



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:35 AM GMT

PDB ID : 2HTE
Title : The crystal structure of spermidine synthase from *p. falciparum* in complex with 5'-methylthioadenosine
Authors : Qiu, W.; Dong, A.; Ren, H.; Wu, H.; Wasney, G.; Vedadi, M.; Lew, J.; Kozieradski, I.; Edwards, A.M.; Arrowsmith, C.H.; Weigelt, J.; Sundstrom, M.; Plotnikov, A.N.; Bochkarev, A.; Hui, R.; Structural Genomics Consortium (SGC)
Deposited on : 2006-07-25
Resolution : 2.00 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 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : trunk26865

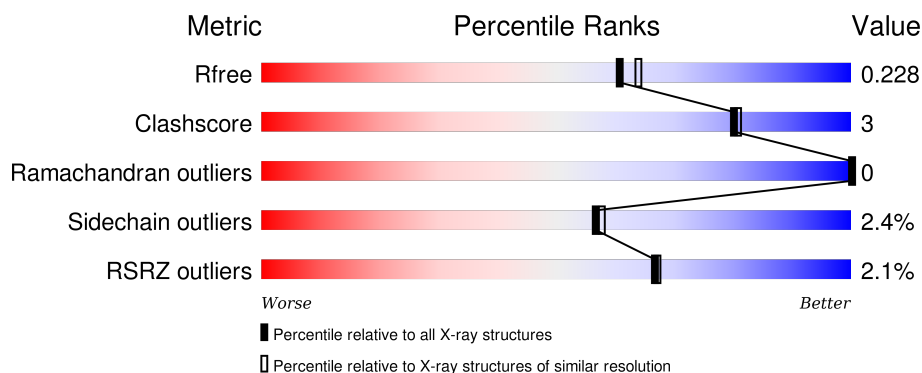
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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}	91344	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (2.00-2.00)

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	283	<div> <div>3%</div> <div>87%</div> <div>8%</div> <div>• •</div> </div>
1	B	283	<div> <div>2%</div> <div>84%</div> <div>11%</div> <div>• 5%</div> </div>
1	C	283	<div> <div>0%</div> <div>89%</div> <div>10%</div> <div>•</div> </div>

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
5	1PG	C	6001	-	-	-	X

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 7404 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 Spermidine synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	272	Total	C	N	O	S	8	6	0
			2220	1444	345	416	15			
1	B	270	Total	C	N	O	S	5	8	0
			2210	1435	343	416	16			
1	C	281	Total	C	N	O	S	1	9	0
			2297	1497	355	430	15			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	cloning artifact	UNP Q8II73
B	1	GLY	-	cloning artifact	UNP Q8II73
C	1	GLY	-	cloning artifact	UNP Q8II73

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

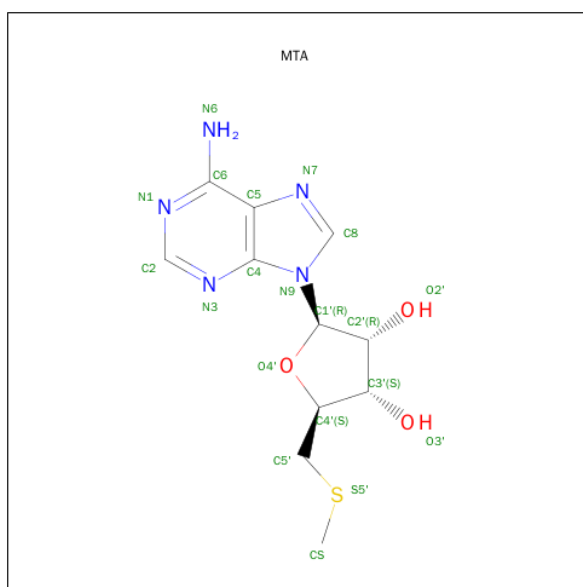


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is UNKNOWN ATOM OR ION (three-letter code: UNX) (formula: X).

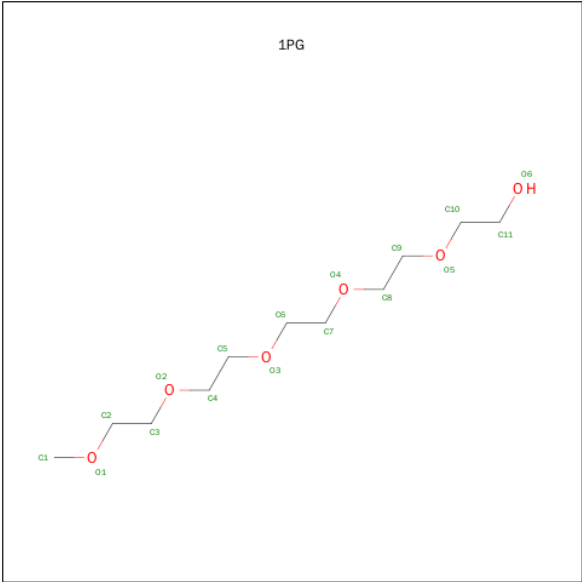
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	5	Total	X	0	0
			5	5		
3	A	5	Total	X	0	0
			5	5		
3	C	3	Total	X	0	0
			3	3		

- Molecule 4 is 5'-DEOXY-5'-METHYLTHIOADENOSINE (three-letter code: MTA) (formula: C₁₁H₁₅N₅O₃S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	S	0	0
			20	11	5	3	1		
4	B	1	Total	C	N	O	S	0	0
			20	11	5	3	1		
4	C	1	Total	C	N	O	S	0	0
			20	11	5	3	1		

- Molecule 5 is 2-(2-{2-[2-(2-METHOXY-ETHOXY)-ETHOXY]-ETHOXY}-ETHOXY)-ETHANOL (three-letter code: 1PG) (formula: C₁₁H₂₄O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			11	7	4		
5	C	1	Total	C	O	0	0
			17	11	6		

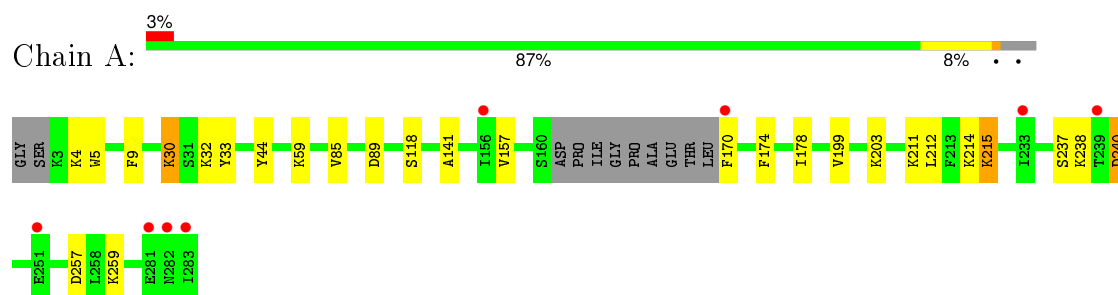
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	141	Total	O	0	0
			141	141		
6	B	159	Total	O	0	0
			159	159		
6	C	271	Total	O	0	0
			271	271		

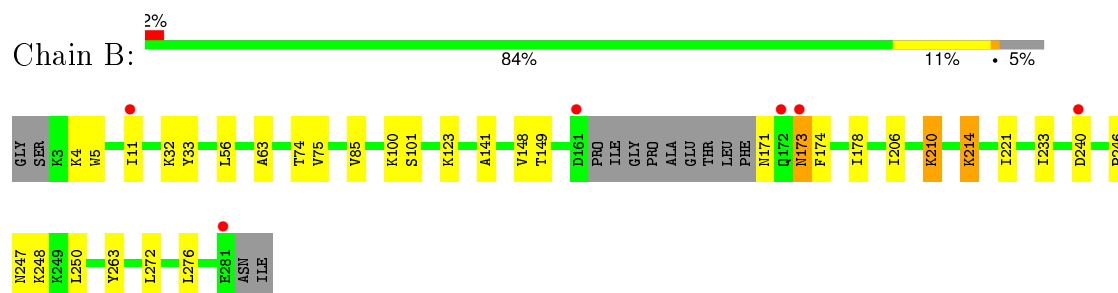
3 Residue-property plots [i](#)

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

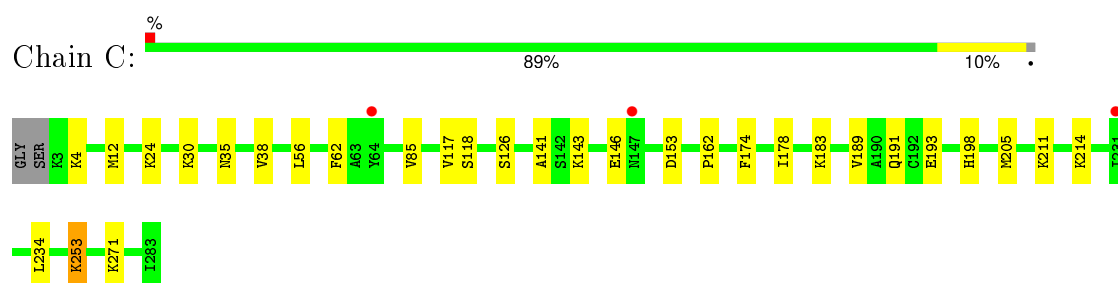
- Molecule 1: Spermidine synthase



- Molecule 1: Spermidine synthase



- Molecule 1: Spermidine synthase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	197.06 Å 134.60 Å 48.34 Å 90.00° 94.66° 90.00°	Depositor
Resolution (Å)	33.08 – 2.00 33.07 – 2.00	Depositor EDS
% Data completeness (in resolution range)	98.5 (33.08-2.00) 98.4 (33.07-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.14 (at 2.00 Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.185 , 0.231 0.183 , 0.228	Depositor DCC
R_{free} test set	4135 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	28.5	Xtriage
Anisotropy	0.076	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 49.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 82754 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7404	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 3.90% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: UNX, MTA, 1PG, 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	0.84	3/2285 (0.1%)	0.77	3/3080 (0.1%)
1	B	0.89	2/2280 (0.1%)	0.78	0/3074
1	C	0.87	0/2374	0.76	0/3204
All	All	0.87	5/6939 (0.1%)	0.77	3/9358 (0.0%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	259	LYS	CE-NZ	-13.69	1.14	1.49
1	A	240	ASP	CB-CG	6.69	1.65	1.51
1	A	30	LYS	CD-CE	-6.64	1.34	1.51
1	B	214	LYS	CD-CE	-6.19	1.35	1.51
1	B	63	ALA	CA-CB	6.03	1.65	1.52

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	240	ASP	CB-CG-OD2	-11.27	108.15	118.30
1	A	240	ASP	CB-CG-OD1	11.08	128.28	118.30
1	A	257	ASP	CB-CG-OD1	6.31	123.97	118.30

There are no chirality outliers.

There are no planarity outliers.

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	2220	0	2246	17	0
1	B	2210	0	2237	13	0
1	C	2297	0	2342	17	0
2	B	5	0	0	0	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
3	C	3	0	0	0	0
4	A	20	0	15	1	0
4	B	20	0	15	0	0
4	C	20	0	15	0	0
5	A	11	0	13	3	0
5	C	17	0	24	0	0
6	A	141	0	0	0	0
6	B	159	0	0	0	0
6	C	271	0	0	4	0
All	All	7404	0	6907	46	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 3.

All (46) 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:C:253:LYS:HD2	6:C:6180:HOH:O	1.53	1.09
1:C:205:MET:HE2	6:C:6272:HOH:O	1.76	0.85
1:C:191:GLN:HE21	1:C:193:GLU:H	1.38	0.71
1:B:272:LEU:HD13	1:B:276:LEU:HD23	1.74	0.69
1:B:174:PHE:CE2	1:B:178:ILE:HD11	2.31	0.66
1:A:174:PHE:CE2	1:A:178[A]:ILE:HD11	2.31	0.66
1:A:157:VAL:HG11	1:A:170:PHE:CZ	2.37	0.60
1:C:174:PHE:CE2	1:C:178:ILE:HD11	2.38	0.58
1:A:9:PHE:HZ	5:A:6000:1PG:H101	1.68	0.58
1:C:193:GLU:HB3	1:C:198:HIS:ND1	2.18	0.57
1:A:9:PHE:CZ	5:A:6000:1PG:H101	2.40	0.56
1:B:171:ASN:HD21	1:B:173:ASN:HB2	1.70	0.56
1:B:75:VAL:O	1:B:247:ASN:ND2	2.41	0.53
1:A:30:LYS:O	1:B:123:LYS:NZ	2.43	0.52
1:A:214:LYS:HD2	1:A:238:LYS:O	2.09	0.52
1:A:89:ASP:OD2	4:A:4000:MTA:HCS3	2.11	0.51
1:C:62:PHE:HD2	6:C:6108:HOH:O	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:32:LYS:HB3	1:B:33:TYR:CD2	2.47	0.50
1:A:215:LYS:HB3	1:A:237:SER:HB2	1.92	0.50
1:A:32:LYS:HE3	1:A:33:TYR:CE1	2.47	0.50
1:B:85:VAL:HG13	1:B:141:ALA:CB	2.43	0.48
1:A:9:PHE:HZ	5:A:6000:1PG:C10	2.27	0.47
1:B:246:PRO:HG2	1:B:263:TYR:CE1	2.51	0.46
1:B:206:ILE:HG22	1:B:210:LYS:HD2	1.97	0.46
1:A:32:LYS:HE3	1:A:33:TYR:CZ	2.52	0.45
1:A:4:LYS:HD2	1:A:5:TRP:H	1.82	0.45
1:C:12[B]:MET:SD	1:C:162:PRO:HD2	2.55	0.45
1:A:199:VAL:O	1:A:203:LYS:HG3	2.17	0.45
1:C:56:LEU:HD12	1:C:56:LEU:C	2.37	0.45
1:C:118[B]:SER:OG	1:C:126:SER:HB3	2.16	0.45
1:C:211:LYS:HE2	1:C:211:LYS:HA	2.00	0.44
1:A:44:TYR:CD2	1:A:59:LYS:HD3	2.52	0.44
1:A:89:ASP:HA	1:A:118[B]:SER:OG	2.17	0.44
1:C:271:LYS:HE3	6:C:6037:HOH:O	2.17	0.44
1:B:221:ILE:HD11	1:B:233:ILE:HD11	2.01	0.43
1:A:157:VAL:CG1	1:A:170:PHE:CZ	3.02	0.42
1:C:30:LYS:HG3	1:C:35:ASN:ND2	2.35	0.42
1:C:38:VAL:HG21	1:C:117[B]:VAL:CG1	2.50	0.42
1:C:143:LYS:HD3	1:C:146:GLU:OE2	2.20	0.41
1:A:85:VAL:HG13	1:A:141:ALA:HB2	2.03	0.41
1:C:189:VAL:HA	1:C:234:LEU:O	2.21	0.41
1:C:85:VAL:HG13	1:C:141:ALA:CB	2.51	0.41
1:C:153:ASP:OD1	1:C:183[A]:LYS:HD2	2.20	0.41
1:B:74:THR:HG22	1:B:250:LEU:HD21	2.02	0.40
1:B:4:LYS:HG3	1:B:5:TRP:N	2.36	0.40
1:B:56:LEU:HD12	1:B:56:LEU:C	2.41	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	274/283 (97%)	267 (97%)	7 (3%)	0	100	100
1	B	274/283 (97%)	268 (98%)	6 (2%)	0	100	100
1	C	288/283 (102%)	281 (98%)	7 (2%)	0	100	100
All	All	836/849 (98%)	816 (98%)	20 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	252/254 (99%)	248 (98%)	4 (2%)	70	73
1	B	252/254 (99%)	242 (96%)	10 (4%)	38	33
1	C	262/254 (103%)	257 (98%)	5 (2%)	65	67
All	All	766/762 (100%)	747 (98%)	19 (2%)	57	55

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

Mol	Chain	Res	Type
1	A	211	LYS
1	A	212	LEU
1	A	215	LYS
1	A	240	ASP
1	B	11	ILE
1	B	100	LYS
1	B	101	SER
1	B	148	VAL
1	B	149	THR
1	B	173	ASN
1	B	210	LYS
1	B	214	LYS
1	B	240	ASP
1	B	248	LYS
1	C	4	LYS

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Mol	Chain	Res	Type
1	C	24[A]	LYS
1	C	24[B]	LYS
1	C	214	LYS
1	C	253	LYS

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

Mol	Chain	Res	Type
1	A	172	GLN
1	B	104	ASN
1	B	135	ASN
1	B	171	ASN
1	B	191	GLN
1	C	35	ASN
1	C	104	ASN
1	C	135	ASN
1	C	172	GLN
1	C	191	GLN
1	C	247	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 19 ligands modelled in this entry, 13 are unknown - 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
4	MTA	A	4000	-	17,22,22	1.22	2 (11%)	17,32,32	3.58	6 (35%)
5	1PG	A	6000	-	10,10,16	1.60	4 (40%)	9,9,15	1.48	2 (22%)
4	MTA	B	4001	-	17,22,22	1.20	1 (5%)	17,32,32	2.71	4 (23%)
2	SO4	B	5000	-	4,4,4	0.20	0	6,6,6	0.56	0
4	MTA	C	4002	-	17,22,22	0.99	1 (5%)	17,32,32	2.72	6 (35%)
5	1PG	C	6001	-	16,16,16	0.96	1 (6%)	15,15,15	0.72	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	MTA	A	4000	-	-	0/3/23/23	0/3/3/3
5	1PG	A	6000	-	-	0/8/8/14	0/0/0/0
4	MTA	B	4001	-	-	0/3/23/23	0/3/3/3
2	SO4	B	5000	-	-	0/0/0/0	0/0/0/0
4	MTA	C	4002	-	-	0/3/23/23	0/3/3/3
5	1PG	C	6001	-	-	0/14/14/14	0/0/0/0

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	6000	1PG	O4-C8	-2.46	1.31	1.42
5	C	6001	1PG	O5-C9	-2.32	1.32	1.42
5	A	6000	1PG	O4-C7	-2.30	1.32	1.42
5	A	6000	1PG	O5-C9	-2.19	1.32	1.42
5	A	6000	1PG	C9-C8	-2.16	1.37	1.48
4	A	4000	MTA	O4'-C1'	2.07	1.43	1.41
4	C	4002	MTA	C5-C4	2.23	1.45	1.40
4	A	4000	MTA	C5-C4	3.25	1.47	1.40
4	B	4001	MTA	C5-C4	3.60	1.48	1.40

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	4000	MTA	N3-C2-N1	-8.07	122.71	128.89
4	C	4002	MTA	N3-C2-N1	-7.90	122.85	128.89
4	B	4001	MTA	N3-C2-N1	-6.92	123.60	128.89
4	A	4000	MTA	C4'-O4'-C1'	-3.90	105.44	109.72
4	C	4002	MTA	C4-C5-N7	-3.10	106.62	109.48
4	B	4001	MTA	C4-C5-N7	-2.62	107.07	109.48
5	A	6000	1PG	O5-C9-C8	-2.55	99.02	110.36
4	B	4001	MTA	C1'-N9-C4	-2.37	123.36	126.94
4	A	4000	MTA	O3'-C3'-C4'	-2.35	104.00	111.05
4	C	4002	MTA	C4'-O4'-C1'	-2.27	107.22	109.72
4	A	4000	MTA	C4-C5-N7	-2.12	107.53	109.48
5	A	6000	1PG	C8-O4-C7	-2.04	104.56	113.31
4	C	4002	MTA	O4'-C1'-N9	2.13	112.55	108.10
4	C	4002	MTA	C2-N1-C6	2.19	122.68	118.77
4	A	4000	MTA	C2-N1-C6	2.35	122.96	118.77
4	C	4002	MTA	CS-S5'-C5'	4.89	109.15	101.21
4	B	4001	MTA	CS-S5'-C5'	7.17	112.87	101.21
4	A	4000	MTA	CS-S5'-C5'	10.29	117.94	101.21

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	4000	MTA	1	0
5	A	6000	1PG	3	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 ⓘ

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	272/283 (96%)	-0.07	8 (2%)	55 56	21, 30, 45, 58	14 (5%)
1	B	270/283 (95%)	-0.12	6 (2%)	65 66	16, 26, 42, 57	9 (3%)
1	C	281/283 (99%)	-0.30	3 (1%)	82 83	15, 23, 35, 48	3 (1%)
All	All	823/849 (96%)	-0.17	17 (2%)	67 67	15, 27, 42, 58	26 (3%)

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	161	ASP	3.6
1	A	170	PHE	3.5
1	A	156	ILE	2.8
1	B	172	GLN	2.5
1	A	233	ILE	2.5
1	B	11	ILE	2.5
1	A	283	ILE	2.4
1	B	240	ASP	2.4
1	C	147	ASN	2.3
1	A	251	GLU	2.2
1	A	282	ASN	2.2
1	B	173	ASN	2.1
1	A	239	THR	2.1
1	A	281	GLU	2.1
1	B	281	GLU	2.0
1	C	64	TYR	2.0
1	C	231	ILE	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 [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(\AA^2)	Q<0.9
5	1PG	C	6001	17/17	0.82	0.25	5.18	31,35,37,38	0
4	MTA	A	4000	20/20	0.94	0.11	0.13	31,34,51,55	0
2	SO4	B	5000	5/5	0.98	0.12	0.12	38,38,42,42	0
4	MTA	B	4001	20/20	0.94	0.10	-0.58	28,35,50,51	0
4	MTA	C	4002	20/20	0.97	0.07	-1.14	22,24,29,34	0
3	UNX	A	3009	1/1	0.89	0.20	-	40,40,40,40	0
3	UNX	B	3013	1/1	0.89	0.34	-	35,35,35,35	0
3	UNX	A	3010	1/1	0.80	0.15	-	31,31,31,31	0
3	UNX	B	3012	1/1	0.62	0.36	-	37,37,37,37	0
3	UNX	B	3006	1/1	0.90	0.26	-	27,27,27,27	0
3	UNX	C	3001	1/1	0.78	0.35	-	37,37,37,37	0
3	UNX	B	3003	1/1	0.91	0.15	-	38,38,38,38	0
3	UNX	C	3007	1/1	0.91	0.20	-	28,28,28,28	0
3	UNX	C	3005	1/1	0.97	0.21	-	21,21,21,21	0
3	UNX	A	3002	1/1	0.92	0.24	-	44,44,44,44	0
5	1PG	A	6000	11/17	0.95	0.11	-	38,40,41,41	11
3	UNX	A	3008	1/1	0.92	0.16	-	22,22,22,22	0
3	UNX	A	3004	1/1	0.96	0.29	-	39,39,39,39	0
3	UNX	B	3011	1/1	0.84	0.33	-	26,26,26,26	0

6.5 Other polymers [i](#)

There are no such residues in this entry.