Full wwPDB X-ray Structure Validation Report (i)

Feb 3, 2024 – 01:32 PM EST

PDB ID	:	1LLD
Title	:	MOLECULAR BASIS OF ALLOSTERIC ACTIVATION OF BACTERIAL L-LACTATE DEHYDROGENASE
Authors	:	Iwata, S.; Ohta, T.
Deposited on	:	1992-06-08
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
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

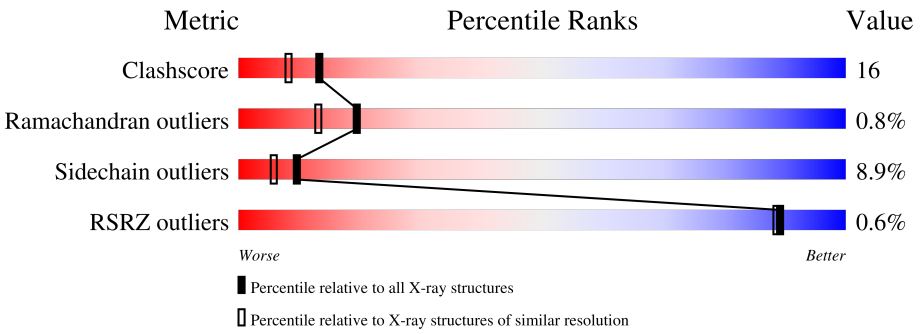
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 2.00 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	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (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 of 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 <=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	319	65% 26% 6% • •
1	B	319	% 66% 24% 7% • •

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 5010 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.

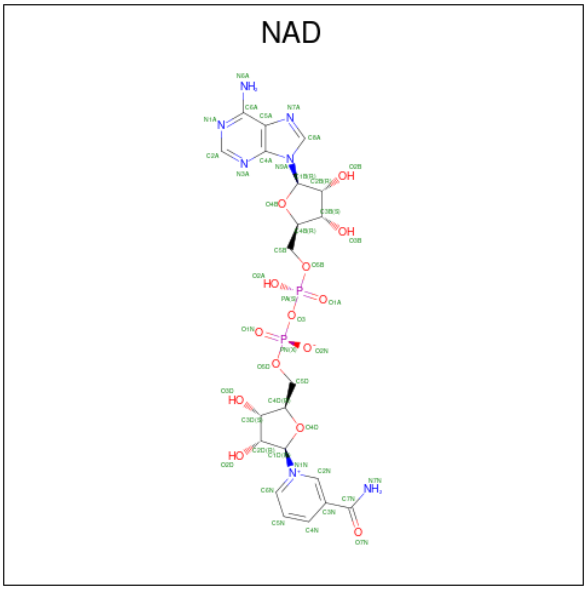
- Molecule 1 is a protein called L-LACTATE DEHYDROGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	313	2350	1474	409	457	10	0	0	0
1	B	313	2350	1474	409	457	10	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	199	SER	CYS	conflict	UNP P19869
B	199	SER	CYS	conflict	UNP P19869

- 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
			Total	C	N	O	P		
2	A	1	44	21	7	14	2	0	0

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

- Molecule 3 is water.

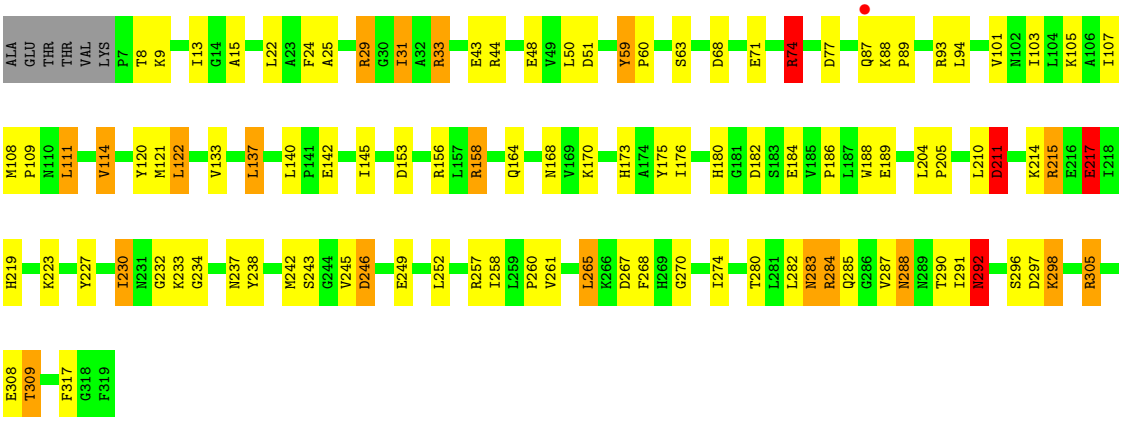
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	112	112	112	0	0
			Total	O		
3	B	110	110	110	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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 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.

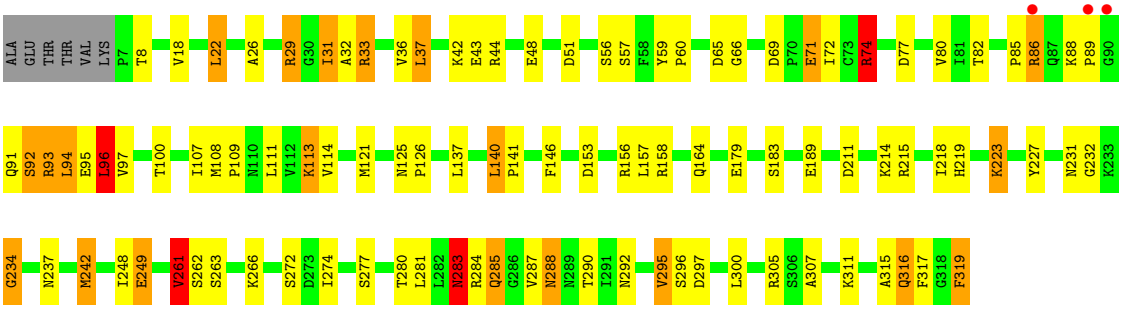
• Molecule 1: L-LACTATE DEHYDROGENASE

Chain A:



• Molecule 1: L-LACTATE DEHYDROGENASE

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α, β, γ	105.80Å 131.40Å 63.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.00 82.41 – 2.00	Depositor EDS
% Data completeness (in resolution range)	(Not available) (10.00-2.00) 79.4 (82.41-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.28 (at 2.00Å)	Xtriage
Refinement program	PROLSQ	Depositor
R, R_{free}	0.179 , (Not available) 0.169 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å^2)	22.0	Xtriage
Anisotropy	0.140	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 83.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	5010	wwPDB-VP
Average B, all atoms (Å^2)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.94% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

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	0.96	0/2390	1.69	51/3248 (1.6%)
1	B	0.95	0/2390	1.61	35/3248 (1.1%)
All	All	0.96	0/4780	1.65	86/6496 (1.3%)

There are no bond length outliers.

All (86) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	44	ARG	NE-CZ-NH1	-13.19	113.71	120.30
1	A	29	ARG	NE-CZ-NH1	12.65	126.63	120.30
1	B	29	ARG	NE-CZ-NH1	12.00	126.30	120.30
1	A	93	ARG	NE-CZ-NH1	11.89	126.24	120.30
1	A	74	ARG	NE-CZ-NH1	11.30	125.95	120.30
1	B	33	ARG	NE-CZ-NH1	10.90	125.75	120.30
1	A	93	ARG	CD-NE-CZ	10.79	138.71	123.60
1	A	158	ARG	NE-CZ-NH1	10.29	125.45	120.30
1	A	211	ASP	CB-CG-OD1	-10.12	109.19	118.30
1	B	71	GLU	OE1-CD-OE2	9.16	134.30	123.30
1	B	305	ARG	NE-CZ-NH1	9.08	124.84	120.30
1	A	74	ARG	NE-CZ-NH2	-8.97	115.81	120.30
1	B	86	ARG	NE-CZ-NH1	8.94	124.77	120.30
1	A	44	ARG	NE-CZ-NH2	8.59	124.60	120.30
1	A	44	ARG	CD-NE-CZ	-8.43	111.80	123.60
1	A	68	ASP	CB-CG-OD1	8.22	125.70	118.30
1	B	215	ARG	NE-CZ-NH2	-7.92	116.34	120.30
1	A	227	TYR	CB-CG-CD1	-7.76	116.34	121.00
1	A	158	ARG	CD-NE-CZ	7.69	134.37	123.60
1	B	153	ASP	CB-CG-OD2	-7.44	111.61	118.30
1	A	93	ARG	NE-CZ-NH2	-7.42	116.59	120.30
1	B	156	ARG	NE-CZ-NH1	7.29	123.94	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	215	ARG	NE-CZ-NH1	7.27	123.93	120.30
1	A	257	ARG	NE-CZ-NH1	7.24	123.92	120.30
1	B	156	ARG	CD-NE-CZ	7.24	133.73	123.60
1	A	210	LEU	C-N-CA	7.16	139.61	121.70
1	A	308	GLU	CA-CB-CG	7.13	129.09	113.40
1	B	74	ARG	NE-CZ-NH2	-7.05	116.77	120.30
1	B	153	ASP	CB-CG-OD1	7.04	124.63	118.30
1	B	249	GLU	OE1-CD-OE2	-6.87	115.06	123.30
1	A	215	ARG	NE-CZ-NH2	-6.87	116.87	120.30
1	A	317	PHE	CB-CG-CD1	-6.79	116.04	120.80
1	B	261	VAL	CA-CB-CG1	6.67	120.90	110.90
1	B	33	ARG	NE-CZ-NH2	-6.65	116.97	120.30
1	A	284	ARG	NE-CZ-NH1	6.60	123.60	120.30
1	A	24	PHE	CB-CG-CD1	-6.44	116.29	120.80
1	A	51	ASP	CB-CG-OD1	6.43	124.08	118.30
1	A	77	ASP	CB-CG-OD2	6.38	124.05	118.30
1	A	153	ASP	CB-CG-OD1	6.32	123.98	118.30
1	A	77	ASP	CB-CG-OD1	-6.26	112.66	118.30
1	B	179	GLU	OE1-CD-OE2	6.11	130.63	123.30
1	B	137	LEU	CA-CB-CG	6.10	129.34	115.30
1	B	249	GLU	CA-CB-CG	6.05	126.71	113.40
1	A	211	ASP	N-CA-CB	-6.00	99.80	110.60
1	A	59	TYR	CB-CG-CD2	-5.98	117.41	121.00
1	B	283	ASN	CB-CA-C	5.96	122.31	110.40
1	A	298	LYS	CA-CB-CG	5.92	126.44	113.40
1	A	121	MET	CA-CB-CG	-5.92	103.23	113.30
1	A	59	TYR	CB-CG-CD1	5.91	124.55	121.00
1	B	263	SER	N-CA-CB	5.91	119.36	110.50
1	A	121	MET	CG-SD-CE	5.90	109.64	100.20
1	A	184	GLU	CG-CD-OE2	-5.88	106.55	118.30
1	B	96	LEU	CB-CA-C	5.86	121.33	110.20
1	A	156	ARG	CD-NE-CZ	5.84	131.78	123.60
1	A	122	LEU	CA-CB-CG	5.80	128.65	115.30
1	A	297	ASP	CB-CG-OD1	5.79	123.51	118.30
1	A	94	LEU	CB-CA-C	5.67	120.97	110.20
1	A	246	ASP	CB-CG-OD2	5.67	123.40	118.30
1	B	51	ASP	CB-CG-OD1	5.60	123.34	118.30
1	B	284	ARG	NE-CZ-NH1	5.59	123.10	120.30
1	A	33	ARG	NE-CZ-NH1	5.53	123.06	120.30
1	B	29	ARG	CD-NE-CZ	5.48	131.27	123.60
1	B	215	ARG	NE-CZ-NH1	5.46	123.03	120.30
1	B	261	VAL	N-CA-CB	-5.40	99.61	111.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	140	LEU	CA-CB-CG	5.40	127.72	115.30
1	B	93	ARG	NE-CZ-NH1	5.39	122.99	120.30
1	B	211	ASP	CB-CG-OD1	-5.38	113.46	118.30
1	A	137	LEU	N-CA-CB	-5.38	99.64	110.40
1	A	211	ASP	CA-CB-CG	-5.35	101.64	113.40
1	B	263	SER	O-C-N	5.31	131.20	122.70
1	A	142	GLU	CB-CA-C	-5.31	99.78	110.40
1	B	22	LEU	CB-CA-C	5.26	120.20	110.20
1	B	305	ARG	CA-CB-CG	5.23	124.90	113.40
1	A	74	ARG	CA-CB-CG	5.22	124.88	113.40
1	A	43	GLU	OE1-CD-OE2	5.13	129.46	123.30
1	B	297	ASP	CB-CG-OD1	5.11	122.90	118.30
1	A	153	ASP	CB-CG-OD2	-5.10	113.71	118.30
1	A	87	GLN	CB-CA-C	5.07	120.55	110.40
1	B	158	ARG	NE-CZ-NH1	5.06	122.83	120.30
1	A	158	ARG	NH1-CZ-NH2	-5.05	113.85	119.40
1	A	184	GLU	CG-CD-OE1	5.04	128.39	118.30
1	A	257	ARG	CD-NE-CZ	5.04	130.65	123.60
1	B	113	LYS	C-N-CA	5.03	134.27	121.70
1	A	217	GLU	CG-CD-OE1	-5.03	108.25	118.30
1	B	29	ARG	CG-CD-NE	5.01	122.32	111.80
1	A	291	ILE	CB-CA-C	5.01	121.62	111.60

There are no chirality outliers.

There are no planarity outliers.

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	2350	0	2366	77	0
1	B	2350	0	2367	78	0
2	A	44	0	26	0	0
2	B	44	0	26	3	0
3	A	112	0	0	4	0
3	B	110	0	0	2	0
All	All	5010	0	4785	155	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 16.

All (155) 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:88:LYS:CG	1:A:89:PRO:HD2	1.69	1.23
1:B:92:SER:OG	1:B:95:GLU:HG3	1.46	1.15
1:B:88:LYS:HB3	1:B:89:PRO:CD	1.76	1.15
1:A:88:LYS:HG2	1:A:89:PRO:CD	1.79	1.12
1:B:88:LYS:HB3	1:B:89:PRO:HD2	1.09	1.06
1:A:71:GLU:OE2	1:A:74:ARG:HD2	1.67	0.95
1:A:305:ARG:NH1	1:A:309:THR:HG23	1.82	0.95
1:B:88:LYS:CB	1:B:89:PRO:HD2	2.00	0.90
1:B:272:SER:O	1:B:311:LYS:NZ	2.05	0.88
1:B:94:LEU:HD11	1:B:316:GLN:OE1	1.75	0.86
1:B:283:ASN:ND2	1:B:285:GLN:HG2	1.92	0.84
1:A:189:GLU:HG3	1:A:296:SER:HB3	1.58	0.84
1:B:283:ASN:HD21	1:B:285:GLN:HG2	1.45	0.82
1:B:26:ALA:HB1	1:B:31:ILE:HD11	1.63	0.80
1:B:189:GLU:HG3	1:B:296:SER:HB3	1.64	0.79
1:A:282:LEU:CD2	1:A:287:VAL:HG12	2.12	0.79
1:A:88:LYS:HG2	1:A:89:PRO:HD2	0.83	0.77
1:B:71:GLU:OE2	1:B:74:ARG:NH1	2.19	0.76
1:A:29:ARG:HB3	1:A:31:ILE:HD11	1.66	0.76
1:A:305:ARG:HH12	1:A:309:THR:HG23	1.50	0.73
1:A:29:ARG:HB3	1:A:31:ILE:CD1	2.18	0.73
1:A:8:THR:HG23	1:A:33:ARG:H	1.53	0.73
1:A:71:GLU:OE2	1:A:74:ARG:NH1	2.18	0.72
1:B:29:ARG:HD3	3:B:428:HOH:O	1.92	0.69
1:A:282:LEU:HD21	1:A:287:VAL:HG12	1.74	0.68
1:A:29:ARG:NH2	1:A:249:GLU:OE1	2.27	0.68
1:A:288:ASN:HD22	1:A:290:THR:H	1.40	0.66
1:A:282:LEU:HD23	1:A:287:VAL:HG12	1.77	0.66
1:A:133:VAL:O	1:A:137:LEU:HB2	1.96	0.65
1:A:15:ALA:HB1	1:A:48:GLU:OE1	1.97	0.65
1:B:92:SER:OG	1:B:95:GLU:CG	2.37	0.63
1:B:44:ARG:NE	1:B:48:GLU:OE2	2.30	0.63
1:B:88:LYS:CB	1:B:89:PRO:CD	2.55	0.63
1:A:182:ASP:O	1:A:223:LYS:HE3	1.99	0.62
1:A:29:ARG:CB	1:A:31:ILE:HD11	2.30	0.62
1:A:211:ASP:HB3	1:A:214:LYS:H	1.65	0.61
1:A:242:MET:HA	1:A:242:MET:CE	2.31	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:288:ASN:HD22	1:A:288:ASN:C	2.03	0.61
1:A:283:ASN:C	1:A:283:ASN:HD22	2.03	0.60
1:A:267:ASP:OD1	1:A:270:GLY:N	2.35	0.60
1:A:288:ASN:ND2	1:A:290:THR:H	2.00	0.59
1:A:175:TYR:HD1	1:A:260:PRO:HG3	1.68	0.59
1:A:284:ARG:O	1:A:284:ARG:HG3	2.02	0.59
1:A:13:ILE:HG13	1:A:111:LEU:HD21	1.84	0.58
1:A:283:ASN:HD22	1:A:285:GLN:H	1.51	0.58
1:A:71:GLU:OE2	1:A:74:ARG:CD	2.49	0.58
1:A:232:GLY:C	1:A:234:GLY:H	2.07	0.58
1:B:22:LEU:C	1:B:22:LEU:HD23	2.23	0.58
1:B:92:SER:HG	1:B:95:GLU:HG3	1.65	0.57
1:B:66:GLY:HA2	1:B:72:ILE:HD13	1.87	0.56
1:B:85:PRO:HG3	1:B:100:THR:OG1	2.06	0.56
1:B:108:MET:N	1:B:109:PRO:HD2	2.22	0.55
1:A:173:HIS:HB3	3:A:347:HOH:O	2.07	0.55
1:A:74:ARG:HA	1:A:114:VAL:HG13	1.89	0.54
1:B:232:GLY:C	1:B:234:GLY:H	2.08	0.54
1:B:86:ARG:NH1	2:B:320:NAD:O1A	2.42	0.53
1:B:261:VAL:O	1:B:277:SER:HA	2.09	0.53
1:B:280:THR:HG21	1:B:287:VAL:HB	1.89	0.53
1:A:285:GLN:NE2	3:A:405:HOH:O	2.42	0.53
1:A:168:ASN:OD1	1:A:170:LYS:HD3	2.09	0.52
1:B:57:SER:O	1:B:60:PRO:HD3	2.10	0.52
1:B:82:THR:O	2:B:320:NAD:C5D	2.58	0.52
1:B:232:GLY:C	1:B:234:GLY:N	2.60	0.52
1:B:288:ASN:C	1:B:288:ASN:HD22	2.11	0.52
1:A:158:ARG:HD2	3:A:322:HOH:O	2.08	0.52
1:B:319:PHE:CD1	1:B:319:PHE:N	2.78	0.51
1:B:8:THR:HA	1:B:77:ASP:OD2	2.11	0.51
1:A:305:ARG:O	1:A:309:THR:OG1	2.28	0.51
1:B:242:MET:HA	1:B:242:MET:CE	2.41	0.51
1:B:26:ALA:CB	1:B:31:ILE:HD11	2.39	0.50
1:B:283:ASN:C	1:B:283:ASN:HD22	2.13	0.50
1:B:26:ALA:O	1:B:31:ILE:HD13	2.11	0.50
1:B:307:ALA:O	1:B:311:LYS:HG3	2.12	0.50
1:B:280:THR:CG2	1:B:287:VAL:HB	2.41	0.50
1:A:283:ASN:ND2	1:A:285:GLN:H	2.10	0.49
1:B:121:MET:HG3	1:B:146:PHE:CE2	2.46	0.49
1:A:219:HIS:CE1	1:A:223:LYS:HD2	2.47	0.49
1:A:242:MET:HA	1:A:242:MET:HE2	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:283:ASN:ND2	1:B:285:GLN:H	2.11	0.49
1:B:80:VAL:HG21	1:B:248:ILE:HD11	1.94	0.49
1:B:44:ARG:NH2	1:B:48:GLU:OE2	2.46	0.48
1:A:232:GLY:C	1:A:234:GLY:N	2.66	0.48
1:A:280:THR:CG2	1:A:287:VAL:HB	2.44	0.48
1:B:37:LEU:O	1:B:66:GLY:HA2	2.13	0.48
1:A:214:LYS:O	1:A:217:GLU:HB3	2.14	0.48
1:B:315:ALA:O	1:B:317:PHE:N	2.46	0.48
1:B:214:LYS:HA	1:B:214:LYS:HD3	1.75	0.47
1:B:29:ARG:HH21	1:B:249:GLU:HG3	1.78	0.47
1:B:74:ARG:HA	1:B:114:VAL:HG13	1.97	0.47
1:A:107:ILE:HG22	1:A:111:LEU:HD22	1.96	0.47
1:B:59:TYR:N	1:B:60:PRO:CD	2.78	0.47
1:A:29:ARG:HB3	1:A:31:ILE:HD13	1.96	0.46
1:A:9:LYS:NZ	1:A:74:ARG:O	2.43	0.46
1:B:227:TYR:CE1	1:B:231:ASN:ND2	2.83	0.46
1:B:315:ALA:C	1:B:317:PHE:N	2.68	0.46
1:A:252:LEU:HD23	1:A:252:LEU:HA	1.72	0.46
1:A:25:ALA:HB3	1:A:245:VAL:HG21	1.98	0.46
1:B:91:GLN:O	1:B:92:SER:O	2.34	0.46
1:A:305:ARG:HD3	1:A:305:ARG:HA	1.34	0.46
1:B:8:THR:HG23	1:B:32:ALA:HA	1.97	0.46
1:A:101:VAL:O	1:A:105:LYS:HG3	2.15	0.46
1:A:88:LYS:CG	1:A:89:PRO:CD	2.63	0.45
1:A:258:ILE:HA	1:A:280:THR:O	2.16	0.45
1:B:183:SER:HB2	3:B:395:HOH:O	2.16	0.45
1:A:243:SER:O	1:A:246:ASP:HB3	2.17	0.45
1:B:26:ALA:HA	1:B:31:ILE:CD1	2.47	0.45
1:B:36:VAL:HA	1:B:65:ASP:O	2.16	0.45
1:A:8:THR:HG23	1:A:33:ARG:N	2.28	0.45
1:A:267:ASP:OD1	1:A:270:GLY:CA	2.65	0.45
1:B:91:GLN:C	1:B:92:SER:O	2.53	0.45
1:A:274:ILE:HD12	1:A:274:ILE:HA	1.72	0.44
1:B:82:THR:O	2:B:320:NAD:H52N	2.17	0.44
1:B:242:MET:HA	1:B:242:MET:HE2	2.00	0.44
1:A:188:TRP:CZ2	1:A:215:ARG:HB3	2.52	0.44
1:A:267:ASP:OD1	1:A:270:GLY:HA2	2.17	0.44
1:B:125:ASN:ND2	1:B:126:PRO:HA	2.33	0.43
1:A:292:ASN:HD22	1:A:292:ASN:HA	1.59	0.43
1:B:26:ALA:HB1	1:B:31:ILE:CD1	2.40	0.43
1:B:274:ILE:HA	1:B:274:ILE:HD12	1.60	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:157:LEU:HD12	1:B:218:ILE:HG22	1.99	0.43
1:B:146:PHE:HB2	1:B:262:SER:O	2.19	0.43
1:A:176:ILE:HA	1:A:186:PRO:HA	2.01	0.43
1:B:189:GLU:HG3	1:B:296:SER:CB	2.41	0.43
1:A:108:MET:HB2	1:A:109:PRO:HD3	1.99	0.43
1:B:69:ASP:O	1:B:72:ILE:HG22	2.18	0.43
1:A:265:LEU:HA	1:A:265:LEU:HD12	1.77	0.43
1:B:93:ARG:O	1:B:97:VAL:HG23	2.18	0.42
1:B:93:ARG:O	1:B:93:ARG:HG2	2.19	0.42
1:A:265:LEU:HG	1:A:268:PHE:HB3	2.01	0.42
1:A:105:LYS:HD3	3:A:411:HOH:O	2.20	0.42
1:A:176:ILE:HD13	1:A:176:ILE:HG21	1.74	0.42
1:B:295:VAL:HG13	1:B:300:LEU:HB2	2.00	0.42
1:B:219:HIS:CE1	1:B:223:LYS:HD3	2.55	0.42
1:A:180:HIS:HE1	1:A:230:ILE:HD11	1.85	0.42
1:A:280:THR:HG21	1:A:287:VAL:HB	2.00	0.41
1:A:25:ALA:CB	1:A:245:VAL:HG21	2.50	0.41
1:B:283:ASN:HD22	1:B:285:GLN:H	1.68	0.41
1:B:44:ARG:CZ	1:B:48:GLU:OE2	2.67	0.41
1:A:204:LEU:HB3	1:A:205:PRO:HD2	2.02	0.41
1:B:288:ASN:HD22	1:B:290:THR:H	1.69	0.41
1:A:103:ILE:HD13	1:A:103:ILE:HG21	1.78	0.41
1:A:238:TYR:O	1:A:242:MET:HG2	2.20	0.41
1:A:107:ILE:CG2	1:A:111:LEU:HD22	2.50	0.41
1:B:37:LEU:O	1:B:66:GLY:CA	2.68	0.41
1:B:44:ARG:HE	1:B:48:GLU:CD	2.22	0.41
1:B:96:LEU:HD21	1:B:126:PRO:HG3	2.03	0.41
1:B:315:ALA:C	1:B:317:PHE:H	2.23	0.41
1:B:94:LEU:CD1	1:B:316:GLN:OE1	2.56	0.41
1:B:113:LYS:HB3	1:B:113:LYS:HE2	1.87	0.41
1:A:214:LYS:HA	1:A:214:LYS:HD3	1.73	0.40
1:B:140:LEU:HA	1:B:141:PRO:HD3	1.91	0.40
1:B:18:VAL:HG22	1:B:237:ASN:HB3	2.02	0.40
1:A:59:TYR:N	1:A:60:PRO:HD3	2.36	0.40
1:A:120:TYR:O	1:A:145:ILE:HA	2.22	0.40
1:A:189:GLU:CG	1:A:296:SER:HB3	2.41	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	311/319 (98%)	298 (96%)	12 (4%)	1 (0%)	41 37
1	B	311/319 (98%)	298 (96%)	9 (3%)	4 (1%)	12 6
All	All	622/638 (98%)	596 (96%)	21 (3%)	5 (1%)	19 13

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	92	SER
1	B	234	GLY
1	A	292	ASN
1	B	316	GLN
1	B	292	ASN

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	254/259 (98%)	231 (91%)	23 (9%)	9 5
1	B	254/259 (98%)	232 (91%)	22 (9%)	10 6
All	All	508/518 (98%)	463 (91%)	45 (9%)	9 6

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	LEU

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Mol	Chain	Res	Type
1	A	31	ILE
1	A	50	LEU
1	A	63	SER
1	A	74	ARG
1	A	111	LEU
1	A	114	VAL
1	A	122	LEU
1	A	140	LEU
1	A	164	GLN
1	A	211	ASP
1	A	217	GLU
1	A	230	ILE
1	A	233	LYS
1	A	237	ASN
1	A	261	VAL
1	A	265	LEU
1	A	283	ASN
1	A	288	ASN
1	A	292	ASN
1	A	298	LYS
1	A	305	ARG
1	A	309	THR
1	B	31	ILE
1	B	33	ARG
1	B	37	LEU
1	B	42	LYS
1	B	43	GLU
1	B	56	SER
1	B	74	ARG
1	B	94	LEU
1	B	96	LEU
1	B	107	ILE
1	B	111	LEU
1	B	164	GLN
1	B	223	LYS
1	B	242	MET
1	B	261	VAL
1	B	266	LYS
1	B	281	LEU
1	B	283	ASN
1	B	285	GLN
1	B	288	ASN

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Mol	Chain	Res	Type
1	B	295	VAL
1	B	319	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	102	ASN
1	A	164	GLN
1	A	173	HIS
1	A	220	GLN
1	A	224	ASN
1	A	231	ASN
1	A	237	ASN
1	A	256	ASN
1	A	283	ASN
1	A	288	ASN
1	A	292	ASN
1	B	87	GLN
1	B	102	ASN
1	B	125	ASN
1	B	164	GLN
1	B	173	HIS
1	B	220	GLN
1	B	224	ASN
1	B	231	ASN
1	B	237	ASN
1	B	256	ASN
1	B	283	ASN
1	B	288	ASN
1	B	292	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

2 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	B	320	-	42,48,48	1.56	5 (11%)	50,73,73	2.92	14 (28%)
2	NAD	A	320	-	42,48,48	1.48	3 (7%)	50,73,73	2.77	15 (30%)

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	B	320	-	-	6/26/62/62	0/5/5/5
2	NAD	A	320	-	-	6/26/62/62	0/5/5/5

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	320	NAD	O4B-C1B	-6.28	1.32	1.41
2	B	320	NAD	O4B-C1B	-5.81	1.33	1.41
2	A	320	NAD	O4D-C1D	-4.15	1.35	1.41
2	B	320	NAD	O4D-C1D	-3.90	1.35	1.41
2	B	320	NAD	C3N-C7N	3.09	1.55	1.50
2	B	320	NAD	C2N-N1N	2.98	1.38	1.35
2	A	320	NAD	C3N-C7N	2.49	1.54	1.50
2	B	320	NAD	C2A-N1A	2.28	1.38	1.33

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	320	NAD	C3N-C7N-N7N	-10.28	105.41	117.75
2	A	320	NAD	C3N-C7N-N7N	-8.69	107.32	117.75
2	A	320	NAD	C5N-C4N-C3N	-7.27	111.74	120.34
2	B	320	NAD	C5N-C4N-C3N	-7.17	111.86	120.34
2	B	320	NAD	O4B-C1B-C2B	-6.22	97.84	106.93
2	A	320	NAD	C6N-C5N-C4N	5.86	127.96	119.44
2	A	320	NAD	C4N-C3N-C7N	-5.74	105.69	121.04
2	B	320	NAD	O7N-C7N-N7N	5.63	130.57	122.58
2	B	320	NAD	C2N-C3N-C4N	5.38	124.35	118.26
2	B	320	NAD	C6N-C5N-C4N	5.32	127.17	119.44
2	A	320	NAD	C2N-C3N-C4N	5.17	124.12	118.26
2	A	320	NAD	C5N-C6N-N1N	-5.05	113.17	120.40
2	A	320	NAD	O4B-C1B-C2B	-4.91	99.75	106.93
2	B	320	NAD	C4N-C3N-C7N	-4.85	108.06	121.04
2	A	320	NAD	C6N-N1N-C2N	4.21	125.82	121.97
2	B	320	NAD	C2N-N1N-C1D	-4.10	110.00	119.14
2	A	320	NAD	O4D-C1D-C2D	-4.10	100.93	106.93
2	B	320	NAD	C5N-C6N-N1N	-3.96	114.73	120.40
2	A	320	NAD	O7N-C7N-C3N	3.69	124.05	119.63
2	A	320	NAD	C2N-C3N-C7N	3.69	130.16	119.46
2	B	320	NAD	O4D-C1D-C2D	-3.53	101.77	106.93
2	A	320	NAD	C3N-C2N-N1N	-3.51	117.00	120.43
2	B	320	NAD	C3N-C2N-N1N	-3.46	117.05	120.43
2	B	320	NAD	C6N-N1N-C2N	3.13	124.83	121.97
2	B	320	NAD	C2N-C3N-C7N	2.79	127.58	119.46
2	A	320	NAD	O7N-C7N-N7N	2.41	126.00	122.58
2	B	320	NAD	O7N-C7N-C3N	2.38	122.48	119.63
2	A	320	NAD	C2N-N1N-C1D	-2.09	114.49	119.14
2	A	320	NAD	O3B-C3B-C4B	2.02	116.90	111.05

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	320	NAD	O4D-C1D-N1N-C6N
2	A	320	NAD	C2D-C1D-N1N-C6N
2	B	320	NAD	O4D-C1D-N1N-C2N
2	B	320	NAD	O4D-C1D-N1N-C6N
2	B	320	NAD	C2D-C1D-N1N-C2N
2	B	320	NAD	C2D-C1D-N1N-C6N
2	B	320	NAD	O4B-C4B-C5B-O5B
2	A	320	NAD	PN-O3-PA-O1A
2	A	320	NAD	C2D-C1D-N1N-C2N

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Mol	Chain	Res	Type	Atoms
2	A	320	NAD	PN-O3-PA-O2A
2	B	320	NAD	C5D-O5D-PN-O1N
2	A	320	NAD	O4B-C4B-C5B-O5B

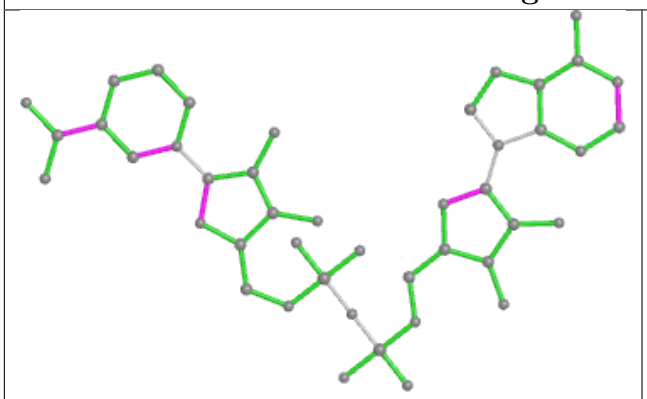
There are no ring outliers.

1 monomer is involved in 3 short contacts:

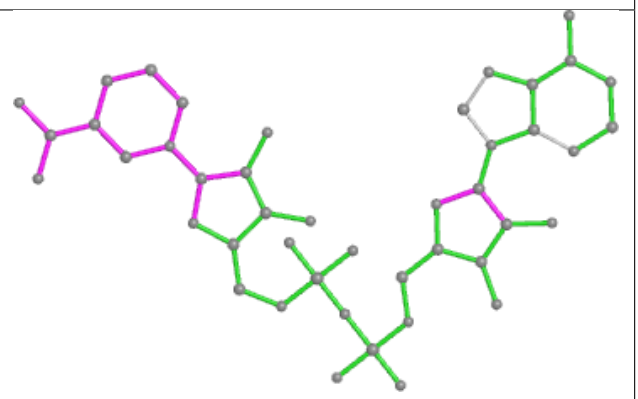
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	320	NAD	3	0

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 then 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.

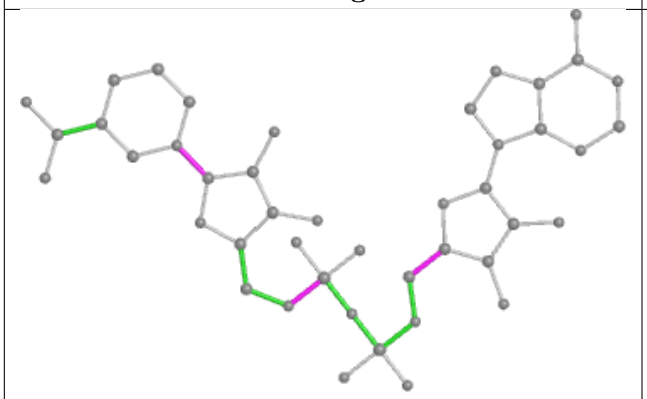
Ligand NAD B 320



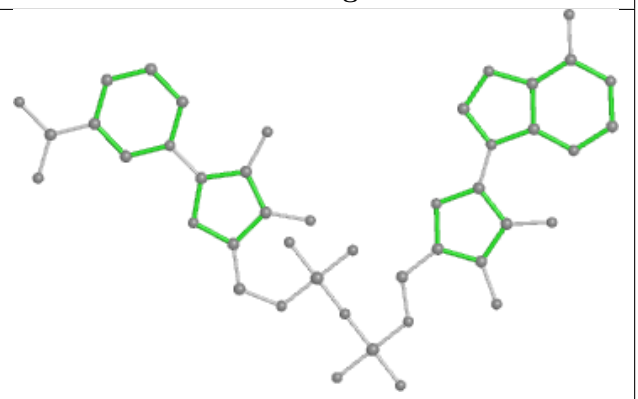
Bond lengths



Bond angles

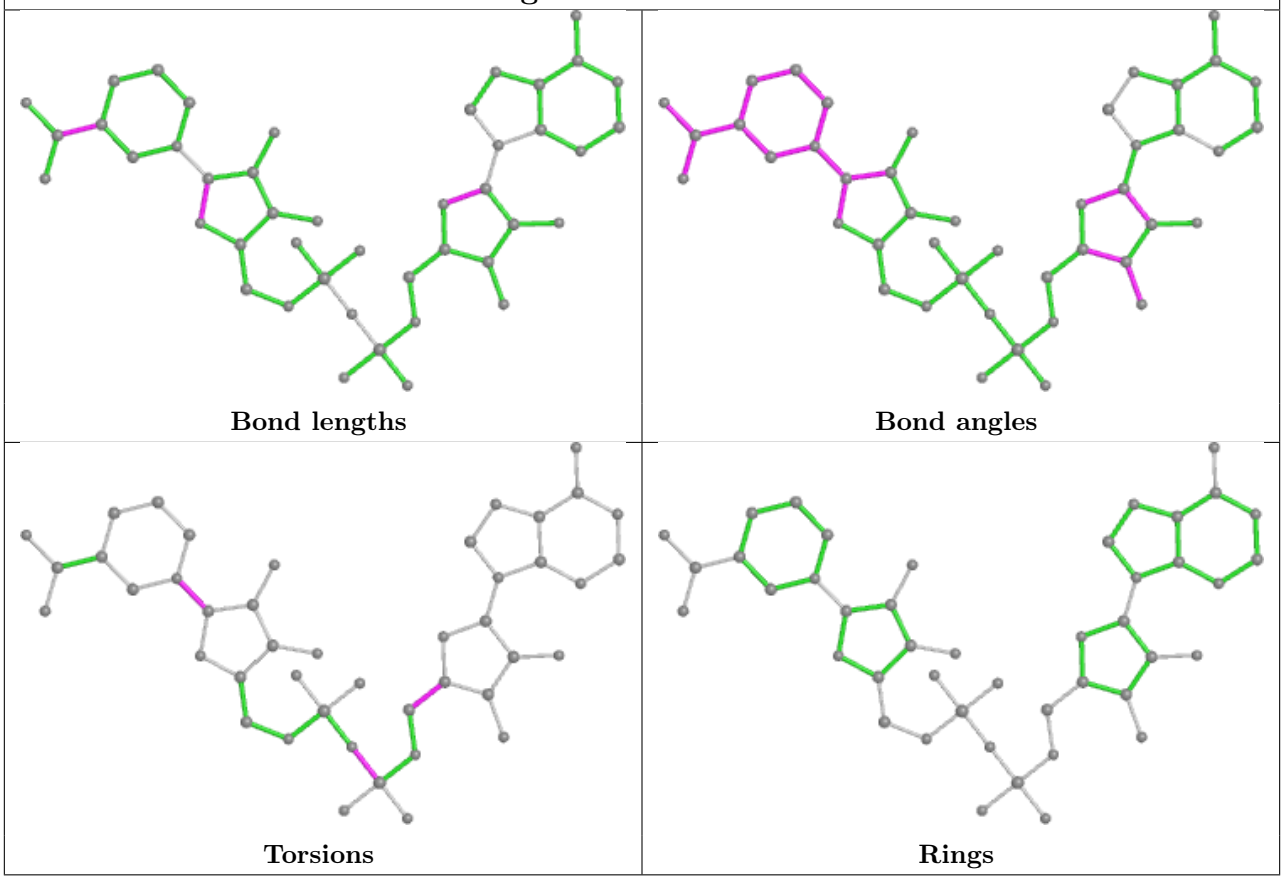


Torsions



Rings

Ligand NAD A 320



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	313/319 (98%)	-0.56	1 (0%)	94	93	9, 23, 44, 66	0
1	B	313/319 (98%)	-0.50	3 (0%)	82	81	8, 25, 51, 67	0
All	All	626/638 (98%)	-0.53	4 (0%)	89	88	8, 24, 48, 67	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	90	GLY	3.9
1	A	87	GLN	3.1
1	B	89	PRO	3.1
1	B	86	ARG	3.1

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 monosaccharides 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. 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	B-factors(Å^2)	Q<0.9
2	NAD	B	320	44/44	0.98	0.08	12,24,28,31	0

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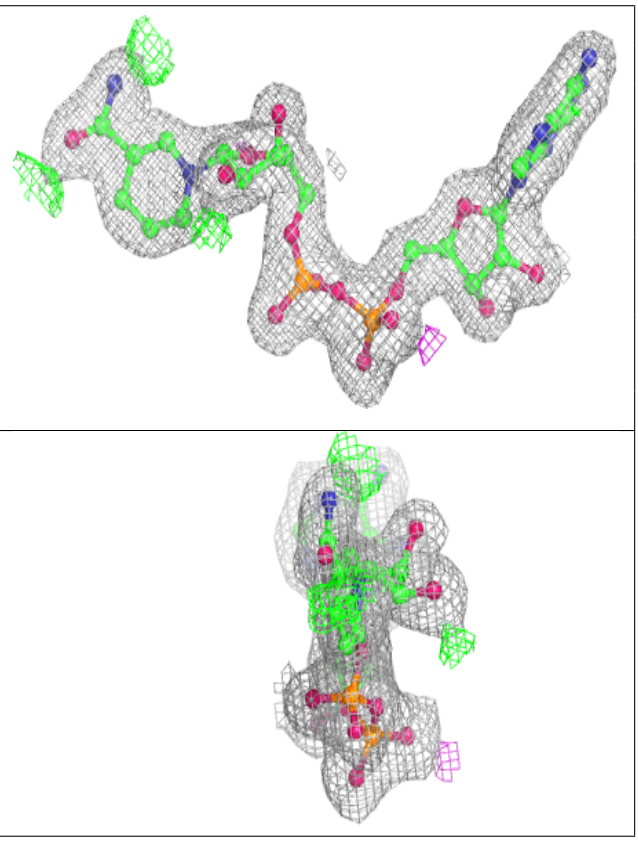
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	NAD	A	320	44/44	0.99	0.08	13,20,28,31	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

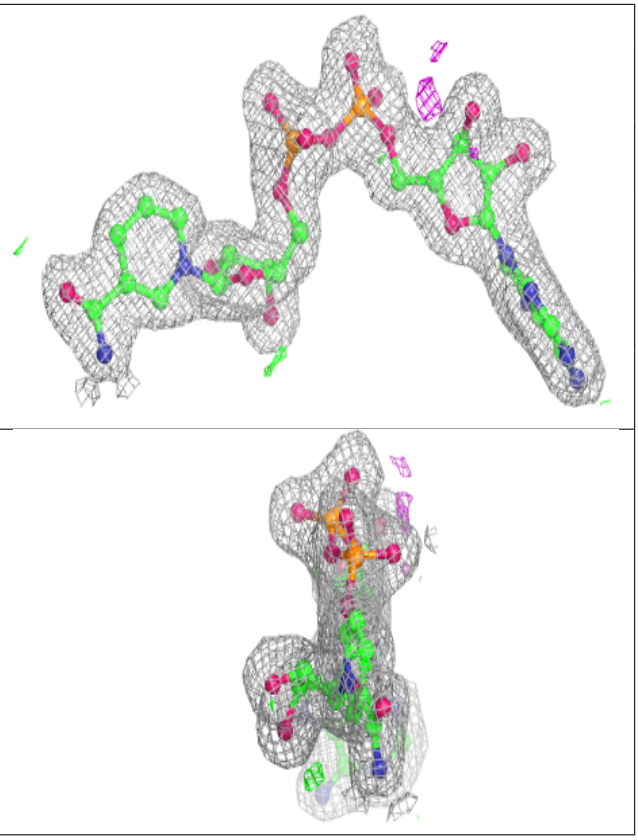
Electron density around NAD B 320:

2mF$_o$-DF$_c$ (at 0.7 rmsd) in gray
mF$_o$-DF$_c$ (at 3 rmsd) in purple (negative) and green (positive)



Electron density around NAD A 320:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers (i)

There are no such residues in this entry.