



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:23 AM GMT

PDB ID : 2QJT  
Title : Crystal structure of a bifunctional NMN adenylyltransferase/ADP ribose pyrophosphatase complexed with AMP and MN ion from Francisella tularensis  
Authors : Huang, N.; Sorci, L.; Zhang, X.; Brautigan, C.; Raffaelli, N.; Magni, G.; Grishin, N.V.; Osterman, A.; Zhang, H.  
Deposited on : 2007-07-09  
Resolution : 2.30 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

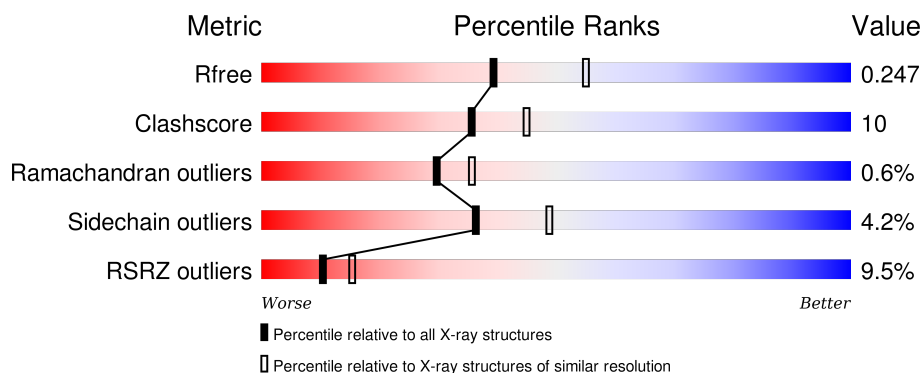
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*


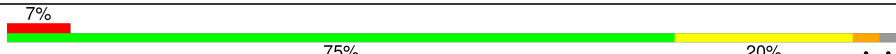
The reported resolution of this entry is 2.30 Å.

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	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

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	352	
1	B	352	

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6073 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 Nicotinamide-nucleotide adenylyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	344	Total	C	N	O	S	0	0	0
			2827	1820	476	519	12			
1	A	340	Total	C	N	O	S	0	0	0
			2792	1798	469	513	12			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-4	GLY	-	EXPRESSION TAG	UNP Q5NHR1
B	-3	ALA	-	EXPRESSION TAG	UNP Q5NHR1
B	-2	MET	-	EXPRESSION TAG	UNP Q5NHR1
B	-1	ASP	-	EXPRESSION TAG	UNP Q5NHR1
B	0	PRO	-	EXPRESSION TAG	UNP Q5NHR1
A	-4	GLY	-	EXPRESSION TAG	UNP Q5NHR1
A	-3	ALA	-	EXPRESSION TAG	UNP Q5NHR1
A	-2	MET	-	EXPRESSION TAG	UNP Q5NHR1
A	-1	ASP	-	EXPRESSION TAG	UNP Q5NHR1
A	0	PRO	-	EXPRESSION TAG	UNP Q5NHR1

- Molecule 2 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	4	Total	Mn	0	0
			4	4		
2	A	5	Total	Mn	0	0
			5	5		

- Molecule 3 is ADENOSINE MONOPHOSPHATE (three-letter code: AMP) (formula: C<sub>10</sub>H<sub>14</sub>N<sub>5</sub>O<sub>7</sub>P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	B	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
3	A	1	Total	C	N	O	P	0	0
			23	10	5	7	1		

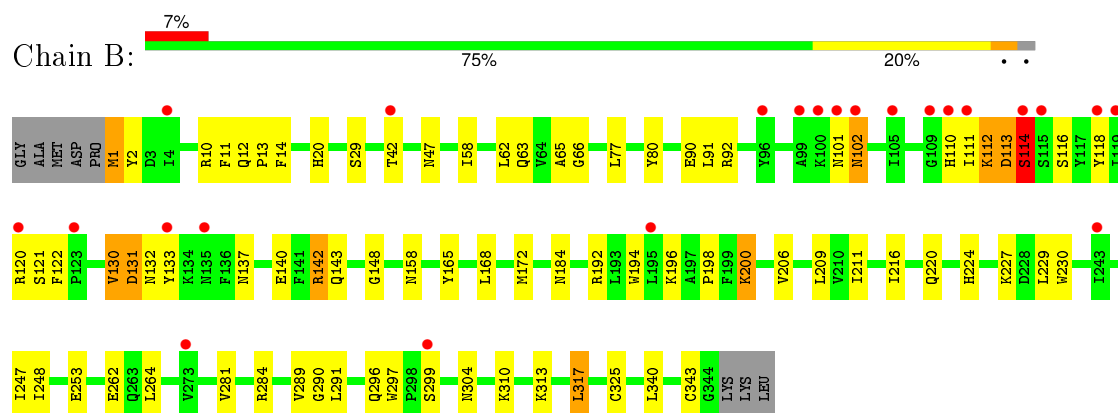
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	203	Total	O	0	0
			203	203		
4	A	196	Total	O	0	0
			196	196		

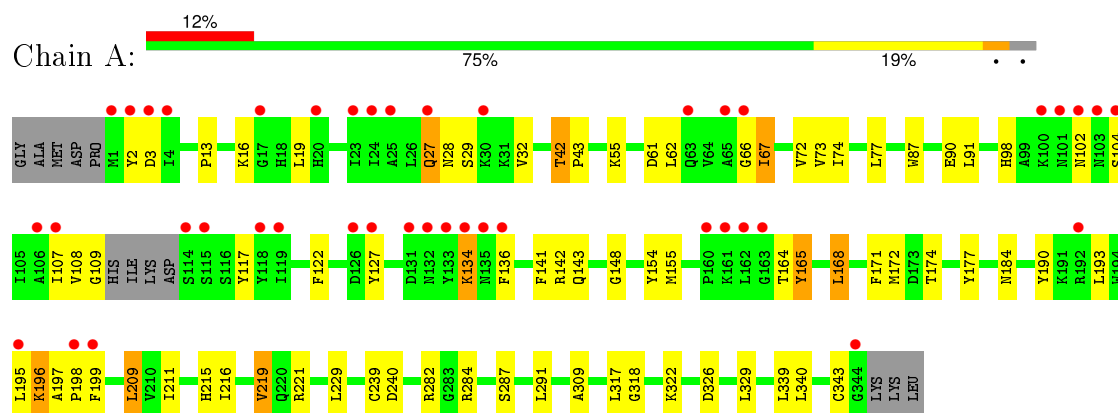
### 3 Residue-property plots

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: Nicotinamide-nucleotide adenylyltransferase



- Molecule 1: Nicotinamide-nucleotide adenylyltransferase



## 4 Data and refinement statistics

Property	Value	Source
Space group	F 2 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	125.64Å 163.65Å 179.97Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	33.61 – 2.30 33.61 – 2.30	Depositor EDS
% Data completeness (in resolution range)	97.0 (33.61-2.30) 96.9 (33.61-2.30)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.49 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.194 , 0.249 0.192 , 0.247	Depositor DCC
$R_{free}$ test set	2003 reflections (5.29%)	DCC
Wilson B-factor (Å <sup>2</sup> )	41.9	Xtriage
Anisotropy	0.020	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 46.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 39854 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	6073	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.56% of the height of the origin peak. No significant pseudotranslation is detected.*

<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 ⓘ

### 5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: AMP, MN

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	0.76	0/2860	0.75	2/3873 (0.1%)
1	B	0.71	0/2897	0.75	3/3924 (0.1%)
All	All	0.74	0/5757	0.75	5/7797 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	4

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	B	142	ARG	NE-CZ-NH2	-7.97	116.32	120.30
1	A	240	ASP	CB-CG-OD2	7.23	124.80	118.30
1	B	114	SER	CB-CA-C	-6.13	98.45	110.10
1	B	142	ARG	NE-CZ-NH1	5.43	123.02	120.30
1	A	240	ASP	CB-CG-OD1	-5.25	113.57	118.30

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	130	VAL	Peptide
1	B	133	TYR	Peptide
1	B	289	VAL	Peptide

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Mol	Chain	Res	Type	Group
1	B	65	ALA	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	2792	0	2742	57	0
1	B	2827	0	2778	56	1
2	A	5	0	0	0	0
2	B	4	0	0	0	0
3	A	23	0	12	0	0
3	B	23	0	12	0	0
4	A	196	0	0	5	0
4	B	203	0	0	8	1
All	All	6073	0	5544	109	1

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 10.

All (109) 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:A:171:PHE:O	1:A:174:THR:HG22	1.57	1.04
1:B:42:THR:HG21	1:B:229:LEU:HD13	1.50	0.94
1:B:42:THR:HG21	1:B:229:LEU:CD1	2.07	0.85
1:A:164:THR:HG22	1:A:168:LEU:HD22	1.61	0.81
1:A:142:ARG:NH1	4:A:641:HOH:O	2.18	0.76
1:A:42:THR:HG22	4:A:615:HOH:O	1.85	0.76
1:A:66:GLY:H	1:A:67:ILE:HG13	1.49	0.76
1:A:73:VAL:HG21	1:A:98:HIS:CE1	2.23	0.73
1:B:1:MET:N	4:B:784:HOH:O	2.21	0.72
1:A:174:THR:HG23	1:A:177:TYR:H	1.54	0.72
1:B:168:LEU:O	1:B:172:MET:HG3	1.92	0.69
1:A:55:LYS:HG3	1:A:74:ILE:HG21	1.76	0.67
1:B:12:GLN:HE21	1:B:47:ASN:HD22	1.41	0.67
1:B:291:LEU:HD23	1:B:343:CYS:HB3	1.77	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:155:MET:HE1	1:A:164:THR:HB	1.78	0.65
1:B:313:LYS:NZ	4:B:791:HOH:O	2.30	0.64
1:B:253:GLU:OE2	4:B:686:HOH:O	2.14	0.64
1:A:16:LYS:HE3	1:A:155:MET:O	1.99	0.62
1:B:211:ILE:HD11	1:B:340:LEU:HD21	1.82	0.61
1:B:196:LYS:HD2	1:A:196:LYS:HB3	1.81	0.61
1:A:196:LYS:NZ	4:A:690:HOH:O	2.32	0.61
1:A:282:ARG:HD3	4:A:723:HOH:O	2.00	0.61
1:B:230:TRP:CH2	1:B:313:LYS:HD3	2.37	0.59
1:B:116:SER:HB2	1:B:120:ARG:HH21	1.68	0.59
1:B:137:ASN:ND2	1:B:140:GLU:HB2	2.18	0.59
1:A:3:ASP:H	1:A:104:SER:HB3	1.68	0.59
1:B:62:LEU:O	1:B:66:GLY:HA2	2.02	0.58
1:B:143:GLN:HE22	1:B:184:ASN:HD21	1.51	0.58
1:A:216:ILE:HD11	1:A:317:LEU:HD21	1.86	0.57
1:B:20:HIS:NE2	1:B:132:ASN:OD1	2.35	0.56
1:B:130:VAL:O	1:B:131:ASP:HB2	2.05	0.56
1:B:1:MET:HG2	1:B:2:TYR:N	2.20	0.56
1:B:110:HIS:HA	1:B:130:VAL:O	2.05	0.56
1:B:304:ASN:HB2	4:B:715:HOH:O	2.06	0.56
1:B:10:ARG:HH12	1:B:47:ASN:CG	2.09	0.55
1:B:248:ILE:HD12	1:B:264:LEU:HD13	1.89	0.55
1:A:211:ILE:HD11	1:A:340:LEU:CD2	2.37	0.55
1:A:66:GLY:N	1:A:67:ILE:HG13	2.21	0.55
1:A:136:PHE:HE1	1:A:154:TYR:HA	1.73	0.54
1:A:155:MET:CE	1:A:164:THR:HB	2.37	0.54
1:A:32:VAL:CG1	1:A:72:VAL:HG22	2.38	0.54
1:A:211:ILE:HD11	1:A:340:LEU:HD21	1.90	0.54
1:B:91:LEU:HD23	1:B:122:PHE:CZ	2.43	0.53
1:A:42:THR:HG21	1:A:229:LEU:HD13	1.90	0.53
1:A:61:ASP:OD2	1:A:164:THR:OG1	2.16	0.53
1:B:224:HIS:CE1	1:B:227:LYS:HE3	2.44	0.52
1:B:192:ARG:HD3	4:B:733:HOH:O	2.09	0.52
1:A:42:THR:HB	1:A:326:ASP:O	2.09	0.52
1:B:216:ILE:HG13	1:B:317:LEU:CD1	2.40	0.52
1:A:155:MET:HE1	1:A:164:THR:CB	2.40	0.52
1:A:143:GLN:HE22	1:A:184:ASN:HD21	1.57	0.52
1:A:13:PRO:HG3	1:A:141:PHE:CE2	2.45	0.51
1:B:220:GLN:NE2	1:B:310:LYS:HD2	2.26	0.51
1:A:87:TRP:CZ3	1:A:117:TYR:HE2	2.27	0.51
1:A:2:TYR:O	1:A:29:SER:HA	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:102:ASN:N	1:B:102:ASN:HD22	2.09	0.51
1:B:148:GLY:HA2	1:B:172:MET:CE	2.40	0.50
1:B:206:VAL:HB	1:B:247:ILE:HG12	1.92	0.50
1:A:209:LEU:HD12	1:A:339:LEU:HD12	1.94	0.50
1:A:193:LEU:HD11	4:A:690:HOH:O	2.12	0.49
1:B:211:ILE:CD1	1:B:340:LEU:HD21	2.42	0.49
1:A:318:GLY:O	1:A:322:LYS:HG3	2.13	0.48
1:B:230:TRP:HH2	1:B:313:LYS:HD3	1.75	0.48
1:B:194:TRP:CE2	1:A:198:PRO:HG2	2.49	0.48
1:B:113:ASP:O	1:B:114:SER:CB	2.61	0.48
1:B:113:ASP:O	1:B:114:SER:HB3	2.14	0.48
1:B:142:ARG:NH2	4:B:629:HOH:O	2.47	0.47
1:B:112:LYS:HE3	1:B:131:ASP:HB3	1.96	0.47
1:B:80:TYR:CE1	1:B:90:GLU:OE1	2.67	0.47
1:B:42:THR:HG21	1:B:229:LEU:HD11	1.93	0.47
1:A:42:THR:OG1	1:A:43:PRO:HD2	2.13	0.47
1:B:10:ARG:NH1	4:B:662:HOH:O	2.47	0.47
1:B:110:HIS:C	1:B:112:LYS:H	2.18	0.46
1:A:2:TYR:HD1	1:A:28:ASN:O	1.98	0.46
1:B:10:ARG:NH1	1:B:47:ASN:OD1	2.43	0.45
1:B:14:PHE:CD1	1:B:58:ILE:HG12	2.51	0.45
1:B:198:PRO:HG3	1:A:193:LEU:HG	1.97	0.45
1:A:219:VAL:HG22	1:A:309:ALA:HB1	1.99	0.44
1:A:134:LYS:HD3	1:A:134:LYS:H	1.82	0.44
1:A:221:ARG:HD2	1:A:329:LEU:HD22	1.99	0.44
1:A:197:ALA:HB1	1:A:199:PHE:O	2.17	0.44
1:A:164:THR:CG2	1:A:168:LEU:HD22	2.40	0.44
1:B:281:VAL:HG22	4:B:693:HOH:O	2.17	0.44
1:B:297:TRP:HD1	1:B:299:SER:O	2.00	0.44
1:A:32:VAL:HG12	1:A:72:VAL:HG22	2.00	0.43
1:B:12:GLN:HE21	1:B:47:ASN:ND2	2.14	0.43
1:B:209:LEU:N	1:B:290:GLY:O	2.29	0.43
1:B:77:LEU:HD23	1:B:90:GLU:HG2	2.00	0.42
1:A:155:MET:HG3	1:A:165:TYR:CD1	2.54	0.42
1:A:77:LEU:HD23	1:A:90:GLU:HG2	2.01	0.42
1:A:19:LEU:HD13	1:A:62:LEU:HD23	2.00	0.42
1:A:190:TYR:CZ	1:A:282:ARG:HD2	2.55	0.42
1:A:107:ILE:HD11	1:A:122:PHE:HB2	2.02	0.42
1:A:291:LEU:HD23	1:A:343:CYS:HB3	2.02	0.42
1:B:2:TYR:O	1:B:29:SER:HA	2.20	0.41
1:A:91:LEU:HD23	1:A:122:PHE:CZ	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:211:ILE:HA	1:A:215:HIS:O	2.20	0.41
1:B:114:SER:O	1:B:118:TYR:CD1	2.74	0.41
1:A:27:GLN:HB3	1:A:27:GLN:HE21	1.69	0.41
1:B:80:TYR:HE1	1:B:90:GLU:OE1	2.03	0.41
1:A:219:VAL:HG22	1:A:309:ALA:CB	2.51	0.41
1:A:55:LYS:HG3	1:A:74:ILE:CG2	2.46	0.41
1:A:193:LEU:HD12	1:A:196:LYS:NZ	2.35	0.41
1:B:148:GLY:HA2	1:B:172:MET:HE3	2.02	0.40
1:A:109:GLY:C	1:A:127:TYR:OH	2.60	0.40
1:B:200:LYS:HB3	1:A:239:CYS:SG	2.61	0.40
1:B:92:ARG:NH2	1:B:121:SER:O	2.55	0.40
1:A:148:GLY:HA2	1:A:172:MET:CE	2.51	0.40
1:B:11:PHE:O	1:B:13:PRO:HA	2.21	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:63:GLN:OE1	4:B:765:HOH:O[2_655]	2.16	0.04

## 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	336/352 (96%)	321 (96%)	14 (4%)	1 (0%)	46	57
1	B	342/352 (97%)	324 (95%)	15 (4%)	3 (1%)	21	24
All	All	678/704 (96%)	645 (95%)	29 (4%)	4 (1%)	30	36

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	114	SER
1	B	131	ASP
1	B	113	ASP
1	A	102	ASN

### 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	307/317 (97%)	294 (96%)	13 (4%)	36	49
1	B	311/317 (98%)	298 (96%)	13 (4%)	36	49
All	All	618/634 (98%)	592 (96%)	26 (4%)	36	49

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	1	MET
1	B	101	ASN
1	B	102	ASN
1	B	111	ILE
1	B	112	LYS
1	B	158	ASN
1	B	165	TYR
1	B	200	LYS
1	B	262	GLU
1	B	284	ARG
1	B	296	GLN
1	B	317	LEU
1	B	325	CYS
1	A	27	GLN
1	A	42	THR
1	A	67	ILE
1	A	108	VAL
1	A	134	LYS
1	A	165	TYR
1	A	168	LEU

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Mol	Chain	Res	Type
1	A	195	LEU
1	A	196	LYS
1	A	209	LEU
1	A	219	VAL
1	A	284	ARG
1	A	287	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	12	GLN
1	B	28	ASN
1	B	63	GLN
1	B	101	ASN
1	B	102	ASN
1	B	137	ASN
1	B	143	GLN
1	B	220	GLN
1	A	27	GLN
1	A	184	ASN
1	A	220	GLN
1	A	323	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](#)

Of 11 ligands modelled in this entry, 9 are monoatomic - leaving 2 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$
3	AMP	A	601	2	20,25,25	1.31	2 (10%)	22,38,38	2.63	4 (18%)
3	AMP	B	602	2	20,25,25	1.28	3 (15%)	22,38,38	2.16	4 (18%)

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
3	AMP	A	601	2	-	0/6/26/26	0/3/3/3
3	AMP	B	602	2	-	0/6/26/26	0/3/3/3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	602	AMP	C2-N3	2.28	1.36	1.32
3	A	601	AMP	C5-C4	2.66	1.46	1.40
3	B	602	AMP	O4'-C1'	2.79	1.44	1.41
3	B	602	AMP	C5-C4	3.31	1.48	1.40
3	A	601	AMP	O4'-C1'	3.75	1.45	1.41

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	601	AMP	N3-C2-N1	-9.73	121.44	128.89
3	B	602	AMP	N3-C2-N1	-7.54	123.12	128.89
3	A	601	AMP	C2'-C1'-N9	-5.26	106.26	114.29
3	B	602	AMP	C2'-C1'-N9	-4.19	107.89	114.29
3	B	602	AMP	O3P-P-O2P	2.01	115.03	107.38
3	A	601	AMP	O4'-C4'-C5'	2.03	116.59	109.32
3	A	601	AMP	C2-N1-C6	2.81	123.78	118.77
3	B	602	AMP	O4'-C1'-N9	2.98	114.33	108.10

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 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	340/352 (96%)	0.45	42 (12%) 5 8	21, 43, 83, 102	0
1	B	344/352 (97%)	0.36	23 (6%) 21 29	26, 42, 76, 99	0
All	All	684/704 (97%)	0.40	65 (9%) 10 15	21, 42, 81, 102	0

All (65) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	133	TYR	10.2
1	A	119	ILE	6.5
1	A	162	LEU	6.3
1	A	101	ASN	5.4
1	B	101	ASN	5.1
1	A	65	ALA	4.7
1	A	134	LYS	4.6
1	B	111	ILE	4.6
1	A	102	ASN	4.4
1	B	102	ASN	4.2
1	B	99	ALA	4.2
1	B	100	LYS	4.0
1	A	135	ASN	4.0
1	A	126	ASP	4.0
1	A	23	ILE	3.8
1	A	127	TYR	3.8
1	B	105	ILE	3.7
1	B	120	ARG	3.6
1	B	135	ASN	3.6
1	B	114	SER	3.6
1	B	96	TYR	3.6
1	A	114	SER	3.5
1	A	198	PRO	3.5
1	B	110	HIS	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	20	HIS	3.3
1	B	195	LEU	3.3
1	A	344	GLY	3.2
1	A	132	ASN	3.1
1	A	1	MET	3.0
1	A	131	ASP	3.0
1	A	63	GLN	3.0
1	A	160	PRO	3.0
1	A	192	ARG	3.0
1	A	161	LYS	2.9
1	B	115	SER	2.8
1	B	133	TYR	2.8
1	A	115	SER	2.7
1	A	199	PHE	2.7
1	A	30	LYS	2.6
1	A	107	ILE	2.5
1	B	118	TYR	2.5
1	A	2	TYR	2.5
1	A	104	SER	2.5
1	B	243	ILE	2.5
1	A	136	PHE	2.5
1	A	195	LEU	2.4
1	A	24	ILE	2.4
1	B	4	ILE	2.4
1	B	299	SER	2.4
1	B	123	PRO	2.4
1	A	17	GLY	2.3
1	A	163	GLY	2.3
1	B	42	THR	2.3
1	A	4	ILE	2.3
1	A	118	TYR	2.2
1	A	66	GLY	2.2
1	A	25	ALA	2.2
1	B	109	GLY	2.2
1	A	27	GLN	2.2
1	B	273	VAL	2.2
1	A	3	ASP	2.2
1	A	106	ALA	2.2
1	A	100	LYS	2.1
1	B	119	ILE	2.1
1	A	103	ASN	2.1

## 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(Å <sup>2</sup> )	Q<0.9
2	MN	A	503	1/1	1.00	0.13	0.61	30,30,30,30	0
2	MN	B	507	1/1	0.99	0.14	0.48	36,36,36,36	0
3	AMP	B	602	23/23	0.95	0.16	0.29	28,48,49,52	0
2	MN	A	504	1/1	0.99	0.12	-0.27	27,27,27,27	0
2	MN	A	501	1/1	0.99	0.12	-0.40	31,31,31,31	0
2	MN	B	506	1/1	0.99	0.12	-0.60	39,39,39,39	0
2	MN	B	505	1/1	1.00	0.12	-0.62	34,34,34,34	0
3	AMP	A	601	23/23	0.98	0.10	-1.66	18,30,34,38	0
2	MN	A	502	1/1	0.96	0.14	-	52,52,52,52	0
2	MN	A	508	1/1	0.84	0.16	-	66,66,66,66	0
2	MN	B	509	1/1	0.82	0.11	-	89,89,89,89	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.