



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 03:22 AM GMT

PDB ID : 3MO5
Title : Human G9a-like (GLP, also known as EHMT1) in complex with inhibitor E72
Authors : Chang, Y.; Horton, J.R.; Cheng, X.
Deposited on : 2010-04-22
Resolution : 2.14 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>

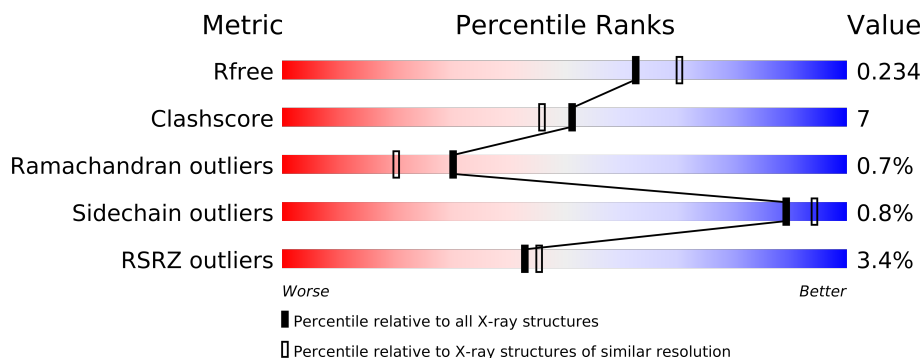
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.14 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	1116 (2.16-2.12)
Clashscore	79885	1302 (2.16-2.12)
Ramachandran outliers	78287	1281 (2.16-2.12)
Sidechain outliers	78261	1281 (2.16-2.12)
RSRZ outliers	66119	1116 (2.16-2.12)

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	
1	C	285	
1	D	285	

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

Mol	Type	Chain	Res	Geometry	Electron density
4	SAH	D	104	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 9067 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.

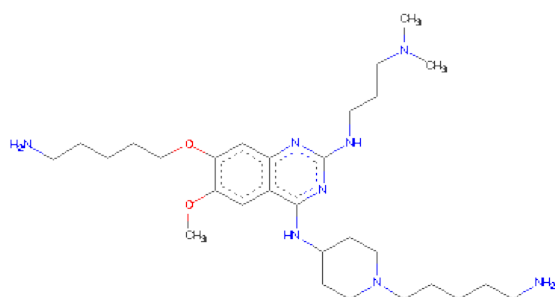
- Molecule 1 is a protein called Histone-lysine N-methyltransferase, H3 lysine-9 specific 5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	255	Total	C	N	O	S	0	0	0
			2022	1262	363	373	24			
1	B	258	Total	C	N	O	S	0	0	0
			2050	1282	366	377	25			
1	C	258	Total	C	N	O	S	0	0	0
			2053	1282	368	378	25			
1	D	256	Total	C	N	O	S	0	0	0
			2019	1263	359	372	25			

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

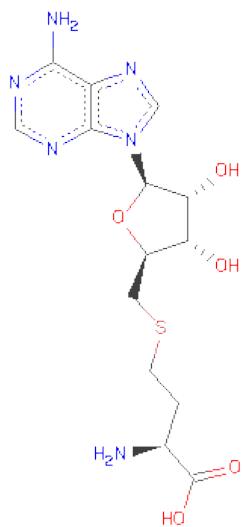
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	4	Total	Zn	0	0
			4	4		
2	A	4	Total	Zn	0	0
			4	4		
2	D	4	Total	Zn	0	0
			4	4		
2	C	4	Total	Zn	0	0
			4	4		

- Molecule 3 is 7-[(5-AMINOPENTYL)OXY]-N 4 -[1-(5-AMINOPENTYL)PIPERIDIN-4-Y L]-N 2 -[3-(DIMETHYLAMINO)PROPYL]-6-METHOXYQUINAZOLINE-2,4-DIAMINE (three-letter code: E72) (formula: C₂₉H₅₂N₈O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			39	29	8	2		
3	B	1	Total	C	N	O	0	0
			39	29	8	2		
3	C	1	Total	C	N	O	0	0
			39	29	8	2		
3	D	1	Total	C	N	O	0	0
			39	29	8	2		

- Molecule 4 is S-ADENOSYL-L-HOMOCYSTEINE (three-letter code: SAH) (formula: $C_{14}H_{20}N_6O_5S$).



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

- Molecule 5 is water.

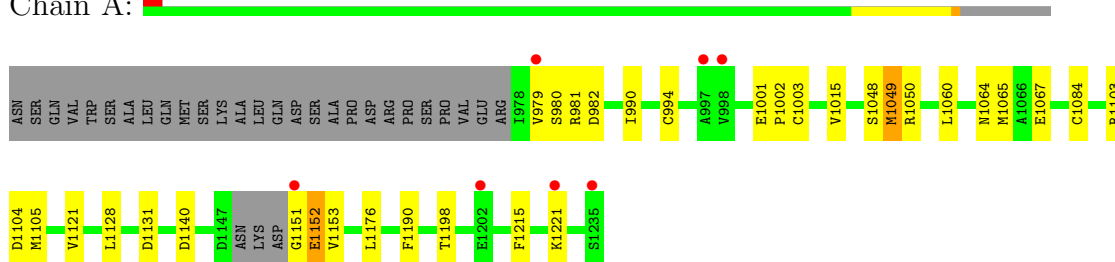
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	193	Total	O	0	0
			193	193		
5	B	201	Total	O	0	0
			201	201		
5	C	182	Total	O	0	0
			182	182		
5	D	123	Total	O	0	0
			123	123		

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.

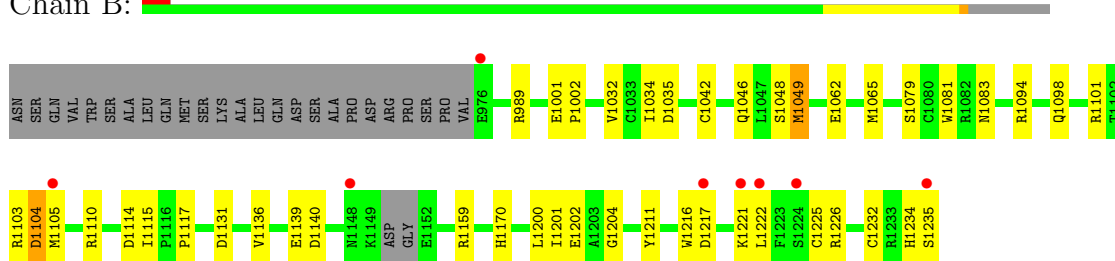
- Molecule 1: Histone-lysine N-methyltransferase, H3 lysine-9 specific 5

Chain A:



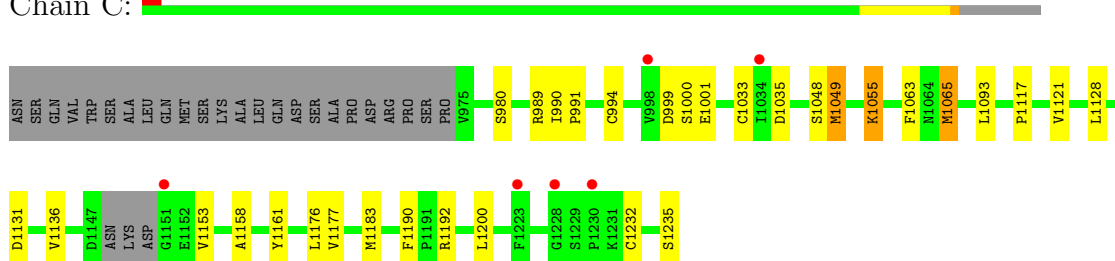
- Molecule 1: Histone-lysine N-methyltransferase, H3 lysine-9 specific 5

Chain B:



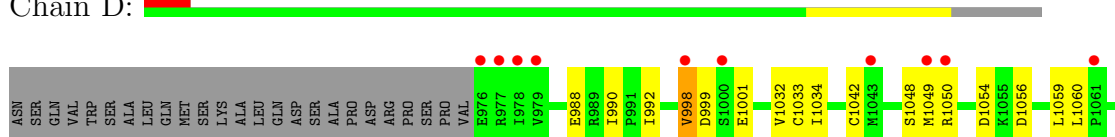
- Molecule 1: Histone-lysine N-methyltransferase, H3 lysine-9 specific 5

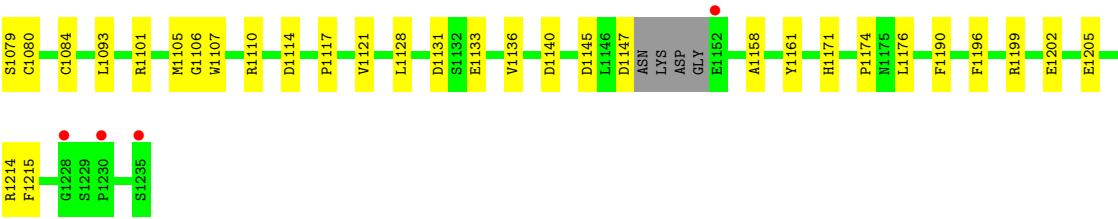
Chain C:



- Molecule 1: Histone-lysine N-methyltransferase, H3 lysine-9 specific 5

Chain D:





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	59.90Å 162.90Å 69.10Å 90.00° 99.00° 90.00°	Depositor
Resolution (Å)	34.14 – 2.14 34.12 – 2.13	Depositor EDS
% Data completeness (in resolution range)	93.3 (34.14-2.14) 92.2 (34.12-2.13)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.10 (at 2.12Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.196 , 0.233 0.205 , 0.234	Depositor DCC
R_{free} test set	3427 reflections (5.11%)	DCC
Wilson B-factor (Å ²)	28.6	Xtriage
Anisotropy	0.309	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 45.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 72203 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9067	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.74% 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: E72, ZN, SAH

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	0.33	0/2072	0.57	0/2807
1	B	0.33	0/2100	0.57	0/2844
1	C	0.33	0/2103	0.57	0/2848
1	D	0.31	0/2069	0.54	0/2807
All	All	0.33	0/8344	0.57	0/11306

There are no bond length outliers.

There are no bond angle outliers.

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	2022	0	1879	30	0
1	B	2050	0	1916	31	0
1	C	2053	0	1916	24	0
1	D	2019	0	1868	35	0
2	A	4	0	0	0	0
2	B	4	0	0	0	0
2	C	4	0	0	0	0
2	D	4	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	39	0	52	3	0
3	B	39	0	52	4	0
3	C	39	0	52	1	0
3	D	39	0	52	3	0
4	C	26	0	18	0	0
4	D	26	0	20	2	0
5	A	193	0	0	0	0
5	B	201	0	0	0	0
5	C	182	0	0	1	0
5	D	123	0	0	1	0
All	All	9067	0	7825	118	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 7.

All (118) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:1050:ARG:HH22	1:D:1060:LEU:HD12	1.30	0.95
1:B:989:ARG:NH1	1:B:1094:ARG:HE	1.71	0.87
1:D:1101:ARG:NH1	1:D:1106:GLY:HA2	1.93	0.83
1:A:1131:ASP:HB2	1:A:1153:VAL:HG23	1.62	0.82
1:D:1080:CYS:SG	5:D:548:HOH:O	2.40	0.78
1:A:1151:GLY:O	1:A:1152:GLU:HG2	1.85	0.76
1:A:1050:ARG:HH12	1:A:1060:LEU:HD12	1.50	0.74
1:D:1050:ARG:NH2	1:D:1060:LEU:HD12	2.05	0.70
1:B:1083:ASN:O	1:C:1055:LYS:HE3	1.94	0.68
1:A:990:ILE:O	1:A:990:ILE:HD12	1.94	0.67
1:A:1050:ARG:NH1	1:A:1060:LEU:HD12	2.09	0.66
1:D:1117:PRO:HB3	1:D:1199:ARG:HA	1.78	0.65
1:B:1034:ILE:HD12	1:B:1035:ASP:N	2.12	0.65
1:D:1101:ARG:HH12	1:D:1106:GLY:HA2	1.62	0.65
1:D:1131:ASP:OD1	3:D:4:E72:HAZA	1.96	0.64
1:D:1048:SER:O	1:D:1049:MET:HB2	1.98	0.63
1:B:1048:SER:O	1:B:1049:MET:HB2	2.00	0.60
1:B:1062:GLU:OE1	1:B:1062:GLU:N	2.31	0.60
1:A:1048:SER:O	1:A:1049:MET:HB2	2.02	0.59
1:C:1065:MET:CE	1:C:1177:VAL:HG11	2.33	0.58
1:B:1139:GLU:HG2	1:B:1159:ARG:HD3	1.84	0.58
1:B:1001:GLU:HG3	1:B:1101:ARG:NH2	2.20	0.57
1:A:1221:LYS:NZ	1:A:1221:LYS:HB2	2.19	0.57
1:D:1133:GLU:O	1:D:1136:VAL:HG22	2.04	0.56
1:C:1128:LEU:HD21	1:C:1190:PHE:CD1	2.41	0.56

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:1131:ASP:OD2	3:B:2:E72:HAZA	2.05	0.56
1:C:1121:VAL:HG21	1:C:1176:LEU:HD21	1.88	0.55
1:D:1101:ARG:HH11	1:D:1106:GLY:HA2	1.73	0.54
1:D:1110:ARG:HB2	1:D:1205:GLU:O	2.07	0.54
1:C:1232:CYS:SG	1:C:1235:SER:HB3	2.47	0.54
1:A:1152:GLU:HG3	1:A:1152:GLU:O	2.08	0.54
1:D:990:ILE:HD11	1:D:1093:LEU:HD13	1.90	0.54
1:B:1140:ASP:O	3:B:2:E72:HAMA	2.07	0.54
1:D:988:GLU:HB3	1:D:1093:LEU:HD12	1.90	0.53
1:B:1232:CYS:SG	1:B:1235:SER:HB3	2.49	0.53
1:B:1042:CYS:O	1:B:1046:GLN:HG3	2.09	0.53
1:C:1048:SER:O	1:C:1049:MET:HB2	2.08	0.53
1:C:989:ARG:HD3	1:C:1093:LEU:O	2.08	0.52
1:B:1217:ASP:O	1:B:1221:LYS:HE2	2.08	0.52
1:D:990:ILE:O	1:D:990:ILE:HD12	2.10	0.52
1:D:1215:PHE:HB2	3:D:4:E72:HBC	1.92	0.52
1:A:979:VAL:HG23	1:A:1003:CYS:HB2	1.91	0.52
1:C:1065:MET:HE2	1:C:1177:VAL:HG11	1.91	0.51
1:B:1103:ARG:C	1:B:1105:MET:H	2.13	0.51
1:D:992:ILE:HD11	1:D:1093:LEU:HD11	1.92	0.51
1:A:1140:ASP:OD2	3:A:1236:E72:HAC	2.11	0.51
1:B:1104:ASP:OD2	1:B:1226:ARG:NE	2.41	0.51
1:C:1131:ASP:HB2	1:C:1153:VAL:HG23	1.92	0.51
1:C:980:SER:HB3	1:C:994:CYS:HB3	1.93	0.51
1:B:989:ARG:HH12	1:B:1094:ARG:HH21	1.57	0.50
1:A:1048:SER:O	1:A:1049:MET:CB	2.60	0.49
1:D:999:ASP:OD1	1:D:1001:GLU:HB2	2.12	0.49
1:D:1121:VAL:HG21	1:D:1176:LEU:HD21	1.94	0.49
1:C:1033:CYS:HB3	1:C:1035:ASP:OD1	2.12	0.49
1:A:1221:LYS:HZ2	1:A:1221:LYS:HB2	1.77	0.49
1:A:1065:MET:CE	1:A:1198:THR:HA	2.43	0.49
1:C:1065:MET:HE2	1:C:1177:VAL:CG1	2.42	0.49
1:C:990:ILE:HG23	1:C:991:PRO:HD2	1.93	0.49
1:A:1103:ARG:C	1:A:1105:MET:H	2.16	0.49
1:B:1221:LYS:C	1:B:1222:LEU:HD12	2.33	0.49
1:A:1152:GLU:CG	1:A:1152:GLU:O	2.60	0.48
1:B:1131:ASP:OD2	3:B:2:E72:HAVA	2.14	0.48
1:D:1033:CYS:SG	1:D:1042:CYS:HA	2.53	0.48
1:D:1128:LEU:HD21	1:D:1190:PHE:CD1	2.48	0.48
1:B:1170:HIS:CE1	1:B:1216:TRP:HE1	2.32	0.48
1:D:1171:HIS:O	1:D:1174:PRO:HD3	2.14	0.47
1:D:1048:SER:C	1:D:1050:ARG:H	2.18	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:1032:VAL:HA	1:B:1079:SER:OG	2.15	0.47
1:B:1225:CYS:HB2	1:B:1234:HIS:HB2	1.97	0.46
1:D:1054:ASP:OD1	1:D:1056:ASP:N	2.49	0.46
1:A:990:ILE:HD12	1:A:990:ILE:C	2.34	0.46
1:C:1063:PHE:CD2	1:C:1065:MET:HE1	2.50	0.46
1:A:982:ASP:HB3	1:B:1081:TRP:CH2	2.52	0.45
1:B:1110:ARG:HD3	1:B:1204:GLY:HA2	1.99	0.45
1:A:1140:ASP:O	3:A:1236:E72:HAMA	2.17	0.44
1:B:1114:ASP:OD1	1:B:1202:GLU:HA	2.18	0.43
1:A:981:ARG:HD2	1:C:1117:PRO:HG2	2.00	0.43
1:D:1117:PRO:HB3	1:D:1199:ARG:CA	2.46	0.43
1:A:1064:ASN:ND2	1:A:1067:GLU:HB3	2.32	0.43
1:B:1034:ILE:C	1:B:1034:ILE:HD12	2.39	0.43
1:D:1158:ALA:HA	1:D:1161:TYR:O	2.19	0.43
1:C:1048:SER:O	1:C:1049:MET:CB	2.66	0.43
1:D:1114:ASP:OD1	1:D:1202:GLU:HA	2.18	0.43
1:C:1065:MET:HG3	1:C:1177:VAL:HG11	2.01	0.43
1:A:1001:GLU:HA	1:A:1002:PRO:HD3	1.91	0.43
1:D:1001:GLU:HG3	1:D:1101:ARG:HH21	1.83	0.42
1:D:1140:ASP:O	3:D:4:E72:HAMA	2.19	0.42
1:B:1170:HIS:HB2	1:B:1211:TYR:CG	2.54	0.42
1:A:1221:LYS:CB	1:A:1221:LYS:NZ	2.83	0.42
1:A:1064:ASN:HD21	1:A:1067:GLU:HB3	1.84	0.42
1:C:1001:GLU:OE1	1:C:1001:GLU:HA	2.20	0.42
1:D:1034:ILE:HD12	1:D:1034:ILE:C	2.40	0.42
1:A:1128:LEU:HD21	1:A:1190:PHE:CD1	2.55	0.42
1:C:999:ASP:HB2	1:C:1000:SER:H	1.71	0.42
1:C:1183:MET:HE3	1:C:1192:ARG:HD3	2.02	0.42
1:B:1115:ILE:HB	1:B:1201:ILE:HB	2.02	0.42
1:D:1049:MET:O	1:D:1050:ARG:HB2	2.20	0.41
1:C:1093:LEU:HD12	5:C:33:HOH:O	2.20	0.41
1:D:998:VAL:HG12	1:D:999:ASP:N	2.35	0.41
4:D:104:SAH:O4'	4:D:104:SAH:HG2	2.21	0.41
1:D:1145:ASP:OD2	1:D:1214:ARG:HD2	2.20	0.41
1:B:1117:PRO:HG3	1:B:1200:LEU:N	2.34	0.41
1:D:1032:VAL:HA	1:D:1079:SER:OG	2.20	0.41
1:A:980:SER:HB3	1:A:994:CYS:HB3	2.02	0.41
1:A:979:VAL:CG2	1:A:1003:CYS:HB2	2.49	0.41
1:A:1015:VAL:HG12	1:A:1128:LEU:HB3	2.03	0.41
1:C:1117:PRO:HG3	1:C:1200:LEU:N	2.36	0.41
1:C:1136:VAL:HA	3:C:1236:E72:HBJ	2.03	0.41
1:A:1121:VAL:HG21	1:A:1176:LEU:HD21	2.03	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:1158:ALA:HA	1:C:1161:TYR:O	2.20	0.41
1:B:1048:SER:O	1:B:1049:MET:CB	2.65	0.41
1:D:1107:TRP:O	4:D:104:SAH:N	2.54	0.41
1:D:1059:LEU:HD11	1:D:1196:PHE:CG	2.55	0.41
1:B:1001:GLU:HA	1:B:1002:PRO:HD3	1.80	0.40
1:B:1136:VAL:HA	3:B:2:E72:HBJ	2.03	0.40
1:A:1065:MET:HE3	1:A:1198:THR:HA	2.03	0.40
1:A:1215:PHE:HB2	3:A:1236:E72:HBC	2.04	0.40
1:B:1098:GLN:OE1	1:B:1110:ARG:NH1	2.52	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	251/285 (88%)	235 (94%)	13 (5%)	3 (1%)	19	9
1	B	254/285 (89%)	239 (94%)	14 (6%)	1 (0%)	43	39
1	C	254/285 (89%)	237 (93%)	16 (6%)	1 (0%)	43	39
1	D	252/285 (88%)	231 (92%)	19 (8%)	2 (1%)	27	16
All	All	1011/1140 (89%)	942 (93%)	62 (6%)	7 (1%)	30	20

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	1105	MET
1	A	1049	MET
1	C	1049	MET
1	A	1104	ASP
1	A	1152	GLU
1	B	1049	MET
1	D	998	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	221/257 (86%)	220 (100%)	1 (0%)	94	97
1	B	225/257 (88%)	223 (99%)	2 (1%)	87	91
1	C	225/257 (88%)	223 (99%)	2 (1%)	87	91
1	D	220/257 (86%)	218 (99%)	2 (1%)	87	91
All	All	891/1028 (87%)	884 (99%)	7 (1%)	89	94

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1084	CYS
1	B	1065	MET
1	B	1104	ASP
1	C	1055	LYS
1	C	1065	MET
1	D	1084	CYS
1	D	1147	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1046	GLN
1	A	1064	ASN
1	A	1086	ASN

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 22 ligands modelled in this entry, 16 are monoatomic - leaving 6 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$
3	E72	A	1236	-	41,41,41	1.17	4 (9%)	52,52,52	1.78	9 (17%)
3	E72	B	2	-	41,41,41	1.19	4 (9%)	52,52,52	1.74	9 (17%)
4	SAH	C	103	-	28,28,28	4.35	21 (75%)	40,40,40	3.05	21 (52%)
3	E72	C	1236	-	41,41,41	1.15	4 (9%)	52,52,52	1.86	9 (17%)
4	SAH	D	104	-	28,28,28	1.26	3 (10%)	40,40,40	2.63	9 (22%)
3	E72	D	4	-	41,41,41	1.21	4 (9%)	52,52,52	1.82	12 (23%)

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
3	E72	A	1236	-	-	0/26/36/36	0/1/3/3
3	E72	B	2	-	-	0/26/36/36	0/1/3/3
4	SAH	C	103	-	-	0/15/31/31	0/1/3/3
3	E72	C	1236	-	-	0/26/36/36	0/1/3/3
4	SAH	D	104	-	-	0/15/31/31	0/1/3/3
3	E72	D	4	-	-	0/26/36/36	0/1/3/3

All (40) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	103	SAH	C4-N9	-10.31	1.22	1.37
4	C	103	SAH	C8-N9	-7.04	1.26	1.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	103	SAH	O3'-C3'	-6.87	1.26	1.43
4	C	103	SAH	O4'-C1'	-6.17	1.31	1.41
4	C	103	SAH	C5'-SD	-6.10	1.68	1.81
4	C	103	SAH	C8-N7	-5.70	1.23	1.34
4	C	103	SAH	C2'-C3'	-5.02	1.39	1.53
4	C	103	SAH	C4-N3	-4.77	1.28	1.35
4	C	103	SAH	C3'-C4'	-4.77	1.39	1.53
4	C	103	SAH	C1'-N9	-4.73	1.33	1.48
4	C	103	SAH	CG-SD	-4.62	1.63	1.81
3	B	2	E72	CAF-CAA	3.86	1.43	1.36
3	C	1236	E72	CAF-CAA	3.70	1.43	1.36
4	C	103	SAH	C2'-C1'	-3.69	1.48	1.53
4	C	103	SAH	C2-N1	-3.66	1.26	1.33
3	A	1236	E72	CAF-CAA	3.64	1.43	1.36
3	D	4	E72	CAC-CAB	3.62	1.43	1.36
4	D	104	SAH	C4-N9	-3.53	1.32	1.37
3	D	4	E72	CAF-CAA	3.44	1.42	1.36
3	B	2	E72	CAC-CAB	3.44	1.42	1.36
4	C	103	SAH	O4'-C4'	-3.43	1.37	1.45
4	C	103	SAH	C5'-C4'	3.42	1.62	1.52
3	A	1236	E72	CAC-CAB	3.34	1.42	1.36
3	C	1236	E72	C6-C5	-3.29	1.40	1.45
4	C	103	SAH	C5-N7	-3.24	1.28	1.40
4	C	103	SAH	C5-C4	-3.21	1.33	1.40
3	A	1236	E72	C6-C5	-3.21	1.41	1.45
4	D	104	SAH	OXT-C	3.17	1.41	1.30
3	C	1236	E72	CAC-CAB	3.03	1.42	1.36
3	B	2	E72	C6-C5	-3.00	1.41	1.45
3	D	4	E72	C6-C5	-2.96	1.41	1.45
3	B	2	E72	C6-N1	2.92	1.37	1.33
3	A	1236	E72	C6-N1	2.92	1.37	1.33
4	C	103	SAH	C6-N6	-2.91	1.25	1.35
3	D	4	E72	C6-N1	2.89	1.37	1.33
3	C	1236	E72	C6-N1	2.45	1.36	1.33
4	C	103	SAH	C2-N3	-2.36	1.27	1.32
4	D	104	SAH	C8-N9	-2.33	1.33	1.36
4	C	103	SAH	O2'-C2'	-2.12	1.37	1.43
4	C	103	SAH	C6-N1	-2.09	1.28	1.37

All (69) bond angle outliers are listed below:

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	104	SAH	N3-C2-N1	-8.64	121.48	128.71
4	C	103	SAH	O4'-C1'-N9	-8.59	100.45	108.44
4	D	104	SAH	OXT-C-CA	8.38	135.68	116.88
4	C	103	SAH	N3-C2-N1	-7.81	122.18	128.71
4	D	104	SAH	OXT-C-O	-7.71	106.64	124.07
3	C	1236	E72	CAC-C5-C6	-6.71	121.38	124.98
3	A	1236	E72	CAC-C5-C6	-6.44	121.52	124.98
3	D	4	E72	CAC-C5-C6	-6.31	121.59	124.98
3	B	2	E72	CAC-C5-C6	-6.26	121.62	124.98
4	C	103	SAH	C4-C5-N7	-5.47	104.84	109.52
4	C	103	SAH	OXT-C-CA	4.55	127.10	116.88
3	C	1236	E72	C6-C5-C4	4.53	118.21	115.63
4	D	104	SAH	N3-C4-N9	4.48	133.53	125.43
3	D	4	E72	C6-C5-C4	4.42	118.14	115.63
4	C	103	SAH	CB-CG-SD	4.40	122.05	113.57
3	A	1236	E72	C6-C5-C4	4.30	118.08	115.63
3	B	2	E72	C6-C5-C4	4.27	118.06	115.63
4	C	103	SAH	O3'-C3'-C2'	-4.23	98.06	111.83
4	C	103	SAH	N6-C6-N1	-4.11	111.30	119.36
3	C	1236	E72	OAL-CAB-CAC	-4.10	120.09	125.25
4	C	103	SAH	N3-C4-N9	4.09	132.81	125.43
3	B	2	E72	N3-C2-N1	-4.03	119.96	126.19
3	D	4	E72	N3-C2-N1	-3.99	120.03	126.19
3	A	1236	E72	N3-C2-N1	-3.92	120.13	126.19
3	C	1236	E72	N3-C2-N1	-3.89	120.18	126.19
3	A	1236	E72	CAM-OAL-CAB	-3.67	112.16	117.59
3	A	1236	E72	OAL-CAB-CAC	-3.64	120.66	125.25
3	D	4	E72	OAL-CAB-CAC	-3.60	120.71	125.25
3	C	1236	E72	C5-C4-N3	-3.60	119.55	122.86
3	B	2	E72	C5-C4-N3	-3.50	119.64	122.86
4	C	103	SAH	OXT-C-O	-3.40	116.38	124.07
3	C	1236	E72	CAM-OAL-CAB	-3.35	112.64	117.59
3	B	2	E72	CAM-OAL-CAB	-3.33	112.66	117.59
3	D	4	E72	C5-C4-N3	-3.30	119.82	122.86
3	A	1236	E72	C5-C4-N3	-3.27	119.85	122.86
4	C	103	SAH	C5-C6-N6	3.12	127.78	120.72
3	B	2	E72	OAL-CAB-CAC	-3.10	121.34	125.25
3	C	1236	E72	OAL-CAB-CAA	3.09	119.90	115.42
3	D	4	E72	CAM-OAL-CAB	-3.06	113.06	117.59
3	D	4	E72	OAL-CAB-CAA	3.03	119.82	115.42
4	C	103	SAH	C2-N3-C4	3.01	122.59	114.01

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	4	E72	OAK-CAA-CAF	-3.01	119.52	125.21
4	C	103	SAH	C5-C4-N3	-2.93	119.32	125.70
4	C	103	SAH	C4'-C5'-SD	2.84	122.29	113.53
4	C	103	SAH	C4'-O4'-C1'	-2.81	106.70	109.75
4	C	103	SAH	O3'-C3'-C4'	-2.80	102.82	111.08
4	D	104	SAH	C5-C4-N3	-2.79	119.62	125.70
4	C	103	SAH	C8-N9-C4	2.75	109.00	106.90
3	A	1236	E72	OAL-CAB-CAA	2.74	119.39	115.42
3	C	1236	E72	OAK-CAA-CAF	-2.70	120.11	125.21
4	C	103	SAH	C5'-C4'-C3'	-2.59	108.28	114.98
3	A	1236	E72	OAK-CAA-CAF	-2.53	120.43	125.21
4	D	104	SAH	C4-C5-N7	-2.50	107.38	109.52
3	D	4	E72	CBB-OAK-CAA	-2.41	111.73	117.65
3	B	2	E72	OAL-CAB-CAA	2.38	118.87	115.42
4	C	103	SAH	O2'-C2'-C3'	-2.30	104.36	111.83
4	D	104	SAH	O4'-C1'-N9	-2.30	106.30	108.44
3	A	1236	E72	C2-N1-C6	2.27	121.69	116.97
4	D	104	SAH	C8-N9-C4	2.25	108.62	106.90
4	D	104	SAH	C2-N3-C4	2.24	120.40	114.01
4	C	103	SAH	O4'-C4'-C5'	2.22	114.84	108.91
3	C	1236	E72	C2-N1-C6	2.21	121.57	116.97
3	D	4	E72	C2-N1-C6	2.16	121.45	116.97
4	C	103	SAH	C6-C5-N7	2.15	139.55	131.34
3	B	2	E72	C2-N1-C6	2.13	121.39	116.97
3	D	4	E72	OAK-CAA-CAB	2.13	120.30	115.80
3	D	4	E72	NAU-C2-N1	2.11	121.15	117.17
3	B	2	E72	NAU-C2-N1	2.08	121.09	117.17
4	C	103	SAH	C2'-C1'-N9	2.03	118.49	113.27

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	255/285 (89%)	0.12	7 (2%) 52 54	17, 32, 54, 59	0
1	B	258/285 (90%)	0.13	8 (3%) 47 49	18, 31, 53, 69	0
1	C	258/285 (90%)	0.07	6 (2%) 57 60	21, 32, 53, 58	0
1	D	256/285 (89%)	0.46	14 (5%) 24 26	26, 41, 61, 71	0
All	All	1027/1140 (90%)	0.19	35 (3%) 43 45	17, 34, 56, 71	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	979	VAL	5.8
1	D	1235	SER	4.6
1	B	1222	LEU	4.3
1	A	979	VAL	4.1
1	D	1228	GLY	4.1
1	A	1235	SER	4.0
1	D	1000	SER	3.7
1	B	1235	SER	3.5
1	D	978	ILE	3.4
1	B	1105	MET	3.3
1	B	1148	ASN	3.3
1	B	976	GLU	3.2
1	D	1049	MET	3.0
1	A	997	ALA	2.9
1	D	976	GLU	2.9
1	C	1034	ILE	2.8
1	D	1043	MET	2.7
1	D	1230	PRO	2.7
1	D	998	VAL	2.6
1	A	1221	LYS	2.6
1	C	998	VAL	2.6

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Mol	Chain	Res	Type	RSRZ
1	C	1151	GLY	2.6
1	A	1151	GLY	2.6
1	A	1202	GLU	2.5
1	A	998	VAL	2.5
1	D	977	ARG	2.4
1	B	1224	SER	2.3
1	C	1228	GLY	2.3
1	D	1050	ARG	2.2
1	C	1230	PRO	2.2
1	B	1217	ASP	2.2
1	D	1061	PRO	2.1
1	B	1221	LYS	2.1
1	D	1152	GLU	2.0
1	C	1223	PHE	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
4	SAH	D	104	26/26	0.23	2.13	62,68,69,70	0
3	E72	A	1236	39/39	0.18	1.60	38,43,51,52	0
2	ZN	B	6	1/1	0.11	1.28	27,27,27,27	0
4	SAH	C	103	26/26	0.20	1.22	50,56,58,59	0
2	ZN	A	1	1/1	0.10	1.18	22,22,22,22	0
3	E72	D	4	39/39	0.19	0.70	44,48,59,61	0
2	ZN	A	3	1/1	0.09	0.50	20,20,20,20	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	E72	C	1236	39/39	0.14	0.45	30,34,45,45	0
2	ZN	D	6	1/1	0.11	0.23	34,34,34,34	0
3	E72	B	2	39/39	0.14	0.15	31,36,45,47	0
2	ZN	B	5	1/1	0.10	0.09	28,28,28,28	0
2	ZN	C	2	1/1	0.09	0.01	28,28,28,28	0
2	ZN	C	3	1/1	0.09	-0.20	28,28,28,28	0
2	ZN	B	7	1/1	0.09	-0.26	28,28,28,28	0
2	ZN	A	2	1/1	0.09	-0.34	18,18,18,18	0
2	ZN	D	7	1/1	0.06	-1.09	39,39,39,39	0
2	ZN	C	1	1/1	0.07	-1.23	27,27,27,27	0
2	ZN	D	5	1/1	0.08	-1.35	35,35,35,35	0
2	ZN	A	4	1/1	0.07	-1.37	41,41,41,41	0
2	ZN	C	4	1/1	0.04	-1.72	39,39,39,39	0
2	ZN	B	8	1/1	0.05	-1.86	41,41,41,41	0
2	ZN	D	8	1/1	0.04	-3.25	49,49,49,49	0

6.5 Other polymers ⓘ

There are no such residues in this entry.