



wwPDB X-ray Structure Validation Summary Report ⓘ

May 15, 2020 – 02:33 pm BST

PDB ID : 3E76
Title : Crystal structure of Wild-type GroEL with bound Thallium ions
Authors : Kiser, P.D.; Lorimer, G.H.; Palczewski, K.
Deposited on : 2008-08-17
Resolution : 3.94 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

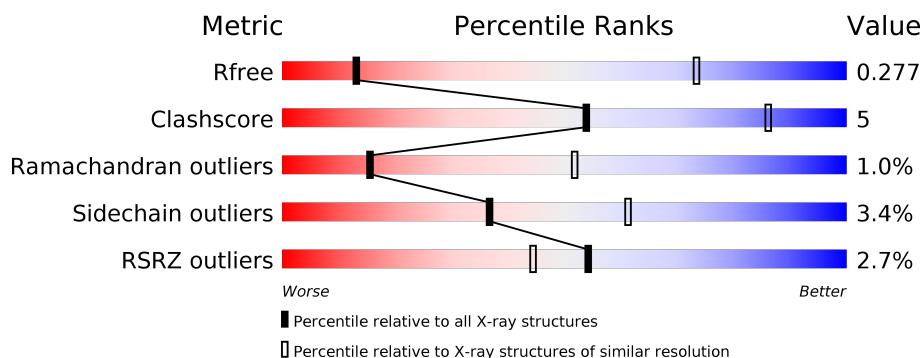
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.94 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}	130704	1036 (4.20-3.68)
Clashscore	141614	1009 (4.18-3.70)
Ramachandran outliers	138981	1057 (4.20-3.68)
Sidechain outliers	138945	1049 (4.20-3.68)
RSRZ outliers	127900	1007 (4.24-3.64)

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 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	547	<div> <div style="width: 78%;"></div> <div style="width: 16%;"></div> <div style="width: 6%;"></div> <div style="width: 2%;"></div> </div>
1	B	547	<div> <div style="width: 82%;"></div> <div style="width: 14%;"></div> <div style="width: 4%;"></div> <div style="width: 2%;"></div> </div>
1	C	547	<div> <div style="width: 2%;"></div> <div style="width: 82%;"></div> <div style="width: 13%;"></div> <div style="width: 3%;"></div> </div>
1	D	547	<div> <div style="width: 1%;"></div> <div style="width: 81%;"></div> <div style="width: 14%;"></div> <div style="width: 4%;"></div> </div>
1	E	547	<div> <div style="width: 2%;"></div> <div style="width: 82%;"></div> <div style="width: 13%;"></div> <div style="width: 3%;"></div> </div>
1	F	547	<div> <div style="width: 2%;"></div> <div style="width: 79%;"></div> <div style="width: 16%;"></div> <div style="width: 3%;"></div> </div>

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Mol	Chain	Length	Quality of chain
1	G	547	<div><div><div></div><div></div><div></div></div><div>%</div><div><div></div><div></div><div></div></div><div>78%</div><div>17%</div><div>.</div><div>.</div></div>
1	H	547	<div><div><div></div><div></div><div></div></div><div>3%</div><div><div></div><div></div><div></div></div><div>82%</div><div>13%</div><div>.</div><div>.</div></div>
1	I	547	<div><div><div></div><div></div><div></div></div><div>2%</div><div><div></div><div></div><div></div></div><div>78%</div><div>17%</div><div>.</div><div>.</div></div>
1	J	547	<div><div><div></div><div></div><div></div></div><div>2%</div><div><div></div><div></div><div></div></div><div>81%</div><div>14%</div><div>.</div><div>.</div></div>
1	K	547	<div><div><div></div><div></div><div></div></div><div>2%</div><div><div></div><div></div><div></div></div><div>79%</div><div>16%</div><div>.</div><div>.</div></div>
1	L	547	<div><div><div></div><div></div><div></div></div><div>4%</div><div><div></div><div></div><div></div></div><div>86%</div><div>10%</div><div>.</div><div>.</div></div>
1	M	547	<div><div><div></div><div></div><div></div></div><div>6%</div><div><div></div><div></div><div></div></div><div>84%</div><div>12%</div><div>.</div><div>.</div></div>
1	N	547	<div><div><div></div><div></div><div></div></div><div>7%</div><div><div></div><div></div><div></div></div><div>86%</div><div>10%</div><div>.</div><div>.</div></div>

2 Entry composition

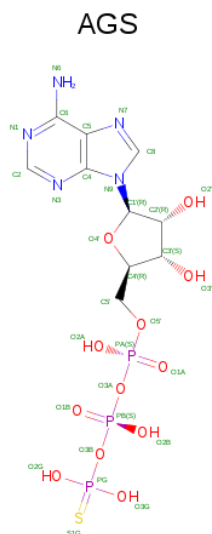
There are 4 unique types of molecules in this entry. The entry contains 54464 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 60 kDa chaperonin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	524	Total	C	N	O	S	0	0	0
			3855	2397	665	773	20			
1	B	524	Total	C	N	O	S	0	0	0
			3855	2397	665	773	20			
1	C	524	Total	C	N	O	S	0	0	0
			3855	2397	665	773	20			
1	D	524	Total	C	N	O	S	0	0	0
			3855	2397	665	773	20			
1	E	524	Total	C	N	O	S	0	0	0
			3855	2397	665	773	20			
1	F	524	Total	C	N	O	S	0	0	0
			3855	2397	665	773	20			
1	G	524	Total	C	N	O	S	0	0	0
			3855	2397	665	773	20			
1	H	524	Total	C	N	O	S	0	0	0
			3855	2397	665	773	20			
1	I	524	Total	C	N	O	S	0	0	0
			3855	2397	665	773	20			
1	J	524	Total	C	N	O	S	0	0	0
			3855	2397	665	773	20			
1	K	524	Total	C	N	O	S	0	0	0
			3855	2397	665	773	20			
1	L	524	Total	C	N	O	S	0	0	0
			3855	2397	665	773	20			
1	M	524	Total	C	N	O	S	0	0	0
			3855	2397	665	773	20			
1	N	524	Total	C	N	O	S	0	0	0
			3855	2397	665	773	20			

- Molecule 2 is PHOSPHOTHIOPHOSPHORIC ACID-ADENYLATE ESTER (three-letter code: AGS) (formula: C₁₀H₁₆N₅O₁₂P₃S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total 31	C 10	N 5	O 12	P 3	S 1	0	0
2	B	1	Total 31	C 10	N 5	O 12	P 3	S 1	0	0
2	C	1	Total 31	C 10	N 5	O 12	P 3	S 1	0	0
2	D	1	Total 31	C 10	N 5	O 12	P 3	S 1	0	0
2	E	1	Total 31	C 10	N 5	O 12	P 3	S 1	0	0
2	F	1	Total 31	C 10	N 5	O 12	P 3	S 1	0	0
2	G	1	Total 31	C 10	N 5	O 12	P 3	S 1	0	0
2	H	1	Total 31	C 10	N 5	O 12	P 3	S 1	0	0
2	I	1	Total 31	C 10	N 5	O 12	P 3	S 1	0	0
2	J	1	Total 31	C 10	N 5	O 12	P 3	S 1	0	0
2	K	1	Total 31	C 10	N 5	O 12	P 3	S 1	0	0
2	L	1	Total 31	C 10	N 5	O 12	P 3	S 1	0	0
2	M	1	Total 31	C 10	N 5	O 12	P 3	S 1	0	0
2	N	1	Total 31	C 10	N 5	O 12	P 3	S 1	0	0

- Molecule 3 is THALLIUM (I) ION (three-letter code: Tl) (formula: Tl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	G	3	Total 3	Tl 3	0	0
3	J	3	Total 3	Tl 3	0	0
3	D	4	Total 4	Tl 4	0	0
3	K	3	Total 3	Tl 3	0	0
3	E	3	Total 3	Tl 3	0	0
3	H	4	Total 4	Tl 4	0	0
3	B	4	Total 4	Tl 4	0	0
3	I	3	Total 3	Tl 3	0	0
3	C	3	Total 3	Tl 3	0	0
3	A	4	Total 4	Tl 4	0	0
3	N	2	Total 2	Tl 2	0	0
3	L	3	Total 3	Tl 3	0	0
3	F	4	Total 4	Tl 4	0	0
3	M	3	Total 3	Tl 3	0	0

- Molecule 4 is MAGNESIUM ION (three-letter code: Mg) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	G	1	Total 1	Mg 1	0	0
4	J	1	Total 1	Mg 1	0	0
4	D	1	Total 1	Mg 1	0	0
4	K	1	Total 1	Mg 1	0	0
4	E	1	Total 1	Mg 1	0	0

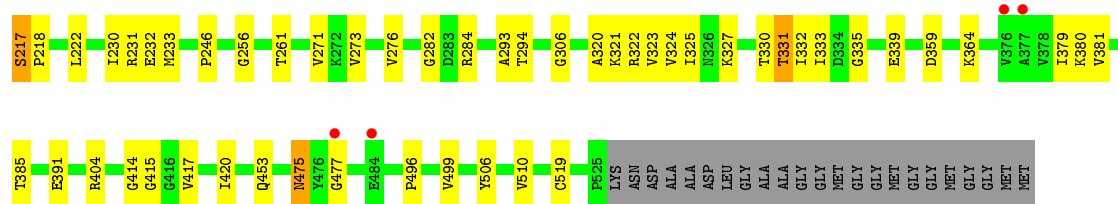
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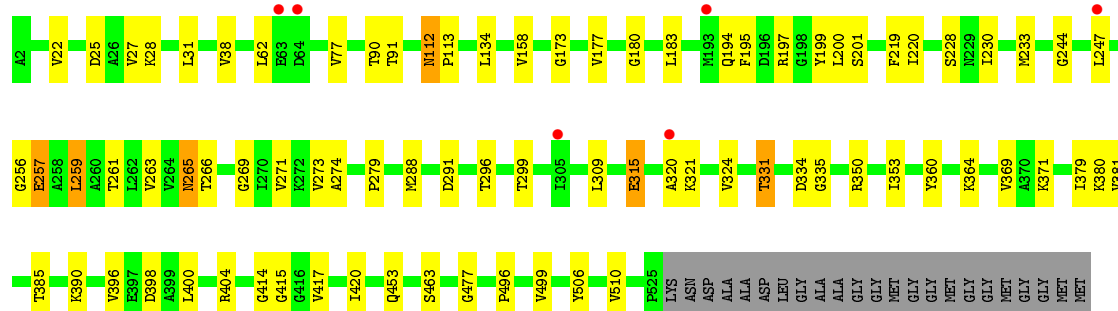
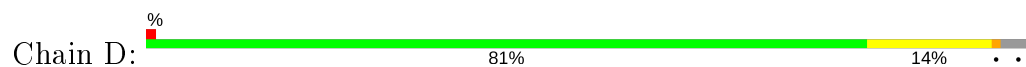
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	H	1	Total 1	Mg 1	0	0
4	B	1	Total 1	Mg 1	0	0
4	I	1	Total 1	Mg 1	0	0
4	C	1	Total 1	Mg 1	0	0
4	A	1	Total 1	Mg 1	0	0
4	N	1	Total 1	Mg 1	0	0
4	L	1	Total 1	Mg 1	0	0
4	F	1	Total 1	Mg 1	0	0
4	M	1	Total 1	Mg 1	0	0

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.

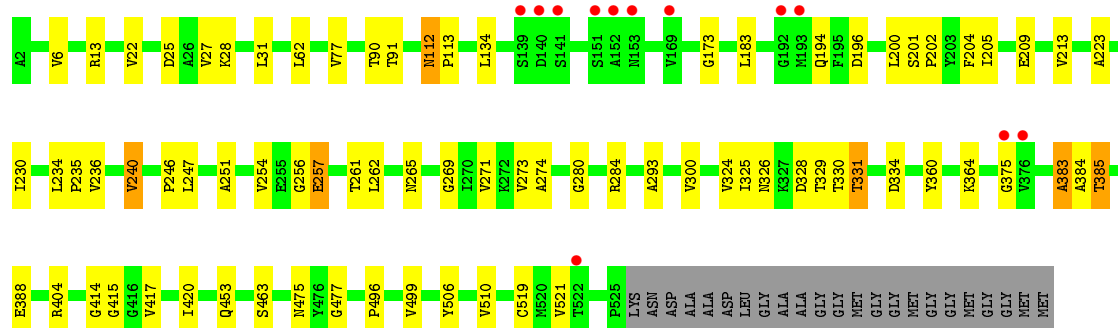
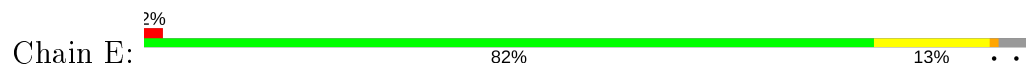
- [illegible]



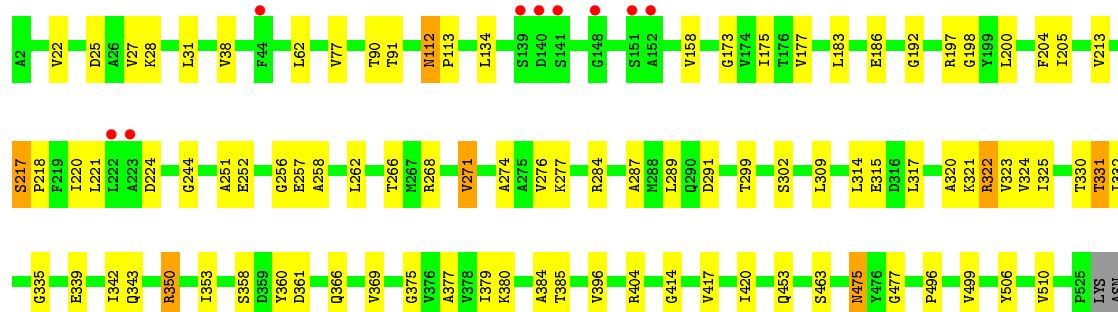
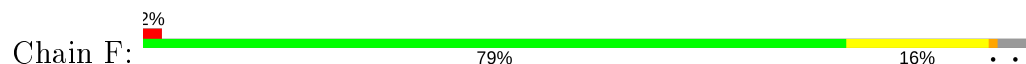
• Molecule 1: 60 kDa chaperonin



• Molecule 1: 60 kDa chaperonin

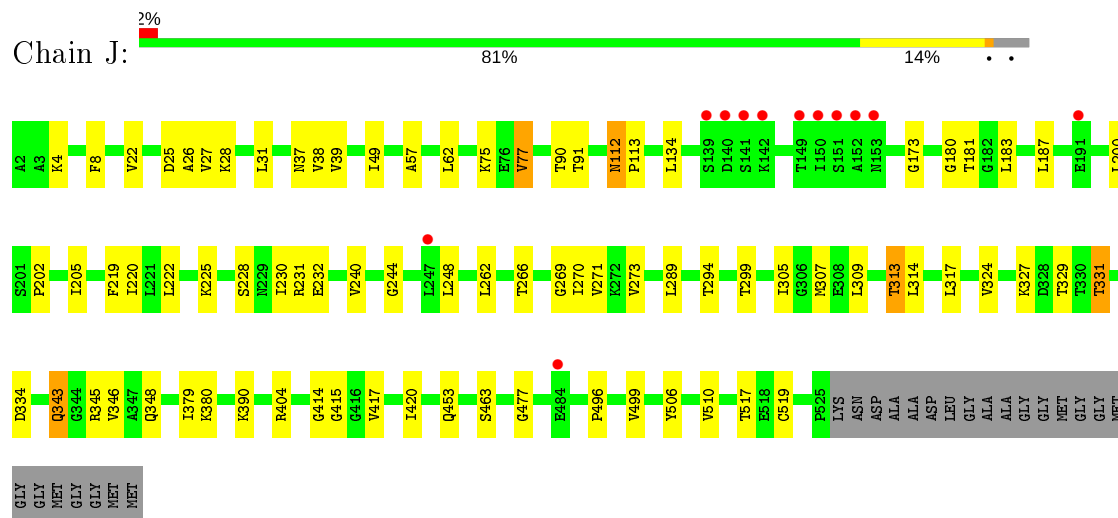


• Molecule 1: 60 kDa chaperonin

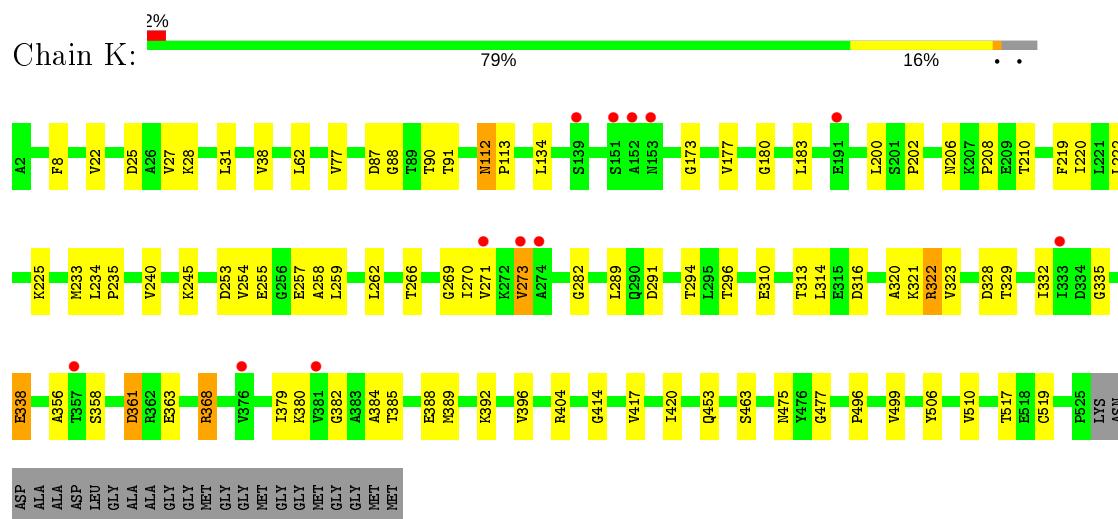


ASN
ASP
ALA
ALA
ASP
LEU
GLY
ALA
ALA
GLY
GLY
MET
MET
GLY
GLY
MET
MET

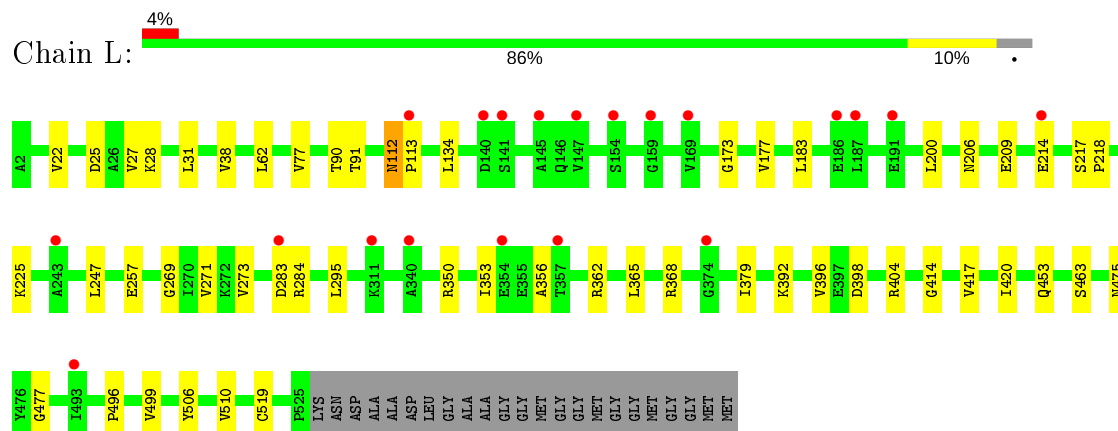
• Molecule 1: 60 kDa chaperonin



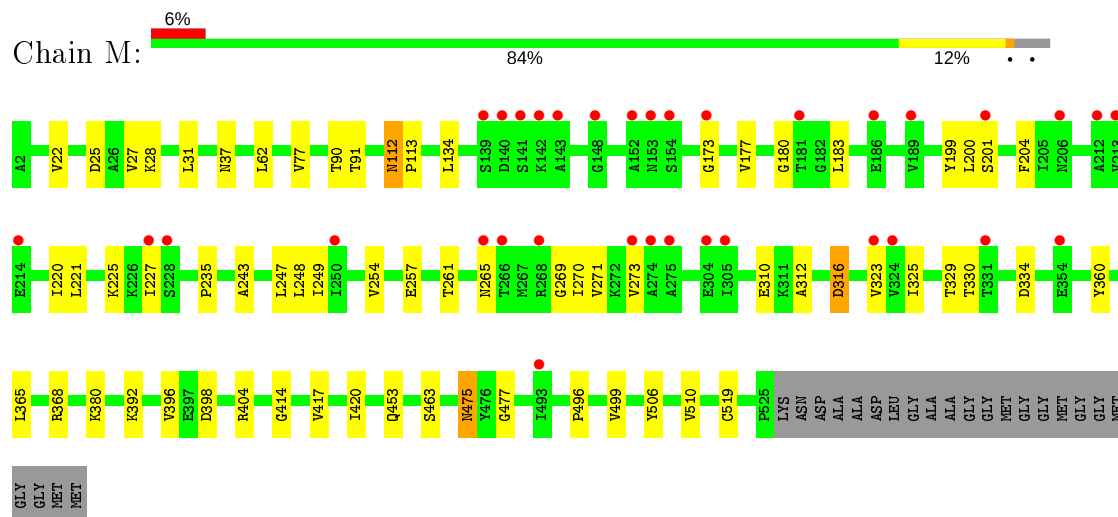
• Molecule 1: 60 kDa chaperonin



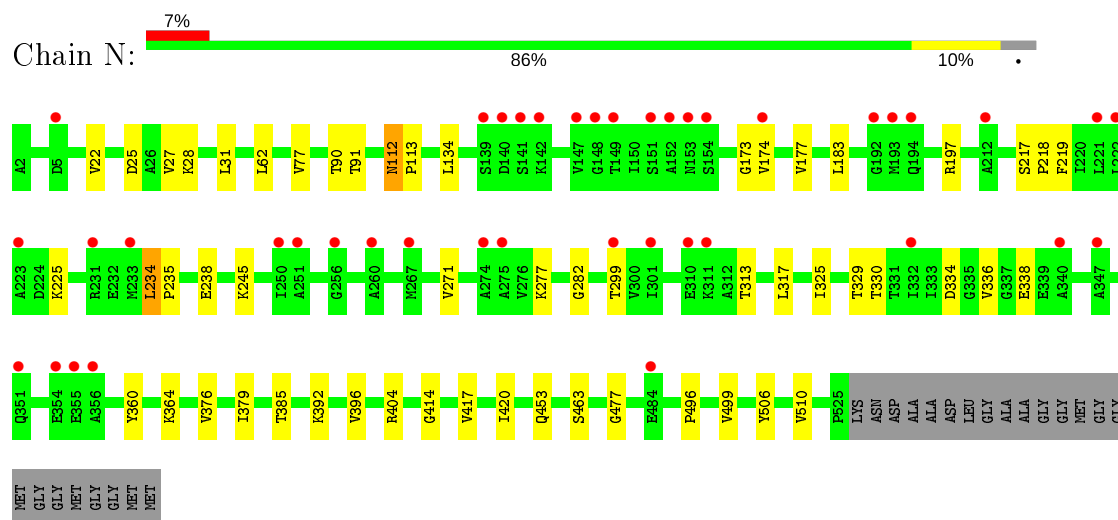
• Molecule 1: 60 kDa chaperonin



• Molecule 1: 60 kDa chaperonin



• Molecule 1: 60 kDa chaperonin



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	135.68Å 260.95Å 287.87Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.39 – 3.94 49.36 – 3.90	Depositor EDS
% Data completeness (in resolution range)	97.7 (49.39-3.94) 96.5 (49.36-3.90)	Depositor EDS
R_{merge}	0.19	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.37 (at 3.88Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.261 , 0.293 0.249 , 0.277	Depositor DCC
R_{free} test set	4528 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	115.0	Xtriage
Anisotropy	0.482	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.24 , 112.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	54464	wwPDB-VP
Average B, all atoms (Å ²)	125.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.99% 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 ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: MG, AGS, TL

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	0.38	0/3883	0.63	4/5243 (0.1%)
1	B	0.36	0/3883	0.48	0/5243
1	C	0.46	4/3883 (0.1%)	0.49	0/5243
1	D	0.43	2/3883 (0.1%)	0.50	0/5243
1	E	0.37	0/3883	0.63	4/5243 (0.1%)
1	F	0.36	0/3883	0.50	0/5243
1	G	0.37	0/3883	0.50	0/5243
1	H	0.43	2/3883 (0.1%)	0.50	0/5243
1	I	0.40	2/3883 (0.1%)	0.49	0/5243
1	J	0.36	0/3883	0.49	0/5243
1	K	0.40	2/3883 (0.1%)	0.49	0/5243
1	L	0.34	0/3883	0.47	0/5243
1	M	0.37	2/3883 (0.1%)	0.47	0/5243
1	N	0.44	3/3883 (0.1%)	0.49	1/5243 (0.0%)
All	All	0.39	17/54362 (0.0%)	0.51	9/73402 (0.0%)

The worst 5 of 17 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	232	GLU	CD-OE2	12.24	1.39	1.25
1	N	338	GLU	CD-OE2	11.57	1.38	1.25
1	H	367	GLU	CD-OE2	10.24	1.36	1.25
1	D	315	GLU	CD-OE1	9.77	1.36	1.25
1	H	367	GLU	CD-OE1	9.42	1.36	1.25

The worst 5 of 9 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	13	ARG	NE-CZ-NH2	-20.52	110.04	120.30
1	A	13	ARG	NE-CZ-NH2	-19.42	110.59	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	13	ARG	NE-CZ-NH1	18.12	129.36	120.30
1	A	13	ARG	NE-CZ-NH1	17.56	129.08	120.30
1	A	13	ARG	CD-NE-CZ	8.14	135.00	123.60

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	3855	0	3976	60	0
1	B	3855	0	3974	40	0
1	C	3855	0	3976	41	0
1	D	3855	0	3976	43	0
1	E	3855	0	3974	43	0
1	F	3855	0	3976	50	0
1	G	3855	0	3975	55	0
1	H	3855	0	3976	44	0
1	I	3855	0	3975	56	0
1	J	3855	0	3974	45	0
1	K	3855	0	3975	49	0
1	L	3855	0	3976	30	0
1	M	3855	0	3976	36	0
1	N	3855	0	3976	34	0
2	A	31	0	12	0	0
2	B	31	0	12	0	0
2	C	31	0	12	1	0
2	D	31	0	12	1	0
2	E	31	0	12	1	0
2	F	31	0	12	0	0
2	G	31	0	12	2	0
2	H	31	0	12	0	0
2	I	31	0	12	1	0
2	J	31	0	12	1	0
2	K	31	0	12	1	0
2	L	31	0	12	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	M	31	0	12	0	0
2	N	31	0	12	0	0
3	A	4	0	0	1	0
3	B	4	0	0	0	0
3	C	3	0	0	0	0
3	D	4	0	0	0	0
3	E	3	0	0	0	0
3	F	4	0	0	0	0
3	G	3	0	0	0	0
3	H	4	0	0	0	0
3	I	3	0	0	1	0
3	J	3	0	0	0	0
3	K	3	0	0	0	0
3	L	3	0	0	0	0
3	M	3	0	0	0	0
3	N	2	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
4	E	1	0	0	0	0
4	F	1	0	0	0	0
4	G	1	0	0	0	0
4	H	1	0	0	0	0
4	I	1	0	0	0	0
4	J	1	0	0	0	0
4	K	1	0	0	0	0
4	L	1	0	0	0	0
4	M	1	0	0	0	0
4	N	1	0	0	0	0
All	All	54464	0	55823	586	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.

The worst 5 of 586 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:M:183:LEU:HD22	1:N:360:TYR:CB	1.96	0.95
1:A:87:ASP:OD2	1:A:499:VAL:HG21	1.69	0.91
1:C:87:ASP:OD2	1:C:499:VAL:HG21	1.75	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:183:LEU:HD22	1:N:360:TYR:HB3	1.54	0.86
1:H:360:TYR:CG	1:N:183:LEU:HD22	2.13	0.84

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	522/547 (95%)	494 (95%)	23 (4%)	5 (1%)	15	52
1	B	522/547 (95%)	490 (94%)	25 (5%)	7 (1%)	12	47
1	C	522/547 (95%)	490 (94%)	27 (5%)	5 (1%)	15	52
1	D	522/547 (95%)	487 (93%)	30 (6%)	5 (1%)	15	52
1	E	522/547 (95%)	493 (94%)	22 (4%)	7 (1%)	12	47
1	F	522/547 (95%)	496 (95%)	19 (4%)	7 (1%)	12	47
1	G	522/547 (95%)	490 (94%)	28 (5%)	4 (1%)	19	57
1	H	522/547 (95%)	496 (95%)	18 (3%)	8 (2%)	10	45
1	I	522/547 (95%)	495 (95%)	25 (5%)	2 (0%)	34	70
1	J	522/547 (95%)	496 (95%)	21 (4%)	5 (1%)	15	52
1	K	522/547 (95%)	485 (93%)	33 (6%)	4 (1%)	19	57
1	L	522/547 (95%)	501 (96%)	20 (4%)	1 (0%)	47	79
1	M	522/547 (95%)	495 (95%)	22 (4%)	5 (1%)	15	52
1	N	522/547 (95%)	483 (92%)	32 (6%)	7 (1%)	12	47
All	All	7308/7658 (95%)	6891 (94%)	345 (5%)	72 (1%)	15	52

5 of 72 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	244	GLY
1	A	375	GLY
1	B	205	ILE
1	H	322	ARG
1	H	336	VAL

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	404/414 (98%)	386 (96%)	18 (4%)	27	54
1	B	404/414 (98%)	392 (97%)	12 (3%)	41	64
1	C	404/414 (98%)	392 (97%)	12 (3%)	41	64
1	D	404/414 (98%)	387 (96%)	17 (4%)	30	56
1	E	404/414 (98%)	392 (97%)	12 (3%)	41	64
1	F	404/414 (98%)	389 (96%)	15 (4%)	34	59
1	G	404/414 (98%)	385 (95%)	19 (5%)	26	53
1	H	404/414 (98%)	393 (97%)	11 (3%)	44	66
1	I	404/414 (98%)	389 (96%)	15 (4%)	34	59
1	J	404/414 (98%)	388 (96%)	16 (4%)	31	57
1	K	404/414 (98%)	385 (95%)	19 (5%)	26	53
1	L	404/414 (98%)	393 (97%)	11 (3%)	44	66
1	M	404/414 (98%)	393 (97%)	11 (3%)	44	66
1	N	404/414 (98%)	397 (98%)	7 (2%)	60	78
All	All	5656/5796 (98%)	5461 (97%)	195 (3%)	37	61

5 of 195 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	G	134	LEU
1	H	217	SER
1	M	199	TYR

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Mol	Chain	Res	Type
1	G	217	SER
1	G	331	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 21 such sidechains are listed below:

Mol	Chain	Res	Type
1	F	112	ASN
1	H	112	ASN
1	J	366	GLN
1	E	351	GLN
1	L	206	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 74 ligands modelled in this entry, 60 are monoatomic - leaving 14 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	AGS	K	549	4	26,33,33	1.50	2 (7%)	26,52,52	1.55	4 (15%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	AGS	H	549	4	26,33,33	1.14	1 (3%)	26,52,52	1.54	4 (15%)
2	AGS	I	549	4	26,33,33	1.26	1 (3%)	26,52,52	1.52	4 (15%)
2	AGS	J	549	4	26,33,33	1.64	1 (3%)	26,52,52	1.56	4 (15%)
2	AGS	C	549	4	26,33,33	1.50	1 (3%)	26,52,52	1.55	4 (15%)
2	AGS	L	549	4	26,33,33	1.47	2 (7%)	26,52,52	1.48	4 (15%)
2	AGS	M	549	4	26,33,33	1.03	3 (11%)	26,52,52	1.51	3 (11%)
2	AGS	N	549	4	26,33,33	1.75	1 (3%)	26,52,52	1.51	4 (15%)
2	AGS	G	549	4	26,33,33	1.43	1 (3%)	26,52,52	1.55	4 (15%)
2	AGS	A	549	4	26,33,33	1.30	1 (3%)	26,52,52	1.55	4 (15%)
2	AGS	B	549	4	26,33,33	1.15	2 (7%)	26,52,52	1.54	4 (15%)
2	AGS	D	549	4	26,33,33	1.20	2 (7%)	26,52,52	1.57	4 (15%)
2	AGS	E	549	4	26,33,33	1.24	2 (7%)	26,52,52	1.55	4 (15%)
2	AGS	F	549	4	26,33,33	1.39	2 (7%)	26,52,52	1.55	4 (15%)

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	AGS	K	549	4	-	4/17/38/38	0/3/3/3
2	AGS	H	549	4	-	4/17/38/38	0/3/3/3
2	AGS	I	549	4	-	5/17/38/38	0/3/3/3
2	AGS	J	549	4	-	4/17/38/38	0/3/3/3
2	AGS	C	549	4	-	5/17/38/38	0/3/3/3
2	AGS	L	549	4	-	4/17/38/38	0/3/3/3
2	AGS	M	549	4	-	5/17/38/38	0/3/3/3
2	AGS	N	549	4	-	4/17/38/38	0/3/3/3
2	AGS	G	549	4	-	4/17/38/38	0/3/3/3
2	AGS	A	549	4	-	5/17/38/38	0/3/3/3
2	AGS	B	549	4	-	4/17/38/38	0/3/3/3
2	AGS	D	549	4	-	4/17/38/38	0/3/3/3
2	AGS	E	549	4	-	5/17/38/38	0/3/3/3
2	AGS	F	549	4	-	4/17/38/38	0/3/3/3

The worst 5 of 22 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	N	549	AGS	PG-S1G	7.36	2.06	1.90
2	J	549	AGS	PG-S1G	6.90	2.05	1.90
2	K	549	AGS	PG-S1G	5.93	2.03	1.90
2	C	549	AGS	PG-S1G	5.72	2.03	1.90
2	G	549	AGS	PG-S1G	5.51	2.02	1.90

The worst 5 of 55 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	549	AGS	N3-C2-N1	-4.78	121.21	128.68
2	K	549	AGS	N3-C2-N1	-4.74	121.27	128.68
2	C	549	AGS	N3-C2-N1	-4.73	121.29	128.68
2	J	549	AGS	N3-C2-N1	-4.72	121.30	128.68
2	E	549	AGS	N3-C2-N1	-4.67	121.38	128.68

There are no chirality outliers.

5 of 61 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	J	549	AGS	C5'-O5'-PA-O2A
2	B	549	AGS	C5'-O5'-PA-O2A
2	G	549	AGS	C5'-O5'-PA-O2A
2	K	549	AGS	C5'-O5'-PA-O2A
2	I	549	AGS	C5'-O5'-PA-O2A

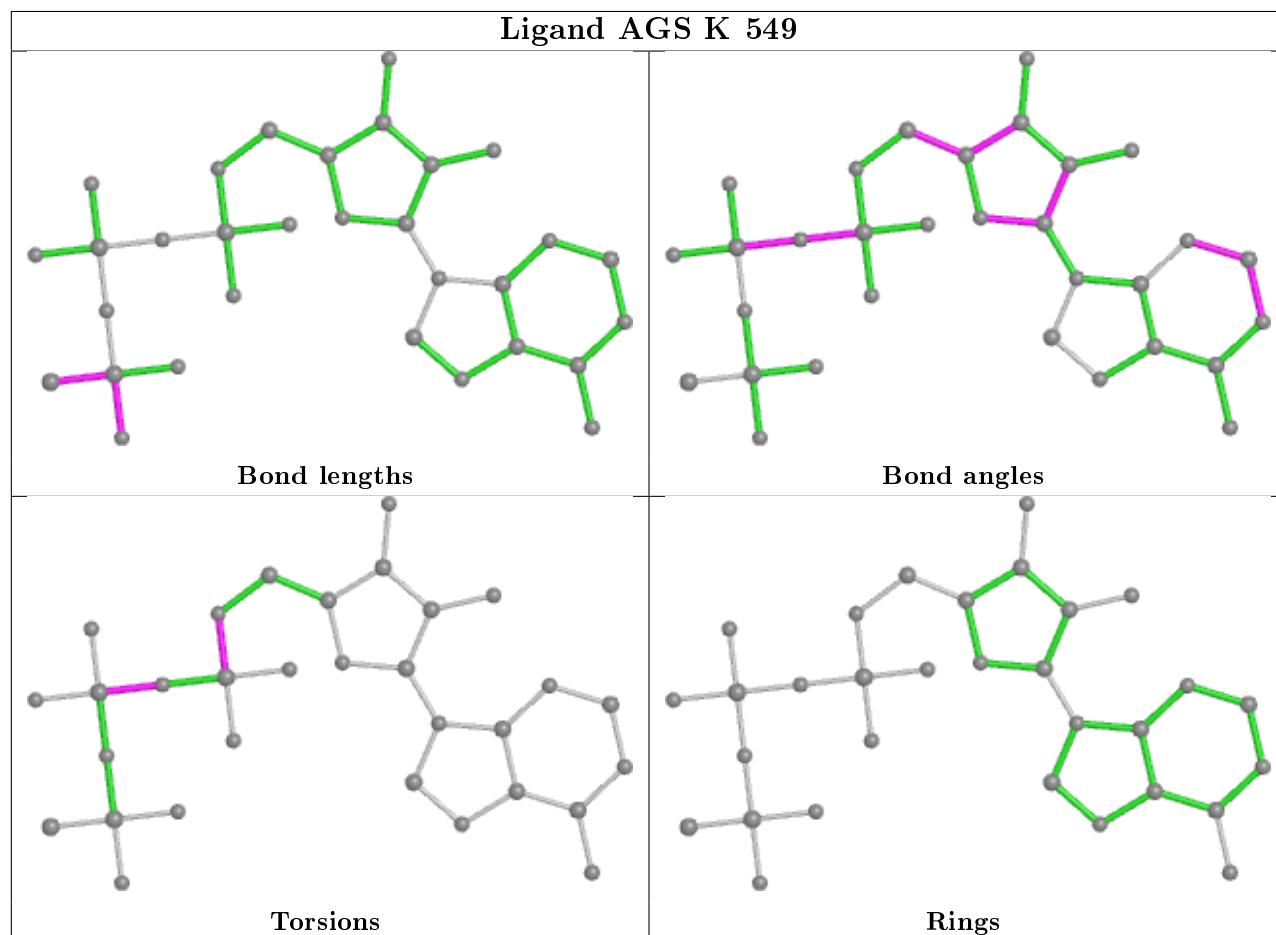
There are no ring outliers.

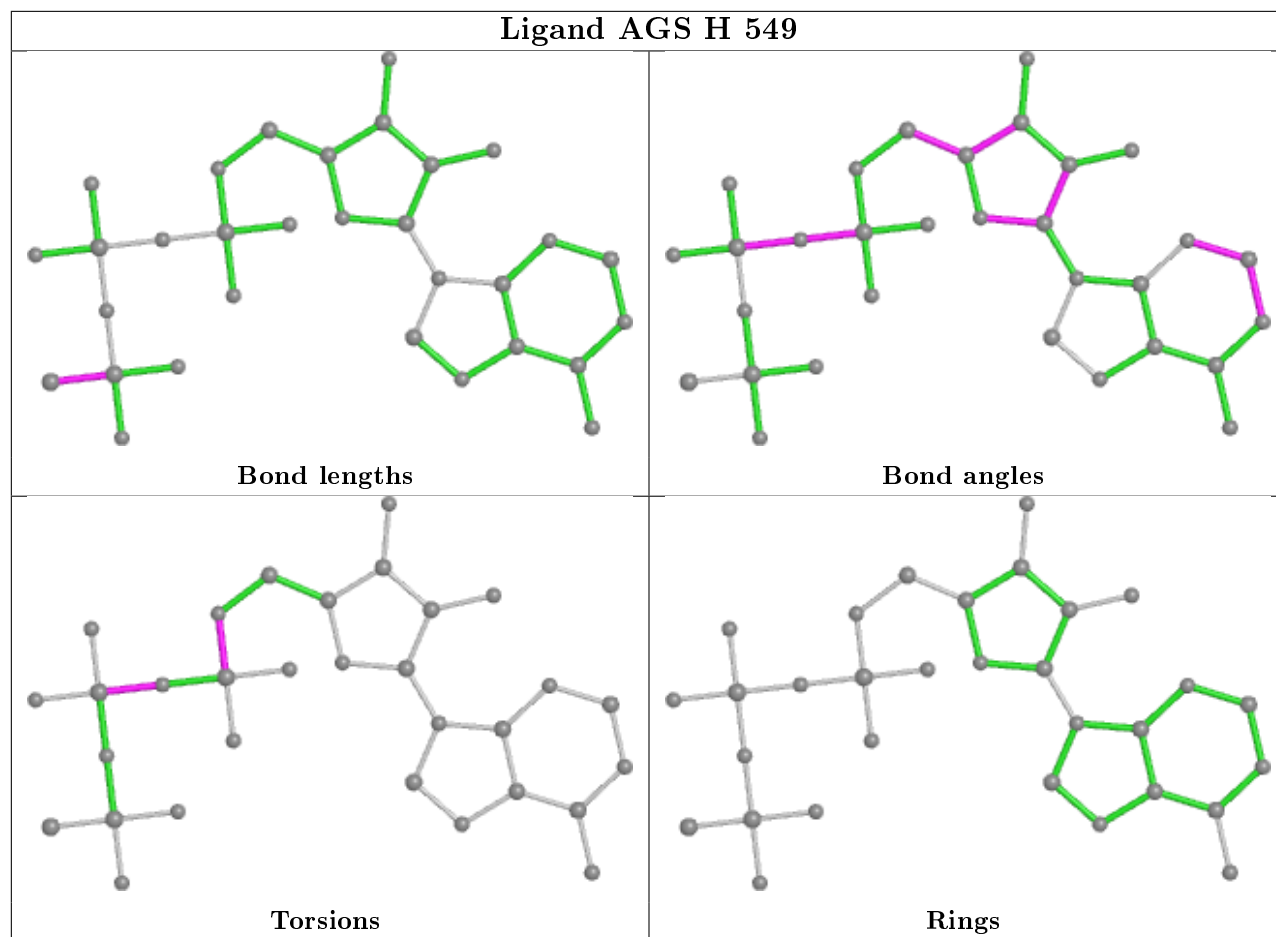
7 monomers are involved in 8 short contacts:

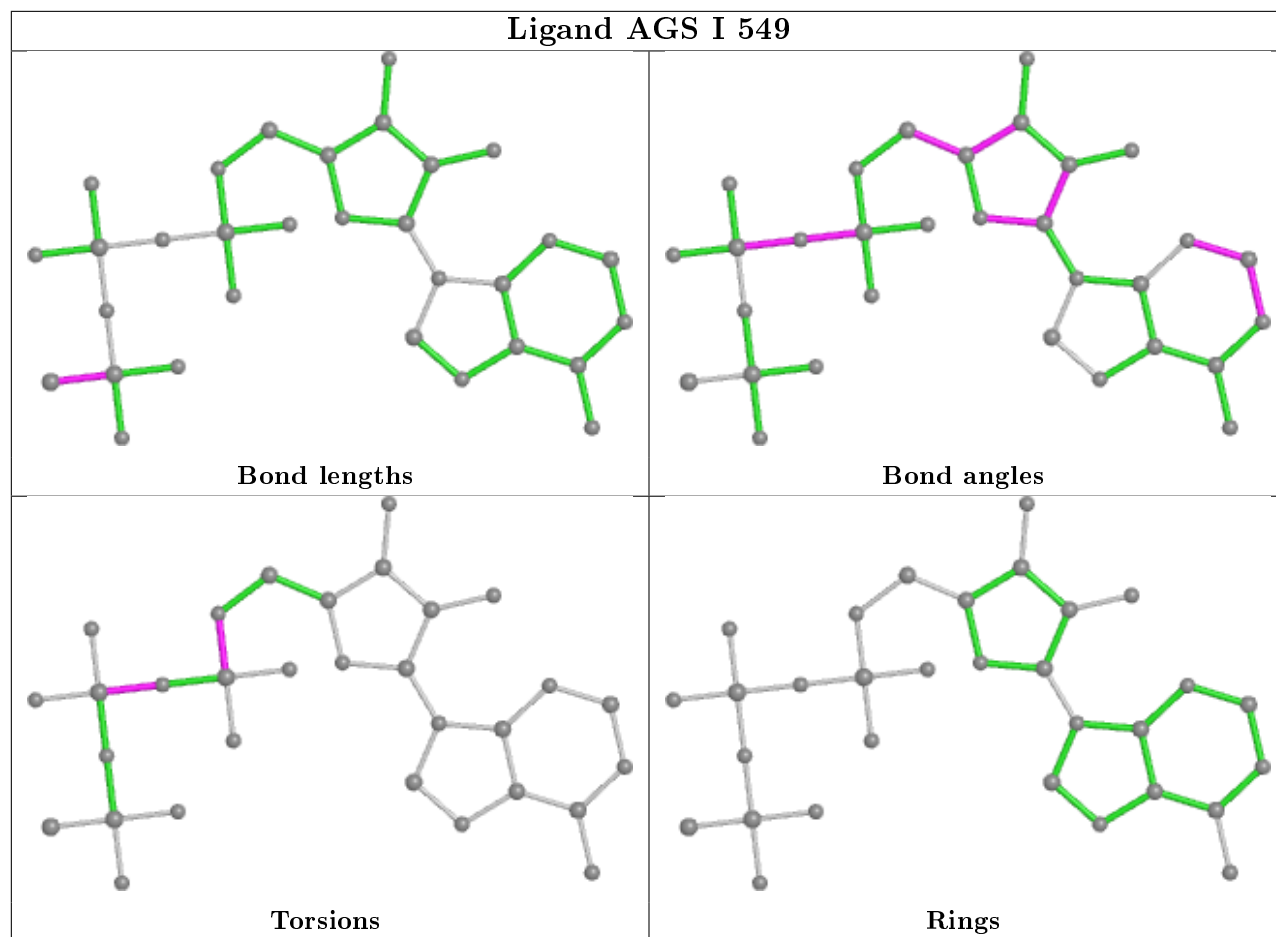
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	K	549	AGS	1	0
2	I	549	AGS	1	0
2	J	549	AGS	1	0
2	C	549	AGS	1	0
2	G	549	AGS	2	0
2	D	549	AGS	1	0
2	E	549	AGS	1	0

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.

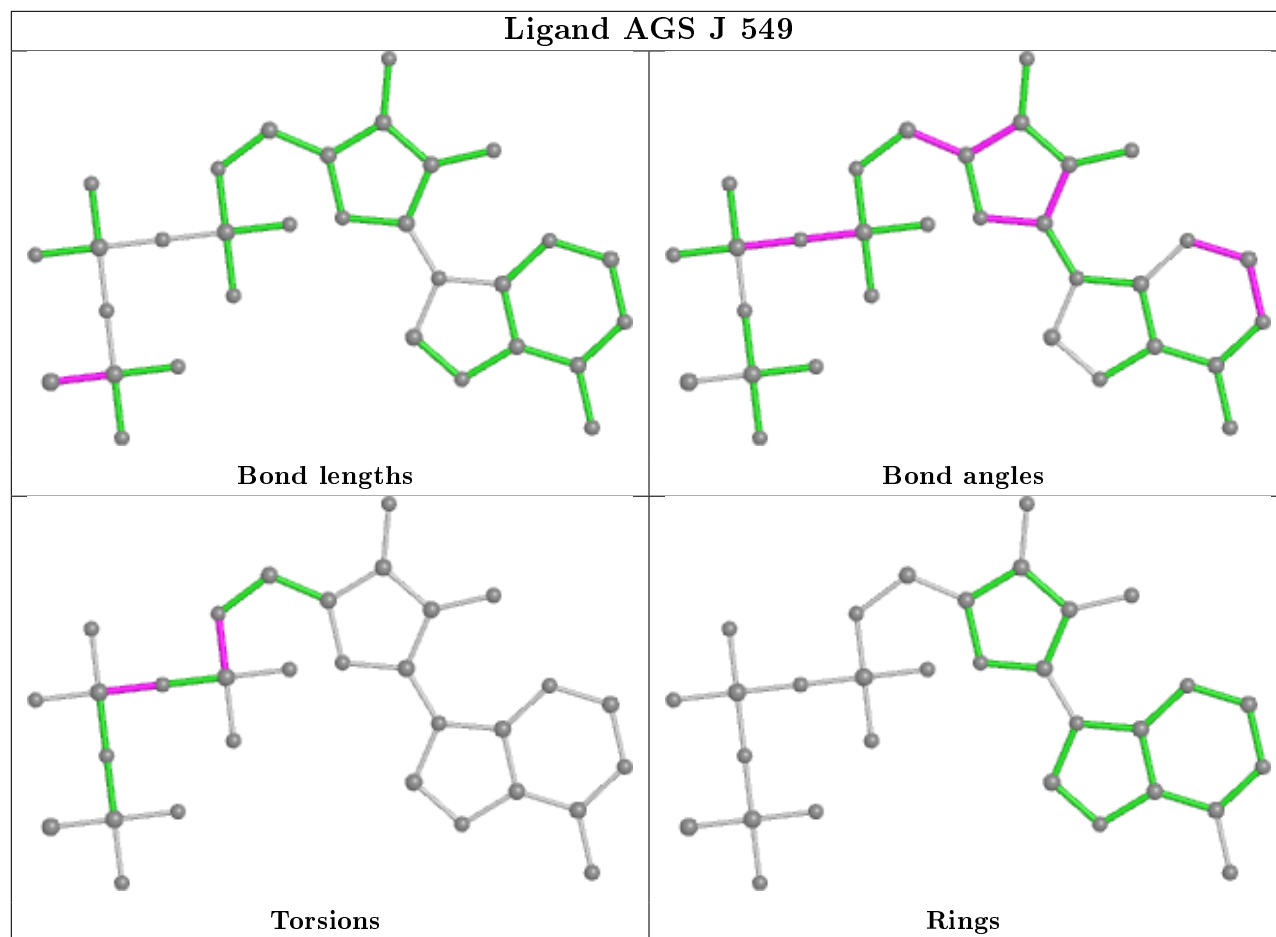
Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD 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.

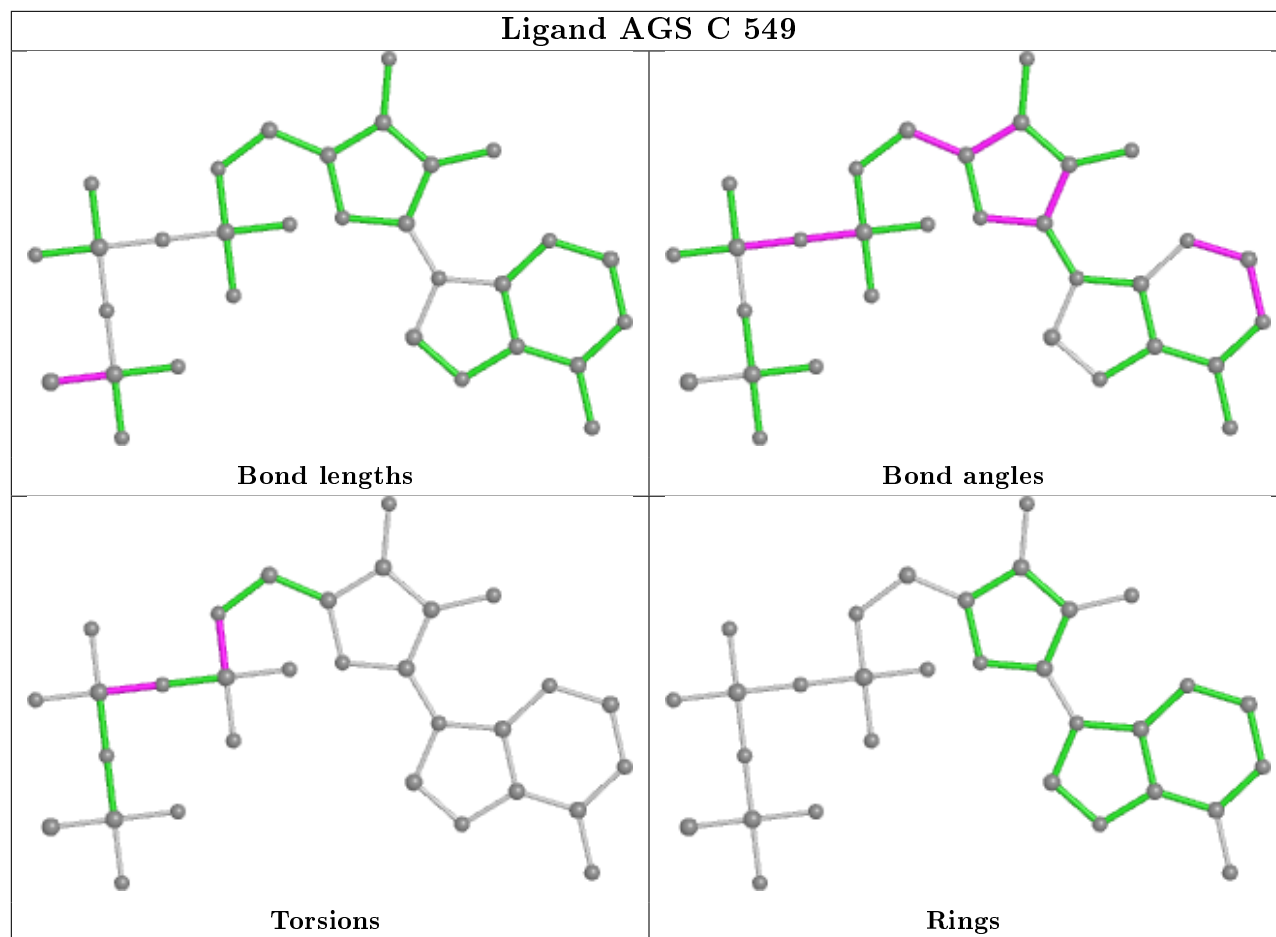


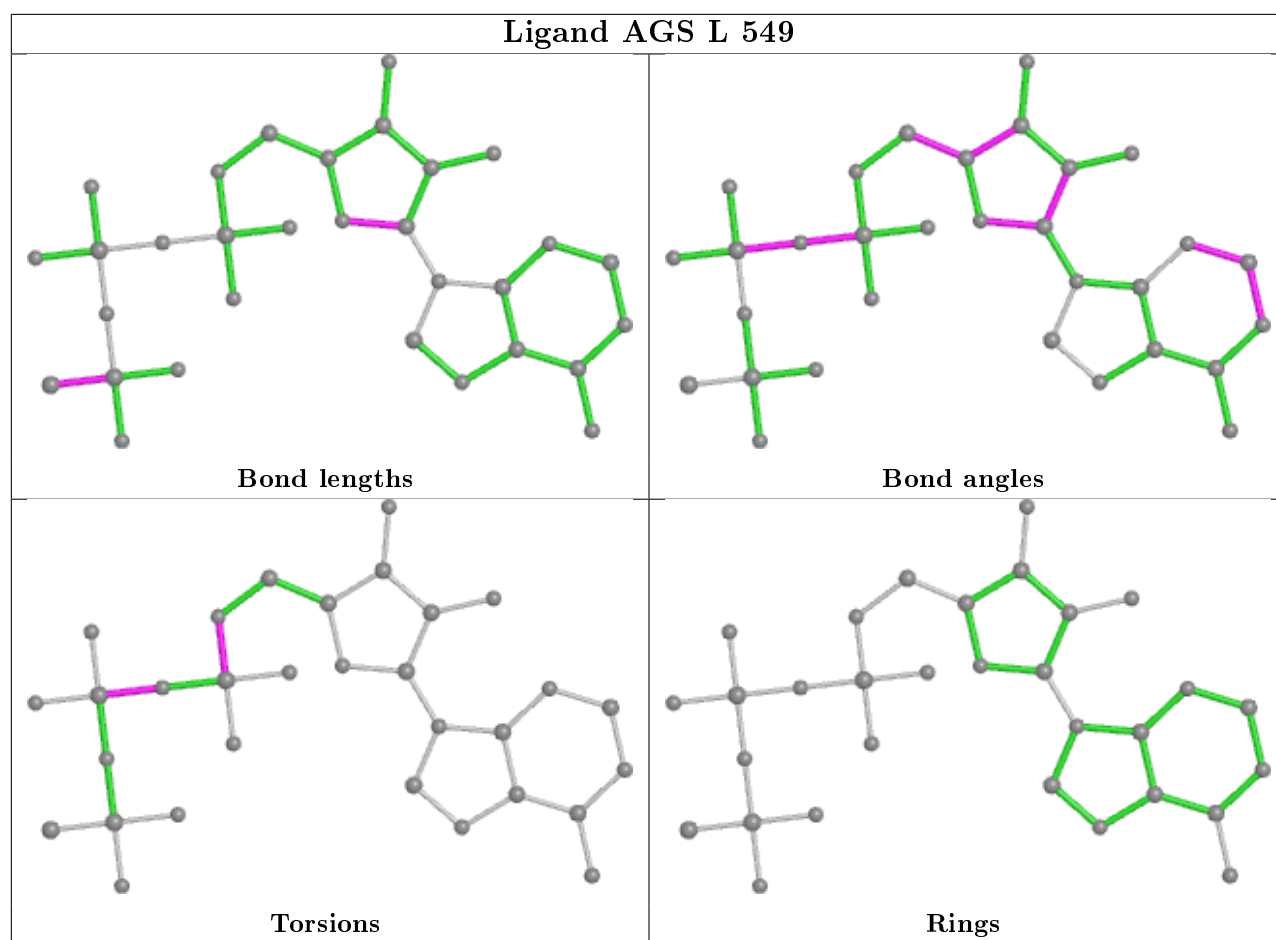


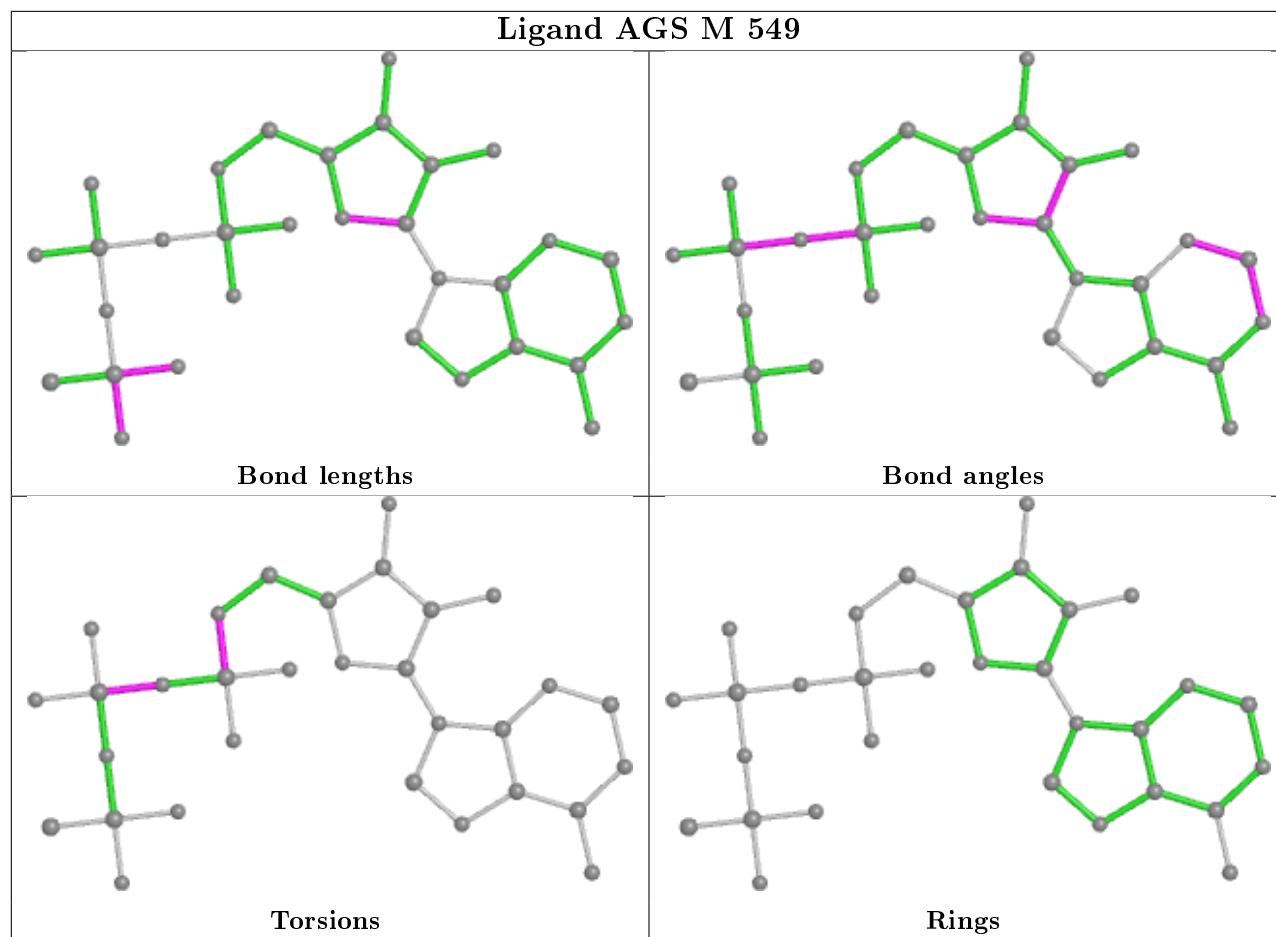


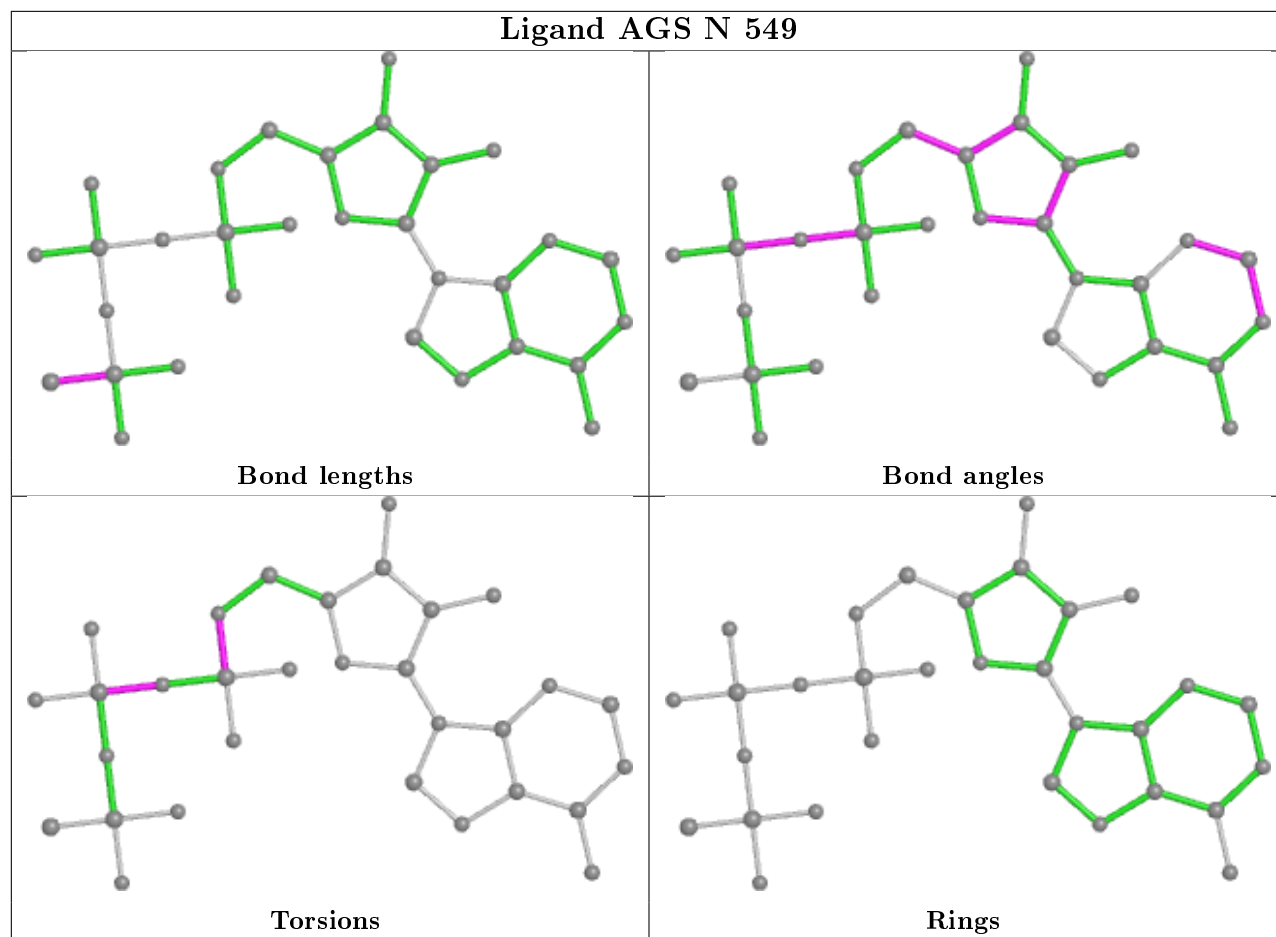
Ligand AGS J 549

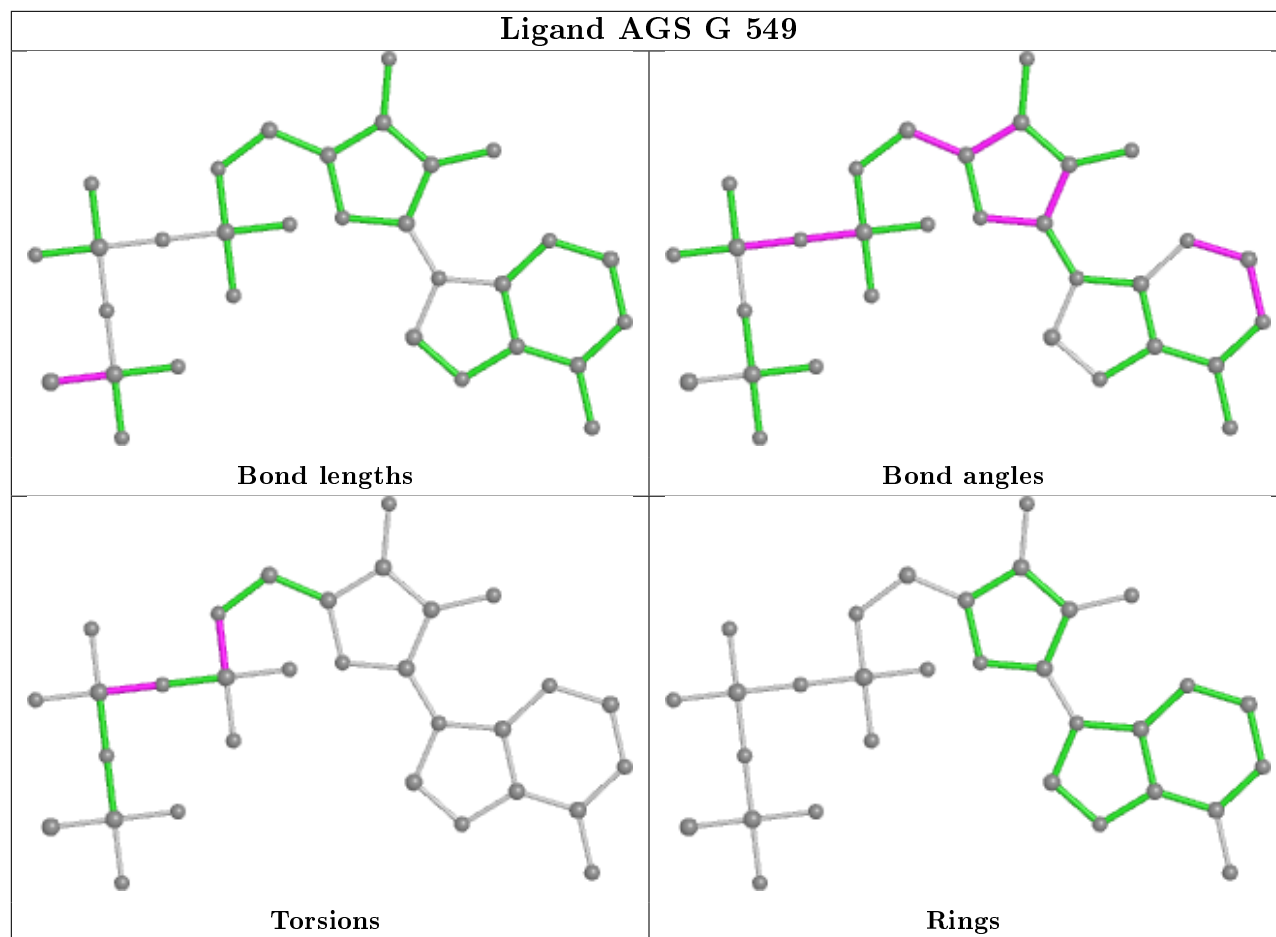


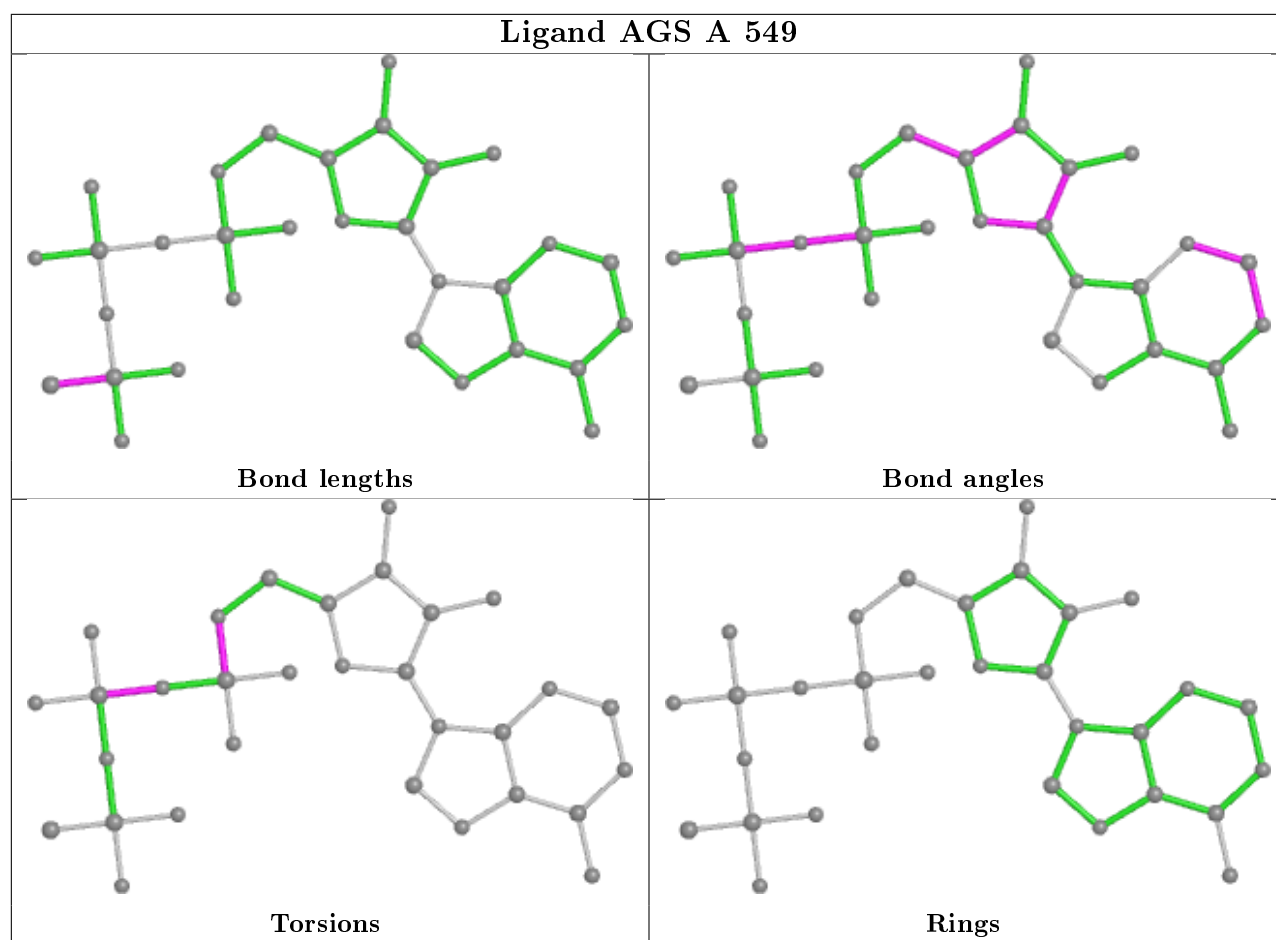


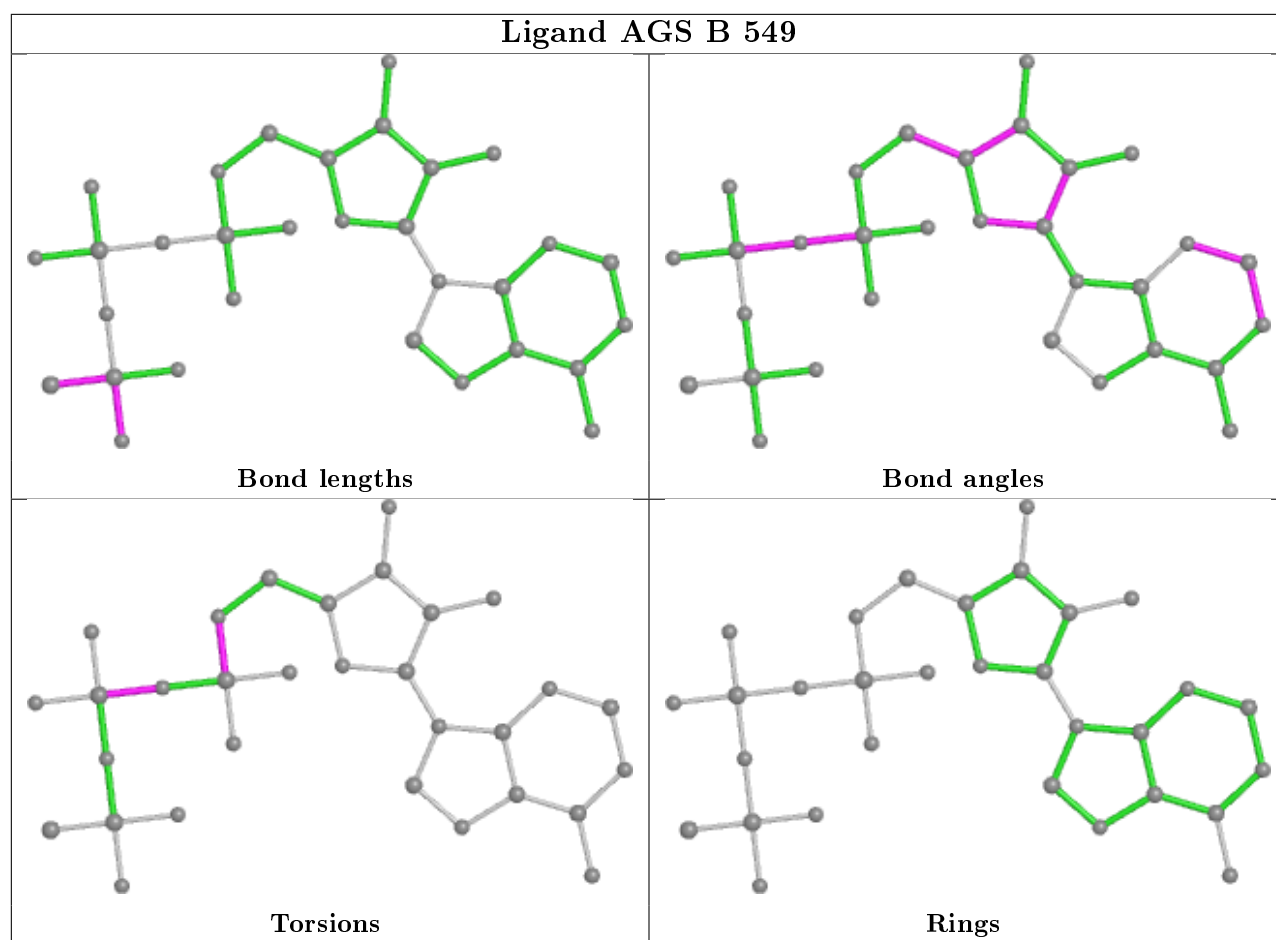


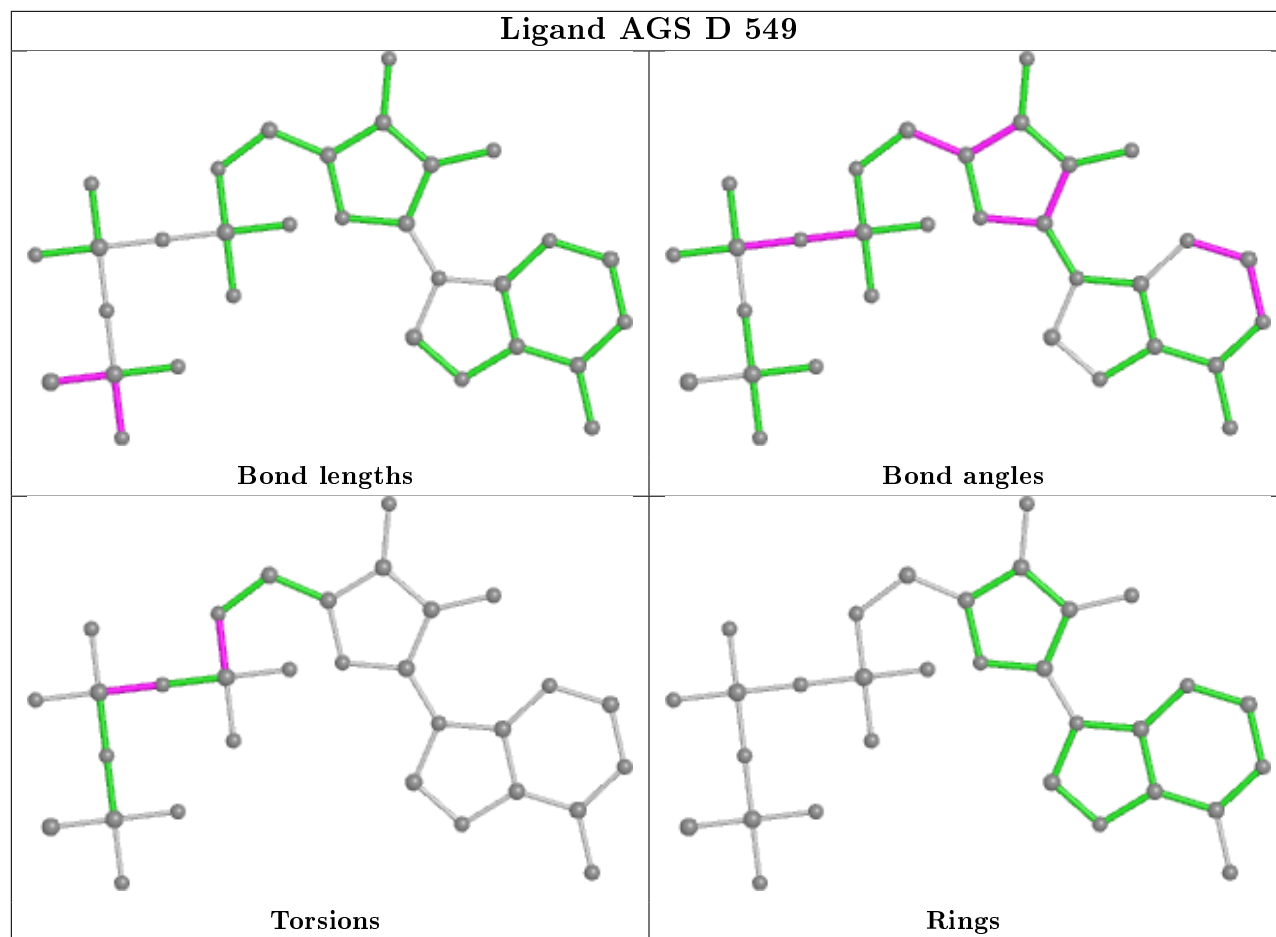


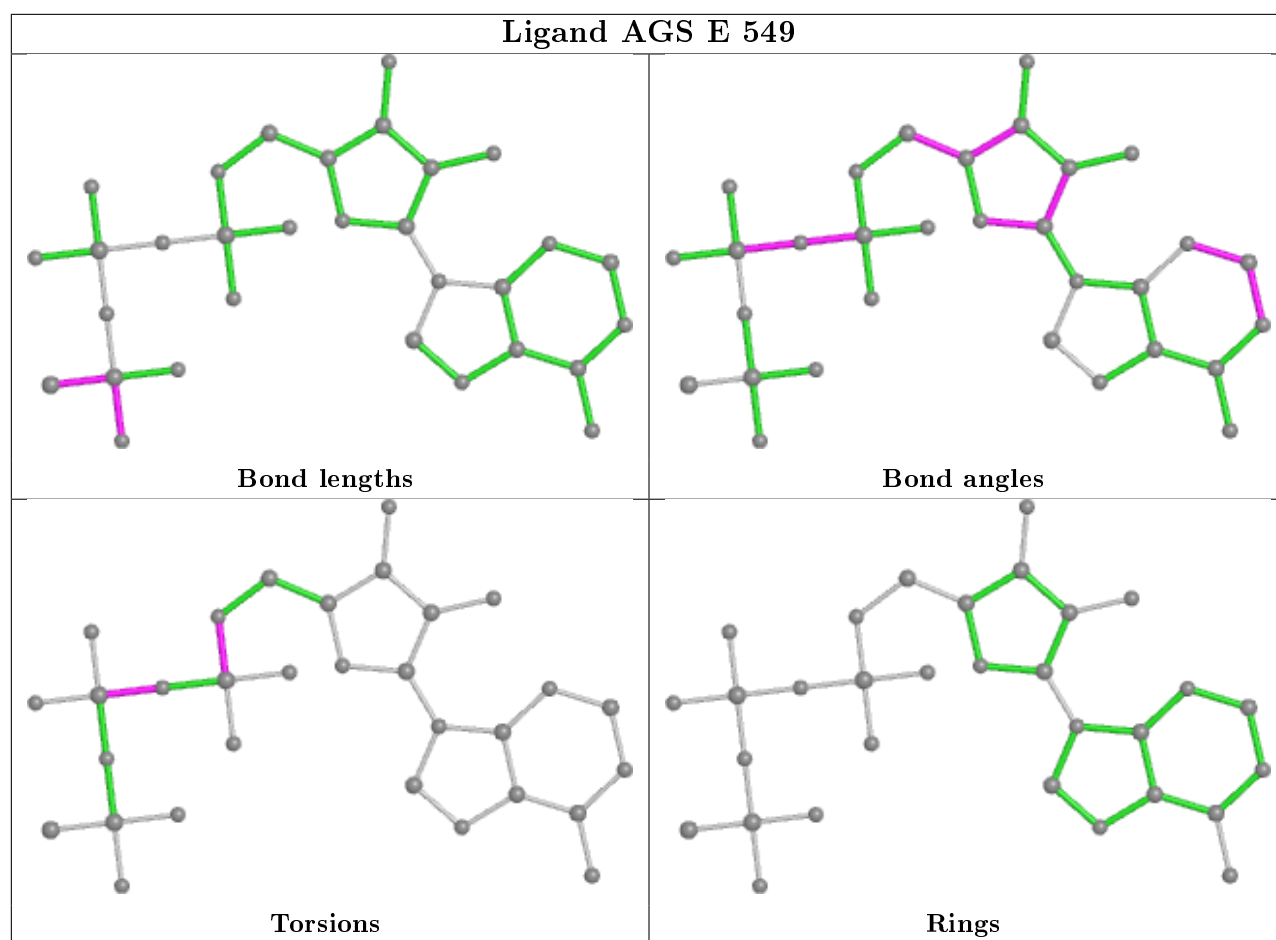


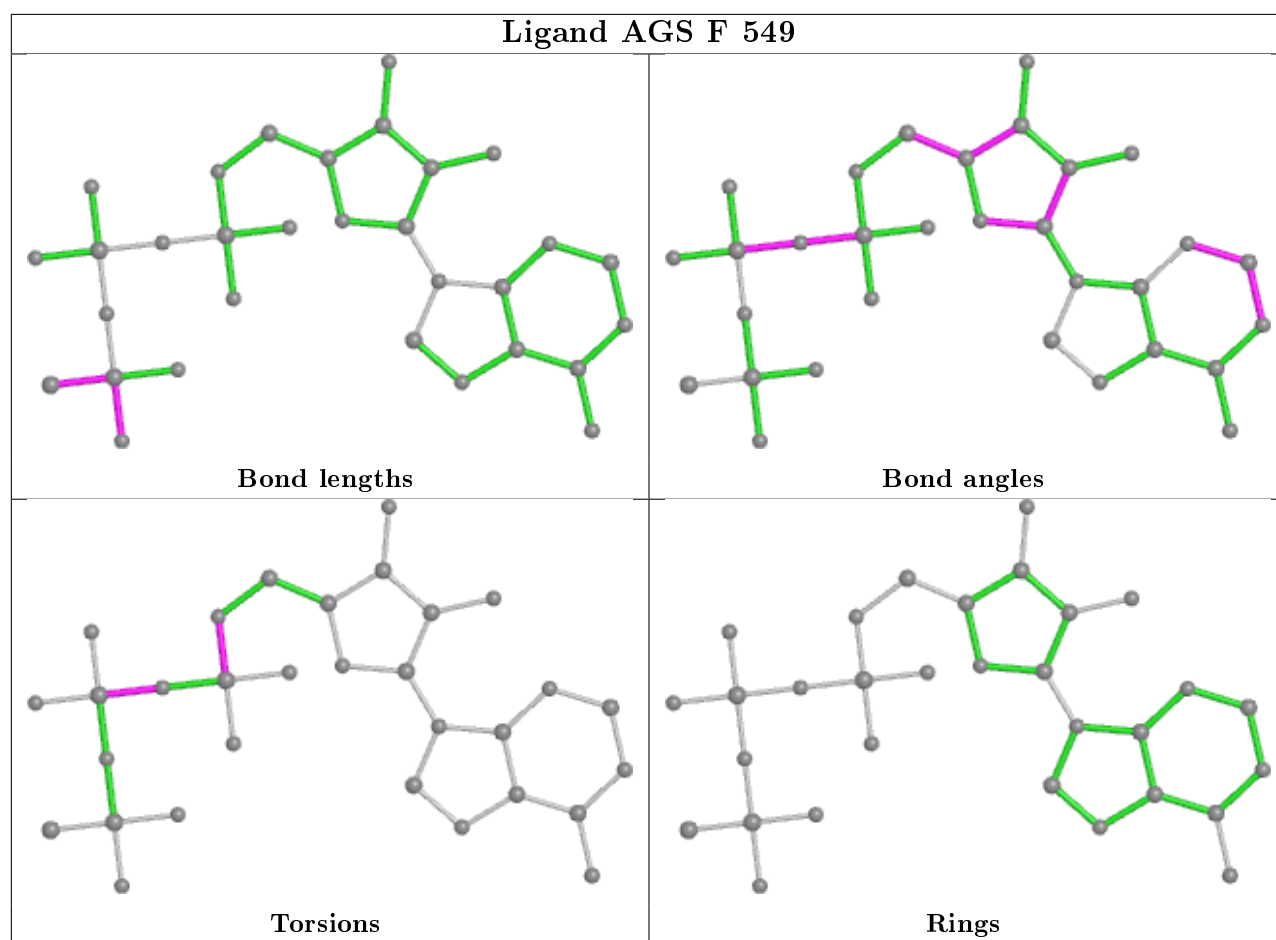












5.7 Other polymers [i](#)

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	524/547 (95%)	-0.19	2 (0%) 92 87	125, 125, 125, 125	0
1	B	524/547 (95%)	-0.17	1 (0%) 95 93	125, 125, 125, 125	0
1	C	524/547 (95%)	-0.08	12 (2%) 60 51	125, 125, 125, 125	0
1	D	524/547 (95%)	-0.12	6 (1%) 80 72	125, 125, 125, 125	0
1	E	524/547 (95%)	-0.02	12 (2%) 60 51	125, 125, 125, 125	0
1	F	524/547 (95%)	-0.11	9 (1%) 70 60	125, 125, 125, 125	0
1	G	524/547 (95%)	-0.14	8 (1%) 73 64	125, 125, 125, 125	0
1	H	524/547 (95%)	0.00	15 (2%) 51 41	125, 125, 125, 125	0
1	I	524/547 (95%)	-0.12	12 (2%) 60 51	125, 125, 125, 125	0
1	J	524/547 (95%)	-0.02	12 (2%) 60 51	125, 125, 125, 125	0
1	K	524/547 (95%)	-0.03	12 (2%) 60 51	125, 125, 125, 125	0
1	L	524/547 (95%)	0.05	20 (3%) 40 32	125, 125, 125, 125	0
1	M	524/547 (95%)	0.23	34 (6%) 18 15	125, 125, 125, 125	0
1	N	524/547 (95%)	0.40	41 (7%) 13 11	125, 125, 125, 125	0
All	All	7336/7658 (95%)	-0.02	196 (2%) 54 44	125, 125, 125, 125	0

The worst 5 of 196 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	N	223	ALA	11.9
1	I	141	SER	7.3
1	N	148	GLY	7.2
1	M	153	ASN	7.1
1	E	152	ALA	6.8

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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(\AA^2)	Q<0.9
3	TL	F	551	1/1	0.73	0.10	125,125,125,125	1
3	TL	G	551	1/1	0.73	0.14	125,125,125,125	1
3	TL	J	551	1/1	0.75	0.18	125,125,125,125	1
2	AGS	E	549	31/31	0.76	0.27	125,125,125,125	0
3	TL	H	554	1/1	0.76	0.15	125,125,125,125	1
4	MG	A	553	1/1	0.79	0.25	125,125,125,125	0
2	AGS	L	549	31/31	0.79	0.25	125,125,125,125	0
2	AGS	C	549	31/31	0.80	0.23	125,125,125,125	0
2	AGS	N	549	31/31	0.80	0.23	125,125,125,125	0
4	MG	C	553	1/1	0.81	0.41	125,125,125,125	0
2	AGS	J	549	31/31	0.82	0.27	125,125,125,125	0
2	AGS	M	549	31/31	0.83	0.21	125,125,125,125	0
2	AGS	H	549	31/31	0.83	0.23	125,125,125,125	0
4	MG	I	553	1/1	0.84	0.23	125,125,125,125	0
2	AGS	F	549	31/31	0.84	0.24	125,125,125,125	0
4	MG	G	553	1/1	0.84	0.22	125,125,125,125	0
3	TL	E	550	1/1	0.85	0.13	125,125,125,125	1
3	TL	L	551	1/1	0.85	0.16	125,125,125,125	1
4	MG	H	553	1/1	0.86	0.24	125,125,125,125	0
2	AGS	A	549	31/31	0.86	0.21	125,125,125,125	0
2	AGS	B	549	31/31	0.86	0.19	125,125,125,125	0
4	MG	N	552	1/1	0.86	0.32	125,125,125,125	0
3	TL	B	552	1/1	0.87	0.18	125,125,125,125	1
3	TL	K	552	1/1	0.87	0.18	125,125,125,125	1
2	AGS	G	549	31/31	0.87	0.22	125,125,125,125	0
3	TL	K	550	1/1	0.88	0.10	125,125,125,125	1
3	TL	A	551	1/1	0.88	0.17	125,125,125,125	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	AGS	D	549	31/31	0.88	0.18	125,125,125,125	0
2	AGS	K	549	31/31	0.88	0.20	125,125,125,125	0
4	MG	F	553	1/1	0.88	0.25	125,125,125,125	0
3	TL	L	550	1/1	0.89	0.15	125,125,125,125	1
3	TL	C	551	1/1	0.89	0.08	125,125,125,125	1
4	MG	D	553	1/1	0.90	0.21	125,125,125,125	0
2	AGS	I	549	31/31	0.90	0.17	125,125,125,125	0
3	TL	D	552	1/1	0.91	0.05	125,125,125,125	1
4	MG	L	553	1/1	0.92	0.21	125,125,125,125	0
3	TL	N	551	1/1	0.93	0.20	125,125,125,125	1
3	TL	H	552	1/1	0.93	0.16	125,125,125,125	1
3	TL	D	551	1/1	0.93	0.12	125,125,125,125	1
3	TL	H	551	1/1	0.93	0.07	125,125,125,125	1
3	TL	G	552	1/1	0.94	0.12	125,125,125,125	1
3	TL	J	550	1/1	0.94	0.14	125,125,125,125	1
4	MG	M	553	1/1	0.94	0.14	125,125,125,125	0
3	TL	N	550	1/1	0.94	0.07	125,125,125,125	1
3	TL	C	550	1/1	0.94	0.09	125,125,125,125	1
3	TL	H	550	1/1	0.94	0.07	125,125,125,125	1
3	TL	E	551	1/1	0.94	0.06	125,125,125,125	1
4	MG	J	553	1/1	0.94	0.13	125,125,125,125	0
3	TL	A	552	1/1	0.94	0.11	125,125,125,125	1
3	TL	L	552	1/1	0.95	0.04	125,125,125,125	1
4	MG	K	553	1/1	0.95	0.12	125,125,125,125	0
3	TL	G	550	1/1	0.95	0.11	125,125,125,125	1
4	MG	E	553	1/1	0.95	0.34	125,125,125,125	0
3	TL	I	551	1/1	0.95	0.13	125,125,125,125	1
3	TL	M	552	1/1	0.95	0.20	125,125,125,125	1
3	TL	B	550	1/1	0.95	0.11	125,125,125,125	1
4	MG	B	553	1/1	0.95	0.16	125,125,125,125	0
3	TL	B	551	1/1	0.96	0.14	125,125,125,125	1
3	TL	A	550	1/1	0.96	0.08	125,125,125,125	1
3	TL	M	550	1/1	0.96	0.10	125,125,125,125	1
3	TL	M	551	1/1	0.96	0.09	125,125,125,125	1
3	TL	D	550	1/1	0.96	0.11	125,125,125,125	1
3	TL	C	552	1/1	0.96	0.25	125,125,125,125	1
3	TL	F	552	1/1	0.97	0.08	125,125,125,125	1
3	TL	K	551	1/1	0.98	0.11	125,125,125,125	1
3	TL	F	554	1/1	0.98	0.17	125,125,125,125	1
3	TL	I	552	1/1	0.98	0.06	125,125,125,125	1
3	TL	I	550	1/1	0.98	0.08	125,125,125,125	1
3	TL	J	552	1/1	0.98	0.04	125,125,125,125	1

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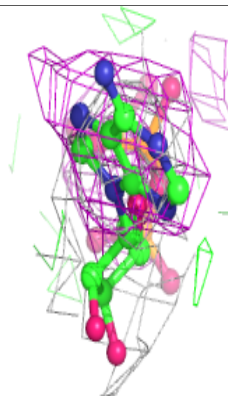
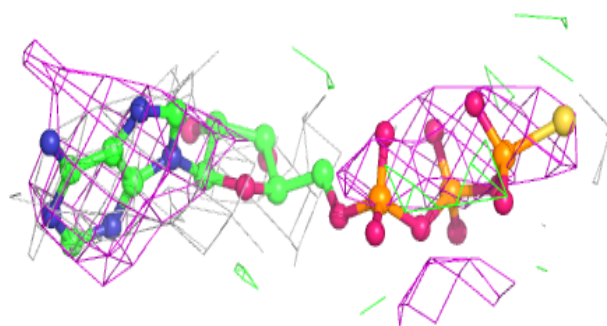
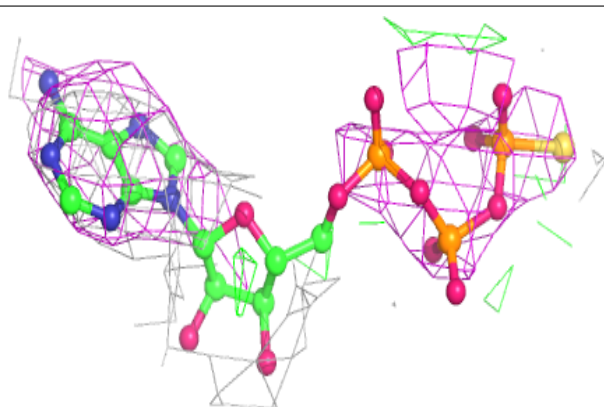
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	TL	A	554	1/1	0.99	0.20	125,125,125,125	1
3	TL	E	552	1/1	0.99	0.25	125,125,125,125	1
3	TL	F	550	1/1	0.99	0.21	125,125,125,125	1
3	TL	D	1	1/1	0.99	0.19	125,125,125,125	1
3	TL	B	554	1/1	1.00	0.19	125,125,125,125	1

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.

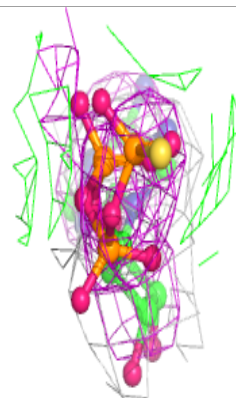
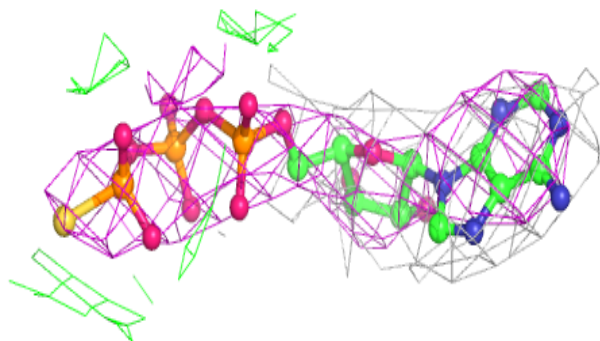
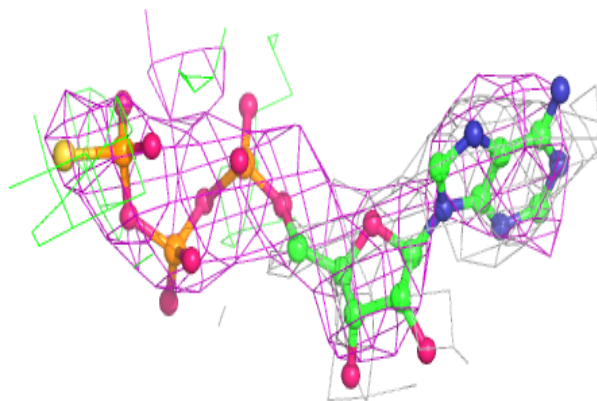
Electron density around AGS E 549:

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

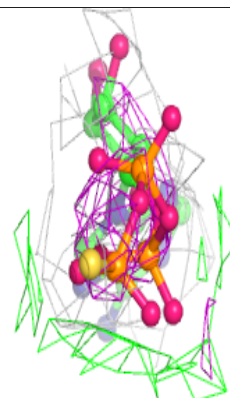
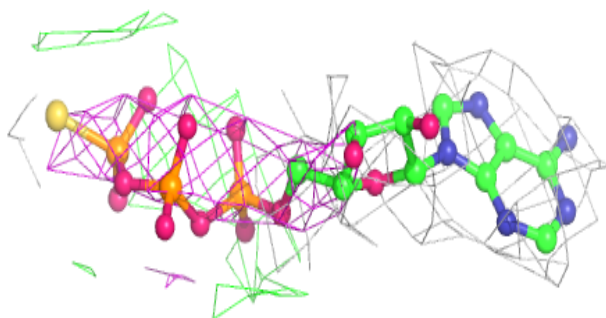
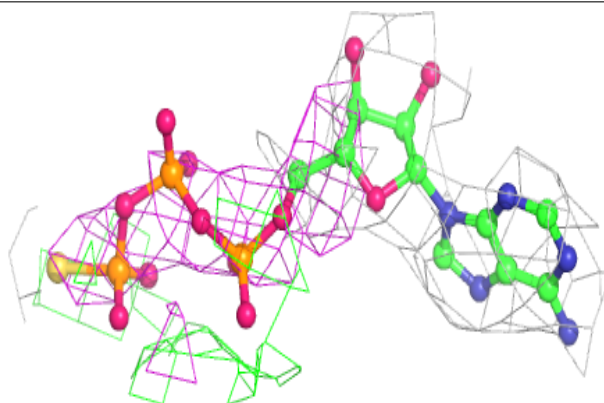


Electron density around AGS L 549:

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

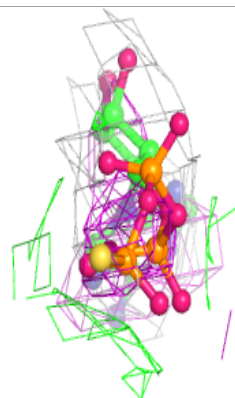
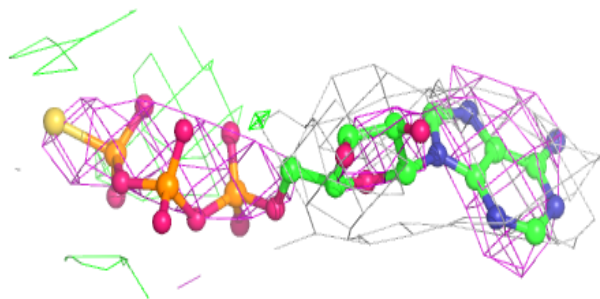
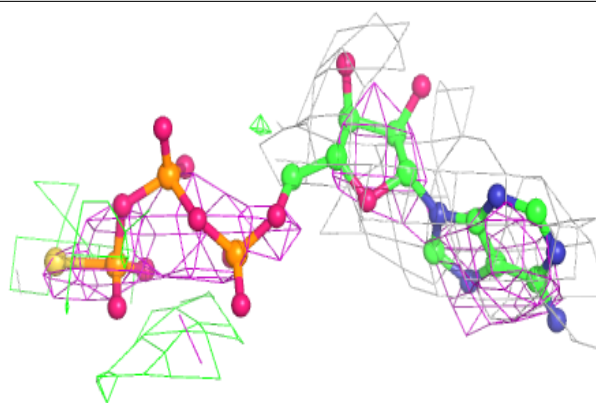
**Electron density around AGS C 549:**

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

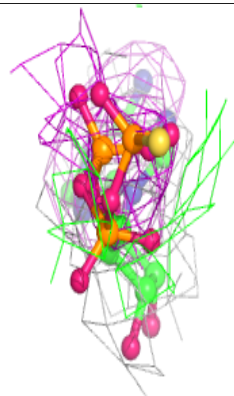
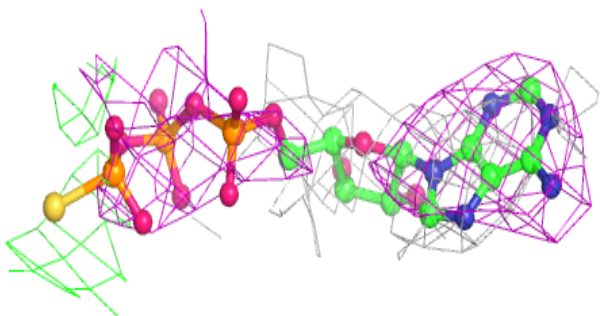
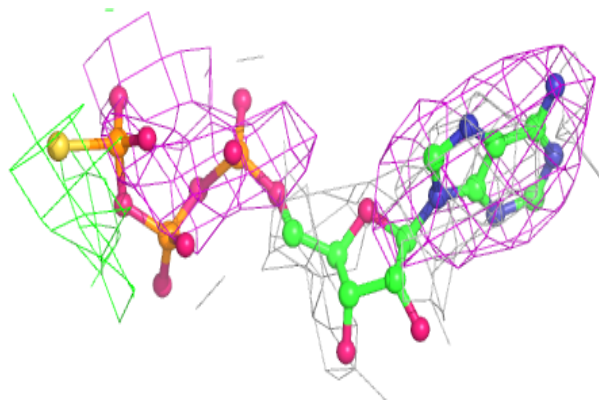


Electron density around AGS N 549:

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

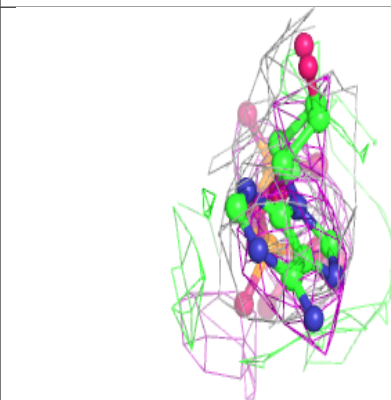
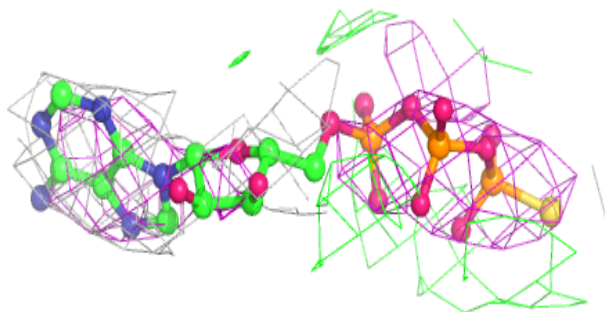
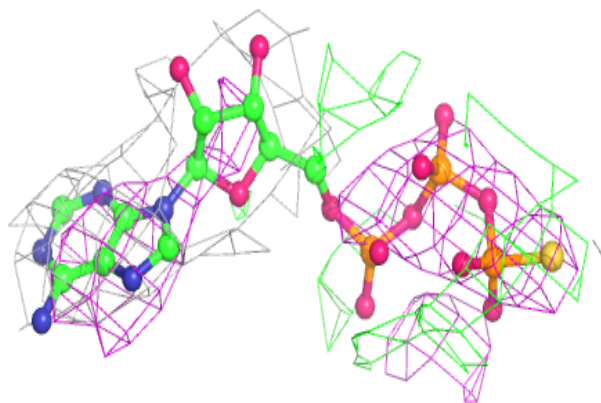
**Electron density around AGS J 549:**

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

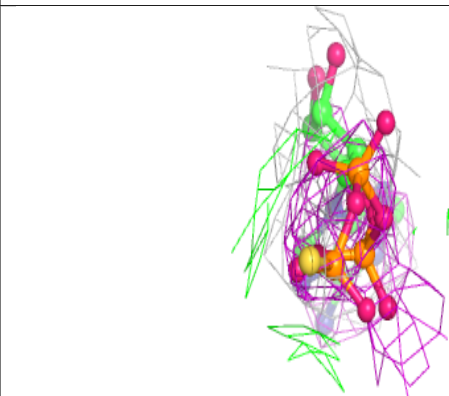
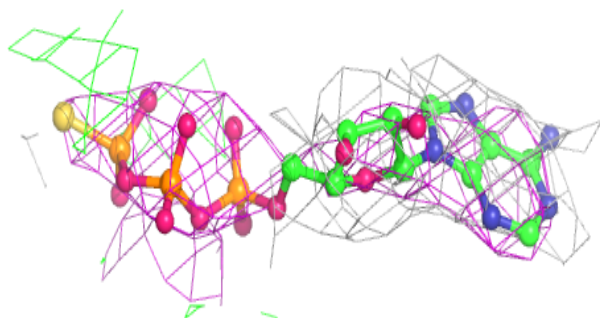
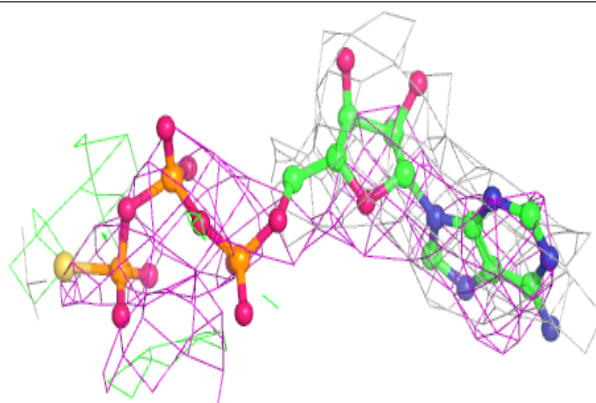


Electron density around AGS M 549:

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

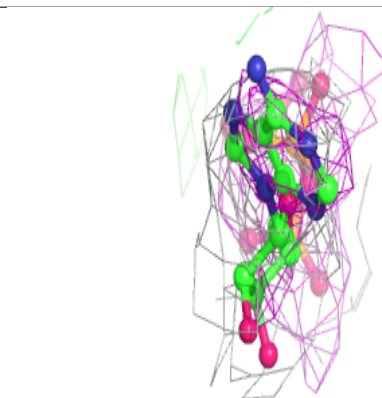
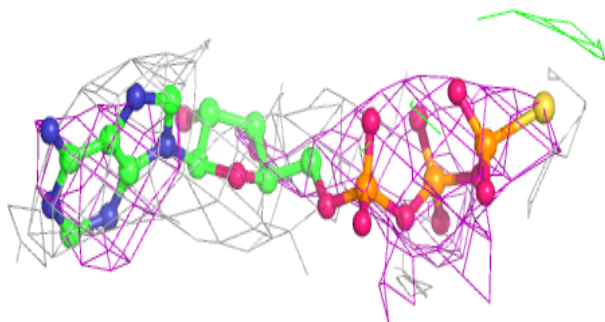
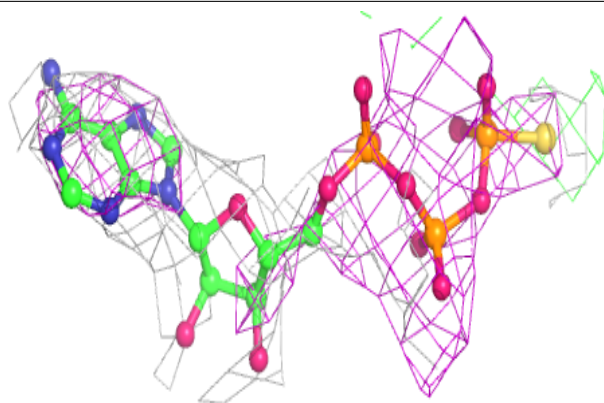
**Electron density around AGS H 549:**

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

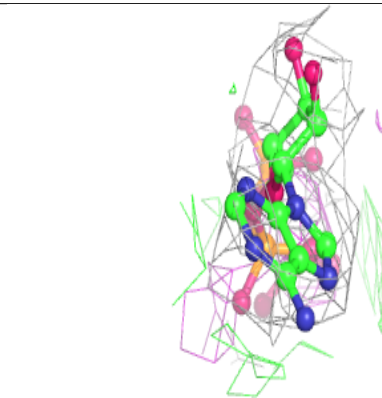
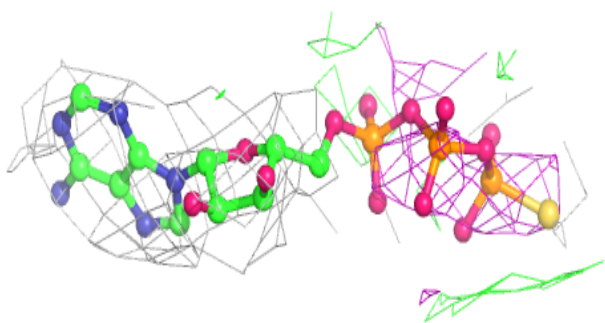
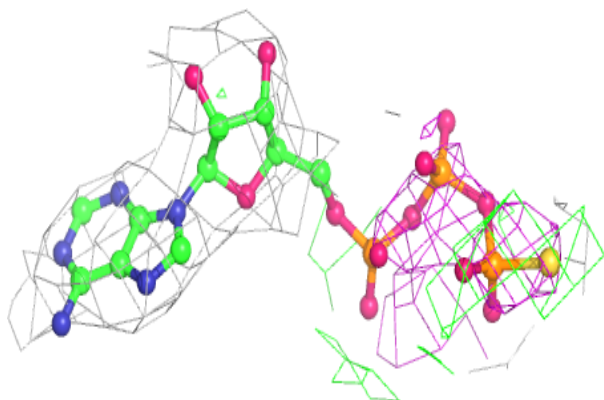


Electron density around AGS F 549:

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

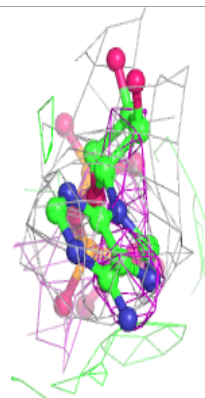
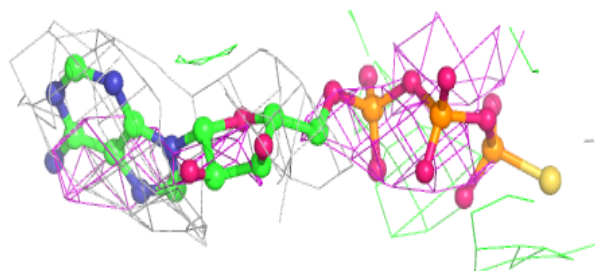
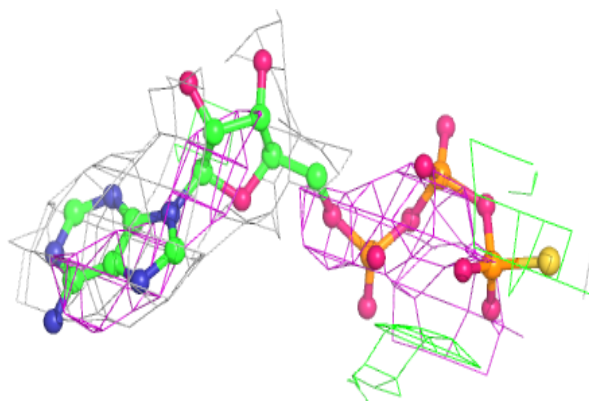
**Electron density around AGS A 549:**

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

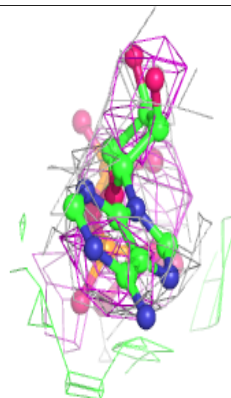
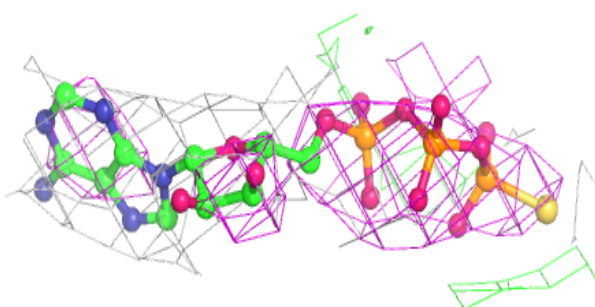
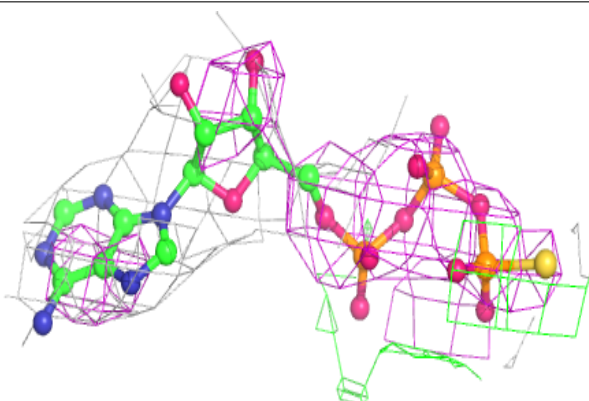


Electron density around AGS B 549:

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

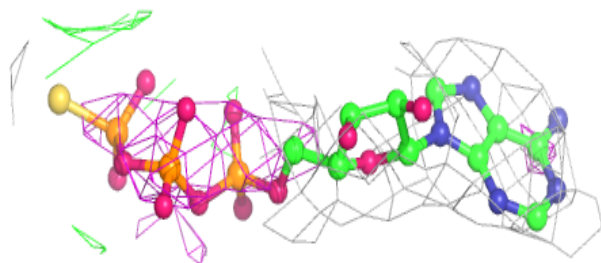
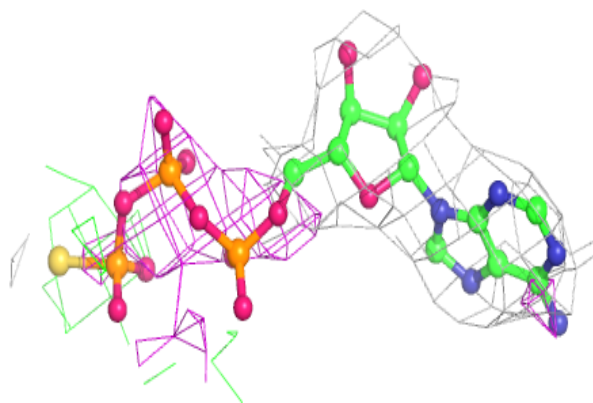
**Electron density around AGS G 549:**

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

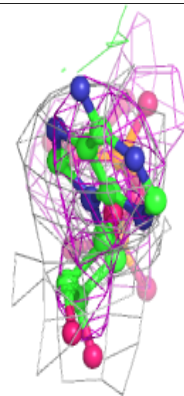
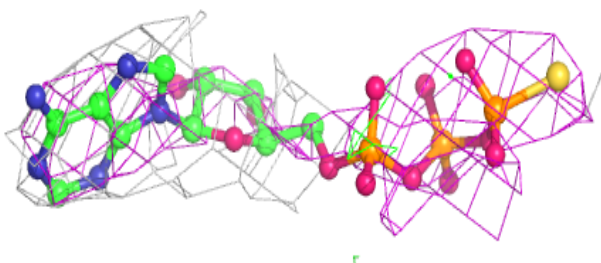
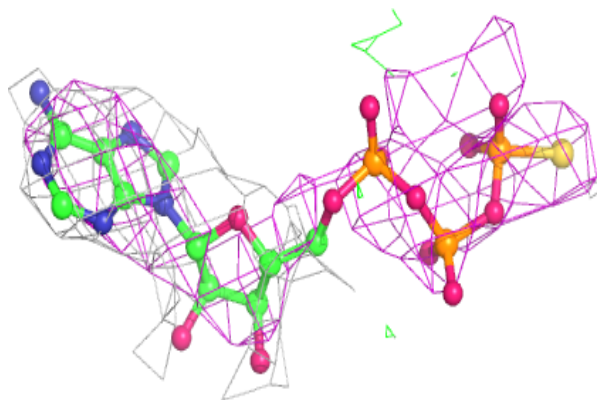


Electron density around AGS D 549:

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

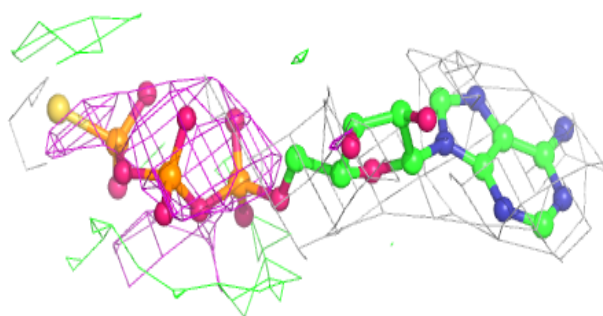
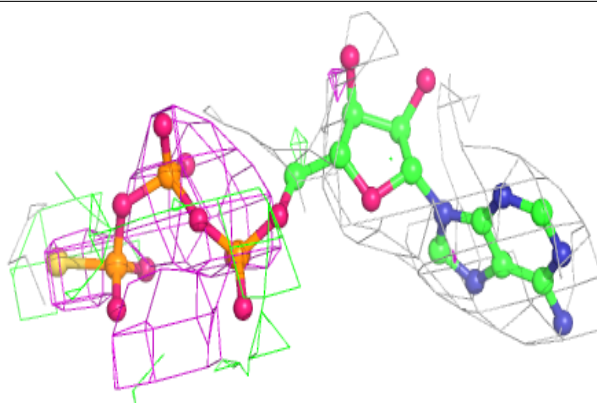
**Electron density around AGS K 549:**

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



Electron density around AGS I 549:

$2mF_o-DF_c$ (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.