



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 10:41 AM GMT

PDB ID : 3MTG
Title : Crystal structure of human S-adenosyl homocysteine hydrolase-like 1 protein
Authors : Wisniewska, M.; Siponen, M.I.; Arrowsmith, C.H.; Berglund, H.; Bountra, C.; Collins, R.; Edwards, A.M.; Flodin, S.; Flores, A.; Graslund, S.; Hammarstrom, M.; Johansson, I.; Karlberg, T.; Kotenyova, T.; Moche, M.; Nordlund, P.; Nyman, T.; Persson, C.; Schutz, P.; Thorsell, A.G.; Tresaugues, L.; van der Berg, S.; Wahlberg, E.; Weigelt, J.; Welin, M.; Schuler, H.; Structural Genomics Consortium (SGC)
Deposited on : 2010-04-30
Resolution : 2.64 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

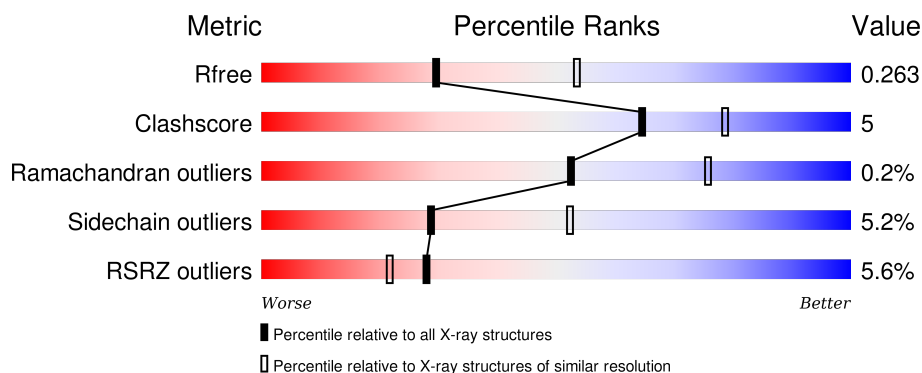
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.64 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	3377 (2.68-2.60)
Clashscore	102246	3781 (2.68-2.60)
Ramachandran outliers	100387	3722 (2.68-2.60)
Sidechain outliers	100360	3722 (2.68-2.60)
RSRZ outliers	91569	3388 (2.68-2.60)

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 ≥ 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	444	<div> <div>5%</div> <div>80%</div> <div>16%</div> <div>• •</div> </div>
1	B	444	<div> <div>5%</div> <div>84%</div> <div>10%</div> <div>• 5%</div> </div>

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 6792 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.

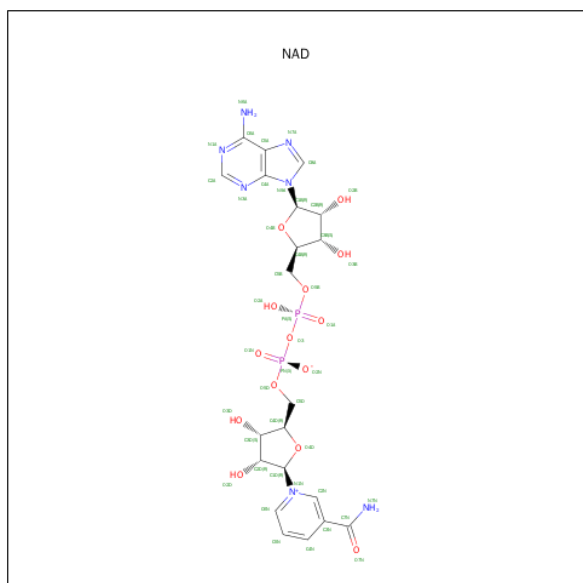
- Molecule 1 is a protein called Putative adenosylhomocysteinase 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	430	Total	C	N	O	S	0	0	0
			3322	2095	576	624	27			
1	B	423	Total	C	N	O	S	0	0	0
			3268	2063	566	613	26			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	87	SER	-	EXPRESSION TAG	UNP O43865
A	88	MET	-	EXPRESSION TAG	UNP O43865
A	508	ALA	THR	ENGINEERED MUTATION	UNP O43865
B	87	SER	-	EXPRESSION TAG	UNP O43865
B	88	MET	-	EXPRESSION TAG	UNP O43865
B	508	ALA	THR	ENGINEERED MUTATION	UNP O43865

- 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		

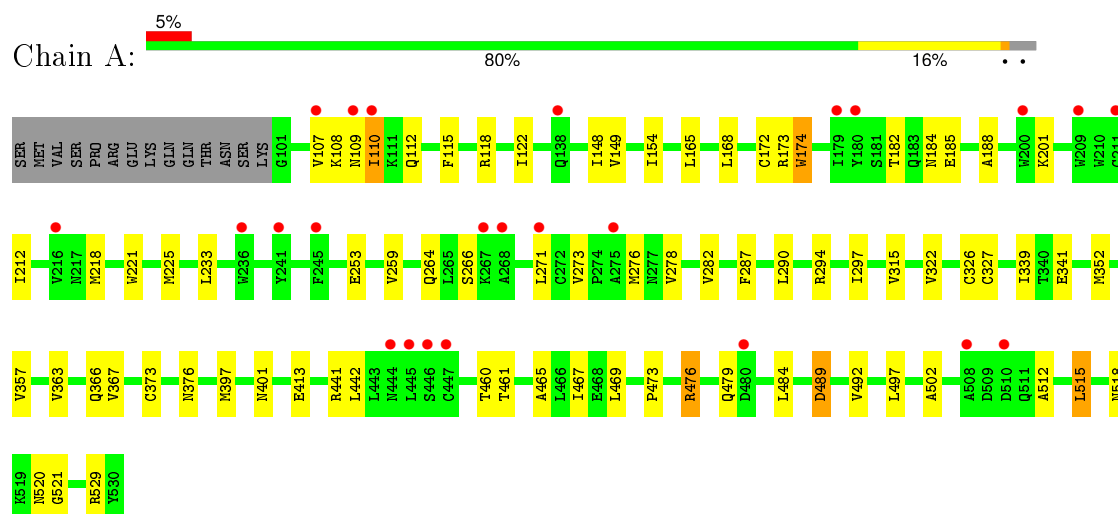
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	55	Total	O	0	0
			55	55		
3	B	59	Total	O	0	0
			59	59		

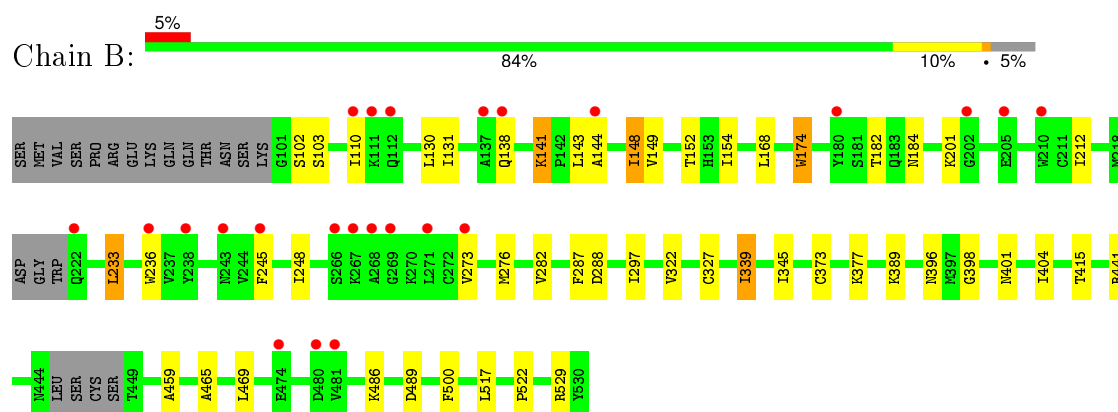
3 Residue-property plots [i](#)

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 ($\text{RSRZ} > 2$). 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: Putative adenosylhomocysteinase 2



• Molecule 1: Putative adenosylhomocysteinase 2



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	186.24Å 68.60Å 90.37Å 90.00° 115.27° 90.00°	Depositor
Resolution (Å)	33.59 – 2.64 32.73 – 2.64	Depositor EDS
% Data completeness (in resolution range)	(Not available) (33.59-2.64) 99.0 (32.73-2.64)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.84 (at 2.65Å)	Xtriage
Refinement program	BUSTER 2.9.2	Depositor
R, R_{free}	0.203 , 0.249 0.208 , 0.263	Depositor DCC
R_{free} test set	628 reflections (2.12%)	DCC
Wilson B-factor (Å ²)	53.1	Xtriage
Anisotropy	0.494	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 58.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 30251 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6792	wwPDB-VP
Average B, all atoms (Å ²)	83.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.66% 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: 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	0.51	0/3383	0.70	0/4592
1	B	0.48	0/3325	0.68	0/4509
All	All	0.50	0/6708	0.69	0/9101

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

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	3322	0	3288	49	0
1	B	3268	0	3243	24	0
2	A	44	0	26	1	0
2	B	44	0	26	1	0
3	A	55	0	0	0	0
3	B	59	0	0	1	0
All	All	6792	0	6583	68	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 5.

All (68) 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:110:ILE:O	1:A:110:ILE:HG22	1.58	1.01
1:A:109:ASN:HD21	1:A:184:ASN:HB3	1.32	0.92
1:A:148:ILE:HG22	1:A:225:MET:HB2	1.67	0.77
1:A:110:ILE:CG2	1:A:110:ILE:O	2.30	0.75
1:A:109:ASN:ND2	1:A:112:GLN:OE1	2.22	0.72
1:B:327:CYS:SG	1:B:339:ILE:HD11	2.32	0.70
1:B:245:PHE:HA	1:B:248:ILE:HD12	1.78	0.65
1:B:149:VAL:HG12	1:B:233:LEU:HD21	1.77	0.65
1:A:109:ASN:HB3	1:A:112:GLN:HB2	1.83	0.61
1:B:152:THR:HG22	1:B:174:TRP:HE1	1.66	0.60
1:A:109:ASN:ND2	1:A:184:ASN:HB3	2.12	0.60
1:A:109:ASN:CB	1:A:188:ALA:HB2	2.33	0.58
1:A:518:ASN:HB2	1:A:521:GLY:HA3	1.86	0.57
1:A:148:ILE:HD12	1:A:174:TRP:HZ3	1.70	0.56
1:B:212:ILE:HG23	1:B:233:LEU:HD12	1.87	0.56
1:A:401:ASN:HB3	1:A:441:ARG:HG2	1.87	0.55
1:A:148:ILE:HD11	1:A:172:CYS:SG	2.47	0.54
1:A:327:CYS:SG	1:A:339:ILE:HD11	2.47	0.54
1:A:352:MET:HE3	1:B:500:PHE:HZ	1.73	0.53
1:A:149:VAL:HG12	1:A:233:LEU:HD21	1.91	0.52
1:A:271:LEU:O	1:A:479:GLN:HB3	2.09	0.52
1:A:276:MET:HB2	1:A:469:LEU:HD11	1.91	0.52
1:A:326:CYS:SG	1:A:397:MET:HE1	2.50	0.52
1:A:276:MET:HG2	1:A:465:ALA:HB1	1.94	0.50
1:A:148:ILE:HD13	1:A:165:LEU:HD13	1.93	0.50
1:A:109:ASN:HB3	1:A:188:ALA:HB2	1.94	0.50
1:B:276:MET:HB2	1:B:469:LEU:HD11	1.94	0.49
1:A:212:ILE:HG23	1:A:233:LEU:HD12	1.94	0.49
1:A:297:ILE:HD11	1:A:322:VAL:HG13	1.94	0.49
1:B:130:LEU:HD11	1:B:459:ALA:HB1	1.95	0.48
1:A:107:VAL:HB	1:A:109:ASN:OD1	2.12	0.48
1:A:352:MET:CE	1:B:500:PHE:CZ	2.97	0.48
1:A:115:PHE:CE2	1:A:185:GLU:HG2	2.48	0.48
1:B:154:ILE:HB	1:B:182:THR:HG23	1.94	0.47
1:A:118:ARG:O	1:A:122:ILE:HG12	2.14	0.47
1:A:489:ASP:HB3	1:B:345:ILE:HD11	1.95	0.47
1:A:115:PHE:HE2	1:A:185:GLU:HG2	1.80	0.47
1:B:143:LEU:HD12	1:B:168:LEU:HB2	1.97	0.46
1:A:352:MET:HE3	1:B:500:PHE:CZ	2.50	0.46
1:A:512:ALA:O	1:A:515:LEU:O	2.34	0.45
1:A:168:LEU:HD13	1:A:467:ILE:HD11	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:290:LEU:HD22	1:A:294:ARG:CZ	2.46	0.45
1:A:185:GLU:H	1:A:185:GLU:CD	2.20	0.45
1:B:141:LYS:HB3	1:B:144:ALA:HB2	1.98	0.44
1:A:109:ASN:HD21	1:A:184:ASN:CB	2.16	0.44
1:B:297:ILE:HD11	1:B:322:VAL:HG13	2.00	0.44
1:A:460:THR:HG22	1:A:492:VAL:HG22	2.00	0.44
1:A:461:THR:HG23	1:A:484:LEU:HD22	1.99	0.44
1:A:154:ILE:HB	1:A:182:THR:HG23	1.99	0.43
1:B:148:ILE:HD12	1:B:174:TRP:HZ3	1.83	0.43
1:B:401:ASN:HB3	1:B:441:ARG:HG3	2.01	0.43
1:B:396:ASN:HB2	1:B:404:ILE:HD12	2.00	0.42
1:A:363:VAL:O	1:A:366:GLN:HG2	2.18	0.42
1:B:184:ASN:ND2	1:B:201:LYS:H	2.18	0.42
1:A:184:ASN:HD21	1:A:201:LYS:H	1.68	0.42
1:A:148:ILE:HG13	1:A:172:CYS:HA	2.02	0.42
1:B:486:LYS:HB3	1:B:522:PRO:HB3	2.01	0.42
1:A:473:PRO:HD2	1:A:476:ARG:HD2	2.02	0.42
1:B:398:GLY:HA2	2:B:1:NAD:O4D	2.21	0.41
1:A:397:MET:O	2:A:1:NAD:H2N	2.21	0.41
1:A:339:ILE:O	1:A:357:VAL:HA	2.21	0.41
1:A:315:VAL:HG12	1:A:367:VAL:HG21	2.01	0.41
1:B:276:MET:HG2	1:B:465:ALA:HB1	2.02	0.41
1:A:529:ARG:HG2	1:B:529:ARG:HA	2.03	0.41
1:A:497:LEU:HB3	1:A:502:ALA:HB3	2.01	0.41
1:A:218:MET:HB3	1:A:221:TRP:HB3	2.01	0.41
1:B:389:LYS:HG2	3:B:59:HOH:O	2.20	0.40
1:A:253:GLU:O	1:A:278:VAL:HG22	2.21	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	428/444 (96%)	406 (95%)	21 (5%)	1 (0%)	52	76
1	B	417/444 (94%)	389 (93%)	27 (6%)	1 (0%)	52	76
All	All	845/888 (95%)	795 (94%)	48 (6%)	2 (0%)	52	76

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	282	VAL
1	A	282	VAL

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	357/378 (94%)	339 (95%)	18 (5%)	30	54
1	B	351/378 (93%)	332 (95%)	19 (5%)	27	50
All	All	708/756 (94%)	671 (95%)	37 (5%)	29	52

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	108	LYS
1	A	110	ILE
1	A	173	ARG
1	A	174	TRP
1	A	259	VAL
1	A	264	GLN
1	A	266	SER
1	A	273	VAL
1	A	287	PHE
1	A	341	GLU
1	A	373	CYS
1	A	376	ASN
1	A	413	GLU
1	A	442	LEU

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Mol	Chain	Res	Type
1	A	476	ARG
1	A	489	ASP
1	A	515	LEU
1	A	520	ASN
1	B	102	SER
1	B	103	SER
1	B	110	ILE
1	B	131	ILE
1	B	138	GLN
1	B	141	LYS
1	B	148	ILE
1	B	174	TRP
1	B	233	LEU
1	B	236	TRP
1	B	273	VAL
1	B	287	PHE
1	B	288	ASP
1	B	339	ILE
1	B	373	CYS
1	B	377	LYS
1	B	415	THR
1	B	489	ASP
1	B	517	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	109	ASN
1	A	224	ASN
1	A	376	ASN
1	A	479	GLN
1	A	503	HIS
1	A	520	ASN
1	B	224	ASN
1	B	376	ASN

5.3.3 RNA ⓘ

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](#)

2 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	1	-	38,48,48	1.66	3 (7%)	47,73,73	2.28	5 (10%)
2	NAD	B	1	-	38,48,48	1.65	3 (7%)	47,73,73	2.13	5 (10%)

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	1	-	-	0/22/62/62	0/5/5/5
2	NAD	B	1	-	-	0/22/62/62	0/5/5/5

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1	NAD	C2A-N1A	2.86	1.39	1.33
2	A	1	NAD	C2A-N1A	2.87	1.39	1.33
2	B	1	NAD	C2A-N3A	3.80	1.38	1.32
2	A	1	NAD	C2A-N3A	3.82	1.39	1.32
2	A	1	NAD	O7N-C7N	7.88	1.41	1.24
2	B	1	NAD	O7N-C7N	7.98	1.41	1.24

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	A	1	NAD	N3A-C2A-N1A	-12.75	119.13	128.89
2	B	1	NAD	N3A-C2A-N1A	-11.83	119.84	128.89
2	A	1	NAD	PN-O3-PA	-2.98	124.36	132.73
2	A	1	NAD	O7N-C7N-C3N	-2.52	116.83	119.59
2	B	1	NAD	O7N-C7N-C3N	-2.21	117.17	119.59
2	B	1	NAD	C2B-C1B-N9A	-2.16	110.99	114.29
2	A	1	NAD	C3N-C7N-N7N	3.08	121.19	117.82
2	B	1	NAD	C3N-C7N-N7N	3.15	121.26	117.82
2	B	1	NAD	O4D-C1D-N1N	3.43	111.90	108.13
2	A	1	NAD	O4D-C1D-N1N	4.81	113.42	108.13

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	A	1	NAD	1	0
2	B	1	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	430/444 (96%)	0.16	24 (5%) 28 22	30, 84, 127, 149	0
1	B	423/444 (95%)	0.20	24 (5%) 27 21	28, 81, 139, 170	0
All	All	853/888 (96%)	0.18	48 (5%) 28 22	28, 82, 134, 170	0

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	222	GLN	8.8
1	B	269	GLY	5.6
1	B	111	LYS	5.4
1	B	137	ALA	5.3
1	A	444	ASN	5.2
1	A	445	LEU	5.1
1	B	267	LYS	4.9
1	A	446	SER	4.8
1	B	474	GLU	4.7
1	B	202	GLY	4.1
1	A	275	ALA	4.1
1	B	480	ASP	4.0
1	A	447	CYS	4.0
1	B	273	VAL	3.9
1	B	245	PHE	3.6
1	B	481	VAL	3.6
1	B	271	LEU	3.6
1	B	236	TRP	3.4
1	A	267	LYS	3.2
1	A	180	TYR	3.2
1	A	510	ASP	3.1
1	B	268	ALA	3.1
1	B	210	TRP	3.0
1	A	109	ASN	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	200	TRP	2.8
1	B	144	ALA	2.7
1	B	266	SER	2.7
1	A	110	ILE	2.6
1	B	243	ASN	2.6
1	B	112	GLN	2.6
1	B	138	GLN	2.6
1	A	236	TRP	2.5
1	A	480	ASP	2.5
1	A	179	ILE	2.5
1	A	209	TRP	2.5
1	A	138	GLN	2.4
1	A	508	ALA	2.3
1	B	110	ILE	2.3
1	A	211	CYS	2.3
1	A	268	ALA	2.3
1	A	271	LEU	2.1
1	B	238	TYR	2.1
1	A	107	VAL	2.1
1	A	216	VAL	2.1
1	A	241	TYR	2.1
1	B	205	GLU	2.1
1	A	245	PHE	2.0
1	B	180	TYR	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 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	B	1	44/44	0.98	0.16	0.04	42,52,59,60	0
2	NAD	A	1	44/44	0.97	0.16	-0.15	47,55,65,67	0

6.5 Other polymers [i](#)

There are no such residues in this entry.