



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 09:41 pm GMT

PDB ID : 4M72
Title : Mutant structure of methyltransferase from *Streptomyces hygrosopicus*
Authors : Liu, Y.C.; Zou, X.W.; Chan, H.C.; Huang, C.J.; Li, T.L.
Deposited on : 2013-08-12
Resolution : 2.10 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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : recalc28949

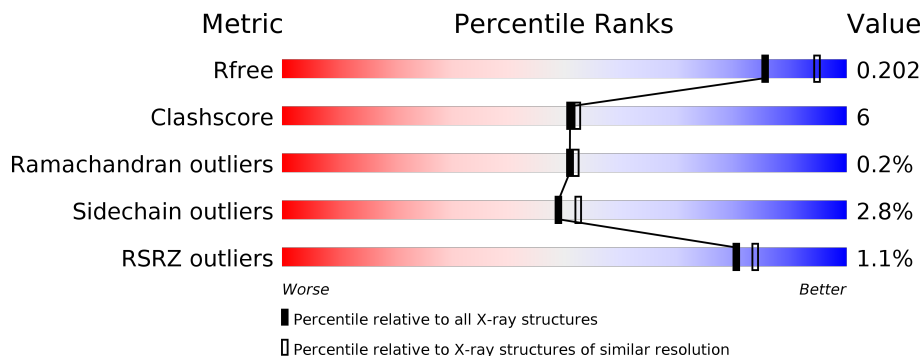
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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}	100719	4243 (2.10-2.10)
Clashscore	112137	4788 (2.10-2.10)
Ramachandran outliers	110173	4740 (2.10-2.10)
Sidechain outliers	110143	4741 (2.10-2.10)
RSRZ outliers	101464	4275 (2.10-2.10)

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	337	
1	B	337	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	M72	A	402	-	-	-	X
3	M72	B	402	-	-	-	X

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 6055 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.

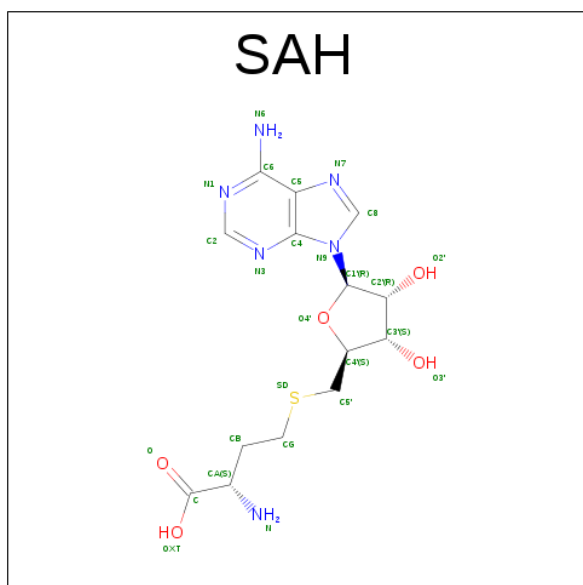
- Molecule 1 is a protein called Methyltransferase MppJ.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	333	Total	C	N	O	S	0	0	0
			2591	1638	459	483	11			
1	B	333	Total	C	N	O	S	0	0	0
			2591	1638	459	483	11			

There are 4 discrepancies between the modelled and reference sequences:

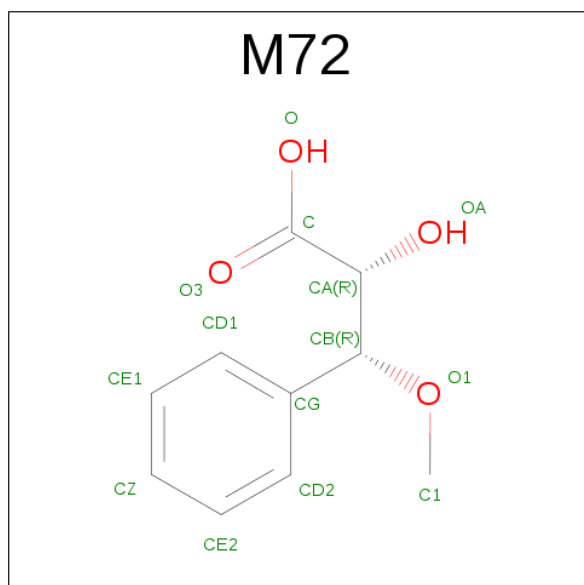
Chain	Residue	Modelled	Actual	Comment	Reference
A	127	LEU	ARG	ENGINEERED MUTATION	UNP Q643C8
A	300	GLU	VAL	ENGINEERED MUTATION	UNP Q643C8
B	127	LEU	ARG	ENGINEERED MUTATION	UNP Q643C8
B	300	GLU	VAL	ENGINEERED MUTATION	UNP Q643C8

- Molecule 2 is S-ADENOSYL-L-HOMOCYSTEINE (three-letter code: SAH) (formula: $C_{14}H_{20}N_6O_5S$).



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

- Molecule 3 is (2R,3R)-2-HYDROXY-3-METHOXY-3-PHENYLPROPANOIC ACID (three-letter code: M72) (formula: C₁₀H₁₂O₄).

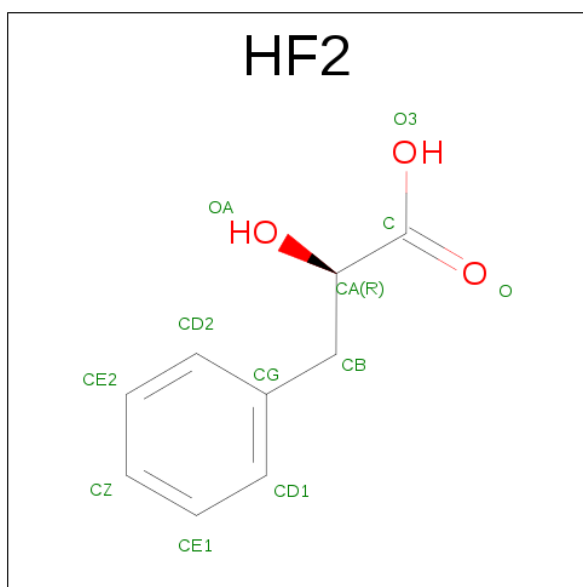


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			14	10	4		
3	B	1	Total	C	O	0	0
			14	10	4		

- Molecule 4 is FE (III) ION (three-letter code: FE) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Fe	0	0
			1	1		
4	A	1	Total	Fe	0	0
			1	1		

- Molecule 5 is (2R)-2-HYDROXY-3-PHENYLPROPANOIC ACID (three-letter code: HF2) (formula: C₉H₁₀O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			12	9	3		
5	B	1	Total	C	O	0	0
			12	9	3		

- Molecule 6 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	3	Total	Ca	0	0
			3	3		
6	A	5	Total	Ca	0	0
			5	5		

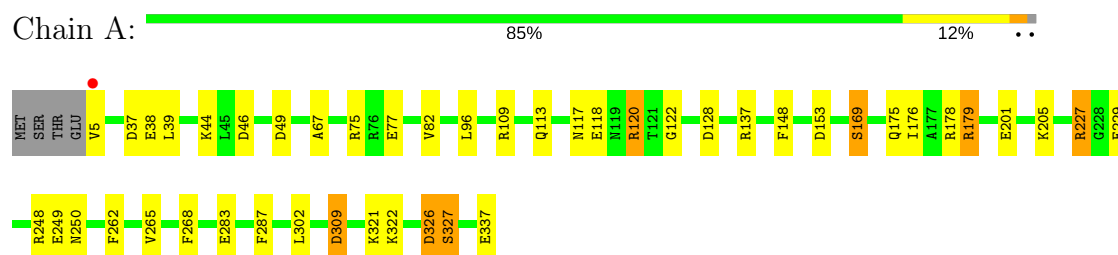
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	411	Total	O	0	0
			411	411		
7	B	348	Total	O	0	0
			348	348		

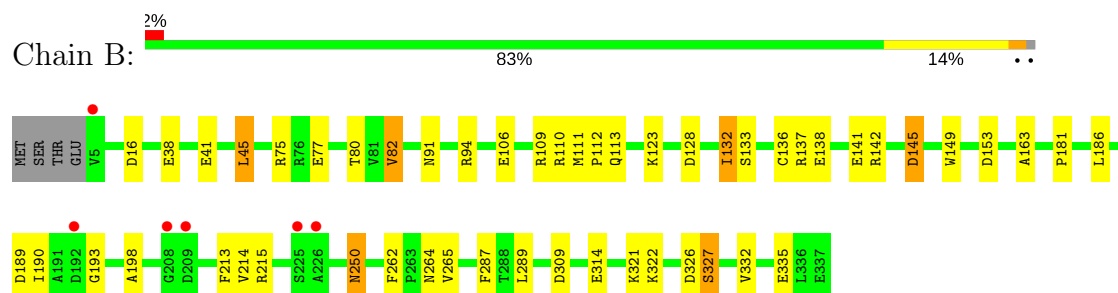
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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: Methyltransferase MppJ



• Molecule 1: Methyltransferase MppJ



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	60.87Å 97.14Å 135.23Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.06 – 2.10 29.04 – 2.10	Depositor EDS
% Data completeness (in resolution range)	95.3 (29.06-2.10) 95.3 (29.04-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.45 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.144 , 0.196 0.156 , 0.202	Depositor DCC
R_{free} test set	2308 reflections (5.35%)	DCC
Wilson B-factor (Å ²)	22.9	Xtriage
Anisotropy	0.058	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 53.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6055	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.17% 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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: M72, CA, SAH, HF2, FE

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	1.00	6/2650 (0.2%)	1.07	9/3587 (0.3%)
1	B	0.95	4/2650 (0.2%)	1.01	10/3587 (0.3%)
All	All	0.97	10/5300 (0.2%)	1.04	19/7174 (0.3%)

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	A	0	1
1	B	0	1
All	All	0	2

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	169	SER	CB-OG	6.75	1.51	1.42
1	B	327	SER	N-CA	-6.42	1.33	1.46
1	A	337	GLU	CD-OE2	-5.74	1.19	1.25
1	A	249	GLU	CG-CD	5.47	1.60	1.51
1	A	120	ARG	CD-NE	-5.45	1.37	1.46
1	B	149	TRP	CB-CG	-5.43	1.40	1.50
1	B	314	GLU	CD-OE1	5.37	1.31	1.25
1	A	283	GLU	CD-OE1	-5.36	1.19	1.25
1	B	138	GLU	CD-OE2	5.36	1.31	1.25
1	A	327	SER	N-CA	-5.30	1.35	1.46

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	120	ARG	NE-CZ-NH2	-18.69	110.95	120.30
1	A	120	ARG	NE-CZ-NH1	14.35	127.47	120.30
1	B	142	ARG	NE-CZ-NH1	8.32	124.46	120.30
1	A	327	SER	N-CA-CB	-7.95	98.58	110.50
1	A	248	ARG	NE-CZ-NH1	7.19	123.90	120.30
1	B	327	SER	N-CA-C	-6.97	92.18	111.00
1	B	45	LEU	CA-CB-CG	6.44	130.12	115.30
1	B	110	ARG	NE-CZ-NH1	6.14	123.37	120.30
1	B	16	ASP	CB-CG-OD1	6.08	123.77	118.30
1	B	142	ARG	NE-CZ-NH2	-5.94	117.33	120.30
1	A	128	ASP	CB-CG-OD1	5.78	123.51	118.30
1	A	109	ARG	NE-CZ-NH1	5.72	123.16	120.30
1	A	309	ASP	CB-CG-OD1	5.57	123.31	118.30
1	B	309	ASP	CB-CG-OD1	5.49	123.24	118.30
1	B	16	ASP	CB-CG-OD2	-5.45	113.39	118.30
1	A	326	ASP	C-N-CA	5.31	134.99	121.70
1	B	332	VAL	CG1-CB-CG2	-5.09	102.75	110.90
1	B	214	VAL	CB-CA-C	-5.03	101.85	111.40
1	A	248	ARG	NE-CZ-NH2	-5.02	117.79	120.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	326	ASP	Peptide
1	B	326	ASP	Peptide

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	2591	0	2503	33	0
1	B	2591	0	2503	30	0
2	A	26	0	19	1	0
2	B	26	0	19	0	0
3	A	14	0	11	0	0
3	B	14	0	11	2	0
4	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	1	0	0	0	0
5	A	12	0	7	2	0
5	B	12	0	7	2	0
6	A	5	0	0	1	0
6	B	3	0	0	0	0
7	A	411	0	0	20	0
7	B	348	0	0	15	0
All	All	6055	0	5080	67	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 6.

All (67) 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:B:153:ASP:HB3	7:B:593:HOH:O	1.45	1.14
1:B:75:ARG:NH1	1:B:77:GLU:OE2	2.03	0.91
6:A:407:CA:CA	7:A:778:HOH:O	1.49	0.86
1:B:41:GLU:OE2	7:B:832:HOH:O	1.95	0.84
1:B:137:ARG:NH1	7:B:810:HOH:O	2.13	0.81
1:A:137:ARG:NH1	7:A:855:HOH:O	2.17	0.78
1:A:175:GLN:HG3	7:A:739:HOH:O	1.84	0.78
1:A:122:GLY:O	7:A:892:HOH:O	2.02	0.77
1:A:321:LYS:HE2	7:A:520:HOH:O	1.84	0.77
1:A:153:ASP:OD1	1:A:179:ARG:NH2	2.18	0.74
1:B:113:GLN:OE1	7:B:806:HOH:O	2.04	0.73
1:A:38:GLU:OE2	7:A:888:HOH:O	2.08	0.71
1:A:169:SER:OG	7:A:898:HOH:O	2.09	0.70
1:A:201:GLU:HG2	7:A:831:HOH:O	1.92	0.69
1:B:106:GLU:OE2	1:B:109:ARG:NH1	2.27	0.68
1:A:37:ASP:OD1	7:A:881:HOH:O	2.14	0.66
1:A:227:ARG:NH1	7:A:900:HOH:O	2.14	0.65
1:B:82:VAL:HG13	7:B:567:HOH:O	1.97	0.64
1:B:38:GLU:OE1	7:B:744:HOH:O	2.15	0.64
1:A:75:ARG:NH1	1:A:77:GLU:OE2	2.28	0.63
1:A:5:VAL:HG22	7:B:812:HOH:O	2.01	0.61
1:A:5:VAL:N	7:A:745:HOH:O	2.36	0.59
3:B:402:M72:OA	7:B:699:HOH:O	1.92	0.59
1:A:113:GLN:OE1	7:A:549:HOH:O	2.16	0.58
1:A:229:GLU:CD	7:A:689:HOH:O	2.43	0.57
1:B:250:ASN:H	1:B:250:ASN:HD22	1.52	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:80:THR:HG22	7:B:808:HOH:O	2.07	0.55
1:B:136:CYS:SG	3:B:402:M72:O1	2.65	0.55
1:A:322:LYS:HD3	5:A:404:HF2:HA	1.88	0.54
1:A:120:ARG:HD3	7:A:632:HOH:O	2.08	0.53
1:B:193:GLY:N	7:B:763:HOH:O	2.10	0.53
1:A:309:ASP:OD2	7:A:827:HOH:O	2.19	0.52
1:B:141:GLU:O	1:B:145:ASP:HB2	2.10	0.52
1:B:91:ASN:O	1:B:94:ARG:HD3	2.10	0.52
1:B:190:ILE:O	1:B:215:ARG:NH1	2.47	0.48
1:A:262:PHE:HB3	1:A:265:VAL:HG23	1.96	0.47
1:A:229:GLU:HG2	7:A:689:HOH:O	2.14	0.47
1:B:322:LYS:HD3	5:B:404:HF2:HA	1.96	0.46
1:A:117:ASN:ND2	7:A:637:HOH:O	2.47	0.46
1:A:148:PHE:HZ	1:A:176:ILE:HD11	1.80	0.46
1:A:118:GLU:HG2	7:B:831:HOH:O	2.16	0.46
1:B:321:LYS:HE3	7:B:546:HOH:O	2.16	0.45
1:B:189:ASP:O	1:B:215:ARG:HA	2.17	0.44
1:A:75:ARG:HD3	7:A:749:HOH:O	2.15	0.44
1:B:181:PRO:HA	7:B:740:HOH:O	2.17	0.44
1:B:198:ALA:HB1	1:B:213:PHE:CZ	2.52	0.44
5:B:404:HF2:C	5:B:404:HF2:CD1	2.96	0.44
1:B:128:ASP:O	1:B:132:ILE:HD12	2.17	0.44
1:B:109:ARG:NH2	7:B:819:HOH:O	2.51	0.43
1:B:163:ALA:HA	1:B:186:LEU:O	2.19	0.43
1:A:175:GLN:NE2	1:A:178:ARG:HH11	2.16	0.43
2:A:401:SAH:HB2	7:A:748:HOH:O	2.17	0.43
1:B:262:PHE:HB3	1:B:265:VAL:HG23	2.00	0.43
1:A:321:LYS:CE	7:A:520:HOH:O	2.57	0.42
1:B:321:LYS:HE2	1:B:335:GLU:OE2	2.19	0.42
1:A:67:ALA:HA	1:B:289:LEU:HD21	2.01	0.42
1:A:46:ASP:HB3	1:A:49:ASP:HB3	2.02	0.42
1:B:264:ASN:ND2	7:B:768:HOH:O	2.49	0.42
1:A:227:ARG:HB2	1:A:227:ARG:HH11	1.84	0.42
1:B:111:MET:HB3	1:B:112:PRO:HD3	2.02	0.41
1:A:250:ASN:H	1:A:250:ASN:HD22	1.68	0.41
1:A:96:LEU:HD13	1:A:287:PHE:HB2	2.01	0.41
1:A:175:GLN:HE22	1:A:178:ARG:HH11	1.69	0.41
1:B:106:GLU:OE2	1:B:109:ARG:CZ	2.69	0.41
1:B:250:ASN:HD22	1:B:250:ASN:N	2.16	0.40
5:A:404:HF2:CD1	5:A:404:HF2:C	2.99	0.40
1:A:39:LEU:HD23	1:A:44:LYS:HA	2.04	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	331/337 (98%)	320 (97%)	11 (3%)	0	100	100
1	B	331/337 (98%)	322 (97%)	8 (2%)	1 (0%)	44	44
All	All	662/674 (98%)	642 (97%)	19 (3%)	1 (0%)	51	52

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	287	PHE

5.3.2 Protein sidechains [i](#)

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	264/268 (98%)	257 (97%)	7 (3%)	50	54
1	B	264/268 (98%)	256 (97%)	8 (3%)	46	49
All	All	528/536 (98%)	513 (97%)	15 (3%)	49	52

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	82	VAL
1	A	179	ARG

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Mol	Chain	Res	Type
1	A	205	LYS
1	A	227	ARG
1	A	268	PHE
1	A	302	LEU
1	A	327	SER
1	B	45	LEU
1	B	82	VAL
1	B	123	LYS
1	B	132	ILE
1	B	133	SER
1	B	145	ASP
1	B	250	ASN
1	B	327	SER

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	113	GLN
1	A	117	ASN
1	A	175	GLN
1	A	250	ASN
1	B	113	GLN
1	B	117	ASN
1	B	250	ASN
1	B	264	ASN

5.3.3 RNA ⓘ

There are no RNA molecules 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 16 ligands modelled in this entry, 10 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$
2	SAH	A	401	-	20,28,28	1.42	3 (15%)	20,40,40	2.49	7 (35%)
3	M72	A	402	4	10,14,14	1.54	2 (20%)	13,18,18	1.69	2 (15%)
5	HF2	A	404	-	9,12,12	0.50	0	11,15,15	1.88	1 (9%)
2	SAH	B	401	-	20,28,28	1.25	3 (15%)	20,40,40	3.01	5 (25%)
3	M72	B	402	4	10,14,14	1.52	2 (20%)	13,18,18	1.98	4 (30%)
5	HF2	B	404	-	9,12,12	0.61	0	11,15,15	3.11	1 (9%)

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	SAH	A	401	-	-	0/7/31/31	0/3/3/3
3	M72	A	402	4	-	1/10/14/14	0/1/1/1
5	HF2	A	404	-	-	0/4/8/8	0/1/1/1
2	SAH	B	401	-	-	0/7/31/31	0/3/3/3
3	M72	B	402	4	-	1/10/14/14	0/1/1/1
5	HF2	B	404	-	-	0/4/8/8	0/1/1/1

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	SAH	C2-N1	2.17	1.38	1.33
2	B	401	SAH	O4'-C1'	2.21	1.44	1.41
3	A	402	M72	O1-CB	2.37	1.47	1.42
2	B	401	SAH	C2-N3	2.41	1.36	1.32
3	B	402	M72	O1-CB	2.64	1.48	1.42
2	B	401	SAH	C4-N3	2.65	1.39	1.35
3	A	402	M72	CD2-CG	2.80	1.43	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	SAH	C2-N3	2.80	1.36	1.32
3	B	402	M72	CD2-CG	2.93	1.43	1.39
2	A	401	SAH	C5-C4	3.44	1.48	1.40

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	SAH	N3-C2-N1	-10.21	119.97	128.86
2	A	401	SAH	N3-C2-N1	-5.65	123.94	128.86
2	B	401	SAH	C1'-N9-C4	-5.35	117.39	126.64
2	A	401	SAH	CB-CG-SD	-5.15	103.65	113.57
3	B	402	M72	CG-CB-CA	-5.02	101.61	112.18
3	A	402	M72	CG-CB-CA	-4.61	102.48	112.18
2	A	401	SAH	O4'-C4'-C5'	-4.52	96.91	108.82
2	B	401	SAH	C5-C6-N6	-3.63	113.08	120.47
2	A	401	SAH	C1'-N9-C4	-3.58	120.44	126.64
2	B	401	SAH	CB-CG-SD	-2.99	107.81	113.57
2	A	401	SAH	C4-C5-N7	-2.76	106.74	109.41
3	B	402	M72	CZ-CE1-CD1	2.00	122.96	120.21
2	A	401	SAH	C2-N1-C6	2.24	122.68	118.77
3	A	402	M72	O1-CB-CA	2.38	119.62	107.50
3	B	402	M72	C1-O1-CB	2.42	121.19	113.58
3	B	402	M72	O1-CB-CA	2.54	120.42	107.50
2	A	401	SAH	O3'-C3'-C4'	2.55	118.53	111.09
2	B	401	SAH	N6-C6-N1	3.18	125.07	118.77
5	A	404	HF2	CB-CA-C	5.93	118.95	111.11
5	B	404	HF2	CB-CA-C	10.07	124.43	111.11

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	402	M72	C1-O1-CB-CA
3	B	402	M72	C1-O1-CB-CA

There are no ring outliers.

4 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	SAH	1	0
5	A	404	HF2	2	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	402	M72	2	0
5	B	404	HF2	2	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	333/337 (98%)	-0.53	1 (0%) 93 94	12, 21, 40, 54	0
1	B	333/337 (98%)	-0.32	6 (1%) 69 73	14, 26, 49, 66	0
All	All	666/674 (98%)	-0.43	7 (1%) 80 84	12, 23, 45, 66	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	5	VAL	5.6
1	A	5	VAL	3.1
1	B	192	ASP	2.7
1	B	209	ASP	2.3
1	B	225	SER	2.3
1	B	226	ALA	2.3
1	B	208	GLY	2.2

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(Å ²)	Q<0.9
3	M72	B	402	14/14	0.81	0.19	5.53	31,49,56,64	0
3	M72	A	402	14/14	0.89	0.16	4.24	24,38,51,52	0
6	CA	A	407	1/1	0.99	0.10	1.19	8,8,8,8	0
2	SAH	B	401	26/26	0.92	0.10	-0.09	29,41,55,60	0
5	HF2	B	404	12/12	0.96	0.09	-0.36	24,26,35,40	0
2	SAH	A	401	26/26	0.97	0.07	-0.52	16,22,46,51	0
5	HF2	A	404	12/12	0.96	0.08	-0.61	26,28,32,35	0
6	CA	A	408	1/1	0.87	0.04	-2.04	67,67,67,67	0
6	CA	B	405	1/1	0.97	0.08	-	34,34,34,34	0
6	CA	A	406	1/1	0.99	0.04	-	38,38,38,38	0
6	CA	B	407	1/1	0.95	0.08	-	58,58,58,58	0
6	CA	B	406	1/1	0.98	0.04	-	49,49,49,49	0
6	CA	A	409	1/1	0.89	0.07	-	45,45,45,45	0
4	FE	A	403	1/1	0.99	0.07	-	16,16,16,16	0
4	FE	B	403	1/1	0.99	0.06	-	19,19,19,19	0
6	CA	A	405	1/1	0.98	0.03	-	42,42,42,42	0

6.5 Other polymers [i](#)

There are no such residues in this entry.