



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 11:45 AM GMT

PDB ID : 3PQF
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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

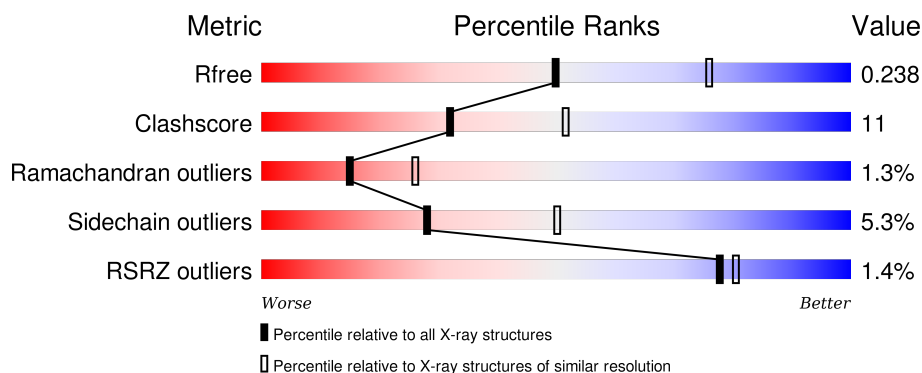
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.49 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}	91344	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	326	<div> <div>85%</div> <div>10% . .</div> </div>
1	B	326	<div> <div>81%</div> <div>13% . .</div> </div>
1	C	326	<div> <div>2%</div> <div>74%</div> <div>18% . . .</div> </div>
1	D	326	<div> <div>2%</div> <div>68%</div> <div>23% . . .</div> </div>

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:

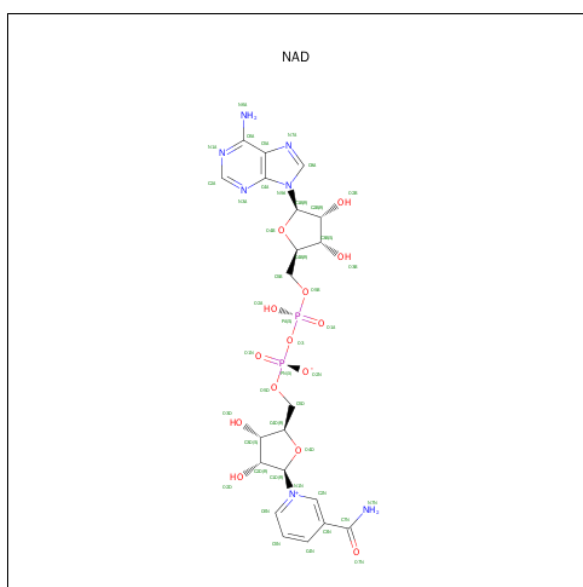
Chain	Residue	Modelled	Actual	Comment	Reference
A	171	CYS	HIS	ENGINEERED MUTATION	UNP P13714
A	321	HIS	-	EXPRESSION TAG	UNP P13714
A	322	HIS	-	EXPRESSION TAG	UNP P13714
A	323	HIS	-	EXPRESSION TAG	UNP P13714
A	324	HIS	-	EXPRESSION TAG	UNP P13714
A	325	HIS	-	EXPRESSION TAG	UNP P13714
A	326	HIS	-	EXPRESSION 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 (three-letter code: NAD) (formula: $C_{21}H_{27}N_7O_{14}P_2$).



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

- Molecule 3 is water.

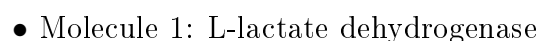
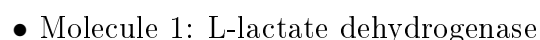
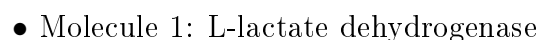
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	82	Total	O	0	0
			82	82		
3	B	94	Total	O	0	0
			94	94		
3	C	17	Total	O	0	0
			17	17		

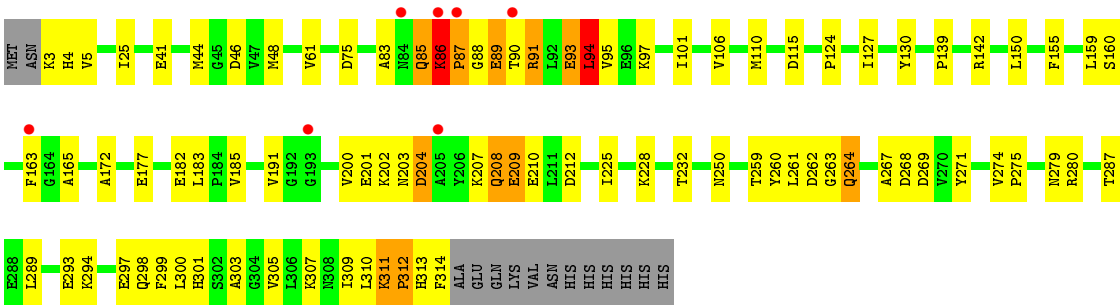
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	24	Total	O	0	0
			24	24		

- 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	133.43 Å 133.43 Å 99.09 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.49 – 2.49 31.76 – 2.49	Depositor EDS
% Data completeness (in resolution range)	99.8 (30.49-2.49) 99.8 (31.76-2.49)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.24 (at 2.48 Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6 _289)	Depositor
R, R_{free}	0.189 , 0.235 0.189 , 0.238	Depositor DCC
R_{free} test set	2062 reflections (3.08%)	DCC
Wilson B-factor (Å ²)	48.6	Xtriage
Anisotropy	0.086	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 29.3	EDS
Estimated twinning fraction	0.478 for -h,-k,l 0.030 for h,-h-k,-l 0.031 for -k,-h,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 68954 reflections	Xtriage
F_o, F_c correlation	0.96	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, 0.375 respectively for untwinned datasets, and 0.333, 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	0.49	0/2443	0.56	0/3311
1	B	0.49	0/2443	0.57	0/3311
1	C	0.39	0/2430	0.53	1/3293 (0.0%)
1	D	0.39	0/2430	0.53	0/3293
All	All	0.44	0/9746	0.55	1/13208 (0.0%)

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 (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	C	86	LYS	C-N-CD	-6.77	105.70	120.60

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	22	0
1	B	2397	0	2384	31	0
1	C	2384	0	2373	78	0
1	D	2384	0	2373	80	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	203	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 (203) 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:B:52:HIS:HD2	1:C:154:ARG:HD2	1.39	0.87
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:C:84:ASN:C	1:C:85:GLN:O	2.12	0.82
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:HD2	1:C:86:LYS:N	2.01	0.76
1:C:86:LYS:NZ	1:C:97:LYS:HZ1	1.84	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:D:93:GLU:C	1:D:95:VAL:H	1.93	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: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:93:GLU:C	1:C:95:VAL:H	1.99	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:B:250:ASN:ND2	1:B:280:ARG:H	1.95	0.64
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
1:C:85:GLN:CG	1:C:87:PRO:HD2	2.18	0.63
1:C:85:GLN:CD	1:C:87:PRO:HD3	2.18	0.63
1:C:86:LYS:HZ3	1:C:97:LYS:HZ1	1.46	0.62
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:OE1	1:B:213:GLN:HA	2.01	0.60
1:C:170:VAL:HG12	1:C:170:VAL:O	2.01	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:C:86:LYS:C	1:C:88:GLY:N	2.53	0.60
1:D:274:VAL:HG23	1:D:275:PRO:HD2	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:D:86:LYS:C	1:D:88:GLY:H	2.05	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:B:156:ARG:C	3:B:417:HOH:O	2.41	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:D:311:LYS:O	1:D:313:HIS:N	2.37	0.58
1:B:2:ASN:HA	3:B:371:HOH:O	2.04	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:D:177:GLU:O	1:D:182:GLU:HB3	2.05	0.57
1:C:40:LYS:HG2	1:C:65:TYR:OH	2.04	0.57
1:C:85:GLN:OE1	1:C:87:PRO:HG3	2.05	0.57
1:C:86:LYS:NZ	1:C:97:LYS:NZ	2.51	0.57
1:B:210:GLU:O	1:B:214:ILE:HG13	2.05	0.56
1:D:89:GLU:HB3	1:D:91:ARG:HG3	1.87	0.56
1:D:90:THR:O	1:D:91:ARG:C	2.44	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:D:94:LEU:O	1:D:94:LEU:HG	2.05	0.54
1:C:89:GLU:HB3	1:C:91:ARG:CG	2.37	0.54
1:D:86:LYS:HE2	1:D:94:LEU:HD12	1.90	0.54
1:A:140:LYS:HD2	1:A:260:TYR:CD2	2.43	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
1:A:44:MET:O	1:A:48:MET:HG3	2.08	0.53
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: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:93:GLU:C	1:D:95:VAL:N	2.62	0.51
1:D:155:PHE:O	1:D:159:LEU:HG	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:D:88:GLY:O	1:D:90:THR:N	2.43	0.51
1:C:209:GLU:HA	1:C:212:ASP:HB2	1.92	0.51
1:C:86:LYS:HD2	1:C:86:LYS:H	1.74	0.51
1:C:93:GLU:C	1:C:95:VAL:N	2.63	0.51
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: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:D:185:VAL:HG11	1:D:289:LEU:HD13	1.96	0.47
1:A:228:LYS:HD2	1:D:48:MET:CE	2.44	0.47
1:C:94:LEU:HG	1:C:94:LEU:O	2.14	0.47
1:D:311:LYS:O	1:D:312:PRO:C	2.53	0.47
1:D:264:GLN:H	1:D:264:GLN:HG2	1.43	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:C:86:LYS:HZ1	1:C:97:LYS:NZ	2.12	0.47
1:D:150:LEU:HB3	3:D:348:HOH:O	2.15	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: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:C:29:ILE:HD13	1:C:245:LYS:HB2	1.98	0.46
1:C:250:ASN:HD22	1:C:279:ASN:HB2	1.80	0.45
1:C:89:GLU:O	1:C:91:ARG:HG3	2.16	0.45
1:D:263:GLY:CA	1:D:267:ALA:O	2.65	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:228:LYS:HD2	1:D:48:MET:HE3	1.98	0.45
1:D:25:ILE:HG12	1:D:61:VAL:HG11	1.97	0.45
1:C:91:ARG:C	1:C:93:GLU:N	2.70	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:C:185:VAL:O	1:C:185:VAL:HG12	2.17	0.45
1:B:77:VAL:HG23	1:B:114:PHE:CE1	2.52	0.44
1:C:162:TYR:HE1	1:C:210:GLU:HG2	1.82	0.44
1:B:194:VAL:HG12	1:B:199:LEU:HG	1.99	0.44
1:D:86:LYS:NZ	1:D:91:ARG:HD3	2.31	0.44
1:B:155:PHE:CE2	1:B:172:ALA:HB1	2.52	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:7:LYS:HE3	1:C:34:VAL:CG2	2.48	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:B:228:LYS:HG3	1:C:48:MET:CE	2.48	0.43
1:B:228:LYS:HD3	1:C:41:GLU:OE1	2.19	0.43
1:A:261:LEU:HD22	1:A:264:GLN:HG2	2.00	0.43
1:B:6:ASN:HB3	1:B:248:LEU:HD22	2.01	0.43
1:D:91:ARG:C	1:D:93:GLU:N	2.71	0.43
1:D:294:LYS:HE2	1:D:298:GLN:HE22	1.84	0.43
1:D:97:LYS:O	1:D:101:ILE:HG13	2.19	0.43
1:D:311:LYS:N	1:D:312:PRO:HD2	2.33	0.43
1:A:41:GLU:OE2	1:D:228:LYS:HE3	2.18	0.43
1:A:40:LYS:HE2	1:A:41:GLU:OE2	2.18	0.43
1:C:262:ASP:H	1:C:264:GLN:HG3	1.84	0.43
1:D:311:LYS:C	1:D:313:HIS:N	2.72	0.42
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:C:86:LYS:HZ3	1:C:97:LYS:NZ	2.14	0.42
1:D:263:GLY:N	1:D:267:ALA:O	2.52	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: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:HG13	1:D:309:ILE:H	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:85:GLN:OE1	1:C:87:PRO:HD3	2.19	0.42
1:A:157:PHE:C	1:A:157:PHE:CD2	2.92	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:191:VAL:HB	1:A:199:LEU:HD11	2.01	0.42
1:C:87:PRO:C	1:C:88:GLY:O	2.58	0.41
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:C:89:GLU:CB	1:C:91:ARG:HG3	2.50	0.41
1:D:208:GLN:CD	1:D:208:GLN:H	2.24	0.41
1:B:229:LYS:HB2	1:C:45:GLY:HA3	2.03	0.41
1:D:93:GLU:O	1:D:95:VAL:N	2.52	0.41
1:A:194:VAL:HG12	1:A:199:LEU:HG	2.01	0.41
1:C:214:ILE:O	1:C:218:VAL:HG23	2.21	0.41
1:B:40:LYS:HE3	1:B:40:LYS:HB3	1.85	0.41
1:D:225:ILE:CG2	1:D:232:THR:HG22	2.51	0.41
1:D:115:ASP:O	1:D:142:ARG:NH2	2.53	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:D:130:TYR:HA	1:D:271:TYR:CD2	2.56	0.40
1:C:202:LYS:HB2	1:C:202:LYS:HE3	1.97	0.40
1:D:185:VAL:O	1:D:185:VAL:HG12	2.21	0.40
1:C:86:LYS:H	1:C:87:PRO:CD	2.34	0.40
1:A:155:PHE:CE2	1:A:172:ALA:HB1	2.56	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%)	46	68
1	C	310/326 (95%)	286 (92%)	16 (5%)	8 (3%)	7	10

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	310/326 (95%)	289 (93%)	14 (4%)	7 (2%)	8	12
All	All	1244/1304 (95%)	1168 (94%)	60 (5%)	16 (1%)	15	26

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
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%)	35	61
1	B	255/267 (96%)	241 (94%)	14 (6%)	27	48
1	C	254/267 (95%)	238 (94%)	16 (6%)	22	40
1	D	254/267 (95%)	241 (95%)	13 (5%)	29	52
All	All	1018/1068 (95%)	964 (95%)	54 (5%)	28	50

All (54) 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	183	LEU
1	A	203	ASN
1	A	204	ASP
1	A	229	LYS
1	A	264	GLN
1	B	61	VAL
1	B	62	LYS
1	B	85	GLN
1	B	152	SER
1	B	156	ARG
1	B	158	MET
1	B	168	GLN
1	B	183	LEU
1	B	190	ASN
1	B	204	ASP
1	B	207	LYS
1	B	208	GLN
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	85	GLN
1	C	86	LYS
1	C	94	LEU
1	C	97	LYS
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	264	GLN
1	C	314	PHE
1	D	5	VAL
1	D	46	ASP

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Mol	Chain	Res	Type
1	D	85	GLN
1	D	86	LYS
1	D	94	LEU
1	D	183	LEU
1	D	203	ASN
1	D	204	ASP
1	D	208	GLN
1	D	209	GLU
1	D	264	GLN
1	D	287	THR
1	D	311	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	51	ASN
1	A	208	GLN
1	A	220	ASN
1	A	250	ASN
1	B	52	HIS
1	B	190	ASN
1	B	208	GLN
1	B	250	ASN
1	C	169	ASN
1	C	188	HIS
1	C	203	ASN
1	C	250	ASN
1	C	264	GLN
1	D	4	HIS
1	D	203	ASN
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 [i](#)

There are no carbohydrates 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	C	352	-	38,48,48	1.69	3 (7%)	47,73,73	2.11	4 (8%)
2	NAD	D	352	-	38,48,48	1.69	3 (7%)	47,73,73	2.07	4 (8%)

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	-	-	0/22/62/62	0/5/5/5
2	NAD	D	352	-	-	0/22/62/62	0/5/5/5

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	352	NAD	C2A-N1A	2.76	1.39	1.33
2	D	352	NAD	C2A-N1A	2.77	1.39	1.33
2	D	352	NAD	C2A-N3A	3.65	1.38	1.32
2	C	352	NAD	C2A-N3A	3.80	1.38	1.32
2	C	352	NAD	O7N-C7N	8.24	1.41	1.24
2	D	352	NAD	O7N-C7N	8.27	1.41	1.24

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	352	NAD	N3A-C2A-N1A	-12.43	119.38	128.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	352	NAD	N3A-C2A-N1A	-12.19	119.56	128.89
2	C	352	NAD	PN-O3-PA	-3.50	122.91	132.73
2	D	352	NAD	PN-O3-PA	-3.12	123.96	132.73
2	D	352	NAD	O4B-C1B-N9A	2.23	112.77	108.10
2	C	352	NAD	O4B-C1B-N9A	2.34	113.00	108.10
2	D	352	NAD	O4D-C1D-N1N	2.58	110.97	108.13
2	C	352	NAD	O4D-C1D-N1N	3.05	111.48	108.13

There are no chirality outliers.

There are no torsion outliers.

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

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 ⓘ

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%)	-0.03	2 (0%) 90 91	29, 41, 70, 88	0
1	B	314/326 (96%)	-0.01	3 (0%) 84 86	29, 42, 71, 85	0
1	C	312/326 (95%)	0.07	5 (1%) 74 78	39, 63, 100, 136	0
1	D	312/326 (95%)	0.08	7 (2%) 65 69	38, 63, 101, 131	0
All	All	1252/1304 (96%)	0.03	17 (1%) 78 80	29, 52, 89, 136	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	205	ALA	5.7
1	D	205	ALA	4.5
1	C	90	THR	4.2
1	D	90	THR	3.4
1	D	87	PRO	3.4
1	B	157	PHE	3.2
1	D	84	ASN	2.9
1	C	163	PHE	2.8
1	A	157	PHE	2.7
1	D	86	LYS	2.6
1	B	168	GLN	2.6
1	C	87	PRO	2.4
1	D	163	PHE	2.3
1	A	205	ALA	2.3
1	C	92	LEU	2.2
1	B	205	ALA	2.1
1	D	193	GLY	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
2	NAD	C	352	44/44	0.95	0.17	0.56	58,74,82,85	0
2	NAD	D	352	44/44	0.96	0.15	-0.09	60,73,82,85	0

6.5 Other polymers [i](#)

There are no such residues in this entry.