



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 12:25 PM GMT

PDB ID : 3SWC
Title : GLP (G9a-like protein) SET domain in complex with Dnmt3aK44me2 peptide
Authors : Chang, Y.; Horton, J.R.; Zhang, X.; Cheng, X.
Deposited on : 2011-07-13
Resolution : 2.33 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>

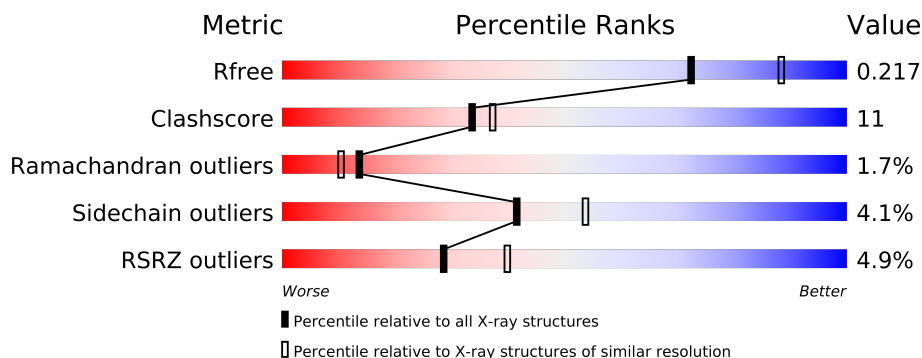
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.33 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	4049 (2.38-2.30)
Clashscore	79885	1094 (2.36-2.32)
Ramachandran outliers	78287	1080 (2.36-2.32)
Sidechain outliers	78261	1081 (2.36-2.32)
RSRZ outliers	66119	4050 (2.38-2.30)

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 ≥ 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	285	
1	B	285	
2	P	12	
2	Q	12	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4528 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 Histone-lysine N-methyltransferase EHMT1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	261	Total	C	N	O	S	0	0	0
			2081	1297	373	386	25			
1	B	261	Total	C	N	O	S	0	1	0
			2096	1307	380	384	25			

- Molecule 2 is a protein called DNA (cytosine-5)-methyltransferase3A.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	P	9	Total	C	N	O	0	0	0
			61	38	13	10			
2	Q	8	Total	C	N	O	0	0	0
			60	37	15	8			

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	4	Total	Zn	0	0
			4	4		
3	A	4	Total	Zn	0	0
			4	4		

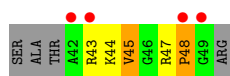
- Molecule 4 is S-ADENOSYL-L-HOMOCYSTEINE (three-letter code: SAH) (formula: C₁₄H₂₀N₆O₅S).



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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	78	Total	O	0	0
			78	78		
5	P	2	Total	O	0	0
			2	2		
5	B	87	Total	O	0	0
			87	87		
5	Q	3	Total	O	0	0
			3	3		



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	80.09Å 90.22Å 95.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.78 – 2.33 47.70 – 2.33	Depositor EDS
% Data completeness (in resolution range)	95.0 (40.78-2.33) 95.0 (47.70-2.33)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.50 (at 2.34Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.289)	Depositor
R, R_{free}	0.187 , 0.227 0.177 , 0.217	Depositor DCC
R_{free} test set	1966 reflections (6.87%)	DCC
Wilson B-factor (Å ²)	38.3	Xtriage
Anisotropy	0.308	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 33.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 29265 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4528	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.95% 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, ZN, MLY

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	0.42	0/2132	0.55	1/2888 (0.0%)
1	B	0.42	0/2150	0.56	0/2909
2	P	0.60	0/50	1.04	0/67
2	Q	0.70	0/49	0.84	0/64
All	All	0.42	0/4381	0.57	1/5928 (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	A	1150	ASP	CB-CG-OD2	5.21	122.99	118.30

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	2081	0	1942	35	0
1	B	2096	0	1975	43	0
2	P	61	0	63	16	0
2	Q	60	0	67	18	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	4	0	0	0	0
3	B	4	0	0	0	0
4	A	26	0	19	0	0
4	B	26	0	19	1	0
5	A	78	0	0	0	0
5	B	87	0	0	2	0
5	P	2	0	0	0	0
5	Q	3	0	0	0	0
All	All	4528	0	4085	93	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 11.

All (93) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:1054:ASP:HB2	1:A:1055:LYS:HE2	1.45	0.96
1:B:1219:LYS:HG2	2:Q:45:VAL:HG11	1.49	0.95
1:A:1058:ARG:HH11	1:A:1058:ARG:HG2	1.44	0.81
2:Q:43:ARG:HG2	2:Q:44:MLY:N	1.98	0.79
2:Q:43:ARG:CG	2:Q:44:MLY:N	2.47	0.77
1:B:1218:ILE:HB	2:Q:45:VAL:HG13	1.68	0.74
1:A:979:VAL:HG23	1:A:1003:CYS:HB2	1.70	0.74
2:P:45:VAL:HG22	2:P:45:VAL:O	1.89	0.71
1:A:1016:THR:O	1:A:1017:SER:HB2	1.90	0.70
1:B:1212:GLY:HA3	5:B:168:HOH:O	1.91	0.68
1:B:1147:ASP:HA	1:B:1153:VAL:HG22	1.78	0.66
2:P:45:VAL:CG2	2:P:45:VAL:O	2.44	0.66
2:Q:47:ARG:HB3	2:Q:48:PRO:HD2	1.79	0.64
1:B:1131:ASP:OD2	2:Q:44:MLY:HH11	1.99	0.63
1:A:1058:ARG:NH1	1:A:1058:ARG:HG2	2.11	0.63
1:B:1215:PHE:CZ	1:B:1219:LYS:HG3	2.34	0.63
1:A:1131:ASP:OD2	2:P:44:MLY:HH11	2.01	0.61
1:A:1048:SER:O	1:A:1049:MET:HB2	2.02	0.59
1:A:1152:GLU:HA	1:A:1152:GLU:OE2	2.02	0.58
1:B:1215:PHE:HA	2:Q:45:VAL:HG12	1.87	0.57
1:B:990:ILE:HD12	1:B:991:PRO:O	2.05	0.56
1:B:1136:VAL:HG23	1:B:1136:VAL:O	2.06	0.56
1:B:1145:ASP:OD1	2:Q:44:MLY:HE3	2.05	0.55
1:B:1059:LEU:HD13	1:B:1063:PHE:CD2	2.41	0.55
1:B:1048:SER:O	1:B:1049:MET:HB2	2.06	0.55
2:Q:43:ARG:NH1	2:Q:43:ARG:HB2	2.22	0.54
2:P:41:THR:O	2:P:42:ALA:HB3	2.07	0.54

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:1149:LYS:C	1:B:1151:GLY:H	2.13	0.52
2:Q:47:ARG:HB3	2:Q:48:PRO:CD	2.39	0.52
1:B:990:ILE:HD13	1:B:1096:ARG:HA	1.91	0.52
1:B:1147:ASP:H	1:B:1153:VAL:HG13	1.75	0.51
1:A:989:ARG:NH1	1:A:1094:ARG:HD3	2.25	0.51
1:A:1016:THR:O	1:A:1017:SER:CB	2.56	0.51
1:A:1082:ARG:O	1:A:1087:ARG:HD2	2.09	0.51
1:A:1016:THR:HG22	1:B:1025:ILE:HG22	1.93	0.51
1:B:1134:ALA:HB3	2:Q:44:MLY:HH13	1.93	0.51
5:B:168:HOH:O	2:Q:48:PRO:HA	2.09	0.51
1:A:978:ILE:HG22	1:A:978:ILE:O	2.11	0.50
1:A:1055:LYS:CD	1:A:1055:LYS:H	2.24	0.50
1:B:1170:HIS:HB2	1:B:1211:TYR:CD1	2.46	0.50
1:B:1048:SER:O	1:B:1049:MET:CB	2.60	0.49
1:B:1135:ASP:OD1	2:Q:44:MLY:HH12	2.12	0.49
1:B:979:VAL:HG23	1:B:1003:CYS:HB2	1.95	0.49
1:A:1054:ASP:HB2	1:A:1055:LYS:CE	2.32	0.49
1:A:1048:SER:O	1:A:1049:MET:CB	2.60	0.49
1:B:1211:TYR:O	2:Q:48:PRO:HD3	2.14	0.48
1:B:1128:LEU:HD21	1:B:1190:PHE:CD1	2.48	0.48
1:B:1139:GLU:HG2	1:B:1159:ARG:HD3	1.96	0.48
1:A:1147:ASP:N	1:A:1180:ARG:NH2	2.61	0.48
1:B:1132:SER:O	1:B:1136:VAL:HG13	2.14	0.48
1:A:979:VAL:CG2	1:A:1003:CYS:HB2	2.41	0.47
1:A:1215:PHE:CZ	1:A:1219:LYS:HG3	2.49	0.47
1:A:1215:PHE:CD1	2:P:45:VAL:HG23	2.50	0.47
1:B:1148:ASN:C	1:B:1150:ASP:H	2.18	0.47
1:A:995:VAL:O	1:A:1100:TYR:HA	2.15	0.47
2:P:41:THR:O	2:P:42:ALA:CB	2.63	0.46
2:P:44:MLY:HE3	2:P:44:MLY:HB2	1.69	0.46
1:B:1085:ARG:O	1:B:1085:ARG:HG2	2.16	0.46
1:A:1147:ASP:C	1:A:1180:ARG:HH22	2.19	0.46
2:P:44:MLY:HD2	2:P:44:MLY:HH13	1.52	0.46
1:A:1043:MET:HA	1:A:1043:MET:HE2	1.98	0.45
1:A:1016:THR:OG1	1:A:1016:THR:O	2.33	0.45
1:A:1081:TRP:O	1:A:1084:CYS:HB3	2.17	0.45
1:B:1152:GLU:HG2	1:B:1154:TYR:CE2	2.52	0.44
1:B:1134:ALA:CB	2:Q:44:MLY:HE2	2.48	0.44
1:A:1145:ASP:OD1	2:P:44:MLY:HE2	2.18	0.43
1:A:1004:PRO:HD3	1:A:1107:TRP:CZ2	2.53	0.43
1:B:1139:GLU:CG	1:B:1159:ARG:HD3	2.49	0.43
1:B:1148:ASN:HD22	1:B:1150:ASP:CB	2.31	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:1199[A]:ARG:NH2	1:B:1205:GLU:OE1	2.46	0.43
2:P:41:THR:OG1	2:P:42:ALA:N	2.51	0.43
1:B:1170:HIS:HB2	1:B:1211:TYR:CG	2.54	0.42
1:B:998:VAL:HG22	1:B:1101:ARG:HG2	2.00	0.42
2:Q:44:MLY:HH12	2:Q:44:MLY:HD2	1.41	0.42
1:A:1232:CYS:SG	1:A:1235:SER:HB3	2.60	0.42
1:A:1211:TYR:O	2:P:48:PRO:HA	2.18	0.42
2:P:41:THR:HG23	2:P:42:ALA:H	1.85	0.42
1:B:1055:LYS:HG3	1:B:1055:LYS:H	1.68	0.42
1:B:1148:ASN:C	1:B:1150:ASP:N	2.73	0.42
1:B:1232:CYS:SG	1:B:1235:SER:HB3	2.60	0.42
1:B:1115:ILE:HB	1:B:1201:ILE:HB	2.01	0.42
1:A:1134:ALA:HB1	2:P:44:MLY:HD3	2.02	0.41
1:A:1054:ASP:OD1	1:A:1056:ASP:N	2.53	0.41
1:B:1218:ILE:CB	2:Q:45:VAL:HG13	2.46	0.41
1:B:1134:ALA:HB1	2:Q:44:MLY:HE2	2.02	0.41
1:A:1145:ASP:OD1	2:P:44:MLY:HH22	2.20	0.41
1:B:1169:ASN:HB3	4:B:201:SAH:N7	2.35	0.41
1:A:1219:LYS:HE2	2:P:45:VAL:HG21	2.03	0.41
1:A:1147:ASP:O	1:A:1180:ARG:NH2	2.53	0.41
1:B:1105:MET:HE3	1:B:1105:MET:HB3	1.86	0.41
1:A:1058:ARG:NH1	1:A:1119:THR:OG1	2.54	0.40
2:P:41:THR:HG23	2:P:42:ALA:N	2.35	0.40
1:B:1158:ALA:HA	1:B:1161:TYR:O	2.22	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	259/285 (91%)	235 (91%)	21 (8%)	3 (1%)	19	18
1	B	260/285 (91%)	236 (91%)	23 (9%)	1 (0%)	43	50
2	P	6/12 (50%)	2 (33%)	0	4 (67%)	0	0

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	Q	5/12 (42%)	1 (20%)	3 (60%)	1 (20%)	0	0
All	All	530/594 (89%)	474 (89%)	47 (9%)	9 (2%)	14	10

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1017	SER
1	B	1147	ASP
2	Q	48	PRO
2	P	42	ALA
1	A	1049	MET
2	P	43	ARG
2	P	48	PRO
1	A	978	ILE
2	P	45	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	229/257 (89%)	222 (97%)	7 (3%)	52	66
1	B	231/257 (90%)	221 (96%)	10 (4%)	40	51
2	P	4/7 (57%)	3 (75%)	1 (25%)	1	0
2	Q	4/7 (57%)	3 (75%)	1 (25%)	1	0
All	All	468/528 (89%)	449 (96%)	19 (4%)	41	54

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1039	SER
1	A	1055	LYS
1	A	1062	GLU
1	A	1084	CYS
1	A	1110	ARG
1	A	1150	ASP

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Mol	Chain	Res	Type
1	A	1152	GLU
2	P	45	VAL
1	B	998	VAL
1	B	1034	ILE
1	B	1055	LYS
1	B	1084	CYS
1	B	1093	LEU
1	B	1105	MET
1	B	1110	ARG
1	B	1131	ASP
1	B	1136	VAL
1	B	1146	LEU
2	Q	45	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1041	ASN
1	B	1148	ASN

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	$\# Z > 2$
2	MLY	P	44	2	10,10,11	5.70	2 (20%)	9,11,13	1.19	1 (11%)
2	MLY	Q	44	2	10,10,11	5.72	2 (20%)	9,11,13	1.50	1 (11%)

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	MLY	P	44	2	-	0/7/9/11	0/0/0/0
2	MLY	Q	44	2	-	0/7/9/11	0/0/0/0

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	Q	44	MLY	O-C	17.78	1.23	1.11
2	P	44	MLY	O-C	17.76	1.23	1.11
2	Q	44	MLY	CA-C	2.71	1.53	1.48
2	P	44	MLY	CA-C	2.54	1.53	1.48

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	Q	44	MLY	C-CA-N	-3.25	110.58	113.83
2	P	44	MLY	C-CA-N	-2.37	111.46	113.83

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

Of 10 ligands modelled in this entry, 8 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
4	SAH	A	202	-	28,28,28	1.03	3 (10%)	40,40,40	2.22	8 (20%)
4	SAH	B	201	-	28,28,28	1.03	2 (7%)	40,40,40	2.19	9 (22%)

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	SAH	A	202	-	-	0/15/31/31	0/1/3/3
4	SAH	B	201	-	-	0/15/31/31	0/1/3/3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	201	SAH	C2-N3	3.15	1.38	1.32
4	A	202	SAH	C2-N3	3.04	1.38	1.32
4	B	201	SAH	C2-N1	2.43	1.38	1.33
4	A	202	SAH	C2'-C1'	-2.33	1.50	1.53
4	A	202	SAH	C2-N1	2.06	1.38	1.33

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	202	SAH	N3-C2-N1	-11.23	119.32	128.71
4	B	201	SAH	N3-C2-N1	-10.00	120.35	128.71
4	A	202	SAH	N3-C4-N9	3.85	132.38	125.43
4	B	201	SAH	N3-C4-N9	3.53	131.81	125.43
4	A	202	SAH	C5'-SD-CG	-3.23	92.79	102.42
4	B	201	SAH	C2'-C1'-N9	-3.20	105.05	113.27
4	B	201	SAH	C4-C5-N7	-3.03	106.93	109.52
4	B	201	SAH	C5'-SD-CG	-3.02	93.40	102.42
4	A	202	SAH	O2'-C2'-C1'	-2.53	103.57	111.23
4	B	201	SAH	C5-C4-N3	-2.52	120.21	125.70
4	B	201	SAH	OXT-C-O	-2.52	118.38	124.07
4	B	201	SAH	O4'-C1'-N9	2.49	110.76	108.44
4	A	202	SAH	C2-N3-C4	2.46	121.01	114.01
4	A	202	SAH	C5-C4-N3	-2.41	120.46	125.70
4	B	201	SAH	C2-N3-C4	2.16	120.17	114.01
4	A	202	SAH	N7-C8-N9	-2.13	108.33	114.36
4	A	202	SAH	OXT-C-CA	2.02	121.42	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	261/285 (91%)	0.19	8 (3%) 47 58	26, 39, 57, 91	0
1	B	261/285 (91%)	0.25	11 (4%) 35 46	24, 36, 56, 87	0
2	P	9/12 (75%)	3.21	4 (44%) 1 0	36, 53, 88, 93	0
2	Q	8/12 (66%)	3.13	4 (50%) 0 0	35, 52, 73, 75	0
All	All	539/594 (90%)	0.31	27 (5%) 28 39	24, 37, 59, 93	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	P	48	PRO	10.3
1	A	975	VAL	7.9
2	Q	42	ALA	7.2
2	P	49	GLY	6.8
2	Q	43	ARG	5.9
2	P	42	ALA	5.5
1	B	1148	ASN	5.5
1	A	977	ARG	5.4
2	Q	49	GLY	5.1
1	B	1151	GLY	5.0
1	B	1222	LEU	4.4
1	B	1150	ASP	4.3
1	A	976	GLU	4.3
2	Q	48	PRO	4.3
1	B	1226	ARG	3.3
2	P	41	THR	3.2
1	B	1136	VAL	3.0
1	B	1214	ARG	2.5
1	A	1148	ASN	2.4
1	B	1220	GLY	2.4
1	A	1150	ASP	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	979	VAL	2.2
1	B	1149	LYS	2.1
1	B	1213	GLU	2.1
1	B	1235	SER	2.1
1	A	1152	GLU	2.0
1	A	1049	MET	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, 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	MLY	P	44	11/12	0.26	1.17	46,52,59,61	0
2	MLY	Q	44	11/12	0.27	0.40	46,51,58,67	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, 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
3	ZN	B	306	1/1	0.17	1.35	31,31,31,31	0
3	ZN	B	307	1/1	0.15	0.64	33,33,33,33	0
3	ZN	A	302	1/1	0.15	0.40	30,30,30,30	0
3	ZN	A	301	1/1	0.14	0.11	29,29,29,29	0
3	ZN	A	303	1/1	0.15	0.05	29,29,29,29	0
4	SAH	B	201	26/26	0.14	-0.18	30,39,45,46	0
3	ZN	B	305	1/1	0.14	-0.20	31,31,31,31	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	ZN	A	304	1/1	0.10	-0.46	39,39,39,39	0
4	SAH	A	202	26/26	0.11	-0.49	31,37,41,44	0
3	ZN	B	308	1/1	0.09	-1.16	40,40,40,40	0

6.5 Other polymers ⓘ

There are no such residues in this entry.