



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 07:13 AM UTC

PDB ID : 1BMD / pdb_00001bmd
Title : DETERMINANTS OF PROTEIN THERMOSTABILITY OBSERVED IN
THE 1.9 ANGSTROMS CRYSTAL STRUCTURE OF MALATE DEHY-
DROGENASE FROM THE THERMOPHILIC BACTERIUM THERMUS
FLAVUS
Authors : Kelly, C.A.; Birktoft, J.J.
Deposited on : 1992-11-10
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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

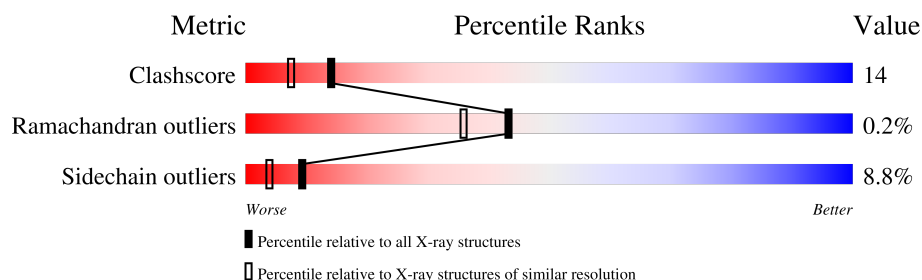
1 Overall quality at a glance

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	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (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 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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	327	 70% 21% 7% .
1	B	327	 76% 19% . .

2 Entry composition i

There are 3 unique types of molecules in this entry. The entry contains 5366 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.

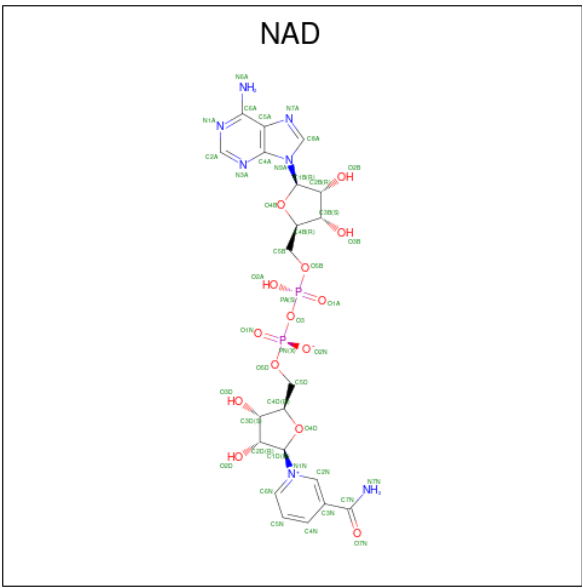
- Molecule 1 is a protein called MALATE DEHYDROGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	327	Total	C	N	O	S	0	1	0
			2493	1577	436	467	13			
1	B	327	Total	C	N	O	S	0	2	0
			2497	1579	436	469	13			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	74	ASP	LYS	conflict	UNP P10584
B	74	ASP	LYS	conflict	UNP P10584

- Molecule 2 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (CCD ID: NAD) (formula: C₂₁H₂₇N₇O₁₄P₂).



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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		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	125	Total	O	0	0
			125	125		
3	B	163	Total	O	0	0
			163	163		

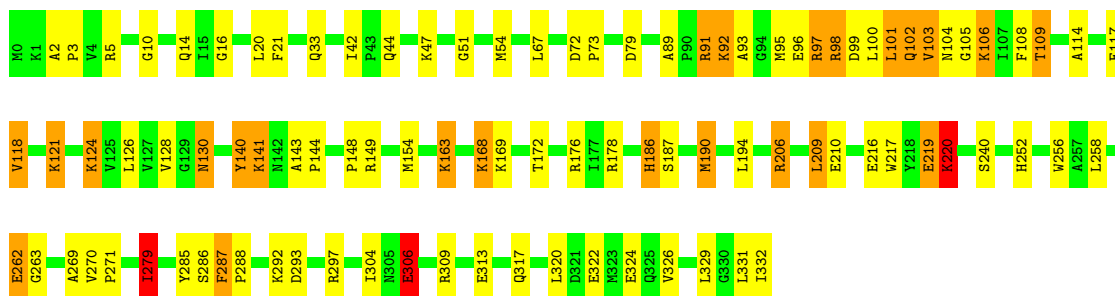
3 Residue-property plots

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. 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.

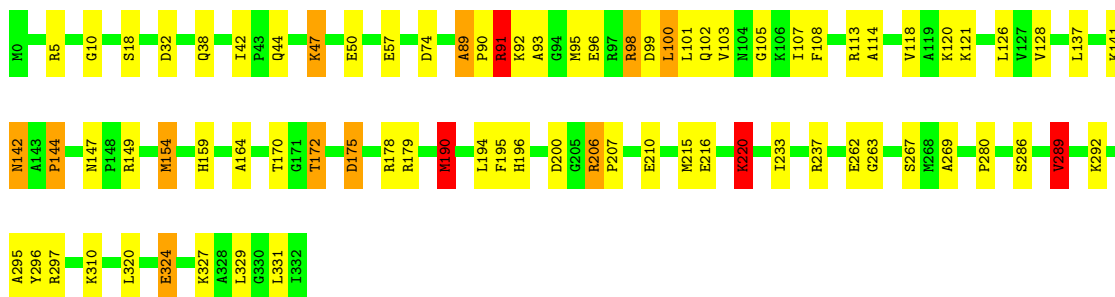
• Molecule 1: MALATE DEHYDROGENASE

Chain A: 



• Molecule 1: MALATE DEHYDROGENASE

Chain B: 



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	71.51Å 87.59Å 118.98Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	(Not available) – 1.90	Depositor
% Data completeness (in resolution range)	(Not available) ((Not available)-1.90)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	TNT, X-PLOR	Depositor
R, R_{free}	0.154 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5366	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: 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	1.23	0/2531	1.80	39/3427 (1.1%)
1	B	1.25	2/2531 (0.1%)	1.79	33/3426 (1.0%)
All	All	1.24	2/5062 (0.0%)	1.80	72/6853 (1.1%)

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

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	200	ASP	CG-OD1	5.20	1.35	1.25
1	B	262	GLU	CD-OE2	5.11	1.35	1.25

All (72) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	96	GLU	N-CA-CB	-13.35	91.36	110.44
1	B	142	ASN	CA-CB-CG	-8.75	103.85	112.60
1	A	279	ILE	CB-CA-C	8.43	118.78	110.94
1	B	175	ASP	CA-CB-CG	-8.41	104.19	112.60
1	B	220	LYS	N-CA-CB	7.72	122.48	110.44
1	A	72	ASP	CB-CA-C	-7.50	101.52	110.15
1	A	130	ASN	CA-CB-CG	-7.31	105.29	112.60
1	B	89	ALA	CA-C-N	7.24	127.67	119.92
1	B	89	ALA	C-N-CA	7.24	127.67	119.92

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	172	THR	CA-CB-OG1	-7.05	99.03	109.60
1	B	200	ASP	CB-CA-C	-7.01	101.93	111.89
1	A	144	PRO	CA-C-N	-6.83	112.58	123.31
1	A	144	PRO	C-N-CA	-6.83	112.58	123.31
1	B	289	VAL	N-CA-CB	-6.77	99.26	111.92
1	A	21	PHE	CA-CB-CG	-6.71	107.09	113.80
1	A	262	GLU	CB-CG-CD	6.69	123.97	112.60
1	B	91	ARG	N-CA-CB	6.68	120.14	110.06
1	A	79	ASP	CA-CB-CG	-6.66	105.94	112.60
1	A	73	PRO	N-CA-CB	6.52	110.43	103.39
1	B	331	LEU	N-CA-C	6.42	119.26	111.82
1	B	99	ASP	CA-C-O	6.25	127.04	120.42
1	A	293	ASP	CA-CB-CG	6.21	118.81	112.60
1	A	5	ARG	NE-CZ-NH2	6.14	124.73	119.20
1	A	297	ARG	NE-CZ-NH1	-6.10	115.40	121.50
1	B	154	MET	N-CA-C	6.08	118.64	108.96
1	A	210	GLU	CA-CB-CG	-6.05	102.01	114.10
1	A	33	GLN	CA-C-O	6.00	125.26	119.62
1	B	144	PRO	CA-C-N	-5.99	111.20	122.05
1	B	144	PRO	C-N-CA	-5.99	111.20	122.05
1	B	105	GLY	CA-C-O	5.99	126.87	120.40
1	A	206	ARG	N-CA-CB	-5.97	102.81	110.03
1	A	270	VAL	N-CA-CB	-5.96	106.60	111.67
1	B	147	ASN	CA-CB-CG	5.95	118.55	112.60
1	B	195	PHE	CA-CB-CG	-5.94	107.86	113.80
1	B	233	ILE	CB-CA-C	-5.89	104.30	112.02
1	A	176	ARG	CA-C-N	-5.83	115.59	122.93
1	A	176	ARG	C-N-CA	-5.83	115.59	122.93
1	A	240	SER	CA-CB-OG	-5.81	99.48	111.10
1	B	215	MET	CB-CA-C	-5.67	98.45	110.31
1	A	324	GLU	N-CA-CB	5.64	118.18	110.01
1	B	237	ARG	CB-CA-C	-5.62	99.87	109.65
1	A	220	LYS	CA-C-N	-5.54	117.28	123.16
1	A	220	LYS	C-N-CA	-5.54	117.28	123.16
1	A	306	GLU	CB-CG-CD	5.51	121.97	112.60
1	A	118	VAL	N-CA-C	5.51	118.44	113.10
1	B	74	ASP	N-CA-CB	5.49	119.50	110.39
1	B	164	ALA	CA-C-N	-5.37	112.66	120.29
1	B	164	ALA	C-N-CA	-5.37	112.66	120.29
1	B	108	PHE	CA-CB-CG	-5.36	108.44	113.80
1	A	258	LEU	N-CA-CB	5.36	118.50	110.14
1	A	33	GLN	CB-CG-CD	-5.31	103.57	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	190	MET	CA-CB-CG	-5.28	103.53	114.10
1	B	107	ILE	CB-CA-C	-5.27	105.11	112.02
1	A	287	PHE	CB-CA-C	-5.26	101.62	109.56
1	A	108	PHE	CA-CB-CG	-5.25	108.55	113.80
1	B	32	ASP	CB-CA-C	-5.22	101.90	110.72
1	A	91	ARG	CA-C-N	-5.21	114.62	122.81
1	A	91	ARG	C-N-CA	-5.21	114.62	122.81
1	A	67	LEU	N-CA-CB	-5.18	102.12	110.71
1	B	100	LEU	CA-C-N	5.17	127.52	120.54
1	B	100	LEU	C-N-CA	5.17	127.52	120.54
1	A	143	ALA	O-C-N	5.17	125.94	121.34
1	A	117	GLU	CA-C-N	-5.13	115.59	122.72
1	A	117	GLU	C-N-CA	-5.13	115.59	122.72
1	B	91	ARG	NE-CZ-NH1	5.13	126.63	121.50
1	A	271	PRO	N-CA-CB	5.08	107.69	103.17
1	A	322	GLU	N-CA-CB	5.06	117.56	110.12
1	A	96	GLU	N-CA-CB	-5.06	102.62	110.46
1	B	262	GLU	CA-C-N	-5.06	113.54	122.09
1	B	262	GLU	C-N-CA	-5.06	113.54	122.09
1	B	324	GLU	CB-CG-CD	5.04	121.17	112.60
1	A	216	GLU	CA-CB-CG	-5.01	104.08	114.10

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	186	HIS	Mainchain

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	2493	0	2520	80	0
1	B	2497	0	2520	65	0
2	A	44	0	26	6	0
2	B	44	0	26	3	0
3	A	125	0	0	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	163	0	0	10	0
All	All	5366	0	5092	145	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 14.

All (145) 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:98:ARG:H	1:A:98:ARG:HD2	1.20	1.05
1:A:89:ALA:H	1:A:104:ASN:HD21	1.13	0.95
1:A:105:GLY:O	1:A:109:THR:HG23	1.74	0.88
1:B:47:LYS:O	1:B:47:LYS:HE2	1.74	0.87
1:B:92:LYS:HG2	1:B:93:ALA:N	1.92	0.83
1:A:306:GLU:HA	1:A:309:ARG:HD2	1.60	0.83
1:A:98:ARG:O	1:A:102:GLN:HG2	1.80	0.81
1:A:121:LYS:N	1:A:121:LYS:HD2	1.95	0.79
1:B:295:ALA:HA	3:B:494:HOH:O	1.83	0.79
1:A:89:ALA:H	1:A:104:ASN:ND2	1.81	0.78
1:A:97:ARG:HB3	1:A:98:ARG:CZ	2.15	0.76
1:B:126:LEU:HD11	1:B:154:MET:HB2	1.66	0.76
1:B:5:ARG:HD3	1:B:38:GLN:HE22	1.48	0.76
1:B:92:LYS:HE2	1:B:93:ALA:O	1.86	0.75
1:A:98:ARG:H	1:A:98:ARG:CD	1.99	0.74
1:A:92:LYS:HG2	1:A:95:MET:HE3	1.70	0.73
1:B:92:LYS:HB3	1:B:95:MET:SD	2.28	0.73
1:B:206:ARG:HD2	1:B:210:GLU:OE1	1.87	0.73
1:B:190:MET:HE2	3:B:474:HOH:O	1.88	0.73
1:B:47:LYS:HE3	1:B:50:GLU:CD	2.14	0.72
1:B:190:MET:O	1:B:190:MET:HG3	1.87	0.72
1:A:326:VAL:HG13	1:A:331:LEU:HB2	1.71	0.71
1:A:89:ALA:HB2	1:A:103:VAL:HG22	1.71	0.71
1:A:140:TYR:CE1	1:A:141:LYS:HD2	2.25	0.70
1:B:10:GLY:HA2	2:B:334:NAD:H1B	1.74	0.70
1:A:92:LYS:HG2	1:A:95:MET:CE	2.21	0.70
1:B:89:ALA:HB2	1:B:103:VAL:HG12	1.73	0.70
1:A:92:LYS:HG2	1:A:95:MET:HB2	1.73	0.68
2:B:334:NAD:H52N	3:B:491:HOH:O	1.92	0.68
1:B:47:LYS:HE2	1:B:47:LYS:C	2.18	0.68
1:B:5:ARG:HD3	1:B:38:GLN:NE2	2.09	0.68
1:A:98:ARG:HD2	1:A:98:ARG:N	2.02	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:92:LYS:HG2	1:A:95:MET:SD	2.35	0.66
1:A:14:GLN:NE2	2:A:334:NAD:O2A	2.28	0.66
1:B:92:LYS:O	1:B:95:MET:HB3	1.94	0.66
1:B:159:HIS:HD2	1:B:267:SER:OG	1.79	0.66
1:A:92:LYS:O	1:A:95:MET:HB3	1.95	0.66
1:B:170:THR:OG1	1:B:172:THR:HG23	1.95	0.66
1:B:128:VAL:O	2:B:334:NAD:H2N	1.96	0.65
1:A:306:GLU:OE1	1:A:306:GLU:N	2.32	0.63
1:A:306:GLU:HA	1:A:309:ARG:CD	2.28	0.63
1:A:92:LYS:HG2	1:A:95:MET:CB	2.28	0.63
1:B:91:ARG:HG3	1:B:91:ARG:HH11	1.63	0.62
1:B:102:GLN:HG2	1:B:329:LEU:HD13	1.80	0.62
1:B:137:LEU:CD1	1:B:141:LYS:HD2	2.32	0.60
1:A:269:ALA:HA	1:A:286:SER:HA	1.84	0.59
1:B:269:ALA:HA	1:B:286:SER:HA	1.84	0.59
1:B:280:PRO:HG2	1:B:320:LEU:HD11	1.84	0.59
1:B:47:LYS:NZ	1:B:50:GLU:OE1	2.33	0.59
1:B:142:ASN:O	1:B:144:PRO:HD3	2.03	0.59
1:A:187:SER:O	1:A:190:MET:HG2	2.02	0.58
1:A:102:GLN:HE22	1:A:329:LEU:HD21	1.67	0.58
1:A:304:ILE:HB	1:A:309:ARG:HH21	1.69	0.58
1:A:99:ASP:O	1:A:103:VAL:HG12	2.04	0.58
1:A:169:LYS:HZ2	1:A:217:TRP:CD1	2.22	0.57
1:B:47:LYS:HE2	1:B:47:LYS:CA	2.34	0.57
1:A:304:ILE:O	1:A:309:ARG:NH2	2.39	0.55
1:A:93:ALA:C	1:A:95:MET:H	2.13	0.55
1:A:102:GLN:OE1	1:A:329:LEU:HD22	2.05	0.55
1:A:89:ALA:CB	1:A:103:VAL:HG22	2.38	0.54
1:B:206:ARG:HG3	1:B:207:PRO:N	2.23	0.54
1:B:289:VAL:HG13	1:B:296:TYR:HB2	1.90	0.54
1:B:47:LYS:HA	1:B:47:LYS:CE	2.38	0.54
1:A:168:LYS:NZ	1:B:57:GLU:OE2	2.41	0.53
1:A:130:ASN:ND2	2:A:334:NAD:O2D	2.37	0.53
1:B:149:ARG:HD2	3:B:450:HOH:O	2.07	0.53
1:A:98:ARG:HA	1:A:101:LEU:HB2	1.90	0.53
1:A:169:LYS:NZ	1:A:217:TRP:CD1	2.78	0.52
1:A:102:GLN:HE22	1:A:329:LEU:CD2	2.23	0.52
1:B:91:ARG:HG3	1:B:91:ARG:NH1	2.23	0.51
1:A:219:GLU:HG2	1:A:220:LYS:HE3	1.92	0.51
1:A:92:LYS:CG	1:A:95:MET:HE3	2.38	0.50
1:A:130:ASN:HB2	2:A:334:NAD:O2D	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:18:SER:O	3:B:425:HOH:O	2.20	0.50
1:B:170:THR:O	1:B:172:THR:HG22	2.11	0.50
1:B:216:GLU:O	1:B:220:LYS:HD2	2.12	0.50
1:B:178:ARG:HH22	1:B:263:GLY:HA3	1.77	0.50
1:B:47:LYS:CA	1:B:47:LYS:CE	2.89	0.50
1:A:92:LYS:CD	1:A:95:MET:HE3	2.42	0.50
1:B:142:ASN:C	1:B:144:PRO:HD3	2.37	0.49
1:A:2:ALA:HB1	1:A:3:PRO:HD2	1.95	0.49
1:A:16:GLY:O	1:A:20:LEU:HG	2.12	0.49
1:A:306:GLU:HA	1:A:309:ARG:CG	2.43	0.49
1:B:196:HIS:HD2	3:B:471:HOH:O	1.96	0.49
1:A:320:LEU:HA	1:A:320:LEU:HD23	1.49	0.48
1:A:126:LEU:HD11	1:A:154:MET:HB2	1.95	0.48
1:B:91:ARG:HH11	1:B:91:ARG:CG	2.24	0.48
1:A:97:ARG:HB3	1:A:98:ARG:NH1	2.28	0.48
1:A:10:GLY:HA2	2:A:334:NAD:H1B	1.95	0.47
1:A:190:MET:HE3	1:A:190:MET:HB2	1.49	0.47
1:B:113:ARG:NH1	1:B:144:PRO:HG3	2.29	0.47
1:A:92:LYS:HD3	1:A:95:MET:HE3	1.96	0.47
1:A:306:GLU:HA	1:A:309:ARG:HG2	1.97	0.47
1:B:149:ARG:HD3	1:B:296:TYR:OH	2.15	0.47
1:A:326:VAL:HG12	1:A:332:ILE:HG12	1.97	0.47
1:A:100:LEU:HA	1:A:100:LEU:HD12	1.73	0.46
1:B:5:ARG:HH11	1:B:38:GLN:NE2	2.14	0.46
1:A:252:HIS:HD2	3:A:354:HOH:O	1.97	0.46
1:A:178:ARG:HH22	1:A:263:GLY:HA3	1.80	0.46
1:A:163:LYS:NZ	3:A:370:HOH:O	2.30	0.45
1:A:92:LYS:HD3	1:A:95:MET:CE	2.47	0.45
1:A:106:LYS:N	1:A:106:LYS:HD2	2.28	0.45
1:B:89:ALA:HB1	1:B:90:PRO:HD2	1.99	0.45
1:B:159:HIS:CD2	1:B:267:SER:OG	2.66	0.45
1:A:97:ARG:O	1:A:100:LEU:HB3	2.17	0.45
1:B:159:HIS:HE1	3:B:354:HOH:O	2.00	0.44
1:B:89:ALA:CB	1:B:90:PRO:HD2	2.42	0.44
1:B:92:LYS:HG2	1:B:93:ALA:H	1.78	0.44
1:B:137:LEU:HD12	1:B:141:LYS:HD2	1.96	0.44
1:A:121:LYS:HD2	1:A:121:LYS:H	1.77	0.44
1:A:287:PHE:HB3	1:A:288:PRO:HD2	1.99	0.44
1:A:42:ILE:HD13	1:A:42:ILE:HG21	1.65	0.43
1:A:178:ARG:NH2	1:A:263:GLY:HA3	2.33	0.43
1:A:100:LEU:HD23	3:A:436:HOH:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:93:ALA:C	1:A:95:MET:N	2.76	0.43
1:A:101:LEU:HD12	1:A:101:LEU:HA	1.82	0.43
1:A:140:TYR:CD1	1:A:148:PRO:HB3	2.53	0.43
1:A:309:ARG:O	1:A:313:GLU:HG2	2.18	0.43
1:B:89:ALA:HB2	1:B:103:VAL:CG1	2.45	0.42
1:B:114:ALA:O	1:B:118:VAL:HG22	2.20	0.42
1:A:92:LYS:CG	1:A:95:MET:CE	2.95	0.42
1:A:190:MET:HE2	3:A:420:HOH:O	2.20	0.42
1:B:175:ASP:OD1	1:B:175:ASP:N	2.42	0.42
1:A:128:VAL:O	2:A:334:NAD:H2N	2.19	0.42
1:B:98:ARG:CZ	1:B:98:ARG:HB3	2.49	0.42
1:A:124:LYS:HB3	1:A:256:TRP:CH2	2.54	0.42
1:B:196:HIS:CD2	3:B:471:HOH:O	2.72	0.42
1:A:279:ILE:HD13	1:A:285:TYR:CE1	2.54	0.42
1:A:209:LEU:HD12	1:A:209:LEU:HA	1.88	0.41
1:B:137:LEU:HD11	1:B:141:LYS:HD2	2.00	0.41
1:B:178:ARG:NH2	1:B:263:GLY:HA3	2.35	0.41
1:A:326:VAL:CG1	1:A:331:LEU:HB2	2.45	0.41
1:B:42:ILE:HG21	1:B:42:ILE:HD13	1.80	0.41
1:B:194:LEU:HA	1:B:194:LEU:HD23	1.87	0.41
1:A:51:GLY:HA2	1:A:54:MET:HE3	2.02	0.41
1:A:306:GLU:H	1:A:306:GLU:CD	2.25	0.41
1:B:100:LEU:HD12	1:B:100:LEU:HA	1.79	0.41
1:B:141:LYS:HZ2	1:B:141:LYS:HG2	1.80	0.41
1:B:179:ARG:HD3	3:B:451:HOH:O	2.21	0.41
1:A:114:ALA:O	1:A:118:VAL:HG22	2.21	0.40
1:A:186:HIS:CD2	2:A:334:NAD:C7N	3.05	0.40
1:B:47:LYS:HE2	1:B:47:LYS:HA	2.02	0.40
1:A:97:ARG:O	1:A:100:LEU:N	2.54	0.40
1:B:120:LYS:HE3	3:B:487:HOH:O	2.20	0.40
1:B:216:GLU:OE2	1:B:220:LYS:HE2	2.22	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	326/327 (100%)	321 (98%)	4 (1%)	1 (0%)	36	29
1	B	327/327 (100%)	325 (99%)	2 (1%)	0	100	100
All	All	653/654 (100%)	646 (99%)	6 (1%)	1 (0%)	43	36

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	97	ARG

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	251/252 (100%)	223 (89%)	28 (11%)	6	2
1	B	250/252 (99%)	234 (94%)	16 (6%)	16	8
All	All	501/504 (99%)	457 (91%)	44 (9%)	9	4

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	44	GLN
1	A	47	LYS
1	A	91	ARG
1	A	92	LYS
1	A	98	ARG
1	A	101	LEU
1	A	102	GLN
1	A	103	VAL
1	A	106	LYS
1	A	109	THR
1	A	121	LYS
1	A	124	LYS

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Mol	Chain	Res	Type
1	A	140	TYR
1	A	141	LYS
1	A	149	ARG
1	A	163	LYS
1	A	168	LYS
1	A	190	MET
1	A	194	LEU
1	A	206	ARG
1	A	209	LEU
1	A	219	GLU
1	A	220	LYS
1	A	262	GLU
1	A	279	ILE
1	A	292	LYS
1	A	306	GLU
1	A	317	GLN
1	B	44	GLN
1	B	47	LYS
1	B	91	ARG
1	B	98	ARG
1	B	101	LEU
1	B	121	LYS
1	B	172	THR
1	B	190	MET
1	B	206	ARG
1	B	220	LYS
1	B	289	VAL
1	B	292	LYS
1	B	297	ARG
1	B	310	LYS
1	B	324	GLU
1	B	327	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	104	ASN
1	A	165	GLN
1	A	252	HIS
1	A	273	GLN
1	A	317	GLN
1	A	325	GLN

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Mol	Chain	Res	Type
1	B	38	GLN
1	B	44	GLN
1	B	102	GLN
1	B	147	ASN
1	B	159	HIS
1	B	273	GLN
1	B	325	GLN

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 oligosaccharides 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	A	334	-	46,48,48	1.30	5 (10%)	64,73,73	1.58	11 (17%)
2	NAD	B	334	-	46,48,48	1.23	5 (10%)	64,73,73	1.54	10 (15%)

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	334	-	-	2/30/62/62	0/5/5/5
2	NAD	B	334	-	-	4/30/62/62	0/5/5/5

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	334	NAD	C4A-N9A	4.37	1.46	1.37
2	A	334	NAD	PN-O3	4.06	1.63	1.59
2	A	334	NAD	C4A-N9A	3.68	1.45	1.37
2	B	334	NAD	PN-O3	-3.27	1.56	1.59
2	B	334	NAD	PA-O3	-3.26	1.56	1.59
2	A	334	NAD	C2N-N1N	3.01	1.38	1.35
2	A	334	NAD	PA-O3	-2.99	1.56	1.59
2	A	334	NAD	C5A-N7A	2.58	1.43	1.39
2	B	334	NAD	C2N-N1N	2.29	1.37	1.35
2	B	334	NAD	O4D-C1D	2.03	1.43	1.40

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	334	NAD	O7N-C7N-C3N	-5.83	112.47	119.60
2	A	334	NAD	O7N-C7N-C3N	-5.60	112.75	119.60
2	A	334	NAD	C4D-O4D-C1D	-5.56	104.83	109.92
2	B	334	NAD	O4B-C1B-C2B	-3.46	99.21	106.62
2	B	334	NAD	C4D-O4D-C1D	-3.38	106.83	109.92
2	A	334	NAD	O7N-C7N-N7N	3.24	127.30	122.62
2	B	334	NAD	C5A-C4A-N9A	-3.01	102.53	105.81
2	B	334	NAD	C3N-C7N-N7N	2.97	121.39	117.74
2	B	334	NAD	C4B-O4B-C1B	-2.91	103.04	109.47
2	A	334	NAD	C6A-C5A-C4A	-2.80	113.36	117.18
2	A	334	NAD	O3D-C3D-C2D	-2.69	103.19	111.82
2	A	334	NAD	C5A-C4A-N3A	2.54	130.21	126.72
2	B	334	NAD	C4A-C5A-N7A	2.47	113.41	110.58
2	B	334	NAD	O7N-C7N-N7N	2.43	126.12	122.62
2	A	334	NAD	C4A-N9A-C1B	-2.32	121.21	126.63
2	A	334	NAD	C1B-N9A-C8A	2.27	132.12	127.09
2	B	334	NAD	O3-PA-O1A	-2.24	103.96	110.70
2	B	334	NAD	O2A-PA-O1A	2.17	122.55	112.44
2	A	334	NAD	O4D-C4D-C3D	-2.12	100.95	105.15
2	A	334	NAD	C5A-C4A-N9A	-2.06	103.56	105.81
2	A	334	NAD	C2D-C3D-C4D	2.00	106.48	102.61

There are no chirality outliers.

All (6) torsion outliers are listed below:

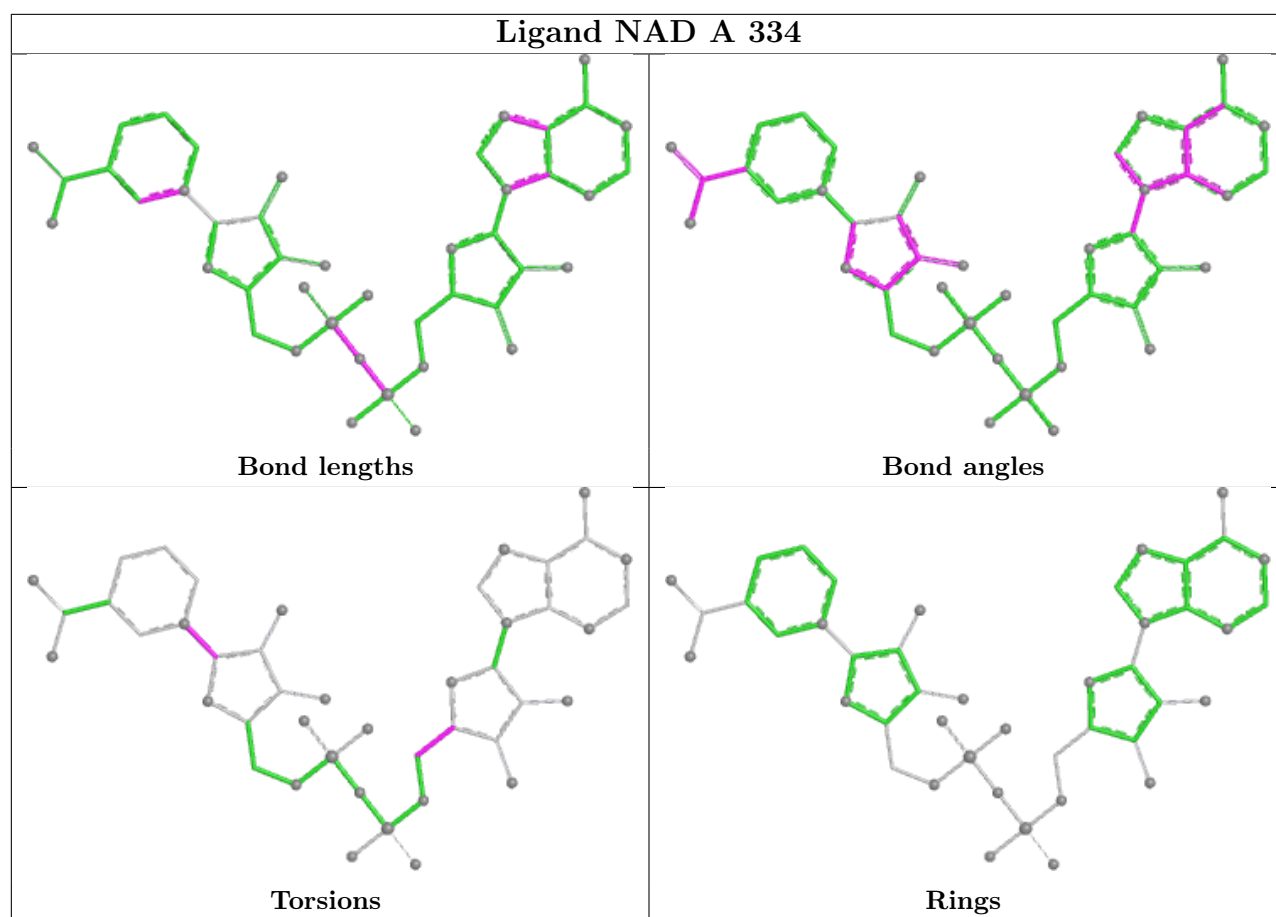
Mol	Chain	Res	Type	Atoms
2	B	334	NAD	O4D-C1D-N1N-C2N
2	B	334	NAD	O4D-C1D-N1N-C6N
2	B	334	NAD	C2D-C1D-N1N-C6N
2	B	334	NAD	C2D-C1D-N1N-C2N
2	A	334	NAD	O4D-C1D-N1N-C2N
2	A	334	NAD	O4B-C4B-C5B-O5B

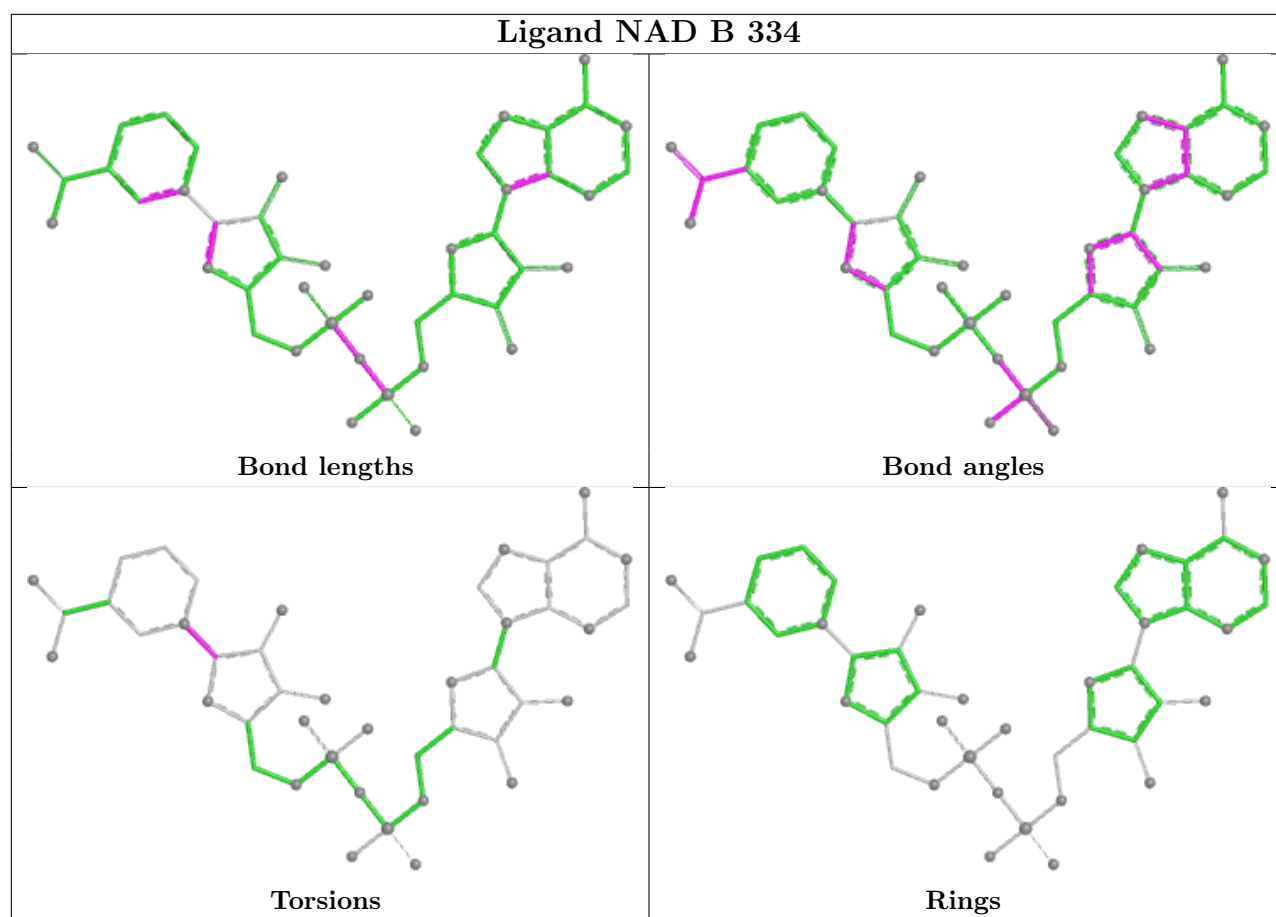
There are no ring outliers.

2 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	334	NAD	6	0
2	B	334	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 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.