



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:14 AM GMT

PDB ID : 2G5C
Title : Crystal Structure of Prephenate Dehydrogenase from Aquifex aeolicus
Authors : Sun, W.; Singh, S.; Zhang, R.; Turnbull, J.L.; Christendat, D.
Deposited on : 2006-02-22
Resolution : 1.90 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<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

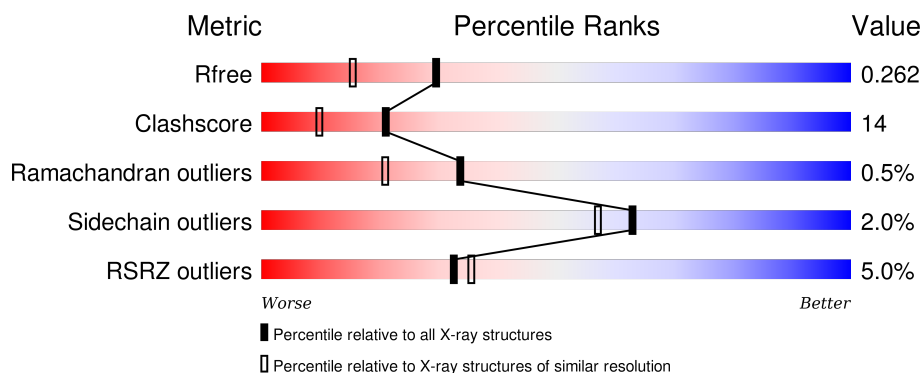
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	4755 (1.90-1.90)
Clashscore	102246	5398 (1.90-1.90)
Ramachandran outliers	100387	5338 (1.90-1.90)
Sidechain outliers	100360	5339 (1.90-1.90)
RSRZ outliers	91569	4766 (1.90-1.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	281	<div> <div>7%</div> <div>67% 30% ..</div> </div>
1	B	281	<div> <div>8%</div> <div>73% 25% .</div> </div>
1	C	281	<div> <div>3%</div> <div>72% 25% ..</div> </div>
1	D	281	<div> <div>%</div> <div>74% 23% ..</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9487 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 prephenate dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	278	Total	C	N	O	Se	0	0	0
			2198	1418	360	412	8			
1	B	280	Total	C	N	O	Se	0	0	0
			2216	1429	365	415	7			
1	C	278	Total	C	N	O	Se	0	0	0
			2198	1418	363	410	7			
1	D	278	Total	C	N	O	Se	0	0	0
			2198	1418	363	410	7			

There are 32 discrepancies between the modelled and reference sequences:

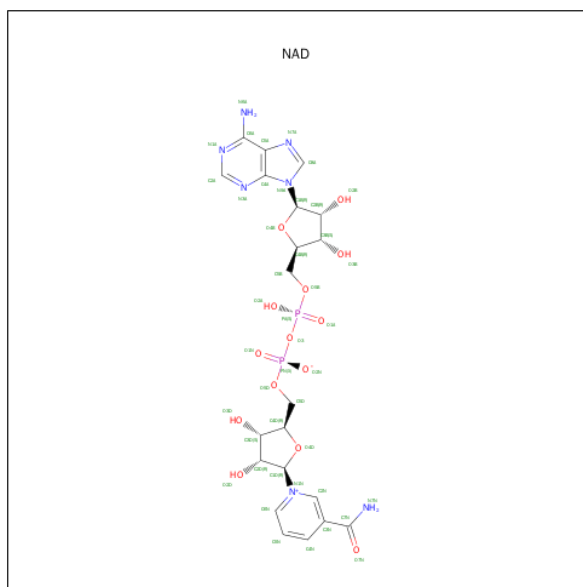
Chain	Residue	Modelled	Actual	Comment	Reference
A	30	MSE	-	INITIATING METHIONINE	UNP 067636
A	41	MSE	MET	MODIFIED RESIDUE	UNP 067636
A	96	MSE	MET	MODIFIED RESIDUE	UNP 067636
A	200	MSE	MET	MODIFIED RESIDUE	UNP 067636
A	230	MSE	MET	MODIFIED RESIDUE	UNP 067636
A	258	MSE	MET	MODIFIED RESIDUE	UNP 067636
A	271	MSE	MET	MODIFIED RESIDUE	UNP 067636
A	308	MSE	MET	MODIFIED RESIDUE	UNP 067636
B	30	MSE	-	INITIATING METHIONINE	UNP 067636
B	41	MSE	MET	MODIFIED RESIDUE	UNP 067636
B	96	MSE	MET	MODIFIED RESIDUE	UNP 067636
B	200	MSE	MET	MODIFIED RESIDUE	UNP 067636
B	230	MSE	MET	MODIFIED RESIDUE	UNP 067636
B	258	MSE	MET	MODIFIED RESIDUE	UNP 067636
B	271	MSE	MET	MODIFIED RESIDUE	UNP 067636
B	308	MSE	MET	MODIFIED RESIDUE	UNP 067636
C	30	MSE	-	INITIATING METHIONINE	UNP 067636
C	41	MSE	MET	MODIFIED RESIDUE	UNP 067636
C	96	MSE	MET	MODIFIED RESIDUE	UNP 067636
C	200	MSE	MET	MODIFIED RESIDUE	UNP 067636
C	230	MSE	MET	MODIFIED RESIDUE	UNP 067636

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Chain	Residue	Modelled	Actual	Comment	Reference
C	258	MSE	MET	MODIFIED RESIDUE	UNP O67636
C	271	MSE	MET	MODIFIED RESIDUE	UNP O67636
C	308	MSE	MET	MODIFIED RESIDUE	UNP O67636
D	30	MSE	-	INITIATING METHIONINE	UNP O67636
D	41	MSE	MET	MODIFIED RESIDUE	UNP O67636
D	96	MSE	MET	MODIFIED RESIDUE	UNP O67636
D	200	MSE	MET	MODIFIED RESIDUE	UNP O67636
D	230	MSE	MET	MODIFIED RESIDUE	UNP O67636
D	258	MSE	MET	MODIFIED RESIDUE	UNP O67636
D	271	MSE	MET	MODIFIED RESIDUE	UNP O67636
D	308	MSE	MET	MODIFIED RESIDUE	UNP O67636

- 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		

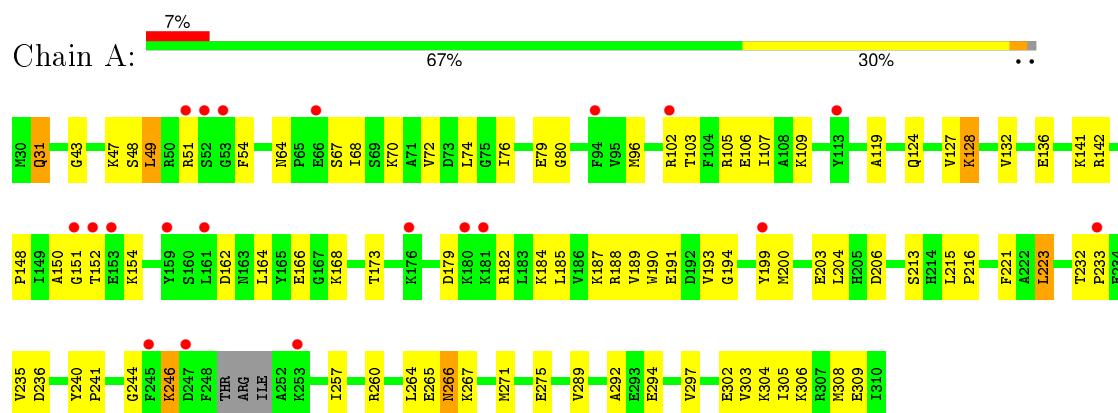
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	84	Total 84	O 84	0	0
3	B	88	Total 88	O 88	0	0
3	C	169	Total 169	O 169	0	0
3	D	160	Total 160	O 160	0	0

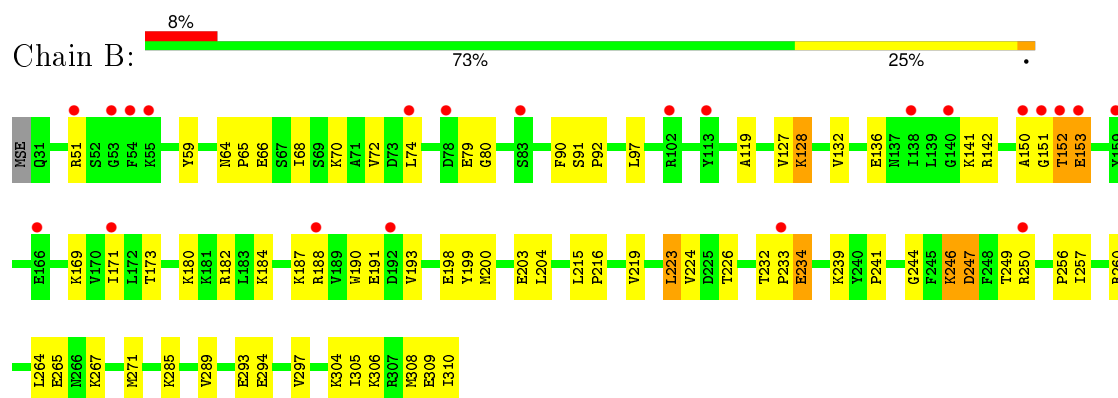
3 Residue-property plots

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.

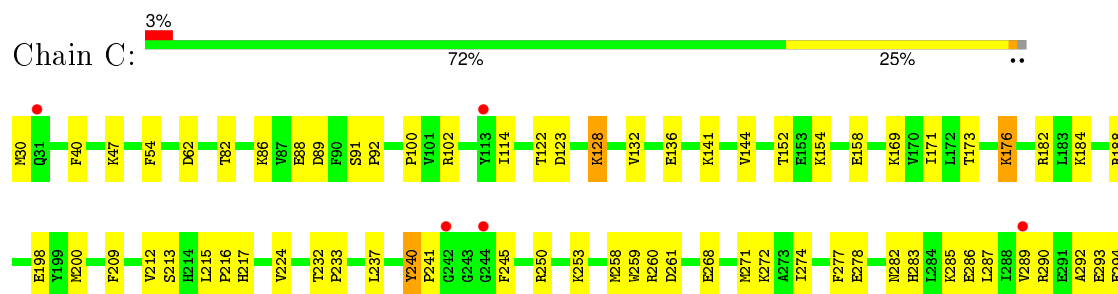
- Molecule 1: prephenate dehydrogenase

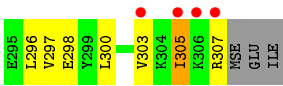


- Molecule 1: prephenate dehydrogenase

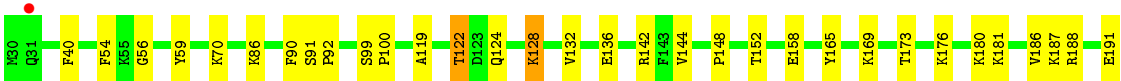


- Molecule 1: prephenate dehydrogenase





● Molecule 1: prephenate dehydrogenase



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4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	42.71 Å 178.95 Å 75.19 Å 90.00° 99.15° 90.00°	Depositor
Resolution (Å)	29.93 – 1.90 34.43 – 1.89	Depositor EDS
% Data completeness (in resolution range)	88.2 (29.93-1.90) 93.4 (34.43-1.89)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.70 (at 1.89 Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.220 , 0.261 0.221 , 0.262	Depositor DCC
R_{free} test set	4738 reflections (5.79%)	DCC
Wilson B-factor (Å ²)	21.7	Xtriage
Anisotropy	0.107	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 53.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Outliers	0 of 160792 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9487	wwPDB-VP
Average B, all atoms (Å ²)	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 74.78 % 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 1.4331e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² 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	0.31	0/2233	0.53	0/2990
1	B	0.31	0/2252	0.53	0/3018
1	C	0.34	0/2235	0.59	0/2998
1	D	0.34	0/2235	0.59	0/2998
All	All	0.32	0/8955	0.56	0/12004

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	2198	0	2241	80	0
1	B	2216	0	2264	75	0
1	C	2198	0	2247	72	0
1	D	2198	0	2247	69	0
2	A	44	0	26	0	0
2	B	44	0	26	0	0
2	C	44	0	26	1	0
2	D	44	0	26	1	0
3	A	84	0	0	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	88	0	0	3	0
3	C	169	0	0	9	0
3	D	160	0	0	7	0
All	All	9487	0	9103	259	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 14.

All (259) 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:305:ILE:HG22	1:D:307:ARG:H	1.32	0.94
1:A:184:LYS:HB3	1:A:188:ARG:HH12	1.32	0.93
1:C:290:ARG:HD2	1:C:292:ALA:HB2	1.51	0.92
1:A:127:VAL:HG21	1:A:265:GLU:HG2	1.50	0.92
1:B:260:ARG:HD2	1:D:293:GLU:HG3	1.55	0.86
1:A:150:ALA:HB2	1:A:168:LYS:HG2	1.56	0.85
1:A:51:ARG:NH1	1:A:51:ARG:HB3	1.91	0.85
1:C:169:LYS:HD3	1:C:198:GLU:OE2	1.77	0.84
1:A:51:ARG:HH11	1:A:51:ARG:HB3	1.43	0.82
1:A:48:SER:HA	1:A:51:ARG:HH12	1.43	0.81
1:C:122:THR:HG23	1:C:144:VAL:O	1.80	0.81
1:B:51:ARG:NH1	1:B:51:ARG:HB3	2.00	0.76
1:D:283:HIS:O	1:D:286:GLU:HG3	1.87	0.75
1:C:305:ILE:HG22	1:C:307:ARG:H	1.52	0.74
1:A:236:ASP:OD2	1:C:307:ARG:HD3	1.87	0.74
1:B:305:ILE:O	1:B:309:GLU:HG2	1.88	0.74
1:D:232:THR:HB	1:D:233:PRO:HD2	1.70	0.73
1:D:237:LEU:O	1:D:241:PRO:HD3	1.89	0.73
1:B:260:ARG:O	1:B:264:LEU:HG	1.89	0.73
1:B:260:ARG:HD3	1:D:297:VAL:CG2	2.19	0.73
1:B:260:ARG:HD2	1:D:293:GLU:CG	2.19	0.72
1:D:119:ALA:O	1:D:142:ARG:HG2	1.90	0.72
1:A:200:MSE:CE	1:A:204:LEU:HG	2.19	0.71
1:D:231:SER:O	1:D:235:VAL:HG13	1.91	0.71
1:A:271:MSE:HE1	1:C:285:LYS:CG	2.21	0.71
1:B:187:LYS:HD2	1:B:199:TYR:OH	1.91	0.70
1:B:150:ALA:O	1:B:152:THR:HG22	1.91	0.70
1:C:158:GLU:CD	1:C:158:GLU:H	1.95	0.70
1:B:51:ARG:HH11	1:B:51:ARG:HB3	1.55	0.69
1:B:271:MSE:HE2	1:D:288:ILE:HG22	1.76	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:48:SER:HA	1:A:51:ARG:NH1	2.07	0.68
1:B:215:LEU:HB3	1:B:216:PRO:HD3	1.74	0.68
1:C:154:LYS:HD2	1:C:154:LYS:N	2.08	0.68
1:A:184:LYS:HB3	1:A:188:ARG:NH1	2.07	0.68
1:B:232:THR:HB	1:B:233:PRO:HD2	1.77	0.67
1:A:200:MSE:HE3	1:A:204:LEU:HG	1.77	0.67
1:B:232:THR:OG1	1:B:234:GLU:HG3	1.95	0.67
1:C:176:LYS:HE3	1:C:176:LYS:HA	1.76	0.66
1:A:128:LYS:HE3	1:A:128:LYS:HA	1.79	0.65
1:B:293:GLU:HG2	1:D:260:ARG:HD2	1.78	0.65
1:A:223:LEU:HD22	1:C:277:PHE:CE1	2.31	0.65
1:A:260:ARG:HD3	1:C:297:VAL:CG2	2.27	0.65
1:A:271:MSE:HE1	1:C:285:LYS:HG2	1.79	0.65
1:C:86:LYS:HD3	1:C:86:LYS:O	1.97	0.65
1:D:132:VAL:O	1:D:136:GLU:HG3	1.97	0.64
1:B:127:VAL:HG21	1:B:265:GLU:HG2	1.79	0.64
1:B:223:LEU:HD22	1:D:277:PHE:CE1	2.31	0.64
1:C:215:LEU:HB3	1:C:216:PRO:HD3	1.78	0.64
1:A:74:LEU:HB2	1:A:76:ILE:HG12	1.79	0.63
1:B:256:PRO:HG2	1:D:301:LYS:HG2	1.79	0.63
1:B:66:GLU:O	1:B:70:LYS:HG2	1.99	0.62
1:A:119:ALA:O	1:A:142:ARG:HG2	2.00	0.62
1:A:132:VAL:O	1:A:136:GLU:HG3	2.00	0.61
1:D:181:LYS:HG2	3:D:6726:HOH:O	2.00	0.61
1:C:290:ARG:CD	1:C:292:ALA:HB2	2.29	0.61
1:A:215:LEU:HB3	1:A:216:PRO:HD3	1.83	0.61
1:B:241:PRO:O	1:B:246:LYS:HG3	2.01	0.60
1:C:88:GLU:HG3	1:C:114:ILE:HG23	1.81	0.60
1:A:305:ILE:O	1:A:309:GLU:HG3	2.01	0.60
1:C:132:VAL:O	1:C:136:GLU:HG3	2.02	0.60
1:B:180:LYS:O	1:B:184:LYS:HG3	2.02	0.60
1:D:215:LEU:HB3	1:D:216:PRO:HD3	1.83	0.60
1:B:187:LYS:HG2	1:B:191:GLU:OE2	2.03	0.59
1:D:306:LYS:CE	1:D:306:LYS:HA	2.32	0.59
1:C:122:THR:HG22	1:C:123:ASP:H	1.67	0.59
1:B:91:SER:N	1:B:92:PRO:HD3	2.18	0.58
1:A:166:GLU:OE1	1:A:194:GLY:HA3	2.02	0.58
1:B:190:TRP:O	1:B:193:VAL:HG22	2.04	0.58
1:C:128:LYS:HA	1:C:128:LYS:HE3	1.86	0.58
1:C:286:GLU:HA	1:C:289:VAL:HG22	1.85	0.58
1:D:122:THR:HG23	1:D:144:VAL:O	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:180:LYS:HE3	3:D:6833:HOH:O	2.04	0.57
1:D:70:LYS:HE2	1:D:158:GLU:HG2	1.86	0.57
1:A:203:GLU:CD	1:A:203:GLU:H	2.06	0.57
1:D:187:LYS:O	1:D:191:GLU:HG3	2.03	0.57
1:A:48:SER:CA	1:A:51:ARG:HH12	2.16	0.57
1:B:271:MSE:HE1	1:D:288:ILE:HB	1.87	0.56
1:B:68:ILE:O	1:B:72:VAL:HG23	2.05	0.56
1:C:213:SER:O	1:C:217:HIS:HD2	1.88	0.56
1:A:294:GLU:O	1:A:297:VAL:HG12	2.05	0.56
1:A:105:ARG:O	1:A:109:LYS:HG3	2.06	0.56
1:D:274:ILE:O	1:D:278:GLU:HG3	2.05	0.56
1:D:128:LYS:HE3	1:D:128:LYS:HA	1.88	0.55
3:C:5855:HOH:O	1:D:180:LYS:HE2	2.06	0.55
1:B:260:ARG:HD3	1:D:297:VAL:HG23	1.87	0.55
1:C:100:PRO:HA	2:C:5686:NAD:O3D	2.06	0.55
1:B:271:MSE:HE3	1:D:289:VAL:CG2	2.37	0.55
1:A:68:ILE:O	1:A:72:VAL:HG23	2.07	0.55
1:A:141:LYS:O	1:A:182:ARG:HD2	2.07	0.55
1:B:70:LYS:O	1:B:74:LEU:HG	2.08	0.54
1:C:250:ARG:HD3	3:C:5832:HOH:O	2.07	0.54
1:B:260:ARG:HG3	1:D:296:LEU:HD23	1.90	0.54
1:A:51:ARG:CB	1:A:51:ARG:HH11	2.17	0.54
1:B:271:MSE:HE1	1:D:289:VAL:N	2.23	0.54
1:D:124:GLN:HE21	1:D:148:PRO:HG3	1.71	0.54
1:A:264:LEU:HB2	3:A:3690:HOH:O	2.07	0.54
1:A:187:LYS:HD2	1:A:199:TYR:OH	2.08	0.54
1:D:297:VAL:HG12	1:D:301:LYS:HE2	1.90	0.54
1:B:203:GLU:HG2	1:B:204:LEU:N	2.23	0.53
1:A:303:VAL:HG13	1:C:224:VAL:HB	1.90	0.53
1:C:289:VAL:HG23	1:C:290:ARG:N	2.24	0.53
1:A:124:GLN:HE21	1:A:148:PRO:HG3	1.73	0.53
1:B:246:LYS:HA	1:B:249:THR:HG23	1.89	0.53
1:D:306:LYS:NZ	1:D:306:LYS:HA	2.22	0.53
1:D:306:LYS:HE2	1:D:306:LYS:HA	1.91	0.53
1:B:271:MSE:CE	1:D:289:VAL:HG23	2.39	0.53
1:B:128:LYS:HE3	1:B:128:LYS:HA	1.90	0.52
1:C:300:LEU:O	1:C:303:VAL:HG22	2.10	0.52
1:A:232:THR:HB	1:A:233:PRO:CD	2.40	0.52
1:A:304:LYS:HE3	1:C:253:LYS:O	2.10	0.51
1:C:307:ARG:HD2	3:C:5743:HOH:O	2.09	0.51
1:C:232:THR:HB	1:C:233:PRO:HD2	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:200:MSE:CE	1:B:204:LEU:HG	2.40	0.51
1:B:294:GLU:O	1:B:297:VAL:HG12	2.10	0.51
1:D:284:LEU:O	1:D:288:ILE:HG13	2.10	0.51
1:C:209:PHE:HA	1:C:212:VAL:HG22	1.92	0.51
1:A:64:ASN:ND2	1:A:67:SER:H	2.09	0.51
1:B:142:ARG:HA	1:B:182:ARG:NE	2.26	0.50
1:D:86:LYS:HD2	3:D:6713:HOH:O	2.10	0.50
1:A:150:ALA:CB	1:A:168:LYS:HG2	2.37	0.50
1:A:106:GLU:HG2	3:A:3764:HOH:O	2.10	0.50
1:A:49:LEU:HD12	1:A:54:PHE:CG	2.46	0.50
1:A:304:LYS:HG3	1:C:253:LYS:HA	1.94	0.50
1:A:102:ARG:HH21	1:A:264:LEU:CD2	2.23	0.50
1:A:187:LYS:O	1:A:191:GLU:HG3	2.11	0.50
1:D:188:ARG:HG3	3:D:6753:HOH:O	2.12	0.50
1:D:287:LEU:HD21	1:D:295:GLU:OE2	2.11	0.50
1:C:294:GLU:O	1:C:298:GLU:HG2	2.12	0.50
1:B:153:GLU:HB3	3:B:4716:HOH:O	2.11	0.50
1:D:122:THR:HG21	1:D:186:VAL:HG11	1.94	0.49
1:C:40:PHE:CD1	1:C:152:THR:HG21	2.47	0.49
1:A:152:THR:OG1	1:A:154:LYS:HG3	2.12	0.49
1:B:271:MSE:HE2	1:D:288:ILE:CG2	2.42	0.49
1:A:260:ARG:HD3	1:C:297:VAL:HG23	1.94	0.49
1:B:119:ALA:O	1:B:142:ARG:HG2	2.13	0.49
1:C:237:LEU:O	1:C:241:PRO:HD3	2.13	0.49
1:B:244:GLY:H	1:B:246:LYS:HZ2	1.61	0.49
1:A:151:GLY:HA2	1:A:164:LEU:HD11	1.94	0.49
1:C:91:SER:N	1:C:92:PRO:HD3	2.28	0.48
1:C:217:HIS:HE1	3:C:5795:HOH:O	1.96	0.48
1:C:258:MSE:HE2	1:C:259:TRP:NE1	2.28	0.48
1:D:59:TYR:CD2	1:D:90:PHE:HB3	2.48	0.48
1:D:209:PHE:HA	1:D:212:VAL:HG22	1.94	0.48
1:B:171:ILE:HD12	1:B:171:ILE:N	2.29	0.48
1:C:122:THR:CG2	3:C:5847:HOH:O	2.61	0.48
1:D:70:LYS:HE2	1:D:158:GLU:CG	2.43	0.48
1:C:240:TYR:N	1:C:241:PRO:HD3	2.28	0.48
1:C:240:TYR:N	1:C:241:PRO:CD	2.77	0.47
1:C:283:HIS:O	1:C:286:GLU:HG2	2.13	0.47
1:A:49:LEU:HD12	1:A:54:PHE:HB2	1.97	0.47
1:B:306:LYS:O	1:B:310:ILE:HG23	2.13	0.47
1:A:70:LYS:O	1:A:74:LEU:HG	2.13	0.47
1:B:267:LYS:O	1:B:271:MSE:HG2	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:122:THR:HG22	3:C:5847:HOH:O	2.15	0.47
1:C:296:LEU:O	1:C:300:LEU:HG	2.14	0.47
1:C:47:LYS:NZ	3:C:5844:HOH:O	2.43	0.47
1:C:102:ARG:NH1	1:C:261:ASP:OD2	2.43	0.47
1:B:132:VAL:O	1:B:136:GLU:HG3	2.14	0.47
1:B:234:GLU:HB2	3:B:4731:HOH:O	2.15	0.47
1:B:169:LYS:HD3	1:B:198:GLU:OE2	2.14	0.47
1:D:54:PHE:CZ	1:D:56:GLY:HA3	2.50	0.47
1:A:257:ILE:N	1:A:257:ILE:HD12	2.31	0.46
1:B:239:LYS:O	1:D:169:LYS:HE3	2.15	0.46
1:B:257:ILE:HD12	1:B:257:ILE:H	1.79	0.46
1:B:257:ILE:HD12	1:B:257:ILE:N	2.29	0.46
1:A:213:SER:C	1:A:216:PRO:HD2	2.36	0.46
1:B:187:LYS:O	1:B:191:GLU:HG3	2.16	0.46
1:A:246:LYS:HD3	1:A:246:LYS:H	1.80	0.46
1:C:173:THR:HA	1:C:200:MSE:O	2.16	0.46
1:C:268:GLU:CD	1:C:272:LYS:HE3	2.36	0.46
1:B:219:VAL:HG11	1:D:219:VAL:CG1	2.46	0.46
1:A:232:THR:HB	1:A:233:PRO:HD2	1.97	0.46
1:B:141:LYS:O	1:B:182:ARG:HD2	2.16	0.45
1:B:224:VAL:HG21	1:D:303:VAL:HG12	1.97	0.45
1:C:141:LYS:O	1:C:182:ARG:HD2	2.16	0.45
1:B:184:LYS:O	1:B:188:ARG:HG3	2.16	0.45
1:D:91:SER:N	1:D:92:PRO:HD3	2.31	0.45
1:B:59:TYR:HB3	1:B:90:PHE:CD1	2.52	0.45
1:B:271:MSE:CE	1:D:289:VAL:N	2.79	0.45
1:A:96:MSE:SE	1:A:124:GLN:HG3	2.66	0.45
1:B:246:LYS:HA	1:B:249:THR:CG2	2.46	0.45
1:A:43:GLY:O	1:A:47:LYS:HG3	2.17	0.45
1:A:179:ASP:OD2	1:A:182:ARG:HG3	2.17	0.45
1:A:102:ARG:HH21	1:A:264:LEU:HD23	1.82	0.45
1:A:275:GLU:OE1	1:A:275:GLU:HA	2.17	0.45
1:B:250:ARG:HB2	1:B:250:ARG:NH1	2.31	0.45
1:A:79:GLU:HG2	1:A:80:GLY:N	2.31	0.45
1:D:40:PHE:CD2	1:D:152:THR:HG21	2.52	0.45
1:C:184:LYS:HE3	3:C:5708:HOH:O	2.16	0.45
1:C:188:ARG:HG2	3:C:5744:HOH:O	2.17	0.45
1:A:240:TYR:HE2	1:C:198:GLU:OE2	1.99	0.44
1:A:244:GLY:N	1:A:246:LYS:HZ2	2.16	0.44
1:A:190:TRP:O	1:A:193:VAL:HG22	2.16	0.44
1:B:285:LYS:O	1:B:289:VAL:HG23	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:31:GLN:OE1	1:A:31:GLN:HA	2.17	0.44
1:C:286:GLU:O	1:C:289:VAL:HG22	2.17	0.44
1:A:271:MSE:HE1	1:C:285:LYS:HG3	1.97	0.44
1:D:208:VAL:O	1:D:212:VAL:HG22	2.17	0.44
1:C:260:ARG:HD2	1:C:260:ARG:C	2.36	0.44
1:C:305:ILE:C	1:C:307:ARG:N	2.70	0.44
1:A:267:LYS:O	1:A:271:MSE:HG2	2.18	0.44
1:C:305:ILE:C	1:C:307:ARG:H	2.21	0.44
1:B:271:MSE:CE	1:D:288:ILE:HB	2.46	0.44
1:A:49:LEU:HD12	1:A:54:PHE:CB	2.48	0.44
1:B:203:GLU:HG2	1:B:204:LEU:H	1.81	0.44
1:A:304:LYS:O	1:A:308:MSE:HG3	2.17	0.44
1:D:240:TYR:N	1:D:241:PRO:CD	2.81	0.43
1:C:154:LYS:CD	1:C:154:LYS:N	2.80	0.43
1:B:250:ARG:HB2	1:B:250:ARG:HH11	1.83	0.43
1:D:124:GLN:O	2:D:6686:NAD:H6N	2.18	0.43
1:C:241:PRO:HB3	1:C:245:PHE:CD2	2.53	0.43
1:C:274:ILE:O	1:C:278:GLU:HG3	2.18	0.43
1:A:271:MSE:O	1:A:275:GLU:HG2	2.18	0.43
1:D:283:HIS:HA	3:D:6728:HOH:O	2.18	0.43
1:B:173:THR:HA	1:B:200:MSE:O	2.18	0.43
1:D:99:SER:HB2	1:D:100:PRO:CD	2.49	0.43
1:B:271:MSE:HE1	1:D:285:LYS:O	2.19	0.43
1:B:226:THR:OG1	1:D:277:PHE:HA	2.18	0.42
1:B:244:GLY:O	1:B:247:ASP:HB2	2.19	0.42
1:A:102:ARG:NH2	1:C:293:GLU:OE2	2.51	0.42
1:D:122:THR:CG2	1:D:144:VAL:O	2.66	0.42
1:B:200:MSE:HE3	1:B:204:LEU:HG	2.01	0.42
1:A:51:ARG:NH1	1:A:162:ASP:O	2.52	0.42
1:A:241:PRO:O	1:A:246:LYS:HD2	2.18	0.42
1:D:173:THR:HA	1:D:200:MSE:O	2.19	0.42
1:C:233:PRO:HB3	3:D:6774:HOH:O	2.19	0.42
1:B:68:ILE:HG23	1:B:80:GLY:HA3	2.00	0.42
1:D:297:VAL:O	1:D:301:LYS:HG3	2.20	0.42
1:C:62:ASP:O	1:C:82:THR:HA	2.19	0.42
1:A:302:GLU:OE2	1:A:306:LYS:HE2	2.19	0.42
1:A:221:PHE:CD1	1:C:303:VAL:HG11	2.55	0.42
1:A:244:GLY:HA2	1:A:246:LYS:HZ3	1.84	0.42
1:C:188:ARG:HH11	1:C:188:ARG:HG3	1.85	0.41
1:C:282:ASN:HA	1:C:282:ASN:HD22	1.67	0.41
1:C:86:LYS:O	1:C:89:ASP:HB2	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:206:ASP:HB3	1:D:266:ASN:HD21	1.83	0.41
1:A:289:VAL:HG22	1:C:271:MSE:SE	2.70	0.41
1:A:185:LEU:O	1:A:189:VAL:HG23	2.21	0.41
1:D:124:GLN:NE2	1:D:165:TYR:OH	2.51	0.41
1:C:122:THR:HG22	1:C:123:ASP:N	2.34	0.41
1:B:271:MSE:HE3	1:D:289:VAL:HG23	2.01	0.41
1:C:171:ILE:HG13	1:C:209:PHE:CE2	2.56	0.41
1:A:265:GLU:HB2	3:A:3761:HOH:O	2.19	0.41
1:A:103:THR:O	1:A:107:ILE:HG13	2.21	0.41
1:A:206:ASP:HB3	1:A:266:ASN:HD21	1.86	0.41
1:D:176:LYS:HE2	3:D:6807:HOH:O	2.20	0.41
1:B:304:LYS:O	1:B:308:MSE:HG3	2.20	0.41
1:D:239:LYS:HG2	1:D:239:LYS:O	2.21	0.41
1:B:271:MSE:SE	1:D:285:LYS:HG3	2.72	0.40
1:B:79:GLU:HG2	1:B:80:GLY:N	2.35	0.40
1:C:268:GLU:OE1	1:C:272:LYS:HE3	2.21	0.40
1:C:30:MSE:HE3	1:C:54:PHE:CE1	2.56	0.40
1:A:64:ASN:ND2	1:A:67:SER:N	2.70	0.40
1:D:240:TYR:H	1:D:241:PRO:HD3	1.86	0.40
1:A:173:THR:HA	1:A:200:MSE:O	2.22	0.40
1:B:152:THR:HA	3:B:4745:HOH:O	2.21	0.40
1:A:292:ALA:HB2	3:A:3700:HOH:O	2.21	0.40
1:D:282:ASN:HA	1:D:282:ASN:HD22	1.70	0.40
1:B:64:ASN:HA	1:B:65:PRO:HD3	1.95	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	274/281 (98%)	254 (93%)	18 (7%)	2 (1%)	26 14

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	278/281 (99%)	259 (93%)	17 (6%)	2 (1%)	26	14
1	C	276/281 (98%)	264 (96%)	10 (4%)	2 (1%)	26	14
1	D	276/281 (98%)	266 (96%)	10 (4%)	0	100	100
All	All	1104/1124 (98%)	1043 (94%)	55 (5%)	6 (0%)	34	21

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	31	GLN
1	B	151	GLY
1	C	305	ILE
1	B	152	THR
1	A	235	VAL
1	C	240	TYR

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	242/237 (102%)	237 (98%)	5 (2%)	61	55
1	B	244/237 (103%)	237 (97%)	7 (3%)	50	40
1	C	242/237 (102%)	239 (99%)	3 (1%)	78	76
1	D	242/237 (102%)	238 (98%)	4 (2%)	68	64
All	All	970/948 (102%)	951 (98%)	19 (2%)	63	57

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	49	LEU
1	A	128	LYS
1	A	223	LEU
1	A	246	LYS
1	A	266	ASN

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Mol	Chain	Res	Type
1	B	97	LEU
1	B	128	LYS
1	B	153	GLU
1	B	223	LEU
1	B	234	GLU
1	B	246	LYS
1	B	247	ASP
1	C	128	LYS
1	C	176	LYS
1	C	287	LEU
1	D	122	THR
1	D	128	LYS
1	D	286	GLU
1	D	306	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	64	ASN
1	A	124	GLN
1	A	266	ASN
1	A	269	ASN
1	A	283	HIS
1	B	31	GLN
1	B	266	ASN
1	B	269	ASN
1	B	282	ASN
1	C	217	HIS
1	C	282	ASN
1	D	31	GLN
1	D	124	GLN
1	D	266	ASN
1	D	269	ASN
1	D	282	ASN
1	D	283	HIS

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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

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	3686	-	38,48,48	1.56	4 (10%)	47,73,73	2.24	9 (19%)
2	NAD	B	4686	-	38,48,48	1.53	3 (7%)	47,73,73	2.22	8 (17%)
2	NAD	C	5686	-	38,48,48	1.51	3 (7%)	47,73,73	2.25	8 (17%)
2	NAD	D	6686	-	38,48,48	1.50	3 (7%)	47,73,73	2.19	9 (19%)

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	3686	-	-	0/22/62/62	0/5/5/5
2	NAD	B	4686	-	-	0/22/62/62	0/5/5/5
2	NAD	C	5686	-	-	0/22/62/62	0/5/5/5
2	NAD	D	6686	-	-	0/22/62/62	0/5/5/5

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	3686	NAD	C4N-C3N	2.00	1.42	1.39
2	A	3686	NAD	C2A-N1A	3.20	1.40	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	5686	NAD	C2A-N1A	3.28	1.40	1.33
2	B	4686	NAD	C2A-N1A	3.30	1.40	1.33
2	D	6686	NAD	C2A-N1A	3.37	1.40	1.33
2	C	5686	NAD	C2A-N3A	3.47	1.38	1.32
2	D	6686	NAD	C2A-N3A	3.73	1.38	1.32
2	B	4686	NAD	C2A-N3A	3.92	1.39	1.32
2	A	3686	NAD	C2A-N3A	3.99	1.39	1.32
2	D	6686	NAD	O7N-C7N	5.95	1.36	1.24
2	C	5686	NAD	O7N-C7N	5.96	1.36	1.24
2	B	4686	NAD	O7N-C7N	6.01	1.37	1.24
2	A	3686	NAD	O7N-C7N	6.11	1.37	1.24

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	3686	NAD	N3A-C2A-N1A	-11.46	120.12	128.89
2	B	4686	NAD	N3A-C2A-N1A	-11.20	120.32	128.89
2	C	5686	NAD	N3A-C2A-N1A	-11.01	120.47	128.89
2	D	6686	NAD	N3A-C2A-N1A	-10.90	120.55	128.89
2	D	6686	NAD	C4B-O4B-C1B	-4.36	104.93	109.72
2	C	5686	NAD	C4B-O4B-C1B	-4.06	105.25	109.72
2	B	4686	NAD	C4B-O4B-C1B	-4.05	105.27	109.72
2	A	3686	NAD	C4B-O4B-C1B	-3.77	105.58	109.72
2	B	4686	NAD	C1B-N9A-C4A	-3.23	122.06	126.94
2	C	5686	NAD	O7N-C7N-N7N	-2.96	118.42	122.59
2	A	3686	NAD	C1B-N9A-C4A	-2.87	122.62	126.94
2	D	6686	NAD	O7N-C7N-N7N	-2.73	118.75	122.59
2	A	3686	NAD	O7N-C7N-N7N	-2.72	118.77	122.59
2	B	4686	NAD	O7N-C7N-N7N	-2.70	118.79	122.59
2	D	6686	NAD	PN-O3-PA	-2.54	125.59	132.73
2	B	4686	NAD	PN-O3-PA	-2.47	125.78	132.73
2	A	3686	NAD	PN-O3-PA	-2.39	126.02	132.73
2	D	6686	NAD	C1B-N9A-C4A	-2.34	123.41	126.94
2	C	5686	NAD	C1B-N9A-C4A	-2.32	123.45	126.94
2	D	6686	NAD	O7N-C7N-C3N	-2.24	117.14	119.59
2	A	3686	NAD	C4A-C5A-N7A	-2.14	107.51	109.48
2	B	4686	NAD	C4A-C5A-N7A	-2.09	107.56	109.48
2	C	5686	NAD	PN-O3-PA	-2.05	126.97	132.73
2	D	6686	NAD	C4A-C5A-N7A	-2.01	107.63	109.48
2	A	3686	NAD	C4D-O4D-C1D	2.16	112.09	109.72
2	D	6686	NAD	C2B-C1B-N9A	2.37	117.91	114.29
2	C	5686	NAD	C2B-C1B-N9A	2.37	117.92	114.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	3686	NAD	C2B-C1B-N9A	3.29	119.31	114.29
2	B	4686	NAD	C2B-C1B-N9A	3.48	119.62	114.29
2	C	5686	NAD	O4D-C1D-N1N	3.97	112.50	108.13
2	A	3686	NAD	C3N-C7N-N7N	5.25	123.56	117.82
2	B	4686	NAD	C3N-C7N-N7N	5.26	123.58	117.82
2	D	6686	NAD	C3N-C7N-N7N	5.75	124.11	117.82
2	C	5686	NAD	C3N-C7N-N7N	5.77	124.14	117.82

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	5686	NAD	1	0
2	D	6686	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	270/281 (96%)	0.56	20 (7%) 17 19	14, 32, 49, 71	0
1	B	273/281 (97%)	0.67	22 (8%) 15 16	17, 35, 51, 68	0
1	C	271/281 (96%)	0.12	9 (3%) 50 53	13, 21, 37, 61	0
1	D	271/281 (96%)	0.13	3 (1%) 82 84	14, 22, 38, 57	0
All	All	1085/1124 (96%)	0.37	54 (4%) 32 35	13, 27, 47, 71	0

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	152	THR	8.5
1	B	151	GLY	6.2
1	B	152	THR	5.8
1	B	113	TYR	5.6
1	C	306	LYS	5.0
1	B	153	GLU	4.6
1	A	151	GLY	4.4
1	A	153	GLU	4.4
1	D	306	LYS	4.1
1	C	307	ARG	3.9
1	D	31	GLN	3.8
1	A	113	TYR	3.7
1	B	188	ARG	3.6
1	C	303	VAL	3.6
1	A	51	ARG	3.5
1	B	138	ILE	3.3
1	B	233	PRO	3.2
1	C	305	ILE	3.0
1	A	53	GLY	2.9
1	C	31	GLN	2.9
1	B	54	PHE	2.8

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Mol	Chain	Res	Type	RSRZ
1	C	289	VAL	2.8
1	B	55	LYS	2.7
1	B	51	ARG	2.6
1	B	150	ALA	2.6
1	B	140	GLY	2.6
1	A	247	ASP	2.5
1	B	102	ARG	2.5
1	A	245	PHE	2.4
1	A	181	LYS	2.4
1	C	244	GLY	2.4
1	B	78	ASP	2.4
1	A	233	PRO	2.4
1	A	159	TYR	2.3
1	B	250	ARG	2.3
1	A	52	SER	2.3
1	B	192	ASP	2.3
1	B	166	GLU	2.3
1	A	102	ARG	2.2
1	A	161	LEU	2.2
1	A	253	LYS	2.2
1	B	53	GLY	2.2
1	A	176	LYS	2.2
1	A	66	GLU	2.1
1	C	113	TYR	2.1
1	B	171	ILE	2.1
1	D	305	ILE	2.1
1	B	159	TYR	2.1
1	A	180	LYS	2.1
1	A	94	PHE	2.1
1	A	199	TYR	2.1
1	B	74	LEU	2.1
1	B	83	SER	2.0
1	C	242	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 ⓘ

There are no carbohydrates in this entry.

6.4 Ligands

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	5686	44/44	0.94	0.11	0.20	15,20,26,31	0
2	NAD	D	6686	44/44	0.95	0.11	0.09	13,18,24,27	0
2	NAD	A	3686	44/44	0.92	0.12	-0.39	27,33,39,41	0
2	NAD	B	4686	44/44	0.92	0.12	-0.46	35,39,41,42	0

6.5 Other polymers

There are no such residues in this entry.