



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:43 PM GMT

PDB ID : 1GUZ  
Title : Structural Basis for Thermophilic Protein Stability: Structures of Thermophilic and Mesophilic Malate Dehydrogenases  
Authors : Dalhus, B.; Sarinen, M.; Sauer, U.H.; Eklund, P.; Johansson, K.; Karlsson, A.; Ramaswamy, S.; Bjork, A.; Synstad, B.; Naterstad, K.; Sirevag, R.; Eklund, H.  
Deposited on : 2002-02-04  
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](mailto: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.

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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

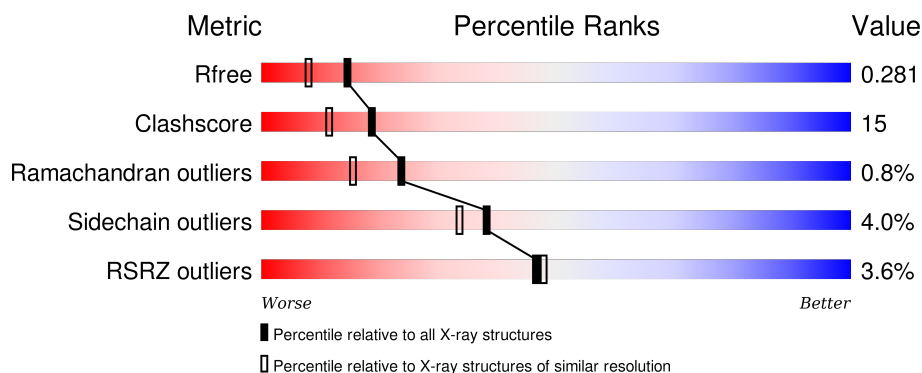
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.00 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))
$R_{free}$	91344	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	310	<div> <div>3%</div> <div>72% 24% . .</div> </div>
1	B	310	<div> <div>4%</div> <div>70% 23% . 5%</div> </div>
1	C	310	<div> <div>4%</div> <div>68% 26% . .</div> </div>
1	D	310	<div> <div>3%</div> <div>64% 28% . 5%</div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10067 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	305	Total	C	N	O	S	0	0	0
			2294	1450	392	437	15			
1	B	296	Total	C	N	O	S	0	0	0
			2221	1406	376	425	14			
1	C	305	Total	C	N	O	S	0	0	0
			2294	1450	392	437	15			
1	D	295	Total	C	N	O	S	0	0	0
			2213	1402	375	422	14			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	227	ALA	SER	CONFLICT	UNP P80039
A	229	ALA	GLY	CONFLICT	UNP P80039
B	227	ALA	SER	CONFLICT	UNP P80039
B	229	ALA	GLY	CONFLICT	UNP P80039
C	227	ALA	SER	CONFLICT	UNP P80039
C	229	ALA	GLY	CONFLICT	UNP P80039
D	227	ALA	SER	CONFLICT	UNP P80039
D	229	ALA	GLY	CONFLICT	UNP P80039

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

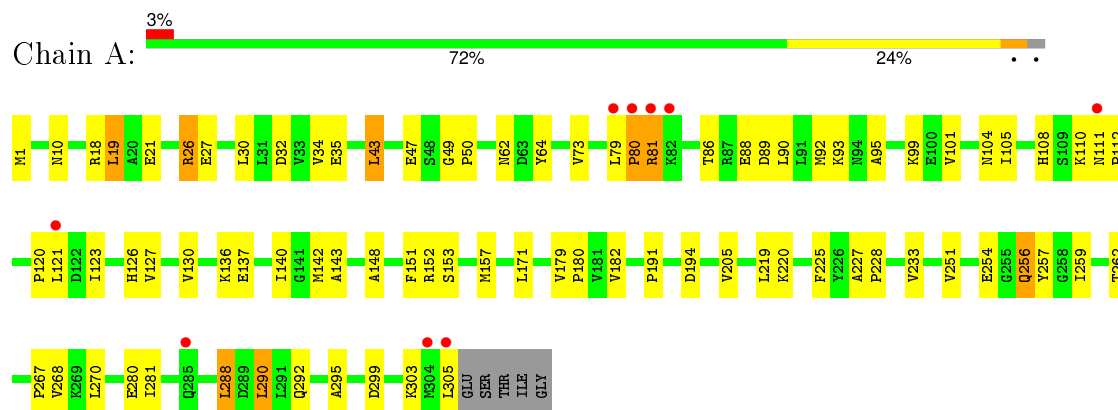
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	194	Total	O	0	0
			194	194		
3	B	243	Total	O	0	0
			243	243		
3	C	185	Total	O	0	0
			185	185		
3	D	247	Total	O	0	0
			247	247		

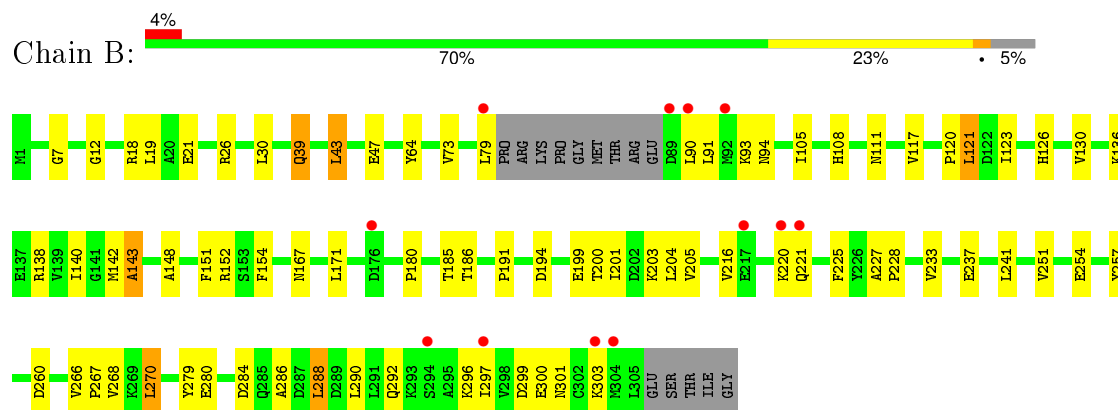
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

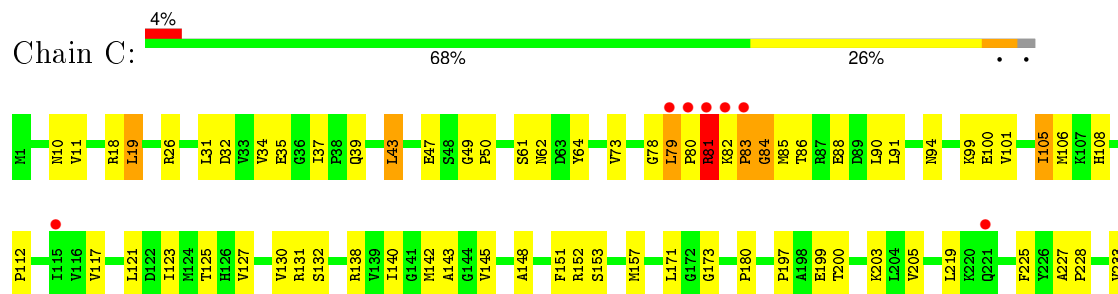
#### • Molecule 1: MALATE DEHYDROGENASE

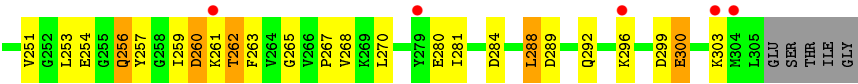


#### • Molecule 1: MALATE DEHYDROGENASE

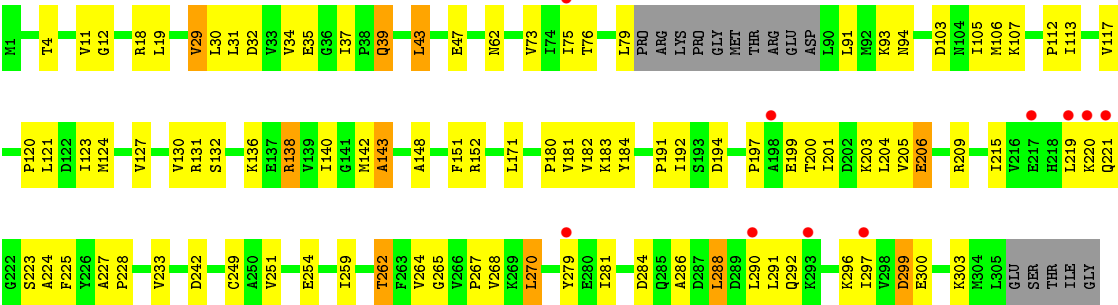


#### • Molecule 1: MALATE DEHYDROGENASE





● Molecule 1: MALATE DEHYDROGENASE



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	83.89Å 117.39Å 125.33Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	26.58 – 2.00 26.58 – 2.00	Depositor EDS
% Data completeness (in resolution range)	98.0 (26.58-2.00) 98.2 (26.58-2.00)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.71 (at 2.01Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.243 , 0.292 0.235 , 0.281	Depositor DCC
$R_{free}$ test set	4114 reflections (5.00%)	DCC
Wilson B-factor (Å <sup>2</sup> )	24.3	Xtriage
Anisotropy	0.418	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 55.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 82230 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	10067	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 49.37 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 7.5318e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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 [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	0.30	0/2324	0.57	0/3145
1	B	0.32	0/2248	0.59	0/3042
1	C	0.30	0/2324	0.58	0/3145
1	D	0.32	0/2240	0.59	0/3031
All	All	0.31	0/9136	0.58	0/12363

There are no bond length outliers.

There are no bond angle outliers.

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	2294	0	2378	63	0
1	B	2221	0	2299	61	0
1	C	2294	0	2378	85	0
1	D	2213	0	2295	80	0
2	A	44	0	25	1	0
2	B	44	0	25	2	0
2	C	44	0	25	4	0
2	D	44	0	25	3	0
3	A	194	0	0	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	243	0	0	4	0
3	C	185	0	0	4	0
3	D	247	0	0	7	0
All	All	10067	0	9450	288	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 15.

All (288) 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:D:121:LEU:HD22	1:D:143:ALA:HB2	1.42	0.99
1:C:80:PRO:HG3	1:C:94:ASN:HD21	1.28	0.99
1:A:121:LEU:HD22	1:A:143:ALA:HB2	1.49	0.94
1:B:18:ARG:HD3	1:B:233:VAL:HG21	1.48	0.94
1:B:121:LEU:HD22	1:B:143:ALA:HB2	1.48	0.94
1:D:34:VAL:HG11	1:D:37:ILE:HD12	1.52	0.90
1:C:121:LEU:HD22	1:C:143:ALA:HB2	1.56	0.87
1:C:81:ARG:HH11	1:C:81:ARG:HB3	1.41	0.84
1:C:11:VAL:HG12	2:C:1306:NAD:O2N	1.77	0.84
1:A:18:ARG:HD3	1:A:233:VAL:HG21	1.60	0.83
1:D:79:LEU:H	1:D:94:ASN:HD21	1.26	0.83
1:C:86:THR:HG22	1:C:88:GLU:H	1.44	0.82
1:C:79:LEU:N	1:C:80:PRO:HD3	1.98	0.78
1:C:81:ARG:NH1	1:C:81:ARG:HB3	1.96	0.78
1:A:171:LEU:HD22	1:A:267:PRO:HD3	1.66	0.78
1:C:81:ARG:HG2	1:C:90:LEU:HD22	1.66	0.78
1:A:153:SER:O	1:A:157:MET:HG3	1.85	0.77
1:C:80:PRO:HG3	1:C:94:ASN:ND2	2.00	0.76
1:D:18:ARG:HD3	1:D:233:VAL:HG21	1.67	0.76
1:A:140:ILE:HG22	1:A:251:VAL:HG12	1.68	0.76
1:D:138:ARG:HH11	1:D:138:ARG:HB3	1.51	0.75
1:C:18:ARG:HD3	1:C:233:VAL:HG21	1.69	0.75
1:C:80:PRO:HA	1:C:90:LEU:HD13	1.68	0.73
1:D:32:ASP:OD1	1:D:34:VAL:HG12	1.88	0.73
1:B:140:ILE:HG22	1:B:251:VAL:HG12	1.71	0.72
1:D:43:LEU:O	1:D:47:GLU:HG3	1.90	0.72
1:C:171:LEU:HD22	1:C:267:PRO:HD3	1.70	0.72
1:A:26:ARG:H	1:A:26:ARG:HD2	1.57	0.69
1:B:180:PRO:HG3	1:B:205:VAL:HG13	1.72	0.69
1:C:79:LEU:H	1:C:80:PRO:HD3	1.55	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:43:LEU:O	1:B:47:GLU:HG3	1.93	0.68
1:A:121:LEU:CD2	1:A:143:ALA:HB2	2.21	0.68
1:C:153:SER:O	1:C:157:MET:HG3	1.93	0.68
1:D:39:GLN:H	1:D:39:GLN:HE21	1.39	0.67
1:C:101:VAL:O	1:C:105:ILE:HG23	1.94	0.67
1:D:39:GLN:H	1:D:39:GLN:NE2	1.92	0.66
1:D:76:THR:HG22	1:D:117:VAL:HG21	1.75	0.66
1:A:179:VAL:HG22	1:A:290:LEU:HD13	1.79	0.65
1:B:79:LEU:HD21	1:B:93:LYS:HB3	1.79	0.65
1:D:113:ILE:HG13	1:D:138:ARG:HH12	1.63	0.64
1:C:121:LEU:CD2	1:C:143:ALA:HB2	2.27	0.64
1:C:292:GLN:O	1:C:296:LYS:HG2	1.97	0.64
1:A:136:LYS:HZ3	1:A:254:GLU:HB3	1.62	0.63
1:B:220:LYS:HD2	1:B:221:GLN:HE21	1.63	0.63
1:B:18:ARG:HD3	1:B:233:VAL:CG2	2.25	0.63
1:B:39:GLN:NE2	1:B:39:GLN:H	1.97	0.62
1:C:180:PRO:HG3	1:C:205:VAL:HG13	1.81	0.62
1:A:73:VAL:HG21	1:A:105:ILE:HD12	1.80	0.62
1:A:136:LYS:NZ	1:A:254:GLU:HB3	2.13	0.62
1:C:127:VAL:O	1:C:130:VAL:HG22	1.99	0.62
1:D:11:VAL:HG22	2:D:1306:NAD:O2N	2.00	0.62
1:A:262:THR:HG21	1:A:295:ALA:HB1	1.81	0.62
1:B:286:ALA:O	1:B:290:LEU:HG	1.98	0.62
1:A:180:PRO:HG3	1:A:205:VAL:HG13	1.81	0.62
1:C:32:ASP:OD1	1:C:34:VAL:HG12	1.99	0.62
1:C:284:ASP:HB3	3:C:2109:HOH:O	2.00	0.61
1:C:91:LEU:HD11	1:C:123:ILE:HG21	1.83	0.61
1:A:81:ARG:HD3	1:A:90:LEU:HB2	1.80	0.61
1:A:81:ARG:HG2	1:A:90:LEU:HD22	1.82	0.60
1:B:299:ASP:O	1:B:303:LYS:HG3	2.01	0.60
1:B:39:GLN:HE21	1:B:39:GLN:H	1.48	0.59
1:A:101:VAL:O	1:A:105:ILE:HG23	2.01	0.59
1:A:81:ARG:CG	1:A:90:LEU:HD22	2.31	0.59
1:B:200:THR:O	1:B:203:LYS:HG2	2.02	0.59
1:B:90:LEU:O	1:B:90:LEU:HD23	2.03	0.59
1:B:167:ASN:O	1:B:185:THR:HA	2.03	0.58
1:A:1:MET:H2	1:A:26:ARG:HD3	1.68	0.58
1:B:73:VAL:HG11	1:B:105:ILE:HD13	1.85	0.58
1:D:113:ILE:HG13	1:D:138:ARG:NH1	2.19	0.57
1:C:19:LEU:HD13	1:C:233:VAL:HG22	1.85	0.57
1:B:79:LEU:HD23	1:B:94:ASN:OD1	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:82:LYS:HB2	1:C:85:MET:HG3	1.85	0.57
1:D:121:LEU:CD2	1:D:143:ALA:HB2	2.26	0.57
1:C:43:LEU:O	1:C:47:GLU:HG3	2.05	0.56
1:A:157:MET:HE1	3:A:2133:HOH:O	2.06	0.56
1:A:43:LEU:O	1:A:47:GLU:HG3	2.05	0.56
1:B:288:LEU:O	1:B:292:GLN:HG3	2.05	0.56
1:A:120:PRO:HB2	3:A:2081:HOH:O	2.05	0.55
1:C:148:ALA:O	1:C:152:ARG:HD3	2.07	0.55
1:D:76:THR:HG22	1:D:117:VAL:CG2	2.37	0.55
1:C:112:PRO:C	1:C:138:ARG:HH12	2.10	0.55
1:D:296:LYS:O	1:D:300:GLU:HG3	2.06	0.55
1:B:216:VAL:HG13	3:B:2178:HOH:O	2.06	0.55
1:D:182:VAL:HG13	1:D:192:ILE:HD11	1.89	0.55
1:B:91:LEU:HD22	1:B:301:ASN:HD22	1.73	0.54
1:B:121:LEU:CD2	1:B:143:ALA:HB2	2.30	0.54
1:B:90:LEU:HD22	1:B:120:PRO:HD3	1.89	0.54
1:C:26:ARG:HH11	1:C:26:ARG:HG2	1.73	0.54
1:C:86:THR:HG22	1:C:88:GLU:N	2.18	0.54
1:B:268:VAL:HG23	1:B:270:LEU:HD13	1.90	0.54
1:D:112:PRO:O	1:D:138:ARG:NH1	2.41	0.54
1:B:79:LEU:O	1:B:79:LEU:HG	2.08	0.53
1:C:262:THR:HB	1:C:299:ASP:OD1	2.09	0.53
1:A:1:MET:N	1:A:26:ARG:HD3	2.23	0.53
1:A:18:ARG:HD3	1:A:233:VAL:CG2	2.35	0.53
1:D:140:ILE:HG22	1:D:251:VAL:HG12	1.90	0.53
1:B:148:ALA:O	1:B:152:ARG:HG3	2.09	0.53
1:C:80:PRO:HA	1:C:90:LEU:CD1	2.37	0.53
1:A:288:LEU:O	1:A:292:GLN:HG3	2.08	0.52
1:C:79:LEU:N	1:C:80:PRO:CD	2.71	0.52
1:B:257:TYR:CZ	1:B:280:GLU:HA	2.44	0.52
1:C:199:GLU:O	1:C:203:LYS:HG3	2.10	0.52
1:B:220:LYS:HD2	1:B:221:GLN:NE2	2.24	0.52
1:C:80:PRO:HD2	2:C:1306:NAD:H3D	1.92	0.52
2:D:1306:NAD:H6N	3:D:2243:HOH:O	2.10	0.52
1:C:253:LEU:HD22	1:C:256:GLN:CG	2.40	0.52
1:C:257:TYR:CZ	1:C:280:GLU:HA	2.45	0.52
1:D:227:ALA:HB3	1:D:228:PRO:HD3	1.92	0.52
1:B:297:ILE:HB	3:B:2226:HOH:O	2.08	0.51
1:C:80:PRO:CG	1:C:94:ASN:HD21	2.11	0.51
1:D:264:VAL:HG11	1:D:291:LEU:HD11	1.92	0.51
1:C:80:PRO:HG3	1:C:90:LEU:HD11	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:227:ALA:HB3	1:C:228:PRO:HD3	1.91	0.51
1:D:279:TYR:HE2	3:D:2215:HOH:O	1.92	0.51
1:A:21:GLU:CD	1:D:18:ARG:HH12	2.13	0.51
1:C:73:VAL:HG21	1:C:105:ILE:HD13	1.93	0.51
1:D:300:GLU:O	1:D:303:LYS:HB3	2.10	0.51
1:C:10:ASN:HB3	1:C:225:PHE:CG	2.45	0.51
1:D:262:THR:HG21	3:D:2230:HOH:O	2.11	0.51
1:D:286:ALA:O	1:D:290:LEU:HG	2.11	0.51
1:B:186:THR:HG22	1:B:191:PRO:HA	1.92	0.51
1:A:148:ALA:O	1:A:152:ARG:HD3	2.11	0.51
1:D:288:LEU:O	1:D:292:GLN:HG3	2.11	0.50
1:D:215:ILE:HG21	1:D:224:ALA:HB2	1.93	0.50
1:D:206:GLU:OE2	1:D:209:ARG:HD2	2.11	0.50
1:A:262:THR:HB	1:A:299:ASP:OD1	2.10	0.50
1:D:148:ALA:O	1:D:152:ARG:HG3	2.11	0.50
1:A:64:TYR:CD2	1:A:104:ASN:HB3	2.46	0.50
1:C:145:VAL:HG23	3:C:2152:HOH:O	2.10	0.50
1:A:268:VAL:HG23	1:A:270:LEU:HD13	1.93	0.50
1:B:79:LEU:H	1:B:79:LEU:HD23	1.76	0.50
1:A:81:ARG:NH2	3:A:2056:HOH:O	2.44	0.50
1:C:268:VAL:HG23	1:C:270:LEU:HD13	1.93	0.50
1:C:79:LEU:HD22	1:C:82:LYS:HE3	1.93	0.50
1:B:191:PRO:O	1:B:194:ASP:HB2	2.12	0.49
1:A:123:ILE:O	1:A:126:HIS:HB3	2.12	0.49
1:D:297:ILE:HB	3:D:2228:HOH:O	2.12	0.49
1:B:200:THR:HG23	1:B:203:LYS:HE2	1.93	0.49
1:C:64:TYR:HB3	1:C:108:HIS:CD2	2.47	0.49
1:C:140:ILE:HG22	1:C:251:VAL:HG12	1.94	0.49
1:B:79:LEU:HD11	1:B:93:LYS:HD3	1.93	0.49
1:D:136:LYS:HE2	1:D:254:GLU:HG2	1.94	0.49
1:B:171:LEU:HD22	1:B:267:PRO:HD3	1.95	0.49
1:D:11:VAL:HG23	1:D:12:GLY:N	2.27	0.49
1:A:127:VAL:O	1:A:130:VAL:HG22	2.13	0.49
1:D:267:PRO:HB2	1:D:279:TYR:HB2	1.95	0.49
1:C:81:ARG:H	1:C:90:LEU:HD22	1.78	0.49
1:C:299:ASP:O	1:C:303:LYS:HG3	2.12	0.49
1:D:262:THR:HB	1:D:299:ASP:OD2	2.13	0.49
1:D:91:LEU:HD11	1:D:123:ILE:HG21	1.95	0.49
1:D:75:ILE:O	1:D:117:VAL:HG22	2.13	0.48
1:D:113:ILE:CG1	1:D:138:ARG:NH1	2.76	0.48
1:B:300:GLU:HA	1:B:303:LYS:NZ	2.27	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:39:GLN:HE22	1:C:62:ASN:ND2	2.10	0.48
1:D:180:PRO:HG3	1:D:205:VAL:HG13	1.94	0.48
1:B:254:GLU:HA	1:B:260:ASP:OD1	2.13	0.48
1:B:199:GLU:H	1:B:199:GLU:CD	2.16	0.48
1:A:88:GLU:O	1:A:92:MET:HG2	2.12	0.48
1:B:117:VAL:O	2:B:1306:NAD:H2N	2.14	0.48
1:D:106:MET:HG3	1:D:132:SER:HA	1.94	0.48
1:D:76:THR:HA	1:D:117:VAL:CG2	2.43	0.48
1:B:171:LEU:CD2	1:B:267:PRO:HD3	2.43	0.48
1:A:303:LYS:C	1:A:305:LEU:H	2.17	0.48
1:B:237:GLU:HG3	1:B:241:LEU:HD12	1.95	0.47
1:A:257:TYR:O	1:A:259:ILE:HG13	2.13	0.47
1:C:18:ARG:HD3	1:C:233:VAL:CG2	2.40	0.47
1:A:32:ASP:OD1	1:A:34:VAL:HG12	2.14	0.47
1:D:34:VAL:CG1	1:D:37:ILE:HD12	2.36	0.47
1:B:21:GLU:CD	1:C:18:ARG:HH12	2.18	0.47
1:B:225:PHE:C	1:B:228:PRO:HD2	2.34	0.47
1:A:257:TYR:CZ	1:A:280:GLU:HA	2.49	0.47
1:D:79:LEU:HD11	1:D:93:LYS:HD3	1.96	0.47
1:B:227:ALA:HB3	1:B:228:PRO:HD3	1.96	0.47
1:A:80:PRO:HA	2:A:1306:NAD:O2D	2.14	0.47
1:B:154:PHE:HB3	1:B:204:LEU:HD22	1.97	0.47
1:D:259:ILE:CG2	1:D:262:THR:HG23	2.45	0.47
1:D:117:VAL:O	2:D:1306:NAD:H2N	2.15	0.46
1:D:223:SER:HB2	3:D:2243:HOH:O	2.16	0.46
1:D:73:VAL:HG11	1:D:105:ILE:HG21	1.97	0.46
1:D:268:VAL:HG23	1:D:270:LEU:HD13	1.97	0.46
1:D:103:ASP:O	1:D:107:LYS:HG3	2.15	0.46
1:C:253:LEU:HD22	1:C:256:GLN:HG3	1.96	0.46
1:D:225:PHE:C	1:D:228:PRO:HD2	2.36	0.46
1:A:225:PHE:C	1:A:228:PRO:HD2	2.36	0.46
1:C:99:LYS:HG3	1:C:100:GLU:N	2.31	0.46
1:A:10:ASN:HB3	1:A:225:PHE:CG	2.50	0.46
1:D:120:PRO:O	1:D:124:MET:HG2	2.15	0.46
1:B:26:ARG:HD3	1:D:242:ASP:OD2	2.15	0.46
1:C:142:MET:C	1:C:142:MET:SD	2.94	0.46
1:A:86:THR:HB	1:A:89:ASP:HB2	1.96	0.46
1:B:90:LEU:HD13	3:B:2113:HOH:O	2.16	0.45
1:B:136:LYS:HE2	1:B:254:GLU:CG	2.46	0.45
1:C:78:GLY:O	1:C:79:LEU:HB2	2.16	0.45
1:D:31:LEU:HD23	1:D:32:ASP:N	2.31	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:80:PRO:O	1:A:81:ARG:HG2	2.16	0.45
1:C:11:VAL:HG12	2:C:1306:NAD:PN	2.56	0.45
1:D:191:PRO:O	1:D:194:ASP:HB2	2.16	0.45
1:A:182:VAL:HG22	3:A:2113:HOH:O	2.17	0.45
1:C:81:ARG:HA	1:C:85:MET:HB2	1.98	0.45
1:B:91:LEU:HD11	1:B:123:ILE:HG21	1.98	0.45
1:C:49:GLY:N	1:C:50:PRO:CD	2.80	0.45
1:A:95:ALA:O	1:A:99:LYS:HG3	2.17	0.45
1:A:35:GLU:HA	1:A:62:ASN:HD21	1.81	0.45
1:A:256:GLN:H	1:A:256:GLN:CD	2.20	0.45
1:D:219:LEU:C	1:D:221:GLN:H	2.20	0.45
1:D:200:THR:O	1:D:203:LYS:HG2	2.17	0.45
1:A:43:LEU:HD22	1:A:47:GLU:HG3	1.99	0.45
1:D:127:VAL:O	1:D:131:ARG:HG2	2.17	0.44
1:B:284:ASP:OD1	1:B:284:ASP:C	2.56	0.44
1:C:106:MET:HG3	1:C:132:SER:HA	1.98	0.44
1:C:37:ILE:N	1:C:37:ILE:HD12	2.32	0.44
1:D:18:ARG:HD3	1:D:233:VAL:CG2	2.42	0.44
1:A:254:GLU:O	1:A:254:GLU:HG3	2.17	0.44
1:C:259:ILE:CG2	1:C:262:THR:HG23	2.47	0.44
1:D:171:LEU:HD22	1:D:267:PRO:HD3	1.99	0.44
1:B:126:HIS:O	1:B:130:VAL:HG12	2.18	0.44
1:C:219:LEU:HD22	3:C:2018:HOH:O	2.18	0.44
1:D:35:GLU:HA	1:D:62:ASN:HD21	1.83	0.44
1:A:79:LEU:HD23	1:A:80:PRO:N	2.33	0.44
1:C:81:ARG:HG2	1:C:90:LEU:CD2	2.41	0.44
1:B:7:GLY:O	1:B:12:GLY:HA3	2.18	0.44
1:A:171:LEU:HD11	1:A:281:ILE:HD11	2.00	0.43
1:D:184:TYR:CD1	1:D:281:ILE:HD12	2.53	0.43
1:D:4:THR:OG1	1:D:29:VAL:HG13	2.19	0.43
1:C:82:LYS:H	1:C:85:MET:HG3	1.83	0.43
1:A:19:LEU:HD13	1:A:233:VAL:HG22	2.01	0.43
1:D:184:TYR:HD1	1:D:281:ILE:HD12	1.83	0.43
1:A:64:TYR:HB3	1:A:108:HIS:CD2	2.54	0.43
1:B:111:ASN:HD21	1:B:138:ARG:CD	2.32	0.43
1:C:81:ARG:HH11	1:C:81:ARG:CB	2.20	0.43
1:C:254:GLU:N	1:C:256:GLN:OE1	2.45	0.43
1:C:82:LYS:HB3	1:C:83:PRO:CD	2.49	0.43
1:D:113:ILE:CG1	1:D:138:ARG:HH12	2.29	0.43
1:D:76:THR:HA	1:D:117:VAL:HG22	2.01	0.43
1:C:253:LEU:C	1:C:260:ASP:O	2.57	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:35:GLU:HA	1:C:62:ASN:HD21	1.84	0.43
1:C:171:LEU:HD11	1:C:281:ILE:HD11	2.01	0.42
1:C:300:GLU:HA	1:C:303:LYS:CE	2.49	0.42
1:A:227:ALA:HB3	1:A:228:PRO:HD3	2.00	0.42
1:B:73:VAL:HG11	1:B:105:ILE:CD1	2.49	0.42
1:D:264:VAL:CG1	1:D:291:LEU:HD11	2.49	0.42
1:B:171:LEU:HD22	1:B:266:VAL:HA	2.00	0.42
1:A:110:LYS:C	1:A:112:PRO:HD3	2.39	0.42
1:C:288:LEU:O	1:C:292:GLN:HG3	2.20	0.42
1:D:201:ILE:O	1:D:205:VAL:HG23	2.20	0.42
1:B:296:LYS:O	1:B:300:GLU:HG3	2.19	0.42
1:C:43:LEU:HD22	1:C:47:GLU:HG3	2.01	0.42
1:A:89:ASP:O	1:A:93:LYS:HG2	2.20	0.42
1:C:31:LEU:HA	1:C:61:SER:O	2.19	0.42
1:D:197:PRO:HA	3:D:2164:HOH:O	2.19	0.42
1:C:171:LEU:HB3	1:C:265:GLY:O	2.20	0.42
1:D:227:ALA:N	1:D:228:PRO:CD	2.83	0.41
1:C:37:ILE:H	1:C:37:ILE:HD12	1.85	0.41
1:B:201:ILE:O	1:B:205:VAL:HG23	2.20	0.41
1:D:130:VAL:HG23	1:D:131:ARG:N	2.34	0.41
1:D:183:LYS:HE2	1:D:184:TYR:CZ	2.55	0.41
1:A:49:GLY:N	1:A:50:PRO:CD	2.83	0.41
1:D:284:ASP:C	1:D:284:ASP:OD1	2.59	0.41
1:C:130:VAL:HG23	1:C:131:ARG:N	2.35	0.41
1:A:79:LEU:C	1:A:79:LEU:HD23	2.41	0.41
1:B:203:LYS:HG3	1:B:204:LEU:N	2.36	0.41
1:C:260:ASP:OD1	1:C:261:LYS:HE2	2.20	0.41
1:D:219:LEU:O	1:D:220:LYS:HB2	2.21	0.41
1:A:219:LEU:O	1:A:220:LYS:HB2	2.20	0.41
1:B:64:TYR:HB3	1:B:108:HIS:CD2	2.56	0.41
1:C:225:PHE:C	1:C:228:PRO:HD2	2.41	0.41
1:B:227:ALA:N	1:B:228:PRO:CD	2.82	0.41
1:D:121:LEU:HD22	1:D:142:MET:O	2.20	0.41
1:D:31:LEU:C	1:D:31:LEU:HD23	2.41	0.41
1:B:267:PRO:HB2	1:B:279:TYR:HB2	2.02	0.41
1:D:181:VAL:HG12	1:D:183:LYS:HG2	2.02	0.41
1:A:26:ARG:HG2	1:A:27:GLU:HG3	2.03	0.41
1:D:249:CYS:O	1:D:265:GLY:HA2	2.20	0.41
1:C:82:LYS:C	1:C:84:GLY:H	2.23	0.41
1:A:136:LYS:HE3	1:A:137:GLU:OE2	2.20	0.41
1:D:183:LYS:HE2	1:D:184:TYR:CE2	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:197:PRO:HG2	1:C:200:THR:OG1	2.21	0.41
1:A:191:PRO:O	1:A:194:ASP:HB2	2.21	0.41
1:C:256:GLN:H	1:C:256:GLN:CD	2.23	0.41
1:A:259:ILE:CG2	1:A:262:THR:CG2	2.98	0.40
2:B:1306:NAD:H6N	3:B:2179:HOH:O	2.21	0.40
1:A:142:MET:C	1:A:142:MET:SD	3.00	0.40
1:B:121:LEU:HD22	1:B:142:MET:O	2.21	0.40
1:D:203:LYS:HG3	1:D:204:LEU:N	2.36	0.40
1:C:117:VAL:O	2:C:1306:NAD:H2N	2.21	0.40
1:D:123:ILE:HG13	3:D:2111:HOH:O	2.22	0.40
1:C:173:GLY:HA2	3:C:2073:HOH:O	2.21	0.40
1:C:125:THR:HG22	1:C:263:PHE:CE2	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	303/310 (98%)	284 (94%)	17 (6%)	2 (1%)	26	19
1	B	292/310 (94%)	286 (98%)	4 (1%)	2 (1%)	26	19
1	C	303/310 (98%)	278 (92%)	20 (7%)	5 (2%)	11	4
1	D	291/310 (94%)	282 (97%)	8 (3%)	1 (0%)	46	41
All	All	1189/1240 (96%)	1130 (95%)	49 (4%)	10 (1%)	24	15

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	81	ARG
1	C	81	ARG
1	A	80	PRO

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Mol	Chain	Res	Type
1	B	143	ALA
1	D	143	ALA
1	C	79	LEU
1	C	83	PRO
1	B	121	LEU
1	C	260	ASP
1	C	84	GLY

### 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	251/255 (98%)	242 (96%)	9 (4%)	42	39
1	B	243/255 (95%)	236 (97%)	7 (3%)	50	49
1	C	251/255 (98%)	241 (96%)	10 (4%)	38	33
1	D	242/255 (95%)	229 (95%)	13 (5%)	27	21
All	All	987/1020 (97%)	948 (96%)	39 (4%)	38	33

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	LEU
1	A	26	ARG
1	A	30	LEU
1	A	43	LEU
1	A	111	ASN
1	A	151	PHE
1	A	256	GLN
1	A	288	LEU
1	A	290	LEU
1	B	19	LEU
1	B	30	LEU
1	B	39	GLN
1	B	43	LEU
1	B	151	PHE

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Mol	Chain	Res	Type
1	B	270	LEU
1	B	288	LEU
1	C	19	LEU
1	C	43	LEU
1	C	81	ARG
1	C	105	ILE
1	C	151	PHE
1	C	256	GLN
1	C	262	THR
1	C	288	LEU
1	C	289	ASP
1	C	300	GLU
1	D	19	LEU
1	D	29	VAL
1	D	30	LEU
1	D	39	GLN
1	D	43	LEU
1	D	138	ARG
1	D	151	PHE
1	D	199	GLU
1	D	206	GLU
1	D	262	THR
1	D	270	LEU
1	D	288	LEU
1	D	299	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	62	ASN
1	A	108	HIS
1	A	111	ASN
1	B	39	GLN
1	B	62	ASN
1	B	111	ASN
1	B	221	GLN
1	B	277	GLN
1	B	301	ASN
1	C	62	ASN
1	C	94	ASN
1	C	108	HIS
1	C	111	ASN

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Mol	Chain	Res	Type
1	C	119	ASN
1	C	301	ASN
1	D	39	GLN
1	D	62	ASN
1	D	94	ASN
1	D	111	ASN
1	D	119	ASN
1	D	282	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

4 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
2	NAD	A	1306	-	38,48,48	2.44	10 (26%)	47,73,73	2.94	16 (34%)
2	NAD	B	1306	-	38,48,48	2.39	11 (28%)	47,73,73	2.93	16 (34%)
2	NAD	C	1306	-	38,48,48	2.47	10 (26%)	47,73,73	2.94	15 (31%)
2	NAD	D	1306	-	38,48,48	2.46	10 (26%)	47,73,73	2.93	16 (34%)

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	1306	-	-	0/22/62/62	0/5/5/5
2	NAD	B	1306	-	-	0/22/62/62	0/5/5/5
2	NAD	C	1306	-	-	0/22/62/62	0/5/5/5
2	NAD	D	1306	-	-	0/22/62/62	0/5/5/5

All (41) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1306	NAD	O2B-C2B	-3.84	1.33	1.43
2	C	1306	NAD	O2B-C2B	-3.38	1.34	1.43
2	A	1306	NAD	O2B-C2B	-3.30	1.35	1.43
2	D	1306	NAD	O2B-C2B	-3.26	1.35	1.43
2	B	1306	NAD	C3B-C4B	2.03	1.58	1.53
2	C	1306	NAD	O4B-C4B	2.22	1.50	1.45
2	B	1306	NAD	O4B-C4B	2.31	1.50	1.45
2	B	1306	NAD	O4D-C4D	2.48	1.50	1.45
2	A	1306	NAD	O4B-C4B	2.49	1.50	1.45
2	D	1306	NAD	O4B-C4B	2.55	1.50	1.45
2	A	1306	NAD	O4D-C4D	2.57	1.50	1.45
2	D	1306	NAD	O4D-C4D	2.58	1.51	1.45
2	C	1306	NAD	O4D-C4D	2.75	1.51	1.45
2	B	1306	NAD	O3B-C3B	3.42	1.51	1.43
2	C	1306	NAD	O3B-C3B	3.43	1.51	1.43
2	B	1306	NAD	C6N-N1N	3.44	1.44	1.35
2	D	1306	NAD	C6N-N1N	3.52	1.44	1.35
2	A	1306	NAD	O3B-C3B	3.54	1.51	1.43
2	B	1306	NAD	C4A-N3A	3.61	1.40	1.35
2	D	1306	NAD	O3B-C3B	3.73	1.51	1.43
2	A	1306	NAD	C4A-N3A	3.73	1.41	1.35
2	C	1306	NAD	C6N-N1N	3.89	1.45	1.35
2	A	1306	NAD	C6N-N1N	3.89	1.45	1.35
2	C	1306	NAD	C4A-N3A	3.98	1.41	1.35
2	D	1306	NAD	C4A-N3A	4.08	1.41	1.35
2	B	1306	NAD	C7N-N7N	4.88	1.42	1.33
2	D	1306	NAD	C7N-N7N	4.93	1.43	1.33
2	C	1306	NAD	C7N-N7N	4.99	1.43	1.33
2	B	1306	NAD	C4N-C3N	5.01	1.47	1.39
2	D	1306	NAD	C4N-C3N	5.11	1.48	1.39
2	A	1306	NAD	C7N-N7N	5.11	1.43	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1306	NAD	C4N-C3N	5.37	1.48	1.39
2	C	1306	NAD	C4N-C3N	5.39	1.48	1.39
2	B	1306	NAD	C2A-N3A	5.55	1.42	1.32
2	D	1306	NAD	C2A-N3A	5.87	1.42	1.32
2	C	1306	NAD	C2A-N3A	6.14	1.43	1.32
2	A	1306	NAD	O4D-C1D	6.16	1.49	1.41
2	B	1306	NAD	O4D-C1D	6.16	1.49	1.41
2	A	1306	NAD	C2A-N3A	6.25	1.43	1.32
2	D	1306	NAD	O4D-C1D	6.48	1.49	1.41
2	C	1306	NAD	O4D-C1D	6.63	1.49	1.41

All (63) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	1306	NAD	N3A-C2A-N1A	-11.38	120.18	128.89
2	B	1306	NAD	N3A-C2A-N1A	-11.10	120.39	128.89
2	C	1306	NAD	N3A-C2A-N1A	-11.03	120.45	128.89
2	A	1306	NAD	N3A-C2A-N1A	-10.96	120.50	128.89
2	A	1306	NAD	C5N-C4N-C3N	-5.72	113.15	120.33
2	B	1306	NAD	C5N-C4N-C3N	-5.70	113.17	120.33
2	C	1306	NAD	C5N-C4N-C3N	-5.65	113.23	120.33
2	D	1306	NAD	C5N-C4N-C3N	-5.60	113.29	120.33
2	A	1306	NAD	C3N-C7N-N7N	-5.17	112.16	117.82
2	B	1306	NAD	C3N-C7N-N7N	-5.14	112.19	117.82
2	C	1306	NAD	C3N-C7N-N7N	-5.12	112.22	117.82
2	D	1306	NAD	C3N-C7N-N7N	-5.00	112.34	117.82
2	D	1306	NAD	C4B-O4B-C1B	-3.63	105.73	109.72
2	B	1306	NAD	C4B-O4B-C1B	-3.62	105.74	109.72
2	C	1306	NAD	C1B-N9A-C4A	-3.52	121.63	126.94
2	A	1306	NAD	C4B-O4B-C1B	-3.48	105.89	109.72
2	A	1306	NAD	C1B-N9A-C4A	-3.37	121.86	126.94
2	B	1306	NAD	C1B-N9A-C4A	-3.32	121.94	126.94
2	C	1306	NAD	C4B-O4B-C1B	-3.26	106.13	109.72
2	D	1306	NAD	C1B-N9A-C4A	-3.05	122.33	126.94
2	D	1306	NAD	C2B-C1B-N9A	-2.12	111.05	114.29
2	B	1306	NAD	C2B-C1B-N9A	-2.05	111.16	114.29
2	B	1306	NAD	C2D-C3D-C4D	2.03	106.78	102.61
2	C	1306	NAD	C2D-C3D-C4D	2.11	106.95	102.61
2	D	1306	NAD	C2D-C3D-C4D	2.13	107.00	102.61
2	A	1306	NAD	C4D-O4D-C1D	2.17	112.11	109.72
2	A	1306	NAD	C2D-C3D-C4D	2.26	107.26	102.61
2	D	1306	NAD	O3-PN-O5D	2.31	109.06	102.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1306	NAD	N6A-C6A-N1A	2.39	124.33	119.20
2	B	1306	NAD	O3-PN-O5D	2.44	109.40	102.94
2	B	1306	NAD	N6A-C6A-N1A	2.48	124.54	119.20
2	C	1306	NAD	O3-PN-O5D	2.49	109.55	102.94
2	C	1306	NAD	N6A-C6A-N1A	2.51	124.58	119.20
2	A	1306	NAD	O3-PN-O5D	2.52	109.62	102.94
2	D	1306	NAD	N6A-C6A-N1A	2.57	124.71	119.20
2	D	1306	NAD	C2B-C3B-C4B	2.68	108.12	102.61
2	B	1306	NAD	C2B-C3B-C4B	2.69	108.15	102.61
2	B	1306	NAD	PN-O3-PA	2.82	140.65	132.73
2	A	1306	NAD	C2B-C3B-C4B	2.89	108.55	102.61
2	C	1306	NAD	C2B-C3B-C4B	2.94	108.67	102.61
2	D	1306	NAD	PN-O3-PA	3.05	141.30	132.73
2	B	1306	NAD	O2B-C2B-C3B	3.07	121.81	111.83
2	C	1306	NAD	O2B-C2B-C3B	3.12	121.99	111.83
2	A	1306	NAD	O2B-C2B-C3B	3.15	122.08	111.83
2	D	1306	NAD	O2B-C2B-C3B	3.18	122.17	111.83
2	A	1306	NAD	PN-O3-PA	3.56	142.74	132.73
2	C	1306	NAD	PN-O3-PA	3.67	143.03	132.73
2	D	1306	NAD	C6N-C5N-C4N	4.03	125.53	119.44
2	C	1306	NAD	C6N-C5N-C4N	4.08	125.61	119.44
2	B	1306	NAD	C6N-C5N-C4N	4.16	125.73	119.44
2	A	1306	NAD	C6N-C5N-C4N	4.17	125.75	119.44
2	B	1306	NAD	C2N-C3N-C4N	4.42	123.21	118.29
2	D	1306	NAD	C2N-C3N-C4N	4.47	123.27	118.29
2	C	1306	NAD	C2N-C3N-C4N	4.63	123.45	118.29
2	A	1306	NAD	C2N-C3N-C4N	4.66	123.48	118.29
2	C	1306	NAD	O7N-C7N-C3N	6.35	126.52	119.59
2	D	1306	NAD	C2A-N1A-C6A	6.38	130.17	118.77
2	A	1306	NAD	C2A-N1A-C6A	6.40	130.21	118.77
2	D	1306	NAD	O7N-C7N-C3N	6.41	126.59	119.59
2	B	1306	NAD	O7N-C7N-C3N	6.42	126.60	119.59
2	A	1306	NAD	O7N-C7N-C3N	6.45	126.63	119.59
2	B	1306	NAD	C2A-N1A-C6A	6.51	130.40	118.77
2	C	1306	NAD	C2A-N1A-C6A	6.52	130.41	118.77

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1306	NAD	1	0
2	B	1306	NAD	2	0
2	C	1306	NAD	4	0
2	D	1306	NAD	3	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	305/310 (98%)	0.23	9 (2%) 54 55	14, 30, 52, 65	0
1	B	296/310 (95%)	0.03	12 (4%) 41 42	12, 22, 42, 63	0
1	C	305/310 (98%)	0.33	12 (3%) 43 45	15, 30, 53, 70	0
1	D	295/310 (95%)	0.12	10 (3%) 49 50	13, 24, 42, 51	0
All	All	1201/1240 (96%)	0.18	43 (3%) 46 48	12, 26, 46, 70	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	80	PRO	7.1
1	C	83	PRO	7.1
1	A	80	PRO	5.3
1	C	79	LEU	4.8
1	B	89	ASP	4.0
1	B	220	LYS	3.8
1	D	221	GLN	3.8
1	B	79	LEU	3.7
1	A	79	LEU	3.5
1	B	90	LEU	3.4
1	C	304	MET	3.3
1	B	221	GLN	3.3
1	A	82	LYS	3.1
1	C	81	ARG	3.1
1	B	303	LYS	2.9
1	B	304	MET	2.9
1	C	221	GLN	2.8
1	A	81	ARG	2.8
1	D	220	LYS	2.7
1	B	294	SER	2.6
1	C	82	LYS	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	217	GLU	2.6
1	D	297	ILE	2.6
1	A	304	MET	2.5
1	C	296	LYS	2.5
1	B	176	ASP	2.5
1	C	303	LYS	2.5
1	B	92	MET	2.5
1	D	279	TYR	2.5
1	D	75	ILE	2.4
1	C	115	ILE	2.3
1	A	285	GLN	2.3
1	C	279	TYR	2.3
1	C	261	LYS	2.2
1	A	121	LEU	2.2
1	A	305	LEU	2.1
1	A	111	ASN	2.1
1	D	290	LEU	2.1
1	B	297	ILE	2.1
1	D	217	GLU	2.1
1	D	293	LYS	2.1
1	D	198	ALA	2.0
1	D	219	LEU	2.0

## 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 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, 95<sup>th</sup> 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( $\text{\AA}^2$ )	Q<0.9
2	NAD	C	1306	44/44	0.84	0.17	0.19	41,47,48,49	0
2	NAD	A	1306	44/44	0.88	0.15	0.01	42,48,51,51	0
2	NAD	B	1306	44/44	0.96	0.09	-0.69	17,22,30,31	0
2	NAD	D	1306	44/44	0.95	0.10	-0.78	15,24,30,32	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.