



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 1, 2025 – 04:17 PM JST

PDB ID : 9W2K / pdb_00009w2k
Title : Structural basis of substrate promiscuity in the archaeal RNA-splicing endonuclease from *Candidatus Micrarchaeum acidiphilum* (ARMAN-2)
Authors : Hirata, A.
Deposited on : 2025-07-28
Resolution : 1.80 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 Phenix2.0rc1
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 2.0rc1
EDS : 3.0
buster-report : 1.1.7 (2018)
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4 : 9.0.006 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.45.1

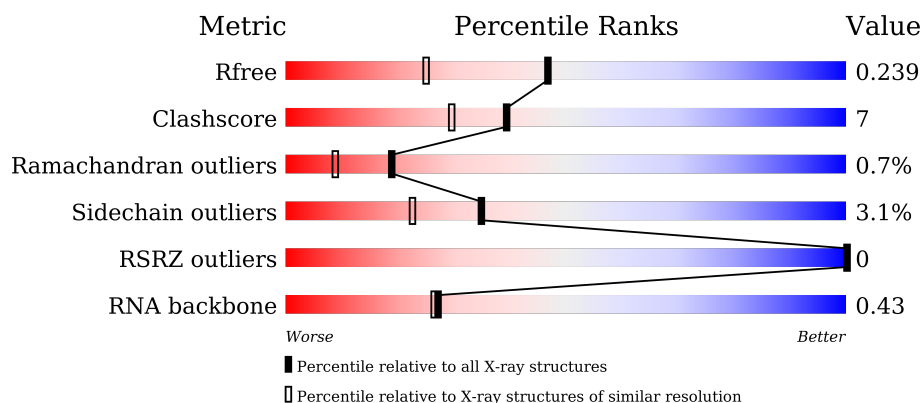
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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}	164625	7108 (1.80-1.80)
Clashscore	180529	8162 (1.80-1.80)
Ramachandran outliers	177936	8077 (1.80-1.80)
Sidechain outliers	177891	8076 (1.80-1.80)
RSRZ outliers	164620	7108 (1.80-1.80)
RNA backbone	3690	1009 (2.24-1.36)

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	395	
1	B	395	
2	I	13	
3	F	8	

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Mol	Chain	Length	Quality of chain
4	H	4	<div>25% 25% 50%</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
5	ACY	A	403	-	-	X	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 7209 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 tRNA intron endonuclease.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	386	Total	C	N	O	S	0	0	0
			3152	2027	545	572	8			
1	B	384	Total	C	N	O	S	0	0	0
			3136	2017	542	569	8			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	388	LEU	-	expression tag	UNP C7DIA5
A	389	GLU	-	expression tag	UNP C7DIA5
A	390	HIS	-	expression tag	UNP C7DIA5
A	391	HIS	-	expression tag	UNP C7DIA5
A	392	HIS	-	expression tag	UNP C7DIA5
A	393	HIS	-	expression tag	UNP C7DIA5
A	394	HIS	-	expression tag	UNP C7DIA5
A	395	HIS	-	expression tag	UNP C7DIA5
B	388	LEU	-	expression tag	UNP C7DIA5
B	389	GLU	-	expression tag	UNP C7DIA5
B	390	HIS	-	expression tag	UNP C7DIA5
B	391	HIS	-	expression tag	UNP C7DIA5
B	392	HIS	-	expression tag	UNP C7DIA5
B	393	HIS	-	expression tag	UNP C7DIA5
B	394	HIS	-	expression tag	UNP C7DIA5
B	395	HIS	-	expression tag	UNP C7DIA5

- Molecule 2 is a RNA chain called DNA/RNA (5'-R(P*AP*CP*CP*GP*AP*CP*CP*A)-D(P*U)-R(P*AP*GP*CP*U)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	I	13	Total	C	N	O	P	0	0	0
			273	123	49	88	13			

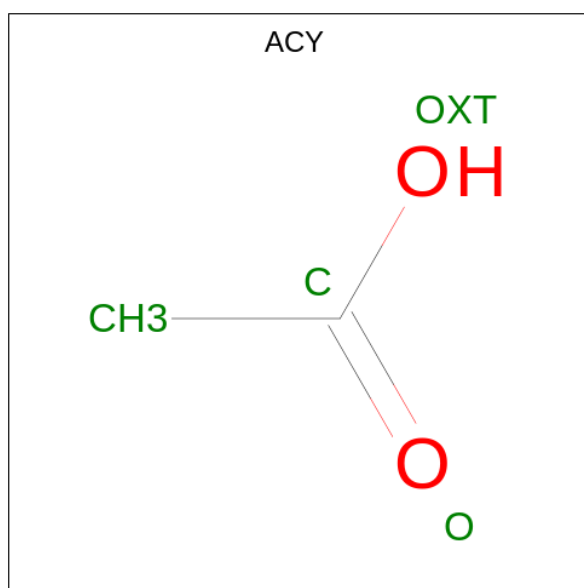
- Molecule 3 is a RNA chain called RNA (5'-R(P*AP*GP*CP*GP*GP*UP*CP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	F	8	Total	C	N	O	P	0	0	0
			173	77	33	55	8			

- Molecule 4 is a RNA chain called RNA (5'-R(*AP*GP*GP*U)-3').

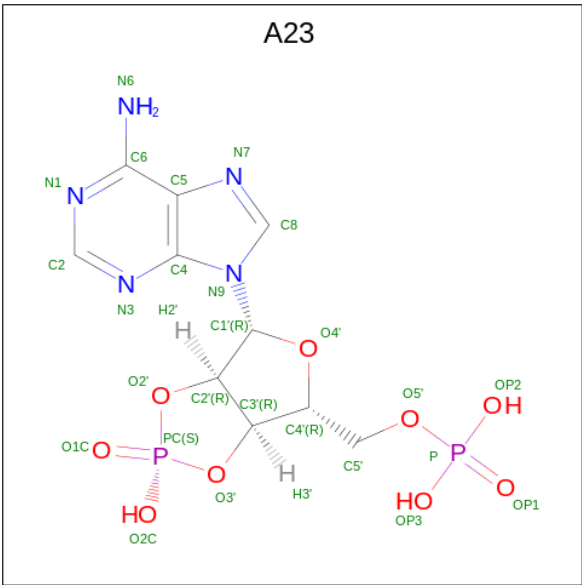
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	H	4	Total	C	N	O	P	0	0	0
			85	39	17	26	3			

- Molecule 5 is ACETIC ACID (CCD ID: ACY) (formula: C₂H₄O₂) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 6 is ADENOSINE-5'-PHOSPHATE-2',3'-CYCLIC PHOSPHATE (CCD ID: A23) (formula: C₁₀H₁₃N₅O₉P₂) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	F	1	Total	C	N	O	P	0	0
			25	10	5	8	2		

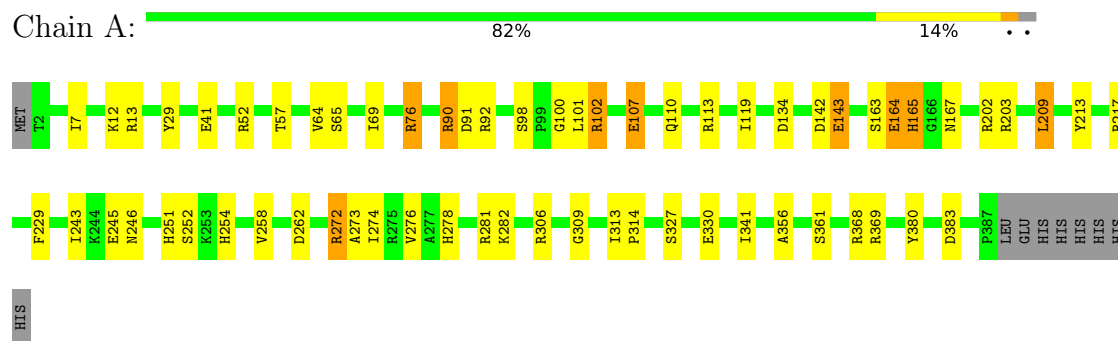
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	168	Total	O	0	0
			168	168		
7	B	159	Total	O	0	0
			159	159		
7	I	12	Total	O	0	0
			12	12		
7	F	7	Total	O	0	0
			7	7		
7	H	3	Total	O	0	0
			3	3		

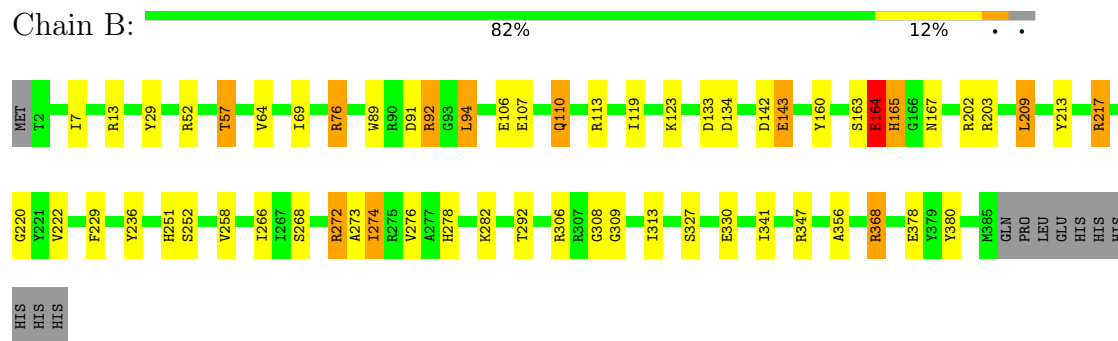
3 Residue-property plots [i](#)

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: tRNA intron endonuclease



- Molecule 1: tRNA intron endonuclease



- Molecule 2: DNA/RNA (5'-R(P*AP*CP*CP*GP*AP*CP*CP*A)-D(P*U)-R(P*AP*GP*CP*U)-3')




- Molecule 3: RNA (5'-R(P*AP*GP*CP*GP*GP*UP*CP*A)-3')





- Molecule 4: RNA (5'-R(*AP*GP*GP*U)-3')

Chain H:  25% 25% 50%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	53.51Å 65.65Å 71.80Å 114.46° 111.89° 90.01°	Depositor
Resolution (Å)	33.53 – 1.80 33.53 – 1.80	Depositor EDS
% Data completeness (in resolution range)	96.8 (33.53-1.80) 96.8 (33.53-1.80)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.83 (at 1.79Å)	Xtriage
Refinement program	REFMAC 5.8.0352	Depositor
R, R_{free}	0.185 , 0.231 0.194 , 0.239	Depositor DCC
R_{free} test set	3707 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	20.3	Xtriage
Anisotropy	0.074	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 32.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.456 for h,-k,-h-l 0.018 for -h,k,-k-l 0.019 for -h,-k,h+k+l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7209	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.26% 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: A23, ACY

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	0.62	1/3221 (0.0%)	1.01	2/4338 (0.0%)
1	B	0.63	0/3204	1.02	2/4314 (0.0%)
2	I	0.83	1/304 (0.3%)	1.19	1/470 (0.2%)
3	F	0.73	0/193	1.21	1/299 (0.3%)
4	H	0.70	0/95	1.10	0/147
All	All	0.64	2/7017 (0.0%)	1.03	6/9568 (0.1%)

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	11
1	B	0	10
All	All	0	21

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	I	13	A	O3'-P	5.49	1.61	1.56
1	A	254	HIS	CE1-NE2	5.18	1.37	1.32

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	274	ILE	N-CA-C	8.05	118.89	112.12
3	F	12	C	O3'-P-O5'	-6.84	93.74	104.00
1	B	143	GLU	CB-CG-CD	6.09	122.95	112.60
2	I	8	C	P-O3'-C3'	-5.75	111.58	120.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	383	ASP	CA-CB-CG	5.52	118.12	112.60
1	A	143	GLU	CB-CG-CD	5.43	121.84	112.60

There are no chirality outliers.

All (21) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	100	GLY	Peptide
1	A	102	ARG	Sidechain
1	A	113	ARG	Sidechain
1	A	13	ARG	Sidechain
1	A	202	ARG	Sidechain
1	A	203	ARG	Sidechain
1	A	272	ARG	Sidechain
1	A	281	ARG	Sidechain
1	A	368	ARG	Sidechain
1	A	76	ARG	Sidechain
1	A	92	ARG	Sidechain
1	B	113	ARG	Sidechain
1	B	13	ARG	Sidechain
1	B	202	ARG	Sidechain
1	B	203	ARG	Sidechain
1	B	217	ARG	Sidechain
1	B	272	ARG	Sidechain
1	B	347	ARG	Sidechain
1	B	368	ARG	Sidechain
1	B	76	ARG	Sidechain
1	B	92	ARG	Sidechain

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	3152	0	3191	42	0
1	B	3136	0	3176	47	0
2	I	273	0	142	3	0
3	F	173	0	87	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	H	85	0	45	6	0
5	A	12	0	9	3	0
5	B	4	0	3	1	0
6	F	25	0	10	6	0
7	A	168	0	0	6	0
7	B	159	0	0	3	0
7	F	7	0	0	0	0
7	H	3	0	0	0	0
7	I	12	0	0	0	0
All	All	7209	0	6663	89	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 7.

All (89) 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:B:92:ARG:HB2	1:B:94:LEU:HD11	1.28	1.06
6:F:101:A23:PC	4:H:15:A:O5'	2.22	0.97
1:B:89:TRP:O	1:B:94:LEU:HD12	1.64	0.95
1:B:306:ARG:NH1	1:B:309:GLY:H	1.65	0.94
1:B:306:ARG:HH11	1:B:309:GLY:H	1.18	0.89
1:B:92:ARG:HB2	1:B:94:LEU:CD1	2.05	0.86
1:A:306:ARG:NH1	1:A:309:GLY:H	1.74	0.86
1:A:306:ARG:NH1	1:A:309:GLY:N	2.28	0.81
1:B:89:TRP:O	1:B:94:LEU:CD1	2.29	0.79
1:A:41:GLU:HG3	7:A:665:HOH:O	1.81	0.79
1:A:91:ASP:OD1	1:B:165:HIS:HE1	1.71	0.74
1:A:369:ARG:HG3	5:A:403:ACY:H3	1.70	0.73
1:B:57:THR:HG22	7:B:610:HOH:O	1.91	0.70
6:F:101:A23:PC	4:H:15:A:HO5'	2.12	0.68
1:A:262:ASP:HB2	7:A:666:HOH:O	1.95	0.66
3:F:12:C:O2'	4:H:16:G:OP2	2.11	0.65
1:A:306:ARG:HH11	1:A:309:GLY:H	1.46	0.63
1:B:306:ARG:NH1	1:B:309:GLY:N	2.43	0.63
1:A:91:ASP:CG	1:B:165:HIS:HE1	2.08	0.62
1:A:165:HIS:HE2	1:B:91:ASP:CG	2.09	0.61
1:A:306:ARG:HH11	1:A:309:GLY:N	1.96	0.61
5:A:403:ACY:O	7:A:501:HOH:O	2.16	0.61
1:A:213:TYR:CE1	1:A:217:ARG:HG3	2.36	0.60
5:A:403:ACY:C	7:A:501:HOH:O	2.48	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:76:ARG:HH11	1:B:76:ARG:HG2	1.67	0.60
1:A:76:ARG:HG2	1:A:76:ARG:HH11	1.67	0.60
1:A:251:HIS:HA	6:F:101:A23:O1C	2.00	0.59
1:B:272:ARG:O	1:B:274:ILE:N	2.37	0.58
1:B:213:TYR:CE1	1:B:217:ARG:HG3	2.39	0.57
1:B:306:ARG:HH11	1:B:306:ARG:HG2	1.71	0.56
1:A:278:HIS:HE1	2:I:12:C:OP1	1.91	0.54
1:B:278:HIS:HE1	3:F:12:C:OP1	1.90	0.54
1:A:306:ARG:HH12	1:A:309:GLY:N	2.03	0.54
1:A:134:ASP:OD1	1:B:330:GLU:OE2	2.26	0.53
1:A:91:ASP:OD1	1:B:165:HIS:CE1	2.57	0.53
1:A:245:GLU:O	1:A:246:ASN:HB2	2.09	0.53
1:A:64:VAL:CG2	1:A:69:ILE:HD11	2.39	0.52
1:A:252:SER:O	1:A:282:LYS:NZ	2.38	0.51
1:B:229:PHE:HA	1:B:276:VAL:HG11	1.92	0.51
1:B:64:VAL:HG21	1:B:69:ILE:HD11	1.93	0.51
1:B:252:SER:O	1:B:282:LYS:NZ	2.40	0.51
1:A:330:GLU:OE2	1:B:134:ASP:OD1	2.29	0.51
1:B:209:LEU:HD22	1:B:258:VAL:HB	1.94	0.50
1:A:119:ILE:HG22	7:A:574:HOH:O	2.10	0.50
1:B:220:GLY:O	5:B:401:ACY:H2	2.11	0.50
1:A:313:ILE:HG22	1:A:314:PRO:O	2.12	0.49
1:A:229:PHE:HA	1:A:276:VAL:HG11	1.94	0.49
1:B:306:ARG:HH12	1:B:308:GLY:N	2.09	0.49
1:B:110:GLN:HA	1:B:110:GLN:OE1	2.12	0.49
1:B:163:SER:C	1:B:164:GLU:HG2	2.38	0.48
1:B:64:VAL:CG2	1:B:69:ILE:HD11	2.43	0.48
1:A:64:VAL:HG21	1:A:69:ILE:HD11	1.94	0.48
1:A:29:TYR:CE2	1:A:313:ILE:HD12	2.49	0.47
1:A:306:ARG:HH11	1:A:309:GLY:CA	2.27	0.47
1:A:107:GLU:OE2	1:A:110:GLN:HG3	2.15	0.47
1:A:361:SER:HB2	1:B:160:TYR:CD2	2.50	0.47
1:B:266:ILE:HG22	1:B:268:SER:OG	2.15	0.47
1:B:106:GLU:HG3	1:B:110:GLN:NE2	2.30	0.46
1:A:163:SER:C	1:A:164:GLU:HG2	2.41	0.46
1:B:119:ILE:HG22	7:B:587:HOH:O	2.15	0.46
1:B:213:TYR:CZ	1:B:217:ARG:HG3	2.50	0.46
1:A:213:TYR:CZ	1:A:217:ARG:HG3	2.50	0.46
1:B:94:LEU:HD12	1:B:94:LEU:H	1.80	0.45
1:B:341:ILE:HG21	1:B:380:TYR:CE1	2.52	0.45
1:B:251:HIS:HA	2:I:15:A:OP2	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:I:9:G:O2'	4:H:15:A:N1	2.49	0.45
1:A:341:ILE:HG21	1:A:380:TYR:CE1	2.52	0.44
1:A:142:ASP:OD1	1:A:167:ASN:HB2	2.17	0.44
1:A:12:LYS:HE3	7:A:549:HOH:O	2.17	0.44
1:A:91:ASP:CG	1:B:165:HIS:CE1	2.93	0.44
1:B:327:SER:HA	1:B:356:ALA:O	2.18	0.44
1:B:92:ARG:CB	1:B:94:LEU:HD11	2.21	0.43
1:B:142:ASP:CG	1:B:167:ASN:HB2	2.44	0.43
1:B:142:ASP:OD1	1:B:167:ASN:HB2	2.18	0.43
6:F:101:A23:O1C	4:H:15:A:O5'	2.37	0.43
1:A:7:ILE:HD12	1:A:69:ILE:HD12	2.00	0.43
1:B:29:TYR:CE2	1:B:313:ILE:HD12	2.54	0.42
1:A:252:SER:N	6:F:101:A23:O1C	2.49	0.42
1:A:272:ARG:O	1:A:274:ILE:N	2.53	0.41
1:A:90:ARG:HD3	1:A:90:ARG:HA	1.93	0.41
6:F:101:A23:O2C	4:H:15:A:O5'	2.38	0.41
1:A:209:LEU:HD22	1:A:258:VAL:HB	2.02	0.41
1:B:368:ARG:NH2	7:B:514:HOH:O	2.53	0.41
1:B:7:ILE:HD12	1:B:69:ILE:HD12	2.03	0.40
1:B:110:GLN:OE1	1:B:110:GLN:CA	2.69	0.40
1:A:327:SER:HA	1:A:356:ALA:O	2.21	0.40
1:A:65:SER:O	1:A:69:ILE:HG12	2.20	0.40
1:B:222:VAL:HB	1:B:236:TYR:HB2	2.03	0.40
1:B:292:THR:HG23	1:B:378:GLU:OE2	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	384/395 (97%)	376 (98%)	6 (2%)	2 (0%)	25 14

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	382/395 (97%)	374 (98%)	5 (1%)	3 (1%)	16	6
All	All	766/790 (97%)	750 (98%)	11 (1%)	5 (1%)	19	9

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	273	ALA
1	B	165	HIS
1	B	164	GLU
1	A	165	HIS
1	A	273	ALA

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	340/349 (97%)	329 (97%)	11 (3%)	34	22
1	B	338/349 (97%)	328 (97%)	10 (3%)	36	24
All	All	678/698 (97%)	657 (97%)	21 (3%)	35	22

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

Mol	Chain	Res	Type
1	A	52	ARG
1	A	57	THR
1	A	90	ARG
1	A	98	SER
1	A	101	LEU
1	A	102	ARG
1	A	107	GLU
1	A	143	GLU
1	A	164	GLU
1	A	209	LEU
1	A	243	ILE

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Mol	Chain	Res	Type
1	B	52	ARG
1	B	57	THR
1	B	94	LEU
1	B	107	GLU
1	B	110	GLN
1	B	123	LYS
1	B	133	ASP
1	B	143	GLU
1	B	164	GLU
1	B	209	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	6	ASN
1	A	108	HIS
1	A	167	ASN
1	A	251	HIS
1	A	278	HIS
1	B	60	ASN
1	B	108	HIS
1	B	165	HIS
1	B	251	HIS
1	B	278	HIS

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	I	10/13 (76%)	2 (20%)	0
3	F	7/8 (87%)	1 (14%)	0
4	H	4/4 (100%)	2 (50%)	1 (25%)
All	All	21/25 (84%)	5 (23%)	1 (4%)

All (5) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	I	13	A
2	I	18	U
3	F	13	A
4	H	16	G

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Mol	Chain	Res	Type
4	H	18	U

All (1) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
4	H	15	A

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](#)

5 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	$\# Z > 2$
5	ACY	A	402	-	3,3,3	1.04	0	3,3,3	0.66	0
5	ACY	A	401	-	3,3,3	1.03	0	3,3,3	0.90	0
5	ACY	A	403	-	3,3,3	1.02	0	3,3,3	0.81	0
6	A23	F	101	3	19,28,29	0.85	0	19,43,46	0.94	0
5	ACY	B	401	-	3,3,3	0.86	0	3,3,3	1.05	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
6	A23	F	101	3	-	0/3/35/36	0/4/4/4

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

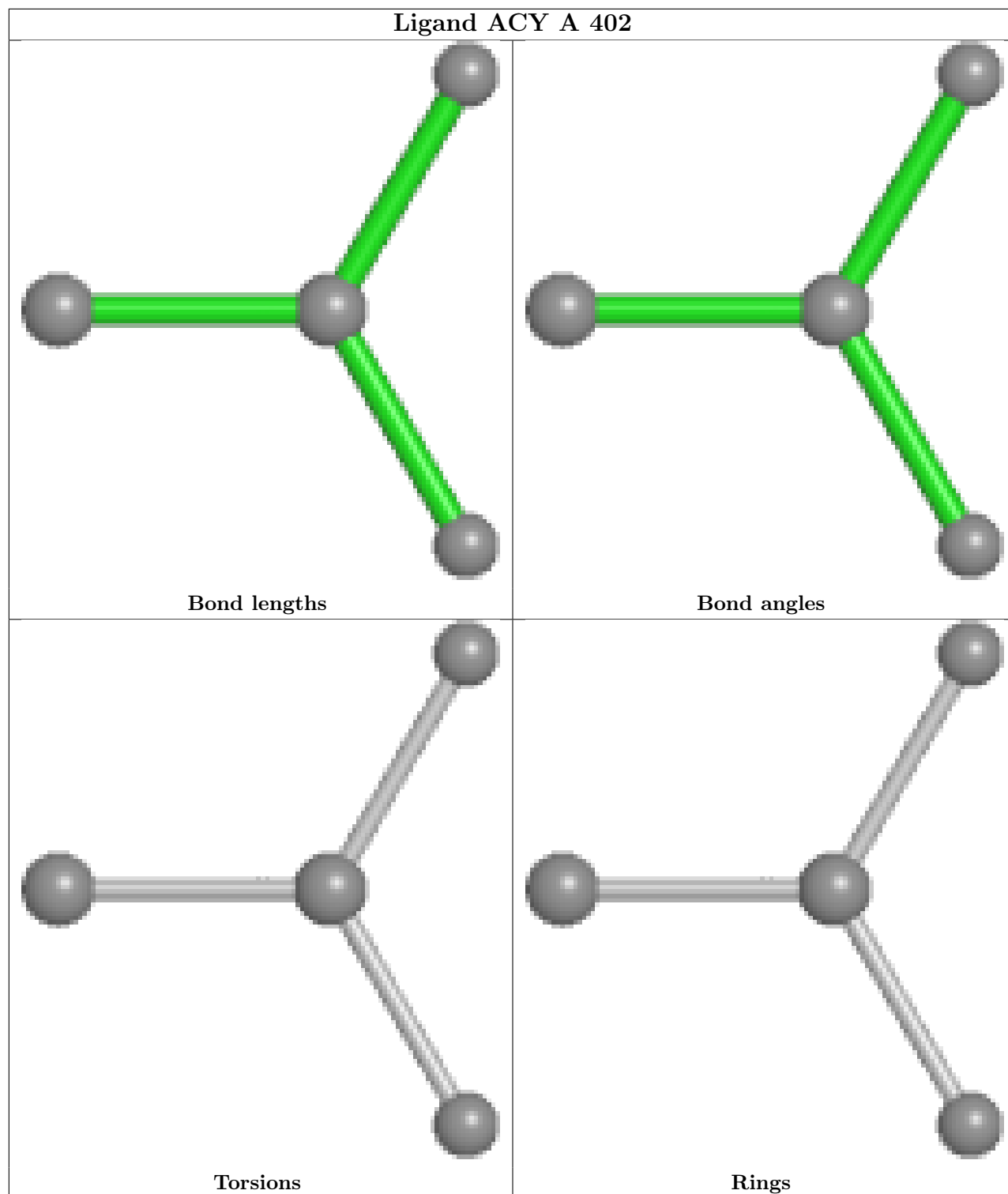
There are no torsion outliers.

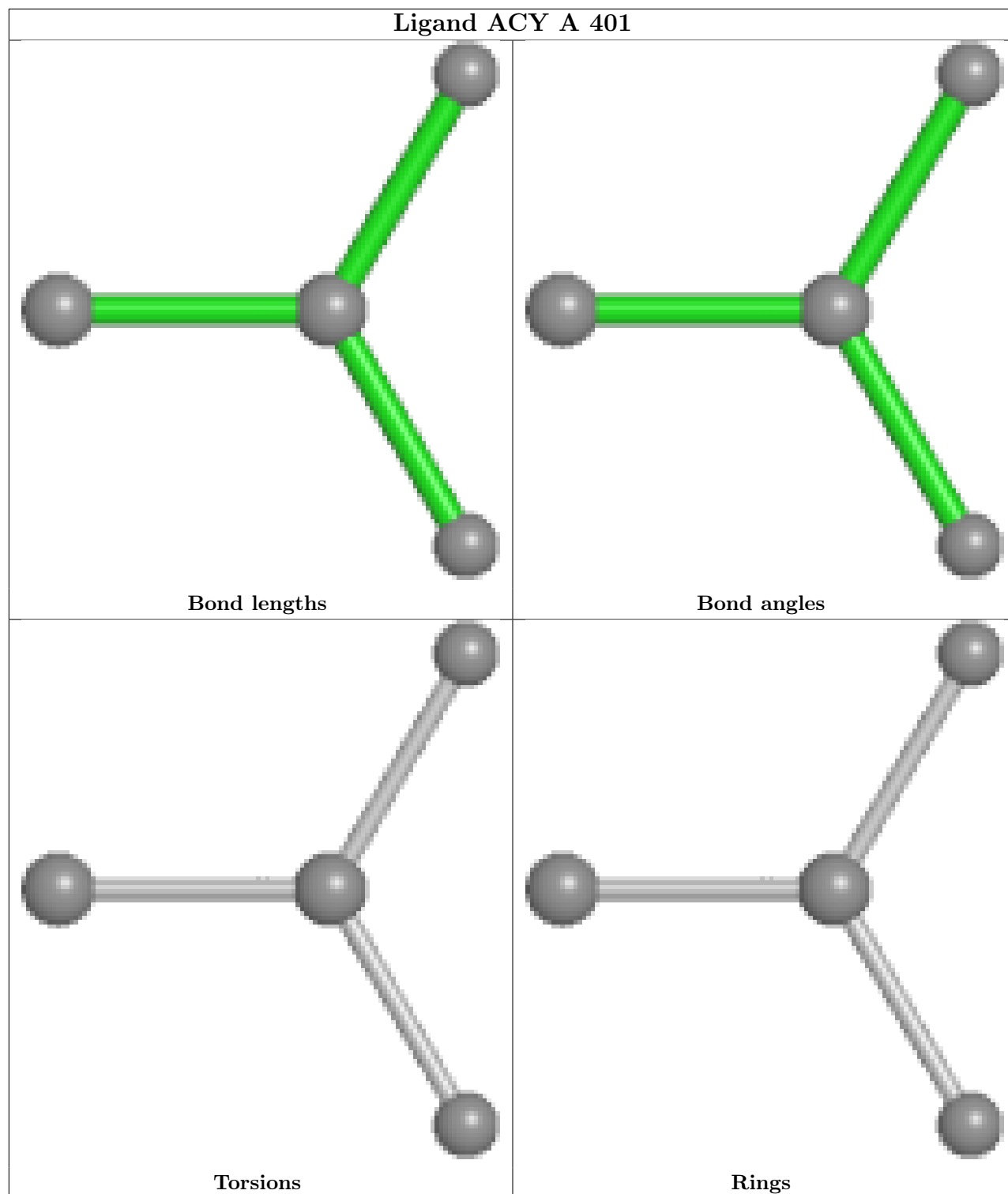
There are no ring outliers.

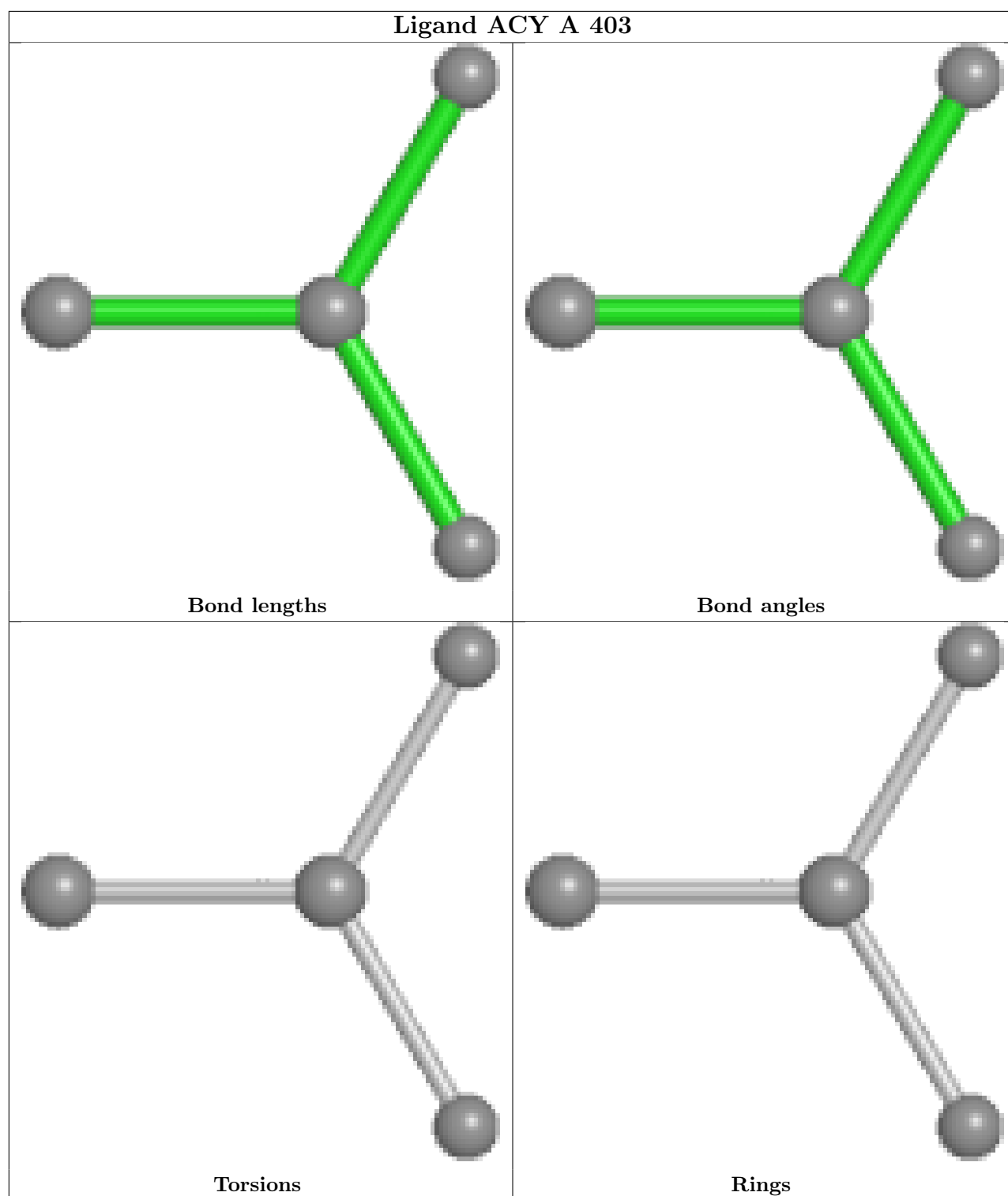
3 monomers are involved in 10 short contacts:

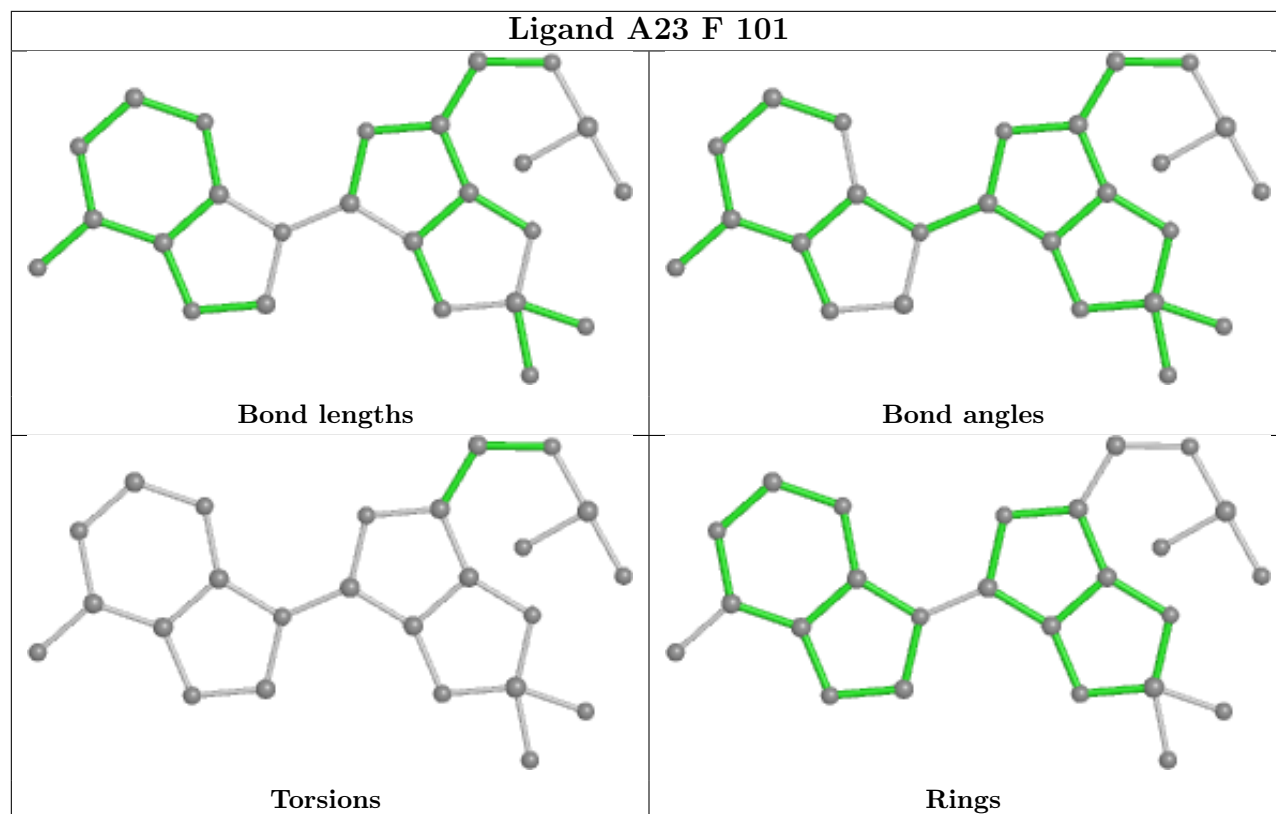
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	403	ACY	3	0
6	F	101	A23	6	0
5	B	401	ACY	1	0

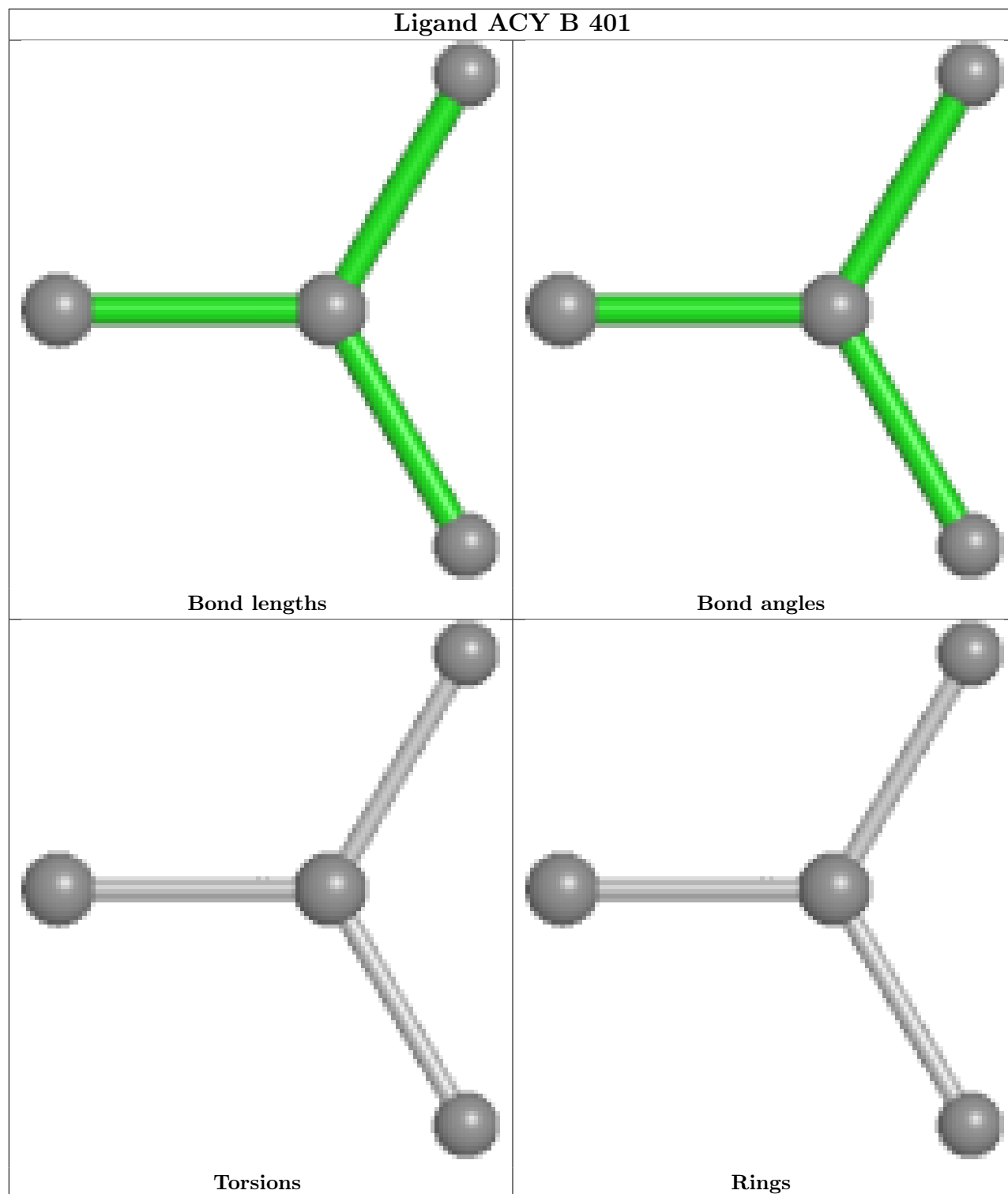
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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











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 [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	386/395 (97%)	-0.90	0 100 100	12, 22, 47, 68	0
1	B	384/395 (97%)	-0.92	0 100 100	12, 21, 45, 72	0
2	I	12/13 (92%)	-0.46	0 100 100	31, 38, 70, 90	0
3	F	8/8 (100%)	-0.63	0 100 100	30, 37, 63, 92	0
4	H	4/4 (100%)	-0.24	0 100 100	37, 38, 60, 73	0
All	All	794/815 (97%)	-0.89	0 100 100	12, 22, 50, 92	0

There are no RSRZ outliers to report.

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
5	ACY	A	402	4/4	0.96	0.05	53,59,60,66	0
5	ACY	B	401	4/4	0.97	0.07	28,35,36,40	0
5	ACY	A	403	4/4	0.98	0.06	20,30,34,38	0

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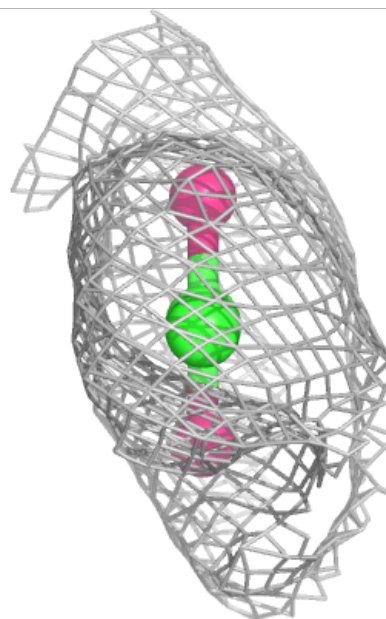
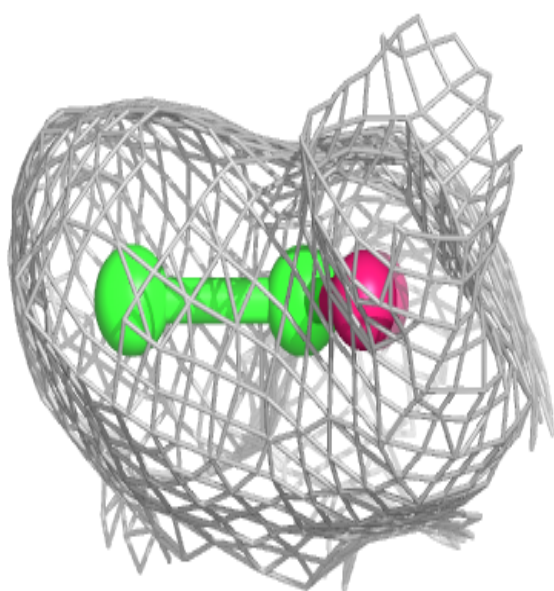
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	ACY	A	401	4/4	0.98	0.08	31,37,39,41	0
6	A23	F	101	25/26	0.99	0.06	29,47,58,69	0

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.

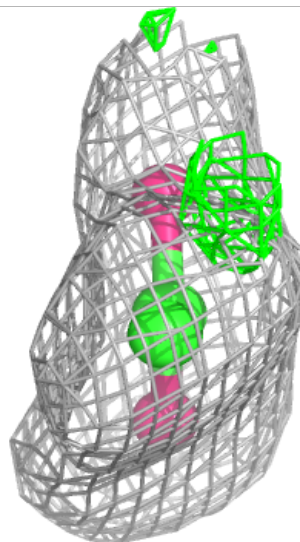
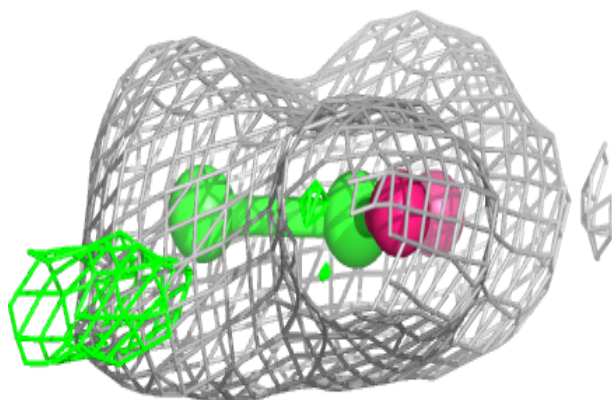
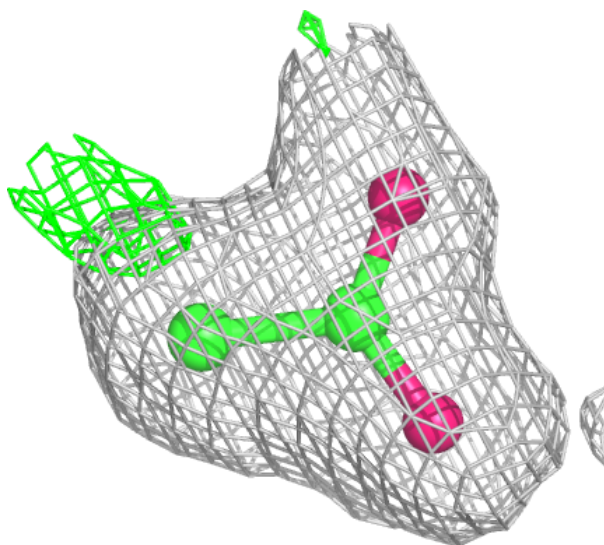
Electron density around ACY A 402:

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



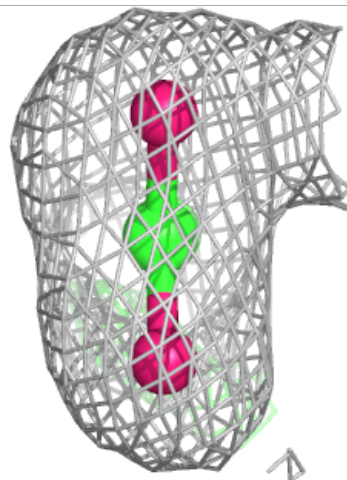
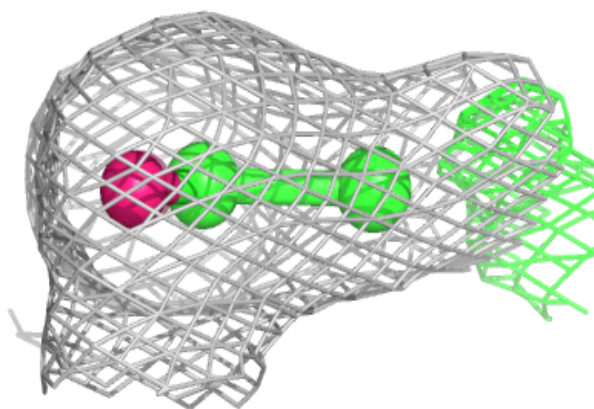
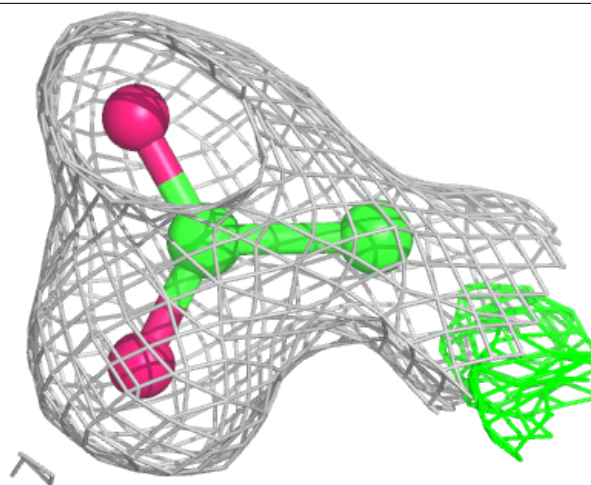
Electron density around ACY B 401:

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



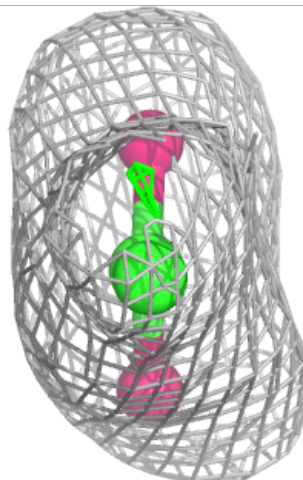
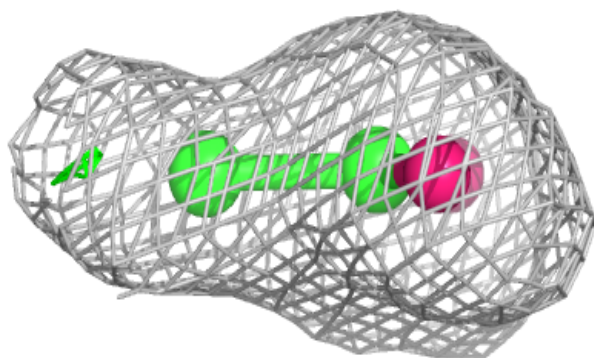
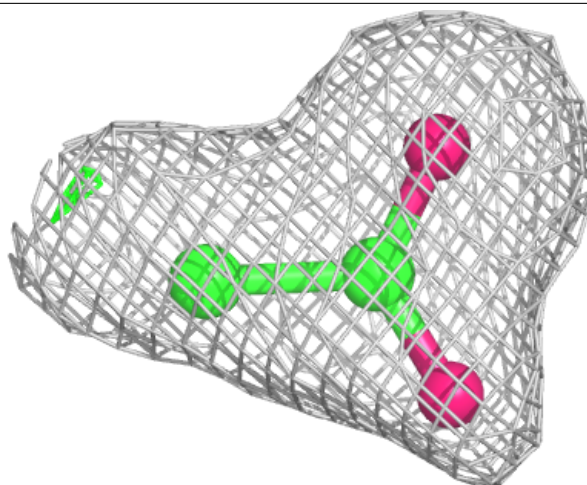
Electron density around ACY A 403:

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



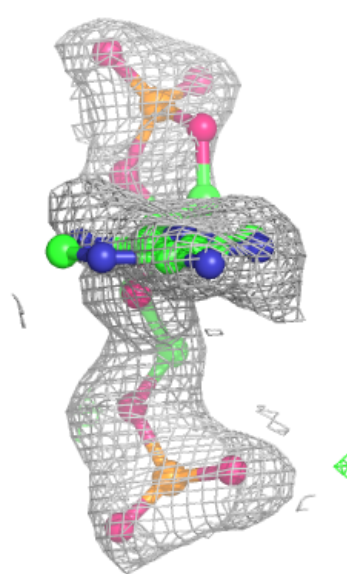
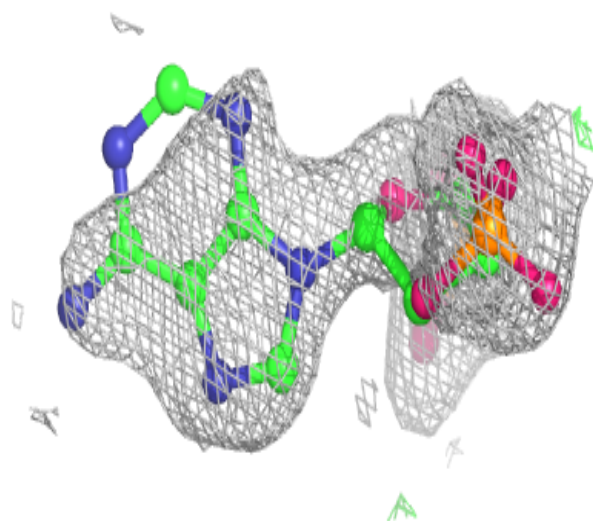
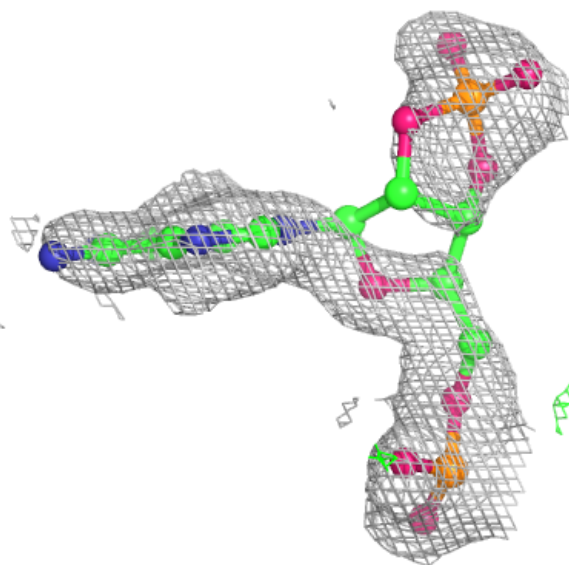
Electron density around ACY A 401:

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



Electron density around A23 F 101:

$2mF_o-DF_c$ (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.