



Full wwPDB EM Validation Report ⓘ

Mar 9, 2026 – 02:15 AM UTC

PDB ID : 8SSB / pdb_00008ssb
EMDB ID : EMD-40750
Title : Structure of LBD-TMD of AMPA receptor GluA2 in complex with auxiliary subunits TARP gamma-5 and cornichon-2 bound to glutamate and channel blocker spermidine (desensitized state)
Authors : Gangwar, S.P.; Yen, L.Y.; Yelshanskaya, M.V.; Sobolevsky, A.I.
Deposited on : 2023-05-08
Resolution : 3.66 Å(reported)

This is a Full wwPDB EM 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/EMValidationReportHelp>
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:

EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

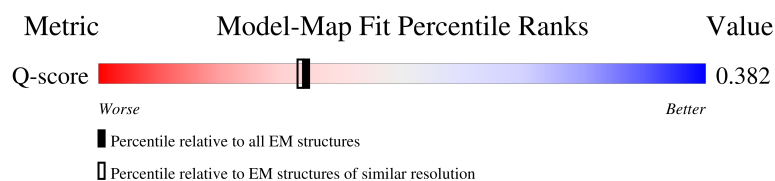
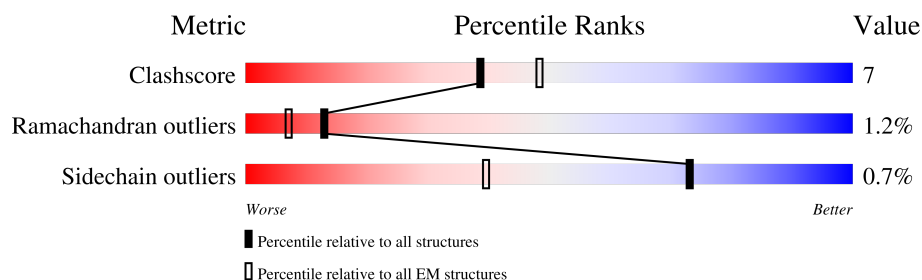
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.66 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)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	11491 (3.16 - 4.16)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the 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 EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	1026	 5% 47% 11% 41%
1	B	1026	 1% 31% 9% 60%
1	C	1026	 5% 47% 11% 41%
1	D	1026	 1% 30% 9% 60%

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Mol	Chain	Length	Quality of chain
2	E	160	15% 71% 15% • 12%
2	F	160	16% 71% 16% • 12%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 19048 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, 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 Glutamate receptor 2, Voltage-dependent calcium channel gamma-5 subunit chimera.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	602	Total	C	N	O	S	0	0
			4698	3062	751	853	32		
1	B	408	Total	C	N	O	S	0	0
			3177	2057	515	584	21		
1	C	602	Total	C	N	O	S	0	0
			4698	3062	751	853	32		
1	D	408	Total	C	N	O	S	0	0
			3177	2057	515	584	21		

There are 64 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	241	GLU	ASN	conflict	UNP P19491
A	382	LEU	VAL	conflict	UNP P19491
A	?	-	LEU	deletion	UNP P19491
A	?	-	THR	deletion	UNP P19491
A	?	-	GLU	deletion	UNP P19491
A	?	-	LEU	deletion	UNP P19491
A	?	-	PRO	deletion	UNP P19491
A	?	-	SER	deletion	UNP P19491
A	384	GLU	GLY	conflict	UNP P19491
A	385	ASP	ASN	conflict	UNP P19491
A	392	GLN	ASN	conflict	UNP P19491
A	827	GLY	-	linker	UNP P19491
A	828	THR	-	linker	UNP P19491
A	829	GLY	-	linker	UNP P19491
A	830	SER	-	linker	UNP P19491
A	831	ALA	-	linker	UNP P19491
B	241	GLU	ASN	conflict	UNP P19491
B	382	LEU	VAL	conflict	UNP P19491
B	?	-	LEU	deletion	UNP P19491
B	?	-	THR	deletion	UNP P19491
B	?	-	GLU	deletion	UNP P19491

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Chain	Residue	Modelled	Actual	Comment	Reference
B	?	-	LEU	deletion	UNP P19491
B	?	-	PRO	deletion	UNP P19491
B	?	-	SER	deletion	UNP P19491
B	384	GLU	GLY	conflict	UNP P19491
B	385	ASP	ASN	conflict	UNP P19491
B	392	GLN	ASN	conflict	UNP P19491
B	827	GLY	-	linker	UNP P19491
B	828	THR	-	linker	UNP P19491
B	829	GLY	-	linker	UNP P19491
B	830	SER	-	linker	UNP P19491
B	831	ALA	-	linker	UNP P19491
C	241	GLU	ASN	conflict	UNP P19491
C	382	LEU	VAL	conflict	UNP P19491
C	?	-	LEU	deletion	UNP P19491
C	?	-	THR	deletion	UNP P19491
C	?	-	GLU	deletion	UNP P19491
C	?	-	LEU	deletion	UNP P19491
C	?	-	PRO	deletion	UNP P19491
C	?	-	SER	deletion	UNP P19491
C	384	GLU	GLY	conflict	UNP P19491
C	385	ASP	ASN	conflict	UNP P19491
C	392	GLN	ASN	conflict	UNP P19491
C	827	GLY	-	linker	UNP P19491
C	828	THR	-	linker	UNP P19491
C	829	GLY	-	linker	UNP P19491
C	830	SER	-	linker	UNP P19491
C	831	ALA	-	linker	UNP P19491
D	241	GLU	ASN	conflict	UNP P19491
D	382	LEU	VAL	conflict	UNP P19491
D	?	-	LEU	deletion	UNP P19491
D	?	-	THR	deletion	UNP P19491
D	?	-	GLU	deletion	UNP P19491
D	?	-	LEU	deletion	UNP P19491
D	?	-	PRO	deletion	UNP P19491
D	?	-	SER	deletion	UNP P19491
D	384	GLU	GLY	conflict	UNP P19491
D	385	ASP	ASN	conflict	UNP P19491
D	392	GLN	ASN	conflict	UNP P19491
D	827	GLY	-	linker	UNP P19491
D	828	THR	-	linker	UNP P19491
D	829	GLY	-	linker	UNP P19491
D	830	SER	-	linker	UNP P19491

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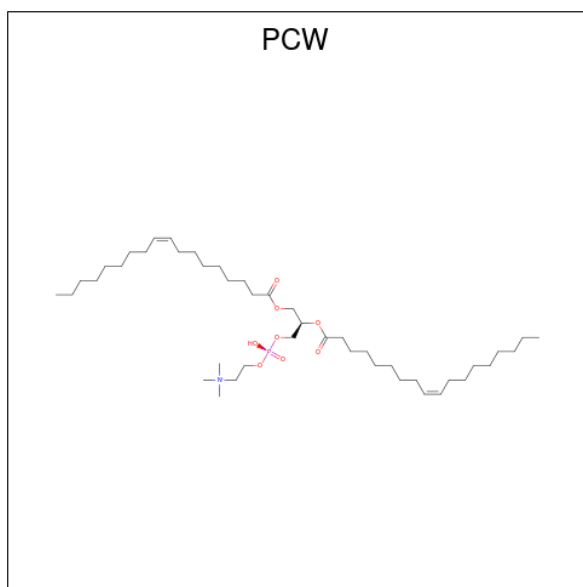
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Chain	Residue	Modelled	Actual	Comment	Reference
D	831	ALA	-	linker	UNP P19491

- Molecule 2 is a protein called Protein cornichon homolog 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	E	140	Total	C	N	O	S	0	0
			1166	787	175	191	13		
2	F	140	Total	C	N	O	S	0	0
			1166	787	175	191	13		

- Molecule 3 is 1,2-DIOLEOYL-SN-GLYCERO-3-PHOSPHOCHOLINE (CCD ID: PCW) (formula: $C_{44}H_{85}NO_8P$).



Mol	Chain	Residues	Atoms					AltConf
3	A	1	Total	C	N	O	P	0
			51	41	1	8	1	
3	A	1	Total	C	N	O	P	0
			51	41	1	8	1	
3	A	1	Total	C	N	O	P	0
			43	33	1	8	1	
3	A	1	Total	C	N	O	P	0
			51	41	1	8	1	
3	A	1	Total	C				0
			11	11				
3	A	1	Total	C				0
			11	11				

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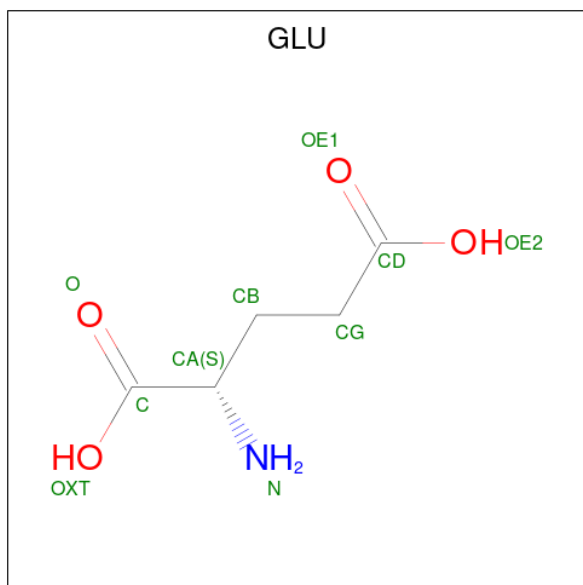
Mol	Chain	Residues	Atoms					AltConf
3	A	1	Total	C				0
			11	11				
3	A	1	Total	C				0
			11	11				
3	B	1	Total	C	N	O	P	0
			41	31	1	8	1	
3	B	1	Total	C	N	O	P	0
			43	33	1	8	1	
3	B	1	Total	C	N	O	P	0
			41	31	1	8	1	
3	B	1	Total	C	N	O	P	0
			51	41	1	8	1	
3	C	1	Total	C	N	O	P	0
			51	41	1	8	1	
3	C	1	Total	C	N	O	P	0
			43	33	1	8	1	
3	C	1	Total	C	N	O	P	0
			51	41	1	8	1	
3	C	1	Total	C				0
			11	11				
3	C	1	Total	C				0
			11	11				
3	C	1	Total	C				0
			11	11				
3	C	1	Total	C				0
			11	11				
3	D	1	Total	C	N	O	P	0
			41	31	1	8	1	
3	D	1	Total	C	N	O	P	0
			43	33	1	8	1	
3	D	1	Total	C	N	O	P	0
			39	29	1	8	1	
3	D	1	Total	C	N	O	P	0
			41	31	1	8	1	
3	E	1	Total	C	N	O	P	0
			39	29	1	8	1	
3	E	1	Total	C				0
			11	11				
3	E	1	Total	C	N	O	P	0
			43	33	1	8	1	
3	F	1	Total	C				0
			11	11				

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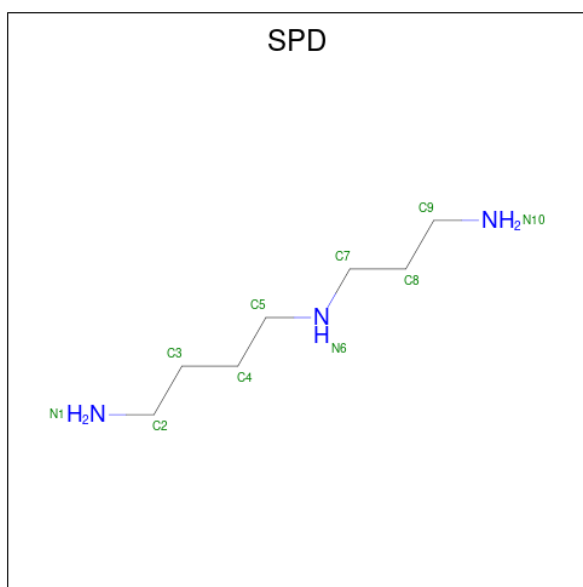
Mol	Chain	Residues	Atoms					AltConf
3	F	1	Total	C	N	O	P	0
			43	33	1	8	1	

- Molecule 4 is GLUTAMIC ACID (CCD ID: GLU) (formula: $C_5H_9NO_4$) (labeled as "Ligand of Interest" by depositor).



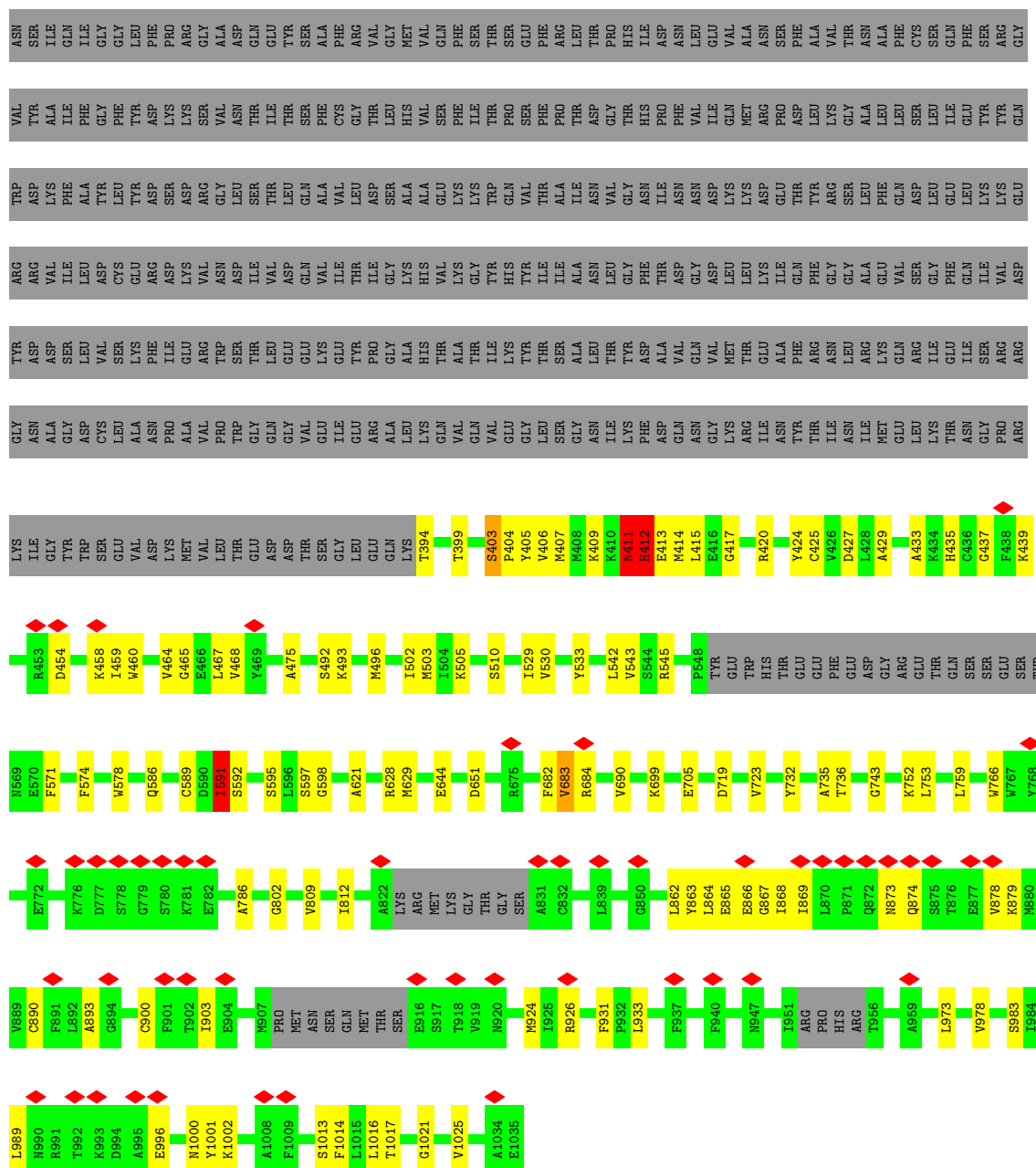
Mol	Chain	Residues	Atoms					AltConf
4	A	1	Total	C	N	O		0
			10	5	1	4		
4	B	1	Total	C	N	O		0
			10	5	1	4		
4	C	1	Total	C	N	O		0
			10	5	1	4		
4	D	1	Total	C	N	O		0
			10	5	1	4		

- Molecule 5 is SPERMIDINE (CCD ID: SPD) (formula: $C_7H_{19}N_3$) (labeled as "Ligand of Interest" by depositor).



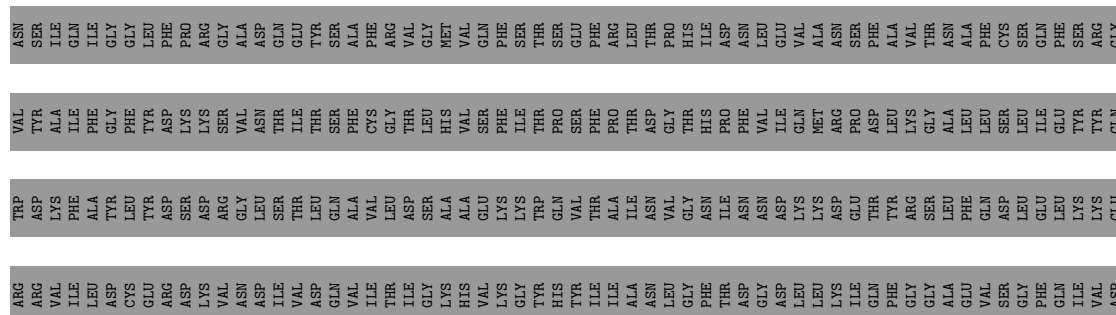
Mol	Chain	Residues	Atoms			AltConf
5	B	1	Total	C	N	0
			10	7	3	



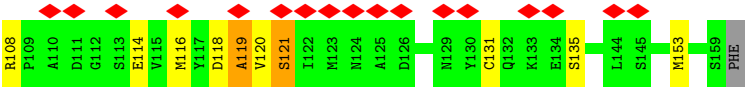


- Molecule 1: Glutamate receptor 2, Voltage-dependent calcium channel gamma-5 subunit chimera

Chain D: 30% 9% 60%







4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C2	Depositor
Number of particles used	48434	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	58	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.903	Depositor
Minimum map value	-0.519	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.016	Depositor
Recommended contour level	0.1	Depositor
Map size (Å)	345.28, 345.28, 345.28	wwPDB
Map dimensions	416, 416, 416	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.83, 0.83, 0.83	Depositor

5 Model quality

5.1 Standard geometry

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

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	0.32	0/4799	0.82	8/6479 (0.1%)
1	B	0.34	0/3245	0.85	1/4378 (0.0%)
1	C	0.32	0/4799	0.82	8/6479 (0.1%)
1	D	0.34	0/3245	0.85	1/4378 (0.0%)
2	E	0.31	0/1203	0.77	2/1636 (0.1%)
2	F	0.31	0/1203	0.78	2/1636 (0.1%)
All	All	0.33	0/18494	0.82	22/24986 (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	10
1	B	0	6
1	C	0	10
1	D	0	6
2	E	0	2
2	F	0	2
All	All	0	36

There are no bond length outliers.

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	411	ASN	CA-C-N	6.57	134.09	121.54
1	A	411	ASN	C-N-CA	6.57	134.09	121.54
1	C	411	ASN	CA-C-N	6.55	134.05	121.54
1	C	411	ASN	C-N-CA	6.55	134.05	121.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	454	ASP	CA-C-N	6.40	129.69	120.79
1	C	454	ASP	C-N-CA	6.40	129.69	120.79
1	A	454	ASP	CA-C-N	6.40	129.69	120.79
1	A	454	ASP	C-N-CA	6.40	129.69	120.79
1	C	591	ILE	CA-C-N	6.33	137.26	121.80
1	C	591	ILE	C-N-CA	6.33	137.26	121.80
1	A	591	ILE	CA-C-N	6.33	137.24	121.80
1	A	591	ILE	C-N-CA	6.33	137.24	121.80
1	A	407	MET	CA-C-N	5.26	131.59	121.54
1	A	407	MET	C-N-CA	5.26	131.59	121.54
1	C	407	MET	CA-C-N	5.26	131.59	121.54
1	C	407	MET	C-N-CA	5.26	131.59	121.54
1	D	784	THR	N-CA-C	5.14	117.54	110.35
2	F	121	SER	CA-C-N	5.12	129.01	120.62
2	F	121	SER	C-N-CA	5.12	129.01	120.62
1	B	784	THR	N-CA-C	5.11	117.51	110.35
2	E	121	SER	CA-C-N	5.01	128.83	120.62
2	E	121	SER	C-N-CA	5.01	128.83	120.62

There are no chirality outliers.

All (36) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	403	SER	Peptide
1	A	411	ASN	Peptide
1	A	412	HIS	Peptide
1	A	458	LYS	Peptide
1	A	510	SER	Peptide
1	A	589	CYS	Peptide
1	A	591	ILE	Peptide
1	A	683	VAL	Peptide
1	A	873	ASN	Peptide
1	A	878	VAL	Peptide
1	B	403	SER	Peptide
1	B	449	LYS	Peptide
1	B	715	ARG	Peptide
1	B	716	LYS	Peptide
1	B	739	GLY	Peptide
1	B	783	LYS	Peptide
1	C	403	SER	Peptide
1	C	411	ASN	Peptide
1	C	412	HIS	Peptide

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Mol	Chain	Res	Type	Group
1	C	458	LYS	Peptide
1	C	510	SER	Peptide
1	C	589	CYS	Peptide
1	C	591	ILE	Peptide
1	C	683	VAL	Peptide
1	C	873	ASN	Peptide
1	C	878	VAL	Peptide
1	D	403	SER	Peptide
1	D	449	LYS	Peptide
1	D	715	ARG	Peptide
1	D	716	LYS	Peptide
1	D	739	GLY	Peptide
1	D	783	LYS	Peptide
2	E	105	TYR	Peptide
2	E	119	ALA	Peptide
2	F	105	TYR	Peptide
2	F	119	ALA	Peptide

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	4698	0	4761	64	0
1	B	3177	0	3221	56	0
1	C	4698	0	4761	60	0
1	D	3177	0	3221	59	0
2	E	1166	0	1152	16	0
2	F	1166	0	1152	15	0
3	A	240	0	346	13	0
3	B	176	0	241	12	0
3	C	189	0	271	13	0
3	D	164	0	216	11	0
3	E	93	0	123	4	0
3	F	54	0	73	3	0
4	A	10	0	5	0	0
4	B	10	0	5	2	0
4	C	10	0	5	0	0
4	D	10	0	5	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	B	10	0	19	2	0
All	All	19048	0	19577	283	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 (283) 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:480:THR:HG1	4:B:1104:GLU:N	1.72	0.86
1:D:480:THR:HG1	4:D:1105:GLU:N	1.79	0.80
3:B:1105:PCW:H121	3:C:1102:PCW:H39	1.68	0.76
1:A:651:ASP:HB2	1:A:682:PHE:HB3	1.76	0.67
1:C:651:ASP:HB2	1:C:682:PHE:HB3	1.76	0.66
2:E:28:ILE:HG22	2:E:135:SER:HB2	1.76	0.66
1:A:863:TYR:HB2	1:A:1002:LYS:HB3	1.79	0.65
1:C:863:TYR:HB2	1:C:1002:LYS:HB3	1.79	0.64
2:F:28:ILE:HG22	2:F:135:SER:HB2	1.80	0.63
1:D:773:CYS:SG	1:D:774:GLY:N	2.71	0.63
1:B:773:CYS:SG	1:B:774:GLY:N	2.71	0.62
1:B:646:ALA:H	1:B:699:LYS:HB3	1.65	0.61
1:D:646:ALA:H	1:D:699:LYS:HB3	1.65	0.61
1:A:888:ARG:HB3	1:A:900:CYS:HB3	1.84	0.60
1:B:467:LEU:HA	1:B:472:ALA:HB3	1.84	0.60
1:C:888:ARG:HB3	1:C:900:CYS:HB3	1.84	0.59
1:D:467:LEU:HA	1:D:472:ALA:HB3	1.84	0.59
1:A:867:GLY:HA2	1:A:879:LYS:HB3	1.85	0.58
1:C:869:ILE:HG22	1:C:996:GLU:HA	1.86	0.58
1:B:715:ARG:HA	1:B:770:LYS:HG3	1.86	0.58
1:B:511:LYS:NZ	1:B:512:PRO:O	2.34	0.58
2:E:104:ARG:NH2	2:E:131:CYS:SG	2.77	0.58
1:C:867:GLY:HA2	1:C:879:LYS:HB3	1.85	0.58
1:D:715:ARG:HA	1:D:770:LYS:HG3	1.86	0.57
1:A:869:ILE:HG22	1:A:996:GLU:HA	1.86	0.57
1:D:493:LYS:HD2	1:D:751:LEU:HD21	1.86	0.57
1:D:511:LYS:NZ	1:D:512:PRO:O	2.34	0.57
1:B:604:VAL:HG21	1:C:802:GLY:HA3	1.87	0.57
1:B:493:LYS:HD2	1:B:751:LEU:HD21	1.86	0.56
3:C:1103:PCW:H121	3:C:1103:PCW:H321	1.88	0.56
1:A:600:ILE:HG12	3:B:1102:PCW:H412	1.86	0.56
1:D:694:ARG:NH1	1:D:719:ASP:OD2	2.39	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:104:ARG:NH2	2:F:131:CYS:SG	2.78	0.56
1:A:586:GLN:HG2	1:B:587:GLN:HE22	1.70	0.56
1:B:694:ARG:NH1	1:B:719:ASP:OD2	2.39	0.56
1:B:406:VAL:HG22	1:B:426:VAL:HG23	1.88	0.56
1:A:1002:LYS:NZ	3:A:1103:PCW:O1P	2.39	0.55
3:A:1104:PCW:H321	3:A:1104:PCW:H121	1.88	0.55
1:D:406:VAL:HG22	1:D:426:VAL:HG23	1.88	0.55
1:A:427:ASP:OD2	1:A:766:TRP:NE1	2.36	0.55
1:C:973:LEU:HD23	1:C:1014:PHE:HB2	1.88	0.55
3:D:1101:PCW:H20	3:D:1102:PCW:H422	1.87	0.55
3:D:1103:PCW:H331	2:F:70:PRO:HA	1.89	0.55
1:B:405:TYR:HA	1:B:424:TYR:HB3	1.89	0.55
1:A:409:LYS:HG3	1:A:411:ASN:H	1.72	0.55
1:A:629:MET:HE2	1:C:629:MET:HE2	1.89	0.55
1:A:644:GLU:OE2	1:A:699:LYS:NZ	2.40	0.55
1:C:644:GLU:OE2	1:C:699:LYS:NZ	2.40	0.55
1:C:475:ALA:HB3	1:C:735:ALA:HB3	1.89	0.55
1:D:403:SER:O	1:D:405:TYR:N	2.39	0.55
1:D:405:TYR:HA	1:D:424:TYR:HB3	1.89	0.55
1:A:475:ALA:HB3	1:A:735:ALA:HB3	1.89	0.54
1:C:409:LYS:HG3	1:C:411:ASN:H	1.72	0.54
1:A:405:TYR:HA	1:A:424:TYR:HB3	1.89	0.54
1:A:973:LEU:HD23	1:A:1014:PHE:HB2	1.88	0.54
1:B:802:GLY:HA2	3:B:1103:PCW:H20	1.89	0.54
1:B:403:SER:O	1:B:405:TYR:N	2.39	0.54
3:B:1101:PCW:H20	3:B:1102:PCW:H422	1.90	0.54
1:C:405:TYR:HA	1:C:424:TYR:HB3	1.88	0.54
1:B:599:ARG:HE	1:C:578:TRP:CG	2.25	0.54
1:A:586:GLN:HB3	5:B:1106:SPD:H81	1.90	0.53
1:C:427:ASP:OD2	1:C:766:TRP:NE1	2.36	0.53
1:C:924:MET:HB3	1:C:983:SER:HB3	1.90	0.53
3:A:1101:PCW:H161	3:A:1103:PCW:H40	1.91	0.53
1:D:694:ARG:HH22	1:D:717:PRO:HG2	1.74	0.53
1:A:492:SER:OG	1:A:493:LYS:N	2.42	0.53
1:A:736:THR:HG21	1:A:743:GLY:HA2	1.91	0.53
1:C:736:THR:HG21	1:C:743:GLY:HA2	1.91	0.53
1:A:924:MET:HB3	1:A:983:SER:HB3	1.90	0.52
1:C:465:GLY:HA2	1:C:468:VAL:HG12	1.92	0.52
1:D:401:LEU:HB3	1:D:444:ILE:HD13	1.92	0.52
1:B:544:SER:O	1:B:569:ASN:ND2	2.42	0.52
1:D:397:VAL:HG23	1:D:441:LYS:H	1.75	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:401:LEU:HB3	1:B:444:ILE:HD13	1.92	0.52
1:A:729:SER:HB3	1:D:497:SER:HB3	1.92	0.52
1:B:694:ARG:HH22	1:B:717:PRO:HG2	1.74	0.52
1:D:710:GLU:OE2	1:D:722:LYS:NZ	2.43	0.52
2:E:108:ARG:HH12	2:E:119:ALA:HB2	1.74	0.52
1:D:544:SER:O	1:D:569:ASN:ND2	2.42	0.52
1:A:399:THR:HB	1:A:406:VAL:HG21	1.92	0.52
1:B:397:VAL:HG23	1:B:441:LYS:H	1.75	0.52
1:B:404:PRO:HA	1:B:407:MET:HE1	1.93	0.51
3:B:1105:PCW:H131	3:C:1101:PCW:H371	1.92	0.51
1:C:989:LEU:O	1:D:508:GLN:NE2	2.43	0.51
3:A:1101:PCW:H121	3:A:1103:PCW:H39	1.92	0.51
1:C:399:THR:HB	1:C:406:VAL:HG21	1.92	0.51
1:C:492:SER:OG	1:C:493:LYS:N	2.42	0.51
1:D:421:TYR:HD2	1:D:442:LEU:HB3	1.76	0.51
1:A:543:VAL:HG23	1:B:810:ALA:HB2	1.93	0.51
1:B:520:PRO:O	1:B:619:ASN:ND2	2.41	0.51
1:D:542:LEU:HD11	2:F:74:ILE:HD11	1.93	0.50
1:A:663:LYS:HE2	1:D:757:GLY:HA3	1.92	0.50
1:B:710:GLU:OE2	1:B:722:LYS:NZ	2.43	0.50
1:A:465:GLY:HA2	1:A:468:VAL:HG12	1.92	0.50
1:C:705:GLU:OE1	1:C:732:TYR:OH	2.28	0.50
1:D:404:PRO:HA	1:D:407:MET:HE1	1.93	0.50
1:B:626:VAL:O	1:C:628:ARG:NE	2.39	0.50
1:A:394:THR:HA	1:A:439:LYS:HB2	1.93	0.50
1:A:417:GLY:O	1:A:420:ARG:NH1	2.43	0.50
3:B:1105:PCW:H161	3:C:1102:PCW:H40	1.92	0.50
1:D:574:PHE:HB2	3:D:1103:PCW:H141	1.93	0.50
1:B:421:TYR:HD2	1:B:442:LEU:HB3	1.76	0.49
2:E:118:ASP:OD1	2:E:118:ASP:N	2.45	0.49
1:C:425:CYS:O	1:C:429:ALA:N	2.40	0.49
2:E:75:HIS:HE1	2:E:94:ASN:HB2	1.78	0.49
1:A:887:TRP:HZ2	1:A:933:LEU:HD11	1.78	0.49
1:C:887:TRP:HZ2	1:C:933:LEU:HD11	1.78	0.49
1:C:394:THR:HA	1:C:439:LYS:HB2	1.93	0.49
1:C:1002:LYS:NZ	3:C:1102:PCW:O1P	2.46	0.49
1:A:578:TRP:CG	1:D:599:ARG:HE	2.31	0.49
1:C:417:GLY:O	1:C:420:ARG:NH1	2.44	0.49
2:E:119:ALA:O	2:E:121:SER:N	2.46	0.49
1:C:862:LEU:HD11	1:C:1001:TYR:HB2	1.94	0.48
2:F:108:ARG:HH12	2:F:119:ALA:HB2	1.79	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:753:LEU:HB3	1:A:759:LEU:HD23	1.95	0.48
2:E:70:PRO:HA	3:E:201:PCW:H331	1.94	0.48
1:A:729:SER:OG	1:D:729:SER:OG	2.26	0.48
1:D:409:LYS:HD2	1:D:422:GLU:HG2	1.96	0.48
3:A:1101:PCW:H122	3:A:1101:PCW:H381	1.95	0.48
2:E:11:MET:HE2	2:E:153:MET:HE3	1.95	0.48
2:F:118:ASP:OD1	2:F:118:ASP:N	2.45	0.48
1:A:502:ILE:HB	1:A:723:VAL:HB	1.96	0.48
1:A:862:LEU:HD11	1:A:1001:TYR:HB2	1.94	0.48
1:B:618:ALA:HA	1:C:621:ALA:HB2	1.96	0.47
1:C:865:GLU:HB2	1:C:1000:ASN:HB2	1.95	0.47
1:C:1013:SER:HA	1:C:1016:LEU:HB2	1.96	0.47
1:A:503:MET:HE1	1:A:690:VAL:HG22	1.95	0.47
1:A:425:CYS:O	1:A:429:ALA:N	2.40	0.47
1:C:753:LEU:HB3	1:C:759:LEU:HD23	1.95	0.47
1:D:520:PRO:O	1:D:619:ASN:ND2	2.41	0.47
1:B:542:LEU:HD11	2:E:74:ILE:HD11	1.95	0.47
1:C:433:ALA:HA	1:C:437:GLY:HA3	1.96	0.47
1:D:399:THR:OG1	1:D:400:ILE:N	2.47	0.47
2:F:119:ALA:O	2:F:121:SER:N	2.48	0.47
1:A:865:GLU:HB2	1:A:1000:ASN:HB2	1.95	0.47
3:B:1101:PCW:H132	3:B:1101:PCW:H162	1.55	0.47
1:A:1013:SER:HA	1:A:1016:LEU:HB2	1.96	0.47
1:A:433:ALA:HA	1:A:437:GLY:HA3	1.96	0.47
1:C:403:SER:O	1:C:405:TYR:N	2.47	0.47
1:C:543:VAL:HG22	1:C:597:SER:HB3	1.97	0.47
3:C:1108:PCW:H211	3:C:1108:PCW:H181	1.74	0.47
3:D:1101:PCW:H162	3:D:1101:PCW:H132	1.52	0.47
2:F:75:HIS:HE1	2:F:94:ASN:HB2	1.80	0.47
1:A:543:VAL:HG22	1:A:597:SER:HB3	1.97	0.46
1:B:409:LYS:HD2	1:B:422:GLU:HG2	1.96	0.46
1:C:503:MET:HE1	1:C:690:VAL:HG22	1.95	0.46
3:B:1105:PCW:H381	3:B:1105:PCW:H122	1.96	0.46
1:C:502:ILE:HB	1:C:723:VAL:HB	1.96	0.46
3:E:203:PCW:H351	3:E:203:PCW:H321	1.73	0.46
1:A:705:GLU:OE1	1:A:732:TYR:OH	2.28	0.46
3:E:203:PCW:H182	3:E:203:PCW:H151	1.73	0.46
2:F:11:MET:HE2	2:F:153:MET:HE3	1.98	0.46
3:A:1109:PCW:H181	3:A:1109:PCW:H211	1.73	0.46
1:B:574:PHE:HB2	3:E:201:PCW:H141	1.97	0.46
3:C:1103:PCW:H131	3:C:1103:PCW:H162	1.73	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:656:LYS:HD3	1:D:660:ARG:HH21	1.81	0.46
1:B:399:THR:OG1	1:B:400:ILE:N	2.47	0.46
1:D:798:ILE:HD13	3:D:1102:PCW:H321	1.98	0.46
2:F:65:ARG:HD3	2:F:114:GLU:HB3	1.97	0.46
3:B:1105:PCW:H412	3:B:1105:PCW:H162	1.99	0.45
1:D:649:THR:HA	1:D:703:LEU:H	1.81	0.45
1:C:545:ARG:NH1	1:C:571:PHE:O	2.50	0.45
2:F:56:ASN:OD1	2:F:59:ARG:NH1	2.50	0.45
1:A:545:ARG:NH1	1:A:571:PHE:O	2.50	0.45
1:B:503:MET:HE1	1:B:690:VAL:HG22	1.99	0.45
1:C:931:PHE:HE1	3:C:1104:PCW:H19	1.82	0.45
2:F:82:PHE:HA	2:F:85:ALA:HB3	1.97	0.45
1:A:606:TRP:CZ3	1:B:582:GLY:HA2	2.52	0.45
1:B:656:LYS:HD3	1:B:660:ARG:HH21	1.81	0.45
1:B:685:THR:HG22	1:B:688:GLU:H	1.82	0.45
1:D:397:VAL:HG13	1:D:476:ILE:HG13	1.99	0.45
1:D:572:GLY:HA3	3:D:1103:PCW:H73	1.97	0.45
1:B:666:VAL:HA	1:B:669:LYS:HE2	1.99	0.45
1:A:596:LEU:HD23	1:B:574:PHE:HE1	1.82	0.44
1:A:1017:THR:O	1:A:1021:GLY:N	2.50	0.44
2:E:65:ARG:HD3	2:E:114:GLU:HB3	1.99	0.44
3:A:1102:PCW:H351	3:A:1102:PCW:H382	1.76	0.44
1:C:881:SER:H	1:C:893:ALA:HB3	1.83	0.44
1:D:671:TRP:O	1:D:675:ARG:N	2.48	0.44
1:A:803:LEU:HD21	1:D:536:VAL:HG22	2.00	0.44
1:B:397:VAL:HG13	1:B:476:ILE:HG13	1.99	0.44
1:D:473:ASP:OD1	1:D:473:ASP:N	2.49	0.44
1:B:619:ASN:HD21	1:C:786:ALA:HA	1.83	0.44
1:C:1017:THR:O	1:C:1021:GLY:N	2.50	0.44
1:B:490:ASP:O	1:B:736:THR:OG1	2.27	0.44
1:D:666:VAL:HA	1:D:669:LYS:HE2	1.99	0.44
1:B:649:THR:HA	1:B:703:LEU:H	1.81	0.44
1:D:685:THR:HG22	1:D:688:GLU:H	1.82	0.44
1:C:586:GLN:HG2	1:D:587:GLN:HE22	1.81	0.44
1:D:394:THR:HG23	1:D:439:LYS:HB2	2.00	0.44
1:D:639:LEU:O	1:D:647:TYR:OH	2.29	0.44
1:D:654:SER:N	4:D:1105:GLU:OE2	2.51	0.44
1:A:403:SER:O	1:A:405:TYR:N	2.47	0.43
1:D:503:MET:HE1	1:D:690:VAL:HG22	1.99	0.43
3:A:1104:PCW:H131	3:A:1104:PCW:H162	1.83	0.43
1:B:671:TRP:O	1:B:675:ARG:N	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:1104:PCW:H181	3:C:1104:PCW:H211	1.82	0.43
1:B:654:SER:N	4:B:1104:GLU:OE1	2.51	0.43
1:A:881:SER:H	1:A:893:ALA:HB3	1.83	0.43
1:B:481:ILE:HD11	1:B:733:GLY:HA3	2.00	0.43
1:C:809:VAL:HA	1:C:812:ILE:HG22	2.01	0.43
1:D:464:VAL:HG11	1:D:485:ARG:HD2	2.00	0.43
1:B:775:ALA:HA	1:B:778:SER:HB2	2.01	0.43
1:C:505:LYS:NZ	1:C:719:ASP:OD2	2.39	0.43
1:C:530:VAL:HG11	3:C:1102:PCW:H182	2.01	0.43
3:C:1101:PCW:H121	3:C:1101:PCW:H152	1.77	0.43
3:C:1101:PCW:H351	3:C:1101:PCW:H382	1.76	0.43
1:D:481:ILE:HD11	1:D:733:GLY:HA3	2.00	0.43
1:D:775:ALA:HA	1:D:778:SER:HB2	2.01	0.43
1:A:529:ILE:O	1:A:533:TYR:N	2.51	0.43
1:A:809:VAL:HA	1:A:812:ILE:HG22	2.01	0.43
1:A:903:ILE:O	1:A:926:ARG:NH2	2.52	0.43
1:B:394:THR:HG23	1:B:439:LYS:HB2	2.00	0.43
1:A:412:HIS:O	1:A:414:MET:N	2.52	0.43
1:A:460:TRP:HB3	1:A:464:VAL:HG23	2.01	0.43
1:A:606:TRP:CD1	1:B:587:GLN:HG3	2.54	0.43
1:C:574:PHE:HB2	3:C:1103:PCW:H142	2.00	0.43
1:A:802:GLY:HA3	1:D:604:VAL:HG21	2.00	0.42
1:C:903:ILE:O	1:C:926:ARG:NH2	2.52	0.42
3:F:202:PCW:H351	3:F:202:PCW:H321	1.77	0.42
1:C:412:HIS:O	1:C:414:MET:N	2.52	0.42
1:C:405:TYR:HE1	1:C:496:MET:HE1	1.84	0.42
1:D:645:ILE:H	1:D:645:ILE:HG13	1.75	0.42
3:A:1101:PCW:H151	3:A:1102:PCW:H371	2.00	0.42
1:B:595:SER:OG	1:B:598:GLY:N	2.49	0.42
1:C:529:ILE:O	1:C:533:TYR:N	2.51	0.42
1:A:628:ARG:NE	1:D:626:VAL:O	2.37	0.42
3:B:1101:PCW:H122	3:B:1102:PCW:H331	2.02	0.42
1:C:435:HIS:CE1	1:C:752:LYS:HE2	2.55	0.42
1:C:542:LEU:HD12	1:C:1025:VAL:HG21	2.02	0.42
1:C:595:SER:OG	1:C:598:GLY:N	2.52	0.42
1:D:577:LEU:HD21	3:D:1101:PCW:H241	2.02	0.42
1:A:542:LEU:HD12	1:A:1025:VAL:HG21	2.02	0.42
1:B:541:PHE:HD2	1:B:542:LEU:HD12	1.85	0.42
3:B:1101:PCW:H51	3:B:1102:PCW:H42	2.01	0.42
3:B:1103:PCW:H181	3:B:1103:PCW:H212	1.77	0.42
1:D:595:SER:OG	1:D:598:GLY:N	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:75:HIS:CE1	2:F:94:ASN:HB2	2.55	0.42
1:B:464:VAL:HG11	1:B:485:ARG:HD2	2.00	0.41
3:D:1102:PCW:H381	3:D:1102:PCW:H351	1.69	0.41
1:B:473:ASP:OD1	1:B:473:ASP:N	2.49	0.41
3:F:202:PCW:H182	3:F:202:PCW:H151	1.73	0.41
1:A:505:LYS:NZ	1:A:719:ASP:OD2	2.39	0.41
1:A:586:GLN:O	5:B:1106:SPD:N6	2.53	0.41
3:A:1102:PCW:H121	3:A:1102:PCW:H152	1.76	0.41
1:C:460:TRP:HB3	1:C:464:VAL:HG23	2.01	0.41
1:D:757:GLY:O	1:D:761:LYS:N	2.51	0.41
3:D:1104:PCW:H212	3:D:1104:PCW:H181	1.72	0.41
2:F:58:GLU:HA	2:F:61:CYS:HB2	2.02	0.41
3:A:1101:PCW:H131	3:A:1102:PCW:H371	2.02	0.41
1:D:541:PHE:HD2	1:D:542:LEU:HD12	1.85	0.41
2:E:108:ARG:HG3	2:E:116:MET:HB2	2.01	0.41
2:F:24:VAL:HA	2:F:27:HIS:HB2	2.02	0.41
1:A:809:VAL:HG12	1:D:597:SER:HB3	2.03	0.41
1:D:527:MET:HE1	3:D:1101:PCW:H2	2.01	0.41
1:A:405:TYR:HE1	1:A:496:MET:HE1	1.84	0.41
1:B:401:LEU:HD23	1:B:444:ILE:HG21	2.02	0.41
1:A:807:MET:HE3	1:A:807:MET:HB2	1.90	0.41
1:B:573:ILE:HD13	1:B:573:ILE:HA	1.91	0.41
2:E:56:ASN:OD1	2:E:59:ARG:NH1	2.54	0.41
1:A:595:SER:OG	1:A:598:GLY:N	2.52	0.41
3:A:1102:PCW:H171	3:A:1102:PCW:H141	1.85	0.41
1:B:716:LYS:NZ	1:B:772:GLU:O	2.54	0.41
1:C:864:LEU:HD13	1:C:866:GLU:H	1.86	0.41
3:F:202:PCW:H381	3:F:202:PCW:H412	1.85	0.41
1:A:435:HIS:CE1	1:A:752:LYS:HE2	2.55	0.41
1:C:411:ASN:O	1:C:415:LEU:HB2	2.21	0.41
1:D:401:LEU:HD23	1:D:444:ILE:HG21	2.02	0.41
1:D:716:LYS:NZ	1:D:772:GLU:O	2.54	0.41
1:A:883:HIS:CE1	1:A:890:CYS:HB2	2.55	0.41
3:A:1103:PCW:H40	3:A:1103:PCW:H431	1.83	0.40
1:B:504:ILE:HD12	1:B:504:ILE:HA	1.93	0.40
2:E:24:VAL:HA	2:E:27:HIS:HB2	2.04	0.40
2:E:132:GLN:O	2:E:135:SER:OG	2.38	0.40
1:B:629:MET:HB3	1:B:629:MET:HE2	1.92	0.40
1:C:883:HIS:CE1	1:C:890:CYS:HB2	2.56	0.40
1:C:985:ASN:HA	1:C:988:MET:HB2	2.03	0.40
1:D:573:ILE:HD13	1:D:573:ILE:HA	1.91	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:978:VAL:HG22	1:D:789:LEU:HD21	2.02	0.40
2:E:58:GLU:HA	2:E:61:CYS:HB2	2.02	0.40
1:B:545:ARG:HH12	2:E:66:LYS:HB3	1.86	0.40
3:D:1101:PCW:H51	3:D:1102:PCW:H42	2.03	0.40
1:A:798:ILE:O	1:A:802:GLY:N	2.55	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 PDB entries followed by that with respect to all EM entries.

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	592/1026 (58%)	521 (88%)	62 (10%)	9 (2%)	8	33
1	B	404/1026 (39%)	349 (86%)	51 (13%)	4 (1%)	12	41
1	C	592/1026 (58%)	522 (88%)	61 (10%)	9 (2%)	8	33
1	D	404/1026 (39%)	349 (86%)	51 (13%)	4 (1%)	12	41
2	E	136/160 (85%)	119 (88%)	16 (12%)	1 (1%)	18	48
2	F	136/160 (85%)	118 (87%)	17 (12%)	1 (1%)	18	48
All	All	2264/4424 (51%)	1978 (87%)	258 (11%)	28 (1%)	13	38

All (28) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	412	HIS
1	A	413	GLU
1	C	412	HIS
1	C	413	GLU
1	A	874	GLN
1	C	874	GLN
2	E	120	VAL

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Mol	Chain	Res	Type
2	F	120	VAL
1	A	404	PRO
1	B	715	ARG
1	C	404	PRO
1	D	715	ARG
1	A	592	SER
1	A	684	ARG
1	B	450	TYR
1	B	636	ALA
1	C	592	SER
1	C	684	ARG
1	D	450	TYR
1	D	636	ALA
1	A	411	ASN
1	C	411	ASN
1	A	591	ILE
1	C	591	ILE
1	B	717	PRO
1	D	717	PRO
1	A	683	VAL
1	C	683	VAL

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 PDB entries followed by that with respect to all EM entries.

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	508/877 (58%)	505 (99%)	3 (1%)	78	78
1	B	343/877 (39%)	340 (99%)	3 (1%)	70	73
1	C	508/877 (58%)	505 (99%)	3 (1%)	78	78
1	D	343/877 (39%)	340 (99%)	3 (1%)	70	73
2	E	126/143 (88%)	125 (99%)	1 (1%)	73	74
2	F	126/143 (88%)	125 (99%)	1 (1%)	73	74
All	All	1954/3794 (52%)	1940 (99%)	14 (1%)	73	76

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	459	ILE
1	A	467	LEU
1	A	868	ILE
1	B	395	VAL
1	B	645	ILE
1	B	683	VAL
1	C	459	ILE
1	C	467	LEU
1	C	868	ILE
1	D	395	VAL
1	D	645	ILE
1	D	683	VAL
2	E	116	MET
2	F	116	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	747	ASN
1	A	920	ASN
1	A	990	ASN
1	B	411	ASN
1	B	435	HIS
1	B	461	ASN
1	B	791	ASN
1	C	747	ASN
1	C	920	ASN
1	C	990	ASN
1	D	411	ASN
1	D	435	HIS
1	D	461	ASN
1	D	756	GLN
1	D	791	ASN
2	E	75	HIS
2	F	75	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

33 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
3	PCW	B	1105	-	50,50,53	1.07	3 (6%)	56,58,61	0.94	4 (7%)
3	PCW	C	1101	-	50,50,53	1.09	3 (6%)	56,58,61	0.90	3 (5%)
4	GLU	C	1107	-	8,9,9	1.05	1 (12%)	8,11,11	1.45	1 (12%)
3	PCW	A	1107	-	10,10,53	0.76	0	9,9,61	0.38	0
3	PCW	E	203	-	42,42,53	1.18	3 (7%)	48,50,61	0.96	3 (6%)
5	SPD	B	1106	-	9,9,9	0.35	0	8,8,8	0.59	0
3	PCW	A	1102	-	50,50,53	1.09	3 (6%)	56,58,61	0.91	3 (5%)
3	PCW	B	1102	-	42,42,53	1.14	3 (7%)	48,50,61	1.04	3 (6%)
3	PCW	D	1101	-	40,40,53	1.17	3 (7%)	46,48,61	0.94	2 (4%)
3	PCW	F	201	-	10,10,53	0.75	0	9,9,61	0.33	0
3	PCW	A	1104	-	50,50,53	1.09	3 (6%)	56,58,61	0.89	3 (5%)
3	PCW	B	1103	-	40,40,53	1.18	3 (7%)	46,48,61	1.09	3 (6%)
3	PCW	C	1106	-	10,10,53	0.75	0	9,9,61	0.38	0
3	PCW	F	202	-	42,42,53	1.17	3 (7%)	48,50,61	0.96	3 (6%)
3	PCW	A	1101	-	50,50,53	1.08	3 (6%)	56,58,61	0.91	2 (3%)
3	PCW	C	1103	-	50,50,53	1.09	3 (6%)	56,58,61	0.89	3 (5%)
3	PCW	A	1109	-	10,10,53	0.76	0	9,9,61	0.35	0
3	PCW	A	1103	-	42,42,53	1.16	3 (7%)	48,50,61	1.05	3 (6%)
3	PCW	C	1105	-	10,10,53	0.75	0	9,9,61	0.32	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	PCW	D	1103	-	38,38,53	1.18	3 (7%)	44,46,61	1.17	4 (9%)
4	GLU	B	1104	-	8,9,9	1.11	1 (12%)	8,11,11	1.29	1 (12%)
3	PCW	E	202	-	10,10,53	0.75	0	9,9,61	0.34	0
3	PCW	A	1105	-	10,10,53	0.76	0	9,9,61	0.31	0
3	PCW	D	1102	-	42,42,53	1.15	3 (7%)	48,50,61	1.07	3 (6%)
3	PCW	C	1104	-	10,10,53	0.75	0	9,9,61	0.33	0
3	PCW	C	1108	-	10,10,53	0.77	0	9,9,61	0.38	0
3	PCW	D	1104	-	40,40,53	1.18	3 (7%)	46,48,61	1.09	3 (6%)
3	PCW	B	1101	-	40,40,53	1.16	3 (7%)	46,48,61	0.94	3 (6%)
3	PCW	A	1106	-	10,10,53	0.75	0	9,9,61	0.31	0
3	PCW	E	201	-	38,38,53	1.19	3 (7%)	44,46,61	1.18	3 (6%)
4	GLU	A	1108	-	8,9,9	1.12	1 (12%)	8,11,11	1.45	1 (12%)
3	PCW	C	1102	-	42,42,53	1.16	3 (7%)	48,50,61	1.10	3 (6%)
4	GLU	D	1105	-	8,9,9	1.09	1 (12%)	8,11,11	1.38	1 (12%)

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	PCW	B	1105	-	-	24/54/54/57	-
3	PCW	C	1101	-	-	29/54/54/57	-
4	GLU	C	1107	-	-	0/9/9/9	-
3	PCW	A	1107	-	-	1/8/8/57	-
3	PCW	E	203	-	-	23/46/46/57	-
5	SPD	B	1106	-	-	1/7/7/7	-
3	PCW	A	1102	-	-	24/54/54/57	-
3	PCW	B	1102	-	-	26/46/46/57	-
3	PCW	D	1101	-	-	22/44/44/57	-
3	PCW	F	201	-	-	4/8/8/57	-
3	PCW	A	1104	-	-	27/54/54/57	-
3	PCW	B	1103	-	-	16/44/44/57	-
3	PCW	C	1106	-	-	1/8/8/57	-
3	PCW	F	202	-	-	20/46/46/57	-
3	PCW	A	1101	-	-	30/54/54/57	-
3	PCW	C	1103	-	-	23/54/54/57	-
3	PCW	A	1109	-	-	2/8/8/57	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	PCW	A	1103	-	-	23/46/46/57	-
3	PCW	C	1105	-	-	5/8/8/57	-
3	PCW	D	1103	-	-	20/42/42/57	-
4	GLU	B	1104	-	-	1/9/9/9	-
3	PCW	E	202	-	-	4/8/8/57	-
3	PCW	A	1105	-	-	3/8/8/57	-
3	PCW	D	1102	-	-	22/46/46/57	-
3	PCW	C	1104	-	-	3/8/8/57	-
3	PCW	C	1108	-	-	1/8/8/57	-
3	PCW	D	1104	-	-	15/44/44/57	-
3	PCW	B	1101	-	-	20/44/44/57	-
3	PCW	A	1106	-	-	5/8/8/57	-
3	PCW	E	201	-	-	19/42/42/57	-
4	GLU	A	1108	-	-	1/9/9/9	-
3	PCW	C	1102	-	-	27/46/46/57	-
4	GLU	D	1105	-	-	1/9/9/9	-

All (58) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	1102	PCW	O2-C31	3.03	1.42	1.34
3	A	1103	PCW	O2-C31	3.00	1.42	1.34
3	C	1101	PCW	O3-C11	2.98	1.42	1.33
3	A	1102	PCW	O3-C11	2.98	1.42	1.33
3	C	1101	PCW	O2-C31	2.97	1.42	1.34
3	A	1102	PCW	O2-C31	2.96	1.42	1.34
3	A	1104	PCW	O3-C11	2.93	1.41	1.33
3	C	1103	PCW	O2-C31	2.91	1.42	1.34
3	A	1104	PCW	O2-C31	2.91	1.42	1.34
3	C	1103	PCW	O3-C11	2.90	1.41	1.33
3	E	203	PCW	O3-C11	2.89	1.41	1.33
3	D	1104	PCW	O3-C11	2.88	1.41	1.33
3	F	202	PCW	O3-C11	2.87	1.41	1.33
3	E	201	PCW	O3-C11	2.84	1.41	1.33
3	B	1105	PCW	O2-C31	2.84	1.42	1.34
3	B	1103	PCW	O3-C11	2.83	1.41	1.33
3	E	201	PCW	O2-C31	2.81	1.42	1.34
3	A	1101	PCW	O3-C11	2.81	1.41	1.33
3	D	1101	PCW	O2-C31	2.80	1.42	1.34
3	D	1103	PCW	O3-C11	2.79	1.41	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1101	PCW	O2-C31	2.78	1.42	1.34
3	B	1101	PCW	O3-C11	2.76	1.41	1.33
3	B	1101	PCW	O2-C31	2.76	1.42	1.34
3	B	1105	PCW	O3-C11	2.76	1.41	1.33
3	C	1102	PCW	O3-C11	2.73	1.41	1.33
3	D	1103	PCW	O2-C31	2.72	1.42	1.34
3	A	1103	PCW	O3-C11	2.72	1.41	1.33
3	B	1102	PCW	O2-C31	2.71	1.41	1.34
3	D	1104	PCW	O2-C31	2.71	1.41	1.34
3	D	1102	PCW	O2-C31	2.71	1.41	1.34
3	B	1103	PCW	O2-C31	2.70	1.41	1.34
3	D	1101	PCW	O3-C11	2.68	1.41	1.33
3	E	203	PCW	O2-C2	-2.68	1.40	1.46
3	D	1102	PCW	O3-C11	2.68	1.41	1.33
3	F	202	PCW	O2-C31	2.67	1.41	1.34
3	E	203	PCW	O2-C31	2.66	1.41	1.34
3	D	1102	PCW	O2-C2	-2.65	1.40	1.46
3	F	202	PCW	O2-C2	-2.63	1.40	1.46
3	B	1102	PCW	O3-C11	2.63	1.41	1.33
3	D	1101	PCW	O2-C2	-2.61	1.40	1.46
3	B	1102	PCW	O2-C2	-2.61	1.40	1.46
3	A	1101	PCW	O2-C2	-2.60	1.40	1.46
3	B	1101	PCW	O2-C2	-2.59	1.40	1.46
3	B	1103	PCW	O2-C2	-2.57	1.40	1.46
3	D	1104	PCW	O2-C2	-2.55	1.40	1.46
3	B	1105	PCW	O2-C2	-2.52	1.40	1.46
4	A	1108	GLU	OXT-C	-2.43	1.22	1.30
4	B	1104	GLU	OXT-C	-2.28	1.23	1.30
4	D	1105	GLU	OXT-C	-2.28	1.23	1.30
4	C	1107	GLU	OXT-C	-2.19	1.23	1.30
3	D	1103	PCW	O2-C2	-2.18	1.41	1.46
3	C	1101	PCW	O2-C2	-2.14	1.41	1.46
3	A	1102	PCW	O2-C2	-2.13	1.41	1.46
3	E	201	PCW	O2-C2	-2.11	1.41	1.46
3	A	1103	PCW	O2-C2	-2.10	1.41	1.46
3	A	1104	PCW	O2-C2	-2.09	1.41	1.46
3	C	1103	PCW	O2-C2	-2.05	1.41	1.46
3	C	1102	PCW	O2-C2	-2.04	1.41	1.46

All (58) bond angle outliers are listed below:

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	201	PCW	O2-C31-C32	4.84	121.96	111.48
3	C	1102	PCW	O2-C31-C32	4.69	121.63	111.48
3	D	1103	PCW	O2-C31-C32	4.59	121.40	111.48
3	A	1103	PCW	O2-C31-C32	4.26	120.69	111.48
3	D	1102	PCW	O2-C31-C32	4.19	120.54	111.48
3	D	1104	PCW	O2-C31-C32	4.11	120.38	111.48
3	B	1103	PCW	O2-C31-C32	4.06	120.26	111.48
3	B	1102	PCW	O2-C31-C32	4.01	120.17	111.48
3	B	1105	PCW	O2-C31-C32	3.82	119.74	111.48
3	C	1103	PCW	O2-C31-C32	3.73	119.56	111.48
3	E	203	PCW	O2-C31-C32	3.71	119.50	111.48
3	A	1104	PCW	O2-C31-C32	3.70	119.48	111.48
3	F	202	PCW	O2-C31-C32	3.69	119.47	111.48
3	A	1101	PCW	O2-C31-C32	3.68	119.45	111.48
3	A	1102	PCW	O2-C31-C32	3.60	119.28	111.48
3	D	1101	PCW	O2-C31-C32	3.60	119.26	111.48
3	C	1101	PCW	O2-C31-C32	3.59	119.25	111.48
3	B	1101	PCW	O2-C31-C32	3.49	119.03	111.48
4	A	1108	GLU	OXT-C-O	-3.47	116.21	124.08
4	C	1107	GLU	OXT-C-O	-3.46	116.22	124.08
4	D	1105	GLU	OXT-C-O	-3.07	117.11	124.08
4	B	1104	GLU	OXT-C-O	-3.06	117.14	124.08
3	A	1102	PCW	O3-C11-C12	2.87	120.57	111.83
3	C	1101	PCW	O3-C11-C12	2.86	120.55	111.83
3	D	1103	PCW	O3-C11-C12	2.60	119.76	111.83
3	F	202	PCW	O3-C11-C12	2.59	119.73	111.83
3	B	1103	PCW	O3-C11-C12	2.58	119.70	111.83
3	E	203	PCW	O3-C11-C12	2.56	119.63	111.83
3	D	1104	PCW	O3-C11-C12	2.55	119.62	111.83
3	A	1103	PCW	O3-C11-C12	2.55	119.60	111.83
3	E	201	PCW	O3-C11-C12	2.55	119.60	111.83
3	A	1101	PCW	O3-C11-C12	2.50	119.47	111.83
3	C	1102	PCW	O3-C11-C12	2.50	119.46	111.83
3	C	1103	PCW	O3-C11-C12	2.46	119.34	111.83
3	A	1104	PCW	O3-C11-C12	2.45	119.32	111.83
3	D	1102	PCW	O3-C11-C12	2.42	119.21	111.83
3	D	1101	PCW	C6-N-C5	2.40	119.44	109.91
3	B	1101	PCW	C6-N-C5	2.37	119.34	109.91
3	B	1102	PCW	O3-C11-C12	2.35	119.00	111.83
3	A	1103	PCW	C6-N-C5	2.33	119.19	109.91
3	D	1102	PCW	C6-N-C5	2.32	119.13	109.91

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1104	PCW	C6-N-C5	2.30	119.05	109.91
3	B	1105	PCW	C6-N-C5	2.28	118.97	109.91
3	B	1105	PCW	O3-C11-C12	2.27	118.77	111.83
3	B	1102	PCW	C6-N-C5	2.27	118.95	109.91
3	C	1102	PCW	C6-N-C5	2.27	118.92	109.91
3	C	1103	PCW	C6-N-C5	2.21	118.68	109.91
3	C	1101	PCW	C6-N-C5	2.20	118.67	109.91
3	A	1102	PCW	C6-N-C5	2.18	118.56	109.91
3	D	1103	PCW	C6-N-C5	2.17	118.54	109.91
3	D	1103	PCW	C2-O2-C31	2.15	122.95	117.80
3	E	201	PCW	C6-N-C5	2.15	118.45	109.91
3	E	203	PCW	C6-N-C5	2.11	118.28	109.91
3	F	202	PCW	C6-N-C5	2.11	118.28	109.91
3	B	1105	PCW	C3-C2-C1	-2.09	106.91	111.78
3	B	1103	PCW	C6-N-C5	2.09	118.21	109.91
3	D	1104	PCW	C6-N-C5	2.02	117.94	109.91
3	B	1101	PCW	O3-C11-C12	2.00	117.94	111.83

There are no chirality outliers.

All (443) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1101	PCW	C4-O4P-P-O1P
3	A	1101	PCW	C4-O4P-P-O2P
3	A	1101	PCW	C4-O4P-P-O3P
3	A	1102	PCW	C12-C11-O3-C3
3	A	1102	PCW	O11-C11-O3-C3
3	A	1102	PCW	O31-C31-O2-C2
3	A	1102	PCW	C4-O4P-P-O1P
3	A	1102	PCW	C4-O4P-P-O2P
3	A	1102	PCW	C4-O4P-P-O3P
3	A	1103	PCW	O3P-C1-C2-O2
3	A	1103	PCW	C1-O3P-P-O2P
3	A	1103	PCW	C4-O4P-P-O3P
3	A	1104	PCW	O31-C31-O2-C2
3	A	1105	PCW	C20-C21-C22-C23
3	B	1101	PCW	C1-O3P-P-O2P
3	B	1101	PCW	C1-O3P-P-O4P
3	B	1101	PCW	C4-O4P-P-O2P
3	B	1101	PCW	C4-O4P-P-O3P
3	B	1102	PCW	C32-C31-O2-C2
3	B	1103	PCW	O31-C31-O2-C2

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Mol	Chain	Res	Type	Atoms
3	B	1105	PCW	O4P-C4-C5-N
3	B	1105	PCW	C1-O3P-P-O1P
3	B	1105	PCW	C1-O3P-P-O2P
3	B	1105	PCW	C1-O3P-P-O4P
3	C	1101	PCW	C12-C11-O3-C3
3	C	1101	PCW	O11-C11-O3-C3
3	C	1101	PCW	C4-O4P-P-O1P
3	C	1101	PCW	C4-O4P-P-O2P
3	C	1101	PCW	C4-O4P-P-O3P
3	C	1102	PCW	O3P-C1-C2-O2
3	C	1102	PCW	C1-O3P-P-O2P
3	C	1102	PCW	C1-O3P-P-O4P
3	C	1102	PCW	C4-O4P-P-O3P
3	C	1103	PCW	O4P-C4-C5-N
3	C	1104	PCW	C20-C21-C22-C23
3	C	1108	PCW	C20-C21-C22-C23
3	D	1101	PCW	C4-O4P-P-O1P
3	D	1101	PCW	C4-O4P-P-O2P
3	D	1101	PCW	C4-O4P-P-O3P
3	D	1102	PCW	C32-C31-O2-C2
3	D	1103	PCW	C32-C31-O2-C2
3	D	1103	PCW	C4-O4P-P-O1P
3	D	1103	PCW	C4-O4P-P-O2P
3	D	1103	PCW	C4-O4P-P-O3P
3	D	1104	PCW	O31-C31-O2-C2
3	E	201	PCW	C32-C31-O2-C2
3	E	201	PCW	C4-O4P-P-O1P
3	E	201	PCW	C4-O4P-P-O2P
3	E	201	PCW	C4-O4P-P-O3P
3	E	202	PCW	C20-C21-C22-C23
3	E	203	PCW	O4P-C4-C5-N
3	E	203	PCW	O31-C31-O2-C2
3	E	203	PCW	C4-O4P-P-O1P
3	E	203	PCW	C4-O4P-P-O3P
3	F	201	PCW	C20-C21-C22-C23
3	F	202	PCW	O4P-C4-C5-N
3	F	202	PCW	O31-C31-O2-C2
3	F	202	PCW	C4-O4P-P-O1P
3	F	202	PCW	C4-O4P-P-O3P
3	C	1103	PCW	O11-C11-O3-C3
3	E	203	PCW	O11-C11-O3-C3
3	F	202	PCW	O11-C11-O3-C3

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Mol	Chain	Res	Type	Atoms
3	C	1103	PCW	C12-C11-O3-C3
3	E	203	PCW	C12-C11-O3-C3
3	F	202	PCW	C12-C11-O3-C3
3	A	1104	PCW	O11-C11-O3-C3
3	B	1102	PCW	O31-C31-O2-C2
3	C	1101	PCW	O31-C31-O2-C2
3	C	1102	PCW	O31-C31-O2-C2
3	C	1103	PCW	O31-C31-O2-C2
3	D	1102	PCW	O31-C31-O2-C2
3	D	1103	PCW	O31-C31-O2-C2
3	E	201	PCW	O31-C31-O2-C2
3	A	1104	PCW	C12-C11-O3-C3
3	A	1102	PCW	C32-C31-O2-C2
3	A	1104	PCW	C32-C31-O2-C2
3	B	1103	PCW	C32-C31-O2-C2
3	C	1101	PCW	C32-C31-O2-C2
3	D	1104	PCW	C32-C31-O2-C2
3	E	203	PCW	C32-C31-O2-C2
3	F	202	PCW	C32-C31-O2-C2
3	A	1103	PCW	C12-C11-O3-C3
3	C	1102	PCW	C12-C11-O3-C3
3	D	1103	PCW	C12-C11-O3-C3
3	A	1103	PCW	C32-C31-O2-C2
3	B	1101	PCW	C32-C31-O2-C2
3	C	1102	PCW	C32-C31-O2-C2
3	C	1103	PCW	C32-C31-O2-C2
3	D	1101	PCW	C32-C31-O2-C2
3	A	1103	PCW	O11-C11-O3-C3
3	C	1102	PCW	O11-C11-O3-C3
3	D	1103	PCW	O11-C11-O3-C3
3	A	1103	PCW	O31-C31-O2-C2
3	D	1102	PCW	C14-C15-C16-C17
3	D	1101	PCW	C13-C14-C15-C16
3	D	1101	PCW	O31-C31-O2-C2
3	D	1101	PCW	C12-C11-O3-C3
3	B	1102	PCW	C31-C32-C33-C34
3	D	1102	PCW	C31-C32-C33-C34
3	D	1103	PCW	C32-C33-C34-C35
3	E	201	PCW	C15-C16-C17-C18
3	D	1101	PCW	O11-C11-O3-C3
3	D	1102	PCW	C12-C11-O3-C3
3	B	1105	PCW	C11-C12-C13-C14

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Mol	Chain	Res	Type	Atoms
3	B	1101	PCW	O31-C31-O2-C2
3	B	1101	PCW	C13-C14-C15-C16
3	D	1104	PCW	C4-C5-N-C8
3	B	1102	PCW	C11-C12-C13-C14
3	C	1102	PCW	C31-C32-C33-C34
3	D	1101	PCW	C31-C32-C33-C34
3	B	1102	PCW	C12-C11-O3-C3
3	E	201	PCW	C12-C11-O3-C3
3	E	201	PCW	C32-C33-C34-C35
3	A	1101	PCW	C4-C5-N-C6
3	A	1101	PCW	C4-C5-N-C7
3	A	1101	PCW	C4-C5-N-C8
3	B	1103	PCW	C4-C5-N-C7
3	B	1103	PCW	C4-C5-N-C8
3	D	1102	PCW	C16-C17-C18-C19
3	C	1106	PCW	C20-C21-C22-C23
3	D	1102	PCW	O11-C11-O3-C3
3	C	1103	PCW	C13-C14-C15-C16
3	B	1103	PCW	C4-C5-N-C6
3	D	1104	PCW	C4-C5-N-C7
3	D	1102	PCW	C32-C33-C34-C35
3	E	202	PCW	C15-C16-C17-C18
3	A	1101	PCW	C12-C13-C14-C15
3	E	202	PCW	C14-C15-C16-C17
3	F	201	PCW	C14-C15-C16-C17
3	A	1104	PCW	C14-C15-C16-C17
3	E	203	PCW	C13-C14-C15-C16
3	C	1101	PCW	C32-C33-C34-C35
3	B	1102	PCW	C16-C17-C18-C19
3	A	1105	PCW	C15-C16-C17-C18
3	B	1101	PCW	C12-C13-C14-C15
3	C	1104	PCW	C15-C16-C17-C18
3	A	1104	PCW	C31-C32-C33-C34
3	B	1102	PCW	C13-C14-C15-C16
3	B	1102	PCW	C32-C33-C34-C35
3	D	1101	PCW	C12-C13-C14-C15
3	E	201	PCW	O11-C11-O3-C3
3	B	1101	PCW	C23-C24-C25-C26
3	D	1103	PCW	C34-C35-C36-C37
3	B	1102	PCW	O11-C11-O3-C3
3	E	203	PCW	C35-C36-C37-C38
3	B	1105	PCW	C31-C32-C33-C34

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Mol	Chain	Res	Type	Atoms
3	A	1103	PCW	C13-C14-C15-C16
3	B	1102	PCW	C35-C36-C37-C38
3	D	1102	PCW	C13-C14-C15-C16
3	F	201	PCW	C15-C16-C17-C18
3	D	1101	PCW	C23-C24-C25-C26
3	B	1105	PCW	C21-C22-C23-C24
3	A	1101	PCW	C31-C32-C33-C34
3	B	1105	PCW	C12-C13-C14-C15
3	F	202	PCW	C13-C14-C15-C16
3	F	202	PCW	C35-C36-C37-C38
3	C	1103	PCW	C33-C34-C35-C36
3	C	1105	PCW	C14-C15-C16-C17
3	A	1106	PCW	C14-C15-C16-C17
3	A	1102	PCW	C41-C42-C43-C44
3	A	1104	PCW	C15-C16-C17-C18
3	B	1102	PCW	C40-C41-C42-C43
3	C	1102	PCW	C36-C37-C38-C39
3	C	1103	PCW	C16-C17-C18-C19
3	D	1104	PCW	C4-C5-N-C6
3	A	1101	PCW	C35-C36-C37-C38
3	B	1105	PCW	C35-C36-C37-C38
3	B	1101	PCW	C31-C32-C33-C34
3	E	203	PCW	O3P-C1-C2-O2
3	C	1103	PCW	C11-C12-C13-C14
3	A	1104	PCW	C16-C17-C18-C19
3	B	1105	PCW	C16-C17-C18-C19
3	B	1105	PCW	C14-C15-C16-C17
3	D	1104	PCW	C33-C34-C35-C36
3	A	1102	PCW	C32-C33-C34-C35
3	A	1102	PCW	C2-C1-O3P-P
3	A	1102	PCW	C21-C22-C23-C24
3	A	1104	PCW	C33-C34-C35-C36
3	A	1104	PCW	C24-C25-C26-C27
3	C	1103	PCW	C24-C25-C26-C27
3	C	1102	PCW	C41-C42-C43-C44
3	D	1103	PCW	C12-C13-C14-C15
3	A	1102	PCW	C36-C37-C38-C39
3	B	1102	PCW	C36-C37-C38-C39
3	B	1105	PCW	C40-C41-C42-C43
3	A	1103	PCW	O3P-C1-C2-C3
3	A	1104	PCW	O3P-C1-C2-C3
3	C	1102	PCW	O3P-C1-C2-C3

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Mol	Chain	Res	Type	Atoms
3	A	1101	PCW	C14-C15-C16-C17
3	A	1102	PCW	C14-C15-C16-C17
3	B	1101	PCW	C12-C11-O3-C3
3	A	1104	PCW	C1-C2-C3-O3
3	B	1103	PCW	C1-C2-C3-O3
3	C	1101	PCW	C1-C2-C3-O3
3	C	1103	PCW	C1-C2-C3-O3
3	D	1102	PCW	C1-C2-C3-O3
3	D	1104	PCW	C1-C2-C3-O3
3	A	1101	PCW	C16-C17-C18-C19
3	A	1102	PCW	C20-C21-C22-C23
3	A	1103	PCW	C16-C17-C18-C19
3	A	1103	PCW	C36-C37-C38-C39
3	C	1101	PCW	C36-C37-C38-C39
3	C	1105	PCW	C16-C17-C18-C19
3	D	1102	PCW	C36-C37-C38-C39
3	E	203	PCW	C16-C17-C18-C19
3	F	202	PCW	C16-C17-C18-C19
3	A	1101	PCW	C21-C22-C23-C24
3	B	1103	PCW	C33-C34-C35-C36
3	F	202	PCW	C32-C33-C34-C35
3	A	1106	PCW	C20-C21-C22-C23
3	A	1109	PCW	C20-C21-C22-C23
3	C	1105	PCW	C20-C21-C22-C23
3	B	1101	PCW	O11-C11-O3-C3
3	C	1101	PCW	C21-C22-C23-C24
3	A	1102	PCW	C1-C2-O2-C31
3	A	1104	PCW	C3-C2-O2-C31
3	C	1101	PCW	C1-C2-O2-C31
3	D	1103	PCW	C1-C2-O2-C31
3	A	1104	PCW	C41-C42-C43-C44
3	C	1101	PCW	C20-C21-C22-C23
3	D	1102	PCW	C40-C41-C42-C43
3	C	1103	PCW	C12-C13-C14-C15
3	D	1102	PCW	C35-C36-C37-C38
3	A	1104	PCW	C11-C12-C13-C14
3	A	1101	PCW	C13-C14-C15-C16
3	B	1101	PCW	O3P-C1-C2-O2
3	F	202	PCW	O3P-C1-C2-O2
3	C	1101	PCW	C14-C15-C16-C17
3	C	1101	PCW	C41-C42-C43-C44
3	A	1101	PCW	C32-C33-C34-C35

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Mol	Chain	Res	Type	Atoms
3	C	1102	PCW	C14-C15-C16-C17
3	B	1102	PCW	C39-C40-C41-C42
3	A	1101	PCW	C20-C21-C22-C23
3	A	1104	PCW	C20-C21-C22-C23
3	A	1106	PCW	C16-C17-C18-C19
3	C	1102	PCW	C42-C43-C44-C45
3	D	1101	PCW	C32-C33-C34-C35
3	D	1101	PCW	C14-C15-C16-C17
3	C	1101	PCW	C2-C1-O3P-P
3	B	1105	PCW	C13-C14-C15-C16
3	E	203	PCW	C15-C16-C17-C18
3	C	1104	PCW	C14-C15-C16-C17
3	D	1102	PCW	C41-C42-C43-C44
3	B	1103	PCW	C11-C12-C13-C14
3	B	1101	PCW	O3P-C1-C2-C3
3	E	203	PCW	O3P-C1-C2-C3
3	F	202	PCW	O3P-C1-C2-C3
3	A	1101	PCW	C12-C11-O3-C3
3	A	1102	PCW	C34-C35-C36-C37
3	A	1104	PCW	C34-C35-C36-C37
3	D	1104	PCW	C37-C38-C39-C40
3	B	1102	PCW	C42-C43-C44-C45
3	C	1103	PCW	C34-C35-C36-C37
3	D	1104	PCW	C35-C36-C37-C38
3	D	1101	PCW	C11-C12-C13-C14
3	E	203	PCW	C32-C33-C34-C35
3	B	1102	PCW	C1-C2-C3-O3
3	B	1103	PCW	C37-C38-C39-C40
3	C	1101	PCW	C22-C23-C24-C25
3	F	202	PCW	C15-C16-C17-C18
3	B	1105	PCW	C32-C33-C34-C35
3	B	1101	PCW	C32-C33-C34-C35
3	A	1102	PCW	O2-C2-C3-O3
3	B	1103	PCW	O2-C2-C3-O3
3	C	1103	PCW	O2-C2-C3-O3
3	D	1101	PCW	O2-C2-C3-O3
3	D	1104	PCW	O2-C2-C3-O3
3	E	203	PCW	O2-C2-C3-O3
3	F	202	PCW	O2-C2-C3-O3
3	A	1104	PCW	C22-C23-C24-C25
3	A	1102	PCW	C12-C13-C14-C15
3	B	1105	PCW	C33-C34-C35-C36

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Mol	Chain	Res	Type	Atoms
3	F	202	PCW	C12-C13-C14-C15
3	D	1103	PCW	C35-C36-C37-C38
3	F	201	PCW	C13-C14-C15-C16
3	D	1103	PCW	C39-C40-C41-C42
3	A	1103	PCW	C32-C33-C34-C35
3	A	1107	PCW	C20-C21-C22-C23
3	C	1103	PCW	C32-C33-C34-C35
3	A	1101	PCW	C11-C12-C13-C14
3	A	1101	PCW	O11-C11-O3-C3
3	E	202	PCW	C13-C14-C15-C16
3	C	1103	PCW	C22-C23-C24-C25
3	D	1101	PCW	C24-C25-C26-C27
3	A	1103	PCW	C3-C2-O2-C31
3	C	1102	PCW	C3-C2-O2-C31
3	C	1103	PCW	C3-C2-O2-C31
3	E	201	PCW	C1-C2-O2-C31
3	A	1101	PCW	C40-C41-C42-C43
3	C	1102	PCW	C33-C34-C35-C36
3	B	1102	PCW	C12-C13-C14-C15
3	C	1103	PCW	O3P-C1-C2-O2
3	D	1102	PCW	O3P-C1-C2-O2
3	E	201	PCW	C12-C13-C14-C15
3	A	1102	PCW	C1-C2-C3-O3
3	E	201	PCW	C1-C2-C3-O3
3	A	1103	PCW	C5-C4-O4P-P
3	B	1103	PCW	C5-C4-O4P-P
3	C	1102	PCW	C5-C4-O4P-P
3	D	1103	PCW	C5-C4-O4P-P
3	D	1104	PCW	C5-C4-O4P-P
3	E	201	PCW	C5-C4-O4P-P
3	B	1102	PCW	C34-C35-C36-C37
3	C	1101	PCW	O2-C2-C3-O3
3	D	1102	PCW	O2-C2-C3-O3
3	E	201	PCW	O2-C2-C3-O3
3	C	1101	PCW	C35-C36-C37-C38
3	D	1102	PCW	C34-C35-C36-C37
3	C	1101	PCW	C12-C13-C14-C15
3	A	1101	PCW	O4P-C4-C5-N
3	A	1102	PCW	O4P-C4-C5-N
3	A	1103	PCW	O4P-C4-C5-N
3	B	1101	PCW	O4P-C4-C5-N
3	B	1102	PCW	O4P-C4-C5-N

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Mol	Chain	Res	Type	Atoms
3	B	1103	PCW	O4P-C4-C5-N
3	C	1101	PCW	O4P-C4-C5-N
3	C	1102	PCW	O4P-C4-C5-N
3	D	1101	PCW	O4P-C4-C5-N
3	D	1102	PCW	O4P-C4-C5-N
3	D	1103	PCW	O4P-C4-C5-N
3	E	201	PCW	O4P-C4-C5-N
3	B	1103	PCW	C15-C16-C17-C18
3	C	1101	PCW	C11-C12-C13-C14
3	C	1103	PCW	C35-C36-C37-C38
3	D	1102	PCW	C11-C12-C13-C14
3	C	1103	PCW	O3P-C1-C2-C3
3	D	1103	PCW	C14-C15-C16-C17
3	E	203	PCW	C12-C13-C14-C15
3	A	1101	PCW	C2-C1-O3P-P
3	A	1104	PCW	O3P-C1-C2-O2
3	B	1102	PCW	O3P-C1-C2-O2
3	C	1102	PCW	C16-C17-C18-C19
3	A	1103	PCW	C12-C13-C14-C15
3	A	1104	PCW	O2-C2-C3-O3
3	B	1102	PCW	O2-C2-C3-O3
3	D	1103	PCW	O2-C2-C3-O3
3	A	1102	PCW	C35-C36-C37-C38
3	A	1101	PCW	C33-C34-C35-C36
3	C	1102	PCW	C1-C2-C3-O3
3	D	1103	PCW	C1-C2-C3-O3
3	A	1104	PCW	C32-C33-C34-C35
3	A	1103	PCW	C4-O4P-P-O2P
3	A	1104	PCW	C1-O3P-P-O2P
3	A	1104	PCW	C1-O3P-P-O4P
3	A	1104	PCW	C4-O4P-P-O2P
3	B	1101	PCW	C4-O4P-P-O1P
3	B	1102	PCW	C1-O3P-P-O2P
3	C	1101	PCW	C1-O3P-P-O2P
3	C	1102	PCW	C4-O4P-P-O2P
3	D	1101	PCW	C1-O3P-P-O2P
3	E	203	PCW	C1-O3P-P-O1P
3	E	203	PCW	C1-O3P-P-O4P
3	A	1109	PCW	C15-C16-C17-C18
3	A	1104	PCW	C40-C41-C42-C43
3	A	1101	PCW	C25-C26-C27-C28
3	D	1101	PCW	C19-C20-C21-C22

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Mol	Chain	Res	Type	Atoms
3	B	1103	PCW	C32-C33-C34-C35
3	D	1103	PCW	C15-C16-C17-C18
3	C	1102	PCW	C34-C35-C36-C37
3	E	201	PCW	C35-C36-C37-C38
3	D	1102	PCW	O3P-C1-C2-C3
3	D	1104	PCW	C12-C13-C14-C15
3	B	1103	PCW	C35-C36-C37-C38
3	D	1102	PCW	C2-C1-O3P-P
3	B	1105	PCW	C32-C31-O2-C2
3	B	1105	PCW	O31-C31-O2-C2
3	A	1103	PCW	C34-C35-C36-C37
3	A	1104	PCW	C13-C14-C15-C16
3	D	1104	PCW	C11-C12-C13-C14
3	B	1102	PCW	C17-C18-C19-C20
3	A	1101	PCW	C15-C16-C17-C18
3	C	1101	PCW	C33-C34-C35-C36
3	C	1101	PCW	C34-C35-C36-C37
3	B	1102	PCW	O3P-C1-C2-C3
3	D	1104	PCW	C15-C16-C17-C18
3	B	1101	PCW	C24-C25-C26-C27
3	A	1101	PCW	C34-C35-C36-C37
3	A	1103	PCW	C41-C42-C43-C44
3	C	1101	PCW	C39-C40-C41-C42
3	D	1102	PCW	C39-C40-C41-C42
3	E	203	PCW	C31-C32-C33-C34
3	B	1103	PCW	C12-C13-C14-C15
3	A	1101	PCW	C37-C38-C39-C40
3	B	1105	PCW	C37-C38-C39-C40
3	B	1105	PCW	C25-C26-C27-C28
3	D	1104	PCW	C32-C33-C34-C35
3	C	1103	PCW	C31-C32-C33-C34
3	C	1103	PCW	C37-C38-C39-C40
3	A	1101	PCW	O2-C2-C3-O3
3	A	1103	PCW	C35-C36-C37-C38
3	A	1103	PCW	C14-C15-C16-C17
3	A	1102	PCW	C11-C12-C13-C14
3	B	1102	PCW	C14-C15-C16-C17
5	B	1106	SPD	C2-C3-C4-C5
3	A	1103	PCW	C1-C2-C3-O3
3	D	1101	PCW	O3P-C1-C2-O2
3	C	1103	PCW	C39-C40-C41-C42
3	A	1106	PCW	C15-C16-C17-C18

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Mol	Chain	Res	Type	Atoms
3	D	1101	PCW	O3P-C1-C2-C3
3	B	1105	PCW	C22-C23-C24-C25
3	C	1102	PCW	C40-C41-C42-C43
3	A	1103	PCW	C37-C38-C39-C40
3	C	1102	PCW	C37-C38-C39-C40
3	D	1101	PCW	C21-C22-C23-C24
3	A	1102	PCW	C39-C40-C41-C42
3	A	1104	PCW	C37-C38-C39-C40
3	E	203	PCW	C1-C2-C3-O3
3	B	1102	PCW	C15-C16-C17-C18
3	D	1103	PCW	C37-C38-C39-C40
3	E	201	PCW	C37-C38-C39-C40
3	A	1101	PCW	C5-C4-O4P-P
3	A	1105	PCW	C14-C15-C16-C17
3	C	1105	PCW	C15-C16-C17-C18
3	B	1101	PCW	C14-C15-C16-C17
3	B	1101	PCW	C19-C20-C21-C22
3	C	1105	PCW	C17-C18-C19-C20
3	F	202	PCW	O2-C31-C32-C33
3	B	1105	PCW	C23-C24-C25-C26
3	B	1102	PCW	C41-C42-C43-C44
3	B	1105	PCW	C34-C35-C36-C37
3	A	1106	PCW	C17-C18-C19-C20
3	E	203	PCW	O2-C31-C32-C33
3	B	1105	PCW	C15-C16-C17-C18
3	F	202	PCW	C1-C2-C3-O3
3	C	1102	PCW	C11-C12-C13-C14
3	C	1101	PCW	O3-C11-C12-C13
3	C	1102	PCW	O3-C11-C12-C13
3	A	1101	PCW	C23-C24-C25-C26
3	E	201	PCW	C39-C40-C41-C42
3	A	1101	PCW	C36-C37-C38-C39
3	E	201	PCW	C34-C35-C36-C37
3	F	202	PCW	O31-C31-C32-C33
4	B	1104	GLU	OXT-C-CA-N
4	D	1105	GLU	OXT-C-CA-N
3	A	1102	PCW	O3-C11-C12-C13
3	B	1105	PCW	C20-C21-C22-C23
4	A	1108	GLU	OXT-C-CA-N
3	E	203	PCW	O31-C31-C32-C33
3	E	203	PCW	O3-C11-C12-C13
3	F	202	PCW	O3-C11-C12-C13

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Mol	Chain	Res	Type	Atoms
3	C	1102	PCW	O11-C11-C12-C13
3	C	1101	PCW	C19-C20-C21-C22
3	C	1101	PCW	O11-C11-C12-C13

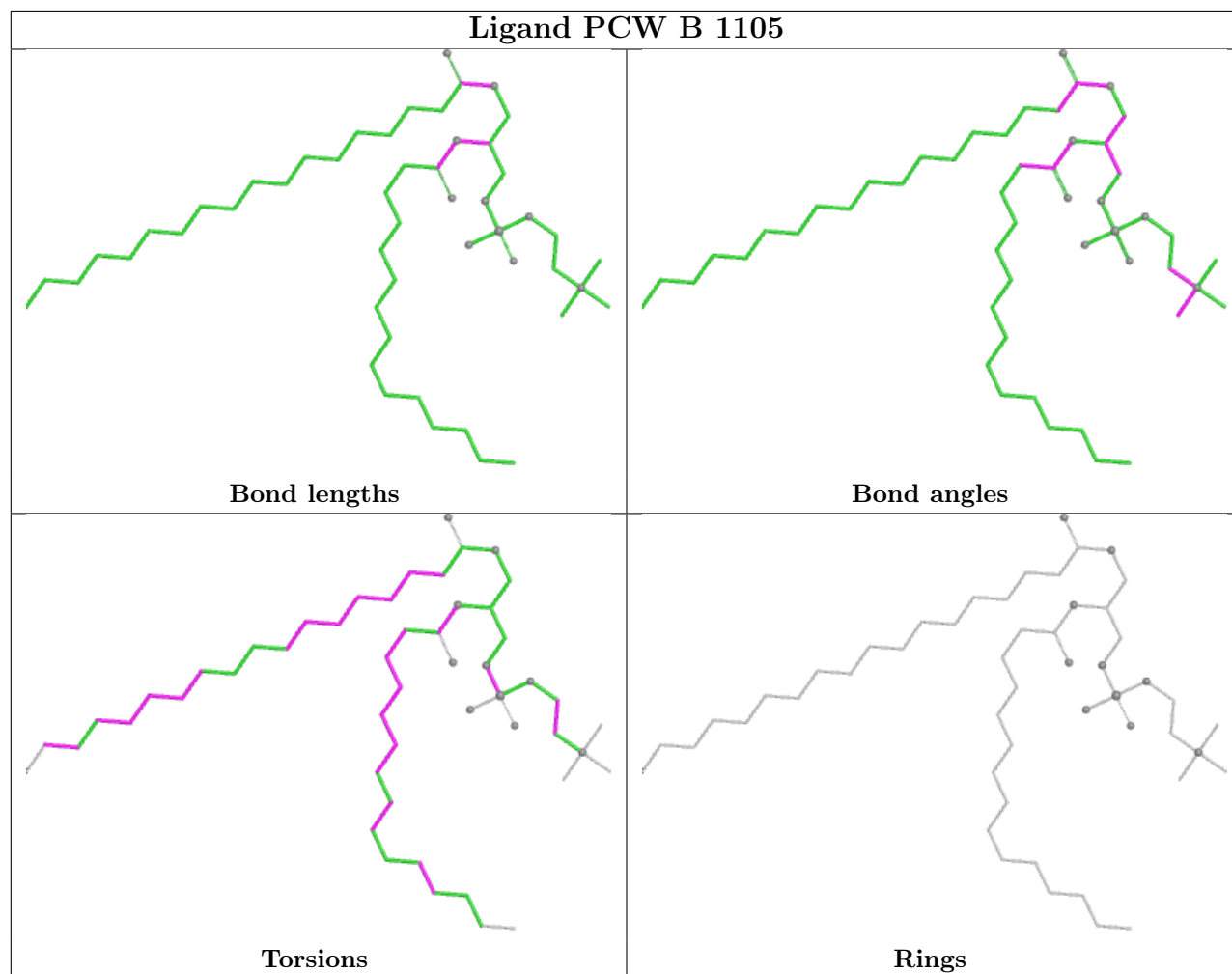
There are no ring outliers.

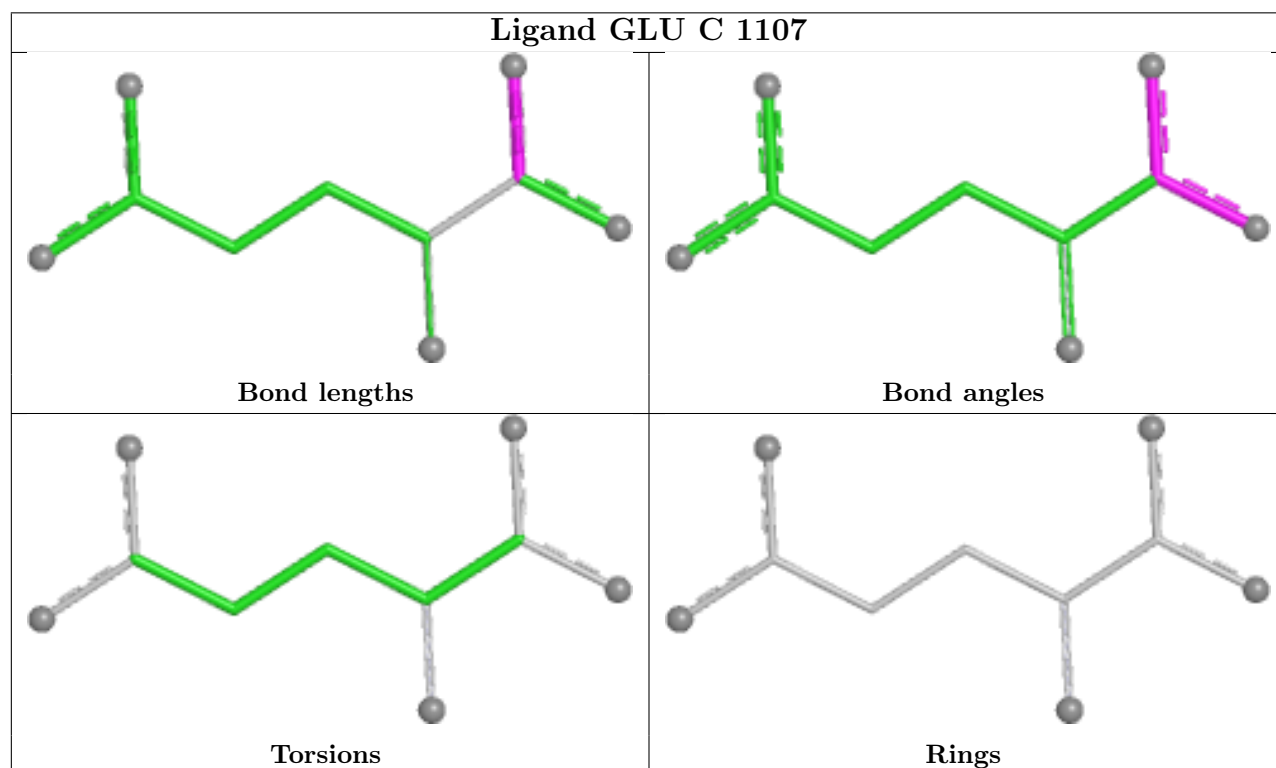
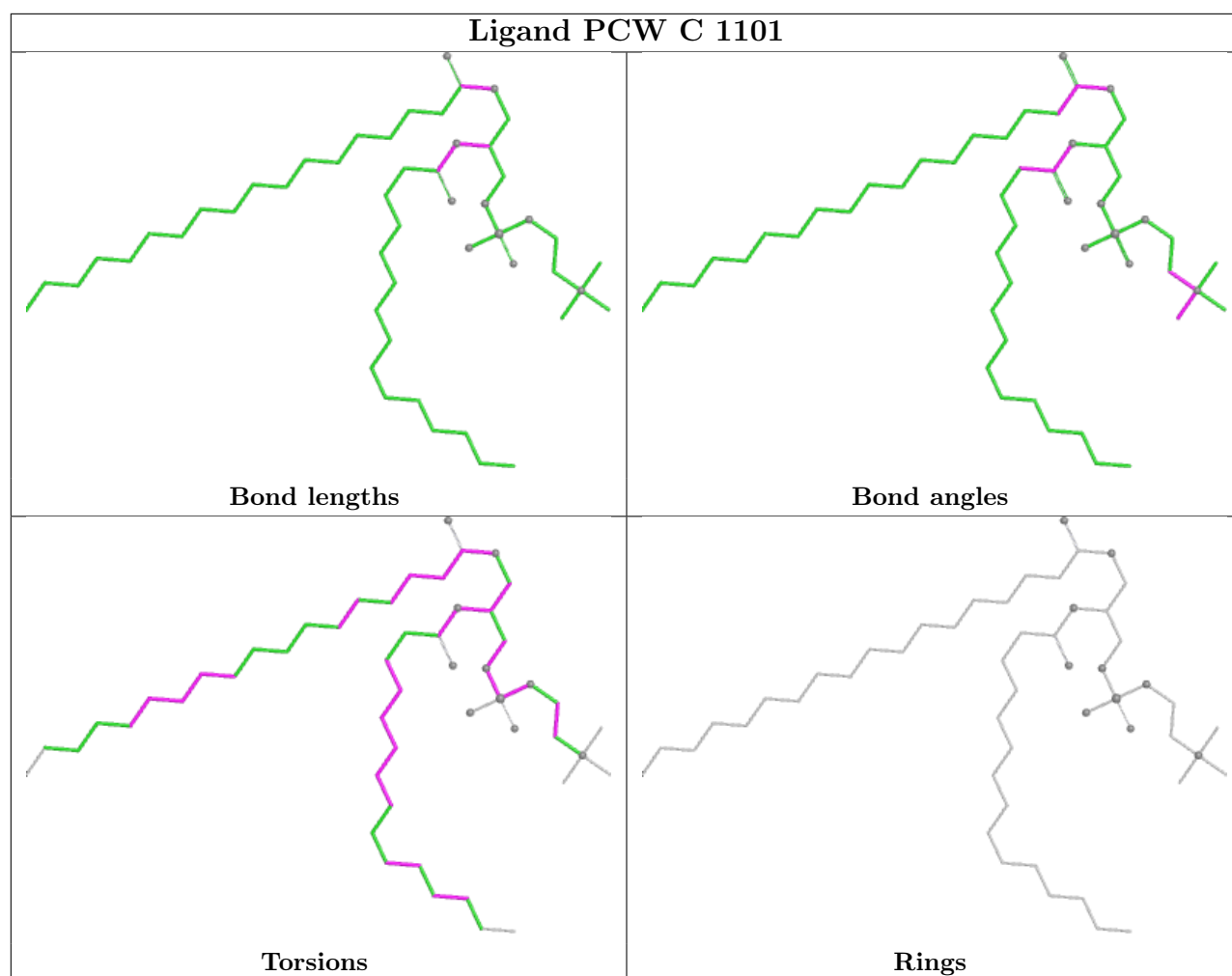
24 monomers are involved in 59 short contacts:

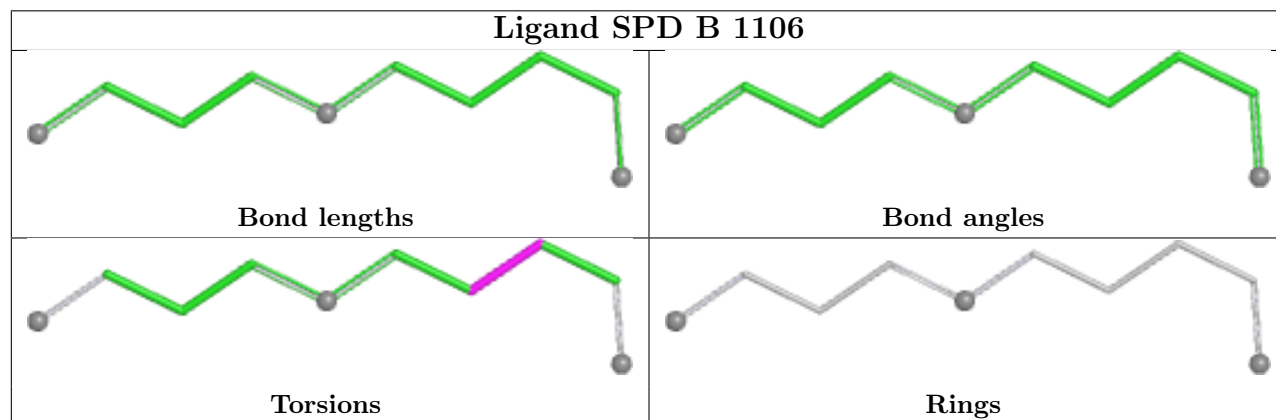
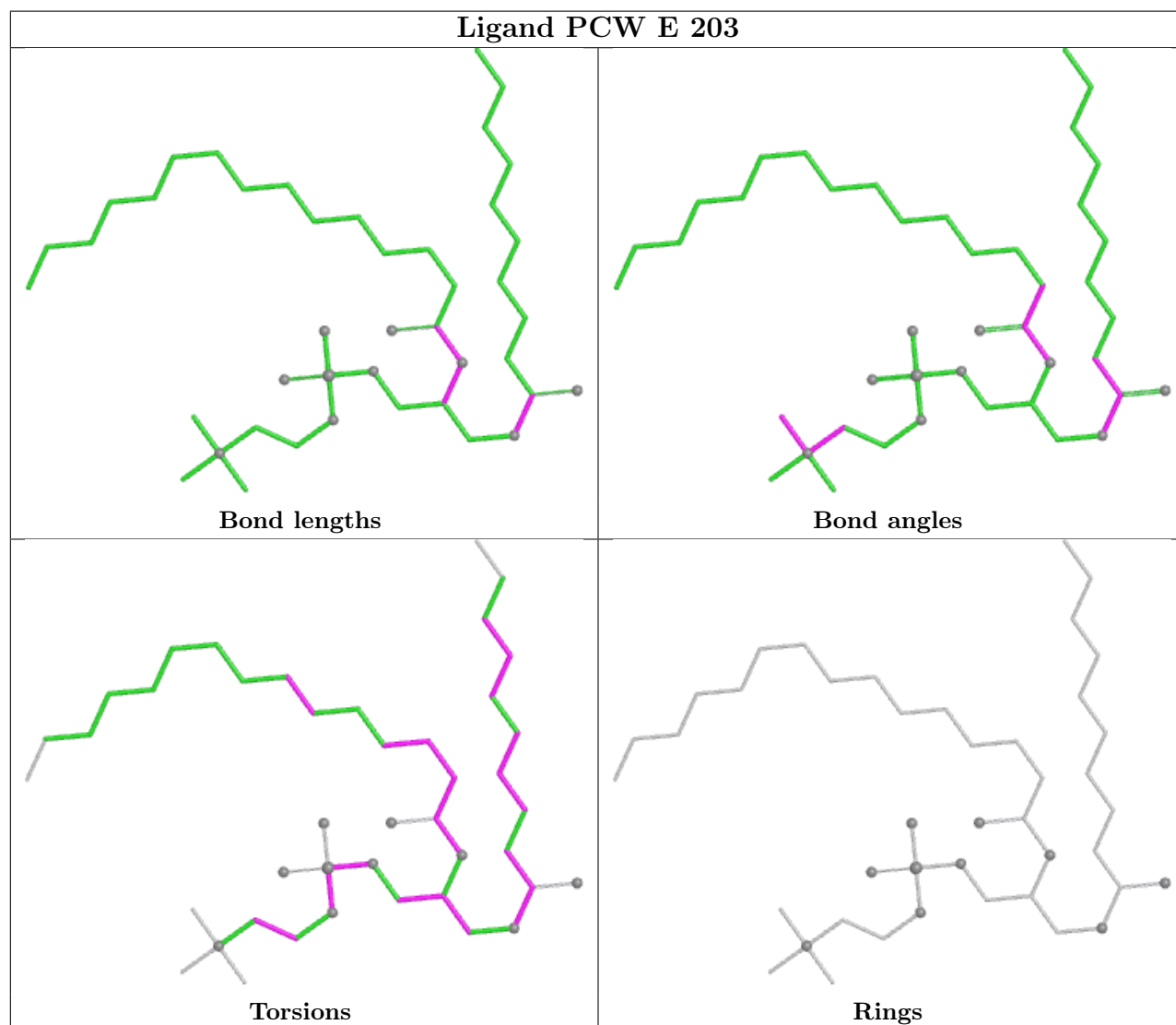
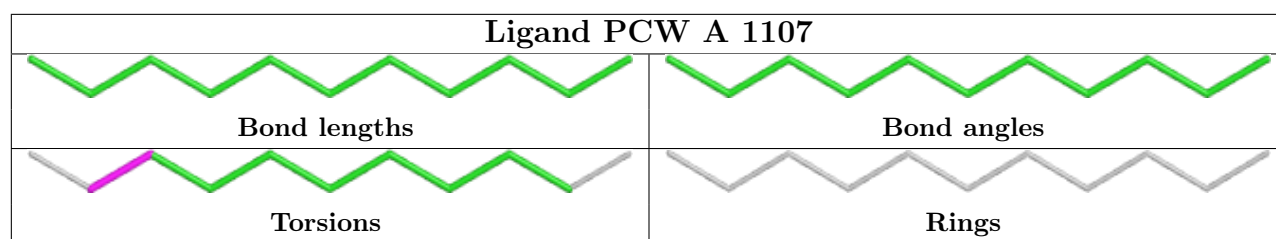
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1105	PCW	5	0
3	C	1101	PCW	3	0
3	E	203	PCW	2	0
5	B	1106	SPD	2	0
3	A	1102	PCW	5	0
3	B	1102	PCW	4	0
3	D	1101	PCW	5	0
3	A	1104	PCW	2	0
3	B	1103	PCW	2	0
3	F	202	PCW	3	0
3	A	1101	PCW	5	0
3	C	1103	PCW	3	0
3	A	1109	PCW	1	0
3	A	1103	PCW	4	0
3	D	1103	PCW	3	0
4	B	1104	GLU	2	0
3	D	1102	PCW	4	0
3	C	1104	PCW	2	0
3	C	1108	PCW	1	0
3	D	1104	PCW	1	0
3	B	1101	PCW	4	0
3	E	201	PCW	2	0
3	C	1102	PCW	4	0
4	D	1105	GLU	2	0

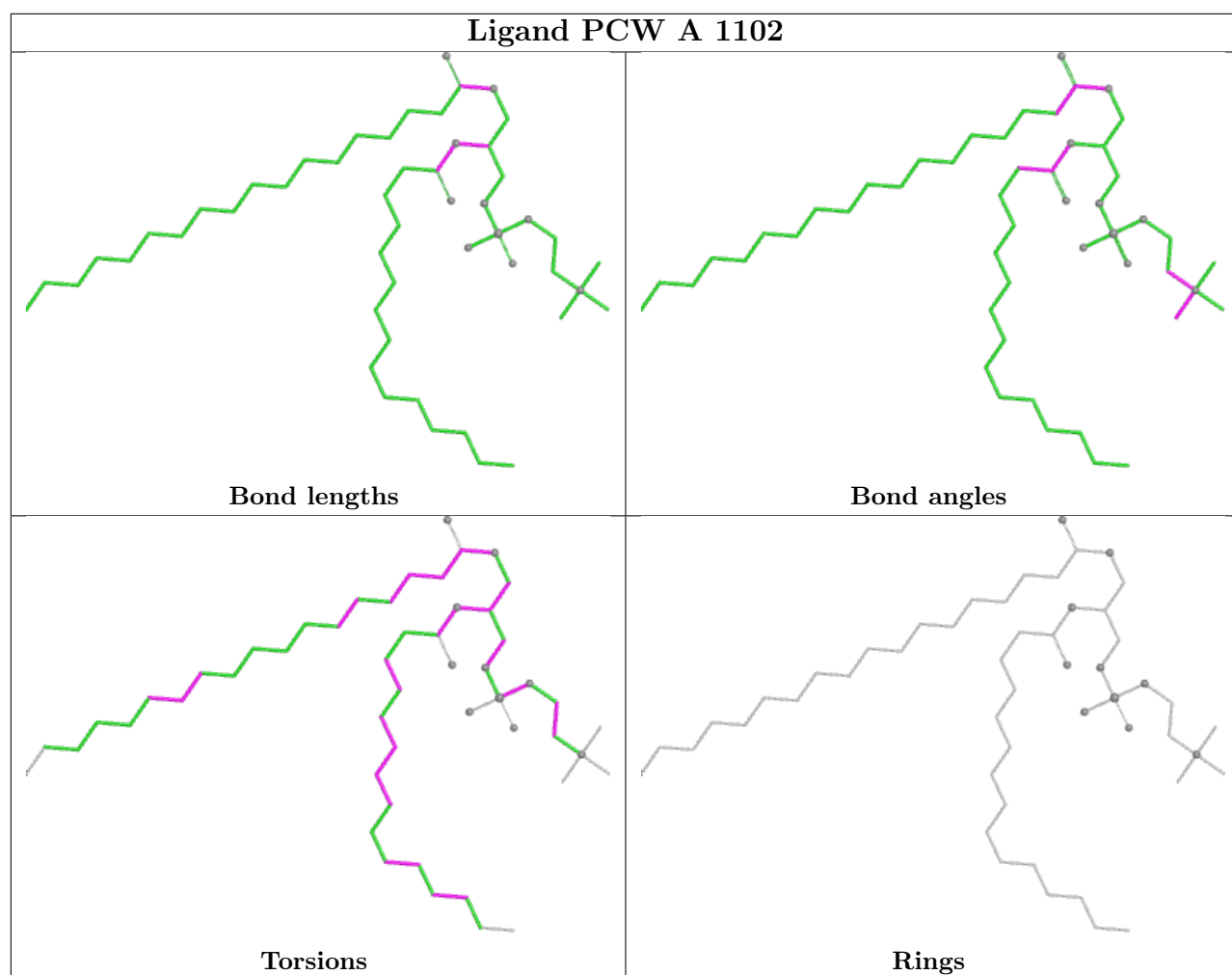
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

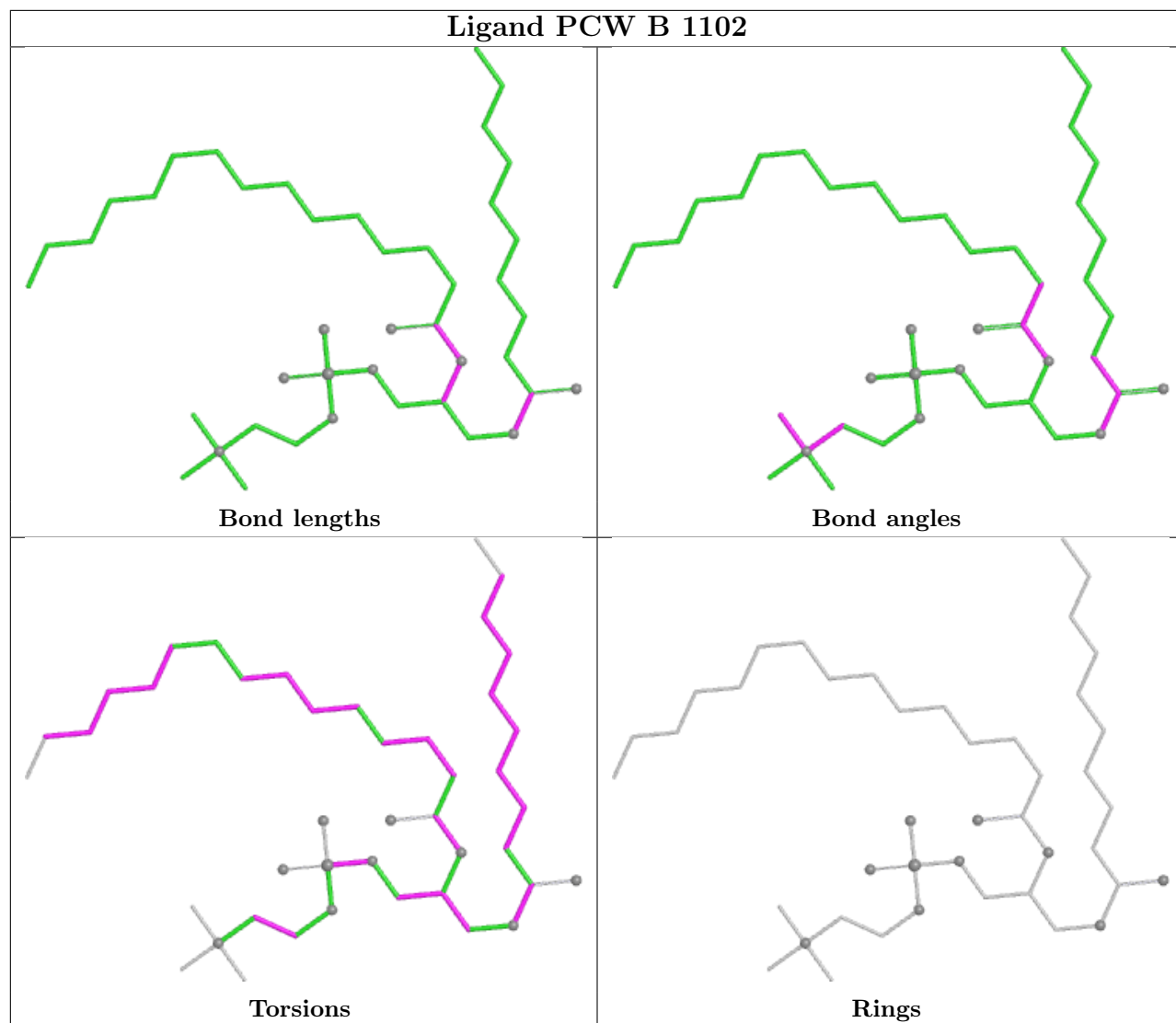
any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

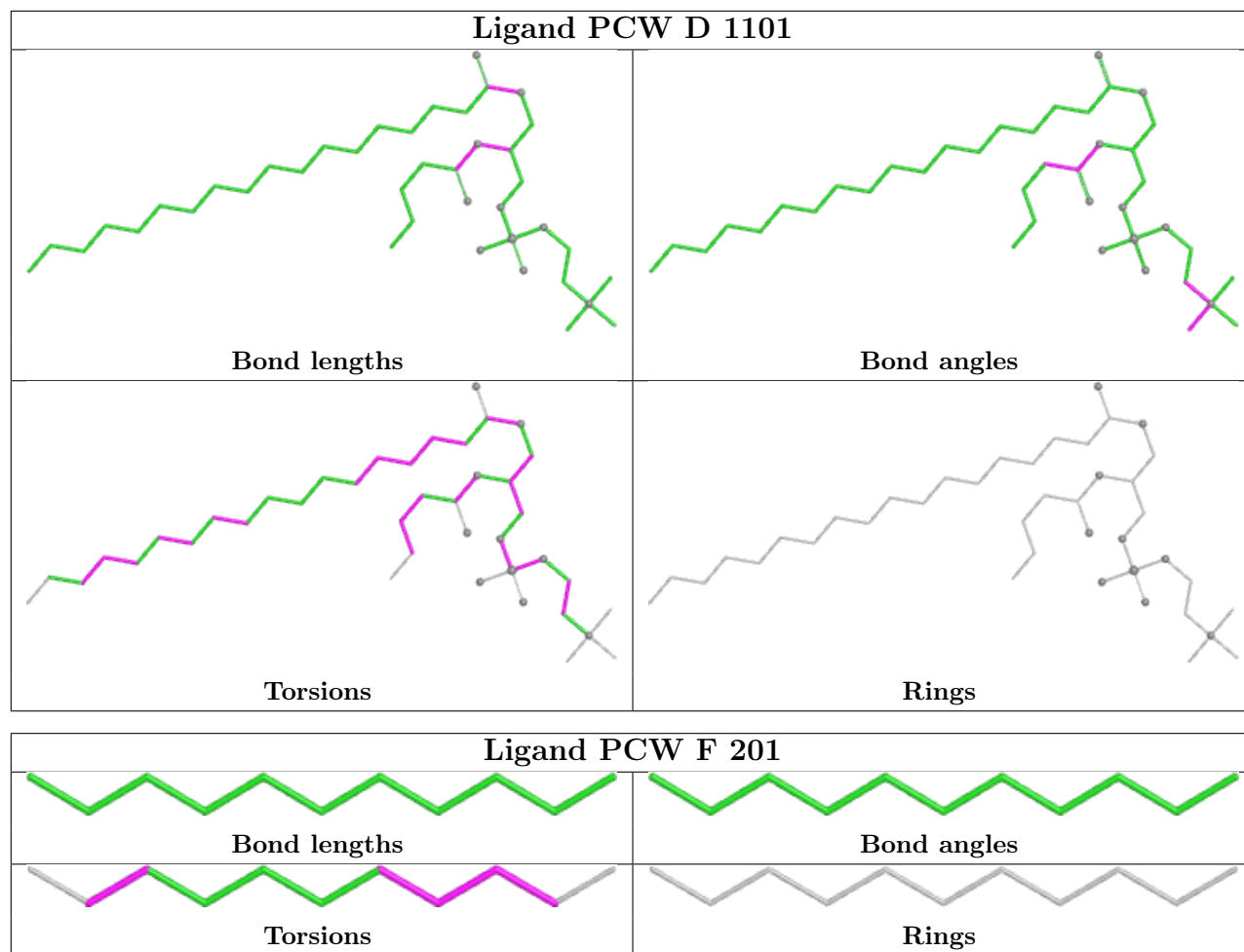


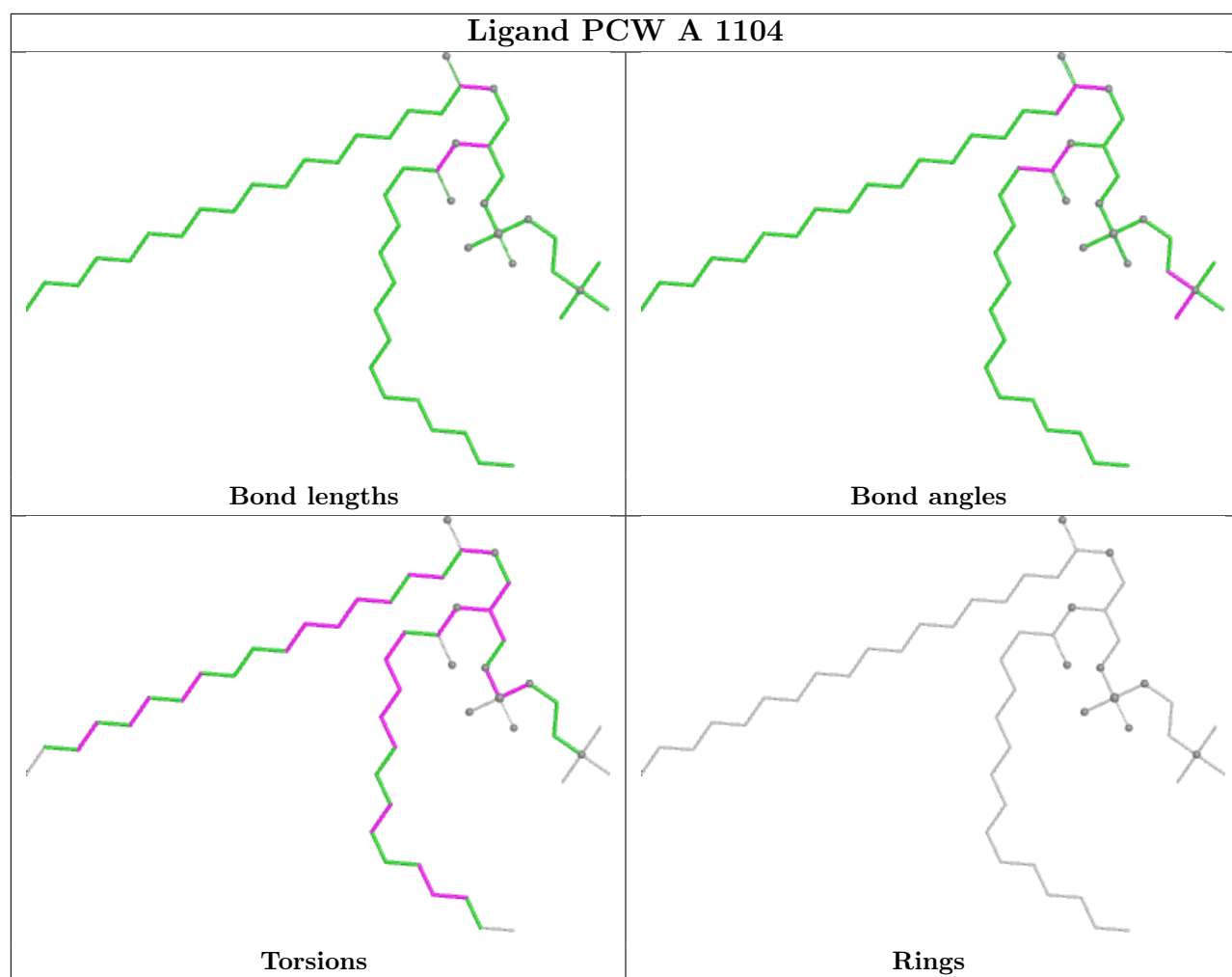


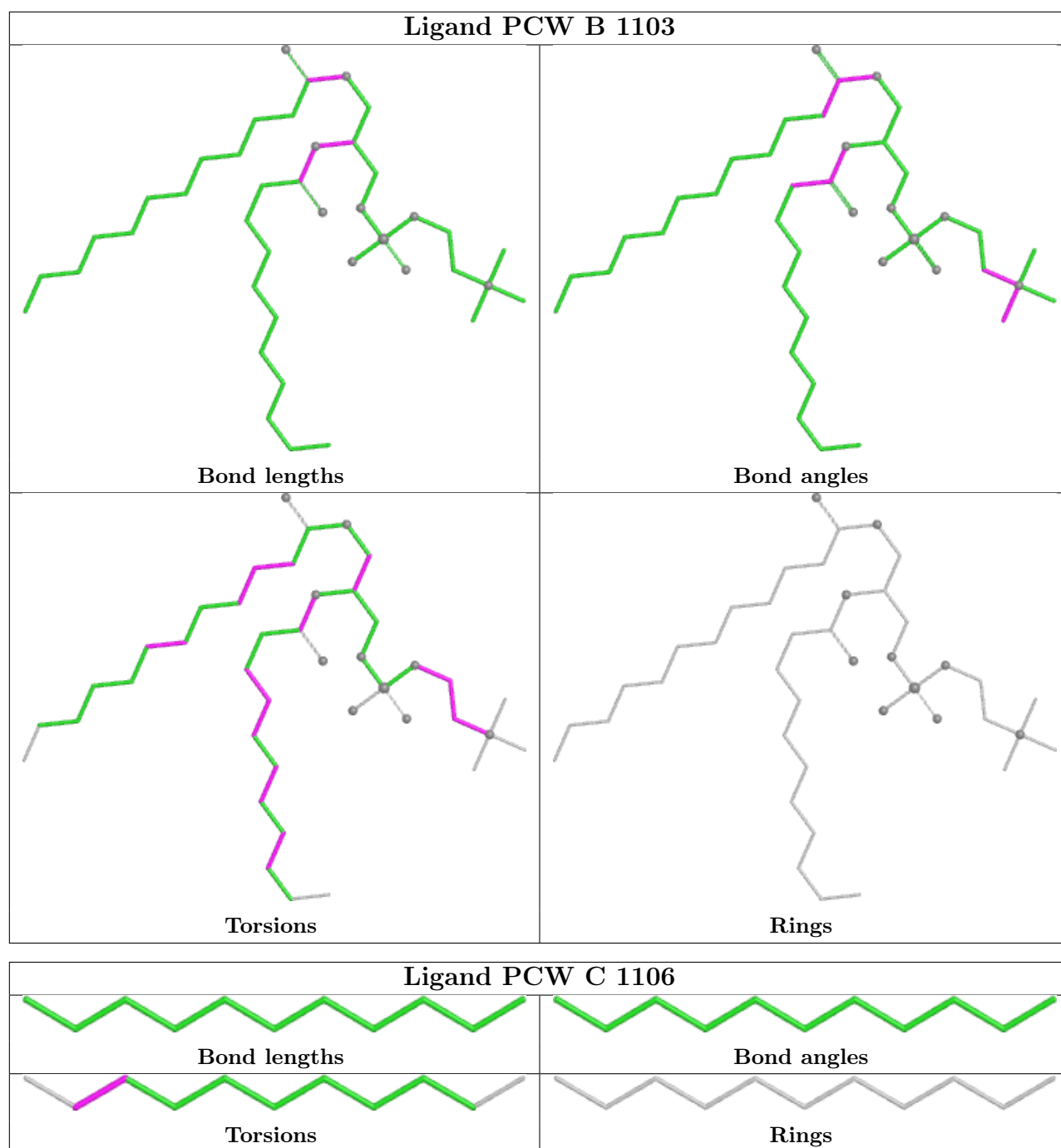


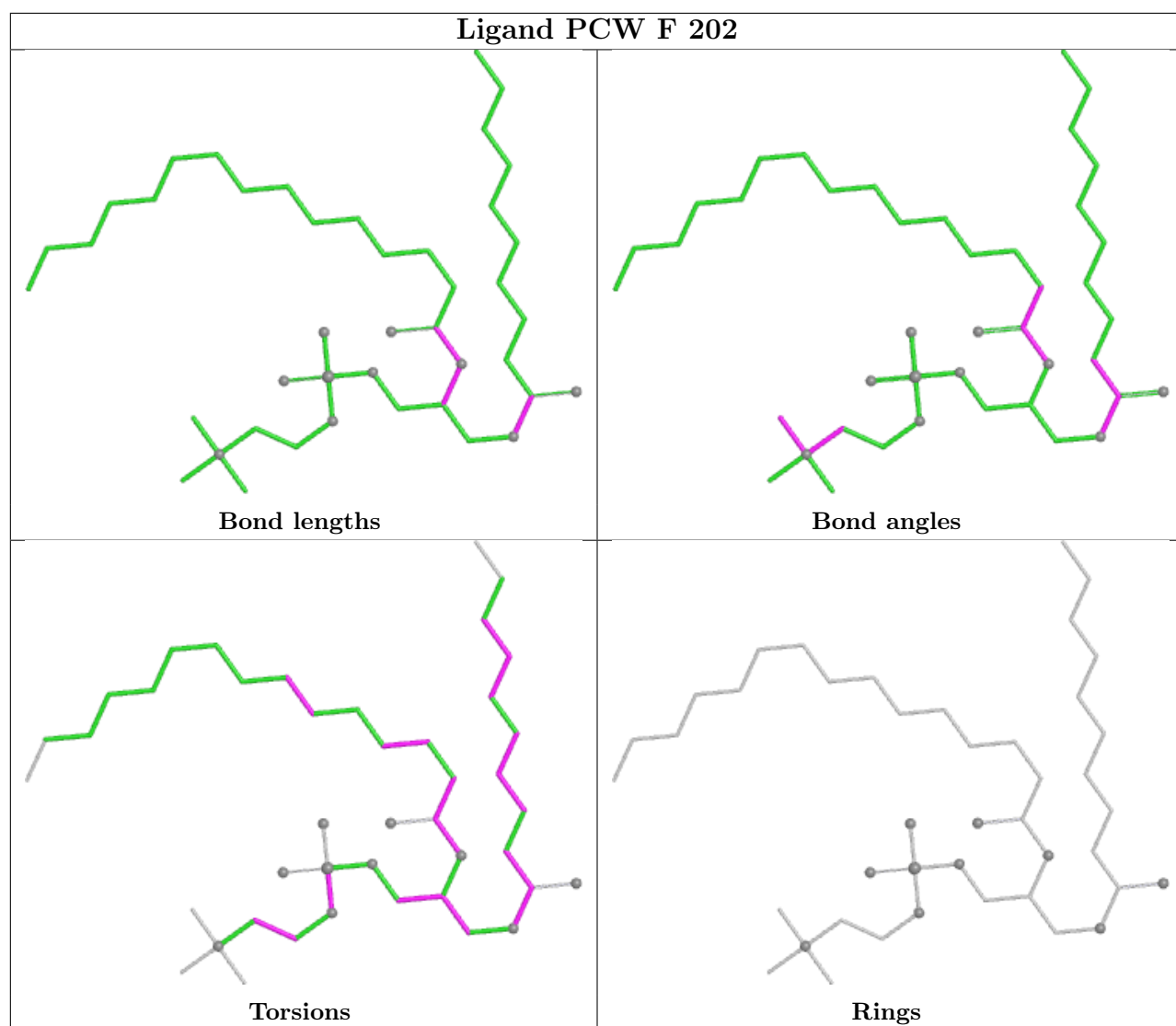


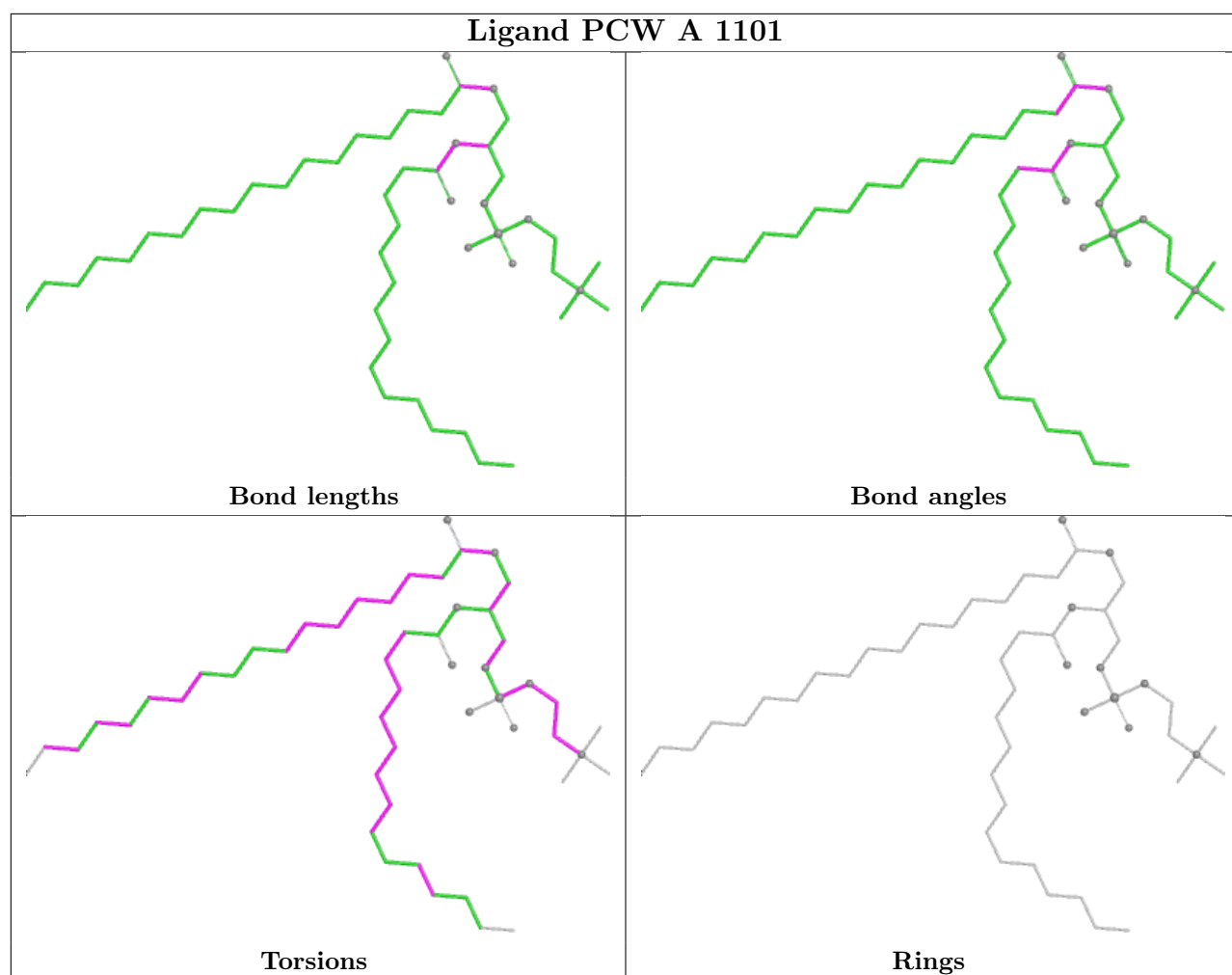


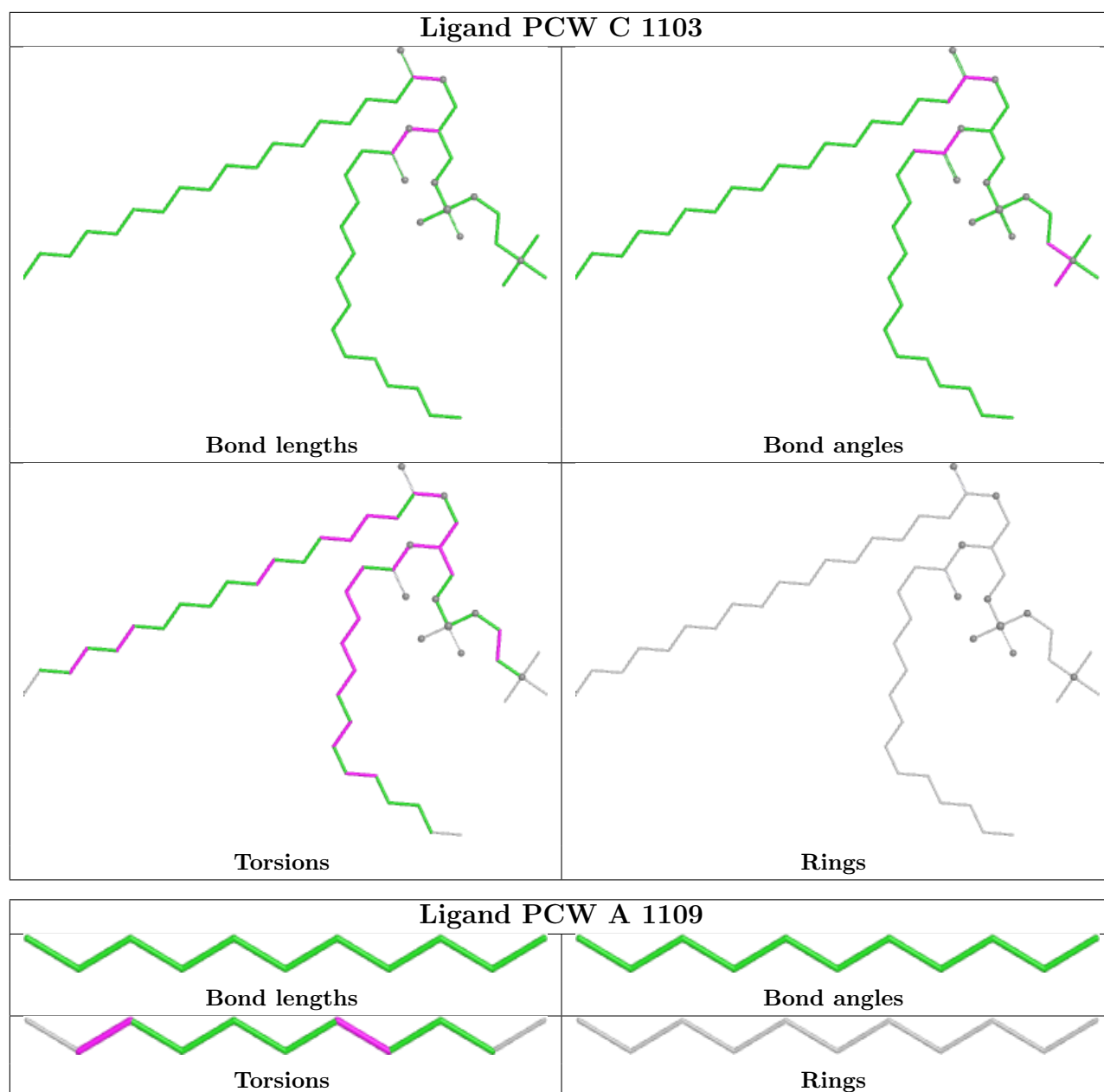


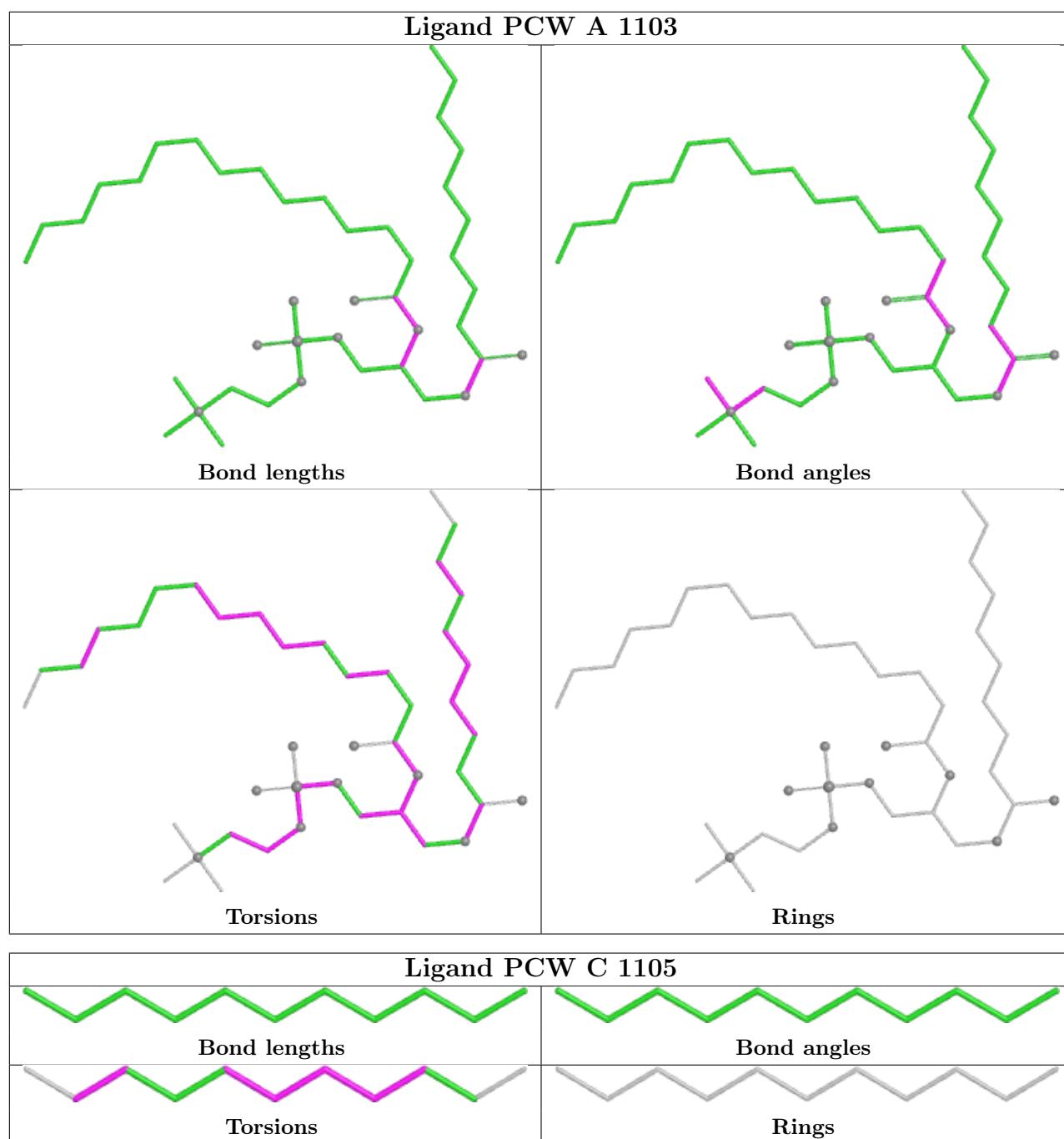


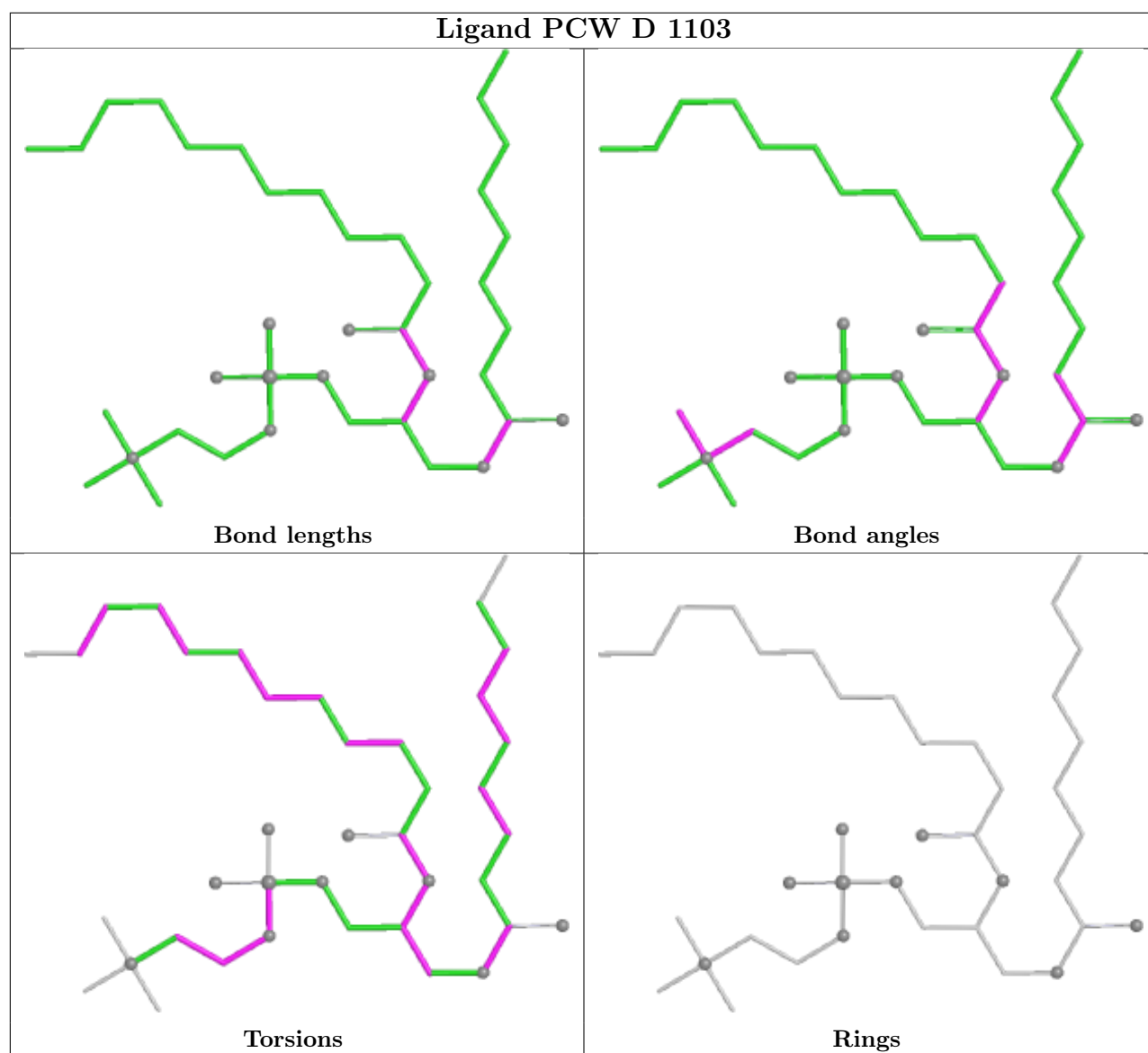


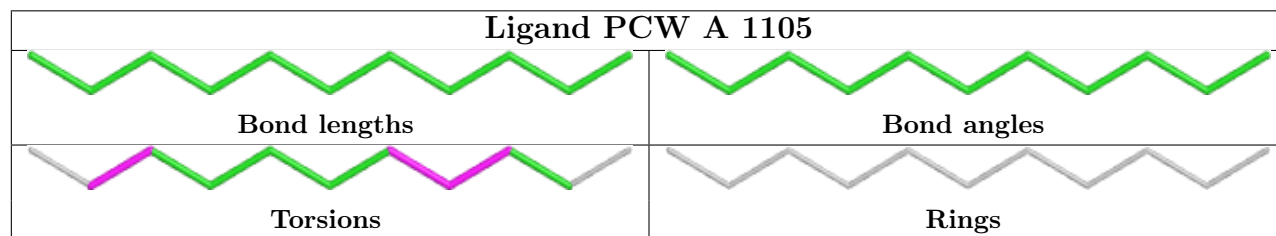
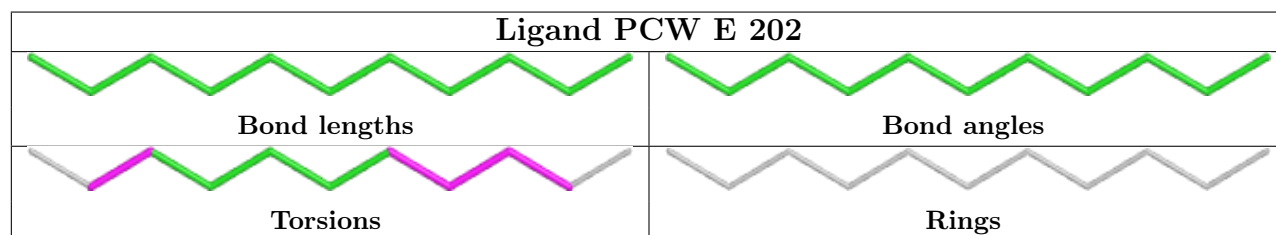
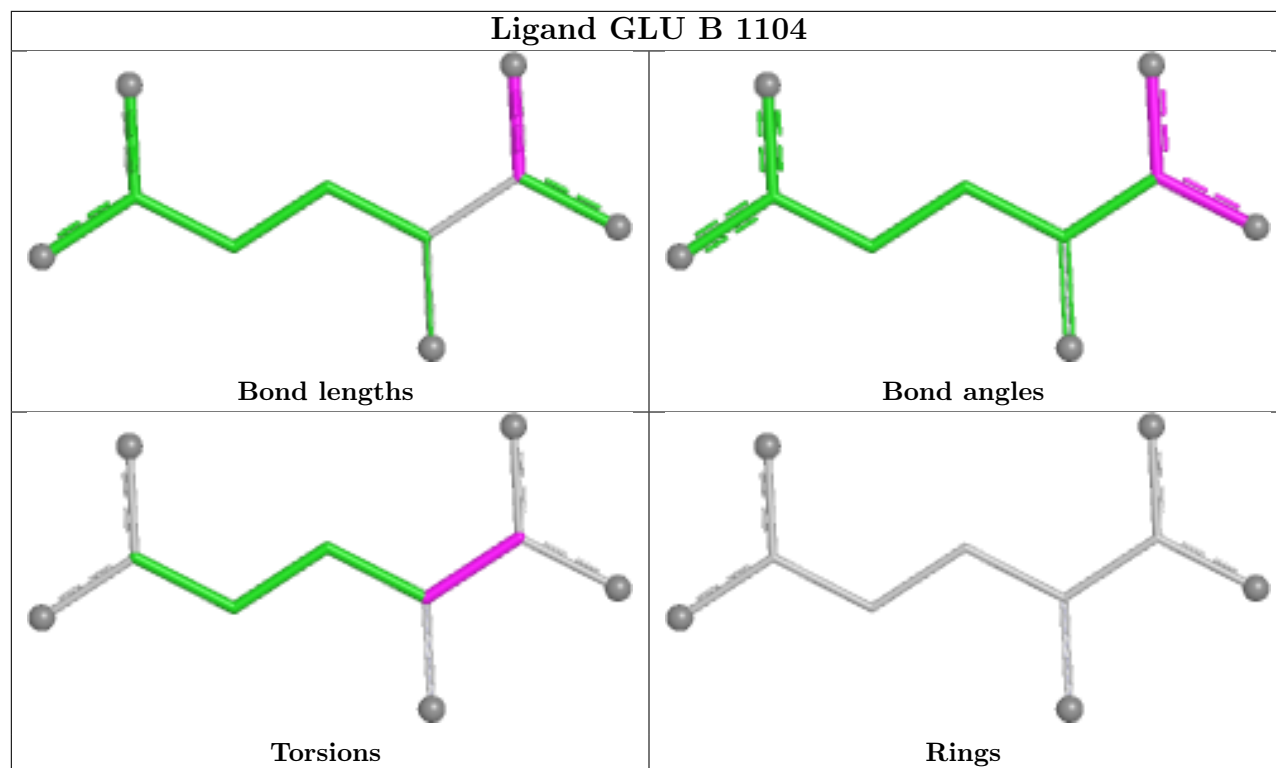


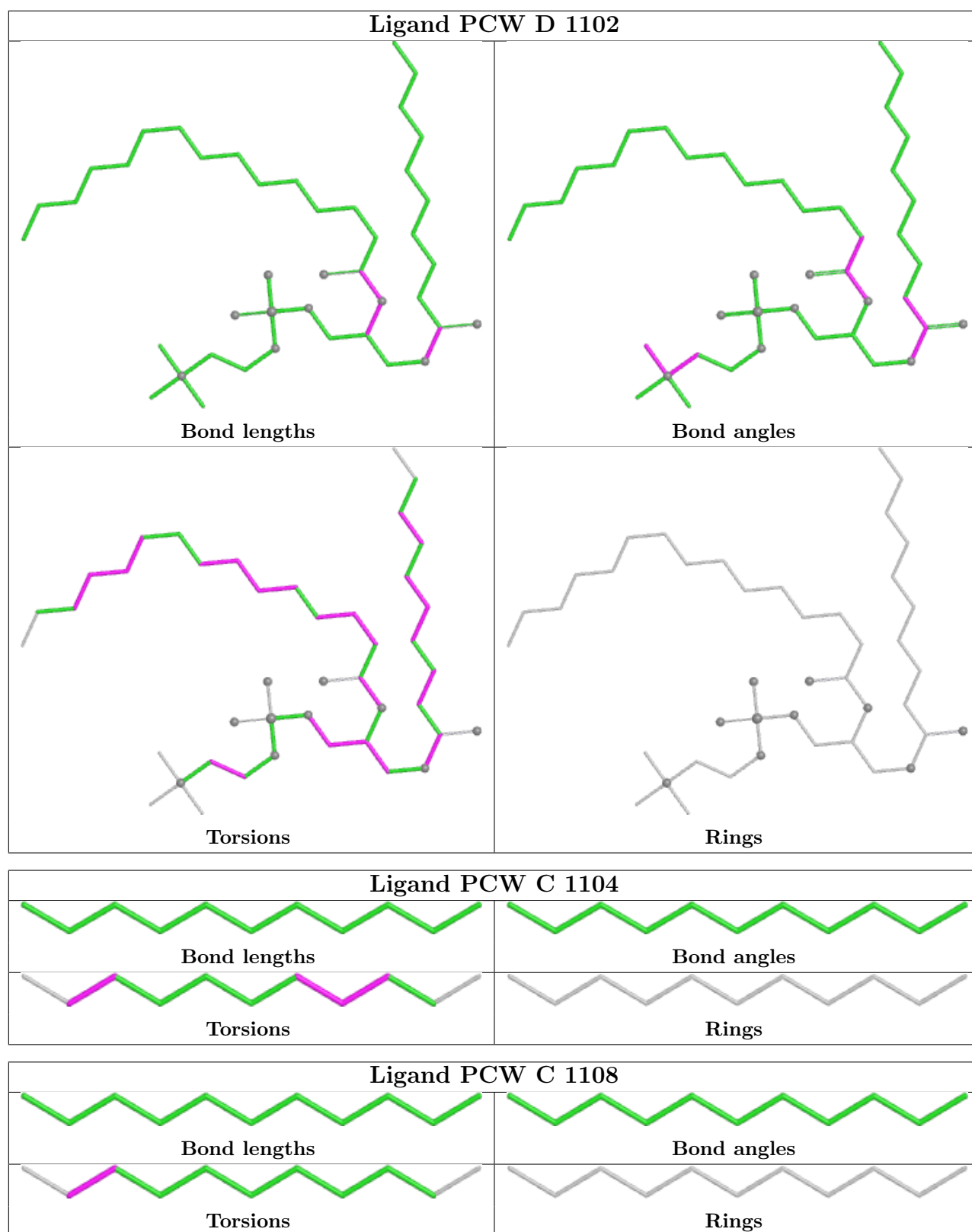


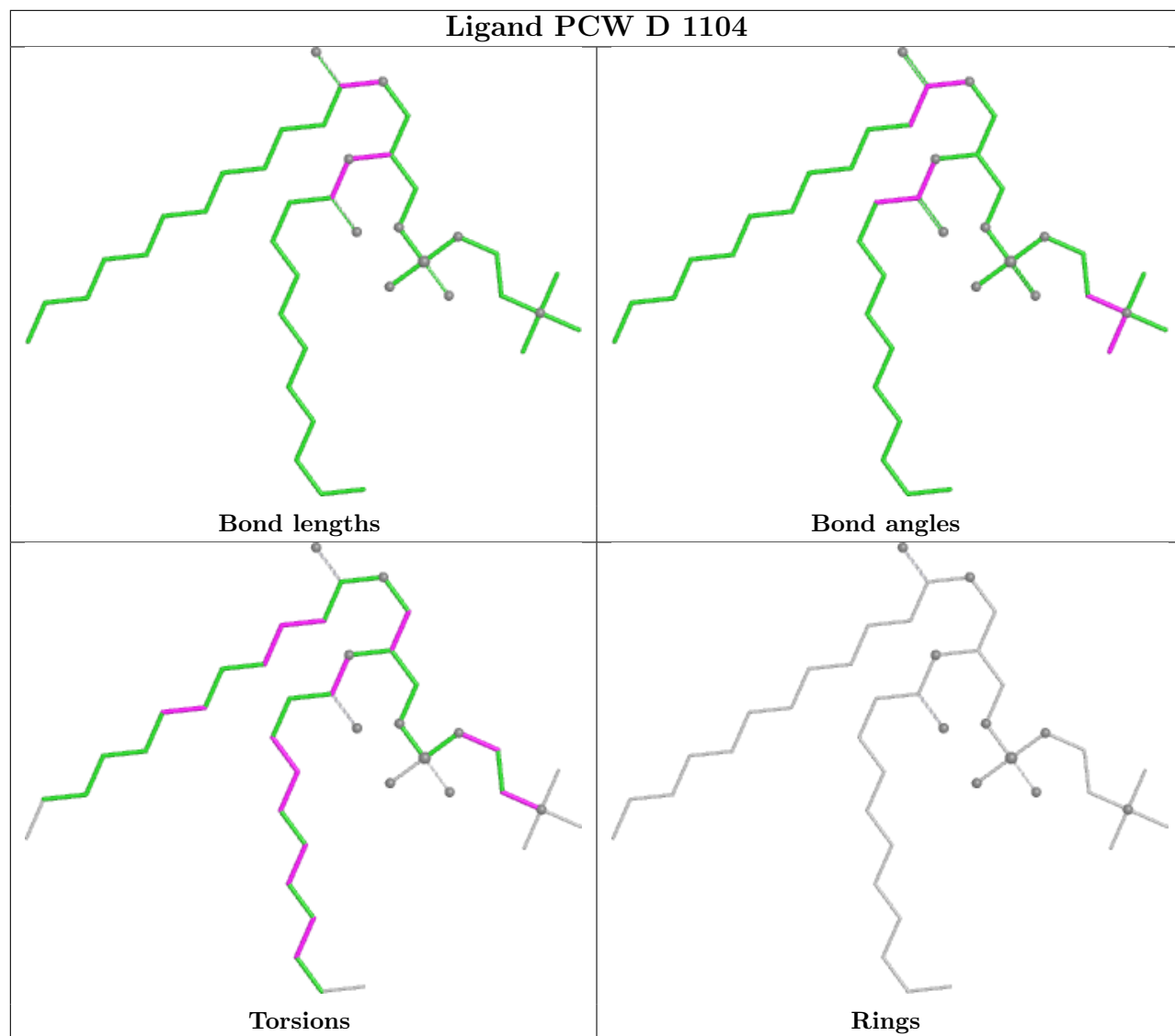


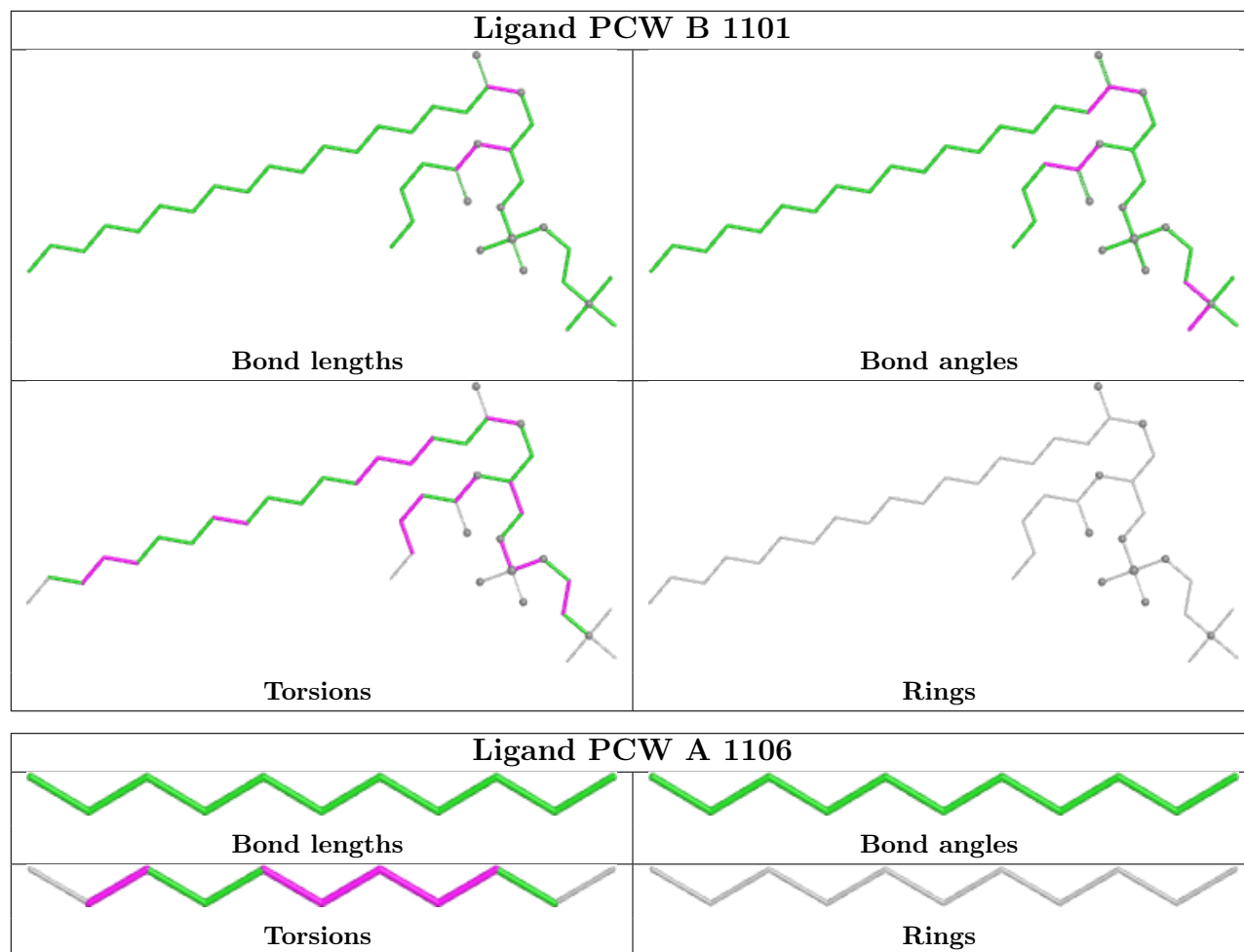


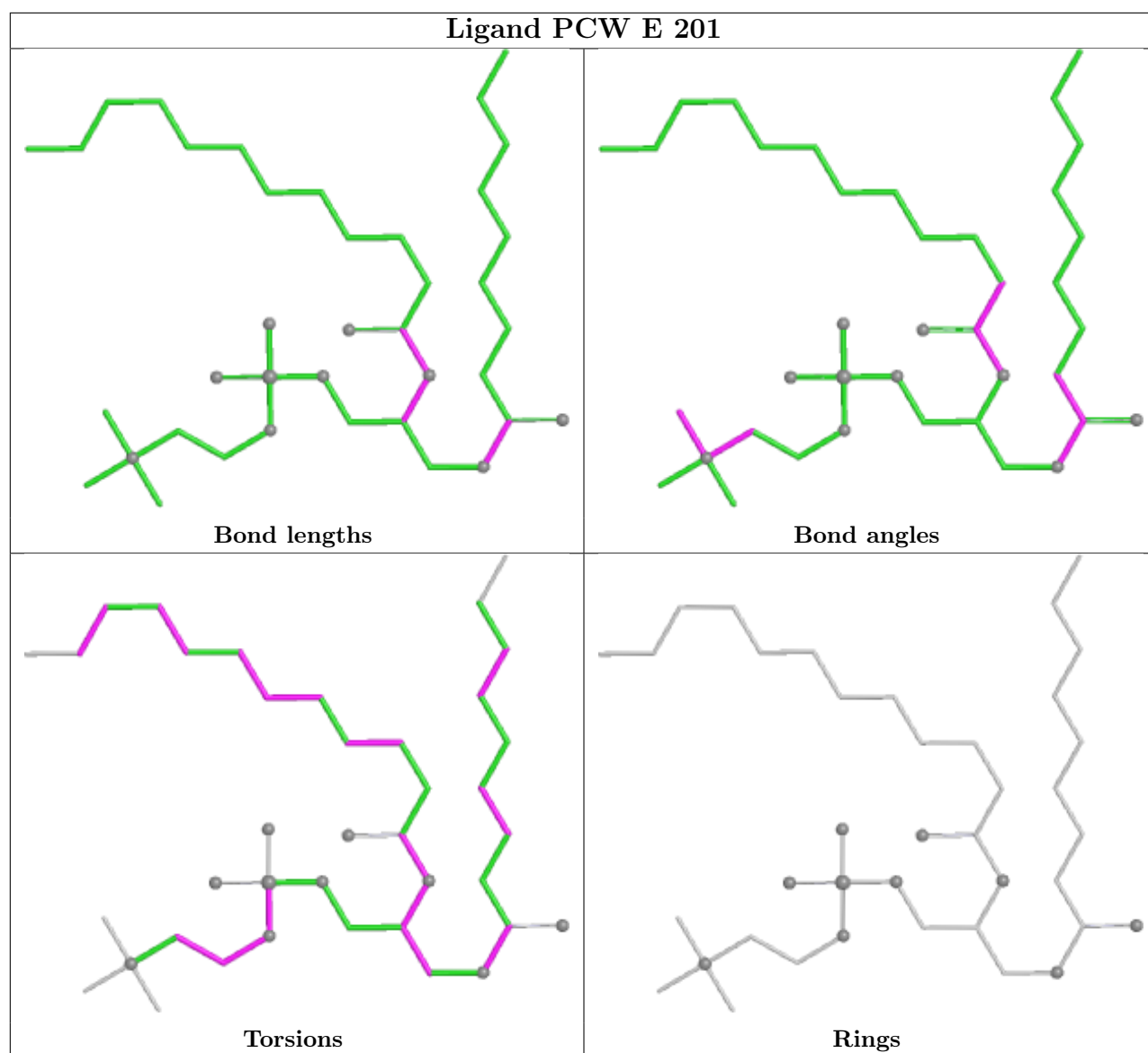


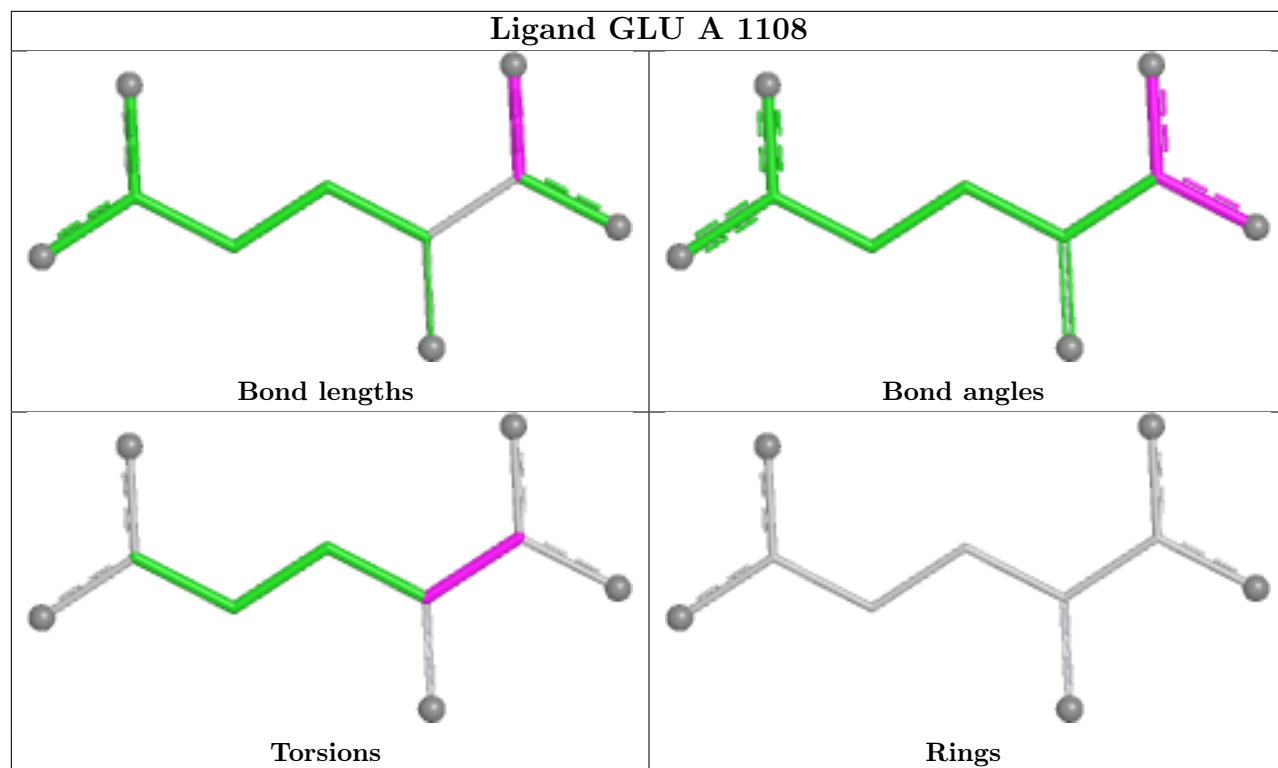


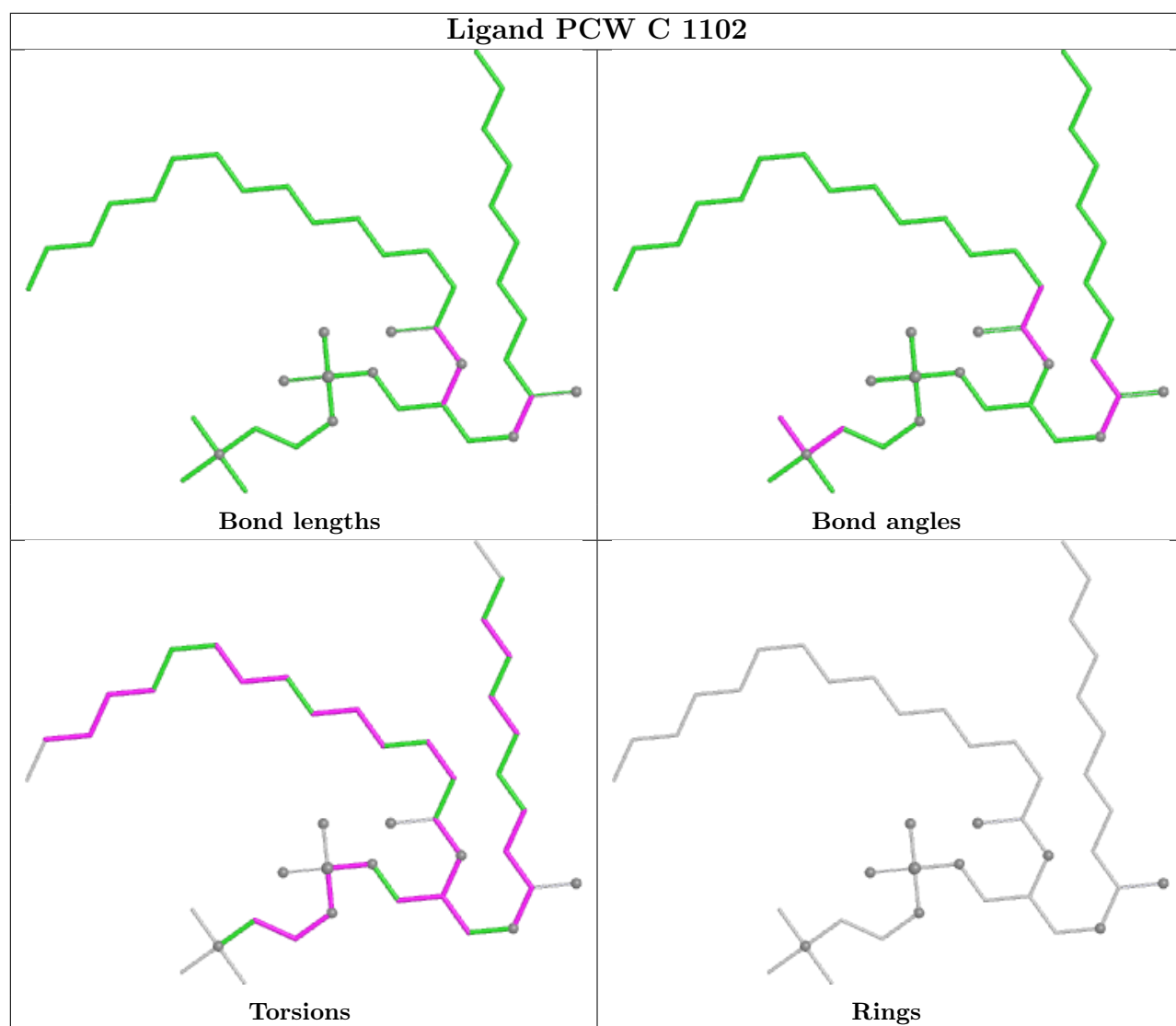


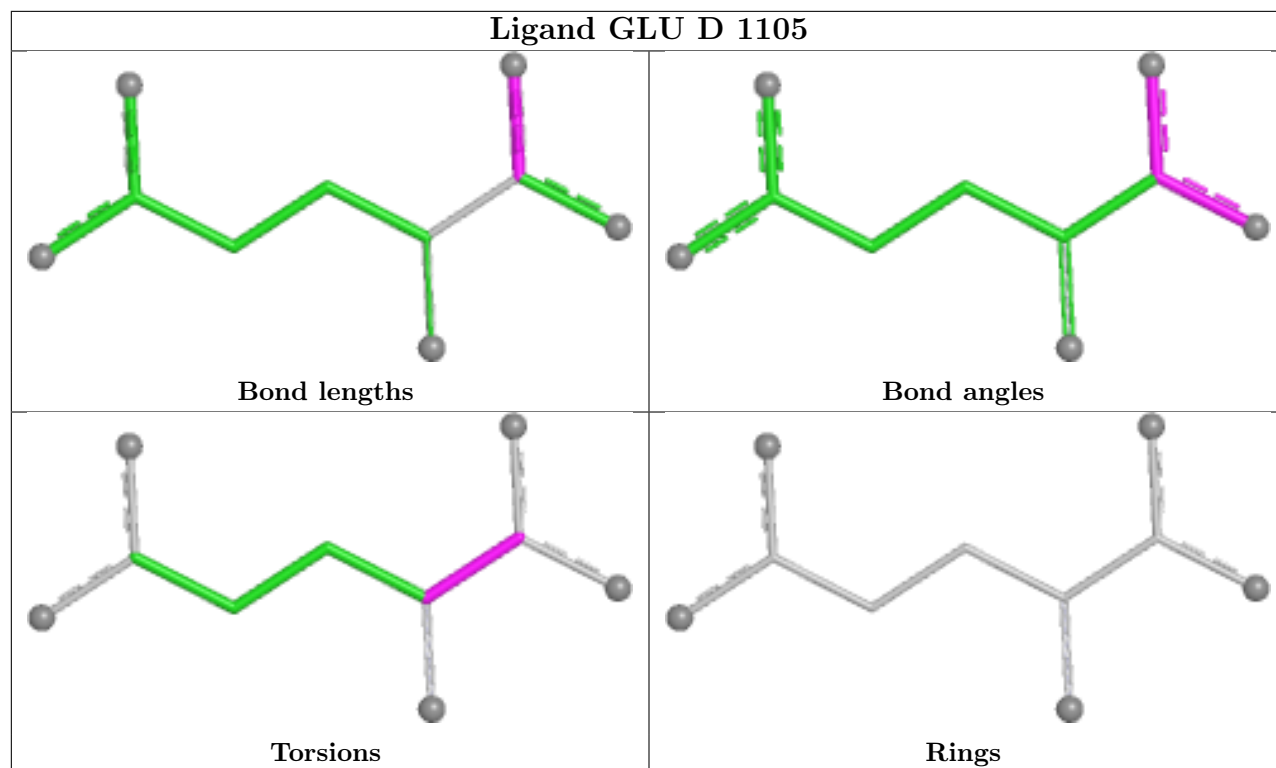












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

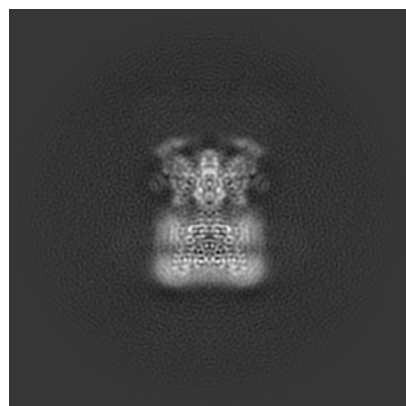
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-40750. These allow visual inspection of the internal detail of the map and identification of artifacts.

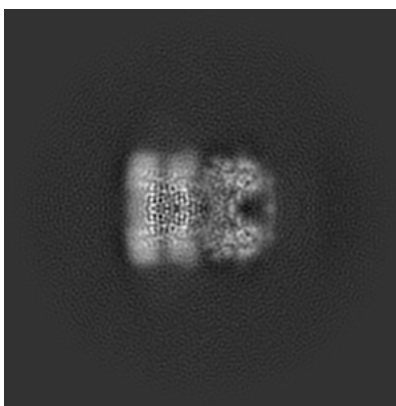
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

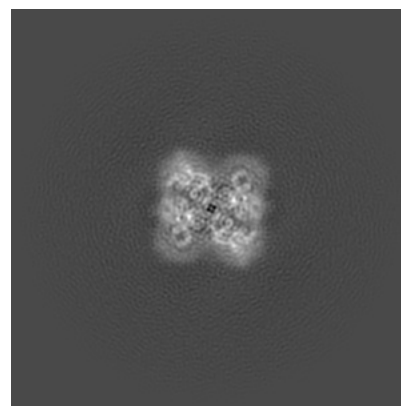
6.1.1 Primary map



X

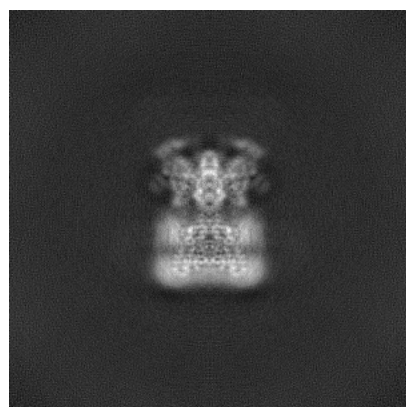


Y

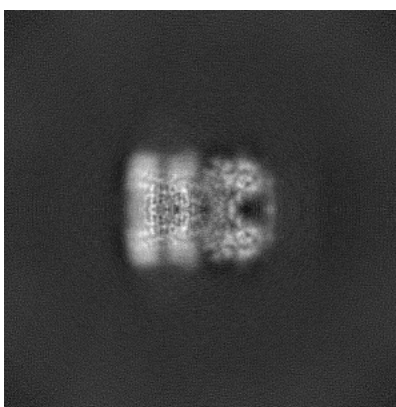


Z

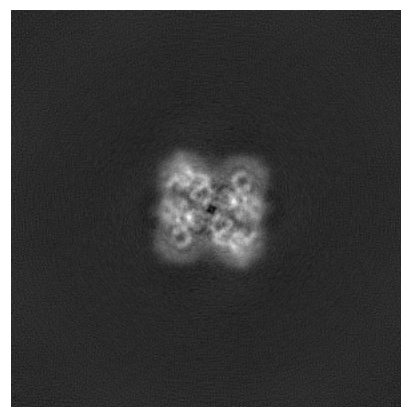
6.1.2 Raw map



X



Y

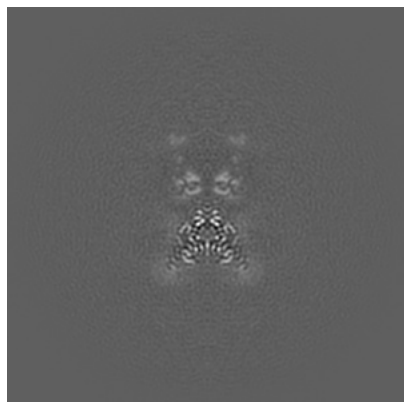


Z

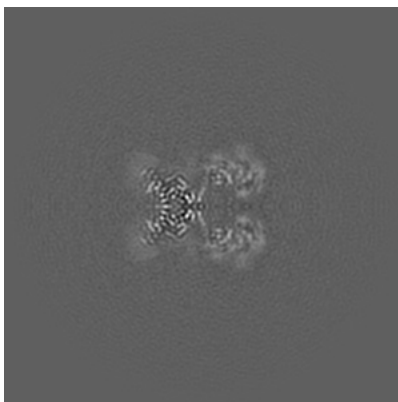
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

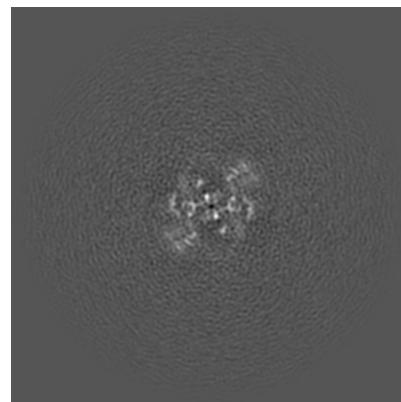
6.2.1 Primary map



X Index: 208

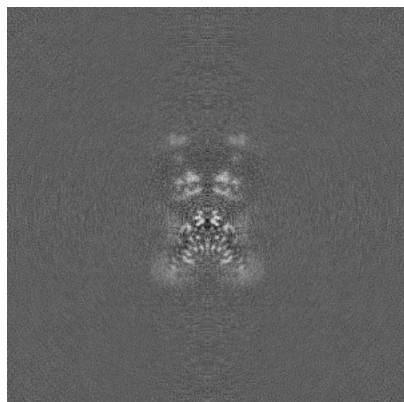


Y Index: 208

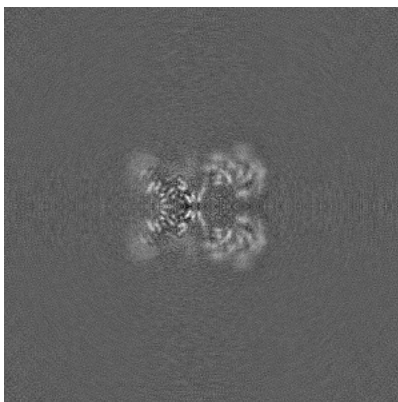


Z Index: 208

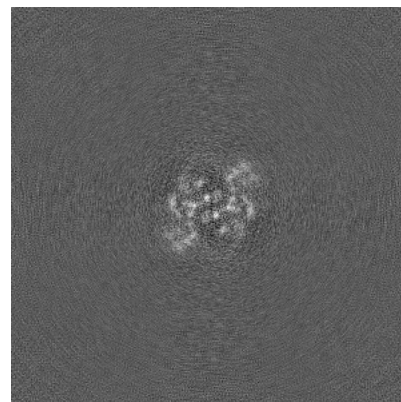
6.2.2 Raw map



X Index: 208



Y Index: 208

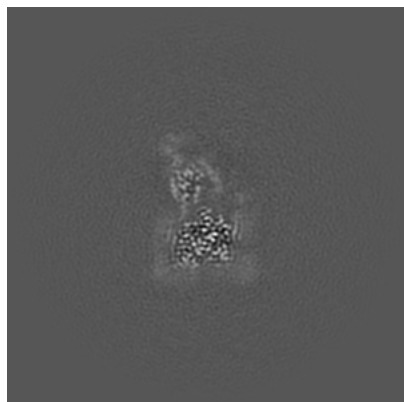


Z Index: 208

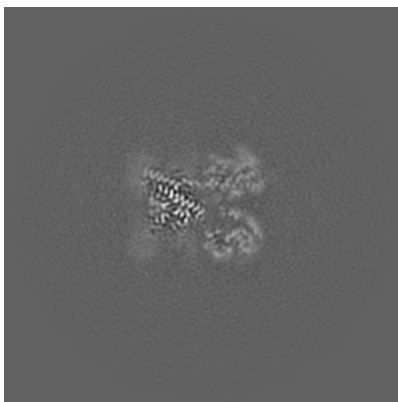
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

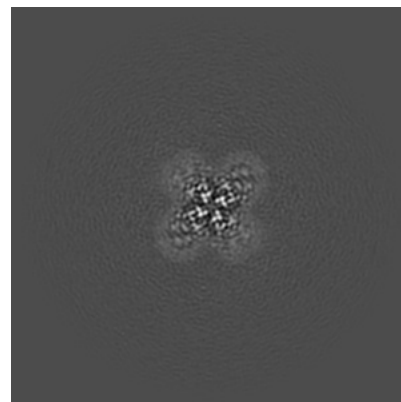
6.3.1 Primary map



X Index: 219

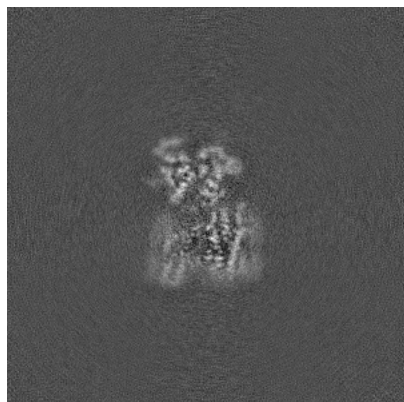


Y Index: 214

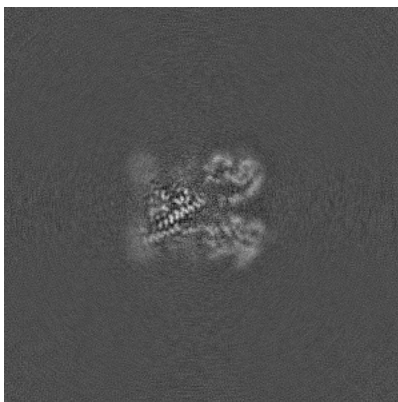


Z Index: 179

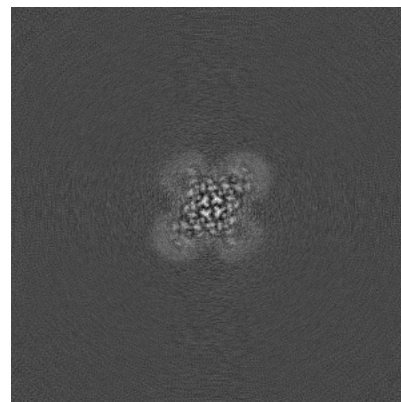
6.3.2 Raw map



X Index: 233



Y Index: 202

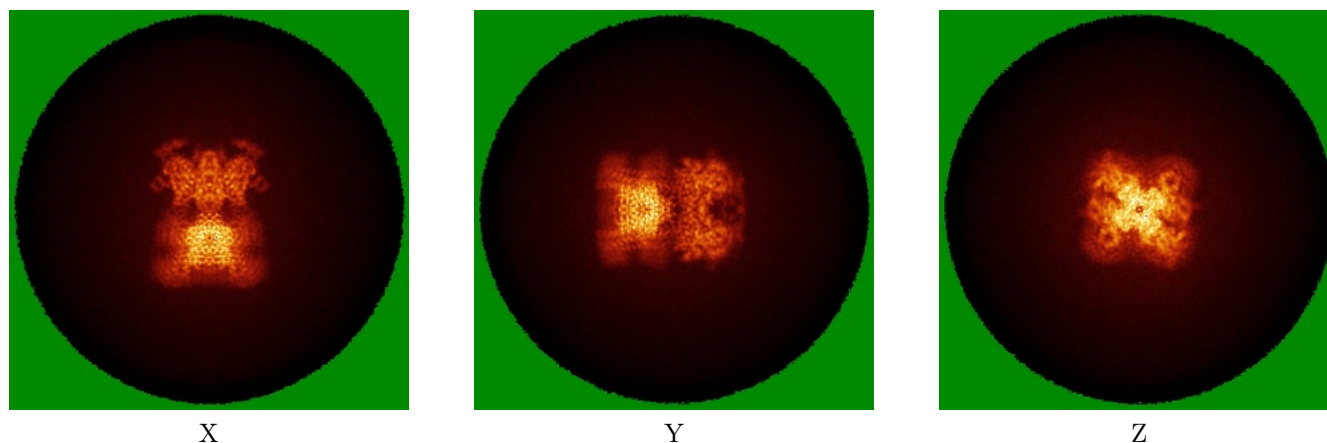


Z Index: 185

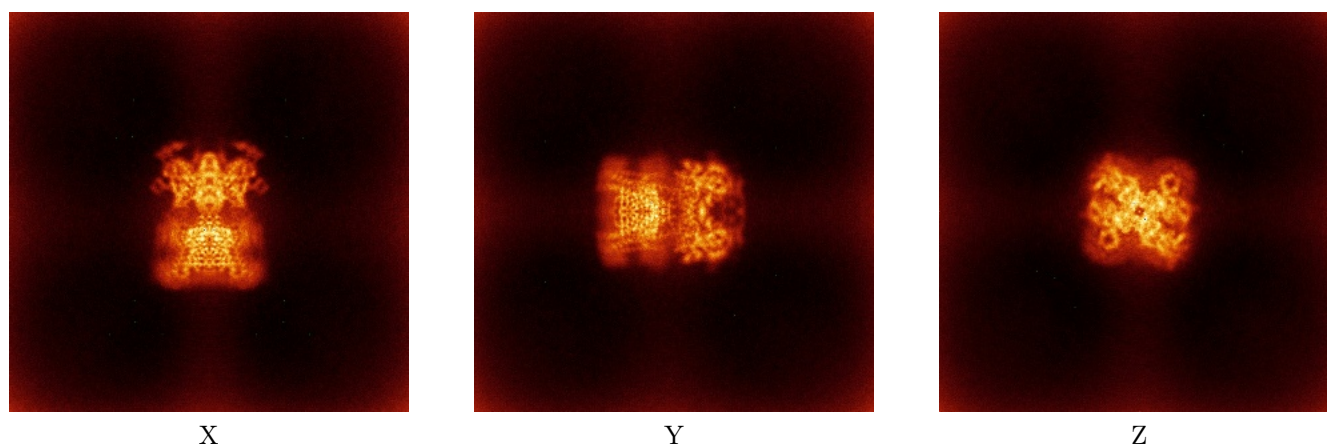
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) [i](#)

6.4.1 Primary map



6.4.2 Raw map



The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views [i](#)

This section was not generated.

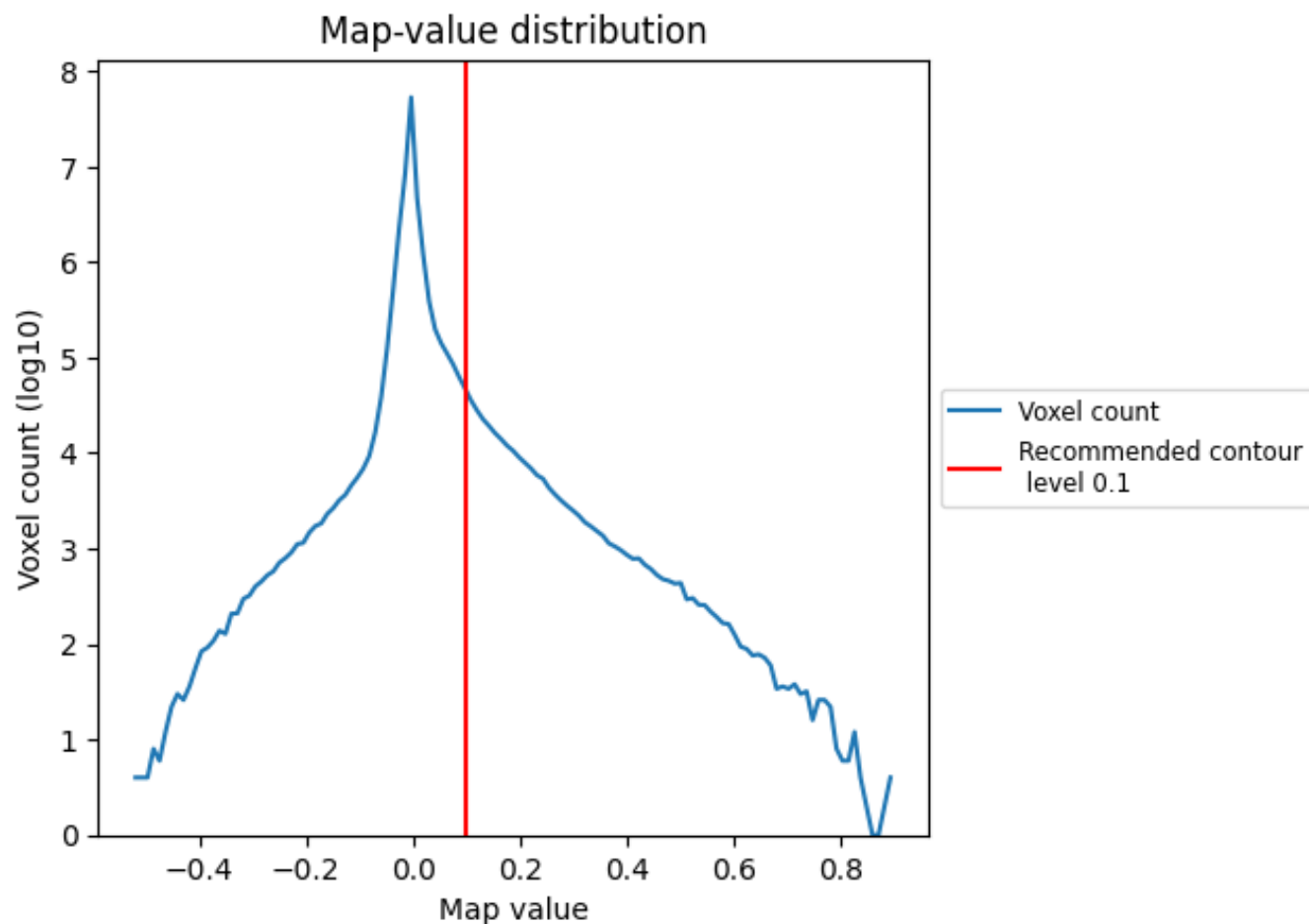
6.6 Mask visualisation [i](#)

This section was not generated. No masks/segmentation were deposited.

7 Map analysis [i](#)

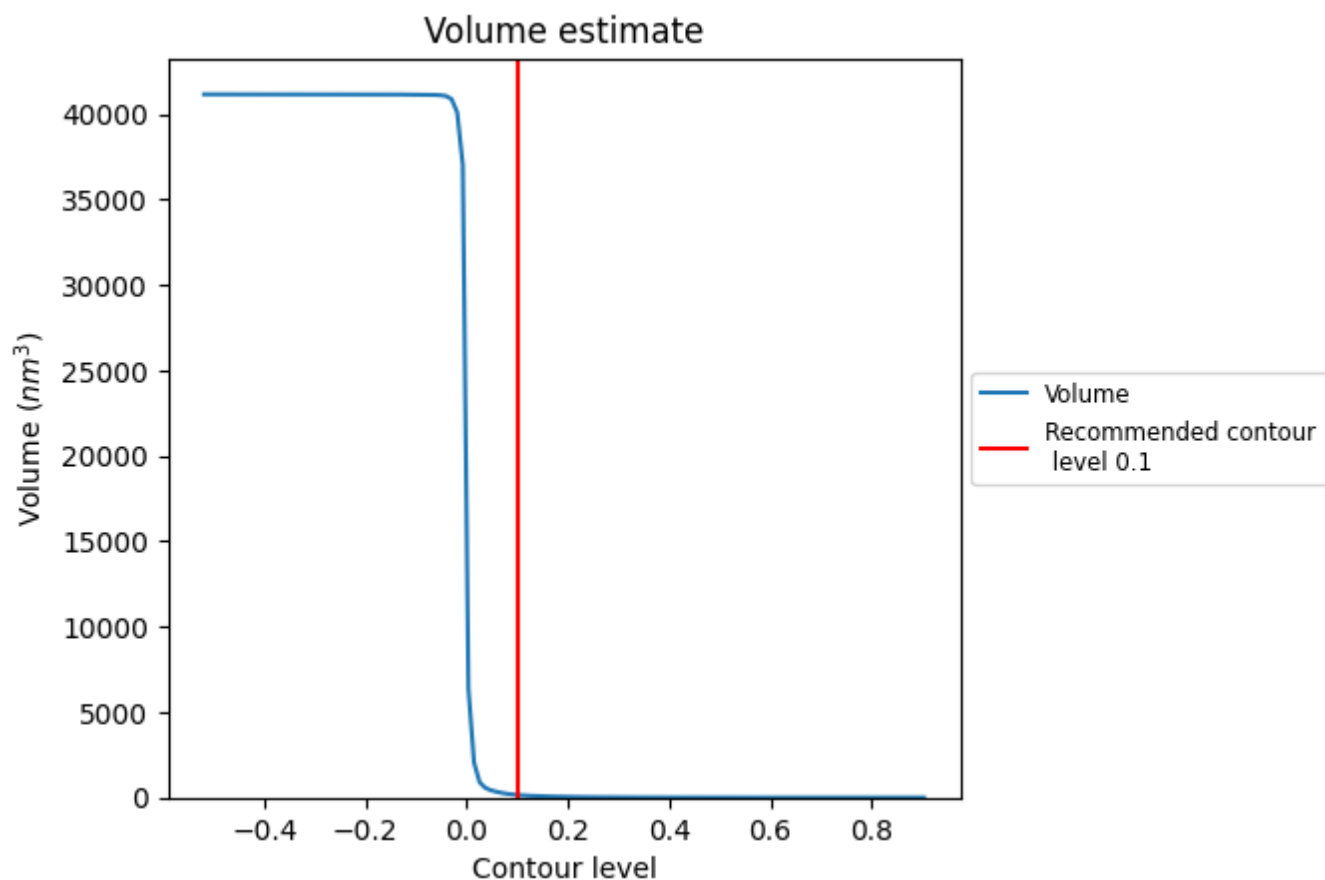
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

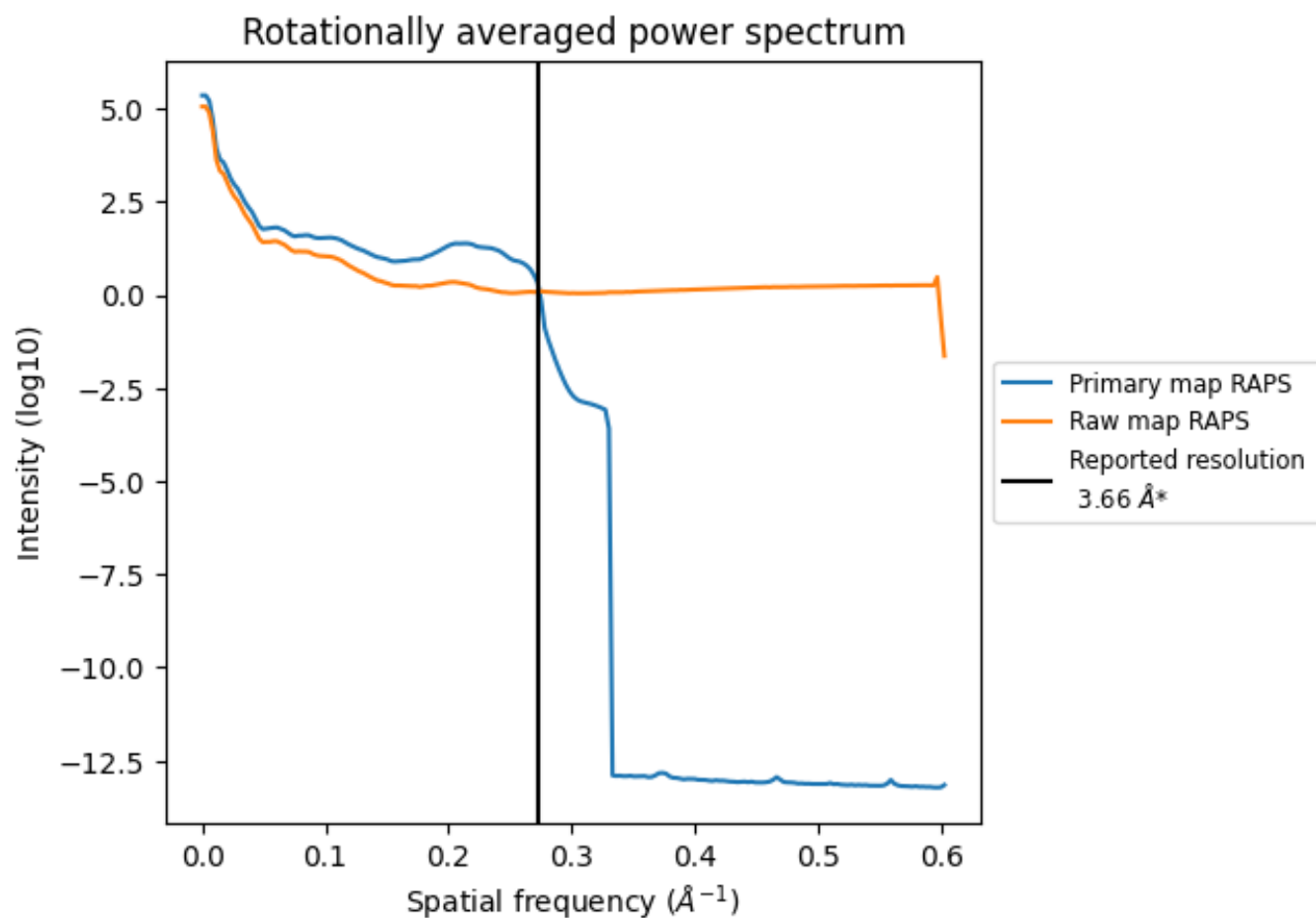
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 153 nm^3 ; this corresponds to an approximate mass of 138 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum ⓘ

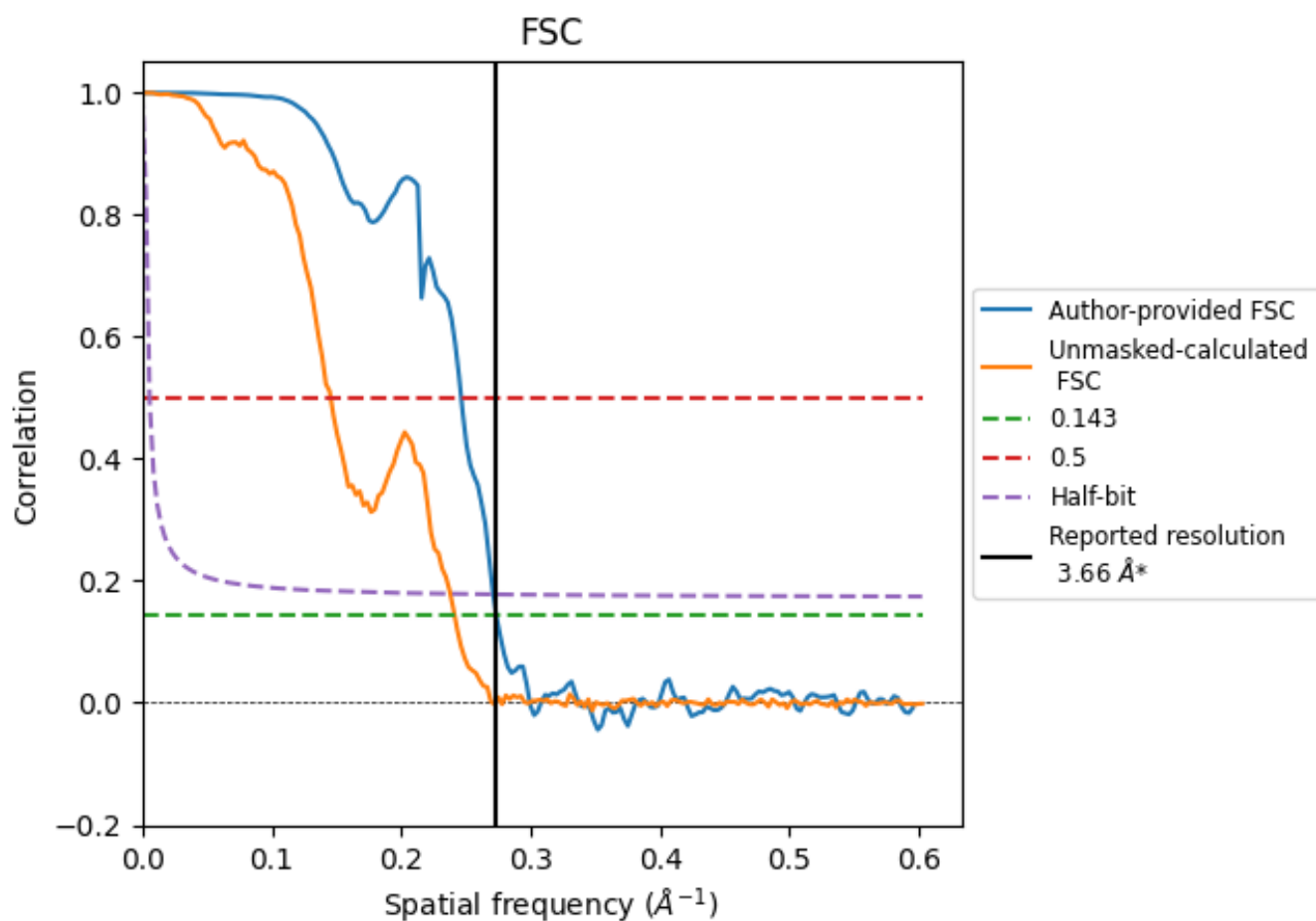


*Reported resolution corresponds to spatial frequency of 0.273 Å⁻¹

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.273 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.66	-	-
Author-provided FSC curve	3.66	4.06	3.69
Unmasked-calculated*	4.14	6.87	4.20

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 4.14 differs from the reported value 3.66 by more than 10 %

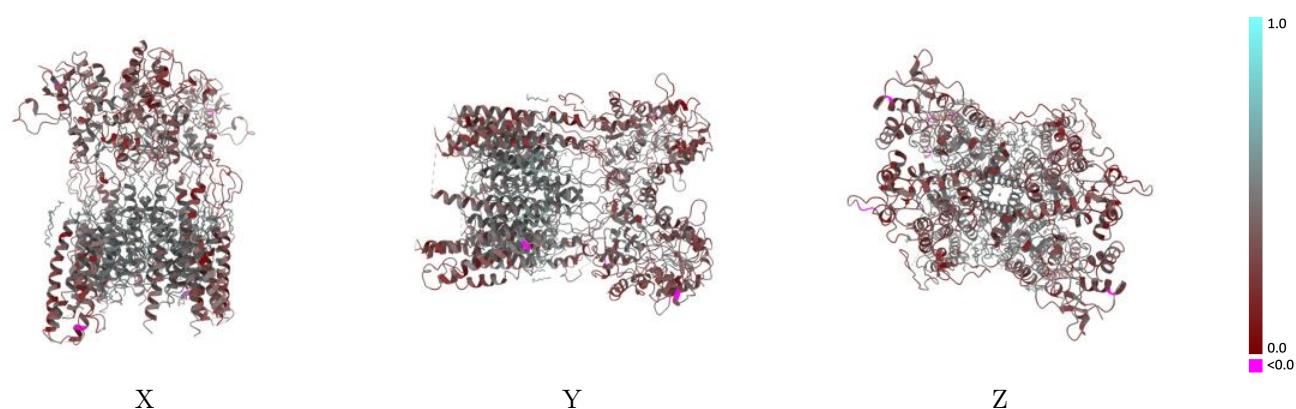
9 Map-model fit [i](#)

This section contains information regarding the fit between EMD map EMD-40750 and PDB model 8SSB. Per-residue inclusion information can be found in section 3 on page 10.

9.1 Map-model overlay [i](#)

This section was not generated.

9.2 Q-score mapped to coordinate model [i](#)

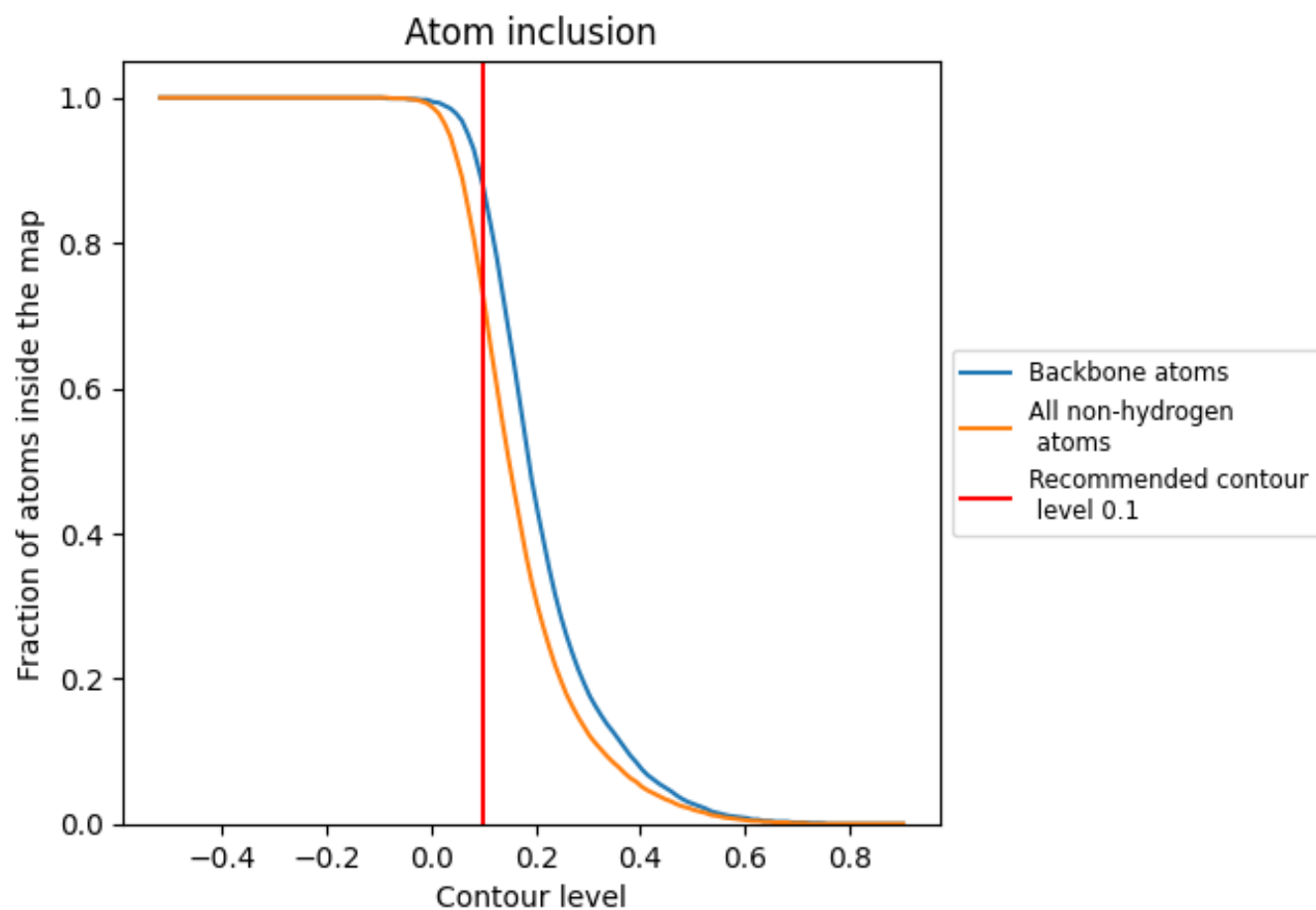


The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model [i](#)

This section was not generated.

9.4 Atom inclusion [i](#)



At the recommended contour level, 88% of all backbone atoms, 72% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary ⓘ

The table lists the average atom inclusion at the recommended contour level (0.1) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7250	0.3820
A	0.7380	0.3970
B	0.7440	0.3780
C	0.7340	0.3870
D	0.7410	0.3750
E	0.6340	0.3600
F	0.6330	0.3490

1.0

0.0

<0.0