	-3.41	105.98	109.72
2	B	301	NAD	C5N-C6N-N1N	-3.39	114.61	120.47
2	D	301	NAD	C2B-C1B-N9A	-2.67	110.21	114.29
2	C	301	NAD	C3N-C7N-N7N	-2.44	115.14	117.82
2	D	301	NAD	C3N-C2N-N1N	-2.35	117.66	120.36
2	A	301	NAD	C3N-C2N-N1N	-2.13	117.91	120.36
2	C	301	NAD	C4B-O4B-C1B	-2.04	107.48	109.72
2	C	301	NAD	C2B-C1B-N9A	-2.02	111.21	114.29
2	A	301	NAD	C4B-O4B-C1B	-2.01	107.51	109.72
2	B	301	NAD	C2N-C3N-C7N	2.10	125.40	119.31
2	C	301	NAD	C2N-C3N-C7N	2.16	125.57	119.31
2	D	301	NAD	C4A-C5A-N7A	2.21	111.52	109.48
2	B	301	NAD	O2A-PA-O3	2.30	115.52	105.09
2	A	301	NAD	C2N-C3N-C7N	2.46	126.45	119.31
2	D	301	NAD	C2N-C3N-C7N	2.66	127.03	119.31
2	B	301	NAD	O7N-C7N-N7N	2.81	126.54	122.59
2	D	301	NAD	O7N-C7N-N7N	4.58	129.05	122.59
2	C	301	NAD	C2N-C3N-C4N	4.81	123.64	118.29
2	A	301	NAD	C2N-C3N-C4N	4.92	123.77	118.29
2	A	301	NAD	C6N-C5N-C4N	5.03	127.05	119.44
2	C	301	NAD	C6N-C5N-C4N	5.42	127.63	119.44
2	B	301	NAD	C6N-C5N-C4N	5.46	127.69	119.44
2	D	301	NAD	C6N-C5N-C4N	5.79	128.19	119.44
2	B	301	NAD	C2N-C3N-C4N	5.88	124.84	118.29
2	D	301	NAD	C2N-C3N-C4N	6.30	125.31	118.29

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 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.