



# Full wwPDB X-ray Structure Validation Report

Nov 10, 2014 – 07:17 PM EST

PDB ID : 4PGF  
Title : The structure of mono-acetylated SAHH  
Authors : Kavran, J.M.; Wang, Y.; Cole, P.A.; Leahy, D.J.  
Deposited on : 2014-05-01  
Resolution : 2.59 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

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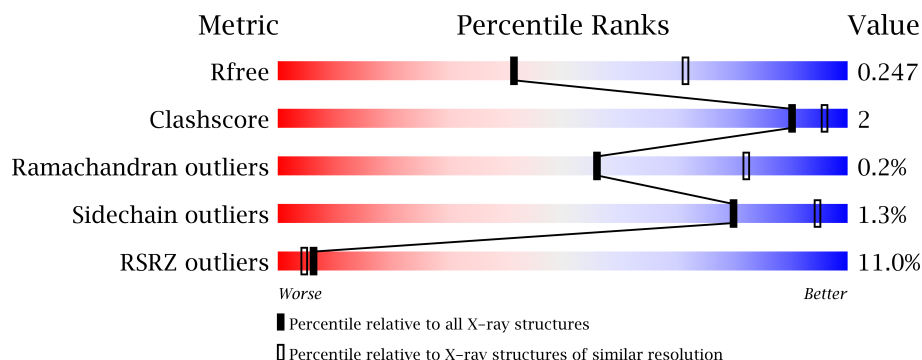
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.16 November 2013  
Xtriage (Phenix) : dev-1439  
EDS : stable24103  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.1.3  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable24103

# 1 Overall quality at a glance



The reported resolution of this entry is 2.59 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))
$R_{free}$	66092	1718 (2.60-2.60)
Clashscore	79885	2154 (2.60-2.60)
Ramachandran outliers	78287	2113 (2.60-2.60)
Sidechain outliers	78261	2113 (2.60-2.60)
RSRZ outliers	66119	1718 (2.60-2.60)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	432	
1	B	432	

## 2 Entry composition i

There are 4 unique types of molecules in this entry. The entry contains 13514 atoms, of which 6715 are hydrogens and 0 are deuterium.

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.

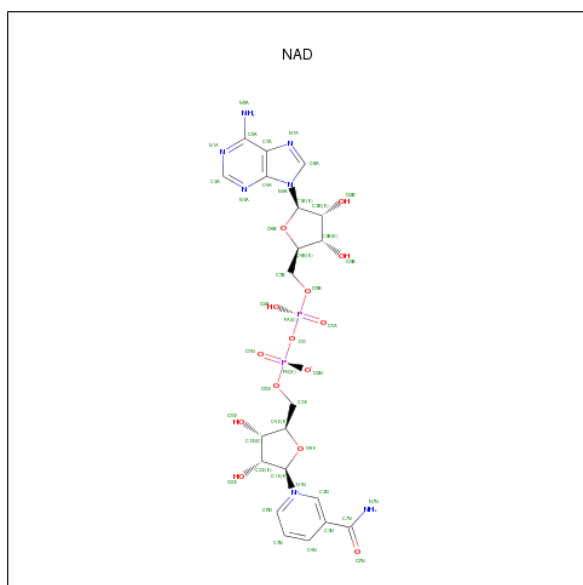
- Molecule 1 is a protein called Adenosylhomocysteinase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	430	Total	C	H	N	O	S	0	1	0
			6671	2116	3335	572	622	26			
1	B	428	Total	C	H	N	O	S	0	0	0
			6615	2102	3302	569	616	26			

There are 4 discrepancies between the modelled and reference sequences:

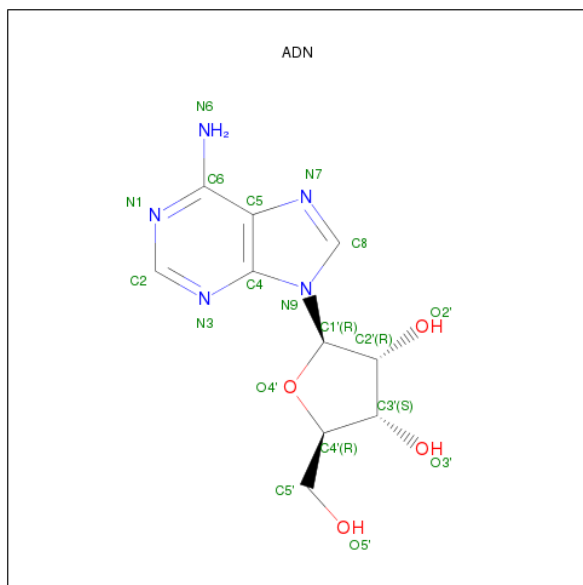
Chain	Residue	Modelled	Actual	Comment	Reference
A	86	ASN	ASP	variant	UNP P23526
A	396	CYS	GLU	engineered mutation	UNP P23526
B	86	ASN	ASP	variant	UNP P23526
B	396	CYS	GLU	engineered mutation	UNP P23526

- 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	H	N	O	0	0
			70	21	26	7	14		
2	B	1	Total	C	H	N	O	0	0
			70	21	26	7	14		

- Molecule 3 is ADENOSINE (three-letter code: ADN) (formula:  $C_{10}H_{13}N_5O_4$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	H	N	O	0	0
			32	10	13	5	4		
3	B	1	Total	C	H	N	O	0	0
			32	10	13	5	4		

- Molecule 4 is water.

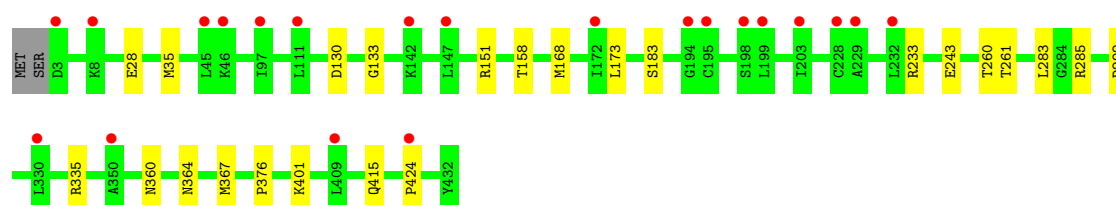
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	13	Total	O	0	0
			13	13		
4	B	11	Total	O	0	0
			11	11		

### 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.

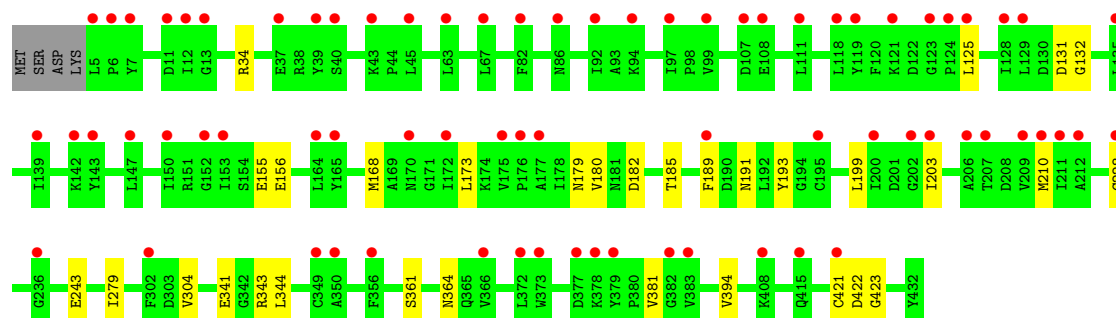
#### • Molecule 1: Adenosylhomocysteinase

Chain A: 



#### • Molecule 1: Adenosylhomocysteinase

Chain B: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	97.60Å 102.74Å 175.02Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.00 – 2.59 44.00 – 2.59	Depositor EDS
% Data completeness (in resolution range)	99.7 (44.00-2.59) 99.7 (44.00-2.59)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.54 (at 2.58Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.4_1496)	Depositor
R, $R_{free}$	0.188 , 0.248 0.188 , 0.247	Depositor DCC
$R_{free}$ test set	1375 reflections (4.98%)	DCC
Wilson B-factor (Å <sup>2</sup> )	61.2	Xtriage
Anisotropy	0.380	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.39 , 37.4	EDS
Estimated twinning fraction	0.012 for k,h,-l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 27636 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	13514	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	73.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.20% of the height of the origin peak. No significant pseudotranslation is detected.*

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

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ADN, NAD, ALY

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	0/3390	0.46	0/4587
1	B	0.28	0/3364	0.45	1/4553 (0.0%)
All	All	0.29	0/6754	0.46	1/9140 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	422	ASP	C-N-CA	8.94	141.07	122.30

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3336	3335	15	11	0
1	B	3313	3302	25	16	0
2	A	44	26	0	2	0
2	B	44	26	0	1	0
3	A	19	13	0	1	0
3	B	19	13	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	13	0	0	1	0
4	B	11	0	0	0	0
All	All	6799	6715	40	28	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 2.

All (28) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:243:GLU:OE2	2:A:501:NAD:O2B	2.20	0.59
1:B:132:GLY:N	1:B:155:GLU:OE2	2.29	0.57
1:A:292:ASP:OD2	1:A:335:ARG:NH2	2.41	0.53
1:B:131:ASP:HA	1:B:155:GLU:OE2	2.10	0.51
1:B:243:GLU:OE2	2:B:501:NAD:O2B	2.30	0.49
1:A:35:MET:CE	1:A:367:MET:HE3	2.43	0.49
1:B:179:ASN:ND2	1:B:182:ASP:OD2	2.46	0.49
1:B:168:MET:HE2	1:B:173:LEU:HB3	1.95	0.47
1:A:283:LEU:HD12	1:A:285:ARG:NH1	2.30	0.47
1:A:151:ARG:HD3	1:A:376:PRO:HG3	1.97	0.47
1:B:279:ILE:HG22	1:B:304:VAL:HB	1.98	0.45
1:B:180:VAL:HG13	1:B:364:ASN:HB3	1.98	0.45
1:B:185:THR:HG21	1:B:394:VAL:HG11	1.99	0.44
1:A:260:THR:OG1	1:A:261:THR:N	2.48	0.44
1:A:28:GLU:OE2	1:A:401:LYS:HE2	2.18	0.43
1:A:360:ASN:O	1:A:364:ASN:ND2	2.49	0.43
1:B:199:LEU:HD22	1:B:228:CYS:SG	2.60	0.42
1:B:156:GLU:OE2	1:B:361:SER:HB3	2.19	0.42
1:B:341:GLU:HB2	1:B:343:ARG:NH1	2.34	0.42
2:A:501:NAD:C4N	3:A:502:ADN:H3'	2.50	0.42
1:B:168:MET:HE1	1:B:381:VAL:HG12	2.02	0.42
1:A:415:GLN:NE2	4:A:612:HOH:O	2.44	0.41
1:B:185:THR:HG21	1:B:394:VAL:CG1	2.50	0.41
1:B:199:LEU:HD11	1:B:203:ILE:HD11	2.02	0.41
1:B:189:PHE:HA	1:B:193:TYR:CD2	2.56	0.41
1:A:130:ASP:OD2	1:A:133:GLY:HA2	2.20	0.40
1:B:131:ASP:CB	1:B:156:GLU:HB3	2.51	0.40
1:A:168:MET:HE2	1:A:173:LEU:HB3	2.04	0.40

There are no symmetry-related clashes.



## 5.3 Torsion angles

### 5.3.1 Protein backbone ⓘ

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	428/432 (99%)	405 (95%)	22 (5%)	1 (0%)	56	82
1	B	425/432 (98%)	402 (95%)	22 (5%)	1 (0%)	56	82
All	All	853/864 (99%)	807 (95%)	44 (5%)	2 (0%)	56	82

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	424	PRO
1	B	423	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	354/355 (100%)	351 (99%)	3 (1%)	89	97
1	B	351/355 (99%)	345 (98%)	6 (2%)	73	92
All	All	705/710 (99%)	696 (99%)	9 (1%)	80	95

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	158	THR
1	A	183	SER
1	A	233	ARG
1	B	34	ARG
1	B	125	LEU

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Mol	Chain	Res	Type
1	B	191	ASN
1	B	210	MET
1	B	344	LEU
1	B	421	CYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 5.3.3 RNA ⓘ

There are no RNA chains in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

2 non-standard protein/DNA/RNA residues 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	$\# Z  > 2$
1	ALY	A	408	1	11,11,12	5.27	2 (18%)	10,12,14	2.27	3 (30%)
1	ALY	B	408	1	11,11,12	5.13	2 (18%)	10,12,14	2.44	4 (40%)

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
1	ALY	A	408	1	-	2/8/10/12	0/0/0/0
1	ALY	B	408	1	-	2/8/10/12	0/0/0/0

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	408	ALY	O-C	17.21	1.23	1.11
1	B	408	ALY	O-C	16.65	1.22	1.11
1	B	408	ALY	CH-NZ	2.58	1.40	1.33
1	A	408	ALY	CH-NZ	2.24	1.39	1.33

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	408	ALY	C-CA-N	-5.78	108.06	113.83
1	A	408	ALY	C-CA-N	-5.56	108.27	113.83
1	B	408	ALY	CE-NZ-CH	3.09	127.41	122.35
1	A	408	ALY	CE-NZ-CH	3.06	127.36	122.35
1	B	408	ALY	CB-CA-N	2.69	118.00	110.35
1	B	408	ALY	CH3-CH-NZ	2.47	119.88	116.15
1	A	408	ALY	CB-CA-N	2.26	116.78	110.35

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	408	ALY	OH-CH-NZ-CE
1	B	408	ALY	OH-CH-NZ-CE
1	B	408	ALY	CH3-CH-NZ-CE
1	A	408	ALY	CH3-CH-NZ-CE

There are no ring outliers.

## 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	501	-	48,48,48	1.57	6 (12%)	73,73,73	1.77	18 (24%)
3	ADN	A	502	-	21,21,21	1.08	3 (14%)	31,31,31	1.04	3 (9%)
2	NAD	B	501	-	48,48,48	1.59	7 (14%)	73,73,73	1.78	16 (21%)
3	ADN	B	502	-	21,21,21	1.23	2 (9%)	31,31,31	0.99	3 (9%)

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	501	-	-	0/30/62/62	0/5/5/5
3	ADN	A	502	-	-	0/6/22/22	0/3/3/3
2	NAD	B	501	-	-	0/30/62/62	0/5/5/5
3	ADN	B	502	-	-	0/6/22/22	0/3/3/3

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	NAD	PN-O5D	6.33	1.80	1.59
2	A	501	NAD	PN-O5D	5.82	1.78	1.59
3	B	502	ADN	C8-N9	4.12	1.42	1.36
2	B	501	NAD	PA-O5B	3.69	1.75	1.59
2	A	501	NAD	PA-O5B	3.60	1.75	1.59
3	A	502	ADN	C8-N9	3.09	1.41	1.36
2	B	501	NAD	C8A-N9A	2.62	1.40	1.36
2	A	501	NAD	C7N-N7N	2.39	1.38	1.33
2	B	501	NAD	C7N-N7N	2.32	1.37	1.33
2	A	501	NAD	C2D-C1D	2.27	1.56	1.53
2	A	501	NAD	C8A-N9A	2.24	1.40	1.36
2	B	501	NAD	C2B-C1B	2.21	1.56	1.53
2	B	501	NAD	O2D-C2D	-2.17	1.37	1.43
2	A	501	NAD	C2N-N1N	2.15	1.38	1.35
3	B	502	ADN	C4-N9	2.14	1.40	1.37
3	A	502	ADN	C2-N3	2.09	1.35	1.32
2	B	501	NAD	C2D-C1D	2.02	1.56	1.53
3	A	502	ADN	C4-N9	2.00	1.40	1.37

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	NAD	PN-O3-PA	-4.85	113.67	133.17
2	A	501	NAD	O3-PN-O1N	4.84	120.47	108.77
2	B	501	NAD	C4B-O4B-C1B	-4.59	104.67	109.72
2	B	501	NAD	O3-PN-O1N	4.29	119.13	108.77
2	A	501	NAD	PN-O3-PA	-4.21	116.22	133.17
2	A	501	NAD	O4D-C1D-N1N	-3.89	103.86	108.13
2	A	501	NAD	C4B-O4B-C1B	-3.84	105.50	109.72
2	B	501	NAD	O2N-PN-O1N	3.53	129.23	118.70
2	A	501	NAD	PN-O5D-C5D	-3.33	108.80	120.35
2	A	501	NAD	O2N-PN-O1N	3.27	128.46	118.70
2	B	501	NAD	O2A-PA-O1A	3.23	130.02	112.14
2	B	501	NAD	PN-O5D-C5D	-3.21	109.24	120.35
2	A	501	NAD	O2A-PA-O1A	3.19	129.78	112.14
2	A	501	NAD	O2N-PN-O5D	-3.05	95.74	108.19
2	B	501	NAD	N3A-C4A-N9A	2.90	130.36	125.39
2	A	501	NAD	N3A-C4A-N9A	2.81	130.21	125.39
2	B	501	NAD	O2N-PN-O3	-2.78	101.17	108.85
2	B	501	NAD	O2N-PN-O5D	-2.72	97.08	108.19
2	A	501	NAD	O2A-PA-O3	2.67	117.83	105.14
2	B	501	NAD	C1B-N9A-C4A	-2.63	122.09	126.64
2	B	501	NAD	C5B-C4B-C3B	-2.63	104.66	115.19
2	A	501	NAD	O2N-PN-O3	-2.63	101.61	108.85
2	B	501	NAD	C8A-N9A-C1B	2.51	130.88	126.15
2	A	501	NAD	C1B-N9A-C4A	-2.50	122.32	126.64
2	B	501	NAD	O2A-PA-O3	2.47	116.88	105.14
2	A	501	NAD	C5B-C4B-C3B	-2.45	105.39	115.19
2	A	501	NAD	O4B-C1B-C2B	-2.43	103.15	106.69
3	A	502	ADN	N3-C4-N9	2.30	129.34	125.39
2	B	501	NAD	O4B-C1B-C2B	-2.30	103.35	106.69
2	A	501	NAD	C3D-C2D-C1D	-2.23	97.43	100.92
2	B	501	NAD	O5D-PN-O1N	-2.22	99.28	108.62
2	A	501	NAD	C8A-N9A-C1B	2.19	130.28	126.15
3	B	502	ADN	N3-C4-N9	2.19	129.14	125.39
2	B	501	NAD	O4D-C1D-N1N	-2.16	105.76	108.13
2	A	501	NAD	O3-PA-O1A	-2.13	95.84	110.43
3	B	502	ADN	C4-C5-N7	2.08	111.42	109.41
2	A	501	NAD	O5D-PN-O1N	-2.08	99.87	108.62
3	A	502	ADN	C8-N9-C1'	2.03	129.97	126.15
3	A	502	ADN	C4'-O4'-C1'	-2.01	107.51	109.72
3	B	502	ADN	C8-N9-C1'	2.01	129.93	126.15

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 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 ⓘ

### 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	430/432 (99%)	0.54	21 (4%) 28 25	44, 58, 72, 106	0
1	B	428/432 (99%)	1.02	73 (17%) 2 1	44, 72, 103, 226	0
All	All	858/864 (99%)	0.78	94 (10%) 6 4	44, 61, 100, 226	0

All (94) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	153	ILE	6.0
1	B	111	LEU	6.0
1	B	378	LYS	5.4
1	B	118	LEU	5.3
1	B	175	VAL	5.2
1	B	124	PRO	4.9
1	B	379	TYR	4.4
1	B	97	ILE	4.0
1	B	86	ASN	3.9
1	B	143	TYR	3.9
1	B	92	ILE	3.8
1	B	129	LEU	3.8
1	B	39	TYR	3.7
1	B	202	GLY	3.7
1	B	150	ILE	3.7
1	B	165	TYR	3.6
1	B	82	PHE	3.6
1	B	177	ALA	3.5
1	B	203	ILE	3.4
1	B	5	LEU	3.4
1	B	408	ALY	3.3
1	A	172	ILE	3.3
1	B	6	PRO	3.3
1	B	99	VAL	3.2

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Mol	Chain	Res	Type	RSRZ
1	B	121	LYS	3.2
1	B	119	TYR	3.2
1	B	372	LEU	3.1
1	B	67	LEU	3.1
1	A	45	LEU	3.0
1	B	377	ASP	3.0
1	B	200	ILE	3.0
1	B	12	ILE	3.0
1	B	11	ASP	3.0
1	B	123	GLY	2.9
1	B	152	GLY	2.9
1	B	45	LEU	2.9
1	B	382	GLY	2.9
1	B	366	VAL	2.9
1	B	108	GLU	2.8
1	A	195	CYS	2.7
1	B	209	VAL	2.7
1	B	13	GLY	2.7
1	A	198	SER	2.7
1	B	40	SER	2.7
1	B	373	TRP	2.7
1	B	164	LEU	2.7
1	B	63	LEU	2.6
1	A	142	LYS	2.6
1	B	172	ILE	2.6
1	A	228	CYS	2.6
1	B	128	ILE	2.6
1	B	349	CYS	2.6
1	B	135	LEU	2.6
1	A	111	LEU	2.5
1	A	3	ASP	2.5
1	B	125	LEU	2.5
1	B	415	GLN	2.5
1	B	356	PHE	2.5
1	B	7	TYR	2.4
1	B	302	PHE	2.4
1	B	212	ALA	2.4
1	A	8	LYS	2.4
1	A	194	GLY	2.4
1	B	421	CYS	2.3
1	A	147	LEU	2.3
1	A	199	LEU	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	206	ALA	2.3
1	B	147	LEU	2.2
1	B	211	ILE	2.2
1	B	350	ALA	2.2
1	B	210	MET	2.2
1	A	424	PRO	2.2
1	B	189	PHE	2.2
1	B	107	ASP	2.2
1	A	409	LEU	2.2
1	A	350	ALA	2.2
1	A	203	ILE	2.1
1	A	330	LEU	2.1
1	B	383	VAL	2.1
1	B	142	LYS	2.1
1	A	46	LYS	2.1
1	B	207	THR	2.1
1	B	139	ILE	2.1
1	B	195	CYS	2.0
1	B	236	GLY	2.0
1	B	94	LYS	2.0
1	B	228	CYS	2.0
1	B	43	LYS	2.0
1	A	229	ALA	2.0
1	B	170	ASN	2.0
1	B	37	GLU	2.0
1	A	232	LEU	2.0
1	A	97	ILE	2.0
1	B	176	PRO	2.0

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

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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
1	ALY	B	408	12/13	0.34	2.02	55,68,79,79	0
1	ALY	A	408	12/13	0.26	0.44	53,73,88,89	0

### 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, 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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	ADN	A	502	19/19	0.22	0.19	56,60,72,73	0
3	ADN	B	502	19/19	0.19	-0.24	71,78,94,96	0
2	NAD	A	501	44/44	0.16	-0.43	47,56,68,70	0
2	NAD	B	501	44/44	0.17	-0.51	53,66,82,87	0

### 6.5 Other polymers ⓘ

There are no such residues in this entry.