



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 07:35 AM UTC

PDB ID : 5YGK / pdb_00005ygk
Title : Crystal structure of a synthase from Streptomyces sp. CL190 with dmaspp
Authors : Gao, J.; Liu, W.D.; Chen, C.C.; Guo, R.T.
Deposited on : 2017-09-23
Resolution : 2.05 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

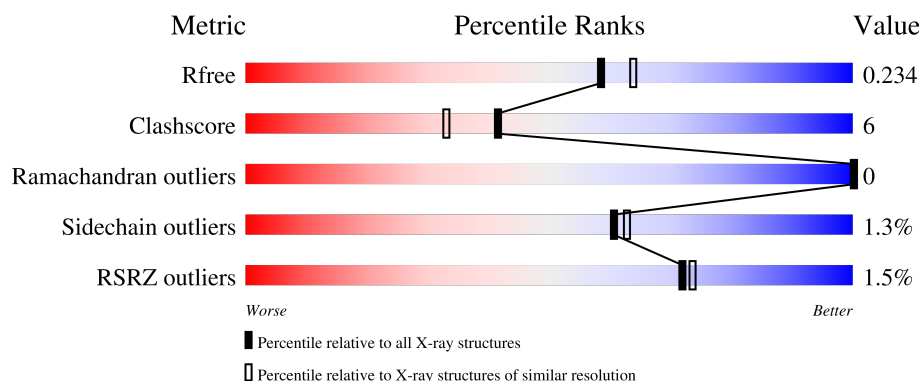
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.05 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	2260 (2.04-2.04)
Clashscore	190562	2333 (2.04-2.04)
Ramachandran outliers	187476	2318 (2.04-2.04)
Sidechain outliers	187428	2318 (2.04-2.04)
RSRZ outliers	180081	2260 (2.04-2.04)

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	217	
1	B	217	
1	C	217	
1	D	217	
1	E	217	

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Mol	Chain	Length	Quality of chain
1	F	217	
1	G	217	
1	H	217	

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	DST	C	301	-	-	X	-

2 Entry composition

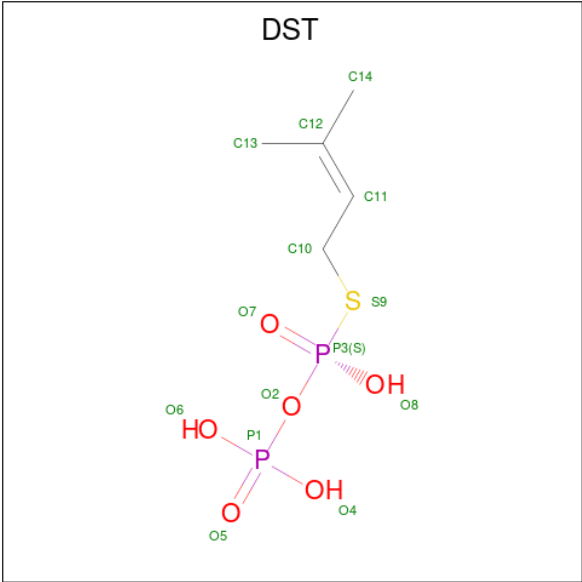
There are 4 unique types of molecules in this entry. The entry contains 14345 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Cyclolavandulyl diphosphate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	210	Total	C	N	O	S	0	0	0
			1657	1051	275	320	11			
1	B	208	Total	C	N	O	S	0	0	0
			1638	1041	270	316	11			
1	C	217	Total	C	N	O	S	0	0	0
			1712	1087	285	329	11			
1	D	208	Total	C	N	O	S	0	0	0
			1638	1041	270	316	11			
1	E	208	Total	C	N	O	S	0	0	0
			1638	1041	270	316	11			
1	F	208	Total	C	N	O	S	0	0	0
			1638	1040	270	318	10			
1	G	208	Total	C	N	O	S	0	0	0
			1638	1040	270	318	10			
1	H	216	Total	C	N	O	S	0	0	0
			1703	1082	284	327	10			

- Molecule 2 is DIMETHYLALLYL S-THIOLODIPHOSPHATE (CCD ID: DST) (formula: C₅H₁₂O₆P₂S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	O	P	S	0	0
			14	5	6	2	1		
2	A	1	Total	C	O	P	S	0	0
			14	5	6	2	1		
2	B	1	Total	C	O	P	S	0	0
			14	5	6	2	1		
2	B	1	Total	C	O	P	S	0	0
			14	5	6	2	1		
2	C	1	Total	C	O	P	S	0	0
			14	5	6	2	1		
2	C	1	Total	C	O	P	S	0	0
			14	5	6	2	1		
2	D	1	Total	C	O	P	S	0	0
			14	5	6	2	1		
2	D	1	Total	C	O	P	S	0	0
			14	5	6	2	1		
2	E	1	Total	C	O	P	S	0	0
			14	5	6	2	1		
2	E	1	Total	C	O	P	S	0	0
			14	5	6	2	1		
2	F	1	Total	C	O	P	S	0	0
			14	5	6	2	1		
2	F	1	Total	C	O	P	S	0	0
			14	5	6	2	1		
2	G	1	Total	C	O	P	S	0	0
			14	5	6	2	1		
2	G	1	Total	C	O	P	S	0	0
			14	5	6	2	1		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	H	1	Total	C	O	P	S	0	0
			14	5	6	2	1		
2	H	1	Total	C	O	P	S	0	0
			14	5	6	2	1		

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		
3	B	1	Total	Mg	0	0
			1	1		
3	C	1	Total	Mg	0	0
			1	1		
3	D	1	Total	Mg	0	0
			1	1		
3	E	1	Total	Mg	0	0
			1	1		
3	F	1	Total	Mg	0	0
			1	1		
3	G	1	Total	Mg	0	0
			1	1		
3	H	1	Total	Mg	0	0
			1	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	177	Total	O	0	0
			177	177		
4	B	77	Total	O	0	0
			77	77		
4	C	150	Total	O	0	0
			150	150		
4	D	89	Total	O	0	0
			89	89		
4	E	107	Total	O	0	0
			107	107		
4	F	37	Total	O	0	0
			37	37		
4	G	89	Total	O	0	0
			89	89		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	H	125	Total 125	O 125	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.

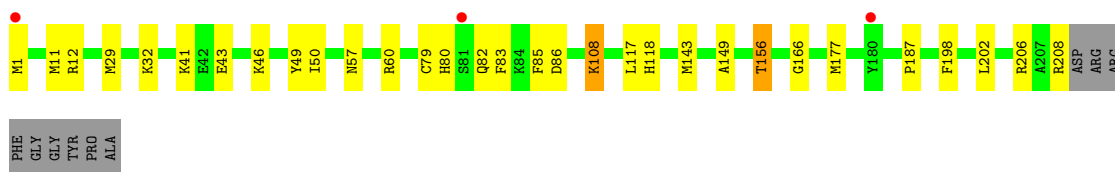
- Molecule 1: Cyclolavandulyl diphosphate synthase

Chain A: 

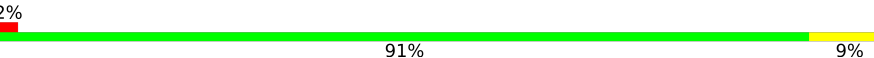


- Molecule 1: Cyclolavandulyl diphosphate synthase

Chain B: 



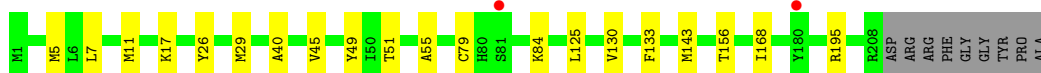
- Molecule 1: Cyclolavandulyl diphosphate synthase

Chain C: 



- Molecule 1: Cyclolavandulyl diphosphate synthase

Chain D: 

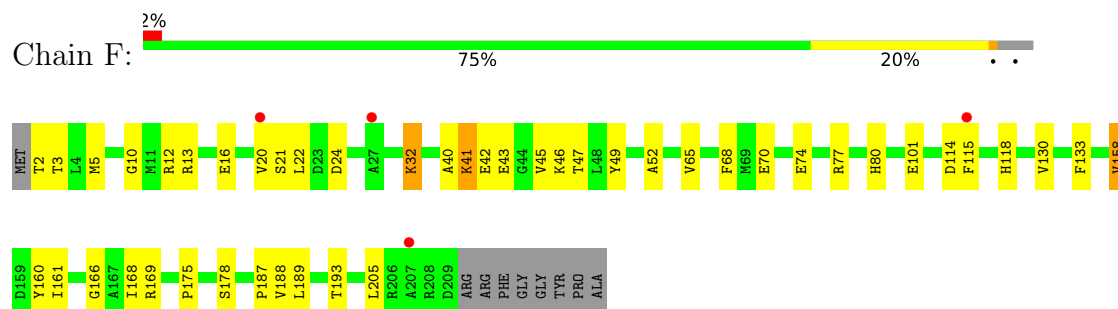


- Molecule 1: Cyclolavandulyl diphosphate synthase

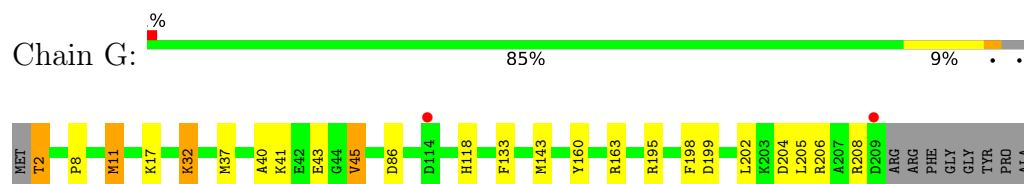
Chain E: 



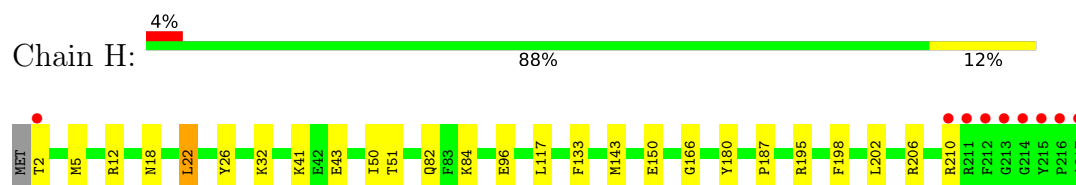
● Molecule 1: Cyclolavandulyl diphosphate synthase



● Molecule 1: Cyclolavandulyl diphosphate synthase



● Molecule 1: Cyclolavandulyl diphosphate synthase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	90.09Å 119.31Å 93.13Å 90.00° 117.35° 90.00°	Depositor
Resolution (Å)	24.87 – 2.05 24.87 – 2.05	Depositor EDS
% Data completeness (in resolution range)	99.4 (24.87-2.05) 96.1 (24.87-2.05)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.85 (at 2.04Å)	Xtriage
Refinement program	PHENIX (1.10_2155: ???)	Depositor
R, R_{free}	0.184 , 0.236 0.185 , 0.234	Depositor DCC
R_{free} test set	2004 reflections (1.83%)	wwPDB-VP
Wilson B-factor (Å ²)	32.6	Xtriage
Anisotropy	0.368	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 47.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.000 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	14345	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 53.04 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 4.4636e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	0.41	0/1693	0.65	0/2287
1	B	0.56	0/1674	0.67	1/2262 (0.0%)
1	C	0.44	0/1751	0.71	1/2364 (0.0%)
1	D	0.70	3/1674 (0.2%)	0.68	1/2262 (0.0%)
1	E	0.57	1/1674 (0.1%)	0.70	0/2262
1	F	0.50	1/1674 (0.1%)	0.66	3/2263 (0.1%)
1	G	0.47	1/1674 (0.1%)	0.69	0/2263
1	H	0.43	1/1742 (0.1%)	0.68	0/2354
All	All	0.52	7/13556 (0.1%)	0.68	6/18317 (0.0%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	55	ALA	C-O	-6.05	1.17	1.24
1	E	125	LEU	C-O	-5.47	1.17	1.24
1	D	125	LEU	C-O	-5.37	1.18	1.24
1	G	11	MET	C-O	-5.25	1.17	1.24
1	F	158	VAL	C-O	-5.14	1.18	1.24
1	H	150	GLU	C-O	-5.08	1.17	1.24
1	D	17	LYS	C-O	-5.07	1.17	1.24

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	79	CYS	N-CA-C	7.69	121.48	112.72
1	C	212	PHE	N-CA-C	7.67	119.33	110.97
1	B	79	CYS	N-CA-C	7.18	121.74	112.41
1	F	80	HIS	N-CA-C	5.83	119.49	112.38
1	F	32	LYS	CA-CB-CG	-5.65	102.81	114.10
1	F	41	LYS	N-CA-C	-5.19	105.54	111.14

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	1657	0	1614	15	0
1	B	1638	0	1597	24	0
1	C	1712	0	1663	16	0
1	D	1638	0	1597	15	0
1	E	1638	0	1597	17	0
1	F	1638	0	1589	33	0
1	G	1638	0	1589	21	0
1	H	1703	0	1651	19	0
2	A	28	0	20	2	0
2	B	28	0	20	5	0
2	C	28	0	20	10	0
2	D	28	0	20	7	0
2	E	28	0	20	5	0
2	F	28	0	20	4	0
2	G	28	0	20	6	0
2	H	28	0	20	3	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0
3	F	1	0	0	0	0
3	G	1	0	0	0	0
3	H	1	0	0	0	0
4	A	177	0	0	0	0
4	B	77	0	0	0	0
4	C	150	0	0	2	0
4	D	89	0	0	1	0
4	E	107	0	0	2	0
4	F	37	0	0	0	0
4	G	89	0	0	1	0
4	H	125	0	0	2	0
All	All	14345	0	13057	170	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 6.

All (170) 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:E:206:ARG:NH1	4:E:401:HOH:O	1.78	1.15
2:D:301:DST:C11	2:D:302:DST:H133	1.90	1.01
2:D:301:DST:H111	2:D:302:DST:H133	1.03	0.99
1:A:43:GLU:HG3	1:A:206:ARG:HH12	1.30	0.97
2:D:301:DST:H111	2:D:302:DST:C13	1.98	0.93
1:B:80:HIS:CD2	1:B:108:LYS:HG3	2.04	0.92
1:C:8:PRO:HD2	2:C:301:DST:H143	1.56	0.85
1:F:74:GLU:OE1	1:F:77:ARG:NH1	2.10	0.85
1:E:43:GLU:HG3	1:E:206:ARG:HH22	1.42	0.84
1:F:169:ARG:NH2	2:F:301:DST:O5	2.09	0.82
1:B:1:MET:H3	1:B:46:LYS:HG3	1.45	0.81
1:F:32:LYS:NZ	1:F:193:THR:O	2.12	0.81
1:A:43:GLU:HG2	1:A:202:LEU:HD13	1.62	0.81
1:H:43:GLU:HG3	1:H:206:ARG:HH12	1.45	0.81
1:G:43:GLU:HG3	1:G:206:ARG:HH12	1.49	0.78
1:G:195:ARG:NH2	1:G:199:ASP:OD1	2.18	0.77
2:G:301:DST:H142	2:G:302:DST:H142	1.66	0.77
1:B:41:LYS:NZ	1:B:82:GLN:O	2.18	0.76
1:B:80:HIS:HD2	1:B:108:LYS:HG3	1.50	0.73
1:G:43:GLU:HG2	1:G:202:LEU:HD13	1.70	0.73
1:E:24:ASP:OD1	4:E:402:HOH:O	2.07	0.70
2:G:301:DST:H111	2:G:302:DST:H101	1.71	0.70
1:A:43:GLU:HG3	1:A:206:ARG:NH1	2.06	0.69
1:C:51:THR:O	2:C:301:DST:H141	1.92	0.69
1:D:40:ALA:HB1	1:D:45:VAL:CG2	2.23	0.68
1:F:20:VAL:HG12	1:F:21:SER:N	2.07	0.67
1:C:8:PRO:HD2	2:C:301:DST:C14	2.25	0.67
1:A:43:GLU:HG2	1:A:202:LEU:CD1	2.25	0.67
2:C:301:DST:H101	2:C:302:DST:H102	1.77	0.66
2:E:301:DST:H101	2:E:302:DST:H112	1.76	0.66
1:B:1:MET:N	1:B:46:LYS:HG3	2.10	0.66
1:G:2:THR:HG22	1:G:45:VAL:HA	1.78	0.66
1:D:40:ALA:HB1	1:D:45:VAL:HG21	1.77	0.65
1:H:180:TYR:O	1:H:210:ARG:NH2	2.27	0.65
2:C:301:DST:H101	2:C:302:DST:C10	2.27	0.64
1:F:13:ARG:NH1	2:F:302:DST:O4	2.31	0.64
1:G:11:MET:HE2	2:G:302:DST:S9	2.38	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:301:DST:H102	2:B:302:DST:H112	1.79	0.63
1:E:202:LEU:O	1:E:206:ARG:HG2	1.99	0.63
1:E:104:ALA:O	1:E:108:LYS:HG3	1.99	0.62
1:E:43:GLU:HG3	1:E:206:ARG:NH2	2.14	0.62
1:A:175:PRO:HG2	1:A:178:SER:HB2	1.82	0.62
1:D:195:ARG:HD2	4:D:477:HOH:O	1.98	0.61
2:G:301:DST:H111	2:G:302:DST:C10	2.30	0.60
1:F:20:VAL:CG1	1:F:24:ASP:HB2	2.31	0.60
2:H:301:DST:H101	2:H:302:DST:H101	1.82	0.59
1:C:8:PRO:CD	2:C:301:DST:H143	2.30	0.59
1:G:8:PRO:HD2	2:G:301:DST:H141	1.83	0.59
1:C:51:THR:OG1	2:C:301:DST:C14	2.50	0.59
1:F:20:VAL:CG1	1:F:21:SER:N	2.65	0.58
1:C:195:ARG:HD2	4:C:525:HOH:O	2.04	0.57
1:E:202:LEU:O	1:E:206:ARG:CG	2.53	0.57
1:E:43:GLU:HG2	1:E:202:LEU:HD13	1.86	0.57
1:B:83:PHE:HE1	1:B:85:PHE:HB2	1.70	0.57
1:F:70:GLU:OE2	1:F:77:ARG:NH1	2.29	0.57
1:H:43:GLU:HG2	1:H:202:LEU:HD13	1.86	0.56
1:H:18:ASN:ND2	4:H:405:HOH:O	2.38	0.56
1:F:77:ARG:NE	1:F:101:GLU:OE2	2.21	0.56
1:G:40:ALA:HB1	1:G:45:VAL:HG21	1.88	0.55
1:B:80:HIS:CD2	1:B:108:LYS:CG	2.84	0.54
1:A:11:MET:HE2	2:A:302:DST:S9	2.47	0.54
1:C:51:THR:OG1	2:C:301:DST:H141	2.08	0.54
2:E:301:DST:H142	2:E:302:DST:H142	1.89	0.54
1:H:43:GLU:HG3	1:H:206:ARG:NH1	2.17	0.54
1:A:5:MET:HE3	1:A:7:LEU:HD11	1.90	0.54
1:F:12:ARG:HG2	1:F:22:LEU:HD13	1.89	0.53
1:G:160:TYR:HB3	1:G:205:LEU:HD13	1.91	0.53
1:F:166:GLY:HA2	1:F:187:PRO:HA	1.91	0.53
1:D:51:THR:OG1	2:D:301:DST:H131	2.09	0.52
1:A:12:ARG:NH2	1:B:12:ARG:HD2	2.25	0.52
1:B:208:ARG:NH1	1:F:168:ILE:HB	2.26	0.51
1:C:143:MET:HB2	1:H:133:PHE:CD2	2.45	0.51
1:E:198:PHE:CZ	1:E:202:LEU:HD11	2.45	0.51
1:B:143:MET:HB2	1:F:133:PHE:CD1	2.46	0.51
1:B:11:MET:HE3	1:B:29:MET:SD	2.51	0.50
1:G:37:MET:HE1	1:G:41:LYS:HE3	1.92	0.50
2:B:301:DST:H101	2:B:302:DST:H101	1.92	0.50
1:H:5:MET:HE2	1:H:51:THR:CG2	2.41	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:40:ALA:HB1	1:G:45:VAL:CG2	2.42	0.50
1:C:11:MET:HE2	2:C:302:DST:S9	2.51	0.50
1:F:114:ASP:OD1	1:F:114:ASP:N	2.32	0.50
2:E:301:DST:C10	2:E:302:DST:H112	2.42	0.50
1:F:20:VAL:HG12	1:F:24:ASP:HB2	1.93	0.49
1:D:26:TYR:HE1	2:D:302:DST:H112	1.77	0.49
1:D:51:THR:O	2:D:301:DST:H141	2.12	0.49
1:F:46:LYS:N	1:F:46:LYS:HD2	2.28	0.49
1:H:12:ARG:HH21	1:H:12:ARG:HG3	1.76	0.49
2:A:301:DST:H101	2:A:302:DST:H112	1.93	0.49
1:F:41:LYS:O	1:F:42:GLU:C	2.51	0.49
2:H:301:DST:H142	2:H:302:DST:H111	1.95	0.49
1:C:51:THR:OG1	2:C:301:DST:H142	2.13	0.49
1:D:11:MET:HE3	1:D:29:MET:SD	2.53	0.49
1:C:166:GLY:HA2	1:C:187:PRO:HA	1.95	0.49
1:G:163:ARG:NH2	2:G:301:DST:O6	2.41	0.48
1:H:43:GLU:HG2	1:H:202:LEU:CD1	2.44	0.48
1:E:40:ALA:O	1:E:45:VAL:HG13	2.14	0.48
1:F:160:TYR:HB3	1:F:205:LEU:HD13	1.96	0.48
1:D:40:ALA:HB1	1:D:45:VAL:HG22	1.95	0.47
1:H:198:PHE:CZ	1:H:202:LEU:HD11	2.49	0.47
2:F:301:DST:H101	2:F:302:DST:H101	1.97	0.47
1:H:84:LYS:NZ	4:H:406:HOH:O	2.48	0.47
1:B:49:TYR:OH	1:B:156:THR:HB	2.14	0.47
1:B:143:MET:HG2	1:F:130:VAL:HA	1.97	0.46
1:C:130:VAL:HG22	1:H:143:MET:HG3	1.97	0.46
1:F:20:VAL:CG1	1:F:21:SER:H	2.28	0.46
1:H:41:LYS:NZ	1:H:82:GLN:O	2.48	0.46
1:E:13:ARG:NH1	2:E:302:DST:O5	2.43	0.46
1:G:43:GLU:HG3	1:G:206:ARG:NH1	2.24	0.46
2:H:301:DST:H101	2:H:302:DST:H133	1.98	0.46
1:A:143:MET:HB2	1:E:133:PHE:CD1	2.51	0.46
1:B:166:GLY:HA2	1:B:187:PRO:HA	1.98	0.45
1:A:131:GLY:O	1:A:135:LYS:HG3	2.16	0.45
1:E:208:ARG:HD3	1:E:208:ARG:HA	1.40	0.45
1:H:22:LEU:HD11	1:H:26:TYR:HE2	1.82	0.45
1:F:10:GLY:HA2	1:F:13:ARG:HH11	1.82	0.44
1:G:32:LYS:HA	1:G:32:LYS:HE2	1.99	0.44
1:B:11:MET:CE	2:B:302:DST:H143	2.46	0.44
1:G:86:ASP:O	1:G:118:HIS:HA	2.18	0.44
1:G:43:GLU:HG2	1:G:202:LEU:CD1	2.42	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:149:ALA:HB2	1:B:177:MET:HE2	1.98	0.44
2:B:301:DST:C10	2:B:302:DST:H112	2.44	0.44
1:F:5:MET:HB3	1:F:161:ILE:HG12	1.99	0.44
1:B:43:GLU:HB3	1:B:206:ARG:HH12	1.82	0.44
1:F:40:ALA:HB1	1:F:45:VAL:CG2	2.48	0.44
1:B:108:LYS:HB3	1:B:108:LYS:HE2	1.54	0.44
1:C:160:TYR:CZ	1:C:205:LEU:HB2	2.53	0.44
2:B:301:DST:S9	2:B:301:DST:H132	2.58	0.43
1:A:60:ARG:HB3	1:A:61:PRO:HD2	2.01	0.43
1:D:11:MET:HE2	2:D:302:DST:S9	2.58	0.43
1:D:49:TYR:OH	1:D:156:THR:HB	2.18	0.43
1:A:166:GLY:HA2	1:A:187:PRO:HA	2.01	0.43
1:A:210:ARG:O	1:A:210:ARG:HG2	2.17	0.43
1:C:84:LYS:HD3	4:C:521:HOH:O	2.18	0.43
1:D:5:MET:HE3	1:D:7:LEU:HD11	1.99	0.43
1:B:57:ASN:O	1:B:60:ARG:HD2	2.18	0.43
1:C:52:ALA:HB3	1:C:121:LEU:HD23	2.00	0.43
1:F:158:VAL:N	1:F:178:SER:OG	2.47	0.43
1:D:130:VAL:HG22	1:G:143:MET:HG3	2.01	0.42
1:D:143:MET:HB2	1:G:133:PHE:CD1	2.54	0.42
1:F:65:VAL:HA	1:F:68:PHE:CE2	2.54	0.42
1:E:49:TYR:CD1	1:E:118:HIS:HB3	2.54	0.42
1:A:43:GLU:CG	1:A:202:LEU:HD13	2.39	0.42
1:B:32:LYS:HA	1:B:32:LYS:HD3	1.79	0.42
1:F:3:THR:HG23	1:F:47:THR:HB	2.01	0.42
1:G:204:ASP:CG	1:G:208:ARG:HH21	2.27	0.42
1:C:133:PHE:CD1	1:H:143:MET:HB2	2.55	0.42
1:D:168:ILE:HB	1:G:208:ARG:NH1	2.35	0.42
1:F:41:LYS:O	1:F:43:GLU:N	2.53	0.42
1:A:198:PHE:CZ	1:A:202:LEU:HD11	2.54	0.42
1:F:160:TYR:CD1	1:F:205:LEU:HB3	2.55	0.42
1:B:83:PHE:CD1	1:B:83:PHE:C	2.98	0.41
1:H:96:GLU:H	1:H:96:GLU:CD	2.28	0.41
1:B:86:ASP:O	1:B:118:HIS:HA	2.20	0.41
1:B:50:ILE:HD12	1:B:117:LEU:HD11	2.03	0.41
1:F:52:ALA:O	2:F:302:DST:H132	2.21	0.41
1:F:175:PRO:HG2	1:F:178:SER:HB2	2.03	0.41
1:F:188:VAL:HG12	1:F:189:LEU:O	2.21	0.41
1:D:133:PHE:CD1	1:D:133:PHE:C	2.99	0.41
1:E:91:LEU:HD22	1:E:99:LEU:HD21	2.02	0.41
1:F:45:VAL:HG23	1:F:115:PHE:HE1	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:49:TYR:CD1	1:F:118:HIS:HB3	2.56	0.41
1:G:198:PHE:CZ	1:G:202:LEU:HD11	2.55	0.41
1:G:208:ARG:HD3	4:G:438:HOH:O	2.19	0.41
1:H:50:ILE:HD12	1:H:117:LEU:HD11	2.02	0.40
1:H:195:ARG:HG2	1:H:195:ARG:HH11	1.86	0.40
1:H:166:GLY:HA2	1:H:187:PRO:HA	2.03	0.40
1:E:43:GLU:HG2	1:E:202:LEU:CD1	2.52	0.40
2:E:301:DST:H101	2:E:302:DST:H101	2.03	0.40
1:B:198:PHE:CZ	1:B:202:LEU:HD11	2.56	0.40
1:E:11:MET:HE3	1:E:29:MET:SD	2.61	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	208/217 (96%)	205 (99%)	3 (1%)	0	100	100
1	B	206/217 (95%)	203 (98%)	3 (2%)	0	100	100
1	C	215/217 (99%)	208 (97%)	7 (3%)	0	100	100
1	D	206/217 (95%)	202 (98%)	4 (2%)	0	100	100
1	E	206/217 (95%)	203 (98%)	3 (2%)	0	100	100
1	F	206/217 (95%)	204 (99%)	2 (1%)	0	100	100
1	G	206/217 (95%)	203 (98%)	3 (2%)	0	100	100
1	H	214/217 (99%)	209 (98%)	5 (2%)	0	100	100
All	All	1667/1736 (96%)	1637 (98%)	30 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	179/183 (98%)	179 (100%)	0	100	100
1	B	177/183 (97%)	175 (99%)	2 (1%)	65	68
1	C	183/183 (100%)	179 (98%)	4 (2%)	45	44
1	D	177/183 (97%)	176 (99%)	1 (1%)	78	83
1	E	177/183 (97%)	175 (99%)	2 (1%)	65	68
1	F	177/183 (97%)	175 (99%)	2 (1%)	65	68
1	G	177/183 (97%)	173 (98%)	4 (2%)	44	42
1	H	182/183 (100%)	179 (98%)	3 (2%)	55	56
All	All	1429/1464 (98%)	1411 (99%)	18 (1%)	61	63

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	108	LYS
1	B	156	THR
1	C	32	LYS
1	C	45	VAL
1	C	172	SER
1	C	210	ARG
1	D	84	LYS
1	E	45	VAL
1	E	206	ARG
1	F	2	THR
1	F	16	GLU
1	G	2	THR
1	G	17	LYS
1	G	32	LYS
1	G	45	VAL
1	H	2	THR
1	H	22	LEU
1	H	32	LYS

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

Mol	Chain	Res	Type
1	A	18	ASN
1	A	57	ASN
1	A	82	GLN
1	B	18	ASN
1	C	18	ASN
1	C	82	GLN
1	E	127	HIS
1	G	57	ASN
1	G	82	GLN
1	H	18	ASN
1	H	82	GLN

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](#)

Of 24 ligands modelled in this entry, 8 are monoatomic - leaving 16 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	DST	E	301	3	10,13,13	2.63	5 (50%)	12,19,19	1.23	1 (8%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	DST	B	301	3	10,13,13	2.38	5 (50%)	12,19,19	1.11	1 (8%)
2	DST	F	302	3	10,13,13	1.54	2 (20%)	12,19,19	1.06	1 (8%)
2	DST	C	301	3	10,13,13	3.00	5 (50%)	12,19,19	1.66	2 (16%)
2	DST	F	301	3	10,13,13	2.34	4 (40%)	12,19,19	1.34	1 (8%)
2	DST	H	302	3	10,13,13	1.74	3 (30%)	12,19,19	1.35	2 (16%)
2	DST	A	301	3	10,13,13	2.96	5 (50%)	12,19,19	1.01	1 (8%)
2	DST	A	302	3	10,13,13	2.55	5 (50%)	12,19,19	1.60	2 (16%)
2	DST	G	302	3	10,13,13	1.55	3 (30%)	12,19,19	1.23	2 (16%)
2	DST	C	302	3	10,13,13	2.87	5 (50%)	12,19,19	1.30	1 (8%)
2	DST	G	301	3	10,13,13	1.62	2 (20%)	12,19,19	1.27	1 (8%)
2	DST	H	301	3	10,13,13	1.54	1 (10%)	12,19,19	1.13	0
2	DST	E	302	3	10,13,13	2.27	4 (40%)	12,19,19	1.54	1 (8%)
2	DST	B	302	3	10,13,13	2.36	4 (40%)	12,19,19	1.59	2 (16%)
2	DST	D	301	3	10,13,13	2.58	4 (40%)	12,19,19	1.64	5 (41%)
2	DST	D	302	3	10,13,13	2.44	4 (40%)	12,19,19	1.40	1 (8%)

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	DST	E	301	3	-	2/7/13/13	-
2	DST	B	301	3	-	0/7/13/13	-
2	DST	F	302	3	-	2/7/13/13	-
2	DST	C	301	3	-	0/7/13/13	-
2	DST	F	301	3	-	0/7/13/13	-
2	DST	H	302	3	-	2/7/13/13	-
2	DST	A	301	3	-	0/7/13/13	-
2	DST	A	302	3	-	1/7/13/13	-
2	DST	G	302	3	-	2/7/13/13	-
2	DST	C	302	3	-	2/7/13/13	-
2	DST	G	301	3	-	0/7/13/13	-
2	DST	H	301	3	-	2/7/13/13	-
2	DST	E	302	3	-	0/7/13/13	-
2	DST	B	302	3	-	2/7/13/13	-
2	DST	D	301	3	-	2/7/13/13	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	DST	D	302	3	-	2/7/13/13	-

All (61) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	301	DST	C10-S9	-6.40	1.76	1.84
2	A	301	DST	C10-S9	-5.86	1.76	1.84
2	D	301	DST	C10-S9	-5.45	1.77	1.84
2	E	301	DST	C10-S9	-5.27	1.77	1.84
2	C	302	DST	C10-S9	-4.81	1.78	1.84
2	D	302	DST	P3-O8	-4.61	1.45	1.56
2	F	301	DST	C10-S9	-4.58	1.78	1.84
2	A	302	DST	P3-O8	-4.54	1.45	1.56
2	B	302	DST	C10-S9	-4.49	1.78	1.84
2	A	301	DST	P3-O8	-4.35	1.45	1.56
2	C	302	DST	P3-O8	-4.34	1.45	1.56
2	B	301	DST	C10-S9	-4.34	1.78	1.84
2	F	301	DST	P3-O8	-4.18	1.46	1.56
2	E	302	DST	P3-O8	-4.16	1.46	1.56
2	A	302	DST	C10-S9	-4.12	1.79	1.84
2	E	301	DST	P3-O8	-3.93	1.46	1.56
2	B	302	DST	P3-O8	-3.80	1.47	1.56
2	D	301	DST	P3-O8	-3.68	1.47	1.56
2	G	301	DST	C10-S9	-3.67	1.79	1.84
2	C	301	DST	P3-O8	-3.67	1.47	1.56
2	D	302	DST	C10-S9	-3.67	1.79	1.84
2	C	302	DST	P1-O5	-3.62	1.39	1.50
2	C	301	DST	P1-O5	-3.55	1.39	1.50
2	C	302	DST	P1-O6	-3.51	1.41	1.54
2	A	301	DST	P1-O6	-3.51	1.41	1.54
2	D	302	DST	P1-O6	-3.45	1.42	1.54
2	B	301	DST	P3-O8	-3.38	1.48	1.56
2	C	301	DST	P1-O6	-3.35	1.42	1.54
2	A	301	DST	P1-O4	-3.35	1.42	1.54
2	B	301	DST	P1-O6	-3.34	1.42	1.54
2	C	302	DST	P1-O4	-3.27	1.42	1.54
2	B	302	DST	P1-O4	-3.27	1.42	1.54
2	E	301	DST	P1-O4	-3.26	1.42	1.54
2	A	302	DST	P1-O4	-3.25	1.42	1.54
2	E	302	DST	C10-S9	-3.25	1.80	1.84
2	D	301	DST	P1-O6	-3.12	1.43	1.54
2	H	302	DST	C10-S9	-3.11	1.80	1.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	301	DST	P1-O6	-3.07	1.43	1.54
2	C	301	DST	P1-O4	-3.03	1.43	1.54
2	A	301	DST	P1-O5	-2.99	1.41	1.50
2	H	301	DST	C10-S9	-2.95	1.80	1.84
2	H	302	DST	P3-O7	2.91	1.53	1.46
2	D	301	DST	P1-O4	-2.87	1.44	1.54
2	E	302	DST	P1-O6	-2.83	1.44	1.54
2	F	301	DST	P1-O6	-2.75	1.44	1.54
2	F	302	DST	C10-S9	-2.74	1.80	1.84
2	D	302	DST	P1-O4	-2.63	1.45	1.54
2	A	302	DST	P1-O6	-2.61	1.45	1.54
2	G	302	DST	C10-S9	-2.60	1.80	1.84
2	B	301	DST	P1-O4	-2.59	1.45	1.54
2	E	302	DST	P1-O4	-2.53	1.45	1.54
2	A	302	DST	P1-O5	-2.46	1.42	1.50
2	F	302	DST	P3-O7	2.36	1.51	1.46
2	B	301	DST	P1-O5	-2.35	1.43	1.50
2	G	301	DST	P3-O7	2.31	1.51	1.46
2	H	302	DST	P3-O8	-2.20	1.51	1.56
2	B	302	DST	P1-O6	-2.19	1.46	1.54
2	F	301	DST	P1-O4	-2.15	1.46	1.54
2	G	302	DST	P3-O8	-2.12	1.51	1.56
2	E	301	DST	P1-O5	-2.10	1.43	1.50
2	G	302	DST	P3-O7	2.06	1.51	1.46

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	302	DST	O8-P3-O7	4.15	119.75	109.86
2	C	301	DST	C11-C10-S9	-3.77	93.47	111.87
2	B	302	DST	O8-P3-O7	3.58	118.38	109.86
2	A	302	DST	C14-C12-C13	-3.40	106.76	114.59
2	G	301	DST	C11-C10-S9	-3.29	95.84	111.87
2	F	301	DST	C14-C12-C13	-3.27	107.06	114.59
2	D	301	DST	O8-P3-O7	3.18	117.44	109.86
2	A	302	DST	O8-P3-O7	3.10	117.25	109.86
2	D	302	DST	C14-C12-C13	-3.06	107.55	114.59
2	H	302	DST	C14-C12-C13	-2.63	108.53	114.59
2	F	302	DST	C14-C12-C13	-2.63	108.53	114.59
2	B	302	DST	O6-P1-O2	2.61	113.40	104.64
2	C	302	DST	C14-C12-C13	-2.51	108.81	114.59
2	G	302	DST	C14-C12-C13	-2.40	109.07	114.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	301	DST	C13-C12-C11	2.37	129.76	122.66
2	D	301	DST	C14-C12-C13	-2.30	109.30	114.59
2	B	301	DST	O8-P3-O7	2.28	115.29	109.86
2	C	301	DST	O8-P3-O7	2.26	115.25	109.86
2	H	302	DST	O4-P1-O2	2.17	111.92	104.64
2	D	301	DST	O4-P1-O2	2.15	111.84	104.64
2	A	301	DST	O8-P3-O2	2.05	112.80	107.27
2	G	302	DST	O6-P1-O2	2.02	111.41	104.64
2	E	301	DST	O6-P1-O4	2.02	115.37	107.80
2	D	301	DST	O6-P1-O2	2.02	111.39	104.64

There are no chirality outliers.

All (19) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	302	DST	P1-O2-P3-O7
2	D	302	DST	P1-O2-P3-O7
2	B	302	DST	P1-O2-P3-O8
2	G	302	DST	P1-O2-P3-O8
2	C	302	DST	P1-O2-P3-O7
2	H	302	DST	P1-O2-P3-O7
2	D	302	DST	P1-O2-P3-O8
2	F	302	DST	P1-O2-P3-O8
2	H	301	DST	P1-O2-P3-O8
2	H	302	DST	P1-O2-P3-O8
2	F	302	DST	P1-O2-P3-O7
2	H	301	DST	P1-O2-P3-O7
2	D	301	DST	P1-O2-P3-O7
2	E	301	DST	P1-O2-P3-O7
2	D	301	DST	P1-O2-P3-O8
2	E	301	DST	P1-O2-P3-O8
2	G	302	DST	P1-O2-P3-O7
2	A	302	DST	P1-O2-P3-O8
2	C	302	DST	P1-O2-P3-O8

There are no ring outliers.

16 monomers are involved in 42 short contacts:

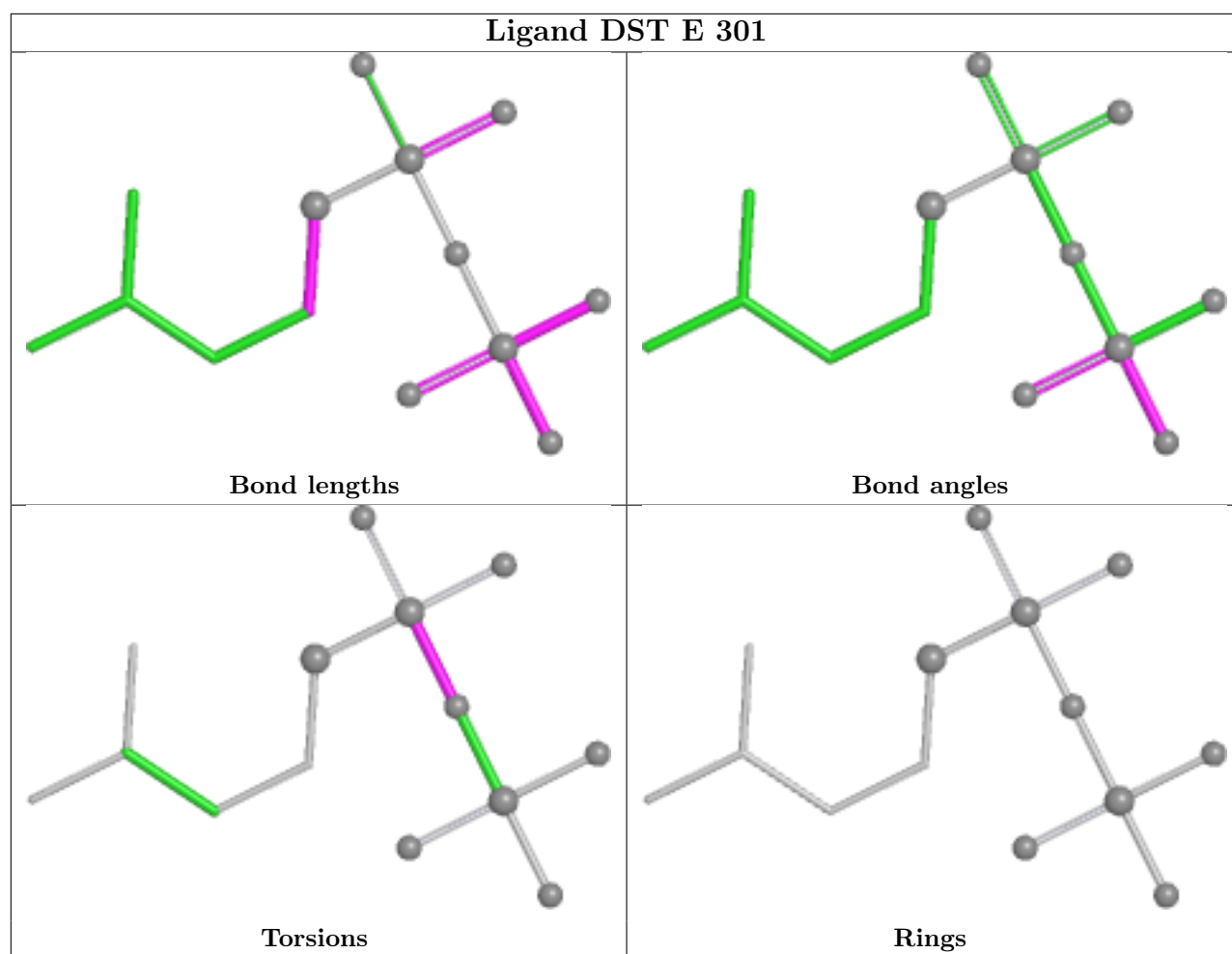
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	301	DST	4	0
2	B	301	DST	4	0

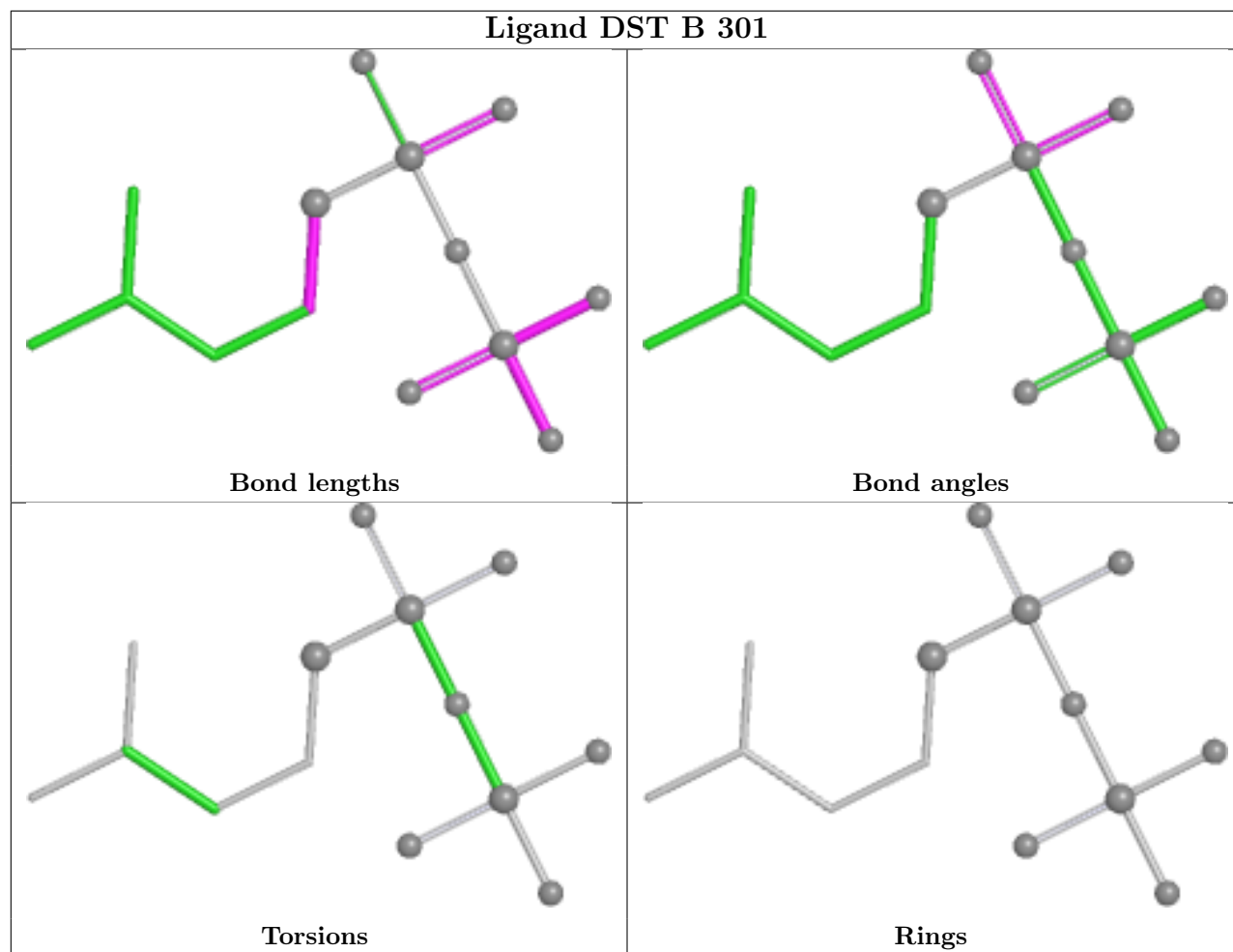
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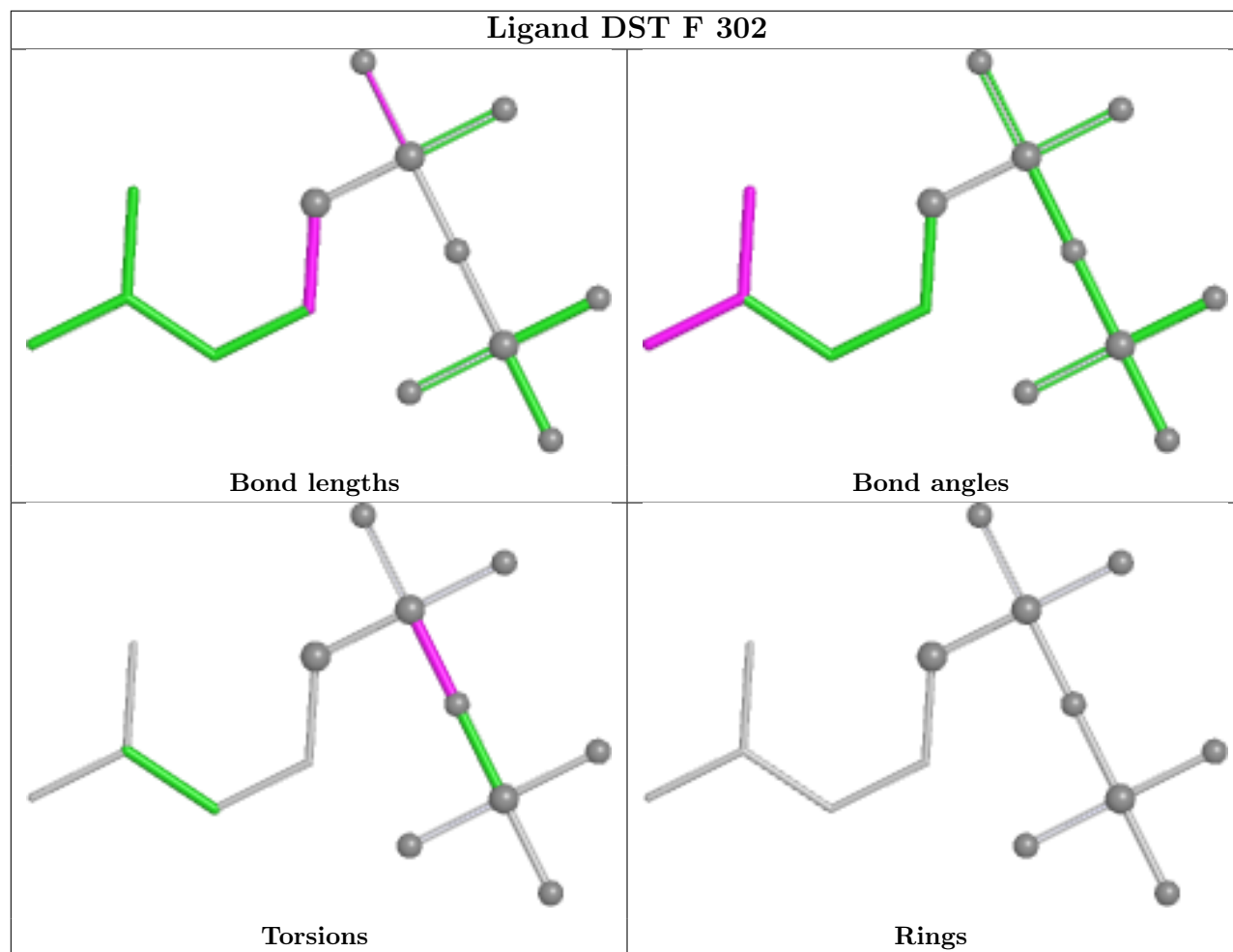
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	F	302	DST	3	0
2	C	301	DST	9	0
2	F	301	DST	2	0
2	H	302	DST	3	0
2	A	301	DST	1	0
2	A	302	DST	2	0
2	G	302	DST	4	0
2	C	302	DST	3	0
2	G	301	DST	5	0
2	H	301	DST	3	0
2	E	302	DST	5	0
2	B	302	DST	4	0
2	D	301	DST	5	0
2	D	302	DST	5	0

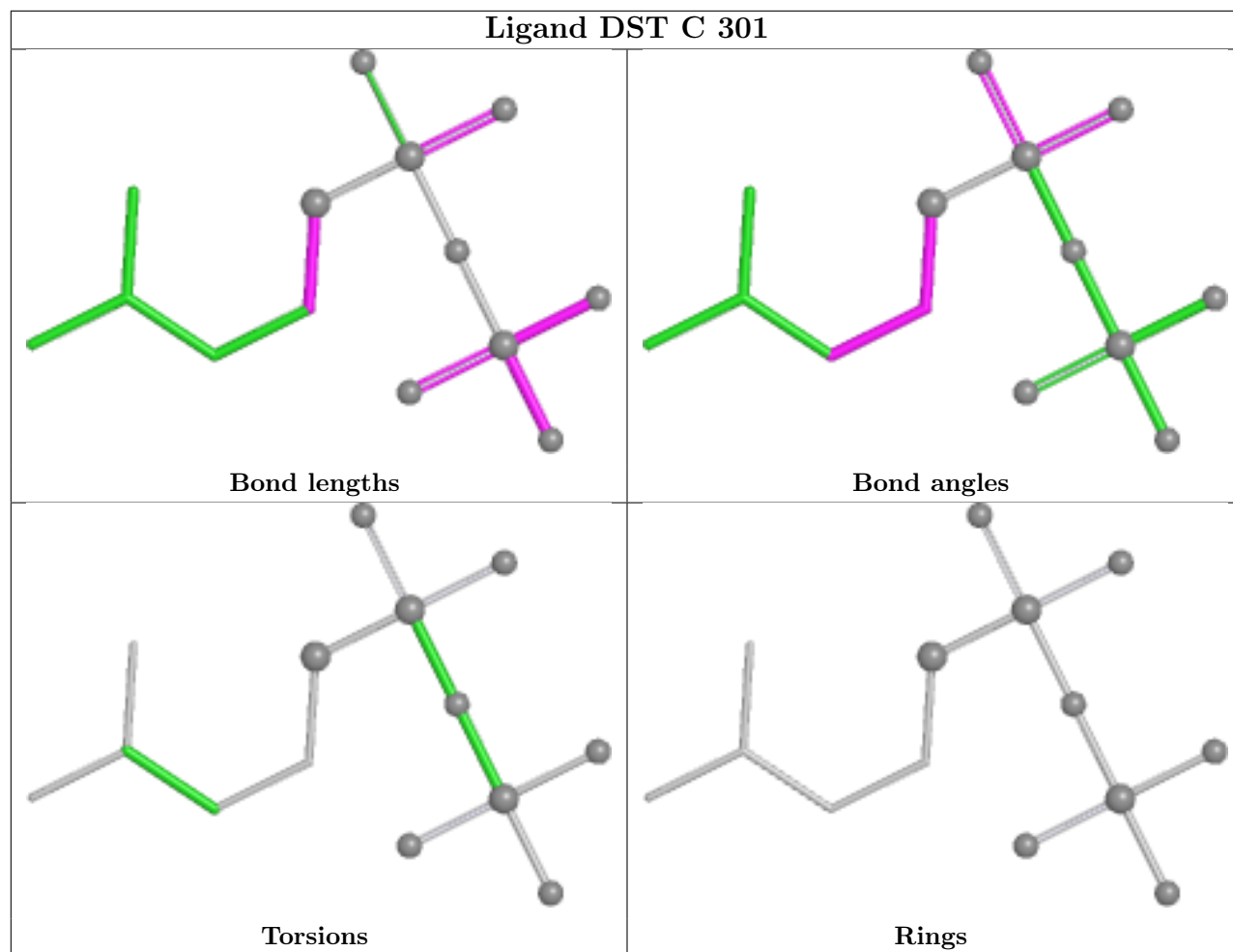
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

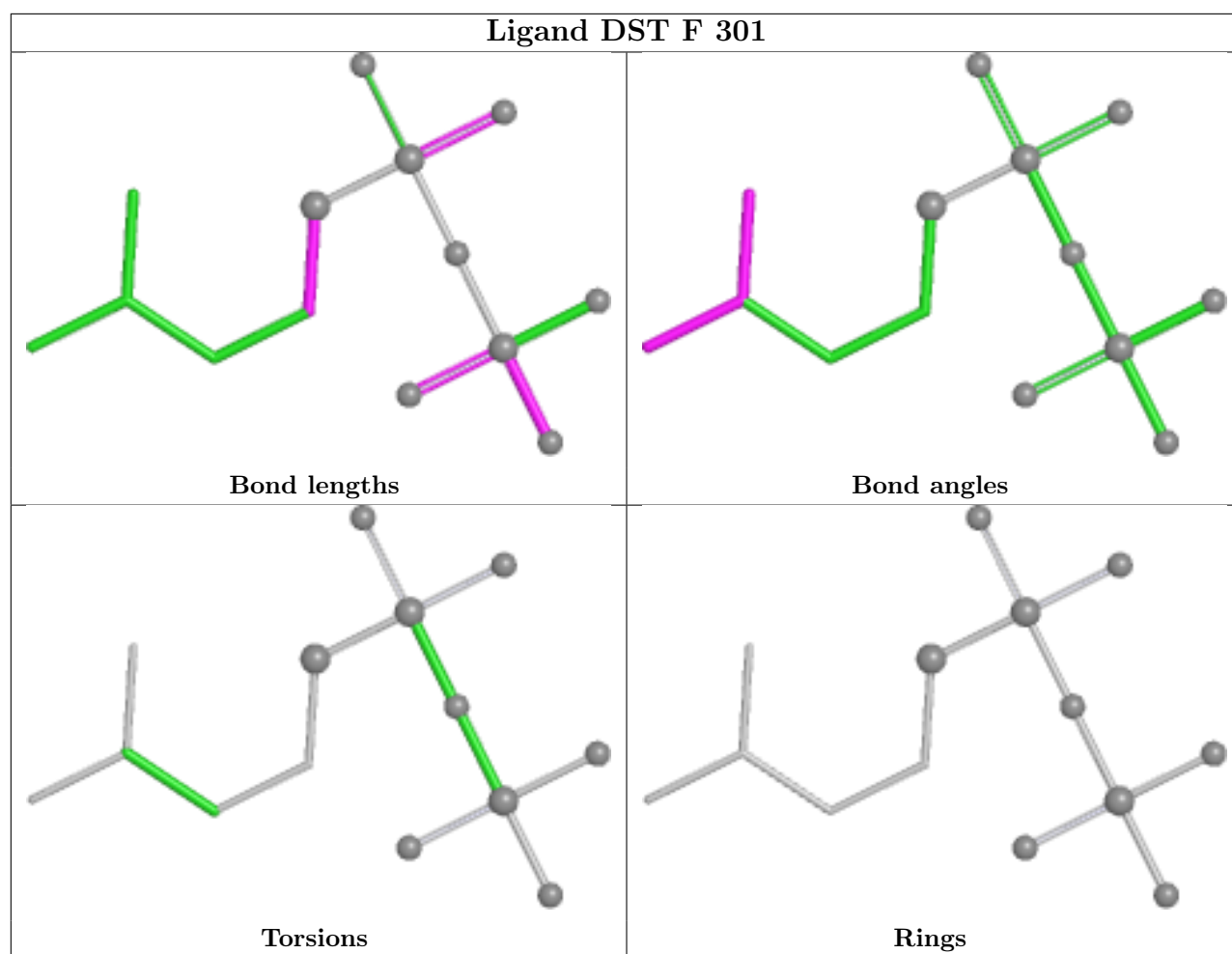


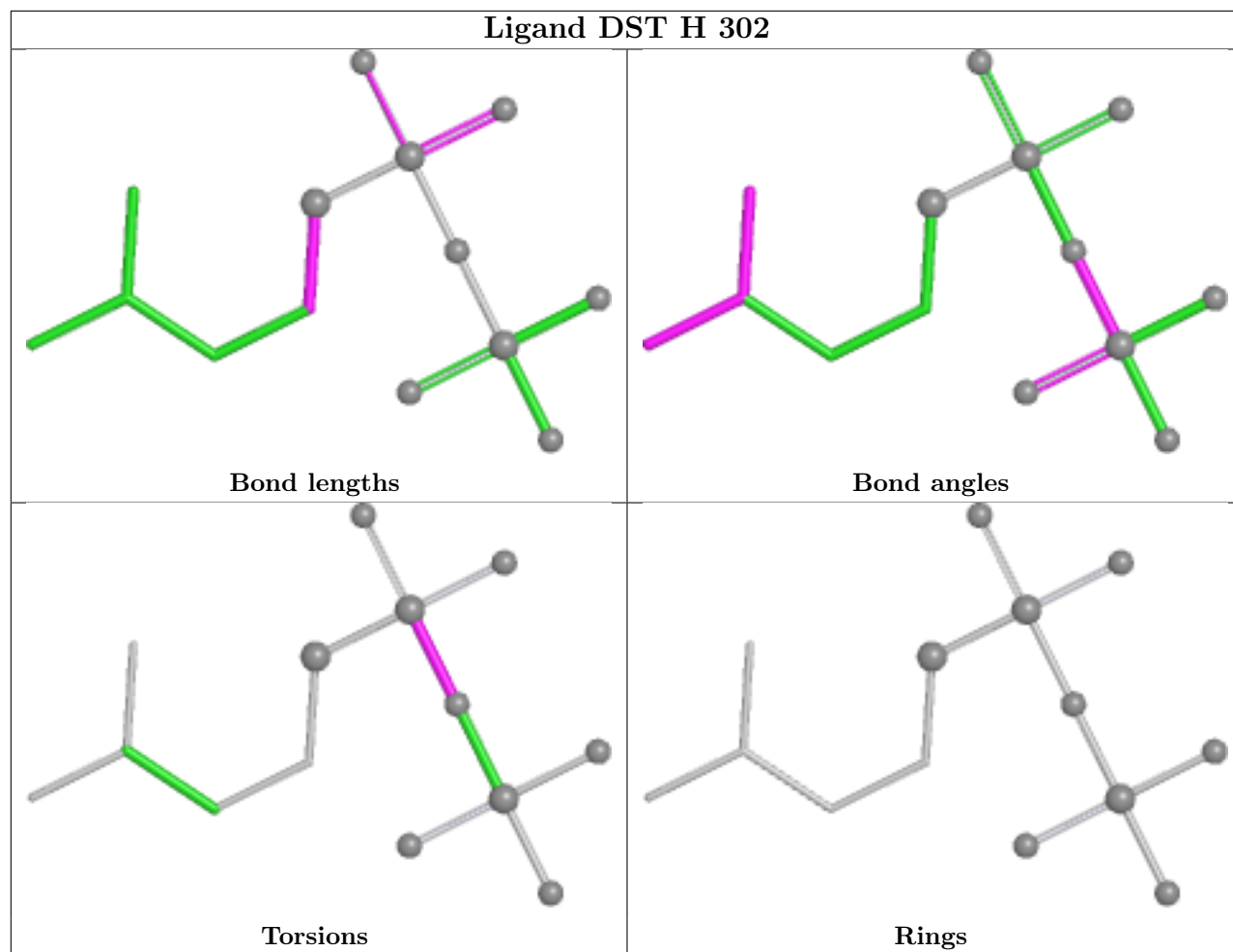


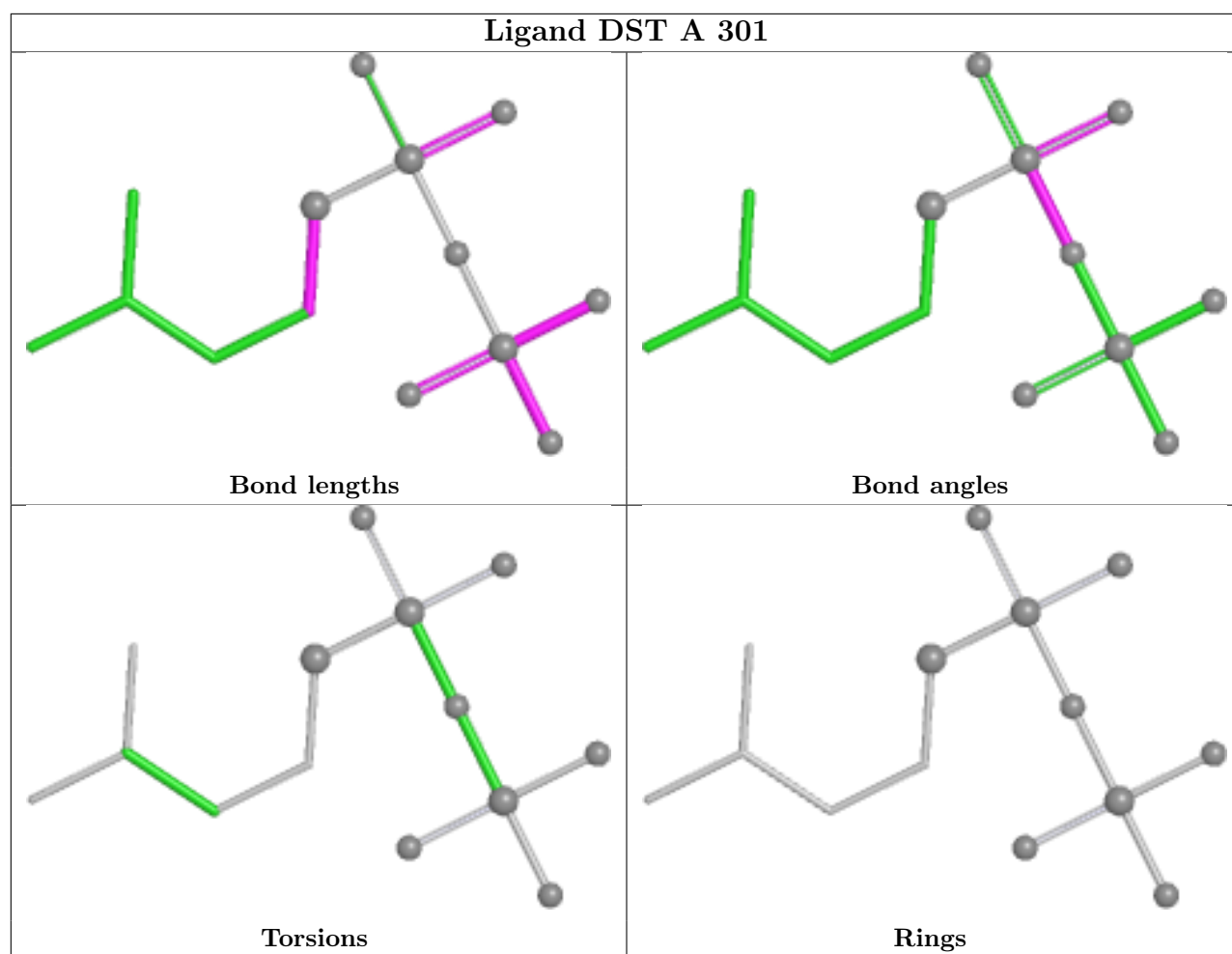
Ligand DST F 302

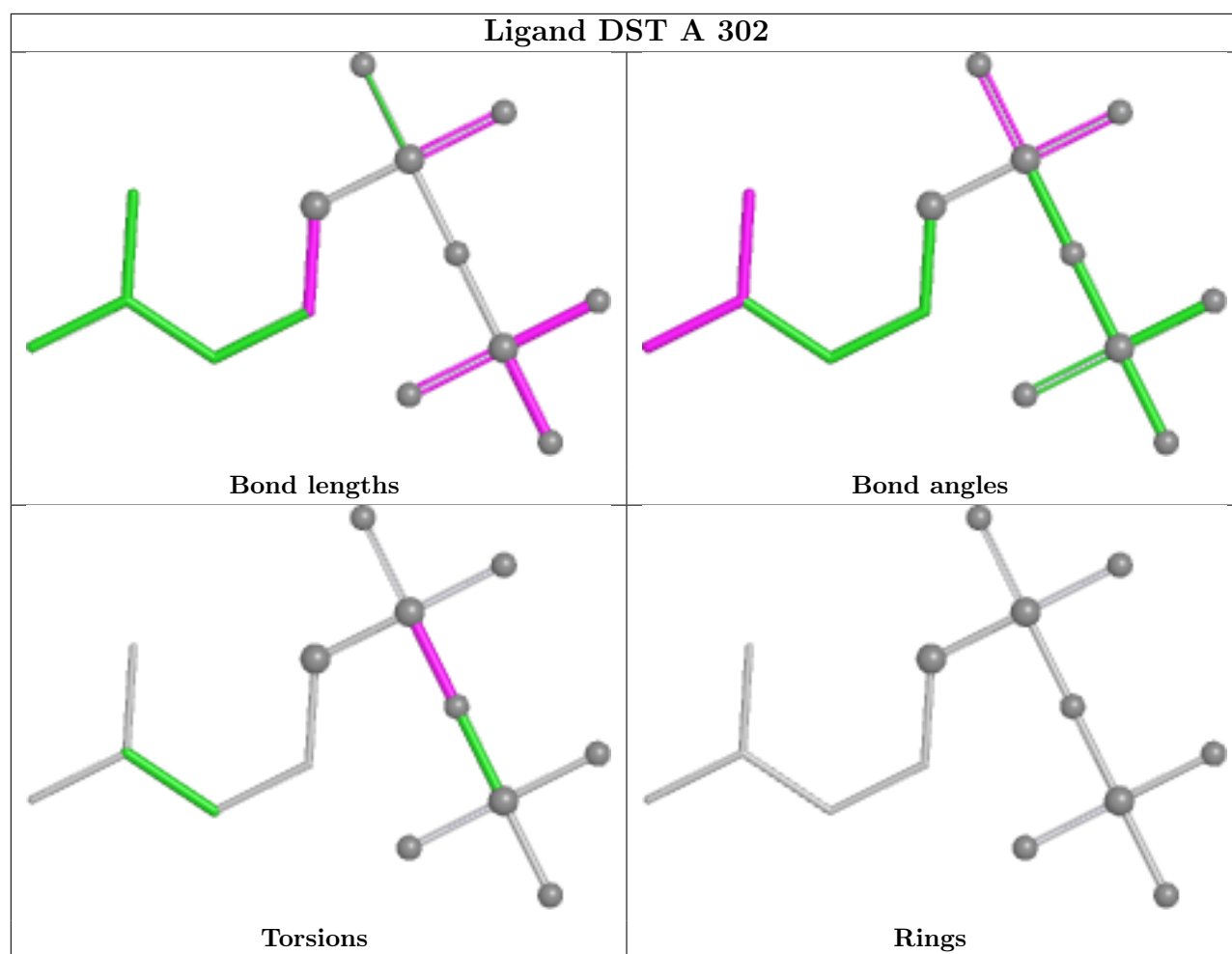


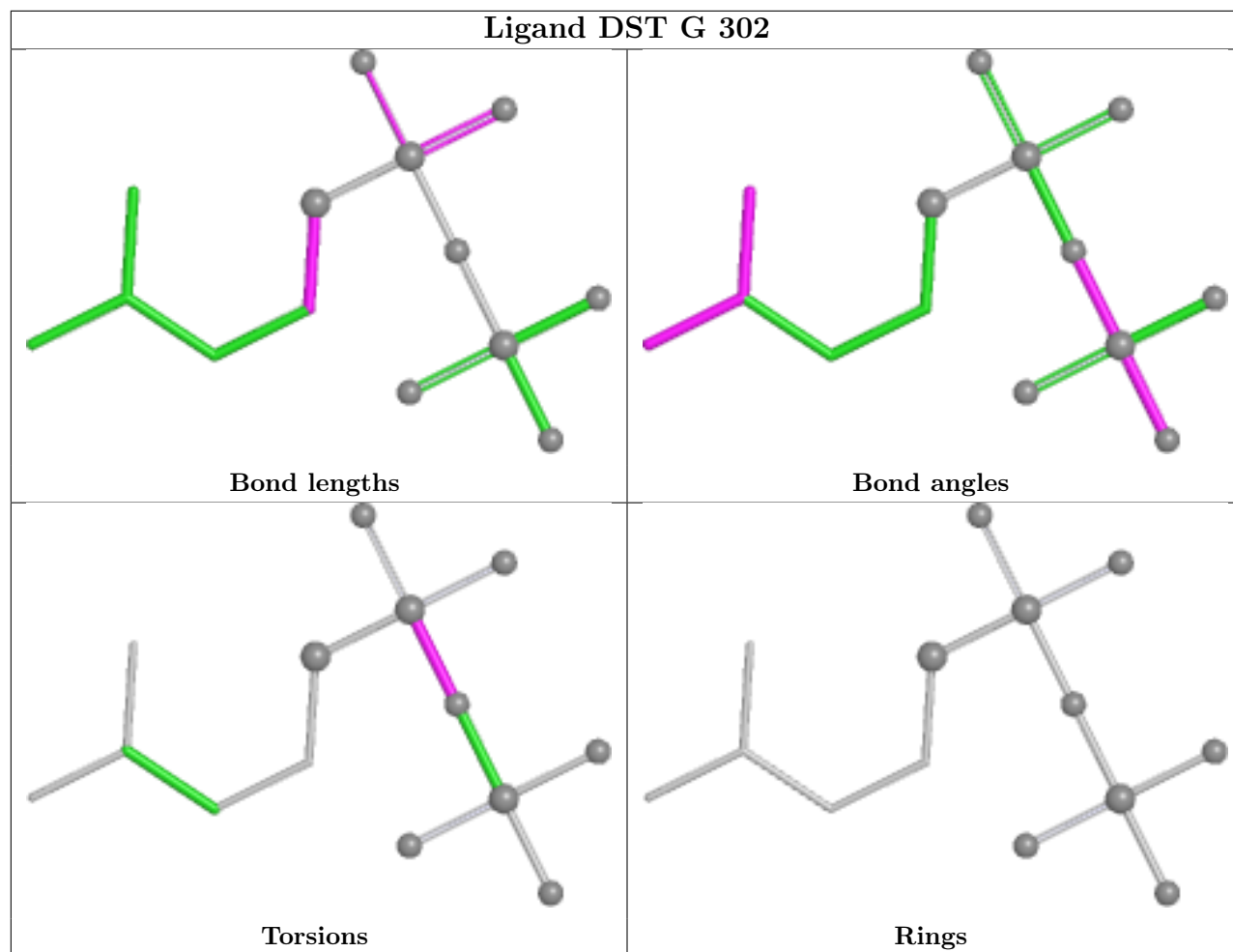


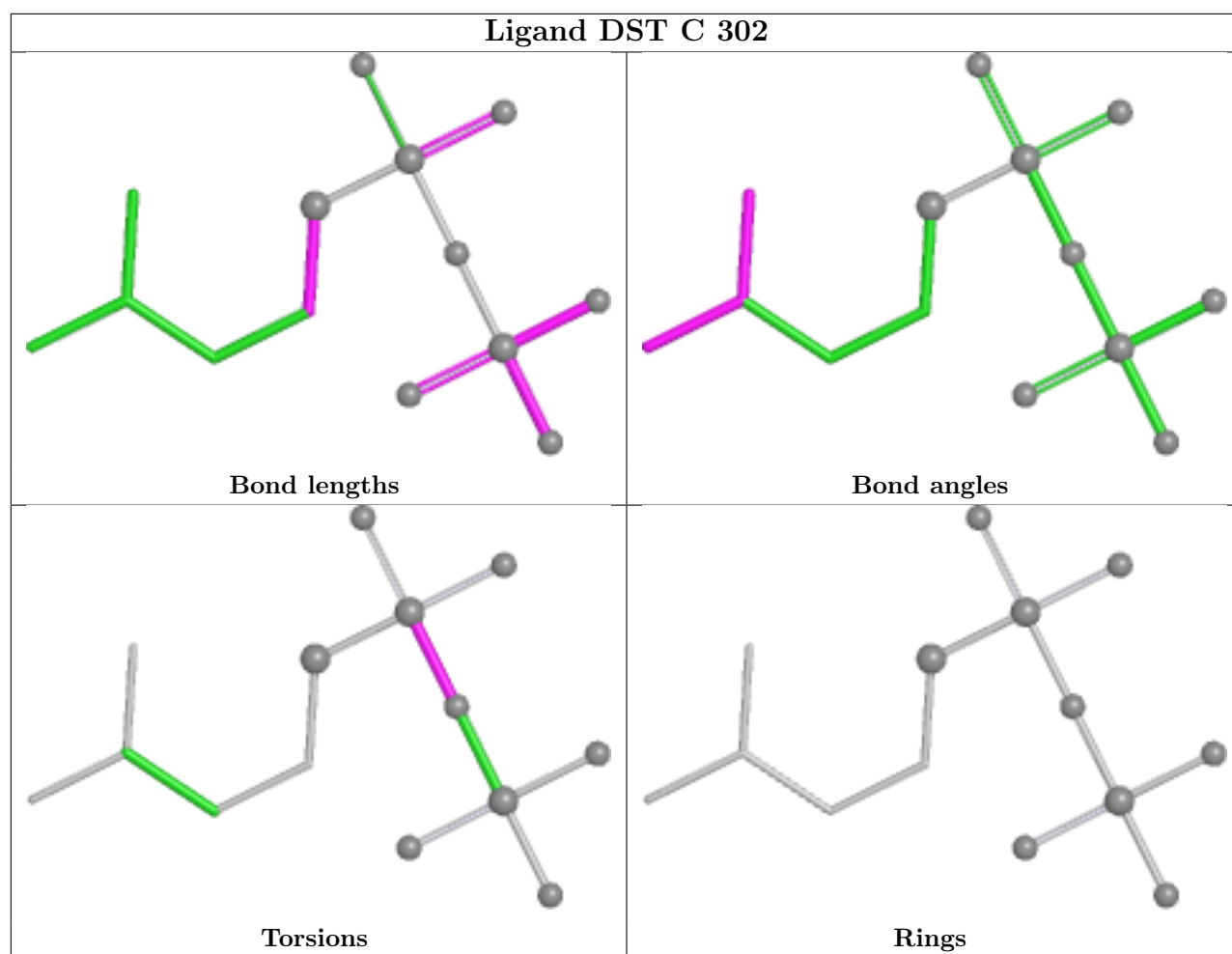


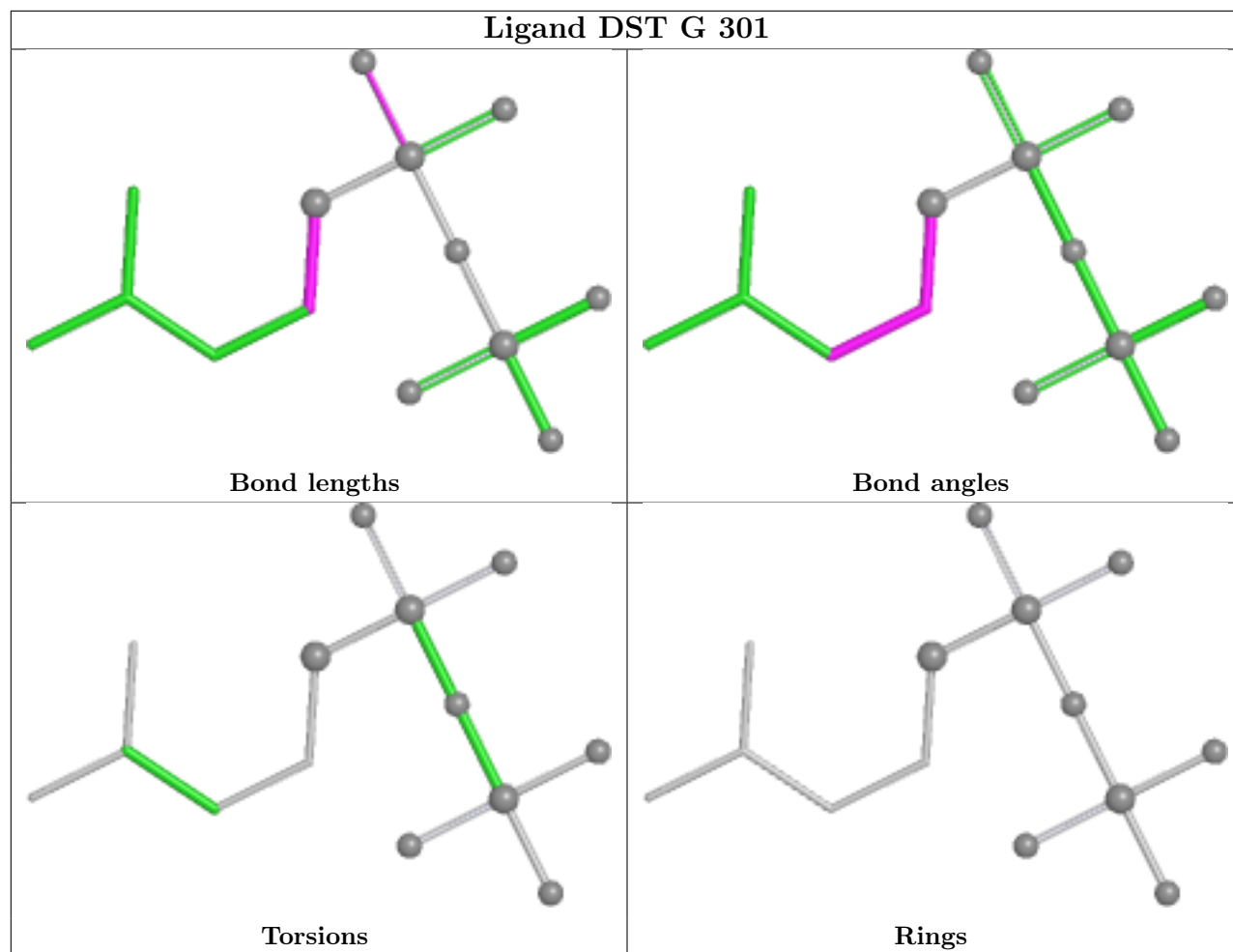


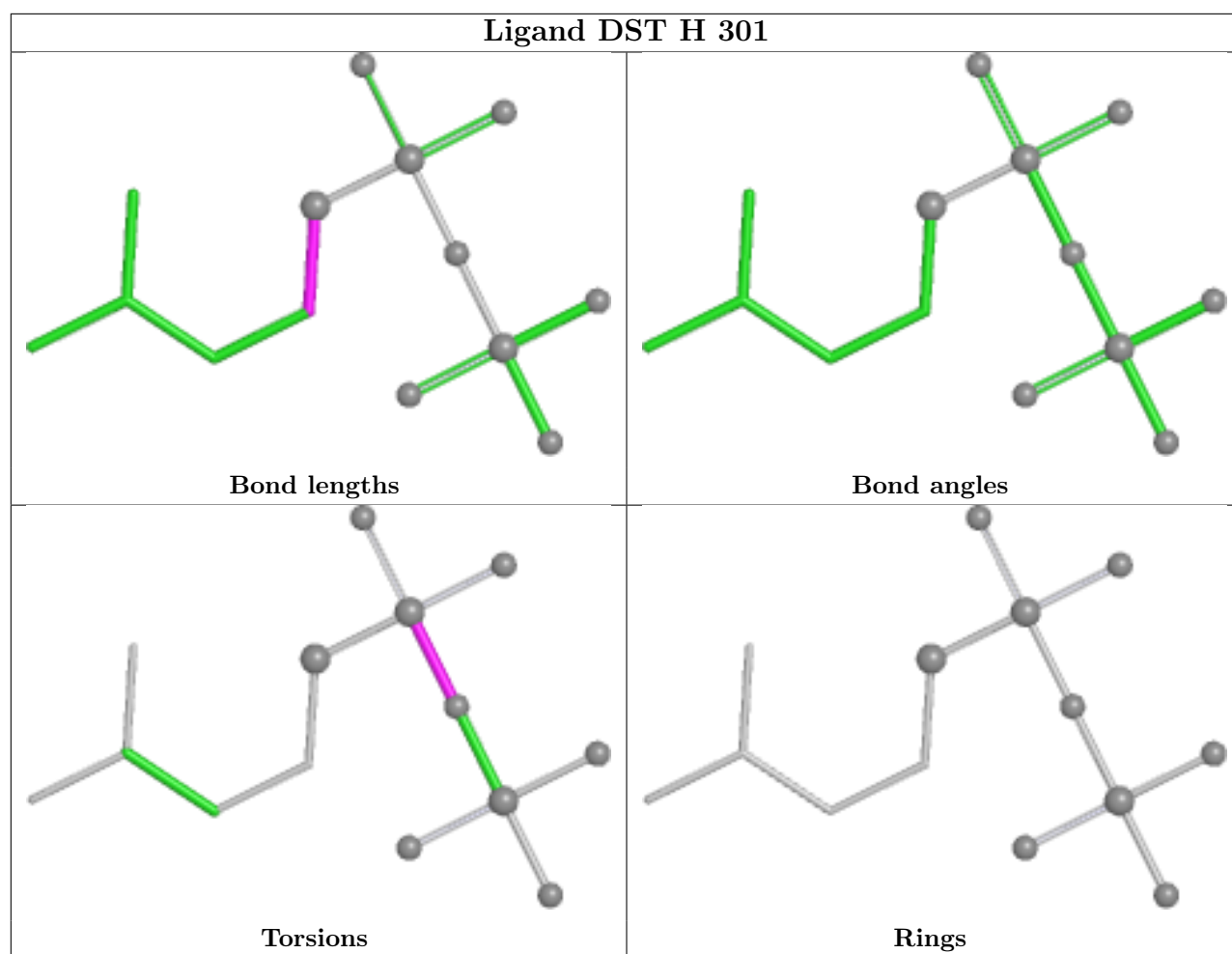




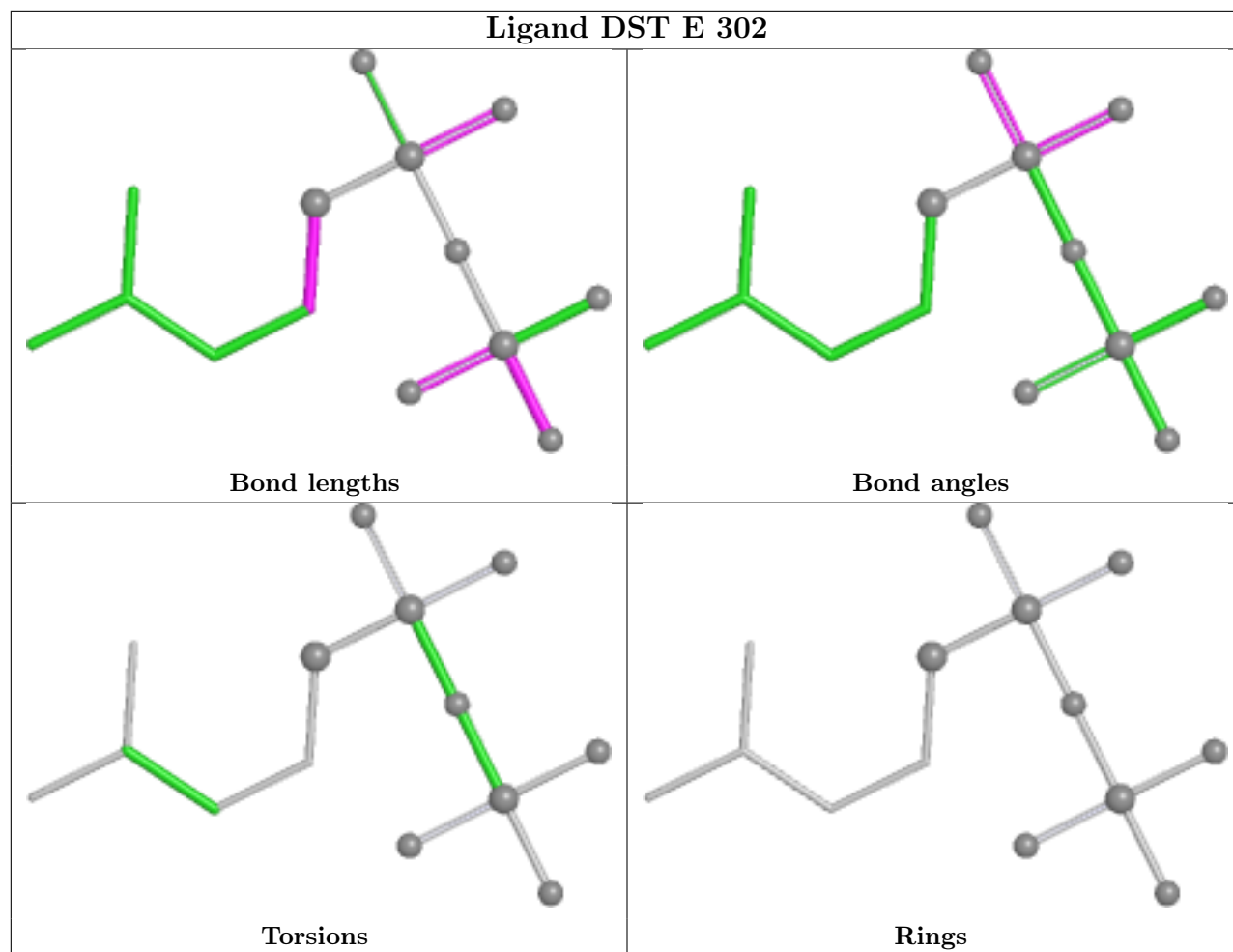


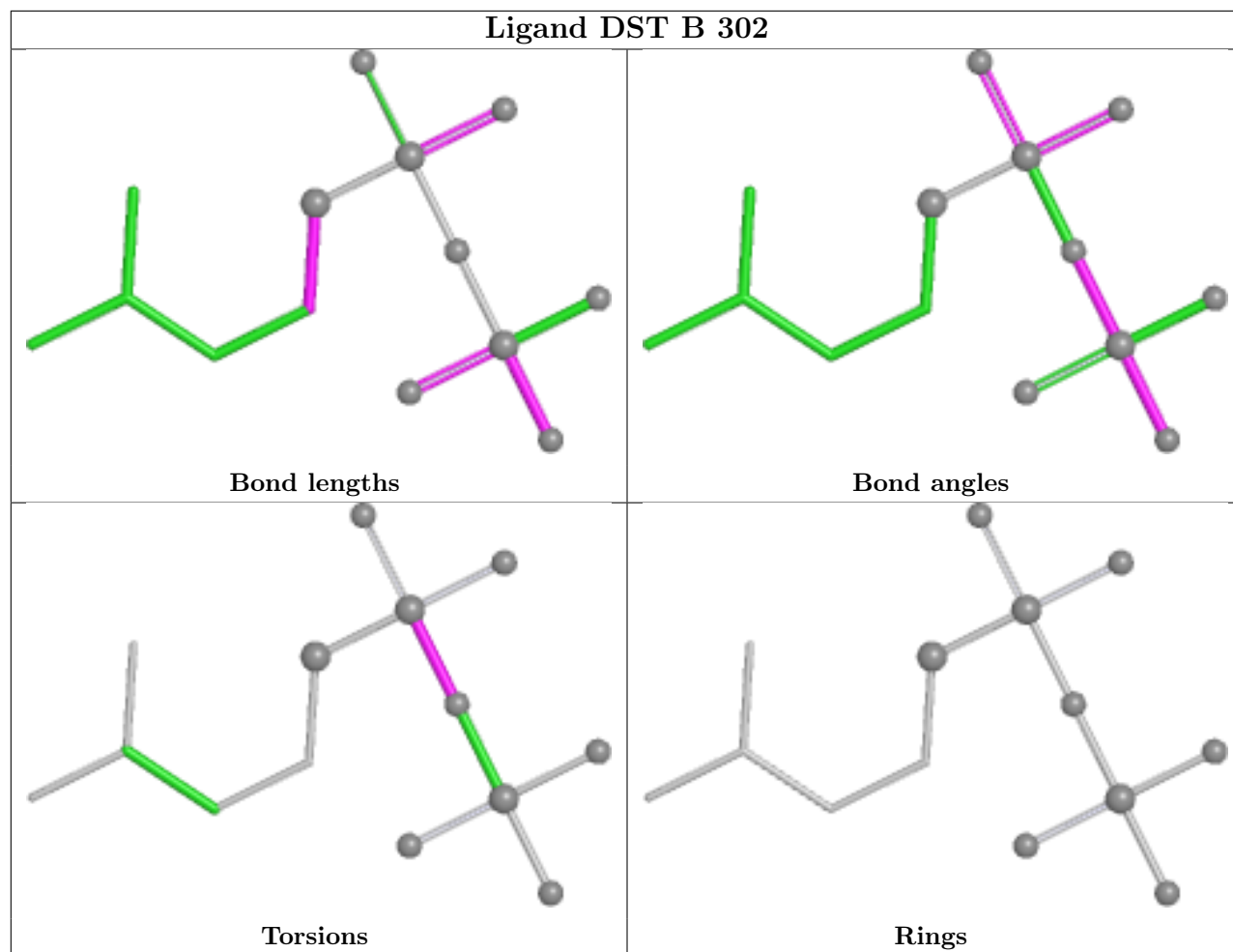


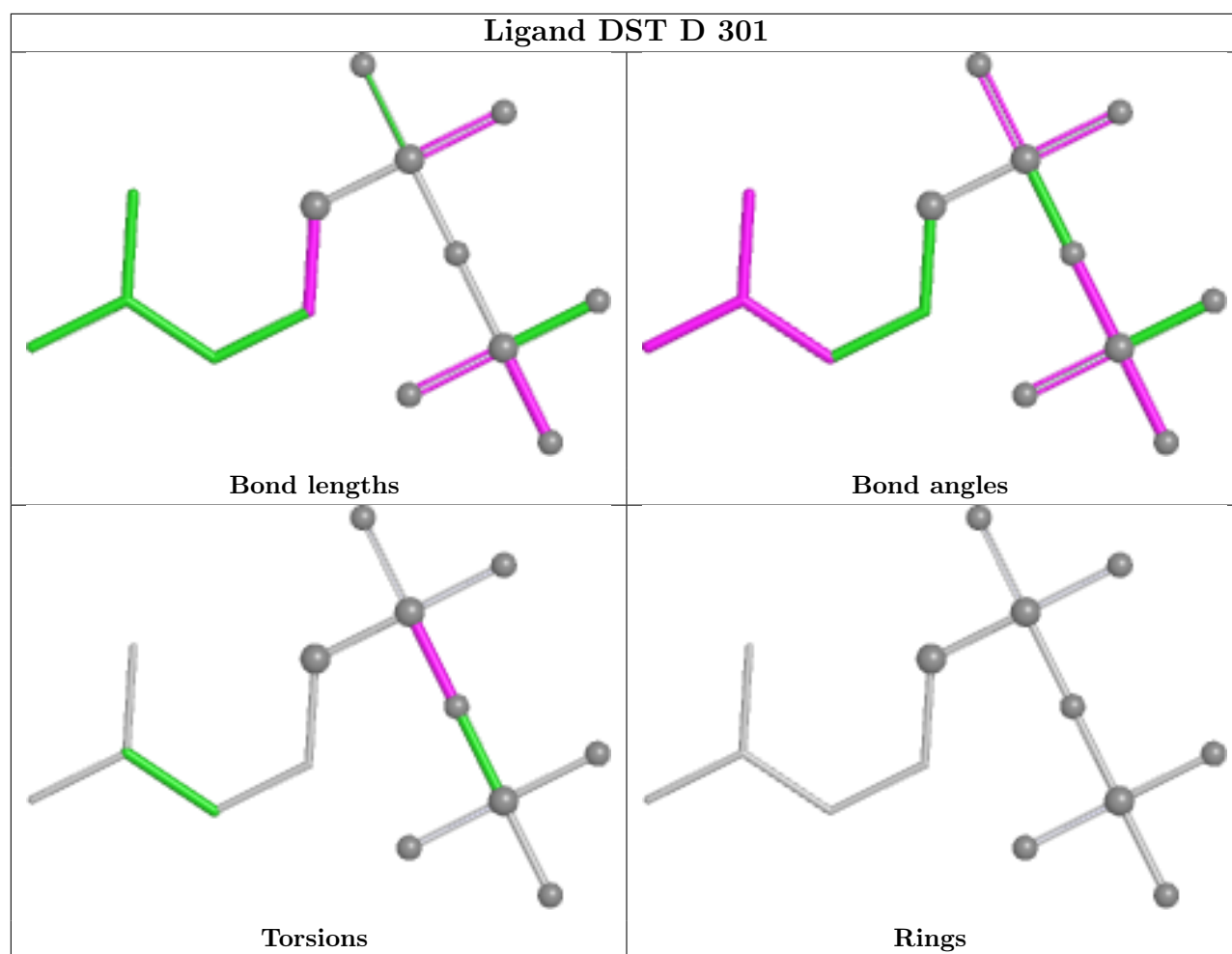


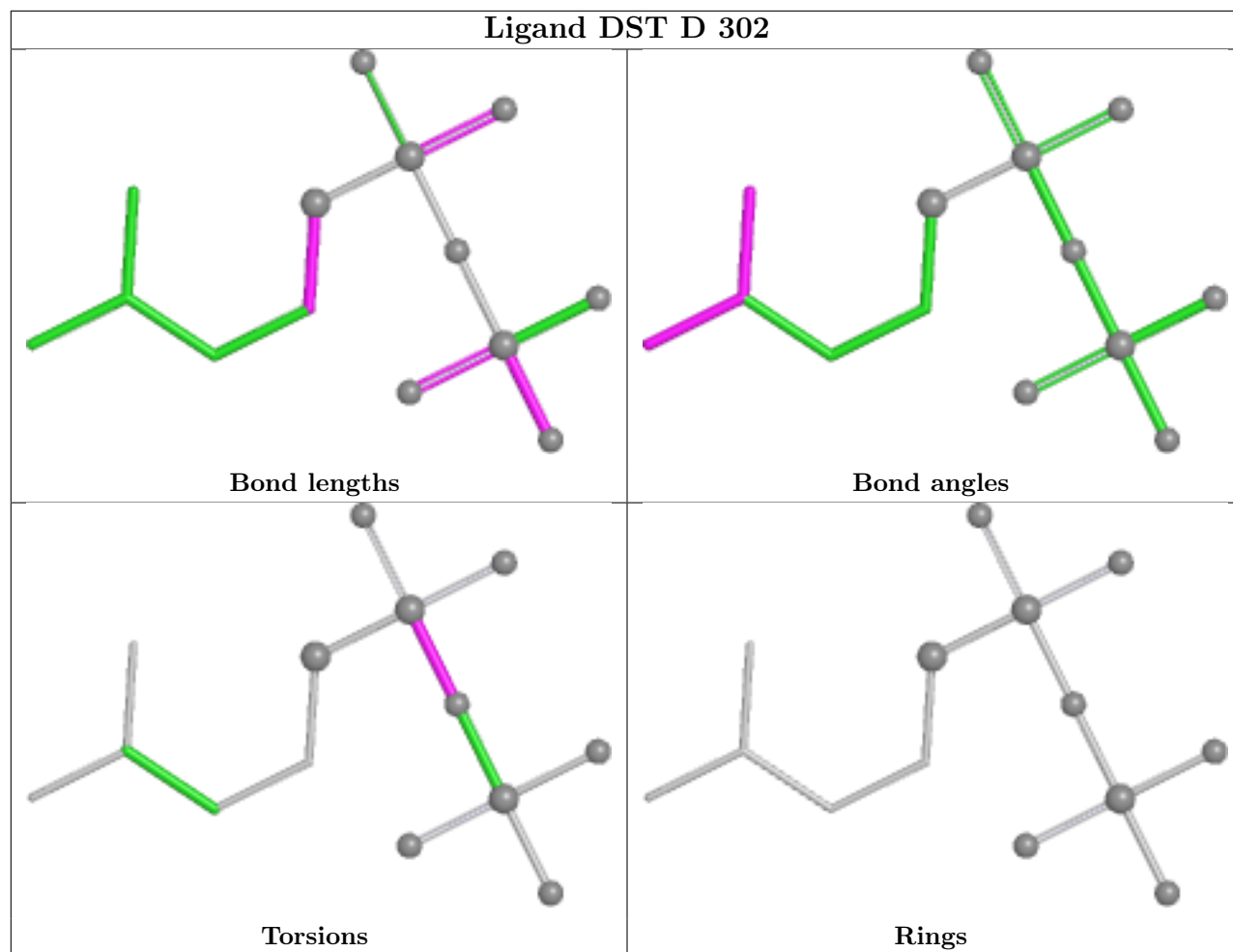


Ligand DST E 302









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	210/217 (96%)	-0.33	0 100 100	22, 31, 49, 83	0
1	B	208/217 (95%)	0.07	3 (1%) 73 75	29, 42, 60, 83	0
1	C	217/217 (100%)	-0.14	5 (2%) 61 63	23, 32, 58, 90	0
1	D	208/217 (95%)	0.04	2 (0%) 79 81	29, 40, 57, 77	0
1	E	208/217 (95%)	-0.04	1 (0%) 87 89	23, 37, 59, 78	0
1	F	208/217 (95%)	0.54	4 (1%) 66 68	33, 56, 77, 92	0
1	G	208/217 (95%)	0.01	2 (0%) 79 81	26, 39, 63, 89	0
1	H	216/217 (99%)	-0.07	9 (4%) 40 40	22, 35, 59, 85	0
All	All	1683/1736 (96%)	0.01	26 (1%) 72 73	22, 38, 67, 92	0

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	217	ALA	5.9
1	H	216	PRO	5.1
1	H	212	PHE	4.6
1	C	212	PHE	3.9
1	H	215	TYR	3.7
1	H	214	GLY	3.7
1	C	217	ALA	3.5
1	C	213	GLY	2.9
1	H	2	THR	2.7
1	F	20	VAL	2.6
1	H	213	GLY	2.6
1	F	27	ALA	2.5
1	C	206	ARG	2.4
1	H	211	ARG	2.3
1	B	81	SER	2.3
1	C	215	TYR	2.3

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Mol	Chain	Res	Type	RSRZ
1	H	210	ARG	2.3
1	G	114	ASP	2.3
1	D	180	TYR	2.3
1	F	115	PHE	2.2
1	E	1	MET	2.2
1	B	180	TYR	2.1
1	B	1	MET	2.1
1	G	209	ASP	2.1
1	D	81	SER	2.0
1	F	207	ALA	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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
3	MG	H	303	1/1	0.89	0.08	36,36,36,36	0
3	MG	B	303	1/1	0.90	0.14	43,43,43,43	0
2	DST	F	301	14/14	0.90	0.09	48,58,72,74	0
2	DST	C	301	14/14	0.91	0.10	35,41,46,52	0
2	DST	B	301	14/14	0.91	0.09	43,52,67,69	0
2	DST	A	301	14/14	0.92	0.09	33,41,60,63	0
2	DST	F	302	14/14	0.93	0.10	55,62,69,70	0
2	DST	G	301	14/14	0.93	0.09	44,52,76,78	0
2	DST	E	302	14/14	0.93	0.09	38,44,49,50	0
2	DST	D	301	14/14	0.93	0.09	37,46,58,59	0
2	DST	B	302	14/14	0.94	0.08	39,46,60,62	0
2	DST	D	302	14/14	0.95	0.09	29,47,58,65	0
2	DST	E	301	14/14	0.95	0.06	30,47,60,63	0

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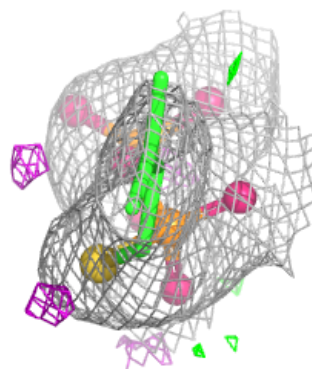
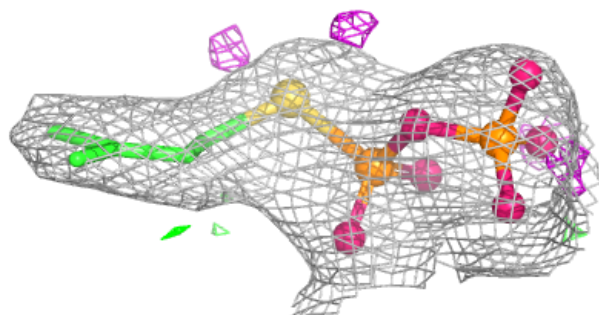
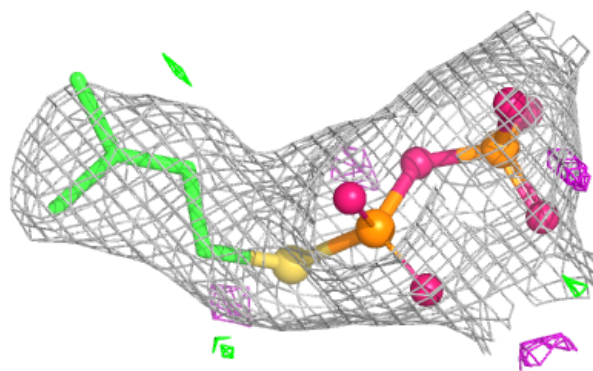
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	DST	A	302	14/14	0.96	0.08	32,43,56,65	0
2	DST	H	302	14/14	0.96	0.08	34,39,46,49	0
2	DST	C	302	14/14	0.97	0.09	27,39,58,58	0
3	MG	A	303	1/1	0.97	0.04	36,36,36,36	0
2	DST	G	302	14/14	0.97	0.07	34,42,50,60	0
3	MG	E	303	1/1	0.97	0.03	42,42,42,42	0
3	MG	F	303	1/1	0.97	0.04	54,54,54,54	0
2	DST	H	301	14/14	0.97	0.06	32,37,44,45	0
3	MG	G	303	1/1	0.98	0.04	48,48,48,48	0
3	MG	D	303	1/1	0.98	0.03	41,41,41,41	0
3	MG	C	303	1/1	0.99	0.04	31,31,31,31	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.

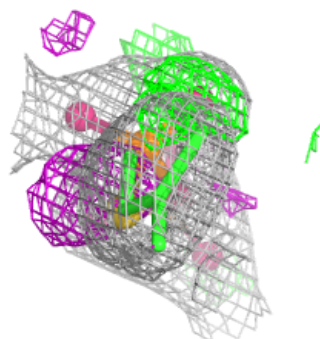
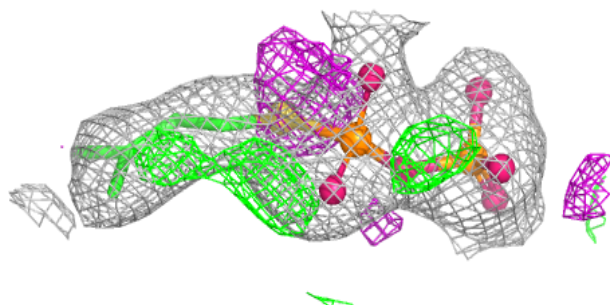
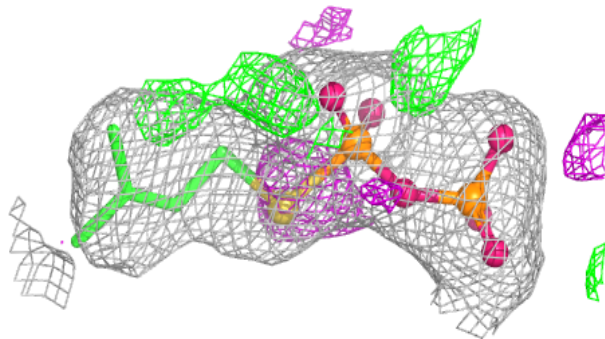
Electron density around DST F 301:

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

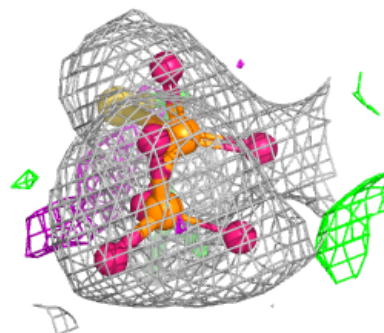
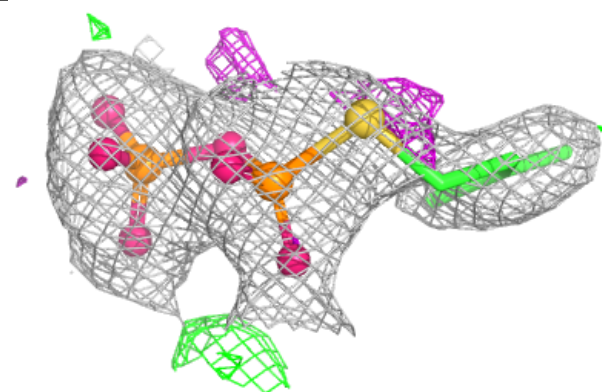
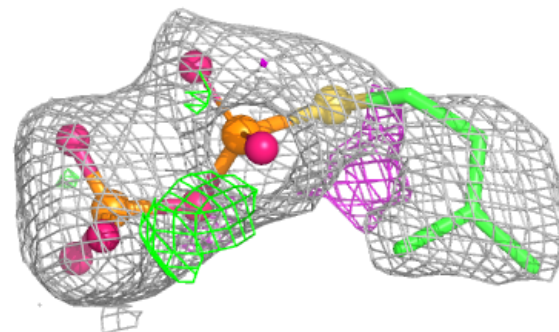


Electron density around DST C 301:

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

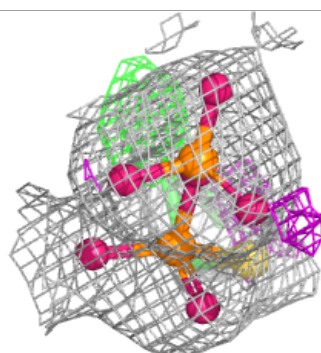
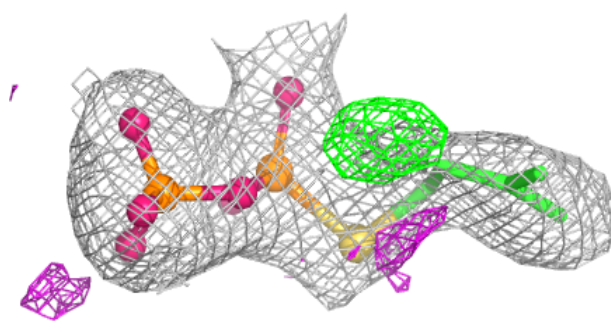
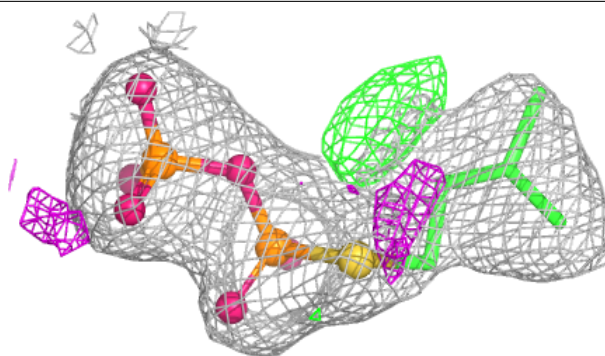
**Electron density around DST B 301:**

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

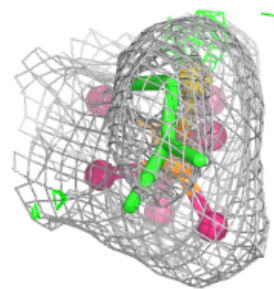
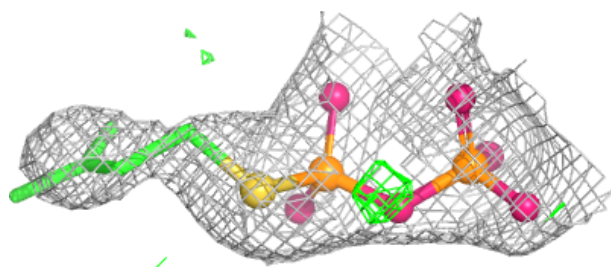
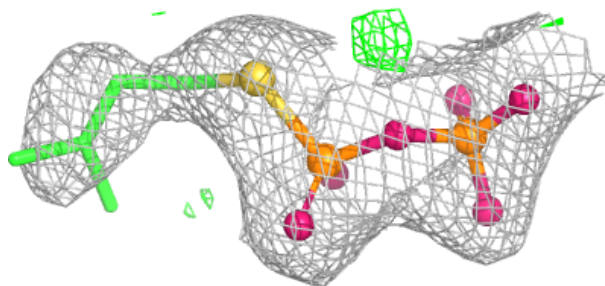


Electron density around DST A 301:

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and green (positive)

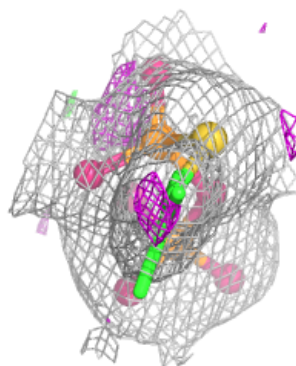
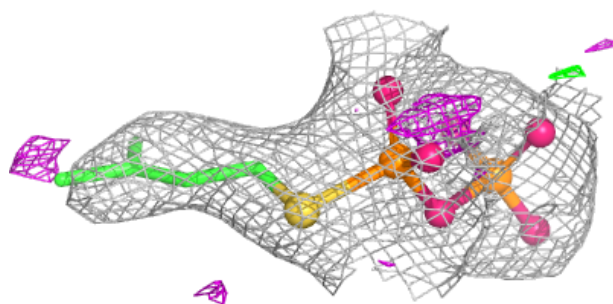
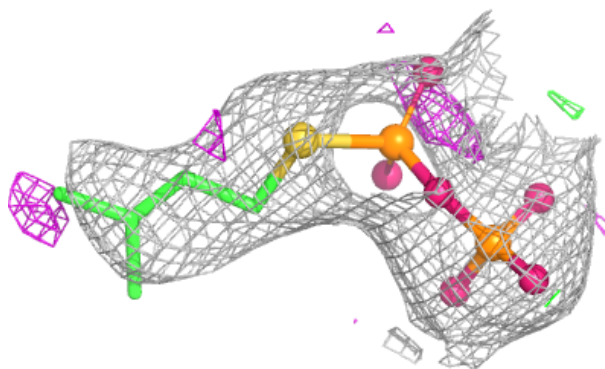
**Electron density around DST F 302:**

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

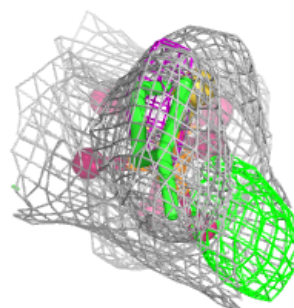
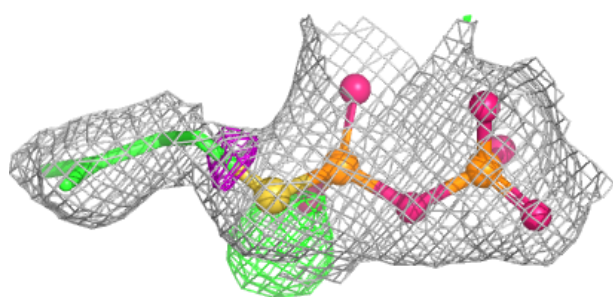
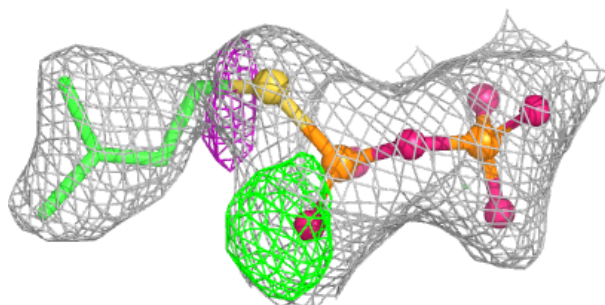


Electron density around DST G 301:

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and green (positive)

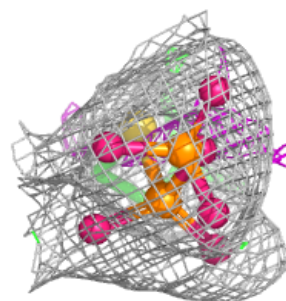
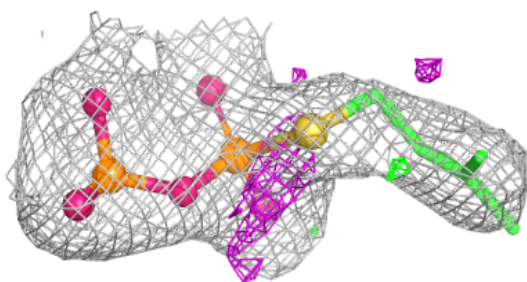
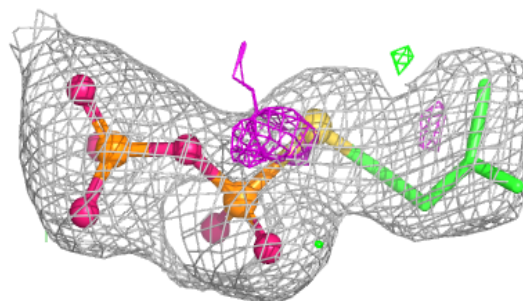
**Electron density around DST E 302:**

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
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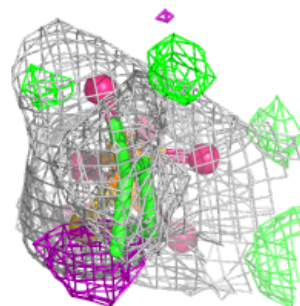
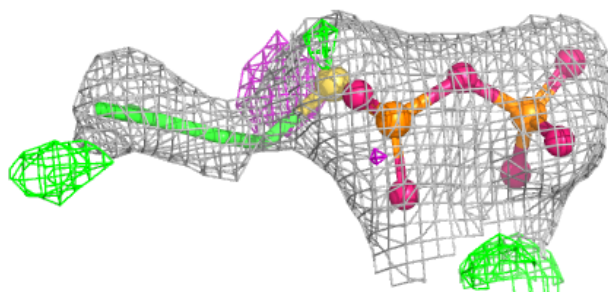
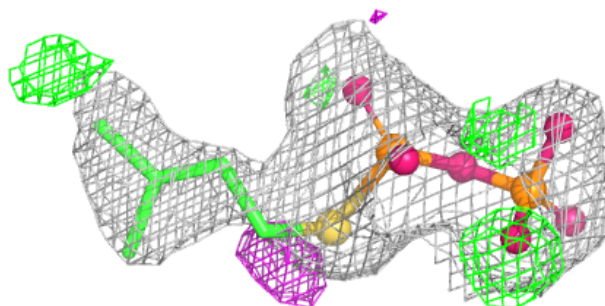


Electron density around DST D 301:

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and green (positive)

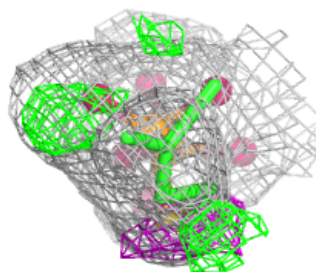
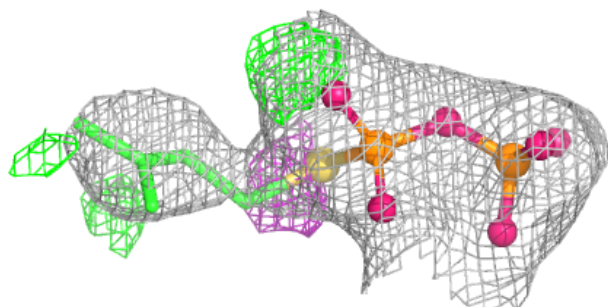
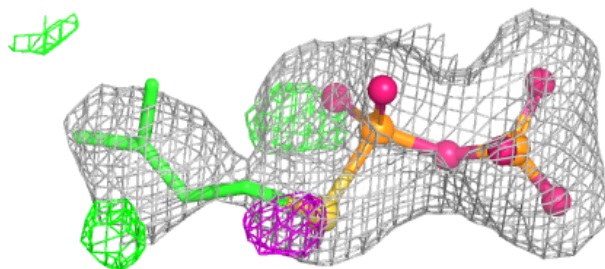
**Electron density around DST B 302:**

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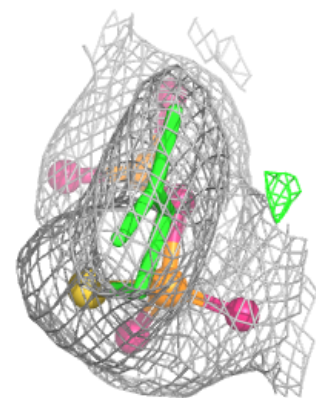
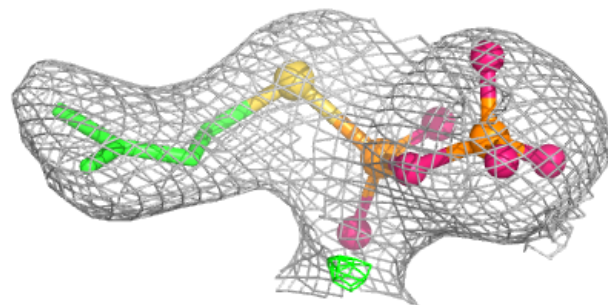
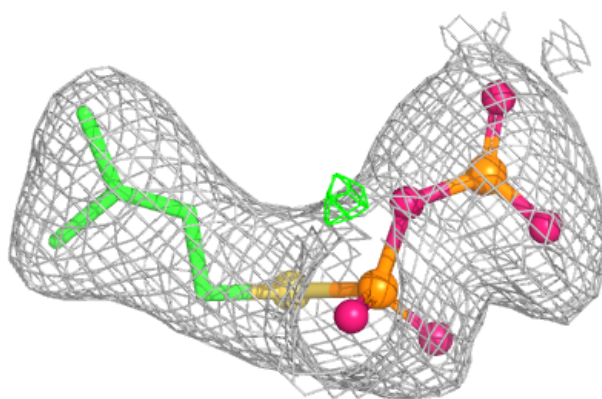


Electron density around DST D 302:

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

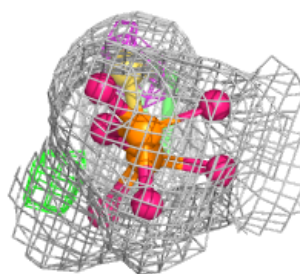
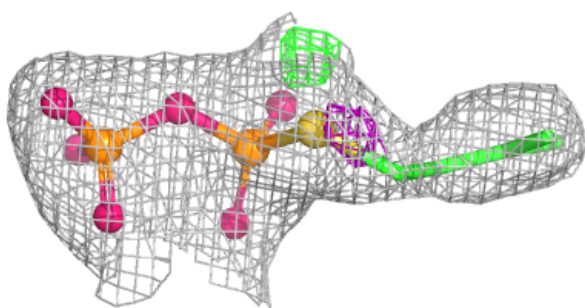
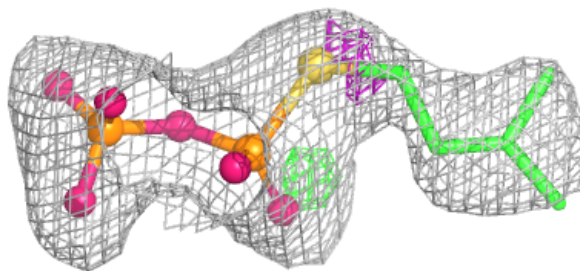
**Electron density around DST E 301:**

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

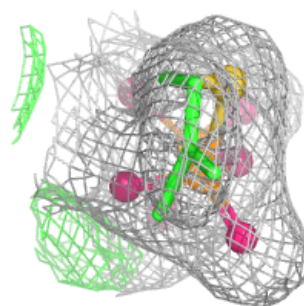
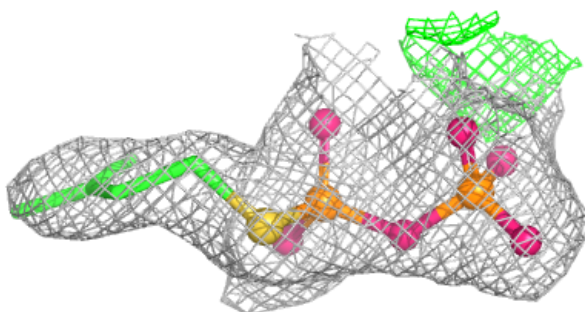
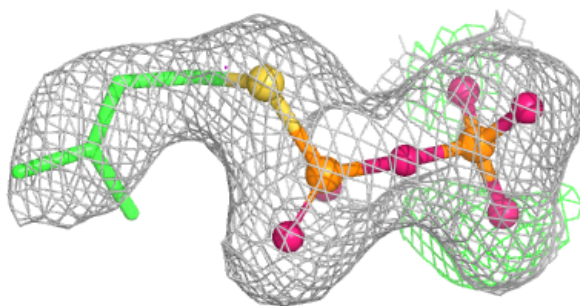


Electron density around DST A 302:

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

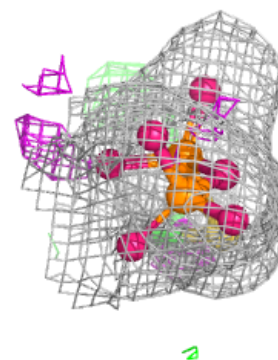
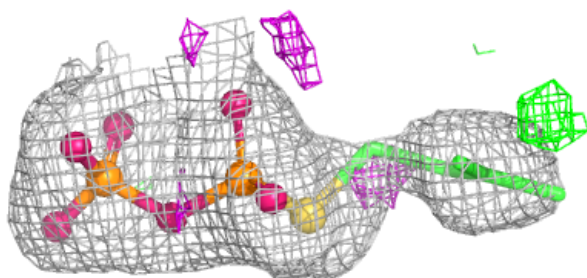
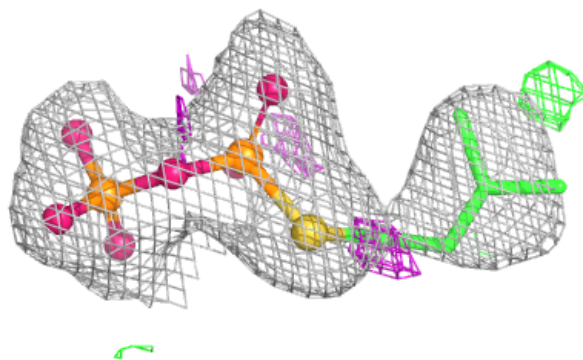
**Electron density around DST H 302:**

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

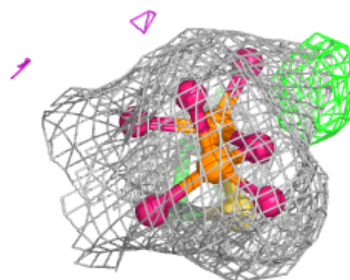
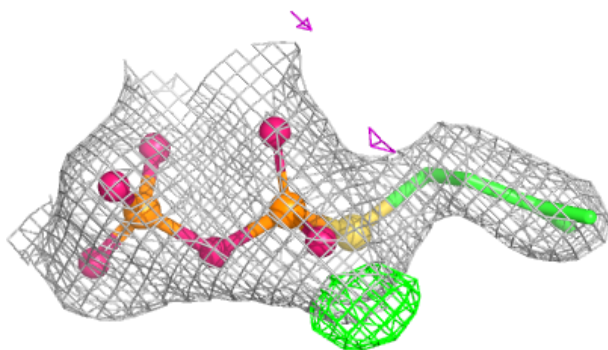
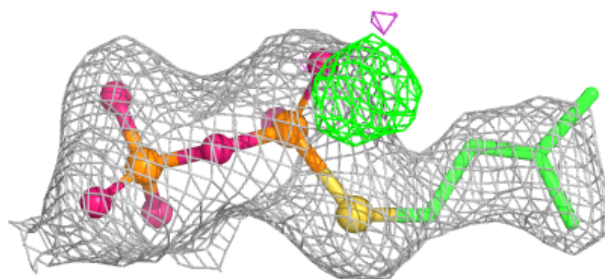


Electron density around DST C 302:

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

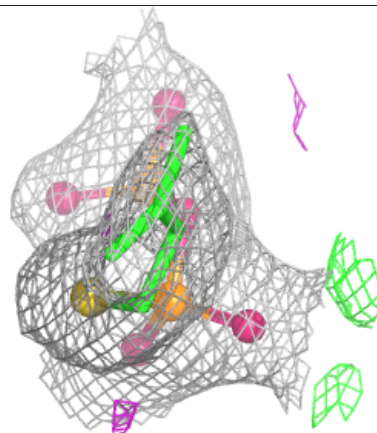
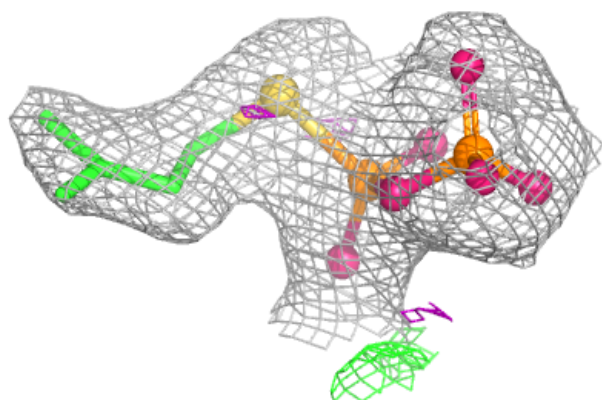
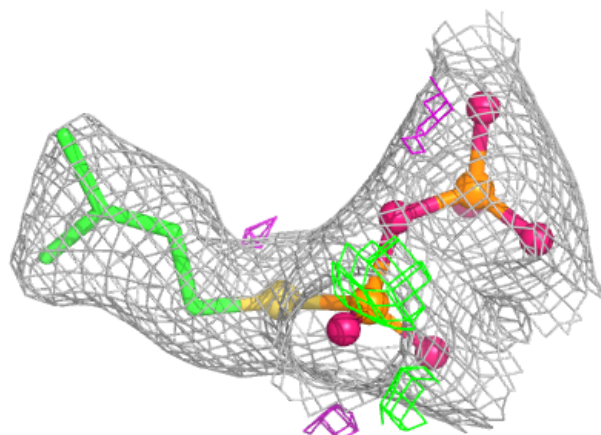
**Electron density around DST G 302:**

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



Electron density around DST H 301:

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



6.5 Other polymers [i](#)

There are no such residues in this entry.