



wwPDB X-ray Structure Validation Summary Report ⓘ

Nov 3, 2025 – 06:15 pm GMT

PDB ID : 9IF3 / pdb_00009if3
Title : Structure of YIUA from Yersinia ruckeri with Iron and DHB-L-Arg-L-Ser
Authors : Thompson, S.; Thomsen, E.; Duhme-Klair, A.; Butler, A.; Grogan, G.
Deposited on : 2025-02-17
Resolution : 2.45 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.46

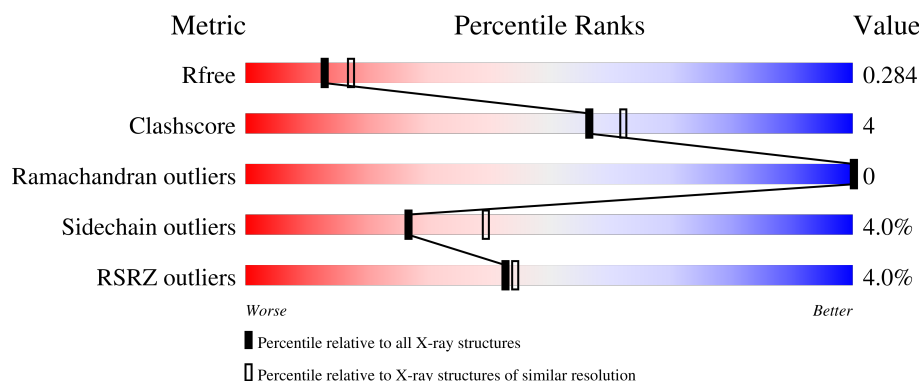
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.45 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.



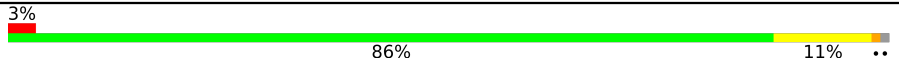

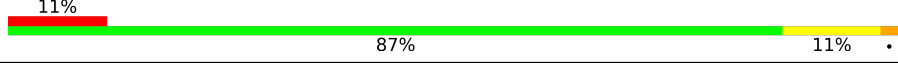
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	164625	1096 (2.46-2.46)
Clashscore	180529	1178 (2.46-2.46)
Ramachandran outliers	177936	1170 (2.46-2.46)
Sidechain outliers	177891	1170 (2.46-2.46)
RSRZ outliers	164620	1096 (2.46-2.46)

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	349	<div> <div>5%</div> <div> <div></div> <div>88%</div> <div>10%</div> <div>..</div> </div> </div>
1	B	349	<div> <div>3%</div> <div> <div></div> <div>88%</div> <div>10%</div> <div>..</div> </div> </div>
1	C	349	<div> <div>2%</div> <div> <div></div> <div>87%</div> <div>11%</div> <div>..</div> </div> </div>
1	D	349	<div> <div>3%</div> <div> <div></div> <div>86%</div> <div>12%</div> <div>..</div> </div> </div>
1	E	349	<div> <div>2%</div> <div> <div></div> <div>89%</div> <div>9%</div> <div>..</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	F	349	 3% 86% 11% ..
1	G	349	 2% 85% 13% ..
1	H	349	 11% 87% 11% ..

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
4	SO4	A	405	-	-	X	-
4	SO4	G	404	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 21896 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 Periplasmic substrate-binding transport protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	346	Total	C	N	O	S	0	0	0
			2609	1660	444	497	8			
1	B	346	Total	C	N	O	S	0	0	0
			2616	1668	441	499	8			
1	C	346	Total	C	N	O	S	0	0	0
			2641	1678	449	506	8			
1	D	347	Total	C	N	O	S	0	0	0
			2609	1661	445	495	8			
1	E	347	Total	C	N	O	S	0	1	0
			2639	1680	449	502	8			
1	F	346	Total	C	N	O	S	0	0	0
			2610	1661	446	495	8			
1	G	347	Total	C	N	O	S	0	0	0
			2611	1661	442	500	8			
1	H	346	Total	C	N	O	S	0	0	0
			2562	1633	430	491	8			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	ALA	conflict	UNP A0A085U4N5
A	12	VAL	ALA	conflict	UNP A0A085U4N5
A	233	THR	ALA	conflict	UNP A0A085U4N5
A	262	GLU	ALA	conflict	UNP A0A085U4N5
A	315	GLN	GLU	conflict	UNP A0A085U4N5
A	324	ASN	LYS	conflict	UNP A0A085U4N5
B	0	SER	ALA	conflict	UNP A0A085U4N5
B	12	VAL	ALA	conflict	UNP A0A085U4N5
B	233	THR	ALA	conflict	UNP A0A085U4N5
B	262	GLU	ALA	conflict	UNP A0A085U4N5
B	315	GLN	GLU	conflict	UNP A0A085U4N5
B	324	ASN	LYS	conflict	UNP A0A085U4N5
C	0	SER	ALA	conflict	UNP A0A085U4N5

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Chain	Residue	Modelled	Actual	Comment	Reference
C	12	VAL	ALA	conflict	UNP A0A085U4N5
C	233	THR	ALA	conflict	UNP A0A085U4N5
C	262	GLU	ALA	conflict	UNP A0A085U4N5
C	315	GLN	GLU	conflict	UNP A0A085U4N5
C	324	ASN	LYS	conflict	UNP A0A085U4N5
D	0	SER	ALA	conflict	UNP A0A085U4N5
D	12	VAL	ALA	conflict	UNP A0A085U4N5
D	233	THR	ALA	conflict	UNP A0A085U4N5
D	262	GLU	ALA	conflict	UNP A0A085U4N5
D	315	GLN	GLU	conflict	UNP A0A085U4N5
D	324	ASN	LYS	conflict	UNP A0A085U4N5
E	0	SER	ALA	conflict	UNP A0A085U4N5
E	12	VAL	ALA	conflict	UNP A0A085U4N5
E	233	THR	ALA	conflict	UNP A0A085U4N5
E	262	GLU	ALA	conflict	UNP A0A085U4N5
E	315	GLN	GLU	conflict	UNP A0A085U4N5
E	324	ASN	LYS	conflict	UNP A0A085U4N5
F	0	SER	ALA	conflict	UNP A0A085U4N5
F	12	VAL	ALA	conflict	UNP A0A085U4N5
F	233	THR	ALA	conflict	UNP A0A085U4N5
F	262	GLU	ALA	conflict	UNP A0A085U4N5
F	315	GLN	GLU	conflict	UNP A0A085U4N5
F	324	ASN	LYS	conflict	UNP A0A085U4N5
G	0	SER	ALA	conflict	UNP A0A085U4N5
G	12	VAL	ALA	conflict	UNP A0A085U4N5
G	233	THR	ALA	conflict	UNP A0A085U4N5
G	262	GLU	ALA	conflict	UNP A0A085U4N5
G	315	GLN	GLU	conflict	UNP A0A085U4N5
G	324	ASN	LYS	conflict	UNP A0A085U4N5
H	0	SER	ALA	conflict	UNP A0A085U4N5
H	12	VAL	ALA	conflict	UNP A0A085U4N5
H	233	THR	ALA	conflict	UNP A0A085U4N5
H	262	GLU	ALA	conflict	UNP A0A085U4N5
H	315	GLN	GLU	conflict	UNP A0A085U4N5
H	324	ASN	LYS	conflict	UNP A0A085U4N5

- Molecule 2 is FE (III) ION (CCD ID: FE) (formula: Fe) (labeled as "Ligand of Interest" by depositor).

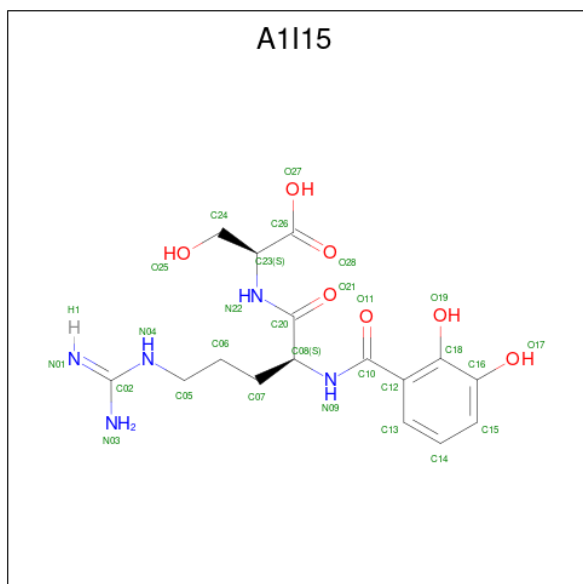
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Fe 1 1	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Fe	0	0
			1	1		
2	C	1	Total	Fe	0	0
			1	1		
2	D	1	Total	Fe	0	0
			1	1		
2	E	1	Total	Fe	0	0
			1	1		
2	F	1	Total	Fe	0	0
			1	1		
2	G	1	Total	Fe	0	0
			1	1		
2	H	1	Total	Fe	0	0
			1	1		

- Molecule 3 is (2 {S})-2-[[[(2 {S})-2-[[2,3-bis(oxidanyl)phenyl]carbonylamino]-5-carbamimidamido-pentanoyl]amino]-3-oxidanyl-propanoic acid (CCD ID: A1I15) (formula: C₁₆H₂₃N₅O₇) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			28	16	5	7		
3	A	1	Total	C	N	O	0	0
			28	16	5	7		
3	B	1	Total	C	N	O	0	0
			28	16	5	7		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	N	O	0	0
			28	16	5	7		
3	C	1	Total	C	N	O	0	0
			28	16	5	7		
3	C	1	Total	C	N	O	0	0
			28	16	5	7		
3	D	1	Total	C	N	O	0	0
			28	16	5	7		
3	D	1	Total	C	N	O	0	0
			28	16	5	7		
3	E	1	Total	C	N	O	0	0
			28	16	5	7		
3	E	1	Total	C	N	O	0	0
			28	16	5	7		
3	F	1	Total	C	N	O	0	0
			28	16	5	7		
3	F	1	Total	C	N	O	0	0
			28	16	5	7		
3	G	1	Total	C	N	O	0	0
			28	16	5	7		
3	G	1	Total	C	N	O	0	0
			28	16	5	7		
3	H	1	Total	C	N	O	0	0
			28	16	5	7		
3	H	1	Total	C	N	O	0	0
			28	16	5	7		

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	E	1	Total	O	S	0	0
			5	4	1		
4	E	1	Total	O	S	0	0
			5	4	1		
4	F	1	Total	O	S	0	0
			5	4	1		
4	F	1	Total	O	S	0	0
			5	4	1		
4	G	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	G	1	Total	O	S	0	0
			5	4	1		
4	G	1	Total	O	S	0	0
			5	4	1		
4	G	1	Total	O	S	0	0
			5	4	1		
4	H	1	Total	O	S	0	0
			5	4	1		

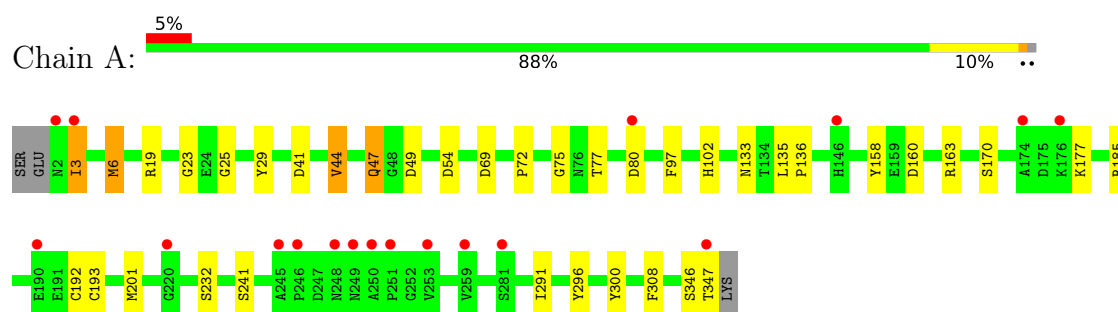
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	60	Total	O	0	0
			60	60		
5	B	50	Total	O	0	0
			50	50		
5	C	60	Total	O	0	0
			60	60		
5	D	61	Total	O	0	0
			61	61		
5	E	82	Total	O	0	0
			82	82		
5	F	45	Total	O	0	0
			45	45		
5	G	65	Total	O	0	0
			65	65		
5	H	30	Total	O	0	0
			30	30		

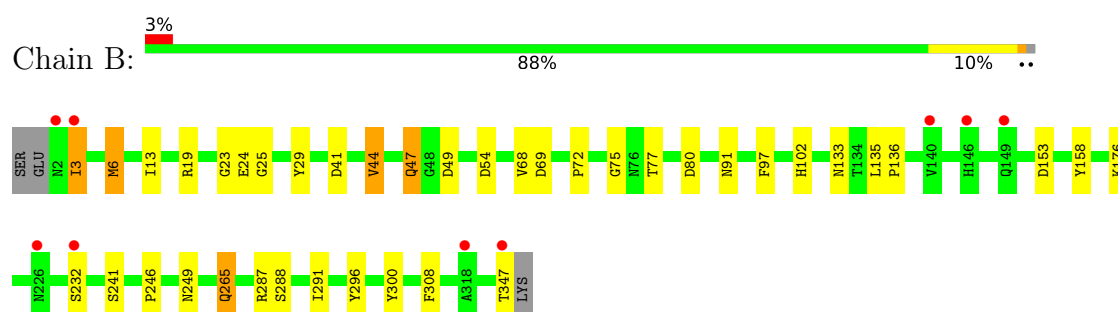
3 Residue-property plots [i](#)

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.

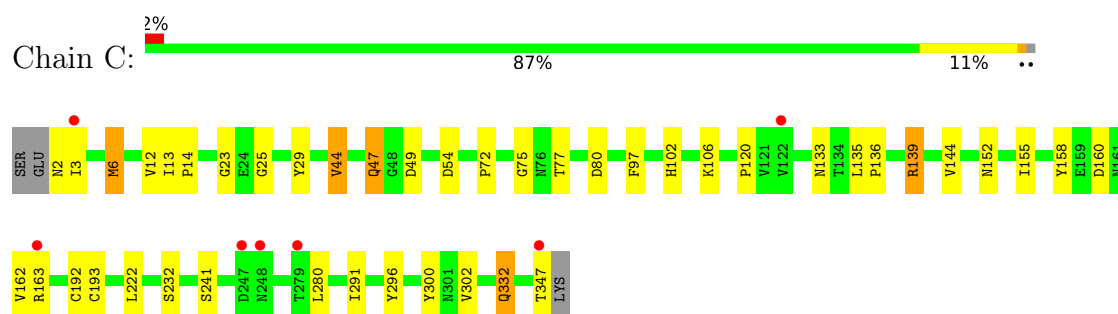
- Molecule 1: Periplasmic substrate-binding transport protein



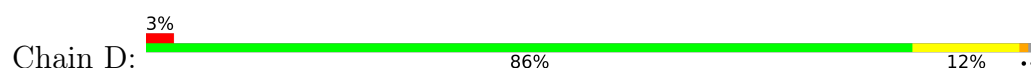
- Molecule 1: Periplasmic substrate-binding transport protein

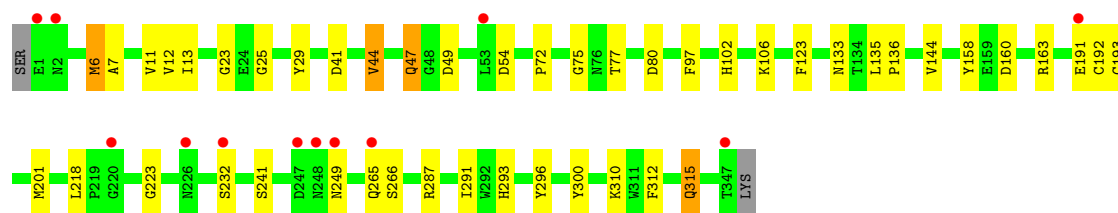


- Molecule 1: Periplasmic substrate-binding transport protein

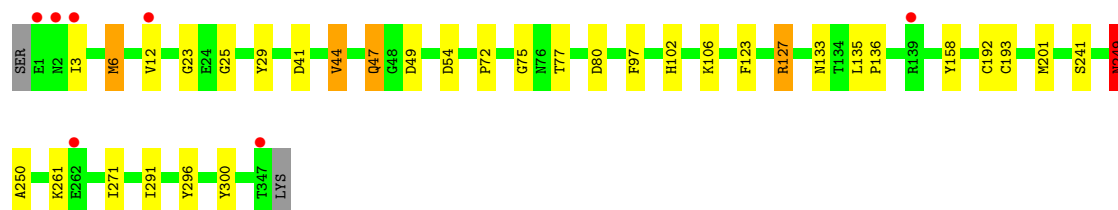
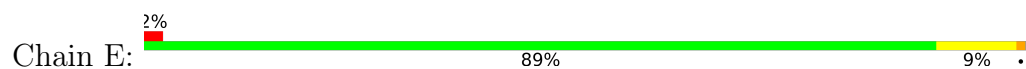


- Molecule 1: Periplasmic substrate-binding transport protein

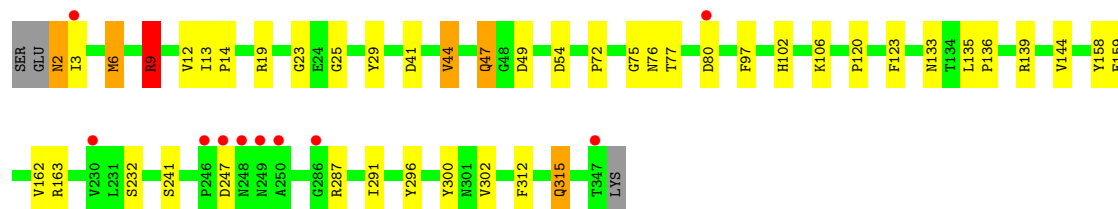
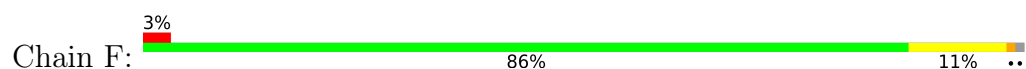




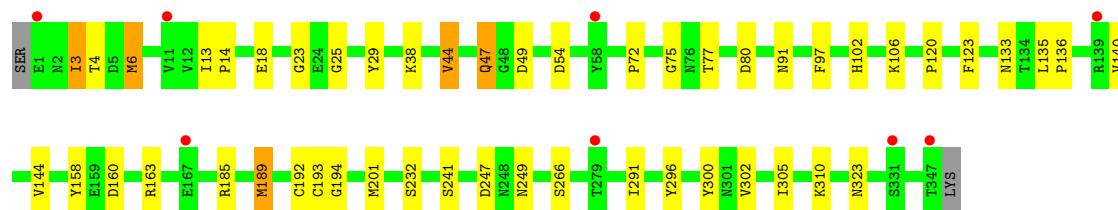
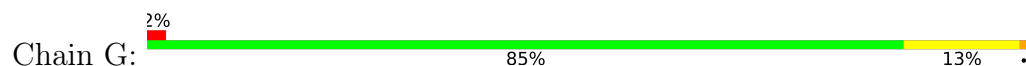
- Molecule 1: Periplasmic substrate-binding transport protein



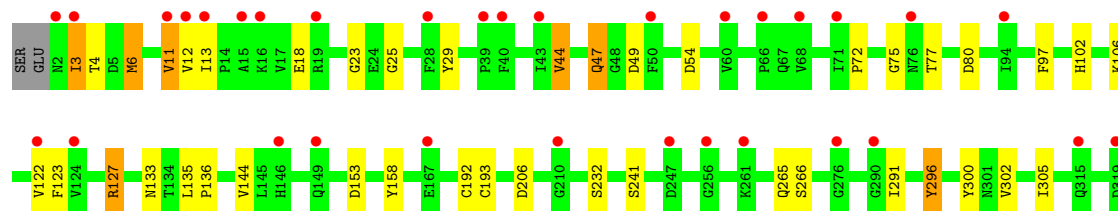
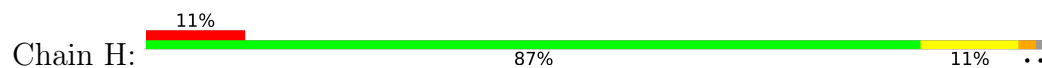
- Molecule 1: Periplasmic substrate-binding transport protein

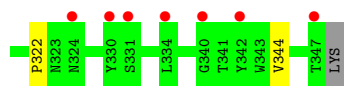


- Molecule 1: Periplasmic substrate-binding transport protein



- Molecule 1: Periplasmic substrate-binding transport protein





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	80.07Å 80.02Å 223.70Å 90.00° 89.86° 90.00°	Depositor
Resolution (Å)	75.56 – 2.45 75.56 – 2.45	Depositor EDS
% Data completeness (in resolution range)	99.4 (75.56-2.45) 99.4 (75.56-2.45)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.19 (at 2.45Å)	Xtriage
Refinement program	REFMAC 5.8.0430	Depositor
R, R_{free}	0.247 , 0.277 0.252 , 0.284	Depositor DCC
R_{free} test set	5183 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	23.9	Xtriage
Anisotropy	1.982	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 34.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.000 for -k,-h,-l 0.000 for k,h,-l 0.078 for h,-k,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	21896	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 34.72 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 6.5282e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, FE, A1I15

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	0.63	0/2666	1.06	3/3634 (0.1%)
1	B	0.63	0/2673	1.08	6/3641 (0.2%)
1	C	0.61	0/2698	1.05	1/3671 (0.0%)
1	D	0.62	0/2666	1.05	2/3634 (0.1%)
1	E	0.64	1/2699 (0.0%)	1.12	6/3673 (0.2%)
1	F	0.62	0/2667	1.07	4/3633 (0.1%)
1	G	0.64	1/2668 (0.0%)	1.06	4/3637 (0.1%)
1	H	0.60	0/2618	1.05	4/3573 (0.1%)
All	All	0.62	2/21355 (0.0%)	1.07	30/29096 (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	1
1	C	0	1
1	E	0	1
1	F	0	2
1	H	0	1
All	All	0	6

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	127	ARG	NE-CZ	-5.88	1.26	1.33
1	G	91	ASN	C-O	-5.18	1.21	1.23

The worst 5 of 30 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	127	ARG	NE-CZ-NH2	-13.29	107.24	119.20
1	E	127	ARG	CD-NE-CZ	9.66	137.93	124.40
1	E	127	ARG	NE-CZ-NH1	9.08	130.58	121.50
1	F	9	ARG	NE-CZ-NH1	6.71	128.21	121.50
1	E	127	ARG	CG-CD-NE	-6.53	97.63	112.00

There are no chirality outliers.

5 of 6 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	19	ARG	Sidechain
1	C	139	ARG	Sidechain
1	E	127	ARG	Sidechain
1	F	19	ARG	Sidechain
1	F	9	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	2609	0	2531	20	0
1	B	2616	0	2546	21	0
1	C	2641	0	2589	28	0
1	D	2609	0	2526	24	0
1	E	2639	0	2588	17	0
1	F	2610	0	2540	26	0
1	G	2611	0	2524	28	0
1	H	2562	0	2433	24	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
2	G	1	0	0	0	0
2	H	1	0	0	0	0
3	A	56	0	0	2	0
3	B	56	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	56	0	0	1	0
3	D	56	0	0	4	0
3	E	56	0	0	1	0
3	F	56	0	0	0	0
3	G	56	0	0	1	0
3	H	56	0	0	0	0
4	A	10	0	0	4	0
4	B	5	0	0	0	0
4	C	20	0	0	0	0
4	D	10	0	0	0	0
4	E	10	0	0	0	0
4	F	10	0	0	0	0
4	G	20	0	0	2	0
4	H	5	0	0	0	0
5	A	60	0	0	1	0
5	B	50	0	0	2	0
5	C	60	0	0	3	0
5	D	61	0	0	1	0
5	E	82	0	0	0	0
5	F	45	0	0	1	0
5	G	65	0	0	3	0
5	H	30	0	0	2	0
All	All	21896	0	20277	182	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 4.

The worst 5 of 182 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:A:170:SER:HB2	1:B:246:PRO:O	1.59	1.00
1:C:139:ARG:HD3	1:G:247:ASP:O	1.68	0.92
1:B:29:TYR:OH	1:B:54:ASP:OD2	2.00	0.80
1:D:29:TYR:OH	1:D:54:ASP:OD2	2.00	0.79
1:A:29:TYR:OH	1:A:54:ASP:OD2	2.01	0.79

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	344/349 (99%)	334 (97%)	10 (3%)	0	100	100
1	B	344/349 (99%)	334 (97%)	10 (3%)	0	100	100
1	C	344/349 (99%)	334 (97%)	10 (3%)	0	100	100
1	D	345/349 (99%)	335 (97%)	10 (3%)	0	100	100
1	E	346/349 (99%)	334 (96%)	12 (4%)	0	100	100
1	F	344/349 (99%)	334 (97%)	10 (3%)	0	100	100
1	G	345/349 (99%)	335 (97%)	10 (3%)	0	100	100
1	H	344/349 (99%)	333 (97%)	11 (3%)	0	100	100
All	All	2756/2792 (99%)	2673 (97%)	83 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	273/293 (93%)	263 (96%)	10 (4%)	29	43
1	B	274/293 (94%)	264 (96%)	10 (4%)	30	44
1	C	282/293 (96%)	272 (96%)	10 (4%)	31	45
1	D	270/293 (92%)	256 (95%)	14 (5%)	19	27
1	E	279/293 (95%)	268 (96%)	11 (4%)	27	41
1	F	273/293 (93%)	261 (96%)	12 (4%)	24	36

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	G	272/293 (93%)	261 (96%)	11 (4%)	27	39
1	H	258/293 (88%)	248 (96%)	10 (4%)	27	41
All	All	2181/2344 (93%)	2093 (96%)	88 (4%)	27	39

5 of 88 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	F	6	MET
1	G	47	GLN
1	F	41	ASP
1	F	300	TYR
1	G	296	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 48 such sidechains are listed below:

Mol	Chain	Res	Type
1	E	76	ASN
1	F	2	ASN
1	E	113	GLN
1	E	226	ASN
1	F	76	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 42 ligands modelled in this entry, 8 are monoatomic - leaving 34 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
4	SO4	G	407	-	4,4,4	0.28	0	6,6,6	0.21	0
4	SO4	C	404	-	4,4,4	0.25	0	6,6,6	0.21	0
4	SO4	D	405	-	4,4,4	0.43	0	6,6,6	0.05	0
3	A1I15	C	402	2	28,28,28	2.26	7 (25%)	35,37,37	1.80	10 (28%)
3	A1I15	C	403	2	28,28,28	2.26	10 (35%)	35,37,37	1.63	7 (20%)
4	SO4	A	405	-	4,4,4	0.31	0	6,6,6	0.11	0
4	SO4	C	406	-	4,4,4	0.28	0	6,6,6	0.30	0
4	SO4	D	404	-	4,4,4	0.30	0	6,6,6	0.15	0
4	SO4	H	404	-	4,4,4	0.30	0	6,6,6	0.16	0
3	A1I15	F	403	2	28,28,28	2.12	8 (28%)	35,37,37	1.42	7 (20%)
4	SO4	B	404	-	4,4,4	0.29	0	6,6,6	0.22	0
3	A1I15	F	402	2	28,28,28	2.54	7 (25%)	35,37,37	1.82	6 (17%)
4	SO4	G	406	-	4,4,4	0.30	0	6,6,6	0.11	0
3	A1I15	B	402	2	28,28,28	2.37	8 (28%)	35,37,37	1.18	4 (11%)
4	SO4	F	405	-	4,4,4	0.32	0	6,6,6	0.15	0
4	SO4	G	405	-	4,4,4	0.39	0	6,6,6	0.31	0
3	A1I15	D	402	2	28,28,28	2.31	8 (28%)	35,37,37	1.26	4 (11%)
3	A1I15	A	403	2	28,28,28	2.24	10 (35%)	35,37,37	1.39	7 (20%)
4	SO4	E	404	-	4,4,4	0.41	0	6,6,6	0.18	0
3	A1I15	E	403	2	28,28,28	2.29	8 (28%)	35,37,37	1.94	12 (34%)
3	A1I15	H	402	2	28,28,28	2.40	7 (25%)	35,37,37	1.53	7 (20%)
3	A1I15	A	402	2	28,28,28	2.32	6 (21%)	35,37,37	1.30	5 (14%)
4	SO4	F	404	-	4,4,4	0.33	0	6,6,6	0.10	0
3	A1I15	H	403	2	28,28,28	2.43	7 (25%)	35,37,37	1.70	7 (20%)
3	A1I15	E	402	2	28,28,28	2.25	8 (28%)	35,37,37	1.35	4 (11%)
4	SO4	E	405	-	4,4,4	0.29	0	6,6,6	0.16	0
3	A1I15	D	403	2	28,28,28	2.32	7 (25%)	35,37,37	1.59	9 (25%)
4	SO4	C	405	-	4,4,4	0.31	0	6,6,6	0.25	0
4	SO4	G	404	-	4,4,4	0.31	0	6,6,6	0.15	0
4	SO4	C	407	-	4,4,4	0.30	0	6,6,6	0.08	0
3	A1I15	G	403	2	28,28,28	2.13	7 (25%)	35,37,37	1.56	5 (14%)
3	A1I15	G	402	2	28,28,28	2.58	6 (21%)	35,37,37	1.23	4 (11%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	A1I15	B	403	2	28,28,28	2.26	8 (28%)	35,37,37	1.66	9 (25%)
4	SO4	A	404	-	4,4,4	0.32	0	6,6,6	0.11	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	A1I15	F	403	2	-	5/29/29/29	0/1/1/1
3	A1I15	F	402	2	-	2/29/29/29	0/1/1/1
3	A1I15	G	403	2	-	7/29/29/29	0/1/1/1
3	A1I15	E	403	2	-	7/29/29/29	0/1/1/1
3	A1I15	C	402	2	-	2/29/29/29	0/1/1/1
3	A1I15	C	403	2	-	2/29/29/29	0/1/1/1
3	A1I15	B	403	2	-	5/29/29/29	0/1/1/1
3	A1I15	G	402	2	-	4/29/29/29	0/1/1/1
3	A1I15	H	402	2	-	2/29/29/29	0/1/1/1
3	A1I15	B	402	2	-	4/29/29/29	0/1/1/1
3	A1I15	A	402	2	-	4/29/29/29	0/1/1/1
3	A1I15	H	403	2	-	2/29/29/29	0/1/1/1
3	A1I15	D	402	2	-	4/29/29/29	0/1/1/1
3	A1I15	E	402	2	-	4/29/29/29	0/1/1/1
3	A1I15	D	403	2	-	3/29/29/29	0/1/1/1
3	A1I15	A	403	2	-	3/29/29/29	0/1/1/1

The worst 5 of 122 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	403	A1I15	C02-N04	7.73	1.48	1.33
3	D	403	A1I15	C02-N04	7.62	1.48	1.33
3	G	402	A1I15	C02-N04	7.50	1.48	1.33
3	A	402	A1I15	C02-N04	7.28	1.47	1.33
3	H	403	A1I15	C02-N04	7.19	1.47	1.33

The worst 5 of 107 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	402	A1I15	C12-C18-C16	6.94	124.39	119.99
3	C	403	A1I15	O21-C20-N22	-4.46	114.68	122.93
3	C	402	A1I15	C12-C18-C16	4.38	122.77	119.99
3	E	403	A1I15	C24-C23-N22	-4.27	100.93	110.56
3	E	403	A1I15	O21-C20-N22	-4.25	115.05	122.93

There are no chirality outliers.

5 of 60 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	402	A1I15	C26-C23-C24-O25
3	B	402	A1I15	C26-C23-C24-O25
3	B	403	A1I15	N22-C23-C24-O25
3	C	403	A1I15	C05-C06-C07-C08
3	E	402	A1I15	C26-C23-C24-O25

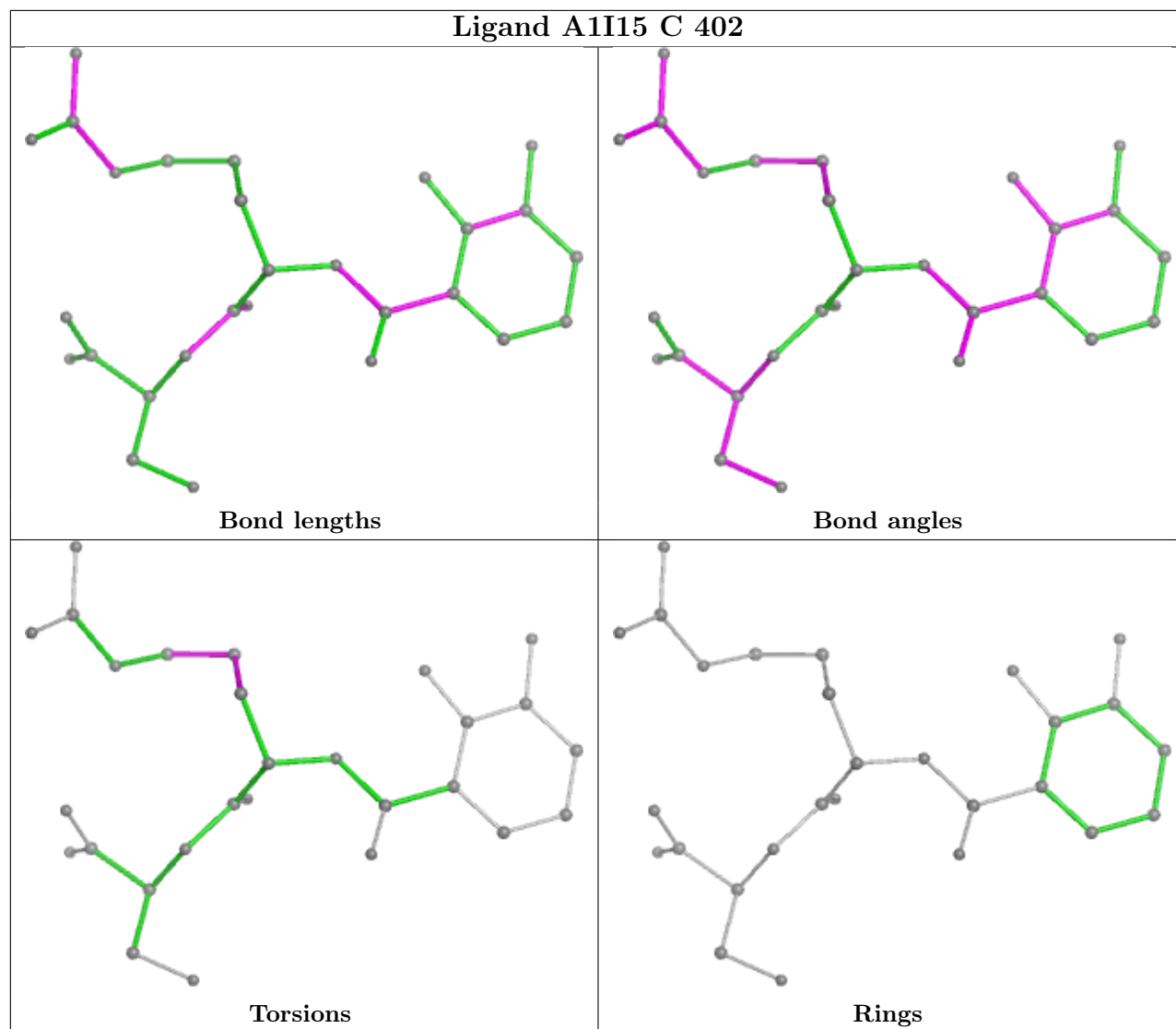
There are no ring outliers.

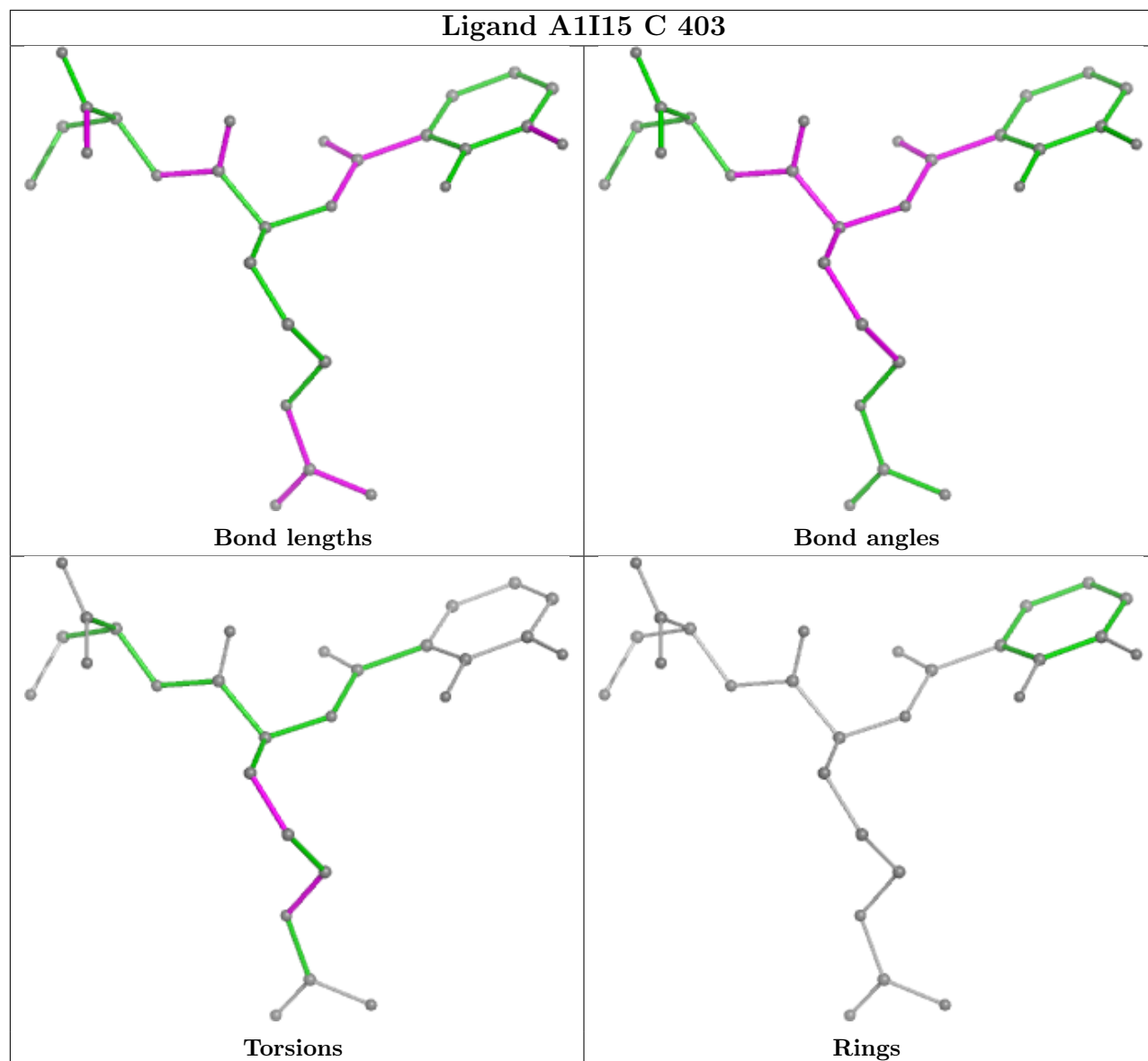
10 monomers are involved in 13 short contacts:

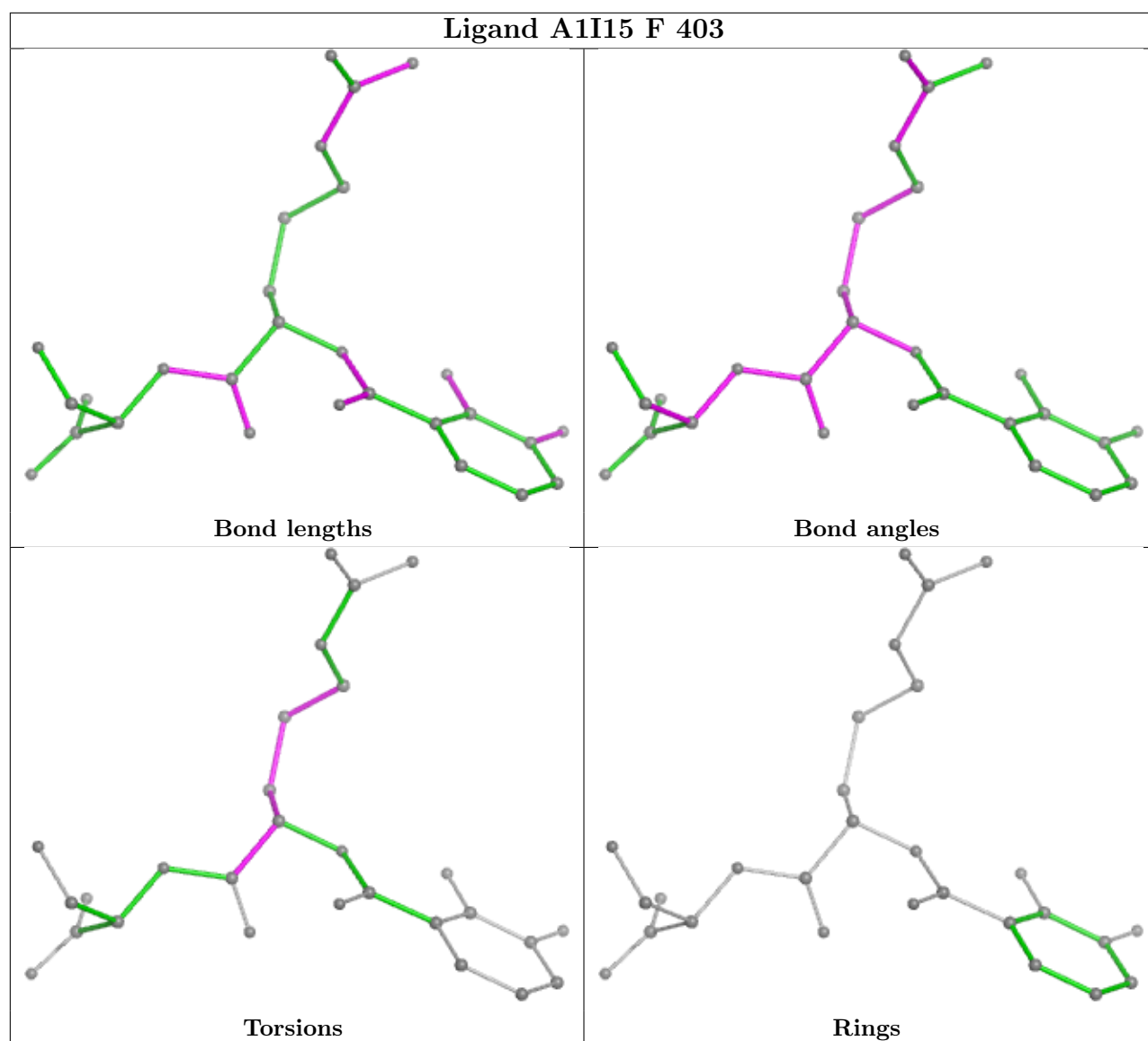
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	403	A1I15	1	0
4	A	405	SO4	3	0
3	B	402	A1I15	1	0
3	D	402	A1I15	2	0
3	A	402	A1I15	2	0
3	E	402	A1I15	1	0
3	D	403	A1I15	2	0
4	G	404	SO4	2	0
3	G	402	A1I15	1	0
4	A	404	SO4	1	0

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

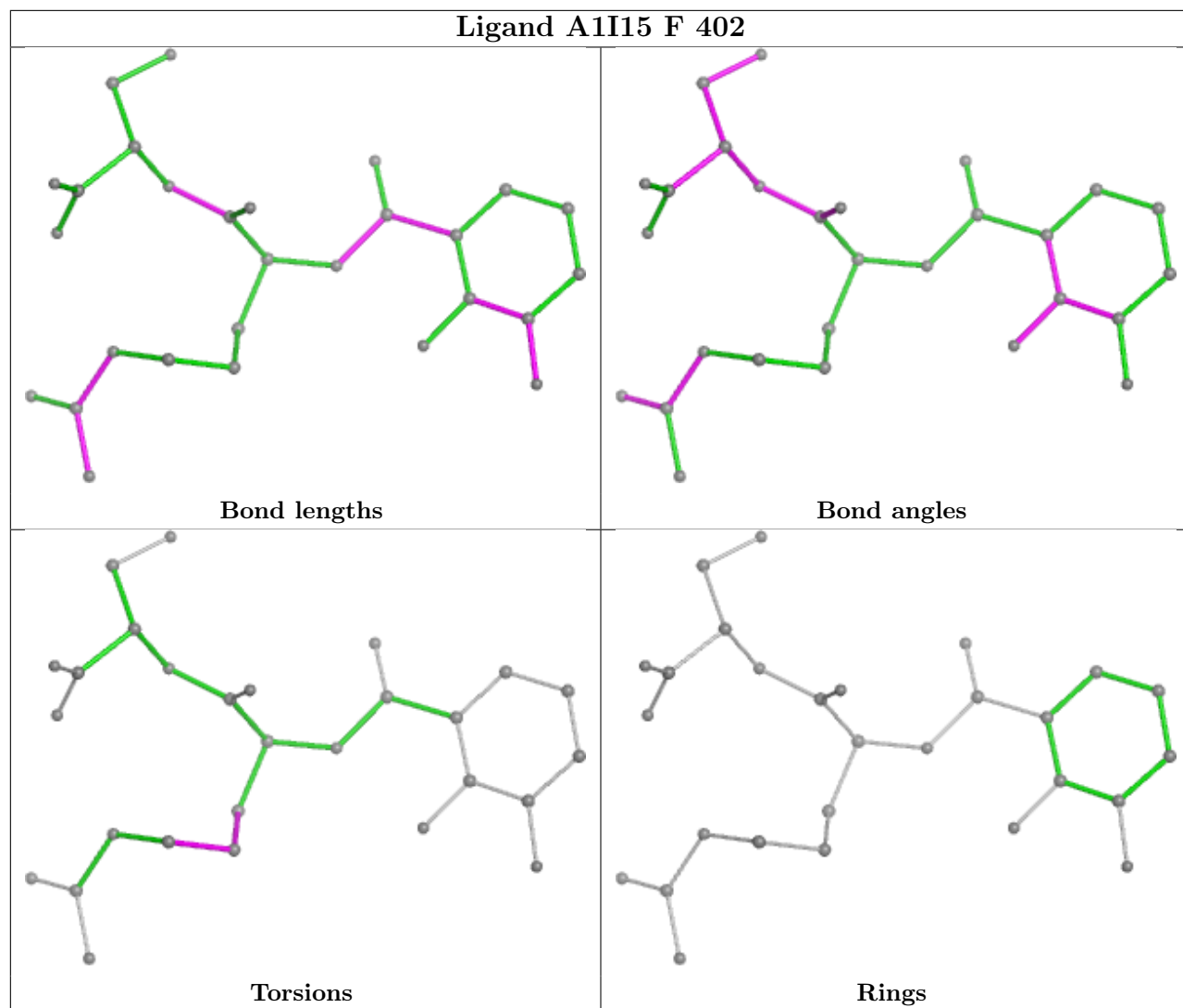
Ligand A1I15 C 402



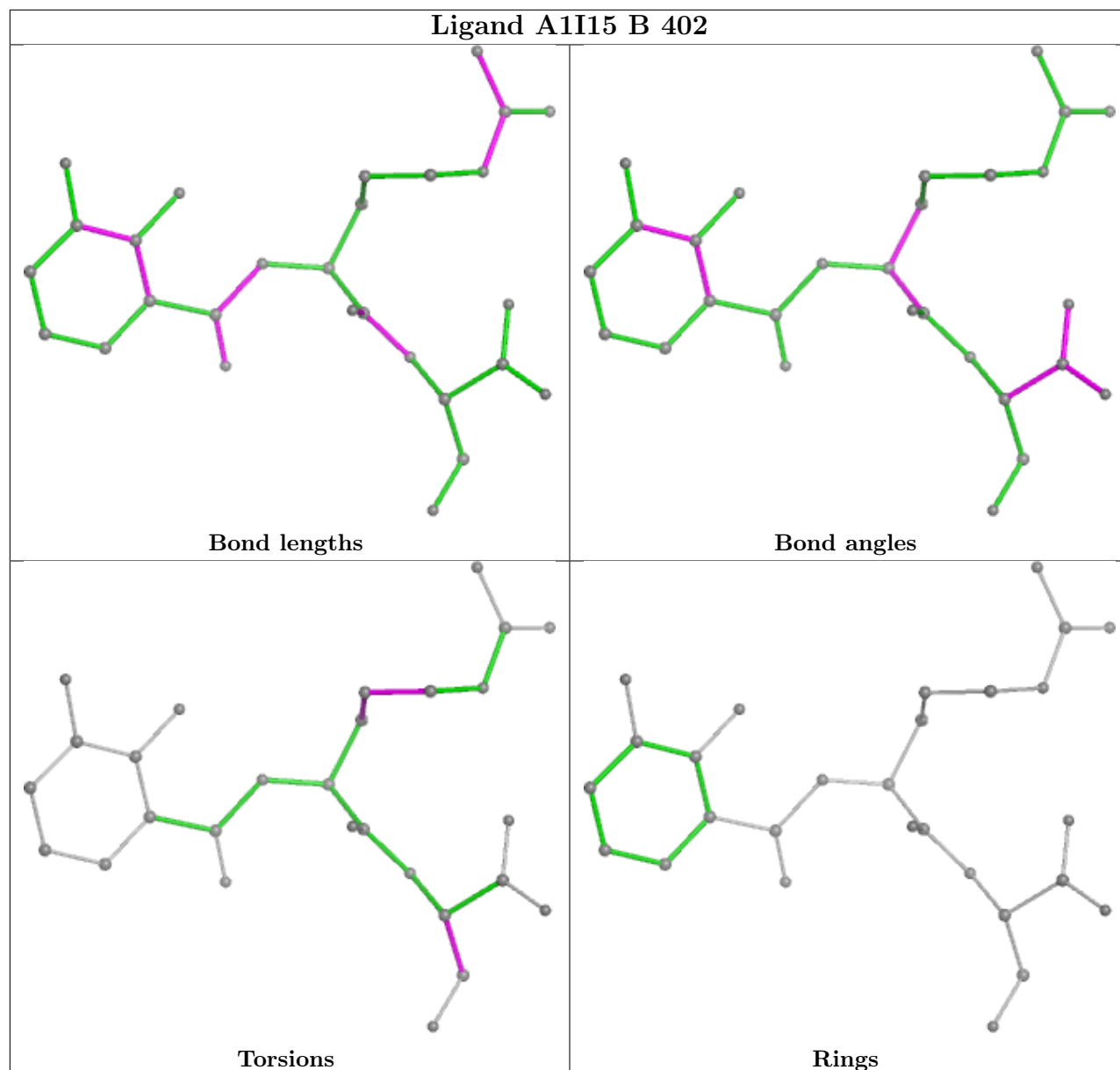




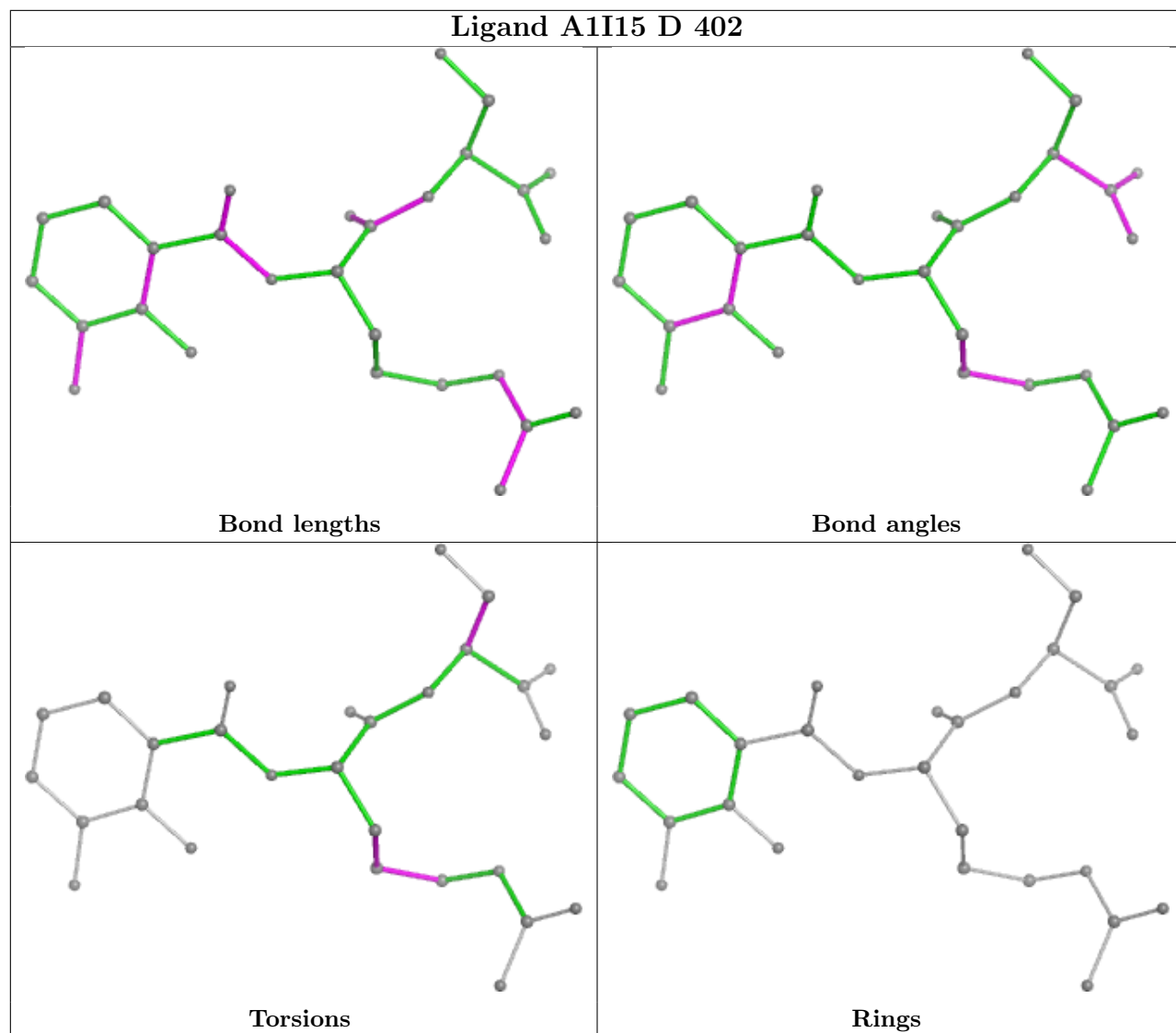
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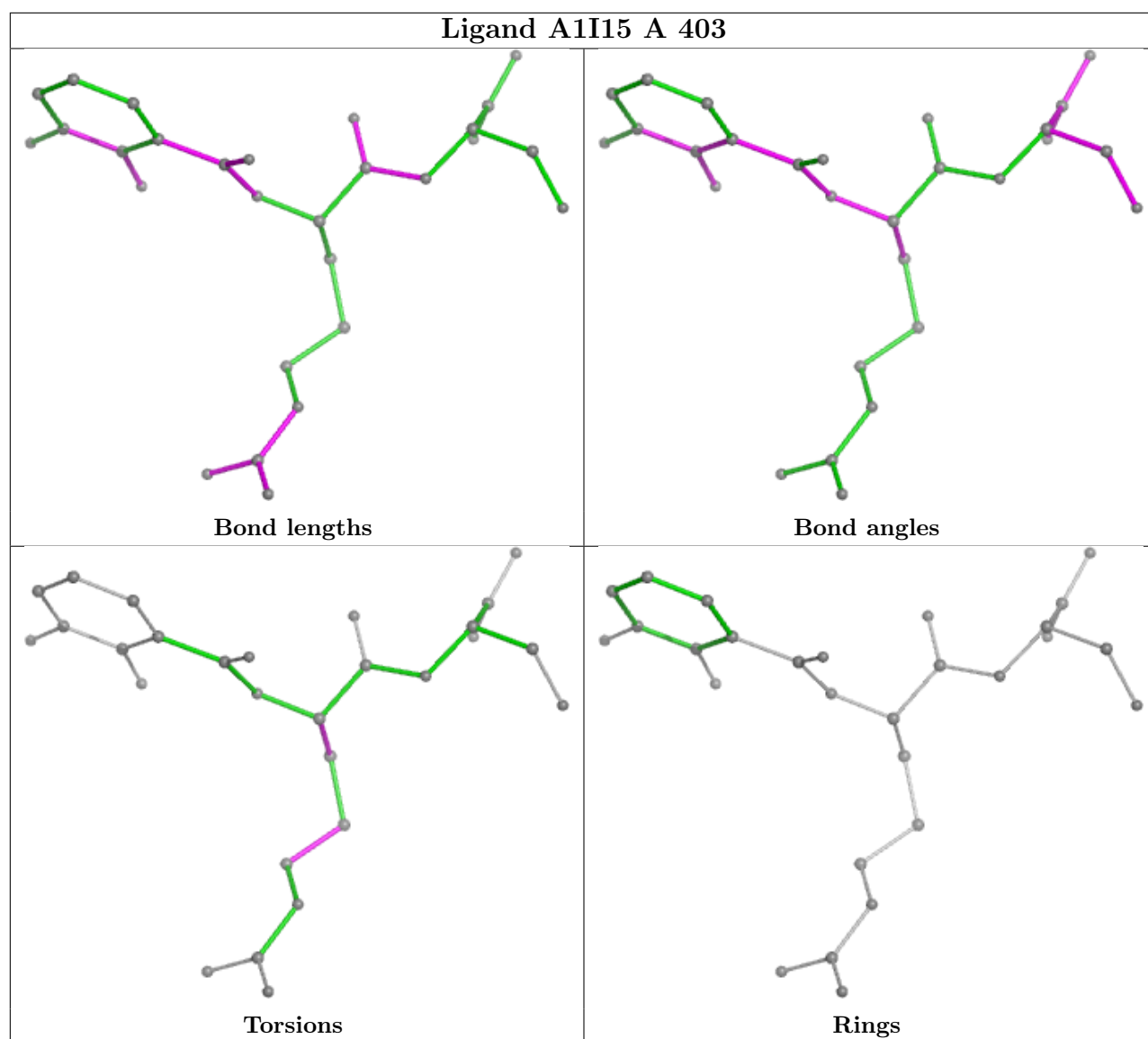


Ligand A1I15 B 402

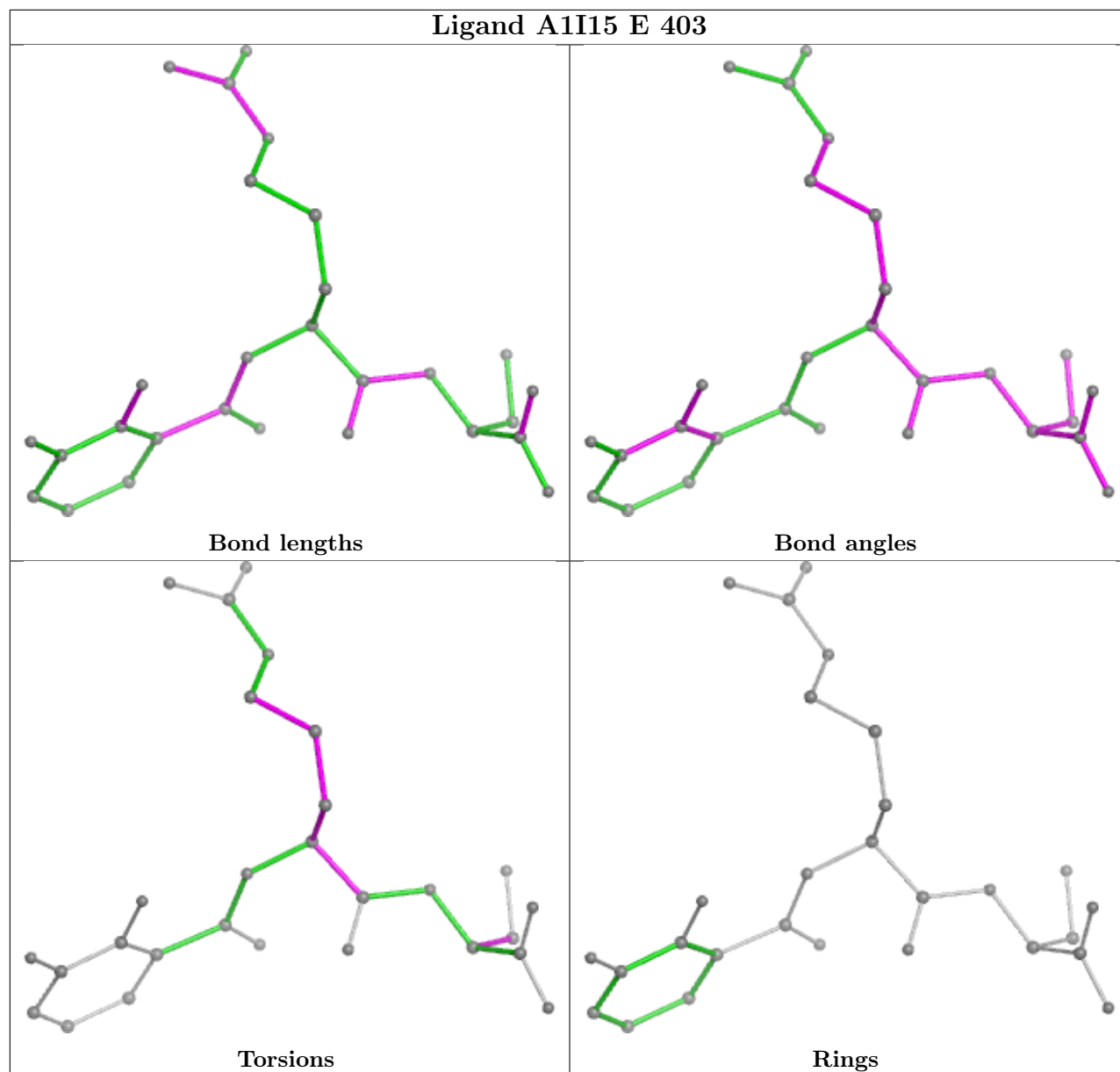


Ligand A1I15 D 402

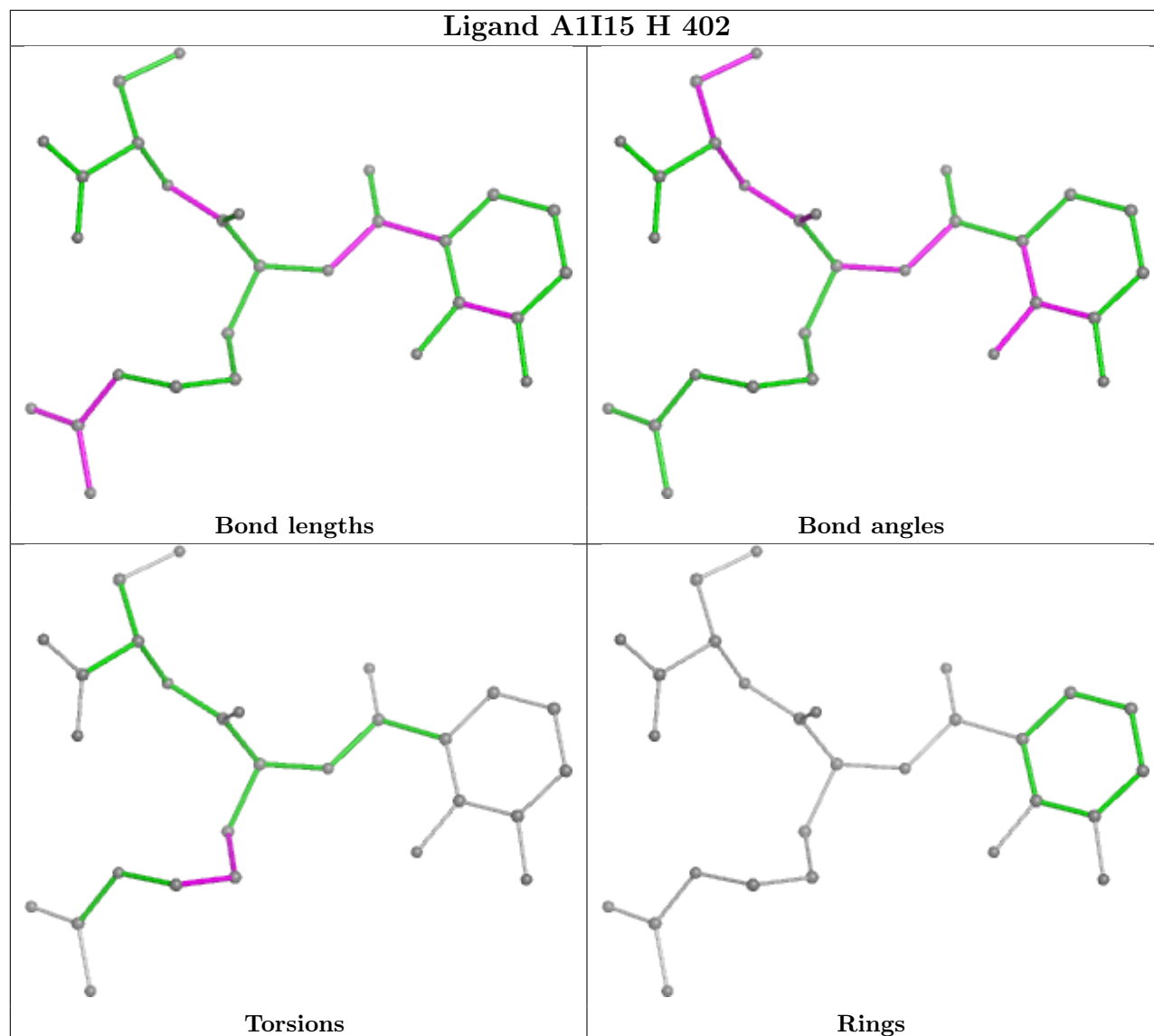




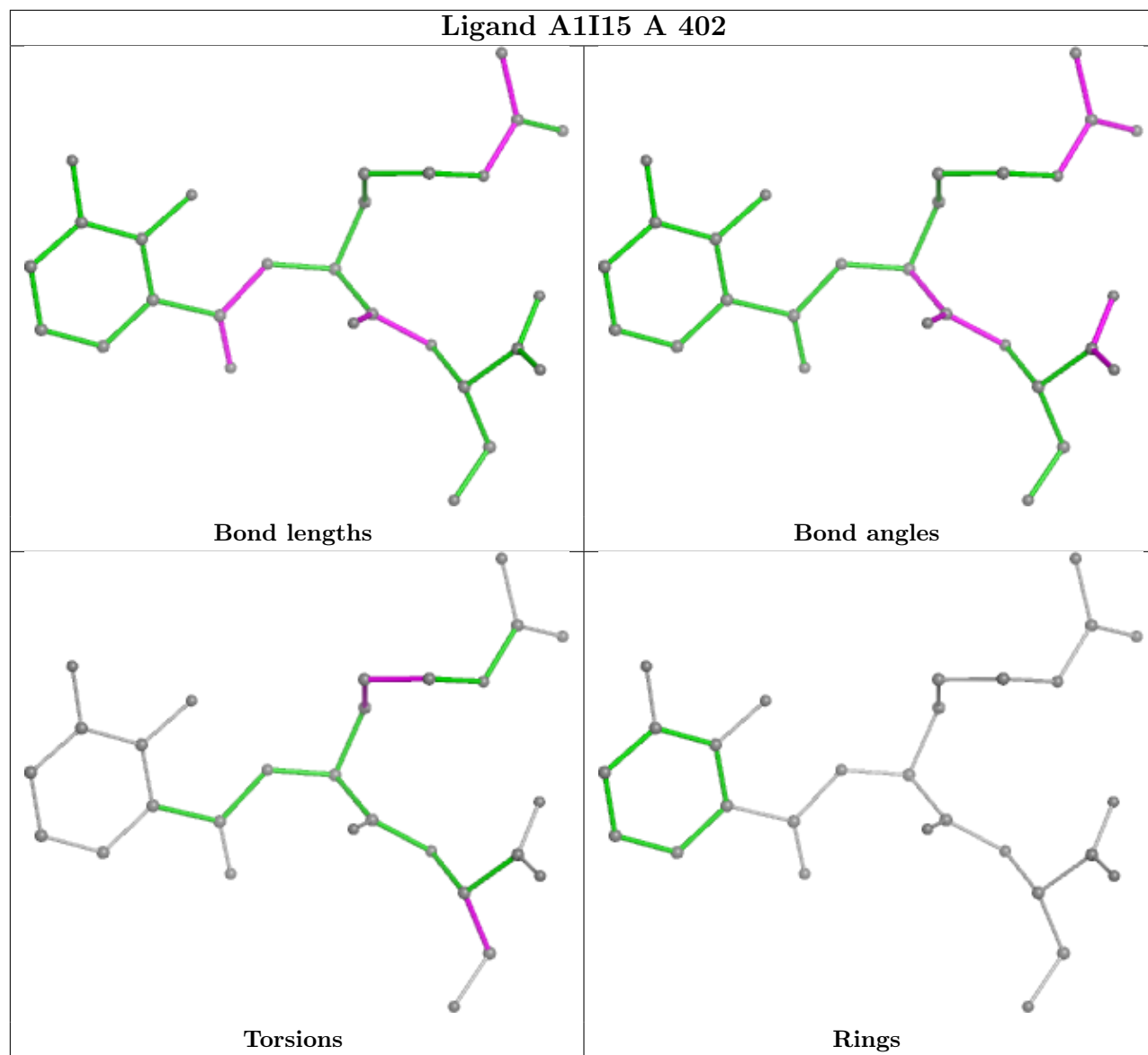
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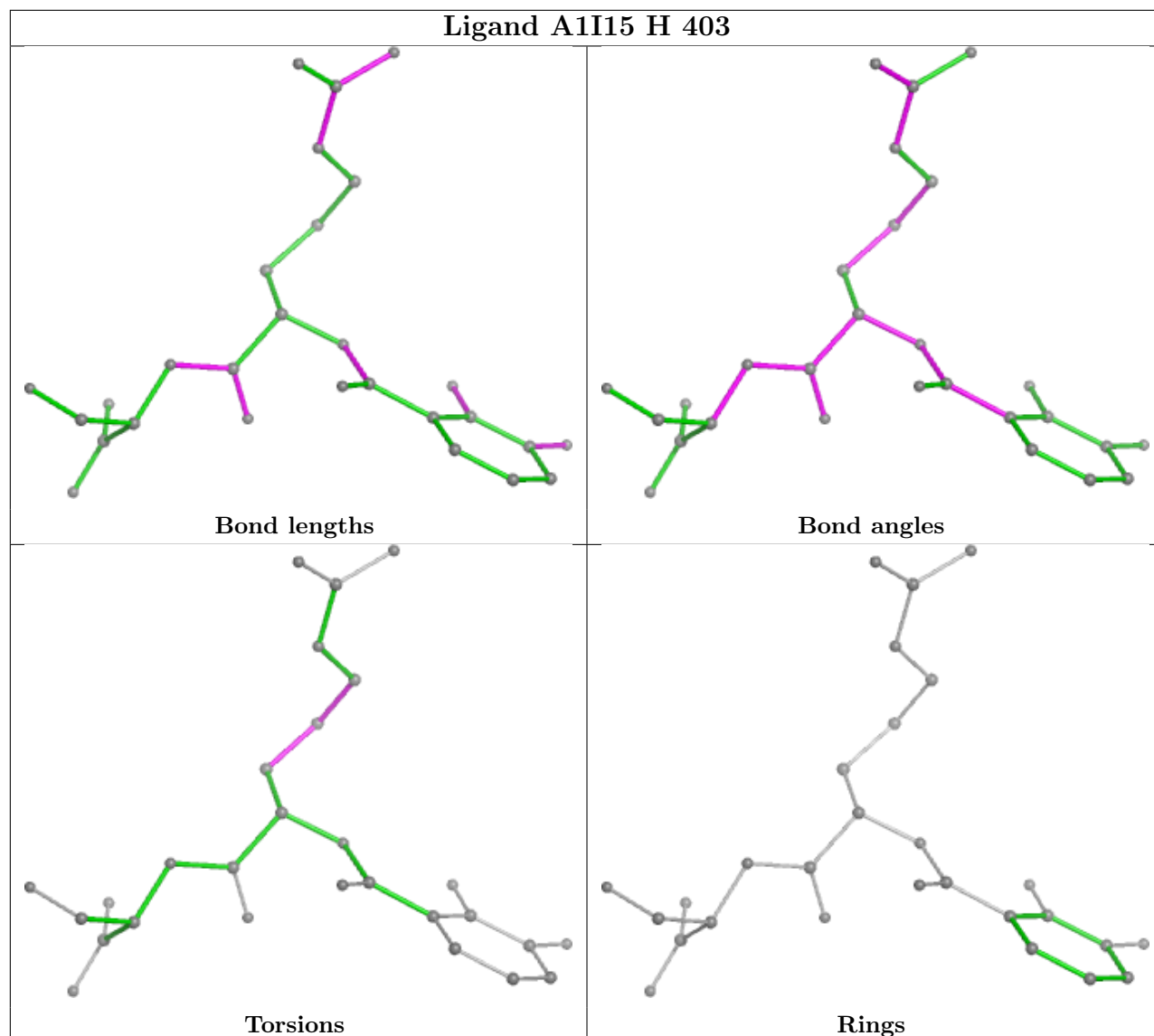
Ligand A1I15 H 402

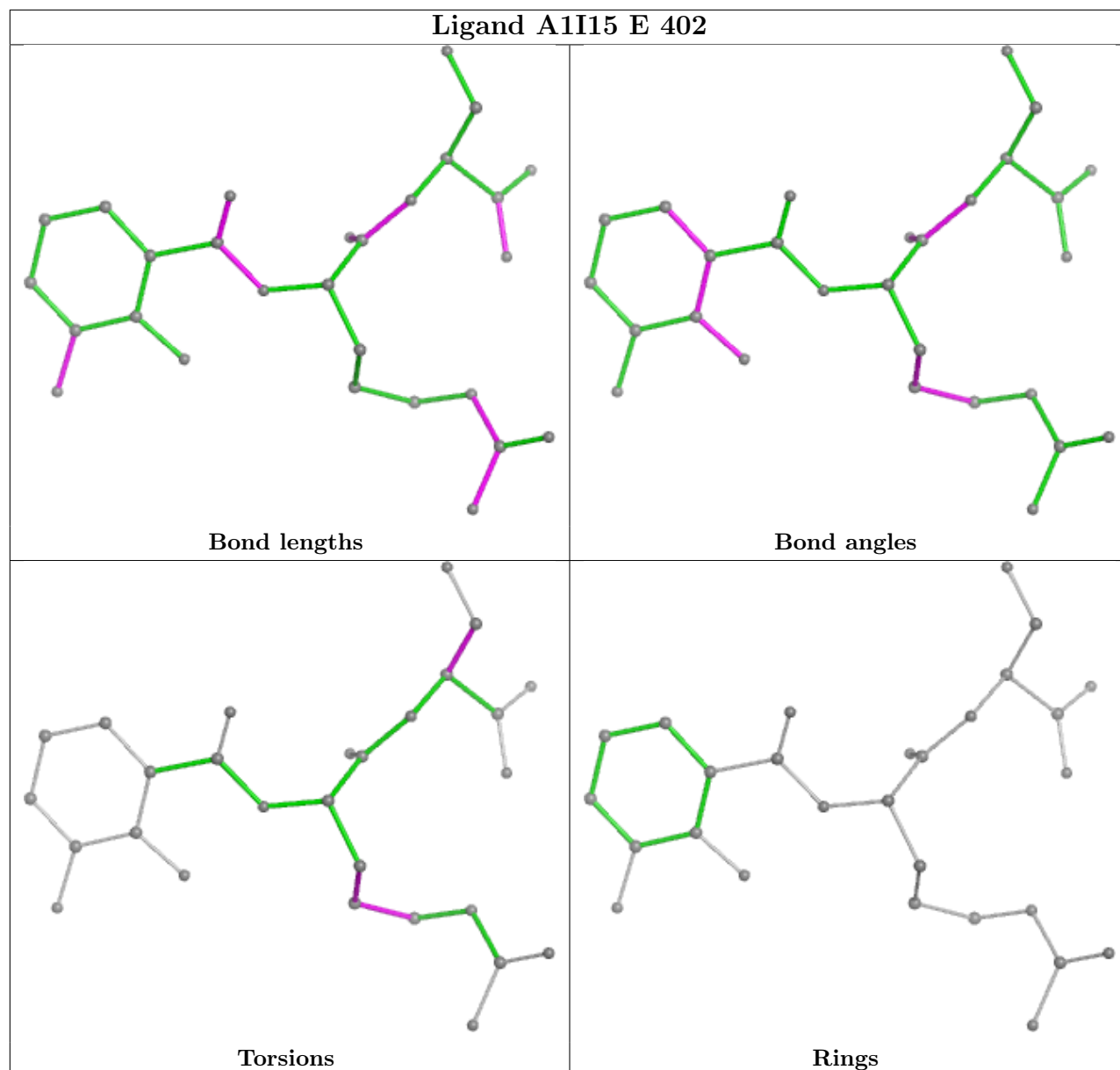


Ligand A1I15 A 402

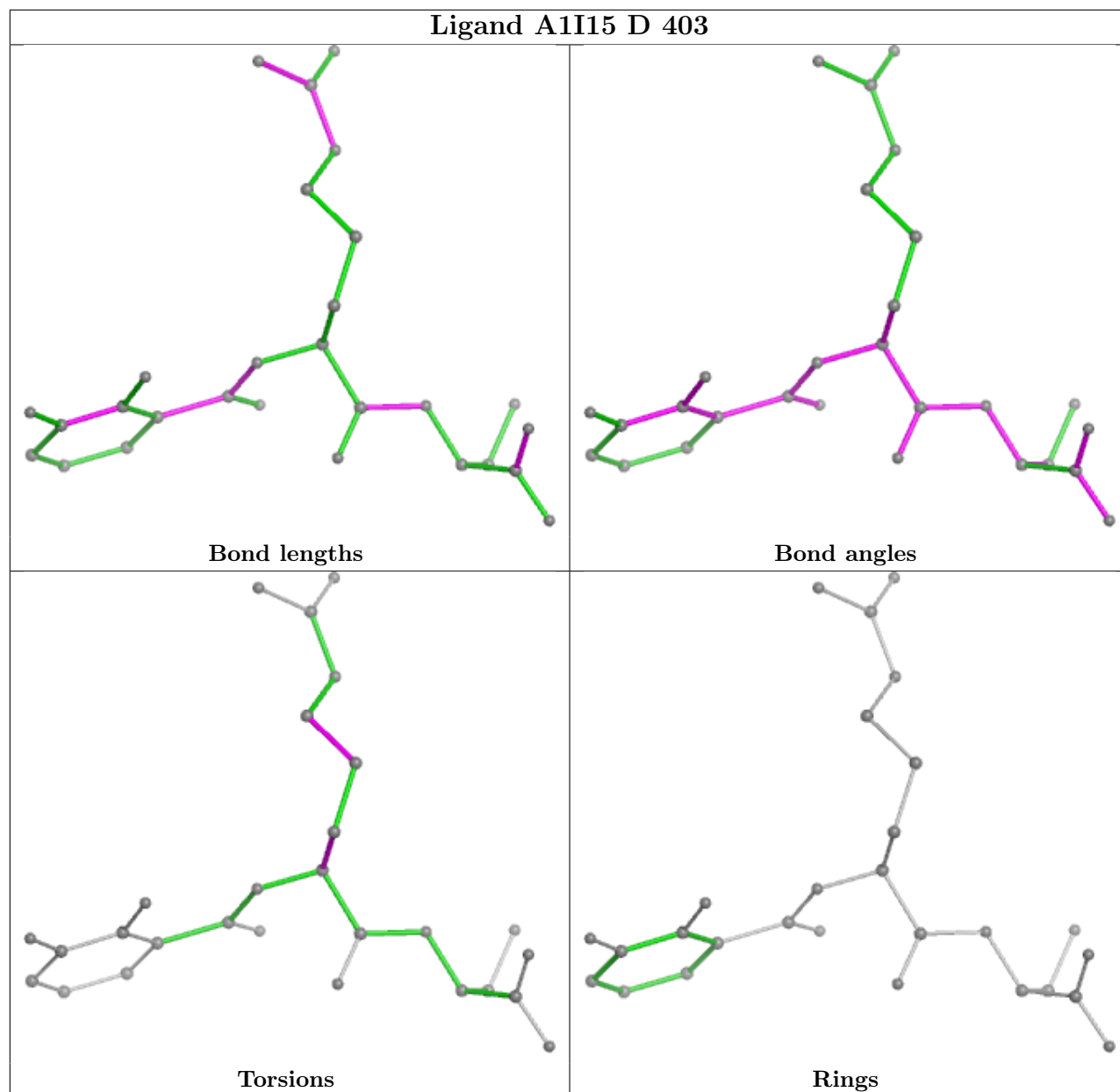


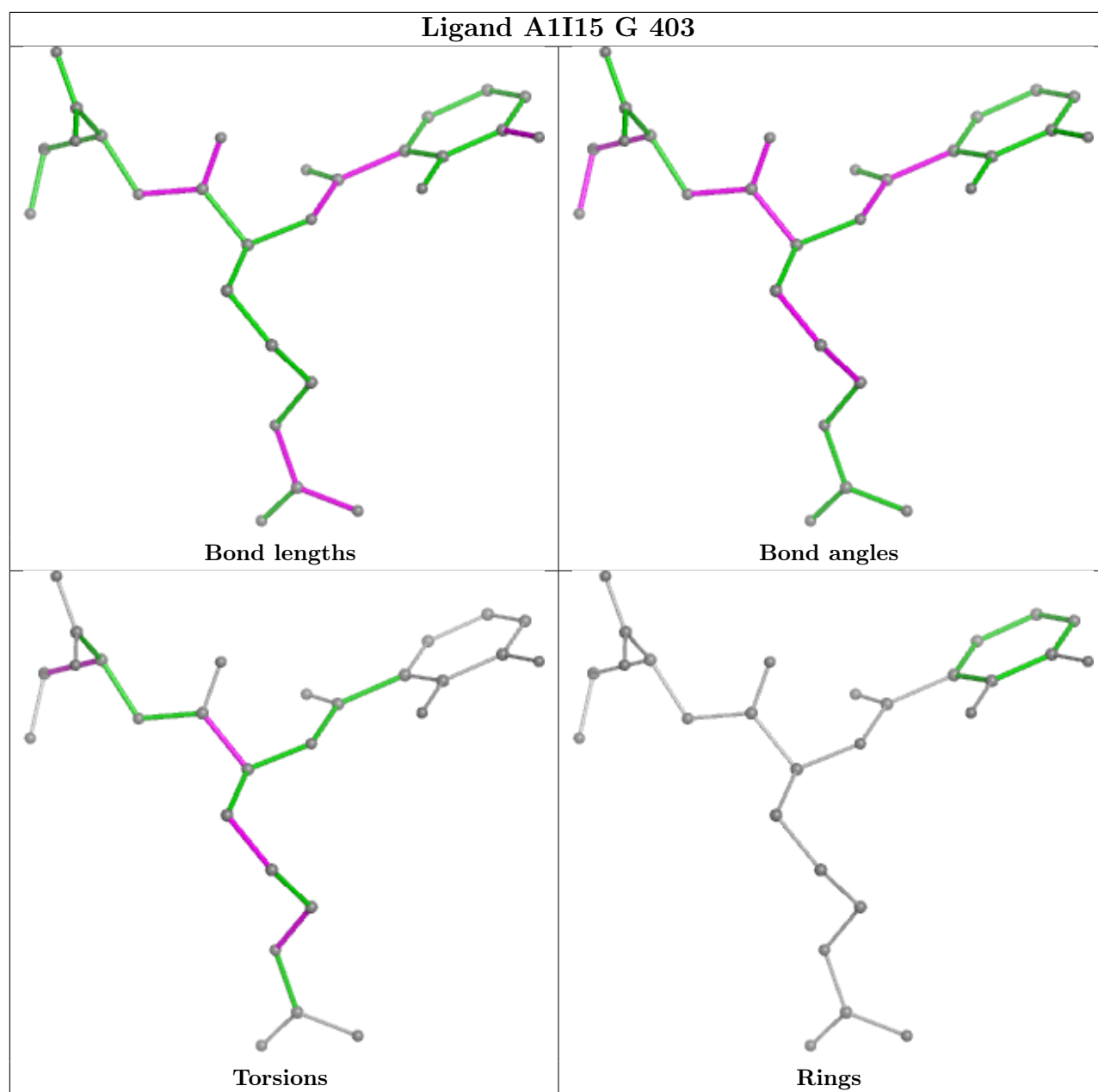
Ligand A1I15 H 403



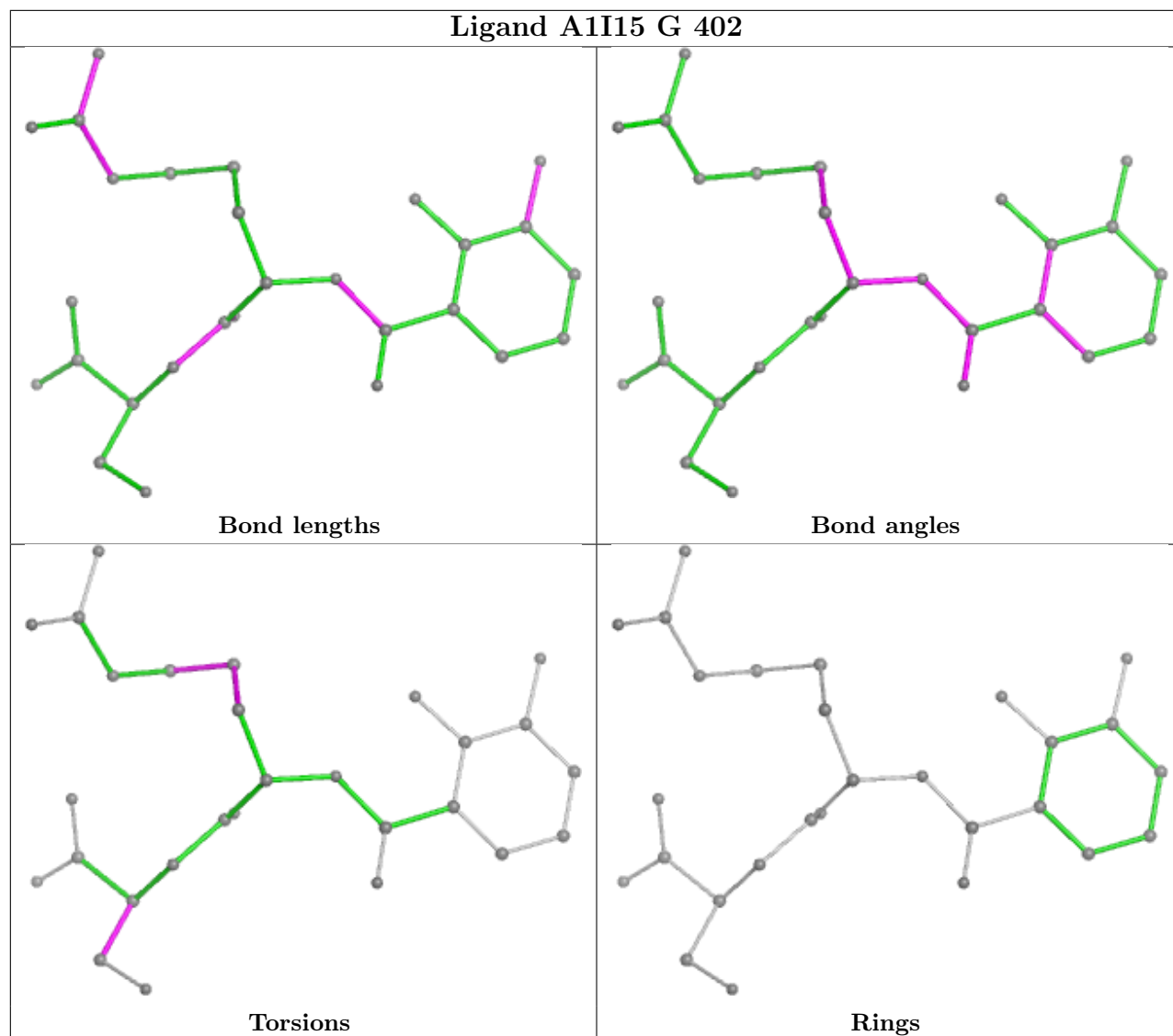


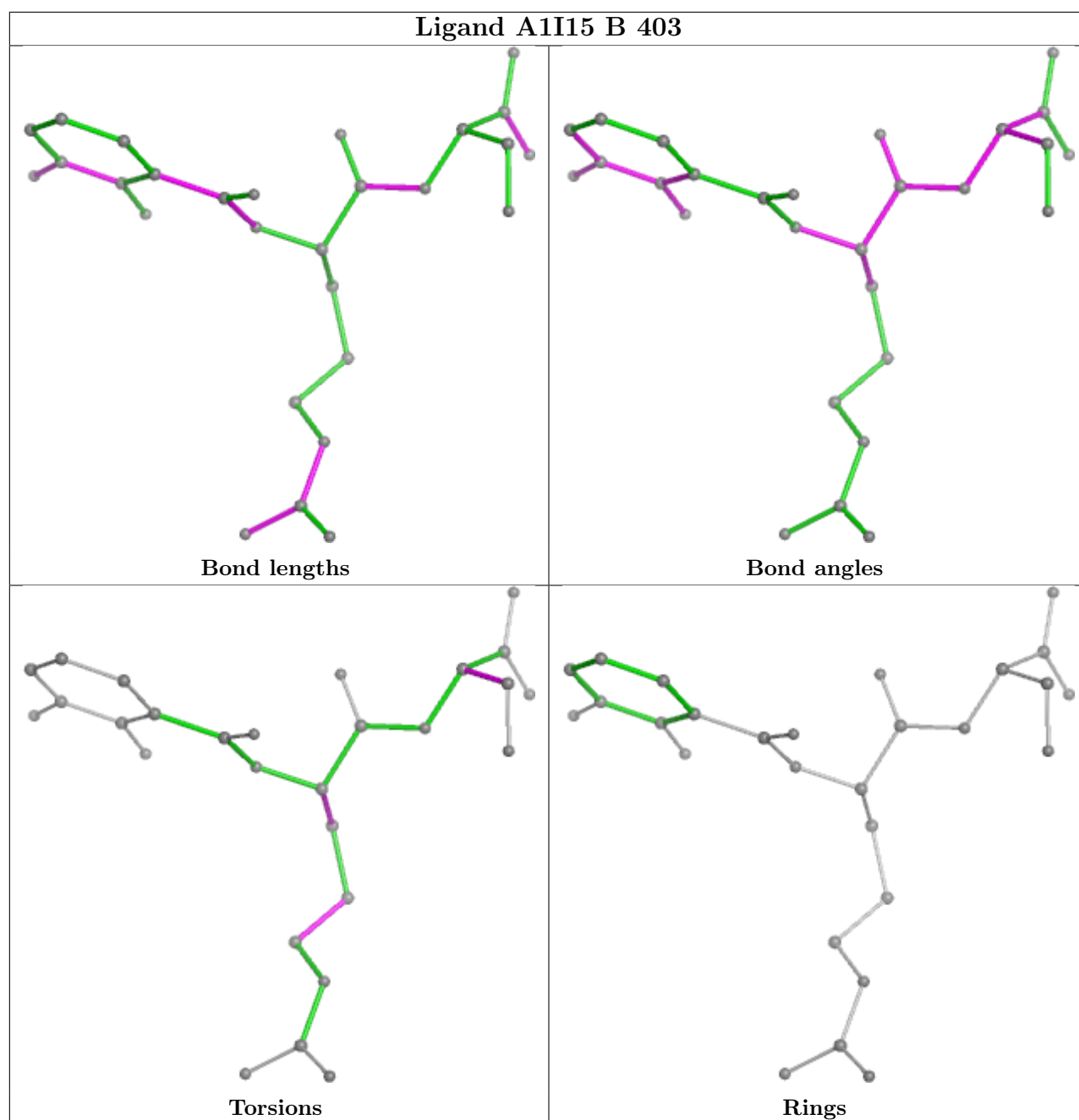
Ligand A1I15 D 403





Ligand A1I15 G 402





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	346/349 (99%)	0.57	18 (5%) 34 34	17, 36, 59, 83	0
1	B	346/349 (99%)	0.54	9 (2%) 57 60	19, 36, 55, 80	0
1	C	346/349 (99%)	0.44	7 (2%) 64 67	22, 34, 52, 76	0
1	D	347/349 (99%)	0.51	12 (3%) 47 49	17, 34, 53, 93	0
1	E	347/349 (99%)	0.34	7 (2%) 64 67	16, 30, 46, 79	1 (0%)
1	F	346/349 (99%)	0.51	10 (2%) 54 56	21, 36, 53, 90	0
1	G	347/349 (99%)	0.49	8 (2%) 61 63	23, 35, 51, 80	0
1	H	346/349 (99%)	0.97	39 (11%) 11 11	28, 46, 68, 91	0
All	All	2771/2792 (99%)	0.55	110 (3%) 43 44	16, 35, 58, 93	1 (0%)

The worst 5 of 110 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	249	ASN	5.5
1	H	2	ASN	5.0
1	H	11	VAL	4.7
1	H	12	VAL	4.6
1	H	3	ILE	4.5

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

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	SO4	D	405	5/5	0.79	0.25	30,30,30,30	0
4	SO4	F	405	5/5	0.82	0.11	72,72,75,77	0
4	SO4	G	405	5/5	0.83	0.11	52,59,62,71	0
4	SO4	G	407	5/5	0.83	0.15	57,59,67,72	0
4	SO4	C	404	5/5	0.84	0.14	47,50,57,58	0
4	SO4	C	407	5/5	0.84	0.10	62,65,75,78	0
4	SO4	C	406	5/5	0.85	0.13	43,47,57,58	0
4	SO4	H	404	5/5	0.85	0.12	54,62,65,67	0
3	A1I15	D	403	28/28	0.86	0.13	30,39,52,59	0
3	A1I15	H	403	28/28	0.86	0.11	30,43,51,59	0
4	SO4	E	405	5/5	0.87	0.14	61,62,71,73	0
3	A1I15	A	402	28/28	0.88	0.12	26,55,64,73	0
4	SO4	G	406	5/5	0.88	0.14	61,62,67,70	0
3	A1I15	B	402	28/28	0.88	0.12	25,60,73,81	0
4	SO4	A	405	5/5	0.88	0.14	53,58,67,76	0
3	A1I15	G	402	28/28	0.89	0.10	18,42,50,54	0
3	A1I15	H	402	28/28	0.89	0.10	26,39,55,59	0
4	SO4	F	404	5/5	0.89	0.10	61,62,64,68	0
3	A1I15	B	403	28/28	0.89	0.11	29,38,52,60	0
3	A1I15	D	402	28/28	0.89	0.11	24,49,65,70	0
3	A1I15	A	403	28/28	0.89	0.12	32,45,51,55	0
3	A1I15	E	402	28/28	0.89	0.12	12,36,65,70	0
3	A1I15	F	402	28/28	0.89	0.12	16,42,50,52	0
3	A1I15	C	402	28/28	0.90	0.11	18,45,57,59	0
3	A1I15	G	403	28/28	0.90	0.10	26,36,48,54	0
4	SO4	A	404	5/5	0.90	0.10	51,53,59,63	0
3	A1I15	F	403	28/28	0.91	0.09	27,36,50,53	0
3	A1I15	C	403	28/28	0.92	0.09	29,34,48,50	0
4	SO4	D	404	5/5	0.92	0.11	49,52,59,59	0
3	A1I15	E	403	28/28	0.93	0.10	16,25,42,49	0
4	SO4	E	404	5/5	0.94	0.08	49,53,59,71	0
4	SO4	G	404	5/5	0.94	0.08	44,46,47,48	0
4	SO4	B	404	5/5	0.95	0.09	41,42,44,44	0
4	SO4	C	405	5/5	0.96	0.07	36,37,39,39	0
2	FE	E	401	1/1	0.97	0.02	15,15,15,15	0
2	FE	G	401	1/1	0.99	0.02	28,28,28,28	0
2	FE	B	401	1/1	0.99	0.03	30,30,30,30	0

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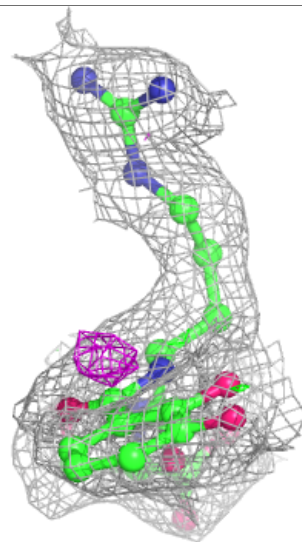
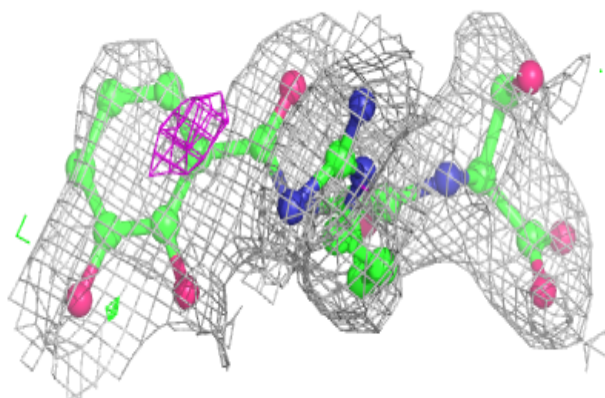
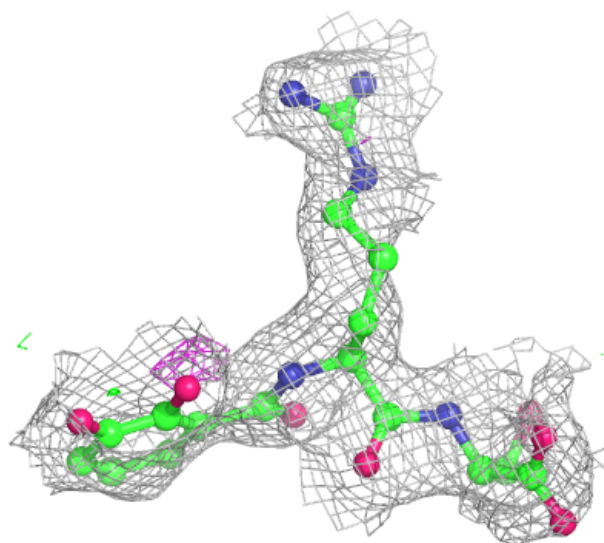
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	FE	D	401	1/1	0.99	0.03	28,28,28,28	0
2	FE	A	401	1/1	0.99	0.03	35,35,35,35	0
2	FE	F	401	1/1	0.99	0.03	27,27,27,27	0
2	FE	C	401	1/1	1.00	0.02	25,25,25,25	0
2	FE	H	401	1/1	1.00	0.02	33,33,33,33	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.

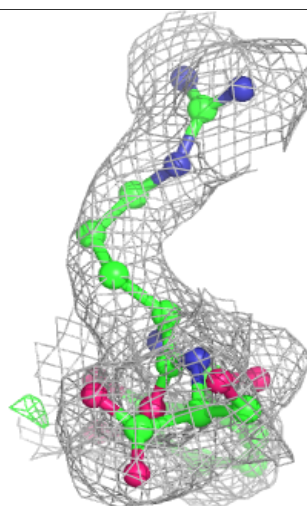
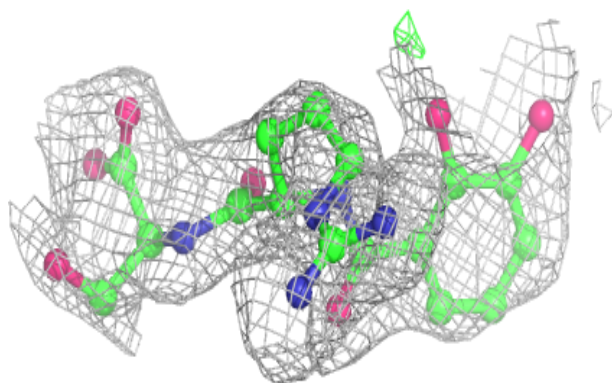
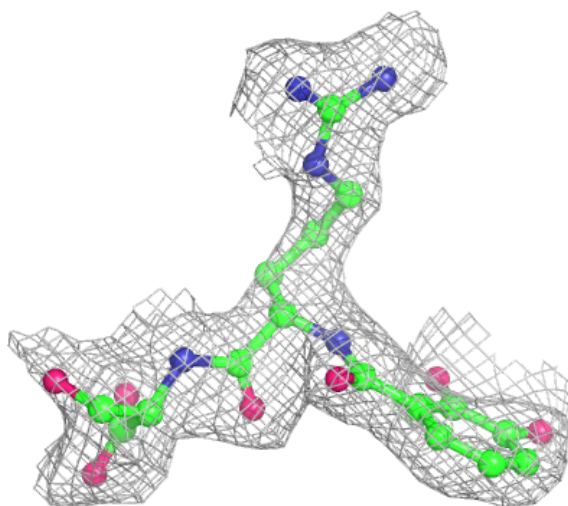
Electron density around A1I15 D 403:

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



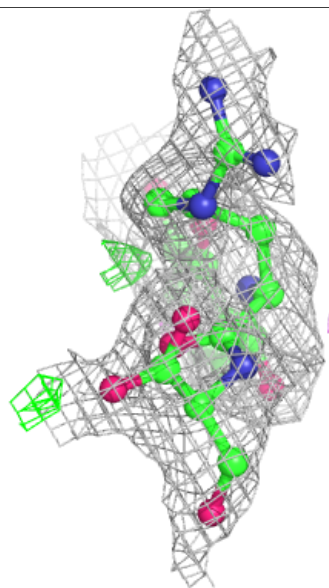
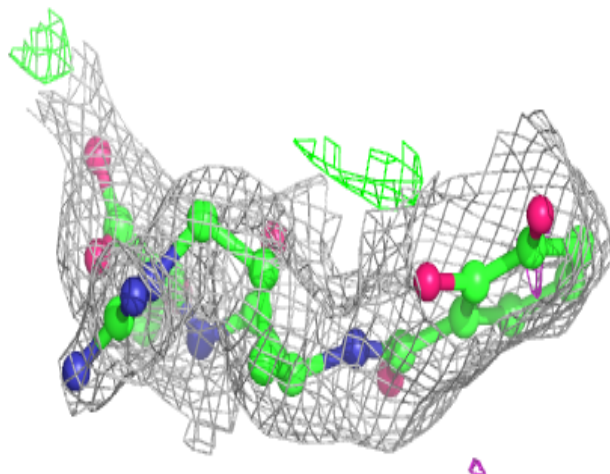
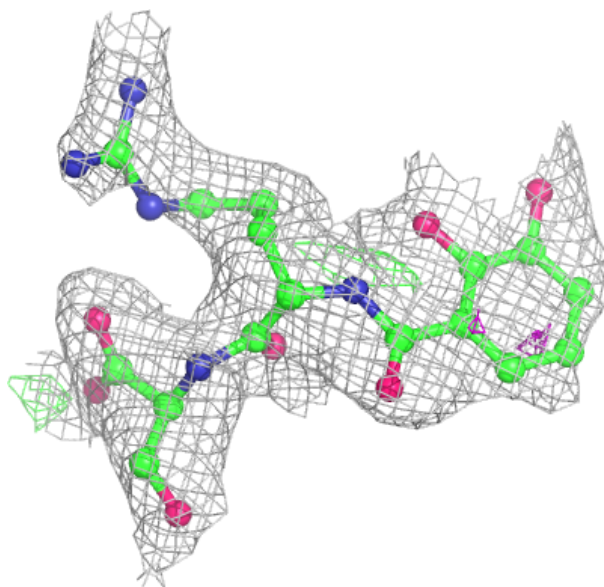
Electron density around A1I15 H 403:

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



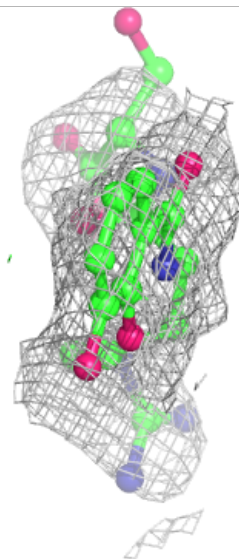
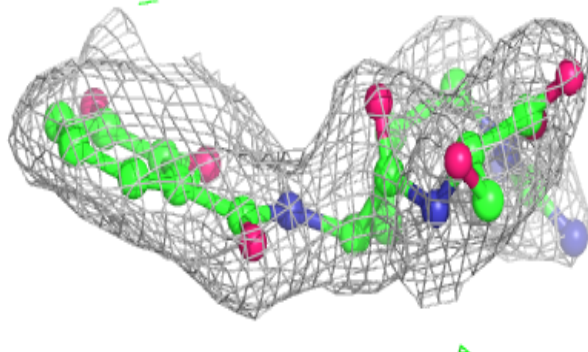
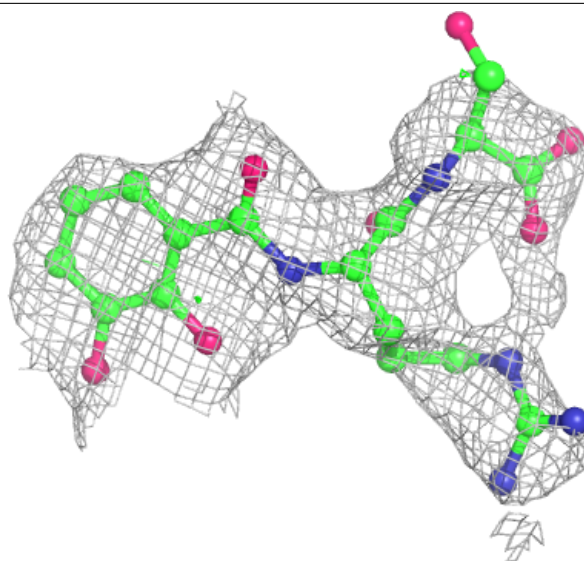
Electron density around A1I15 A 402:

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and green (positive)



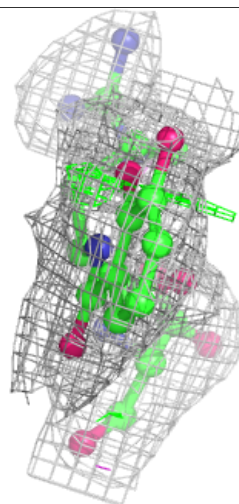
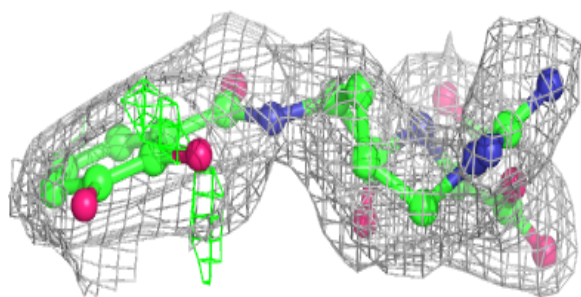
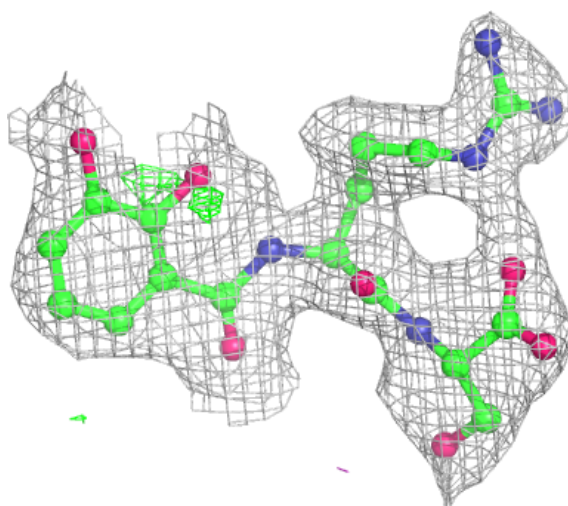
Electron density around A1I15 B 402:

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and green (positive)



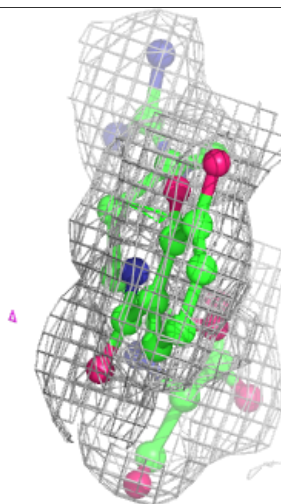
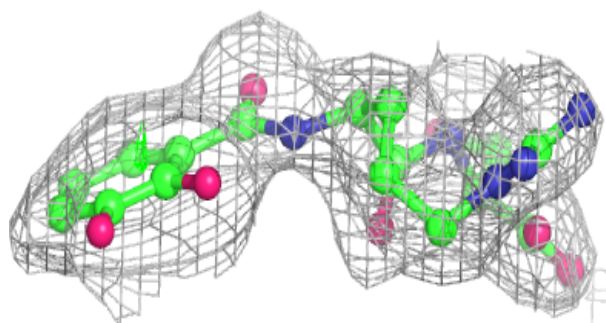
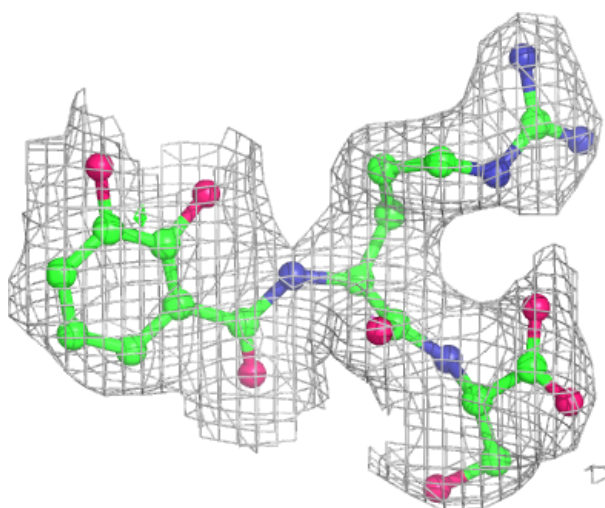
Electron density around A1I15 G 402:

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



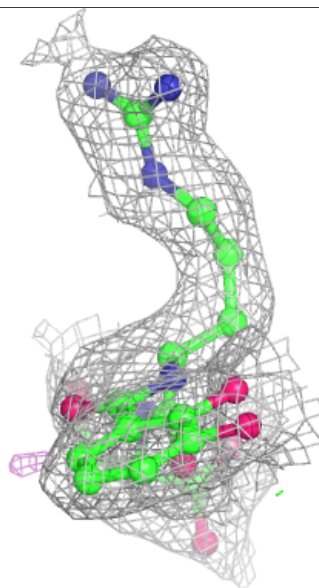
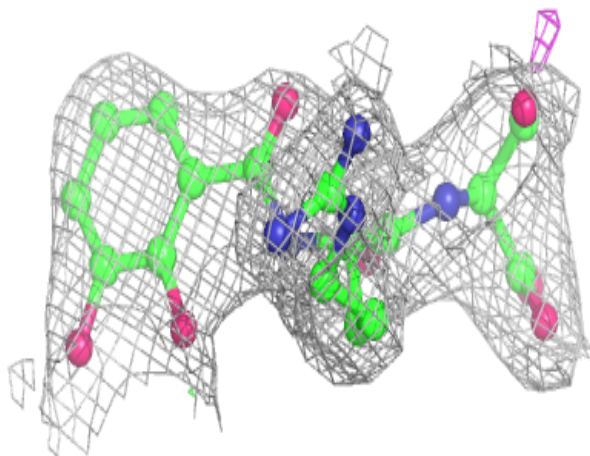
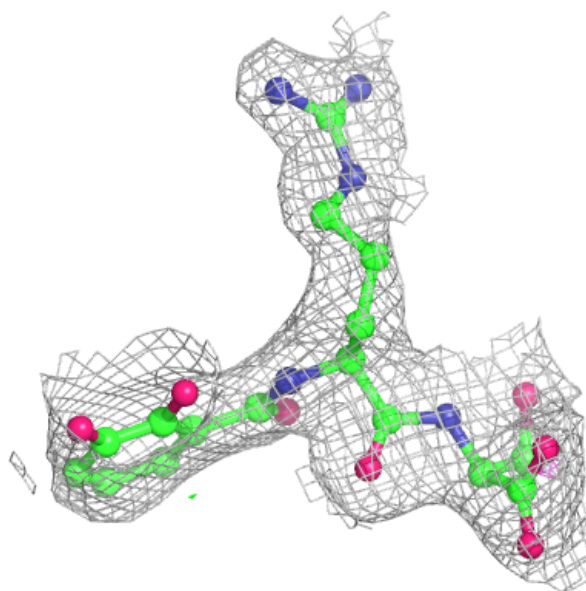
Electron density around A1I15 H 402:

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



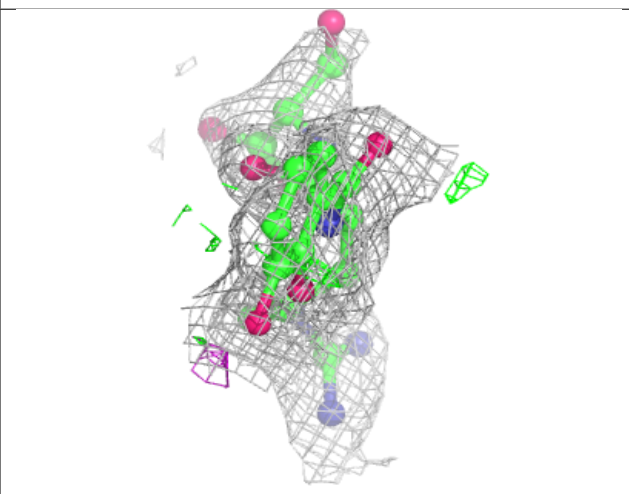
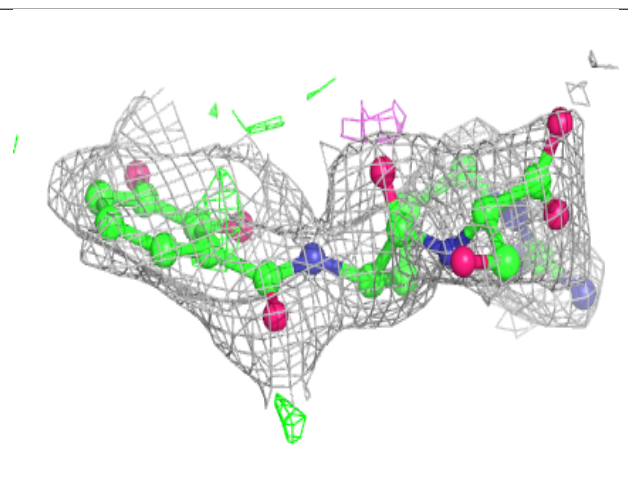
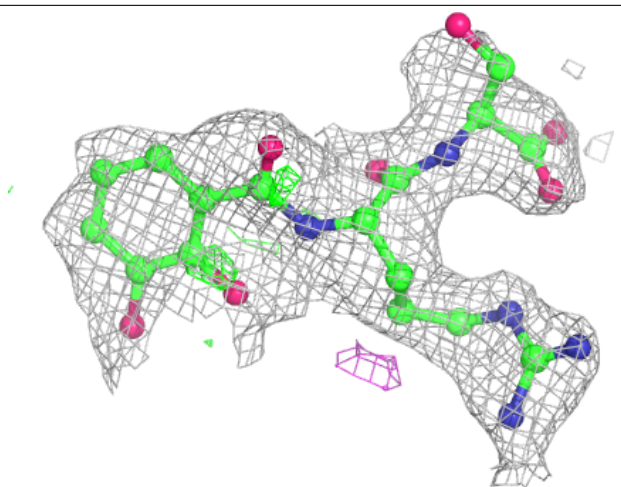
Electron density around A1I15 B 403:

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



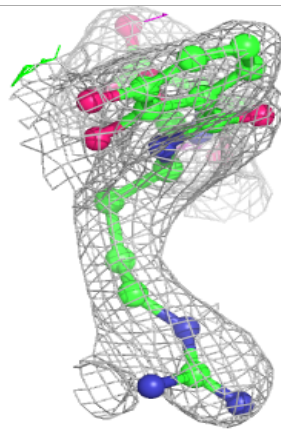
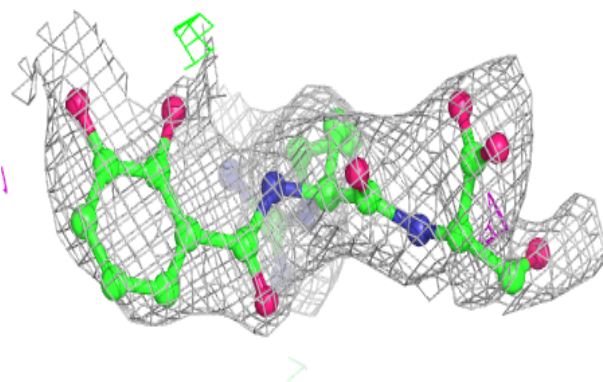
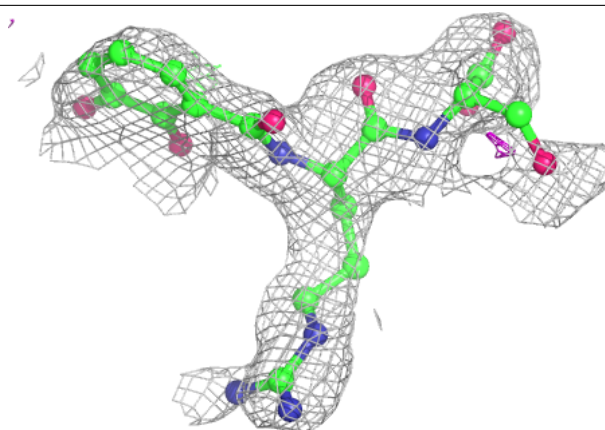
Electron density around A1I15 D 402:

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



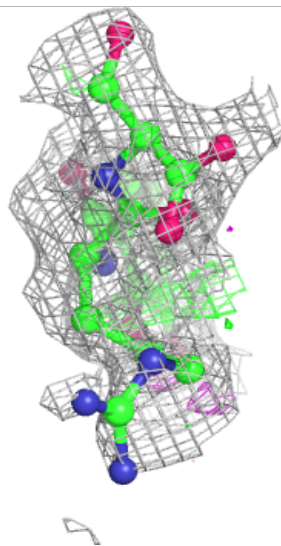
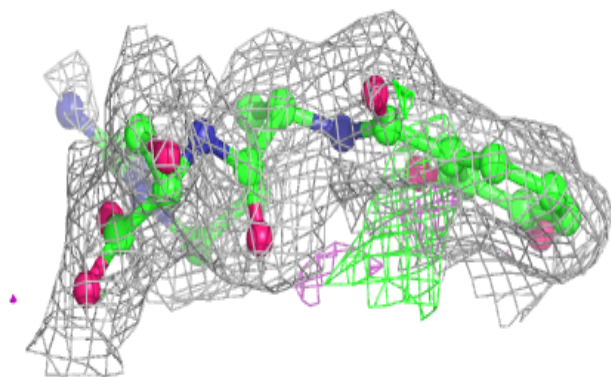
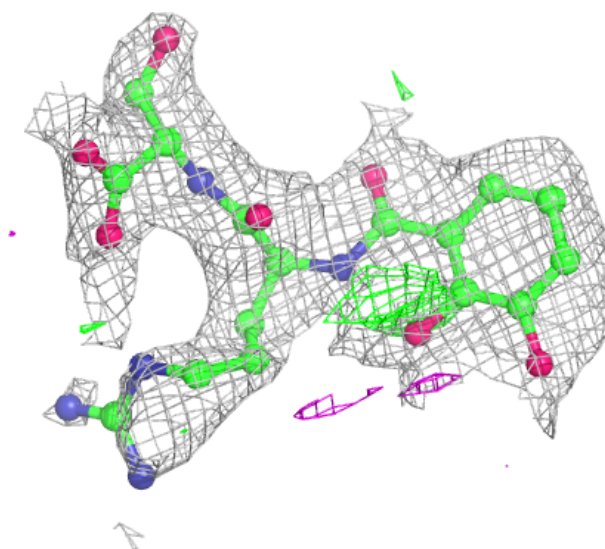
Electron density around A1I15 A 403:

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



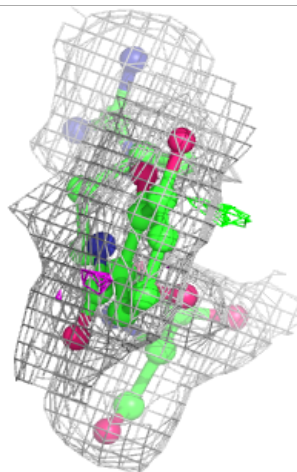
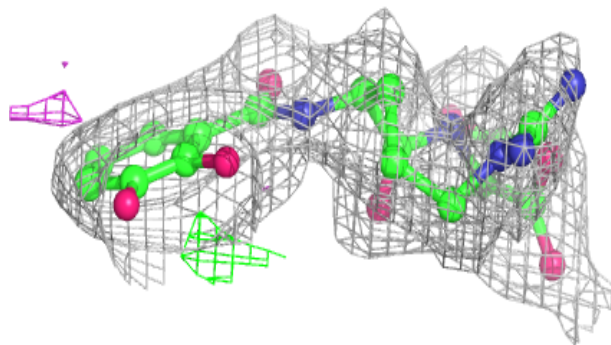
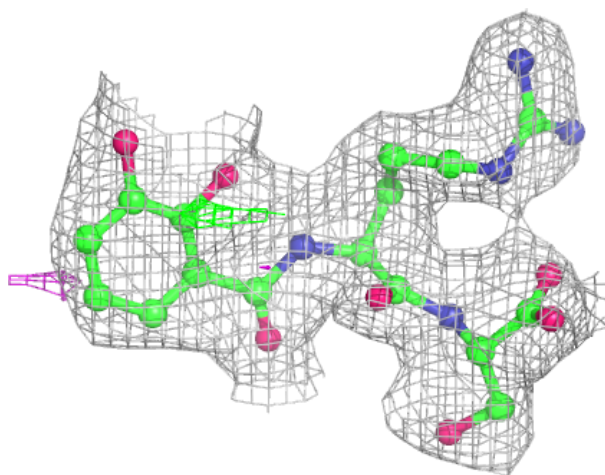
Electron density around A1I15 E 402:

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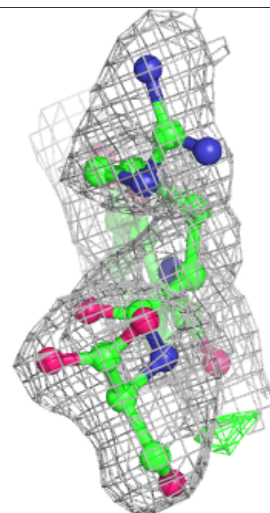
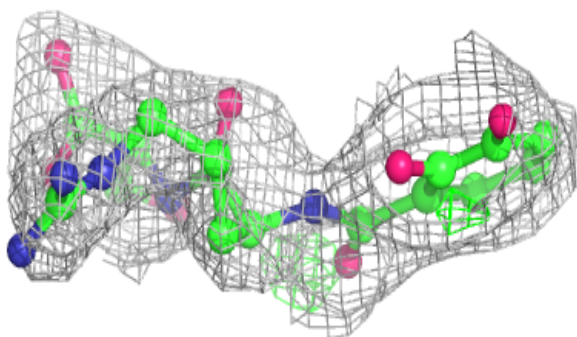
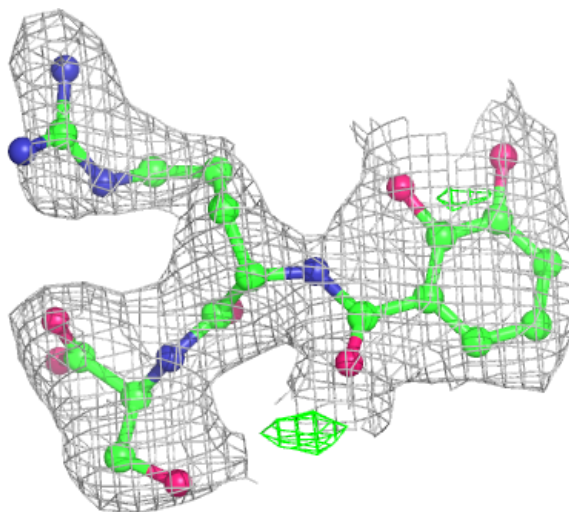
Electron density around A1I15 F 402:

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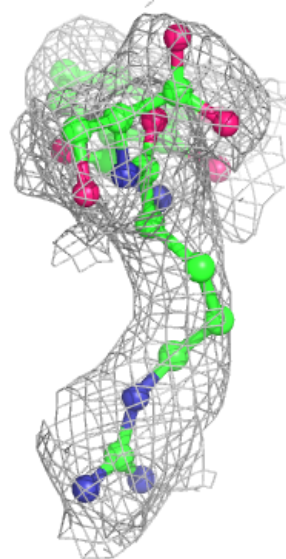
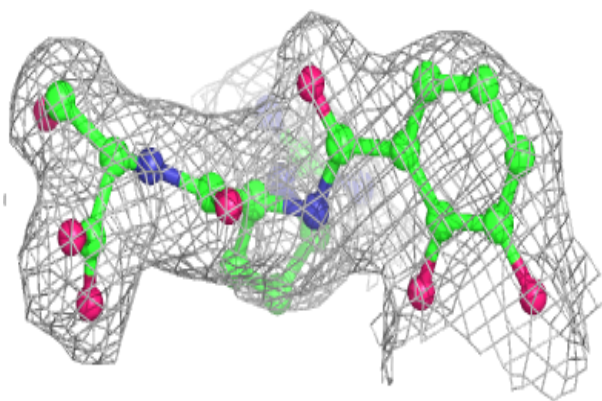
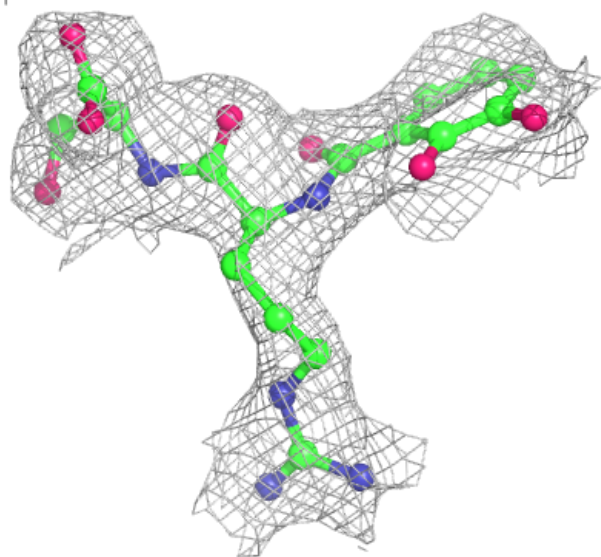
Electron density around A1I15 C 402:

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and green (positive)



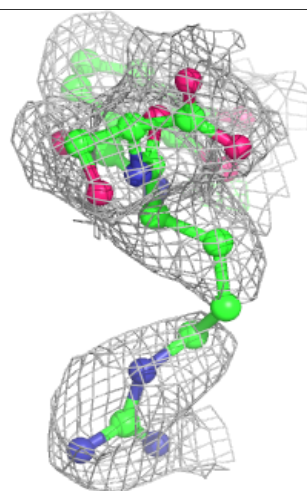
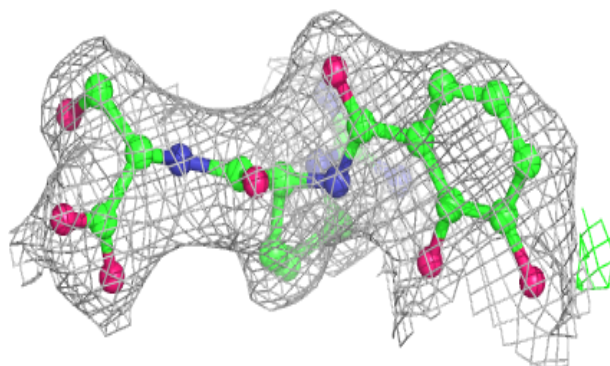
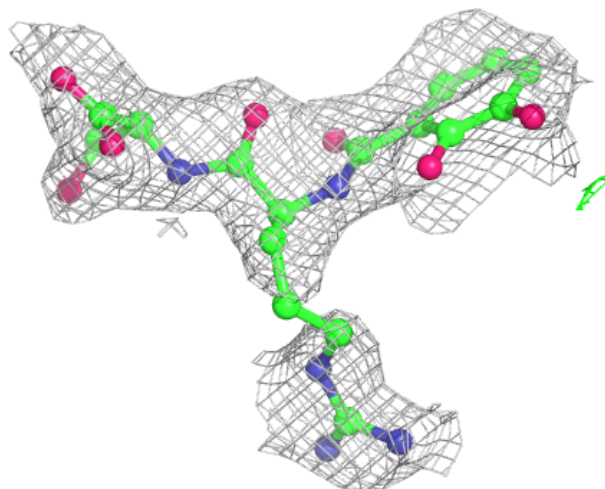
Electron density around A1I15 G 403:

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and green (positive)



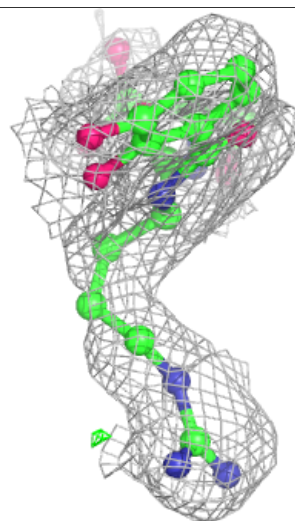
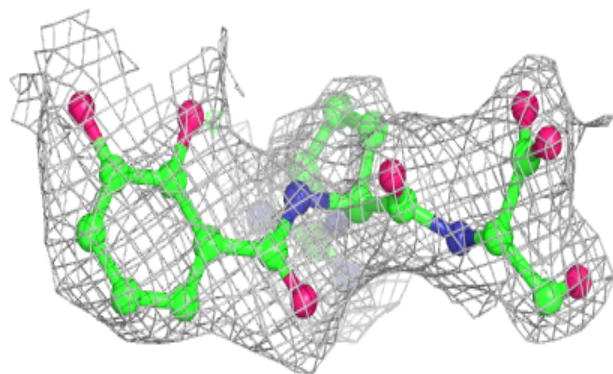
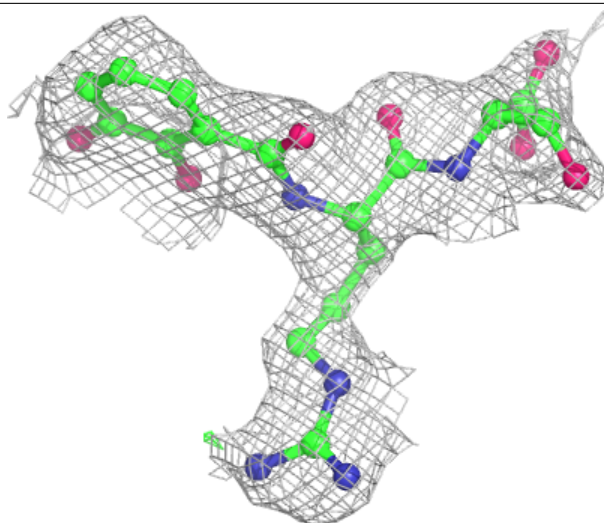
Electron density around A1I15 F 403:

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and green (positive)



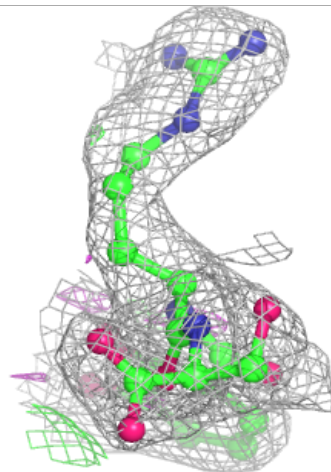
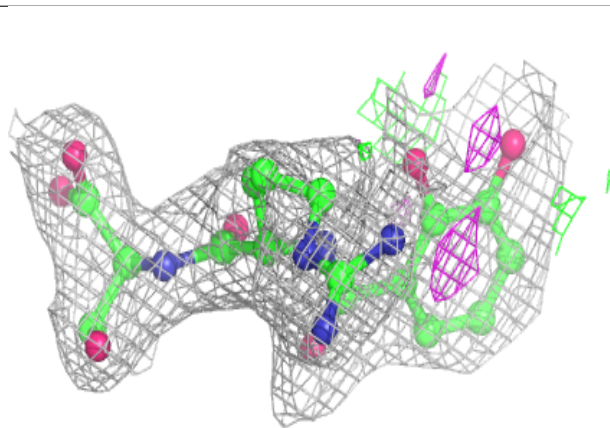
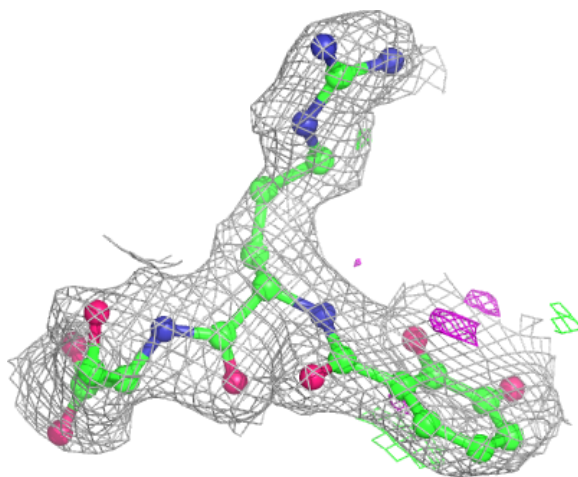
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and green (positive)



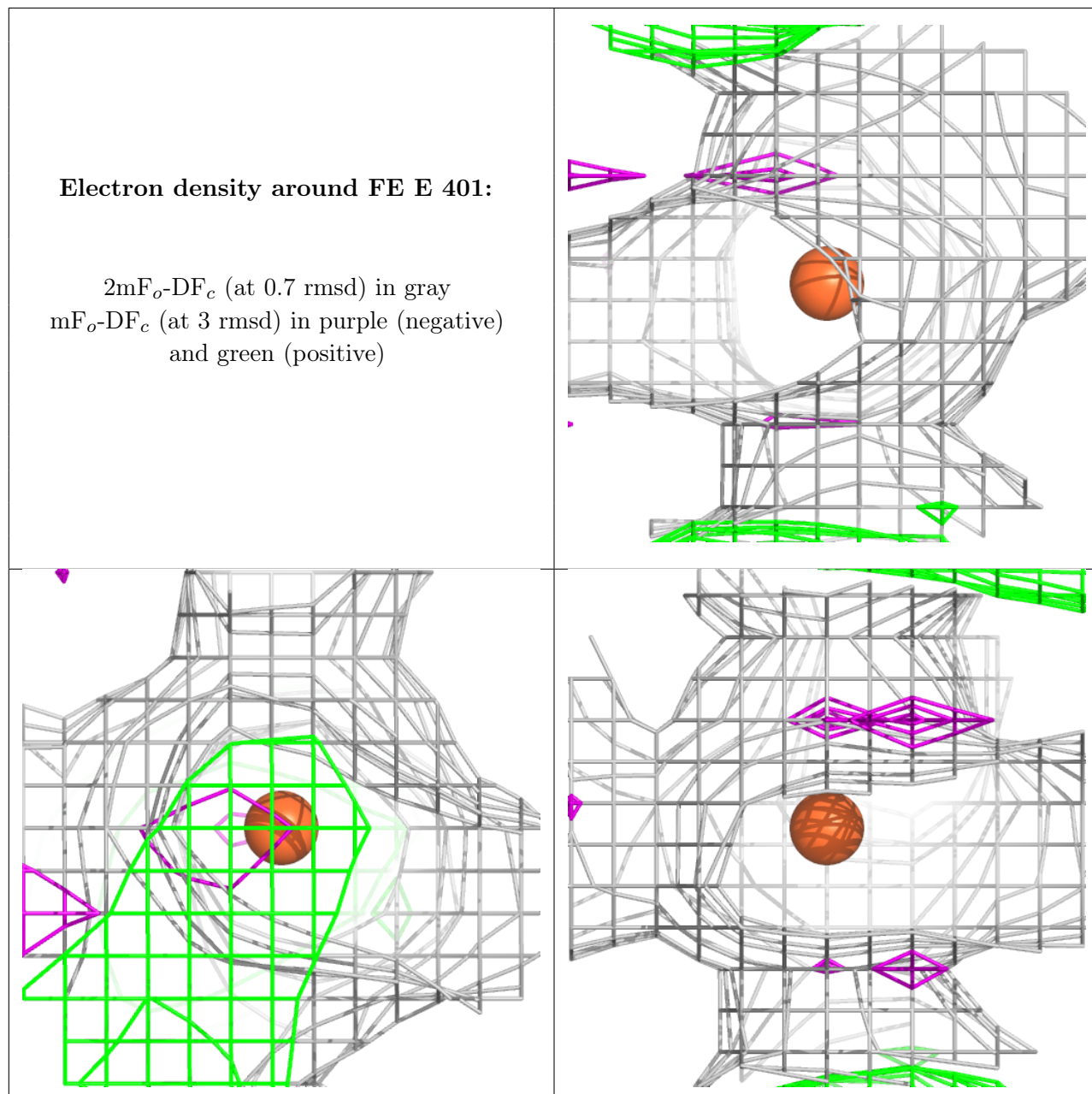
Electron density around A1I15 E 403:

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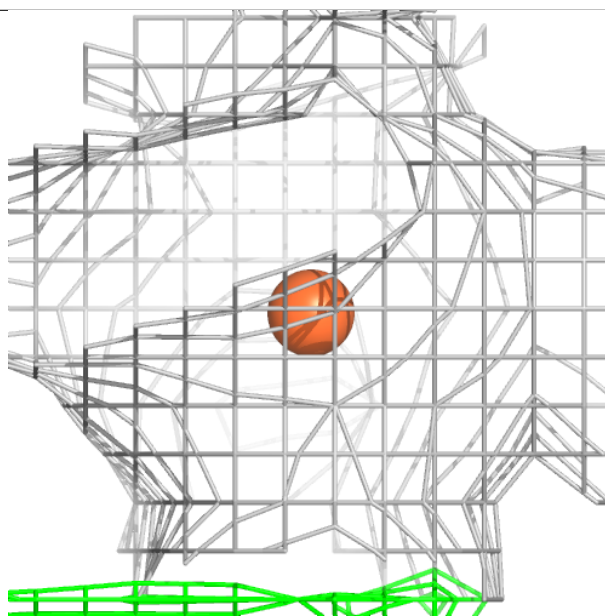
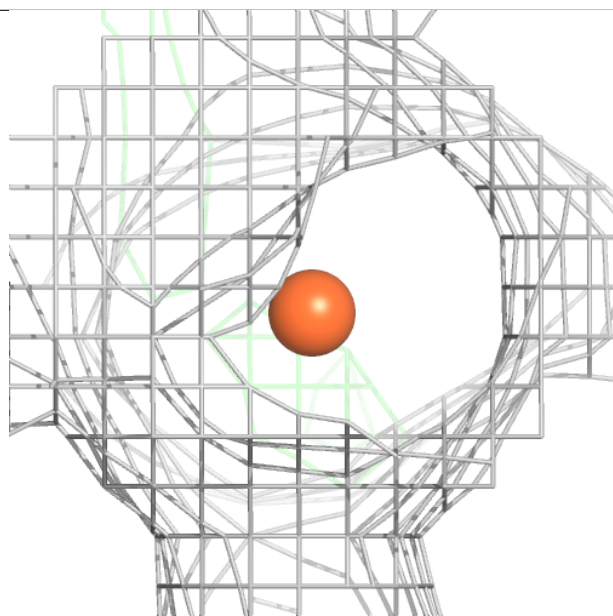
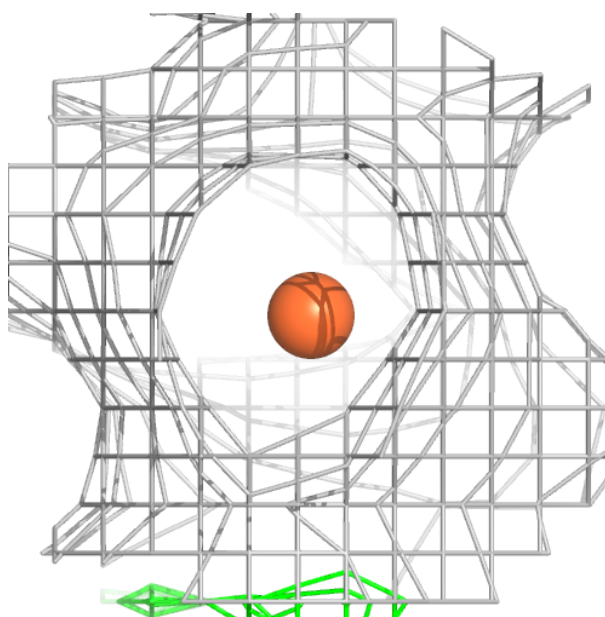
Electron density around FE E 401:

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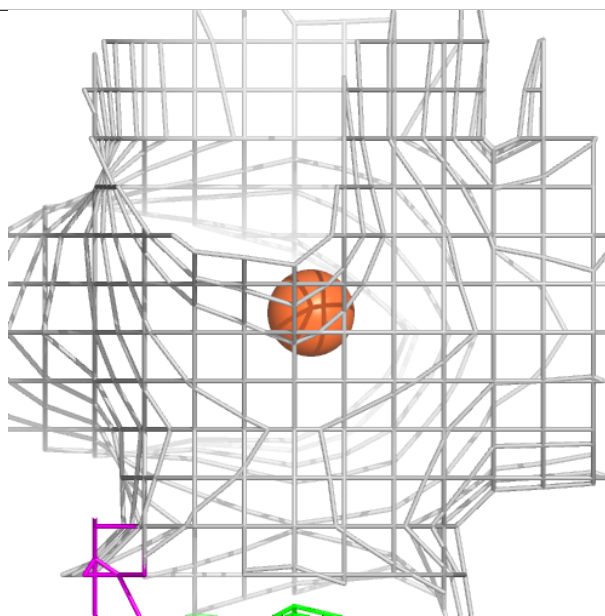
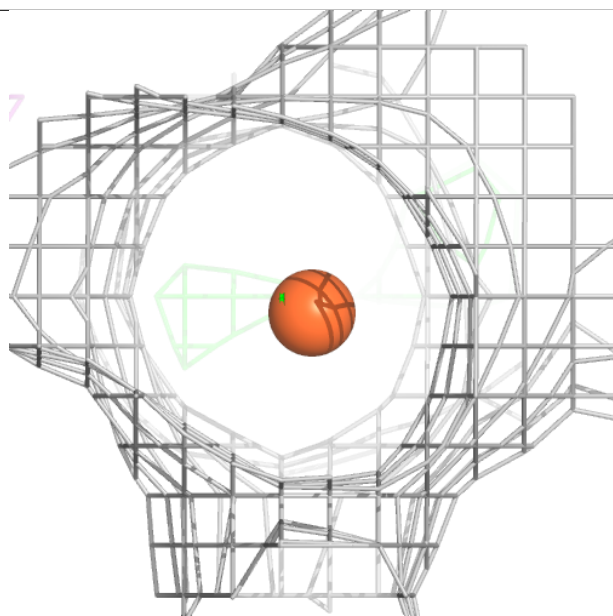
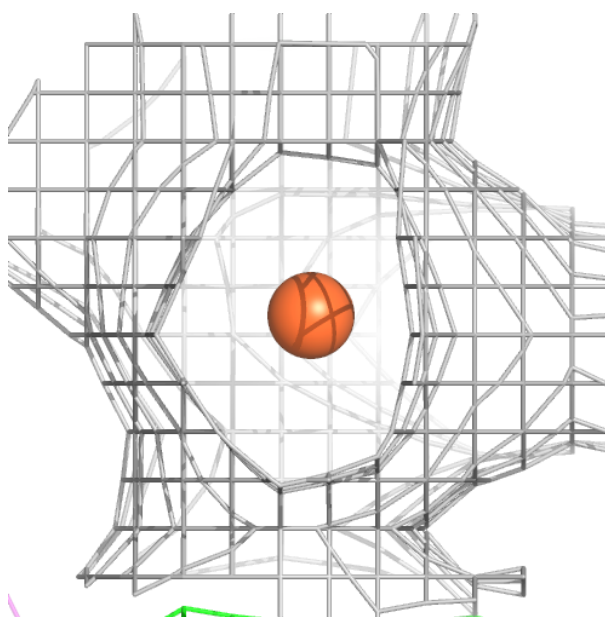
Electron density around FE G 401:

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and green (positive)



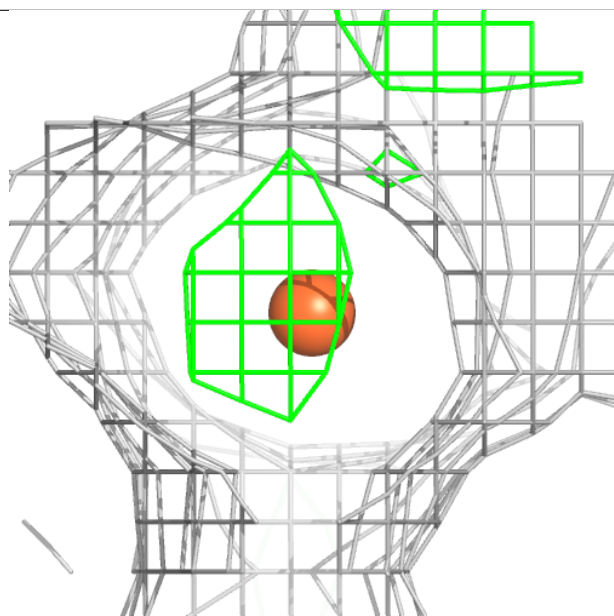
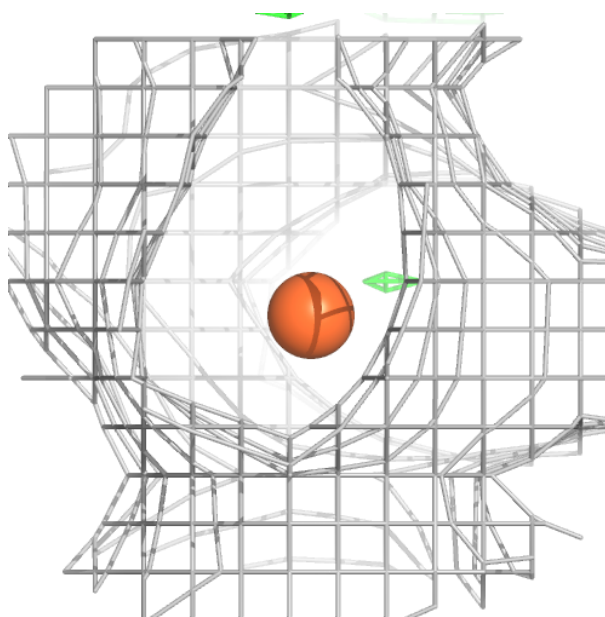
Electron density around FE B 401:

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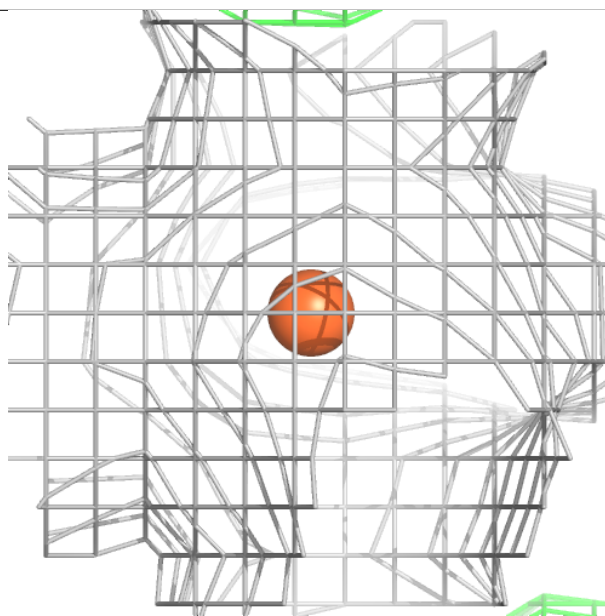
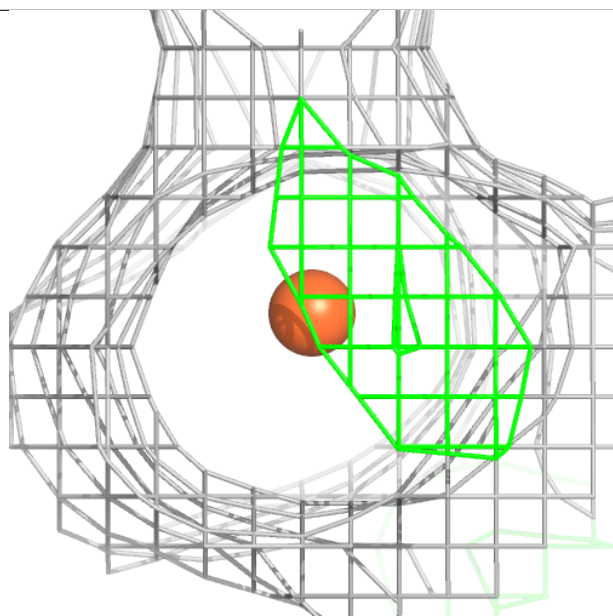
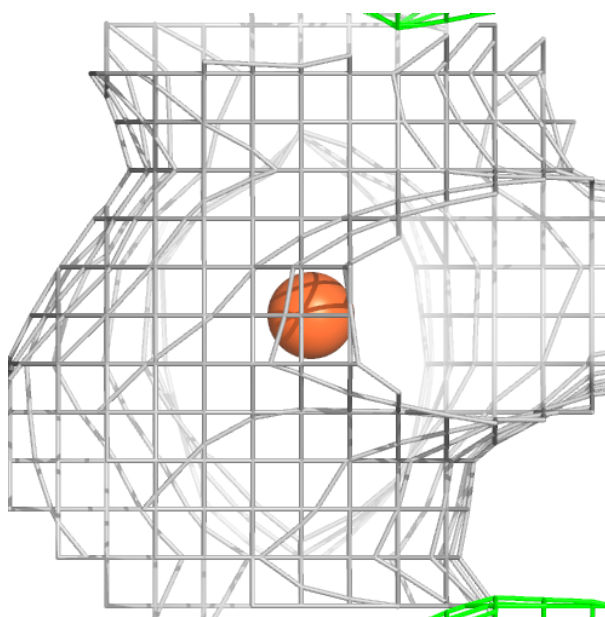
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and green (positive)



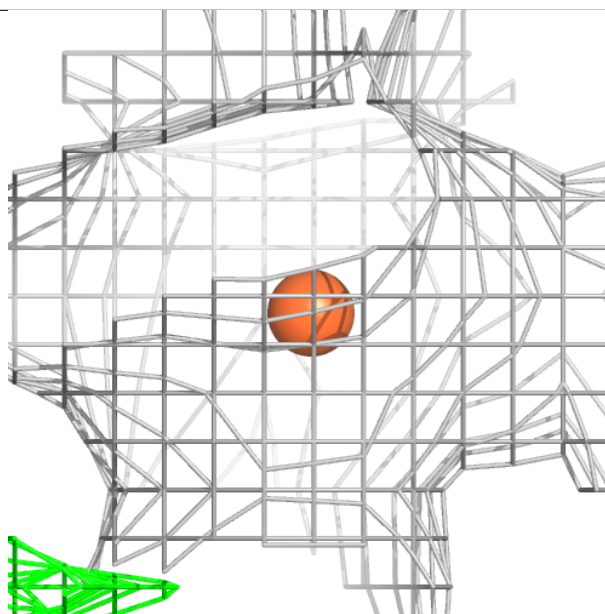
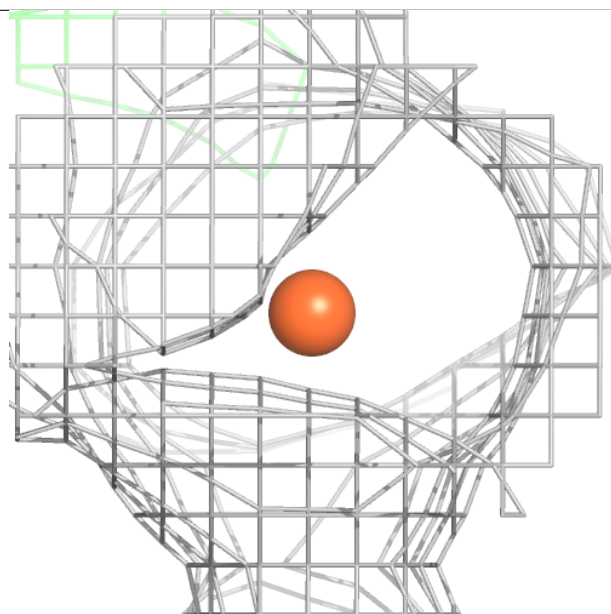
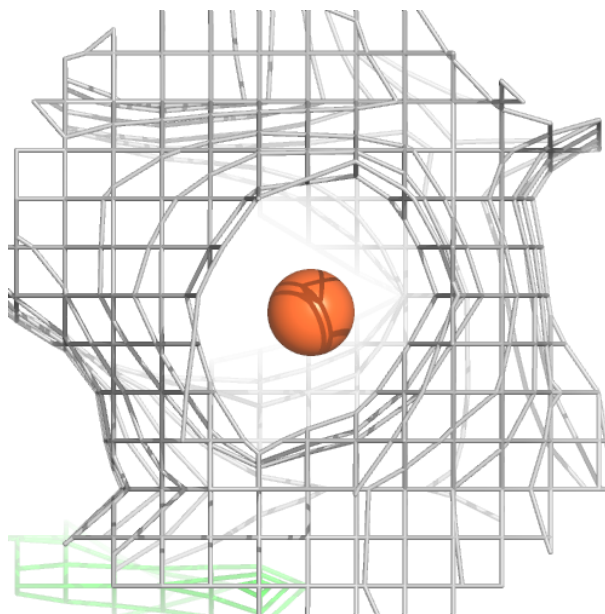
Electron density around FE A 401:

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and green (positive)



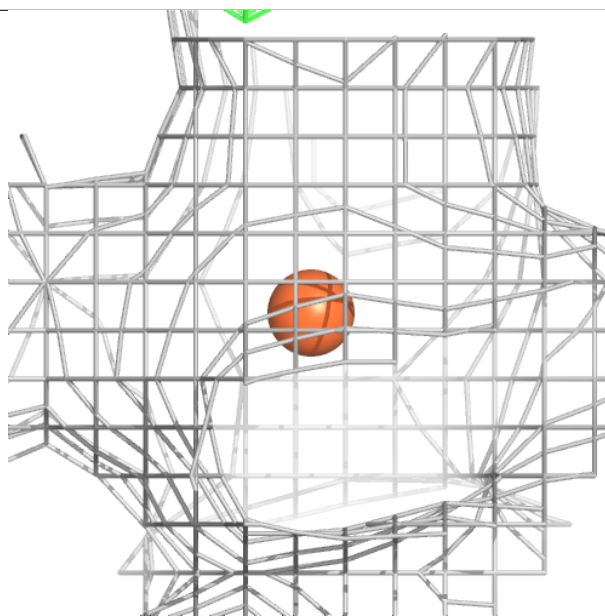
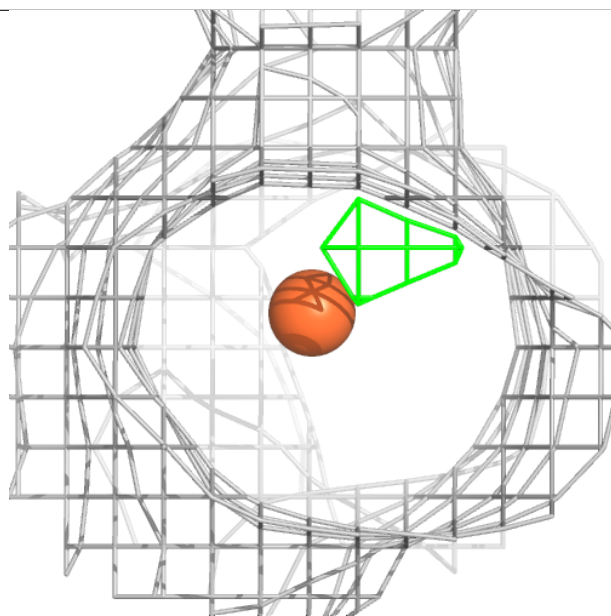
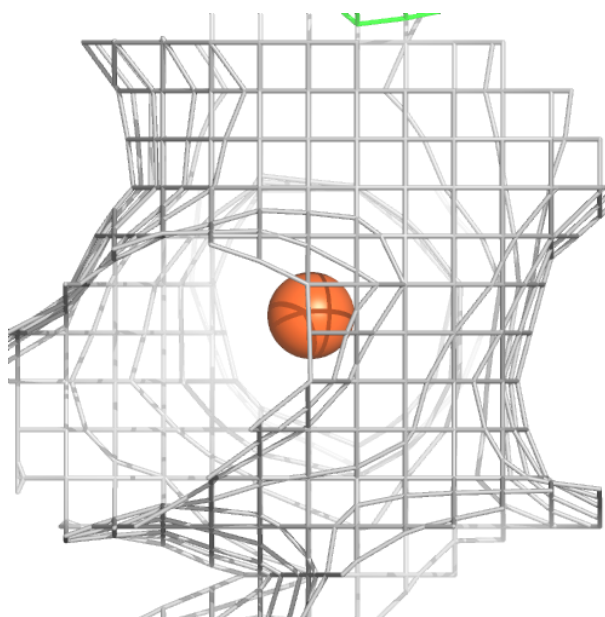
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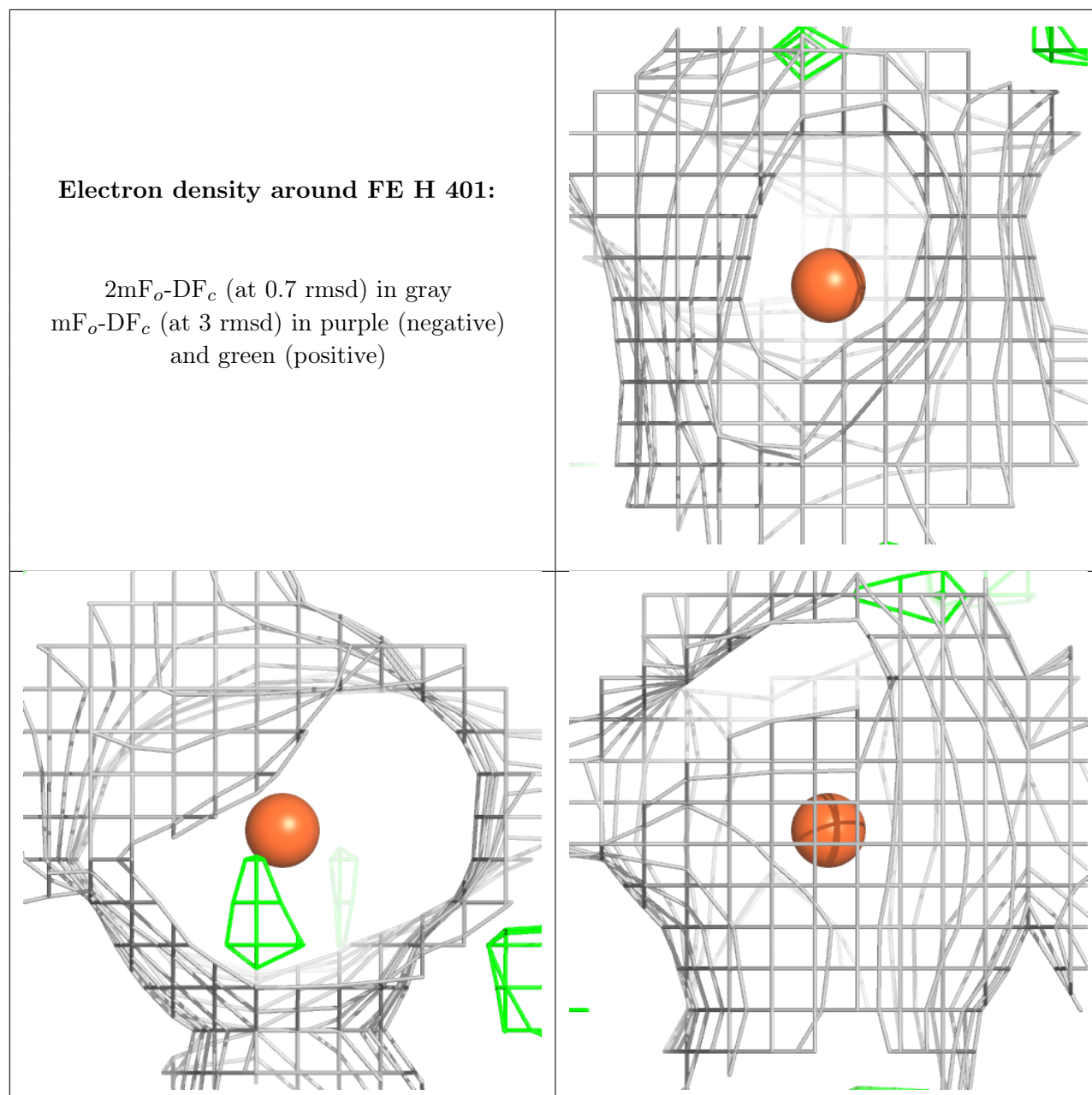
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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around FE 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)





6.5 Other polymers ⓘ

There are no such residues in this entry.