



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 25, 2024 – 12:33 AM EDT

PDB ID : 6NXA
Title : ECAII(D90T,K162T) MUTANT AT PH 7
Authors : Lubkowski, J.; Wlodawer, A.
Deposited on : 2019-02-08
Resolution : 1.93 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.37.1
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.37.1

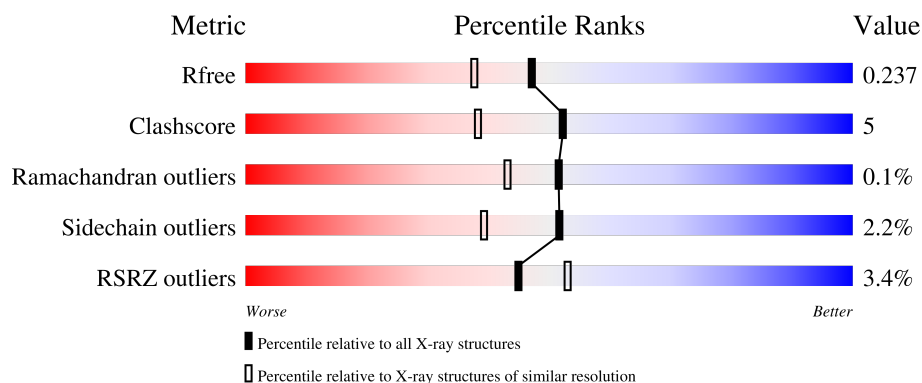
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.93 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	4310 (1.96-1.92)
Clashscore	141614	1023 (1.94-1.94)
Ramachandran outliers	138981	1007 (1.94-1.94)
Sidechain outliers	138945	1007 (1.94-1.94)
RSRZ outliers	127900	4250 (1.96-1.92)

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	333	<div> <div>5%</div> <div> <div></div> <div>79%</div> <div>16%</div> <div>• •</div> </div> </div>
1	B	333	<div> <div>5%</div> <div> <div></div> <div>76%</div> <div>17%</div> <div>• 7%</div> </div> </div>
1	C	333	<div> <div>2%</div> <div> <div></div> <div>82%</div> <div>11%</div> <div>• 6%</div> </div> </div>
1	D	333	<div> <div>%</div> <div> <div></div> <div>83%</div> <div>10%</div> <div>• 6%</div> </div> </div>

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	ACY	B	502	-	-	X	-
2	ACY	D	401	-	-	X	-
3	GOL	B	501	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10658 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 L-asparaginase 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	319	Total	C	N	O	S	0	3	0
			2401	1502	408	483	8			
1	B	311	Total	C	N	O	S	0	1	0
			2336	1460	398	470	8			
1	C	313	Total	C	N	O	S	0	2	0
			2355	1471	403	473	8			
1	D	313	Total	C	N	O	S	0	1	0
			2348	1467	402	471	8			

There are 36 discrepancies between the modelled and reference sequences:

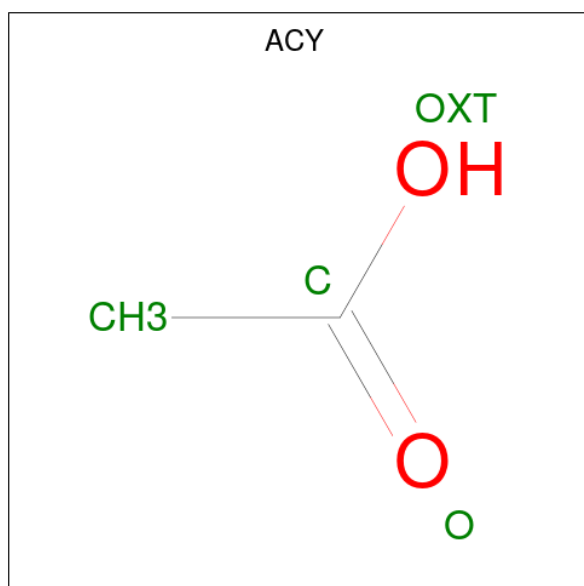
Chain	Residue	Modelled	Actual	Comment	Reference
A	-6	MET	-	initiating methionine	UNP P00805
A	-5	HIS	-	expression tag	UNP P00805
A	-4	HIS	-	expression tag	UNP P00805
A	-3	HIS	-	expression tag	UNP P00805
A	-2	HIS	-	expression tag	UNP P00805
A	-1	HIS	-	expression tag	UNP P00805
A	0	HIS	-	expression tag	UNP P00805
A	90	THR	ASP	conflict	UNP P00805
A	162	THR	LYS	engineered mutation	UNP P00805
B	-6	MET	-	initiating methionine	UNP P00805
B	-5	HIS	-	expression tag	UNP P00805
B	-4	HIS	-	expression tag	UNP P00805
B	-3	HIS	-	expression tag	UNP P00805
B	-2	HIS	-	expression tag	UNP P00805
B	-1	HIS	-	expression tag	UNP P00805
B	0	HIS	-	expression tag	UNP P00805
B	90	THR	ASP	conflict	UNP P00805
B	162	THR	LYS	engineered mutation	UNP P00805
C	-6	MET	-	initiating methionine	UNP P00805
C	-5	HIS	-	expression tag	UNP P00805
C	-4	HIS	-	expression tag	UNP P00805

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-3	HIS	-	expression tag	UNP P00805
C	-2	HIS	-	expression tag	UNP P00805
C	-1	HIS	-	expression tag	UNP P00805
C	0	HIS	-	expression tag	UNP P00805
C	90	THR	ASP	conflict	UNP P00805
C	162	THR	LYS	engineered mutation	UNP P00805
D	-6	MET	-	initiating methionine	UNP P00805
D	-5	HIS	-	expression tag	UNP P00805
D	-4	HIS	-	expression tag	UNP P00805
D	-3	HIS	-	expression tag	UNP P00805
D	-2	HIS	-	expression tag	UNP P00805
D	-1	HIS	-	expression tag	UNP P00805
D	0	HIS	-	expression tag	UNP P00805
D	90	THR	ASP	conflict	UNP P00805
D	162	THR	LYS	engineered mutation	UNP P00805

- Molecule 2 is ACETIC ACID (three-letter code: ACY) (formula: C₂H₄O₂) (labeled as "Ligand of Interest" by depositor).



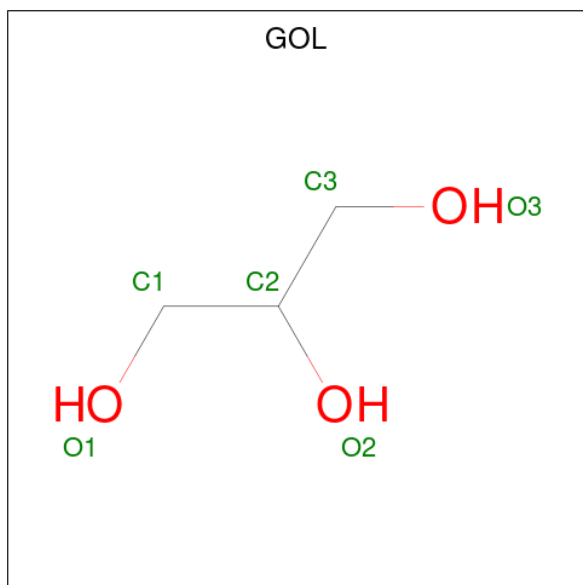
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 4 2 2	0	0
2	B	1	Total C O 4 2 2	0	0
2	C	1	Total C O 4 2 2	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	D	1	Total	C	O	0	0
			4	2	2		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			6	3	3		
3	C	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		

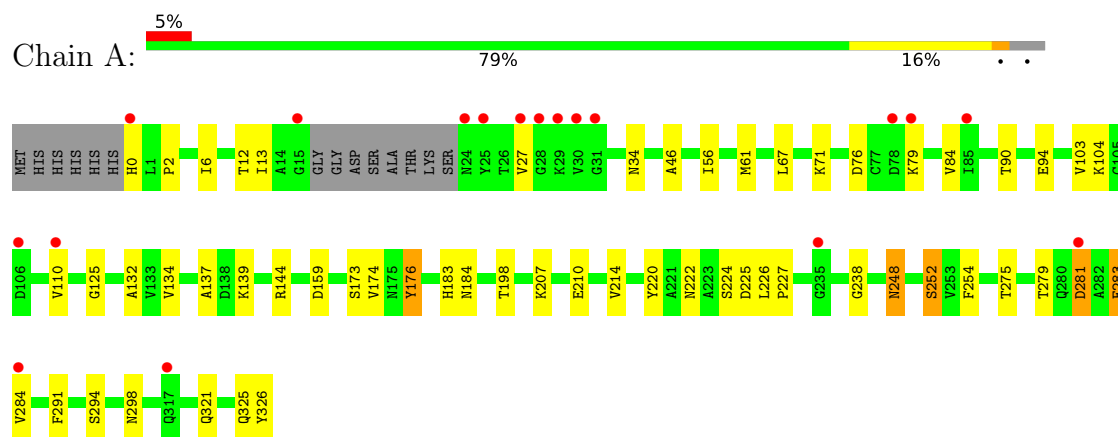
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	298	Total	O	0	1
			299	299		
4	B	247	Total	O	0	0
			247	247		
4	C	307	Total	O	0	2
			309	309		
4	D	329	Total	O	0	0
			329	329		

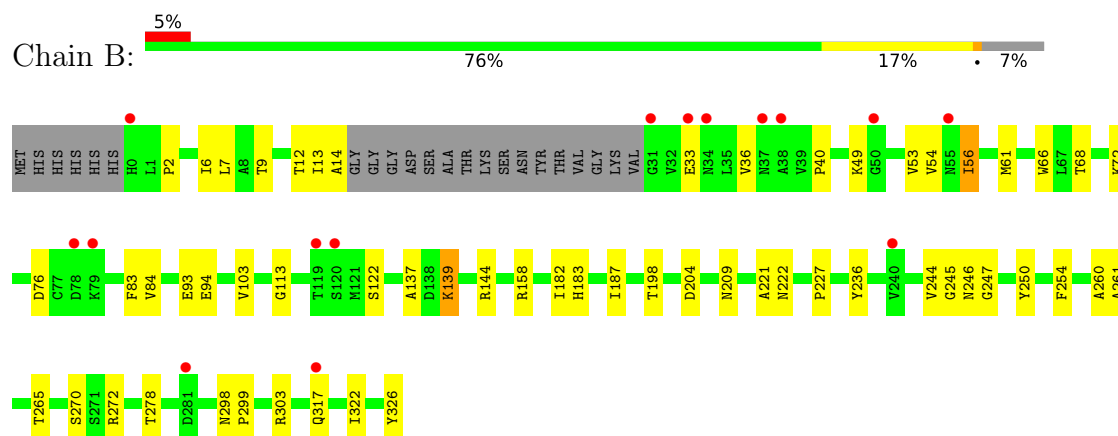
3 Residue-property plots [i](#)

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.

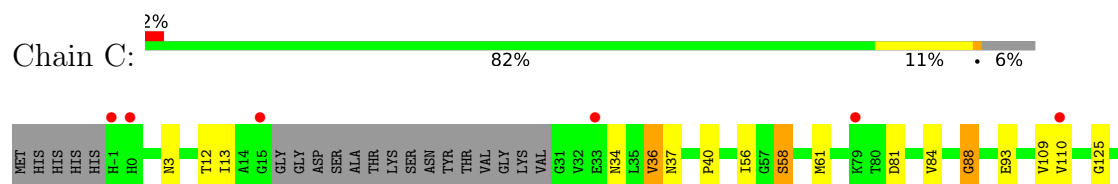
• Molecule 1: L-asparaginase 2

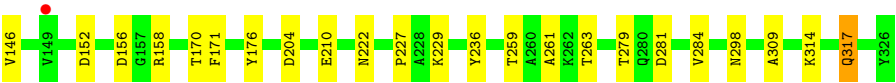


• Molecule 1: L-asparaginase 2

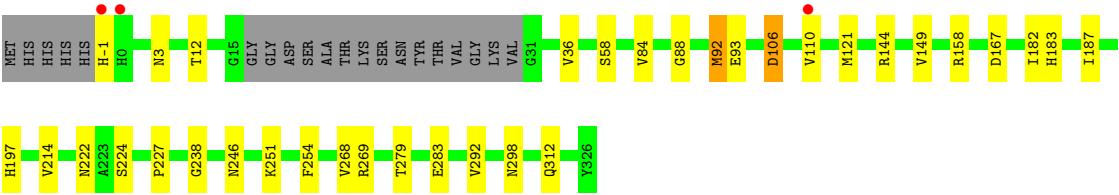
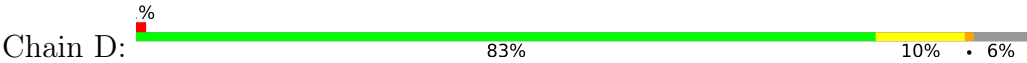


• Molecule 1: L-asparaginase 2





● Molecule 1: L-asparaginase 2



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	151.88Å 62.51Å 143.16Å 90.00° 118.19° 90.00°	Depositor
Resolution (Å)	26.54 – 1.93 26.52 – 1.93	Depositor EDS
% Data completeness (in resolution range)	96.0 (26.54-1.93) 96.0 (26.52-1.93)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.29 (at 1.93Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.152 , 0.227 0.167 , 0.237	Depositor DCC
R_{free} test set	2649 reflections (3.09%)	wwPDB-VP
Wilson B-factor (Å ²)	21.1	Xtriage
Anisotropy	0.051	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 57.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	10658	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.77% 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: ACY, GOL

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	1.11	6/2447 (0.2%)	1.21	5/3334 (0.1%)
1	B	1.14	11/2375 (0.5%)	1.21	10/3237 (0.3%)
1	C	1.10	7/2398 (0.3%)	1.18	5/3269 (0.2%)
1	D	1.14	4/2388 (0.2%)	1.22	8/3254 (0.2%)
All	All	1.12	28/9608 (0.3%)	1.21	28/13094 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	D	0	1
All	All	0	3

All (28) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	210	GLU	CD-OE1	8.34	1.34	1.25
1	A	94	GLU	CD-OE2	7.56	1.33	1.25
1	C	93	GLU	CD-OE2	6.89	1.33	1.25
1	D	93	GLU	CD-OE2	6.32	1.32	1.25
1	A	134	VAL	C-O	6.22	1.35	1.23
1	A	283	GLU	CD-OE1	6.07	1.32	1.25
1	D	121	MET	C-O	5.97	1.34	1.23
1	B	83	PHE	C-O	5.89	1.34	1.23
1	B	326	TYR	C-O	5.75	1.34	1.23
1	B	270	SER	C-O	5.67	1.34	1.23
1	B	94	GLU	CD-OE2	5.67	1.31	1.25
1	B	66	TRP	C-O	5.63	1.34	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	122	SER	C-O	5.59	1.33	1.23
1	B	326	TYR	C-OXT	5.55	1.33	1.23
1	B	221	ALA	C-O	5.53	1.33	1.23
1	D	92	MET	CG-SD	-5.46	1.67	1.81
1	C	309	ALA	C-O	5.42	1.33	1.23
1	A	132	ALA	C-O	5.35	1.33	1.23
1	C	229	LYS	C-O	5.30	1.33	1.23
1	B	245	GLY	C-O	5.25	1.32	1.23
1	C	88	GLY	C-O	5.24	1.32	1.23
1	B	113	GLY	C-O	5.23	1.32	1.23
1	C	58	SER	CB-OG	5.16	1.49	1.42
1	B	93	GLU	C-O	5.14	1.33	1.23
1	C	261	ALA	C-O	5.10	1.33	1.23
1	A	294	SER	C-O	5.09	1.33	1.23
1	C	210	GLU	C-O	5.02	1.32	1.23
1	D	197	HIS	C-O	5.01	1.32	1.23

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	144	ARG	NE-CZ-NH1	8.24	124.42	120.30
1	B	144	ARG	NE-CZ-NH1	7.93	124.26	120.30
1	D	269	ARG	NE-CZ-NH2	-7.78	116.41	120.30
1	B	272	ARG	NE-CZ-NH1	-7.71	116.44	120.30
1	C	204	ASP	CB-CG-OD2	-7.67	111.40	118.30
1	A	176	TYR	CB-CG-CD2	-7.64	116.42	121.00
1	D	144	ARG	NE-CZ-NH2	-7.15	116.72	120.30
1	A	144	ARG	NE-CZ-NH2	-7.09	116.75	120.30
1	D	167	ASP	CB-CG-OD2	-6.91	112.08	118.30
1	B	12	THR	N-CA-CB	6.89	123.39	110.30
1	A	176	TYR	CB-CG-CD1	6.49	124.89	121.00
1	D	158	ARG	NE-CZ-NH2	-6.09	117.26	120.30
1	B	12	THR	CB-CA-C	-6.07	95.22	111.60
1	C	158	ARG	NE-CZ-NH2	-6.02	117.29	120.30
1	C	158	ARG	NE-CZ-NH1	6.00	123.30	120.30
1	A	298	ASN	CB-CA-C	5.91	122.23	110.40
1	B	158	ARG	NE-CZ-NH2	-5.90	117.35	120.30
1	B	68	THR	CA-CB-OG1	-5.86	96.69	109.00
1	B	303	ARG	NE-CZ-NH2	-5.64	117.48	120.30
1	B	250	TYR	CB-CG-CD1	-5.60	117.64	121.00
1	D	269	ARG	NE-CZ-NH1	5.58	123.09	120.30
1	B	14	ALA	CA-C-O	-5.54	108.46	120.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	204	ASP	CB-CG-OD1	5.51	123.26	118.30
1	D	92	MET	CG-SD-CE	5.47	108.95	100.20
1	A	291	PHE	CB-CG-CD2	-5.23	117.14	120.80
1	D	167	ASP	CB-CG-OD1	5.17	122.95	118.30
1	C	298	ASN	CB-CA-C	5.12	120.63	110.40
1	B	139	LYS	CB-CA-C	-5.03	100.34	110.40

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	176	TYR	Mainchain
1	A	46	ALA	Peptide
1	D	106	ASP	Mainchain

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	2401	0	2398	30	0
1	B	2336	0	2327	26	0
1	C	2355	0	2343	26	0
1	D	2348	0	2338	19	0
2	A	4	0	3	1	0
2	B	4	0	3	3	0
2	C	4	0	3	0	0
2	D	4	0	3	2	0
3	B	6	0	8	4	0
3	C	6	0	8	0	0
3	D	6	0	8	1	0
4	A	299	0	0	8	0
4	B	247	0	0	9	0
4	C	309	0	0	8	1
4	D	329	0	0	12	1
All	All	10658	0	9442	103	2

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including

hydrogen atoms). The all-atom clashscore for this structure is 5.

All (103) 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:D:3:ASN:HB3	3:D:402:GOL:H2	1.46	0.97
1:C:317:GLN:NE2	4:C:501:HOH:O	2.02	0.92
1:B:209:ASN:O	4:B:601:HOH:O	1.89	0.90
4:A:562:HOH:O	3:B:501:GOL:H32	1.70	0.90
1:C:279:THR:HG22	4:C:803:HOH:O	1.74	0.88
1:A:248:ASN:ND2	1:A:283:GLU:HB3	1.92	0.85
1:A:248:ASN:HD21	1:A:283:GLU:HB3	1.42	0.84
1:A:279:THR:HG23	4:A:635:HOH:O	1.81	0.80
2:A:401:ACY:H1	4:A:690:HOH:O	1.80	0.79
1:C:227:PRO:HB3	1:D:227:PRO:HB3	1.67	0.76
1:A:225:ASP:HB3	1:A:252:SER:OG	1.86	0.75
1:A:279:THR:CG2	4:A:635:HOH:O	2.34	0.74
1:D:279:THR:HG23	4:D:671:HOH:O	1.88	0.71
1:C:279:THR:HG23	4:C:608:HOH:O	1.95	0.67
2:D:401:ACY:CH3	4:D:768:HOH:O	2.42	0.66
1:C:34:ASN:ND2	1:C:37:ASN:HD22	1.94	0.66
1:D:312:GLN:NE2	4:D:503:HOH:O	2.22	0.66
1:D:279:THR:HG22	4:D:826:HOH:O	1.96	0.66
1:A:104:LYS:HE2	4:A:515:HOH:O	1.94	0.65
1:C:58:SER:HB3	1:C:88:GLY:HA3	1.80	0.63
1:B:246:ASN:H	3:B:501:GOL:H31	1.63	0.62
1:A:104:LYS:CE	4:A:515:HOH:O	2.48	0.62
1:D:183:HIS:HD2	4:D:659:HOH:O	1.84	0.60
1:B:246:ASN:N	3:B:501:GOL:H31	2.17	0.60
2:D:401:ACY:H1	4:D:768:HOH:O	2.01	0.60
1:A:13:ILE:O	1:A:125:GLY:HA3	2.02	0.59
1:D:279:THR:CG2	4:D:671:HOH:O	2.46	0.59
1:B:13:ILE:HA	4:B:762:HOH:O	2.02	0.58
1:A:159:ASP:HB3	1:A:173:SER:HB3	1.84	0.58
1:C:84:VAL:HA	1:C:110:VAL:O	2.03	0.58
1:C:170:THR:HG23	1:C:171:PHE:N	2.17	0.58
1:D:106:ASP:N	4:D:501:HOH:O	1.76	0.56
1:C:12:THR:HG22	4:C:771:HOH:O	2.05	0.56
1:A:56:ILE:HD11	1:A:61:MET:HE2	1.89	0.54
2:B:502:ACY:H3	4:B:842:HOH:O	2.09	0.53
1:B:322:ILE:HD11	4:B:813:HOH:O	2.07	0.53
1:A:67:LEU:O	1:A:71:LYS:HG3	2.09	0.52
1:B:246:ASN:H	3:B:501:GOL:C3	2.22	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:12:THR:HG22	4:D:786:HOH:O	2.10	0.52
1:C:259:THR:O	1:C:263:THR:HG23	2.09	0.52
1:C:281:ASP:HA	1:C:284:VAL:O	2.10	0.52
1:B:182:ILE:HG12	1:B:187:ILE:HG12	1.91	0.52
1:A:27:VAL:HG13	1:A:34:ASN:ND2	2.26	0.51
1:B:72:LYS:NZ	1:B:76:ASP:OD2	2.43	0.51
1:A:76:ASP:HA	1:A:79:LYS:HD3	1.92	0.51
1:A:104:LYS:NZ	4:A:515:HOH:O	2.44	0.51
1:D:182:ILE:HG12	1:D:187:ILE:HG12	1.92	0.51
1:D:246:ASN:HB3	1:D:283:GLU:HG3	1.93	0.50
1:A:90:THR:HB	1:B:244:VAL:HB	1.92	0.50
1:A:174:VAL:HB	1:A:275:THR:HG22	1.94	0.50
1:A:227:PRO:HB3	1:B:227:PRO:HB3	1.94	0.50
1:A:6:ILE:HA	1:A:84:VAL:O	2.13	0.49
1:A:281:ASP:HA	1:A:284:VAL:O	2.13	0.49
1:B:56:ILE:CD1	1:B:61:MET:HE2	2.43	0.49
1:A:214:VAL:HA	1:A:238:GLY:O	2.13	0.49
1:A:248:ASN:ND2	1:A:283:GLU:CB	2.72	0.49
1:B:139:LYS:HG3	4:B:731:HOH:O	2.11	0.49
1:C:36:VAL:O	1:C:40:PRO:HA	2.14	0.48
1:B:103:VAL:O	1:B:198:THR:HA	2.14	0.47
1:A:2:PRO:HG3	1:A:137:ALA:HB1	1.96	0.47
1:B:33:GLU:HG2	4:B:705:HOH:O	2.14	0.46
1:C:13:ILE:O	1:C:125:GLY:HA3	2.16	0.46
1:D:58:SER:HB3	1:D:88:GLY:HA3	1.98	0.46
2:B:502:ACY:CH3	4:B:726:HOH:O	2.64	0.46
1:B:2:PRO:HG2	1:B:137:ALA:HB1	1.97	0.46
1:A:224:SER:HB2	1:B:236:TYR:OH	2.17	0.45
1:C:34:ASN:HA	1:C:37:ASN:ND2	2.31	0.45
1:B:9:THR:HA	1:B:54:VAL:HG23	1.98	0.45
1:C:34:ASN:ND2	1:C:37:ASN:ND2	2.64	0.45
1:B:260:ALA:HB1	1:B:265:THR:HB	1.99	0.45
1:C:152:ASP:OD2	4:C:503:HOH:O	2.21	0.45
1:C:279:THR:CG2	4:C:608:HOH:O	2.60	0.45
1:A:183:HIS:O	1:A:184:ASN:C	2.55	0.44
1:C:170:THR:CG2	1:C:171:PHE:N	2.80	0.44
1:C:56:ILE:HD11	1:C:61:MET:HE1	1.99	0.44
1:D:214:VAL:HA	1:D:238:GLY:O	2.18	0.44
1:B:36:VAL:O	1:B:40:PRO:HA	2.18	0.44
1:B:183:HIS:HD2	4:B:796:HOH:O	1.99	0.44
1:C:176:TYR:OH	4:C:502:HOH:O	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:247:GLY:HA3	1:B:278:THR:HG23	1.99	0.43
1:A:226:LEU:HB2	1:A:227:PRO:HD3	2.01	0.43
1:C:236:TYR:OH	1:D:224:SER:HB2	2.19	0.42
1:B:261:ALA:HA	1:B:265:THR:O	2.18	0.42
1:A:248:ASN:ND2	4:A:531:HOH:O	2.52	0.42
1:A:220:TYR:CD1	1:A:220:TYR:N	2.86	0.42
1:C:34:ASN:HD22	1:C:37:ASN:HD22	1.66	0.41
1:D:149:VAL:HG12	4:D:564:HOH:O	2.21	0.41
2:B:502:ACY:H2	4:B:726:HOH:O	2.21	0.41
1:D:-1:HIS:N	4:D:535:HOH:O	2.52	0.41
1:C:314:LYS:HE3	4:C:743:HOH:O	2.21	0.41
1:D:84:VAL:HA	1:D:110:VAL:O	2.20	0.41
1:D:268:VAL:HG22	1:D:292:VAL:HB	2.03	0.41
1:B:204:ASP:C	1:B:204:ASP:OD1	2.58	0.41
1:A:103:VAL:O	1:A:198:THR:HA	2.21	0.41
1:B:6:ILE:HA	1:B:84:VAL:O	2.21	0.41
1:A:325:GLN:HG3	1:A:326:TYR:CE2	2.56	0.40
1:B:7:LEU:HD22	1:B:53:VAL:CG2	2.51	0.40
1:C:3:ASN:O	1:C:81:ASP:HB2	2.21	0.40
1:C:34:ASN:HD22	1:C:37:ASN:ND2	2.19	0.40
1:A:84:VAL:HA	1:A:110:VAL:O	2.21	0.40
1:B:298:ASN:HB2	1:B:299:PRO:CD	2.51	0.40
1:C:109:VAL:HB	1:C:146:VAL:HG22	2.04	0.40
1:D:298:ASN:HA	4:D:693:HOH:O	2.21	0.40

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:619:HOH:O	4:D:619:HOH:O[2_556]	0.62	1.58
4:C:654:HOH:O	4:C:654:HOH:O[2_556]	0.82	1.38

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	318/333 (96%)	309 (97%)	8 (2%)	1 (0%)	41	32
1	B	308/333 (92%)	297 (96%)	11 (4%)	0	100	100
1	C	311/333 (93%)	302 (97%)	9 (3%)	0	100	100
1	D	310/333 (93%)	304 (98%)	6 (2%)	0	100	100
All	All	1247/1332 (94%)	1212 (97%)	34 (3%)	1 (0%)	51	43

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	281	ASP

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	265/273 (97%)	256 (97%)	9 (3%)	37	22
1	B	257/273 (94%)	252 (98%)	5 (2%)	57	45
1	C	259/273 (95%)	255 (98%)	4 (2%)	65	56
1	D	258/273 (94%)	253 (98%)	5 (2%)	57	45
All	All	1039/1092 (95%)	1016 (98%)	23 (2%)	52	39

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	0	HIS
1	A	12	THR
1	A	139	LYS
1	A	207	LYS
1	A	222	ASN
1	A	248	ASN
1	A	252	SER

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Mol	Chain	Res	Type
1	A	254	PHE
1	A	321	GLN
1	B	49	LYS
1	B	56	ILE
1	B	222	ASN
1	B	254	PHE
1	B	317	GLN
1	C	36	VAL
1	C	156	ASP
1	C	222	ASN
1	C	317	GLN
1	D	36	VAL
1	D	92	MET
1	D	222	ASN
1	D	251	LYS
1	D	254	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	280	GLN
1	B	64	ASN
1	B	280	GLN
1	B	324	ASN
1	C	34	ASN
1	C	64	ASN
1	C	183	HIS
1	C	209	ASN
1	C	317	GLN
1	D	64	ASN
1	D	143	ASN
1	D	183	HIS
1	D	318	GLN
1	D	324	ASN

5.3.3 RNA ⓘ

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

7 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	GOL	B	501	-	5,5,5	0.19	0	5,5,5	0.38	0
2	ACY	A	401	-	3,3,3	0.70	0	3,3,3	0.87	0
3	GOL	C	402	-	5,5,5	0.18	0	5,5,5	0.63	0
3	GOL	D	402	-	5,5,5	0.37	0	5,5,5	0.60	0
2	ACY	D	401	-	3,3,3	1.28	0	3,3,3	1.22	0
2	ACY	C	401	-	3,3,3	1.01	0	3,3,3	0.90	0
2	ACY	B	502	-	3,3,3	0.97	0	3,3,3	0.64	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
3	GOL	C	402	-	-	2/4/4/4	-
3	GOL	B	501	-	-	2/4/4/4	-
3	GOL	D	402	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (6) torsion outliers are listed below:

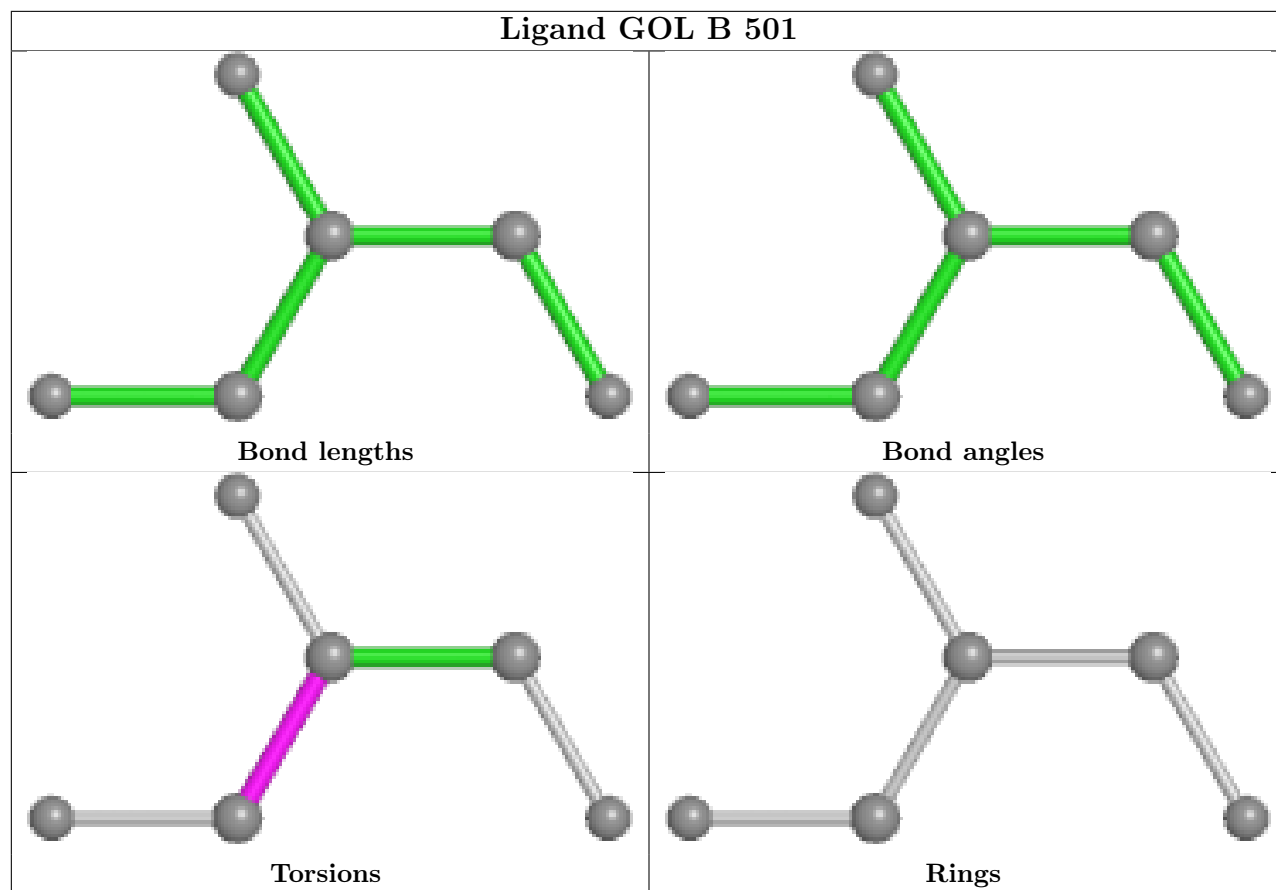
Mol	Chain	Res	Type	Atoms
3	C	402	GOL	C1-C2-C3-O3
3	D	402	GOL	O1-C1-C2-O2
3	B	501	GOL	C1-C2-C3-O3
3	D	402	GOL	O1-C1-C2-C3
3	C	402	GOL	O2-C2-C3-O3
3	B	501	GOL	O2-C2-C3-O3

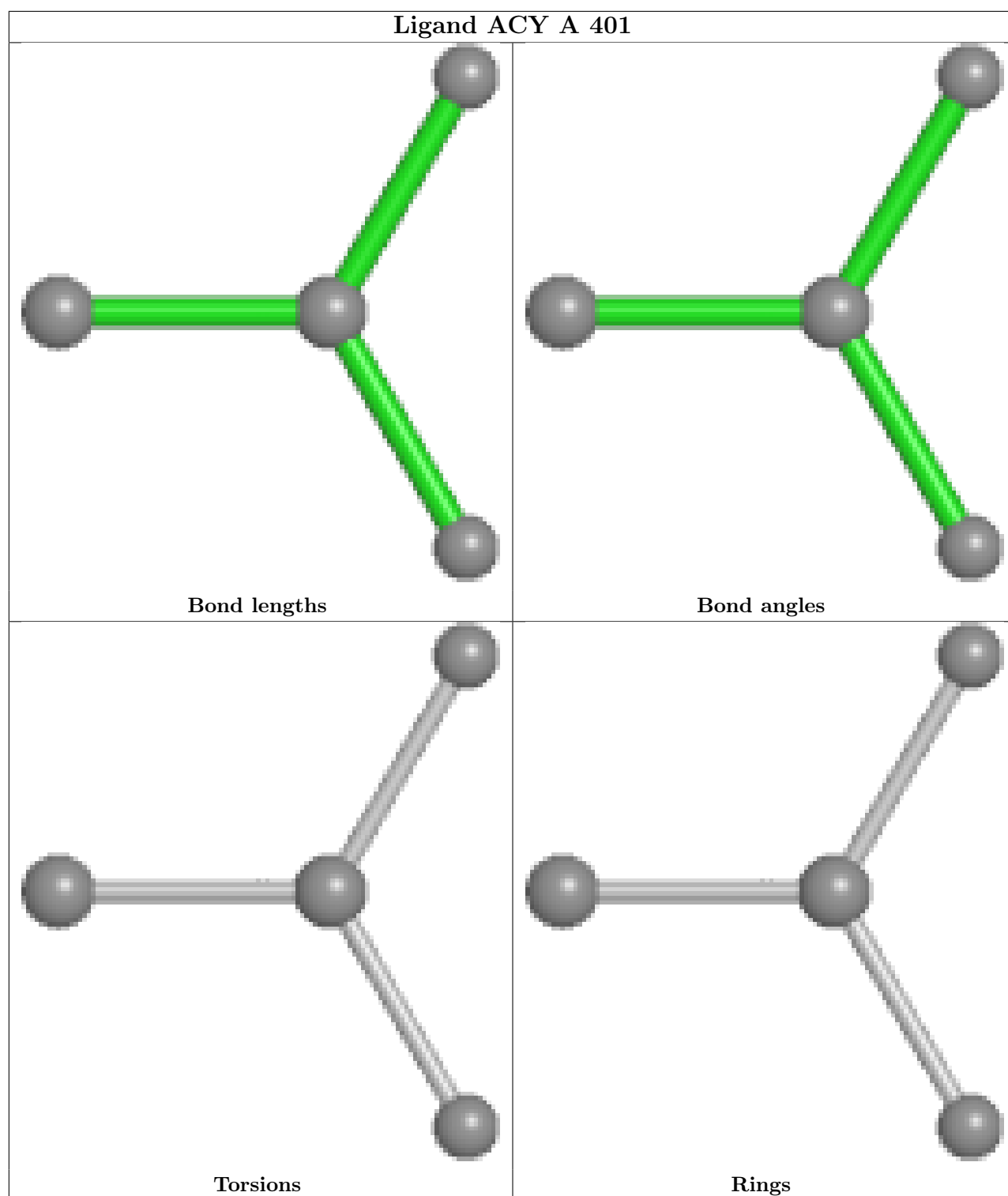
There are no ring outliers.

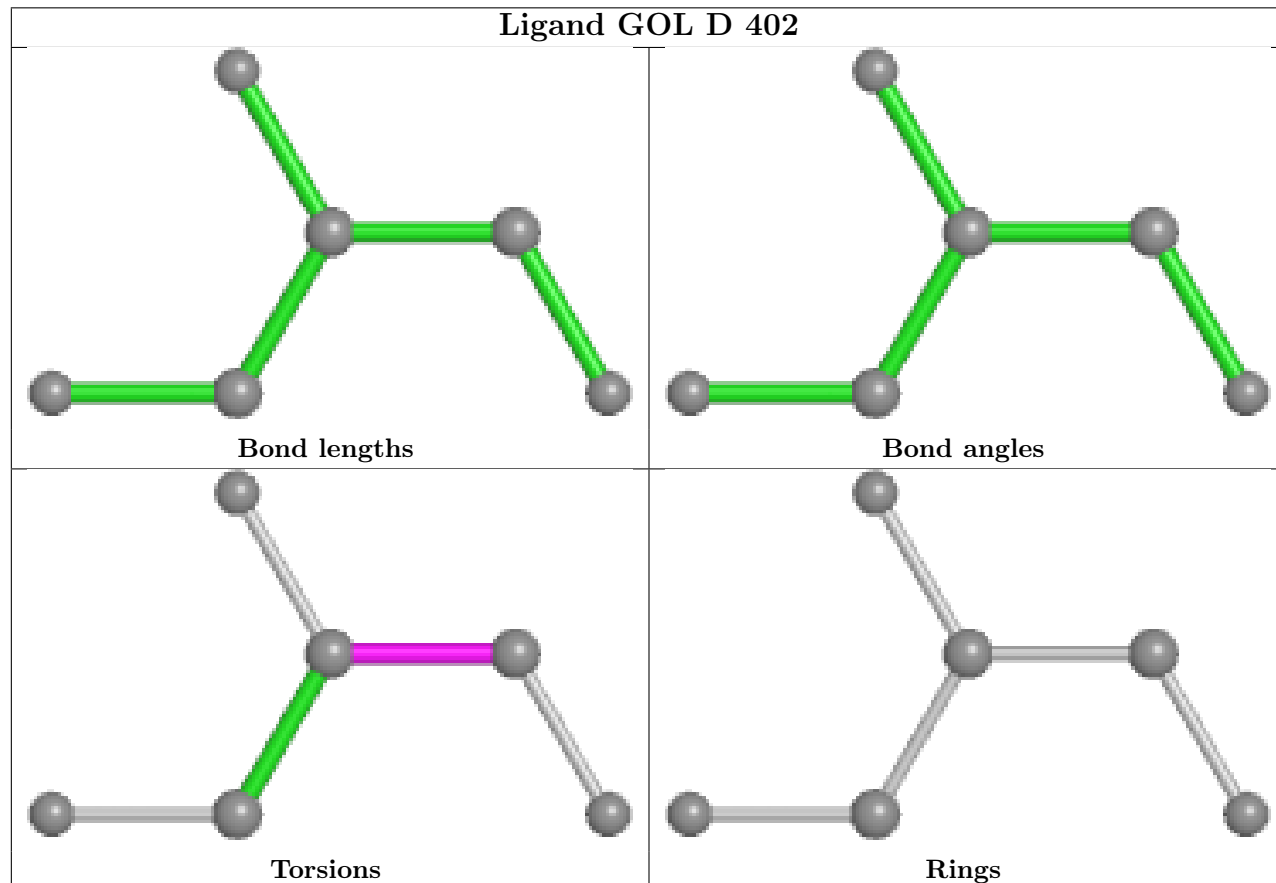
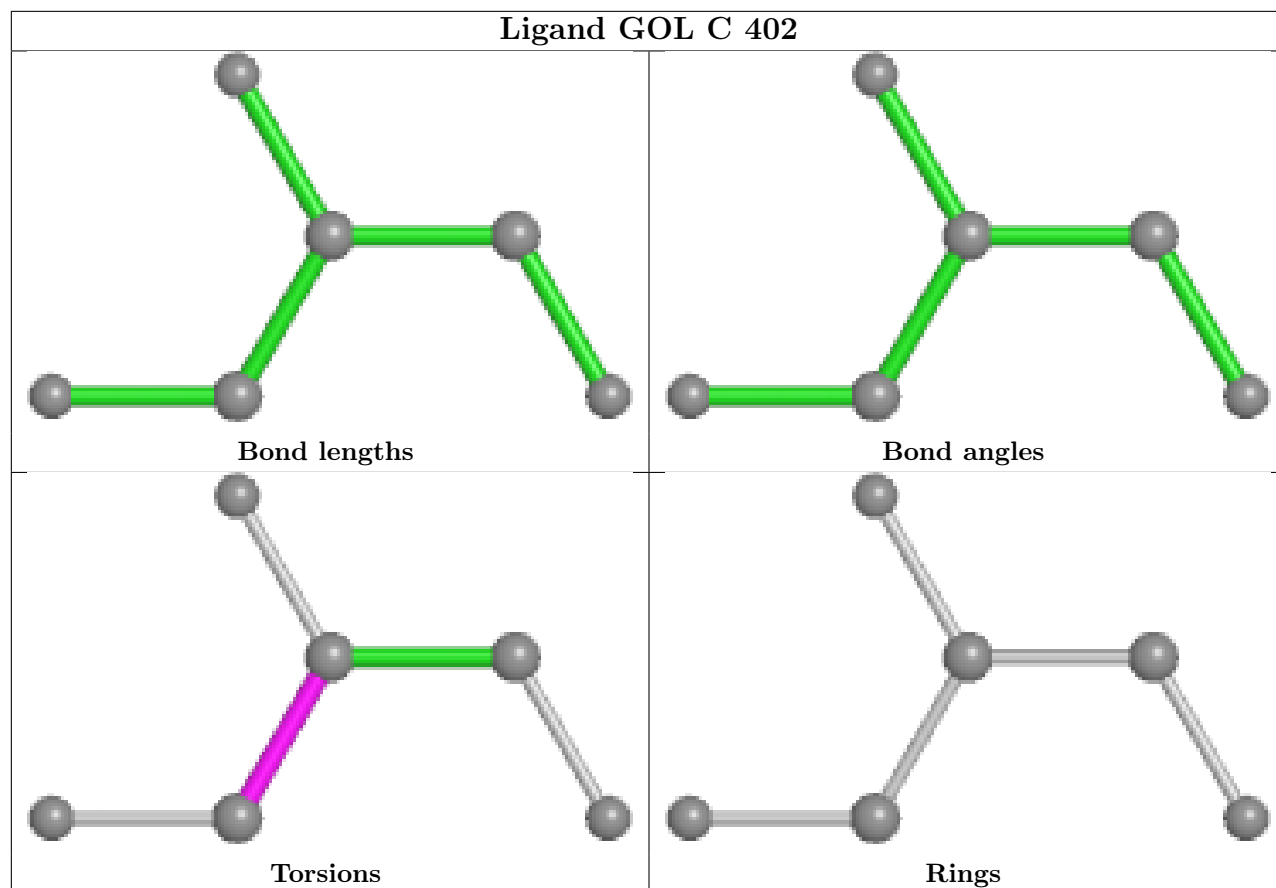
5 monomers are involved in 11 short contacts:

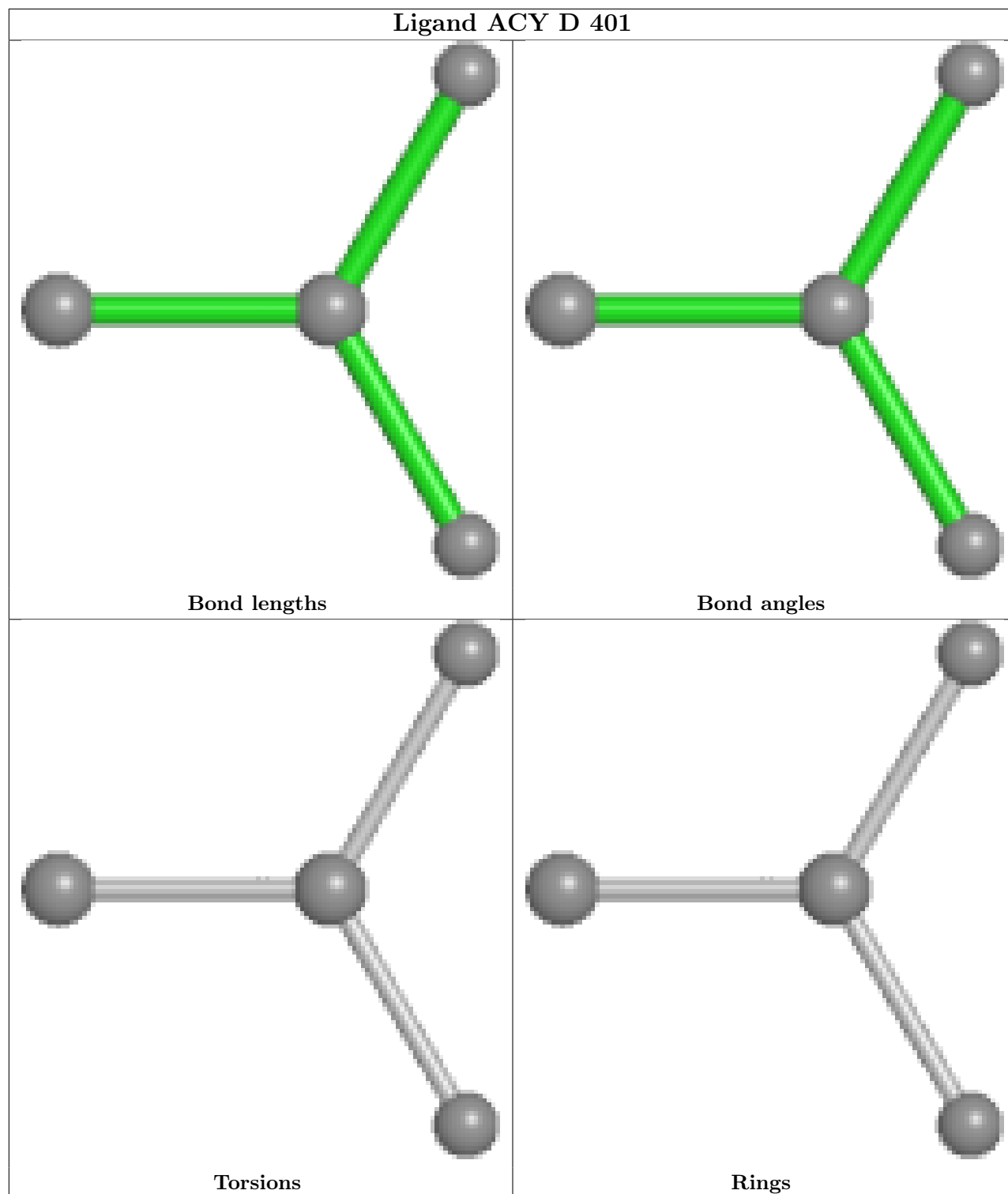
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	501	GOL	4	0
2	A	401	ACY	1	0
3	D	402	GOL	1	0
2	D	401	ACY	2	0
2	B	502	ACY	3	0

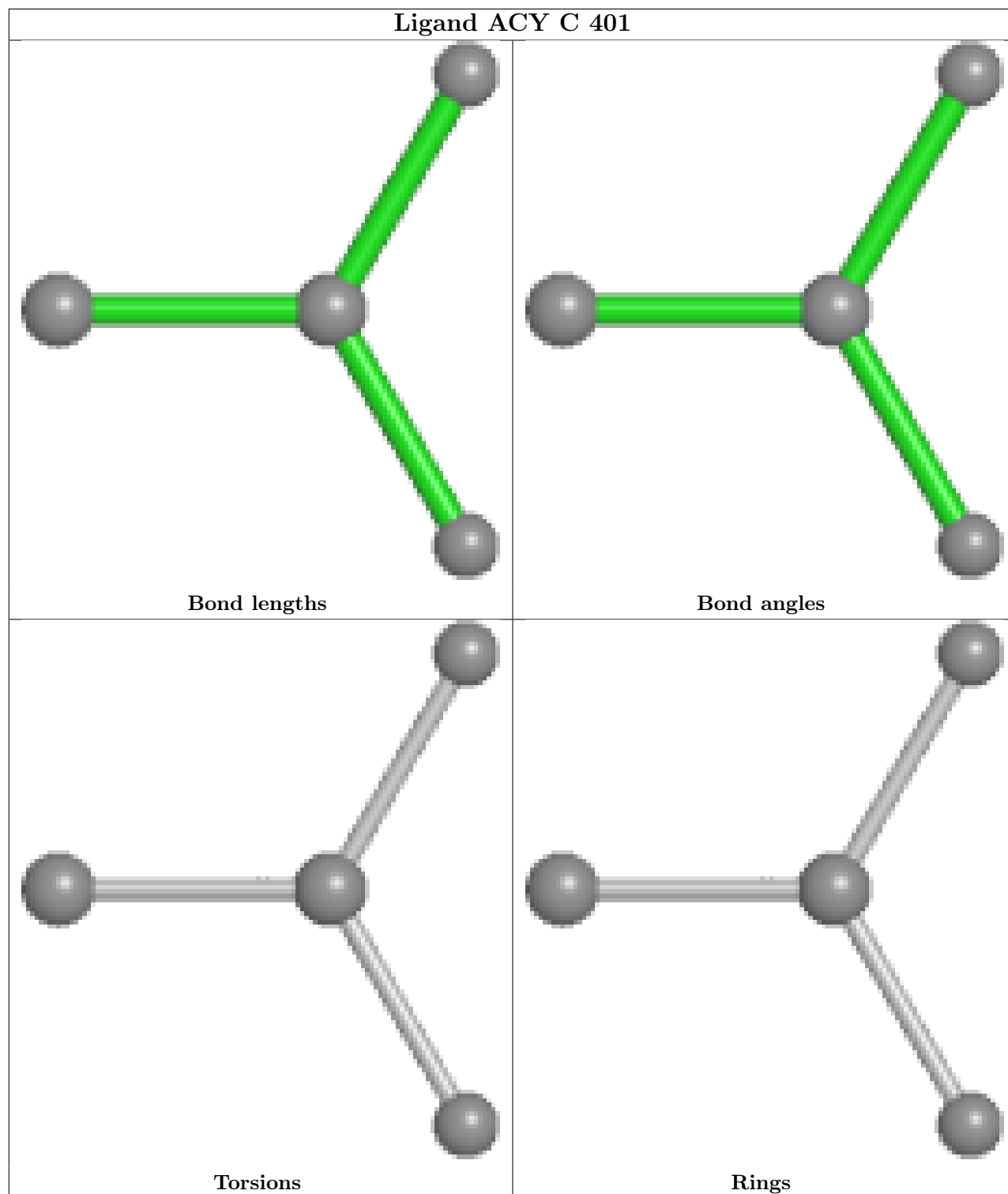
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

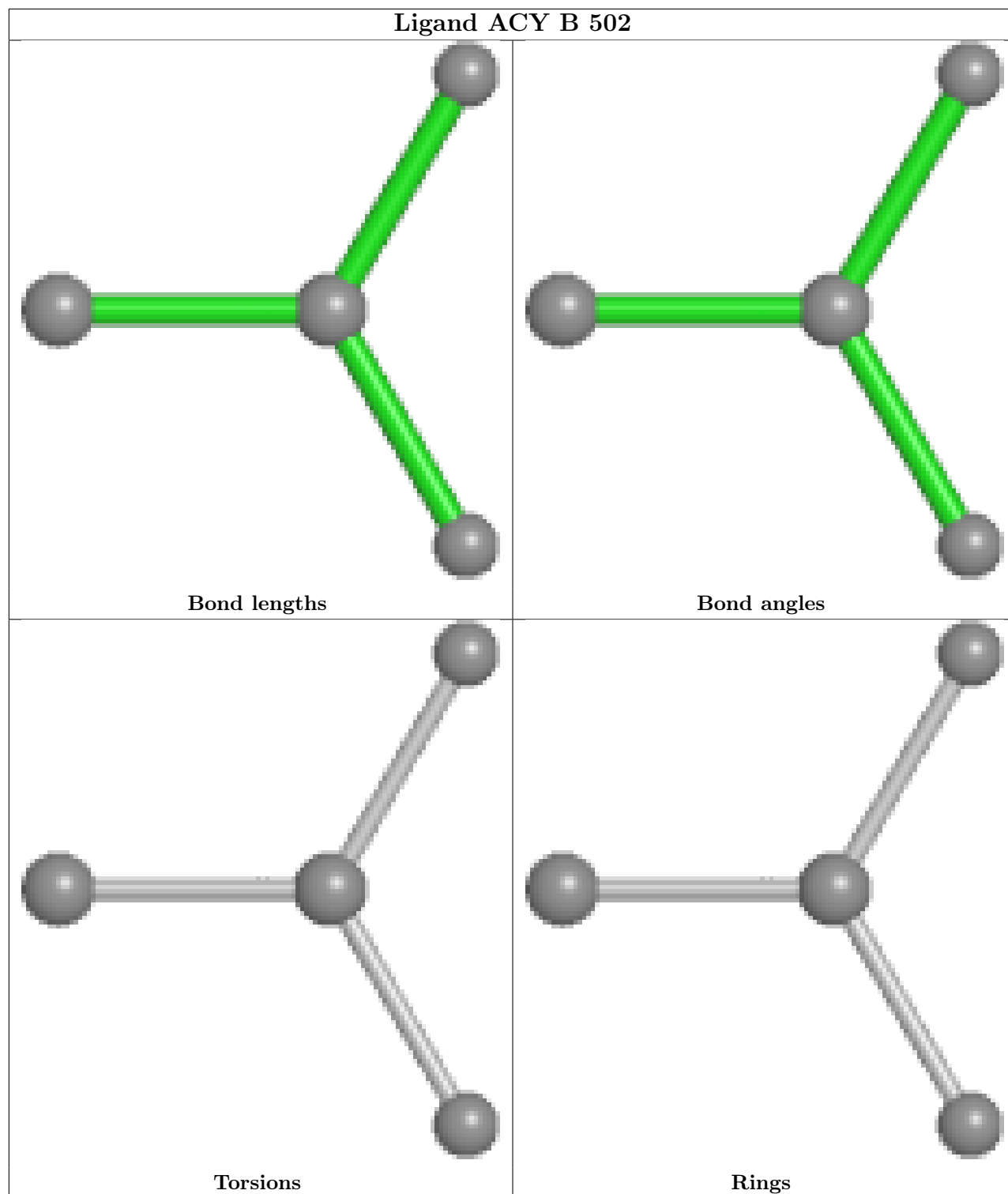












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	319/333 (95%)	0.25	18 (5%) 24 31	27, 34, 53, 87	0
1	B	311/333 (93%)	0.36	15 (4%) 30 38	28, 38, 56, 78	0
1	C	313/333 (93%)	0.09	7 (2%) 62 69	24, 31, 49, 86	0
1	D	313/333 (93%)	0.07	3 (0%) 82 86	23, 30, 44, 84	0
All	All	1256/1332 (94%)	0.19	43 (3%) 45 53	23, 34, 52, 87	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	-1	HIS	7.3
1	D	-1	HIS	5.4
1	A	28	GLY	5.3
1	A	29	LYS	4.7
1	A	0	HIS	4.6
1	D	0	HIS	4.2
1	A	106	ASP	3.4
1	A	30	VAL	3.4
1	B	38	ALA	3.4
1	B	0	HIS	3.4
1	B	281	ASP	3.1
1	A	25	TYR	3.1
1	B	120	SER	3.1
1	C	110	VAL	3.1
1	B	78[A]	ASP	3.0
1	B	79	LYS	3.0
1	B	55	ASN	2.9
1	B	34	ASN	2.7
1	A	31	GLY	2.7
1	B	119	THR	2.6
1	A	110	VAL	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	31	GLY	2.5
1	C	0	HIS	2.5
1	A	15	GLY	2.5
1	A	317	GLN	2.4
1	C	149	VAL	2.4
1	A	235	GLY	2.4
1	A	78	ASP	2.4
1	B	33	GLU	2.3
1	A	24	ASN	2.3
1	A	281	ASP	2.2
1	C	33	GLU	2.2
1	B	240	VAL	2.2
1	A	27	VAL	2.2
1	A	79	LYS	2.1
1	B	37	ASN	2.1
1	A	284	VAL	2.1
1	B	317	GLN	2.1
1	B	50	GLY	2.1
1	D	110	VAL	2.1
1	C	15	GLY	2.0
1	C	79	LYS	2.0
1	A	85	ILE	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 monosaccharides 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	GOL	D	402	6/6	0.89	0.21	41,70,85,88	0

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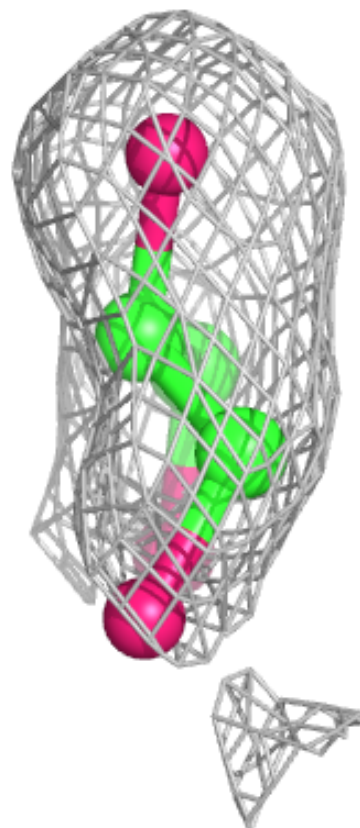
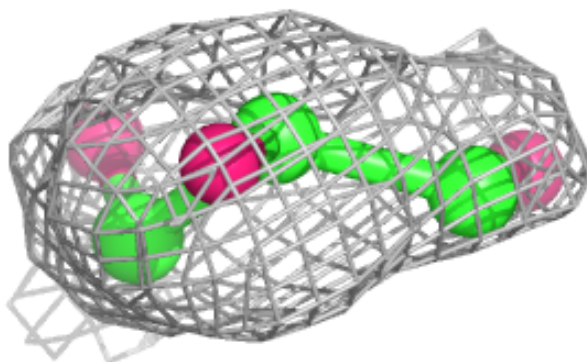
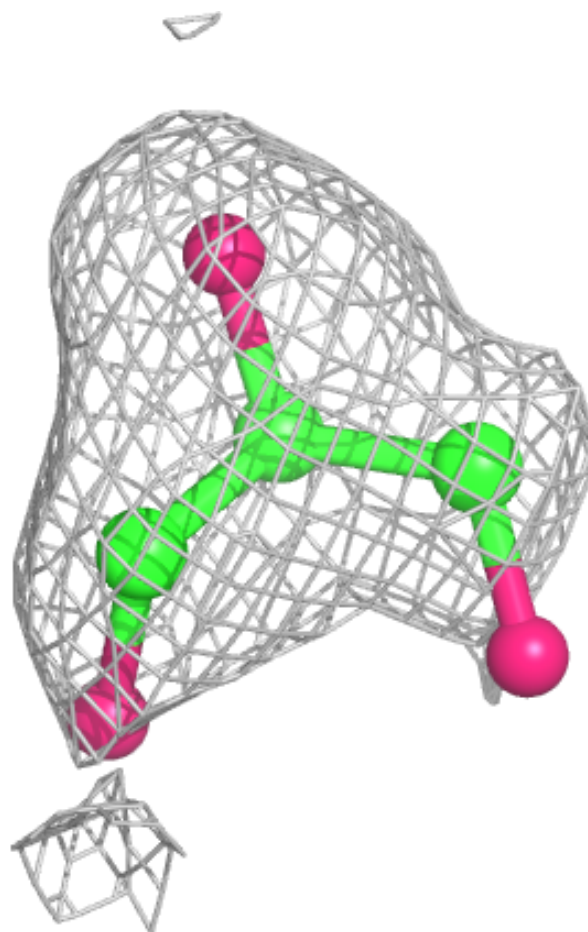
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	GOL	C	402	6/6	0.90	0.20	49,60,61,65	0
3	GOL	B	501	6/6	0.90	0.18	38,46,51,56	0
2	ACY	C	401	4/4	0.94	0.09	42,44,45,46	0
2	ACY	D	401	4/4	0.95	0.10	35,51,51,52	0
2	ACY	A	401	4/4	0.95	0.10	36,38,45,46	0
2	ACY	B	502	4/4	0.96	0.09	36,55,57,69	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.

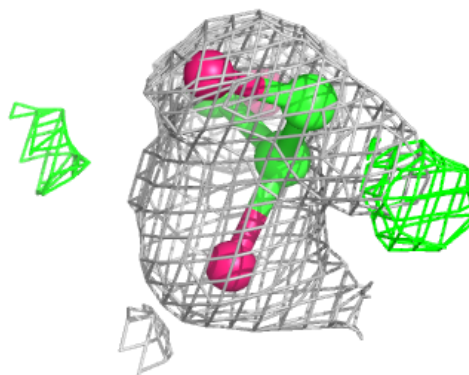
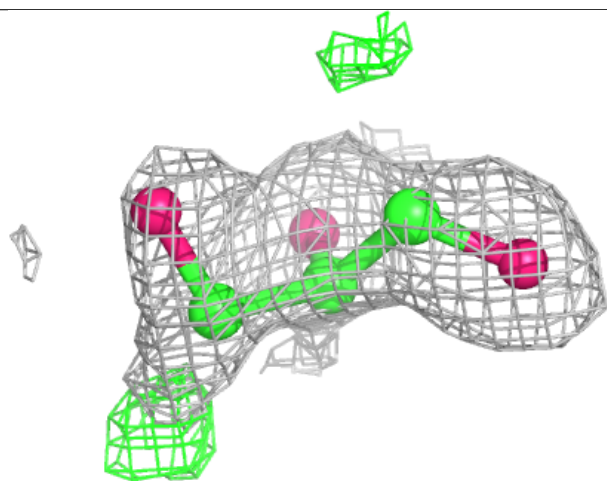
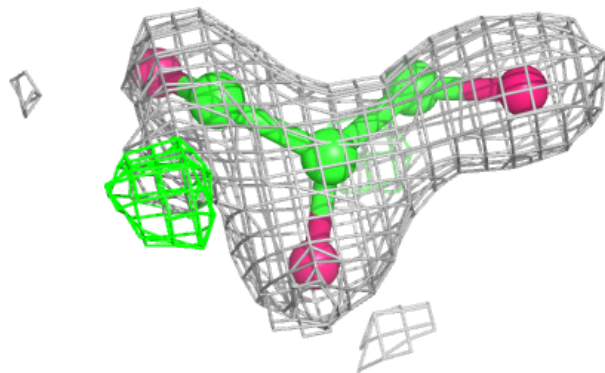
Electron density around GOL D 402:

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



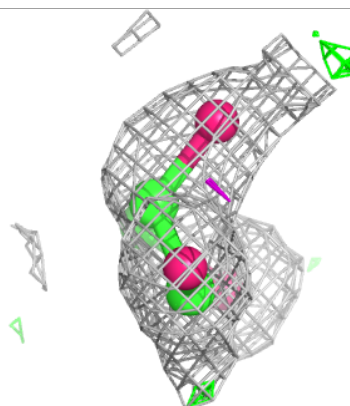
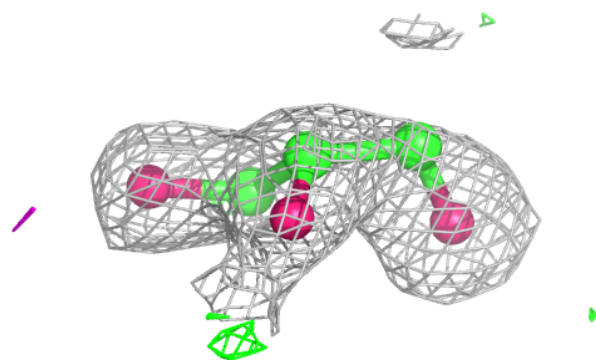
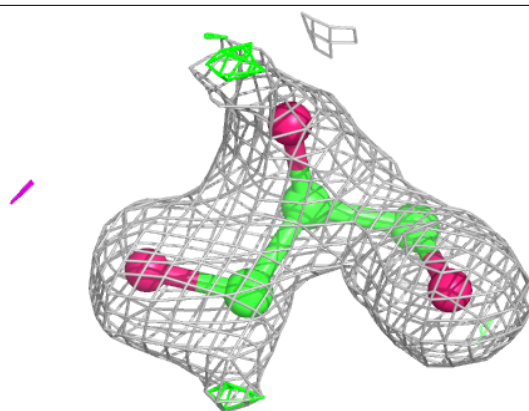
Electron density around GOL C 402:

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



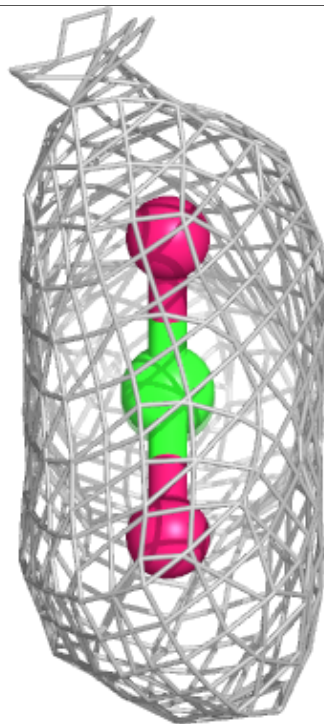
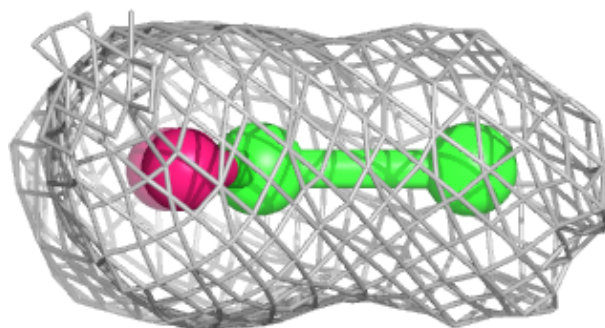
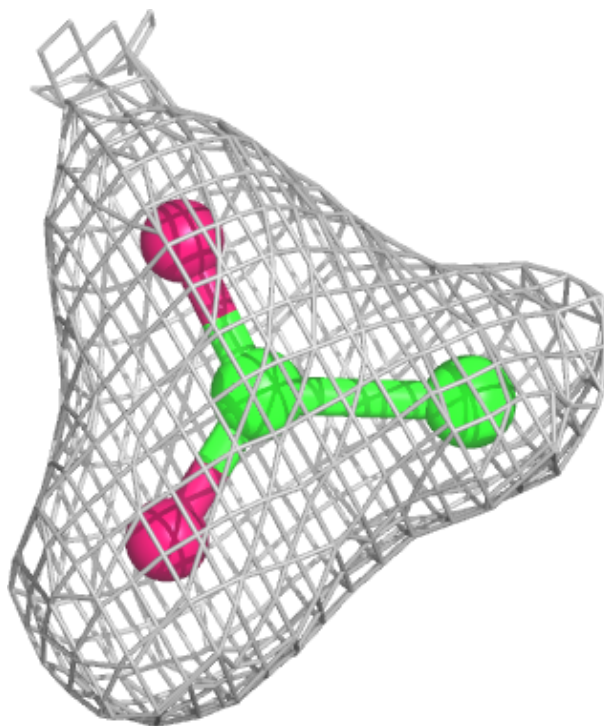
Electron density around GOL B 501:

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



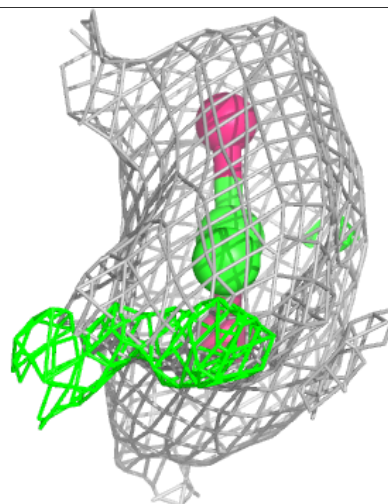
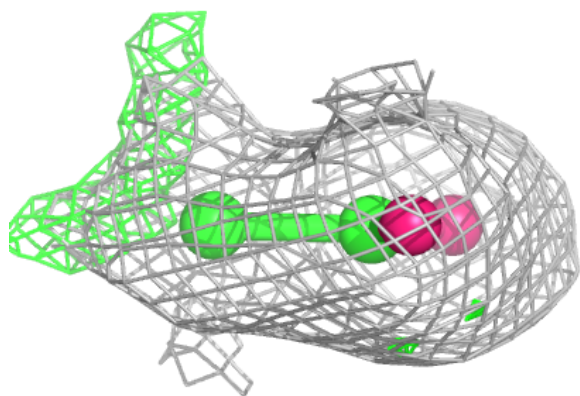
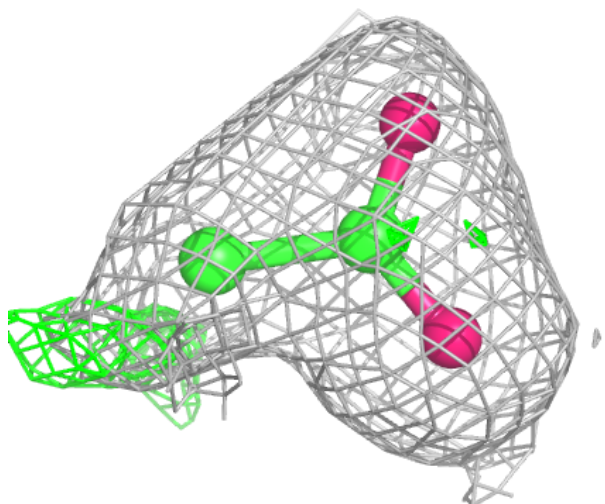
Electron density around ACY C 401:

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



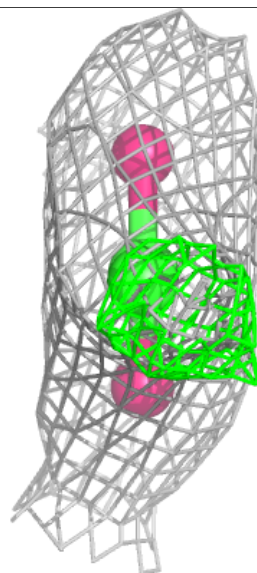
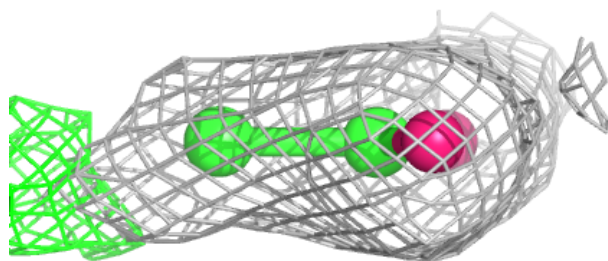
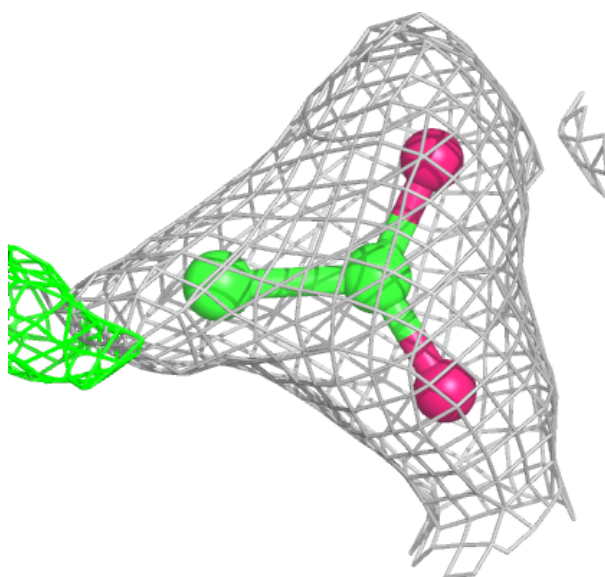
Electron density around ACY D 401:

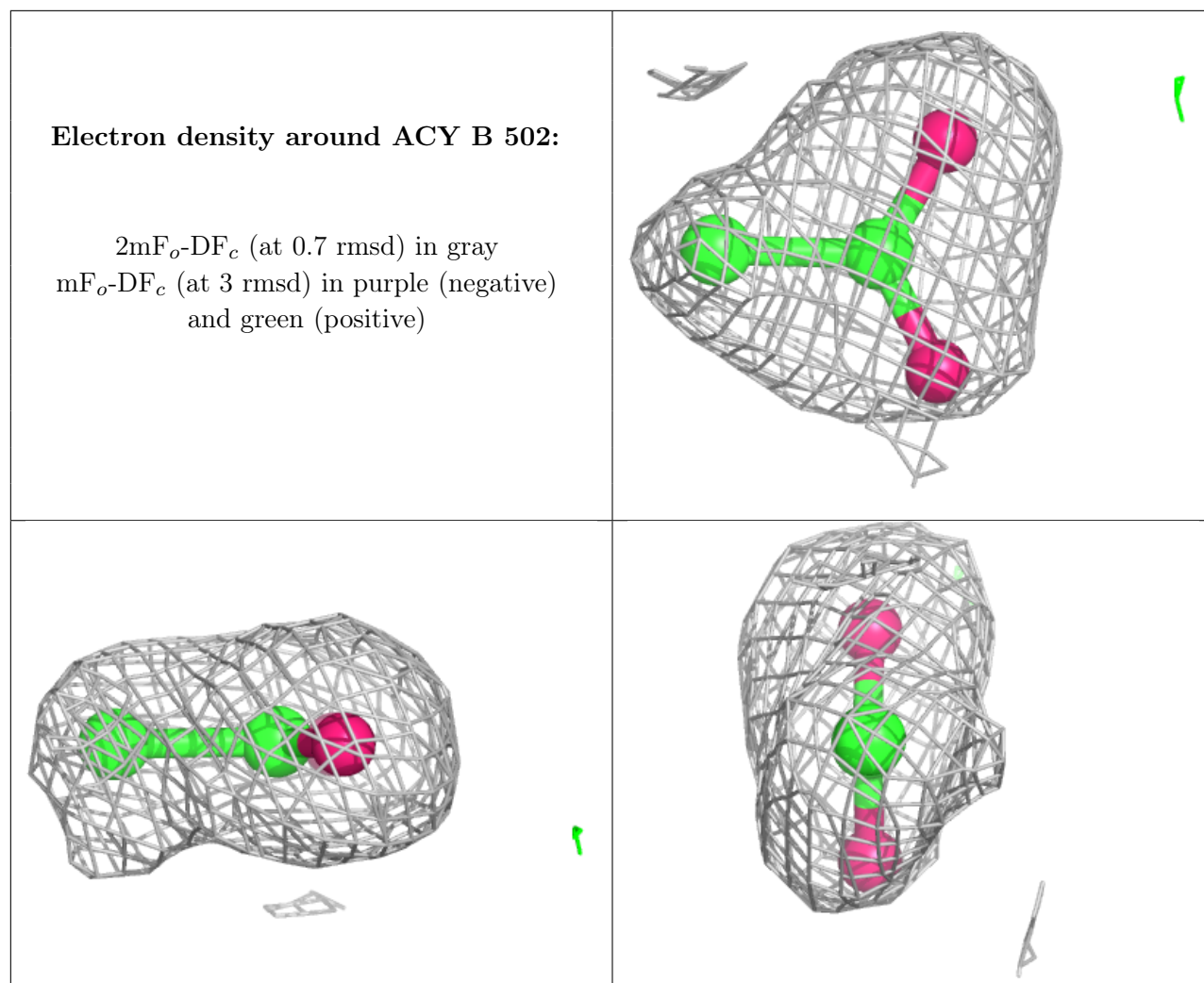
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around ACY A 401:

$2mF_o-DF_c$ (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.