



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 27, 2014 – 03:21 PM GMT

PDB ID : 2CDA
Title : SULFOLOBUS SOLFATARICUS GLUCOSE DEHYDROGENASE 1 IN
COMPLEX WITH NADP
Authors : Milburn, C.C.; Lamble, H.J.; Theodossis, A.; Hough, D.W.; Danson, M.J.;
Taylor, G.L.
Deposited on : 2006-01-23
Resolution : 2.28 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

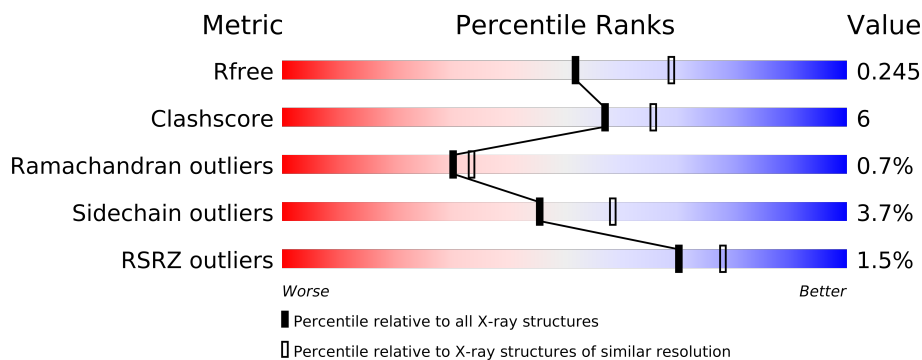
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.28 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	3861 (2.30-2.26)
Clashscore	79885	4801 (2.30-2.26)
Ramachandran outliers	78287	4729 (2.30-2.26)
Sidechain outliers	78261	4728 (2.30-2.26)
RSRZ outliers	66119	3864 (2.30-2.26)

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. 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	366	
1	B	366	

2 Entry composition i

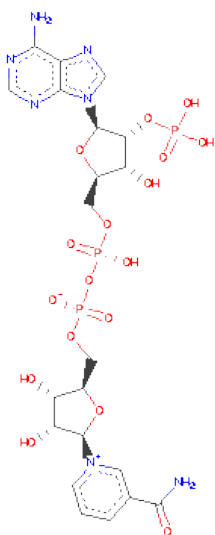
There are 4 unique types of molecules in this entry. The entry contains 5913 atoms, of which 0 are hydrogen and 0 are deuterium.

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 GLUCOSE DEHYDROGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	360	Total	C	N	O	S	0	0	0
			2833	1813	482	525	13			
1	B	359	Total	C	N	O	S	0	0	0
			2824	1807	480	524	13			

- Molecule 2 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDEPHOSPHATE (three-letter code: NAP) (formula: $C_{21}H_{28}N_7O_{17}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	B	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	2	Total 2	Zn 2	0	0
3	A	2	Total 2	Zn 2	0	0

- Molecule 4 is water.

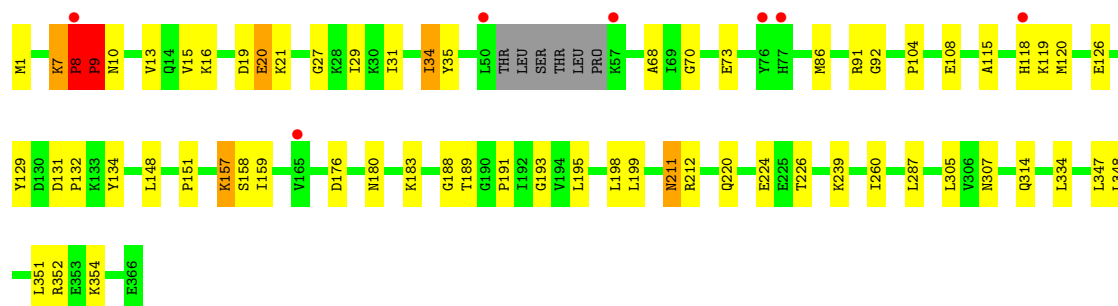
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	75	Total 75	O 75	0	0
4	B	81	Total 81	O 81	0	0

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.

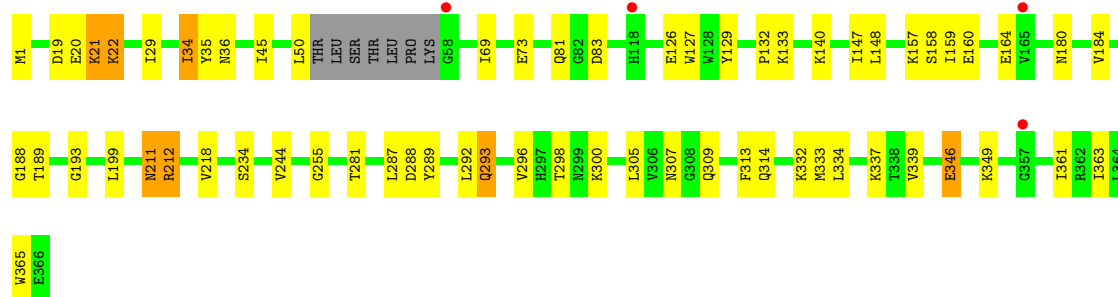
• Molecule 1: GLUCOSE DEHYDROGENASE

Chain A: 



• Molecule 1: GLUCOSE DEHYDROGENASE

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	68.62Å 92.30Å 138.26Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	138.68 – 2.28 48.70 – 2.28	Depositor EDS
% Data completeness (in resolution range)	98.1 (138.68-2.28) 98.2 (48.70-2.28)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.37 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.196 , 0.246 0.196 , 0.245	Depositor DCC
R_{free} test set	2008 reflections (5.27%)	DCC
Wilson B-factor (Å ²)	19.5	Xtriage
Anisotropy	0.842	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 15.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtriage
Outliers	0 of 40085 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5913	wwPDB-VP
Average B, all atoms (Å ²)	20.0	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 22.87 % 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 5.2223e-03.*

¹Intensities 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: ZN, NAP

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	0.91	1/2886 (0.0%)	0.87	6/3899 (0.2%)
1	B	0.87	1/2877 (0.0%)	0.86	1/3888 (0.0%)
All	All	0.89	2/5763 (0.0%)	0.87	7/7787 (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	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	108	GLU	CD-OE2	5.27	1.31	1.25
1	B	184	VAL	CB-CG1	5.21	1.63	1.52

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	83	ASP	CB-CG-OD1	7.04	124.64	118.30
1	A	176	ASP	CB-CG-OD1	6.58	124.22	118.30
1	A	9	PRO	CA-N-CD	-5.66	103.58	111.50
1	A	9	PRO	N-CA-CB	5.50	109.91	103.30
1	A	8	PRO	N-CA-C	5.47	126.33	112.10
1	A	189	THR	C-N-CA	-5.13	111.52	122.30
1	A	7	LYS	O-C-N	5.10	130.80	121.10

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	255	GLY	Peptide

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	2833	0	2887	36	0
1	B	2824	0	2874	43	0
2	A	48	0	25	0	0
2	B	48	0	25	0	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
4	A	75	0	0	1	0
4	B	81	0	0	2	0
All	All	5913	0	5811	74	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 6.

All (74) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:29:ILE:HD11	1:A:132:PRO:HG3	1.37	1.04
1:A:180:ASN:HD21	1:B:314:GLN:HE22	1.05	0.98
1:A:183:LYS:HE3	4:A:2047:HOH:O	1.63	0.97
1:A:211:ASN:HD22	1:A:212:ARG:H	1.25	0.82
1:A:314:GLN:HE22	1:B:180:ASN:HD21	1.26	0.80
1:B:19:ASP:HB3	1:B:22:LYS:HD3	1.67	0.77
1:B:148:LEU:HD11	1:B:334:LEU:HB3	1.66	0.76
1:B:36:ASN:ND2	1:B:365:TRP:HE1	1.93	0.66
1:A:159:ILE:HD12	1:A:199:LEU:HD23	1.80	0.64
1:A:180:ASN:HD21	1:B:314:GLN:NE2	1.88	0.63
1:A:211:ASN:HD22	1:A:212:ARG:N	1.96	0.63
1:B:34:ILE:HG12	1:B:35:TYR:CE1	2.34	0.63
1:B:19:ASP:HB3	1:B:22:LYS:CD	2.29	0.63
1:A:148:LEU:HD11	1:A:334:LEU:HB3	1.80	0.62

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:1:MET:HG3	1:A:126:GLU:HB2	1.81	0.62
1:B:332:LYS:HE3	4:B:2071:HOH:O	1.99	0.62
1:B:158:SER:HB3	1:B:305:LEU:HD13	1.83	0.60
1:B:189:THR:OG1	1:B:211:ASN:ND2	2.38	0.56
1:A:92:GLY:HA2	1:A:104:PRO:HB3	1.88	0.56
1:B:35:TYR:HA	1:B:363:ILE:O	2.05	0.55
1:B:29:ILE:HG13	1:B:132:PRO:HD3	1.89	0.54
1:A:15:VAL:HG21	1:A:347:LEU:HD23	1.89	0.54
1:B:212:ARG:HG2	1:B:234:SER:HA	1.89	0.53
1:B:339:VAL:HG13	1:B:346:GLU:CG	2.38	0.52
1:B:36:ASN:HD22	1:B:365:TRP:HE1	1.57	0.51
1:A:158:SER:HB3	1:A:305:LEU:HD13	1.93	0.50
1:A:131:ASP:OD1	1:A:132:PRO:HD2	2.12	0.50
1:B:292:LEU:O	1:B:296:VAL:HG23	2.13	0.49
1:A:348:LEU:O	1:A:352:ARG:HD2	2.13	0.49
1:A:188:GLY:O	1:A:193:GLY:HA3	2.13	0.49
1:B:160:GLU:O	1:B:164:GLU:HG3	2.12	0.49
1:B:339:VAL:HG13	1:B:346:GLU:HG2	1.95	0.48
1:A:191:PRO:O	1:A:195:LEU:HD12	2.13	0.48
1:A:29:ILE:O	1:A:129:TYR:HA	2.12	0.48
1:A:180:ASN:ND2	1:B:314:GLN:HE22	1.90	0.48
1:A:9:PRO:HG2	1:A:10:ASN:OD1	2.14	0.48
1:A:91:ARG:HG2	1:A:134:TYR:CE1	2.49	0.48
1:A:115:ALA:HA	1:A:120:MET:HB3	1.95	0.48
1:B:19:ASP:O	1:B:22:LYS:HG2	2.14	0.47
1:B:1:MET:HG2	1:B:126:GLU:HB2	1.97	0.47
1:B:73:GLU:O	1:B:81:GLN:HG3	2.15	0.47
1:A:19:ASP:OD1	1:A:21:LYS:HG3	2.15	0.47
1:A:157:LYS:HD3	1:A:157:LYS:C	2.36	0.46
1:B:69:ILE:HD11	1:B:147:ILE:N	2.31	0.46
1:A:20:GLU:H	1:A:20:GLU:HG2	1.40	0.45
1:A:7:LYS:HA	1:A:8:PRO:HD2	1.77	0.45
1:B:298:THR:OG1	1:B:300:LYS:HE3	2.17	0.45
1:B:157:LYS:HD3	1:B:157:LYS:C	2.36	0.45
1:B:337:LYS:HG2	1:B:361:ILE:HD12	1.98	0.44
1:A:29:ILE:CD1	1:A:132:PRO:HG3	2.28	0.44
1:B:188:GLY:O	1:B:193:GLY:HA3	2.17	0.44
1:A:118:HIS:NE2	1:A:119:LYS:HE2	2.32	0.44
1:A:34:ILE:HG12	1:A:35:TYR:CE2	2.51	0.44
1:A:198:LEU:HD23	1:A:226:THR:HG21	1.99	0.44
1:B:21:LYS:HE3	4:B:2004:HOH:O	2.18	0.44
1:B:339:VAL:CG1	1:B:346:GLU:HG3	2.48	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:314:GLN:NE2	1:B:180:ASN:HD21	2.05	0.43
1:A:13:VAL:HG21	1:A:351:LEU:HD13	2.00	0.43
1:B:45:ILE:HG13	1:B:50:LEU:HD12	1.99	0.43
1:B:159:ILE:HD12	1:B:199:LEU:HD23	2.01	0.43
1:B:1:MET:CE	1:B:20:GLU:HB2	2.48	0.43
1:B:189:THR:H	1:B:211:ASN:HD21	1.68	0.42
1:B:289:TYR:O	1:B:293:GLN:HB2	2.20	0.42
1:A:27:GLY:O	1:A:129:TYR:HB3	2.19	0.42
1:B:50:LEU:HA	1:B:50:LEU:HD23	1.94	0.42
1:B:127:TRP:CZ3	1:B:129:TYR:CE2	3.08	0.41
1:B:45:ILE:CG1	1:B:50:LEU:HD12	2.51	0.41
1:B:148:LEU:CD1	1:B:334:LEU:HD13	2.51	0.41
1:B:218:VAL:HG22	1:B:333:MET:HB3	2.02	0.41
1:B:309:GLN:O	1:B:313:PHE:HD1	2.03	0.41
1:B:349:LYS:HA	1:B:349:LYS:HD2	1.89	0.41
1:A:31:ILE:HG22	1:A:70:GLY:HA3	2.03	0.40
1:A:68:ALA:O	1:A:86:MET:HG3	2.22	0.40
1:A:220:GLN:HG3	1:A:224:GLU:OE2	2.21	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	356/366 (97%)	339 (95%)	13 (4%)	4 (1%)	21	19
1	B	355/366 (97%)	337 (95%)	17 (5%)	1 (0%)	50	59
All	All	711/732 (97%)	676 (95%)	30 (4%)	5 (1%)	30	33

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	8	PRO
1	A	9	PRO

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Mol	Chain	Res	Type
1	A	307	ASN
1	B	307	ASN
1	A	260	ILE

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	312/318 (98%)	302 (97%)	10 (3%)	51	66
1	B	311/318 (98%)	298 (96%)	13 (4%)	40	51
All	All	623/636 (98%)	600 (96%)	23 (4%)	45	58

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	LYS
1	A	20	GLU
1	A	34	ILE
1	A	73	GLU
1	A	151	PRO
1	A	157	LYS
1	A	211	ASN
1	A	239	LYS
1	A	287	LEU
1	A	354	LYS
1	B	21	LYS
1	B	22	LYS
1	B	34	ILE
1	B	133	LYS
1	B	140	LYS
1	B	211	ASN
1	B	212	ARG
1	B	244	VAL
1	B	281	THR
1	B	287	LEU
1	B	288	ASP
1	B	293	GLN

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Mol	Chain	Res	Type
1	B	346	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	47	ASN
1	A	211	ASN
1	B	14	GLN
1	B	36	ASN
1	B	180	ASN
1	B	211	ASN
1	B	259	ASN
1	B	263	ASN
1	B	293	GLN
1	B	297	HIS
1	B	314	GLN

5.3.3 RNA ⓘ

There are no RNA chains 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 6 ligands modelled in this entry, 4 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
2	NAP	A	1367	-	52,52,52	2.03	10 (19%)	80,80,80	1.93	14 (17%)
2	NAP	B	1367	-	52,52,52	1.77	9 (17%)	80,80,80	1.93	17 (21%)

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	NAP	A	1367	-	-	0/35/67/67	0/3/5/5
2	NAP	B	1367	-	-	0/35/67/67	0/3/5/5

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1367	NAP	C2N-N1N	6.29	1.43	1.35
2	B	1367	NAP	P2B-O2B	6.22	1.79	1.59
2	B	1367	NAP	P2B-O1X	5.16	1.68	1.51
2	A	1367	NAP	P2B-O1X	4.89	1.67	1.51
2	B	1367	NAP	C2N-N1N	4.77	1.41	1.35
2	A	1367	NAP	P2B-O2B	4.45	1.73	1.59
2	A	1367	NAP	PA-O3	4.25	1.67	1.59
2	B	1367	NAP	O4D-C1D	4.25	1.47	1.41
2	A	1367	NAP	PN-O3	3.73	1.68	1.60
2	A	1367	NAP	C4A-N9A	-3.63	1.32	1.37
2	A	1367	NAP	C3N-C7N	3.54	1.56	1.50
2	A	1367	NAP	P2B-O2X	3.06	1.65	1.54
2	A	1367	NAP	C6N-N1N	3.05	1.44	1.35
2	B	1367	NAP	PN-O5D	2.94	1.69	1.60
2	B	1367	NAP	C4A-N9A	-2.78	1.33	1.37
2	B	1367	NAP	P2B-O2X	2.69	1.64	1.54
2	A	1367	NAP	O4D-C1D	2.51	1.45	1.41
2	B	1367	NAP	C6N-N1N	2.12	1.41	1.35
2	B	1367	NAP	PA-O3	2.06	1.63	1.59

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1367	NAP	N3A-C2A-N1A	-9.39	120.86	128.71
2	A	1367	NAP	N3A-C2A-N1A	-9.14	121.07	128.71
2	A	1367	NAP	O4B-C1B-N9A	-6.87	102.05	108.44
2	B	1367	NAP	O4B-C1B-N9A	-6.26	102.62	108.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1367	NAP	C4B-O4B-C1B	-4.52	104.84	109.75
2	B	1367	NAP	N3A-C4A-N9A	4.03	132.71	125.43
2	A	1367	NAP	N3A-C4A-N9A	4.02	132.70	125.43
2	B	1367	NAP	C4B-O4B-C1B	-3.57	105.87	109.75
2	A	1367	NAP	C8A-N9A-C4A	3.37	109.47	106.90
2	A	1367	NAP	C6N-N1N-C2N	-3.26	118.36	122.04
2	B	1367	NAP	C4A-C5A-N7A	-2.95	106.99	109.52
2	B	1367	NAP	O4D-C1D-N1N	2.95	110.97	107.95
2	A	1367	NAP	C3B-C2B-C1B	-2.92	97.05	102.73
2	B	1367	NAP	O4D-C1D-C2D	-2.85	102.41	106.77
2	A	1367	NAP	C3N-C7N-N7N	-2.84	114.54	117.77
2	A	1367	NAP	O3X-P2B-O2B	2.82	115.22	107.09
2	B	1367	NAP	O3X-P2B-O2B	2.80	115.15	107.09
2	B	1367	NAP	O4D-C4D-C5D	-2.72	99.66	109.36
2	B	1367	NAP	C5A-C4A-N3A	-2.71	119.79	125.70
2	B	1367	NAP	C2A-N3A-C4A	2.66	121.59	114.01
2	A	1367	NAP	P2B-O2B-C2B	2.47	127.16	121.96
2	A	1367	NAP	C2A-N1A-C6A	2.34	123.00	118.77
2	B	1367	NAP	C4D-O4D-C1D	-2.26	107.30	109.75
2	B	1367	NAP	C1B-N9A-C4A	-2.20	122.83	126.64
2	A	1367	NAP	O4D-C1D-C2D	-2.17	103.44	106.77
2	A	1367	NAP	N7A-C8A-N9A	-2.15	108.27	114.36
2	B	1367	NAP	C5B-C4B-C3B	-2.15	106.60	115.21
2	B	1367	NAP	O2X-P2B-O1X	-2.06	103.72	110.44
2	A	1367	NAP	C5A-C4A-N3A	-2.06	121.22	125.70
2	B	1367	NAP	O3-PN-O5D	-2.04	93.19	101.36
2	B	1367	NAP	N7A-C8A-N9A	-2.02	108.63	114.36

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, 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	360/366 (98%)	0.02	7 (1%) 64 72	4, 19, 37, 57	1 (0%)
1	B	359/366 (98%)	-0.01	4 (1%) 77 85	5, 19, 36, 44	0
All	All	719/732 (98%)	0.00	11 (1%) 70 78	4, 19, 36, 57	1 (0%)

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	118	HIS	4.4
1	A	8	PRO	3.9
1	A	76	TYR	3.9
1	A	50	LEU	3.5
1	B	357	GLY	3.1
1	B	58	GLY	3.0
1	A	77	HIS	2.6
1	A	57	LYS	2.4
1	A	165	VAL	2.2
1	B	165	VAL	2.1
1	B	118	HIS	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	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	NAP	A	1367	48/48	0.12	0.07	11,21,50,50	0
2	NAP	B	1367	48/48	0.12	-0.33	16,24,46,48	0
3	ZN	A	1369	1/1	0.08	-0.92	21,21,21,21	0
3	ZN	B	1369	1/1	0.05	-1.72	19,19,19,19	0
3	ZN	A	1368	1/1	0.06	-2.29	35,35,35,35	1
3	ZN	B	1368	1/1	0.06	-3.25	30,30,30,30	1

6.5 Other polymers

There are no such residues in this entry.