



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 15, 2026 – 02:41 PM UTC

PDB ID : 8EJ0 / pdb_00008ej0
Title : Crystal structure of Fe-S cluster-dependent dehydratase from *Paralcaligenes ureilyticus* in complex with Mg
Authors : Bayaraa, T.; Lonhienne, T.; Guddat, L.W.
Deposited on : 2022-09-16
Resolution : 2.59 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

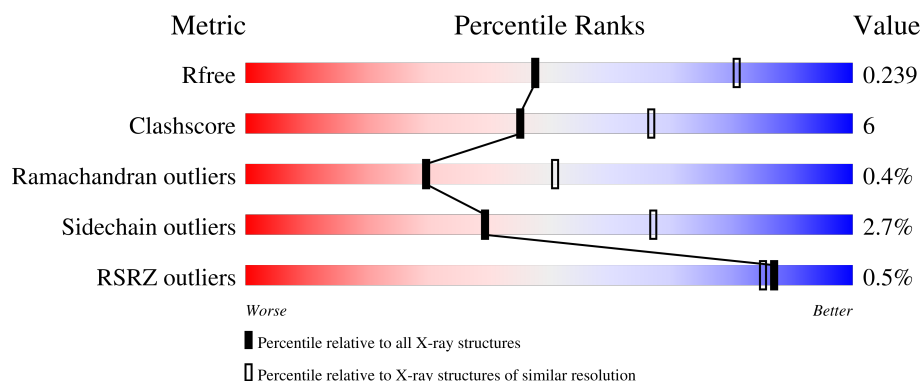
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.59 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	575	 84% 14% ..
1	C	575	 79% 18% ..
1	E	575	 86% 12% ..
1	G	575	 79% 18% ..

2 Entry composition [i](#)

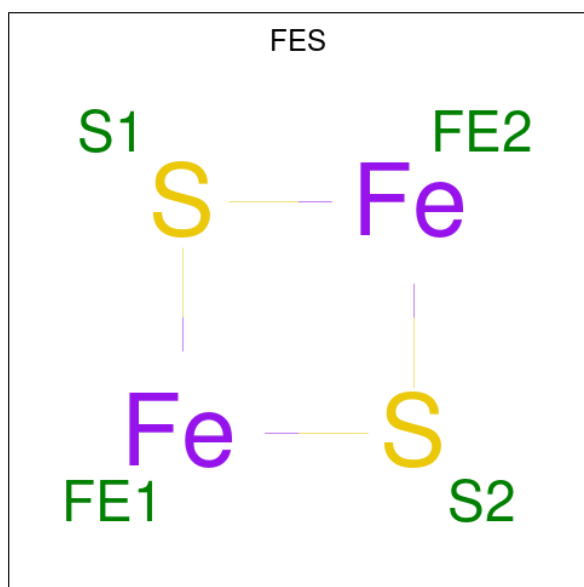
There are 6 unique types of molecules in this entry. The entry contains 17475 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Dihydroxyacid dehydratase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	568	Total	C	N	O	S	0	1	0
			4322	2699	784	801	38			
1	C	568	Total	C	N	O	S	0	3	0
			4344	2711	791	804	38			
1	E	568	Total	C	N	O	S	0	2	0
			4337	2707	788	804	38			
1	G	568	Total	C	N	O	S	0	1	0
			4294	2682	774	800	38			

- Molecule 2 is FE2/S2 (INORGANIC) CLUSTER (CCD ID: FES) (formula: Fe_2S_2) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	Fe	S	0	0
			4	2	2		
2	C	1	Total	Fe	S	0	0
			4	2	2		

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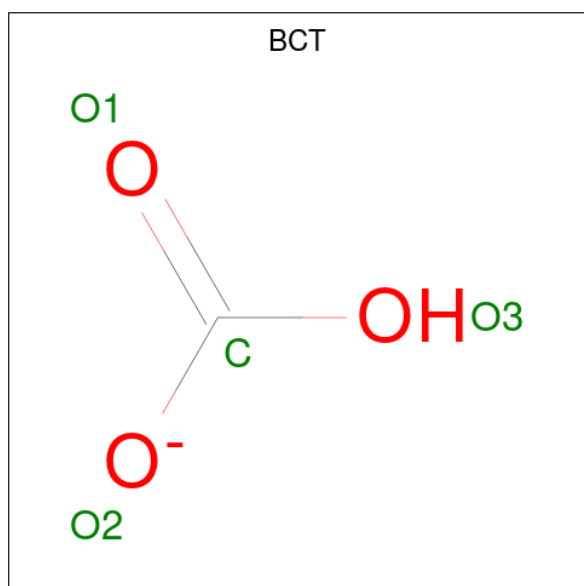
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	E	1	Total	Fe	S	0	0
			4	2	2		
2	G	1	Total	Fe	S	0	0
			4	2	2		

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		
3	C	1	Total	Mg	0	0
			1	1		
3	E	1	Total	Mg	0	0
			1	1		
3	G	1	Total	Mg	0	0
			1	1		

- Molecule 4 is BICARBONATE ION (CCD ID: BCT) (formula: CHO₃) (labeled as "Ligand of Interest" by depositor).



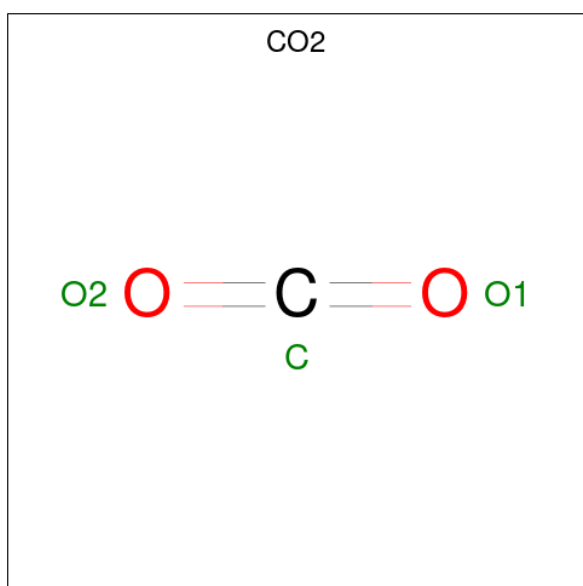
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			4	1	3		
4	A	1	Total	C	O	0	0
			4	1	3		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	C	1	Total	C	O	0	0
			4	1	3		
4	E	1	Total	C	O	0	0
			4	1	3		
4	E	1	Total	C	O	0	0
			4	1	3		
4	E	1	Total	C	O	0	0
			4	1	3		

- Molecule 5 is CARBON DIOXIDE (CCD ID: CO2) (formula: CO₂) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	G	1	Total	C	O	0	0
			3	1	2		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	31	Total	O	0	0
			31	31		
6	C	36	Total	O	0	0
			36	36		
6	E	42	Total	O	0	0
			42	42		
6	G	22	Total	O	0	0
			22	22		

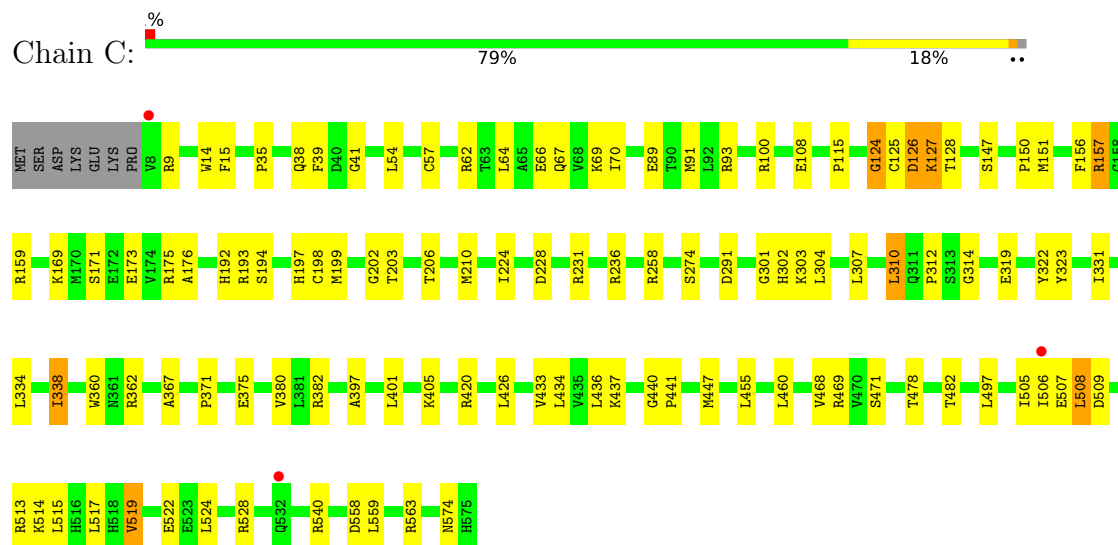
3 Residue-property plots

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.

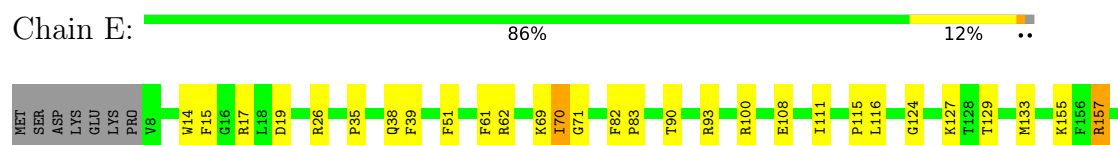
• Molecule 1: Dihydroxyacid dehydratase

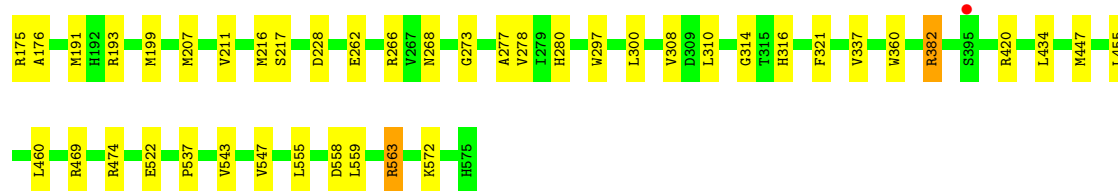


• Molecule 1: Dihydroxyacid dehydratase

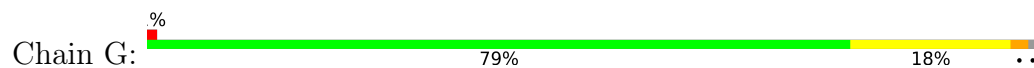


• Molecule 1: Dihydroxyacid dehydratase





• Molecule 1: Dihydroxyacid dehydratase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	69.91Å 178.07Å 92.86Å 90.00° 101.72° 90.00°	Depositor
Resolution (Å)	44.05 – 2.59 44.05 – 2.59	Depositor EDS
% Data completeness (in resolution range)	99.3 (44.05-2.59) 99.3 (44.05-2.59)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.09 (at 2.58Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.194 , 0.239 0.195 , 0.239	Depositor DCC
R_{free} test set	2000 reflections (2.91%)	wwPDB-VP
Wilson B-factor (Å ²)	45.4	Xtriage
Anisotropy	0.646	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 25.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	17475	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.90% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 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: CO2, KCX, BCT, MG, FES

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	0.49	2/4400 (0.0%)	0.38	5/5957 (0.1%)
1	C	0.52	3/4422 (0.1%)	0.39	2/5985 (0.0%)
1	E	0.49	2/4415 (0.0%)	0.33	2/5976 (0.0%)
1	G	0.17	0/4369	0.33	0/5922
All	All	0.44	7/17606 (0.0%)	0.36	9/23840 (0.0%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	69	LYS	C-N	-30.87	0.94	1.33
1	E	69	LYS	C-N	-28.72	0.98	1.33
1	C	124	GLY	C-N	23.29	1.64	1.33
1	C	69	LYS	C-N	-19.97	1.09	1.33
1	C	70	ILE	C-N	-13.49	1.14	1.33
1	E	70	ILE	C-N	-13.24	1.14	1.33
1	A	70	ILE	C-N	-6.87	1.24	1.33

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	124	GLY	CA-C-N	-9.46	103.75	120.97
1	C	124	GLY	C-N-CA	-9.46	103.75	120.97
1	A	70	ILE	CA-C-N	-7.29	111.86	120.03
1	A	70	ILE	C-N-CA	-7.29	111.86	120.03
1	A	69	LYS	CA-C-N	7.25	129.84	120.56
1	A	69	LYS	C-N-CA	7.25	129.84	120.56
1	A	69	LYS	O-C-N	-5.91	115.30	122.22
1	E	70	ILE	CA-C-N	-5.67	113.68	120.03
1	E	70	ILE	C-N-CA	-5.67	113.68	120.03

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	4322	0	4284	53	0
1	C	4344	0	4306	69	0
1	E	4337	0	4300	47	0
1	G	4294	0	4219	64	0
2	A	4	0	0	0	0
2	C	4	0	0	0	0
2	E	4	0	0	0	0
2	G	4	0	0	0	0
3	A	1	0	0	0	0
3	C	1	0	0	0	0
3	E	1	0	0	0	0
3	G	1	0	0	0	0
4	A	8	0	1	0	0
4	C	4	0	0	0	0
4	E	12	0	1	0	0
5	G	3	0	0	0	0
6	A	31	0	0	0	0
6	C	36	0	0	0	0
6	E	42	0	0	0	0
6	G	22	0	0	2	0
All	All	17475	0	17111	203	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 (203) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:397:ALA:HB1	1:C:482:THR:HG22	1.67	0.76
1:C:441:PRO:HG3	1:C:447:MET:HE2	1.69	0.74
1:C:206:THR:HG21	1:C:274:SER:H	1.52	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:175:ARG:NH2	1:C:455:LEU:O	2.22	0.73
1:A:163:SER:OG	1:A:199:MET:SD	2.47	0.71
1:C:9:ARG:NH2	1:C:41:GLY:O	2.26	0.69
1:E:108:GLU:OE1	1:G:540:ARG:NH2	2.27	0.68
1:E:17:ARG:O	1:E:26:ARG:NH2	2.27	0.67
1:A:37:ASP:O	1:A:42:ARG:NH1	2.29	0.64
1:A:108:GLU:OE1	1:C:540:ARG:NH2	2.31	0.64
1:C:362:ARG:HH21	1:C:367:ALA:HB2	1.63	0.64
1:G:397:ALA:HB1	1:G:482:THR:HG22	1.81	0.63
1:A:225:PRO:HG2	1:A:228:ASP:HB2	1.80	0.63
1:A:447:MET:HG2	1:A:474:ARG:HG3	1.80	0.62
1:C:331:ILE:HG23	1:C:338:ILE:HD13	1.81	0.62
1:E:14:TRP:CD2	1:E:115:PRO:HB3	2.34	0.62
1:E:266:ARG:NH1	1:E:337:VAL:O	2.33	0.61
1:G:57:CYS:HB3	1:G:197:HIS:HA	1.82	0.61
1:G:111:ILE:HA	1:G:116:LEU:HD12	1.82	0.61
1:E:310:LEU:HD11	1:E:321:PHE:HB2	1.83	0.60
1:G:225:PRO:HG2	1:G:228:ASP:HB2	1.82	0.60
1:G:32:ARG:NH1	6:G:703:HOH:O	2.34	0.60
1:E:15:PHE:HB3	1:E:39:PHE:HB3	1.82	0.60
1:G:14:TRP:CD2	1:G:115:PRO:HB3	2.36	0.60
1:A:510:VAL:O	1:A:513:ARG:NH1	2.35	0.59
1:E:175:ARG:NH2	1:E:455:LEU:O	2.35	0.59
1:C:258:ARG:NH1	1:C:291:ASP:O	2.36	0.59
1:G:121:LEU:HD22	1:G:132:LEU:HB3	1.84	0.59
1:A:258:ARG:NH2	1:A:291:ASP:O	2.35	0.59
1:A:176:ALA:HB2	1:A:460:LEU:HB3	1.86	0.58
1:C:401:LEU:HB2	1:C:482:THR:HG21	1.86	0.58
1:C:57:CYS:HB3	1:C:197:HIS:HA	1.86	0.57
1:E:572:LYS:NZ	1:G:414:SER:OG	2.27	0.57
1:A:154:GLY:H	1:A:199:MET:HE3	1.68	0.57
1:A:15:PHE:HB3	1:A:39:PHE:HB3	1.86	0.57
1:A:452:ASN:HB3	1:A:475:MET:HE2	1.86	0.57
1:E:100:ARG:HH11	1:E:280:HIS:HD1	1.53	0.57
1:A:75:SER:HB2	1:A:243:VAL:HG11	1.86	0.56
1:C:405:LYS:HE2	1:C:507:GLU:HG3	1.87	0.56
1:A:82:PHE:HB3	1:C:54:LEU:HD12	1.87	0.56
1:E:129:THR:HG21	1:E:207:MET:HB3	1.87	0.56
1:E:308:VAL:HG12	1:E:310:LEU:HG	1.86	0.56
1:G:517:LEU:O	1:G:519:VAL:N	2.38	0.56
1:C:14:TRP:CG	1:C:115:PRO:HB3	2.41	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:133:MET:HE1	1:E:216:MET:HG3	1.88	0.55
1:E:555:LEU:O	1:E:563:ARG:NH2	2.39	0.55
1:C:517:LEU:HG	1:C:519:VAL:HG13	1.88	0.55
1:C:14:TRP:CD2	1:C:115:PRO:HB3	2.41	0.55
1:C:67:GLN:HG3	1:C:236:ARG:HB2	1.89	0.54
1:C:126:ASP:OD2	1:C:203:THR:HB	2.07	0.54
1:C:434:LEU:HB2	1:C:469:ARG:HG2	1.90	0.54
1:G:298:ASP:OD1	1:G:513:ARG:NE	2.36	0.54
1:E:310:LEU:O	1:E:314:GLY:N	2.36	0.54
1:C:440:GLY:HA2	1:C:447:MET:HG3	1.90	0.54
1:A:175:ARG:NH2	1:A:422:ASP:OD2	2.40	0.54
1:A:308:VAL:HG12	1:A:310:LEU:HB2	1.89	0.53
1:A:360:TRP:CH2	1:G:157:ARG:HD2	2.43	0.53
1:E:19:ASP:OD2	1:E:19:ASP:N	2.41	0.53
1:C:433:VAL:HG21	1:C:506:ILE:HB	1.90	0.53
1:C:15:PHE:HB3	1:C:39:PHE:HB3	1.91	0.53
1:A:133:MET:HE1	1:A:216:MET:HG3	1.90	0.53
1:A:170:MET:HB3	1:A:179:MET:HE1	1.90	0.52
1:G:176:ALA:HB2	1:G:460:LEU:HB3	1.92	0.52
1:G:310:LEU:O	1:G:314:GLY:N	2.40	0.52
1:A:14:TRP:CD2	1:A:115:PRO:HB3	2.45	0.52
1:G:563:ARG:NH1	6:G:705:HOH:O	2.39	0.52
1:A:155[B]:LYS:HD2	1:G:323:TYR:HB3	1.91	0.52
1:G:434:LEU:HB2	1:G:469:ARG:HG2	1.90	0.51
1:A:155[B]:LYS:HE2	1:A:158:GLY:C	2.35	0.51
1:A:10:ARG:HD3	1:A:116:LEU:O	2.11	0.51
1:A:302:HIS:CD2	1:A:513:ARG:HH22	2.29	0.51
1:A:302:HIS:HA	1:A:377:GLY:HA3	1.92	0.51
1:C:157:ARG:HD2	1:E:360:TRP:CH2	2.46	0.51
1:A:331:ILE:HD13	1:A:350:LEU:HD11	1.93	0.51
1:E:447:MET:HG2	1:E:474:ARG:HG3	1.93	0.51
1:G:14:TRP:CG	1:G:115:PRO:HB3	2.46	0.51
1:G:424:GLU:OE2	1:G:461:ARG:NH2	2.43	0.51
1:A:15:PHE:HZ	1:C:91:MET:HE1	1.76	0.50
1:A:540:ARG:NH2	1:C:108:GLU:OE1	2.42	0.50
1:E:93:ARG:HD3	1:G:14:TRP:CE2	2.46	0.50
1:E:572:LYS:HZ1	1:G:414:SER:HG	1.57	0.50
1:G:170:MET:HE2	1:G:184:PHE:HD1	1.76	0.50
1:A:540:ARG:HH22	1:C:108:GLU:CD	2.19	0.50
1:G:344:THR:OG1	1:G:346:ASN:OD1	2.29	0.50
1:A:157:ARG:HD2	1:G:360:TRP:CH2	2.47	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:360:TRP:CH2	1:E:157:ARG:HD2	2.47	0.50
1:C:380:VAL:O	1:C:513:ARG:HG3	2.11	0.49
1:G:128:THR:O	1:G:132:LEU:HG	2.11	0.49
1:C:524:LEU:O	1:C:528:ARG:HG3	2.12	0.49
1:C:558:ASP:OD1	1:C:559:LEU:N	2.46	0.48
1:G:409:VAL:HG21	1:G:428:VAL:CG1	2.43	0.48
1:A:14:TRP:CE2	1:C:93:ARG:HD3	2.48	0.48
1:A:93:ARG:HD3	1:C:14:TRP:CE2	2.48	0.48
1:C:100:ARG:HB2	1:C:127:KCX:HG2	1.95	0.48
1:C:35:PRO:HG2	1:C:38:GLN:HG2	1.95	0.48
1:G:508:LEU:HD23	1:G:515:LEU:HD13	1.95	0.48
1:A:162:GLY:N	1:A:166:ASP:OD2	2.47	0.47
1:G:175:ARG:NH2	1:G:422:ASP:OD1	2.46	0.47
1:C:303:LYS:HA	1:C:375:GLU:HG2	1.97	0.47
1:G:376:ALA:HB1	1:G:395:SER:HB3	1.97	0.47
1:G:378:ILE:H	1:G:378:ILE:HD12	1.80	0.47
1:E:207:MET:HA	1:E:207:MET:HE2	1.97	0.47
1:E:262:GLU:HG2	1:E:300:LEU:HD11	1.97	0.47
1:C:171:SER:O	1:C:175:ARG:HG3	2.15	0.46
1:A:310:LEU:O	1:A:314:GLY:N	2.39	0.46
1:C:169:LYS:HE2	1:C:173:GLU:OE2	2.16	0.46
1:E:211:VAL:HG12	1:E:217:SER:HB3	1.96	0.46
1:G:457:PRO:O	1:G:461:ARG:HG3	2.16	0.46
1:E:90:THR:HG21	1:G:28:TRP:CD2	2.50	0.46
1:E:310:LEU:HD23	1:E:316:HIS:HB2	1.98	0.46
1:C:206:THR:HG21	1:C:274:SER:N	2.25	0.46
1:G:558:ASP:OD1	1:G:559:LEU:N	2.46	0.46
1:A:380:VAL:O	1:A:513:ARG:HD2	2.15	0.46
1:A:470:VAL:HA	1:A:484:VAL:HG22	1.98	0.46
1:E:14:TRP:CE2	1:G:93:ARG:HD3	2.51	0.46
1:G:89:GLU:HG3	1:G:126:ASP:HB2	1.97	0.46
1:C:436:LEU:HB3	1:C:471:SER:HB3	1.98	0.46
1:G:199:MET:HE3	1:G:199:MET:HB2	1.75	0.46
1:A:10:ARG:HD2	1:A:141:LEU:HD21	1.98	0.45
1:A:28:TRP:NE1	1:C:198:CYS:SG	2.85	0.45
1:G:163:SER:H	1:G:199:MET:HE2	1.82	0.45
1:G:303:LYS:HA	1:G:375:GLU:HG2	1.98	0.45
1:C:319:GLU:HG2	1:C:323:TYR:CE2	2.52	0.45
1:C:206:THR:O	1:C:210:MET:N	2.43	0.45
1:C:508:LEU:HD23	1:C:515:LEU:HD13	1.99	0.45
1:G:100:ARG:HB2	1:G:127:KCX:HG3	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:129:THR:HG21	1:A:207:MET:HB3	1.99	0.45
1:C:125:CYS:O	1:C:128:THR:OG1	2.29	0.45
1:C:202:GLY:O	1:C:206:THR:HG23	2.16	0.45
1:C:301:GLY:HA2	1:C:304:LEU:HD12	1.99	0.45
1:G:44:VAL:HG22	1:G:78:PHE:HD2	1.81	0.45
1:E:14:TRP:CG	1:E:115:PRO:HB3	2.51	0.44
1:C:310:LEU:O	1:C:314:GLY:N	2.46	0.44
1:E:522:GLU:OE1	1:E:522:GLU:N	2.48	0.44
1:C:147:SER:O	1:C:231:ARG:NH1	2.50	0.44
1:C:519:VAL:HG22	1:C:524:LEU:HG	1.99	0.44
1:E:82:PHE:HB3	1:G:54:LEU:HD12	2.00	0.44
1:C:176:ALA:HB2	1:C:460:LEU:HB3	1.98	0.44
1:A:193:ARG:NH1	1:G:323:TYR:OH	2.51	0.44
1:G:51:PHE:HB2	1:G:61:PHE:HB2	2.00	0.44
1:A:108:GLU:CD	1:C:540:ARG:HH22	2.24	0.43
1:C:307:LEU:HD23	1:C:371:PRO:HB3	1.99	0.43
1:E:176:ALA:HB2	1:E:460:LEU:HB3	1.99	0.43
1:G:458:LYS:HG2	1:G:459:ILE:HD12	2.00	0.43
1:C:509:ASP:HB3	1:C:514:LYS:HB3	2.00	0.43
1:E:191:MET:O	1:E:199:MET:HE1	2.18	0.43
1:E:35:PRO:HG2	1:E:38:GLN:HG2	2.00	0.43
1:G:437:LYS:HD2	1:G:498:ALA:HA	2.01	0.43
1:G:298:ASP:HB2	1:G:380:VAL:HB	2.00	0.43
1:A:93:ARG:HD3	1:C:14:TRP:NE1	2.34	0.43
1:C:62[B]:ARG:O	1:C:66[B]:GLU:HG2	2.19	0.43
1:C:323:TYR:CE1	1:E:155:LYS:HG3	2.53	0.43
1:C:433:VAL:HA	1:C:468:VAL:HG23	2.01	0.43
1:G:499:LEU:HD22	1:G:527:ARG:HB2	2.00	0.43
1:E:382:ARG:HE	1:E:382:ARG:HB3	1.57	0.43
1:G:15:PHE:HB3	1:G:39:PHE:HB3	2.00	0.43
1:C:228:ASP:OD2	1:E:193:ARG:HD3	2.18	0.43
1:G:268:ASN:OD1	1:G:273:GLY:HA3	2.19	0.43
1:A:14:TRP:CG	1:A:115:PRO:HB3	2.54	0.42
1:A:224:ILE:HG12	1:A:322:TYR:CE1	2.54	0.42
1:G:522:GLU:H	1:G:522:GLU:HG2	1.50	0.42
1:A:64:LEU:HD21	1:A:231:ARG:HG3	2.01	0.42
1:E:558:ASP:OD1	1:E:559:LEU:N	2.51	0.42
1:G:352:ASP:HA	1:G:355:LYS:HE2	2.02	0.42
1:C:420:ARG:HG2	1:C:426:LEU:HD13	2.01	0.42
1:G:51:PHE:CE2	1:G:62:ARG:HD2	2.55	0.42
1:C:312:PRO:HG3	1:C:478:THR:HA	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:51:PHE:HB2	1:E:61:PHE:HB2	2.01	0.42
1:G:411:PHE:CD1	1:G:417:MET:HA	2.54	0.42
1:A:94:PRO:O	1:C:574:ASN:HA	2.19	0.42
1:E:83:PRO:HB2	1:G:83:PRO:HB2	2.02	0.42
1:A:213:ALA:HB2	1:A:267:VAL:HG21	2.02	0.42
1:E:108:GLU:CD	1:G:540:ARG:HH22	2.27	0.42
1:C:156:PHE:O	1:C:159:ARG:HG2	2.19	0.41
1:A:211:VAL:HG12	1:A:217:SER:HB3	2.02	0.41
1:A:28:TRP:CD2	1:C:192:HIS:CE1	3.08	0.41
1:E:268:ASN:OD1	1:E:273:GLY:HA3	2.20	0.41
1:E:278:VAL:HA	1:E:297:TRP:CZ2	2.55	0.41
1:C:302:HIS:CD2	1:C:303:LYS:HG2	2.55	0.41
1:C:150:PRO:HD2	1:C:224:ILE:O	2.20	0.41
1:E:207:MET:HE3	1:E:277:ALA:HA	2.02	0.41
1:E:537:PRO:HG2	1:E:547:VAL:HG21	2.01	0.41
1:G:150:PRO:HD2	1:G:224:ILE:O	2.20	0.41
1:C:437:LYS:HD3	1:C:497:LEU:O	2.21	0.41
1:G:426:LEU:HG	1:G:428:VAL:HG22	2.01	0.41
1:A:151:MET:HG3	1:A:196:GLY:C	2.46	0.41
1:G:199:MET:HE1	1:G:317:LEU:HD13	2.01	0.41
1:A:540:ARG:HG2	1:A:541:GLY:N	2.36	0.41
1:C:224:ILE:HG12	1:C:322:TYR:CE1	2.56	0.41
1:G:151:MET:HG3	1:G:196:GLY:C	2.46	0.41
1:G:333:GLU:OE1	1:G:362:ARG:NH2	2.54	0.41
1:E:70:ILE:O	1:E:71:GLY:C	2.59	0.41
1:E:111:ILE:HA	1:E:116:LEU:HD12	2.03	0.41
1:E:434:LEU:HB2	1:E:469:ARG:HG2	2.03	0.41
1:G:142:PRO:HD3	1:G:251:VAL:HG12	2.03	0.41
1:A:424:GLU:OE2	1:A:461:ARG:NH2	2.33	0.40
1:A:475:MET:HE1	1:A:483:VAL:HB	2.03	0.40
1:G:447:MET:HG2	1:G:474:ARG:HG3	2.02	0.40
1:G:539:ALA:C	1:G:544:LYS:HD2	2.47	0.40
1:A:35:PRO:HG2	1:A:38:GLN:HG2	2.04	0.40
1:C:151:MET:HE3	1:C:194:SER:O	2.21	0.40
1:C:193:ARG:HD2	1:E:228:ASP:OD2	2.22	0.40
1:G:129:THR:HG21	1:G:207:MET:HB3	2.03	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	566/575 (98%)	552 (98%)	12 (2%)	2 (0%)	30	51
1	C	568/575 (99%)	552 (97%)	14 (2%)	2 (0%)	30	51
1	E	567/575 (99%)	549 (97%)	17 (3%)	1 (0%)	43	66
1	G	566/575 (98%)	538 (95%)	25 (4%)	3 (0%)	24	46
All	All	2267/2300 (99%)	2191 (97%)	68 (3%)	8 (0%)	30	51

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	126	ASP
1	G	126	ASP
1	G	518	HIS
1	A	124	GLY
1	E	124	GLY
1	C	124	GLY
1	C	126	ASP
1	G	124	GLY

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	454/461 (98%)	444 (98%)	10 (2%)	45	72
1	C	456/461 (99%)	442 (97%)	14 (3%)	35	63

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	E	456/461 (99%)	449 (98%)	7 (2%)	57 80
1	G	447/461 (97%)	426 (95%)	21 (5%)	23 48
All	All	1813/1844 (98%)	1761 (97%)	52 (3%)	39 65

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	ARG
1	A	10	ARG
1	A	11	SER
1	A	62	ARG
1	A	66	GLU
1	A	155[A]	LYS
1	A	155[B]	LYS
1	A	157	ARG
1	A	189	SER
1	A	374	THR
1	C	64	LEU
1	C	89	GLU
1	C	157	ARG
1	C	199	MET
1	C	310	LEU
1	C	334	LEU
1	C	338	ILE
1	C	382[A]	ARG
1	C	382[B]	ARG
1	C	505	ILE
1	C	508	LEU
1	C	519	VAL
1	C	522	GLU
1	C	563	ARG
1	E	62	ARG
1	E	157	ARG
1	E	382	ARG
1	E	420[A]	ARG
1	E	420[B]	ARG
1	E	543	VAL
1	E	563	ARG
1	G	8	VAL
1	G	62	ARG
1	G	66[A]	GLU

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Mol	Chain	Res	Type
1	G	66[B]	GLU
1	G	89	GLU
1	G	157	ARG
1	G	159	ARG
1	G	181	GLN
1	G	199	MET
1	G	278	VAL
1	G	370	GLU
1	G	428	VAL
1	G	431	ASN
1	G	458	LYS
1	G	505	ILE
1	G	507	GLU
1	G	510	VAL
1	G	520	SER
1	G	522	GLU
1	G	523	GLU
1	G	527	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	244	GLN
1	A	316	HIS
1	C	302	HIS
1	C	404	HIS
1	E	316	HIS
1	E	431	ASN
1	G	38	GLN
1	G	67	GLN
1	G	311	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 non-standard protein/DNA/RNA residues 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$
1	KCX	E	127	3,1	10,11,12	2.08	2 (20%)	6,12,14	1.32	1 (16%)
1	KCX	A	127	3,1	10,11,12	2.05	2 (20%)	6,12,14	1.45	1 (16%)
1	KCX	G	127	3,1	10,11,12	2.03	2 (20%)	6,12,14	1.65	1 (16%)
1	KCX	C	127	3,1	10,11,12	2.00	2 (20%)	6,12,14	1.46	1 (16%)

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
1	KCX	E	127	3,1	-	4/9/10/12	-
1	KCX	A	127	3,1	-	4/9/10/12	-
1	KCX	G	127	3,1	-	7/9/10/12	-
1	KCX	C	127	3,1	-	6/9/10/12	-

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	127	KCX	CX-NZ	5.90	1.45	1.35
1	A	127	KCX	CX-NZ	5.79	1.45	1.35
1	G	127	KCX	CX-NZ	5.78	1.45	1.35
1	C	127	KCX	CX-NZ	5.76	1.45	1.35
1	E	127	KCX	OQ1-CX	2.55	1.26	1.21
1	A	127	KCX	OQ1-CX	2.47	1.26	1.21
1	G	127	KCX	OQ1-CX	2.43	1.26	1.21
1	C	127	KCX	OQ1-CX	2.19	1.25	1.21

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	127	KCX	OQ1-CX-NZ	-3.54	119.54	124.92
1	C	127	KCX	OQ1-CX-NZ	-3.28	119.94	124.92

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	127	KCX	OQ1-CX-NZ	-3.18	120.09	124.92
1	E	127	KCX	OQ1-CX-NZ	-3.02	120.34	124.92

There are no chirality outliers.

All (21) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	127	KCX	C-CA-CB-CG
1	C	127	KCX	N-CA-CB-CG
1	C	127	KCX	C-CA-CB-CG
1	C	127	KCX	OQ1-CX-NZ-CE
1	E	127	KCX	C-CA-CB-CG
1	G	127	KCX	C-CA-CB-CG
1	E	127	KCX	CG-CD-CE-NZ
1	A	127	KCX	CA-CB-CG-CD
1	C	127	KCX	CA-CB-CG-CD
1	G	127	KCX	CA-CB-CG-CD
1	G	127	KCX	CE-CD-CG-CB
1	C	127	KCX	OQ2-CX-NZ-CE
1	G	127	KCX	OQ1-CX-NZ-CE
1	G	127	KCX	OQ2-CX-NZ-CE
1	C	127	KCX	CE-CD-CG-CB
1	A	127	KCX	N-CA-CB-CG
1	E	127	KCX	N-CA-CB-CG
1	G	127	KCX	N-CA-CB-CG
1	A	127	KCX	CE-CD-CG-CB
1	G	127	KCX	CG-CD-CE-NZ
1	E	127	KCX	CE-CD-CG-CB

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	G	127	KCX	1	0
1	C	127	KCX	1	0

5.5 Carbohydrates

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 15 ligands modelled in this entry, 4 are monoatomic - leaving 11 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	FES	E	601	1	0,4,4	-	-	-		
4	BCT	E	605	-	3,3,3	2.21	1 (33%)	2,3,3	0.12	0
4	BCT	A	604	3	3,3,3	2.20	1 (33%)	2,3,3	0.07	0
2	FES	C	601	1	0,4,4	-	-	-		
4	BCT	E	604	3	3,3,3	2.23	1 (33%)	2,3,3	0.14	0
5	CO2	G	603	3	2,2,2	1.07	0	1,1,1	0.51	0
4	BCT	E	603	-	3,3,3	2.19	1 (33%)	2,3,3	0.10	0
2	FES	G	601	1	0,4,4	-	-	-		
2	FES	A	601	1	0,4,4	-	-	-		
4	BCT	C	603	3	3,3,3	2.18	1 (33%)	2,3,3	0.03	0
4	BCT	A	603	-	3,3,3	2.21	1 (33%)	2,3,3	0.12	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	FES	G	601	1	-	-	0/1/1/1
2	FES	A	601	1	-	-	0/1/1/1
2	FES	C	601	1	-	-	0/1/1/1
2	FES	E	601	1	-	-	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	E	604	BCT	O1-C	3.61	1.38	1.25
4	A	603	BCT	O1-C	3.57	1.38	1.25
4	E	605	BCT	O1-C	3.57	1.38	1.25
4	A	604	BCT	O1-C	3.54	1.38	1.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	E	603	BCT	O1-C	3.54	1.37	1.25
4	C	603	BCT	O1-C	3.53	1.37	1.25

There are no bond angle outliers.

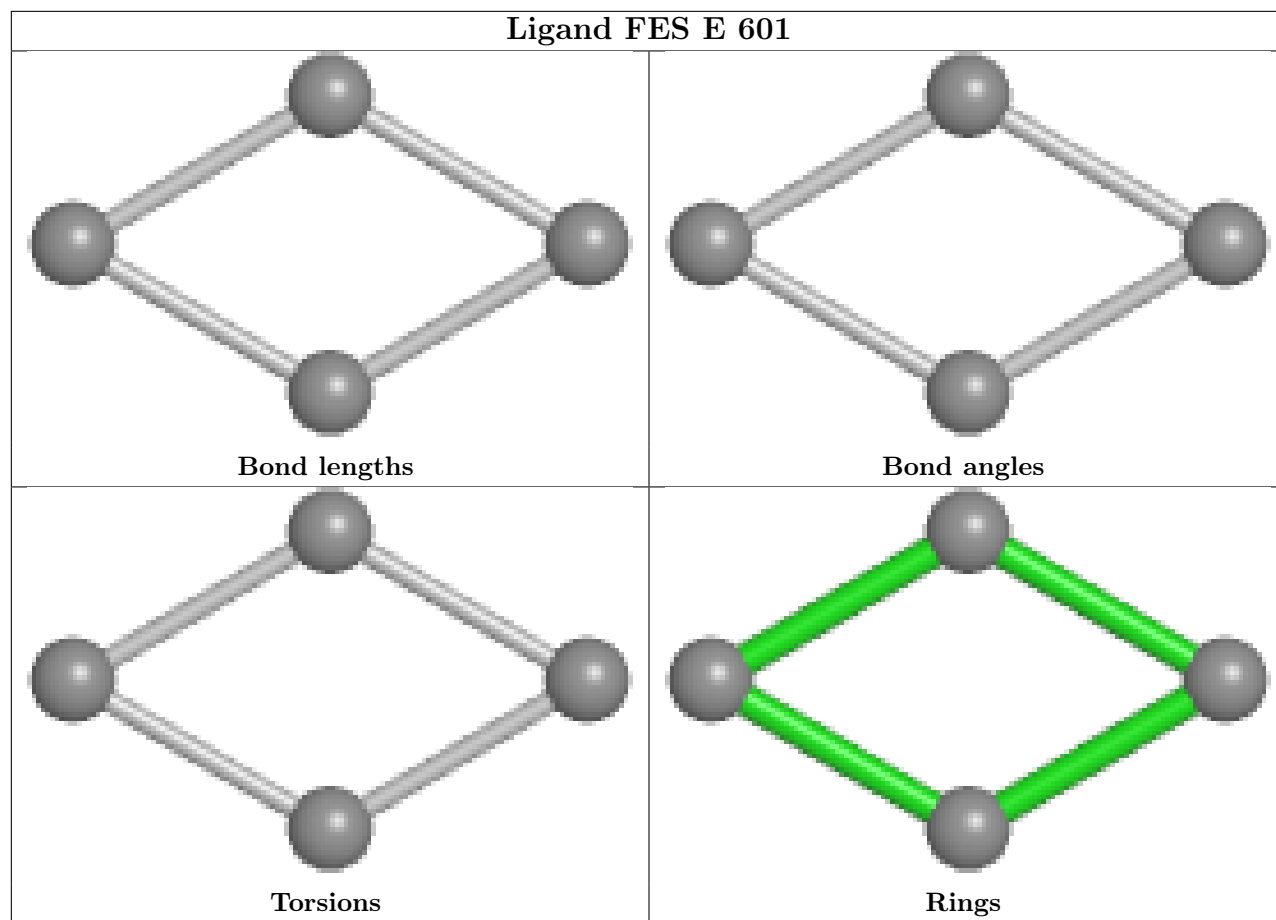
There are no chirality outliers.

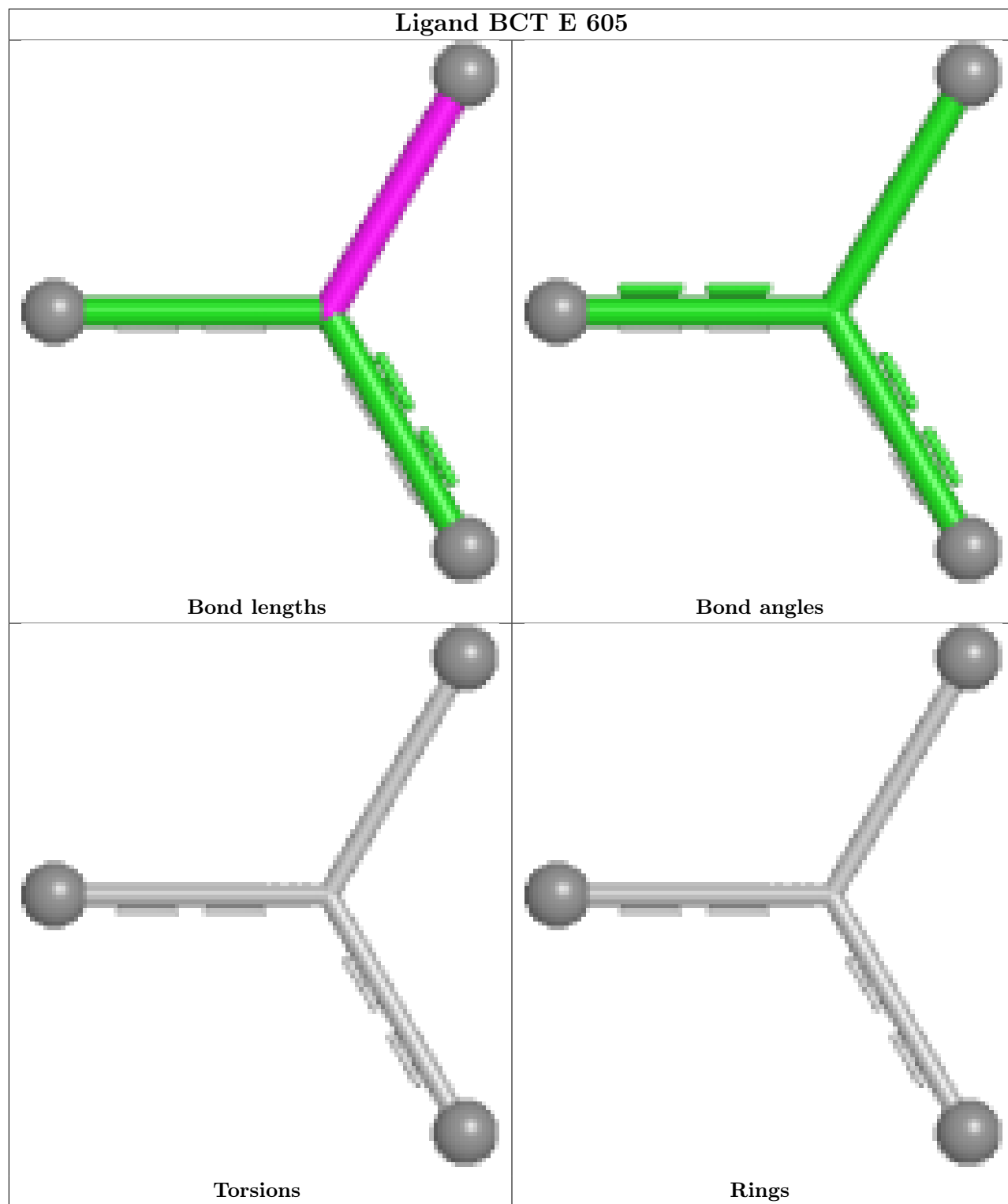
There are no torsion outliers.

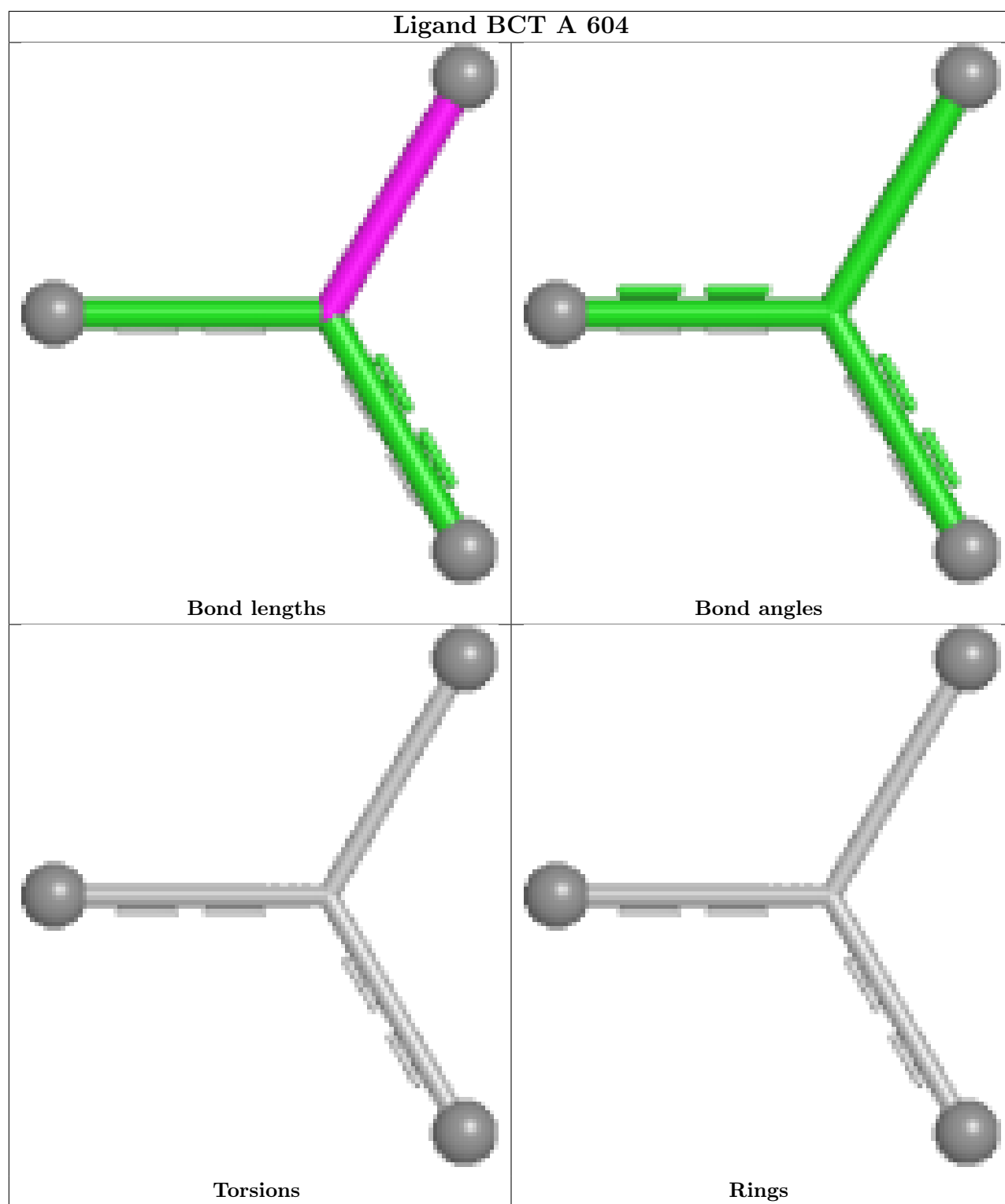
There are no ring outliers.

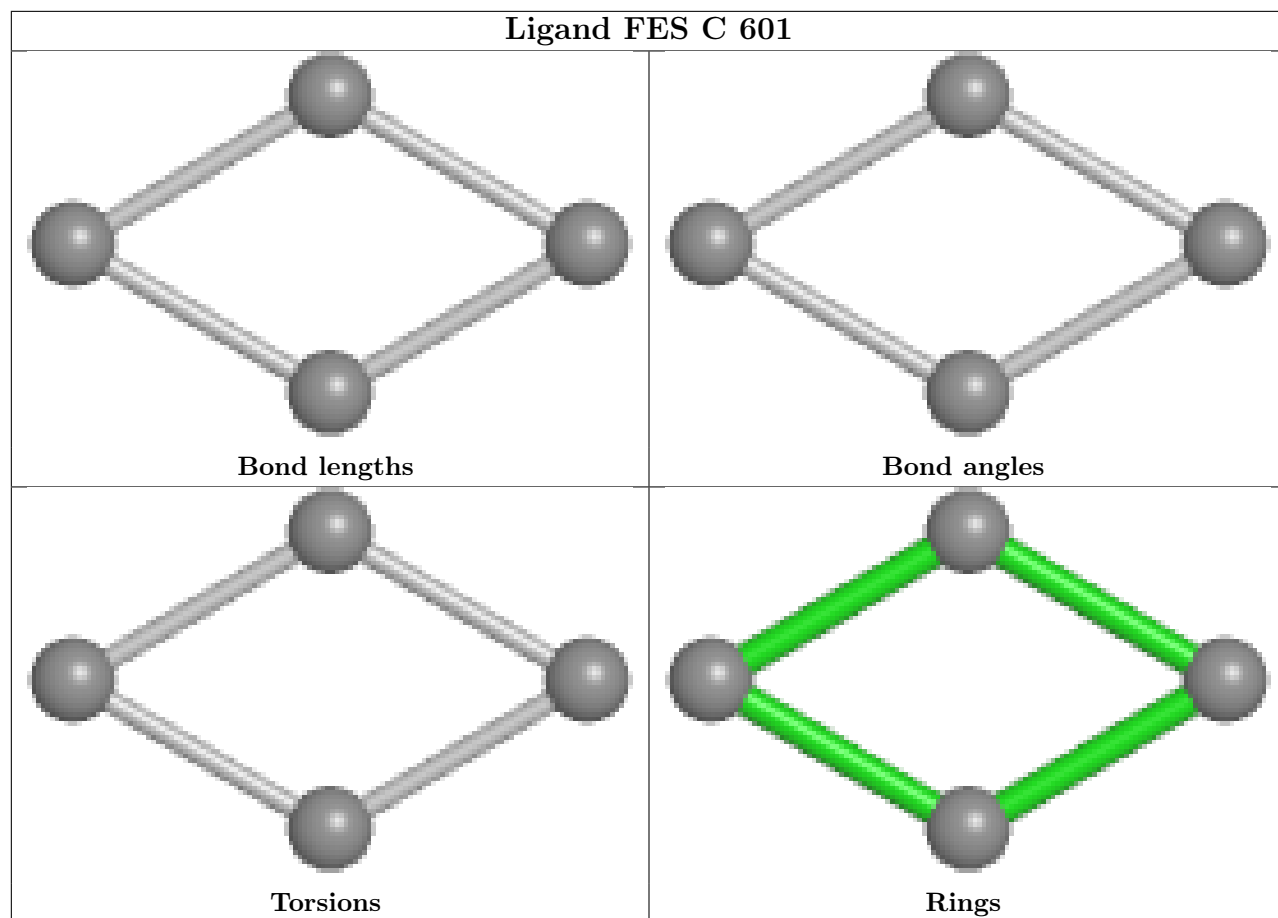
No monomer is involved in short contacts.

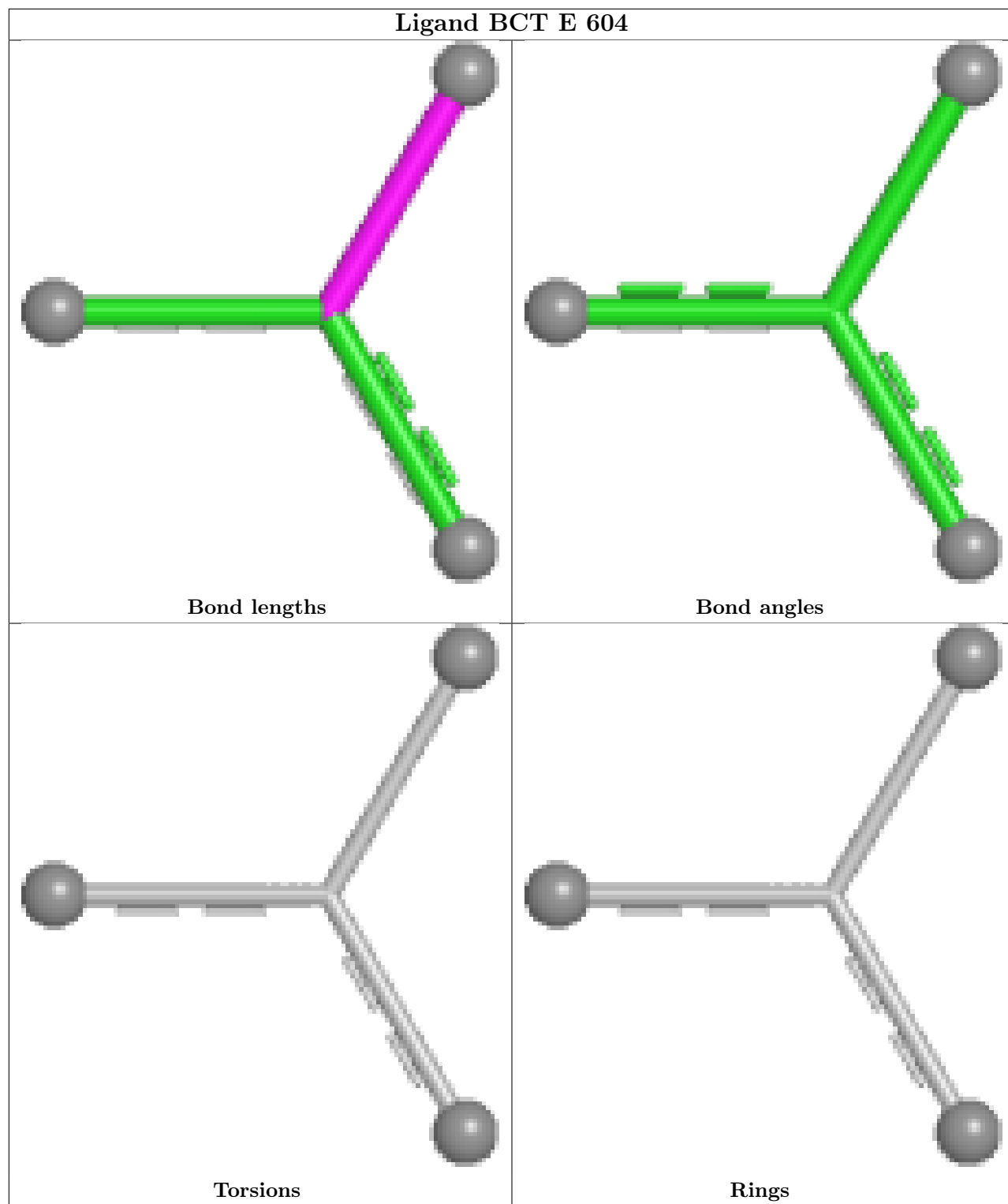
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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



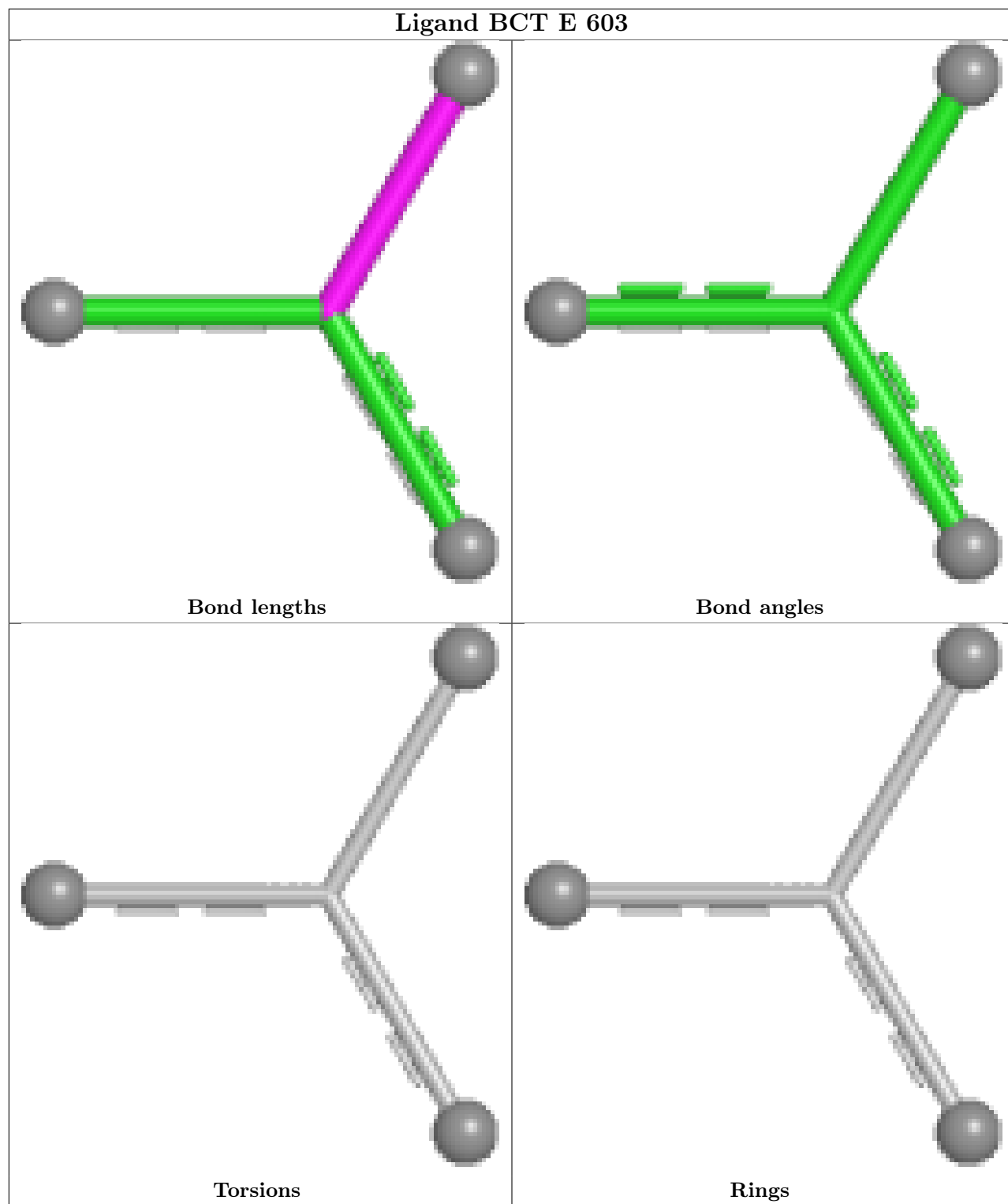


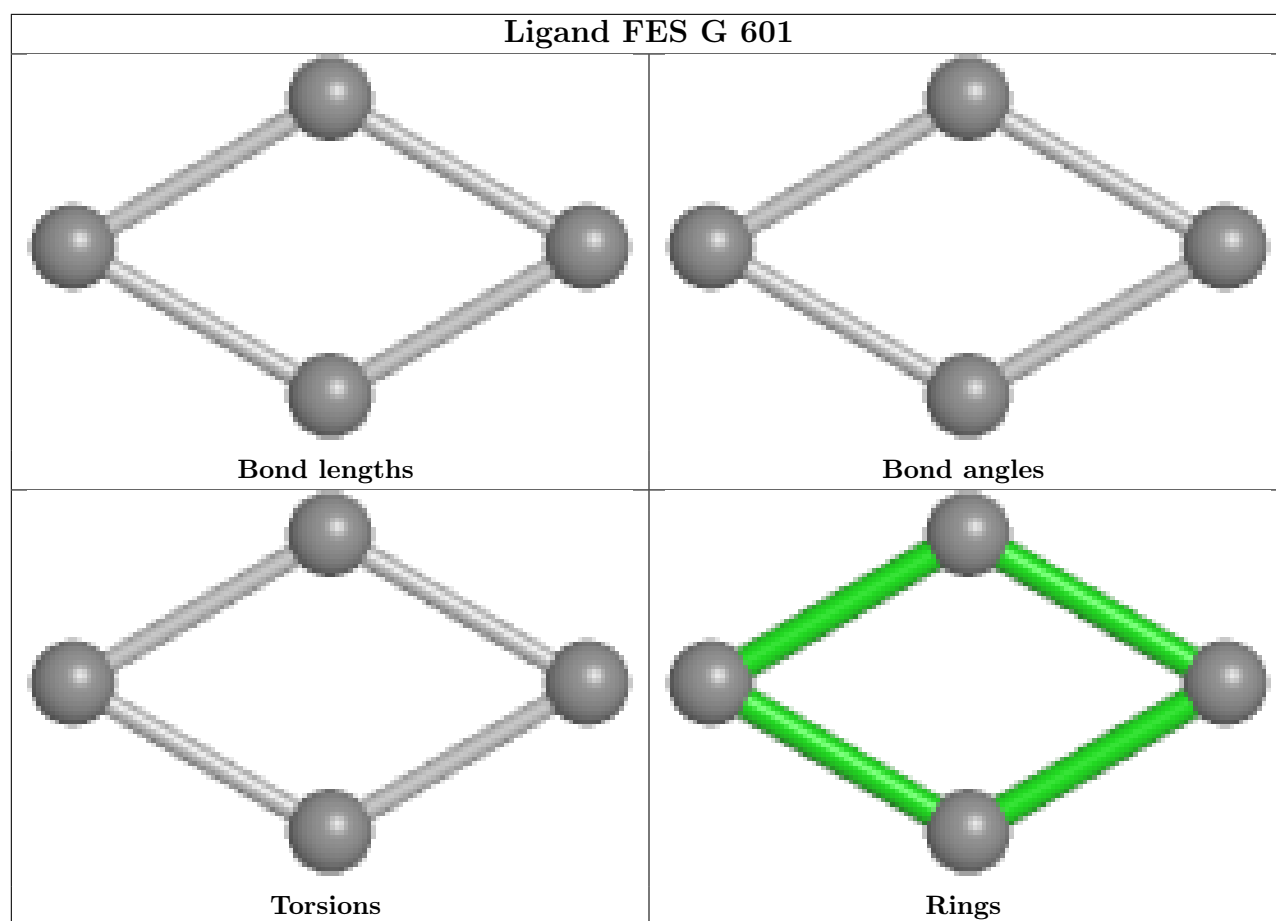


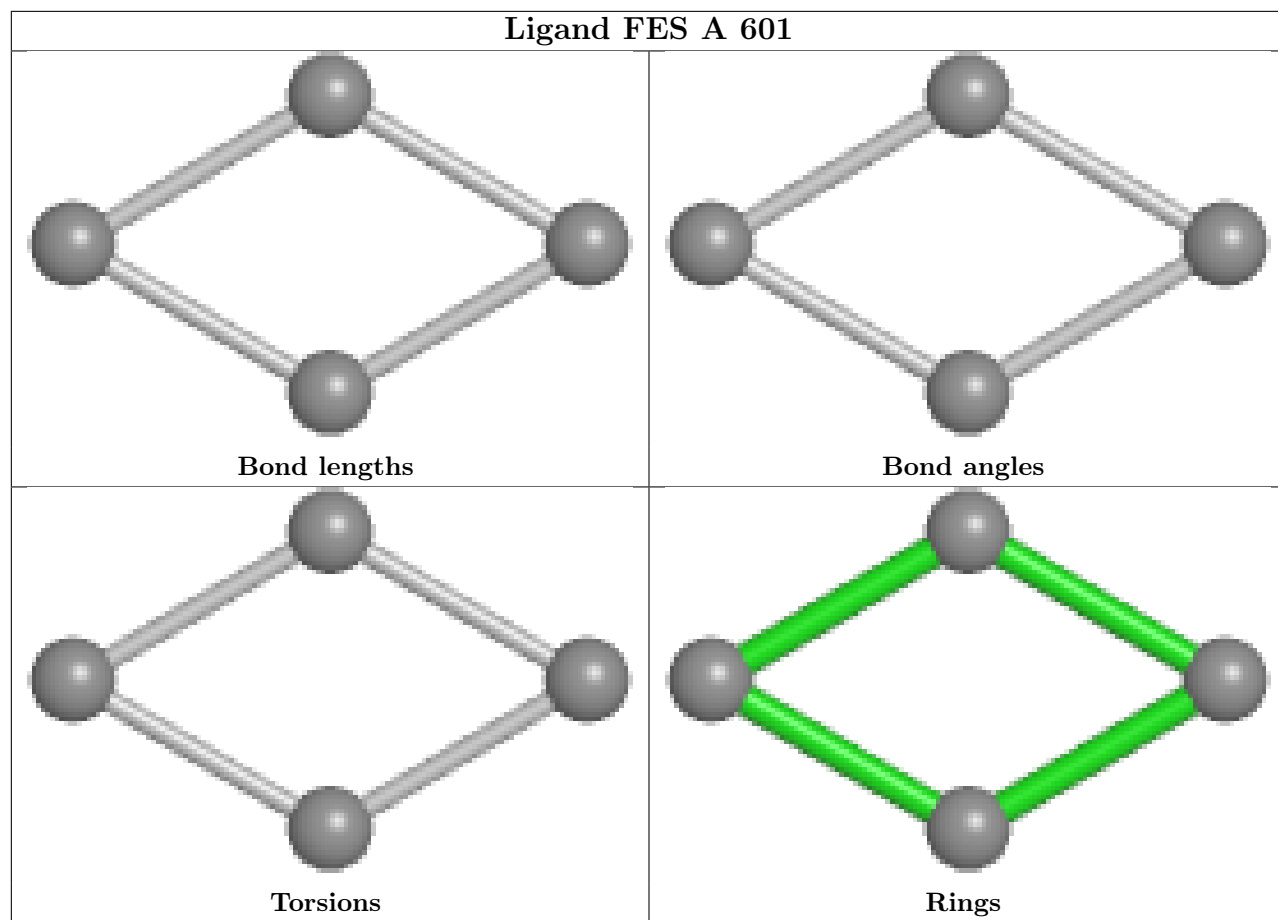


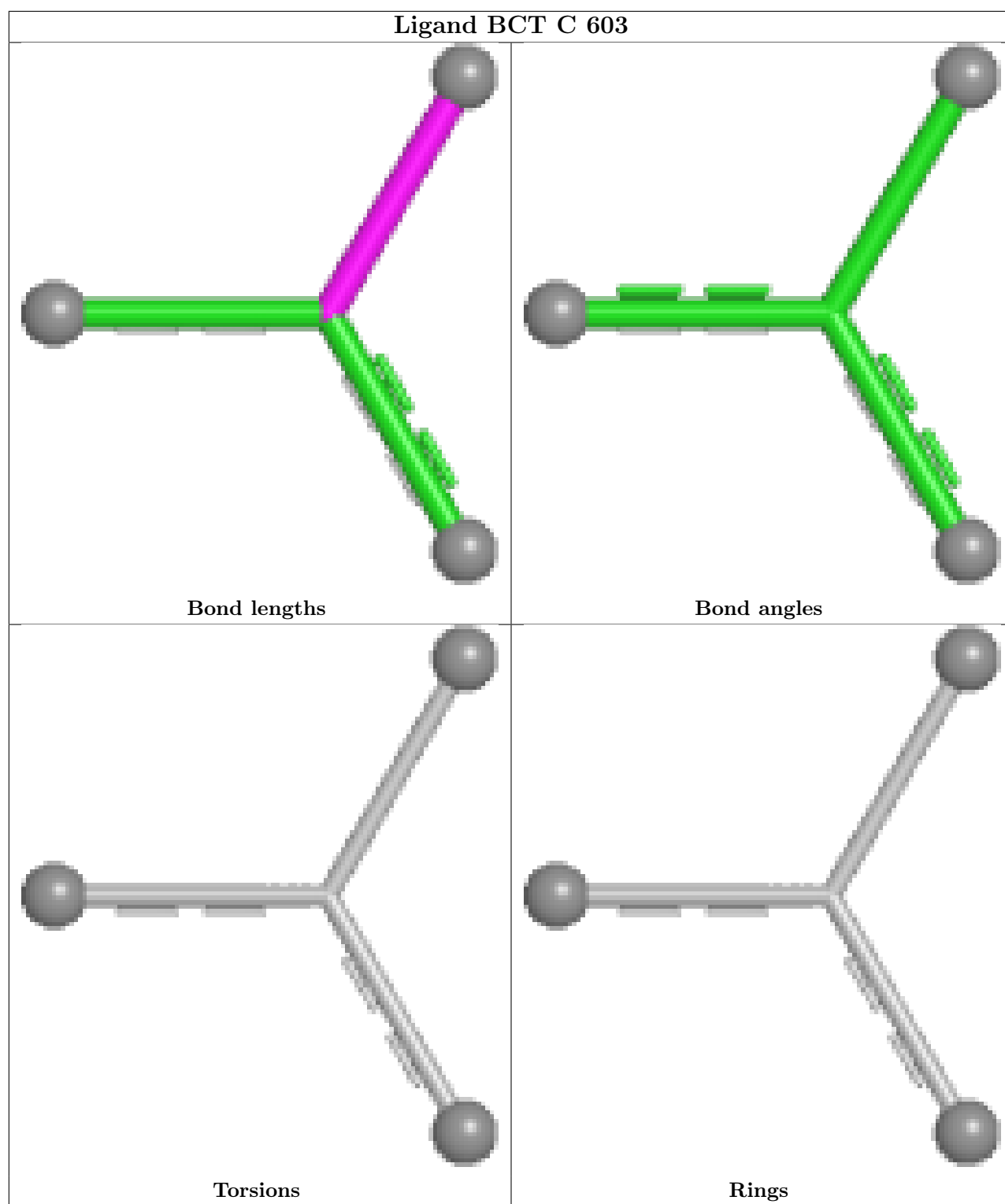


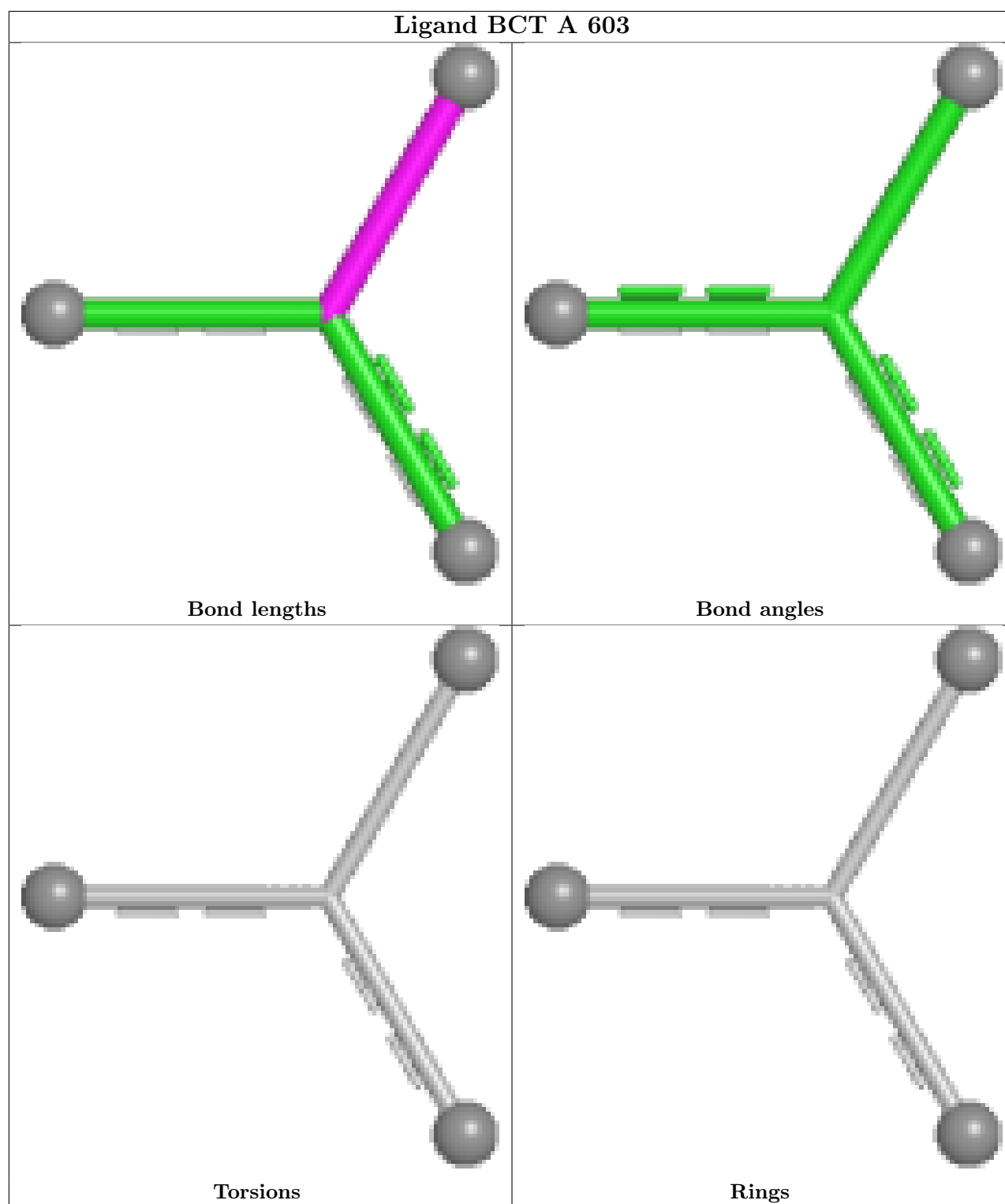
Ligand CO2 G 603			
			
Bond lengths		Bond angles	
			
Torsions		Rings	











5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	C	3
1	E	2
1	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	C	124:GLY	C	125:CYS	N	1.64
1	C	70:ILE	C	71:GLY	N	1.14
1	E	70:ILE	C	71:GLY	N	1.14
1	C	69:LYS	C	70:ILE	N	1.09
1	E	69:LYS	C	70:ILE	N	0.98
1	A	69:LYS	C	70:ILE	N	0.94

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	567/575 (98%)	-0.11	0 100 100	32, 46, 65, 81	1 (0%)
1	C	567/575 (98%)	0.04	3 (0%) 87 85	18, 51, 74, 86	3 (0%)
1	E	567/575 (98%)	-0.12	1 (0%) 91 90	18, 45, 66, 79	2 (0%)
1	G	567/575 (98%)	0.17	7 (1%) 76 73	20, 54, 87, 100	1 (0%)
All	All	2268/2300 (98%)	-0.00	11 (0%) 87 85	18, 49, 76, 100	7 (0%)

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	455	LEU	3.0
1	G	378	ILE	3.0
1	G	506	ILE	2.9
1	C	532	GLN	2.4
1	G	525	ALA	2.4
1	C	8	VAL	2.3
1	G	399	PRO	2.3
1	E	395	SER	2.2
1	G	526	ARG	2.1
1	C	506	ILE	2.1
1	G	486	HIS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	KCX	G	127	12/13	0.93	0.09	42,46,52,53	0

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
1	KCX	E	127	12/13	0.94	0.10	36,41,43,46	0
1	KCX	C	127	12/13	0.94	0.10	39,43,49,50	0
1	KCX	A	127	12/13	0.96	0.07	37,42,44,44	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

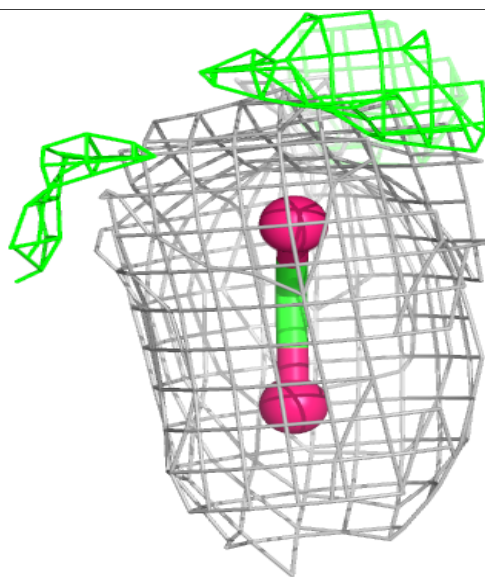
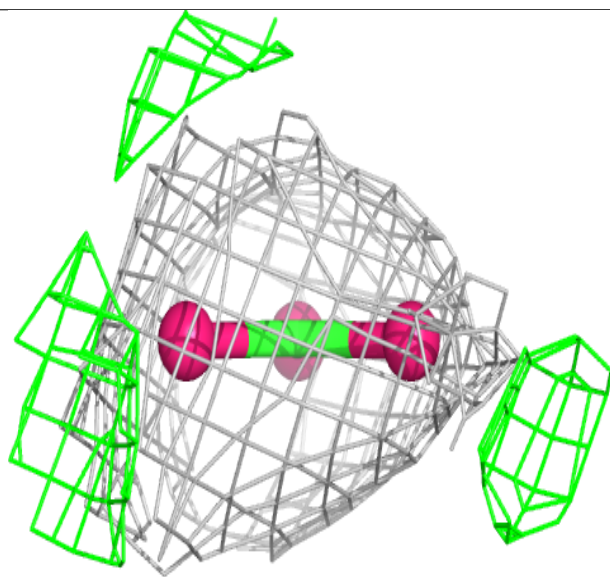
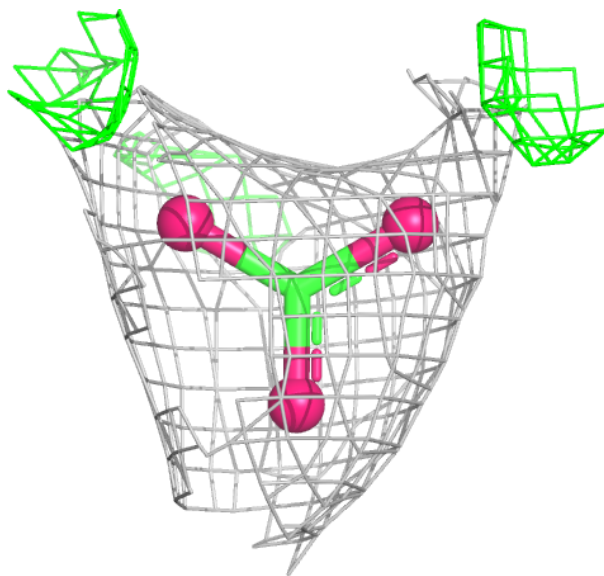
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	BCT	E	603	4/4	0.75	0.16	57,62,66,67	0
4	BCT	E	605	4/4	0.83	0.15	61,65,68,73	0
4	BCT	E	604	4/4	0.85	0.14	51,52,54,57	0
4	BCT	A	604	4/4	0.85	0.10	49,50,54,56	0
4	BCT	A	603	4/4	0.88	0.12	62,63,63,64	0
5	CO2	G	603	3/3	0.88	0.09	57,57,58,60	0
4	BCT	C	603	4/4	0.90	0.09	52,53,54,55	0
3	MG	C	602	1/1	0.95	0.07	46,46,46,46	0
2	FES	C	601	4/4	0.95	0.06	40,44,46,67	4
2	FES	G	601	4/4	0.96	0.06	45,46,48,64	4
2	FES	E	601	4/4	0.97	0.04	36,42,45,54	4
2	FES	A	601	4/4	0.97	0.05	39,42,47,62	4
3	MG	E	602	1/1	0.98	0.07	31,31,31,31	0
3	MG	A	602	1/1	0.98	0.07	36,36,36,36	0
3	MG	G	602	1/1	0.99	0.06	41,41,41,41	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.

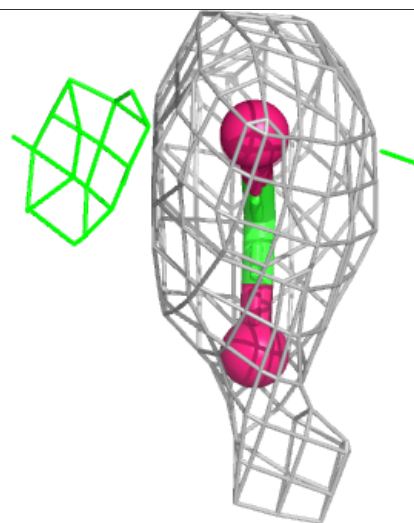
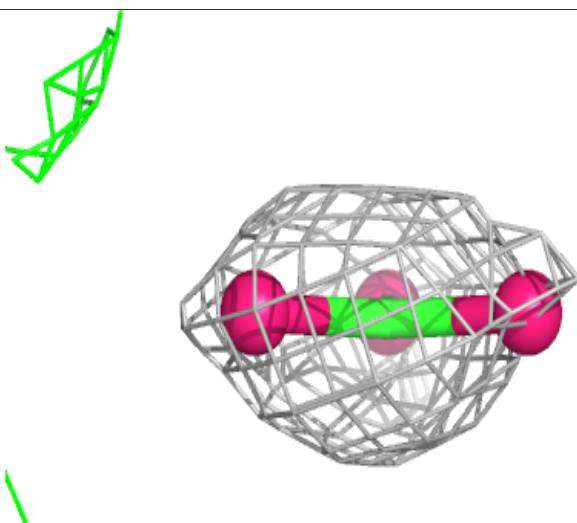
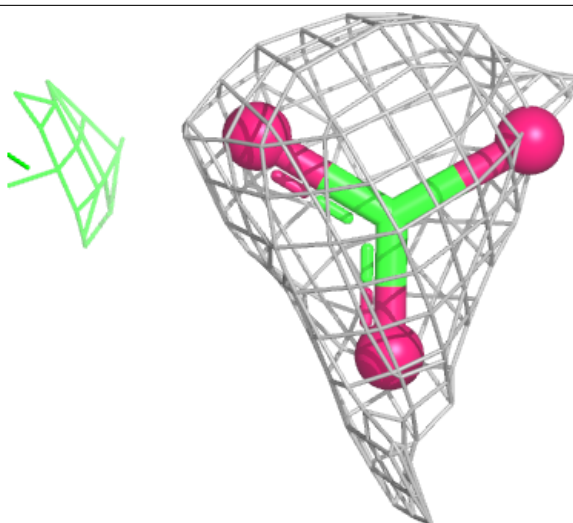
Electron density around BCT E 603:

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



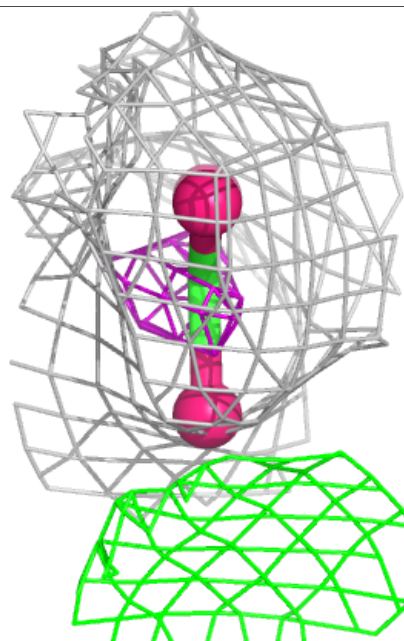
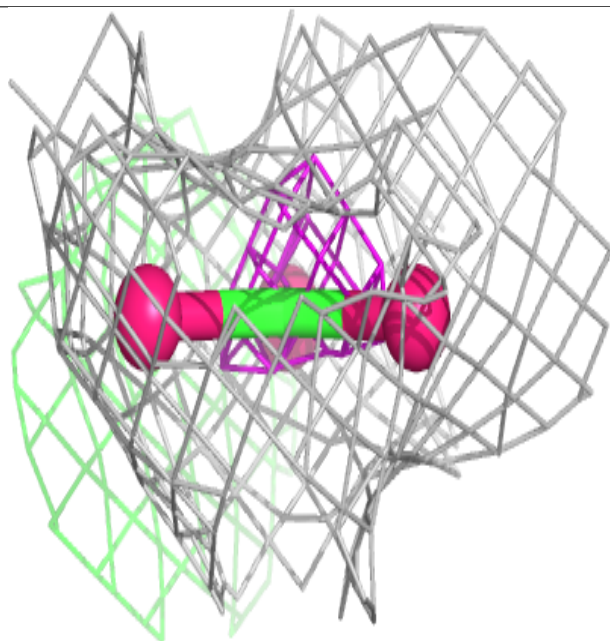
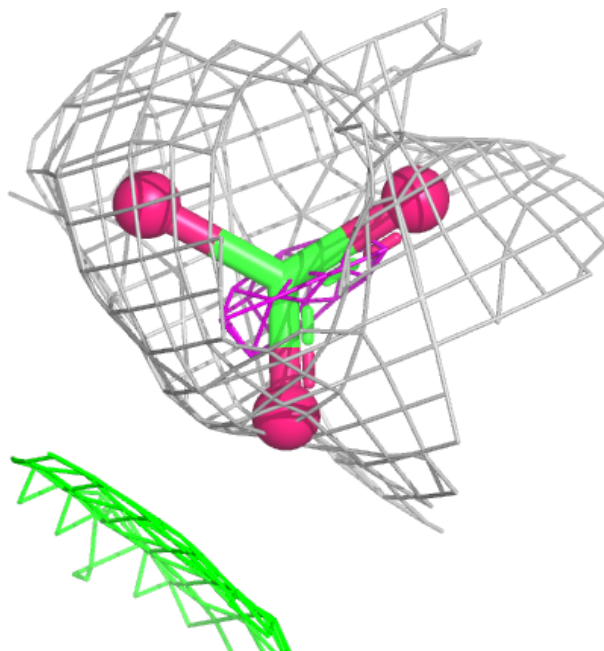
Electron density around BCT E 605:

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



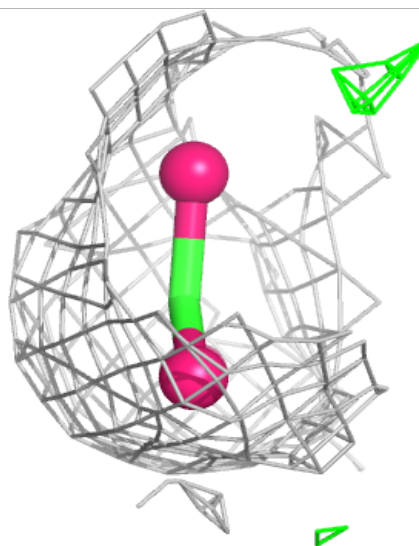
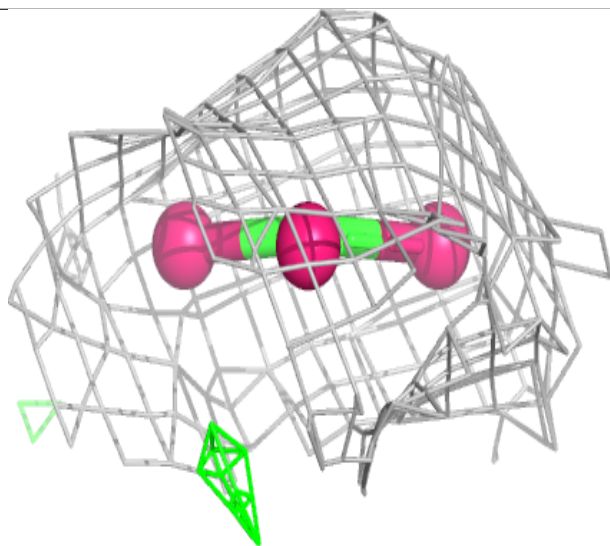
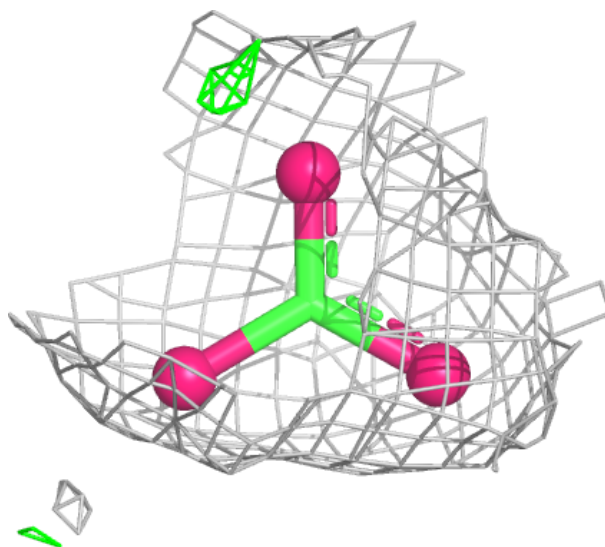
Electron density around BCT E 604:

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



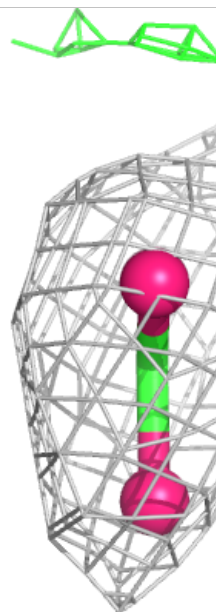
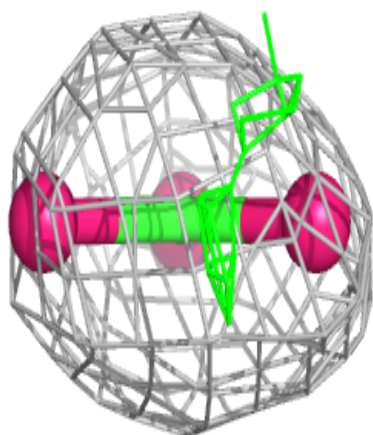
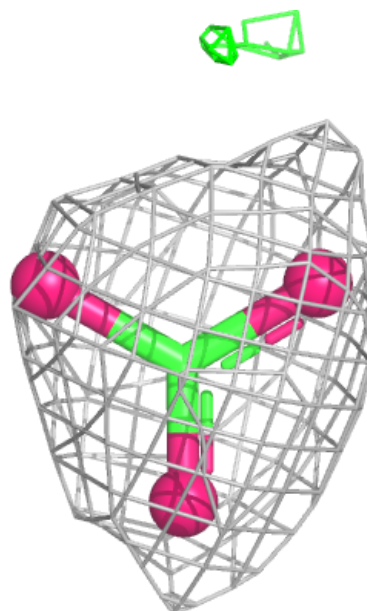
Electron density around BCT A 604:

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



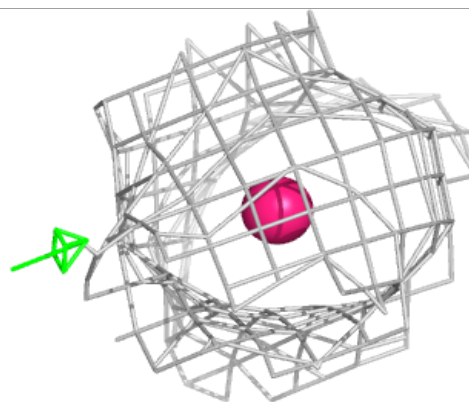
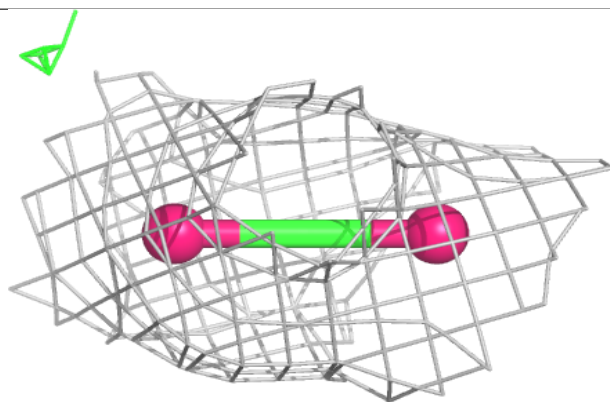
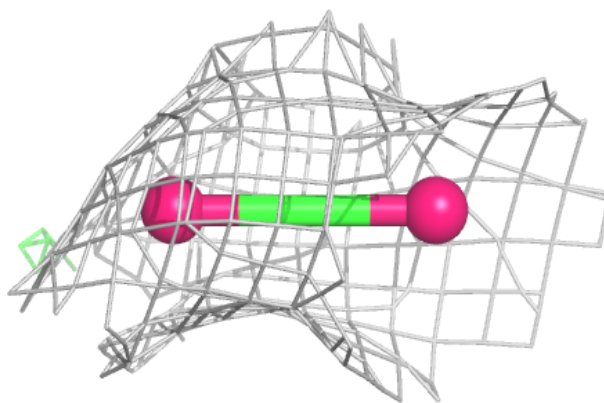
Electron density around BCT A 603:

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and green (positive)



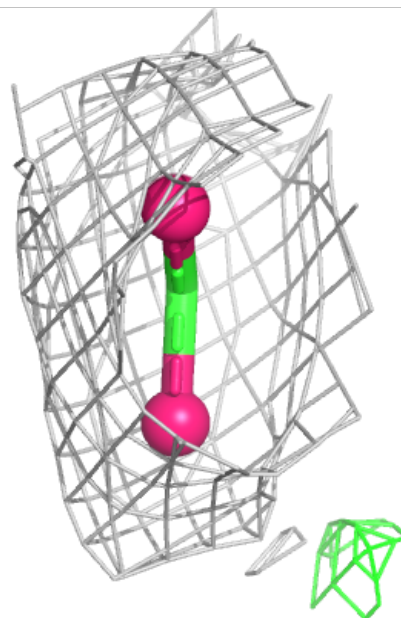
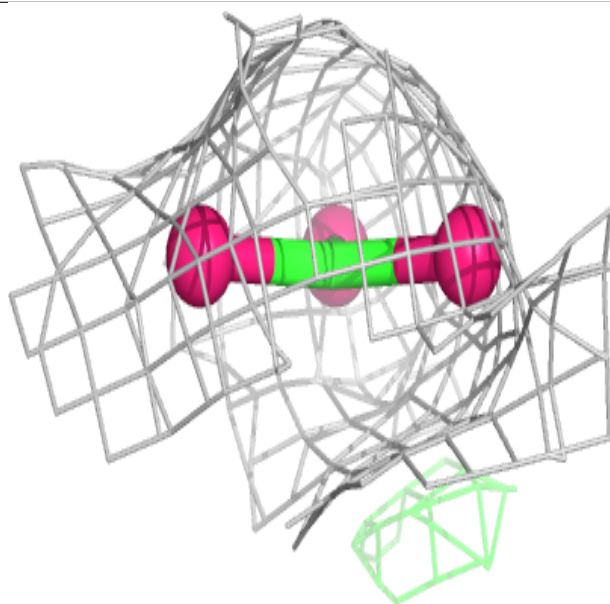
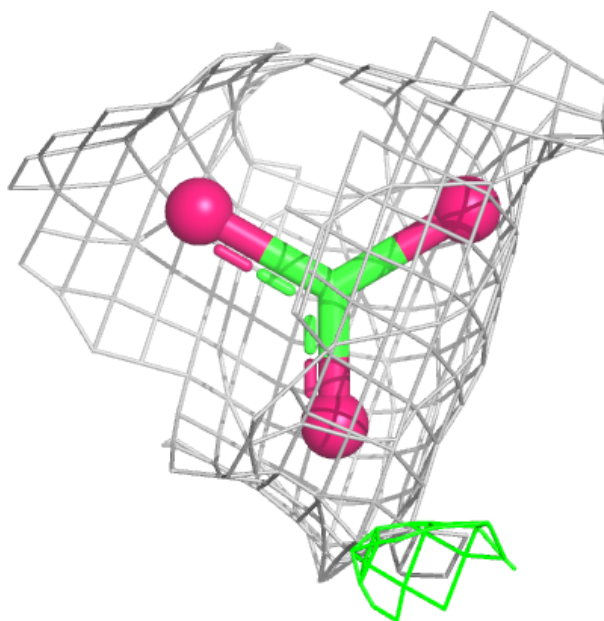
Electron density around CO2 G 603:

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and green (positive)



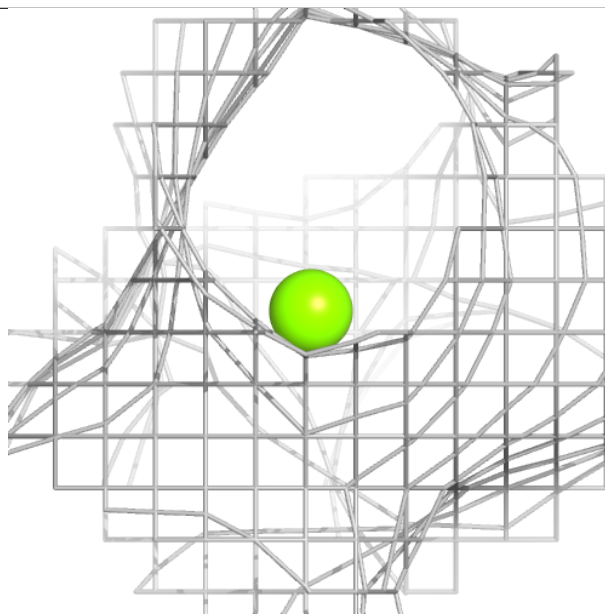
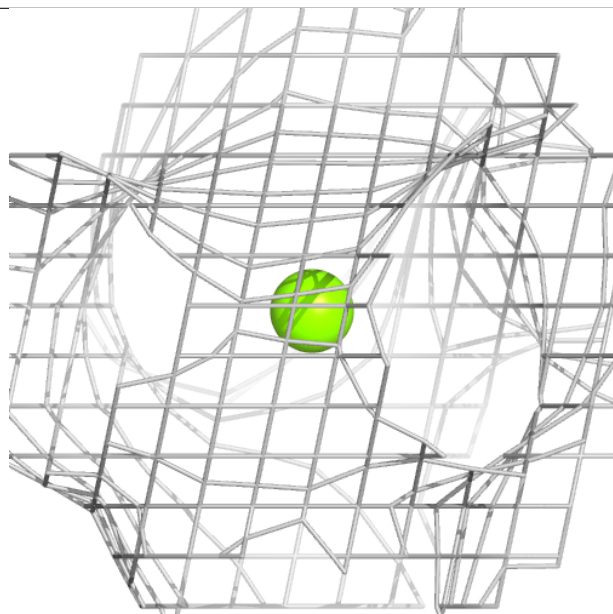
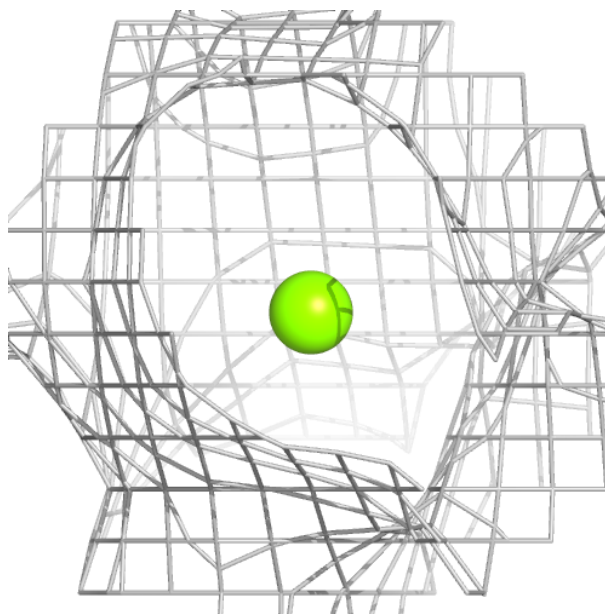
Electron density around BCT C 603:

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and green (positive)



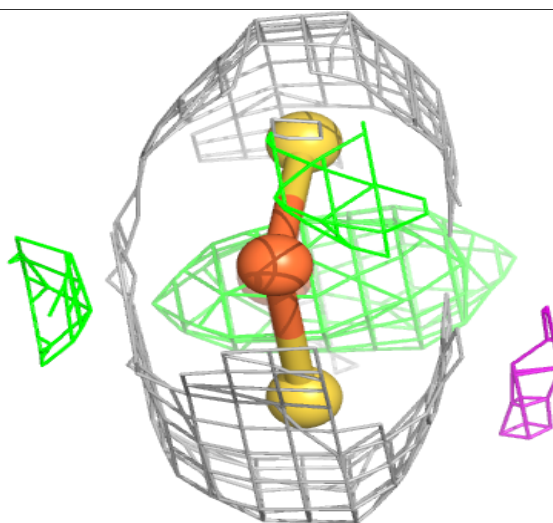
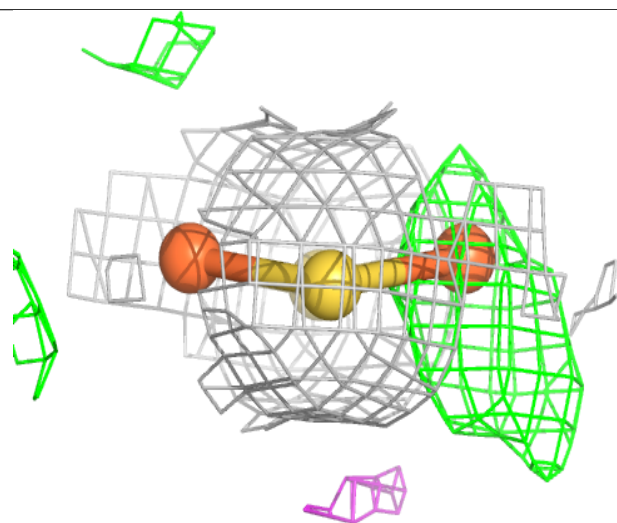
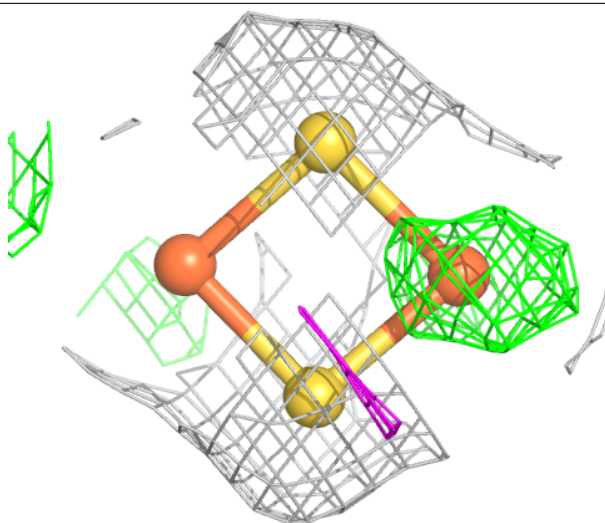
Electron density around MG C 602:

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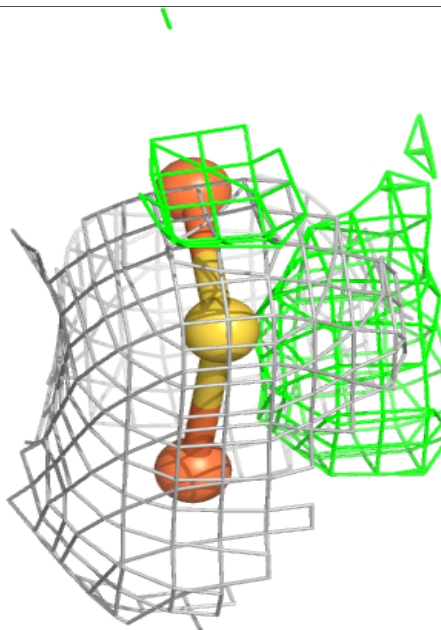
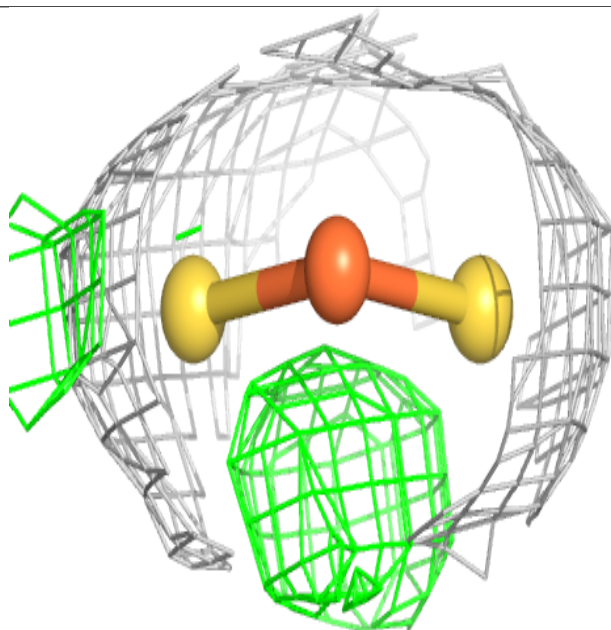
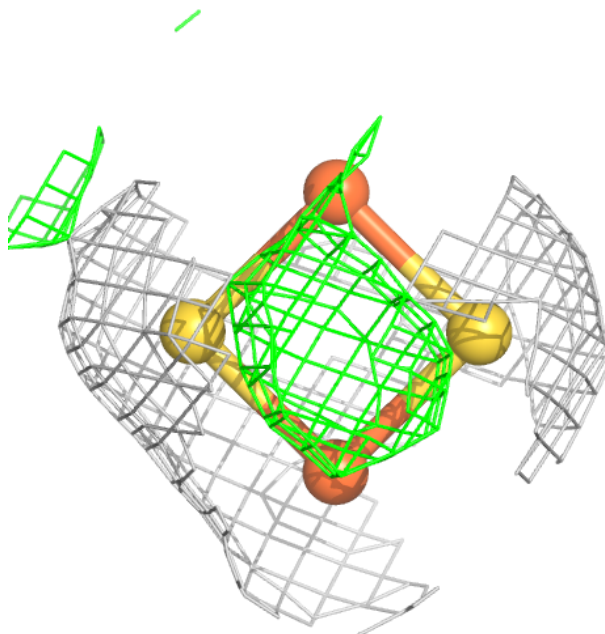
Electron density around FES C 601:

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and green (positive)



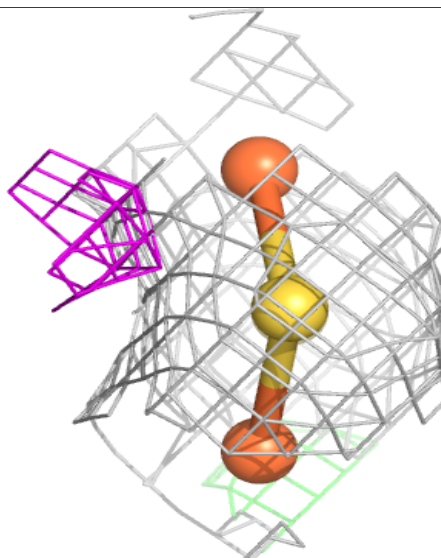
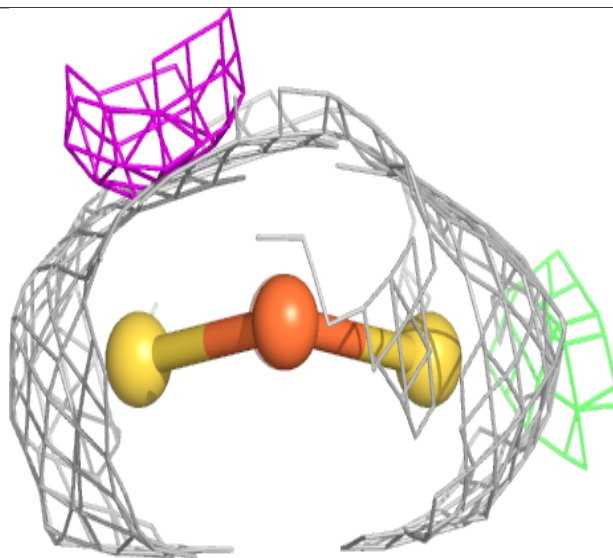
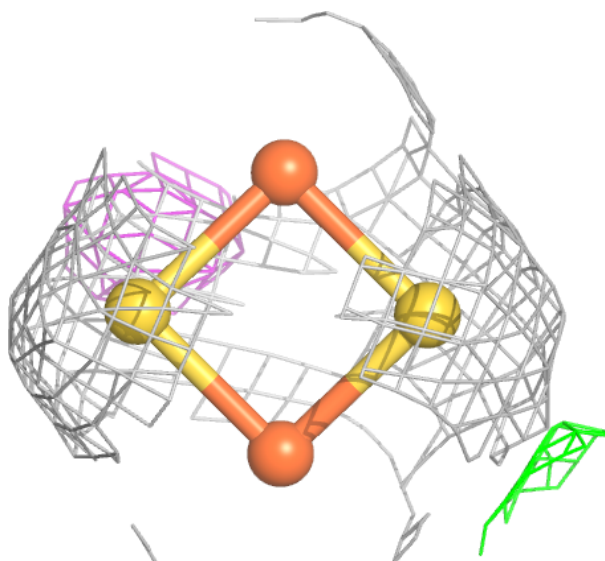
Electron density around FES G 601:

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and green (positive)



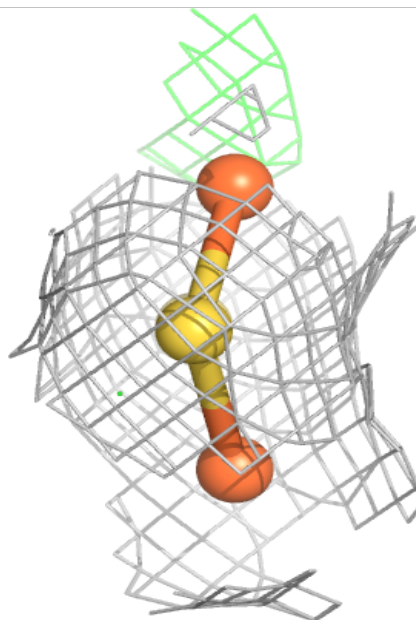
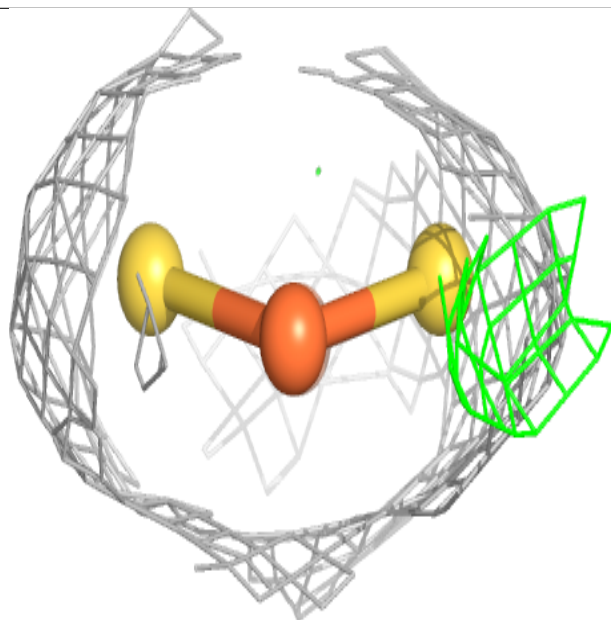
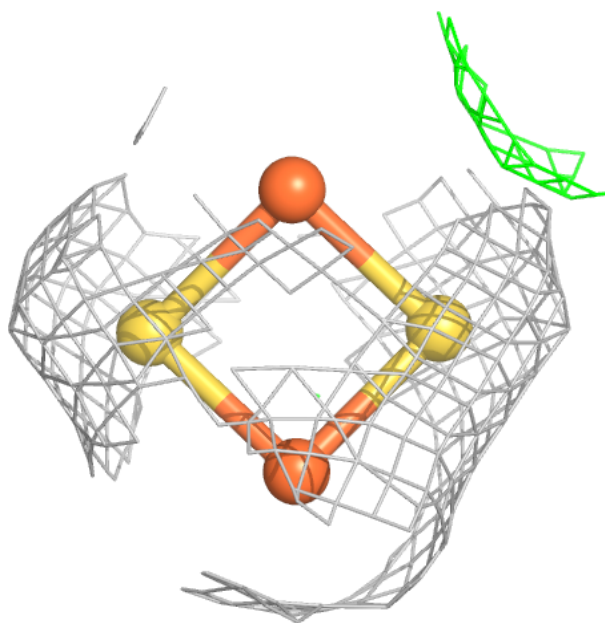
Electron density around FES E 601:

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and green (positive)



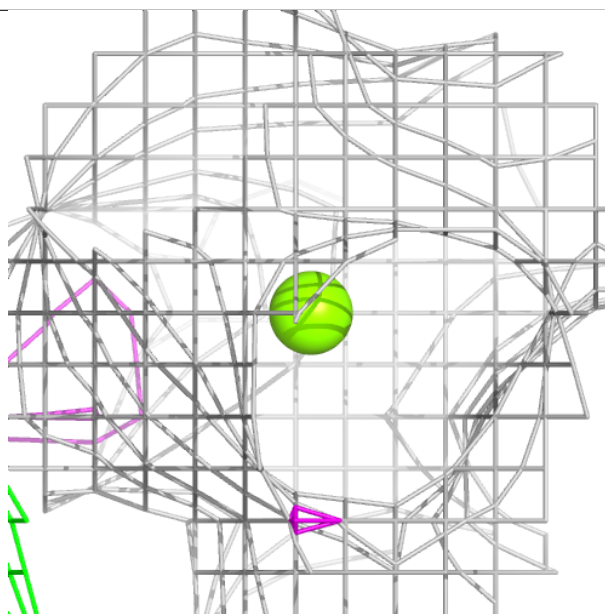
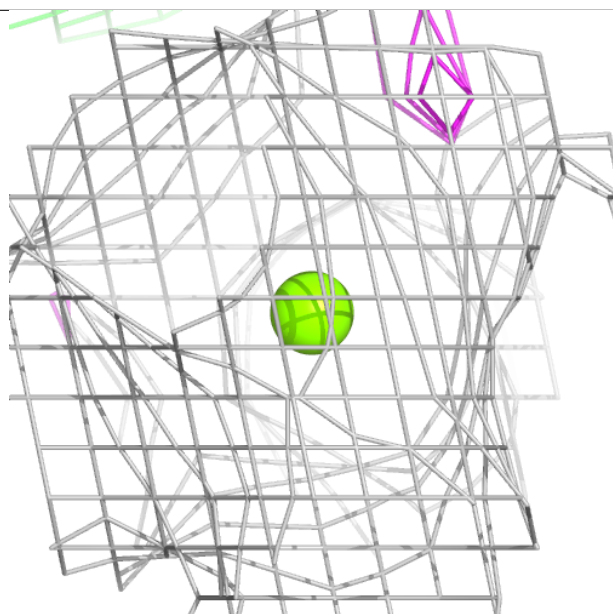
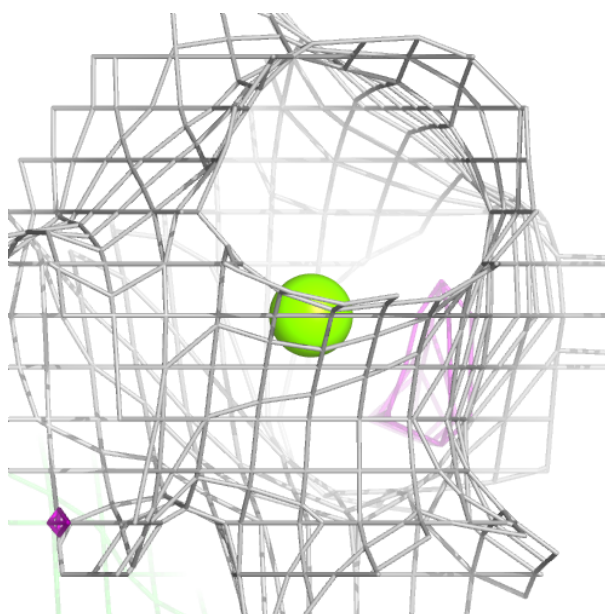
Electron density around FES A 601:

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and green (positive)



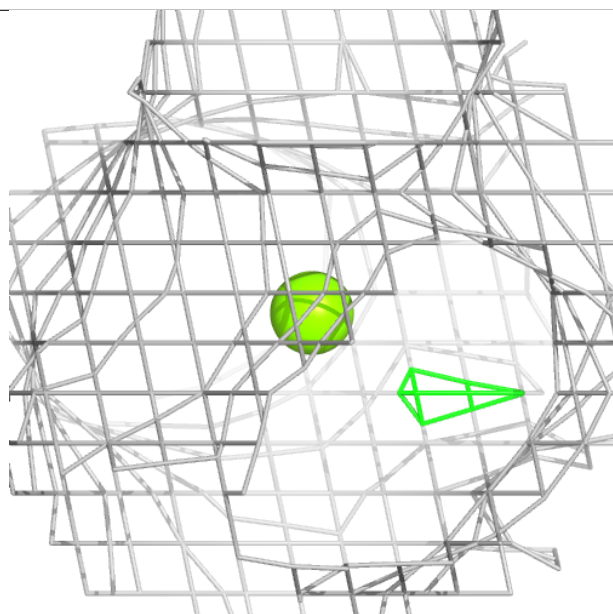
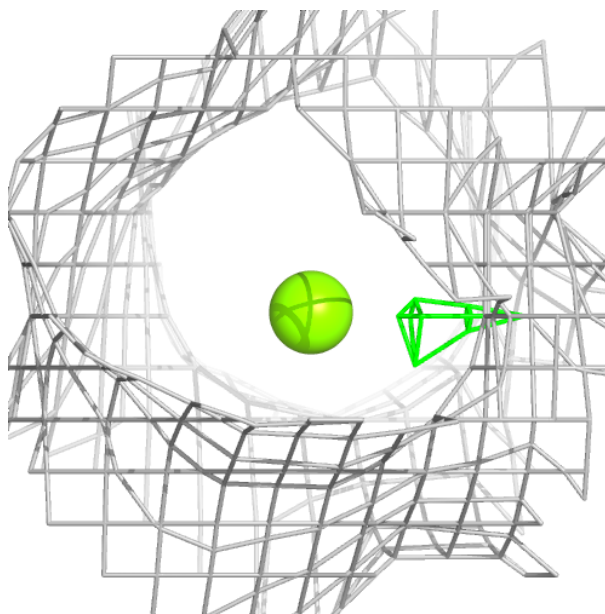
Electron density around MG E 602:

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and green (positive)



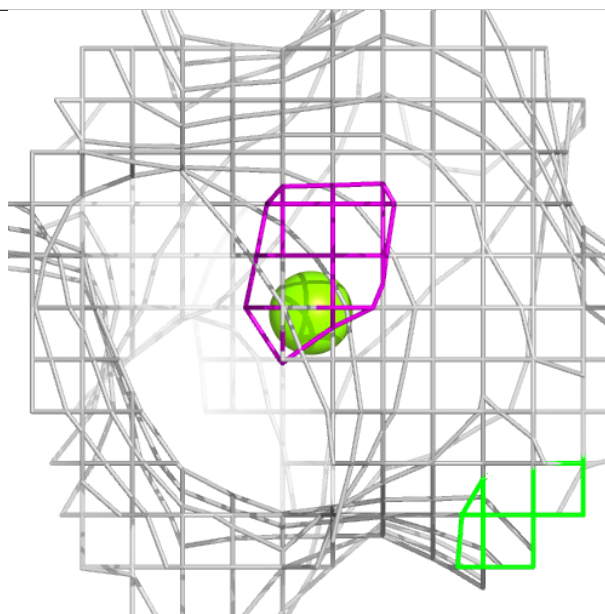
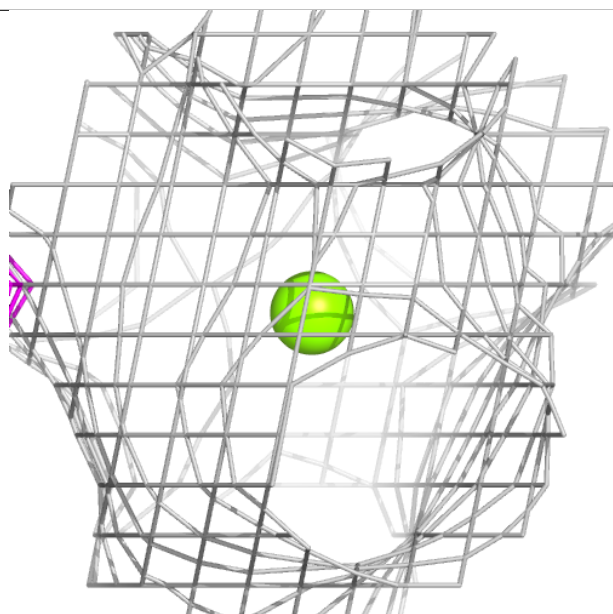
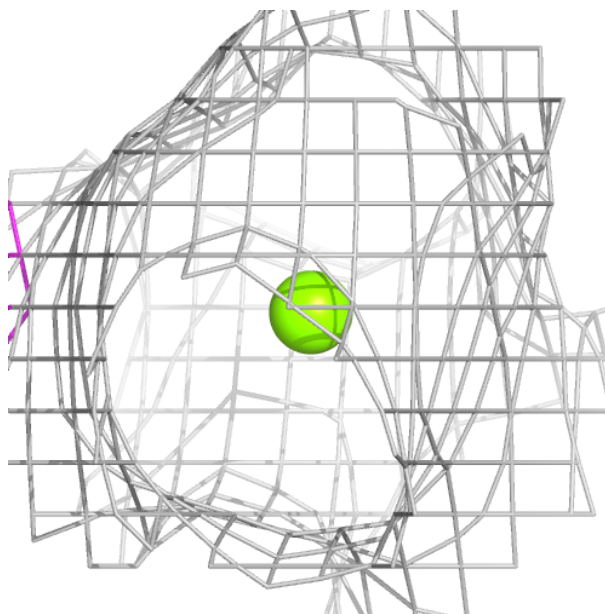
Electron density around MG A 602:

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



Electron density around MG G 602:

$2mF_o-DF_c$ (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.