



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 10:06 PM GMT

PDB ID : 1R6A
Title : Structure of the dengue virus 2'O methyltransferase in complex with s-adenosyl
homocysteine and ribavirin 5' triphosphate
Authors : Benarroch, D.; Egloff, M.P.; Mulard, L.; Romette, J.L.; Canard, B.
Deposited on : 2003-10-15
Resolution : 2.60 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>

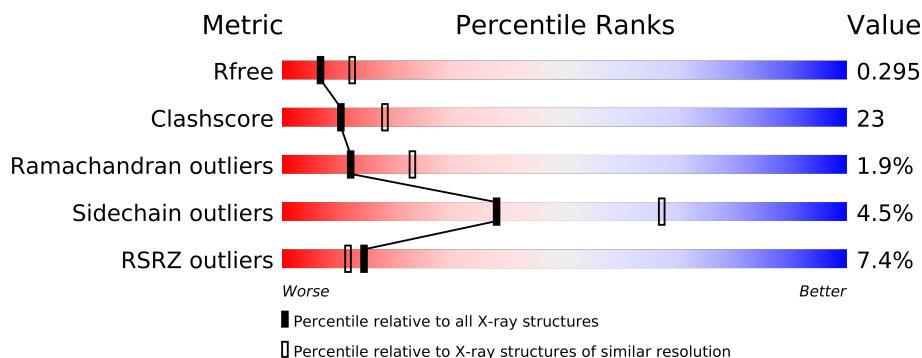
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.60 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 ≥ 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	295	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
2	SO4	A	902	-	X
2	SO4	A	905	-	X
2	SO4	A	906	-	X
2	SO4	A	907	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 2197 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.

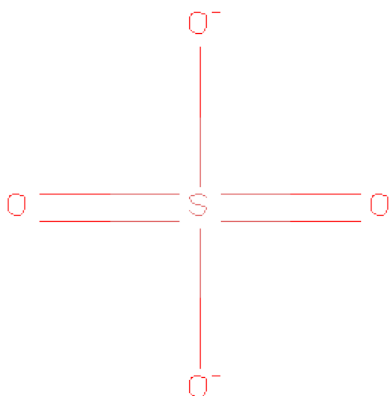
- Molecule 1 is a protein called Genome polyprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	262	2049	1282	376	379	12	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

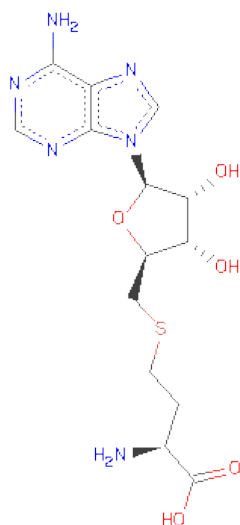
Chain	Residue	Modelled	Actual	Comment	Reference
A	2	GLY	-	CLONING ARTIFACT	UNP P12823
A	3	SER	-	CLONING ARTIFACT	UNP P12823
A	135	ILE	VAL	CONFLICT	UNP P12823
A	139	ARG	LYS	CONFLICT	UNP P12823
A	173	SER	ASN	CONFLICT	UNP P12823
A	193	LYS	ARG	CONFLICT	UNP P12823
A	196	ALA	THR	CONFLICT	UNP P12823
A	201	HIS	TYR	CONFLICT	UNP P12823

- Molecule 2 is SULFATE ION (three-letter code: SO₄) (formula: O₄S).



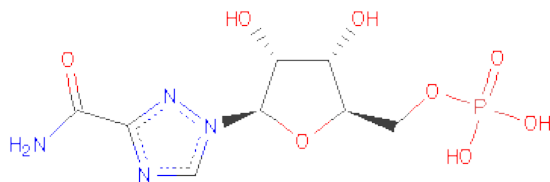
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is S-ADENOSYL-L-HOMOCYSTEINE (three-letter code: SAH) (formula: $C_{14}H_{20}N_6O_5S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			26	14	6	5	1		

- Molecule 4 is RIBAVIRIN MONOPHOSPHATE (three-letter code: RVP) (formula: $C_8H_{13}N_4O_8P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			21	8	4	8	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	66	Total	O	0	0
			66	66		

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- Molecule 1: Genome polypeptide

[illegible]

4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	111.99Å 111.99Å 56.27Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	25.00 – 2.60 28.03 – 2.40	Depositor EDS
% Data completeness (in resolution range)	(Not available) (25.00-2.60) 94.2 (28.03-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.14 (at 2.39Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.218 , 0.243 0.295 , 0.295	Depositor DCC
R_{free} test set	613 reflections (5.18%)	DCC
Wilson B-factor (Å ²)	52.3	Xtriage
Anisotropy	0.467	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 25.8	EDS
Estimated twinning fraction	0.049 for -h,-k,l	Xtriage
L-test for twinning	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 15239 reflections	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	2197	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.99% of the height of the origin peak. No significant pseudotranslation is detected.*

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: SAH, RVP, SO4

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	1/2091 (0.0%)	0.83	6/2823 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	188	PRO	C-N	-9.05	1.13	1.34

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	96	ASN	N-CA-CB	-14.14	85.15	110.60
1	A	188	PRO	O-C-N	-13.29	101.44	122.70
1	A	139	ARG	NE-CZ-NH2	7.36	123.98	120.30
1	A	95	LYS	N-CA-C	6.53	128.63	111.00
1	A	188	PRO	CA-C-N	5.16	128.54	117.20
1	A	187	MET	CG-SD-CE	5.08	108.33	100.20

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	2049	0	2045	98	0
2	A	35	0	0	1	0
3	A	26	0	19	3	0
4	A	21	0	10	0	0
5	A	66	0	0	4	0
All	All	2197	0	2074	98	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 23.

All (98) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:7:GLU:HB2	1:A:188:PRO:HG3	1.16	1.14
1:A:7:GLU:OE1	1:A:188:PRO:HG2	1.54	1.08
1:A:7:GLU:CB	1:A:188:PRO:HG3	1.94	0.96
1:A:83:GLY:H	1:A:104:THR:HG21	1.29	0.94
1:A:7:GLU:HB2	1:A:188:PRO:CG	1.99	0.91
1:A:74:GLU:H	1:A:142:THR:HG21	1.32	0.91
1:A:208:ASN:HD22	1:A:210:LEU:H	1.19	0.90
1:A:83:GLY:H	1:A:104:THR:CG2	1.88	0.86
1:A:234:MET:HG2	5:A:945:HOH:O	1.80	0.82
1:A:72:THR:H	1:A:177:GLN:NE2	1.78	0.81
1:A:83:GLY:N	1:A:104:THR:HG21	2.00	0.76
1:A:98:ARG:HH11	1:A:265:ILE:HG23	1.53	0.73
1:A:265:ILE:HG22	1:A:266:GLY:H	1.52	0.72
1:A:194:MET:HB3	1:A:228:ILE:HD13	1.71	0.71
1:A:177:GLN:HE22	1:A:223:ASN:HD21	1.39	0.69
1:A:72:THR:HB	1:A:177:GLN:HE21	1.56	0.69
1:A:51:ASP:OD1	1:A:260:SER:HB2	1.92	0.69
1:A:31:SER:HB3	1:A:245:MET:HE3	1.75	0.68
1:A:208:ASN:ND2	1:A:210:LEU:H	1.90	0.68
1:A:9:LEU:O	1:A:9:LEU:HD23	1.94	0.67
1:A:245:MET:O	1:A:248:LYS:HE3	1.96	0.66
1:A:74:GLU:H	1:A:142:THR:CG2	2.09	0.63
1:A:72:THR:HB	1:A:177:GLN:NE2	2.15	0.61
1:A:73:PRO:HA	1:A:142:THR:HG21	1.82	0.60
1:A:74:GLU:N	1:A:142:THR:HG21	2.13	0.59
1:A:73:PRO:HG2	1:A:94:LEU:CD1	2.31	0.59
1:A:92:GLY:O	1:A:120:GLY:HA3	2.02	0.59
1:A:73:PRO:HG2	1:A:94:LEU:HD12	1.82	0.59
1:A:133:PHE:CE2	1:A:163:ARG:HB3	2.38	0.59
1:A:246:ARG:O	1:A:247:HIS:HB2	2.03	0.59
1:A:7:GLU:OE1	1:A:188:PRO:CG	2.40	0.59

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:72:THR:N	1:A:177:GLN:NE2	2.52	0.57
1:A:122:ASN:OD1	1:A:262:THR:HB	2.04	0.56
1:A:71:VAL:HG21	1:A:179:CYS:HB2	1.88	0.55
1:A:122:ASN:N	1:A:122:ASN:HD22	2.04	0.55
1:A:58:GLY:HA3	3:A:887:SAH:O	2.07	0.55
1:A:191:ILE:O	1:A:195:GLU:HG3	2.05	0.54
1:A:38:ARG:HD3	2:A:903:SO4:O4	2.06	0.54
1:A:38:ARG:NH2	1:A:54:ALA:O	2.40	0.53
1:A:133:PHE:HE2	1:A:163:ARG:HB3	1.72	0.53
1:A:212:ARG:HG2	1:A:212:ARG:HH11	1.73	0.53
1:A:204:ALA:HB2	1:A:225:SER:HA	1.91	0.53
1:A:72:THR:CB	1:A:177:GLN:HE21	2.22	0.51
1:A:142:THR:HG22	1:A:177:GLN:HB2	1.91	0.50
1:A:164:VAL:O	1:A:168:VAL:HG13	2.11	0.50
1:A:181:LYS:HB2	1:A:219:TYR:CE2	2.48	0.49
1:A:74:GLU:O	1:A:142:THR:HG23	2.13	0.49
1:A:119:TYR:CZ	1:A:263:ARG:HG2	2.48	0.49
1:A:122:ASN:N	1:A:122:ASN:ND2	2.61	0.48
1:A:82:CYS:N	1:A:104:THR:HG23	2.28	0.48
1:A:166:ASN:OD1	1:A:193:LYS:NZ	2.47	0.48
1:A:195:GLU:HA	1:A:228:ILE:HD12	1.96	0.47
1:A:63:ARG:CZ	1:A:256:VAL:HG22	2.44	0.47
1:A:55:VAL:HG22	5:A:953:HOH:O	2.13	0.47
1:A:77:VAL:HG22	1:A:142:THR:OG1	2.14	0.47
1:A:142:THR:HA	1:A:177:GLN:O	2.15	0.46
1:A:246:ARG:NH2	5:A:927:HOH:O	2.49	0.46
1:A:122:ASN:HD22	1:A:123:LEU:H	1.64	0.46
1:A:212:ARG:HG2	1:A:212:ARG:NH1	2.31	0.46
1:A:72:THR:H	1:A:177:GLN:HE21	1.57	0.45
1:A:83:GLY:CA	1:A:104:THR:HG21	2.46	0.45
1:A:122:ASN:H	1:A:122:ASN:ND2	2.15	0.45
1:A:186:TYR:HA	1:A:232:VAL:HG11	1.98	0.45
1:A:246:ARG:O	1:A:247:HIS:CB	2.64	0.45
1:A:265:ILE:HG22	1:A:266:GLY:N	2.29	0.44
1:A:122:ASN:HD22	1:A:123:LEU:N	2.14	0.44
1:A:17:LEU:HD11	1:A:25:PHE:HE1	1.83	0.44
1:A:102:GLY:O	1:A:103:LEU:HD23	2.16	0.44
1:A:264:ASN:C	1:A:265:ILE:HG13	2.37	0.44
1:A:27:ILE:CG2	1:A:248:LYS:HD3	2.48	0.44
1:A:100:VAL:HB	1:A:124:VAL:HG22	1.98	0.44
1:A:31:SER:CB	1:A:245:MET:HE3	2.43	0.44
1:A:200:LYS:HG2	1:A:201:HIS:CE1	2.53	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:14:LYS:HD3	5:A:928:HOH:O	2.18	0.43
1:A:34:GLN:HB3	1:A:210:LEU:HD23	2.00	0.43
1:A:47:ARG:HD2	1:A:49:GLU:OE2	2.19	0.43
1:A:7:GLU:HB2	1:A:188:PRO:CB	2.47	0.43
1:A:136:PRO:HA	1:A:137:PRO:HD3	1.85	0.43
1:A:78:VAL:HG12	1:A:80:LEU:HD13	2.00	0.43
1:A:82:CYS:O	1:A:85:GLY:N	2.52	0.43
1:A:17:LEU:HD11	1:A:25:PHE:CE1	2.54	0.42
1:A:98:ARG:HD2	1:A:265:ILE:CG2	2.49	0.42
1:A:72:THR:CA	1:A:177:GLN:HE21	2.33	0.42
1:A:188:PRO:O	1:A:189:SER:C	2.37	0.42
1:A:72:THR:N	1:A:177:GLN:HE21	2.15	0.42
1:A:16:ARG:O	1:A:20:LEU:HD13	2.19	0.41
1:A:66:VAL:C	1:A:68:ARG:H	2.23	0.41
1:A:146:ASP:OD1	3:A:887:SAH:HB2	2.20	0.41
1:A:104:THR:HG22	3:A:887:SAH:H1'	2.01	0.41
1:A:62:LEU:HD22	1:A:66:VAL:HG23	2.02	0.41
1:A:8:THR:HG21	1:A:186:TYR:O	2.21	0.41
1:A:7:GLU:N	1:A:188:PRO:HG3	2.36	0.41
1:A:68:ARG:HH11	1:A:68:ARG:HG3	1.85	0.41
1:A:153:ASN:HA	1:A:154:PRO:HD2	1.95	0.41
1:A:208:ASN:HD22	1:A:208:ASN:C	2.24	0.41
1:A:132:VAL:HG23	1:A:167:LEU:HD11	2.03	0.41
1:A:227:ASN:ND2	1:A:230:SER:HB2	2.36	0.41
1:A:84:ARG:HA	1:A:113:PRO:HA	2.03	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	260/295 (88%)	240 (92%)	15 (6%)	5 (2%)	12	23

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	174	ASN
1	A	267	ILE
1	A	95	LYS
1	A	265	ILE
1	A	260	SER

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	224/256 (88%)	214 (96%)	10 (4%)	38 67

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	LEU
1	A	62	LEU
1	A	104	THR
1	A	122	ASN
1	A	127	GLN
1	A	167	LEU
1	A	174	ASN
1	A	188	PRO
1	A	208	ASN
1	A	212	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	122	ASN
1	A	177	GLN
1	A	198	GLN
1	A	208	ASN
1	A	227	ASN
1	A	247	HIS

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 ⓘ

9 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	$\# Z > 2$
4	RVP	A	300	-	22,22,22	4.10	14 (63%)	32,33,33	4.77	14 (43%)
3	SAH	A	887	-	28,28,28	0.86	1 (3%)	40,40,40	1.35	2 (5%)
2	SO4	A	901	-	4,4,4	0.31	0	6,6,6	0.08	0
2	SO4	A	902	-	4,4,4	0.25	0	6,6,6	0.14	0
2	SO4	A	903	-	4,4,4	0.32	0	6,6,6	0.20	0
2	SO4	A	904	-	4,4,4	0.20	0	6,6,6	0.09	0
2	SO4	A	905	-	4,4,4	0.35	0	6,6,6	0.15	0
2	SO4	A	906	-	4,4,4	0.32	0	6,6,6	0.13	0
2	SO4	A	907	-	4,4,4	0.28	0	6,6,6	0.10	0

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
4	RVP	A	300	-	-	0/8/30/30	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	SAH	A	887	-	-	0/15/31/31	0/1/3/3
2	SO4	A	901	-	-	0/0/0/0	0/0/0/0
2	SO4	A	902	-	-	0/0/0/0	0/0/0/0
2	SO4	A	903	-	-	0/0/0/0	0/0/0/0
2	SO4	A	904	-	-	0/0/0/0	0/0/0/0
2	SO4	A	905	-	-	0/0/0/0	0/0/0/0
2	SO4	A	906	-	-	0/0/0/0	0/0/0/0
2	SO4	A	907	-	-	0/0/0/0	0/0/0/0

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	300	RVP	C6-N1	10.72	1.57	1.33
4	A	300	RVP	C2'-C1'	7.18	1.63	1.53
4	A	300	RVP	C5-C6	-6.56	1.33	1.49
4	A	300	RVP	C1'-N9	4.55	1.61	1.48
4	A	300	RVP	C8-N9	4.47	1.44	1.34
4	A	300	RVP	P-O1P	4.29	1.70	1.54
4	A	300	RVP	O4'-C1'	4.10	1.47	1.41
4	A	300	RVP	C5'-C4'	3.97	1.64	1.51
4	A	300	RVP	C5-N7	3.33	1.41	1.36
4	A	300	RVP	C5-N4	3.29	1.37	1.33
4	A	300	RVP	O2'-C2'	-3.28	1.35	1.43
4	A	300	RVP	P-O5'	2.62	1.69	1.60
4	A	300	RVP	P-O3P	2.56	1.59	1.51
4	A	300	RVP	C8-N7	2.15	1.39	1.35
3	A	887	SAH	C2-N3	2.07	1.36	1.32

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	300	RVP	N4-C5-N7	-14.20	103.59	115.09
4	A	300	RVP	O6-C6-N1	-10.68	107.16	122.59
4	A	300	RVP	O4'-C1'-N9	9.26	116.28	108.61
4	A	300	RVP	O6-C6-C5	7.92	139.87	119.87
4	A	300	RVP	N7-C8-N9	-7.05	99.90	111.40
4	A	300	RVP	C4'-O4'-C1'	6.32	116.61	109.75
4	A	300	RVP	C8-N7-C5	5.90	115.72	103.79
4	A	300	RVP	O4'-C4'-C5'	5.56	129.21	109.36
3	A	887	SAH	OXT-C-O	-5.05	112.65	124.07
4	A	300	RVP	O4'-C1'-C2'	-4.64	99.66	106.77
4	A	300	RVP	C5-N4-N9	4.51	109.56	102.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	300	RVP	O1P-P-O5'	4.47	118.99	106.65
4	A	300	RVP	O1P-P-O3P	-3.56	98.79	110.44
4	A	300	RVP	C2'-C3'-C4'	2.81	108.25	102.65
4	A	300	RVP	O2P-P-O3P	2.60	118.95	110.44
3	A	887	SAH	OXT-C-CA	-2.39	111.53	116.88

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	262/295 (88%)	0.11	16 (6%) 21 17	20, 55, 71, 94	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	7	GLU	8.6
1	A	268	GLU	6.2
1	A	266	GLY	5.4
1	A	267	ILE	4.0
1	A	139	ARG	3.2
1	A	179	CYS	2.7
1	A	265	ILE	2.5
1	A	109	GLY	2.4
1	A	144	LEU	2.4
1	A	219	TYR	2.3
1	A	146	ASP	2.3
1	A	181	LYS	2.1
1	A	180	VAL	2.1
1	A	98	ARG	2.0
1	A	147	ILE	2.0
1	A	145	CYS	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	SO4	A	907	5/5	0.47	12.14	0,0,0,0	0
2	SO4	A	906	5/5	0.47	11.98	0,0,0,0	0
2	SO4	A	905	5/5	0.46	10.92	0,0,0,0	0
2	SO4	A	902	5/5	0.40	4.23	0,0,0,99	0
4	RVP	A	300	21/21	0.26	0.92	71,74,84,84	0
2	SO4	A	904	5/5	0.22	0.52	92,92,92,93	0
2	SO4	A	903	5/5	0.15	0.28	58,60,61,62	0
3	SAH	A	887	26/26	0.19	0.24	56,63,65,67	0
2	SO4	A	901	5/5	0.15	-1.69	91,91,91,92	0

6.5 Other polymers ⓘ

There are no such residues in this entry.