



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 08:17 AM GMT

PDB ID : 3E27
Title : Nicotinic acid mononucleotide (NaMN) adenylyltransferase from *Bacillus anthracis*: product complex
Authors : Martynowski, D.; Eyobo, Y.; Zhang, H.
Deposited on : 2008-08-05
Resolution : 2.20 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

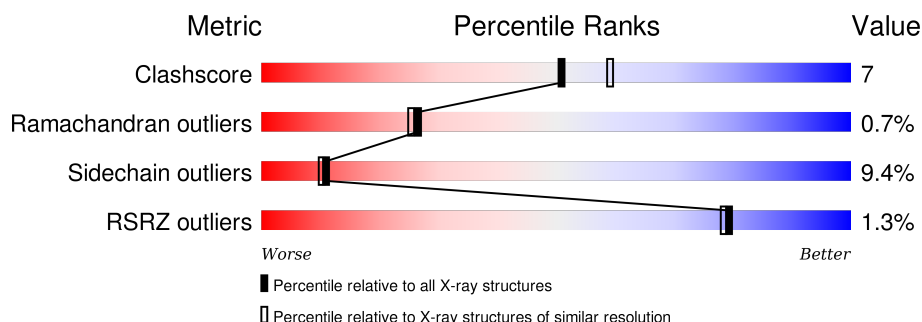
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	102246	4477 (2.20-2.20)
Ramachandran outliers	100387	4404 (2.20-2.20)
Sidechain outliers	100360	4405 (2.20-2.20)
RSRZ outliers	91569	3781 (2.20-2.20)

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	189	<div> <div>3%</div> <div>75%</div> <div>20%</div> <div>• •</div> </div>
1	B	189	<div> <div>%</div> <div>78%</div> <div>19%</div> <div>• •</div> </div>
1	C	189	<div> <div>%</div> <div>84%</div> <div>13%</div> <div>•</div> </div>
1	D	189	<div> <div>%</div> <div>76%</div> <div>19%</div> <div>• •</div> </div>

2 Entry composition [i](#)

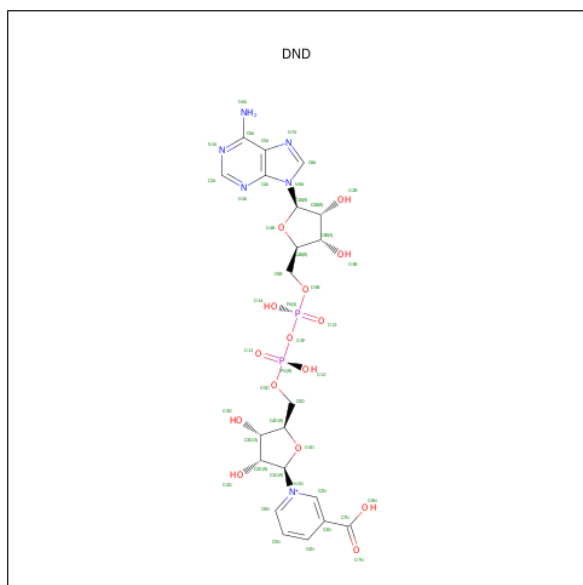
There are 4 unique types of molecules in this entry. The entry contains 6870 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Nicotinate (Nicotinamide) nucleotide adenylyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	188	Total	C	N	O	S	0	1	0
			1547	1005	253	284	5			
1	B	188	Total	C	N	O	S	0	2	0
			1556	1012	255	283	6			
1	C	189	Total	C	N	O	S	0	1	0
			1557	1012	256	283	6			
1	D	187	Total	C	N	O	S	0	0	0
			1536	1000	251	280	5			

- Molecule 2 is NICOTINIC ACID ADENINE DINUCLEOTIDE (three-letter code: DND) (formula: $C_{21}H_{27}N_6O_{15}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			44	21	6	15	2		
2	B	1	Total	C	N	O	P	0	0
			44	21	6	15	2		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	C	1	Total	C	N	O	P	0	0
			44	21	6	15	2		
2	D	1	Total	C	N	O	P	0	0
			44	21	6	15	2		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Mg	0	0
			1	1		
3	A	1	Total	Mg	0	0
			1	1		
3	D	1	Total	Mg	0	0
			1	1		
3	C	1	Total	Mg	0	0
			1	1		

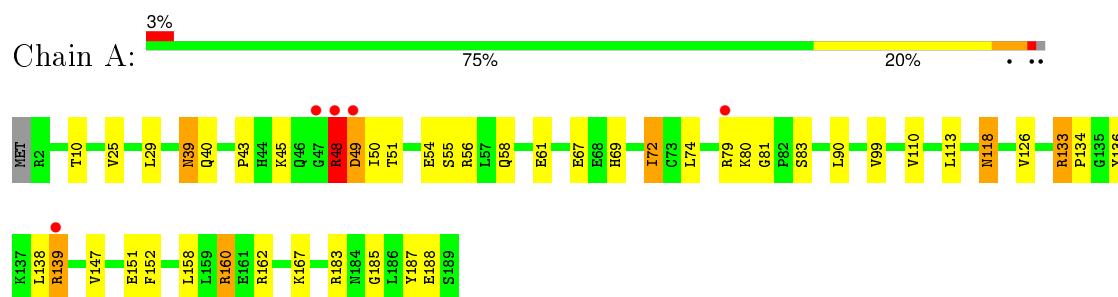
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	117	Total	O	0	0
			117	117		
4	B	141	Total	O	0	0
			141	141		
4	C	126	Total	O	0	0
			126	126		
4	D	110	Total	O	0	0
			110	110		

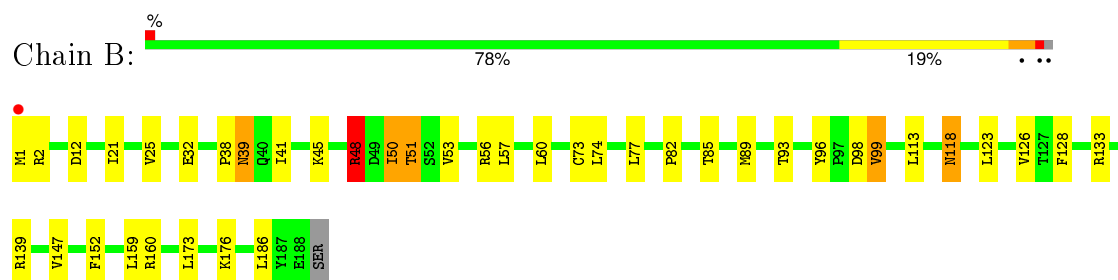
3 Residue-property plots [i](#)

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.

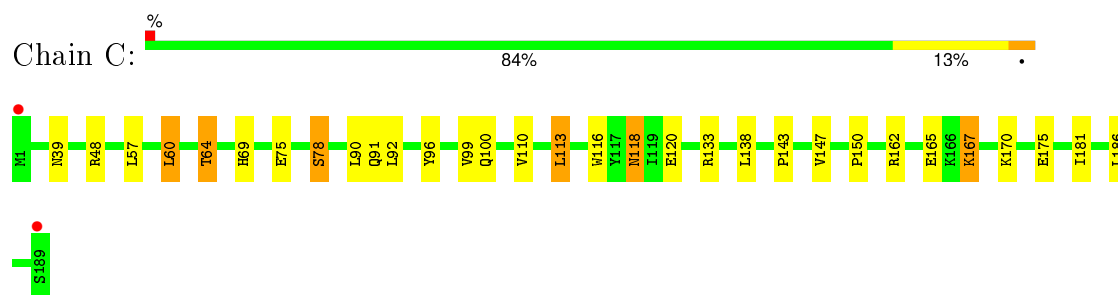
- Molecule 1: Nicotinate (Nicotinamide) nucleotide adenylyltransferase



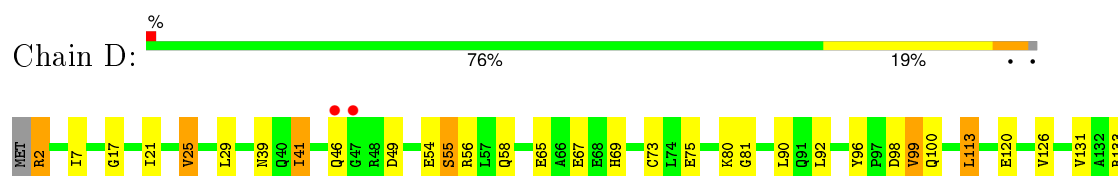
- Molecule 1: Nicotinate (Nicotinamide) nucleotide adenylyltransferase



- Molecule 1: Nicotinate (Nicotinamide) nucleotide adenylyltransferase



- Molecule 1: Nicotinate (Nicotinamide) nucleotide adenylyltransferase



P134	G135	Y136	K137
I149	P150	E151	F152
L158	L159	R160	K167
K176	Y180	E188	SER

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	41.86Å 137.41Å 143.97Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.20 49.70 – 2.19	Depositor EDS
% Data completeness (in resolution range)	99.4 (50.00-2.20) 99.3 (49.70-2.19)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.15 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.201 , 0.273 0.204 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	31.7	Xtriage
Anisotropy	0.497	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 41.9	EDS
Estimated twinning fraction	0.007 for -h,l,k	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	1 of 43274 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6870	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.57% 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: DND, MG

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	0.65	0/1593	0.74	1/2162 (0.0%)
1	B	0.63	0/1606	0.71	2/2179 (0.1%)
1	C	0.63	0/1603	0.70	0/2175
1	D	0.66	0/1578	0.72	2/2143 (0.1%)
All	All	0.64	0/6380	0.72	5/8659 (0.1%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	56	ARG	NE-CZ-NH2	-6.44	117.08	120.30
1	B	56	ARG	NE-CZ-NH2	-5.75	117.43	120.30
1	A	160	ARG	NE-CZ-NH2	-5.66	117.47	120.30
1	B	186	LEU	CA-CB-CG	5.60	128.18	115.30
1	D	113	LEU	CA-CB-CG	5.32	127.54	115.30

There are no chirality outliers.

There are no planarity outliers.

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	1547	0	1543	23	0
1	B	1556	0	1559	22	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	1557	0	1562	19	0
1	D	1536	0	1536	29	0
2	A	44	0	25	0	0
2	B	44	0	24	1	0
2	C	44	0	24	0	0
2	D	44	0	24	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	117	0	0	1	0
4	B	141	0	0	2	0
4	C	126	0	0	2	0
4	D	110	0	0	3	0
All	All	6870	0	6297	88	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 7.

All (88) 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:B:1:MET:HB2	1:B:32[B]:GLU:OE1	1.79	0.81
1:B:25:VAL:HG22	1:B:147:VAL:HG21	1.64	0.79
1:D:41:ILE:HD11	1:D:46:GLN:HG2	1.67	0.76
1:B:96:TYR:HB3	1:B:99:VAL:CG1	2.25	0.66
1:C:170:LYS:HE2	1:D:67:GLU:OE2	1.95	0.65
1:B:96:TYR:HB3	1:B:99:VAL:HG13	1.79	0.64
1:C:96:TYR:HB3	1:C:99:VAL:HG13	1.79	0.64
1:A:69:HIS:H	1:A:69:HIS:HD1	1.45	0.63
1:A:67:GLU:OE2	1:A:69:HIS:HE1	1.82	0.62
1:B:89:MET:O	1:B:93:THR:HG22	2.02	0.60
1:D:80:LYS:HE2	4:D:274:HOH:O	2.01	0.60
1:D:75:GLU:HG3	1:D:92:LEU:HD21	1.84	0.58
1:A:90:LEU:HD11	1:C:78:SER:HB3	1.84	0.58
1:B:12:ASP:OD1	1:B:51:THR:HG23	2.02	0.58
1:D:96:TYR:HB3	1:D:99:VAL:CG1	2.33	0.58
1:C:167:LYS:NZ	4:C:300:HOH:O	2.36	0.58
1:C:75:GLU:HG3	1:C:92:LEU:HD21	1.86	0.57
1:D:98:ASP:HB2	4:D:228:HOH:O	2.04	0.57
1:D:131:VAL:HG11	1:D:149:ILE:HD13	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:45:LYS:HB3	1:B:48:ARG:CD	2.35	0.56
1:A:133:ARG:HG2	1:A:136:TYR:CD1	2.42	0.55
1:A:160:ARG:HG2	1:A:187:TYR:CD1	2.42	0.55
1:C:60:LEU:O	1:C:64:THR:HB	2.07	0.55
1:D:65:GLU:OE2	1:D:176:LYS:NZ	2.40	0.55
1:D:96:TYR:HB3	1:D:99:VAL:HG13	1.89	0.54
1:B:45:LYS:HB3	1:B:48:ARG:HD2	1.89	0.54
1:D:21:ILE:O	1:D:25:VAL:CG1	2.56	0.54
1:A:90:LEU:CD1	1:C:78:SER:HB3	2.37	0.53
1:D:17:GLY:O	1:D:21:ILE:HG23	2.09	0.53
1:C:69:HIS:HD1	1:C:69:HIS:H	1.56	0.53
1:D:69:HIS:H	1:D:69:HIS:HD1	1.58	0.52
1:A:54:GLU:O	1:A:58:GLN:HG3	2.10	0.52
1:D:131:VAL:CG1	1:D:149:ILE:HD13	2.40	0.52
1:B:21:ILE:O	1:B:25:VAL:HG23	2.10	0.52
1:B:74:LEU:HD22	1:B:77:LEU:HD12	1.93	0.51
1:D:67:GLU:HB3	1:D:69:HIS:CE1	2.44	0.51
1:D:21:ILE:O	1:D:25:VAL:HG13	2.10	0.51
1:B:41:ILE:HB	1:B:82:PRO:HA	1.93	0.51
1:C:162:ARG:HG2	1:C:167:LYS:HB2	1.93	0.50
1:C:181:ILE:HG23	1:C:186:LEU:HB2	1.93	0.50
1:B:123:LEU:HD23	1:B:128:PHE:HD1	1.76	0.50
1:A:185:GLY:O	1:A:188:GLU:HG2	2.12	0.50
1:A:45:LYS:HB3	1:A:48:ARG:HG3	1.93	0.49
1:A:80:LYS:HG2	1:A:81:GLY:H	1.78	0.49
1:D:73:CYS:SG	1:D:75:GLU:HG2	2.53	0.48
1:A:51:THR:O	1:A:56:ARG:HD3	2.13	0.47
1:C:165:GLU:OE1	1:C:167:LYS:HE3	2.15	0.47
1:D:134:PRO:HD3	1:D:151:GLU:HB3	1.96	0.47
1:D:160:ARG:NH2	4:D:259:HOH:O	2.47	0.47
1:A:43:PRO:HG3	1:A:83:SER:O	2.14	0.47
1:C:150:PRO:HG3	1:D:158:LEU:HD21	1.97	0.47
1:D:80:LYS:HD2	1:D:81:GLY:H	1.78	0.46
1:C:118:ASN:HD22	1:C:118:ASN:HA	1.58	0.46
1:D:58:GLN:NE2	1:D:180:TYR:OH	2.49	0.46
1:D:133:ARG:HG2	1:D:136:TYR:HD1	1.81	0.45
1:A:80:LYS:CG	1:A:81:GLY:H	2.29	0.45
1:D:2:ARG:O	1:D:100:GLN:N	2.46	0.45
1:D:55:SER:HA	1:D:58:GLN:HE21	1.82	0.45
1:B:96:TYR:CB	1:B:99:VAL:HG13	2.45	0.44
1:D:54:GLU:O	1:D:58:GLN:HG3	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:139:ARG:H	1:A:139:ARG:HD2	1.83	0.44
1:B:45:LYS:HD2	1:B:48:ARG:CZ	2.47	0.44
1:C:110:VAL:O	1:C:113:LEU:HB2	2.18	0.43
1:D:41:ILE:HG23	1:D:81:GLY:O	2.18	0.43
1:A:134:PRO:HD3	1:A:151:GLU:HB3	2.00	0.43
1:A:162:ARG:HG2	1:A:167:LYS:HB2	2.00	0.43
1:A:110:VAL:HG21	1:A:138:LEU:HD23	2.01	0.43
1:D:7:ILE:CD1	1:D:21:ILE:HD11	2.48	0.43
1:A:133:ARG:HD2	4:A:274:HOH:O	2.19	0.43
1:B:38:PRO:HD2	1:B:73:CYS:O	2.18	0.43
1:B:85:THR:OG1	2:B:191:DND:H13	2.19	0.43
1:D:41:ILE:HD11	1:D:46:GLN:CG	2.45	0.42
1:C:48:ARG:HD2	4:C:214:HOH:O	2.18	0.42
1:B:98:ASP:HB2	4:B:306:HOH:O	2.19	0.42
1:B:39:ASN:HD22	1:B:39:ASN:C	2.22	0.42
1:A:133:ARG:HG2	1:A:136:TYR:HD1	1.82	0.42
1:B:160:ARG:NH1	4:B:322:HOH:O	2.50	0.41
1:B:118:ASN:HD22	1:B:118:ASN:HA	1.73	0.41
1:A:51:THR:HG22	1:A:55:SER:HB3	2.02	0.41
1:B:159:LEU:HD13	1:B:173:LEU:HD13	2.01	0.41
1:D:167:LYS:HE3	1:D:167:LYS:HB2	1.92	0.41
1:B:123:LEU:HD23	1:B:128:PHE:CD1	2.55	0.41
1:C:138:LEU:HD23	1:C:138:LEU:HA	1.84	0.41
1:A:72:ILE:HD11	1:A:74:LEU:CD2	2.51	0.41
1:C:113:LEU:HA	1:C:116:TRP:CD1	2.56	0.40
1:A:118:ASN:H	1:C:91:GLN:NE2	2.18	0.40
1:A:10:THR:HA	1:A:39:ASN:HD21	1.87	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	187/189 (99%)	177 (95%)	7 (4%)	3 (2%)	12	8
1	B	188/189 (100%)	180 (96%)	6 (3%)	2 (1%)	17	14
1	C	188/189 (100%)	181 (96%)	7 (4%)	0	100	100
1	D	185/189 (98%)	183 (99%)	2 (1%)	0	100	100
All	All	748/756 (99%)	721 (96%)	22 (3%)	5 (1%)	26	25

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	49	ASP
1	B	48	ARG
1	B	50	ILE
1	A	48	ARG
1	A	79	ARG

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	171/171 (100%)	152 (89%)	19 (11%)	8	6
1	B	172/171 (101%)	155 (90%)	17 (10%)	10	9
1	C	172/171 (101%)	158 (92%)	14 (8%)	15	14
1	D	169/171 (99%)	154 (91%)	15 (9%)	12	11
All	All	684/684 (100%)	619 (90%)	65 (10%)	11	10

All (65) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	VAL
1	A	29	LEU
1	A	39	ASN
1	A	40	GLN
1	A	48	ARG
1	A	49	ASP

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Mol	Chain	Res	Type
1	A	50	ILE
1	A	61	GLU
1	A	72	ILE
1	A	99	VAL
1	A	113	LEU
1	A	118	ASN
1	A	126	VAL
1	A	133	ARG
1	A	139	ARG
1	A	147	VAL
1	A	152	PHE
1	A	158	LEU
1	A	183	ARG
1	B	2[A]	ARG
1	B	2[B]	ARG
1	B	39	ASN
1	B	48	ARG
1	B	50	ILE
1	B	51	THR
1	B	53	VAL
1	B	57	LEU
1	B	60	LEU
1	B	99	VAL
1	B	113	LEU
1	B	118	ASN
1	B	126	VAL
1	B	133	ARG
1	B	139	ARG
1	B	152	PHE
1	B	176	LYS
1	C	39	ASN
1	C	57	LEU
1	C	60	LEU
1	C	64	THR
1	C	78	SER
1	C	90	LEU
1	C	113	LEU
1	C	118	ASN
1	C	120	GLU
1	C	133	ARG
1	C	143	PRO
1	C	147	VAL

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Mol	Chain	Res	Type
1	C	167	LYS
1	C	175	GLU
1	D	2	ARG
1	D	25	VAL
1	D	29	LEU
1	D	39	ASN
1	D	41	ILE
1	D	49	ASP
1	D	55	SER
1	D	90	LEU
1	D	99	VAL
1	D	113	LEU
1	D	120	GLU
1	D	126	VAL
1	D	137	LYS
1	D	152	PHE
1	D	176	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	39	ASN
1	A	46	GLN
1	A	118	ASN
1	B	39	ASN
1	B	100	GLN
1	B	102	HIS
1	B	118	ASN
1	C	23	ASN
1	C	39	ASN
1	C	91	GLN
1	C	118	ASN
1	D	23	ASN
1	D	39	ASN
1	D	46	GLN
1	D	58	GLN

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 ⓘ

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 for Mogul analysis.

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	DND	A	191	-	35,48,48	1.13	2 (5%)	44,73,73	2.31	6 (13%)
2	DND	B	191	3	35,48,48	1.10	2 (5%)	44,73,73	2.09	4 (9%)
2	DND	C	191	3	35,48,48	1.06	2 (5%)	44,73,73	2.05	3 (6%)
2	DND	D	191	3	35,48,48	1.18	3 (8%)	44,73,73	2.14	5 (11%)

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	DND	A	191	-	-	0/18/62/62	0/5/5/5
2	DND	B	191	3	-	0/18/62/62	0/5/5/5
2	DND	C	191	3	-	0/18/62/62	0/5/5/5
2	DND	D	191	3	-	0/18/62/62	0/5/5/5

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	191	DND	PN-O12	-2.05	1.46	1.54
2	A	191	DND	C2A-N1A	2.34	1.38	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	191	DND	C2A-N1A	2.47	1.38	1.33
2	C	191	DND	C2A-N1A	2.69	1.39	1.33
2	D	191	DND	C2A-N1A	2.88	1.39	1.33
2	B	191	DND	C2A-N3A	3.42	1.38	1.32
2	C	191	DND	C2A-N3A	3.56	1.38	1.32
2	D	191	DND	C2A-N3A	3.64	1.38	1.32
2	A	191	DND	C2A-N3A	3.76	1.38	1.32

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	191	DND	N3A-C2A-N1A	-13.16	118.81	128.89
2	D	191	DND	N3A-C2A-N1A	-11.84	119.83	128.89
2	B	191	DND	N3A-C2A-N1A	-11.75	119.90	128.89
2	C	191	DND	N3A-C2A-N1A	-11.44	120.13	128.89
2	A	191	DND	C1B-N9A-C4A	-3.10	122.27	126.94
2	D	191	DND	O3P-PN-O5D	-2.87	95.33	102.94
2	B	191	DND	C1B-N9A-C4A	-2.62	122.98	126.94
2	A	191	DND	O3P-PN-O5D	-2.51	96.28	102.94
2	C	191	DND	C2B-C1B-N9A	-2.19	110.94	114.29
2	A	191	DND	O3D-C3D-C2D	-2.02	105.27	111.83
2	B	191	DND	O3D-C3D-C2D	-2.01	105.27	111.83
2	D	191	DND	C4N-C3N-C7N	2.09	123.26	120.45
2	D	191	DND	O14-PA-O3P	2.54	116.61	105.09
2	A	191	DND	O4D-C1D-N1N	2.55	110.93	108.13
2	D	191	DND	O4D-C1D-N1N	2.60	110.99	108.13
2	A	191	DND	O14-PA-O3P	2.84	117.96	105.09
2	B	191	DND	O4D-C1D-N1N	3.14	111.59	108.13
2	C	191	DND	O4D-C1D-N1N	3.61	112.10	108.13

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	191	DND	1	0

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	188/189 (99%)	-0.27	5 (2%) 58 57	18, 33, 49, 62	0
1	B	188/189 (99%)	-0.41	1 (0%) 91 91	18, 29, 46, 58	0
1	C	189/189 (100%)	-0.34	2 (1%) 82 82	18, 31, 50, 57	0
1	D	187/189 (98%)	-0.47	2 (1%) 82 82	19, 30, 46, 58	0
All	All	752/756 (99%)	-0.37	10 (1%) 79 78	18, 31, 48, 62	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	79	ARG	4.5
1	D	47	GLY	3.6
1	C	1	MET	3.5
1	A	49	ASP	3.5
1	B	1	MET	3.1
1	A	47	GLY	2.8
1	A	48	ARG	2.7
1	A	139	ARG	2.4
1	C	189	SER	2.4
1	D	46	GLN	2.2

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, 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(Å ²)	Q<0.9
2	DND	C	191	44/44	0.98	0.10	0.01	16,21,25,26	0
2	DND	B	191	44/44	0.97	0.09	-0.59	18,25,28,30	0
2	DND	A	191	44/44	0.98	0.09	-0.78	21,24,28,31	0
2	DND	D	191	44/44	0.98	0.08	-1.19	18,23,29,32	0
3	MG	C	190	1/1	0.94	0.10	-	40,40,40,40	0
3	MG	B	190	1/1	0.96	0.10	-	36,36,36,36	0
3	MG	A	190	1/1	0.95	0.09	-	40,40,40,40	0
3	MG	D	190	1/1	0.94	0.18	-	39,39,39,39	0

6.5 Other polymers [i](#)

There are no such residues in this entry.