



Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 08:06 pm BST

PDB ID : 2DC1
Title : Crystal Structure Of L-Aspartate Dehydrogenase From Hyperthermophilic Archaeon Archaeoglobus fulgidus
Authors : Yoneda, K.; Sakuraba, H.; Tsuge, H.; Ohshima, T.
Deposited on : 2005-12-19
Resolution : 1.90 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtrriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
buster-report : 1.1.7 (2018)
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) : 2.11

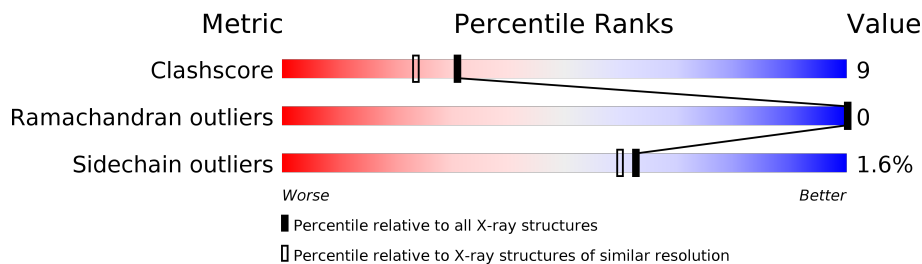
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)

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 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	236	
1	B	236	

2 Entry composition [i](#)

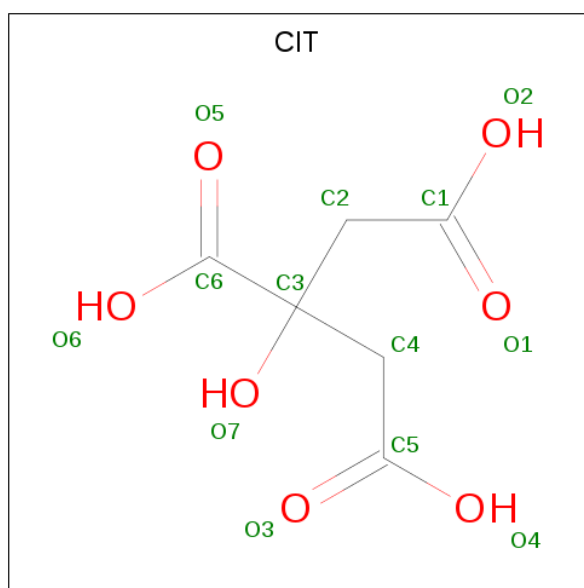
There are 4 unique types of molecules in this entry. The entry contains 3936 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 L-aspartate dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	236	Total 1843	C 1165	N 332	O 340	S 6	0	0	0
1	B	236	Total 1843	C 1165	N 332	O 340	S 6	0	0	0

- Molecule 2 is CITRIC ACID (three-letter code: CIT) (formula: C₆H₈O₇).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	A	1	Total 13	C 6	O 7	0	0
2	B	1	Total 13	C 6	O 7	0	0

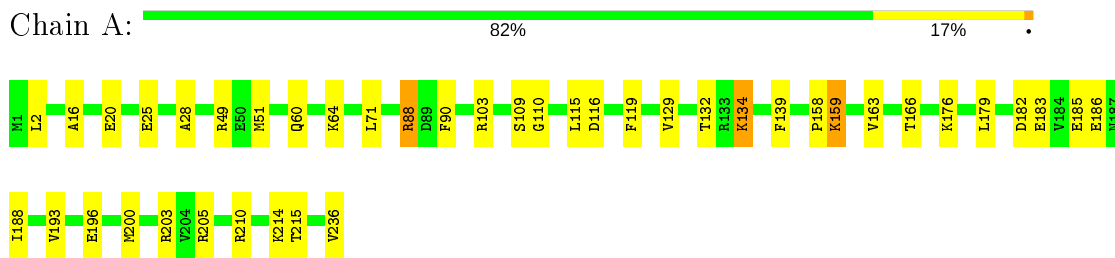
- Molecule 3 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: C₂₁H₂₇N₇O₁₄P₂).

3 Residue-property plots [i](#)

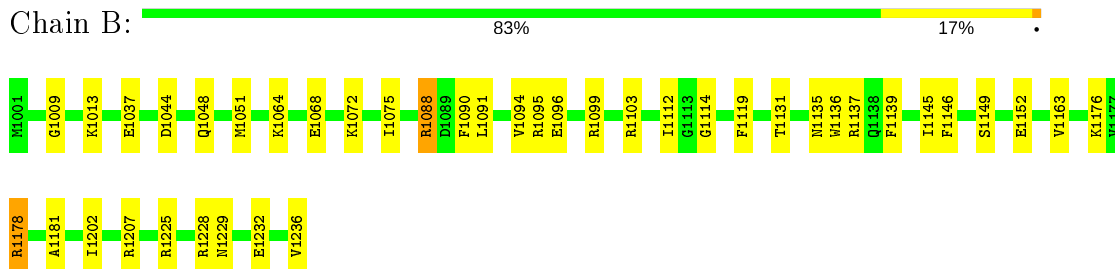
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded according 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: L-aspartate dehydrogenase



- Molecule 1: L-aspartate dehydrogenase



4 Data and refinement statistics

Xtrriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	47.52Å 89.58Å 100.49Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	32.60 – 1.90	Depositor
% Data completeness (in resolution range)	(Not available) (32.60-1.90)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.217 , 0.226	Depositor
Estimated twinning fraction	No twinning to report.	Xtrriage
Total number of atoms	3936	wwPDB-VP
Average B, all atoms (Å ²)	29.0	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, CIT

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	0.30	0/1865	0.57	0/2505
1	B	0.32	0/1865	0.57	0/2505
All	All	0.31	0/3730	0.57	0/5010

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	A	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	203	ARG	Sidechain

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	1843	0	1903	33	0
1	B	1843	0	1900	36	0
2	A	13	0	5	3	0
2	B	13	0	5	1	0
3	A	44	0	26	4	0
3	B	44	0	26	2	0
4	A	70	0	0	1	0
4	B	66	0	0	0	0
All	All	3936	0	3865	70	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 9.

All (70) 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:1145:ILE:HD11	1:B:1181:ALA:HB2	1.27	1.13
1:B:1096:GLU:HG2	1:B:1099:ARG:HH22	1.26	0.99
1:A:119:PHE:HD2	1:B:1225:ARG:HG3	1.33	0.93
1:A:119:PHE:CD2	1:B:1225:ARG:HG3	2.14	0.82
1:A:88:ARG:H	1:A:88:ARG:HD3	1.45	0.82
1:B:1112:ILE:HD13	1:B:1163:VAL:HG13	1.62	0.81
1:A:186:GLU:OE1	1:A:205:ARG:HG2	1.83	0.78
1:B:1131:THR:HG23	1:B:1178:ARG:HD2	1.72	0.71
1:B:1146:PHE:HZ	1:B:1152:GLU:HG3	1.58	0.68
1:A:134:LYS:HD3	1:A:179:LEU:HD11	1.76	0.67
1:B:1131:THR:HG23	1:B:1178:ARG:CD	2.25	0.66
1:A:158:PRO:O	1:A:159:LYS:HG2	1.96	0.65
1:A:88:ARG:CD	1:A:88:ARG:H	2.08	0.65
1:B:1068:GLU:O	1:B:1072:LYS:HG2	1.97	0.64
1:B:1096:GLU:HG2	1:B:1099:ARG:NH2	2.06	0.64
1:A:109:SER:HA	1:A:115:LEU:HD12	1.82	0.61
1:B:1088:ARG:HD2	1:B:1088:ARG:H	1.66	0.60
1:B:1145:ILE:HD11	1:B:1181:ALA:CB	2.18	0.60
1:B:1136:TRP:HA	1:B:1181:ALA:HB1	1.85	0.59
1:A:115:LEU:HD11	1:A:166:THR:HG21	1.87	0.57
1:B:1114:GLY:HA3	1:B:1202:ILE:CD1	2.36	0.55
1:B:1090:PHE:O	1:B:1094:VAL:HG23	2.07	0.55
2:A:1585:CIT:H42	3:A:242:NAD:C4N	2.38	0.54
1:B:1131:THR:HA	1:B:1178:ARG:HG3	1.89	0.53
1:A:236:VAL:HG13	1:B:1119:PHE:CD2	2.44	0.52
3:A:242:NAD:H52N	3:A:242:NAD:H6N	1.90	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1135:ASN:HD21	1:B:1137:ARG:NH1	2.07	0.52
1:A:196:GLU:O	1:B:1207:ARG:HD2	2.11	0.51
3:B:1242:NAD:H6N	3:B:1242:NAD:H52N	1.92	0.51
1:A:134:LYS:NZ	1:A:134:LYS:HB3	2.25	0.51
1:A:129:VAL:HG22	1:A:176:LYS:HB2	1.93	0.51
1:B:1044:ASP:O	1:B:1048:GLN:HG3	2.11	0.50
2:B:2585:CIT:H42	3:B:1242:NAD:C4N	2.41	0.50
1:B:1064:LYS:HA	1:B:1090:PHE:CZ	2.46	0.50
1:B:1178:ARG:HG3	1:B:1178:ARG:O	2.12	0.49
1:A:215:THR:HG23	3:A:242:NAD:H5N	1.94	0.49
1:A:28:ALA:HB2	1:A:49:ARG:HG3	1.95	0.49
1:B:1091:LEU:HD21	1:B:1095:ARG:NH2	2.28	0.48
1:B:1139:PHE:CG	1:B:1145:ILE:HD13	2.48	0.48
1:A:132:THR:HG22	1:A:134:LYS:HD2	1.95	0.48
1:B:1139:PHE:CD1	1:B:1145:ILE:HD13	2.49	0.48
1:A:2:LEU:HB3	1:A:51:MET:HA	1.96	0.47
1:B:1135:ASN:HD21	1:B:1137:ARG:HH12	1.61	0.47
1:A:163:VAL:HG23	2:A:1585:CIT:H41	1.97	0.46
1:A:210:ARG:HG3	1:A:210:ARG:HH11	1.80	0.46
1:B:1103:ARG:HA	1:B:1103:ARG:NE	2.31	0.46
1:B:1131:THR:CG2	1:B:1178:ARG:HD2	2.45	0.46
1:A:71:LEU:HD22	1:A:103:ARG:HB3	1.97	0.45
1:A:110:GLY:HA3	3:A:242:NAD:O7N	2.19	0.43
1:A:193:VAL:HB	1:A:200:MET:HG2	2.00	0.43
1:B:1229:ASN:HD21	1:B:1236:VAL:H	1.66	0.43
1:B:1037:GLU:N	1:B:1037:GLU:OE1	2.37	0.42
1:B:1176:LYS:HB3	1:B:1176:LYS:HE2	1.88	0.42
1:A:134:LYS:HG3	1:A:139:PHE:HE1	1.84	0.41
1:A:188:ILE:HG12	1:A:205:ARG:HG3	2.02	0.41
1:A:214:LYS:HA	4:A:1607:HOH:O	2.19	0.41
1:B:1009:GLY:O	1:B:1013:LYS:HG3	2.20	0.41
1:A:2:LEU:HD12	1:A:25:GLU:HB3	2.03	0.41
1:A:182:ASP:OD1	1:A:183:GLU:N	2.48	0.41
1:A:60:GLN:HG3	1:A:159:LYS:HA	2.03	0.41
1:B:1135:ASN:ND2	1:B:1137:ARG:NH1	2.69	0.41
1:B:1149:SER:OG	1:B:1152:GLU:HG2	2.21	0.41
1:A:16:ALA:O	1:A:20:GLU:HG3	2.21	0.41
1:B:1228:ARG:O	1:B:1232:GLU:HG2	2.21	0.41
1:A:64:LYS:HA	1:A:90:PHE:CZ	2.56	0.41
1:A:134:LYS:HZ3	1:A:134:LYS:HB3	1.85	0.40
2:A:1585:CIT:C5	2:A:1585:CIT:C1	2.98	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:2:LEU:HA	1:A:2:LEU:HD12	1.95	0.40
1:A:185:GLU:O	1:A:185:GLU:CD	2.60	0.40
1:B:1051:MET:CG	1:B:1075:ILE:HD13	2.52	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	234/236 (99%)	227 (97%)	7 (3%)	0	100	100
1	B	234/236 (99%)	226 (97%)	8 (3%)	0	100	100
All	All	468/472 (99%)	453 (97%)	15 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	191/191 (100%)	187 (98%)	4 (2%)	53	48
1	B	191/191 (100%)	189 (99%)	2 (1%)	76	76
All	All	382/382 (100%)	376 (98%)	6 (2%)	62	60

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	88	ARG
1	A	116	ASP
1	A	134	LYS
1	A	159	LYS
1	B	1088	ARG
1	B	1178	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	48	GLN
1	A	60	GLN
1	A	155	GLN
1	A	160	ASN
1	A	206	ASN
1	B	1229	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](#)

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	CIT	A	1585	-	3,12,12	3.52	2 (66%)	3,17,17	5.30	3 (100%)
3	NAD	A	242	-	42,48,48	2.31	10 (23%)	50,73,73	2.14	16 (32%)
2	CIT	B	2585	-	3,12,12	5.59	2 (66%)	3,17,17	4.89	2 (66%)
3	NAD	B	1242	-	42,48,48	2.70	10 (23%)	50,73,73	2.21	16 (32%)

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	CIT	A	1585	-	-	2/6/16/16	-
3	NAD	A	242	-	-	4/26/62/62	0/5/5/5
2	CIT	B	2585	-	-	3/6/16/16	-
3	NAD	B	1242	-	-	4/26/62/62	0/5/5/5

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1242	NAD	O7N-C7N	13.41	1.49	1.24
3	A	242	NAD	O7N-C7N	10.83	1.44	1.24
2	B	2585	CIT	C4-C3	-9.19	1.41	1.54
2	A	1585	CIT	O7-C3	5.17	1.51	1.43
3	A	242	NAD	O5D-C5D	-4.67	1.26	1.44
3	B	1242	NAD	C2A-N3A	4.35	1.39	1.32
3	B	1242	NAD	C2D-C1D	4.28	1.60	1.53
3	B	1242	NAD	O5D-C5D	-3.70	1.30	1.44
3	B	1242	NAD	C4N-C3N	3.12	1.44	1.39
2	A	1585	CIT	C4-C3	-3.06	1.50	1.54
3	A	242	NAD	C2A-N3A	3.00	1.36	1.32
3	B	1242	NAD	C7N-N7N	2.96	1.38	1.33
3	B	1242	NAD	C2A-N1A	2.93	1.39	1.33
3	A	242	NAD	O4B-C1B	2.86	1.45	1.41
2	B	2585	CIT	O7-C3	2.85	1.47	1.43
3	A	242	NAD	C6A-C5A	-2.44	1.34	1.43
3	B	1242	NAD	C2D-C3D	-2.40	1.46	1.53
3	B	1242	NAD	PN-O1N	-2.33	1.42	1.50
3	A	242	NAD	O4B-C4B	-2.20	1.40	1.45
3	A	242	NAD	C2D-C3D	-2.19	1.47	1.53
3	A	242	NAD	C4N-C3N	2.19	1.43	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	242	NAD	C2N-N1N	2.15	1.37	1.35
3	B	1242	NAD	O4B-C1B	-2.11	1.38	1.41
3	A	242	NAD	C5A-N7A	-2.04	1.32	1.39

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	2585	CIT	C3-C4-C5	-7.09	103.63	114.98
2	A	1585	CIT	C3-C2-C1	7.01	126.20	114.98
3	B	1242	NAD	C5D-C4D-C3D	-6.82	89.60	115.18
3	A	242	NAD	C5D-C4D-C3D	-6.29	91.62	115.18
2	A	1585	CIT	C3-C4-C5	-5.41	106.32	114.98
3	B	1242	NAD	N3A-C2A-N1A	-5.26	120.45	128.68
3	B	1242	NAD	O2D-C2D-C3D	-4.64	96.80	111.82
3	B	1242	NAD	C3N-C2N-N1N	4.53	124.86	120.43
3	A	242	NAD	C6N-N1N-C2N	-4.38	117.98	121.97
3	A	242	NAD	N3A-C2A-N1A	-4.37	121.85	128.68
2	B	2585	CIT	C3-C2-C1	4.37	121.98	114.98
3	A	242	NAD	C3N-C7N-N7N	3.87	122.40	117.75
3	B	1242	NAD	O2N-PN-O1N	3.64	130.22	112.24
3	A	242	NAD	O2D-C2D-C3D	-3.55	100.33	111.82
3	A	242	NAD	C3N-C2N-N1N	3.35	123.70	120.43
3	A	242	NAD	C3D-C2D-C1D	3.33	105.99	100.98
3	A	242	NAD	O2N-PN-O1N	3.25	128.33	112.24
3	B	1242	NAD	C3N-C7N-N7N	3.23	121.62	117.75
3	A	242	NAD	O7N-C7N-C3N	-3.17	115.84	119.63
3	A	242	NAD	PN-O3-PA	-3.15	122.03	132.83
3	B	1242	NAD	O4B-C1B-C2B	-2.95	102.62	106.93
3	A	242	NAD	O2N-PN-O5D	-2.92	94.18	107.75
3	B	1242	NAD	C6N-N1N-C2N	-2.91	119.32	121.97
3	A	242	NAD	O4D-C4D-C3D	2.76	110.57	105.11
3	B	1242	NAD	O3D-C3D-C2D	-2.69	103.13	111.82
3	B	1242	NAD	C1B-N9A-C4A	-2.46	122.32	126.64
3	B	1242	NAD	O4D-C4D-C5D	2.43	117.38	109.37
2	A	1585	CIT	C4-C3-C2	-2.40	102.91	109.33
3	B	1242	NAD	C4A-C5A-N7A	-2.37	106.93	109.40
3	B	1242	NAD	O2N-PN-O5D	-2.34	96.87	107.75
3	A	242	NAD	C5B-C4B-C3B	-2.33	106.44	115.18
3	A	242	NAD	C4A-C5A-N7A	-2.33	106.97	109.40
3	B	1242	NAD	O2B-C2B-C3B	2.29	119.24	111.82
3	A	242	NAD	C1B-N9A-C4A	-2.28	122.64	126.64
3	B	1242	NAD	N6A-C6A-N1A	2.20	123.15	118.57

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	242	NAD	C2A-N1A-C6A	2.13	122.40	118.75
3	B	1242	NAD	O4D-C4D-C3D	2.01	109.10	105.11

There are no chirality outliers.

All (13) torsion outliers are listed below:

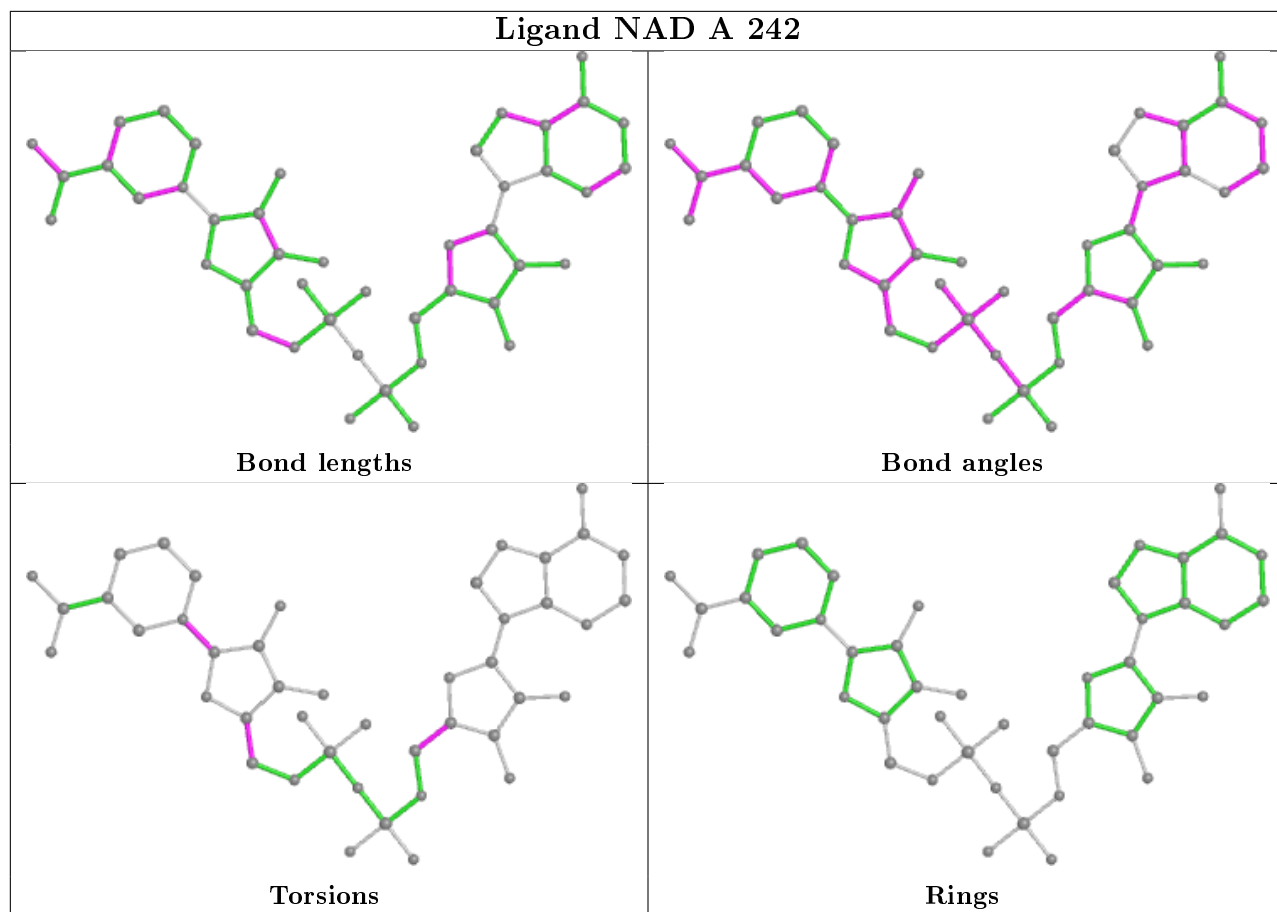
Mol	Chain	Res	Type	Atoms
3	A	242	NAD	O4D-C1D-N1N-C6N
2	B	2585	CIT	C1-C2-C3-O7
2	B	2585	CIT	C1-C2-C3-C4
3	B	1242	NAD	O4D-C1D-N1N-C6N
2	A	1585	CIT	C1-C2-C3-C4
2	A	1585	CIT	C1-C2-C3-O7
3	A	242	NAD	O4D-C4D-C5D-O5D
3	A	242	NAD	C3D-C4D-C5D-O5D
3	B	1242	NAD	C5D-O5D-PN-O3
2	B	2585	CIT	C1-C2-C3-C6
3	A	242	NAD	O4B-C4B-C5B-O5B
3	B	1242	NAD	O4B-C4B-C5B-O5B
3	B	1242	NAD	C5D-O5D-PN-O1N

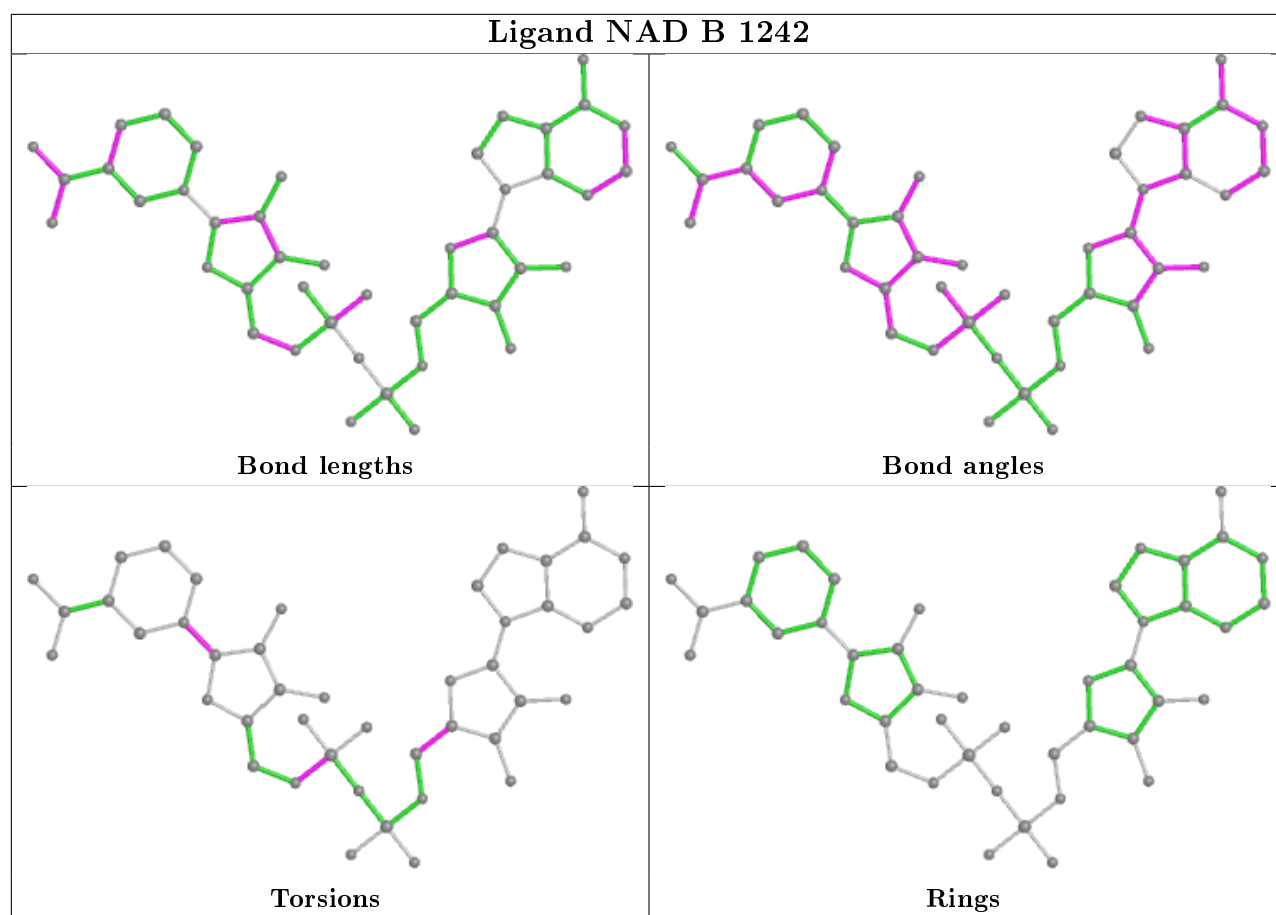
There are no ring outliers.

4 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1585	CIT	3	0
3	A	242	NAD	4	0
2	B	2585	CIT	1	0
3	B	1242	NAD	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





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.