



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 12, 2026 – 03:22 AM UTC

PDB ID : 4TWF / pdb_00004twf
Title : X-ray structure of a pentameric ligand gated ion channel from *Erwinia chrysanthemi* (ELIC) in complex with bromomemantine
Authors : Ulens, C.; Spurny, R.; Thompson, A.J.; Alqazzaz, M.; Debaveye, S.; Lu, H.; Price, K.; Villalgordo, J.M.; Tresadern, G.; Lynch, J.W.; Lummis, S.C.R.
Deposited on : 2014-06-30
Resolution : 3.90 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 : 2022.3.0, CSD as543be (2022)
Xtrriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

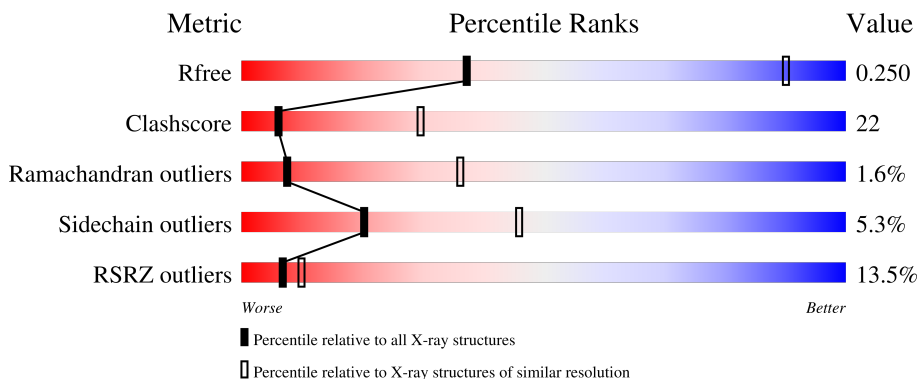
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.90 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1270 (4.10-3.70)
Clashscore	190562	1034 (4.08-3.72)
Ramachandran outliers	187476	1251 (4.10-3.70)
Sidechain outliers	187428	1243 (4.10-3.70)
RSRZ outliers	180081	1269 (4.10-3.70)

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	307	
1	B	307	
1	C	307	
1	D	307	

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Mol	Chain	Length	Quality of chain
1	E	307	14% 56% 37% 6% .
1	F	307	12% 57% 36% 5% .
1	G	307	19% 55% 39% 5% .
1	H	307	13% 53% 39% 6% .
1	I	307	13% 55% 40% . .
1	J	307	13% 60% 33% 6% .

2 Entry composition i

There are 2 unique types of molecules in this entry. The entry contains 25138 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 Cys-loop ligand-gated ion channel.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	307	2497	1624	416	451	6	0	0	0
1	B	307	2497	1624	416	451	6	0	0	0
1	C	307	2497	1624	416	451	6	0	0	0
1	D	307	2497	1624	416	451	6	0	0	0
1	E	307	2497	1624	416	451	6	0	0	0
1	F	307	2497	1624	416	451	6	0	0	0
1	G	307	2497	1624	416	451	6	0	0	0
1	H	307	2497	1624	416	451	6	0	0	0
1	I	307	2497	1624	416	451	6	0	0	0
1	J	307	2497	1624	416	451	6	0	0	0

There are 40 discrepancies between the modelled and reference sequences:

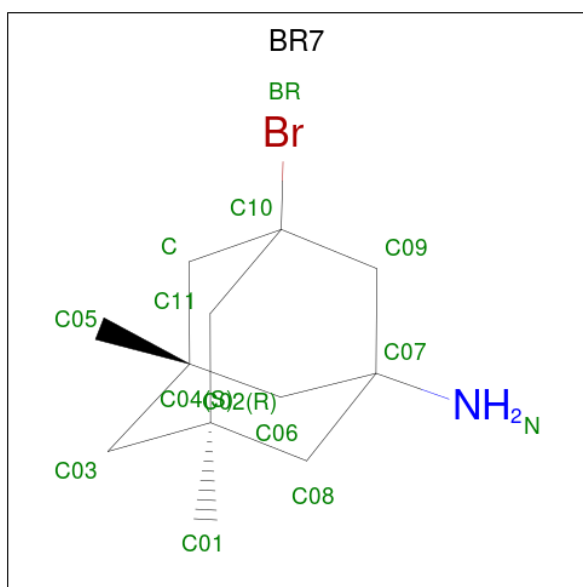
Chain	Residue	Modelled	Actual	Comment	Reference
A	152	ALA	ILE	conflict	UNP P0C7B7
A	164	GLY	-	insertion	UNP P0C7B7
A	247	SER	PHE	engineered mutation	UNP P0C7B7
A	289	ASN	MET	conflict	UNP P0C7B7
B	152	ALA	ILE	conflict	UNP P0C7B7
B	164	GLY	-	insertion	UNP P0C7B7
B	247	SER	PHE	engineered mutation	UNP P0C7B7
B	289	ASN	MET	conflict	UNP P0C7B7
C	152	ALA	ILE	conflict	UNP P0C7B7

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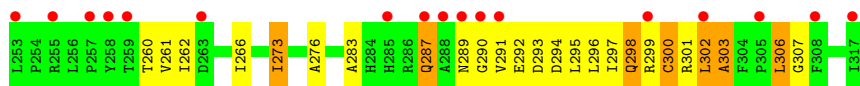
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Chain	Residue	Modelled	Actual	Comment	Reference
C	164	GLY	-	insertion	UNP P0C7B7
C	247	SER	PHE	engineered mutation	UNP P0C7B7
C	289	ASN	MET	conflict	UNP P0C7B7
D	152	ALA	ILE	conflict	UNP P0C7B7
D	164	GLY	-	insertion	UNP P0C7B7
D	247	SER	PHE	engineered mutation	UNP P0C7B7
D	289	ASN	MET	conflict	UNP P0C7B7
E	152	ALA	ILE	conflict	UNP P0C7B7
E	164	GLY	-	insertion	UNP P0C7B7
E	247	SER	PHE	engineered mutation	UNP P0C7B7
E	289	ASN	MET	conflict	UNP P0C7B7
F	152	ALA	ILE	conflict	UNP P0C7B7
F	164	GLY	-	insertion	UNP P0C7B7
F	247	SER	PHE	engineered mutation	UNP P0C7B7
F	289	ASN	MET	conflict	UNP P0C7B7
G	152	ALA	ILE	conflict	UNP P0C7B7
G	164	GLY	-	insertion	UNP P0C7B7
G	247	SER	PHE	engineered mutation	UNP P0C7B7
G	289	ASN	MET	conflict	UNP P0C7B7
H	152	ALA	ILE	conflict	UNP P0C7B7
H	164	GLY	-	insertion	UNP P0C7B7
H	247	SER	PHE	engineered mutation	UNP P0C7B7
H	289	ASN	MET	conflict	UNP P0C7B7
I	152	ALA	ILE	conflict	UNP P0C7B7
I	164	GLY	-	insertion	UNP P0C7B7
I	247	SER	PHE	engineered mutation	UNP P0C7B7
I	289	ASN	MET	conflict	UNP P0C7B7
J	152	ALA	ILE	conflict	UNP P0C7B7
J	164	GLY	-	insertion	UNP P0C7B7
J	247	SER	PHE	engineered mutation	UNP P0C7B7
J	289	ASN	MET	conflict	UNP P0C7B7

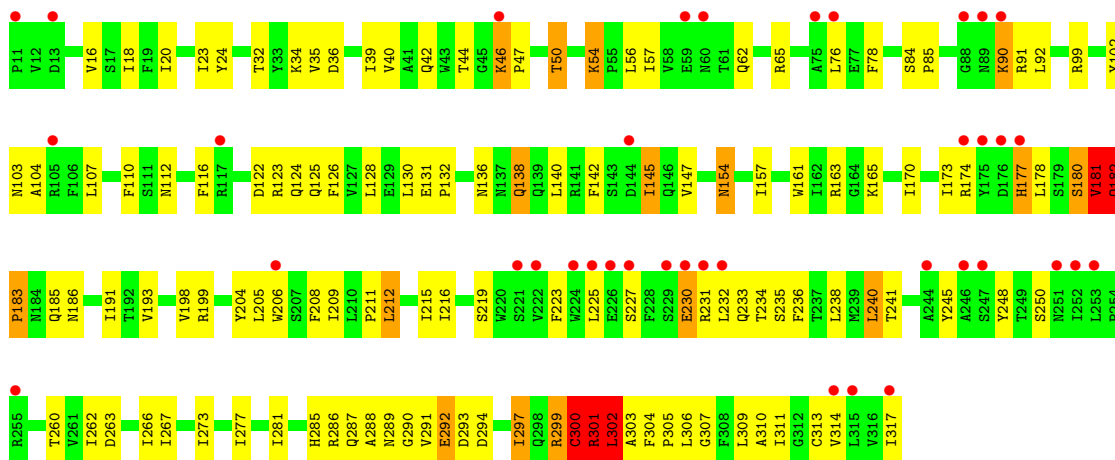
- Molecule 2 is Bromomemantine (CCD ID: BR7) (formula: C₁₂H₂₀BrN).



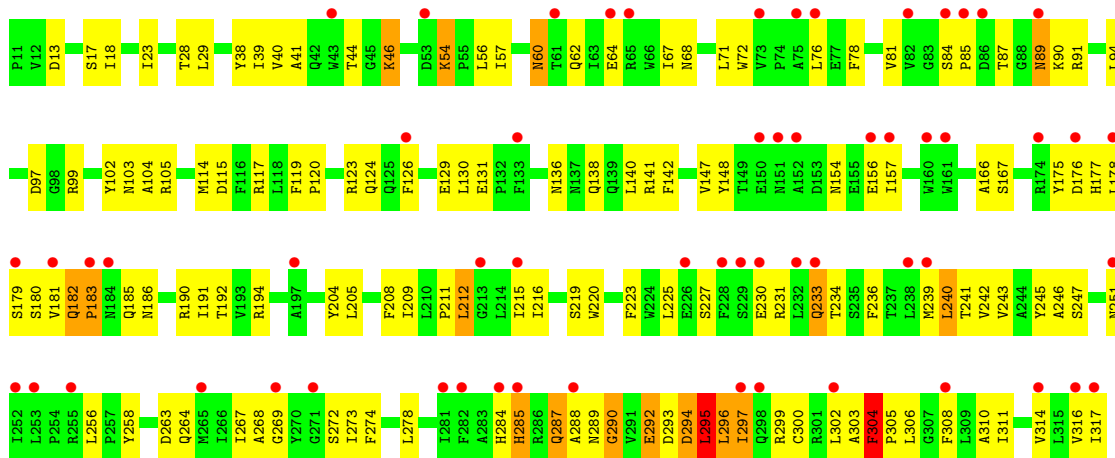
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	Br	C	N		
2	A	1	14	1	12	1	0	0
2	A	1	14	1	12	1	0	0
2	B	1	14	1	12	1	0	0
2	C	1	14	1	12	1	0	0
2	D	1	14	1	12	1	0	0
2	E	1	14	1	12	1	0	0
2	F	1	14	1	12	1	0	0
2	F	1	14	1	12	1	0	0
2	G	1	14	1	12	1	0	0
2	H	1	14	1	12	1	0	0
2	I	1	14	1	12	1	0	0
2	J	1	14	1	12	1	0	0



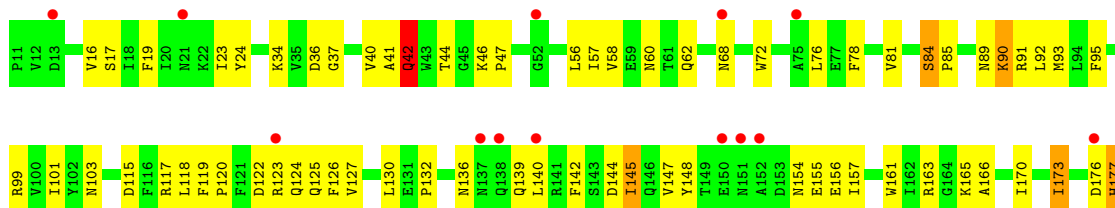
• Molecule 1: Cys-loop ligand-gated ion channel

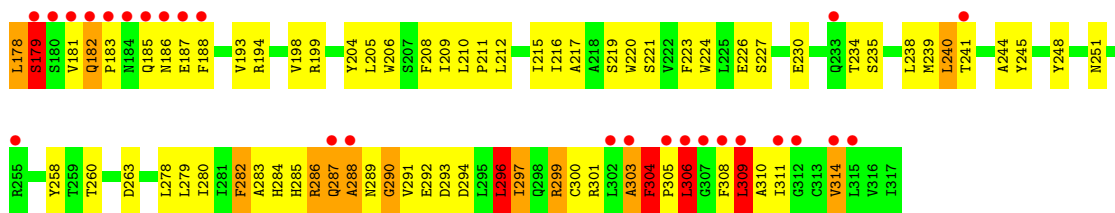


• Molecule 1: Cys-loop ligand-gated ion channel

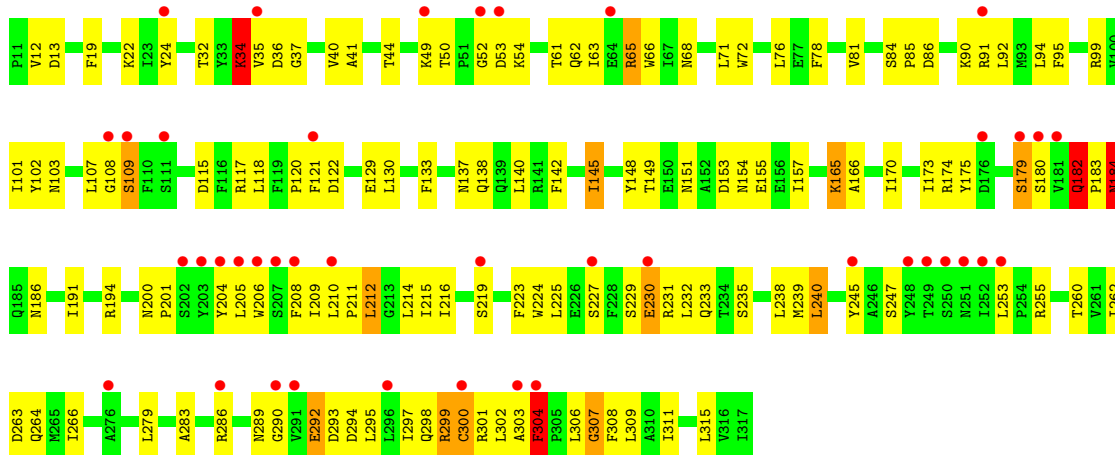


• Molecule 1: Cys-loop ligand-gated ion channel

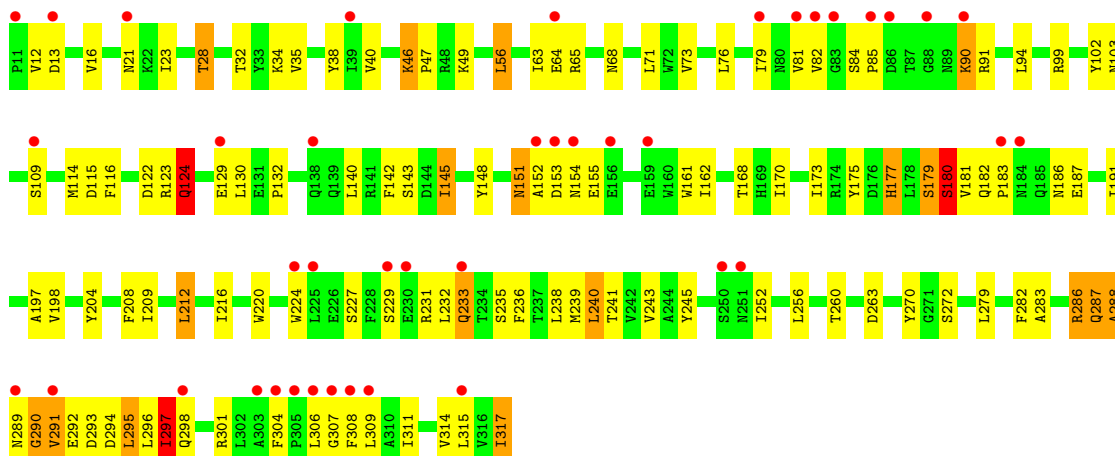




● Molecule 1: Cys-loop ligand-gated ion channel



● Molecule 1: Cys-loop ligand-gated ion channel



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	106.50Å 266.09Å 112.17Å 90.00° 107.70° 90.00°	Depositor
Resolution (Å)	49.83 – 3.90 49.83 – 3.90	Depositor EDS
% Data completeness (in resolution range)	99.2 (49.83-3.90) 91.0 (49.83-3.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.95 (at 3.88Å)	Xtrriage
Refinement program	PHENIX (phenix.refine: 1.8.4_1496)	Depositor
R, R_{free}	0.202 , 0.250 0.207 , 0.250	Depositor DCC
R_{free} test set	2719 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	106.1	Xtrriage
Anisotropy	0.404	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.20 , 74.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	25138	wwPDB-VP
Average B, all atoms (Å ²)	142.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.96% 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 i

5.1 Standard geometry i

Bond lengths and bond angles in the following residue types are not validated in this section: BR7

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	0.55	0/2564	1.07	16/3495 (0.5%)
1	B	0.54	0/2564	1.10	16/3495 (0.5%)
1	C	0.55	0/2564	1.09	16/3495 (0.5%)
1	D	0.50	0/2564	1.08	13/3495 (0.4%)
1	E	0.68	4/2564 (0.2%)	1.19	21/3495 (0.6%)
1	F	0.57	0/2564	1.14	12/3495 (0.3%)
1	G	0.52	0/2564	1.08	8/3495 (0.2%)
1	H	0.56	0/2564	1.17	21/3495 (0.6%)
1	I	0.52	0/2564	1.11	12/3495 (0.3%)
1	J	0.57	0/2564	1.10	13/3495 (0.4%)
All	All	0.56	4/25640 (0.0%)	1.11	148/34950 (0.4%)

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	B	0	7
1	C	0	3
1	D	0	2
1	E	0	3
1	F	0	5
1	G	0	2
1	H	0	3
1	I	0	2
1	J	0	2
All	All	0	31

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	273	ILE	CA-CB	-11.76	1.38	1.54
1	E	239	MET	SD-CE	6.97	1.97	1.79
1	E	240	LEU	N-CA	-6.42	1.38	1.46
1	E	239	MET	CA-C	-5.05	1.46	1.52

The worst 5 of 148 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	54	LYS	CA-C-N	-10.96	108.78	119.85
1	F	54	LYS	C-N-CA	-10.96	108.78	119.85
1	E	45	GLY	N-CA-C	-10.77	97.20	112.60
1	G	182	GLN	CA-C-N	10.16	132.54	119.84
1	G	182	GLN	C-N-CA	10.16	132.54	119.84

There are no chirality outliers.

5 of 31 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	177	HIS	Peptide
1	A	178	LEU	Peptide
1	B	177	HIS	Peptide
1	B	178	LEU	Peptide
1	B	179	SER	Peptide

5.2 Too-close contacts [i](#)

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	2497	0	2468	120	0
1	B	2497	0	2468	119	0
1	C	2497	0	2468	123	0
1	D	2497	0	2468	83	0
1	E	2497	0	2468	123	0
1	F	2497	0	2468	128	0
1	G	2497	0	2468	126	0
1	H	2497	0	2468	135	0
1	I	2497	0	2468	140	0
1	J	2497	0	2468	119	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	28	0	40	2	0
2	B	14	0	20	4	0
2	C	14	0	20	5	0
2	D	14	0	20	0	0
2	E	14	0	20	2	0
2	F	28	0	40	1	0
2	G	14	0	20	3	0
2	H	14	0	20	0	0
2	I	14	0	20	1	0
2	J	14	0	20	2	0
All	All	25138	0	24920	1126	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 22.

The worst 5 of 1126 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:221:SER:HA	1:A:301:ARG:HH22	1.15	1.07
1:I:63:ILE:HD13	1:I:90:LYS:HE3	1.40	1.04
1:I:299:ARG:HH12	1:I:302:LEU:HD23	1.21	1.01
1:E:239:MET:HE1	1:E:273:ILE:HD13	1.46	0.98
1:C:298:GLN:O	1:C:301:ARG:NH1	1.98	0.97

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	305/307 (99%)	278 (91%)	23 (8%)	4 (1%)	9	40
1	B	305/307 (99%)	280 (92%)	19 (6%)	6 (2%)	6	32

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	305/307 (99%)	281 (92%)	20 (7%)	4 (1%)	9	40
1	D	305/307 (99%)	273 (90%)	23 (8%)	9 (3%)	3	26
1	E	305/307 (99%)	277 (91%)	26 (8%)	2 (1%)	18	53
1	F	305/307 (99%)	280 (92%)	20 (7%)	5 (2%)	7	36
1	G	305/307 (99%)	280 (92%)	18 (6%)	7 (2%)	5	30
1	H	305/307 (99%)	280 (92%)	19 (6%)	6 (2%)	6	32
1	I	305/307 (99%)	276 (90%)	25 (8%)	4 (1%)	9	40
1	J	305/307 (99%)	279 (92%)	24 (8%)	2 (1%)	18	53
All	All	3050/3070 (99%)	2784 (91%)	217 (7%)	49 (2%)	7	36

5 of 49 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	150	GLU
1	B	154	ASN
1	B	179	SER
1	B	185	GLN
1	C	152	ALA

5.3.2 Protein sidechains [i](#)

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	274/274 (100%)	254 (93%)	20 (7%)	13	39
1	B	274/274 (100%)	266 (97%)	8 (3%)	37	58
1	C	274/274 (100%)	262 (96%)	12 (4%)	25	49
1	D	274/274 (100%)	259 (94%)	15 (6%)	19	45
1	E	274/274 (100%)	261 (95%)	13 (5%)	23	48
1	F	274/274 (100%)	259 (94%)	15 (6%)	19	45
1	G	274/274 (100%)	263 (96%)	11 (4%)	28	51
1	H	274/274 (100%)	261 (95%)	13 (5%)	23	48

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	I	274/274 (100%)	257 (94%)	17 (6%)	16	43
1	J	274/274 (100%)	253 (92%)	21 (8%)	12	37
All	All	2740/2740 (100%)	2595 (95%)	145 (5%)	20	46

5 of 145 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	I	184	ASN
1	J	309	LEU
1	I	239	MET
1	J	145	ILE
1	D	167	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 47 such sidechains are listed below:

Mol	Chain	Res	Type
1	G	138	GLN
1	H	139	GLN
1	G	139	GLN
1	G	285	HIS
1	H	289	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](#)

12 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
2	BR7	H	401	-	13,16,16	3.70	9 (69%)	23,30,30	3.70	14 (60%)
2	BR7	A	402	-	13,16,16	3.99	10 (76%)	23,30,30	3.74	14 (60%)
2	BR7	J	401	-	13,16,16	3.64	8 (61%)	23,30,30	3.51	13 (56%)
2	BR7	F	401	-	13,16,16	3.66	8 (61%)	23,30,30	3.73	15 (65%)
2	BR7	A	401	-	13,16,16	3.63	9 (69%)	23,30,30	3.53	14 (60%)
2	BR7	E	401	-	13,16,16	3.72	11 (84%)	23,30,30	4.26	12 (52%)
2	BR7	G	401	-	13,16,16	3.81	9 (69%)	23,30,30	3.58	15 (65%)
2	BR7	I	401	-	13,16,16	3.81	9 (69%)	23,30,30	3.40	10 (43%)
2	BR7	D	401	-	13,16,16	3.82	9 (69%)	23,30,30	4.11	13 (56%)
2	BR7	C	401	-	13,16,16	3.90	10 (76%)	23,30,30	4.07	14 (60%)
2	BR7	F	402	-	13,16,16	3.88	10 (76%)	23,30,30	3.73	14 (60%)
2	BR7	B	401	-	13,16,16	3.86	11 (84%)	23,30,30	4.38	12 (52%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	BR7	H	401	-	-	-	0/4/3/3
2	BR7	A	402	-	-	-	0/4/3/3
2	BR7	J	401	-	-	-	0/4/3/3
2	BR7	F	401	-	-	-	0/4/3/3
2	BR7	A	401	-	-	-	0/4/3/3
2	BR7	E	401	-	-	-	0/4/3/3
2	BR7	G	401	-	-	-	0/4/3/3
2	BR7	I	401	-	-	-	0/4/3/3
2	BR7	D	401	-	-	-	0/4/3/3
2	BR7	C	401	-	-	-	0/4/3/3
2	BR7	F	402	-	-	-	0/4/3/3
2	BR7	B	401	-	-	-	0/4/3/3

The worst 5 of 113 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	402	BR7	C09-C10	-10.49	1.37	1.53
2	C	401	BR7	C09-C10	-10.37	1.38	1.53
2	F	402	BR7	C09-C10	-10.34	1.38	1.53
2	I	401	BR7	C09-C10	-10.31	1.38	1.53
2	D	401	BR7	C09-C10	-10.30	1.38	1.53

The worst 5 of 160 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	401	BR7	C-C10-C11	-14.02	102.70	110.55
2	E	401	BR7	C-C10-C11	-13.47	103.01	110.55
2	B	401	BR7	C-C10-C11	-13.33	103.09	110.55
2	C	401	BR7	C-C10-C11	-12.71	103.44	110.55
2	H	401	BR7	C-C10-C11	-11.91	103.88	110.55

There are no chirality outliers.

There are no torsion outliers.

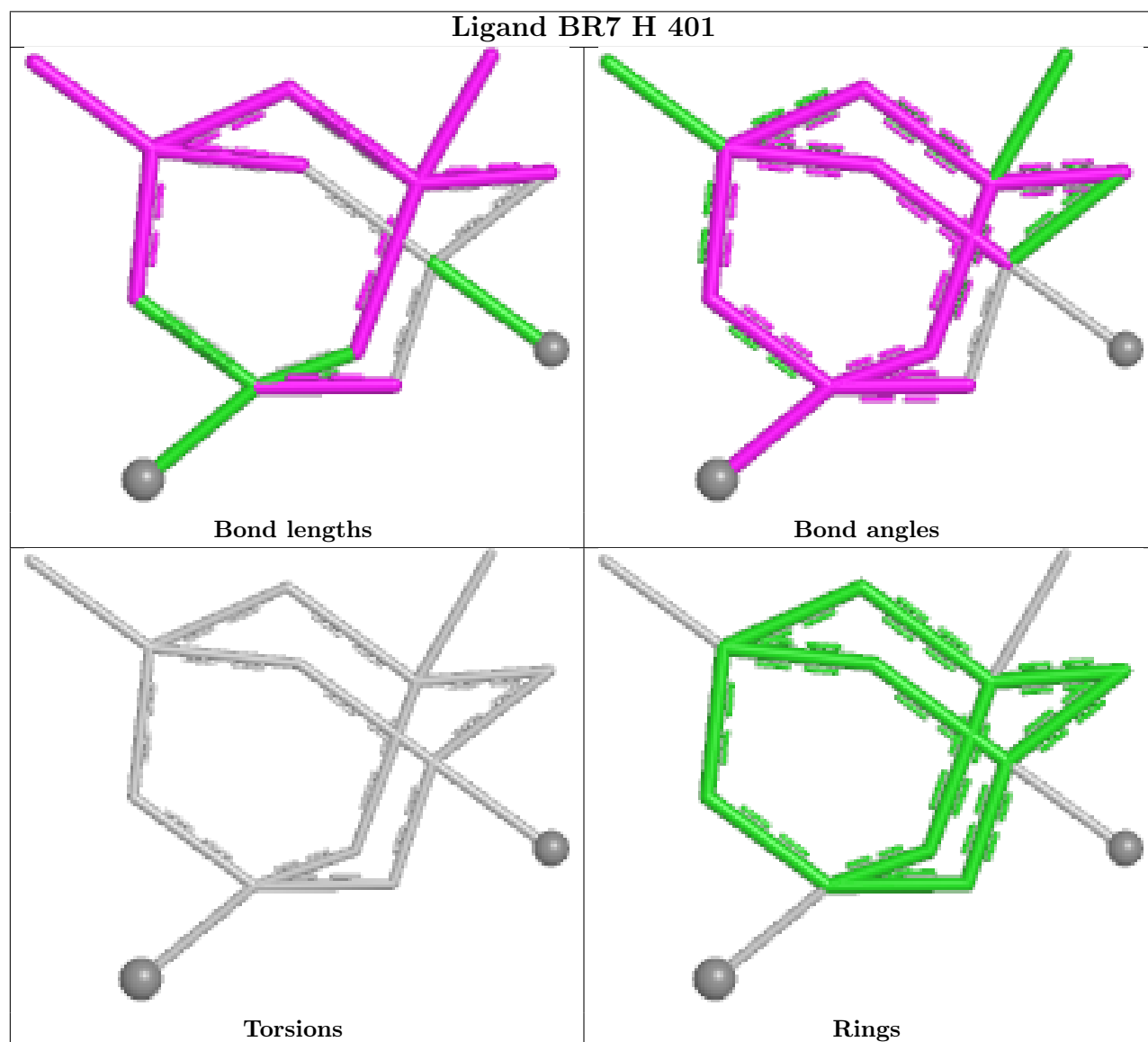
There are no ring outliers.

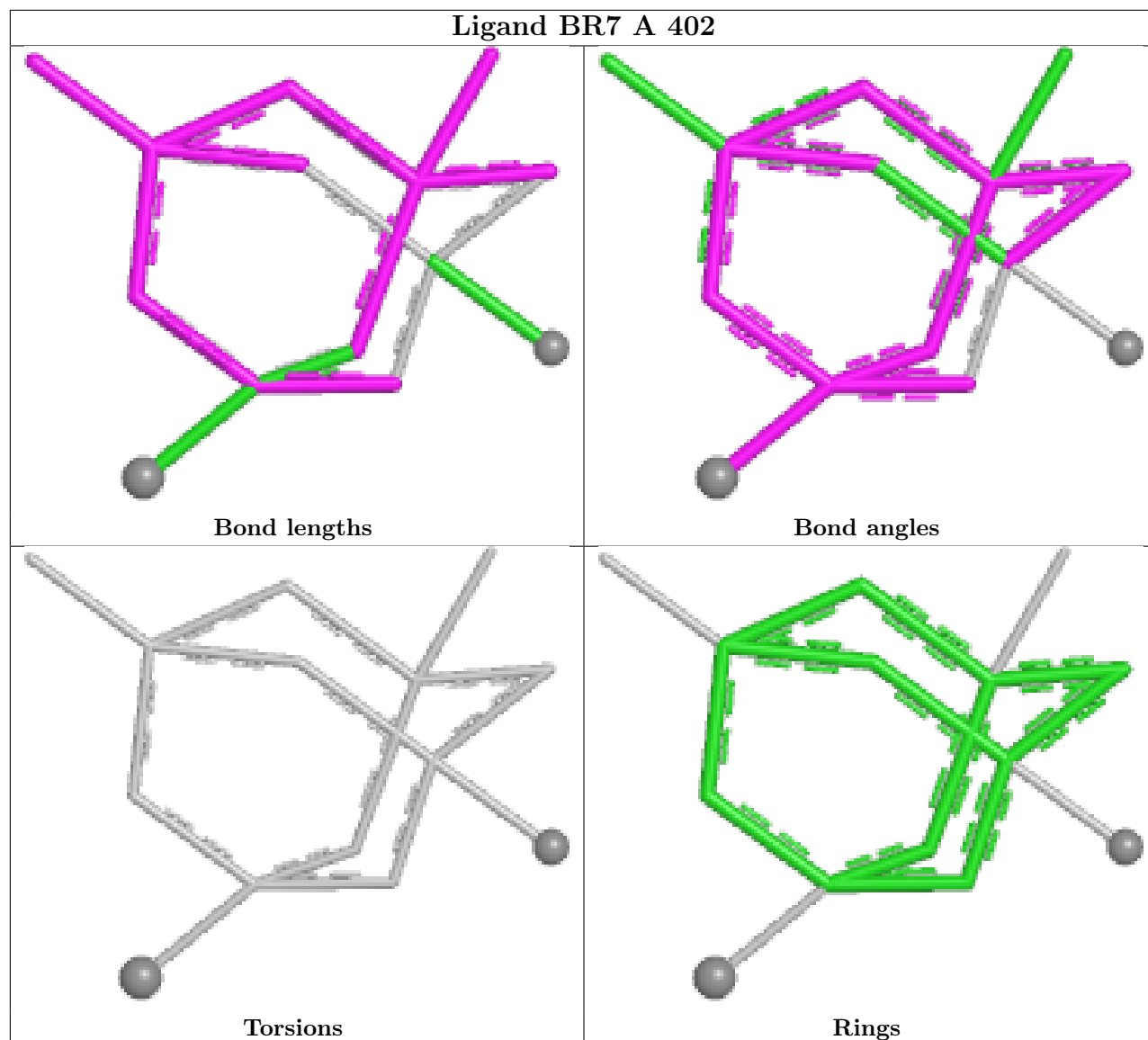
8 monomers are involved in 20 short contacts:

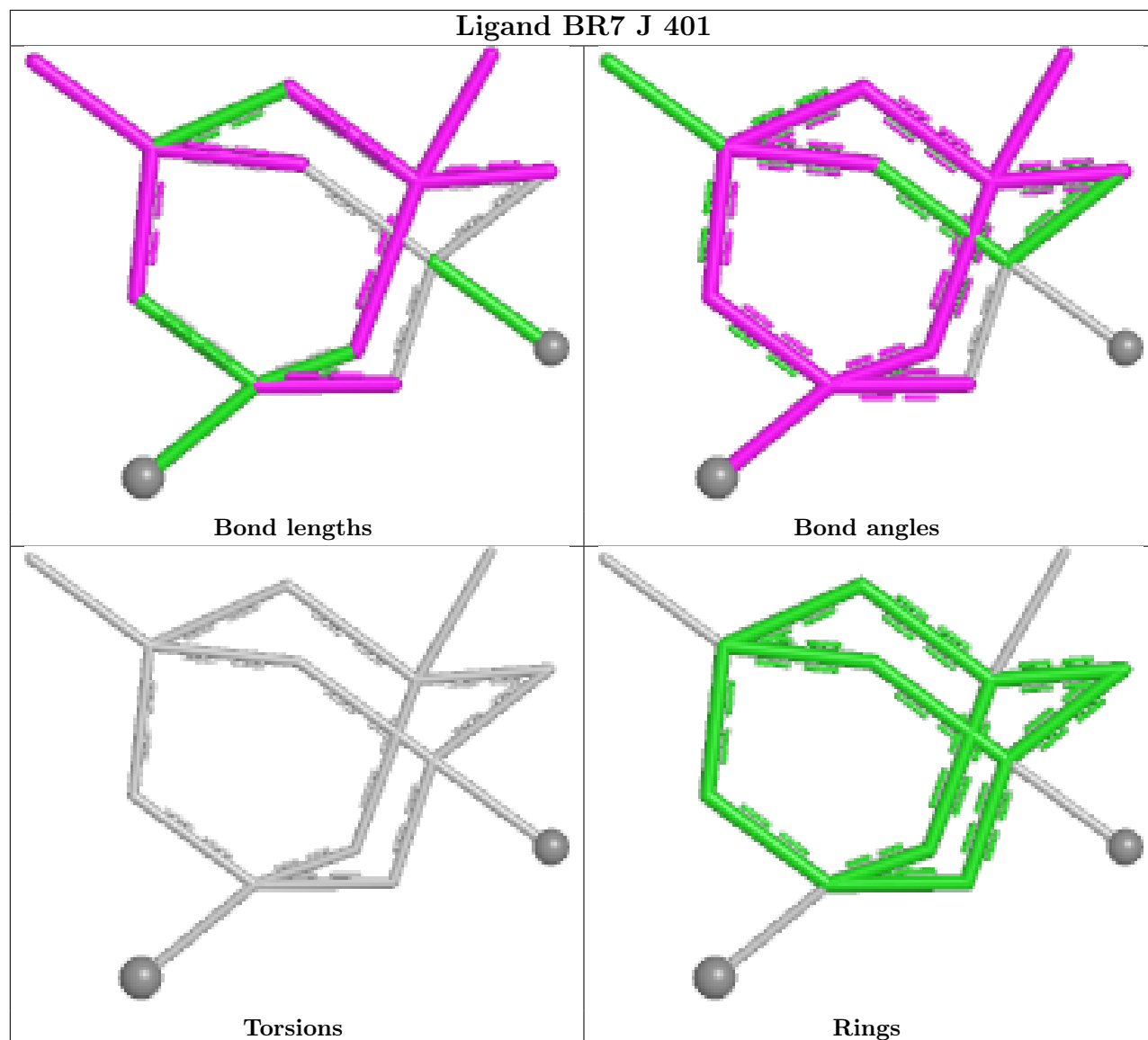
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	402	BR7	2	0
2	J	401	BR7	2	0
2	E	401	BR7	2	0
2	G	401	BR7	3	0
2	I	401	BR7	1	0
2	C	401	BR7	5	0
2	F	402	BR7	1	0
2	B	401	BR7	4	0

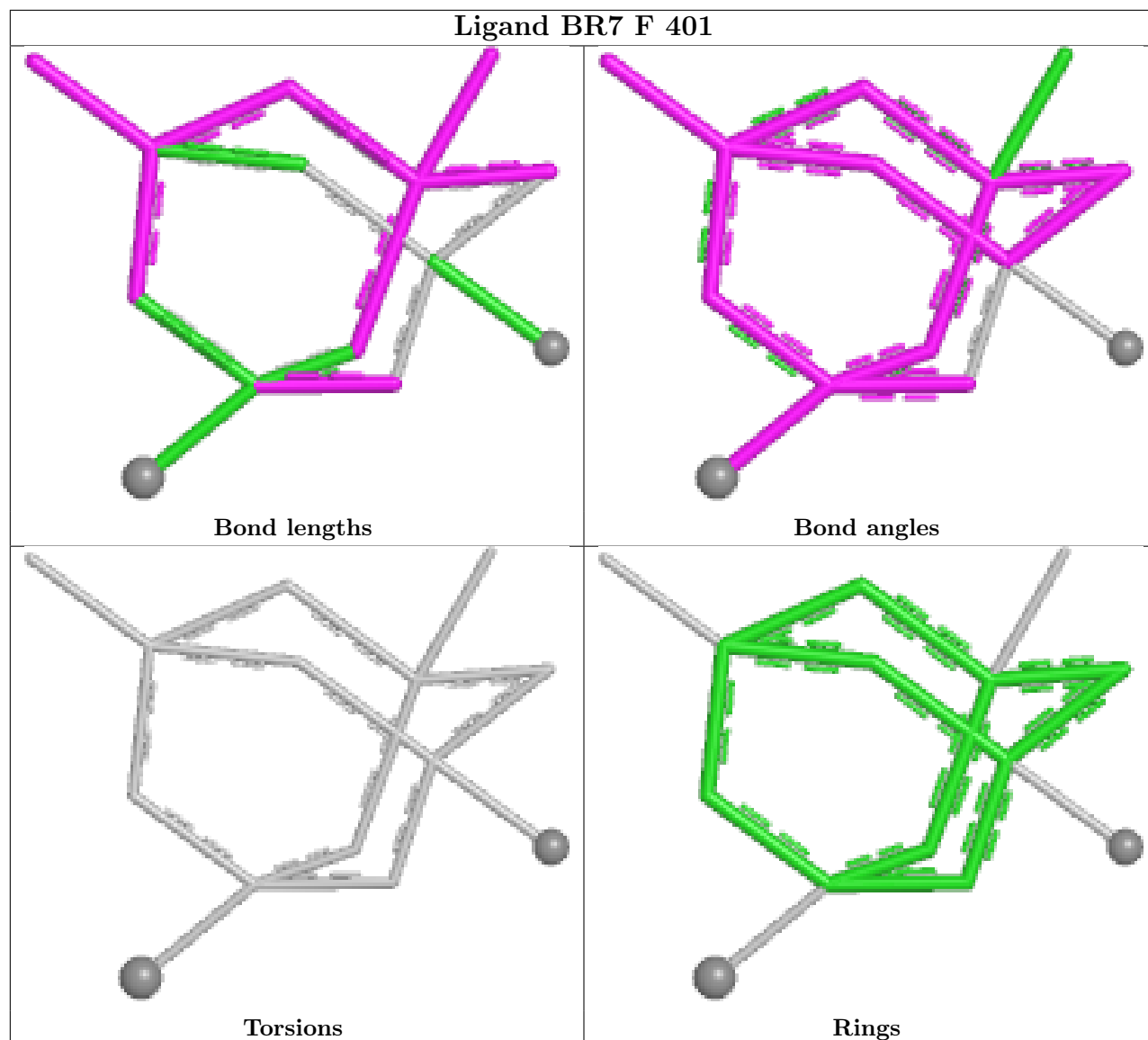
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

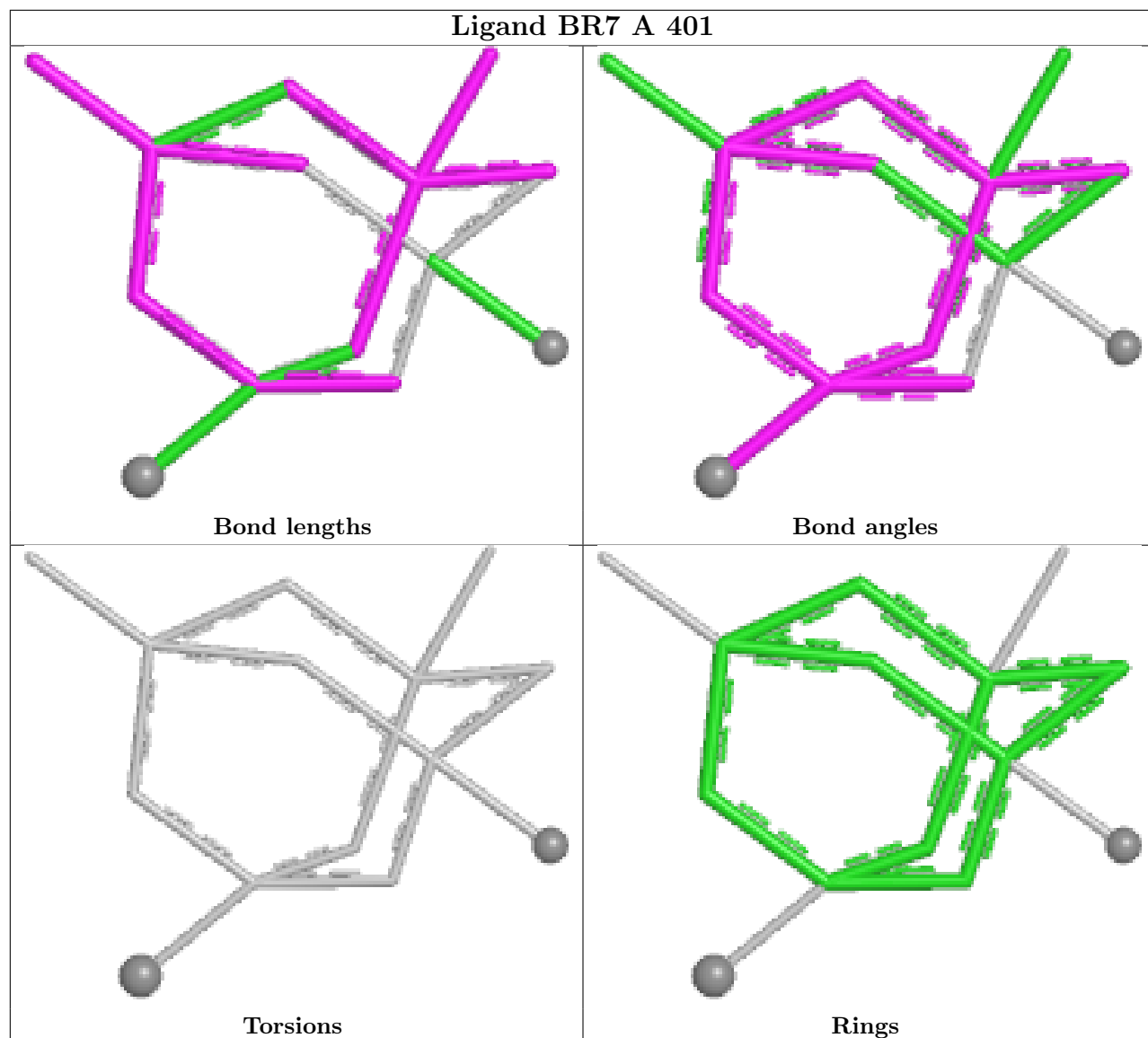
The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

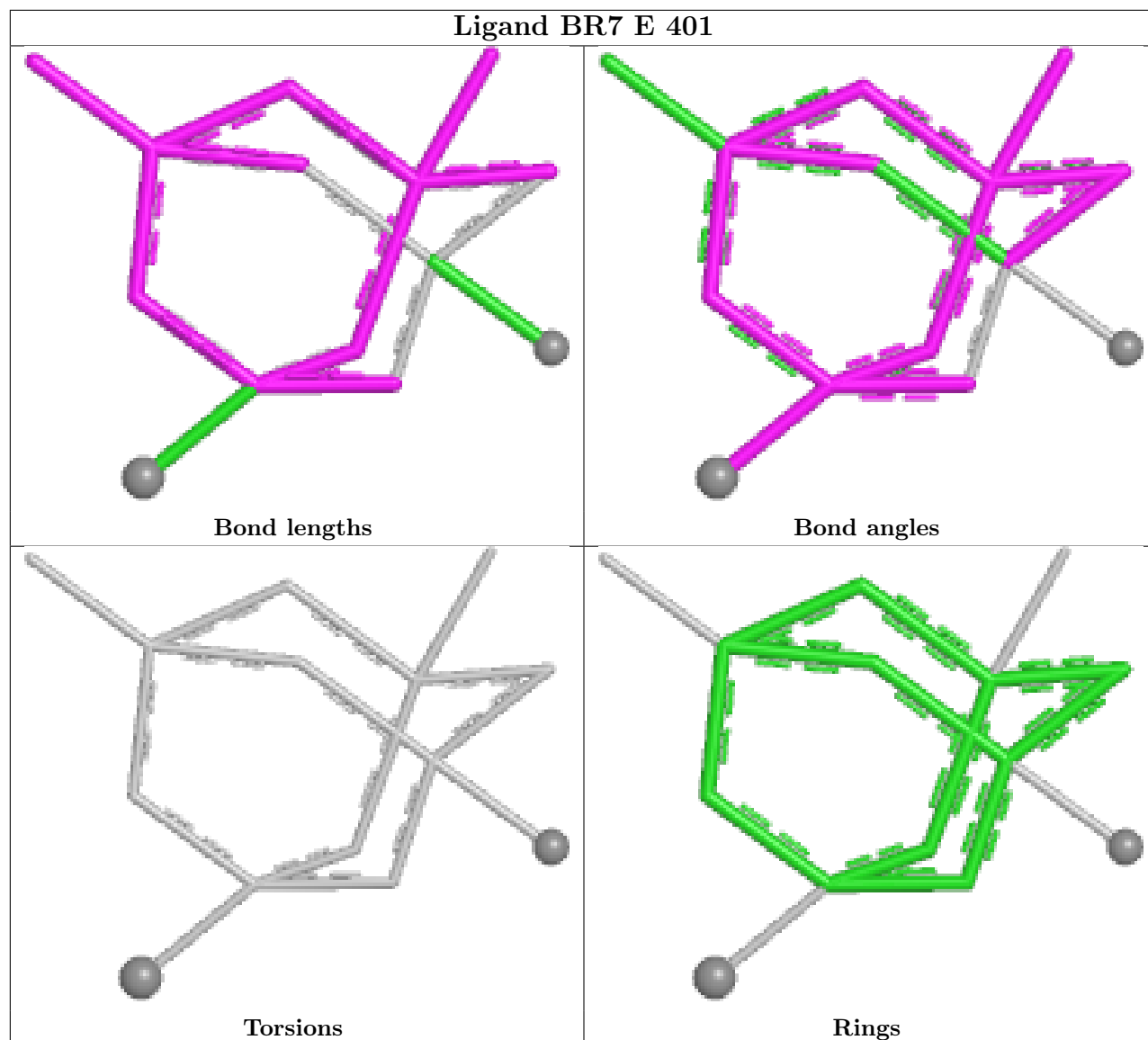


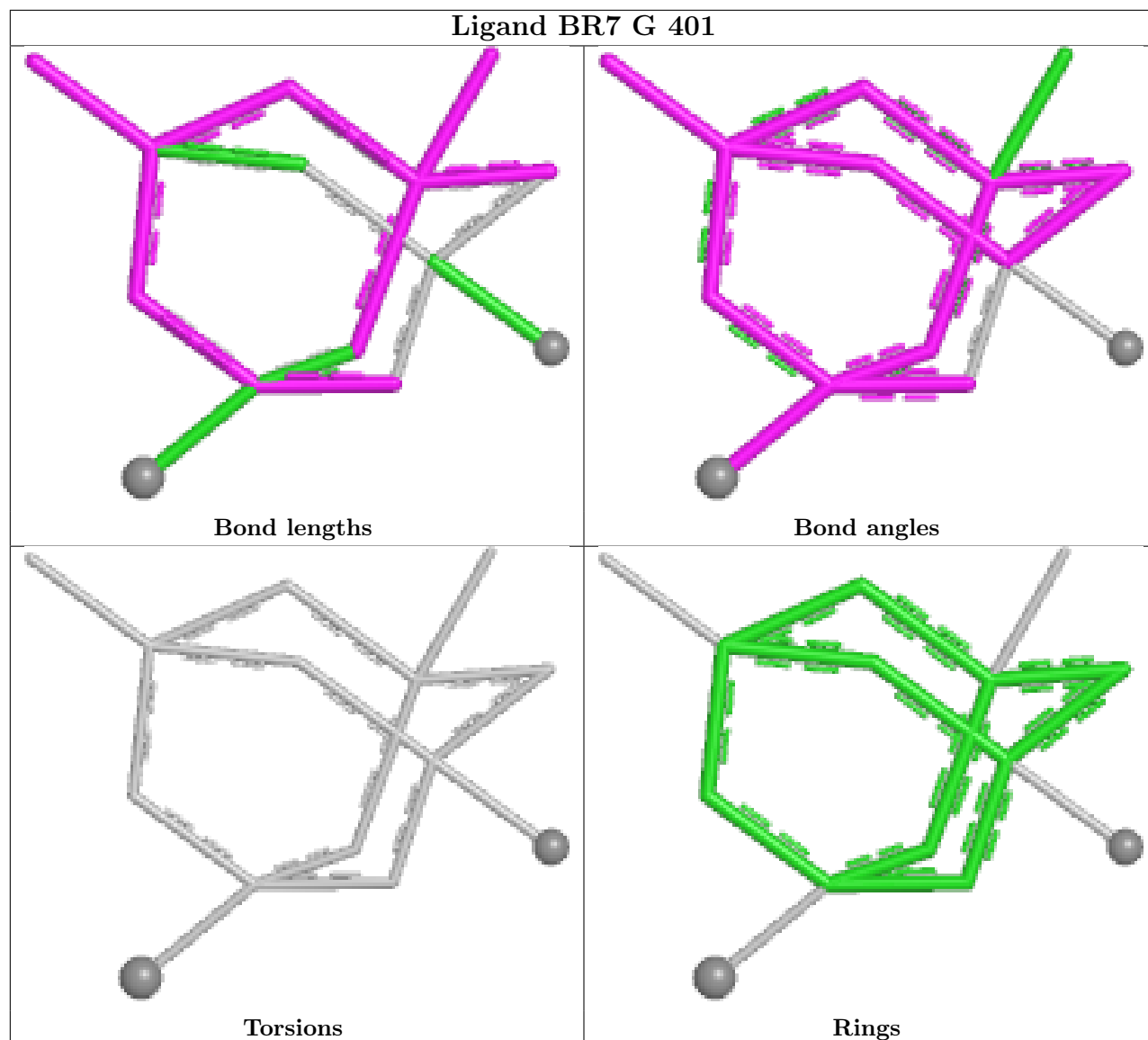


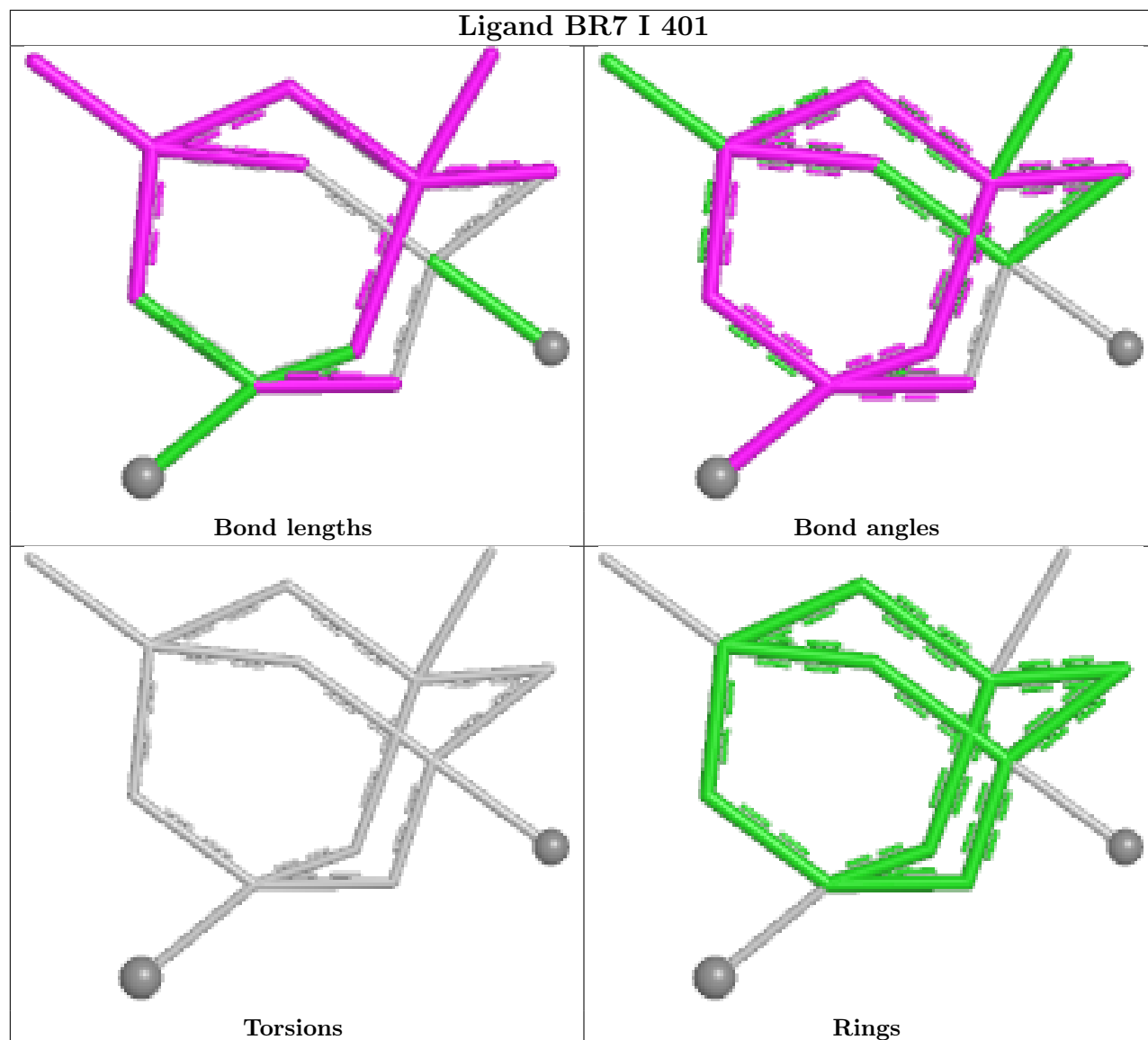


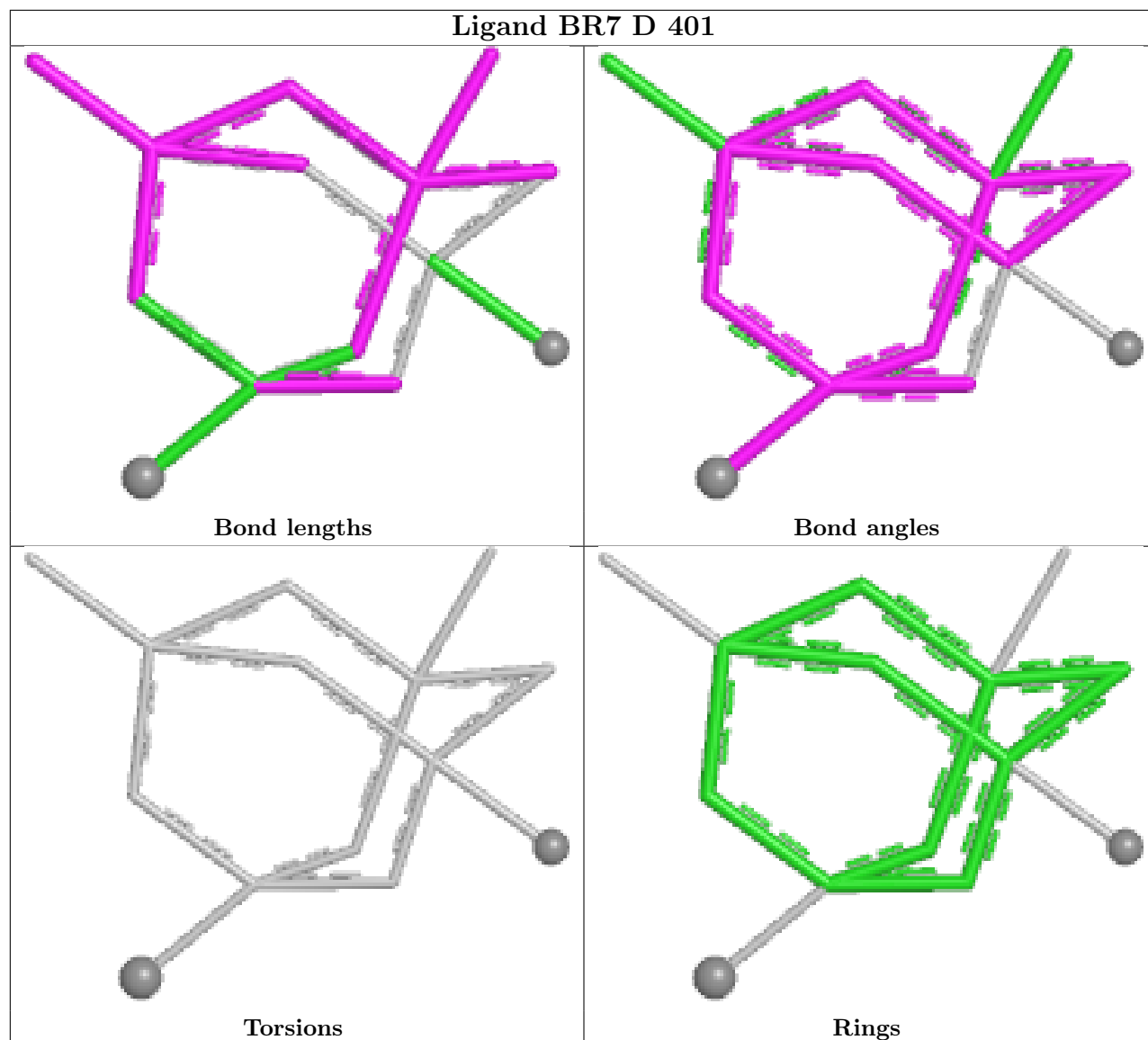


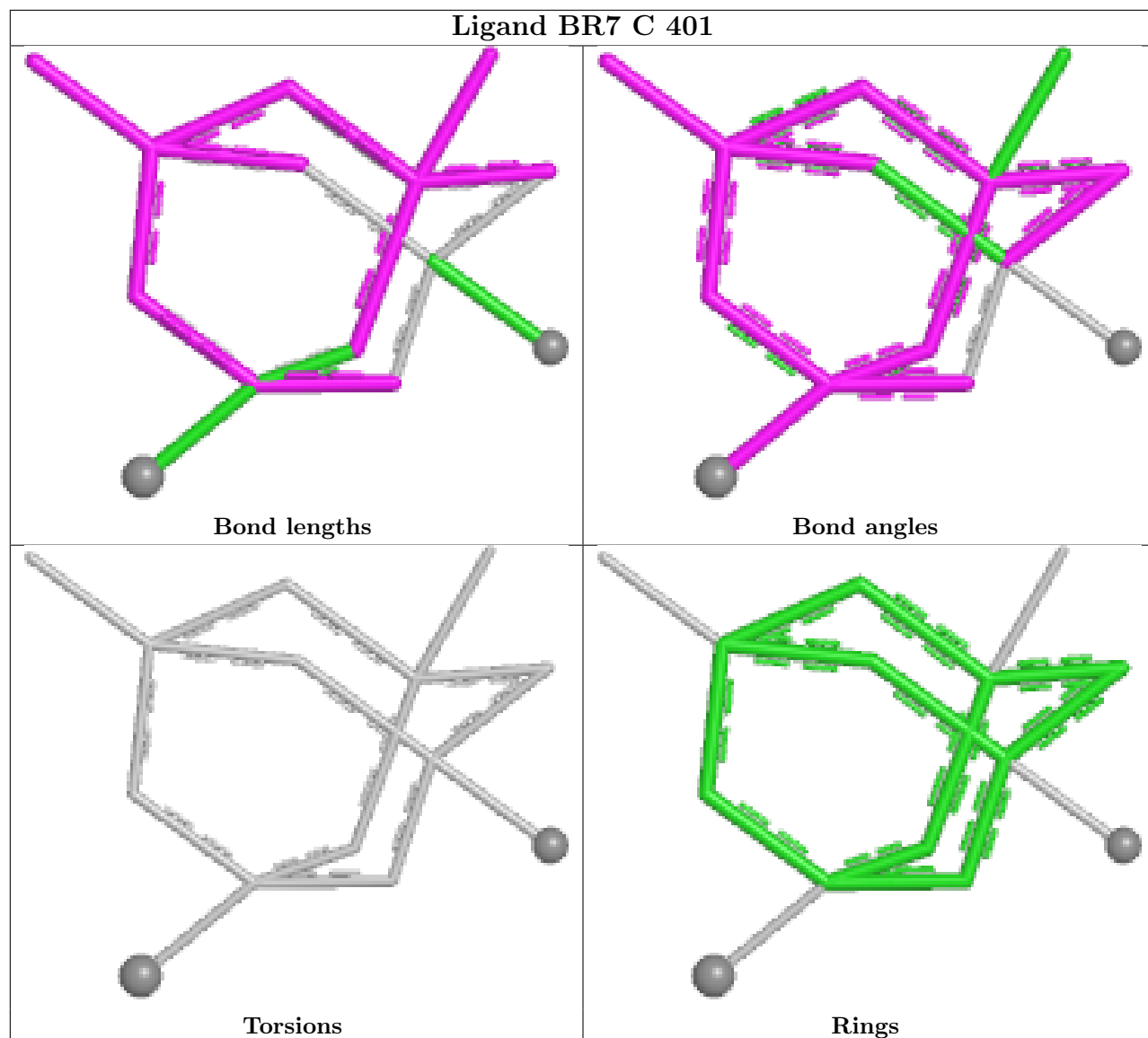


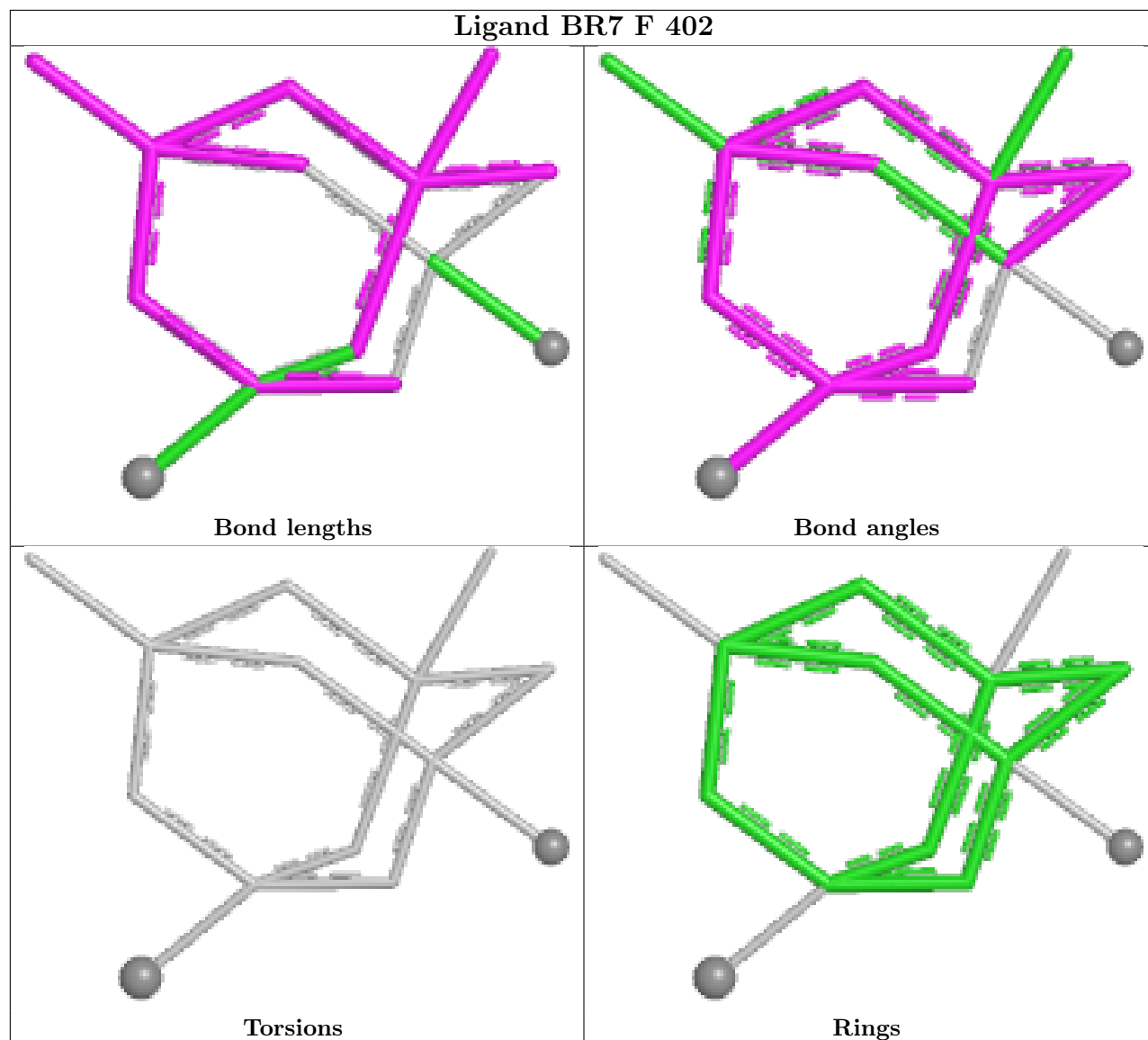


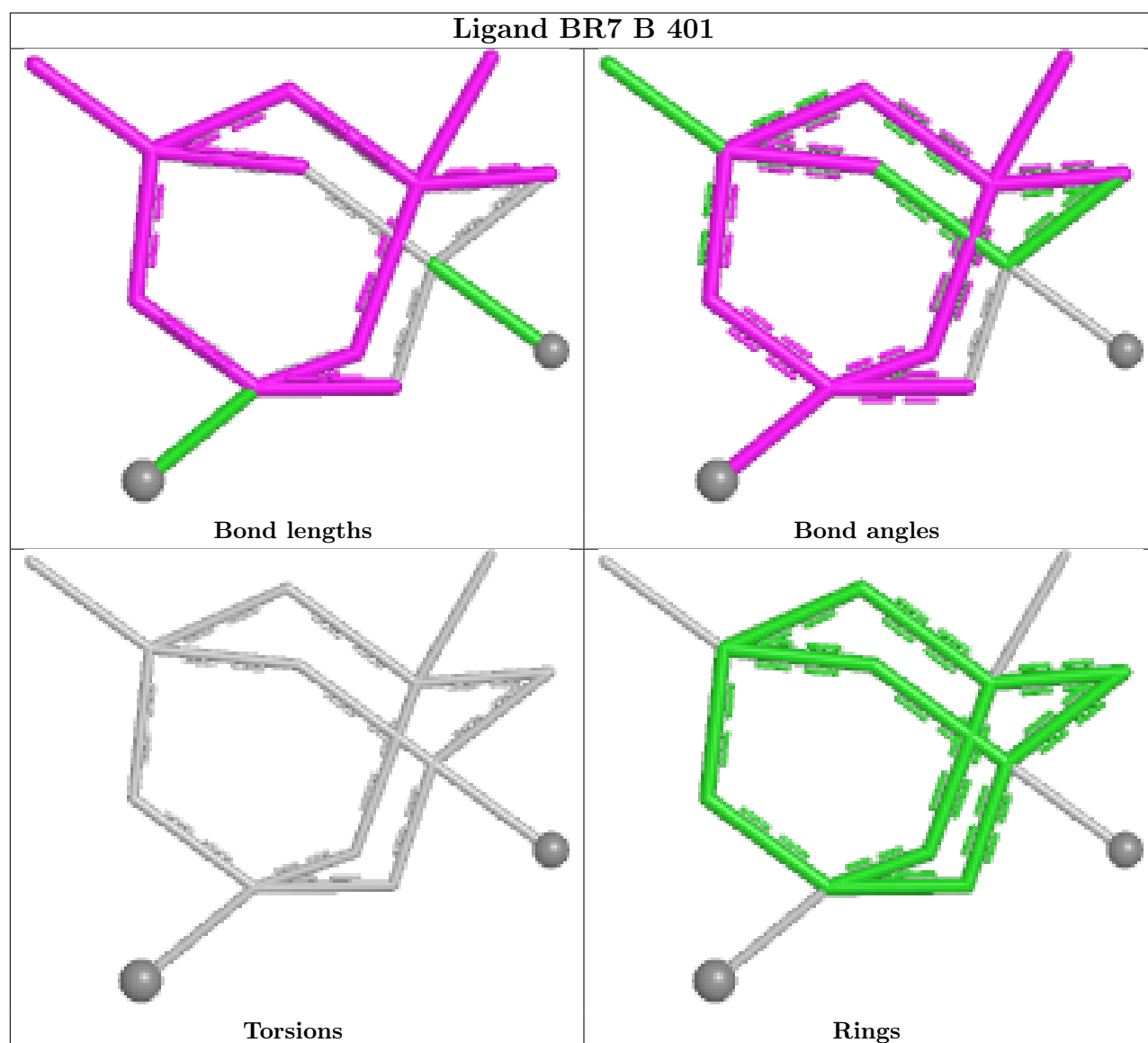












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

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	307/307 (100%)	0.53	23 (7%) 20 19	89, 136, 232, 324	0
1	B	307/307 (100%)	1.05	59 (19%) 3 5	78, 123, 237, 352	0
1	C	307/307 (100%)	0.67	31 (10%) 12 15	71, 119, 234, 291	0
1	D	307/307 (100%)	0.70	40 (13%) 7 10	73, 121, 228, 325	0
1	E	307/307 (100%)	0.78	43 (14%) 6 9	83, 129, 235, 381	0
1	F	307/307 (100%)	0.74	38 (12%) 8 11	84, 138, 250, 372	0
1	G	307/307 (100%)	0.98	59 (19%) 3 5	78, 122, 221, 324	0
1	H	307/307 (100%)	0.80	39 (12%) 8 11	79, 133, 250, 378	0
1	I	307/307 (100%)	0.85	41 (13%) 7 10	76, 131, 234, 327	0
1	J	307/307 (100%)	0.71	41 (13%) 7 10	91, 139, 272, 347	0
All	All	3070/3070 (100%)	0.78	414 (13%) 7 10	71, 129, 243, 381	0

The worst 5 of 414 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	290	GLY	11.2
1	C	312	GLY	10.6
1	E	291	VAL	10.4
1	C	313	CYS	10.3
1	E	289	ASN	8.7

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](#)

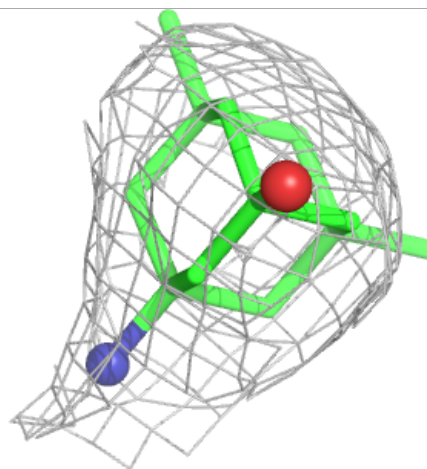
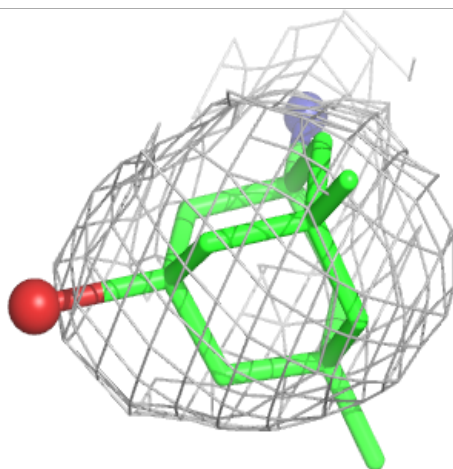
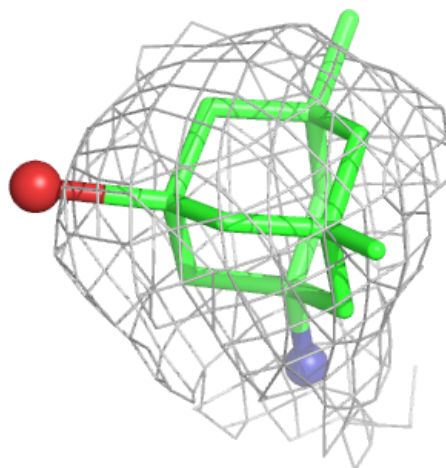
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	BR7	F	401	14/14	0.88	0.14	99,129,135,272	0
2	BR7	J	401	14/14	0.90	0.11	98,112,125,194	0
2	BR7	C	401	14/14	0.91	0.12	111,127,145,217	0
2	BR7	B	401	14/14	0.92	0.12	104,108,119,265	0
2	BR7	A	401	14/14	0.92	0.13	112,142,147,354	0
2	BR7	G	401	14/14	0.94	0.10	77,115,131,268	0
2	BR7	F	402	14/14	0.94	0.17	110,131,138,389	0
2	BR7	D	401	14/14	0.95	0.10	98,111,119,277	0
2	BR7	H	401	14/14	0.95	0.09	129,141,153,250	0
2	BR7	I	401	14/14	0.95	0.10	86,100,119,321	0
2	BR7	E	401	14/14	0.95	0.11	84,110,120,309	0
2	BR7	A	402	14/14	0.96	0.13	98,109,129,224	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.

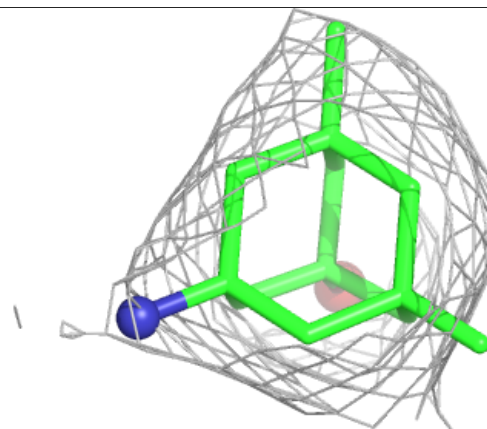
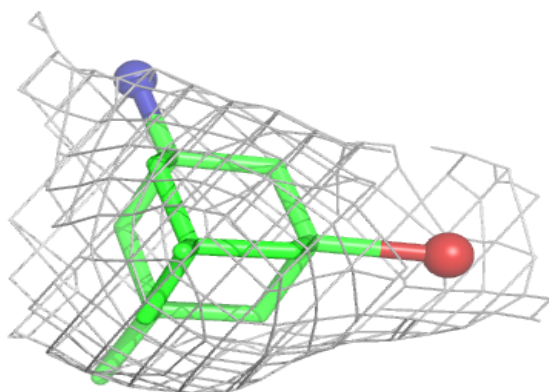
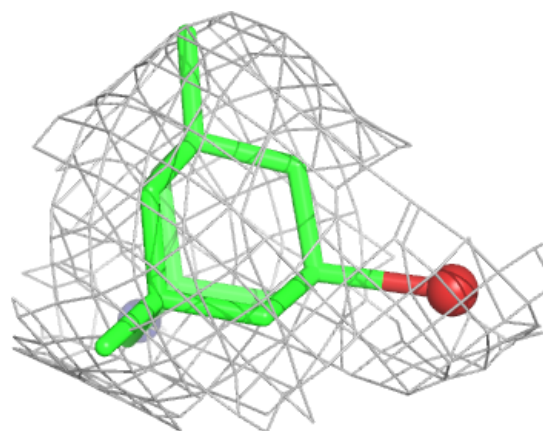
Electron density around BR7 F 401:

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



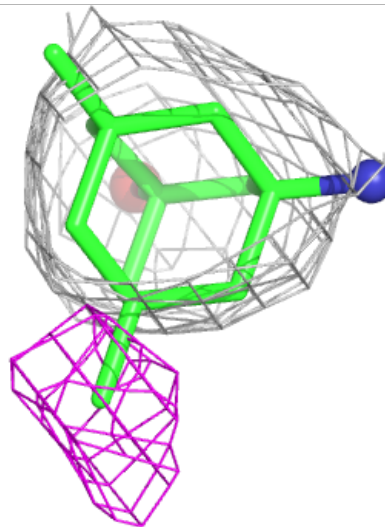
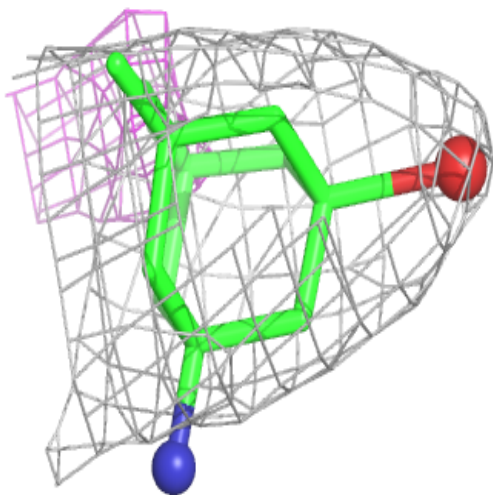
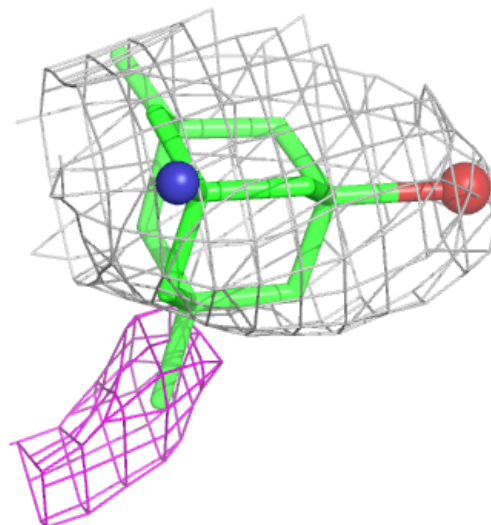
Electron density around BR7 J 401:

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



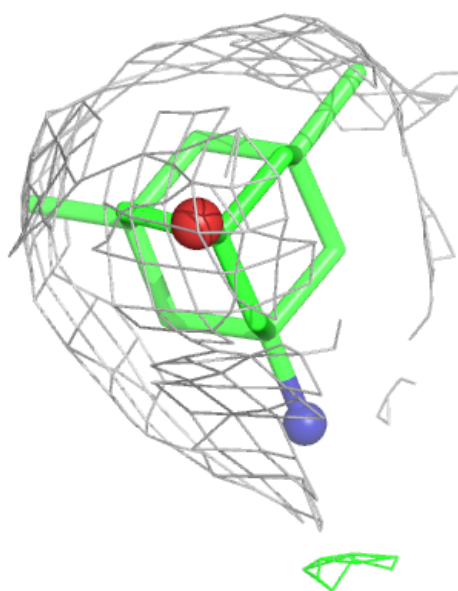
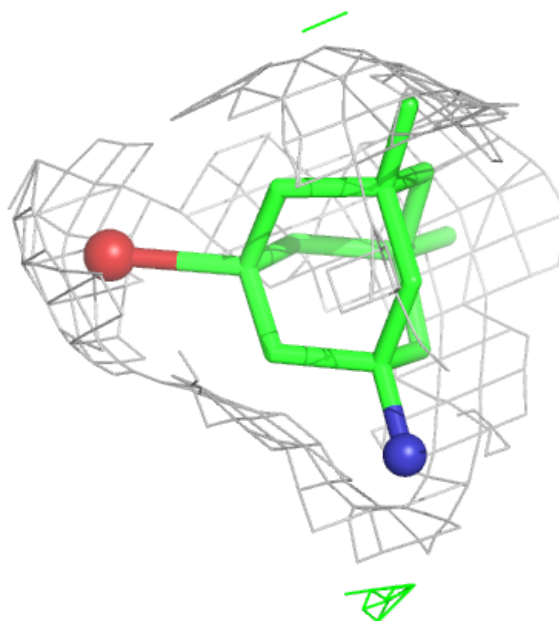
Electron density around BR7 C 401:

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



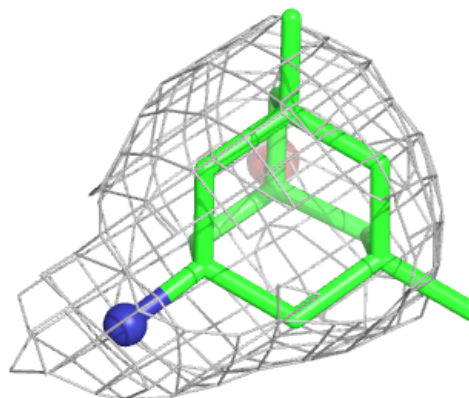
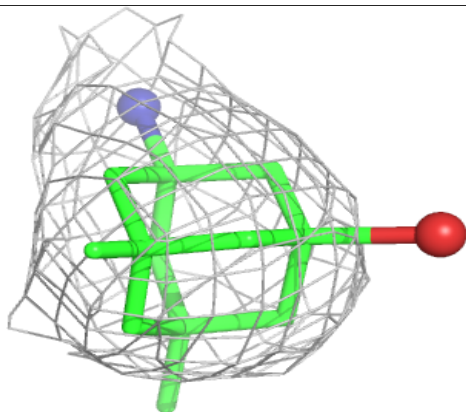
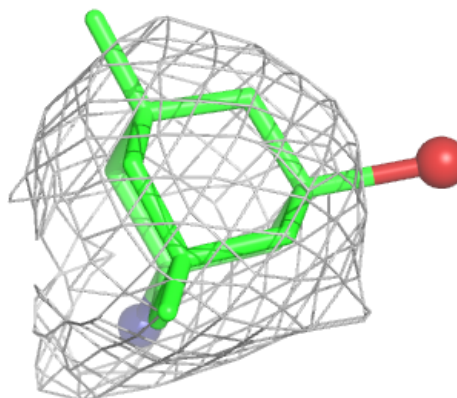
Electron density around BR7 B 401:

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and green (positive)



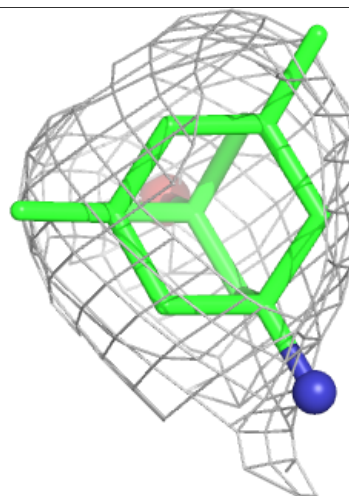
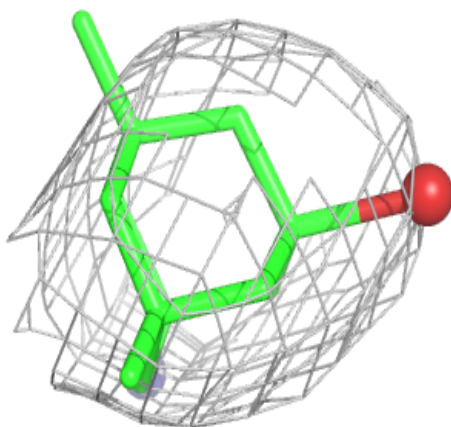
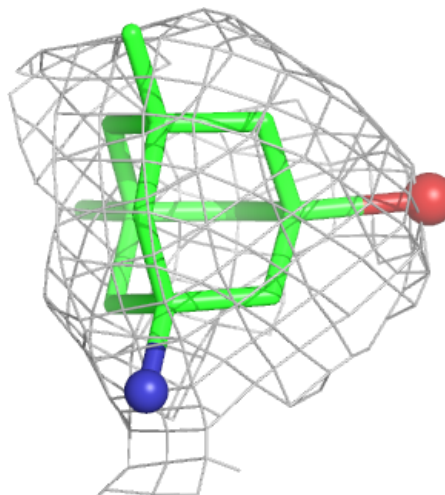
Electron density around BR7 A 401:

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



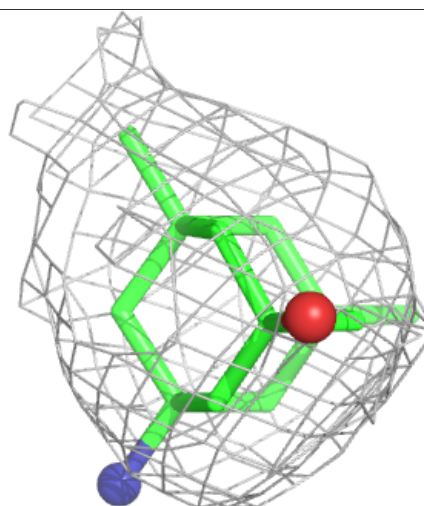
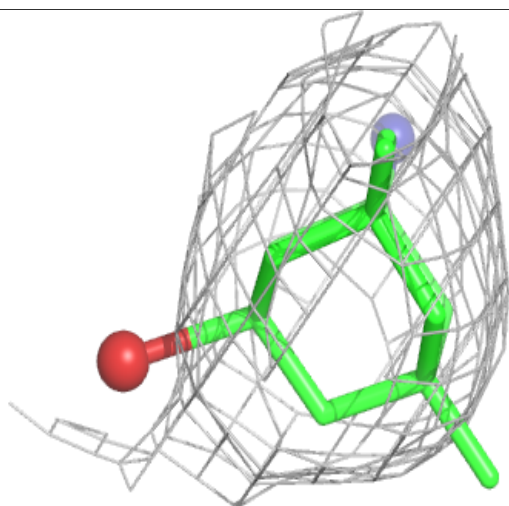
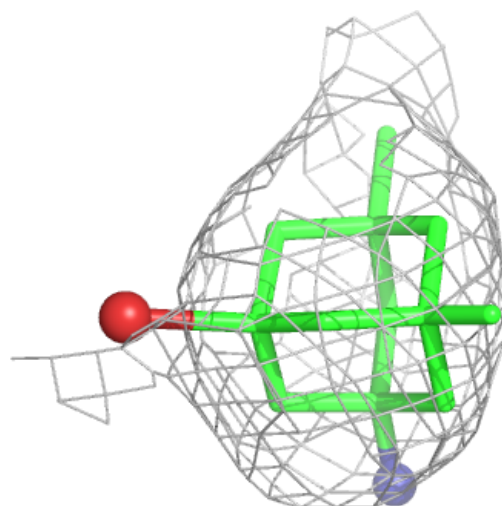
Electron density around BR7 G 401:

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



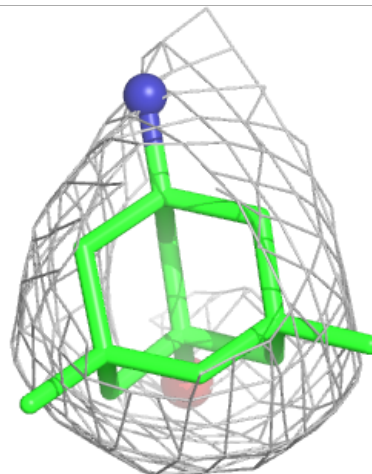
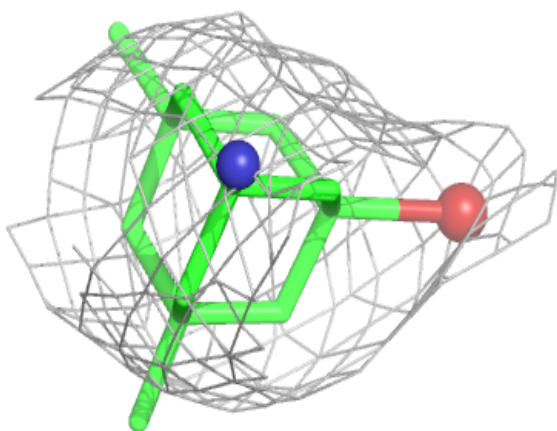
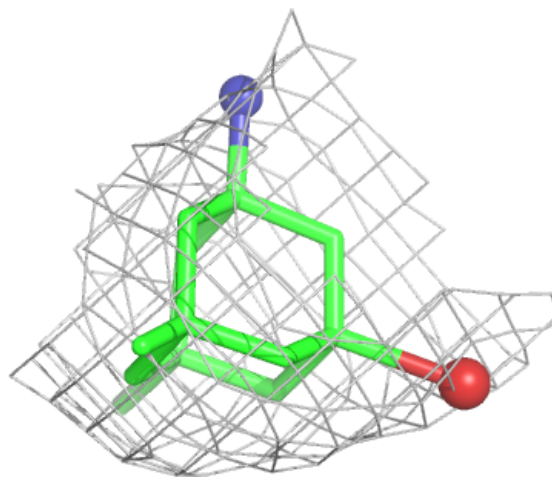
Electron density around BR7 F 402:

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



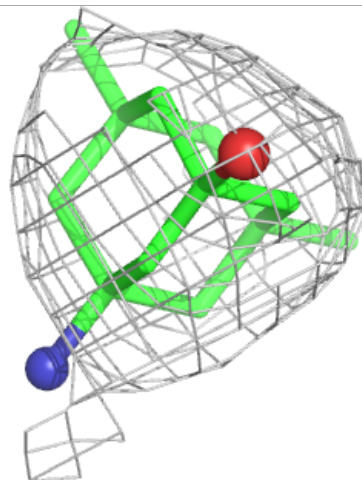
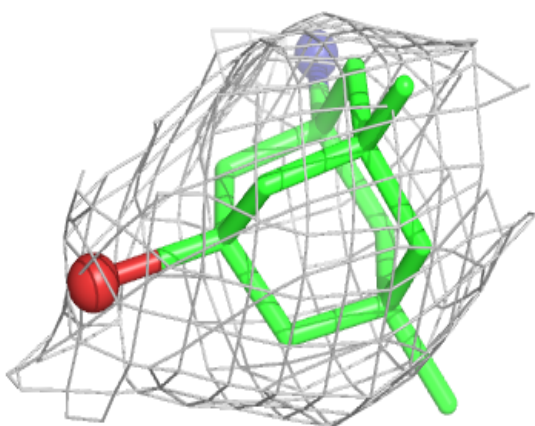
Electron density around BR7 D 401:

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and green (positive)



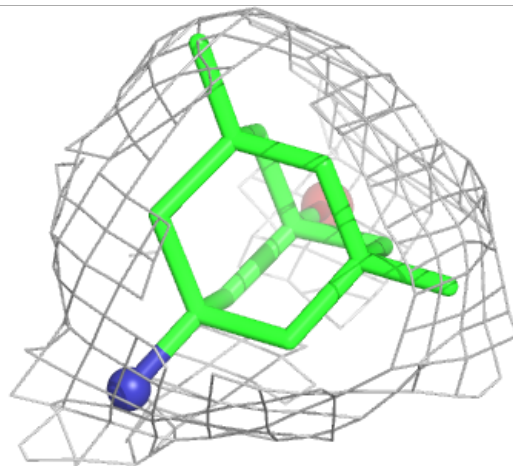
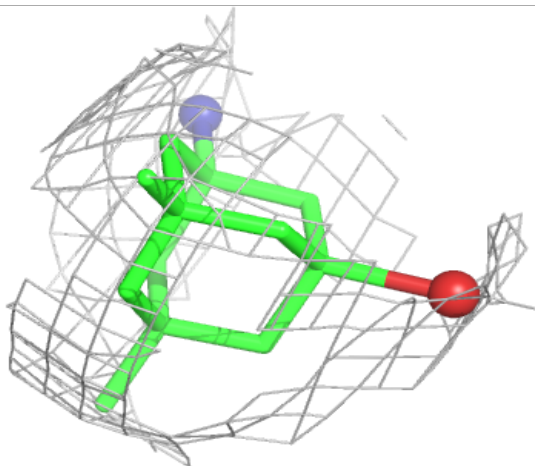
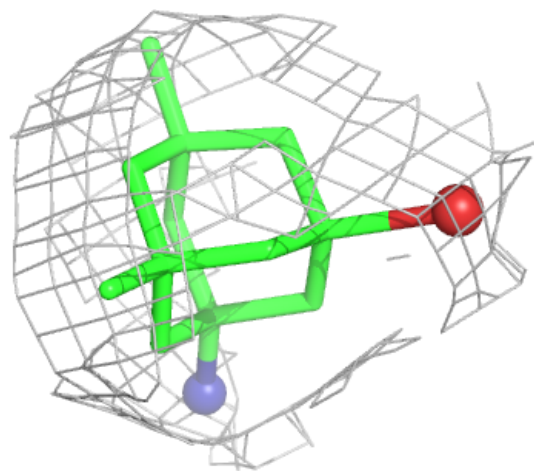
Electron density around BR7 H 401:

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and green (positive)



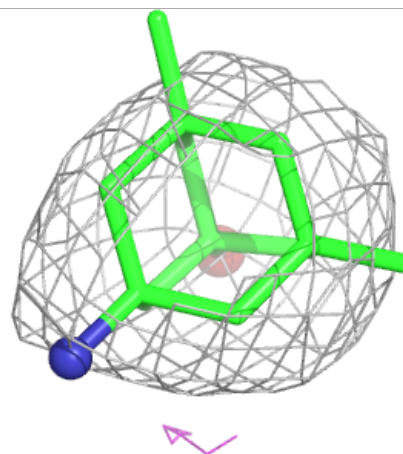
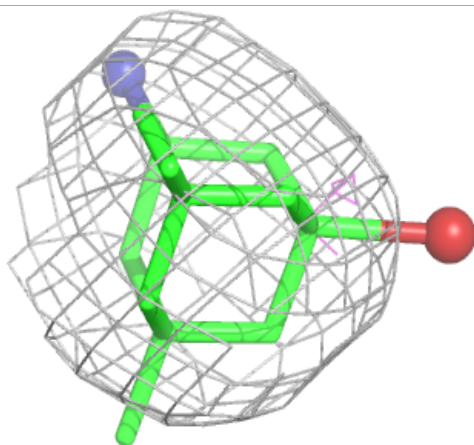
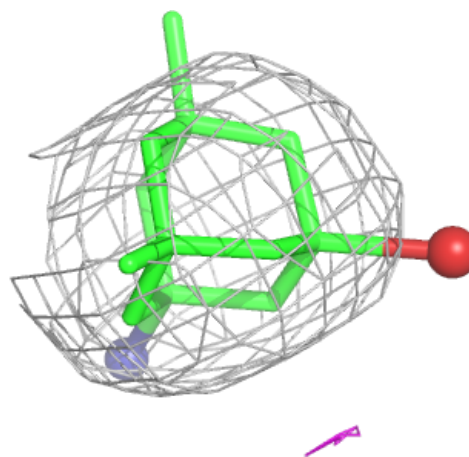
Electron density around BR7 I 401:

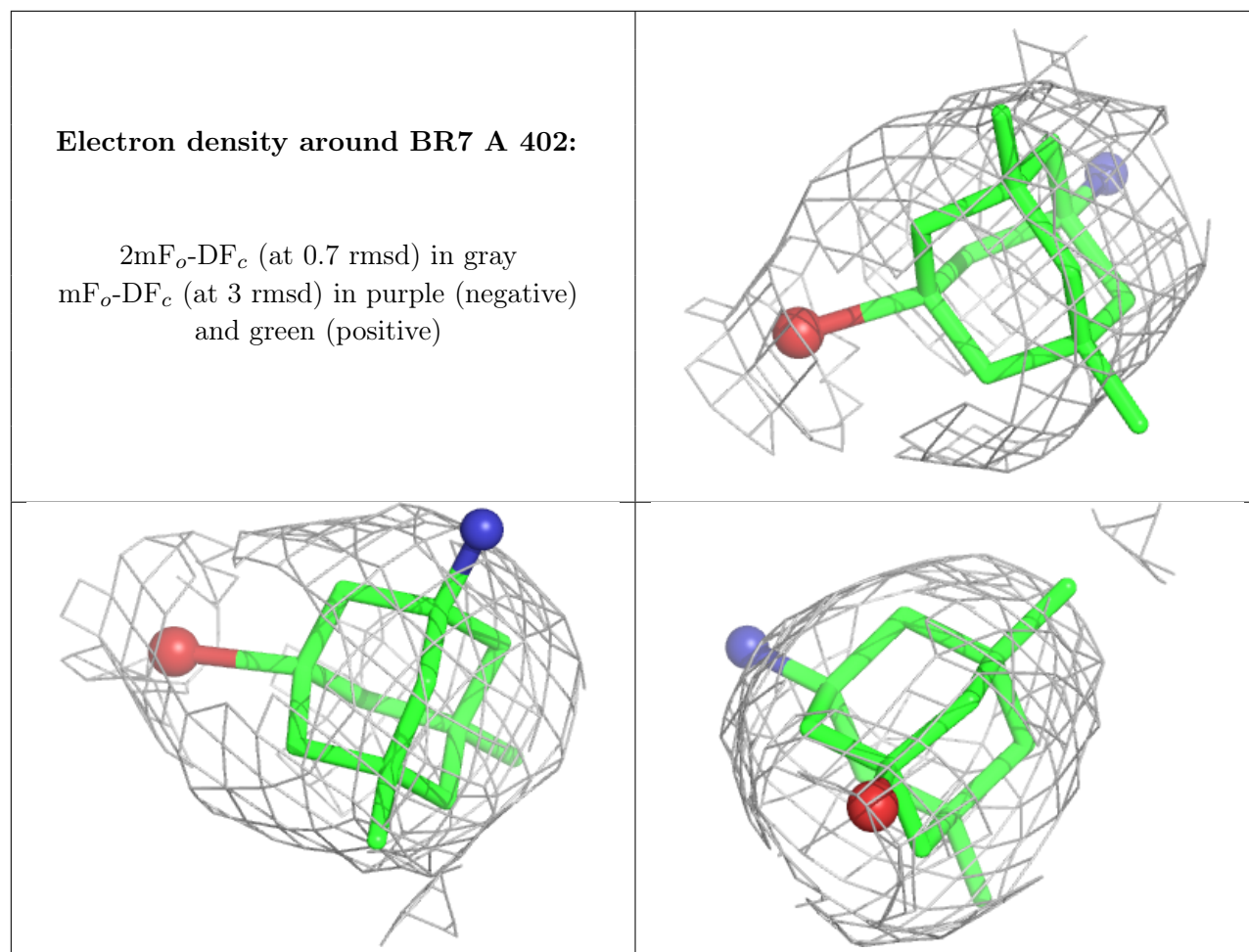
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around BR7 E 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.