



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:15 AM UTC

PDB ID : 8BLL / pdb_00008bll
Title : Structure of RutB
Authors : Rajendran, C.
Deposited on : 2022-11-09
Resolution : 1.54 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

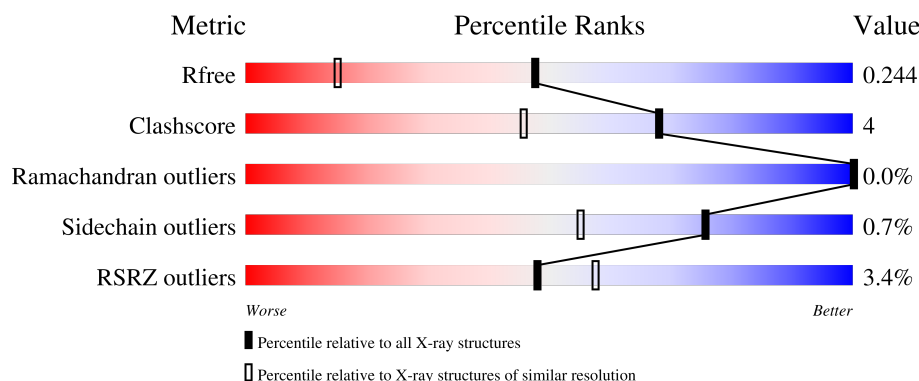
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.54 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	1003 (1.54-1.54)
Clashscore	190562	1025 (1.54-1.54)
Ramachandran outliers	187476	1007 (1.54-1.54)
Sidechain outliers	187428	1007 (1.54-1.54)
RSRZ outliers	180081	1002 (1.54-1.54)

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	230	
1	B	230	
1	C	230	
1	D	230	
1	E	230	

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Mol	Chain	Length	Quality of chain
1	F	230	
1	G	230	
1	H	230	
1	I	230	
1	J	230	
1	K	230	
1	L	230	
1	M	230	
1	N	230	
1	O	230	
1	P	230	

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
2	ACT	C	301	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 57341 atoms, of which 26479 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.

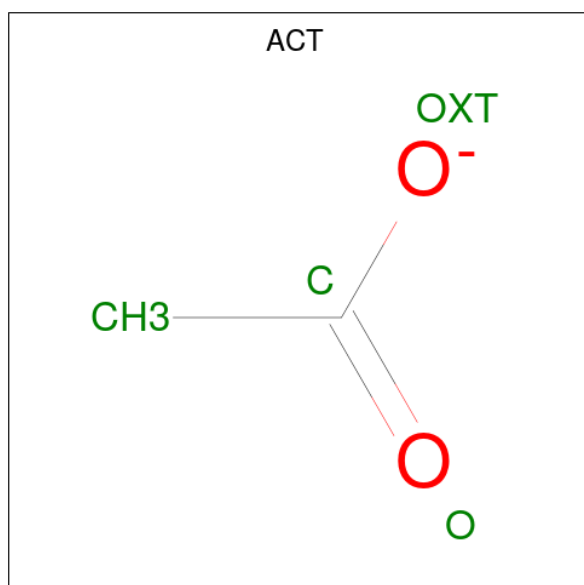
- Molecule 1 is a protein called Ureidoacrylate amidohydrolase RutB.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	222	Total	C	H	N	O	S	0	0	0
			3369	1104	1654	285	322	4			
1	B	222	Total	C	H	N	O	S	0	0	0
			3369	1104	1654	285	322	4			
1	C	222	Total	C	H	N	O	S	0	0	0
			3359	1104	1644	285	322	4			
1	D	222	Total	C	H	N	O	S	0	0	0
			3369	1104	1654	285	322	4			
1	E	222	Total	C	H	N	O	S	0	0	0
			3359	1104	1644	285	322	4			
1	F	222	Total	C	H	N	O	S	0	0	0
			3369	1104	1654	285	322	4			
1	G	222	Total	C	H	N	O	S	0	0	0
			3369	1104	1654	285	322	4			
1	H	222	Total	C	H	N	O	S	0	0	0
			3369	1104	1654	285	322	4			
1	I	222	Total	C	H	N	O	S	0	0	0
			3356	1104	1641	285	322	4			
1	J	222	Total	C	H	N	O	S	0	0	0
			3369	1104	1654	285	322	4			
1	K	222	Total	C	H	N	O	S	0	0	0
			3369	1104	1654	285	322	4			
1	L	222	Total	C	H	N	O	S	0	0	0
			3369	1104	1654	285	322	4			
1	M	222	Total	C	H	N	O	S	0	0	0
			3344	1099	1639	282	320	4			
1	N	222	Total	C	H	N	O	S	0	0	0
			3354	1101	1643	284	322	4			
1	O	222	Total	C	H	N	O	S	0	0	0
			3358	1104	1643	285	322	4			
1	P	222	Total	C	H	N	O	S	0	0	0
			3352	1101	1643	282	322	4			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	166	SER	CYS	conflict	UNP C9QZ65
B	166	SER	CYS	conflict	UNP C9QZ65
C	166	SER	CYS	conflict	UNP C9QZ65
D	166	SER	CYS	conflict	UNP C9QZ65
E	166	SER	CYS	conflict	UNP C9QZ65
F	166	SER	CYS	conflict	UNP C9QZ65
G	166	SER	CYS	conflict	UNP C9QZ65
H	166	SER	CYS	conflict	UNP C9QZ65
I	166	SER	CYS	conflict	UNP C9QZ65
J	166	SER	CYS	conflict	UNP C9QZ65
K	166	SER	CYS	conflict	UNP C9QZ65
L	166	SER	CYS	conflict	UNP C9QZ65
M	166	SER	CYS	conflict	UNP C9QZ65
N	166	SER	CYS	conflict	UNP C9QZ65
O	166	SER	CYS	conflict	UNP C9QZ65
P	166	SER	CYS	conflict	UNP C9QZ65

- Molecule 2 is ACETATE ION (CCD ID: ACT) (formula: $C_2H_3O_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	H	O	0	0
			7	2	3	2		
2	A	1	Total	C	H	O	0	0
			7	2	3	2		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	B	1	Total 7	C 2	H 3	O 2	0	0
2	B	1	Total 7	C 2	H 3	O 2	0	0
2	C	1	Total 7	C 2	H 3	O 2	0	0
2	C	1	Total 7	C 2	H 3	O 2	0	0
2	D	1	Total 7	C 2	H 3	O 2	0	0
2	D	1	Total 7	C 2	H 3	O 2	0	0
2	E	1	Total 7	C 2	H 3	O 2	0	0
2	E	1	Total 7	C 2	H 3	O 2	0	0
2	F	1	Total 7	C 2	H 3	O 2	0	0
2	F	1	Total 7	C 2	H 3	O 2	0	0
2	G	1	Total 7	C 2	H 3	O 2	0	0
2	G	1	Total 7	C 2	H 3	O 2	0	0
2	H	1	Total 7	C 2	H 3	O 2	0	0
2	H	1	Total 7	C 2	H 3	O 2	0	0
2	I	1	Total 7	C 2	H 3	O 2	0	0
2	I	1	Total 7	C 2	H 3	O 2	0	0
2	J	1	Total 7	C 2	H 3	O 2	0	0
2	J	1	Total 7	C 2	H 3	O 2	0	0
2	K	1	Total 7	C 2	H 3	O 2	0	0
2	K	1	Total 7	C 2	H 3	O 2	0	0
2	L	1	Total 7	C 2	H 3	O 2	0	0

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	L	1	Total	C	H	O	0	0
			7	2	3	2		
2	M	1	Total	C	H	O	0	0
			7	2	3	2		
2	M	1	Total	C	H	O	0	0
			7	2	3	2		
2	N	1	Total	C	H	O	0	0
			7	2	3	2		
2	N	1	Total	C	H	O	0	0
			7	2	3	2		
2	O	1	Total	C	H	O	0	0
			7	2	3	2		
2	O	1	Total	C	H	O	0	0
			7	2	3	2		
2	P	1	Total	C	H	O	0	0
			7	2	3	2		
2	P	1	Total	C	H	O	0	0
			7	2	3	2		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	238	Total	O	0	0
			238	238		
3	B	252	Total	O	0	0
			252	252		
3	C	233	Total	O	0	0
			233	233		
3	D	255	Total	O	0	0
			255	255		
3	E	239	Total	O	0	0
			239	239		
3	F	228	Total	O	0	0
			228	228		
3	G	233	Total	O	0	0
			233	233		
3	H	210	Total	O	0	0
			210	210		
3	I	173	Total	O	0	0
			173	173		
3	J	246	Total	O	0	0
			246	246		

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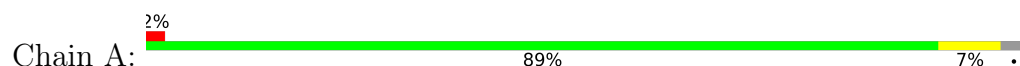
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	K	189	Total 189	O 189	0	0
3	L	208	Total 208	O 208	0	0
3	M	148	Total 148	O 148	0	0
3	N	155	Total 155	O 155	0	0
3	O	168	Total 168	O 168	0	0
3	P	139	Total 139	O 139	0	0

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: Ureidoacrylate amidohydrolase RutB



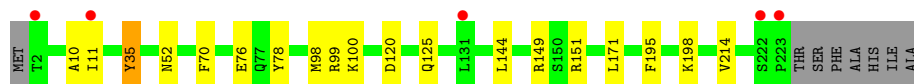
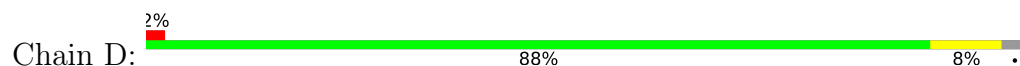
- Molecule 1: Ureidoacrylate amidohydrolase RutB



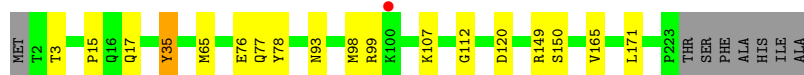
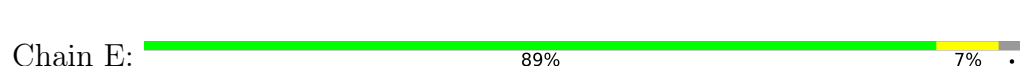
- Molecule 1: Ureidoacrylate amidohydrolase RutB



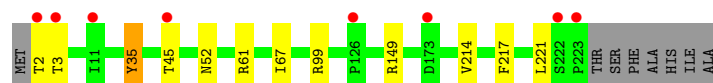
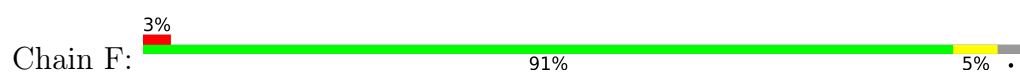
- Molecule 1: Ureidoacrylate amidohydrolase RutB



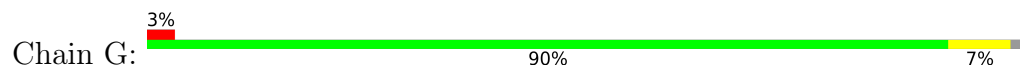
- Molecule 1: Ureidoacrylate amidohydrolase RutB



- Molecule 1: Ureidoacrylate amidohydrolase RutB



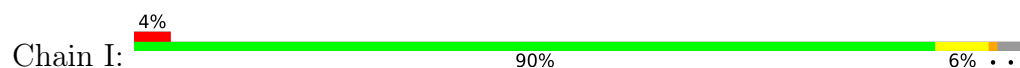
• Molecule 1: Ureidoacrylate amidohydrolase RutB



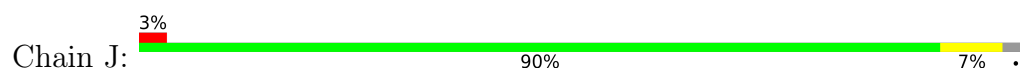
• Molecule 1: Ureidoacrylate amidohydrolase RutB



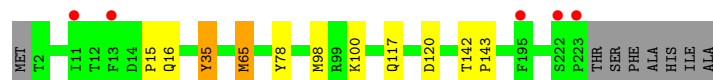
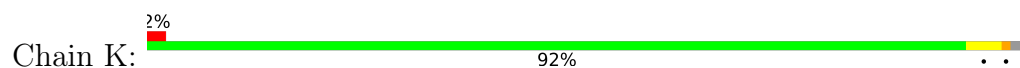
• Molecule 1: Ureidoacrylate amidohydrolase RutB



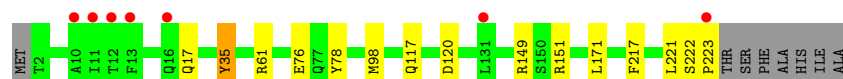
• Molecule 1: Ureidoacrylate amidohydrolase RutB



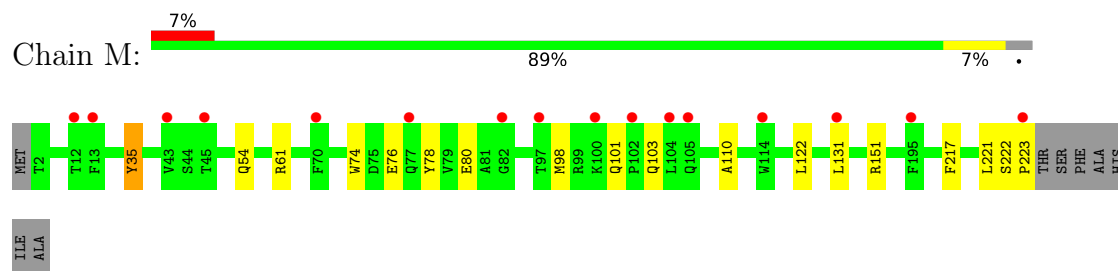
• Molecule 1: Ureidoacrylate amidohydrolase RutB



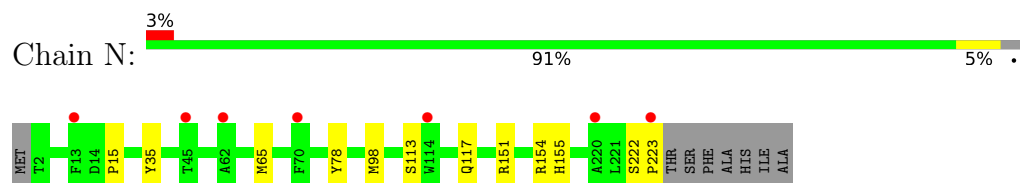
• Molecule 1: Ureidoacrylate amidohydrolase RutB



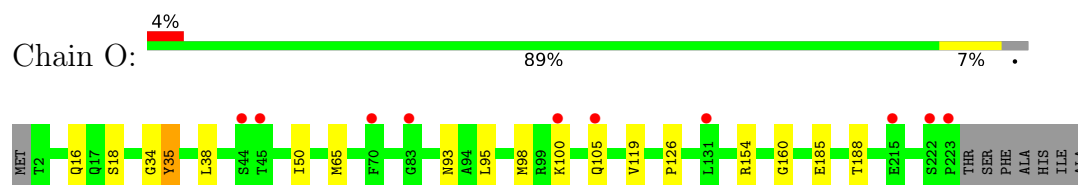
- Molecule 1: Ureidoacrylate amidohydrolase RutB



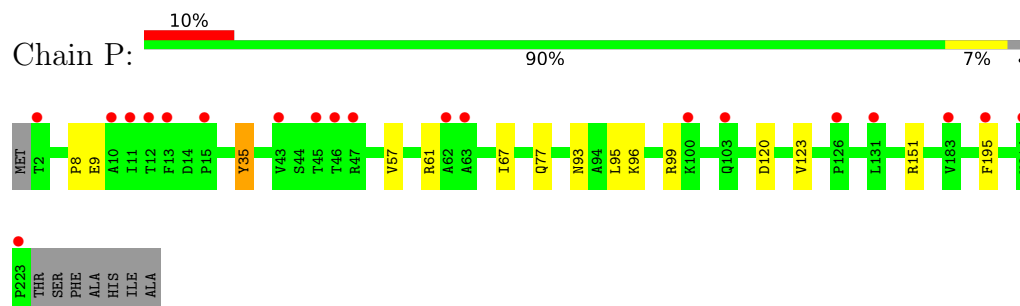
- Molecule 1: Ureidoacrylate amidohydrolase RutB



- Molecule 1: Ureidoacrylate amidohydrolase RutB



- Molecule 1: Ureidoacrylate amidohydrolase RutB



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	69.46Å 108.61Å 124.34Å 85.66° 83.22° 74.34°	Depositor
Resolution (Å)	47.00 – 1.54 47.00 – 1.54	Depositor EDS
% Data completeness (in resolution range)	83.7 (47.00-1.54) 85.5 (47.00-1.54)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.11 (at 1.54Å)	Xtriage
Refinement program	PHENIX 1.18rc2_3793	Depositor
R, R_{free}	0.207 , 0.243 0.207 , 0.244	Depositor DCC
R_{free} test set	23015 reflections (4.49%)	wwPDB-VP
Wilson B-factor (Å ²)	17.0	Xtriage
Anisotropy	0.047	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 38.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.79	EDS
Total number of atoms	57341	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 14.99% 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: ACT

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	0.48	0/1760	0.67	0/2401
1	B	0.48	0/1760	0.63	0/2401
1	C	0.47	0/1760	0.63	0/2401
1	D	0.51	0/1760	0.68	0/2401
1	E	0.47	0/1760	0.64	0/2401
1	F	0.47	0/1760	0.62	0/2401
1	G	0.49	0/1760	0.63	0/2401
1	H	0.45	0/1760	0.59	0/2401
1	I	0.43	0/1760	0.58	0/2401
1	J	0.51	1/1760 (0.1%)	0.66	0/2401
1	K	0.42	0/1760	0.58	0/2401
1	L	0.44	0/1760	0.62	0/2401
1	M	0.42	0/1750	0.57	0/2389
1	N	0.42	0/1756	0.58	0/2397
1	O	0.42	0/1760	0.59	0/2401
1	P	0.42	0/1754	0.57	0/2394
All	All	0.46	1/28140 (0.0%)	0.62	0/38393

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	J	181	GLY	CA-C	-5.80	1.48	1.52

There are no bond angle outliers.

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	1715	1654	1675	15	0
1	B	1715	1654	1675	10	0
1	C	1715	1644	1675	14	0
1	D	1715	1654	1675	16	0
1	E	1715	1644	1675	16	0
1	F	1715	1654	1675	9	0
1	G	1715	1654	1675	16	0
1	H	1715	1654	1675	13	0
1	I	1715	1641	1675	11	0
1	J	1715	1654	1675	13	0
1	K	1715	1654	1675	8	0
1	L	1715	1654	1675	13	0
1	M	1705	1639	1660	11	0
1	N	1711	1643	1664	8	0
1	O	1715	1643	1675	12	0
1	P	1709	1643	1664	14	0
2	A	8	6	6	1	0
2	B	8	6	6	1	0
2	C	8	6	6	2	0
2	D	8	6	6	0	0
2	E	8	6	6	0	0
2	F	8	6	6	0	0
2	G	8	6	6	0	0
2	H	8	6	6	1	0
2	I	8	6	6	0	0
2	J	8	6	6	0	0
2	K	8	6	6	0	0
2	L	8	6	6	0	0
2	M	8	6	6	0	0
2	N	8	6	6	0	0
2	O	8	6	6	0	0
2	P	8	6	6	0	0
3	A	238	0	0	5	2
3	B	252	0	0	2	0
3	C	233	0	0	7	1
3	D	255	0	0	8	0
3	E	239	0	0	7	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	F	228	0	0	3	1
3	G	233	0	0	7	0
3	H	210	0	0	10	2
3	I	173	0	0	6	0
3	J	246	0	0	4	3
3	K	189	0	0	4	2
3	L	208	0	0	7	0
3	M	148	0	0	2	0
3	N	155	0	0	3	0
3	O	168	0	0	4	1
3	P	139	0	0	8	0
All	All	30862	26479	26859	193	6

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 4.

All (193) 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:H:215:GLU:OE2	3:H:401:HOH:O	1.58	1.20
1:H:151:ARG:NH2	3:H:402:HOH:O	1.96	0.97
1:E:149:ARG:NH2	3:E:401:HOH:O	1.98	0.96
1:E:3:THR:HG23	1:G:3:THR:H	1.30	0.94
1:A:125:GLN:NE2	3:A:401:HOH:O	2.00	0.93
1:B:166:SER:OG	2:B:302:ACT:H1	1.71	0.90
1:D:149:ARG:NH2	3:D:401:HOH:O	2.06	0.89
1:A:151:ARG:NH2	3:A:402:HOH:O	2.07	0.87
1:A:166:SER:OG	2:A:301:ACT:H1	1.75	0.87
1:C:147:ILE:O	3:C:401:HOH:O	1.94	0.86
1:D:151:ARG:NH2	3:D:405:HOH:O	2.13	0.81
1:H:223:PRO:O	3:H:403:HOH:O	1.98	0.80
1:I:223:PRO:O	3:I:401:HOH:O	2.00	0.79
1:B:7:ARG:NH1	3:B:401:HOH:O	2.15	0.79
1:E:112:GLY:O	3:E:402:HOH:O	2.01	0.79
1:L:117:GLN:NE2	3:L:403:HOH:O	2.15	0.78
1:O:185:GLU:OE2	3:O:401:HOH:O	2.02	0.76
1:E:120:ASP:OD1	3:E:403:HOH:O	2.03	0.76
1:K:16:GLN:OE1	3:K:401:HOH:O	2.06	0.73
1:C:154:ARG:NH1	3:C:402:HOH:O	2.12	0.73
1:D:10:ALA:O	3:D:402:HOH:O	2.08	0.72
1:I:154:ARG:NH1	3:I:403:HOH:O	2.23	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:120:ASP:OD1	3:D:403:HOH:O	2.08	0.71
1:G:71:GLN:OE1	3:G:401:HOH:O	2.07	0.71
1:H:125:GLN:NE2	3:H:404:HOH:O	2.00	0.70
1:O:126:PRO:O	3:O:402:HOH:O	2.09	0.70
1:F:3:THR:OG1	3:F:401:HOH:O	2.08	0.69
1:A:149:ARG:NH2	3:A:404:HOH:O	2.23	0.69
1:M:151:ARG:NH2	3:M:401:HOH:O	2.25	0.68
1:D:76:GLU:OE2	3:D:404:HOH:O	2.10	0.68
1:P:151:ARG:NH2	3:P:403:HOH:O	2.26	0.68
1:O:154:ARG:NH1	3:O:404:HOH:O	2.28	0.67
1:E:3:THR:HG23	1:G:3:THR:N	2.08	0.67
1:K:120:ASP:OD1	3:K:402:HOH:O	2.13	0.66
1:L:76:GLU:OE2	3:L:401:HOH:O	2.13	0.66
1:L:120:ASP:OD1	3:L:402:HOH:O	2.15	0.65
1:H:114:TRP:N	3:H:409:HOH:O	2.31	0.64
1:I:149:ARG:HG2	1:I:149:ARG:HH11	1.62	0.64
1:C:147:ILE:HA	3:C:401:HOH:O	1.99	0.62
1:B:100:LYS:NZ	1:B:101:GLN:OE1	2.33	0.61
1:J:151:ARG:NH2	3:J:405:HOH:O	2.33	0.61
1:L:76:GLU:OE1	3:L:404:HOH:O	2.16	0.61
1:E:17:GLN:OE1	3:E:404:HOH:O	2.16	0.61
1:I:198:LYS:NZ	3:I:406:HOH:O	2.34	0.60
1:P:120:ASP:O	3:P:401:HOH:O	2.17	0.60
1:E:107:LYS:NZ	3:E:407:HOH:O	2.35	0.59
1:G:99:ARG:NH2	3:G:403:HOH:O	2.33	0.59
1:N:15:PRO:O	1:N:65:MET:HG2	2.02	0.59
1:J:99:ARG:NH2	3:J:406:HOH:O	2.33	0.59
1:N:155:HIS:ND1	3:N:404:HOH:O	2.32	0.59
1:F:99:ARG:NH2	3:F:405:HOH:O	2.33	0.57
1:P:9:GLU:OE2	3:P:402:HOH:O	2.17	0.57
1:K:100:LYS:NZ	3:K:404:HOH:O	2.37	0.57
1:G:154:ARG:NH1	3:G:405:HOH:O	2.38	0.57
1:K:78:TYR:CZ	1:K:98:MET:HE1	2.42	0.55
2:H:301:ACT:H2	3:H:511:HOH:O	2.06	0.55
1:A:99:ARG:NH2	3:A:406:HOH:O	2.32	0.55
1:C:150:SER:HB2	3:C:401:HOH:O	2.07	0.55
1:O:100:LYS:N	1:O:100:LYS:HD3	2.22	0.54
1:J:70:PHE:HZ	1:J:144:LEU:HD22	1.73	0.54
1:I:42:ASP:O	3:I:402:HOH:O	2.18	0.53
1:H:99:ARG:NH2	3:H:413:HOH:O	2.40	0.53
1:A:61:ARG:HG3	3:A:414:HOH:O	2.08	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:151:ARG:NH2	3:L:412:HOH:O	2.43	0.52
1:P:77:GLN:NE2	3:P:412:HOH:O	2.41	0.52
3:E:401:HOH:O	1:F:149:ARG:NE	2.43	0.52
1:G:78:TYR:CZ	1:G:98:MET:HE1	2.44	0.52
1:E:35:TYR:C	1:E:35:TYR:CD1	2.88	0.51
1:A:195:PHE:CD2	1:C:198:LYS:HE3	2.46	0.51
1:H:35:TYR:OH	1:H:93:ASN:HB2	2.10	0.51
1:K:15:PRO:O	1:K:65:MET:HG3	2.10	0.51
1:B:100:LYS:HG3	1:B:101:GLN:HG3	1.92	0.51
1:K:117:GLN:HG3	3:K:497:HOH:O	2.11	0.51
1:F:52:ASN:HB3	1:F:214:VAL:HG21	1.93	0.51
1:O:50:ILE:HD13	1:O:119:VAL:HG21	1.93	0.51
1:G:222:SER:OG	1:G:223:PRO:HD3	2.11	0.50
1:I:99:ARG:NH2	3:I:409:HOH:O	2.39	0.50
1:I:2:THR:CG2	1:I:220:ALA:O	2.59	0.50
1:L:35:TYR:CD1	1:L:35:TYR:C	2.90	0.49
1:N:113:SER:O	1:N:117:GLN:HG3	2.13	0.49
1:P:123:VAL:N	3:P:407:HOH:O	2.34	0.49
1:C:120:ASP:OD2	3:C:403:HOH:O	2.19	0.49
1:F:35:TYR:C	1:F:35:TYR:CD1	2.91	0.49
1:M:35:TYR:C	1:M:35:TYR:CD1	2.90	0.49
1:D:35:TYR:CD1	1:D:35:TYR:C	2.91	0.48
1:D:78:TYR:CZ	1:D:98:MET:HE1	2.48	0.48
1:P:35:TYR:CD1	1:P:35:TYR:C	2.91	0.48
1:P:57:VAL:HG13	1:P:67:ILE:HD13	1.96	0.48
1:A:35:TYR:C	1:A:35:TYR:CD1	2.91	0.47
1:J:113:SER:O	1:J:117:GLN:HG3	2.13	0.47
1:C:35:TYR:OH	1:C:93:ASN:HB2	2.14	0.47
1:A:11:ILE:HD11	1:A:182:VAL:HG21	1.96	0.47
1:C:35:TYR:C	1:C:35:TYR:CD1	2.92	0.47
1:O:160:GLY:O	1:O:188:THR:HA	2.15	0.47
1:P:61:ARG:HB2	1:P:67:ILE:HD11	1.95	0.47
1:B:78:TYR:CZ	1:B:98:MET:HE1	2.50	0.47
1:B:149:ARG:HD3	1:D:149:ARG:CZ	2.45	0.47
1:M:54:GLN:HG3	1:M:122:LEU:HD21	1.95	0.47
1:M:101:GLN:HB3	1:M:103:GLN:NE2	2.30	0.47
1:E:3:THR:CG2	1:G:3:THR:H	2.15	0.47
1:M:217:PHE:CZ	1:M:221:LEU:HD11	2.50	0.47
1:A:78:TYR:CZ	1:A:98:MET:HE1	2.50	0.46
1:E:76:GLU:OE1	1:E:77:GLN:OE1	2.34	0.46
1:G:7:ARG:NH1	3:G:410:HOH:O	2.44	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:35:TYR:CD1	1:G:35:TYR:C	2.93	0.46
1:G:154:ARG:NH1	3:G:412:HOH:O	2.46	0.46
1:H:107:LYS:HE2	3:H:459:HOH:O	2.15	0.46
1:O:18:SER:HB3	1:O:65:MET:HE3	1.98	0.46
1:L:17:GLN:OE1	3:L:405:HOH:O	2.20	0.46
1:O:35:TYR:C	1:O:35:TYR:CD1	2.94	0.46
1:L:78:TYR:CZ	1:L:98:MET:HE1	2.50	0.46
1:D:99:ARG:NH2	3:D:415:HOH:O	2.43	0.46
1:K:35:TYR:C	1:K:35:TYR:CD1	2.93	0.46
1:H:78:TYR:CZ	1:H:98:MET:HE1	2.50	0.45
1:I:52:ASN:HB3	1:I:214:VAL:HG21	1.98	0.45
1:N:35:TYR:C	1:N:35:TYR:CD1	2.94	0.45
1:F:217:PHE:CZ	1:F:221:LEU:HD11	2.51	0.45
1:C:136:TYR:HD1	2:C:301:ACT:H2	1.82	0.45
1:C:166:SER:OG	2:C:301:ACT:CH3	2.64	0.45
1:D:70:PHE:HZ	1:D:144:LEU:HD22	1.81	0.45
1:D:171:LEU:C	1:D:171:LEU:HD23	2.41	0.45
1:P:95:LEU:O	1:P:99:ARG:HG3	2.15	0.45
1:A:171:LEU:C	1:A:171:LEU:HD23	2.42	0.45
1:J:171:LEU:C	1:J:171:LEU:HD23	2.42	0.45
1:N:222:SER:OG	1:N:223:PRO:HD3	2.16	0.45
1:D:52:ASN:HB3	1:D:214:VAL:HG21	1.99	0.45
1:O:35:TYR:OH	1:O:93:ASN:HB2	2.17	0.45
1:F:61:ARG:HB2	1:F:67:ILE:HD11	1.99	0.44
1:H:149:ARG:NE	3:H:410:HOH:O	2.32	0.44
1:J:2:THR:HG23	1:J:220:ALA:O	2.18	0.44
1:J:70:PHE:CZ	1:J:144:LEU:HD22	2.52	0.44
1:A:195:PHE:CD1	1:A:195:PHE:C	2.96	0.44
1:I:35:TYR:C	1:I:35:TYR:CD1	2.96	0.44
1:P:96:LYS:HG3	3:P:421:HOH:O	2.17	0.44
1:J:151:ARG:NH2	3:J:401:HOH:O	2.22	0.44
1:E:15:PRO:O	1:E:65:MET:HG2	2.18	0.44
1:E:171:LEU:C	1:E:171:LEU:HD23	2.43	0.43
1:A:217:PHE:CZ	1:A:221:LEU:HD11	2.54	0.43
1:H:151:ARG:NH2	3:H:411:HOH:O	2.34	0.43
1:C:198:LYS:HB3	1:C:198:LYS:HE2	1.70	0.43
1:M:74:TRP:HB3	1:M:80:GLU:HB2	2.00	0.43
1:G:142:THR:HB	1:G:143:PRO:HD2	2.00	0.43
1:G:193:PRO:HG2	1:G:195:PHE:CE2	2.54	0.43
3:I:566:HOH:O	1:J:147:ILE:HD11	2.19	0.43
1:L:217:PHE:CZ	1:L:221:LEU:HD11	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:65:MET:HE2	1:A:65:MET:HB3	1.86	0.43
1:E:99:ARG:NH1	3:E:415:HOH:O	2.48	0.43
1:J:195:PHE:C	1:J:195:PHE:CD1	2.97	0.43
1:J:35:TYR:C	1:J:35:TYR:CD1	2.96	0.42
1:P:35:TYR:OH	1:P:93:ASN:HB2	2.19	0.42
1:E:78:TYR:CZ	1:E:98:MET:HE1	2.53	0.42
1:H:35:TYR:CD1	1:H:35:TYR:C	2.97	0.42
1:K:142:THR:HB	1:K:143:PRO:HD2	2.01	0.42
1:O:105:GLN:NE2	3:O:406:HOH:O	2.38	0.42
1:P:8:PRO:HB3	3:P:524:HOH:O	2.19	0.42
1:A:70:PHE:HZ	1:A:144:LEU:HD22	1.84	0.42
1:B:86:SER:HB2	1:E:150:SER:OG	2.19	0.42
1:C:117:GLN:NE2	3:C:412:HOH:O	2.52	0.42
1:C:155:HIS:NE2	3:C:404:HOH:O	2.36	0.42
1:D:171:LEU:HD23	1:D:171:LEU:O	2.19	0.42
1:M:78:TYR:CZ	1:M:98:MET:HE1	2.54	0.42
1:D:195:PHE:CD1	1:D:195:PHE:C	2.97	0.42
1:J:154:ARG:NH1	3:J:414:HOH:O	2.51	0.42
1:M:61:ARG:HG3	3:M:455:HOH:O	2.19	0.42
1:B:161:ILE:HA	1:B:162:ALA:HA	1.83	0.42
1:D:198:LYS:NZ	3:D:419:HOH:O	2.46	0.42
1:P:99:ARG:NH2	3:P:421:HOH:O	2.53	0.42
1:B:35:TYR:CD1	1:B:35:TYR:C	2.98	0.42
1:B:125:GLN:NE2	3:B:402:HOH:O	2.32	0.41
1:L:171:LEU:C	1:L:171:LEU:HD23	2.45	0.41
1:N:154:ARG:NH1	3:N:415:HOH:O	2.52	0.41
1:P:195:PHE:C	1:P:195:PHE:CD1	2.98	0.41
1:L:222:SER:HB3	1:L:223:PRO:HD3	2.01	0.41
1:M:131:LEU:HD23	1:M:131:LEU:HA	1.96	0.41
1:E:35:TYR:OH	1:E:93:ASN:HB2	2.21	0.41
1:N:78:TYR:CZ	1:N:98:MET:HE1	2.55	0.41
1:N:151:ARG:NH2	3:N:417:HOH:O	2.53	0.41
1:J:217:PHE:CZ	1:J:221:LEU:HD11	2.56	0.41
1:I:35:TYR:OH	1:I:93:ASN:HB2	2.21	0.41
1:D:125:GLN:NE2	3:D:406:HOH:O	2.21	0.41
1:F:2:THR:N	3:F:416:HOH:O	2.53	0.41
1:G:154:ARG:NH2	1:H:85:GLY:O	2.54	0.41
1:I:15:PRO:O	1:I:65:MET:HG2	2.21	0.41
1:M:76:GLU:HG3	1:M:110:ALA:HB1	2.03	0.41
1:O:95:LEU:HD23	1:O:98:MET:CE	2.51	0.41
1:M:222:SER:OG	1:M:223:PRO:HD3	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:O:34:GLY:O	1:O:38:LEU:HD23	2.21	0.41
1:C:154:ARG:HE	1:C:154:ARG:HB3	1.69	0.40
1:G:194:LYS:NZ	3:G:419:HOH:O	2.54	0.40
1:L:61:ARG:HG3	1:L:61:ARG:HH11	1.86	0.40
1:L:149:ARG:HD2	3:L:428:HOH:O	2.22	0.40
1:F:61:ARG:CB	1:F:67:ILE:HD11	2.52	0.40
1:G:37:ASP:O	3:G:402:HOH:O	2.22	0.40

All (6) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:557:HOH:O	3:H:558:HOH:O[1_655]	2.02	0.18
3:C:620:HOH:O	3:F:593:HOH:O[1_465]	2.03	0.17
3:J:577:HOH:O	3:K:457:HOH:O[1_556]	2.08	0.12
3:A:437:HOH:O	3:H:421:HOH:O[1_655]	2.14	0.06
3:J:577:HOH:O	3:K:566:HOH:O[1_556]	2.18	0.02
3:J:542:HOH:O	3:O:537:HOH:O[1_456]	2.19	0.01

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	220/230 (96%)	215 (98%)	5 (2%)	0	100	100
1	B	220/230 (96%)	215 (98%)	5 (2%)	0	100	100
1	C	220/230 (96%)	214 (97%)	6 (3%)	0	100	100
1	D	220/230 (96%)	216 (98%)	4 (2%)	0	100	100
1	E	220/230 (96%)	213 (97%)	6 (3%)	1 (0%)	24	8
1	F	220/230 (96%)	216 (98%)	4 (2%)	0	100	100
1	G	220/230 (96%)	216 (98%)	4 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	H	220/230 (96%)	216 (98%)	4 (2%)	0	100	100
1	I	220/230 (96%)	217 (99%)	3 (1%)	0	100	100
1	J	220/230 (96%)	214 (97%)	6 (3%)	0	100	100
1	K	220/230 (96%)	213 (97%)	7 (3%)	0	100	100
1	L	220/230 (96%)	215 (98%)	5 (2%)	0	100	100
1	M	220/230 (96%)	215 (98%)	5 (2%)	0	100	100
1	N	220/230 (96%)	216 (98%)	4 (2%)	0	100	100
1	O	220/230 (96%)	215 (98%)	5 (2%)	0	100	100
1	P	220/230 (96%)	217 (99%)	3 (1%)	0	100	100
All	All	3520/3680 (96%)	3443 (98%)	76 (2%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	165	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	181/188 (96%)	180 (99%)	1 (1%)	78	59
1	B	181/188 (96%)	180 (99%)	1 (1%)	78	59
1	C	181/188 (96%)	180 (99%)	1 (1%)	78	59
1	D	181/188 (96%)	178 (98%)	3 (2%)	53	23
1	E	181/188 (96%)	180 (99%)	1 (1%)	78	59
1	F	181/188 (96%)	179 (99%)	2 (1%)	65	38
1	G	181/188 (96%)	180 (99%)	1 (1%)	78	59
1	H	181/188 (96%)	181 (100%)	0	100	100
1	I	181/188 (96%)	178 (98%)	3 (2%)	53	23

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	J	181/188 (96%)	181 (100%)	0	100	100
1	K	181/188 (96%)	179 (99%)	2 (1%)	65	38
1	L	181/188 (96%)	180 (99%)	1 (1%)	78	59
1	M	179/188 (95%)	178 (99%)	1 (1%)	78	59
1	N	180/188 (96%)	180 (100%)	0	100	100
1	O	181/188 (96%)	179 (99%)	2 (1%)	65	38
1	P	180/188 (96%)	178 (99%)	2 (1%)	65	38
All	All	2892/3008 (96%)	2871 (99%)	21 (1%)	76	55

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	35	TYR
1	B	35	TYR
1	C	35	TYR
1	D	11	ILE
1	D	35	TYR
1	D	100	LYS
1	E	35	TYR
1	F	35	TYR
1	F	45	THR
1	G	35	TYR
1	I	2	THR
1	I	35	TYR
1	I	222	SER
1	K	35	TYR
1	K	65	MET
1	L	35	TYR
1	M	35	TYR
1	O	16	GLN
1	O	35	TYR
1	P	35	TYR
1	P	222	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	54	GLN

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Mol	Chain	Res	Type
1	A	77	GLN
1	A	189	HIS
1	D	17	GLN
1	D	117	GLN
1	E	101	GLN
1	G	105	GLN
1	H	17	GLN
1	H	105	GLN
1	I	17	GLN
1	I	77	GLN
1	J	189	HIS
1	K	117	GLN
1	L	117	GLN
1	M	16	GLN
1	M	17	GLN
1	M	189	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](#)

32 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
2	ACT	N	301	-	3,3,3	1.17	0	3,3,3	1.77	2 (66%)
2	ACT	I	301	-	3,3,3	1.10	0	3,3,3	1.46	0
2	ACT	L	302	-	3,3,3	1.20	0	3,3,3	1.32	0
2	ACT	M	301	-	3,3,3	0.93	0	3,3,3	1.46	1 (33%)
2	ACT	O	301	-	3,3,3	0.80	0	3,3,3	1.77	2 (66%)
2	ACT	E	302	-	3,3,3	0.84	0	3,3,3	1.70	2 (66%)
2	ACT	O	302	-	3,3,3	1.46	0	3,3,3	1.36	0
2	ACT	K	302	-	3,3,3	1.09	0	3,3,3	1.59	0
2	ACT	B	302	-	3,3,3	1.31	0	3,3,3	1.47	1 (33%)
2	ACT	A	301	-	3,3,3	1.49	0	3,3,3	1.87	1 (33%)
2	ACT	F	302	-	3,3,3	1.05	0	3,3,3	1.36	0
2	ACT	B	301	-	3,3,3	0.94	0	3,3,3	1.39	0
2	ACT	A	302	-	3,3,3	0.64	0	3,3,3	1.32	0
2	ACT	J	302	-	3,3,3	0.64	0	3,3,3	1.41	0
2	ACT	P	301	-	3,3,3	1.27	0	3,3,3	1.67	1 (33%)
2	ACT	D	301	-	3,3,3	1.12	0	3,3,3	1.28	0
2	ACT	C	301	-	3,3,3	1.57	1 (33%)	3,3,3	1.49	0
2	ACT	N	302	-	3,3,3	1.05	0	3,3,3	1.49	0
2	ACT	L	301	-	3,3,3	1.14	0	3,3,3	1.09	0
2	ACT	J	301	-	3,3,3	0.48	0	3,3,3	1.85	2 (66%)
2	ACT	P	302	-	3,3,3	0.88	0	3,3,3	1.67	1 (33%)
2	ACT	E	301	-	3,3,3	1.25	0	3,3,3	1.66	1 (33%)
2	ACT	K	301	-	3,3,3	0.82	0	3,3,3	1.31	0
2	ACT	G	302	-	3,3,3	0.83	0	3,3,3	1.52	0
2	ACT	G	301	-	3,3,3	1.10	0	3,3,3	1.61	1 (33%)
2	ACT	I	302	-	3,3,3	1.47	0	3,3,3	1.29	0
2	ACT	H	302	-	3,3,3	1.18	0	3,3,3	1.59	0
2	ACT	H	301	-	3,3,3	1.31	1 (33%)	3,3,3	1.29	0
2	ACT	F	301	-	3,3,3	0.97	0	3,3,3	1.52	0
2	ACT	C	302	-	3,3,3	1.27	0	3,3,3	1.53	1 (33%)
2	ACT	D	302	-	3,3,3	0.90	0	3,3,3	1.51	0
2	ACT	M	302	-	3,3,3	0.81	0	3,3,3	1.66	1 (33%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	301	ACT	CH3-C	-2.53	1.39	1.49
2	H	301	ACT	O-C	2.19	1.31	1.22

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301	ACT	OXT-C-O	2.54	131.44	122.03
2	J	301	ACT	OXT-C-O	2.27	130.47	122.03
2	N	301	ACT	O-C-CH3	-2.25	113.30	122.53
2	J	301	ACT	O-C-CH3	-2.23	113.39	122.53
2	O	301	ACT	OXT-C-O	2.21	130.22	122.03
2	M	301	ACT	OXT-C-O	2.15	129.99	122.03
2	G	301	ACT	O-C-CH3	-2.13	113.78	122.53
2	P	302	ACT	OXT-C-O	2.12	129.90	122.03
2	P	301	ACT	OXT-C-O	2.12	129.88	122.03
2	E	302	ACT	OXT-C-O	2.11	129.85	122.03
2	O	301	ACT	O-C-CH3	-2.11	113.89	122.53
2	C	302	ACT	OXT-C-O	2.10	129.82	122.03
2	E	301	ACT	O-C-CH3	-2.10	113.93	122.53
2	B	302	ACT	OXT-C-O	2.09	129.79	122.03
2	M	302	ACT	O-C-CH3	-2.09	113.97	122.53
2	E	302	ACT	O-C-CH3	-2.04	114.17	122.53
2	N	301	ACT	OXT-C-O	2.01	129.47	122.03

There are no chirality outliers.

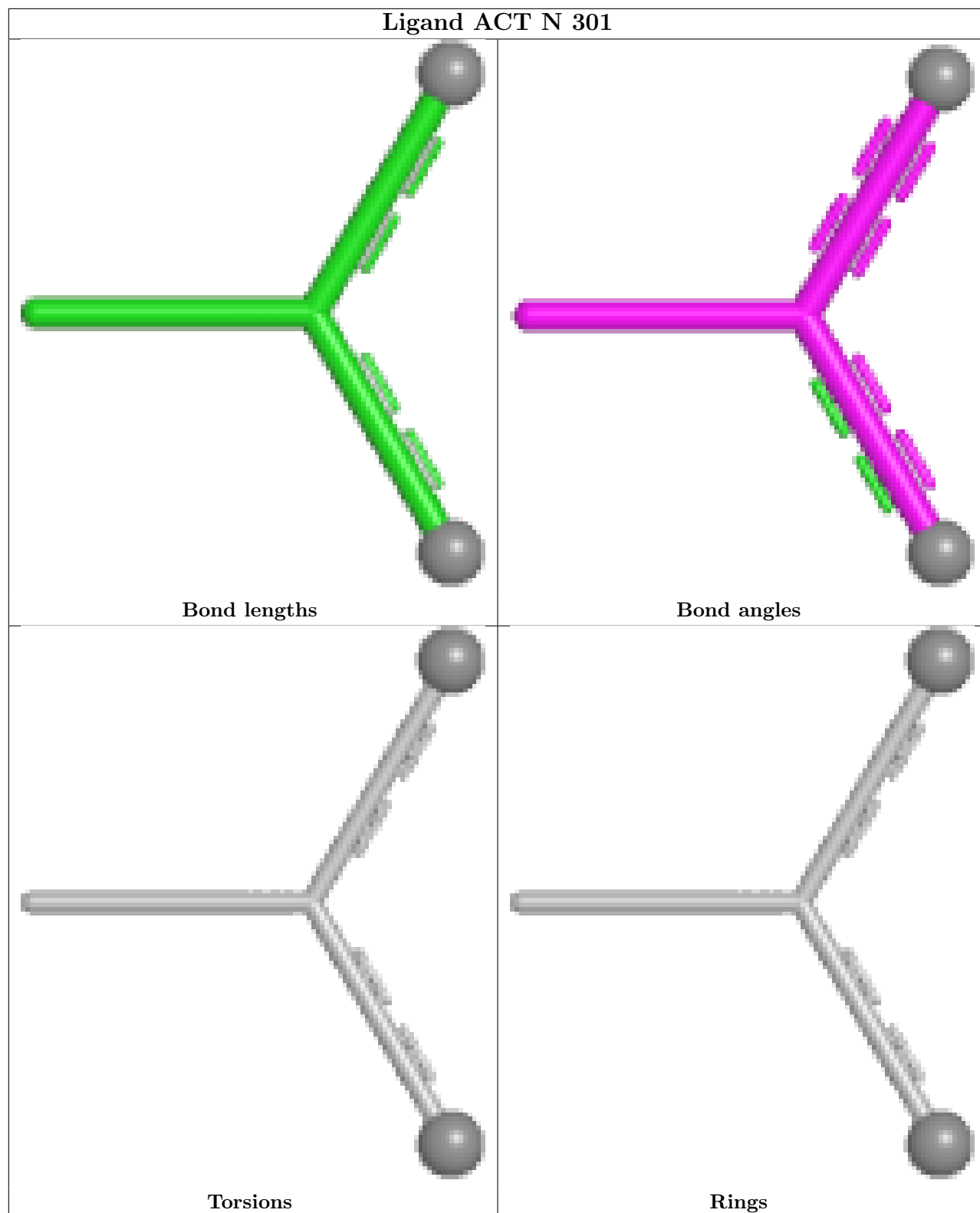
There are no torsion outliers.

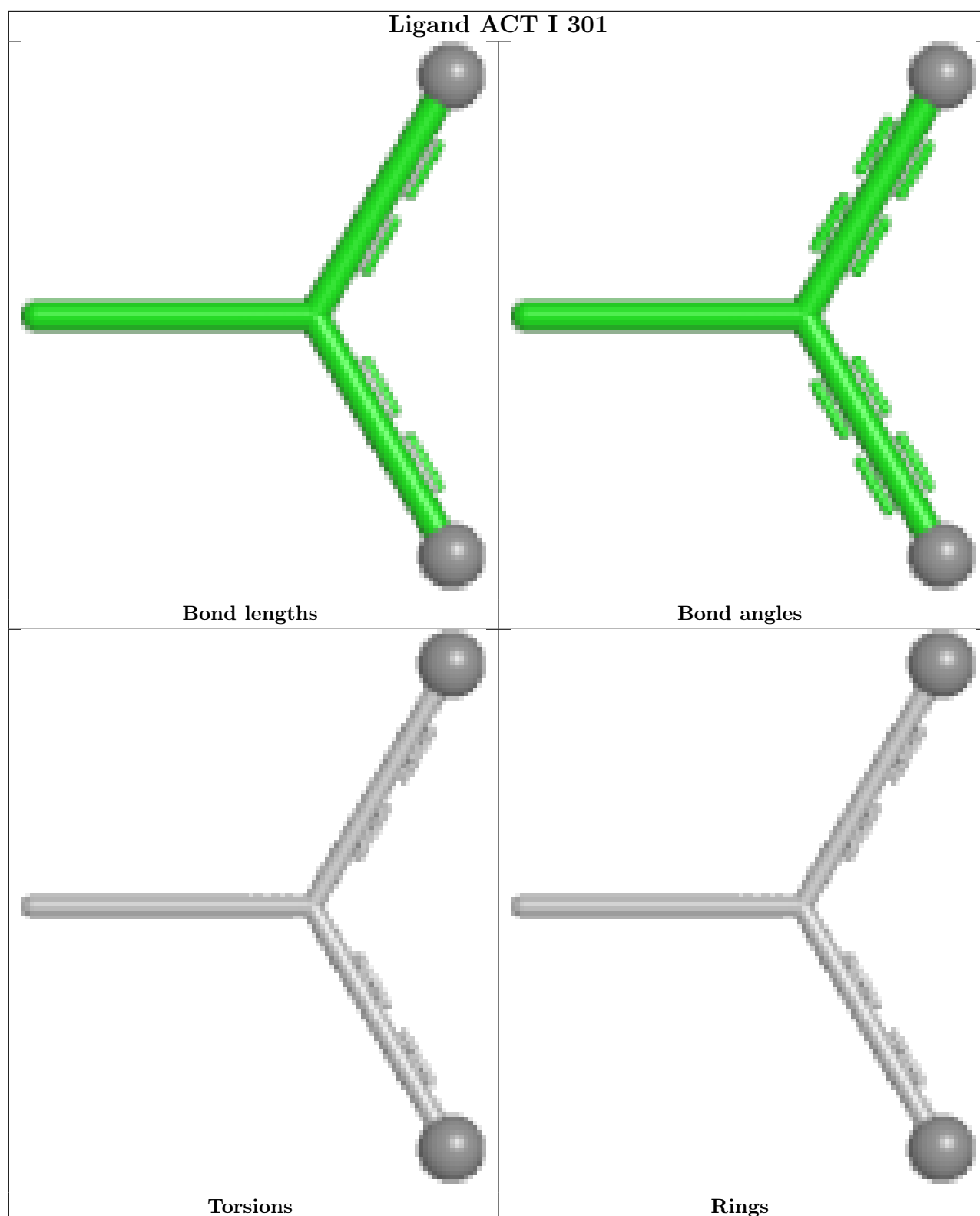
There are no ring outliers.

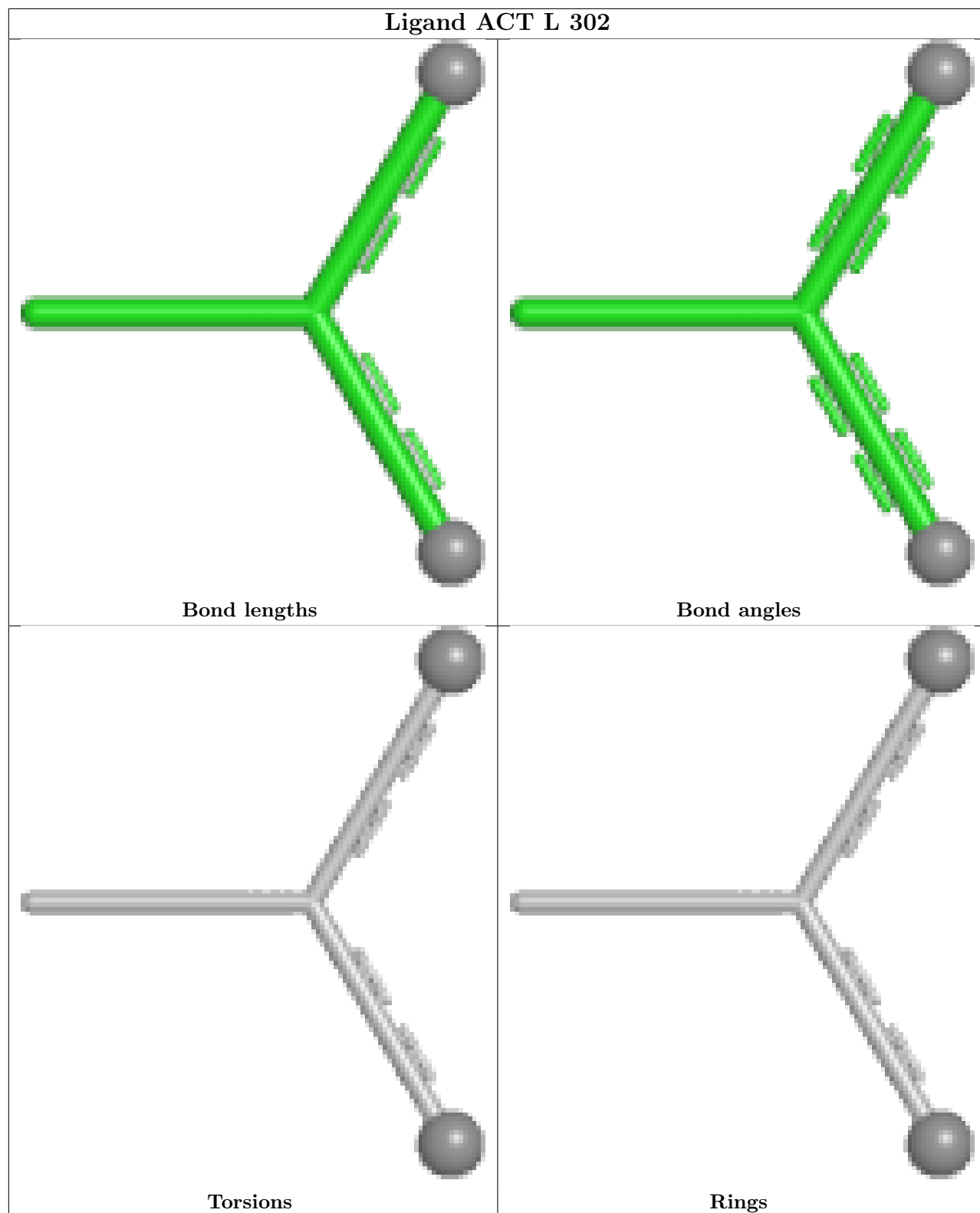
4 monomers are involved in 5 short contacts:

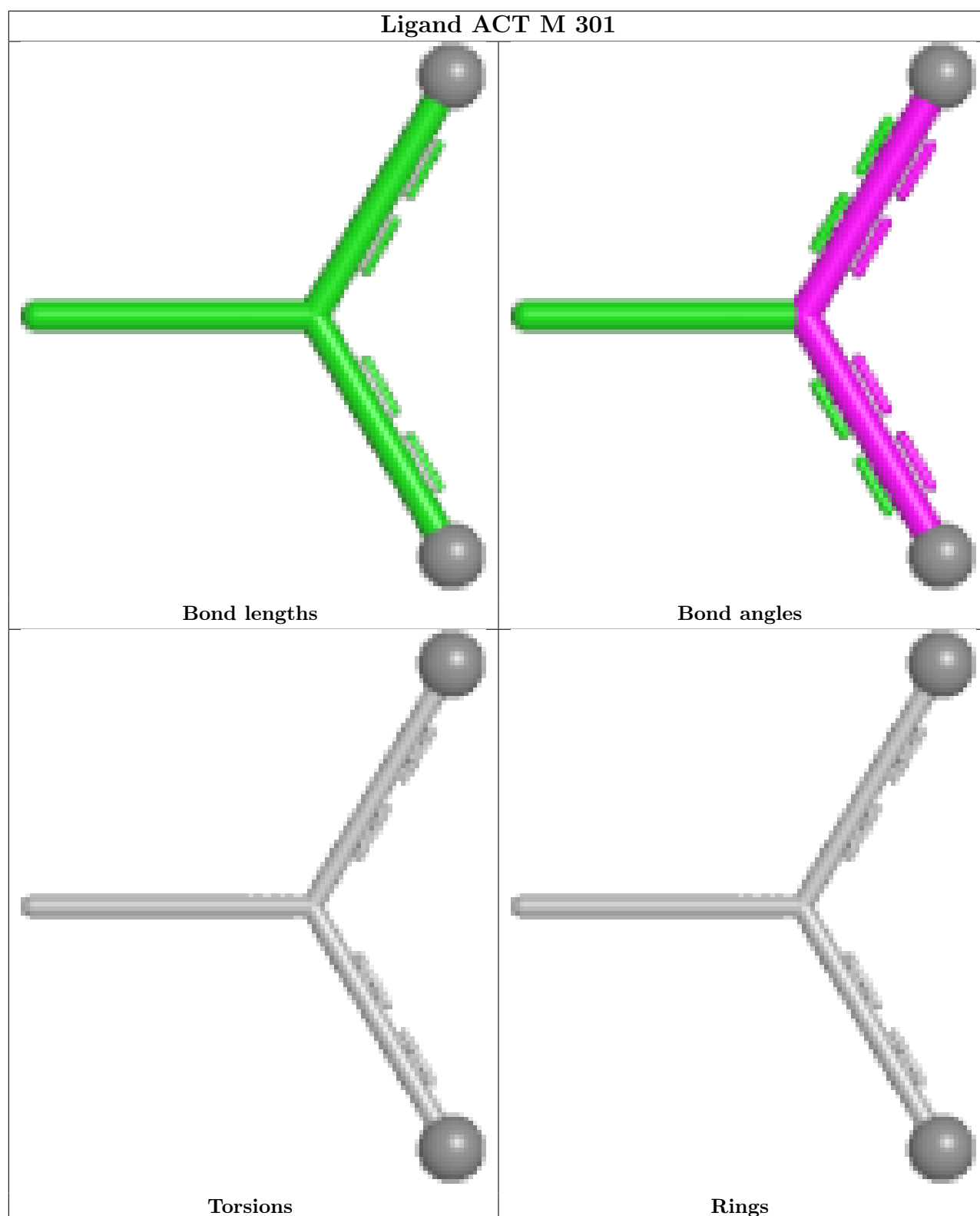
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	302	ACT	1	0
2	A	301	ACT	1	0
2	C	301	ACT	2	0
2	H	301	ACT	1	0

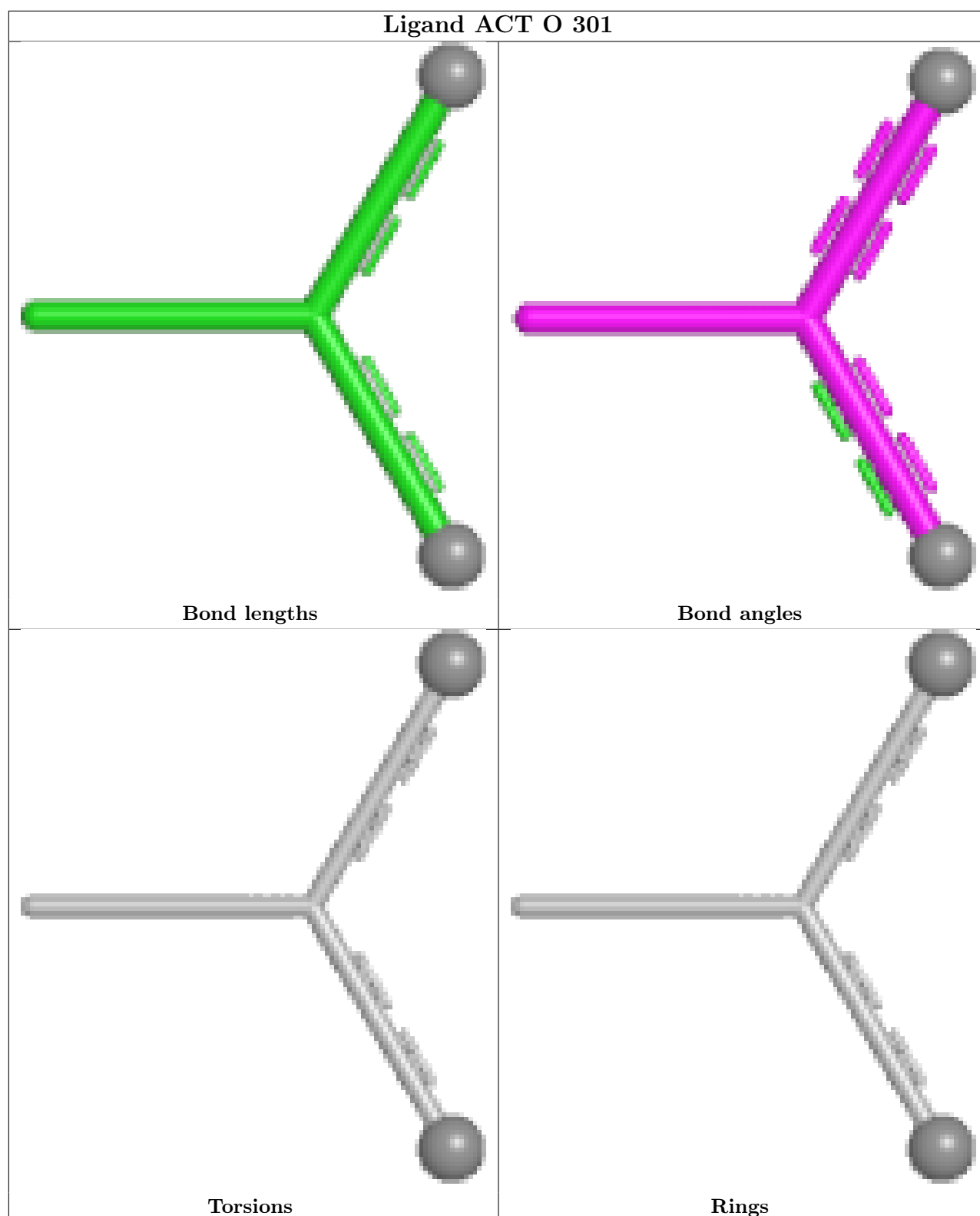
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

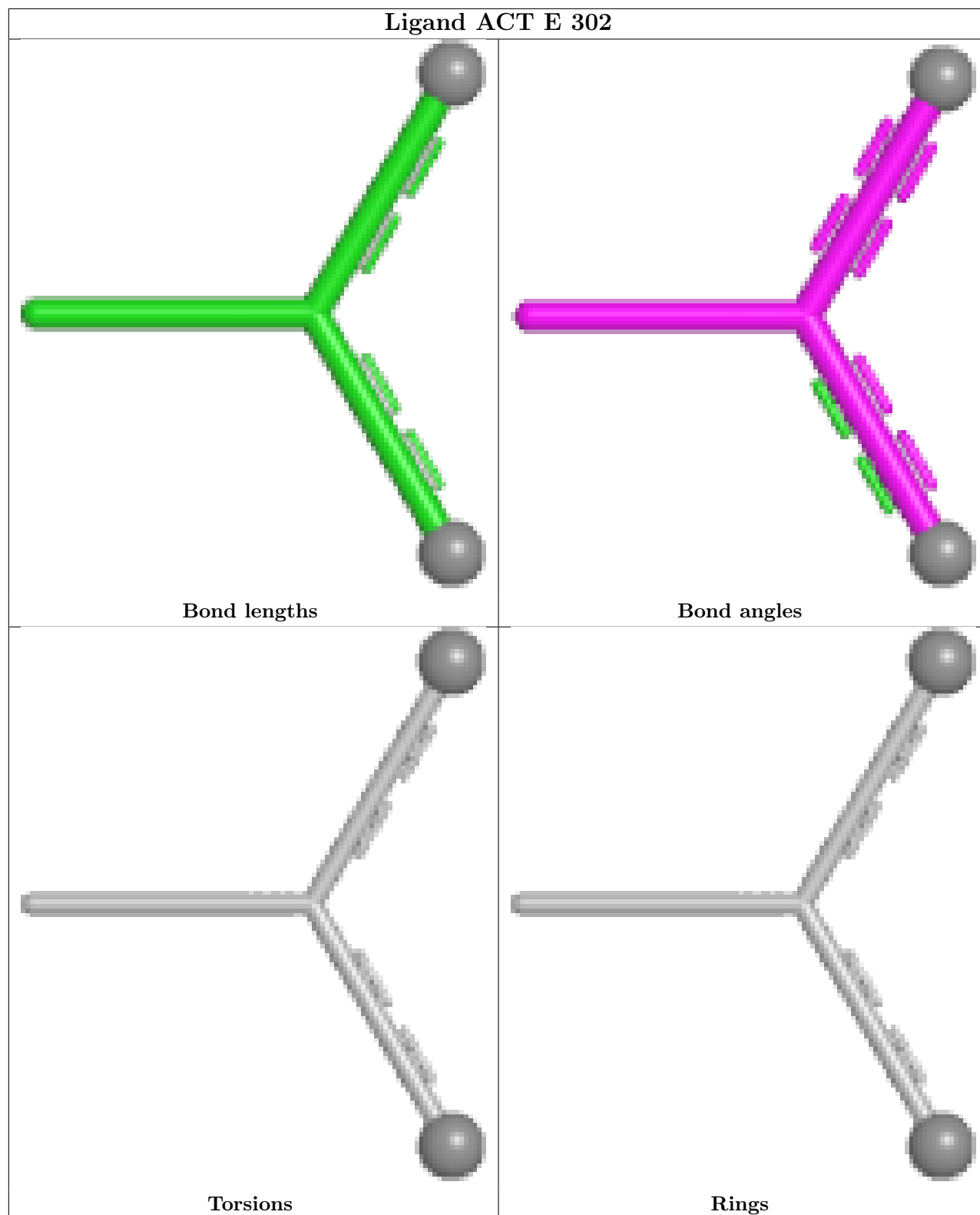


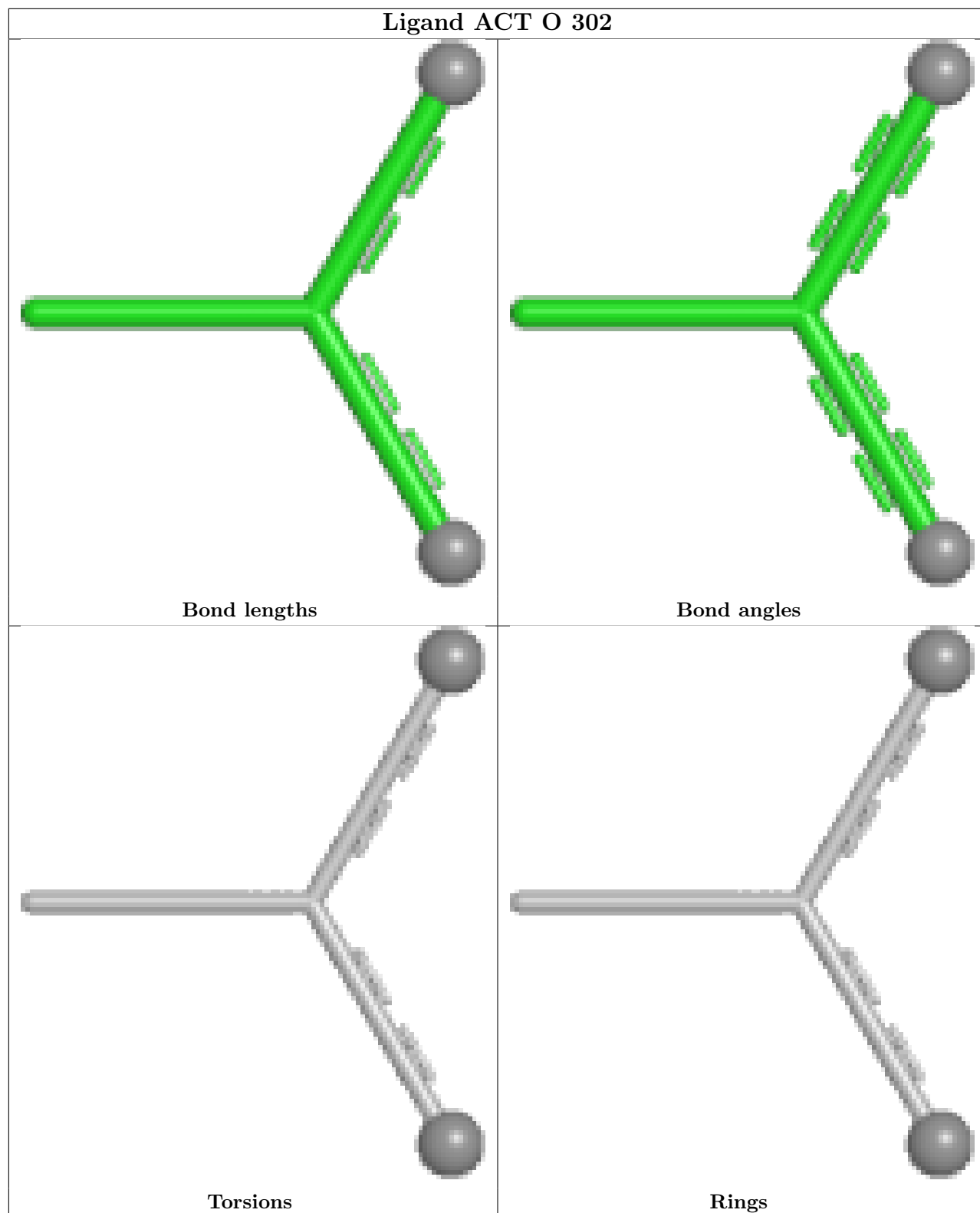


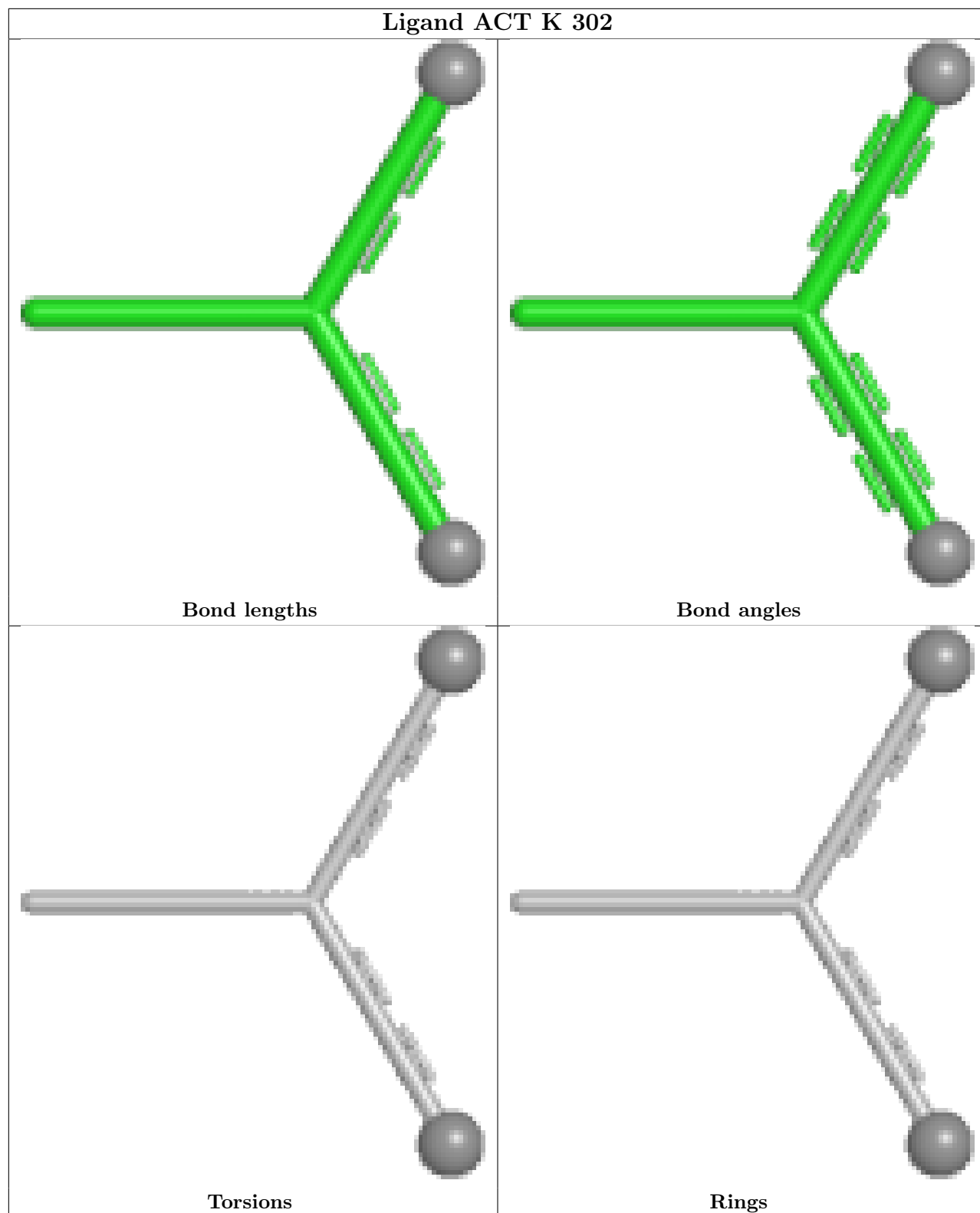


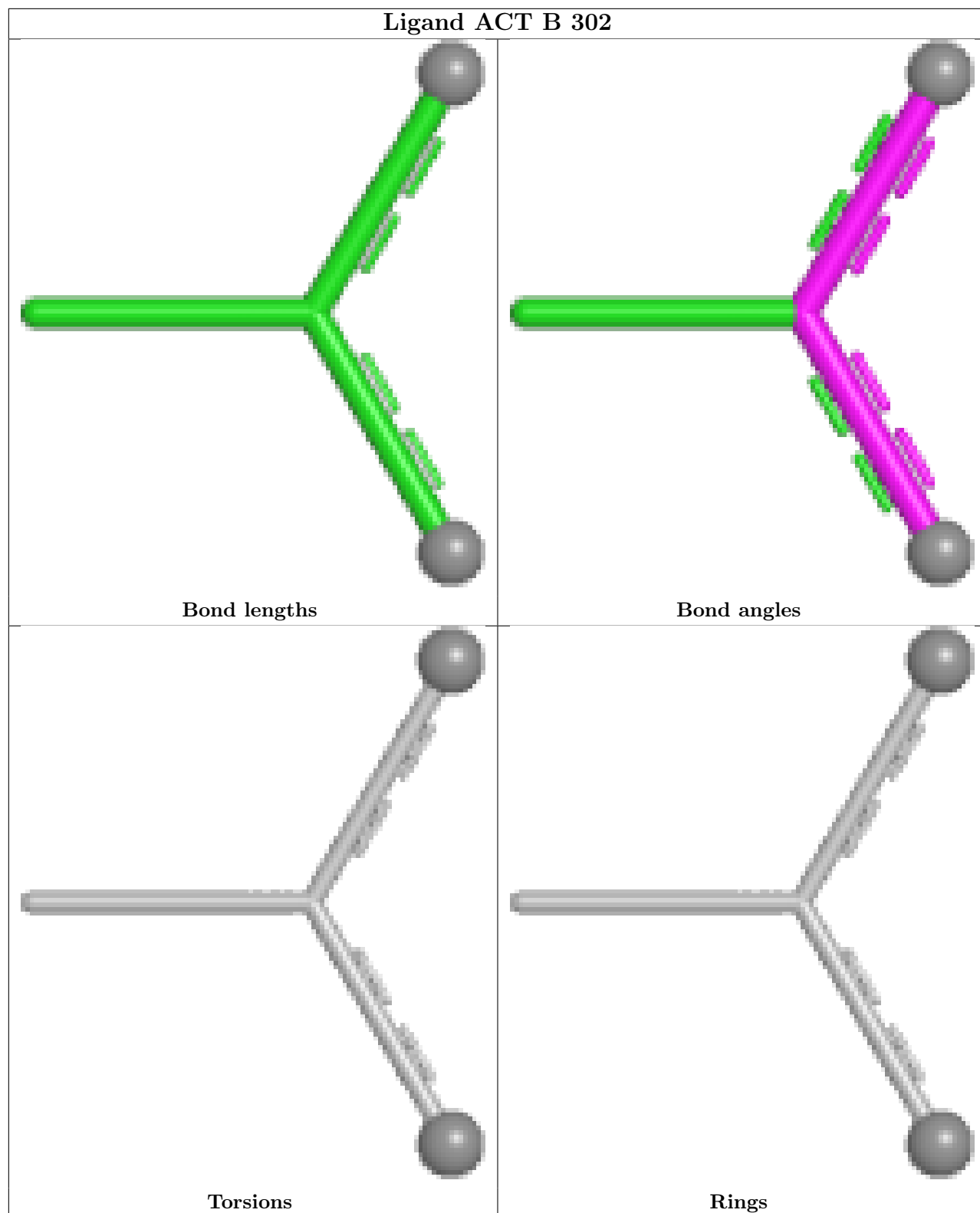


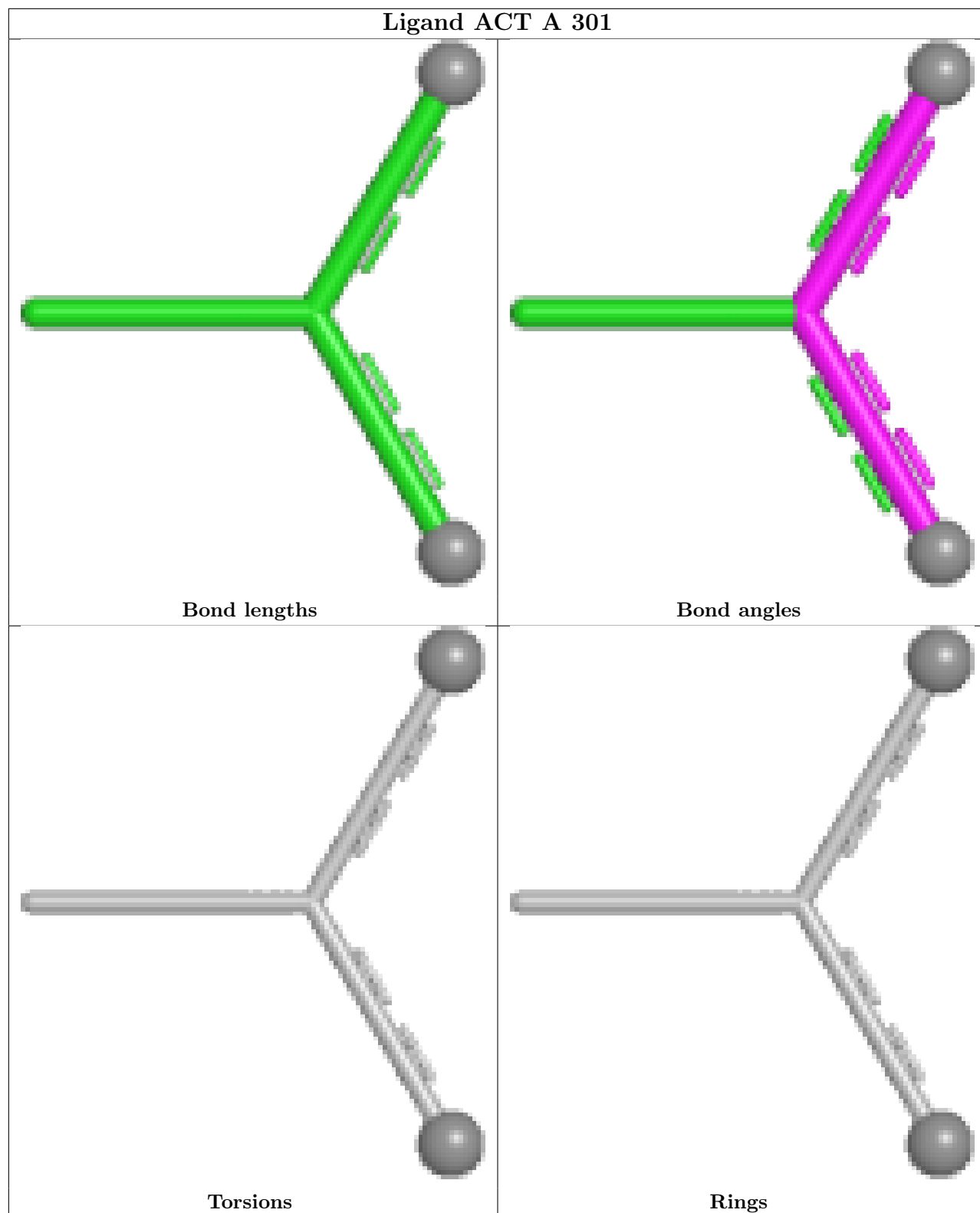


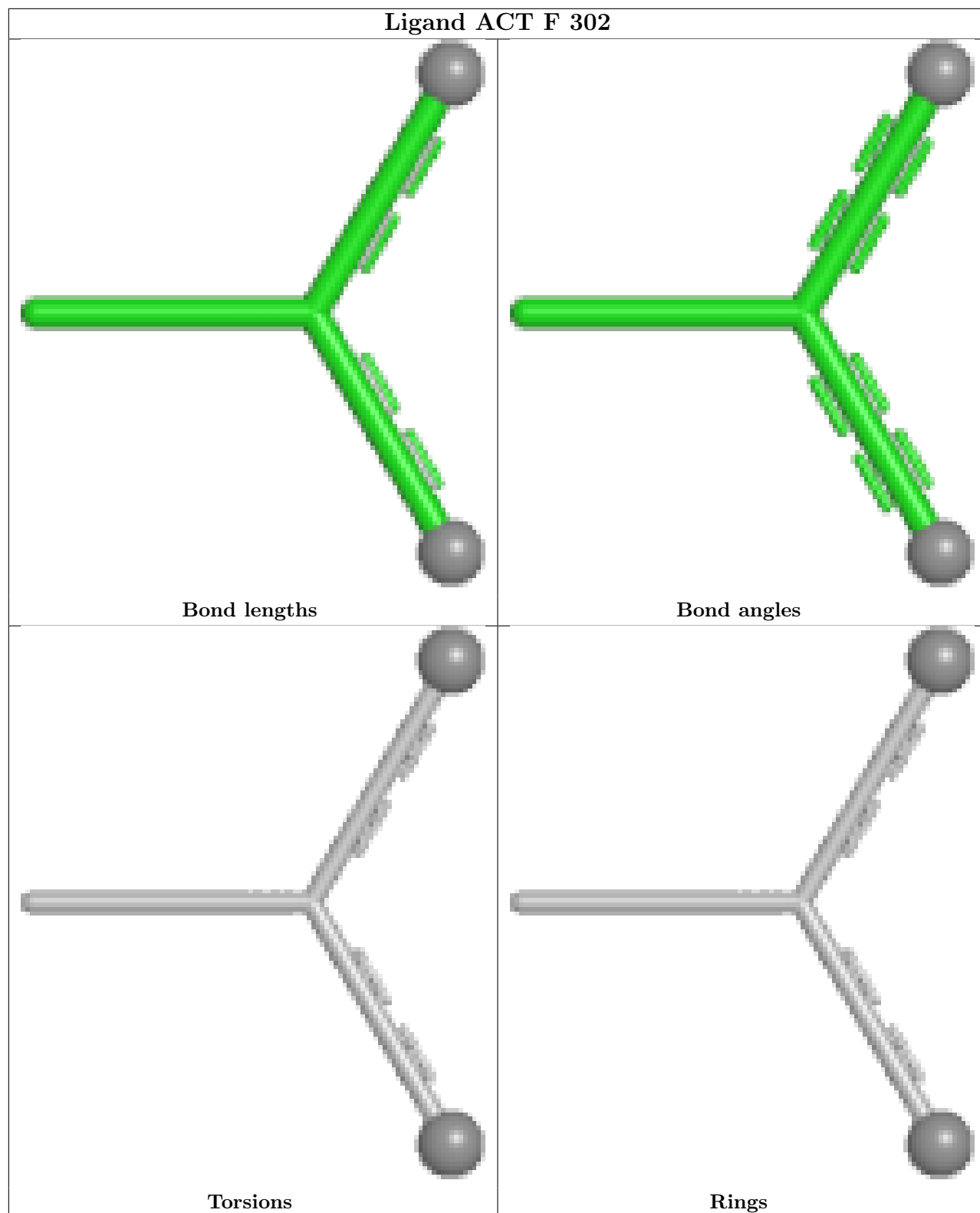


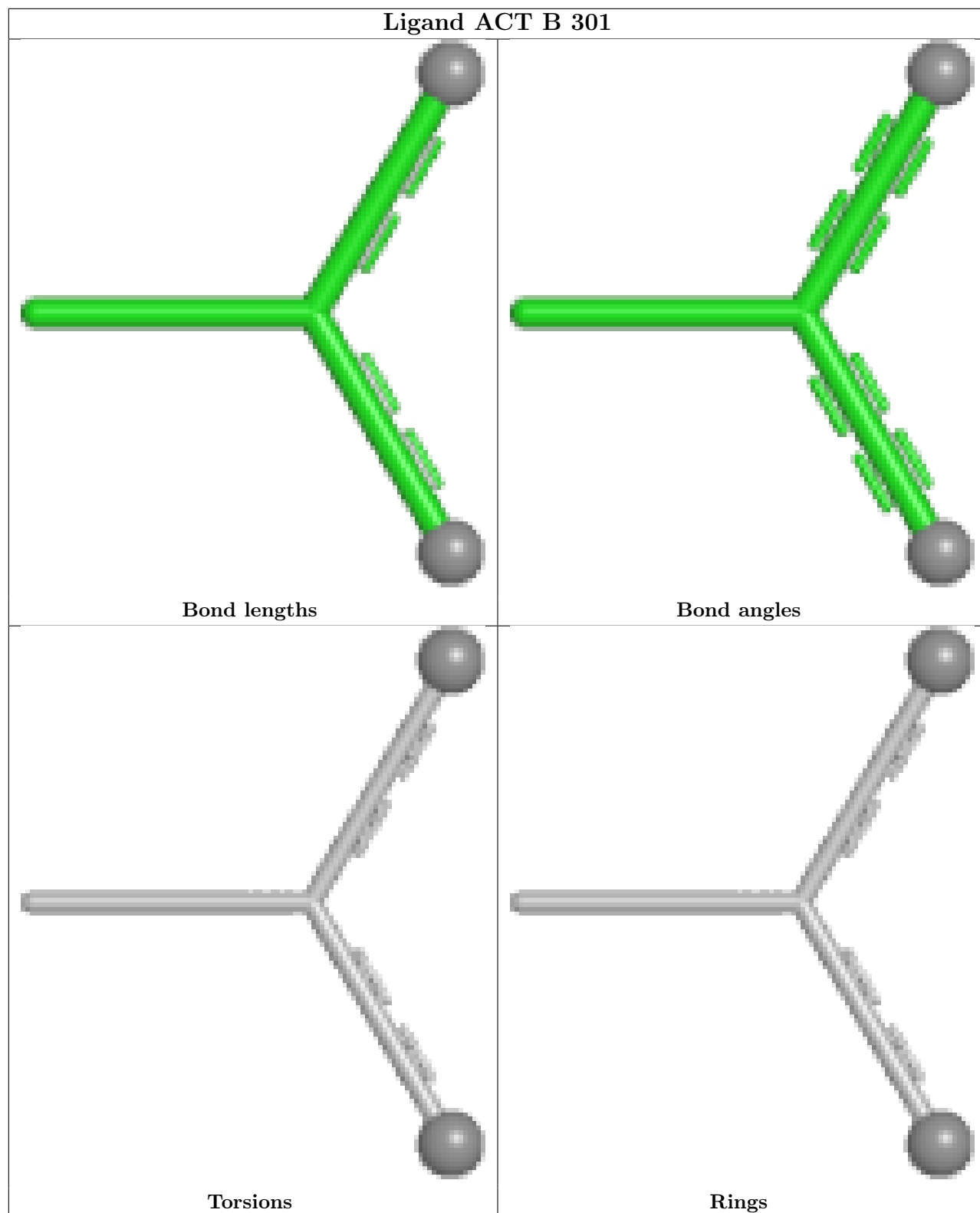


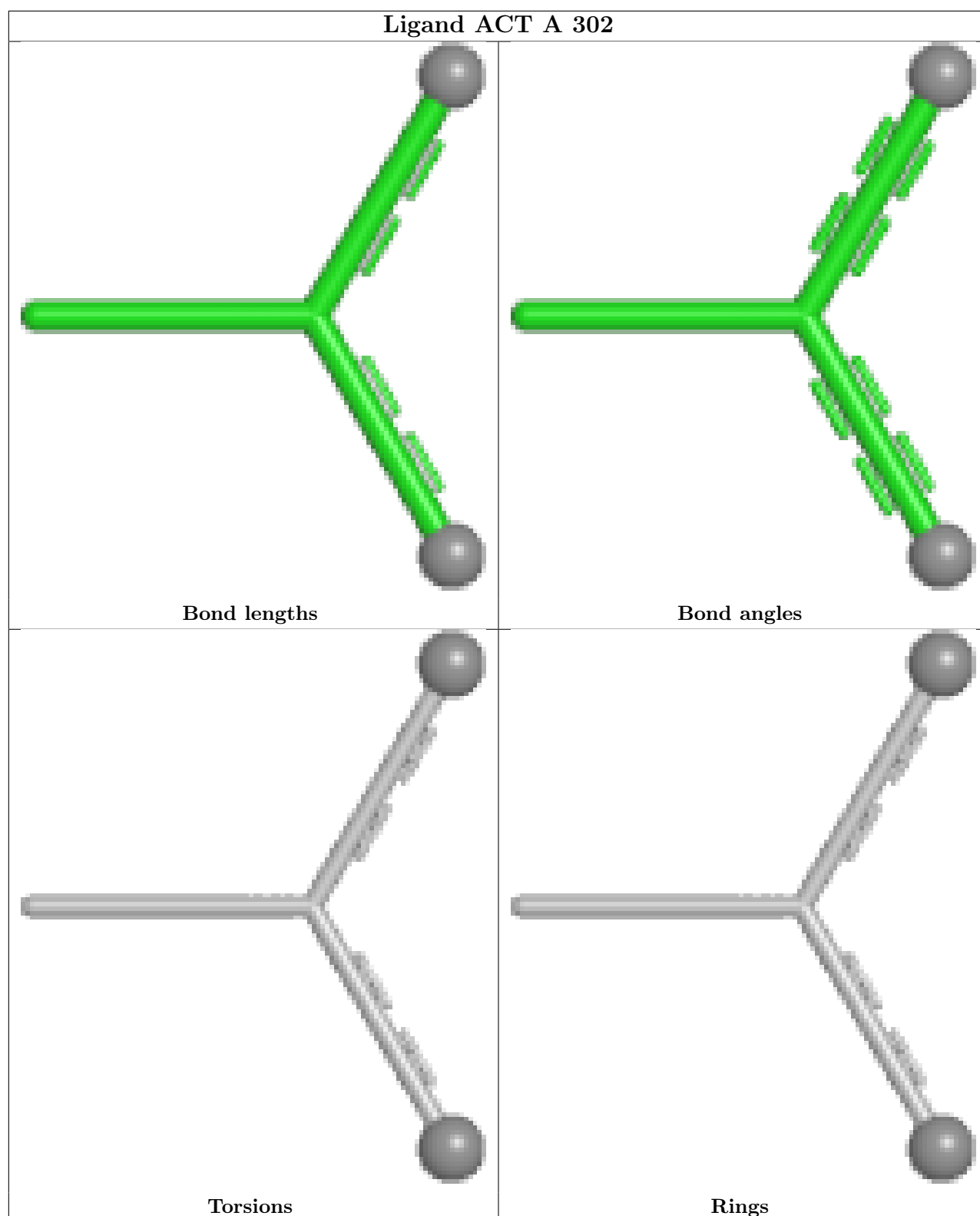


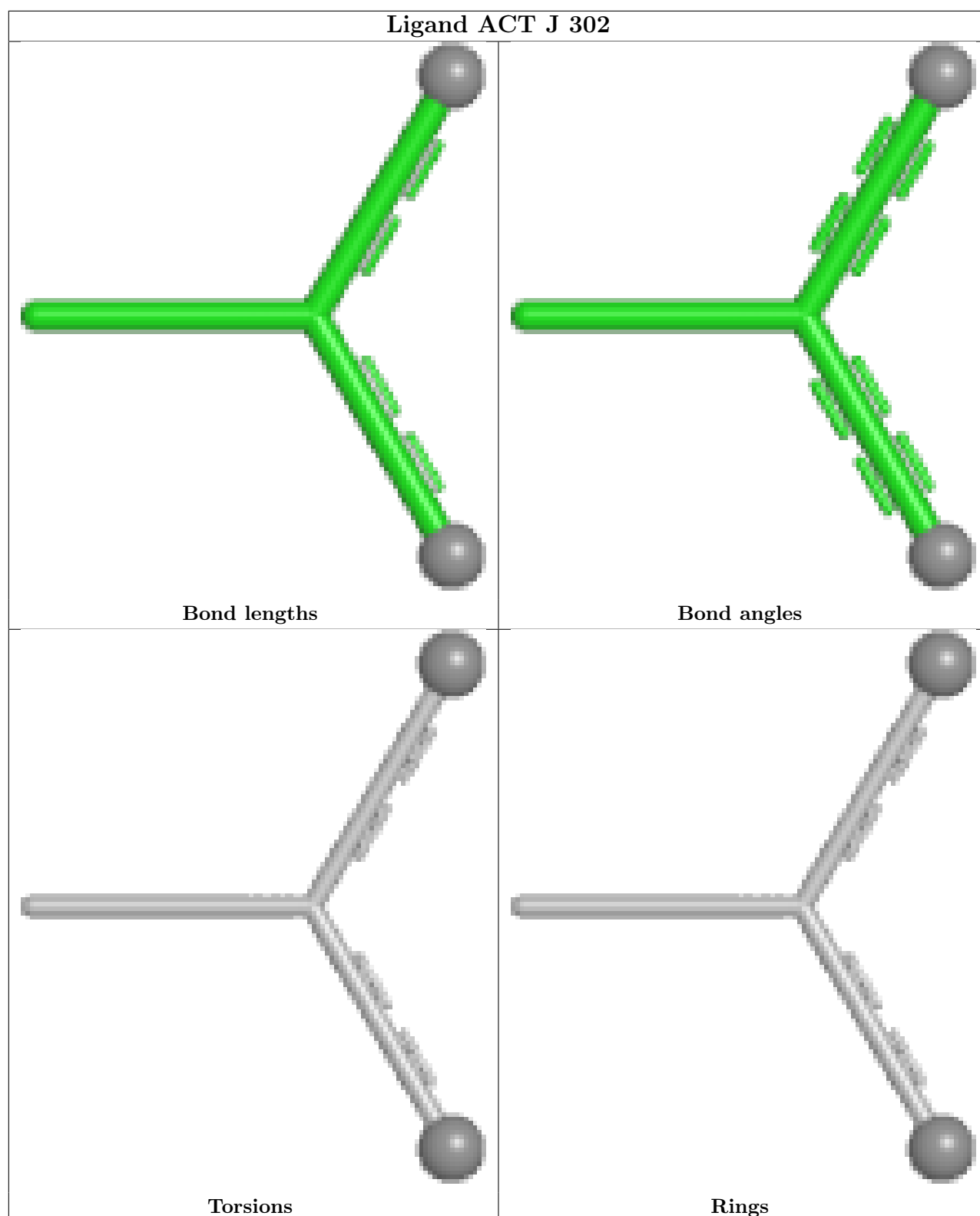


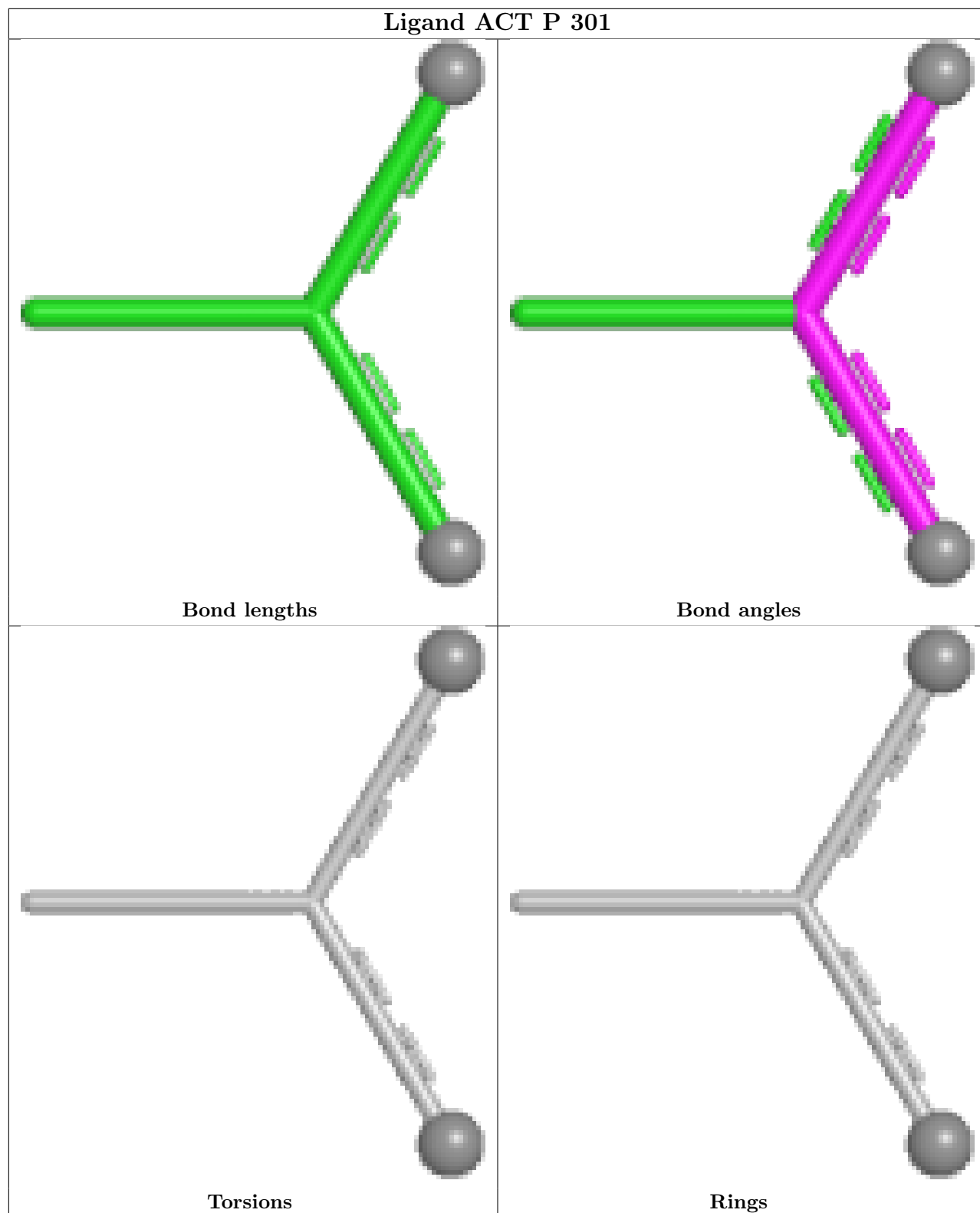


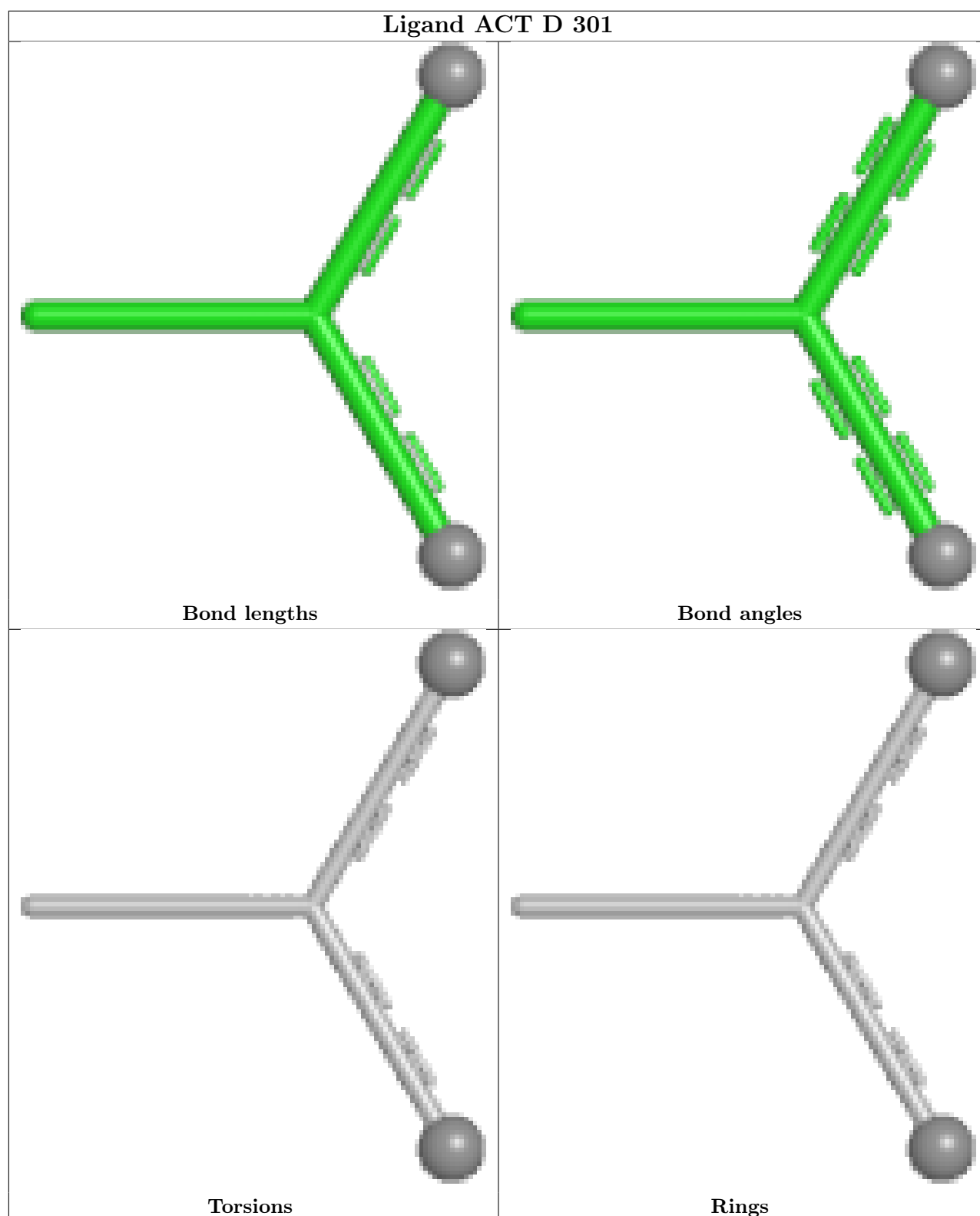


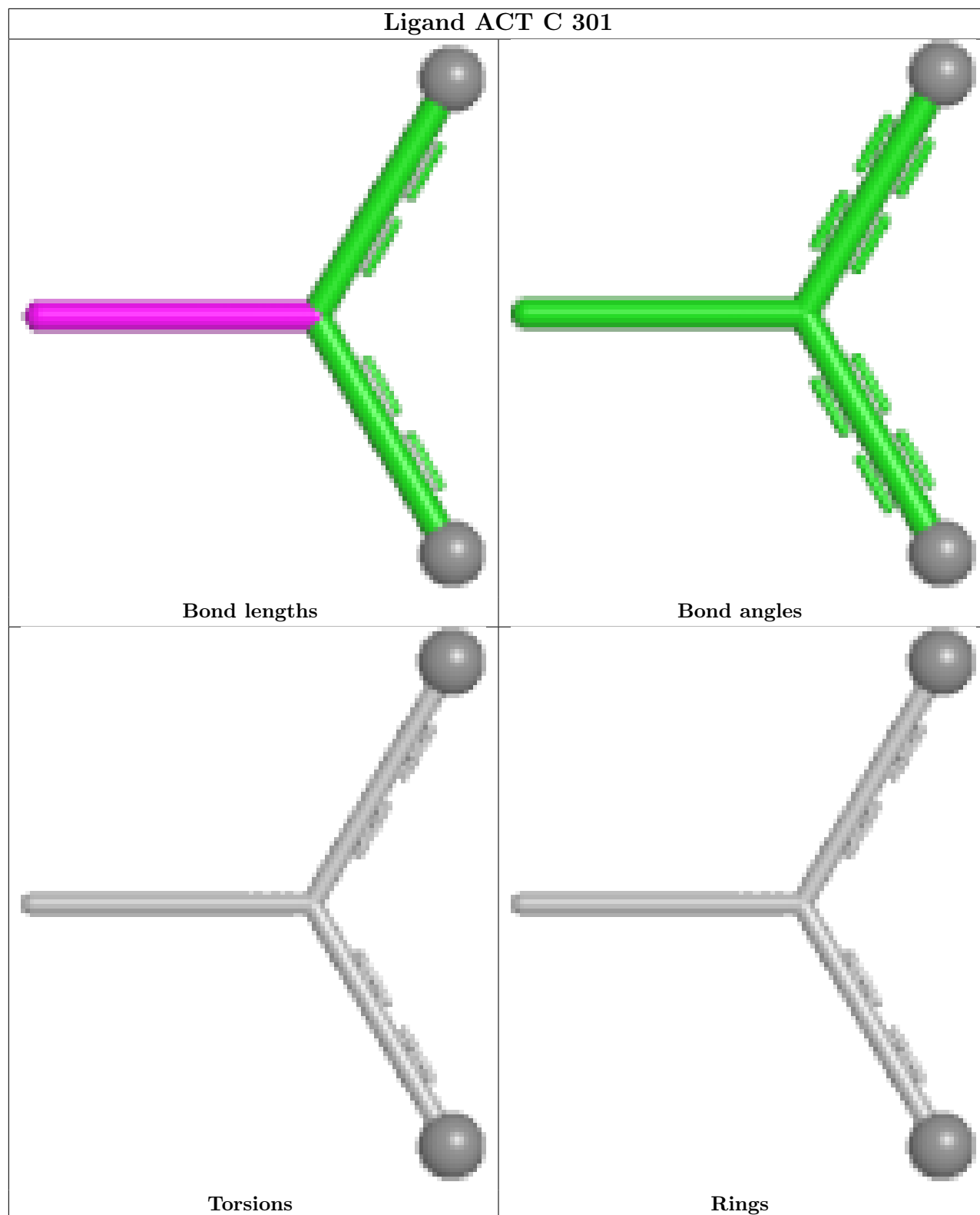


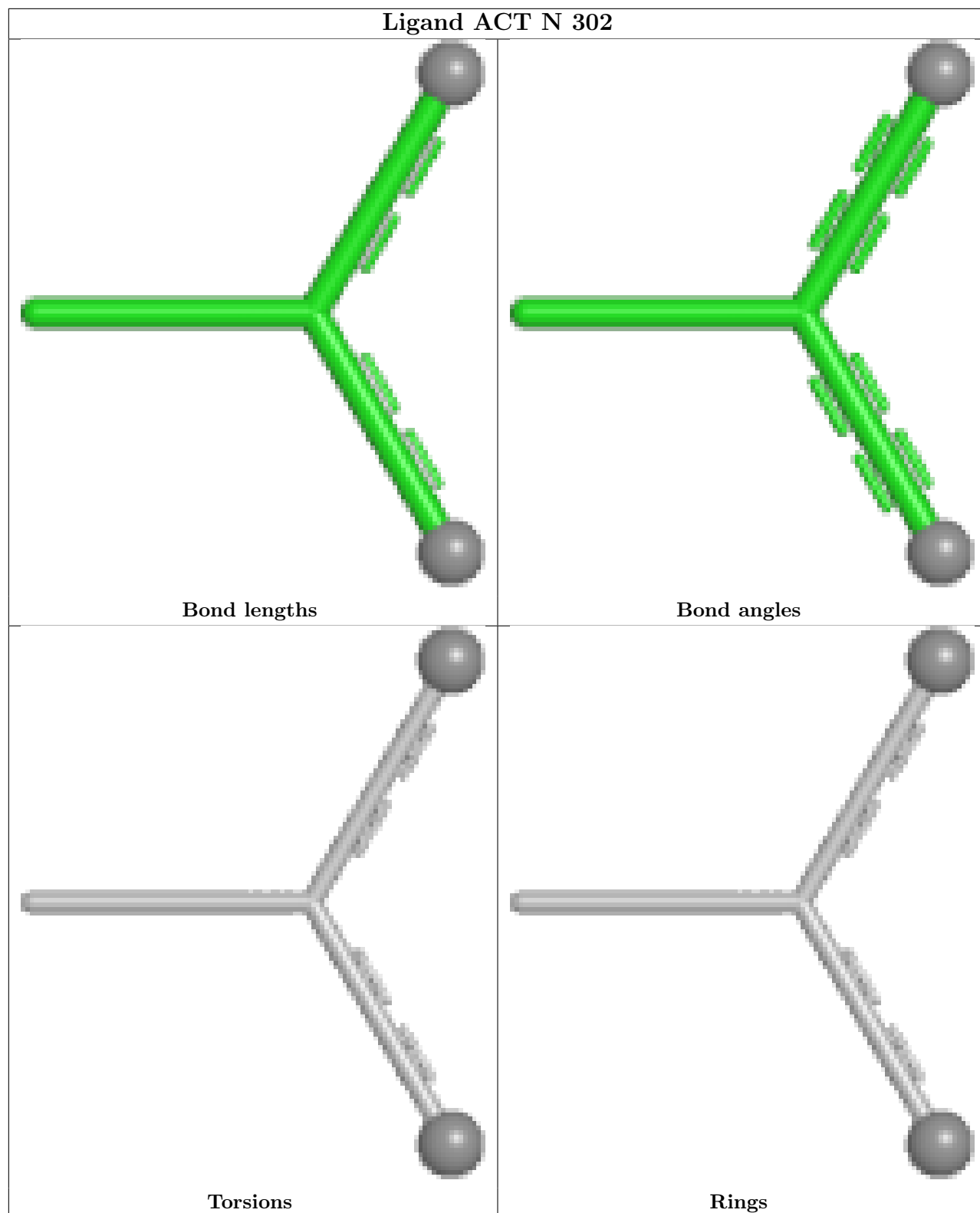


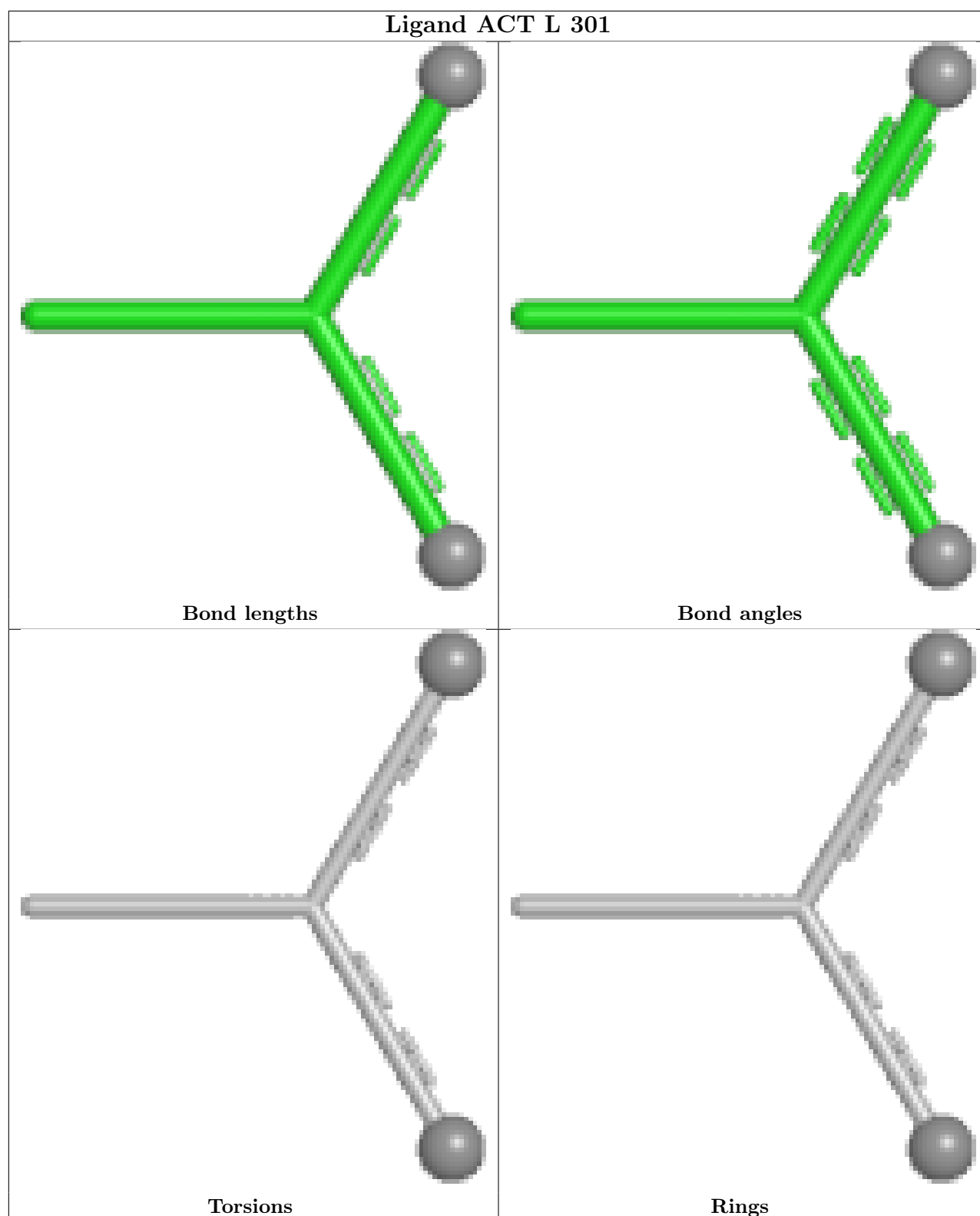


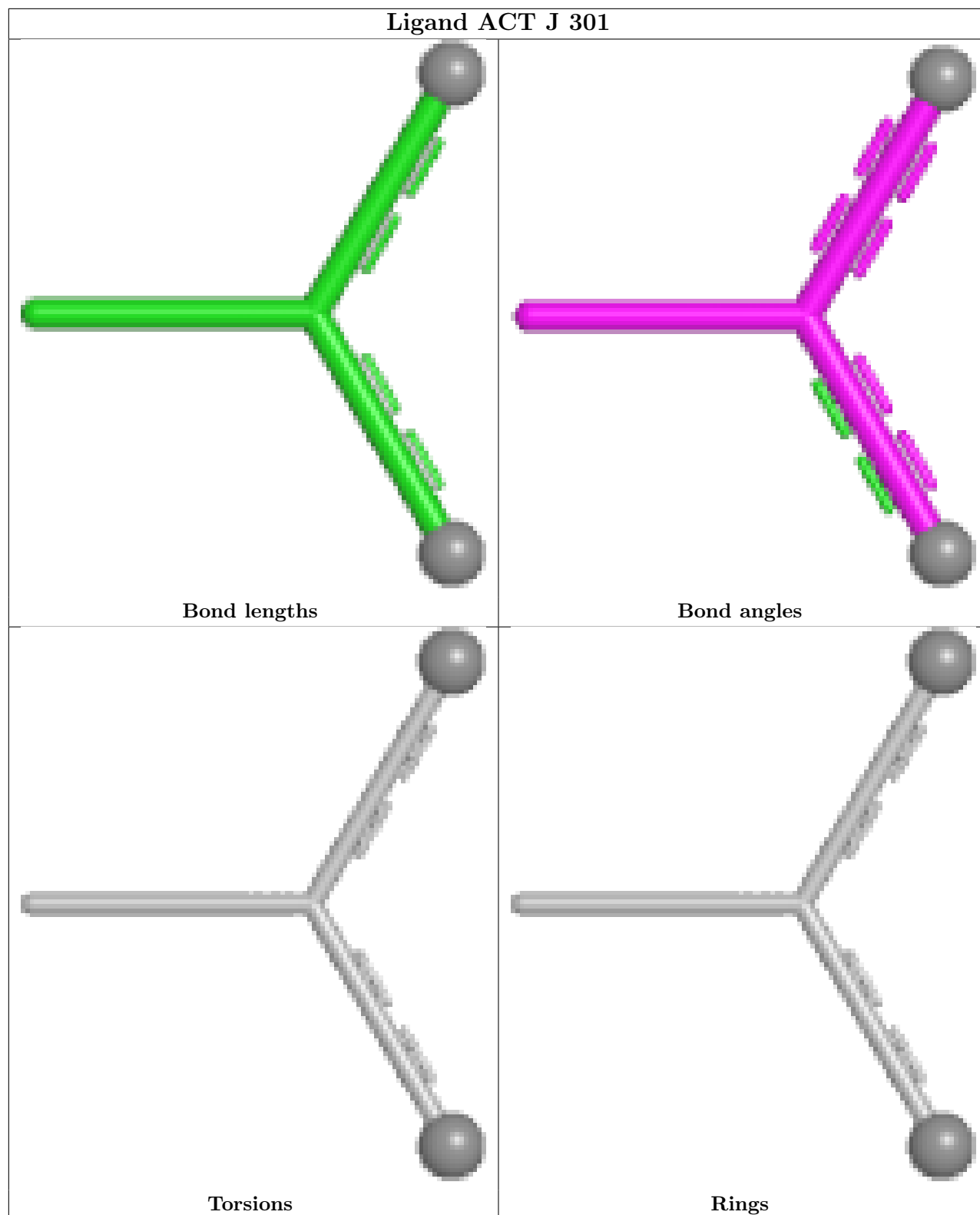


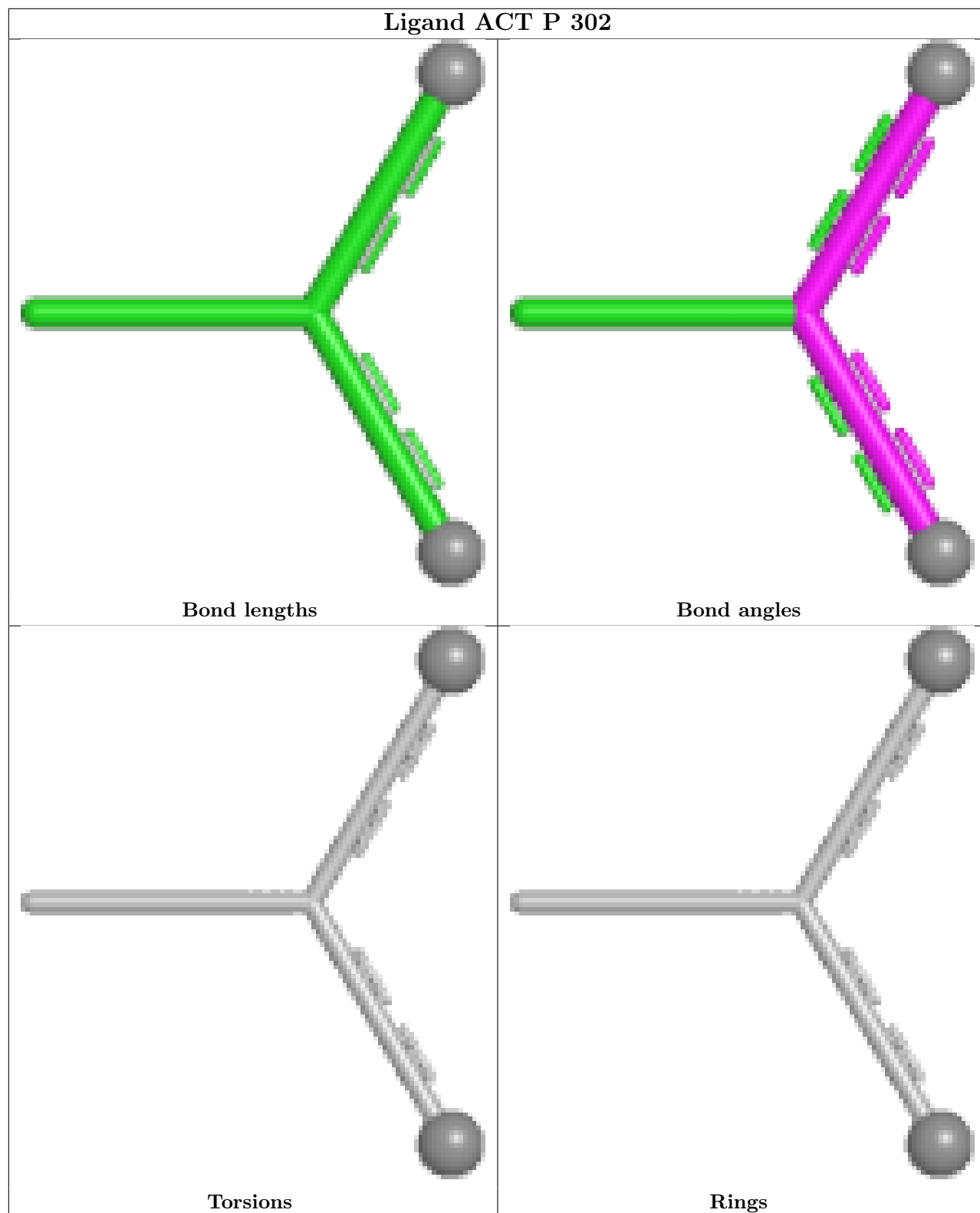


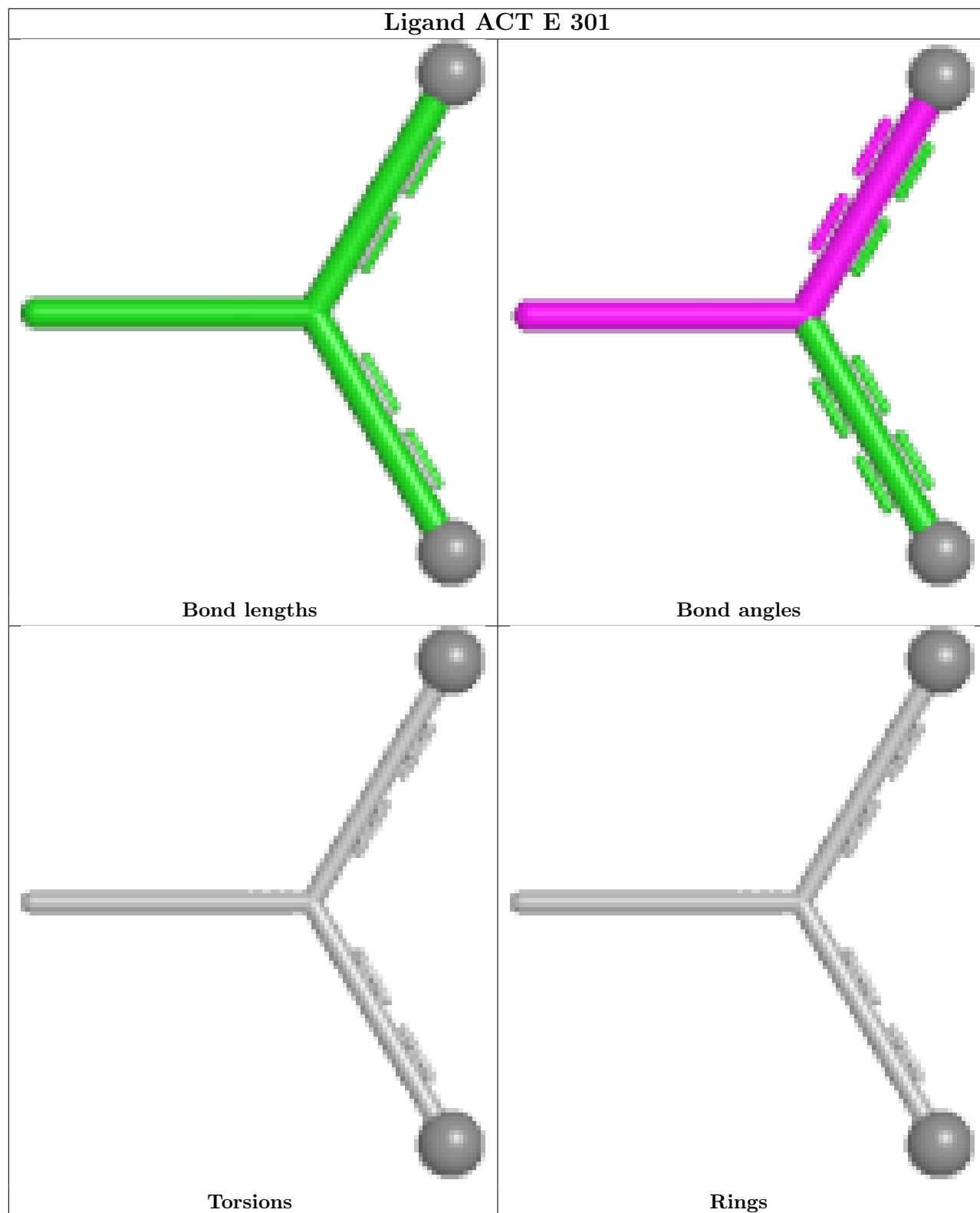


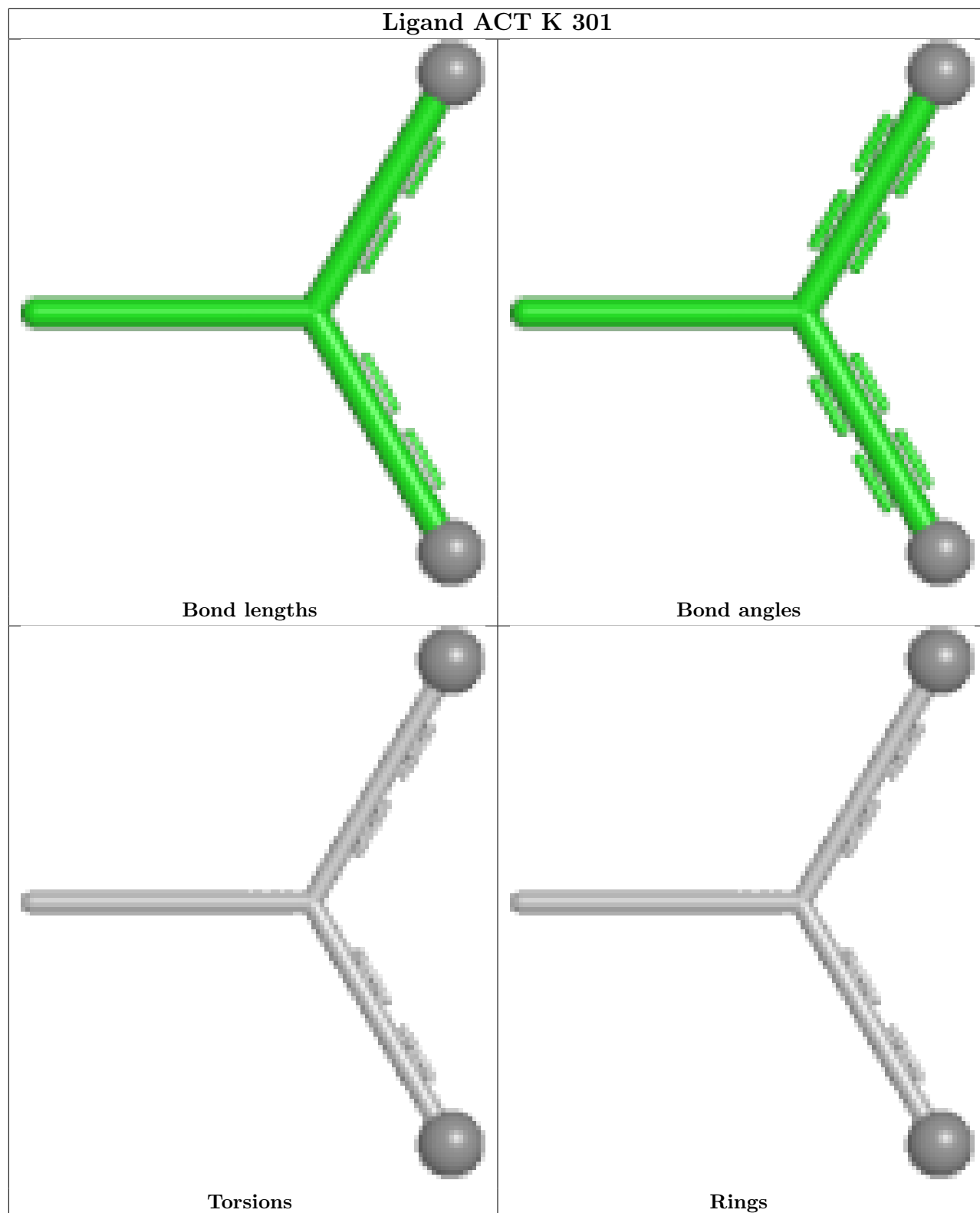


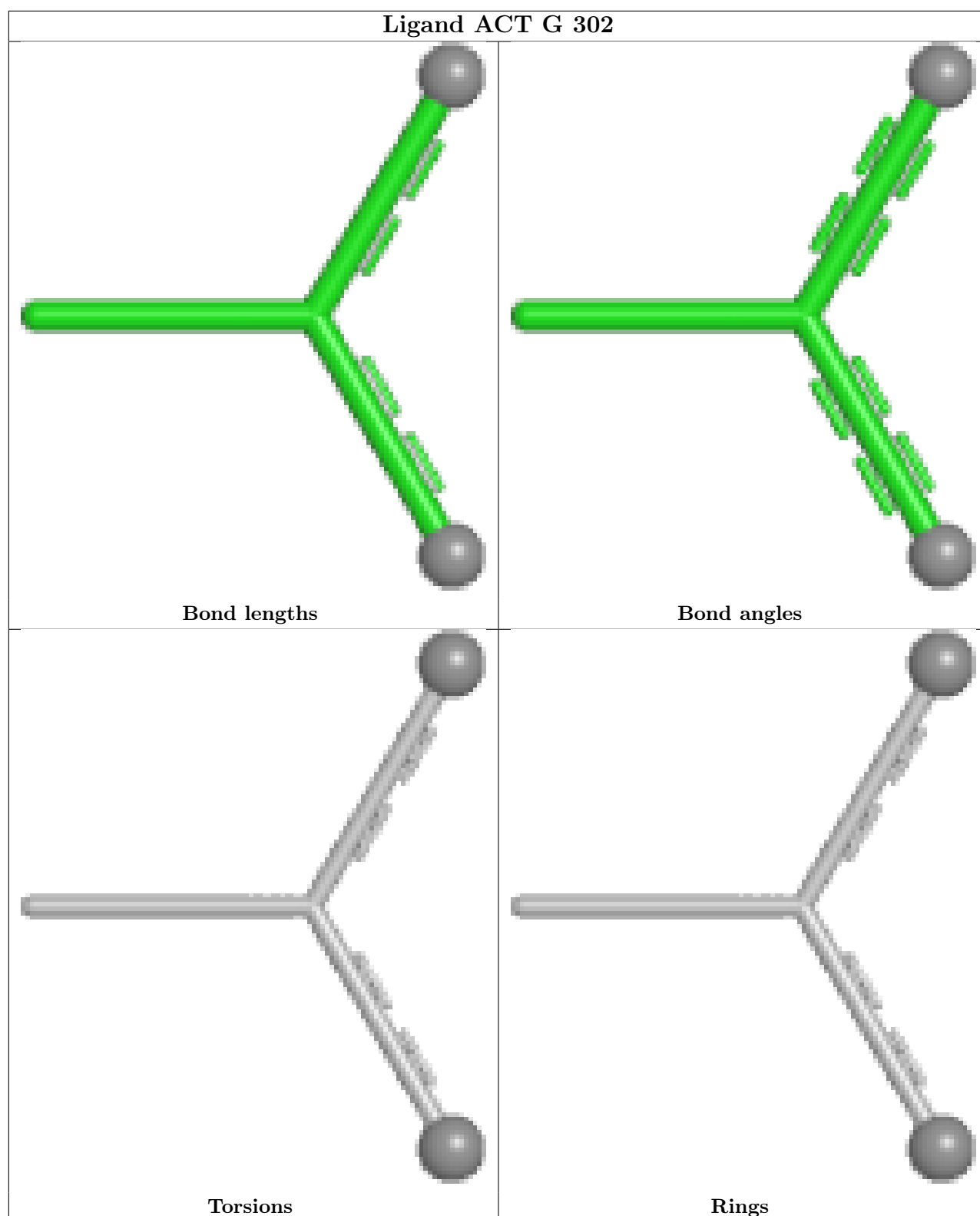


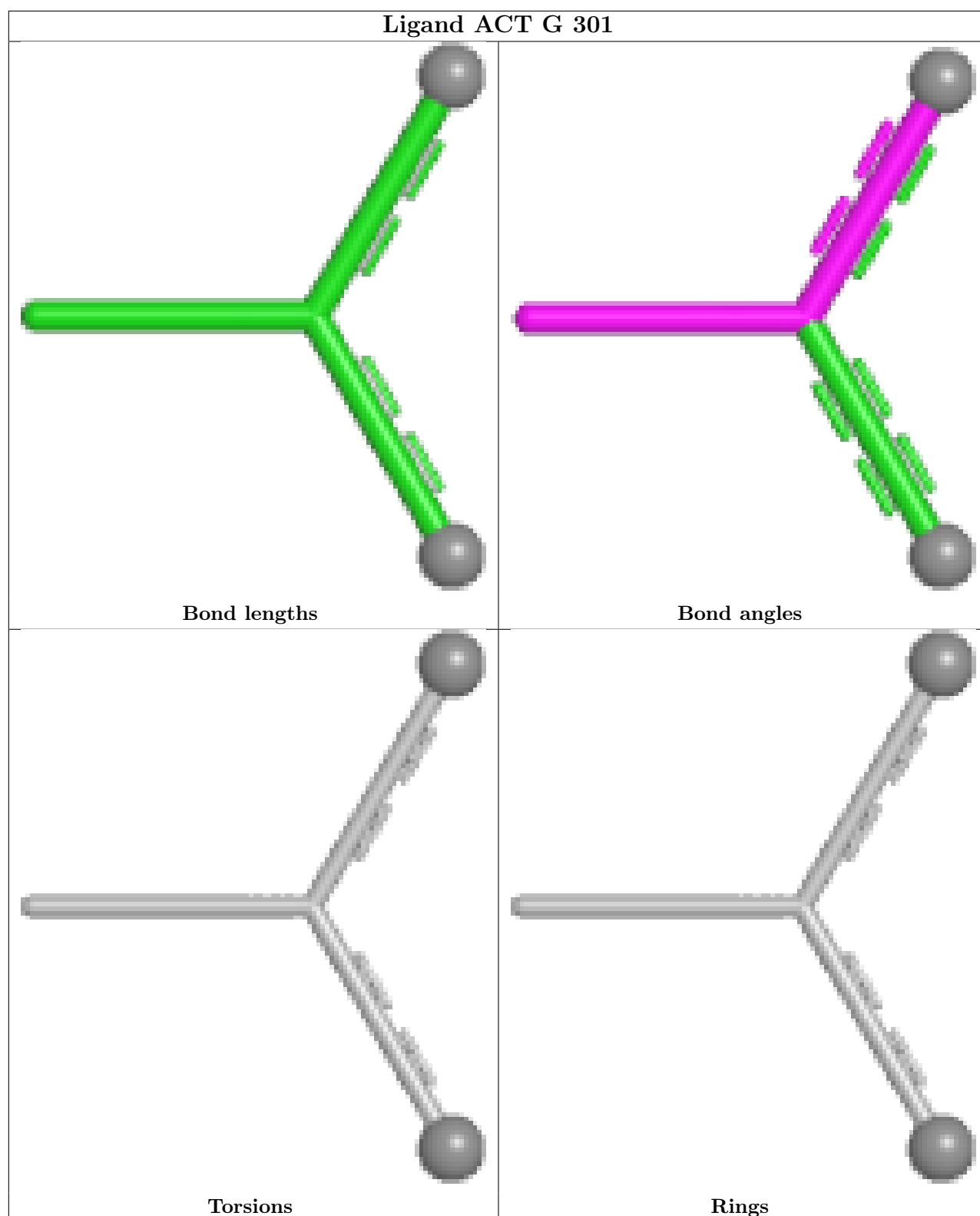




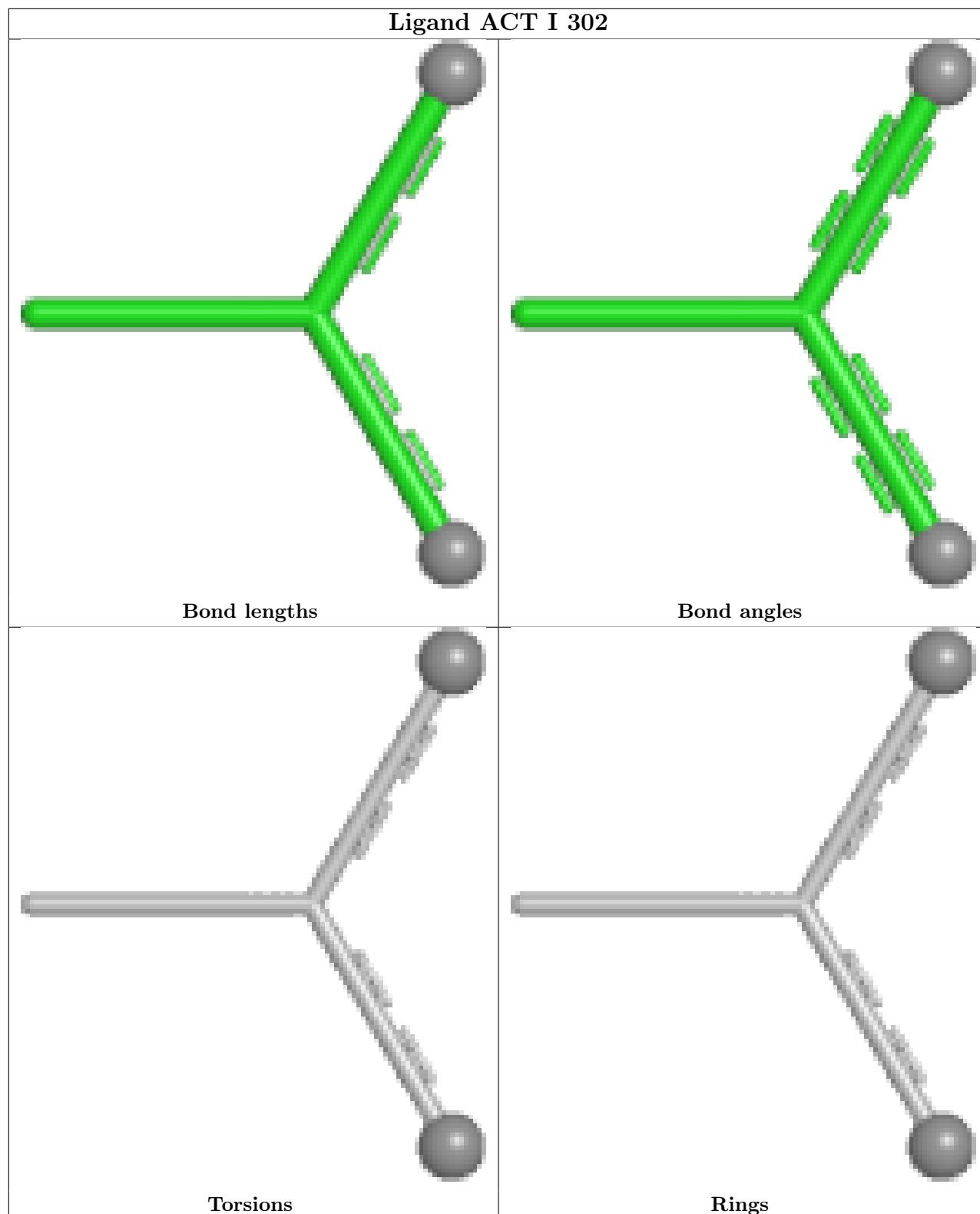


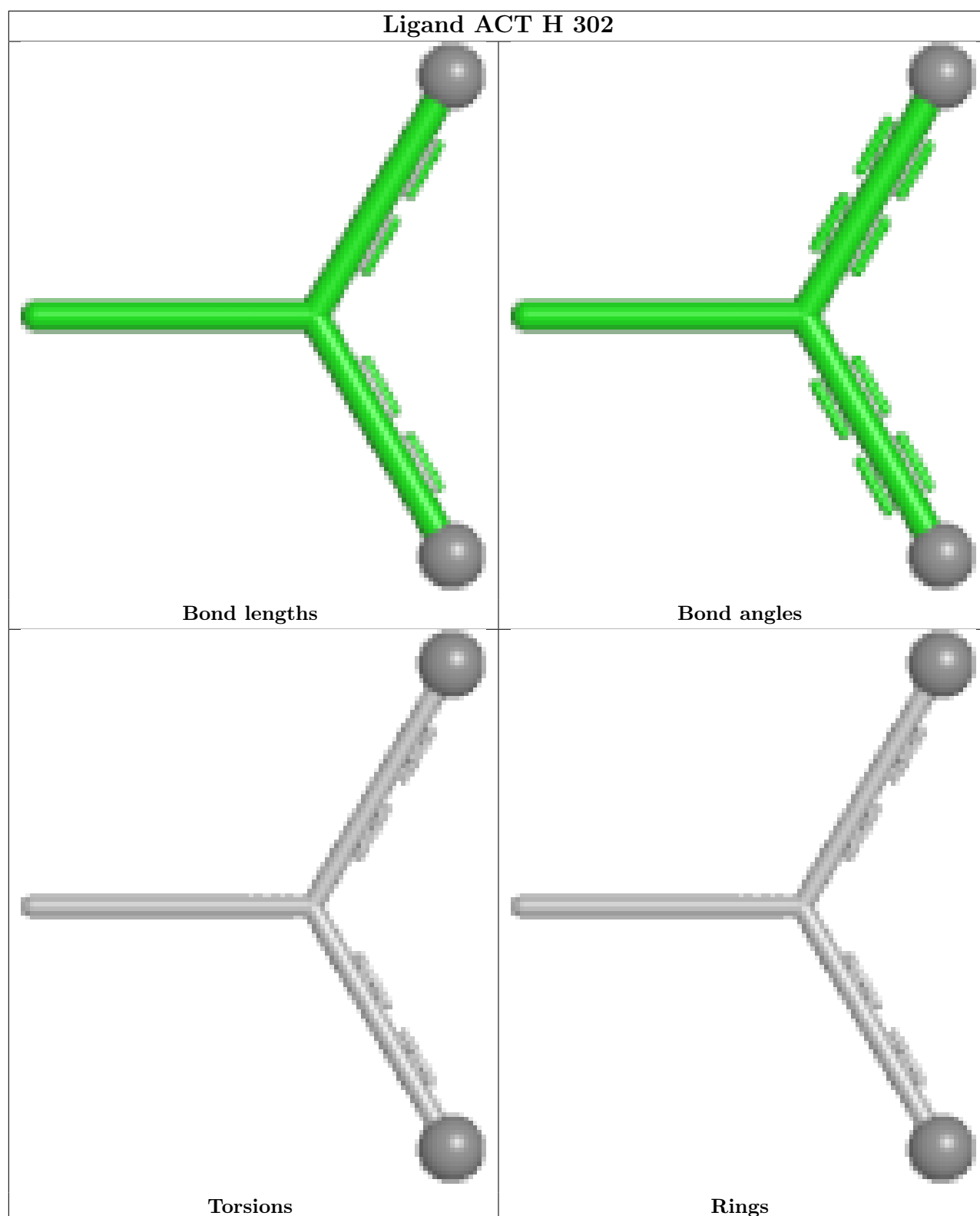


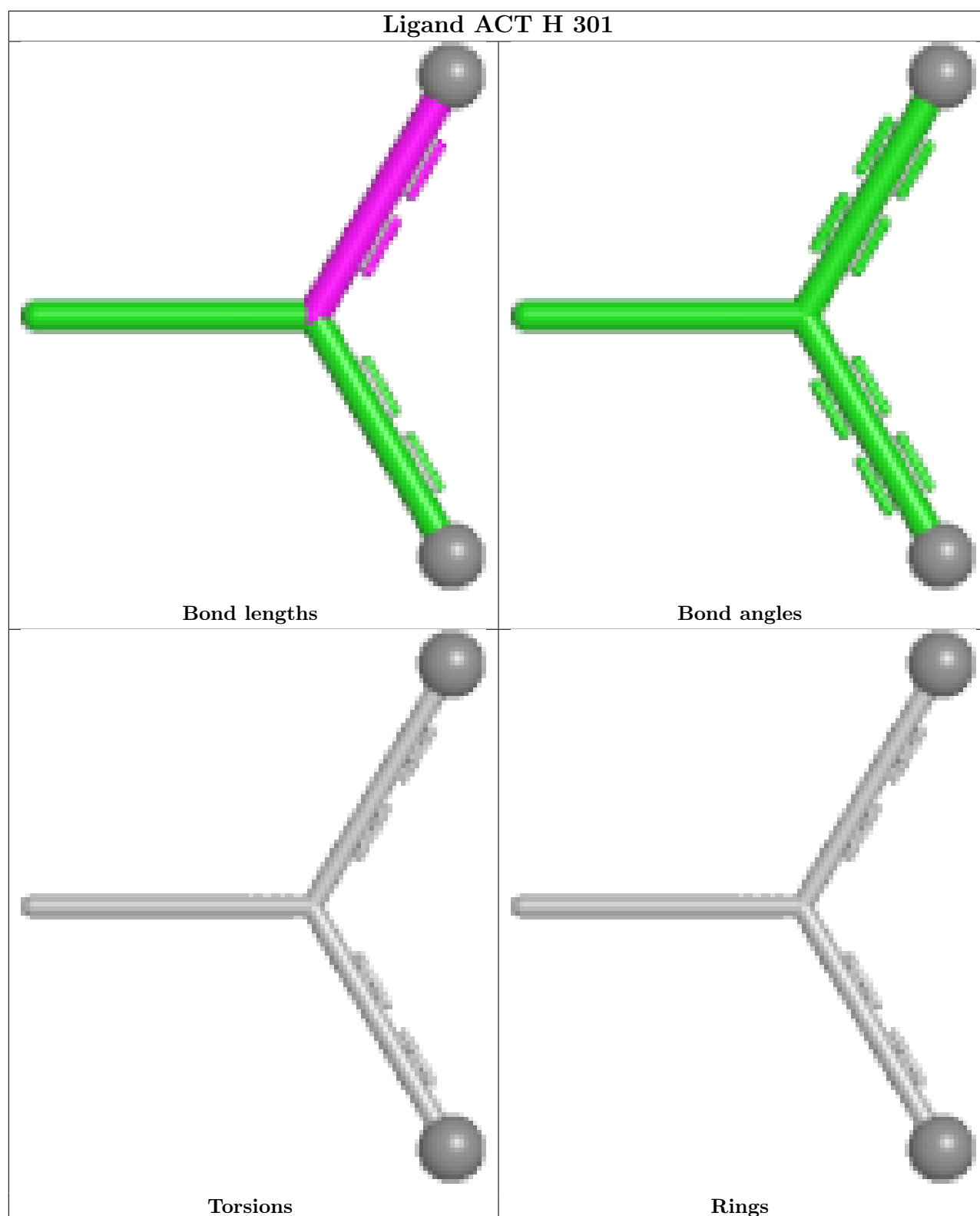


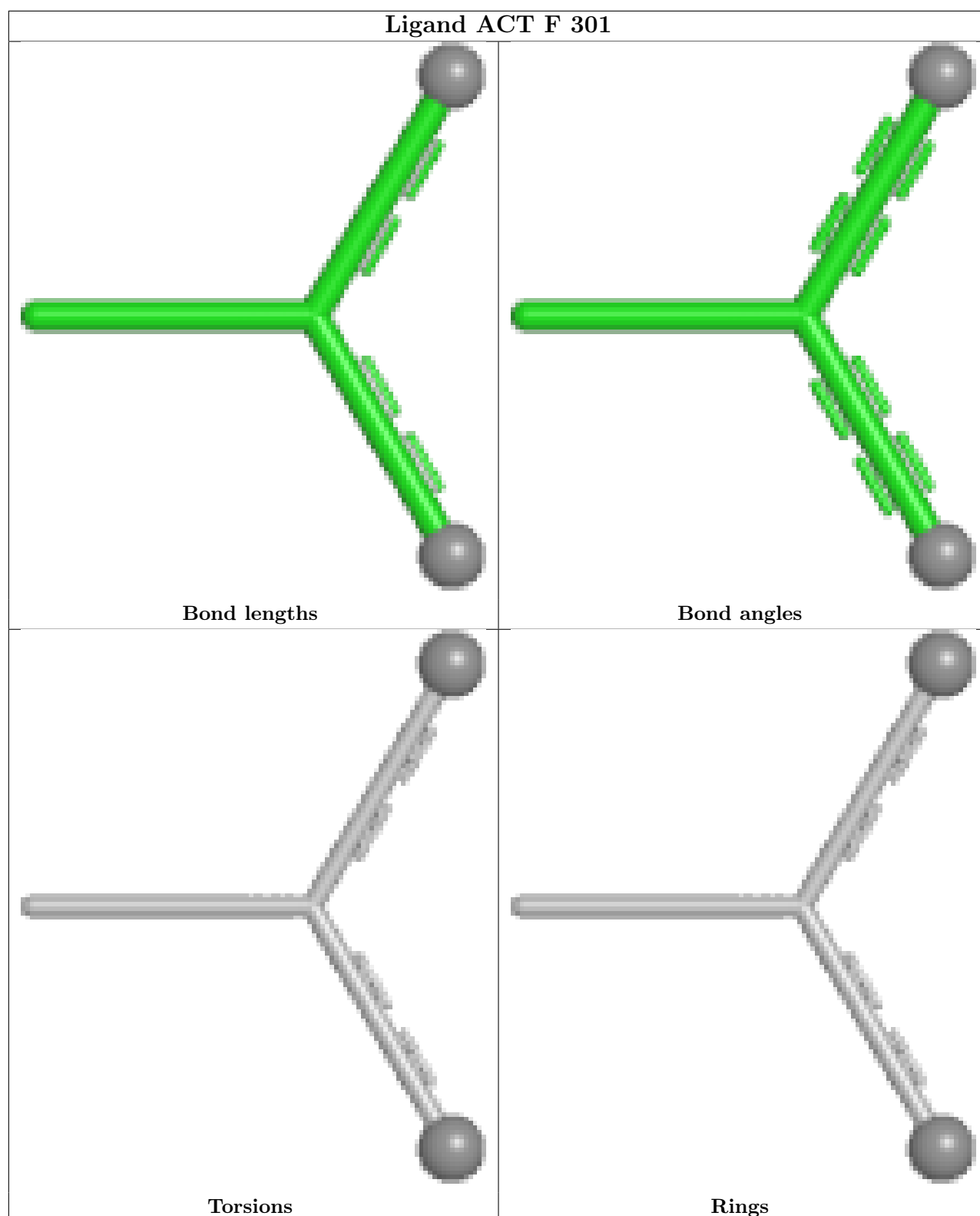


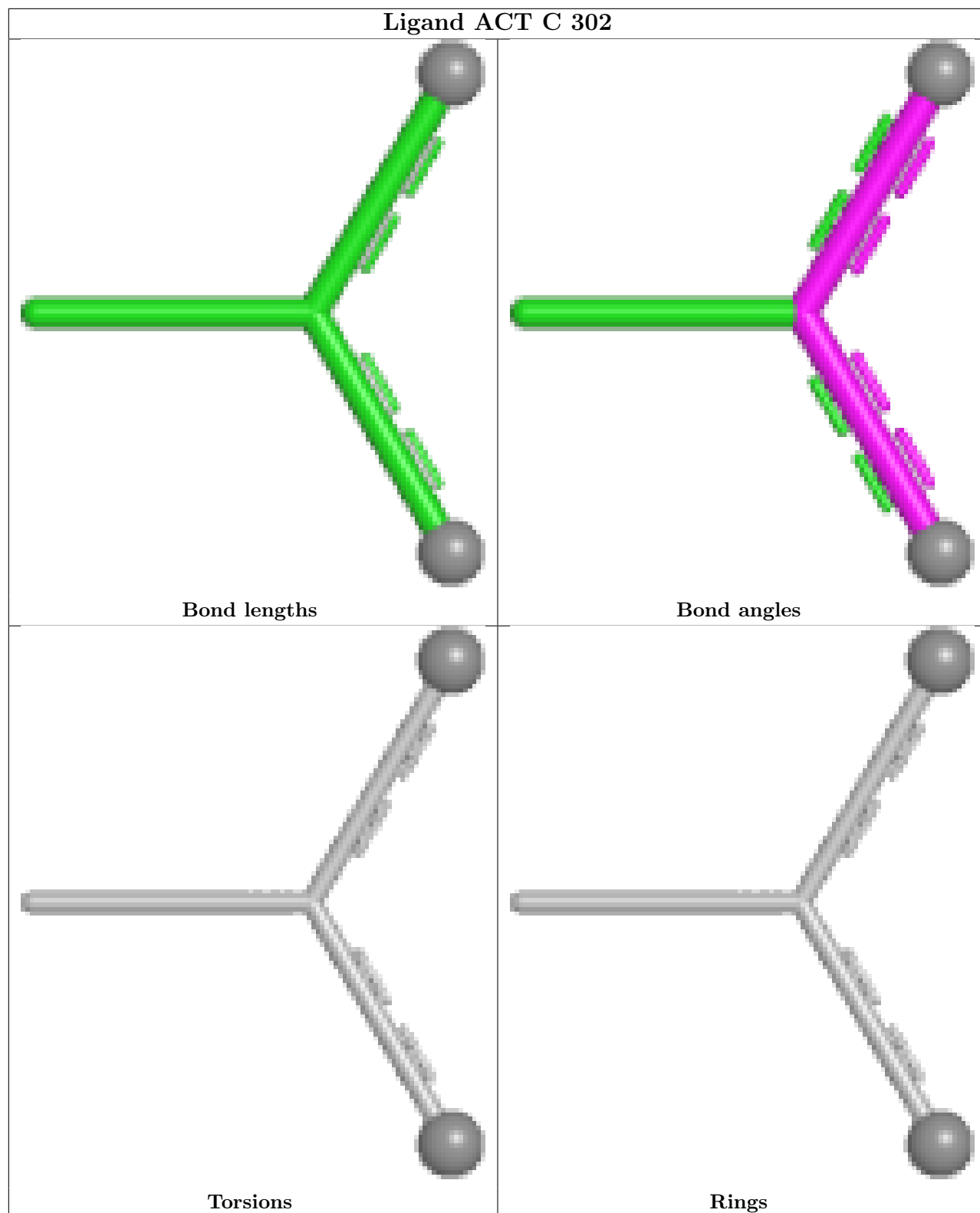
Ligand ACT I 302

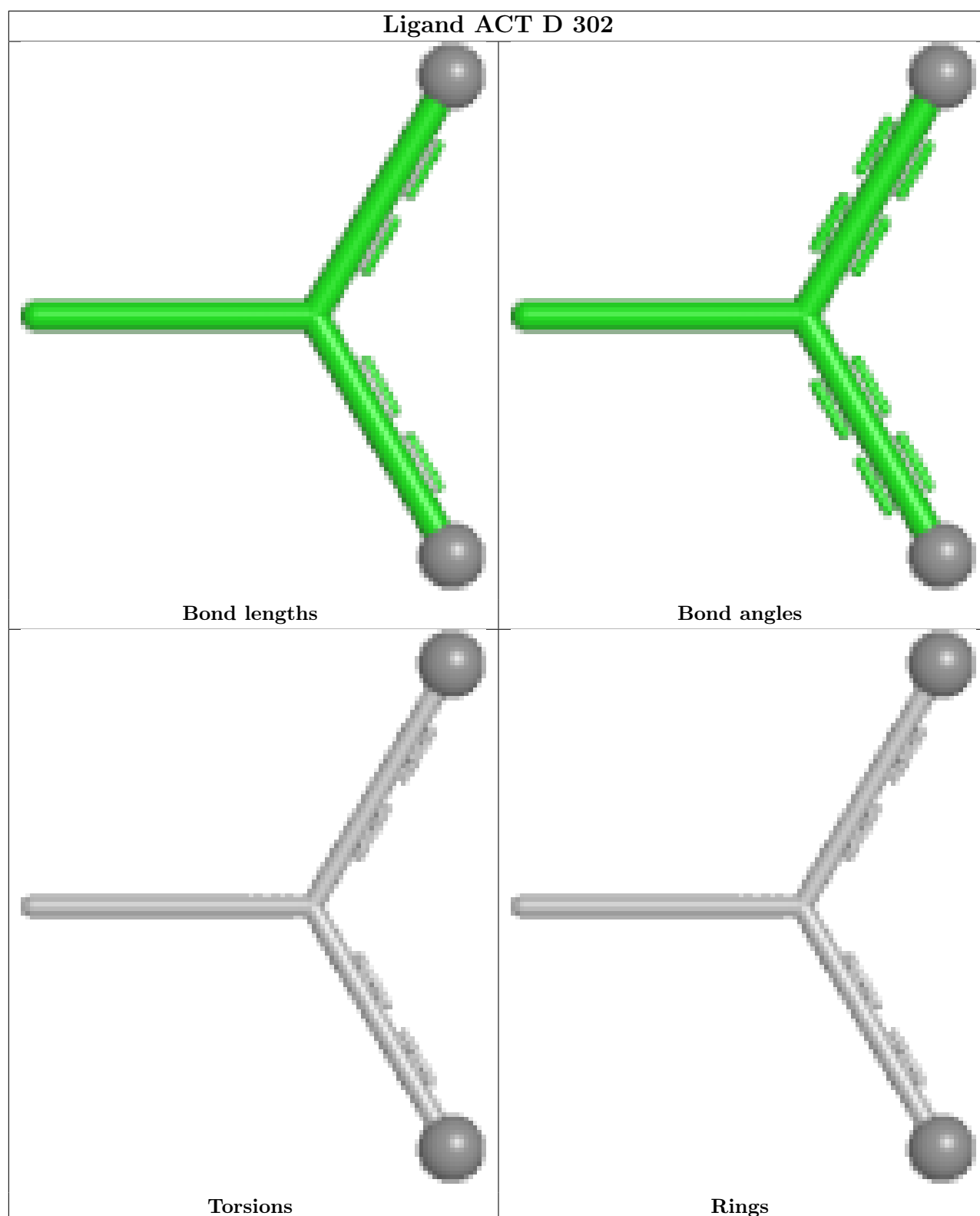


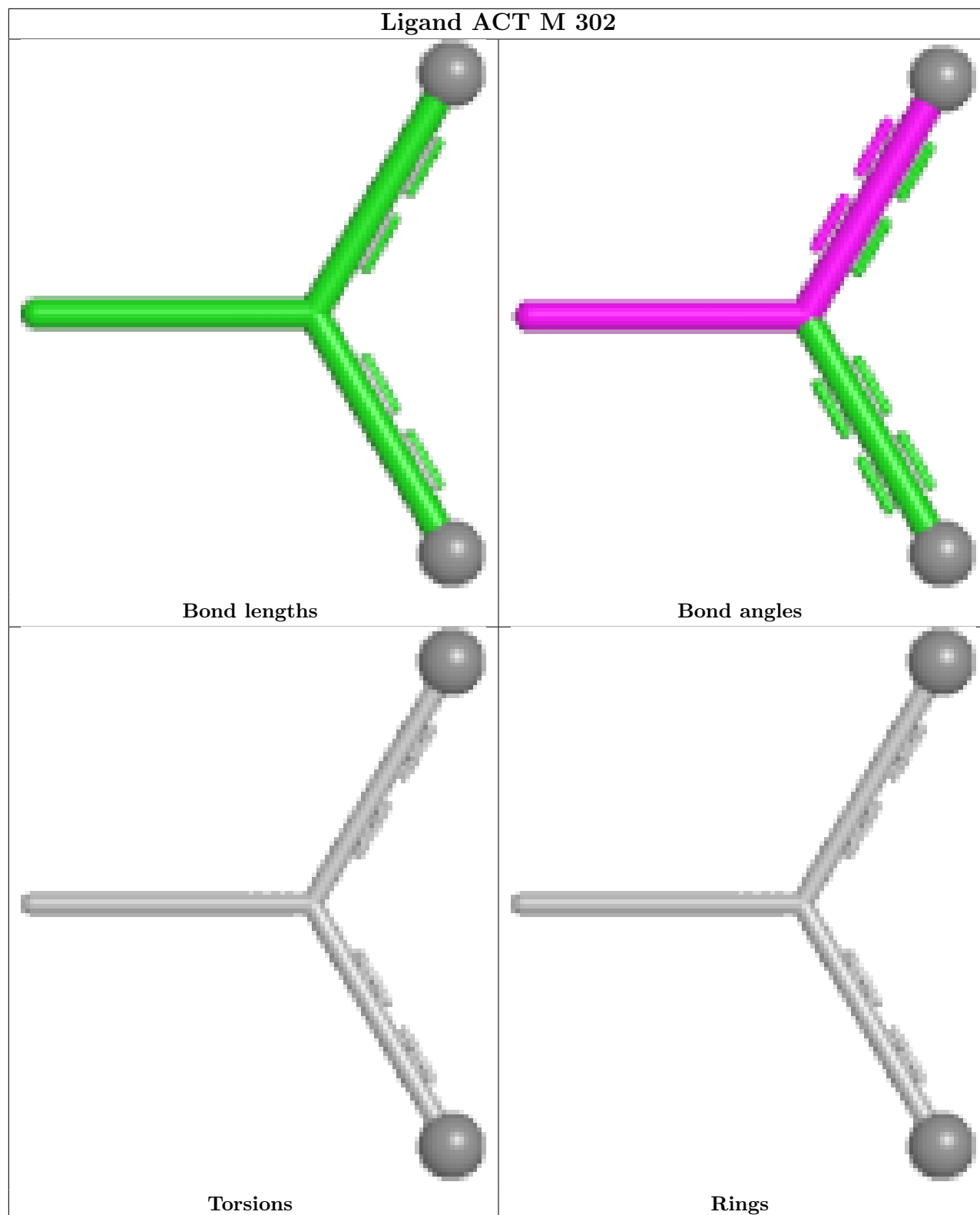












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	222/230 (96%)	0.23	5 (2%) 61 71	12, 18, 28, 40	0
1	B	222/230 (96%)	0.16	3 (1%) 73 80	12, 18, 29, 43	0
1	C	222/230 (96%)	0.30	2 (0%) 81 86	11, 19, 31, 45	0
1	D	222/230 (96%)	0.22	5 (2%) 61 71	11, 18, 31, 36	0
1	E	222/230 (96%)	0.29	1 (0%) 87 90	11, 19, 29, 44	0
1	F	222/230 (96%)	0.39	8 (3%) 46 54	11, 20, 34, 47	0
1	G	222/230 (96%)	0.35	6 (2%) 56 65	13, 20, 30, 49	0
1	H	222/230 (96%)	0.50	9 (4%) 41 50	12, 22, 37, 49	0
1	I	222/230 (96%)	0.57	9 (4%) 41 50	14, 23, 38, 55	0
1	J	222/230 (96%)	0.29	6 (2%) 56 65	13, 18, 30, 42	0
1	K	222/230 (96%)	0.53	5 (2%) 61 71	14, 22, 33, 44	0
1	L	222/230 (96%)	0.41	7 (3%) 50 59	14, 20, 34, 44	0
1	M	222/230 (96%)	0.80	16 (7%) 21 27	14, 26, 40, 58	0
1	N	222/230 (96%)	0.63	7 (3%) 50 59	16, 24, 34, 44	0
1	O	222/230 (96%)	0.71	10 (4%) 38 46	13, 24, 39, 47	0
1	P	222/230 (96%)	0.94	23 (10%) 11 16	15, 28, 43, 57	0
All	All	3552/3680 (96%)	0.46	122 (3%) 48 57	11, 21, 36, 58	0

All (122) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	L	223	PRO	4.7
1	P	11	ILE	4.7
1	G	223	PRO	4.6
1	P	223	PRO	4.3
1	K	223	PRO	4.1

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Mol	Chain	Res	Type	RSRZ
1	F	223	PRO	4.0
1	L	11	ILE	3.9
1	N	45	THR	3.8
1	A	222	SER	3.7
1	P	45	THR	3.5
1	P	43	VAL	3.5
1	M	43	VAL	3.5
1	I	223	PRO	3.4
1	A	223	PRO	3.3
1	D	223	PRO	3.2
1	A	11	ILE	3.1
1	F	11	ILE	3.1
1	P	62	ALA	3.1
1	M	13	PHE	3.0
1	J	11	ILE	3.0
1	I	2	THR	3.0
1	B	223	PRO	3.0
1	K	13	PHE	3.0
1	O	222	SER	3.0
1	J	131	LEU	2.9
1	O	215	GLU	2.9
1	P	2	THR	2.9
1	P	103	GLN	2.9
1	I	11	ILE	2.8
1	H	223	PRO	2.8
1	O	223	PRO	2.8
1	G	13	PHE	2.8
1	M	100	LYS	2.8
1	P	220	ALA	2.8
1	J	70	PHE	2.8
1	P	13	PHE	2.8
1	M	131	LEU	2.8
1	M	195	PHE	2.7
1	P	183	VAL	2.7
1	N	223	PRO	2.7
1	P	10	ALA	2.6
1	M	12	THR	2.6
1	K	222	SER	2.6
1	I	149	ARG	2.6
1	A	131	LEU	2.5
1	I	221	LEU	2.5
1	L	13	PHE	2.5

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Mol	Chain	Res	Type	RSRZ
1	H	49	VAL	2.5
1	M	223	PRO	2.5
1	P	126	PRO	2.5
1	D	222	SER	2.5
1	D	11	ILE	2.5
1	P	221	LEU	2.5
1	P	195	PHE	2.5
1	I	222	SER	2.5
1	F	2	THR	2.5
1	D	131	LEU	2.5
1	J	223	PRO	2.5
1	H	45	THR	2.5
1	G	45	THR	2.4
1	C	222	SER	2.4
1	M	77	GLN	2.4
1	O	131	LEU	2.4
1	A	2	THR	2.4
1	F	222	SER	2.4
1	H	11	ILE	2.4
1	G	222	SER	2.4
1	O	83	GLY	2.4
1	P	12	THR	2.3
1	K	11	ILE	2.3
1	P	63	ALA	2.3
1	J	194	LYS	2.3
1	O	105	GLN	2.3
1	N	62	ALA	2.3
1	K	195	PHE	2.3
1	O	100	LYS	2.3
1	B	131	LEU	2.3
1	I	104	LEU	2.3
1	P	131	LEU	2.3
1	L	10	ALA	2.2
1	N	220	ALA	2.2
1	H	2	THR	2.2
1	L	12	THR	2.2
1	N	70	PHE	2.2
1	P	46	THR	2.2
1	P	47	ARG	2.2
1	C	223	PRO	2.2
1	M	102	PRO	2.2
1	D	2	THR	2.2

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Mol	Chain	Res	Type	RSRZ
1	F	3	THR	2.2
1	M	70	PHE	2.2
1	N	13	PHE	2.2
1	I	43	VAL	2.2
1	L	131	LEU	2.1
1	F	126	PRO	2.1
1	G	11	ILE	2.1
1	H	35	TYR	2.1
1	M	97	THR	2.1
1	O	44	SER	2.1
1	G	77	GLN	2.1
1	E	100	LYS	2.1
1	J	13	PHE	2.1
1	B	222	SER	2.1
1	M	82	GLY	2.1
1	M	104	LEU	2.1
1	I	12	THR	2.1
1	F	173	ASP	2.1
1	P	100	LYS	2.1
1	H	103	GLN	2.1
1	L	16	GLN	2.1
1	O	70	PHE	2.1
1	P	214	VAL	2.1
1	P	15	PRO	2.1
1	M	114	TRP	2.1
1	N	114	TRP	2.1
1	F	45	THR	2.0
1	M	45	THR	2.0
1	H	149	ARG	2.0
1	H	221	LEU	2.0
1	M	105	GLN	2.0
1	P	219	ASP	2.0
1	O	45	THR	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 ⓘ

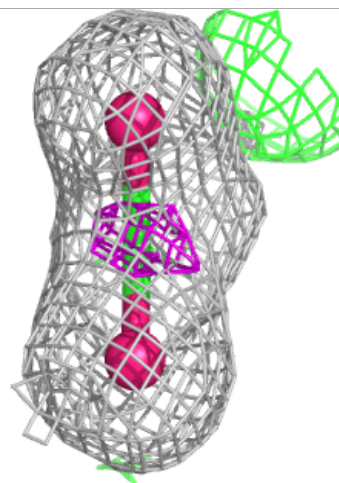
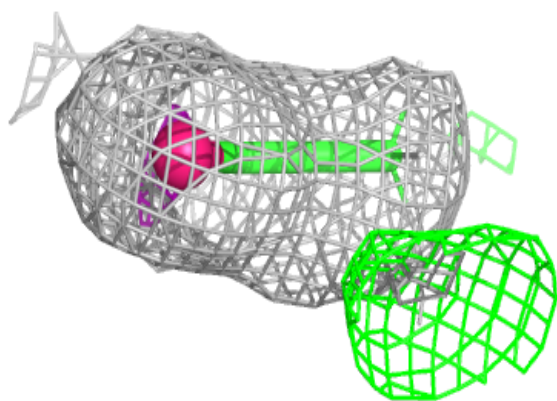
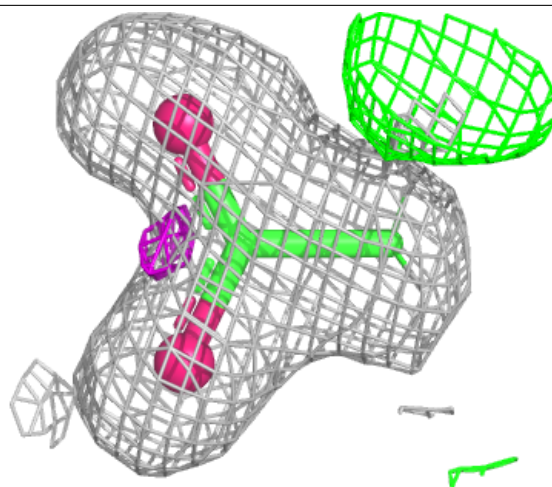
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	ACT	H	301	4/4	0.92	0.09	14,17,17,24	0
2	ACT	P	302	4/4	0.92	0.09	19,21,23,23	0
2	ACT	I	302	4/4	0.94	0.08	16,21,22,22	0
2	ACT	M	302	4/4	0.94	0.08	15,18,24,29	0
2	ACT	C	302	4/4	0.94	0.07	14,17,20,20	0
2	ACT	I	301	4/4	0.95	0.07	16,21,26,26	0
2	ACT	C	301	4/4	0.95	0.08	9,11,12,28	0
2	ACT	A	301	4/4	0.95	0.08	7,9,13,23	0
2	ACT	N	302	4/4	0.95	0.07	16,22,28,28	0
2	ACT	O	301	4/4	0.95	0.07	19,20,28,28	0
2	ACT	P	301	4/4	0.95	0.07	20,23,24,24	0
2	ACT	B	302	4/4	0.95	0.07	8,9,15,22	0
2	ACT	K	301	4/4	0.96	0.07	15,18,24,24	0
2	ACT	L	301	4/4	0.96	0.06	13,15,18,18	0
2	ACT	L	302	4/4	0.96	0.06	13,16,23,23	0
2	ACT	M	301	4/4	0.96	0.07	13,20,24,24	0
2	ACT	B	301	4/4	0.96	0.07	12,15,18,18	0
2	ACT	H	302	4/4	0.96	0.07	18,19,26,26	0
2	ACT	A	302	4/4	0.96	0.07	13,14,18,18	0
2	ACT	O	302	4/4	0.96	0.07	14,17,23,23	0
2	ACT	F	302	4/4	0.96	0.06	11,14,19,19	0
2	ACT	J	301	4/4	0.96	0.06	13,15,19,19	0
2	ACT	E	302	4/4	0.97	0.06	14,16,20,20	0
2	ACT	D	301	4/4	0.97	0.06	14,16,19,19	0
2	ACT	J	302	4/4	0.97	0.06	13,15,19,19	0
2	ACT	G	302	4/4	0.97	0.08	14,17,23,23	0
2	ACT	N	301	4/4	0.97	0.07	16,17,30,30	0
2	ACT	D	302	4/4	0.98	0.05	10,13,18,18	0
2	ACT	F	301	4/4	0.98	0.06	15,16,21,21	0
2	ACT	E	301	4/4	0.98	0.06	13,14,24,24	0
2	ACT	G	301	4/4	0.98	0.05	13,15,25,25	0
2	ACT	K	302	4/4	0.98	0.04	14,17,27,27	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.

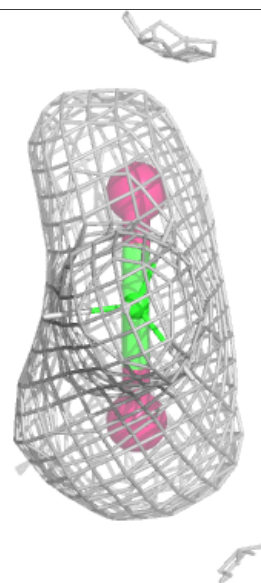
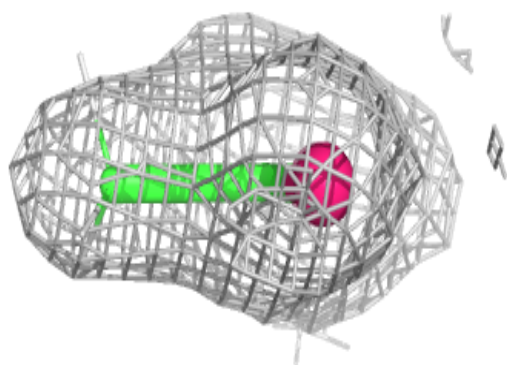
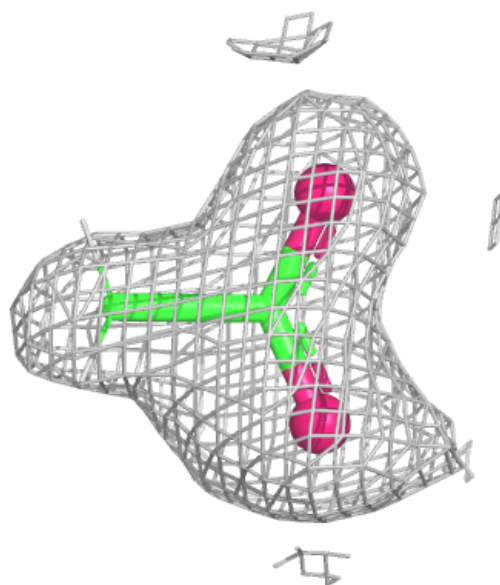
Electron density around ACT H 301:

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



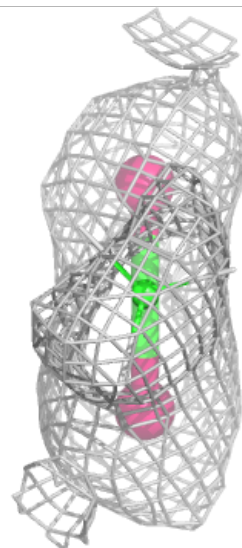
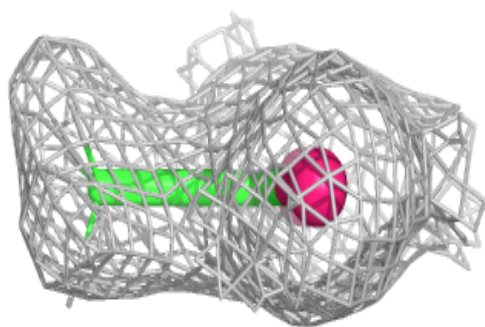
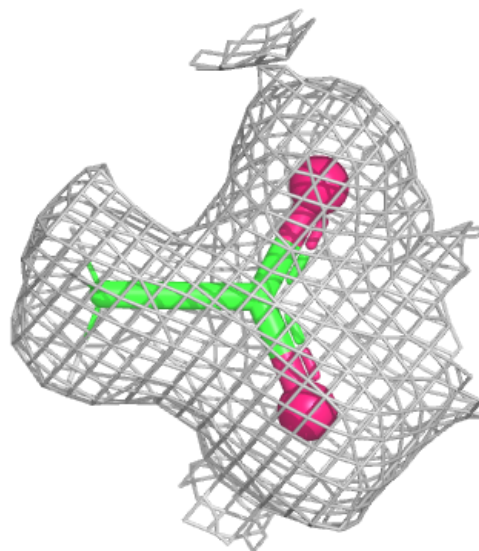
Electron density around ACT P 302:

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and green (positive)



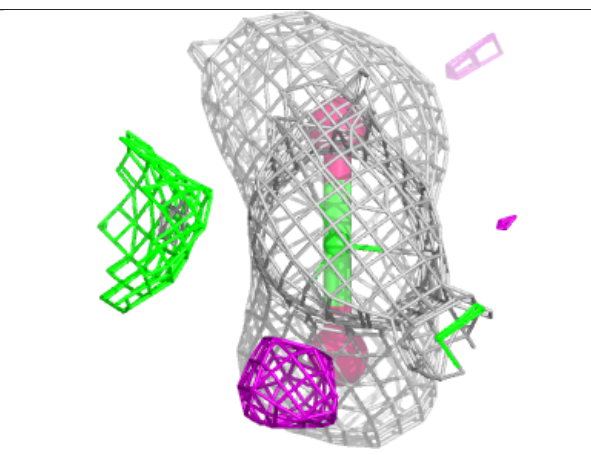
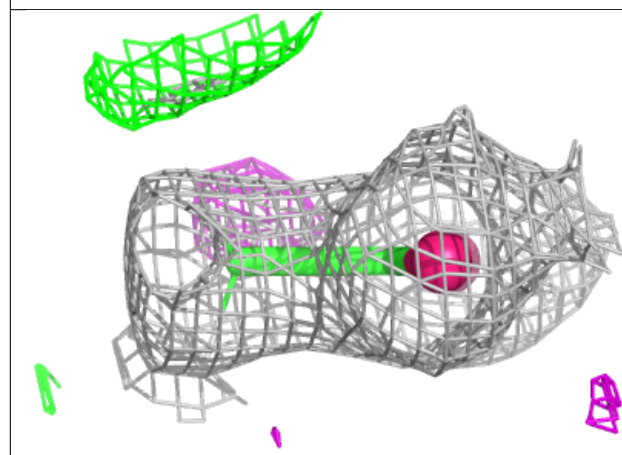
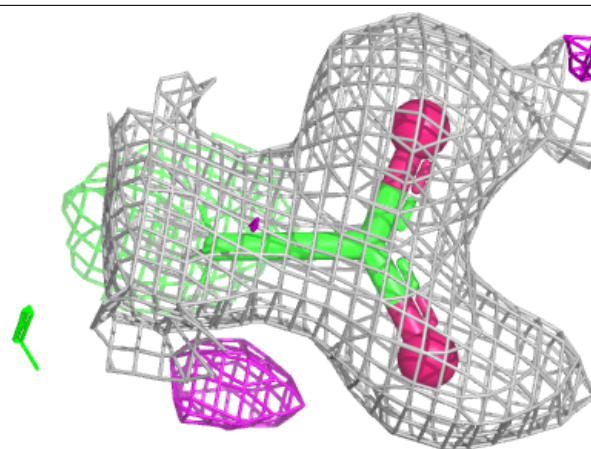
Electron density around ACT I 302:

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and green (positive)



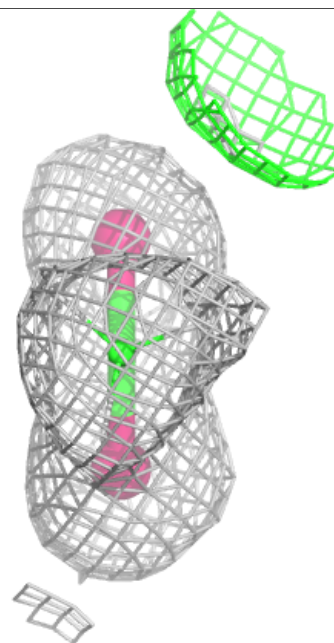
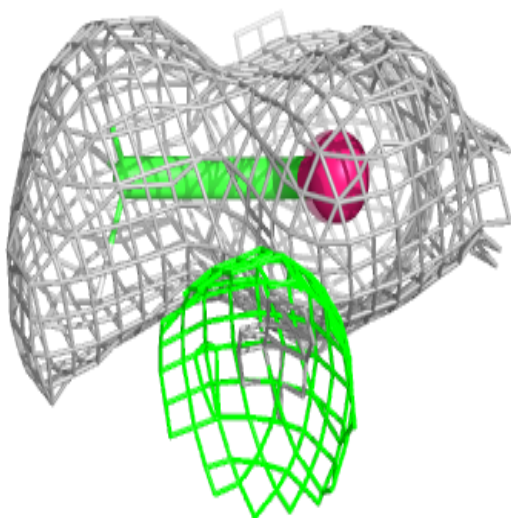
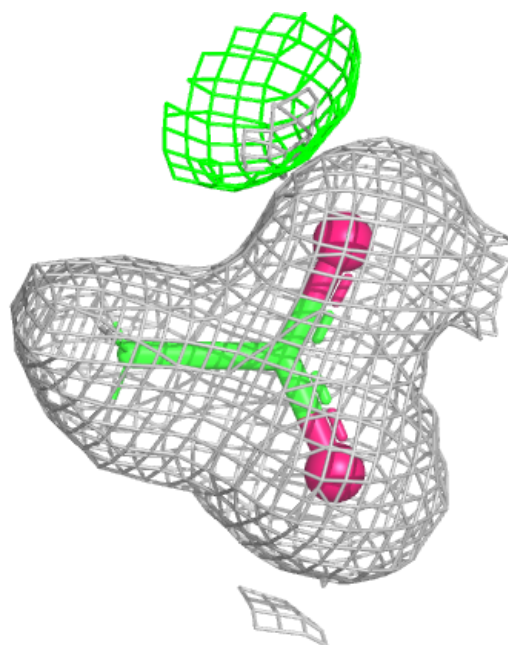
Electron density around ACT M 302:

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and green (positive)



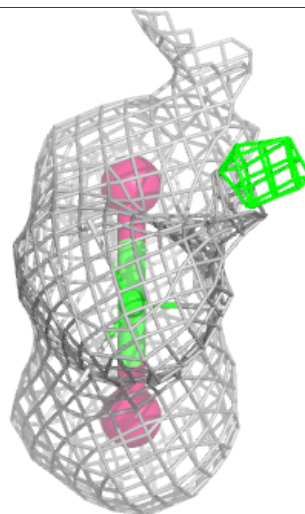
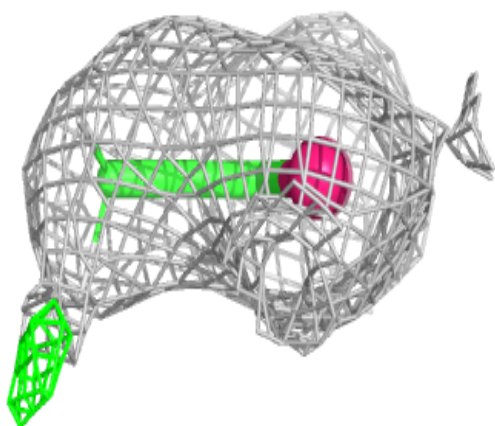
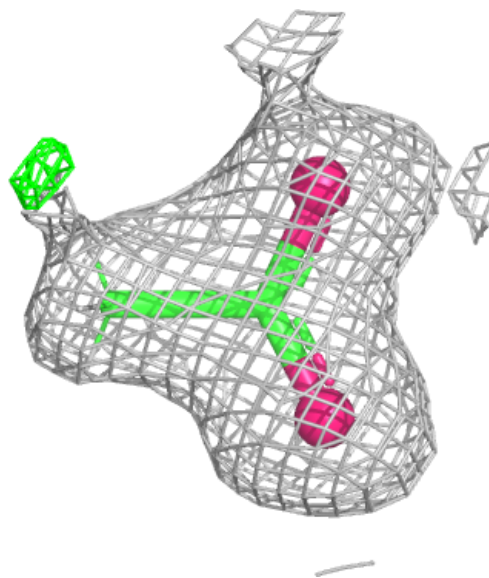
Electron density around ACT C 302:

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and green (positive)



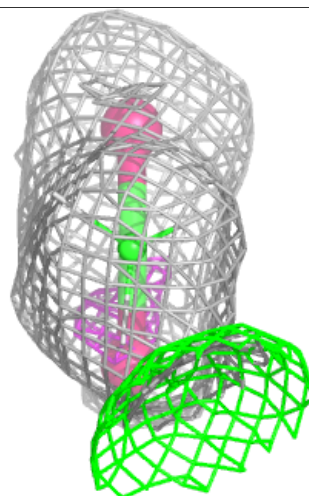
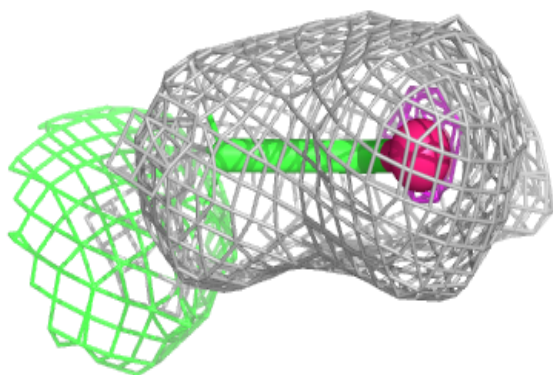
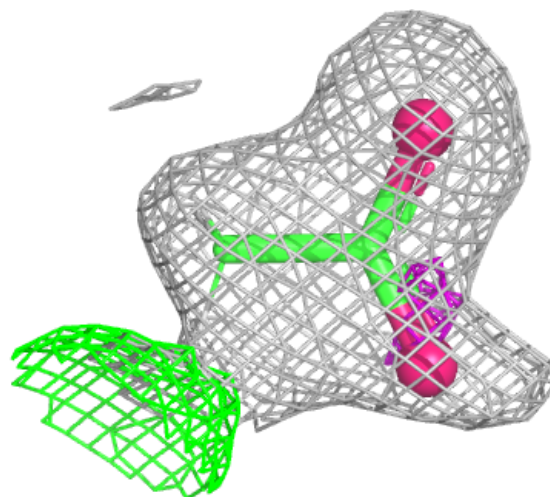
Electron density around ACT I 301:

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and green (positive)



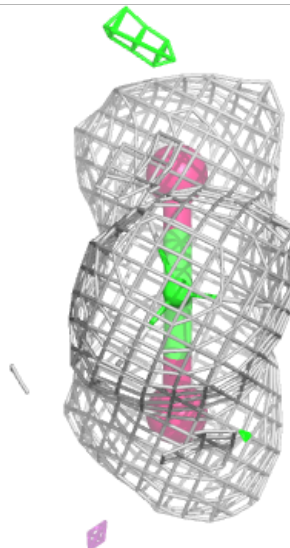
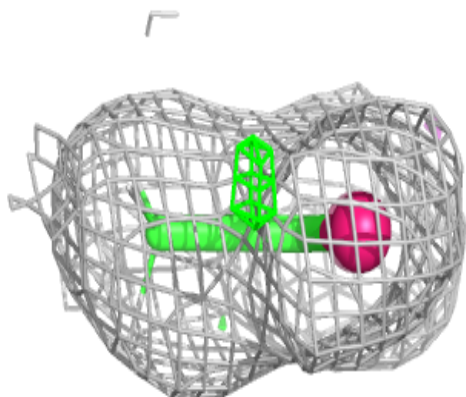
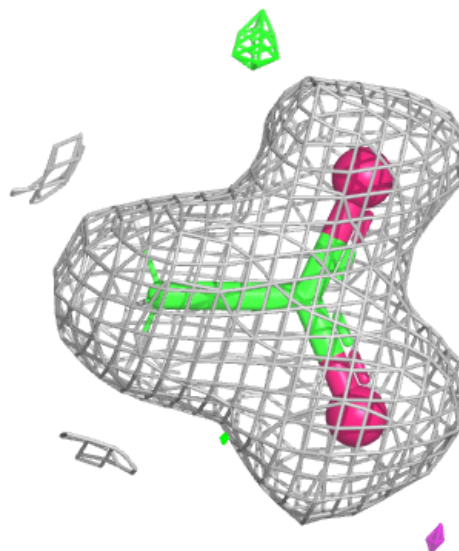
Electron density around ACT C 301:

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



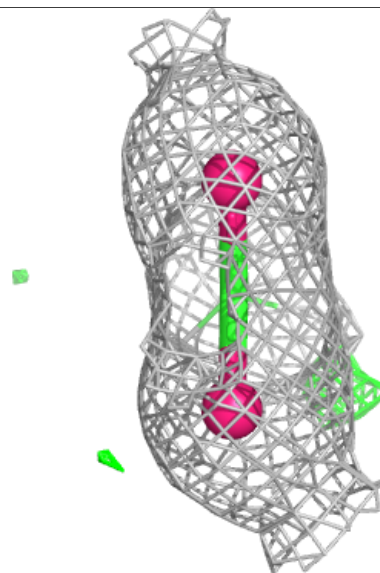
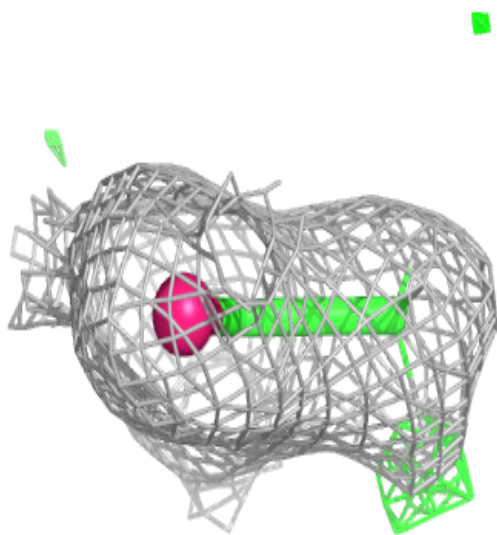
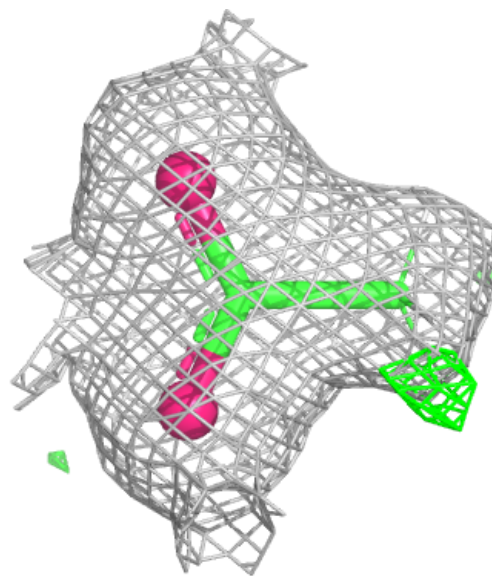
Electron density around ACT A 301:

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and green (positive)



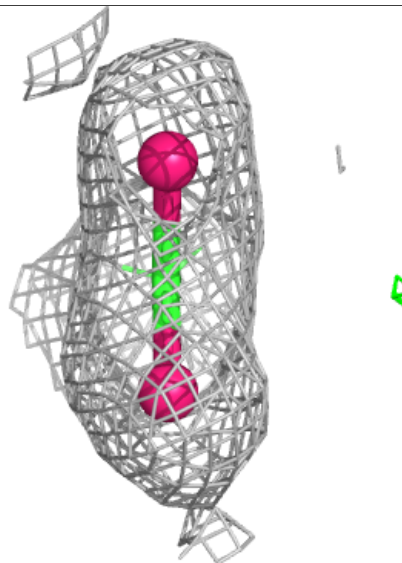
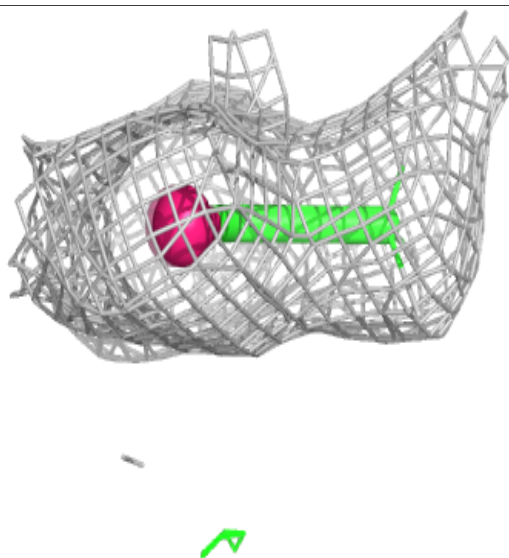
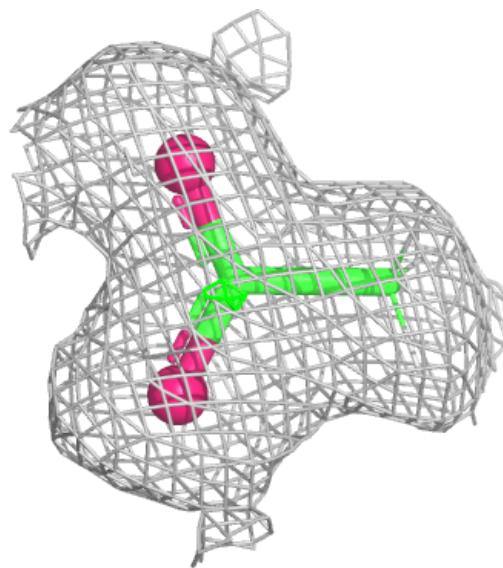
Electron density around ACT N 302:

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and green (positive)



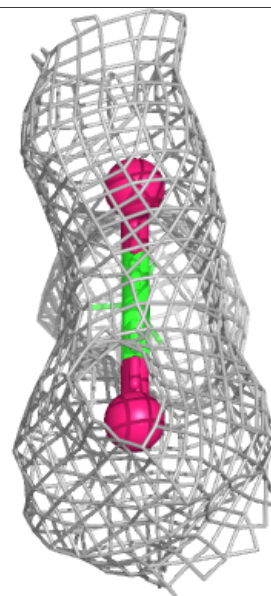
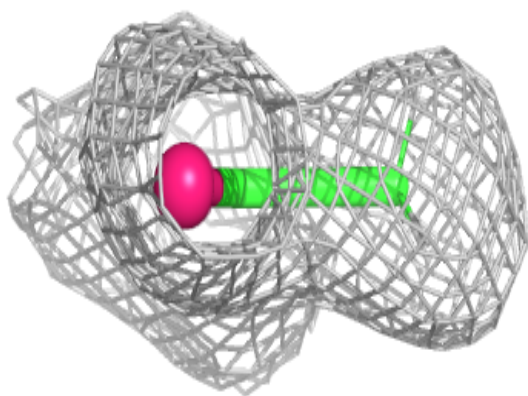
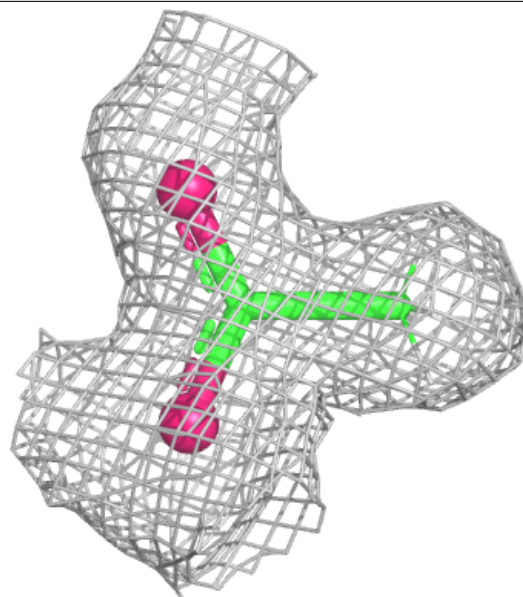
Electron density around ACT O 301:

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and green (positive)



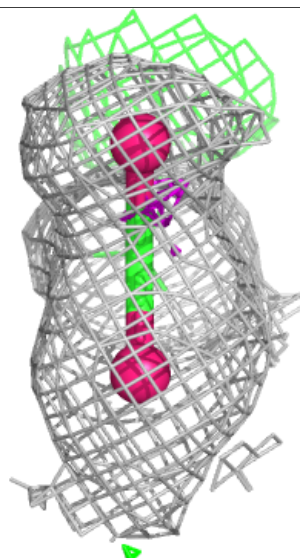
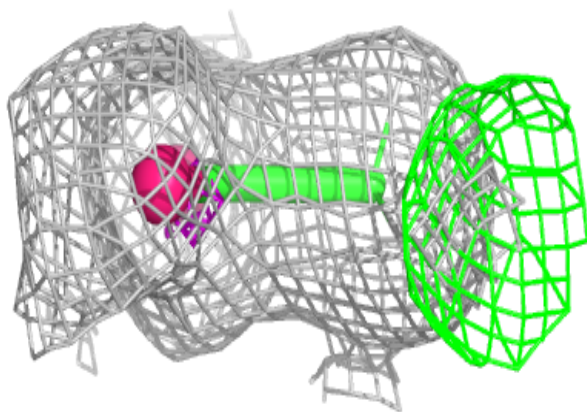
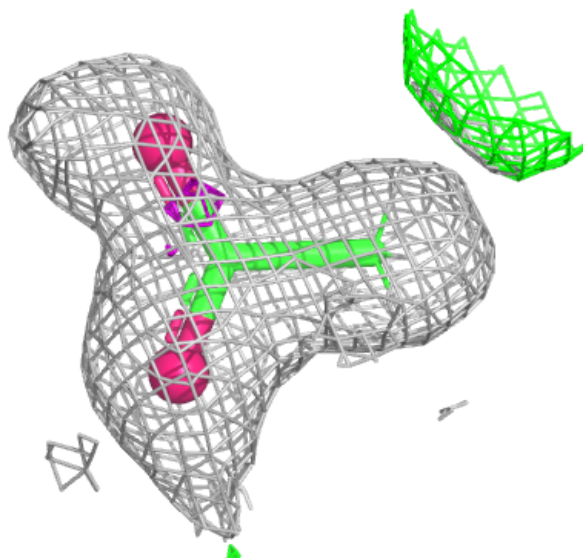
Electron density around ACT P 301:

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and green (positive)



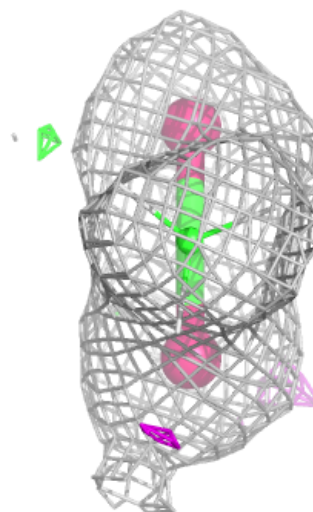
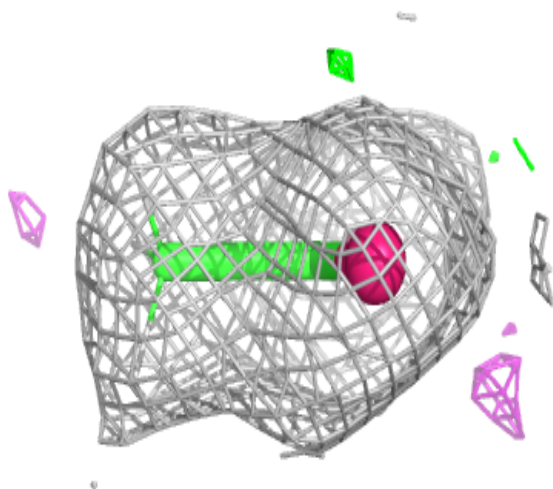
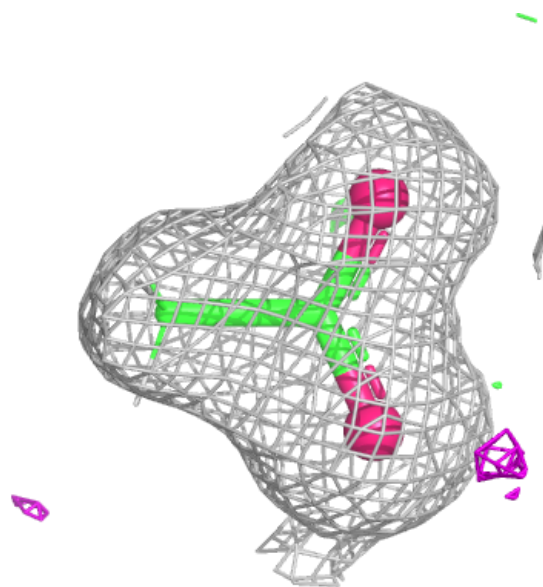
Electron density around ACT B 302:

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and green (positive)



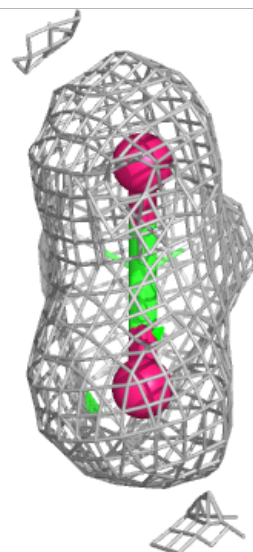
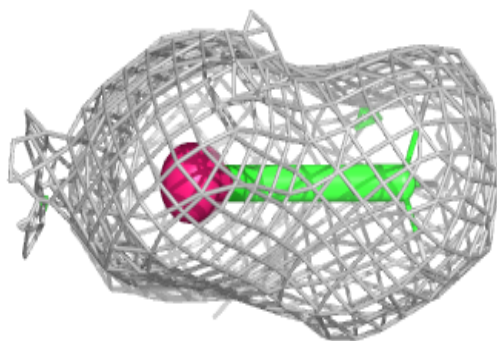
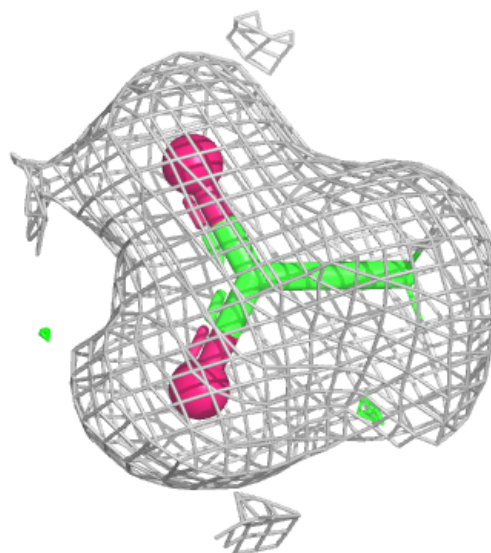
Electron density around ACT K 301:

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and green (positive)



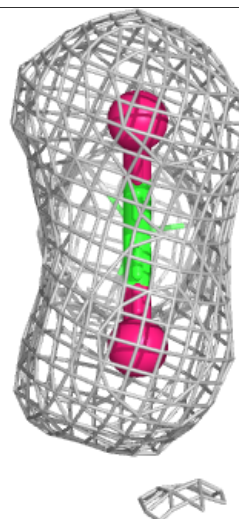
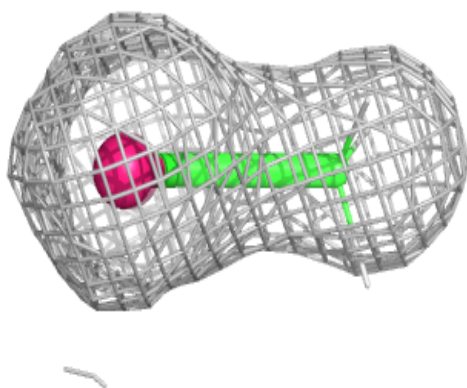
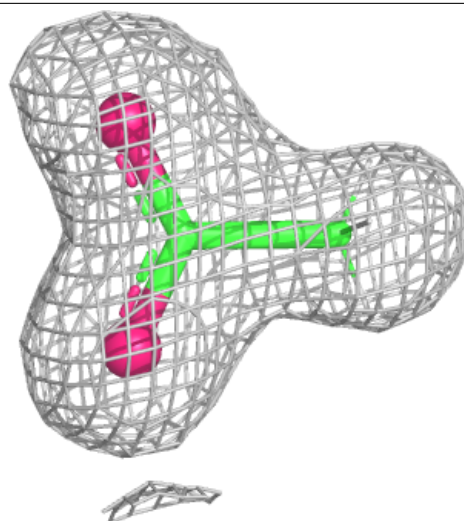
Electron density around ACT L 301:

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and green (positive)



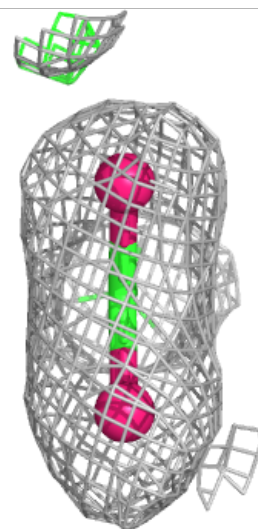
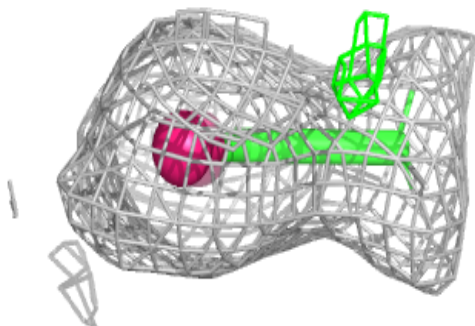
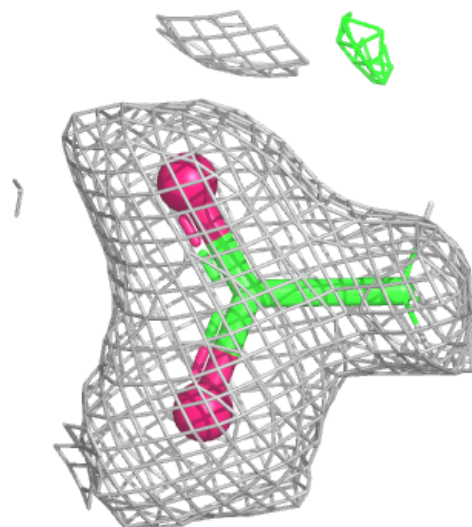
Electron density around ACT L 302:

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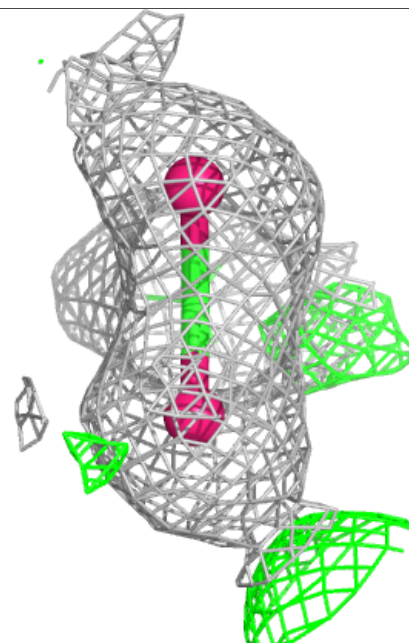
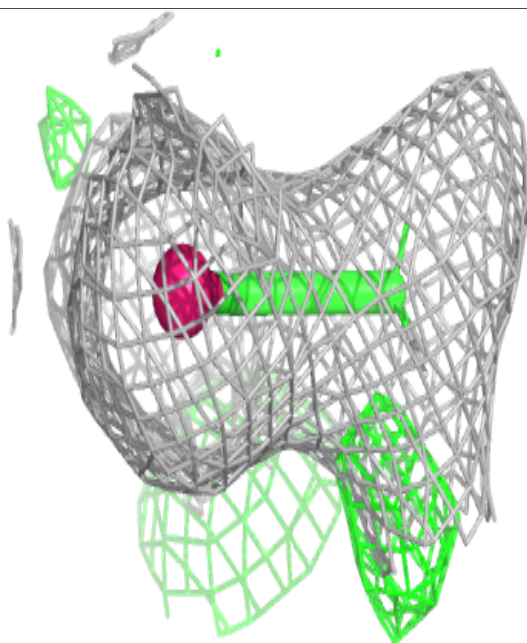
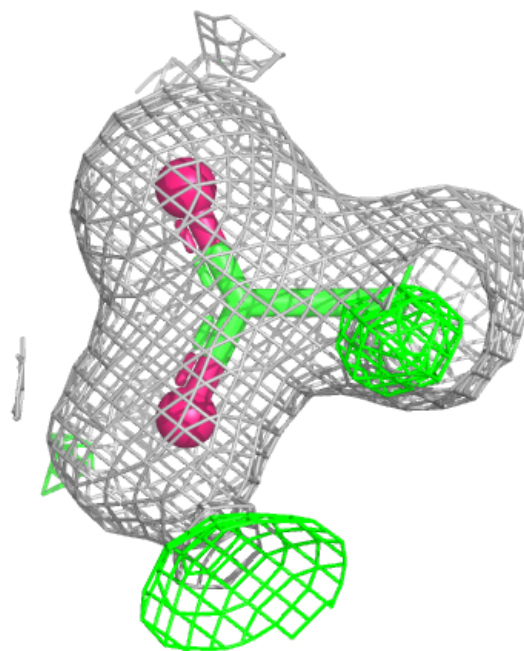
Electron density around ACT M 301:

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and green (positive)



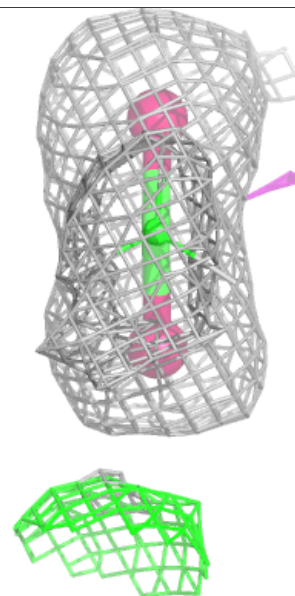
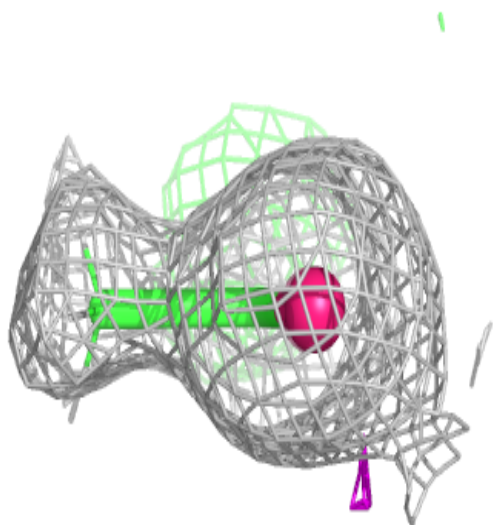
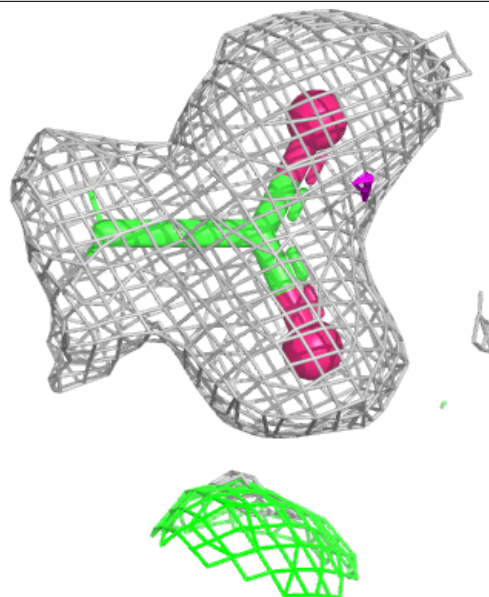
Electron density around ACT B 301:

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and green (positive)



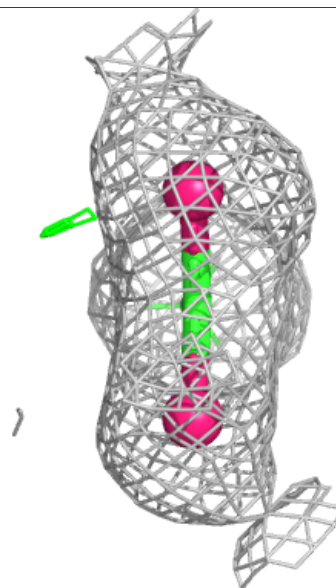
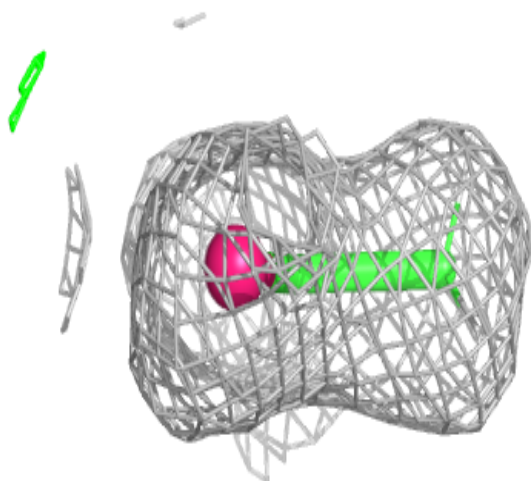
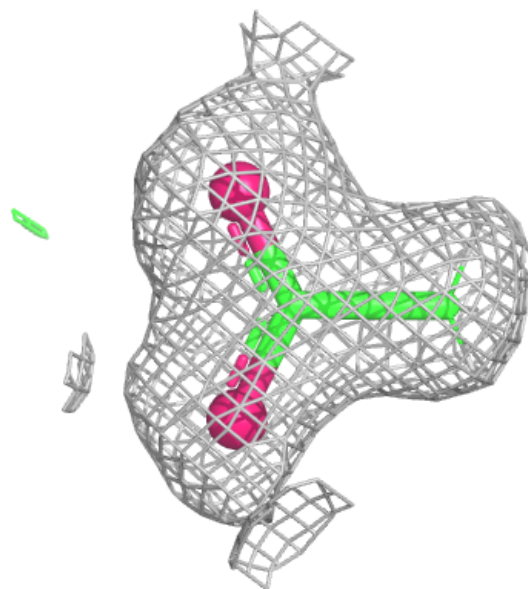
Electron density around ACT H 302:

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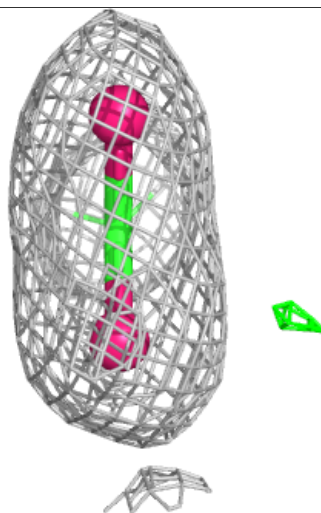
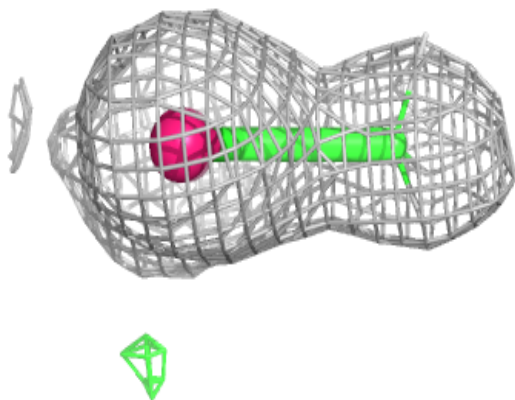
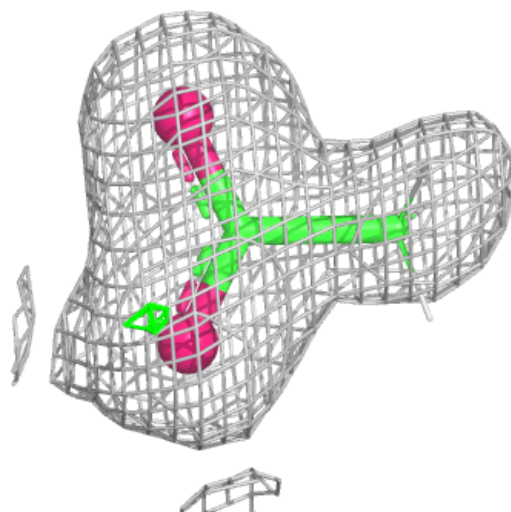
Electron density around ACT A 302:

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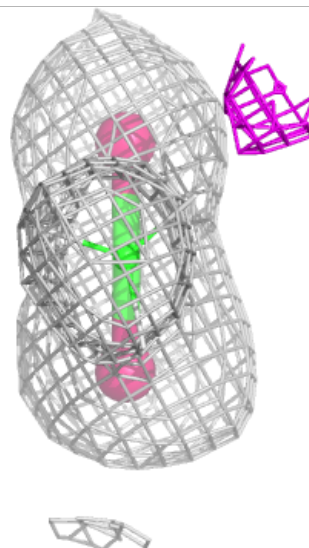
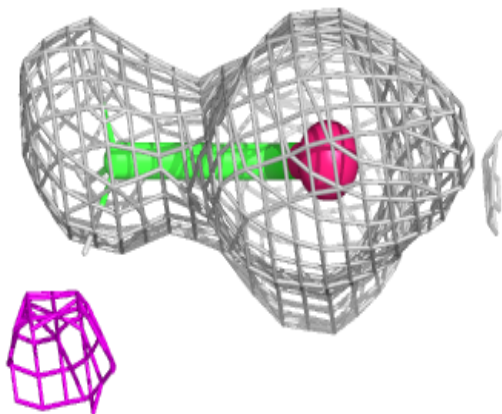
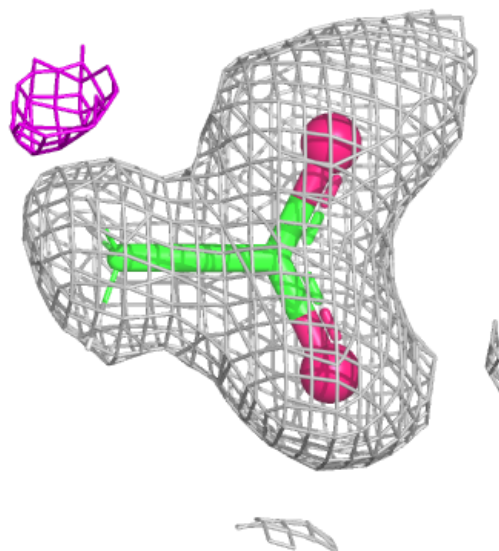
Electron density around ACT O 302:

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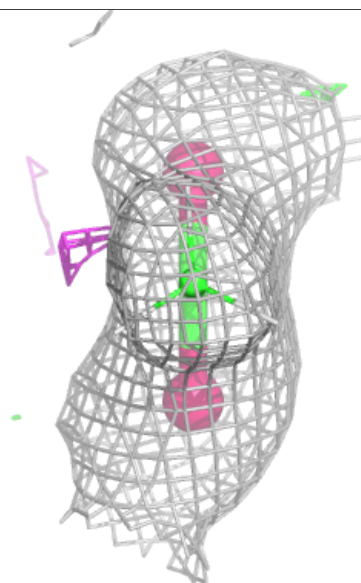
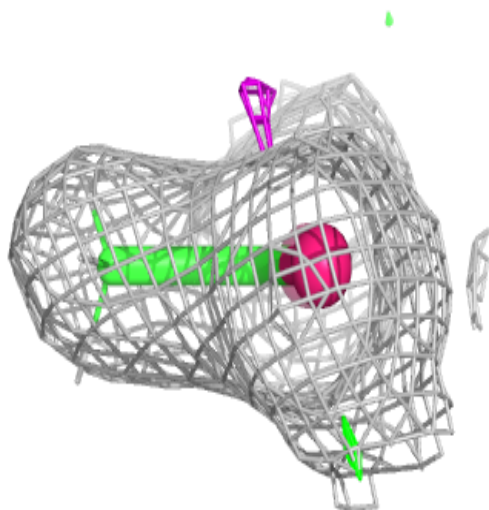
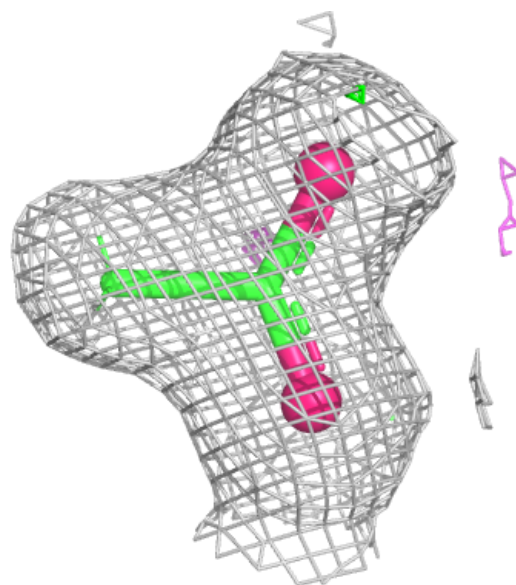
Electron density around ACT F 302:

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and green (positive)



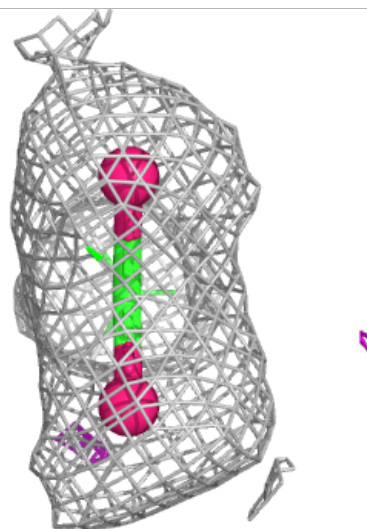
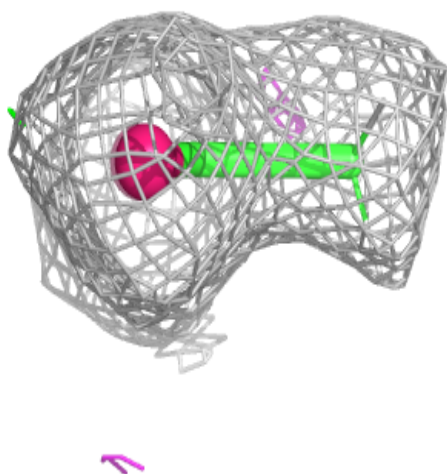
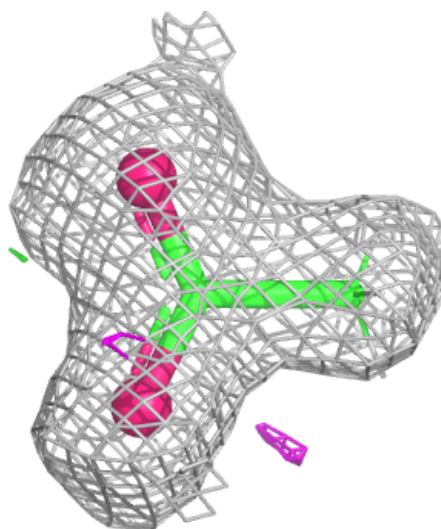
Electron density around ACT J 301:

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and green (positive)



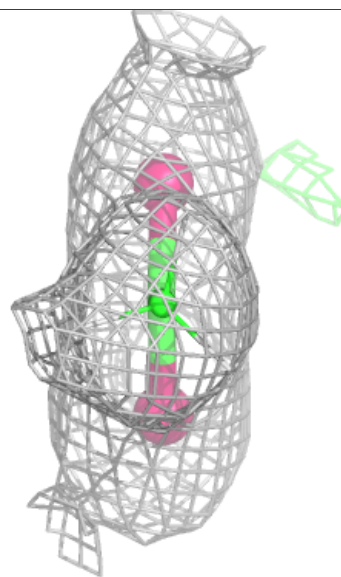
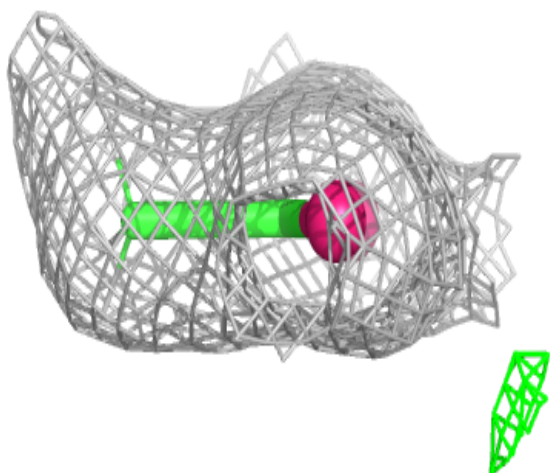
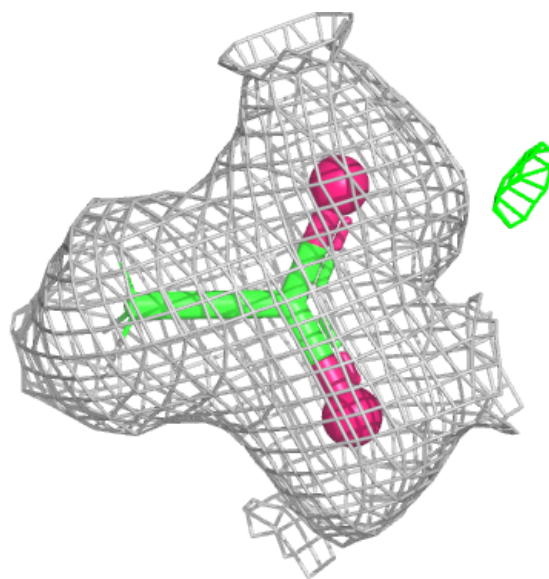
Electron density around ACT E 302:

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and green (positive)



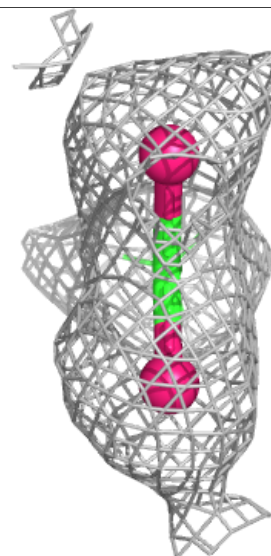
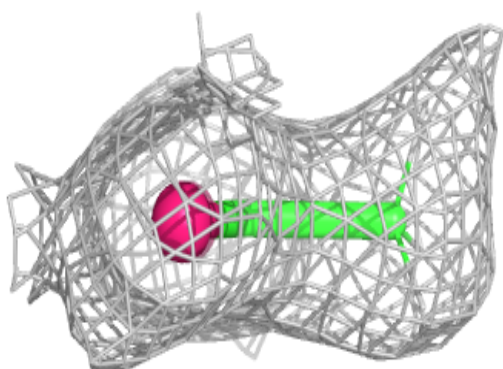
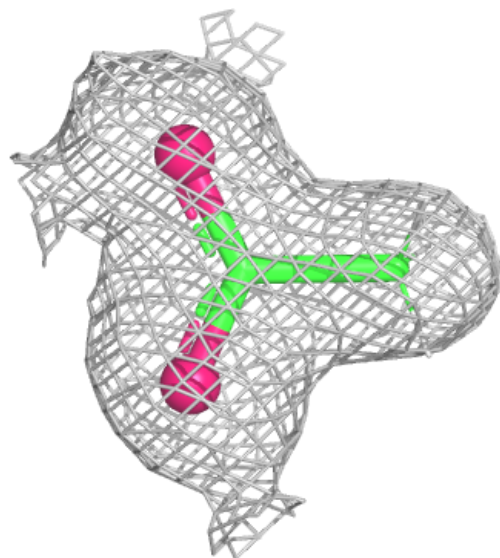
Electron density around ACT D 301:

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



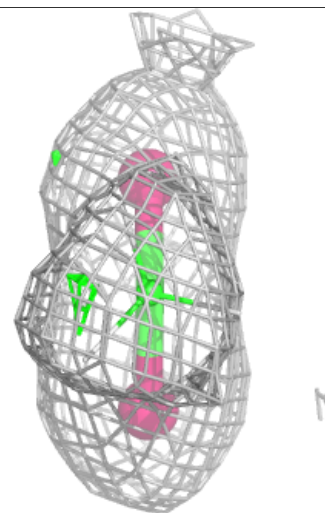
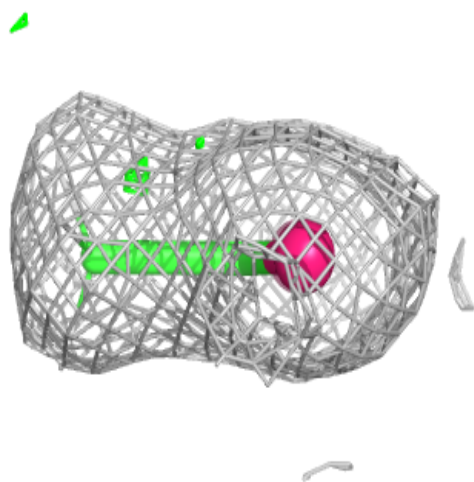
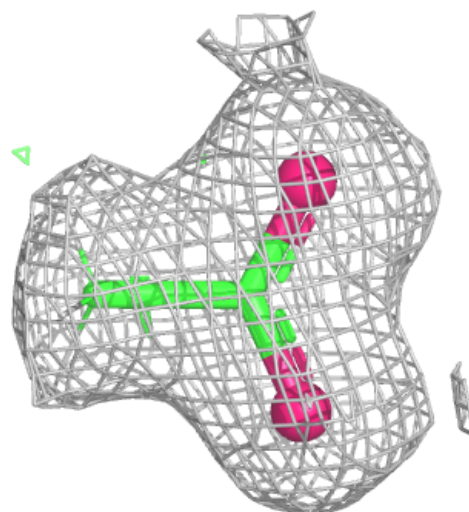
Electron density around ACT J 302:

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



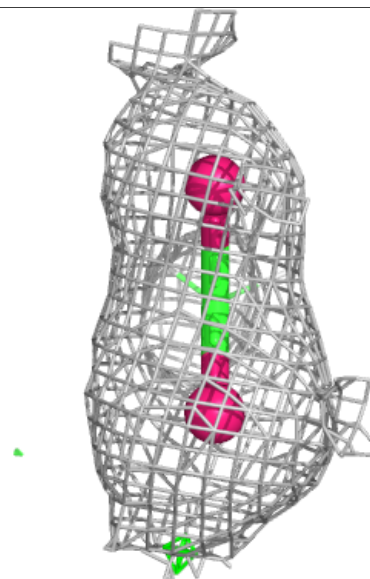
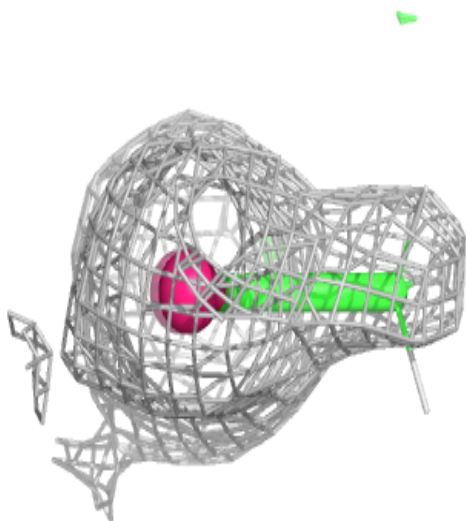
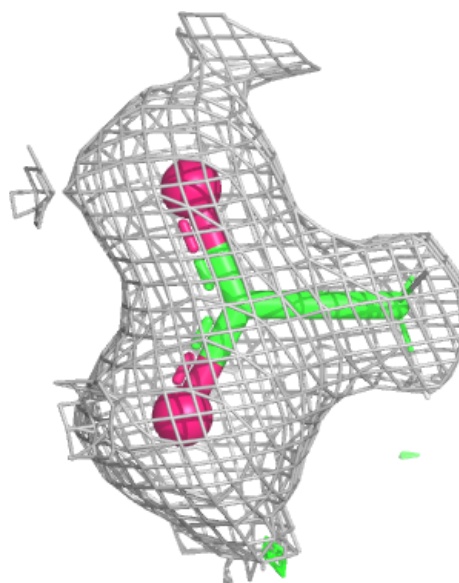
Electron density around ACT G 302:

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



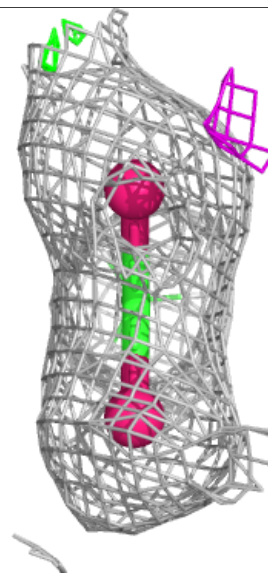
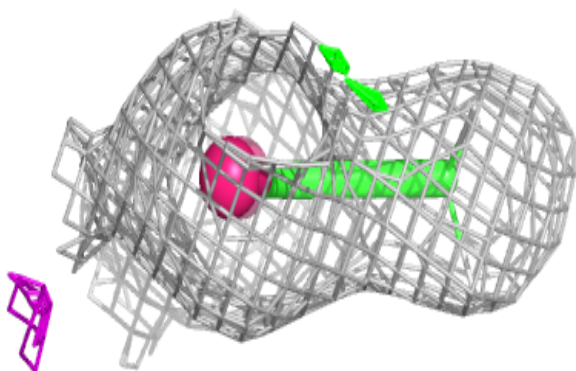
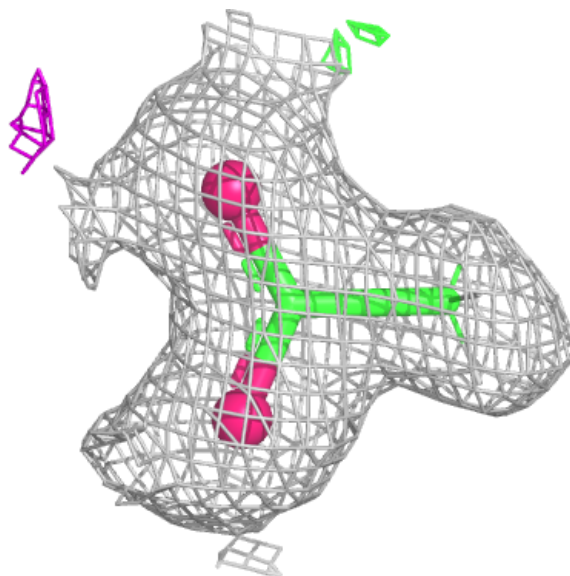
Electron density around ACT N 301:

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



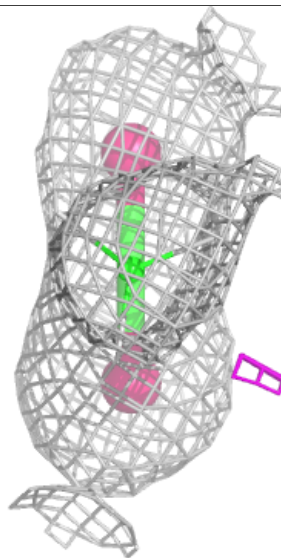
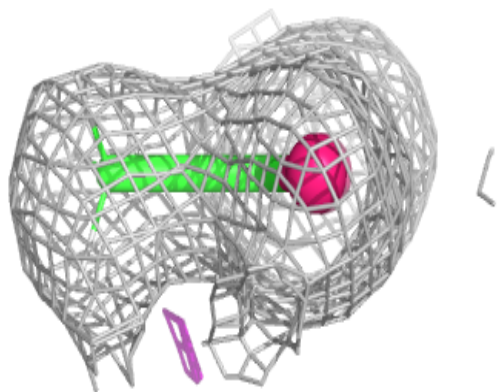
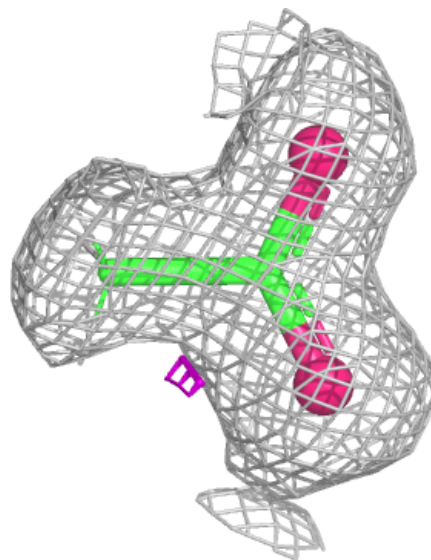
Electron density around ACT D 302:

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



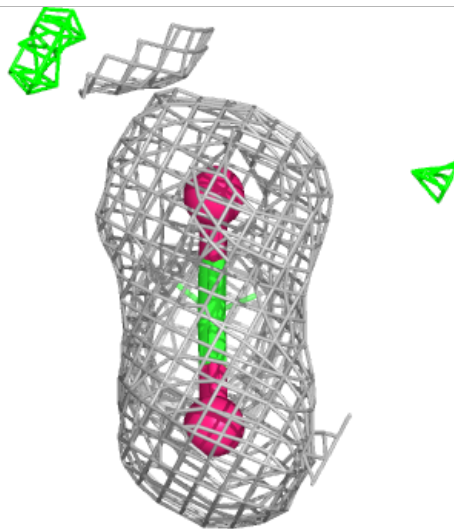
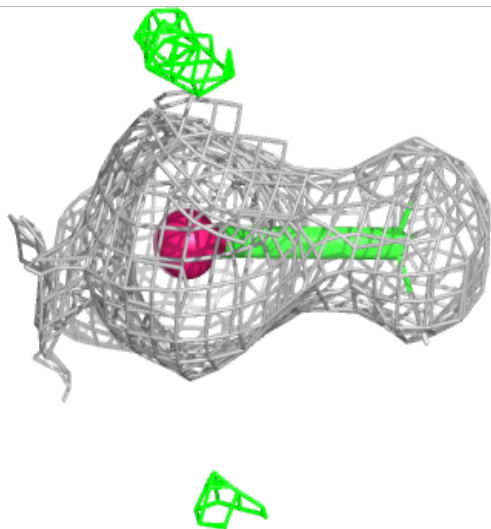
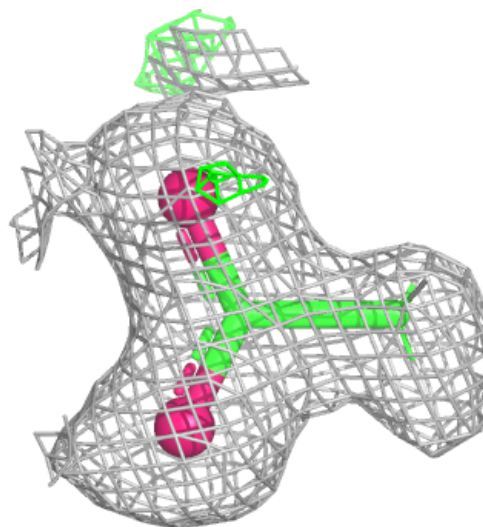
Electron density around ACT F 301:

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



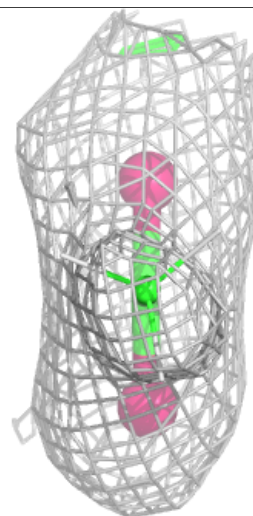
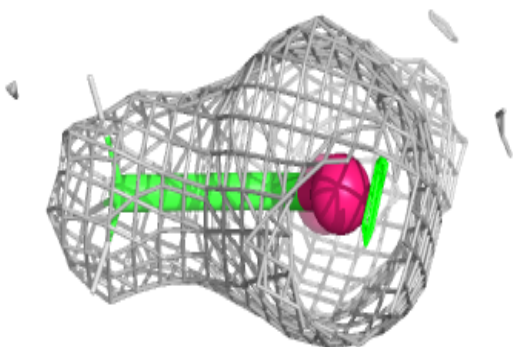
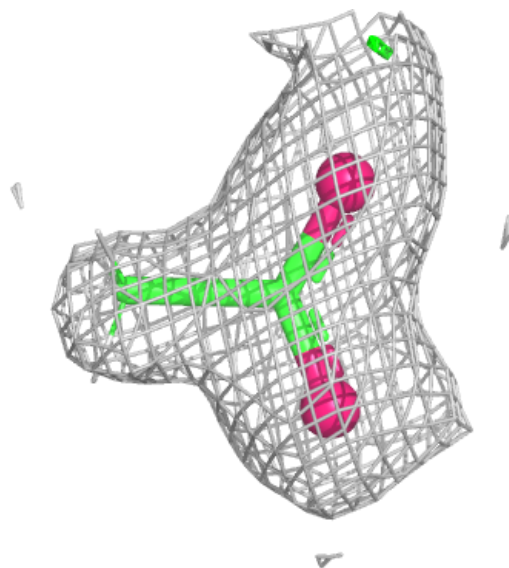
Electron density around ACT E 301:

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



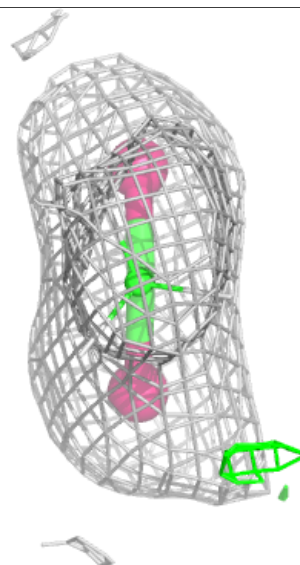
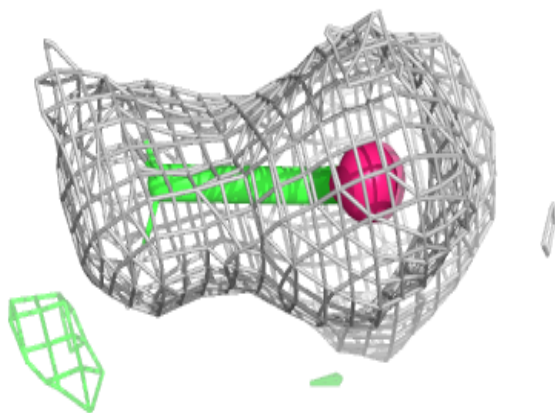
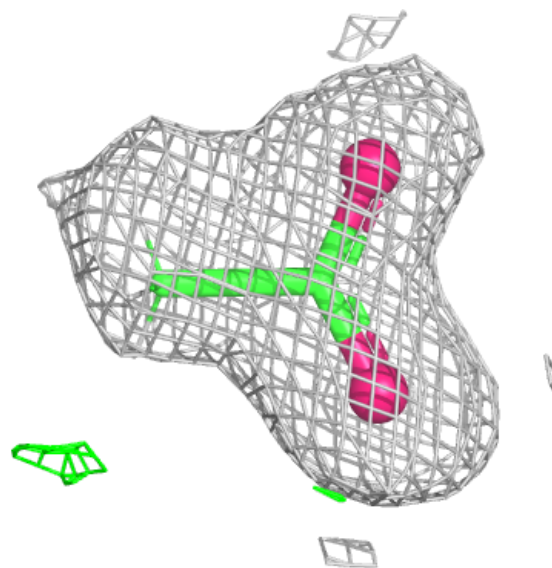
Electron density around ACT G 301:

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



Electron density around ACT K 302:

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



6.5 Other polymers ⓘ

There are no such residues in this entry.