



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 09:54 PM UTC

PDB ID : 7XIG / pdb_00007xig
Title : Crystal structure of the aminopropyltransferase, SpeE from hyperthermophilic crenarchaeon, Pyrobaculum caldifontis in complex with 5'-methylthioadenosine (MTA) and spermine
Authors : Mizohata, E.; Yasuda, Y.
Deposited on : 2022-04-13
Resolution : 2.25 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

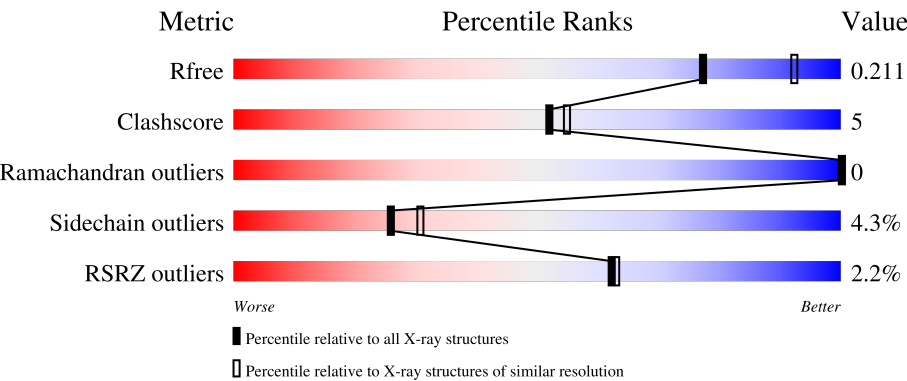
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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}	180053	1898 (2.26-2.26)
Clashscore	190562	2005 (2.26-2.26)
Ramachandran outliers	187476	1965 (2.26-2.26)
Sidechain outliers	187428	1966 (2.26-2.26)
RSRZ outliers	180081	1898 (2.26-2.26)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	309	% 81%11%7%
1	B	309	2% 80%13%7%
1	C	309	% 82%11%7%
1	D	309	3% 79%13%7%

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Mol	Chain	Length	Quality of chain
1	E	309	 % 80% 11% • 7%
1	F	309	 4% 81% 10% • 7%
1	G	309	 2% 83% 9% • 7%
1	H	309	 2% 80% 12% • 7%

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	SPM	H	302	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 19343 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 Polyamine aminopropyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	286	Total	C	N	O	S	0	0	0
			2276	1460	385	424	7			
1	B	288	Total	C	N	O	S	0	0	0
			2296	1472	391	426	7			
1	C	286	Total	C	N	O	S	0	0	0
			2276	1460	385	424	7			
1	D	288	Total	C	N	O	S	0	0	0
			2296	1472	391	426	7			
1	E	286	Total	C	N	O	S	0	1	0
			2284	1465	388	424	7			
1	F	286	Total	C	N	O	S	0	0	0
			2276	1460	385	424	7			
1	G	286	Total	C	N	O	S	0	0	0
			2276	1460	385	424	7			
1	H	288	Total	C	N	O	S	0	0	0
			2296	1472	391	426	7			

There are 160 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP A3MU81
A	-18	GLY	-	expression tag	UNP A3MU81
A	-17	SER	-	expression tag	UNP A3MU81
A	-16	SER	-	expression tag	UNP A3MU81
A	-15	HIS	-	expression tag	UNP A3MU81
A	-14	HIS	-	expression tag	UNP A3MU81
A	-13	HIS	-	expression tag	UNP A3MU81
A	-12	HIS	-	expression tag	UNP A3MU81
A	-11	HIS	-	expression tag	UNP A3MU81
A	-10	HIS	-	expression tag	UNP A3MU81
A	-9	SER	-	expression tag	UNP A3MU81
A	-8	SER	-	expression tag	UNP A3MU81
A	-7	GLY	-	expression tag	UNP A3MU81

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-6	LEU	-	expression tag	UNP A3MU81
A	-5	VAL	-	expression tag	UNP A3MU81
A	-4	PRO	-	expression tag	UNP A3MU81
A	-3	ARG	-	expression tag	UNP A3MU81
A	-2	GLY	-	expression tag	UNP A3MU81
A	-1	SER	-	expression tag	UNP A3MU81
A	0	HIS	-	expression tag	UNP A3MU81
B	-19	MET	-	initiating methionine	UNP A3MU81
B	-18	GLY	-	expression tag	UNP A3MU81
B	-17	SER	-	expression tag	UNP A3MU81
B	-16	SER	-	expression tag	UNP A3MU81
B	-15	HIS	-	expression tag	UNP A3MU81
B	-14	HIS	-	expression tag	UNP A3MU81
B	-13	HIS	-	expression tag	UNP A3MU81
B	-12	HIS	-	expression tag	UNP A3MU81
B	-11	HIS	-	expression tag	UNP A3MU81
B	-10	HIS	-	expression tag	UNP A3MU81
B	-9	SER	-	expression tag	UNP A3MU81
B	-8	SER	-	expression tag	UNP A3MU81
B	-7	GLY	-	expression tag	UNP A3MU81
B	-6	LEU	-	expression tag	UNP A3MU81
B	-5	VAL	-	expression tag	UNP A3MU81
B	-4	PRO	-	expression tag	UNP A3MU81
B	-3	ARG	-	expression tag	UNP A3MU81
B	-2	GLY	-	expression tag	UNP A3MU81
B	-1	SER	-	expression tag	UNP A3MU81
B	0	HIS	-	expression tag	UNP A3MU81
C	-19	MET	-	initiating methionine	UNP A3MU81
C	-18	GLY	-	expression tag	UNP A3MU81
C	-17	SER	-	expression tag	UNP A3MU81
C	-16	SER	-	expression tag	UNP A3MU81
C	-15	HIS	-	expression tag	UNP A3MU81
C	-14	HIS	-	expression tag	UNP A3MU81
C	-13	HIS	-	expression tag	UNP A3MU81
C	-12	HIS	-	expression tag	UNP A3MU81
C	-11	HIS	-	expression tag	UNP A3MU81
C	-10	HIS	-	expression tag	UNP A3MU81
C	-9	SER	-	expression tag	UNP A3MU81
C	-8	SER	-	expression tag	UNP A3MU81
C	-7	GLY	-	expression tag	UNP A3MU81
C	-6	LEU	-	expression tag	UNP A3MU81
C	-5	VAL	-	expression tag	UNP A3MU81

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-4	PRO	-	expression tag	UNP A3MU81
C	-3	ARG	-	expression tag	UNP A3MU81
C	-2	GLY	-	expression tag	UNP A3MU81
C	-1	SER	-	expression tag	UNP A3MU81
C	0	HIS	-	expression tag	UNP A3MU81
D	-19	MET	-	initiating methionine	UNP A3MU81
D	-18	GLY	-	expression tag	UNP A3MU81
D	-17	SER	-	expression tag	UNP A3MU81
D	-16	SER	-	expression tag	UNP A3MU81
D	-15	HIS	-	expression tag	UNP A3MU81
D	-14	HIS	-	expression tag	UNP A3MU81
D	-13	HIS	-	expression tag	UNP A3MU81
D	-12	HIS	-	expression tag	UNP A3MU81
D	-11	HIS	-	expression tag	UNP A3MU81
D	-10	HIS	-	expression tag	UNP A3MU81
D	-9	SER	-	expression tag	UNP A3MU81
D	-8	SER	-	expression tag	UNP A3MU81
D	-7	GLY	-	expression tag	UNP A3MU81
D	-6	LEU	-	expression tag	UNP A3MU81
D	-5	VAL	-	expression tag	UNP A3MU81
D	-4	PRO	-	expression tag	UNP A3MU81
D	-3	ARG	-	expression tag	UNP A3MU81
D	-2	GLY	-	expression tag	UNP A3MU81
D	-1	SER	-	expression tag	UNP A3MU81
D	0	HIS	-	expression tag	UNP A3MU81
E	-19	MET	-	initiating methionine	UNP A3MU81
E	-18	GLY	-	expression tag	UNP A3MU81
E	-17	SER	-	expression tag	UNP A3MU81
E	-16	SER	-	expression tag	UNP A3MU81
E	-15	HIS	-	expression tag	UNP A3MU81
E	-14	HIS	-	expression tag	UNP A3MU81
E	-13	HIS	-	expression tag	UNP A3MU81
E	-12	HIS	-	expression tag	UNP A3MU81
E	-11	HIS	-	expression tag	UNP A3MU81
E	-10	HIS	-	expression tag	UNP A3MU81
E	-9	SER	-	expression tag	UNP A3MU81
E	-8	SER	-	expression tag	UNP A3MU81
E	-7	GLY	-	expression tag	UNP A3MU81
E	-6	LEU	-	expression tag	UNP A3MU81
E	-5	VAL	-	expression tag	UNP A3MU81
E	-4	PRO	-	expression tag	UNP A3MU81
E	-3	ARG	-	expression tag	UNP A3MU81

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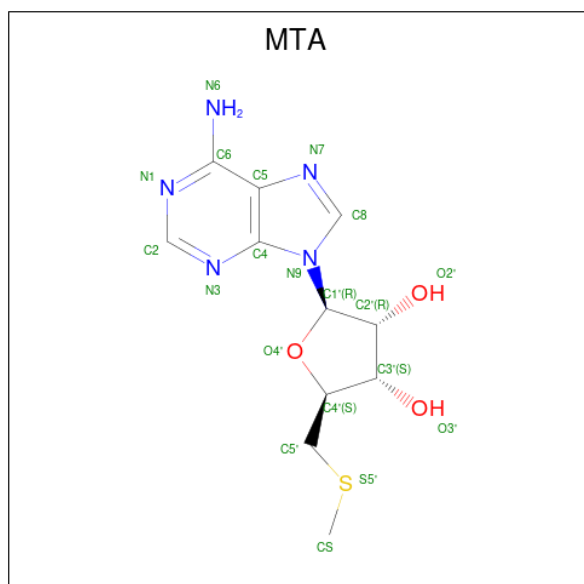
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E	-2	GLY	-	expression tag	UNP A3MU81
E	-1	SER	-	expression tag	UNP A3MU81
E	0	HIS	-	expression tag	UNP A3MU81
F	-19	MET	-	initiating methionine	UNP A3MU81
F	-18	GLY	-	expression tag	UNP A3MU81
F	-17	SER	-	expression tag	UNP A3MU81
F	-16	SER	-	expression tag	UNP A3MU81
F	-15	HIS	-	expression tag	UNP A3MU81
F	-14	HIS	-	expression tag	UNP A3MU81
F	-13	HIS	-	expression tag	UNP A3MU81
F	-12	HIS	-	expression tag	UNP A3MU81
F	-11	HIS	-	expression tag	UNP A3MU81
F	-10	HIS	-	expression tag	UNP A3MU81
F	-9	SER	-	expression tag	UNP A3MU81
F	-8	SER	-	expression tag	UNP A3MU81
F	-7	GLY	-	expression tag	UNP A3MU81
F	-6	LEU	-	expression tag	UNP A3MU81
F	-5	VAL	-	expression tag	UNP A3MU81
F	-4	PRO	-	expression tag	UNP A3MU81
F	-3	ARG	-	expression tag	UNP A3MU81
F	-2	GLY	-	expression tag	UNP A3MU81
F	-1	SER	-	expression tag	UNP A3MU81
F	0	HIS	-	expression tag	UNP A3MU81
G	-19	MET	-	initiating methionine	UNP A3MU81
G	-18	GLY	-	expression tag	UNP A3MU81
G	-17	SER	-	expression tag	UNP A3MU81
G	-16	SER	-	expression tag	UNP A3MU81
G	-15	HIS	-	expression tag	UNP A3MU81
G	-14	HIS	-	expression tag	UNP A3MU81
G	-13	HIS	-	expression tag	UNP A3MU81
G	-12	HIS	-	expression tag	UNP A3MU81
G	-11	HIS	-	expression tag	UNP A3MU81
G	-10	HIS	-	expression tag	UNP A3MU81
G	-9	SER	-	expression tag	UNP A3MU81
G	-8	SER	-	expression tag	UNP A3MU81
G	-7	GLY	-	expression tag	UNP A3MU81
G	-6	LEU	-	expression tag	UNP A3MU81
G	-5	VAL	-	expression tag	UNP A3MU81
G	-4	PRO	-	expression tag	UNP A3MU81
G	-3	ARG	-	expression tag	UNP A3MU81
G	-2	GLY	-	expression tag	UNP A3MU81
G	-1	SER	-	expression tag	UNP A3MU81

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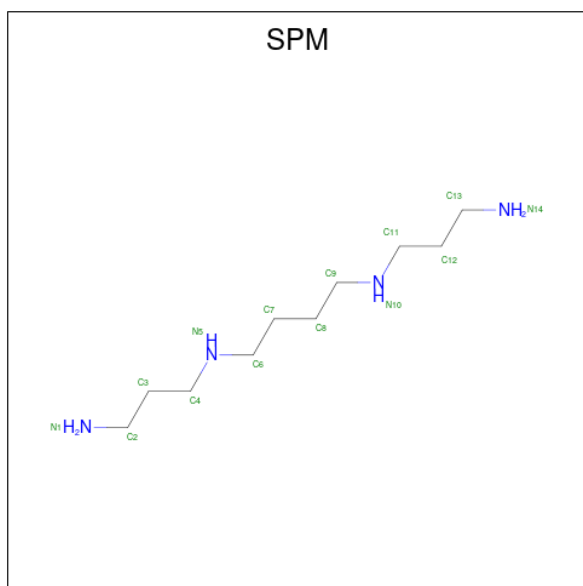
Chain	Residue	Modelled	Actual	Comment	Reference
G	0	HIS	-	expression tag	UNP A3MU81
H	-19	MET	-	initiating methionine	UNP A3MU81
H	-18	GLY	-	expression tag	UNP A3MU81
H	-17	SER	-	expression tag	UNP A3MU81
H	-16	SER	-	expression tag	UNP A3MU81
H	-15	HIS	-	expression tag	UNP A3MU81
H	-14	HIS	-	expression tag	UNP A3MU81
H	-13	HIS	-	expression tag	UNP A3MU81
H	-12	HIS	-	expression tag	UNP A3MU81
H	-11	HIS	-	expression tag	UNP A3MU81
H	-10	HIS	-	expression tag	UNP A3MU81
H	-9	SER	-	expression tag	UNP A3MU81
H	-8	SER	-	expression tag	UNP A3MU81
H	-7	GLY	-	expression tag	UNP A3MU81
H	-6	LEU	-	expression tag	UNP A3MU81
H	-5	VAL	-	expression tag	UNP A3MU81
H	-4	PRO	-	expression tag	UNP A3MU81
H	-3	ARG	-	expression tag	UNP A3MU81
H	-2	GLY	-	expression tag	UNP A3MU81
H	-1	SER	-	expression tag	UNP A3MU81
H	0	HIS	-	expression tag	UNP A3MU81

- Molecule 2 is 5'-DEOXY-5'-METHYLTHIOADENOSINE (CCD ID: MTA) (formula: $C_{11}H_{15}N_5O_3S$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			20	11	5	3	1		
2	B	1	Total	C	N	O	S	0	0
			20	11	5	3	1		
2	C	1	Total	C	N	O	S	0	0
			20	11	5	3	1		
2	D	1	Total	C	N	O	S	0	0
			20	11	5	3	1		
2	E	1	Total	C	N	O	S	0	0
			20	11	5	3	1		
2	F	1	Total	C	N	O	S	0	0
			20	11	5	3	1		
2	G	1	Total	C	N	O	S	0	0
			20	11	5	3	1		
2	H	1	Total	C	N	O	S	0	0
			20	11	5	3	1		

- Molecule 3 is SPERMINE (CCD ID: SPM) (formula: $C_{10}H_{26}N_4$) (labeled as "Ligand of Interest" by depositor).



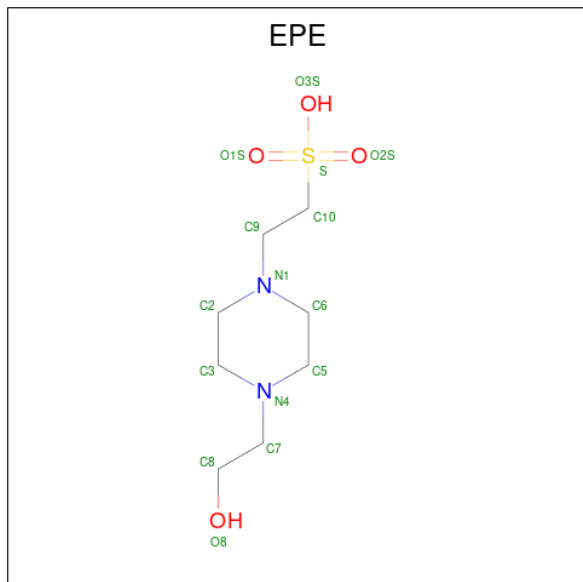
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	N	0	0
			14	10	4		
3	B	1	Total	C	N	0	0
			14	10	4		
3	C	1	Total	C	N	0	0
			14	10	4		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	D	1	Total	C	N	0	0
			14	10	4		
3	E	1	Total	C	N	0	0
			14	10	4		
3	F	1	Total	C	N	0	0
			14	10	4		
3	G	1	Total	C	N	0	0
			14	10	4		
3	H	1	Total	C	N	0	0
			14	10	4		

- Molecule 4 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (CCD ID: EPE) (formula: $C_8H_{18}N_2O_4S$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	S	0	0
			15	8	2	4	1		
4	G	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	114	Total	O	0	0
			114	114		
5	B	79	Total	O	0	0
			79	79		

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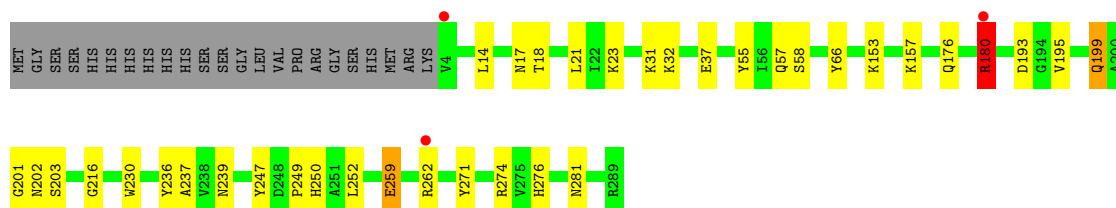
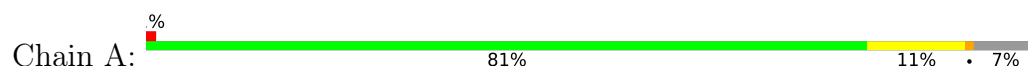
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	C	118	Total 118	O 118	0	0
5	D	68	Total 68	O 68	0	0
5	E	99	Total 99	O 99	0	0
5	F	76	Total 76	O 76	0	0
5	G	113	Total 113	O 113	0	0
5	H	98	Total 98	O 98	0	0

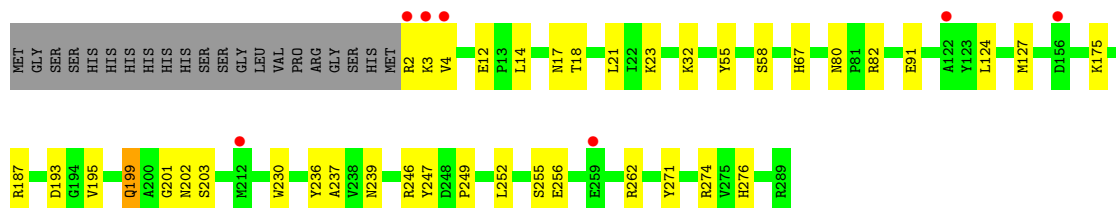
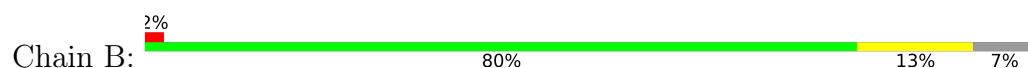
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

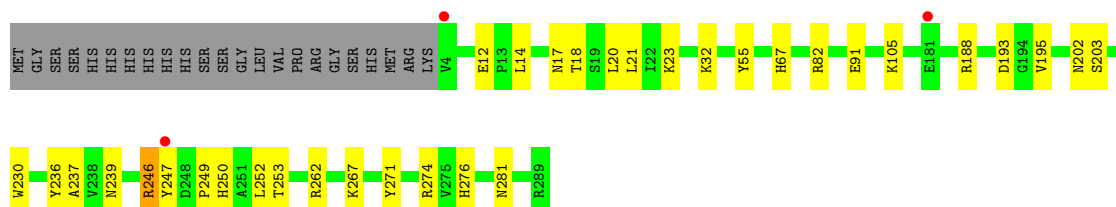
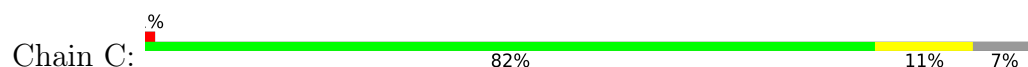
• Molecule 1: Polyamine aminopropyltransferase



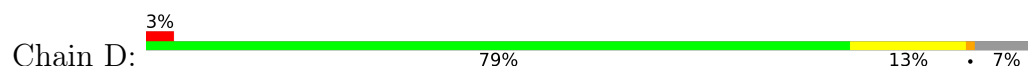
• Molecule 1: Polyamine aminopropyltransferase

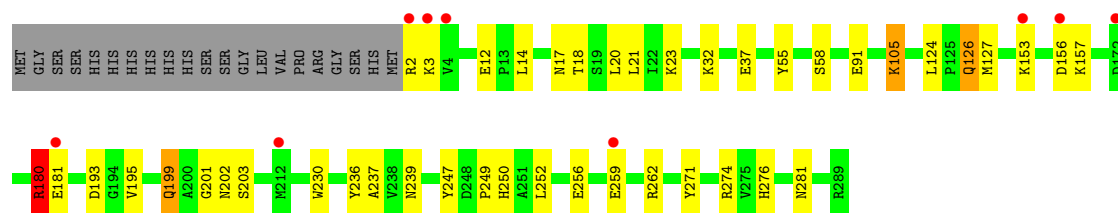


• Molecule 1: Polyamine aminopropyltransferase

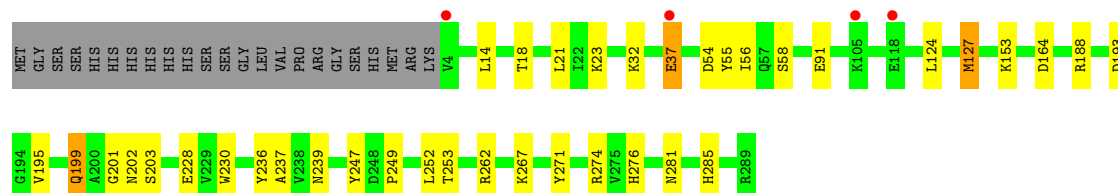
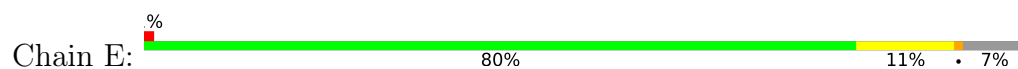


• Molecule 1: Polyamine aminopropyltransferase

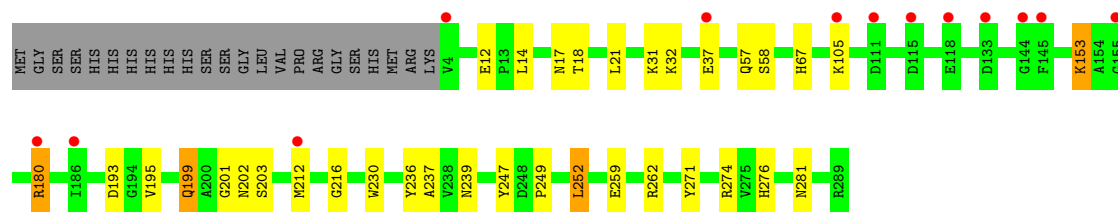
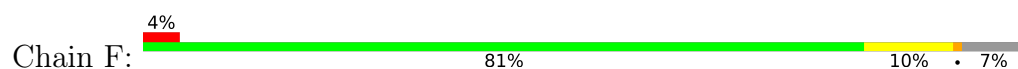




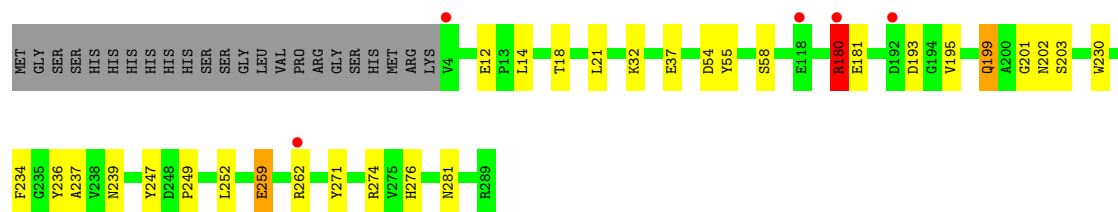
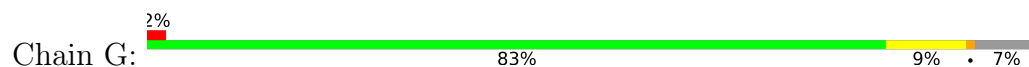
• Molecule 1: Polyamine aminopropyltransferase



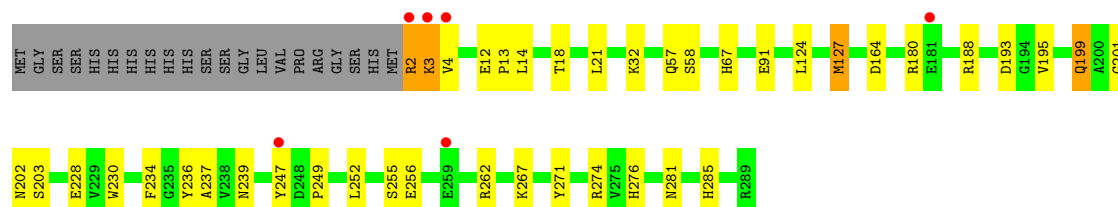
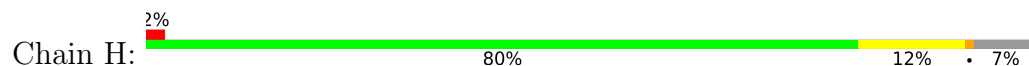
• Molecule 1: Polyamine aminopropyltransferase



• Molecule 1: Polyamine aminopropyltransferase



• Molecule 1: Polyamine aminopropyltransferase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	77.83Å 80.39Å 97.84Å 72.46° 89.13° 88.08°	Depositor
Resolution (Å)	46.69 – 2.25 46.69 – 2.25	Depositor EDS
% Data completeness (in resolution range)	98.9 (46.69-2.25) 98.9 (46.69-2.25)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.72 (at 2.24Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.173 , 0.207 0.176 , 0.211	Depositor DCC
R_{free} test set	5275 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	25.5	Xtriage
Anisotropy	0.043	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 34.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.006 for h,-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	19343	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 75.27 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.3061e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MTA, EPE, SPM

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	0.58	0/2329	0.98	3/3164 (0.1%)
1	B	0.57	0/2349	0.96	1/3189 (0.0%)
1	C	0.58	1/2329 (0.0%)	0.96	3/3164 (0.1%)
1	D	0.56	0/2349	0.97	3/3189 (0.1%)
1	E	0.58	0/2340	0.97	3/3178 (0.1%)
1	F	0.56	0/2329	0.97	0/3164
1	G	0.58	0/2329	0.95	1/3164 (0.0%)
1	H	0.56	0/2349	0.96	2/3189 (0.1%)
All	All	0.57	1/18703 (0.0%)	0.97	16/25401 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	67	HIS	CE1-NE2	5.11	1.37	1.32

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	180	ARG	CG-CD-NE	9.05	131.91	112.00
1	D	180	ARG	CB-CG-CD	8.03	129.78	111.30
1	A	274	ARG	CB-CG-CD	-7.58	93.87	111.30
1	D	256	GLU	CB-CG-CD	6.16	123.08	112.60
1	C	246	ARG	CB-CG-CD	6.11	125.35	111.30
1	E	37	GLU	CB-CG-CD	5.96	122.73	112.60
1	H	188	ARG	CB-CG-CD	5.94	124.97	111.30
1	A	259	GLU	CB-CG-CD	5.90	122.62	112.60
1	C	82	ARG	CB-CG-CD	5.75	124.52	111.30
1	C	188	ARG	CB-CG-CD	5.55	124.08	111.30
1	G	180	ARG	CG-CD-NE	5.43	123.94	112.00
1	D	2	ARG	CG-CD-NE	-5.42	100.07	112.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	246	ARG	CB-CG-CD	5.34	123.58	111.30
1	E	54	ASP	CA-CB-CG	5.33	117.93	112.60
1	E	188	ARG	CB-CG-CD	5.33	123.57	111.30
1	H	180	ARG	CG-CD-NE	-5.14	100.69	112.00

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	2276	0	2265	27	0
1	B	2296	0	2291	25	0
1	C	2276	0	2265	31	0
1	D	2296	0	2291	43	0
1	E	2284	0	2278	25	0
1	F	2276	0	2265	27	0
1	G	2276	0	2265	20	0
1	H	2296	0	2291	32	0
2	A	20	0	15	0	0
2	B	20	0	15	0	0
2	C	20	0	15	1	0
2	D	20	0	15	0	0
2	E	20	0	15	0	0
2	F	20	0	15	0	0
2	G	20	0	15	1	0
2	H	20	0	15	1	0
3	A	14	0	26	5	0
3	B	14	0	26	3	0
3	C	14	0	26	4	0
3	D	14	0	26	5	0
3	E	14	0	26	4	0
3	F	14	0	26	3	0
3	G	14	0	26	4	0
3	H	14	0	26	9	0
4	A	15	0	18	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	G	15	0	17	0	0
5	A	114	0	0	1	0
5	B	79	0	0	2	0
5	C	118	0	0	1	0
5	D	68	0	0	1	0
5	E	99	0	0	0	0
5	F	76	0	0	1	0
5	G	113	0	0	0	0
5	H	98	0	0	0	0
All	All	19343	0	18574	191	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 5.

All (191) 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:2:ARG:HD2	1:B:4:VAL:O	1.62	1.00
1:D:180:ARG:HG2	1:D:180:ARG:HH21	1.26	0.97
1:A:259:GLU:HB2	1:C:253:THR:HG21	1.46	0.95
1:D:126:GLN:HE21	1:D:126:GLN:HA	1.32	0.92
1:F:259:GLU:OE1	1:H:255:SER:HB2	1.68	0.92
1:F:259:GLU:OE1	1:H:255:SER:CB	2.19	0.91
1:C:20:LEU:CD2	1:D:20:LEU:CD2	2.51	0.89
1:D:180:ARG:HG2	1:D:180:ARG:NH2	1.86	0.86
1:A:176:GLN:HG2	5:A:461:HOH:O	1.76	0.84
1:D:180:ARG:HH21	1:D:180:ARG:CG	1.91	0.83
1:A:14:LEU:HD11	1:A:236:TYR:HD2	1.44	0.81
1:C:14:LEU:HD11	1:C:236:TYR:HD2	1.46	0.80
1:G:14:LEU:HD11	1:G:236:TYR:HD2	1.47	0.79
1:B:14:LEU:HD11	1:B:236:TYR:HD2	1.48	0.78
1:H:14:LEU:HD11	1:H:236:TYR:HD2	1.48	0.78
1:F:14:LEU:HD11	1:F:236:TYR:HD2	1.49	0.77
1:D:14:LEU:HD11	1:D:236:TYR:HD2	1.48	0.77
1:E:14:LEU:HD11	1:E:236:TYR:HD2	1.51	0.74
1:F:259:GLU:CD	1:H:255:SER:HB2	2.12	0.74
1:E:124:LEU:HD13	1:E:127:MET:HE2	1.70	0.73
1:D:124:LEU:HD13	1:D:127:MET:CE	2.19	0.72
1:H:124:LEU:HD13	1:H:127:MET:HE2	1.72	0.71
1:H:67:HIS:NE2	3:H:302:SPM:N1	2.38	0.71
1:F:259:GLU:OE1	1:H:255:SER:HB3	1.90	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:124:LEU:HD13	1:E:127:MET:CE	2.23	0.69
1:H:124:LEU:HD13	1:H:127:MET:CE	2.23	0.69
1:C:250:HIS:HD2	5:D:421:HOH:O	1.77	0.68
1:H:234:PHE:O	3:H:302:SPM:H132	1.94	0.68
1:B:124:LEU:HD13	1:B:127:MET:CE	2.24	0.67
1:E:55:TYR:HD1	3:E:302:SPM:H31	1.59	0.67
1:D:12:GLU:OE2	3:D:302:SPM:H131	1.97	0.65
1:D:124:LEU:HD13	1:D:127:MET:HE2	1.78	0.64
1:H:91:GLU:HG3	1:H:127:MET:HE3	1.80	0.64
1:C:91:GLU:OE2	3:C:302:SPM:H132	1.99	0.63
1:F:12:GLU:OE2	3:F:302:SPM:H131	1.98	0.62
1:E:91:GLU:HG3	1:E:127:MET:HE3	1.83	0.61
2:H:301:MTA:S5'	3:H:302:SPM:H22	2.41	0.61
1:B:80:ASN:ND2	1:B:82:ARG:HD2	2.16	0.60
1:C:281:ASN:OD1	1:D:274:ARG:HD2	2.02	0.60
5:C:495:HOH:O	1:D:250:HIS:HD2	1.83	0.60
1:H:164:ASP:OD1	3:H:302:SPM:H21	2.01	0.60
1:D:12:GLU:CD	3:D:302:SPM:H131	2.27	0.60
1:H:164:ASP:CG	3:H:302:SPM:H21	2.27	0.60
1:C:20:LEU:CD2	1:D:20:LEU:HD23	2.33	0.59
1:D:55:TYR:HD1	3:D:302:SPM:H122	1.66	0.58
1:E:202:ASN:HA	1:E:239:ASN:HD22	1.69	0.58
1:D:202:ASN:HA	1:D:239:ASN:HD22	1.69	0.57
1:C:20:LEU:HD23	1:D:20:LEU:HD22	1.86	0.57
1:H:202:ASN:HA	1:H:239:ASN:HD22	1.69	0.57
1:G:202:ASN:HA	1:G:239:ASN:HD22	1.69	0.57
1:A:180:ARG:HG3	1:A:216:GLY:HA2	1.87	0.56
1:A:259:GLU:HB2	1:C:253:THR:CG2	2.29	0.56
1:A:202:ASN:HA	1:A:239:ASN:HD22	1.70	0.56
1:C:20:LEU:HD23	1:D:20:LEU:CD2	2.33	0.56
1:C:202:ASN:HA	1:C:239:ASN:HD22	1.70	0.56
1:F:202:ASN:HA	1:F:239:ASN:HD22	1.71	0.56
1:B:202:ASN:HA	1:B:239:ASN:HD22	1.70	0.56
1:F:14:LEU:CD1	1:F:236:TYR:HD2	2.19	0.56
1:G:14:LEU:CD1	1:G:236:TYR:HD2	2.19	0.56
1:E:55:TYR:CD1	3:E:302:SPM:H31	2.41	0.55
1:E:281:ASN:OD1	1:F:274:ARG:HD2	2.05	0.55
1:D:14:LEU:CD1	1:D:236:TYR:HD2	2.19	0.55
1:D:126:GLN:HA	1:D:126:GLN:NE2	2.11	0.54
1:H:14:LEU:CD1	1:H:236:TYR:HD2	2.19	0.54
1:C:20:LEU:HD22	1:D:20:LEU:HD23	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:55:TYR:HB3	3:B:302:SPM:H111	1.88	0.54
1:C:14:LEU:CD1	1:C:236:TYR:HD2	2.19	0.54
1:B:124:LEU:HD13	1:B:127:MET:HE2	1.89	0.53
1:B:14:LEU:CD1	1:B:236:TYR:HD2	2.20	0.53
1:C:20:LEU:CD2	1:D:20:LEU:HD22	2.36	0.53
1:A:55:TYR:HD2	3:A:302:SPM:H21	1.73	0.53
1:G:12:GLU:OE2	3:G:302:SPM:H21	2.08	0.53
1:A:14:LEU:CD1	1:A:236:TYR:HD2	2.18	0.53
1:B:193:ASP:HB2	1:B:247:TYR:CE2	2.44	0.52
1:B:91:GLU:HB2	1:B:127:MET:HE3	1.91	0.52
1:A:281:ASN:OD1	1:B:274:ARG:HD2	2.10	0.52
1:B:12:GLU:OE2	3:B:302:SPM:H131	2.10	0.51
1:F:193:ASP:HB2	1:F:247:TYR:CE2	2.46	0.51
1:A:21:LEU:HD21	1:B:21:LEU:HD11	1.92	0.51
1:G:274:ARG:HD2	1:H:281:ASN:OD1	2.09	0.51
1:E:14:LEU:CD1	1:E:236:TYR:HD2	2.22	0.51
1:F:203:SER:H	1:F:239:ASN:ND2	2.09	0.51
1:H:193:ASP:HB2	1:H:247:TYR:CE2	2.46	0.51
1:C:274:ARG:HD2	1:D:281:ASN:OD1	2.11	0.51
1:D:203:SER:H	1:D:239:ASN:ND2	2.09	0.51
1:G:203:SER:H	1:G:239:ASN:ND2	2.09	0.50
1:D:91:GLU:HB2	1:D:127:MET:HE3	1.94	0.50
1:F:180:ARG:HD2	5:F:468:HOH:O	2.11	0.50
1:A:203:SER:H	1:A:239:ASN:ND2	2.09	0.50
1:C:193:ASP:HB2	1:C:247:TYR:CE2	2.46	0.50
1:C:203:SER:H	1:C:239:ASN:ND2	2.10	0.50
1:F:180:ARG:HG3	1:F:216:GLY:HA2	1.92	0.50
1:G:281:ASN:OD1	1:H:274:ARG:HD2	2.11	0.50
1:A:17:ASN:HD22	1:B:23:LYS:HD2	1.77	0.50
1:D:55:TYR:CD1	3:D:302:SPM:H122	2.46	0.50
1:D:193:ASP:HB2	1:D:247:TYR:CE2	2.47	0.49
1:E:249:PRO:O	1:E:276:HIS:HE1	1.94	0.49
1:A:193:ASP:HB2	1:A:247:TYR:CE2	2.46	0.49
1:E:203:SER:H	1:E:239:ASN:ND2	2.10	0.49
1:F:57:GLN:NE2	3:F:302:SPM:H42	2.27	0.49
1:E:199:GLN:NE2	1:E:201:GLY:H	2.11	0.49
1:G:54:ASP:OD2	1:H:3:LYS:HE2	2.12	0.49
1:H:203:SER:H	1:H:239:ASN:ND2	2.09	0.49
1:B:203:SER:H	1:B:239:ASN:ND2	2.10	0.49
1:D:126:GLN:HE21	1:D:126:GLN:CA	2.05	0.49
1:E:164:ASP:OD1	3:E:302:SPM:H112	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:199:GLN:NE2	1:F:201:GLY:H	2.11	0.48
1:G:193:ASP:HB2	1:G:247:TYR:CE2	2.48	0.48
1:C:249:PRO:O	1:C:276:HIS:HE1	1.95	0.48
1:C:55:TYR:HB3	3:C:302:SPM:H32	1.95	0.48
1:H:199:GLN:NE2	1:H:201:GLY:H	2.11	0.48
1:D:199:GLN:NE2	1:D:201:GLY:H	2.11	0.48
1:B:199:GLN:NE2	1:B:201:GLY:H	2.11	0.48
1:B:249:PRO:O	1:B:276:HIS:HE1	1.97	0.48
1:G:199:GLN:NE2	1:G:201:GLY:H	2.11	0.48
2:C:301:MTA:S5'	3:C:302:SPM:H112	2.53	0.48
1:A:249:PRO:O	1:A:276:HIS:HE1	1.96	0.48
1:B:67:HIS:NE2	3:B:302:SPM:N1	2.62	0.48
1:A:199:GLN:NE2	1:A:201:GLY:H	2.11	0.47
1:G:18:THR:HA	1:H:21:LEU:O	2.14	0.47
1:G:180:ARG:HB2	1:G:180:ARG:HH11	1.79	0.47
1:C:20:LEU:CD2	1:D:20:LEU:HD21	2.42	0.46
1:D:249:PRO:O	1:D:276:HIS:HE1	1.98	0.46
1:F:249:PRO:O	1:F:276:HIS:HE1	1.97	0.46
1:G:249:PRO:O	1:G:276:HIS:HE1	1.97	0.46
1:H:249:PRO:O	1:H:276:HIS:HE1	1.97	0.46
1:E:253:THR:HG21	1:G:259:GLU:HB2	1.98	0.46
1:D:55:TYR:HB3	3:D:302:SPM:H111	1.98	0.46
1:A:180:ARG:HH11	1:A:180:ARG:HB2	1.80	0.45
1:E:193:ASP:HB2	1:E:247:TYR:HE2	1.81	0.45
1:C:20:LEU:HD21	1:D:20:LEU:HD21	1.97	0.45
1:E:23:LYS:HD2	1:F:17:ASN:HD22	1.81	0.45
1:E:193:ASP:HB2	1:E:247:TYR:CE2	2.51	0.45
1:E:253:THR:HG21	1:G:259:GLU:CB	2.47	0.45
1:E:274:ARG:HD2	1:F:281:ASN:OD1	2.15	0.45
1:G:21:LEU:O	1:H:18:THR:HA	2.16	0.45
1:C:230:TRP:HA	1:C:237:ALA:HA	1.99	0.45
1:G:271:TYR:OH	1:G:276:HIS:CD2	2.70	0.45
1:A:18:THR:HA	1:B:21:LEU:O	2.17	0.44
1:A:57:GLN:HE22	3:A:302:SPM:H91	1.81	0.44
1:C:17:ASN:HD22	1:D:23:LYS:HD2	1.82	0.44
1:C:20:LEU:HD21	1:D:20:LEU:CD2	2.40	0.44
1:C:21:LEU:O	1:D:18:THR:HA	2.17	0.44
1:H:57:GLN:NE2	3:H:302:SPM:H42	2.32	0.44
1:D:271:TYR:OH	1:D:276:HIS:CD2	2.71	0.44
1:A:21:LEU:O	1:B:18:THR:HA	2.17	0.44
1:A:271:TYR:OH	1:A:276:HIS:CD2	2.71	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:255:SER:HB3	1:D:259:GLU:OE1	2.18	0.44
1:E:230:TRP:HA	1:E:237:ALA:HA	2.00	0.44
1:F:230:TRP:HA	1:F:237:ALA:HA	1.99	0.44
1:B:271:TYR:OH	1:B:276:HIS:CD2	2.71	0.43
1:C:271:TYR:OH	1:C:276:HIS:CD2	2.71	0.43
1:D:230:TRP:HA	1:D:237:ALA:HA	2.00	0.43
1:F:67:HIS:NE2	3:F:302:SPM:N1	2.66	0.43
1:A:23:LYS:HD2	1:B:17:ASN:HD22	1.84	0.43
1:H:271:TYR:OH	1:H:276:HIS:CD2	2.72	0.43
1:H:12:GLU:OE2	3:H:302:SPM:H131	2.19	0.43
1:A:57:GLN:NE2	3:A:302:SPM:H91	2.34	0.43
1:F:252:LEU:HD12	1:F:252:LEU:HA	1.88	0.43
1:F:271:TYR:OH	1:F:276:HIS:CD2	2.71	0.43
1:G:230:TRP:HA	1:G:237:ALA:HA	1.99	0.43
1:C:23:LYS:HD2	1:D:17:ASN:HD22	1.84	0.42
1:E:21:LEU:O	1:F:18:THR:HA	2.20	0.42
1:A:14:LEU:HD11	1:A:236:TYR:CD2	2.36	0.42
1:A:230:TRP:HA	1:A:237:ALA:HA	1.99	0.42
2:G:301:MTA:S5'	3:G:302:SPM:H122	2.59	0.42
1:B:187:ARG:HD3	5:B:423:HOH:O	2.18	0.42
1:E:271:TYR:OH	1:E:276:HIS:CD2	2.72	0.42
1:C:12:GLU:OE2	3:C:302:SPM:H31	2.20	0.42
1:A:250:HIS:HD2	5:B:460:HOH:O	2.01	0.42
1:C:21:LEU:HD21	1:D:21:LEU:HD11	2.02	0.42
1:H:2:ARG:NE	1:H:4:VAL:O	2.39	0.42
1:F:153:LYS:HD3	1:F:153:LYS:HA	1.81	0.41
1:G:234:PHE:CE2	3:G:302:SPM:H72	2.55	0.41
1:H:230:TRP:HA	1:H:237:ALA:HA	2.00	0.41
1:A:66:TYR:CE2	3:A:302:SPM:H131	2.55	0.41
1:A:55:TYR:HD2	3:A:302:SPM:C2	2.32	0.41
1:C:271:TYR:OH	1:C:276:HIS:HD2	2.03	0.41
1:H:228:GLU:OE2	1:H:285:HIS:ND1	2.52	0.41
1:B:230:TRP:HA	1:B:237:ALA:HA	2.01	0.41
1:F:180:ARG:H	1:F:180:ARG:NH1	2.18	0.41
1:F:259:GLU:OE2	1:H:255:SER:HB2	2.20	0.41
1:C:18:THR:HA	1:D:21:LEU:O	2.21	0.41
1:E:228:GLU:OE2	1:E:285:HIS:ND1	2.52	0.41
1:D:105:LYS:HA	1:D:105:LYS:HD2	1.72	0.41
1:E:18:THR:HA	1:F:21:LEU:O	2.21	0.41
1:G:55:TYR:HD1	3:G:302:SPM:H22	1.86	0.41
1:H:13:PRO:O	3:H:302:SPM:N14	2.54	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:164:ASP:OD2	3:H:302:SPM:H21	2.21	0.40
1:D:271:TYR:OH	1:D:276:HIS:HD2	2.04	0.40
1:E:56:ILE:O	3:E:302:SPM:H42	2.21	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	284/309 (92%)	276 (97%)	8 (3%)	0	100	100
1	B	286/309 (93%)	279 (98%)	7 (2%)	0	100	100
1	C	284/309 (92%)	276 (97%)	8 (3%)	0	100	100
1	D	286/309 (93%)	278 (97%)	8 (3%)	0	100	100
1	E	285/309 (92%)	277 (97%)	8 (3%)	0	100	100
1	F	284/309 (92%)	276 (97%)	8 (3%)	0	100	100
1	G	284/309 (92%)	276 (97%)	8 (3%)	0	100	100
1	H	286/309 (93%)	278 (97%)	8 (3%)	0	100	100
All	All	2279/2472 (92%)	2216 (97%)	63 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	245/265 (92%)	234 (96%)	11 (4%)	24	29
1	B	247/265 (93%)	238 (96%)	9 (4%)	31	39
1	C	245/265 (92%)	238 (97%)	7 (3%)	37	47
1	D	247/265 (93%)	232 (94%)	15 (6%)	17	17
1	E	246/265 (93%)	236 (96%)	10 (4%)	27	33
1	F	245/265 (92%)	233 (95%)	12 (5%)	22	24
1	G	245/265 (92%)	235 (96%)	10 (4%)	27	33
1	H	247/265 (93%)	236 (96%)	11 (4%)	24	29
All	All	1967/2120 (93%)	1882 (96%)	85 (4%)	26	31

All (85) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	31	LYS
1	A	32	LYS
1	A	37	GLU
1	A	58	SER
1	A	153	LYS
1	A	157	LYS
1	A	180	ARG
1	A	195	VAL
1	A	199	GLN
1	A	252	LEU
1	A	262	ARG
1	B	3	LYS
1	B	32	LYS
1	B	58	SER
1	B	175	LYS
1	B	195	VAL
1	B	199	GLN
1	B	252	LEU
1	B	256	GLU
1	B	262	ARG
1	C	32	LYS
1	C	105	LYS
1	C	195	VAL
1	C	246	ARG
1	C	252	LEU
1	C	262	ARG
1	C	267	LYS

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Mol	Chain	Res	Type
1	D	3	LYS
1	D	32	LYS
1	D	37	GLU
1	D	58	SER
1	D	105	LYS
1	D	126	GLN
1	D	153	LYS
1	D	156	ASP
1	D	157	LYS
1	D	180	ARG
1	D	181	GLU
1	D	195	VAL
1	D	199	GLN
1	D	252	LEU
1	D	262	ARG
1	E	32	LYS
1	E	37	GLU
1	E	58	SER
1	E	127	MET
1	E	153	LYS
1	E	195	VAL
1	E	199	GLN
1	E	252	LEU
1	E	262	ARG
1	E	267	LYS
1	F	31	LYS
1	F	32	LYS
1	F	37	GLU
1	F	58	SER
1	F	105	LYS
1	F	153	LYS
1	F	180	ARG
1	F	195	VAL
1	F	199	GLN
1	F	212	MET
1	F	252	LEU
1	F	262	ARG
1	G	32	LYS
1	G	37	GLU
1	G	58	SER
1	G	180	ARG
1	G	181	GLU

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Mol	Chain	Res	Type
1	G	195	VAL
1	G	199	GLN
1	G	252	LEU
1	G	259	GLU
1	G	262	ARG
1	H	2	ARG
1	H	3	LYS
1	H	32	LYS
1	H	58	SER
1	H	127	MET
1	H	195	VAL
1	H	199	GLN
1	H	252	LEU
1	H	256	GLU
1	H	262	ARG
1	H	267	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (67) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	17	ASN
1	A	57	GLN
1	A	129	GLN
1	A	142	GLN
1	A	191	ASN
1	A	199	GLN
1	A	239	ASN
1	A	276	HIS
1	B	17	ASN
1	B	57	GLN
1	B	80	ASN
1	B	129	GLN
1	B	142	GLN
1	B	191	ASN
1	B	199	GLN
1	B	239	ASN
1	B	276	HIS
1	C	17	ASN
1	C	57	GLN
1	C	129	GLN
1	C	142	GLN
1	C	191	ASN

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Mol	Chain	Res	Type
1	C	239	ASN
1	C	276	HIS
1	D	17	ASN
1	D	36	GLN
1	D	126	GLN
1	D	142	GLN
1	D	191	ASN
1	D	199	GLN
1	D	239	ASN
1	D	276	HIS
1	E	17	ASN
1	E	129	GLN
1	E	142	GLN
1	E	191	ASN
1	E	199	GLN
1	E	239	ASN
1	E	276	HIS
1	F	17	ASN
1	F	36	GLN
1	F	57	GLN
1	F	80	ASN
1	F	129	GLN
1	F	142	GLN
1	F	191	ASN
1	F	199	GLN
1	F	239	ASN
1	F	276	HIS
1	G	17	ASN
1	G	36	GLN
1	G	57	GLN
1	G	129	GLN
1	G	142	GLN
1	G	191	ASN
1	G	199	GLN
1	G	239	ASN
1	G	276	HIS
1	H	17	ASN
1	H	36	GLN
1	H	57	GLN
1	H	129	GLN
1	H	142	GLN
1	H	191	ASN

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Mol	Chain	Res	Type
1	H	199	GLN
1	H	239	ASN
1	H	276	HIS

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

18 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	SPM	B	302	-	13,13,13	0.36	0	12,12,12	0.50	0
3	SPM	A	302	-	13,13,13	0.35	0	12,12,12	0.77	1 (8%)
3	SPM	C	302	-	13,13,13	0.37	0	12,12,12	0.68	0
2	MTA	C	301	-	22,22,22	1.68	4 (18%)	32,32,32	1.99	7 (21%)
2	MTA	D	301	-	22,22,22	1.36	3 (13%)	32,32,32	2.27	13 (40%)
3	SPM	G	302	-	13,13,13	0.45	0	12,12,12	0.73	0
2	MTA	B	301	-	22,22,22	1.58	3 (13%)	32,32,32	1.93	10 (31%)
2	MTA	G	301	-	22,22,22	1.45	3 (13%)	32,32,32	2.13	7 (21%)
3	SPM	F	302	-	13,13,13	0.29	0	12,12,12	0.52	0
3	SPM	H	302	-	13,13,13	0.29	0	12,12,12	0.62	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	MTA	A	301	-	22,22,22	1.54	4 (18%)	32,32,32	2.07	11 (34%)
3	SPM	E	302	-	13,13,13	0.32	0	12,12,12	0.68	0
4	EPE	A	303	-	15,15,15	2.24	1 (6%)	19,20,20	1.46	2 (10%)
3	SPM	D	302	-	13,13,13	0.24	0	12,12,12	0.46	0
2	MTA	H	301	-	22,22,22	1.59	4 (18%)	32,32,32	2.05	12 (37%)
2	MTA	E	301	-	22,22,22	1.79	4 (18%)	32,32,32	2.07	9 (28%)
4	EPE	G	303	-	15,15,15	1.97	1 (6%)	19,20,20	1.33	2 (10%)
2	MTA	F	301	-	22,22,22	1.54	6 (27%)	32,32,32	2.19	12 (37%)

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	SPM	B	302	-	-	4/11/11/11	-
3	SPM	A	302	-	-	8/11/11/11	-
3	SPM	C	302	-	-	7/11/11/11	-
2	MTA	C	301	-	-	2/7/23/23	0/3/3/3
2	MTA	D	301	-	-	2/7/23/23	0/3/3/3
3	SPM	G	302	-	-	6/11/11/11	-
2	MTA	B	301	-	-	0/7/23/23	0/3/3/3
2	MTA	G	301	-	-	1/7/23/23	0/3/3/3
3	SPM	F	302	-	-	9/11/11/11	-
3	SPM	H	302	-	-	5/11/11/11	-
2	MTA	A	301	-	-	2/7/23/23	0/3/3/3
3	SPM	E	302	-	-	8/11/11/11	-
4	EPE	A	303	-	-	3/9/19/19	0/1/1/1
3	SPM	D	302	-	-	4/11/11/11	-
2	MTA	H	301	-	-	2/7/23/23	0/3/3/3
2	MTA	E	301	-	-	2/7/23/23	0/3/3/3
4	EPE	G	303	-	-	3/9/19/19	0/1/1/1
2	MTA	F	301	-	-	2/7/23/23	0/3/3/3

All (33) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	303	EPE	C10-S	-8.32	1.65	1.77
4	G	303	EPE	C10-S	-7.15	1.67	1.77
2	E	301	MTA	C5-C4	5.47	1.48	1.39
2	C	301	MTA	C5-C4	4.80	1.47	1.39
2	H	301	MTA	C5-C4	4.76	1.47	1.39
2	B	301	MTA	C5-C4	4.45	1.47	1.39
2	A	301	MTA	C5-C4	4.16	1.46	1.39
2	F	301	MTA	C5-C4	4.05	1.46	1.39
2	G	301	MTA	C5-C4	3.87	1.46	1.39
2	E	301	MTA	C8-N7	3.73	1.38	1.31
2	A	301	MTA	C8-N7	3.69	1.38	1.31
2	B	301	MTA	C4-N9	-3.46	1.30	1.37
2	C	301	MTA	C5-N7	-3.37	1.32	1.39
2	D	301	MTA	C4-N9	-3.08	1.31	1.37
2	H	301	MTA	C5-C6	2.95	1.49	1.41
2	H	301	MTA	C4-N9	-2.93	1.31	1.37
2	A	301	MTA	C5-N7	-2.91	1.33	1.39
2	D	301	MTA	C5-C4	2.86	1.44	1.39
2	C	301	MTA	C5-C6	2.78	1.48	1.41
2	D	301	MTA	C5-C6	2.74	1.48	1.41
2	B	301	MTA	C5-C6	2.72	1.48	1.41
2	F	301	MTA	C8-N7	2.64	1.36	1.31
2	G	301	MTA	C5-C6	2.63	1.48	1.41
2	G	301	MTA	C4-N9	-2.52	1.32	1.37
2	E	301	MTA	C5-N7	-2.46	1.34	1.39
2	F	301	MTA	C4-N9	-2.46	1.32	1.37
2	H	301	MTA	C8-N7	2.26	1.36	1.31
2	C	301	MTA	C4-N9	-2.24	1.33	1.37
2	E	301	MTA	C5'-S5'	-2.24	1.75	1.81
2	F	301	MTA	C5-C6	2.21	1.47	1.41
2	F	301	MTA	O4'-C4'	-2.17	1.40	1.45
2	F	301	MTA	C5-N7	-2.13	1.35	1.39
2	A	301	MTA	C5-C6	2.13	1.46	1.41

All (86) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	301	MTA	C5-C4-N3	-5.52	119.11	126.72
2	D	301	MTA	C4-N9-C8	5.28	111.29	105.74
2	A	301	MTA	C5-C4-N3	-5.26	119.47	126.72
2	F	301	MTA	N3-C2-N1	-5.10	120.86	128.58
2	D	301	MTA	N3-C2-N1	-5.09	120.87	128.58
2	G	301	MTA	C5-C4-N3	-4.90	119.97	126.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	301	MTA	C5-C4-N3	-4.55	120.45	126.72
2	G	301	MTA	N3-C4-N9	4.45	134.74	127.17
2	C	301	MTA	N3-C4-N9	4.43	134.69	127.17
2	G	301	MTA	N3-C2-N1	-4.30	122.07	128.58
2	E	301	MTA	N3-C4-N9	4.21	134.33	127.17
2	F	301	MTA	C5-C4-N3	-4.10	121.07	126.72
2	E	301	MTA	N3-C2-N1	-3.95	122.61	128.58
2	B	301	MTA	N3-C2-N1	-3.84	122.77	128.58
2	B	301	MTA	C5-C4-N3	-3.83	121.45	126.72
2	A	301	MTA	N3-C4-N9	3.76	133.57	127.17
2	F	301	MTA	C4-N9-C8	3.72	109.65	105.74
2	D	301	MTA	C5-C4-N3	-3.71	121.61	126.72
2	H	301	MTA	C4-N9-C8	3.71	109.63	105.74
2	G	301	MTA	C2-N3-C4	3.71	120.88	111.83
2	D	301	MTA	N3-C4-N9	3.67	133.40	127.17
2	C	301	MTA	C2-N3-C4	3.64	120.73	111.83
2	G	301	MTA	C4'-O4'-C1'	-3.64	101.43	109.47
2	H	301	MTA	C5-C4-N3	-3.63	121.72	126.72
2	F	301	MTA	C2-N3-C4	3.63	120.69	111.83
2	F	301	MTA	C2-N1-C6	3.60	124.64	118.73
2	H	301	MTA	C4-C5-N7	-3.60	106.47	110.58
2	D	301	MTA	N9-C8-N7	-3.56	108.89	113.94
2	D	301	MTA	C2-N3-C4	3.53	120.45	111.83
2	F	301	MTA	N3-C4-N9	3.49	133.11	127.17
2	D	301	MTA	C2-N1-C6	3.45	124.40	118.73
2	A	301	MTA	C2-N3-C4	3.42	120.19	111.83
2	G	301	MTA	C4-N9-C8	3.42	109.33	105.74
2	H	301	MTA	C4-N9-C1'	-3.27	119.00	126.63
2	B	301	MTA	C2-N1-C6	3.26	124.08	118.73
2	A	301	MTA	N3-C2-N1	-3.25	123.66	128.58
2	C	301	MTA	N3-C2-N1	-3.25	123.66	128.58
2	E	301	MTA	C4-N9-C8	3.25	109.15	105.74
2	B	301	MTA	C4-N9-C8	3.22	109.12	105.74
4	G	303	EPE	O2S-S-C10	3.20	111.56	106.73
2	C	301	MTA	C4-C5-N7	-3.15	106.98	110.58
2	H	301	MTA	C6-C5-N7	3.13	138.12	132.09
4	A	303	EPE	C6-N1-C2	3.10	115.51	108.84
2	B	301	MTA	N3-C4-N9	3.09	132.42	127.17
2	E	301	MTA	C2-N1-C6	3.06	123.77	118.73
2	A	301	MTA	C4-C5-N7	-3.05	107.10	110.58
2	E	301	MTA	C4'-C5'-S5'	-3.04	103.77	113.89
2	B	301	MTA	C2-N3-C4	3.03	119.23	111.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301	MTA	C4'-O4'-C1'	-3.03	102.78	109.47
2	B	301	MTA	C4-C5-N7	-2.93	107.23	110.58
2	E	301	MTA	C2-N3-C4	2.88	118.87	111.83
2	H	301	MTA	N3-C2-N1	-2.87	124.23	128.58
2	H	301	MTA	C2-N3-C4	2.85	118.80	111.83
2	H	301	MTA	N3-C4-N9	2.85	132.02	127.17
2	G	301	MTA	C4-C5-N7	-2.84	107.33	110.58
2	F	301	MTA	C4-N9-C1'	-2.83	120.02	126.63
4	A	303	EPE	O1S-S-C10	2.83	111.00	106.73
2	F	301	MTA	C4-C5-N7	-2.75	107.44	110.58
2	A	301	MTA	N9-C8-N7	-2.72	110.08	113.94
2	H	301	MTA	N9-C8-N7	-2.68	110.14	113.94
2	F	301	MTA	C6-C5-N7	2.63	137.16	132.09
2	B	301	MTA	C6-C5-N7	2.61	137.12	132.09
2	A	301	MTA	C5-N7-C8	2.59	107.52	103.45
2	H	301	MTA	C5-N7-C8	2.56	107.48	103.45
2	D	301	MTA	C4-N9-C1'	-2.55	120.66	126.63
2	B	301	MTA	C2'-C3'-C4'	2.53	107.50	102.61
2	F	301	MTA	N9-C8-N7	-2.53	110.35	113.94
2	A	301	MTA	C4-N9-C8	2.47	108.33	105.74
2	D	301	MTA	C4-C5-N7	-2.47	107.76	110.58
2	E	301	MTA	N6-C6-N1	2.46	123.87	118.38
4	G	303	EPE	O1S-S-C10	2.46	110.44	106.73
2	D	301	MTA	C6-C5-N7	2.40	136.71	132.09
2	C	301	MTA	C5-N7-C8	2.39	107.21	103.45
2	H	301	MTA	C4'-C5'-S5'	-2.34	106.09	113.89
2	D	301	MTA	C5-N7-C8	2.31	107.08	103.45
2	F	301	MTA	C4'-C5'-S5'	-2.29	106.25	113.89
2	B	301	MTA	C4-N9-C1'	-2.27	121.33	126.63
2	E	301	MTA	C4-N9-C1'	-2.25	121.38	126.63
2	A	301	MTA	C4'-C5'-S5'	-2.23	106.48	113.89
3	A	302	SPM	C3-C4-N5	-2.12	106.39	112.07
2	H	301	MTA	O2'-C2'-C3'	2.10	118.55	111.82
2	F	301	MTA	C5-N7-C8	2.09	106.74	103.45
2	C	301	MTA	C4-N9-C8	2.04	107.89	105.74
2	A	301	MTA	C6-C5-N7	2.02	135.98	132.09
2	D	301	MTA	C4'-O4'-C1'	-2.01	105.02	109.47
2	D	301	MTA	O3'-C3'-C4'	-2.00	105.33	111.08

There are no chirality outliers.

All (70) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	301	MTA	O4'-C4'-C5'-S5'
2	A	301	MTA	C3'-C4'-C5'-S5'
2	C	301	MTA	C3'-C4'-C5'-S5'
2	D	301	MTA	O4'-C4'-C5'-S5'
2	D	301	MTA	C3'-C4'-C5'-S5'
2	E	301	MTA	O4'-C4'-C5'-S5'
2	E	301	MTA	C3'-C4'-C5'-S5'
2	F	301	MTA	O4'-C4'-C5'-S5'
2	F	301	MTA	C3'-C4'-C5'-S5'
3	B	302	SPM	C3-C4-N5-C6
3	E	302	SPM	C12-C11-N10-C9
4	A	303	EPE	C10-C9-N1-C2
4	A	303	EPE	C10-C9-N1-C6
3	G	302	SPM	C7-C8-C9-N10
3	F	302	SPM	N5-C6-C7-C8
3	F	302	SPM	C7-C8-C9-N10
3	E	302	SPM	N5-C6-C7-C8
3	G	302	SPM	N5-C6-C7-C8
3	A	302	SPM	C8-C9-N10-C11
3	E	302	SPM	C7-C8-C9-N10
3	A	302	SPM	C2-C3-C4-N5
3	D	302	SPM	N10-C11-C12-C13
3	H	302	SPM	C2-C3-C4-N5
3	E	302	SPM	C2-C3-C4-N5
3	A	302	SPM	C7-C6-N5-C4
3	C	302	SPM	C3-C4-N5-C6
3	D	302	SPM	C7-C6-N5-C4
3	E	302	SPM	C7-C6-N5-C4
3	H	302	SPM	C3-C4-N5-C6
4	A	303	EPE	N4-C7-C8-O8
3	B	302	SPM	N10-C11-C12-C13
3	D	302	SPM	C7-C8-C9-N10
3	G	302	SPM	C7-C6-N5-C4
3	D	302	SPM	C6-C7-C8-C9
3	A	302	SPM	C7-C8-C9-N10
3	H	302	SPM	C6-C7-C8-C9
3	C	302	SPM	N5-C6-C7-C8
4	G	303	EPE	C10-C9-N1-C2
4	G	303	EPE	C10-C9-N1-C6
3	F	302	SPM	C6-C7-C8-C9
4	G	303	EPE	N4-C7-C8-O8
3	G	302	SPM	C6-C7-C8-C9
3	B	302	SPM	C2-C3-C4-N5

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Mol	Chain	Res	Type	Atoms
3	A	302	SPM	N1-C2-C3-C4
3	B	302	SPM	N1-C2-C3-C4
3	C	302	SPM	C12-C11-N10-C9
3	F	302	SPM	C3-C4-N5-C6
3	G	302	SPM	C11-C12-C13-N14
3	H	302	SPM	N10-C11-C12-C13
3	C	302	SPM	C7-C6-N5-C4
3	F	302	SPM	C2-C3-C4-N5
2	C	301	MTA	O4'-C4'-C5'-S5'
2	G	301	MTA	O4'-C4'-C5'-S5'
2	H	301	MTA	O4'-C4'-C5'-S5'
3	C	302	SPM	N10-C11-C12-C13
3	A	302	SPM	C3-C4-N5-C6
3	C	302	SPM	N1-C2-C3-C4
3	E	302	SPM	N1-C2-C3-C4
3	F	302	SPM	N1-C2-C3-C4
3	F	302	SPM	C12-C11-N10-C9
3	H	302	SPM	C8-C9-N10-C11
3	E	302	SPM	C3-C4-N5-C6
3	E	302	SPM	C8-C9-N10-C11
3	F	302	SPM	C8-C9-N10-C11
3	A	302	SPM	C11-C12-C13-N14
3	F	302	SPM	C7-C6-N5-C4
3	G	302	SPM	C12-C11-N10-C9
3	A	302	SPM	C12-C11-N10-C9
2	H	301	MTA	C3'-C4'-C5'-S5'
3	C	302	SPM	C2-C3-C4-N5

There are no ring outliers.

11 monomers are involved in 37 short contacts:

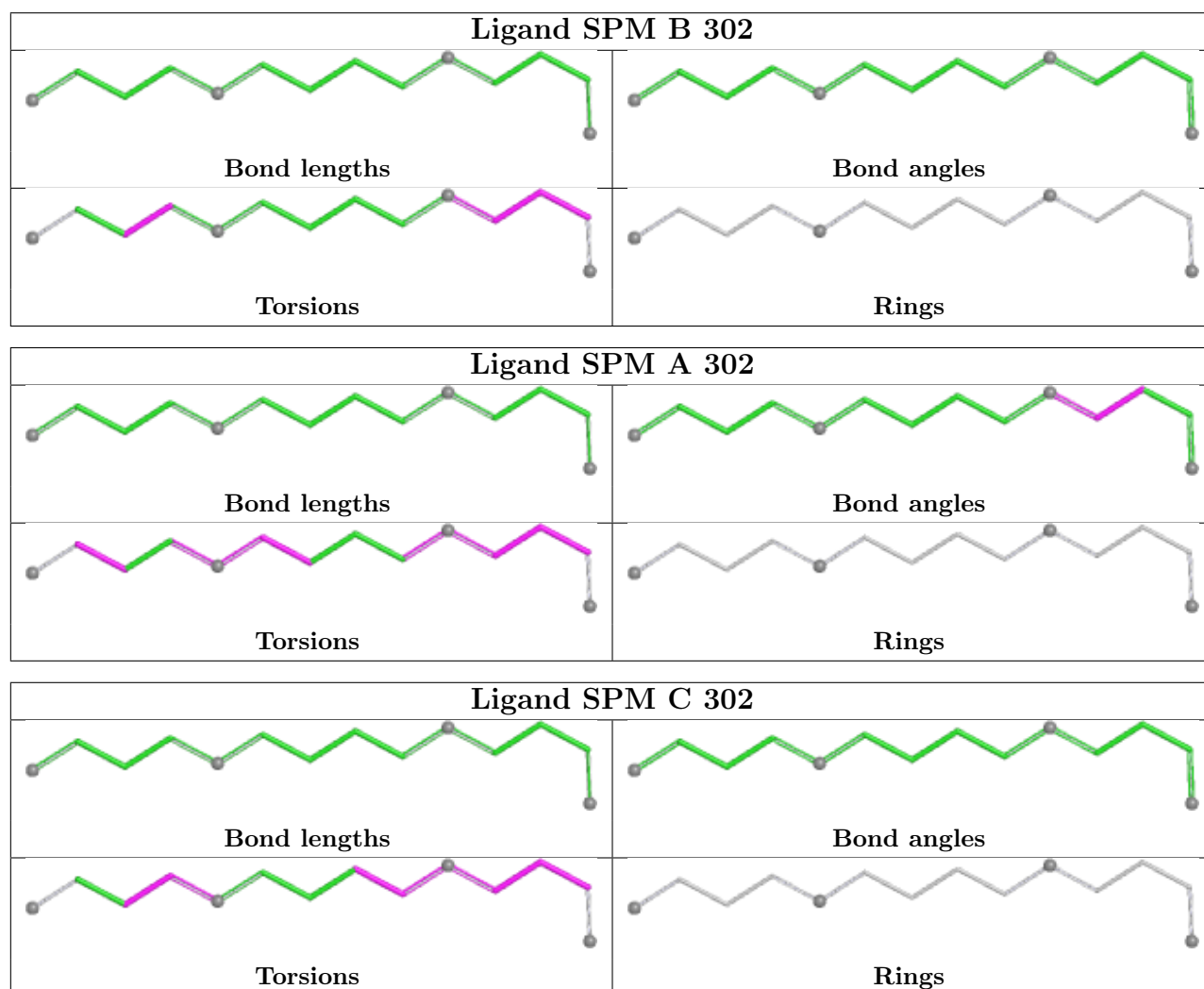
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	302	SPM	3	0
3	A	302	SPM	5	0
3	C	302	SPM	4	0
2	C	301	MTA	1	0
3	G	302	SPM	4	0
2	G	301	MTA	1	0
3	F	302	SPM	3	0
3	H	302	SPM	9	0
3	E	302	SPM	4	0
3	D	302	SPM	5	0

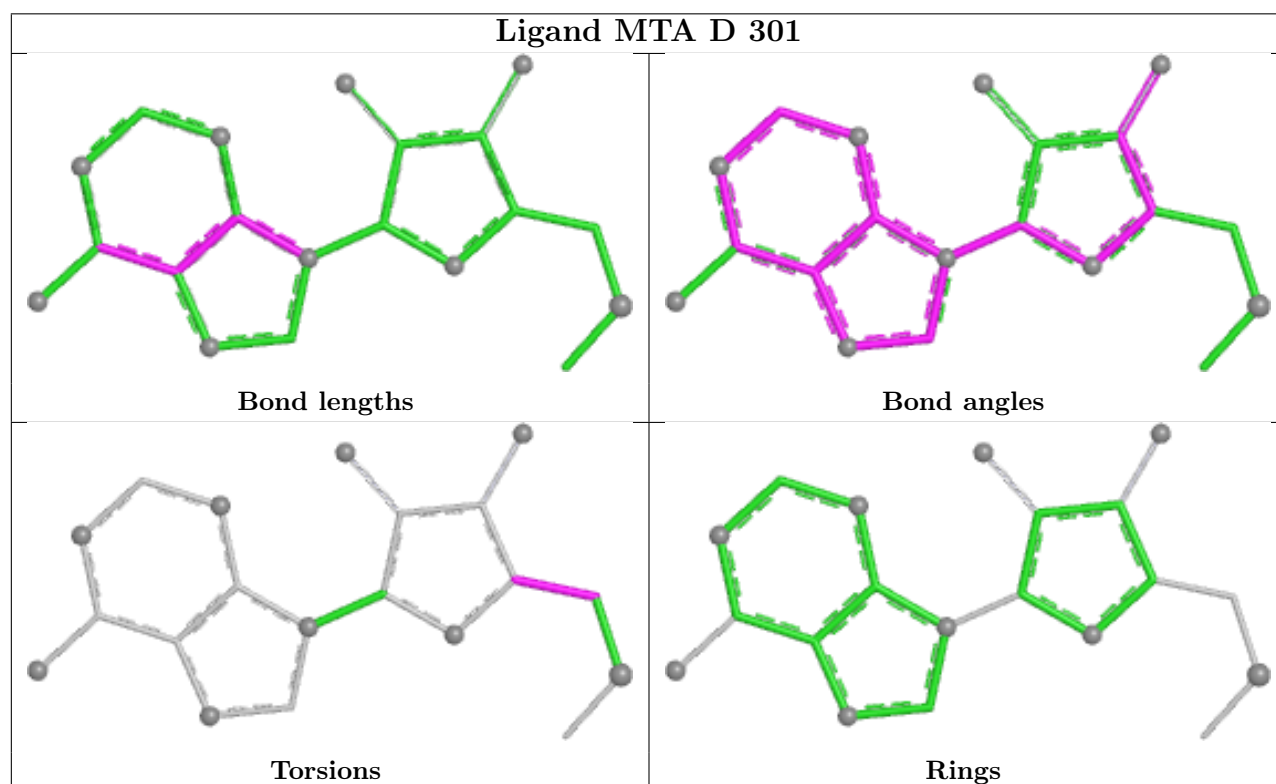
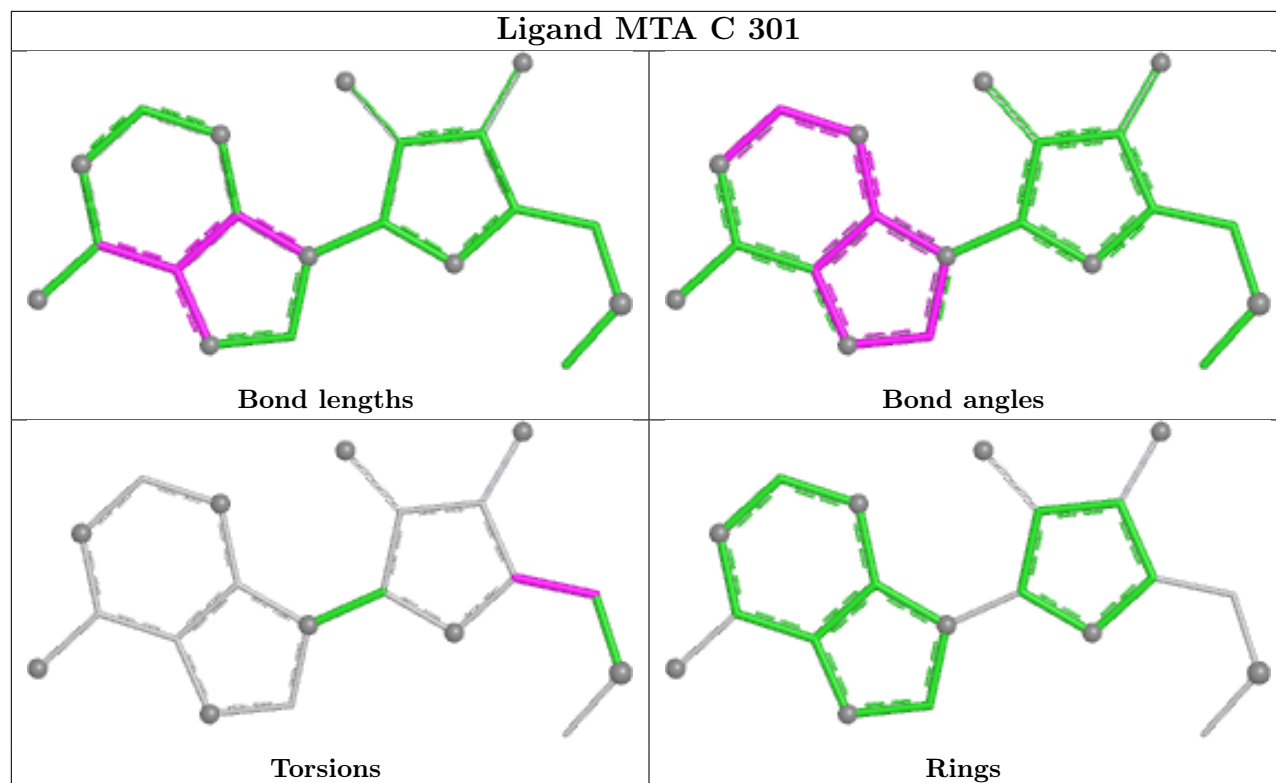
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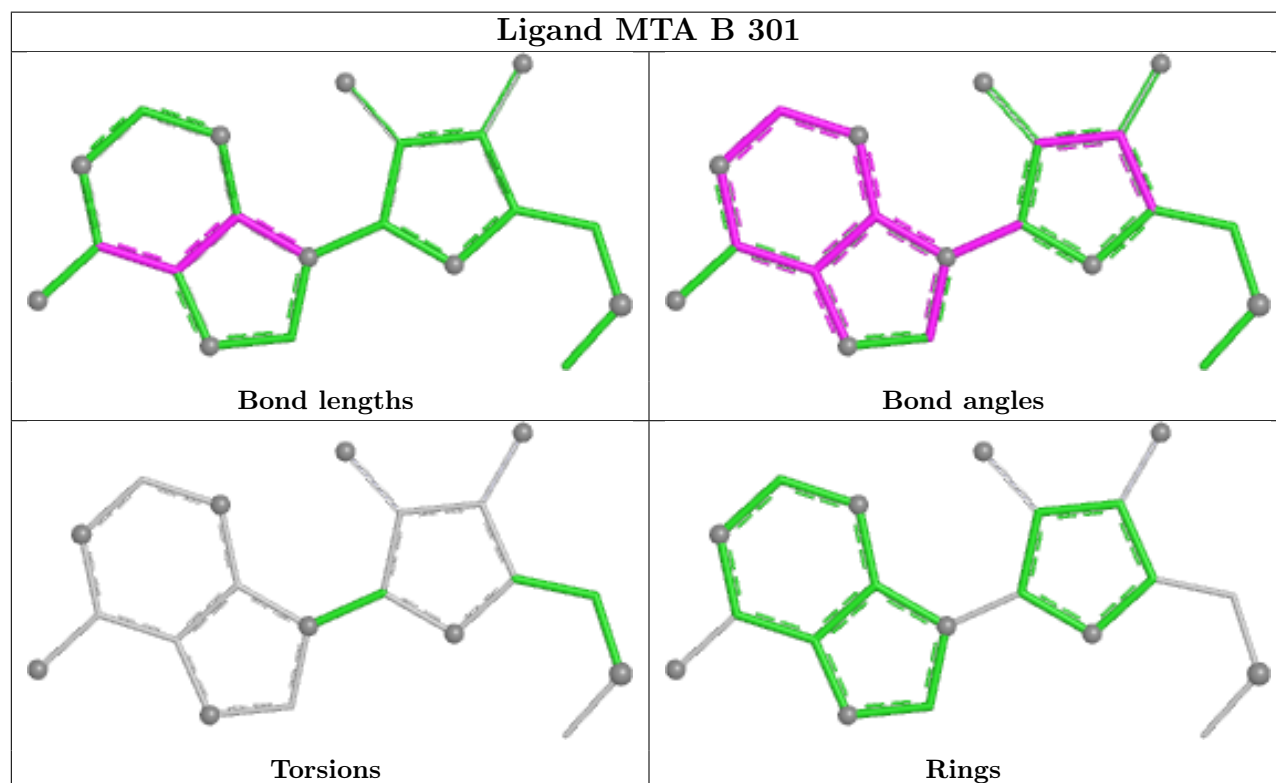
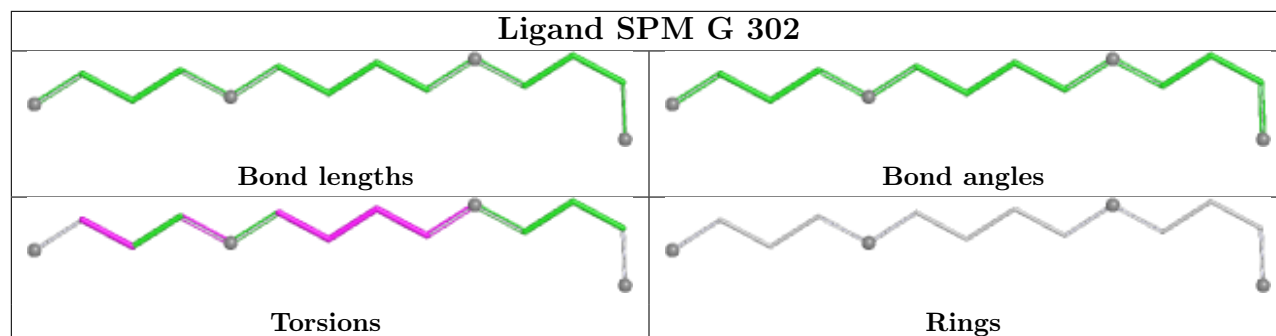
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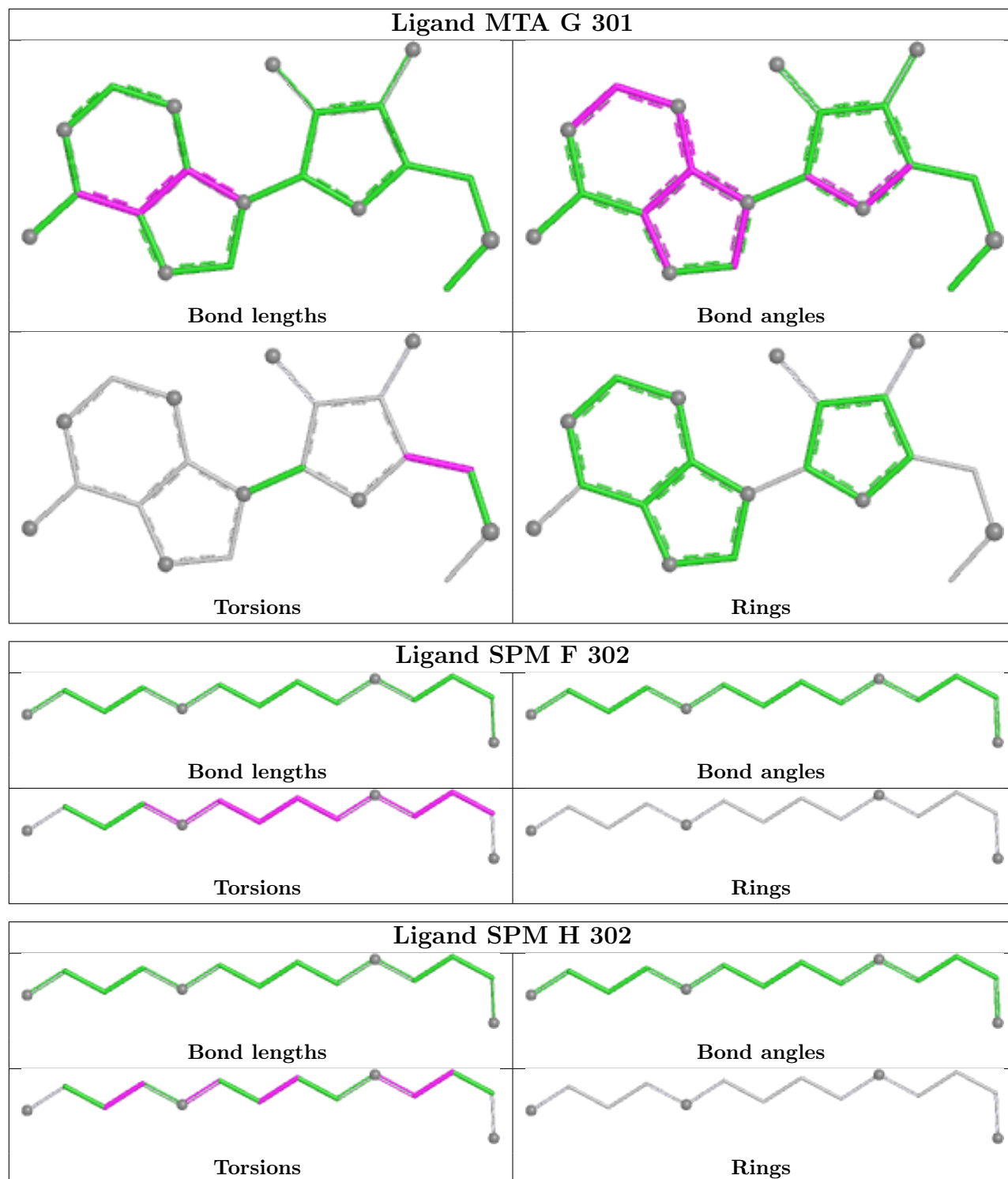
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	H	301	MTA	1	0

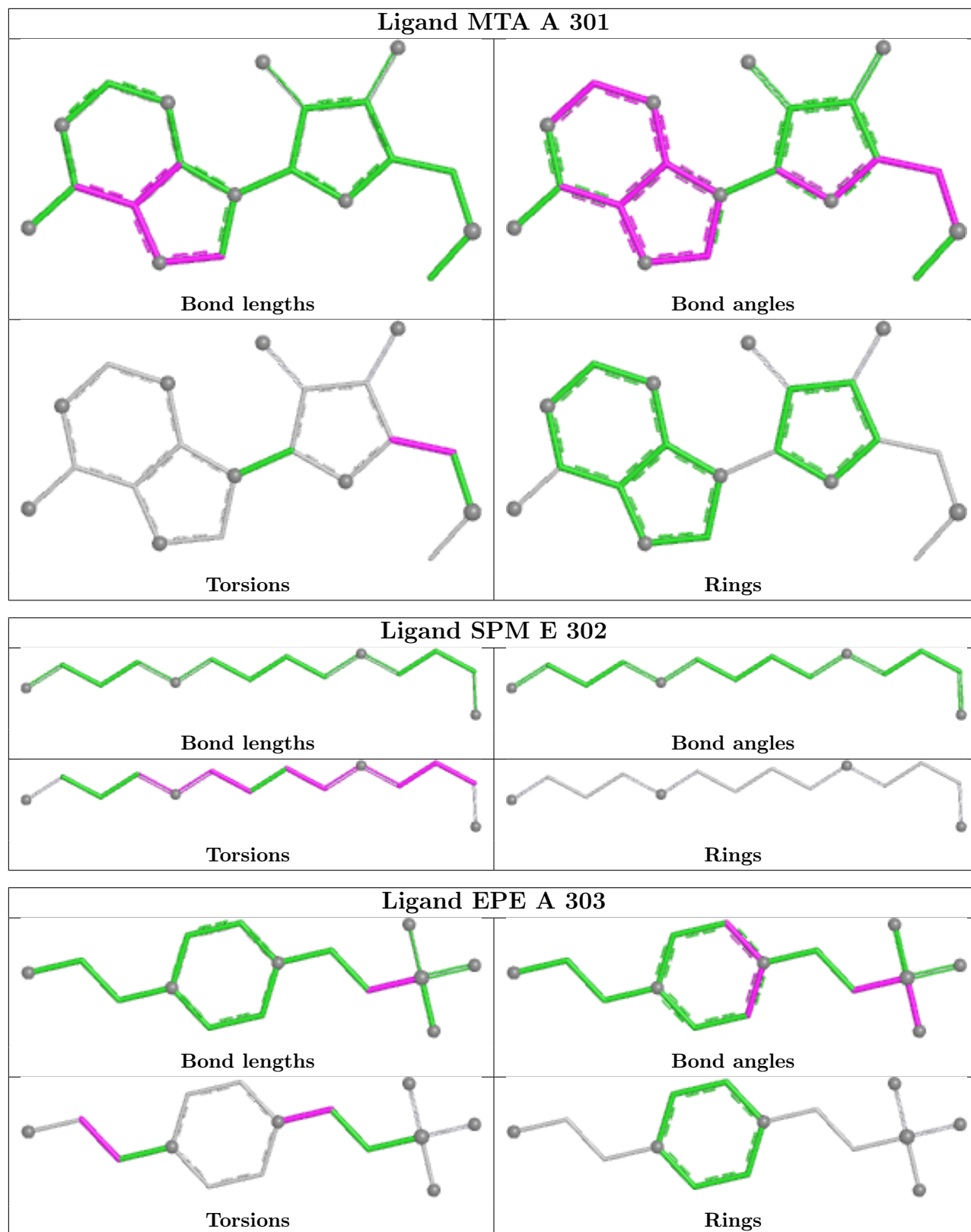
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

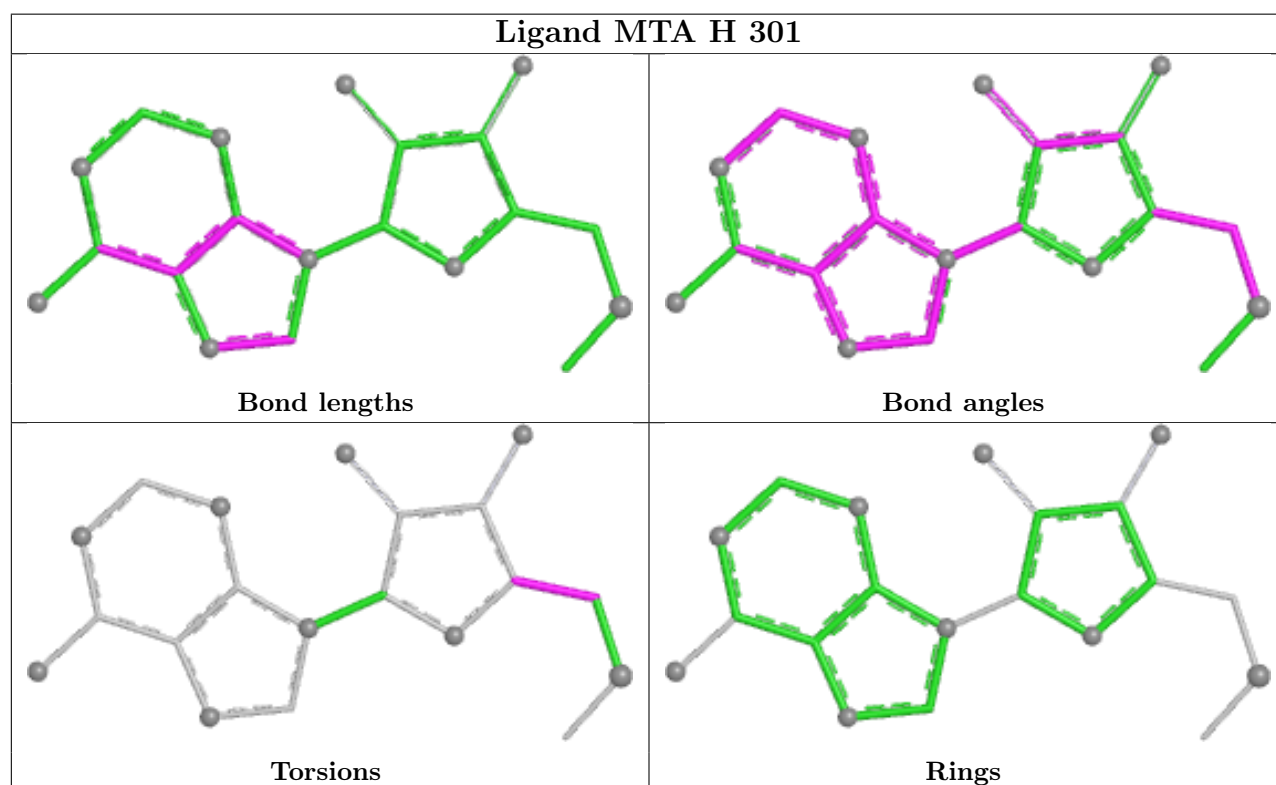
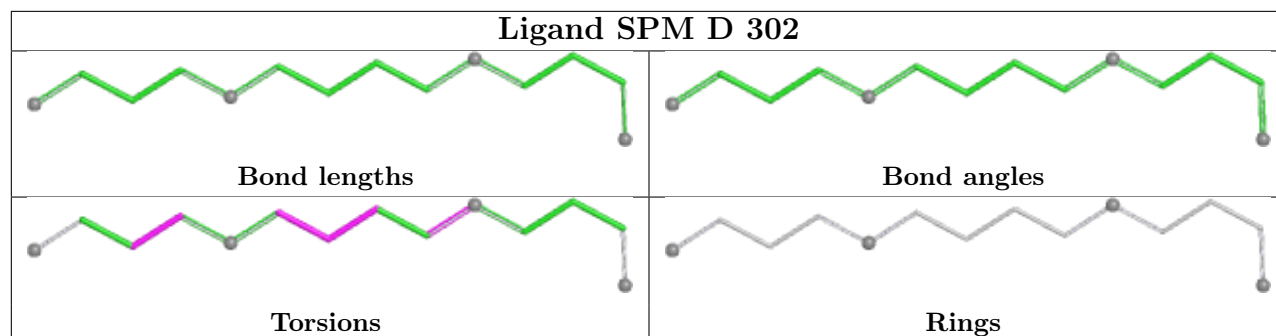


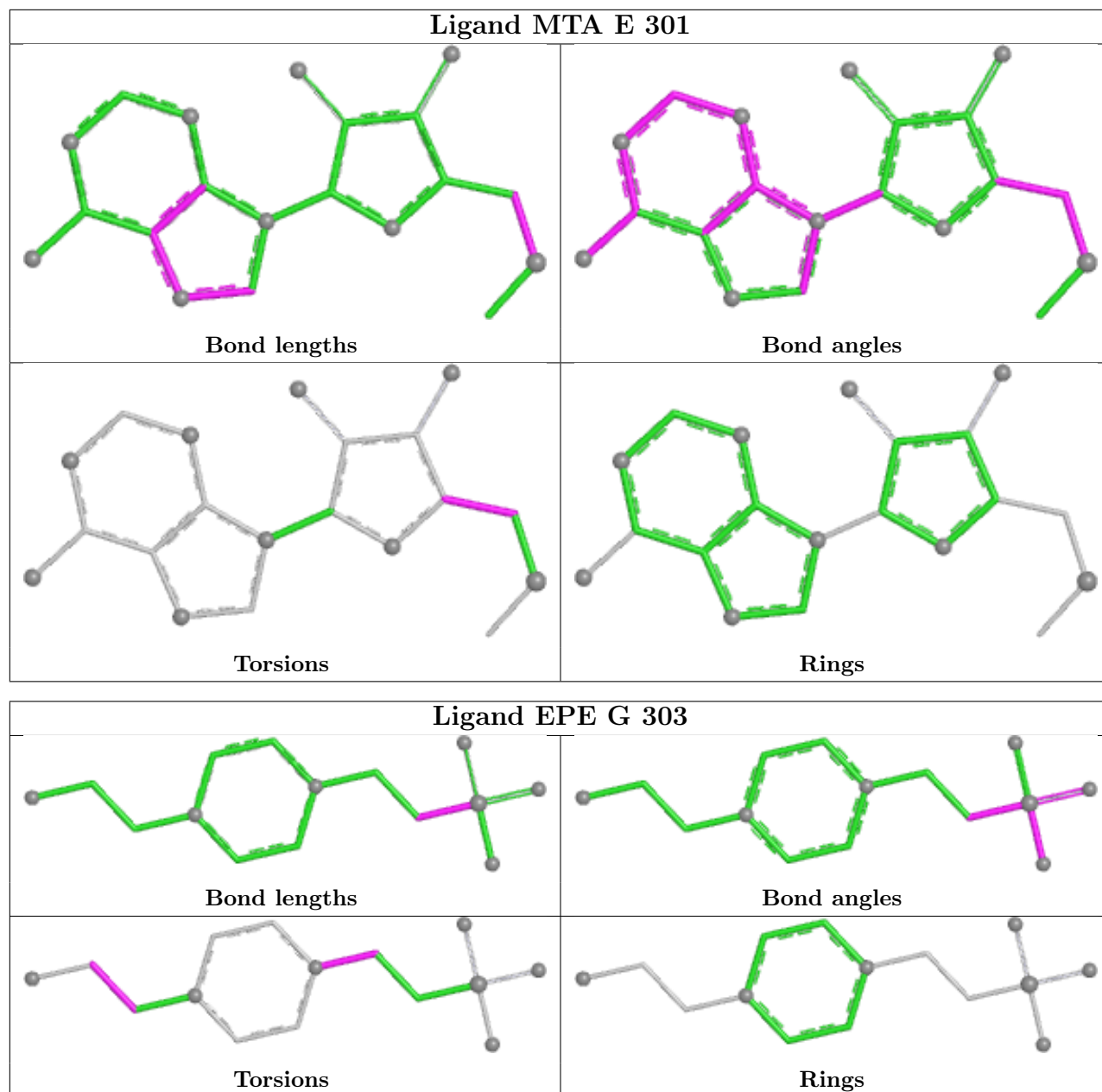


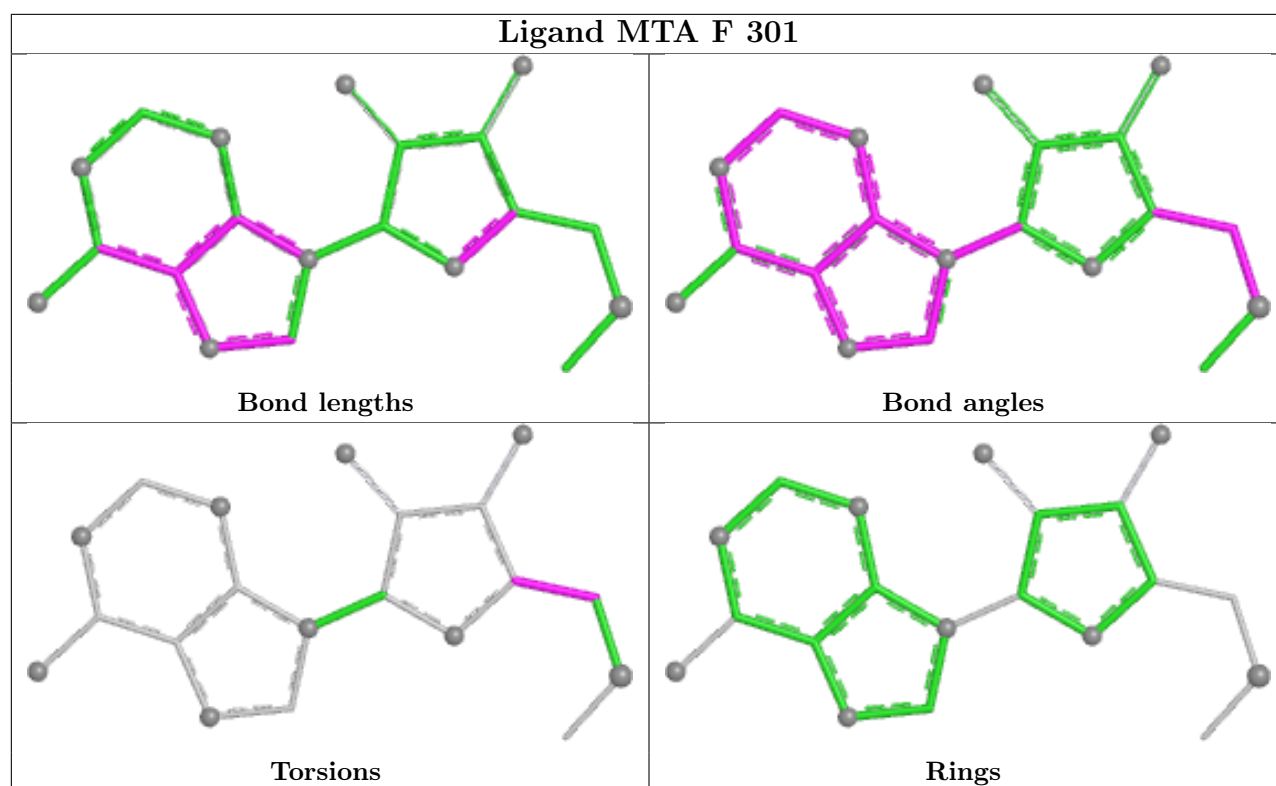












5.7 Other polymers [i](#)

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	286/309 (92%)	0.00	3 (1%) 79 81	29, 41, 61, 92	0
1	B	288/309 (93%)	0.28	7 (2%) 59 61	29, 44, 71, 86	0
1	C	286/309 (92%)	0.19	3 (1%) 79 81	28, 43, 65, 85	0
1	D	288/309 (93%)	0.41	9 (3%) 51 51	28, 46, 75, 88	0
1	E	286/309 (92%)	0.47	4 (1%) 73 74	28, 46, 71, 97	1 (0%)
1	F	286/309 (92%)	0.57	13 (4%) 38 37	29, 48, 79, 90	0
1	G	286/309 (92%)	0.06	5 (1%) 69 70	29, 41, 62, 102	0
1	H	288/309 (93%)	0.21	6 (2%) 63 64	28, 44, 70, 90	0
All	All	2294/2472 (92%)	0.27	50 (2%) 62 63	28, 44, 71, 102	1 (0%)

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	4	VAL	6.3
1	A	4	VAL	6.2
1	E	4	VAL	5.9
1	C	4	VAL	5.1
1	F	212	MET	4.3
1	F	155	GLY	3.9
1	H	4	VAL	3.6
1	F	133	ASP	3.6
1	F	4	VAL	3.5
1	D	212	MET	3.3
1	B	156	ASP	3.2
1	F	180	ARG	3.1
1	B	3	LYS	3.0
1	D	3	LYS	3.0
1	E	105	LYS	3.0
1	H	2	ARG	3.0

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Mol	Chain	Res	Type	RSRZ
1	D	2	ARG	2.9
1	F	105	LYS	2.7
1	B	259	GLU	2.6
1	H	3	LYS	2.6
1	C	181	GLU	2.5
1	E	118	GLU	2.5
1	D	181	GLU	2.5
1	G	180	ARG	2.4
1	D	4	VAL	2.4
1	F	145	PHE	2.4
1	F	144	GLY	2.3
1	B	122	ALA	2.3
1	A	180	ARG	2.3
1	H	259	GLU	2.3
1	B	4	VAL	2.3
1	F	111	ASP	2.2
1	B	2	ARG	2.2
1	D	156	ASP	2.2
1	F	115	ASP	2.2
1	B	212	MET	2.1
1	F	186	ILE	2.1
1	F	118	GLU	2.1
1	G	118	GLU	2.1
1	H	181	GLU	2.1
1	A	262	ARG	2.1
1	G	262	ARG	2.1
1	F	37	GLU	2.1
1	D	172	ASP	2.1
1	G	192	ASP	2.1
1	C	247	TYR	2.1
1	H	247	TYR	2.0
1	D	259	GLU	2.0
1	E	37	GLU	2.0
1	D	153	LYS	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 oligosaccharides in this entry.

6.4 Ligands ⓘ

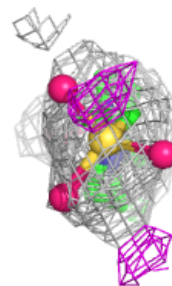
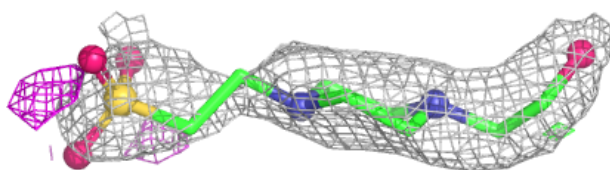
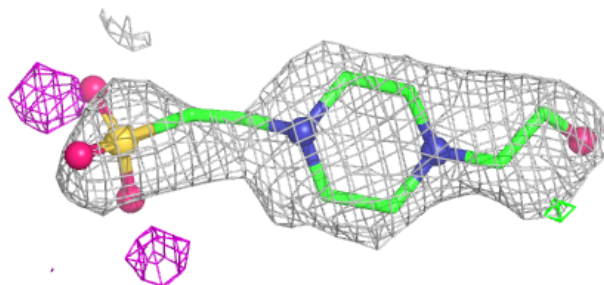
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	EPE	A	303	15/15	0.79	0.19	73,85,135,141	0
4	EPE	G	303	15/15	0.89	0.14	58,69,93,97	0
3	SPM	F	302	14/14	0.90	0.15	45,60,68,69	0
3	SPM	G	302	14/14	0.90	0.14	32,54,59,61	0
3	SPM	C	302	14/14	0.92	0.14	41,48,53,65	0
3	SPM	E	302	14/14	0.92	0.13	38,49,57,63	0
3	SPM	A	302	14/14	0.92	0.12	33,47,55,56	0
3	SPM	B	302	14/14	0.93	0.12	44,52,59,60	0
3	SPM	H	302	14/14	0.93	0.12	39,52,58,58	0
2	MTA	F	301	20/20	0.95	0.08	43,48,59,67	0
3	SPM	D	302	14/14	0.95	0.10	40,55,62,66	0
2	MTA	D	301	20/20	0.95	0.08	41,46,52,53	0
2	MTA	E	301	20/20	0.95	0.08	38,43,48,56	0
2	MTA	A	301	20/20	0.97	0.07	33,36,42,45	0
2	MTA	B	301	20/20	0.97	0.07	39,42,50,59	0
2	MTA	C	301	20/20	0.97	0.06	38,40,49,55	0
2	MTA	G	301	20/20	0.97	0.07	33,37,42,50	0
2	MTA	H	301	20/20	0.97	0.06	42,44,54,57	0

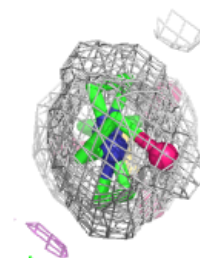
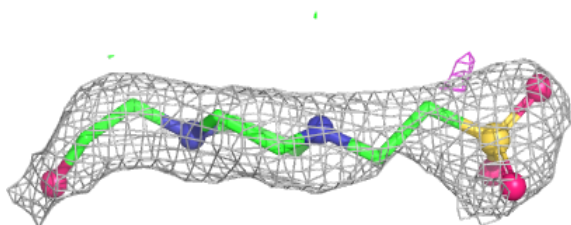
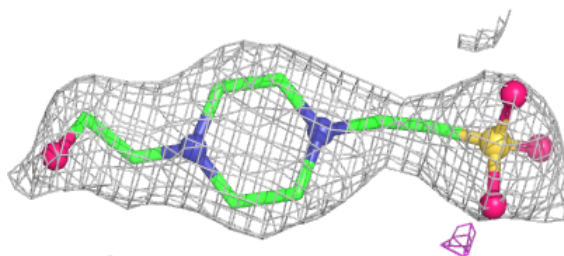
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around EPE A 303:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

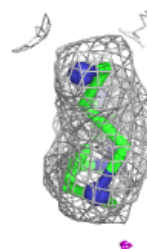
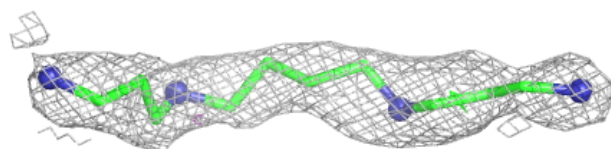
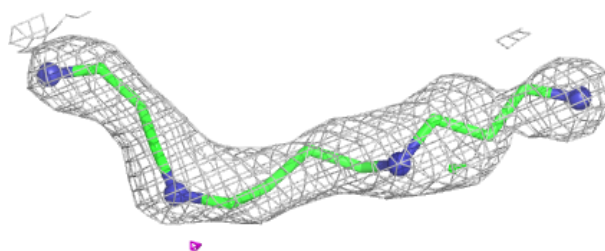
**Electron density around EPE G 303:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

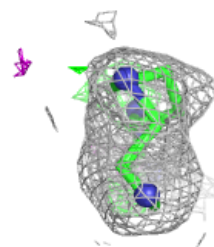
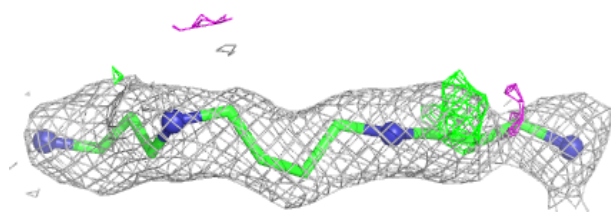
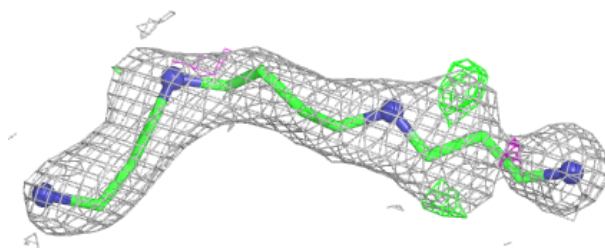


Electron density around SPM F 302:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

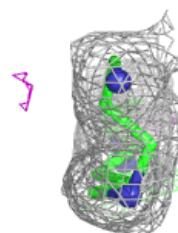
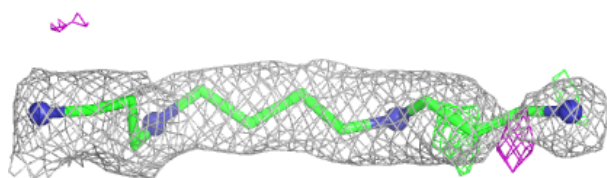
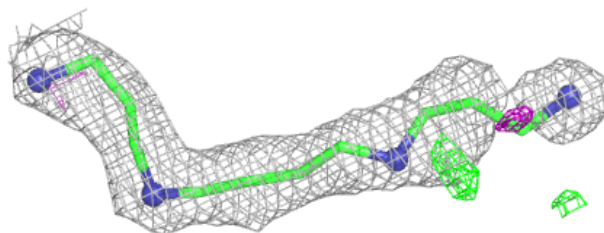
**Electron density around SPM G 302:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

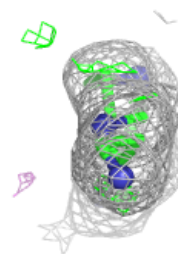
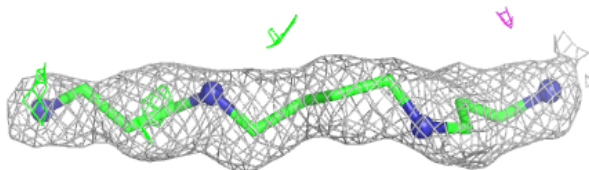
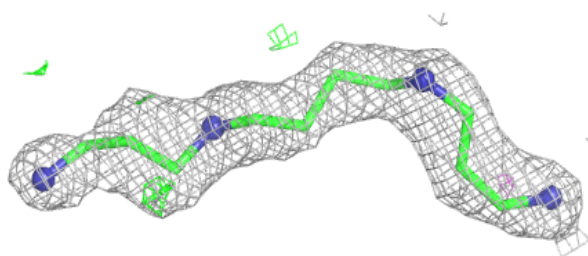


Electron density around SPM C 302:

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and green (positive)

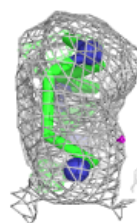
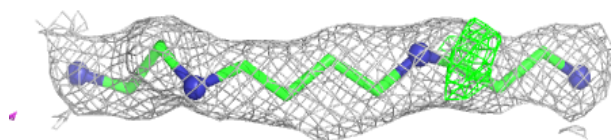
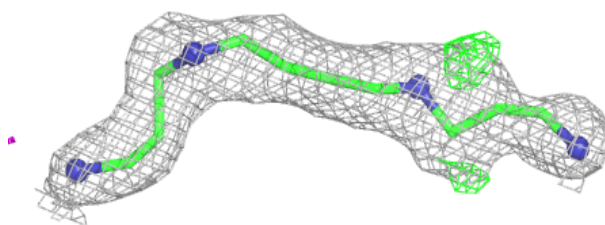
**Electron density around SPM E 302:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

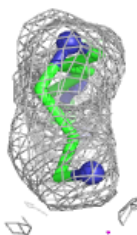
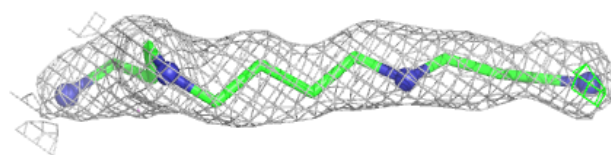
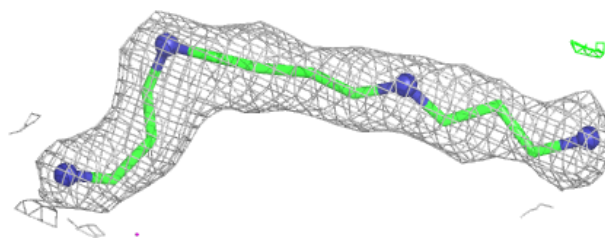


Electron density around SPM A 302:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

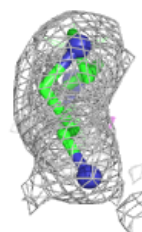
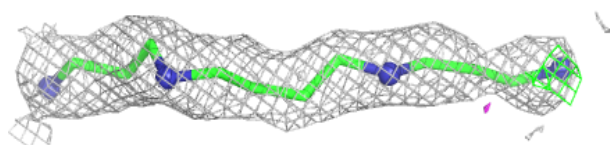
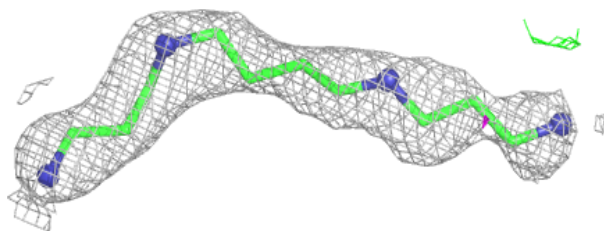
**Electron density around SPM B 302:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

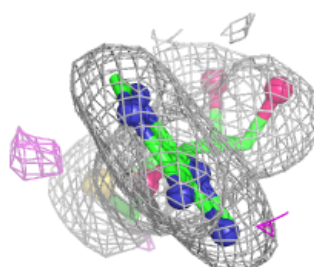
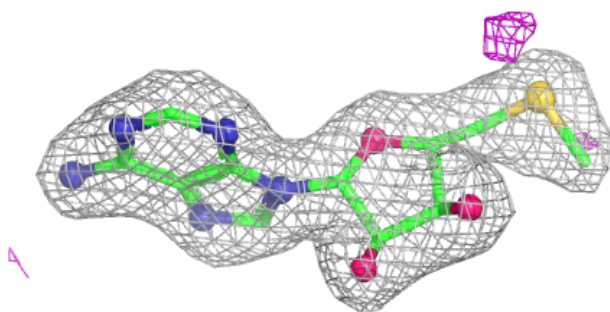
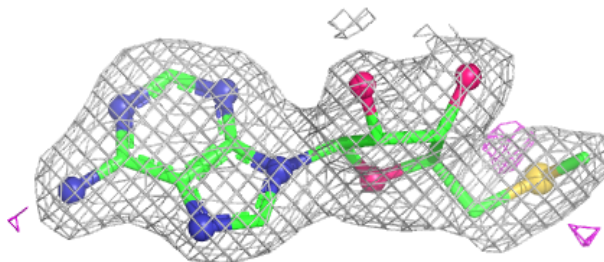


Electron density around SPM H 302:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

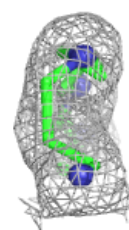
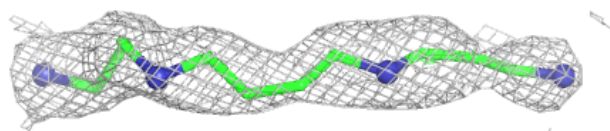
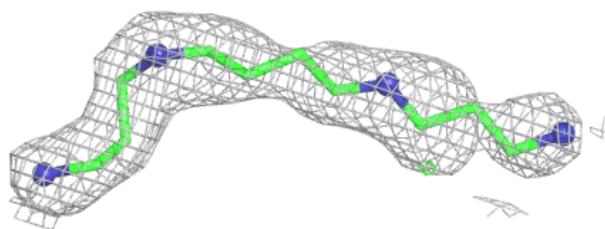
**Electron density around MTA F 301:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

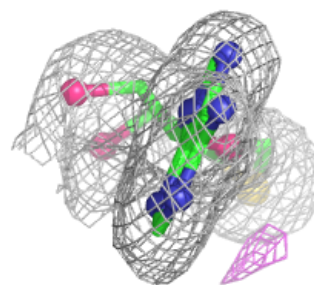
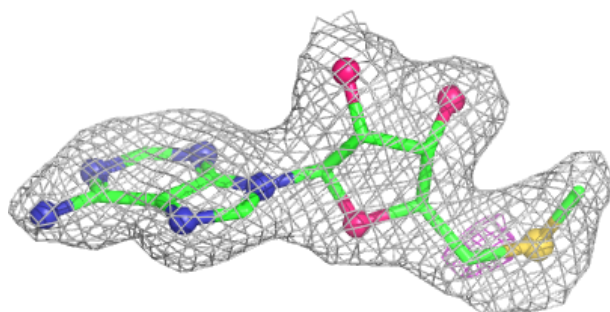
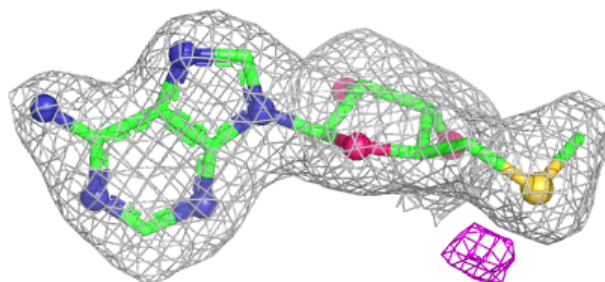


Electron density around SPM D 302:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

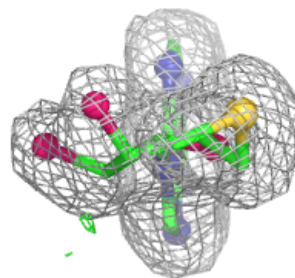
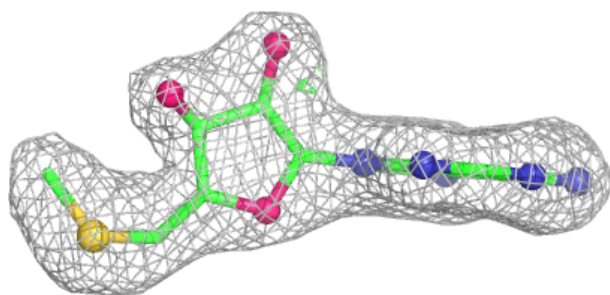
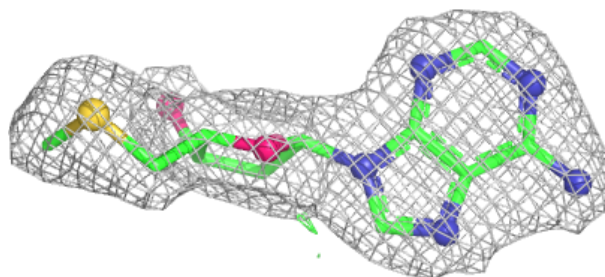
**Electron density around MTA D 301:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

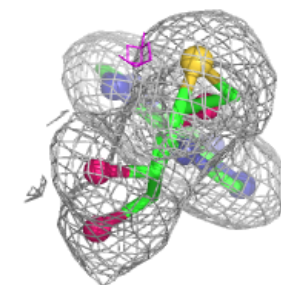
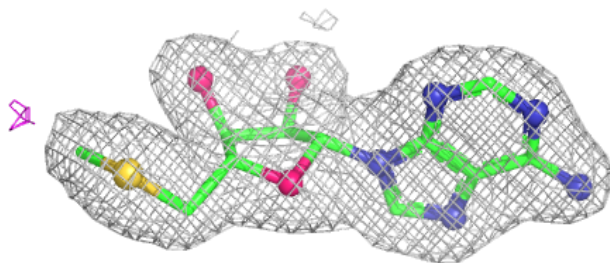
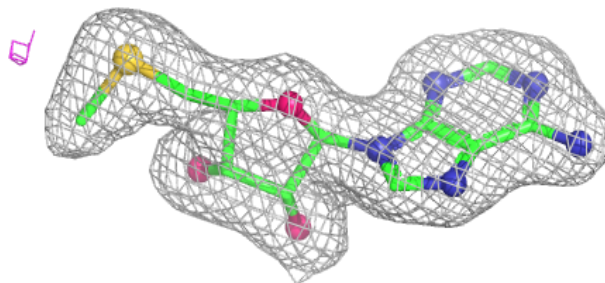


Electron density around MTA E 301:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

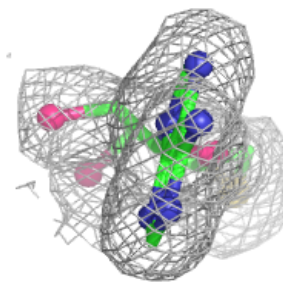
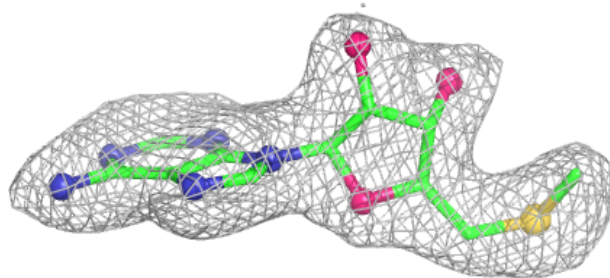
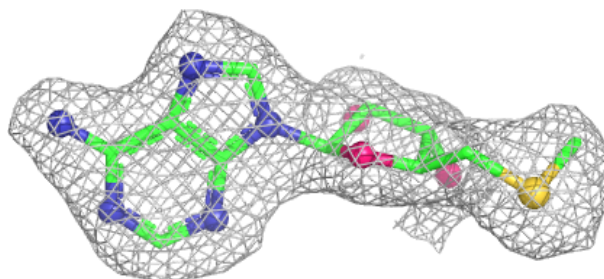
**Electron density around MTA A 301:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

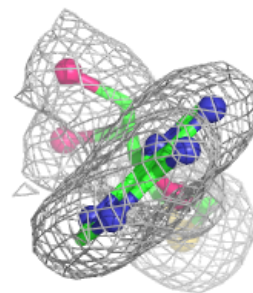
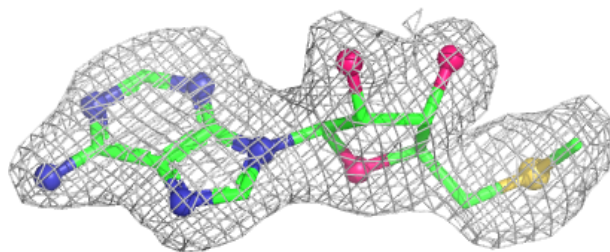
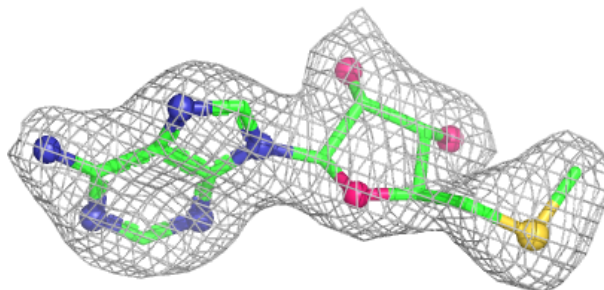


Electron density around MTA B 301:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

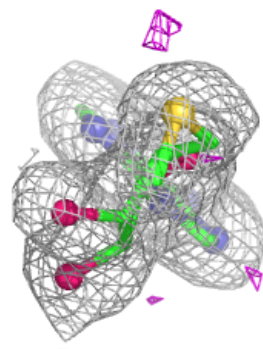
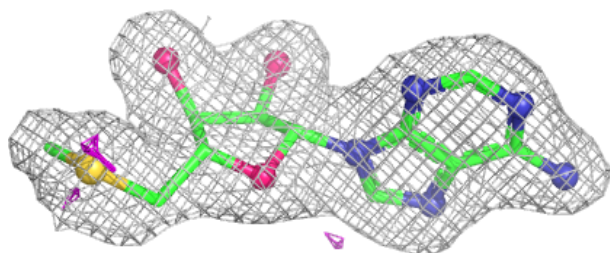
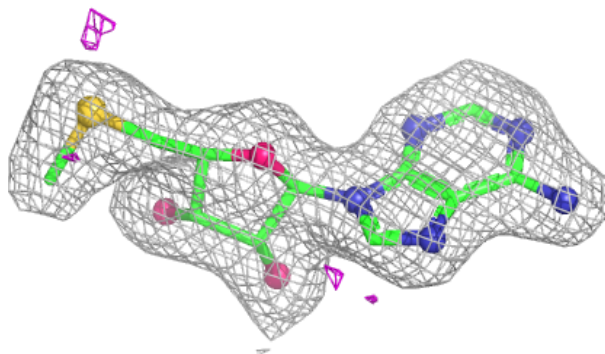
**Electron density around MTA C 301:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

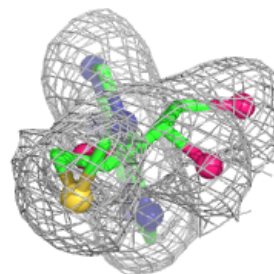
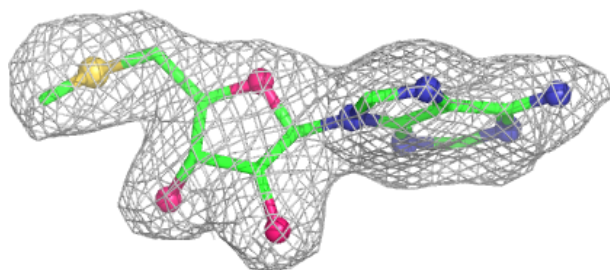
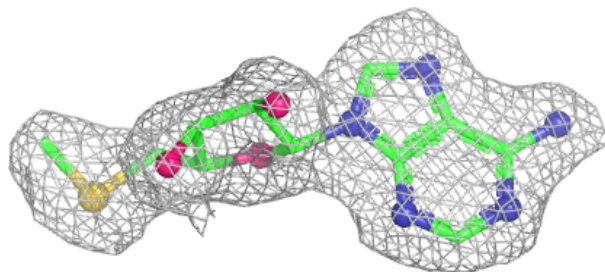


Electron density around MTA G 301:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around MTA H 301:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.