



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 12:14 AM UTC

PDB ID : 3PQF / pdb_00003pqf
Title : Crystal structure of L-lactate dehydrogenase from *Bacillus subtilis* mutation H171C complexed with NAD⁺
Authors : Zhang, Y.; Garavito, R.M.
Deposited on : 2010-11-26
Resolution : 2.49 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)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

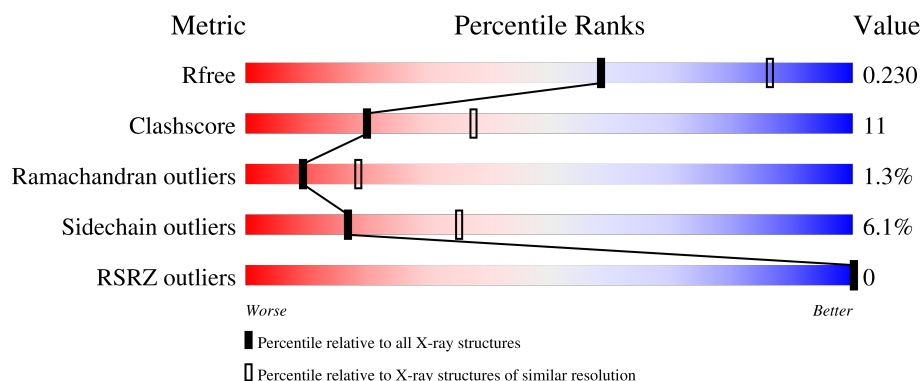
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.49 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))
R_{free}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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 $\leq 5\%$. 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	326	
1	B	326	
1	C	326	
1	D	326	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9867 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
1	A	314	Total	C	N	O	S	0	0	0
			2397	1527	401	460	9			
1	B	314	Total	C	N	O	S	0	0	0
			2397	1527	401	460	9			
1	C	312	Total	C	N	O	S	0	0	0
			2384	1520	398	457	9			
1	D	312	Total	C	N	O	S	0	0	0
			2384	1520	398	457	9			

There are 28 discrepancies between the modelled and reference sequences:

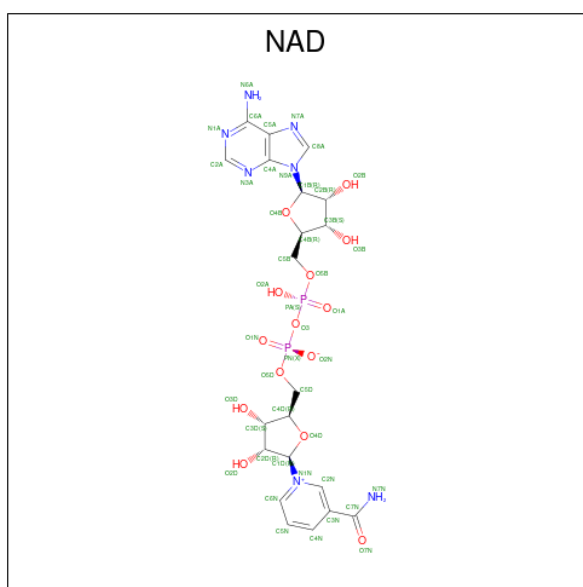
Chain	Residue	Modelled	Actual	Comment	Reference
A	171	CYS	HIS	engineered mutation	UNP P13714
A	321	HIS	-	expression tag	UNP P13714
A	322	HIS	-	expression tag	UNP P13714
A	323	HIS	-	expression tag	UNP P13714
A	324	HIS	-	expression tag	UNP P13714
A	325	HIS	-	expression tag	UNP P13714
A	326	HIS	-	expression tag	UNP P13714
B	171	CYS	HIS	engineered mutation	UNP P13714
B	321	HIS	-	expression tag	UNP P13714
B	322	HIS	-	expression tag	UNP P13714
B	323	HIS	-	expression tag	UNP P13714
B	324	HIS	-	expression tag	UNP P13714
B	325	HIS	-	expression tag	UNP P13714
B	326	HIS	-	expression tag	UNP P13714
C	171	CYS	HIS	engineered mutation	UNP P13714
C	321	HIS	-	expression tag	UNP P13714
C	322	HIS	-	expression tag	UNP P13714
C	323	HIS	-	expression tag	UNP P13714
C	324	HIS	-	expression tag	UNP P13714
C	325	HIS	-	expression tag	UNP P13714
C	326	HIS	-	expression tag	UNP P13714

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Chain	Residue	Modelled	Actual	Comment	Reference
D	171	CYS	HIS	engineered mutation	UNP P13714
D	321	HIS	-	expression tag	UNP P13714
D	322	HIS	-	expression tag	UNP P13714
D	323	HIS	-	expression tag	UNP P13714
D	324	HIS	-	expression tag	UNP P13714
D	325	HIS	-	expression tag	UNP P13714
D	326	HIS	-	expression tag	UNP P13714

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



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

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	82	Total O 82 82	0	0
3	B	94	Total O 94 94	0	0
3	C	17	Total O 17 17	0	0

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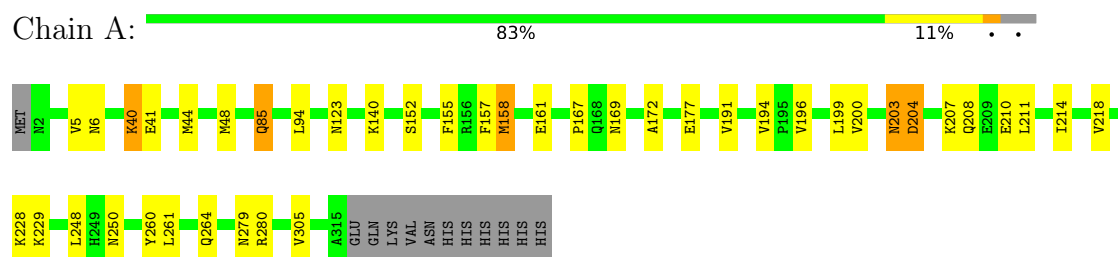
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	24	Total	O	0	0
			24	24		

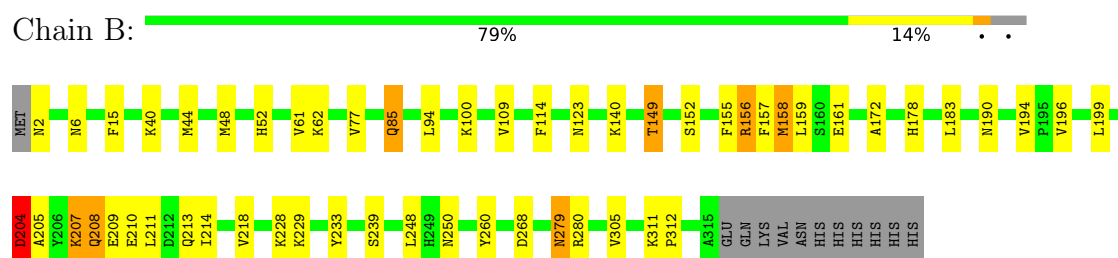
3 Residue-property plots

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 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.

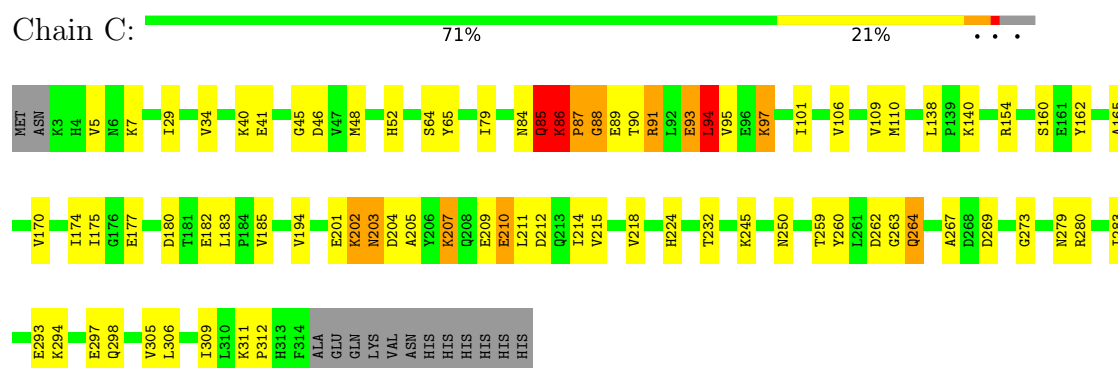
- Molecule 1: L-lactate dehydrogenase



- Molecule 1: L-lactate dehydrogenase

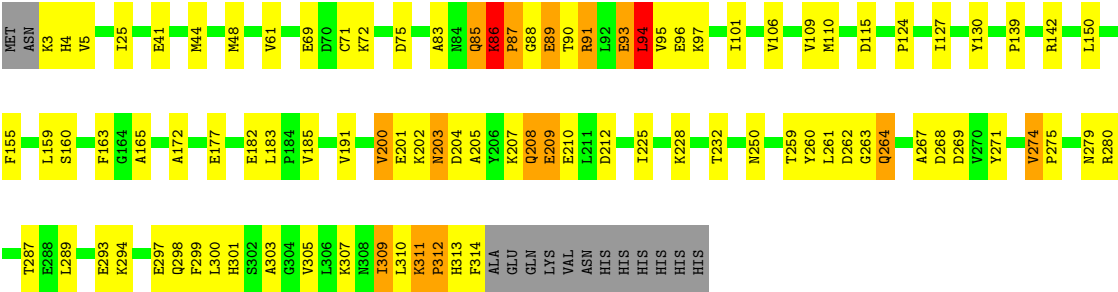


- Molecule 1: L-lactate dehydrogenase



- Molecule 1: L-lactate dehydrogenase





4 Data and refinement statistics

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, α , β , γ	133.43Å 133.43Å 99.09Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.49 – 2.49 30.49 – 2.49	Depositor EDS
% Data completeness (in resolution range)	99.8 (30.49-2.49) 99.8 (30.49-2.49)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.24 (at 2.48Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6_289)	Depositor
R, R_{free}	0.189 , 0.235 0.185 , 0.230	Depositor DCC
R_{free} test set	2062 reflections (2.99%)	wwPDB-VP
Wilson B-factor (Å ²)	48.6	Xtriage
Anisotropy	0.086	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 27.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.478 for -h,-k,l 0.030 for h,-h-k,-l 0.031 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	9867	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.36% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

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	0.63	0/2443	0.87	2/3311 (0.1%)
1	B	0.62	0/2443	0.86	3/3311 (0.1%)
1	C	0.49	0/2430	0.83	4/3293 (0.1%)
1	D	0.48	0/2430	0.85	3/3293 (0.1%)
All	All	0.56	0/9746	0.85	12/13208 (0.1%)

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	B	0	1

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	86	LYS	CA-C-N	7.60	129.34	119.84
1	D	86	LYS	C-N-CA	7.60	129.34	119.84
1	D	205	ALA	N-CA-C	-6.48	105.34	113.18
1	C	205	ALA	N-CA-C	-5.78	106.07	113.01
1	C	194	VAL	N-CA-C	5.63	113.44	107.76
1	A	5	VAL	N-CA-C	5.33	116.34	108.45
1	C	85	GLN	N-CA-C	-5.31	99.50	110.80
1	B	123	ASN	N-CA-C	5.26	117.93	110.24
1	A	123	ASN	N-CA-C	5.10	117.68	110.24
1	B	205	ALA	N-CA-C	-5.02	107.13	113.20
1	C	79	ILE	N-CA-C	5.01	114.79	107.37
1	B	279	ASN	N-CA-C	5.00	114.97	108.07

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	204	ASP	Peptide

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	2397	0	2384	24	0
1	B	2397	0	2384	34	0
1	C	2384	0	2373	85	0
1	D	2384	0	2373	85	0
2	C	44	0	26	1	0
2	D	44	0	26	1	0
3	A	82	0	0	0	0
3	B	94	0	0	4	0
3	C	17	0	0	0	0
3	D	24	0	0	1	0
All	All	9867	0	9566	219	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 11.

All (219) 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:C:85:GLN:HG3	1:C:87:PRO:CD	1.43	1.44
1:C:85:GLN:CG	1:C:87:PRO:HD3	1.62	1.29
1:C:86:LYS:O	1:C:88:GLY:N	1.67	1.26
1:C:311:LYS:HG3	1:C:312:PRO:HD3	1.25	1.08
1:D:250:ASN:HD22	1:D:280:ARG:H	1.11	0.99
1:C:85:GLN:CG	1:C:87:PRO:CD	2.30	0.97
1:C:204:ASP:O	1:C:207:LYS:HE3	1.73	0.88
1:C:85:GLN:HG3	1:C:87:PRO:HD2	1.53	0.87
1:C:84:ASN:C	1:C:85:GLN:O	2.12	0.87
1:B:52:HIS:HD2	1:C:154:ARG:HD2	1.39	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:228:LYS:HG3	1:C:48:MET:HE1	1.59	0.85
1:B:250:ASN:HD22	1:B:280:ARG:H	1.26	0.84
1:C:250:ASN:HD22	1:C:280:ARG:H	1.23	0.83
1:A:250:ASN:HD22	1:A:280:ARG:H	1.23	0.82
1:D:86:LYS:O	1:D:88:GLY:N	2.12	0.82
1:A:250:ASN:ND2	1:A:279:ASN:HB2	1.93	0.82
1:C:84:ASN:O	1:C:85:GLN:O	1.98	0.81
1:C:85:GLN:HG3	1:C:87:PRO:HD3	0.81	0.80
1:D:93:GLU:HG3	1:D:94:LEU:H	1.45	0.79
1:D:207:LYS:HB2	1:D:210:GLU:HG3	1.65	0.79
1:C:87:PRO:O	1:C:88:GLY:O	2.03	0.77
1:C:86:LYS:N	1:C:86:LYS:HD2	2.01	0.76
1:D:93:GLU:C	1:D:95:VAL:H	1.93	0.75
1:B:85:GLN:HB2	1:B:94:LEU:HD22	1.71	0.72
1:B:250:ASN:ND2	1:B:279:ASN:HB2	2.05	0.71
1:C:93:GLU:C	1:C:95:VAL:H	1.99	0.71
1:D:303:ALA:O	1:D:307:LYS:HG3	1.91	0.70
1:B:159:LEU:N	3:B:417:HOH:O	2.23	0.69
1:B:52:HIS:CD2	1:C:154:ARG:HD2	2.25	0.69
1:D:263:GLY:N	1:D:268:ASP:HB3	2.07	0.69
1:D:85:GLN:HB2	1:D:87:PRO:HD3	1.75	0.68
1:C:85:GLN:CD	1:C:87:PRO:HD3	2.18	0.68
1:A:204:ASP:O	1:A:207:LYS:HE2	1.94	0.67
1:D:250:ASN:ND2	1:D:280:ARG:H	1.88	0.67
1:A:250:ASN:HD22	1:A:279:ASN:HB2	1.57	0.67
1:C:250:ASN:ND2	1:C:279:ASN:HB2	2.09	0.67
1:C:259:THR:HG23	1:C:283:ILE:HD13	1.77	0.67
1:D:311:LYS:CG	1:D:312:PRO:HD3	2.24	0.66
1:D:232:THR:H	2:D:352:NAD:H72N	1.42	0.66
1:C:86:LYS:N	1:C:87:PRO:CD	2.57	0.66
1:D:204:ASP:O	1:D:207:LYS:HD3	1.94	0.66
1:D:263:GLY:H	1:D:268:ASP:HB3	1.61	0.65
1:C:250:ASN:ND2	1:C:280:ARG:H	1.93	0.65
1:C:263:GLY:HA2	1:C:267:ALA:O	1.97	0.65
1:D:311:LYS:HG3	1:D:312:PRO:HD3	1.77	0.65
1:C:86:LYS:C	1:C:88:GLY:N	2.53	0.65
1:C:86:LYS:NZ	1:C:97:LYS:HZ1	1.96	0.64
1:B:250:ASN:ND2	1:B:280:ARG:H	1.95	0.64
1:D:86:LYS:C	1:D:88:GLY:H	2.05	0.63
1:B:140:LYS:HD2	1:B:260:TYR:CD2	2.33	0.63
1:D:250:ASN:ND2	1:D:279:ASN:HB2	2.14	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:85:GLN:CG	1:C:87:PRO:HD2	2.18	0.63
1:B:156:ARG:C	3:B:417:HOH:O	2.41	0.63
1:A:250:ASN:ND2	1:A:280:ARG:H	1.92	0.62
1:C:232:THR:H	2:C:352:NAD:H72N	1.47	0.61
1:B:213:GLN:HA	1:B:213:GLN:OE1	2.01	0.60
1:C:170:VAL:HG12	1:C:170:VAL:O	2.01	0.60
1:D:90:THR:O	1:D:91:ARG:C	2.44	0.60
1:A:210:GLU:O	1:A:214:ILE:HG13	2.01	0.60
1:D:301:HIS:O	1:D:305:VAL:HG23	2.02	0.60
1:D:274:VAL:HG23	1:D:275:PRO:HD2	1.83	0.60
1:C:140:LYS:HD2	1:C:260:TYR:CD2	2.36	0.60
1:B:214:ILE:O	1:B:218:VAL:HG23	2.02	0.59
1:B:250:ASN:HD22	1:B:279:ASN:HB2	1.66	0.59
1:C:86:LYS:O	1:C:87:PRO:C	2.44	0.59
1:D:260:TYR:OH	1:D:269:ASP:HA	2.03	0.58
1:C:203:ASN:HD22	1:C:204:ASP:H	1.51	0.58
1:C:86:LYS:NZ	1:C:97:LYS:NZ	2.51	0.58
1:A:85:GLN:HB2	1:A:94:LEU:HD22	1.86	0.58
1:D:208:GLN:H	1:D:208:GLN:NE2	2.02	0.58
1:B:2:ASN:HA	3:B:371:HOH:O	2.04	0.58
1:D:311:LYS:O	1:D:313:HIS:N	2.37	0.58
1:D:201:GLU:HG3	1:D:202:LYS:N	2.19	0.57
1:B:207:LYS:HG2	1:B:210:GLU:HG3	1.86	0.57
1:C:40:LYS:HG2	1:C:65:TYR:OH	2.04	0.57
1:D:177:GLU:O	1:D:182:GLU:HB3	2.05	0.57
1:C:85:GLN:OE1	1:C:87:PRO:HG3	2.05	0.57
1:B:210:GLU:O	1:B:214:ILE:HG13	2.05	0.56
1:D:93:GLU:C	1:D:95:VAL:N	2.62	0.56
1:D:89:GLU:HB3	1:D:91:ARG:HG3	1.87	0.56
1:C:93:GLU:C	1:C:95:VAL:N	2.63	0.56
1:C:86:LYS:N	1:C:87:PRO:HD2	2.21	0.55
1:C:85:GLN:HG3	1:C:86:LYS:N	2.22	0.55
1:A:228:LYS:HD2	1:D:48:MET:HE1	1.89	0.55
1:C:89:GLU:HB3	1:C:91:ARG:CG	2.37	0.54
1:A:140:LYS:HD2	1:A:260:TYR:CD2	2.43	0.54
1:D:86:LYS:HE2	1:D:94:LEU:HD12	1.90	0.54
1:D:94:LEU:O	1:D:94:LEU:HG	2.05	0.54
1:C:311:LYS:HG3	1:C:312:PRO:CD	2.18	0.53
1:D:41:GLU:HA	1:D:41:GLU:OE2	2.07	0.53
1:C:89:GLU:HB3	1:C:91:ARG:HG3	1.89	0.53
1:D:307:LYS:HA	1:D:310:LEU:HD12	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:44:MET:O	1:A:48:MET:HG3	2.08	0.53
1:C:86:LYS:HZ3	1:C:97:LYS:NZ	2.07	0.52
1:C:263:GLY:CA	1:C:267:ALA:O	2.57	0.52
1:C:85:GLN:CG	1:C:86:LYS:N	2.73	0.52
1:D:209:GLU:OE1	1:D:212:ASP:HB3	2.09	0.52
1:D:311:LYS:O	1:D:312:PRO:C	2.53	0.52
1:C:260:TYR:CZ	1:C:269:ASP:HA	2.44	0.52
1:C:260:TYR:OH	1:C:269:ASP:HA	2.09	0.52
1:D:93:GLU:HG3	1:D:94:LEU:N	2.19	0.52
1:D:261:LEU:HD22	1:D:264:GLN:HG3	1.91	0.52
1:C:207:LYS:N	1:C:207:LYS:HD2	2.25	0.51
1:D:155:PHE:O	1:D:159:LEU:HG	2.10	0.51
1:D:311:LYS:O	1:D:314:PHE:N	2.44	0.51
1:C:97:LYS:O	1:C:101:ILE:HG13	2.11	0.51
1:C:209:GLU:HA	1:C:212:ASP:HB2	1.92	0.51
1:D:88:GLY:O	1:D:90:THR:N	2.43	0.51
1:C:86:LYS:HD2	1:C:86:LYS:H	1.74	0.51
1:C:86:LYS:HZ3	1:C:97:LYS:HZ1	1.57	0.50
1:D:263:GLY:HA2	1:D:267:ALA:O	2.11	0.50
1:B:208:GLN:O	1:B:211:LEU:HB2	2.11	0.50
1:B:15:PHE:HB3	1:B:233:TYR:CD1	2.48	0.49
1:C:84:ASN:O	1:C:85:GLN:C	2.54	0.49
1:B:48:MET:HE1	1:C:224:HIS:HB3	1.95	0.49
1:D:260:TYR:CZ	1:D:269:ASP:HA	2.48	0.49
1:D:200:VAL:HG13	1:D:208:GLN:OE1	2.11	0.49
1:D:293:GLU:O	1:D:297:GLU:HG3	2.12	0.49
1:C:311:LYS:CG	1:C:312:PRO:HD3	2.17	0.49
1:C:262:ASP:H	1:C:264:GLN:CG	2.26	0.48
1:C:293:GLU:O	1:C:297:GLU:HG3	2.13	0.48
1:B:44:MET:O	1:B:48:MET:HG3	2.14	0.48
1:D:259:THR:O	1:D:271:TYR:HA	2.13	0.48
1:B:149:THR:HG23	1:B:239:SER:HB2	1.95	0.48
1:D:139:PRO:HG2	1:D:142:ARG:HD3	1.96	0.47
1:C:106:VAL:O	1:C:110:MET:HG2	2.14	0.47
1:A:228:LYS:HD2	1:D:48:MET:CE	2.44	0.47
1:D:185:VAL:HG11	1:D:289:LEU:HD13	1.96	0.47
1:C:94:LEU:O	1:C:94:LEU:HG	2.14	0.47
1:D:311:LYS:C	1:D:313:HIS:N	2.72	0.47
1:A:208:GLN:O	1:A:211:LEU:HB2	2.14	0.47
1:D:124:PRO:HG2	1:D:127:ILE:HB	1.95	0.47
1:D:264:GLN:H	1:D:264:GLN:HG2	1.43	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:150:LEU:HB3	3:D:348:HOH:O	2.15	0.47
1:A:157:PHE:CD2	1:A:157:PHE:C	2.92	0.47
1:D:86:LYS:HE2	1:D:97:LYS:NZ	2.29	0.47
1:A:214:ILE:O	1:A:218:VAL:HG23	2.15	0.46
1:C:87:PRO:C	1:C:88:GLY:O	2.58	0.46
1:D:44:MET:O	1:D:48:MET:HG3	2.14	0.46
1:D:83:ALA:HB1	1:D:97:LYS:HE2	1.97	0.46
1:D:163:PHE:CG	1:D:191:VAL:HG21	2.50	0.46
1:A:158:MET:HE2	1:A:214:ILE:HG23	1.96	0.46
1:C:29:ILE:HD13	1:C:245:LYS:HB2	1.98	0.46
1:C:91:ARG:C	1:C:93:GLU:N	2.70	0.46
1:D:208:GLN:H	1:D:208:GLN:CD	2.24	0.46
1:C:89:GLU:O	1:C:91:ARG:HG3	2.16	0.45
1:C:250:ASN:HD22	1:C:279:ASN:HB2	1.80	0.45
1:D:263:GLY:CA	1:D:267:ALA:O	2.65	0.45
1:D:91:ARG:C	1:D:93:GLU:N	2.71	0.45
1:A:6:ASN:HB3	1:A:248:LEU:HD22	1.98	0.45
1:C:177:GLU:O	1:C:182:GLU:HB3	2.16	0.45
1:C:294:LYS:HE2	1:C:298:GLN:HE22	1.81	0.45
1:D:25:ILE:HG12	1:D:61:VAL:HG11	1.97	0.45
1:C:201:GLU:C	1:C:203:ASN:H	2.25	0.45
1:D:4:HIS:HE1	1:D:75:ASP:OD2	2.00	0.45
1:D:262:ASP:HB2	1:D:264:GLN:OE1	2.17	0.45
1:B:77:VAL:HG23	1:B:114:PHE:CE1	2.52	0.44
1:B:194:VAL:HG12	1:B:199:LEU:HG	1.99	0.44
1:C:90:THR:O	1:C:91:ARG:C	2.60	0.44
1:C:162:TYR:HE1	1:C:210:GLU:HG2	1.82	0.44
1:B:155:PHE:CE2	1:B:172:ALA:HB1	2.52	0.44
1:D:86:LYS:NZ	1:D:91:ARG:HD3	2.31	0.44
1:C:160:SER:HB2	1:C:165:ALA:O	2.17	0.44
1:D:155:PHE:CE2	1:D:172:ALA:HB1	2.53	0.44
1:C:185:VAL:O	1:C:185:VAL:HG12	2.17	0.44
1:C:7:LYS:HE3	1:C:34:VAL:CG2	2.48	0.44
1:C:110:MET:HE1	1:C:138:LEU:HD21	1.98	0.44
1:C:94:LEU:O	1:C:97:LYS:HB2	2.18	0.43
1:B:204:ASP:O	1:B:207:LYS:HE3	2.18	0.43
1:A:261:LEU:HD22	1:A:264:GLN:HG2	2.00	0.43
1:B:228:LYS:HD3	1:C:41:GLU:OE1	2.19	0.43
1:B:6:ASN:HB3	1:B:248:LEU:HD22	2.01	0.43
1:B:157:PHE:CD2	1:B:157:PHE:C	2.96	0.43
1:D:294:LYS:HE2	1:D:298:GLN:HE22	1.84	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:41:GLU:OE2	1:D:228:LYS:HE3	2.18	0.43
1:D:97:LYS:O	1:D:101:ILE:HG13	2.19	0.43
1:D:208:GLN:CD	1:D:208:GLN:N	2.77	0.43
1:D:311:LYS:N	1:D:312:PRO:HD2	2.33	0.43
1:A:40:LYS:HE2	1:A:41:GLU:OE2	2.18	0.43
1:D:201:GLU:C	1:D:203:ASN:H	2.27	0.43
1:C:262:ASP:H	1:C:264:GLN:HG3	1.84	0.43
1:B:156:ARG:O	1:B:156:ARG:HG2	2.19	0.42
1:C:177:GLU:HG3	1:C:306:LEU:HD21	2.01	0.42
1:B:311:LYS:HB3	1:B:312:PRO:HD3	2.01	0.42
1:D:160:SER:HB2	1:D:165:ALA:O	2.19	0.42
1:D:263:GLY:N	1:D:267:ALA:O	2.52	0.42
1:B:158:MET:O	1:B:161:GLU:HG2	2.19	0.42
1:D:299:PHE:HD2	1:D:300:LEU:HD12	1.84	0.42
1:D:309:ILE:H	1:D:309:ILE:HG13	1.57	0.42
1:D:4:HIS:CE1	1:D:75:ASP:OD2	2.72	0.42
1:A:200:VAL:O	1:A:203:ASN:O	2.37	0.42
1:C:88:GLY:O	1:C:89:GLU:C	2.60	0.42
1:C:97:LYS:HZ3	1:C:97:LYS:HG3	1.46	0.42
1:D:71:CYS:O	1:D:72:LYS:C	2.62	0.42
1:D:212:ASP:OD1	1:D:212:ASP:C	2.63	0.42
1:A:191:VAL:HB	1:A:199:LEU:HD11	2.01	0.42
1:D:163:PHE:CD2	1:D:191:VAL:HG21	2.55	0.41
1:D:311:LYS:CG	1:D:312:PRO:CD	2.95	0.41
3:B:389:HOH:O	1:D:3:LYS:HE2	2.20	0.41
1:D:106:VAL:O	1:D:110:MET:HG2	2.21	0.41
1:B:229:LYS:HB2	1:C:45:GLY:HA3	2.03	0.41
1:C:89:GLU:CB	1:C:91:ARG:HG3	2.50	0.41
1:A:167:PRO:C	1:A:169:ASN:H	2.27	0.41
1:A:194:VAL:HG12	1:A:199:LEU:HG	2.01	0.41
1:B:40:LYS:HE3	1:B:40:LYS:HB3	1.85	0.41
1:C:214:ILE:O	1:C:218:VAL:HG23	2.21	0.41
1:D:93:GLU:O	1:D:95:VAL:N	2.52	0.41
1:C:85:GLN:OE1	1:C:87:PRO:HD3	2.19	0.41
1:D:115:ASP:O	1:D:142:ARG:NH2	2.53	0.41
1:D:225:ILE:CG2	1:D:232:THR:HG22	2.51	0.41
1:C:305:VAL:O	1:C:309:ILE:HG13	2.21	0.40
1:C:175:ILE:HB	1:C:273:GLY:O	2.21	0.40
1:C:202:LYS:HE3	1:C:202:LYS:HB2	1.97	0.40
1:D:94:LEU:C	1:D:96:GLU:N	2.78	0.40
1:D:130:TYR:HA	1:D:271:TYR:CD2	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:185:VAL:HG12	1:D:185:VAL:O	2.21	0.40
1:A:155:PHE:CE2	1:A:172:ALA:HB1	2.56	0.40
1:B:158:MET:HE2	1:B:214:ILE:HG23	2.04	0.40
1:C:86:LYS:H	1:C:87:PRO:CD	2.34	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	312/326 (96%)	297 (95%)	15 (5%)	0	100	100
1	B	312/326 (96%)	296 (95%)	15 (5%)	1 (0%)	36	55
1	C	310/326 (95%)	286 (92%)	16 (5%)	8 (3%)	4	7
1	D	310/326 (95%)	289 (93%)	14 (4%)	7 (2%)	5	8
All	All	1244/1304 (95%)	1168 (94%)	60 (5%)	16 (1%)	9	18

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	85	GLN
1	C	86	LYS
1	C	87	PRO
1	C	88	GLY
1	C	93	GLU
1	D	87	PRO
1	C	91	ARG
1	D	89	GLU
1	D	91	ARG
1	D	93	GLU
1	C	94	LEU

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Mol	Chain	Res	Type
1	C	202	LYS
1	D	94	LEU
1	B	178	HIS
1	D	86	LYS
1	D	312	PRO

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	255/267 (96%)	244 (96%)	11 (4%)	26	51
1	B	255/267 (96%)	237 (93%)	18 (7%)	13	29
1	C	254/267 (95%)	237 (93%)	17 (7%)	15	31
1	D	254/267 (95%)	238 (94%)	16 (6%)	16	34
All	All	1018/1068 (95%)	956 (94%)	62 (6%)	17	35

All (62) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	40	LYS
1	A	85	GLN
1	A	152	SER
1	A	158	MET
1	A	161	GLU
1	A	177	GLU
1	A	196	VAL
1	A	203	ASN
1	A	204	ASP
1	A	229	LYS
1	A	305	VAL
1	B	61	VAL
1	B	62	LYS
1	B	85	GLN
1	B	100	LYS
1	B	109	VAL

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Mol	Chain	Res	Type
1	B	149	THR
1	B	152	SER
1	B	156	ARG
1	B	158	MET
1	B	183	LEU
1	B	190	ASN
1	B	196	VAL
1	B	204	ASP
1	B	207	LYS
1	B	208	GLN
1	B	209	GLU
1	B	268	ASP
1	B	305	VAL
1	C	5	VAL
1	C	46	ASP
1	C	52	HIS
1	C	64	SER
1	C	86	LYS
1	C	94	LEU
1	C	97	LYS
1	C	109	VAL
1	C	174	ILE
1	C	180	ASP
1	C	183	LEU
1	C	203	ASN
1	C	207	LYS
1	C	210	GLU
1	C	211	LEU
1	C	215	VAL
1	C	264	GLN
1	D	5	VAL
1	D	69	GLU
1	D	85	GLN
1	D	86	LYS
1	D	94	LEU
1	D	109	VAL
1	D	183	LEU
1	D	200	VAL
1	D	203	ASN
1	D	208	GLN
1	D	209	GLU
1	D	264	GLN

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Mol	Chain	Res	Type
1	D	274	VAL
1	D	287	THR
1	D	309	ILE
1	D	311	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	51	ASN
1	A	123	ASN
1	A	208	GLN
1	A	220	ASN
1	A	250	ASN
1	B	4	HIS
1	B	52	HIS
1	B	85	GLN
1	B	123	ASN
1	B	190	ASN
1	B	208	GLN
1	B	250	ASN
1	C	59	GLN
1	C	123	ASN
1	C	169	ASN
1	C	178	HIS
1	C	188	HIS
1	C	203	ASN
1	C	250	ASN
1	C	264	GLN
1	C	301	HIS
1	D	4	HIS
1	D	52	HIS
1	D	59	GLN
1	D	123	ASN
1	D	178	HIS
1	D	250	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

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	C	352	-	46,48,48	1.79	5 (10%)	64,73,73	1.56	8 (12%)
2	NAD	D	352	-	46,48,48	1.80	5 (10%)	64,73,73	1.56	8 (12%)

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	C	352	-	-	2/30/62/62	0/5/5/5
2	NAD	D	352	-	-	2/30/62/62	0/5/5/5

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	352	NAD	O7N-C7N	9.49	1.41	1.24
2	C	352	NAD	O7N-C7N	9.46	1.41	1.24
2	D	352	NAD	C2A-N1A	2.95	1.39	1.33
2	C	352	NAD	C2A-N1A	2.93	1.39	1.33
2	C	352	NAD	C2A-N3A	2.80	1.38	1.33
2	D	352	NAD	C2A-N3A	2.65	1.38	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	352	NAD	C2N-N1N	2.35	1.37	1.35
2	C	352	NAD	C2N-N1N	2.30	1.37	1.35
2	C	352	NAD	C8A-N7A	2.27	1.36	1.31
2	D	352	NAD	C8A-N7A	2.19	1.35	1.31

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	352	NAD	N3A-C2A-N1A	-6.08	119.38	128.58
2	D	352	NAD	N3A-C2A-N1A	-5.96	119.56	128.58
2	C	352	NAD	C5A-C4A-N3A	-4.21	120.92	126.72
2	D	352	NAD	C5A-C4A-N3A	-4.15	121.01	126.72
2	C	352	NAD	N9A-C8A-N7A	-4.13	108.08	113.94
2	D	352	NAD	N9A-C8A-N7A	-4.04	108.21	113.94
2	C	352	NAD	C2A-N3A-C4A	3.54	120.48	111.83
2	D	352	NAD	C2A-N3A-C4A	3.43	120.21	111.83
2	C	352	NAD	C5A-N7A-C8A	3.36	108.74	103.45
2	D	352	NAD	C5A-N7A-C8A	3.28	108.61	103.45
2	D	352	NAD	N3A-C4A-N9A	3.21	132.62	127.17
2	C	352	NAD	N3A-C4A-N9A	3.12	132.47	127.17
2	C	352	NAD	O4B-C1B-N9A	2.56	113.00	108.09
2	D	352	NAD	O4B-C1B-N9A	2.44	112.77	108.09
2	D	352	NAD	C4A-N9A-C8A	2.38	108.23	105.74
2	C	352	NAD	C4A-N9A-C8A	2.37	108.22	105.74

There are no chirality outliers.

All (4) torsion outliers are listed below:

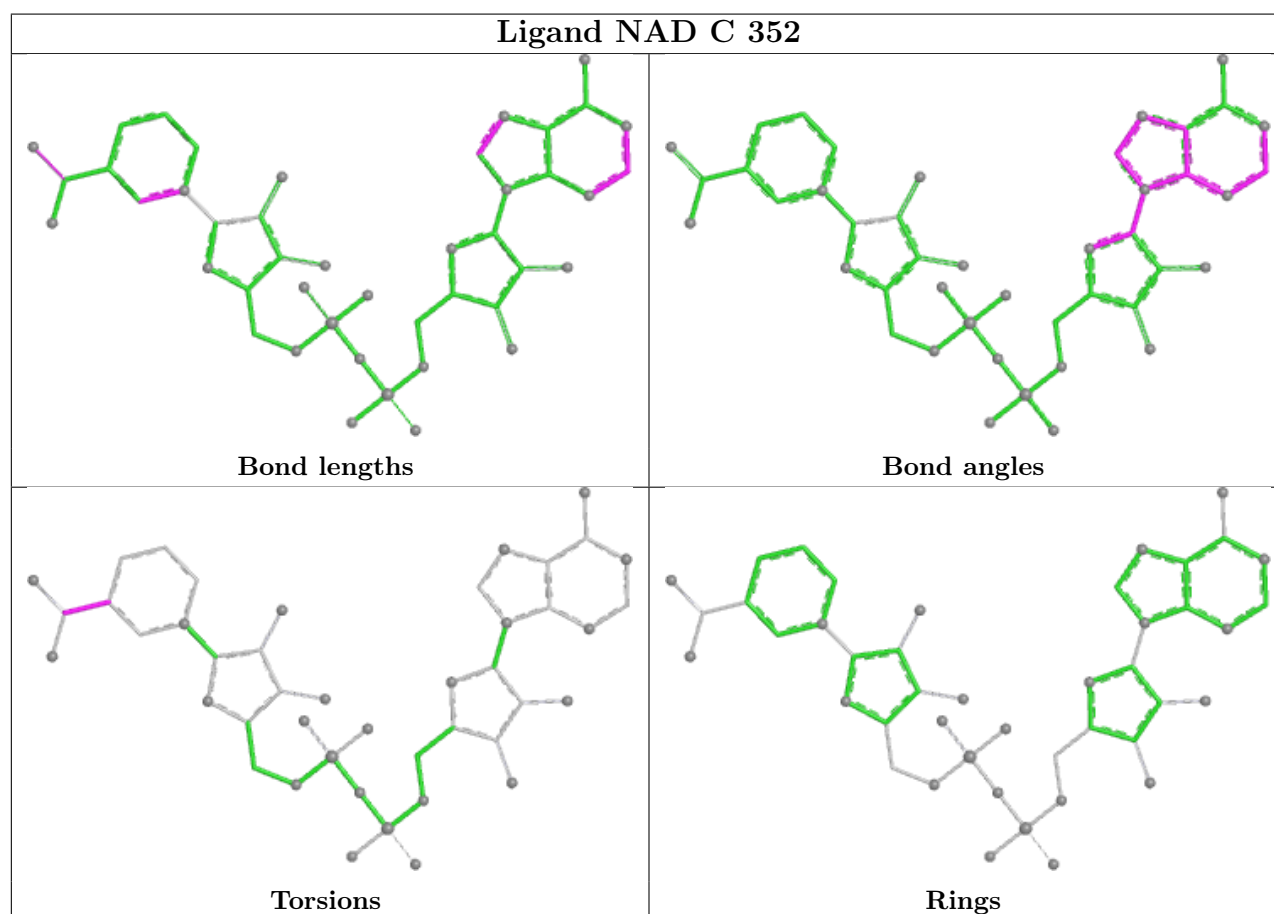
Mol	Chain	Res	Type	Atoms
2	C	352	NAD	C2N-C3N-C7N-O7N
2	C	352	NAD	C2N-C3N-C7N-N7N
2	D	352	NAD	C2N-C3N-C7N-O7N
2	D	352	NAD	C2N-C3N-C7N-N7N

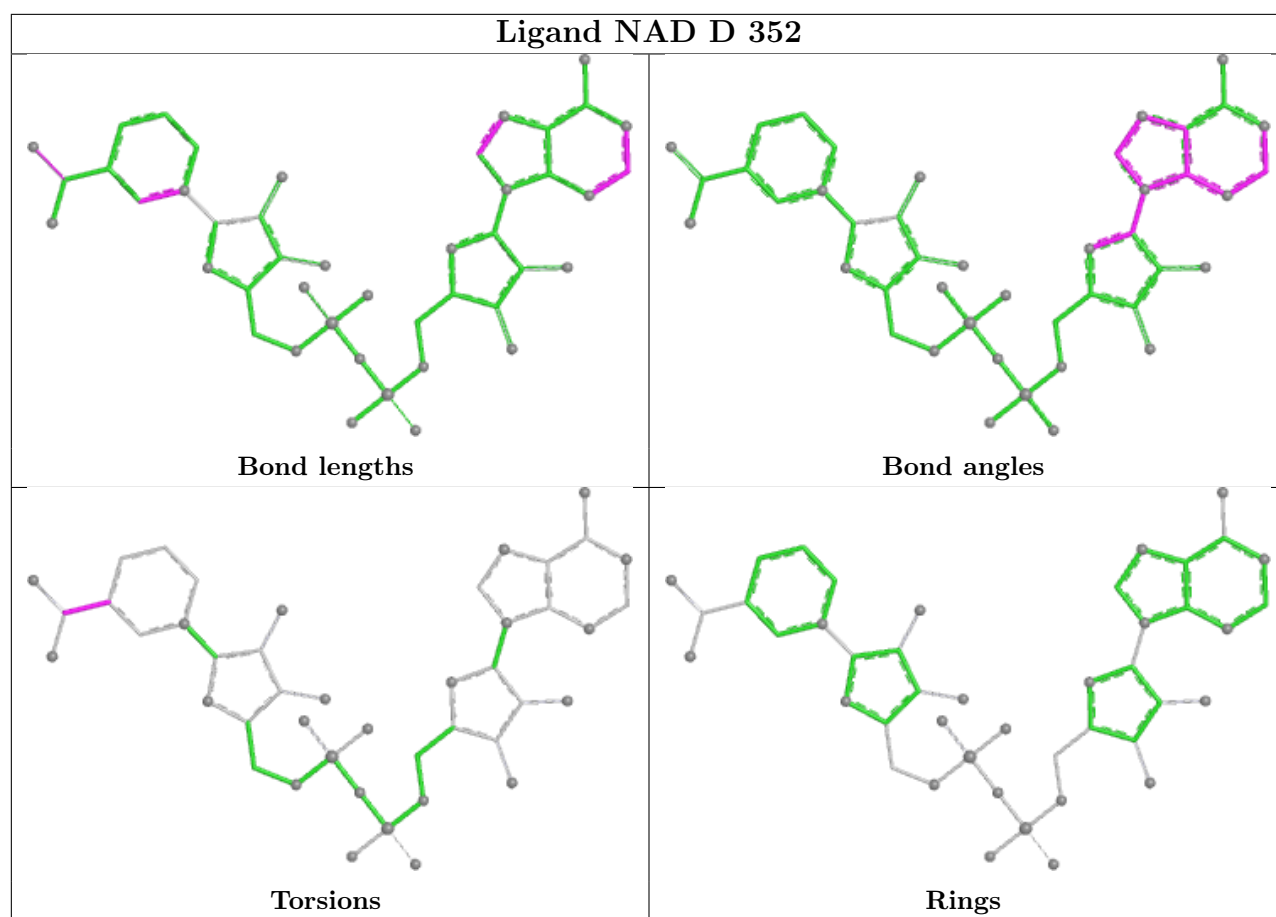
There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	352	NAD	1	0
2	D	352	NAD	1	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 [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	314/326 (96%)	-1.56	0 100 100	29, 41, 70, 88	0
1	B	314/326 (96%)	-1.56	0 100 100	29, 42, 71, 85	0
1	C	312/326 (95%)	-1.38	0 100 100	39, 63, 100, 136	0
1	D	312/326 (95%)	-1.38	0 100 100	38, 63, 101, 131	0
All	All	1252/1304 (96%)	-1.47	0 100 100	29, 52, 89, 136	0

There are no RSRZ outliers to report.

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 oligosaccharides in this entry.

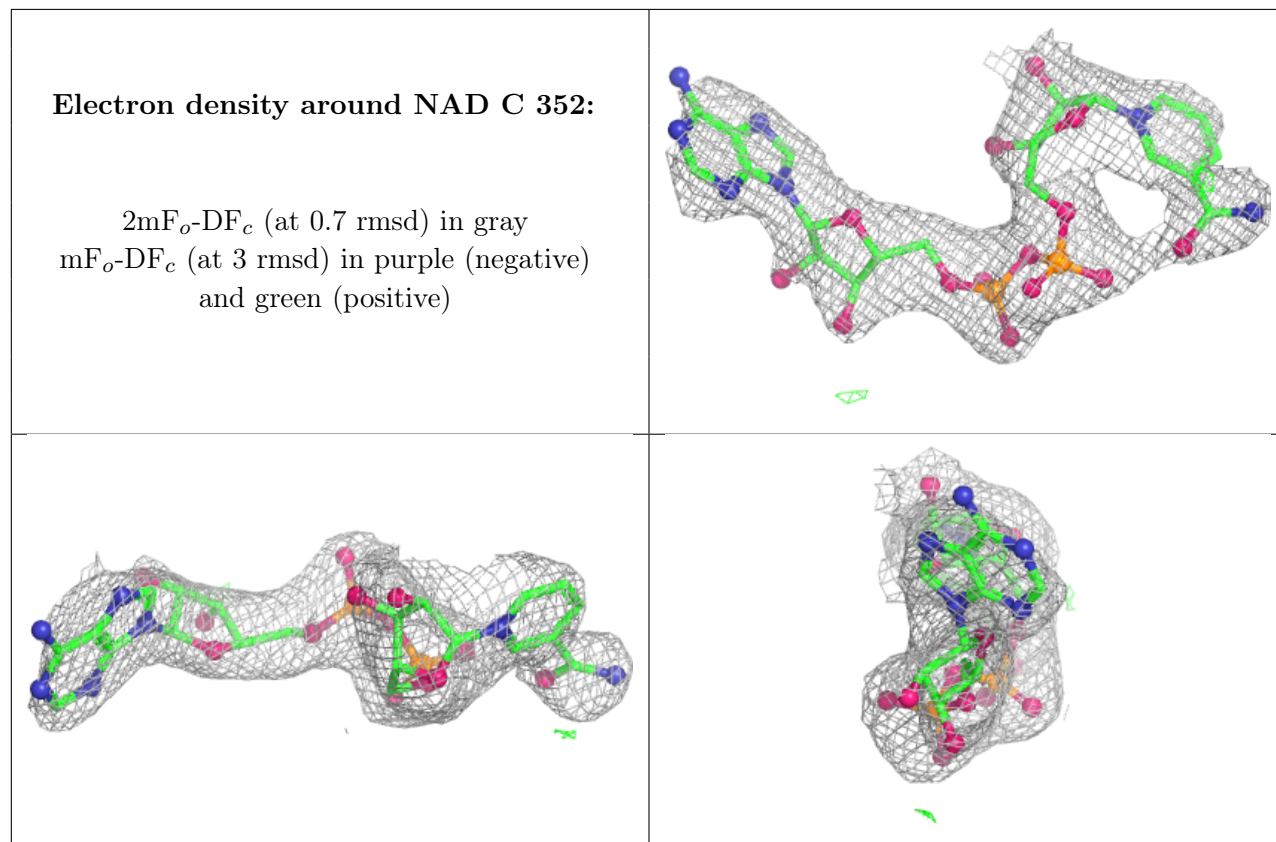
6.4 Ligands [i](#)

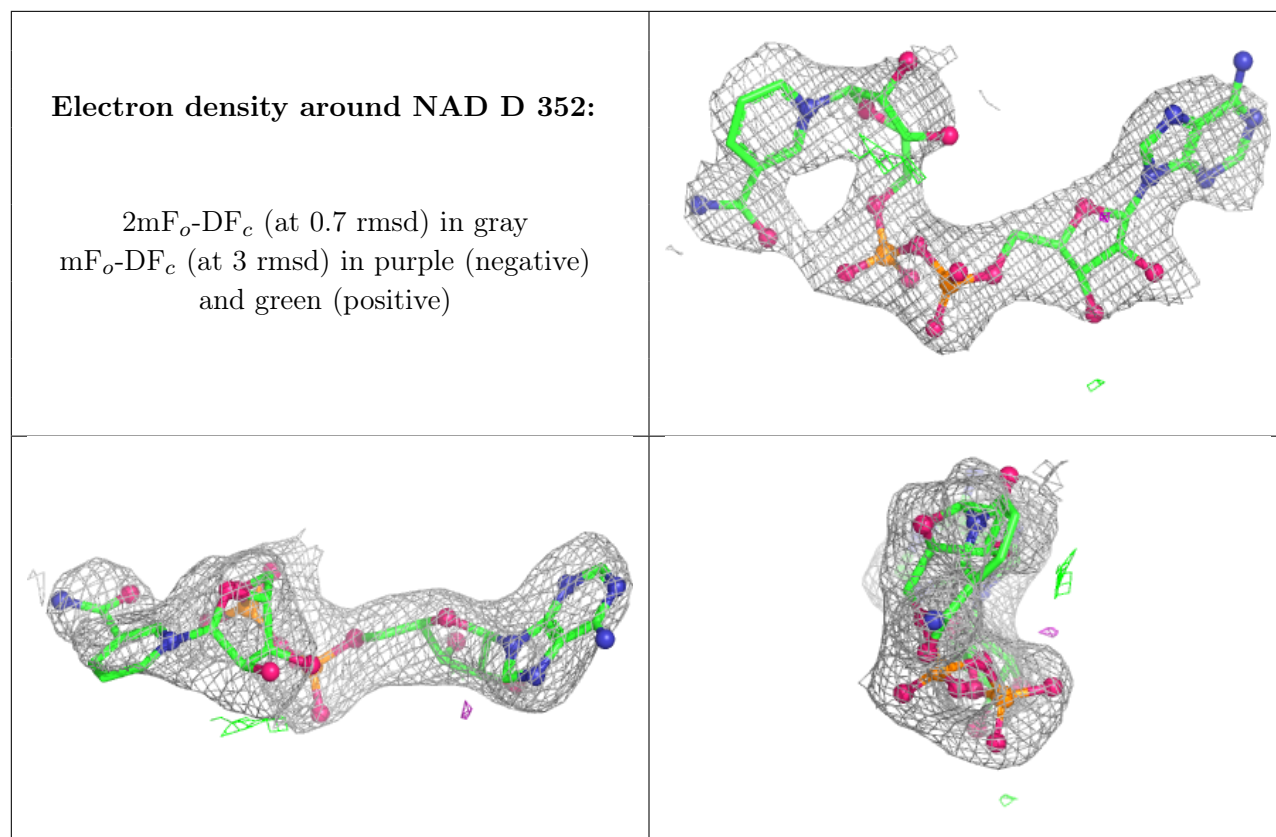
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists 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	C	352	44/44	0.99	0.04	58,74,82,85	0
2	NAD	D	352	44/44	0.99	0.04	60,73,82,85	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.





6.5 Other polymers [i](#)

There are no such residues in this entry.