



Full wwPDB EM Validation Report ⓘ

Mar 19, 2026 – 11:52 PM UTC

PDB ID : 7N0M / pdb_00007n0m
EMDB ID : EMD-24109
Title : Inactivated state of 2-APB-bound wildtype rat TRPV2 in nanodiscs
Authors : Pumroy, R.A.; Protopopova, A.D.; Gallo, P.N.; Moiseenkova-Bell, V.Y.
Deposited on : 2021-05-25
Resolution : 3.50 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 : **NOT EXECUTED**
MapQ : **FAILED**
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

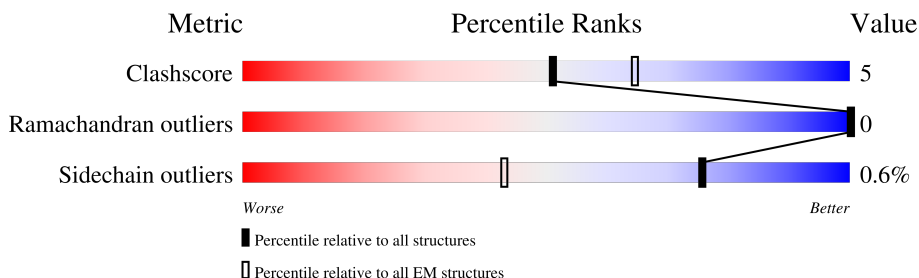
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.50 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)
Clashscore	229148	23984
Ramachandran outliers	224038	23583
Sidechain outliers	223484	23102

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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	761	69% 12% 19%
1	B	761	69% 13% 19%
1	C	761	69% 12% 19%
1	D	761	70% 12% 19%

2 Entry composition [i](#)

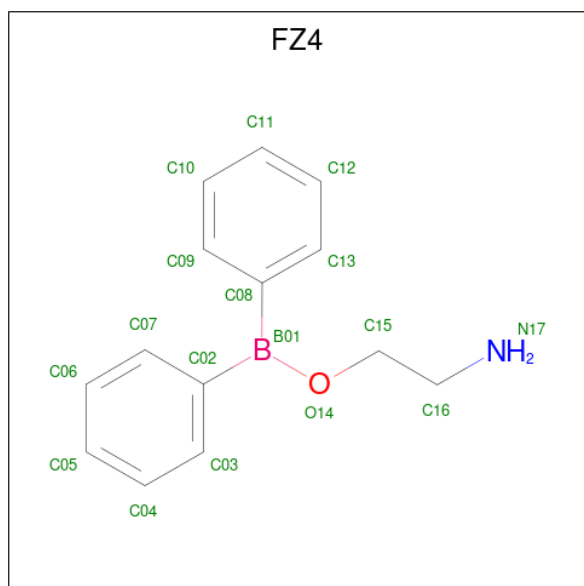
There are 2 unique types of molecules in this entry. The entry contains 19780 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Transient receptor potential cation channel subfamily V member 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	619	Total	C	N	O	S	0	0
			4928	3233	825	843	27		
1	B	619	Total	C	N	O	S	0	0
			4928	3233	825	843	27		
1	C	619	Total	C	N	O	S	0	0
			4928	3233	825	843	27		
1	D	619	Total	C	N	O	S	0	0
			4928	3233	825	843	27		

- Molecule 2 is 2-aminoethyl diphenylborinate (CCD ID: FZ4) (formula: C₁₄H₁₆BNO) (labeled as "Ligand of Interest" by depositor).

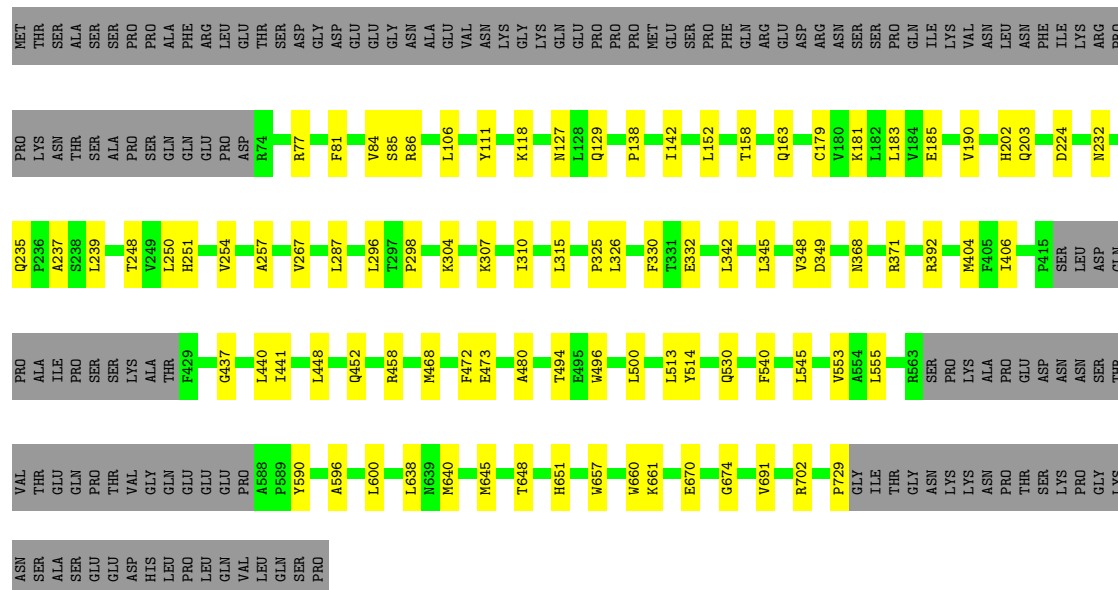


Mol	Chain	Residues	Atoms					AltConf
2	A	1	Total	B	C	N	O	0
			17	1	14	1	1	
2	A	1	Total	B	C	N	O	0
			17	1	14	1	1	

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Mol	Chain	Residues	Atoms					AltConf
2	C	1	Total	B	C	N	O	0
			17	1	14	1	1	
2	C	1	Total	B	C	N	O	0
			17	1	14	1	1	



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C4	Depositor
Number of particles used	92948	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 ($e^-/\text{\AA}^2$)	42	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: FZ4

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	0.36	0/5054	0.56	2/6872 (0.0%)
1	B	0.36	0/5054	0.56	2/6872 (0.0%)
1	C	0.36	0/5054	0.56	2/6872 (0.0%)
1	D	0.36	0/5054	0.56	2/6872 (0.0%)
All	All	0.36	0/20216	0.56	8/27488 (0.0%)

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	348	VAL	N-CA-C	-5.99	107.66	113.53
1	B	348	VAL	N-CA-C	-5.99	107.66	113.53
1	A	348	VAL	N-CA-C	-5.98	107.67	113.53
1	D	348	VAL	N-CA-C	-5.96	107.69	113.53
1	B	404	MET	CB-CG-SD	-5.02	97.64	112.70
1	D	404	MET	CB-CG-SD	-5.02	97.65	112.70
1	A	404	MET	CB-CG-SD	-5.01	97.67	112.70
1	C	404	MET	CB-CG-SD	-5.01	97.68	112.70

There are no chirality outliers.

There are no planarity outliers.

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	4928	0	4936	59	0
1	B	4928	0	4936	61	0
1	C	4928	0	4936	58	0
1	D	4928	0	4936	54	0
2	A	34	0	0	1	0
2	C	34	0	0	1	0
All	All	19780	0	19744	216	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 5.

All (216) 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:645:MET:HG2	1:D:645:MET:HE1	1.72	0.72
1:A:648:THR:HG21	1:B:645:MET:HG3	1.73	0.70
1:A:158:THR:HA	1:A:163:GLN:HG3	1.77	0.67
1:C:158:THR:HA	1:C:163:GLN:HG3	1.77	0.66
1:B:158:THR:HA	1:B:163:GLN:HG3	1.77	0.66
1:D:158:THR:HA	1:D:163:GLN:HG3	1.77	0.65
1:D:257:ALA:HB3	1:D:307:LYS:HG2	1.80	0.64
1:A:257:ALA:HB3	1:A:307:LYS:HG2	1.80	0.64
1:D:555:LEU:HD12	1:D:600:LEU:HD21	1.79	0.64
1:B:555:LEU:HD12	1:B:600:LEU:HD21	1.79	0.63
1:A:555:LEU:HD12	1:A:600:LEU:HD21	1.79	0.63
1:B:494:THR:HG23	1:B:496:TRP:H	1.64	0.63
1:C:555:LEU:HD12	1:C:600:LEU:HD21	1.80	0.63
1:B:257:ALA:HB3	1:B:307:LYS:HG2	1.80	0.62
1:C:494:THR:HG23	1:C:496:TRP:H	1.64	0.62
1:D:494:THR:HG23	1:D:496:TRP:H	1.64	0.62
1:A:494:THR:HG23	1:A:496:TRP:H	1.64	0.62
1:C:257:ALA:HB3	1:C:307:LYS:HG2	1.80	0.62
1:C:540:PHE:HA	1:C:640:MET:HE3	1.84	0.60
1:A:540:PHE:HA	1:A:640:MET:HE3	1.84	0.60
1:D:540:PHE:HA	1:D:640:MET:HE3	1.84	0.59
1:B:540:PHE:HA	1:B:640:MET:HE3	1.84	0.58
1:C:648:THR:HG21	1:D:645:MET:HG3	1.86	0.57
1:B:458:ARG:NH2	1:B:670:GLU:OE1	2.38	0.57
1:C:458:ARG:NH2	1:C:670:GLU:OE1	2.38	0.57
1:A:458:ARG:NH2	1:A:670:GLU:OE1	2.38	0.56
1:A:106:LEU:HD22	1:A:152:LEU:HD11	1.87	0.56
1:B:106:LEU:HD22	1:B:152:LEU:HD11	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:648:THR:HG21	1:C:645:MET:HG3	1.87	0.56
1:D:458:ARG:NH2	1:D:670:GLU:OE1	2.38	0.56
1:A:533:ILE:HD13	1:B:639:ASN:HB2	1.88	0.56
1:C:106:LEU:HD22	1:C:152:LEU:HD11	1.88	0.55
1:A:267:VAL:HG22	1:A:310:ILE:HD12	1.88	0.55
1:D:267:VAL:HG22	1:D:310:ILE:HD12	1.88	0.55
1:D:106:LEU:HD22	1:D:152:LEU:HD11	1.88	0.55
1:A:325:PRO:O	1:A:702:ARG:NH2	2.40	0.55
1:C:267:VAL:HG22	1:C:310:ILE:HD12	1.89	0.55
1:C:325:PRO:O	1:C:702:ARG:NH2	2.40	0.55
1:D:325:PRO:O	1:D:702:ARG:NH2	2.40	0.55
1:C:590:TYR:HB3	1:C:596:ALA:HB2	1.90	0.54
1:B:645:MET:HE1	1:C:645:MET:HG2	1.89	0.54
1:C:452:GLN:NE2	1:C:473:GLU:OE2	2.34	0.54
1:A:645:MET:HE1	1:B:645:MET:HG2	1.89	0.54
1:B:590:TYR:HB3	1:B:596:ALA:HB2	1.90	0.54
1:A:452:GLN:NE2	1:A:473:GLU:OE2	2.34	0.53
1:B:267:VAL:HG22	1:B:310:ILE:HD12	1.89	0.53
1:B:325:PRO:O	1:B:702:ARG:NH2	2.40	0.53
1:D:468:MET:O	1:D:530:GLN:NE2	2.40	0.53
1:B:468:MET:O	1:B:530:GLN:NE2	2.40	0.53
1:C:645:MET:HE1	1:D:645:MET:HG2	1.90	0.53
1:C:232:ASN:OD1	1:C:235:GLN:N	2.38	0.53
1:A:232:ASN:OD1	1:A:235:GLN:N	2.38	0.53
1:A:590:TYR:HB3	1:A:596:ALA:HB2	1.90	0.53
1:D:590:TYR:HB3	1:D:596:ALA:HB2	1.90	0.53
1:A:224:ASP:N	1:A:224:ASP:OD1	2.42	0.52
1:C:468:MET:O	1:C:530:GLN:NE2	2.40	0.52
1:C:224:ASP:N	1:C:224:ASP:OD1	2.42	0.52
1:D:224:ASP:OD1	1:D:224:ASP:N	2.42	0.52
1:B:224:ASP:N	1:B:224:ASP:OD1	2.42	0.52
1:D:452:GLN:NE2	1:D:473:GLU:OE2	2.34	0.52
1:C:202:HIS:ND1	1:C:203:GLN:O	2.44	0.51
1:D:472:PHE:CE2	1:D:514:TYR:HB2	2.46	0.50
1:C:472:PHE:CE2	1:C:514:TYR:HB2	2.46	0.50
1:A:472:PHE:CE2	1:A:514:TYR:HB2	2.46	0.50
1:B:472:PHE:CE2	1:B:514:TYR:HB2	2.46	0.50
1:D:202:HIS:ND1	1:D:203:GLN:O	2.44	0.50
1:D:232:ASN:OD1	1:D:235:GLN:N	2.38	0.49
1:B:202:HIS:ND1	1:B:203:GLN:O	2.44	0.49
1:B:452:GLN:NE2	1:B:473:GLU:OE2	2.34	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:408:THR:HG23	1:B:553:VAL:HG11	1.94	0.49
1:A:502:VAL:HG13	1:B:554:ALA:HB1	1.94	0.49
1:C:345:LEU:HD23	1:C:349:ASP:HB3	1.95	0.49
1:A:729:PRO:HD2	1:B:205:THR:HG21	1.95	0.49
1:A:202:HIS:ND1	1:A:203:GLN:O	2.44	0.48
1:D:345:LEU:HD23	1:D:349:ASP:HB3	1.95	0.48
1:A:510:LEU:HD11	1:B:632:LEU:HD11	1.95	0.48
1:A:296:LEU:HD11	1:A:304:LYS:HD3	1.96	0.48
1:D:296:LEU:HD11	1:D:304:LYS:HD3	1.96	0.48
1:A:345:LEU:HD23	1:A:349:ASP:HB3	1.95	0.48
1:A:645:MET:HG3	1:D:648:THR:HG21	1.96	0.48
1:B:345:LEU:HD23	1:B:349:ASP:HB3	1.95	0.47
1:C:296:LEU:HD11	1:C:304:LYS:HD3	1.96	0.47
1:A:77:ARG:HG3	1:A:111:TYR:HA	1.97	0.47
1:D:118:LYS:HB2	1:D:118:LYS:HE2	1.64	0.47
1:B:296:LEU:HD11	1:B:304:LYS:HD3	1.96	0.47
1:D:77:ARG:HG3	1:D:111:TYR:HA	1.97	0.47
1:D:190:VAL:HG11	1:D:237:ALA:HB2	1.97	0.47
1:B:190:VAL:HG11	1:B:237:ALA:HB2	1.96	0.47
1:B:232:ASN:OD1	1:B:235:GLN:N	2.38	0.47
1:C:190:VAL:HG11	1:C:237:ALA:HB2	1.96	0.47
1:D:86:ARG:O	1:D:129:GLN:NE2	2.46	0.47
1:A:190:VAL:HG11	1:A:237:ALA:HB2	1.96	0.46
1:C:86:ARG:O	1:C:129:GLN:NE2	2.46	0.46
1:A:118:LYS:HB2	1:A:118:LYS:HE2	1.64	0.46
1:C:77:ARG:HG3	1:C:111:TYR:HA	1.97	0.46
1:C:406:ILE:HD11	1:C:440:LEU:HD22	1.98	0.46
1:A:332:GLU:HB2	1:A:342:LEU:HD13	1.98	0.46
1:B:118:LYS:HE2	1:B:118:LYS:HB2	1.64	0.46
1:D:406:ILE:HD11	1:D:440:LEU:HD22	1.98	0.46
1:B:332:GLU:HB2	1:B:342:LEU:HD13	1.98	0.46
1:B:406:ILE:HD11	1:B:440:LEU:HD22	1.98	0.46
1:A:468:MET:O	1:A:530:GLN:NE2	2.40	0.46
1:B:77:ARG:HG3	1:B:111:TYR:HA	1.97	0.46
1:C:81:PHE:HA	1:C:84:VAL:HG12	1.98	0.46
1:D:81:PHE:HA	1:D:84:VAL:HG12	1.98	0.46
1:A:406:ILE:HD11	1:A:440:LEU:HD22	1.98	0.45
1:C:254:VAL:HG22	1:C:310:ILE:HD13	1.99	0.45
1:B:81:PHE:HA	1:B:84:VAL:HG12	1.98	0.45
1:D:138:PRO:O	1:D:142:ILE:HG12	2.17	0.45
1:D:254:VAL:HG22	1:D:310:ILE:HD13	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:392:ARG:HH22	1:A:674:GLY:HA2	1.81	0.45
1:A:478:LEU:HD23	1:A:478:LEU:HA	1.85	0.45
1:B:406:ILE:HG13	1:B:441:ILE:HD11	1.99	0.45
1:C:392:ARG:HH22	1:C:674:GLY:HA2	1.81	0.45
1:C:138:PRO:O	1:C:142:ILE:HG12	2.17	0.45
1:C:332:GLU:HB2	1:C:342:LEU:HD13	1.98	0.45
1:D:392:ARG:HH22	1:D:674:GLY:HA2	1.81	0.45
1:A:205:THR:HG21	1:D:729:PRO:HD2	1.97	0.45
1:D:332:GLU:HB2	1:D:342:LEU:HD13	1.98	0.45
1:A:406:ILE:HG13	1:A:441:ILE:HD11	1.99	0.44
1:B:86:ARG:O	1:B:129:GLN:NE2	2.46	0.44
1:B:254:VAL:HG22	1:B:310:ILE:HD13	1.99	0.44
1:B:392:ARG:HH22	1:B:674:GLY:HA2	1.81	0.44
1:A:138:PRO:O	1:A:142:ILE:HG12	2.17	0.44
1:A:81:PHE:HA	1:A:84:VAL:HG12	1.98	0.44
1:C:406:ILE:HG13	1:C:441:ILE:HD11	1.99	0.44
1:B:209:PHE:O	1:B:215:SER:OG	2.29	0.44
1:A:254:VAL:HG22	1:A:310:ILE:HD13	1.99	0.44
1:C:118:LYS:HB2	1:C:118:LYS:HE2	1.64	0.44
1:A:181:LYS:O	1:A:185:GLU:HG2	2.18	0.44
1:B:138:PRO:O	1:B:142:ILE:HG12	2.17	0.44
1:C:209:PHE:O	1:C:215:SER:OG	2.29	0.44
1:B:408:THR:HG23	1:C:553:VAL:HG11	2.00	0.43
1:B:181:LYS:O	1:B:185:GLU:HG2	2.18	0.43
1:D:406:ILE:HG13	1:D:441:ILE:HD11	1.99	0.43
1:B:251:HIS:NE2	1:B:298:PRO:HD3	2.34	0.43
1:C:181:LYS:O	1:C:185:GLU:HG2	2.18	0.43
1:D:251:HIS:NE2	1:D:298:PRO:HD3	2.34	0.43
1:C:326:LEU:H	1:C:326:LEU:HG	1.62	0.43
1:C:626:LEU:HD23	1:C:626:LEU:HA	1.84	0.43
1:A:86:ARG:O	1:A:129:GLN:NE2	2.46	0.43
1:B:85:SER:HA	1:B:127:ASN:HD21	1.84	0.43
1:C:251:HIS:NE2	1:C:298:PRO:HD3	2.34	0.43
1:C:408:THR:HG23	1:D:553:VAL:HG11	2.01	0.43
1:D:181:LYS:O	1:D:185:GLU:HG2	2.18	0.43
1:A:251:HIS:NE2	1:A:298:PRO:HD3	2.34	0.43
1:C:85:SER:HA	1:C:127:ASN:HD21	1.84	0.43
1:A:239:LEU:HD22	1:A:287:LEU:HD13	2.00	0.43
1:A:357:LEU:HD23	1:A:357:LEU:HA	1.85	0.43
1:A:245:LEU:HD12	1:A:245:LEU:HA	1.90	0.42
1:A:248:THR:HG22	1:A:251:HIS:CD2	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:181:LYS:HE3	1:C:181:LYS:HB2	1.86	0.42
1:A:85:SER:HA	1:A:127:ASN:HD21	1.84	0.42
1:A:209:PHE:O	1:A:215:SER:OG	2.29	0.42
1:B:248:THR:HG22	1:B:251:HIS:CD2	2.54	0.42
1:B:368:ASN:OD1	1:B:371:ARG:NH1	2.43	0.42
1:B:239:LEU:HD22	1:B:287:LEU:HD13	2.00	0.42
1:D:368:ASN:OD1	1:D:371:ARG:NH1	2.43	0.42
1:B:507:LEU:HD12	1:B:507:LEU:HA	1.88	0.42
1:D:248:THR:HG22	1:D:251:HIS:CD2	2.54	0.42
1:A:335:TYR:CE1	1:B:173:GLU:HG3	2.55	0.42
1:B:330:PHE:CZ	1:B:691:VAL:HG11	2.55	0.42
1:D:239:LEU:HD22	1:D:287:LEU:HD13	2.00	0.42
1:A:315:LEU:HA	1:A:315:LEU:HD12	1.86	0.42
1:C:239:LEU:HD22	1:C:287:LEU:HD13	2.00	0.42
1:A:248:THR:HG23	1:A:250:LEU:H	1.85	0.42
1:A:330:PHE:CZ	1:A:691:VAL:HG11	2.55	0.41
1:A:221:LYS:HE2	1:A:221:LYS:HB3	1.88	0.41
1:B:248:THR:HG23	1:B:250:LEU:H	1.85	0.41
1:B:251:HIS:CD2	1:B:298:PRO:HD3	2.55	0.41
1:C:248:THR:HG22	1:C:251:HIS:CD2	2.55	0.41
1:C:478:LEU:HD23	1:C:478:LEU:HA	1.85	0.41
1:D:85:SER:HA	1:D:127:ASN:HD21	1.84	0.41
1:C:251:HIS:CD2	1:C:298:PRO:HD3	2.55	0.41
1:C:330:PHE:CZ	1:C:691:VAL:HG11	2.55	0.41
1:D:248:THR:HG23	1:D:250:LEU:H	1.85	0.41
1:C:248:THR:HG23	1:C:250:LEU:H	1.85	0.41
1:C:437:GLY:O	1:C:441:ILE:HG12	2.21	0.41
1:C:507:LEU:HD12	1:C:507:LEU:HA	1.88	0.41
2:C:802:FZ4:C07	2:C:802:FZ4:C08	2.97	0.41
1:D:437:GLY:O	1:D:441:ILE:HG12	2.20	0.41
1:D:540:PHE:HD1	1:D:640:MET:HG3	1.86	0.41
1:B:245:LEU:HD12	1:B:245:LEU:HA	1.90	0.41
1:B:555:LEU:HD23	1:B:555:LEU:HA	1.87	0.41
1:C:448:LEU:HD22	1:C:480:ALA:HB2	2.03	0.41
1:D:330:PHE:CZ	1:D:691:VAL:HG11	2.55	0.41
1:D:448:LEU:HD22	1:D:480:ALA:HB2	2.03	0.41
1:B:513:LEU:HD23	1:B:513:LEU:HA	1.91	0.41
1:B:540:PHE:HD1	1:B:640:MET:HG3	1.86	0.41
1:B:657:TRP:HA	1:B:660:TRP:CD1	2.56	0.41
1:C:513:LEU:HD23	1:C:513:LEU:HA	1.91	0.41
1:C:634:TYR:HD1	1:C:634:TYR:HA	1.73	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:251:HIS:CD2	1:D:298:PRO:HD3	2.55	0.41
1:D:513:LEU:HD23	1:D:513:LEU:HA	1.91	0.41
1:A:437:GLY:O	1:A:441:ILE:HG12	2.21	0.41
2:A:801:FZ4:C07	2:A:801:FZ4:C08	2.97	0.41
1:B:533:ILE:HD13	1:C:639:ASN:HB2	2.03	0.41
1:C:97:GLU:O	1:C:101:TRP:HB2	2.21	0.41
1:C:657:TRP:HA	1:C:660:TRP:CD1	2.56	0.41
1:D:315:LEU:HD12	1:D:315:LEU:HA	1.86	0.41
1:B:453:LEU:HD23	1:B:453:LEU:HA	1.93	0.41
1:A:657:TRP:HA	1:A:660:TRP:CD1	2.56	0.40
1:D:500:LEU:HD23	1:D:500:LEU:HA	1.93	0.40
1:B:97:GLU:O	1:B:101:TRP:HB2	2.21	0.40
1:C:386:TRP:CE2	1:C:664:LYS:HE3	2.57	0.40
1:D:179:CYS:O	1:D:183:LEU:HB2	2.22	0.40
1:A:179:CYS:O	1:A:183:LEU:HB2	2.22	0.40
1:A:251:HIS:CD2	1:A:298:PRO:HD3	2.55	0.40
1:B:386:TRP:CE2	1:B:664:LYS:HE3	2.57	0.40
1:B:528:MET:O	1:B:532:VAL:N	2.46	0.40
1:C:694:ARG:HD2	1:C:694:ARG:HA	1.93	0.40
1:D:545:LEU:HD12	1:D:545:LEU:HA	1.88	0.40
1:D:657:TRP:HA	1:D:660:TRP:CD1	2.56	0.40
1:D:661:LYS:HA	1:D:661:LYS:HD3	1.92	0.40
1:A:386:TRP:CE2	1:A:664:LYS:HE3	2.57	0.40
1:A:725:ASP:OD1	1:A:725:ASP: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	613/761 (81%)	579 (94%)	34 (6%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	613/761 (81%)	580 (95%)	33 (5%)	0	100	100
1	C	613/761 (81%)	580 (95%)	33 (5%)	0	100	100
1	D	613/761 (81%)	580 (95%)	33 (5%)	0	100	100
All	All	2452/3044 (81%)	2319 (95%)	133 (5%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all 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	523/676 (77%)	520 (99%)	3 (1%)	78	79
1	B	523/676 (77%)	520 (99%)	3 (1%)	78	79
1	C	523/676 (77%)	520 (99%)	3 (1%)	78	79
1	D	523/676 (77%)	520 (99%)	3 (1%)	78	79
All	All	2092/2704 (77%)	2080 (99%)	12 (1%)	76	79

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	326	LEU
1	A	638	LEU
1	A	651	HIS
1	B	326	LEU
1	B	638	LEU
1	B	651	HIS
1	C	326	LEU
1	C	638	LEU
1	C	651	HIS
1	D	326	LEU
1	D	638	LEU
1	D	651	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (33) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	141	GLN
1	A	186	ASN
1	A	251	HIS
1	A	276	GLN
1	A	293	HIS
1	A	324	GLN
1	A	383	GLN
1	A	413	HIS
1	A	521	HIS
1	B	141	GLN
1	B	186	ASN
1	B	251	HIS
1	B	276	GLN
1	B	293	HIS
1	B	324	GLN
1	B	383	GLN
1	B	413	HIS
1	C	141	GLN
1	C	186	ASN
1	C	251	HIS
1	C	276	GLN
1	C	293	HIS
1	C	324	GLN
1	C	383	GLN
1	C	413	HIS
1	D	186	ASN
1	D	251	HIS
1	D	276	GLN
1	D	293	HIS
1	D	324	GLN
1	D	383	GLN
1	D	413	HIS
1	D	414	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

4 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	FZ4	C	801	-	17,18,18	0.63	0	18,22,22	1.15	2 (11%)
2	FZ4	C	802	-	17,18,18	0.61	0	18,22,22	1.14	1 (5%)
2	FZ4	A	801	-	17,18,18	0.62	0	18,22,22	1.15	2 (11%)
2	FZ4	A	802	-	17,18,18	0.63	0	18,22,22	1.15	2 (11%)

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	FZ4	C	801	-	-	9/12/12/12	0/2/2/2
2	FZ4	C	802	-	-	9/12/12/12	0/2/2/2
2	FZ4	A	801	-	-	9/12/12/12	0/2/2/2
2	FZ4	A	802	-	-	9/12/12/12	0/2/2/2

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	801	FZ4	B01-C02-C07	-2.39	118.05	121.51
2	A	801	FZ4	B01-C02-C07	-2.38	118.07	121.51
2	C	802	FZ4	B01-C02-C07	-2.37	118.08	121.51
2	A	802	FZ4	B01-C02-C07	-2.36	118.10	121.51
2	A	801	FZ4	C07-C02-C03	2.03	119.20	116.86
2	A	802	FZ4	C07-C02-C03	2.02	119.17	116.86
2	C	801	FZ4	C07-C02-C03	2.02	119.17	116.86

There are no chirality outliers.

All (36) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	801	FZ4	C02-B01-O14-C15
2	A	801	FZ4	O14-C15-C16-N17
2	A	801	FZ4	C16-C15-O14-B01
2	A	802	FZ4	C02-B01-O14-C15
2	A	802	FZ4	O14-C15-C16-N17
2	A	802	FZ4	C16-C15-O14-B01
2	C	801	FZ4	C02-B01-O14-C15
2	C	801	FZ4	O14-C15-C16-N17
2	C	801	FZ4	C16-C15-O14-B01
2	C	802	FZ4	C02-B01-O14-C15
2	C	802	FZ4	O14-C15-C16-N17
2	C	802	FZ4	C16-C15-O14-B01
2	A	801	FZ4	C02-B01-C08-C09
2	A	801	FZ4	C02-B01-C08-C13
2	A	802	FZ4	C02-B01-C08-C09
2	A	802	FZ4	C02-B01-C08-C13
2	C	801	FZ4	C02-B01-C08-C09
2	C	801	FZ4	C02-B01-C08-C13
2	C	802	FZ4	C02-B01-C08-C09
2	C	802	FZ4	C02-B01-C08-C13
2	A	801	FZ4	O14-B01-C02-C03
2	A	801	FZ4	O14-B01-C02-C07
2	A	801	FZ4	O14-B01-C08-C09
2	A	801	FZ4	O14-B01-C08-C13
2	A	802	FZ4	O14-B01-C02-C03
2	A	802	FZ4	O14-B01-C02-C07
2	A	802	FZ4	O14-B01-C08-C09
2	A	802	FZ4	O14-B01-C08-C13
2	C	801	FZ4	O14-B01-C02-C03
2	C	801	FZ4	O14-B01-C02-C07
2	C	801	FZ4	O14-B01-C08-C09

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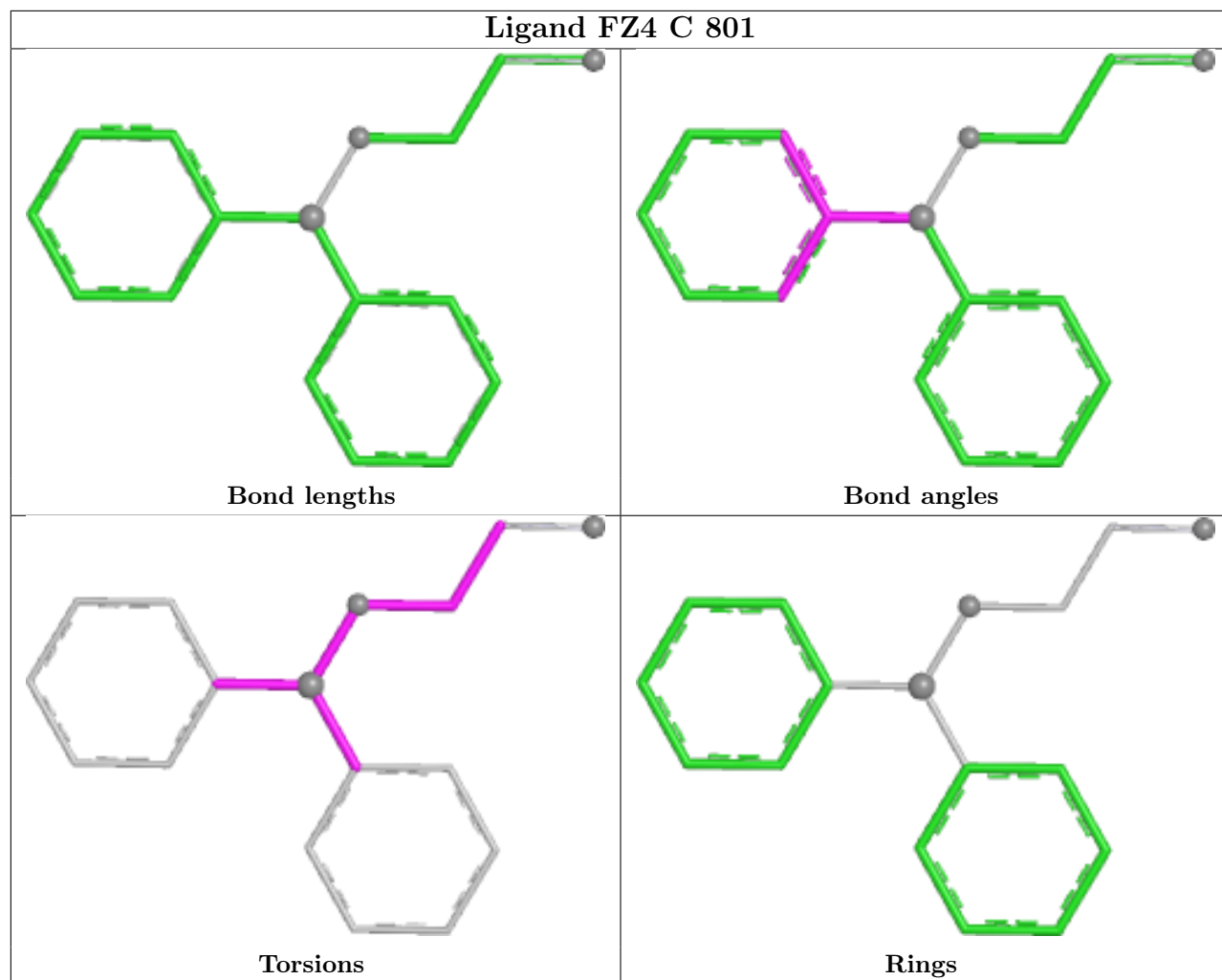
Mol	Chain	Res	Type	Atoms
2	C	801	FZ4	O14-B01-C08-C13
2	C	802	FZ4	O14-B01-C02-C03
2	C	802	FZ4	O14-B01-C02-C07
2	C	802	FZ4	O14-B01-C08-C09
2	C	802	FZ4	O14-B01-C08-C13

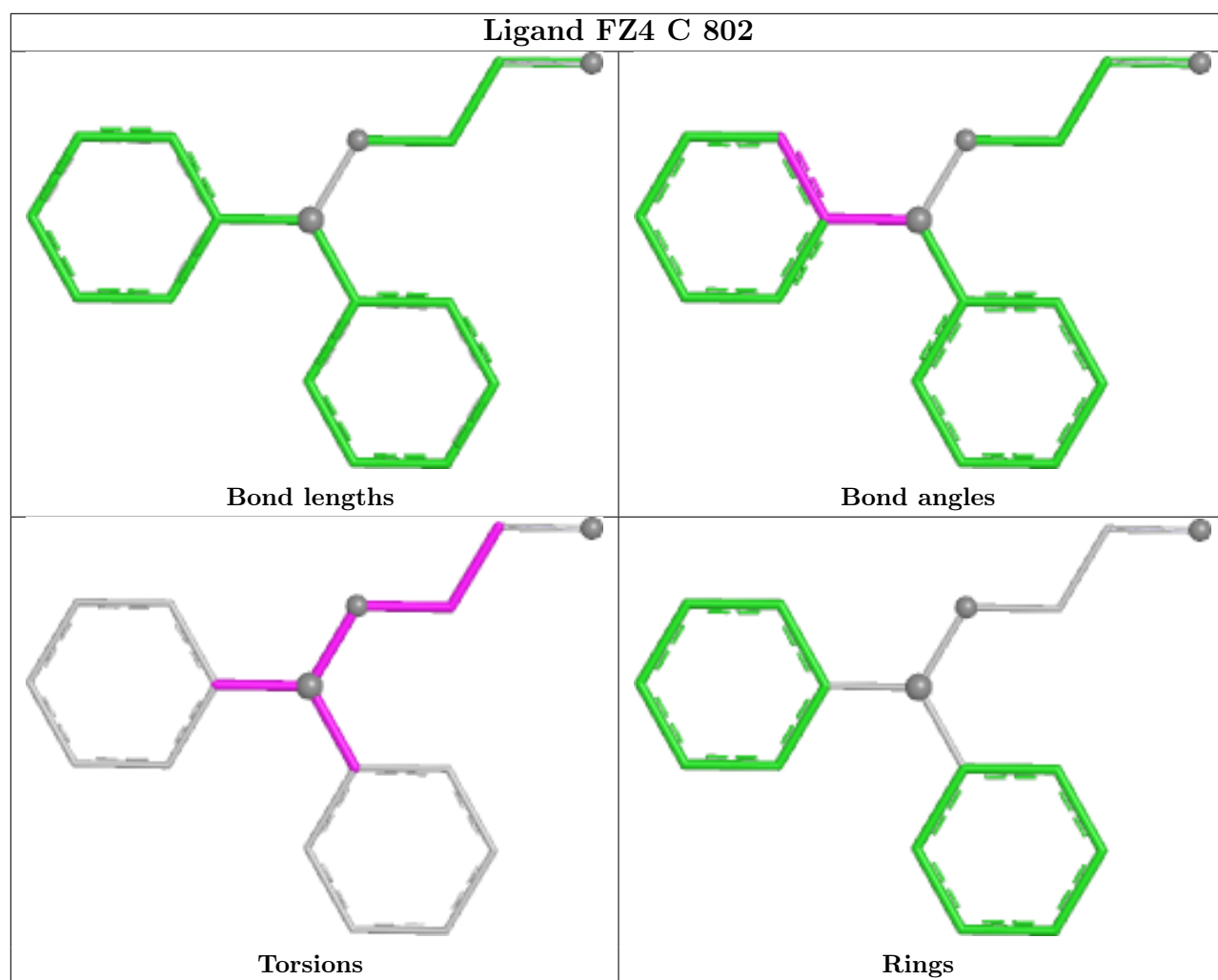
There are no ring outliers.

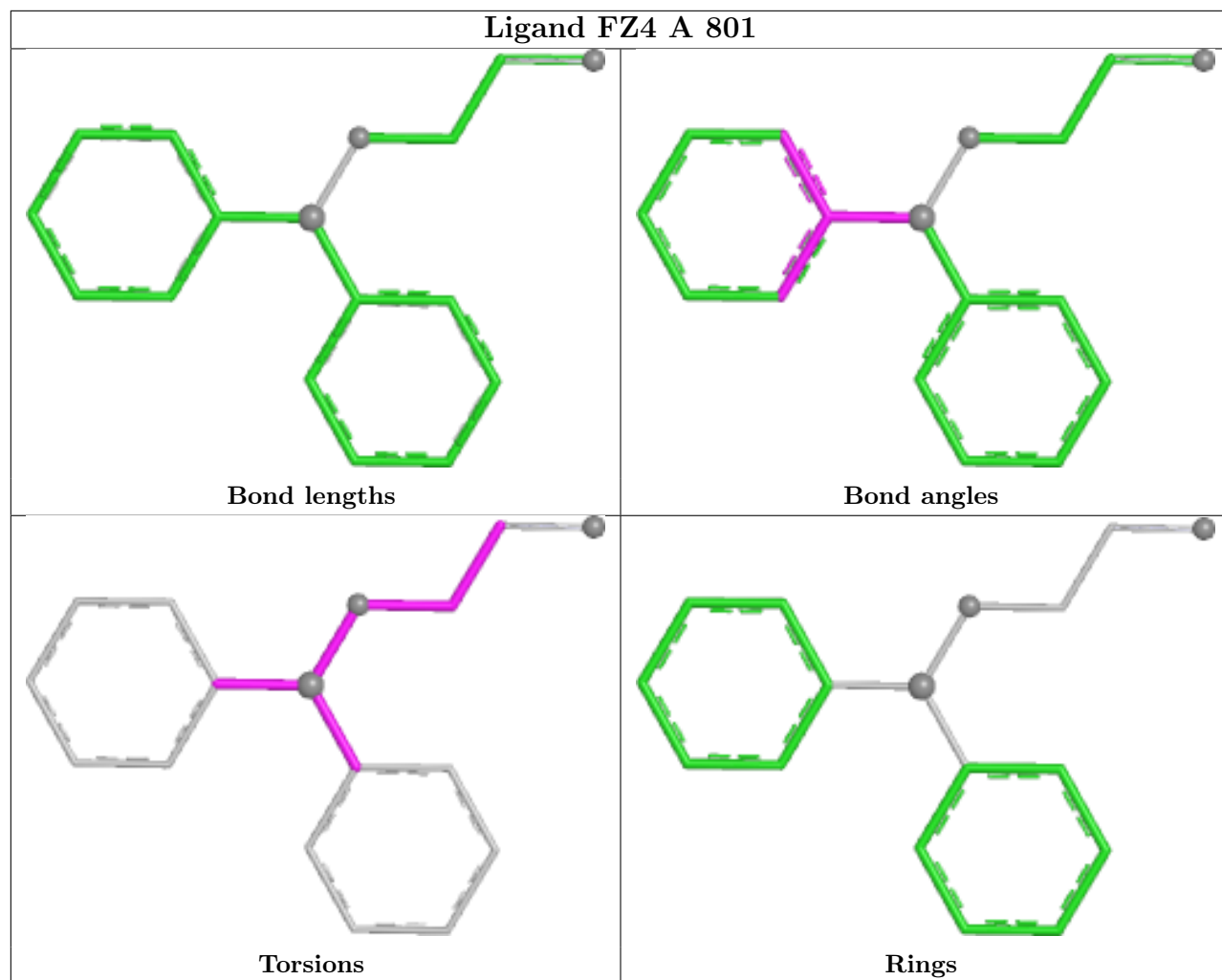
2 monomers are involved in 2 short contacts:

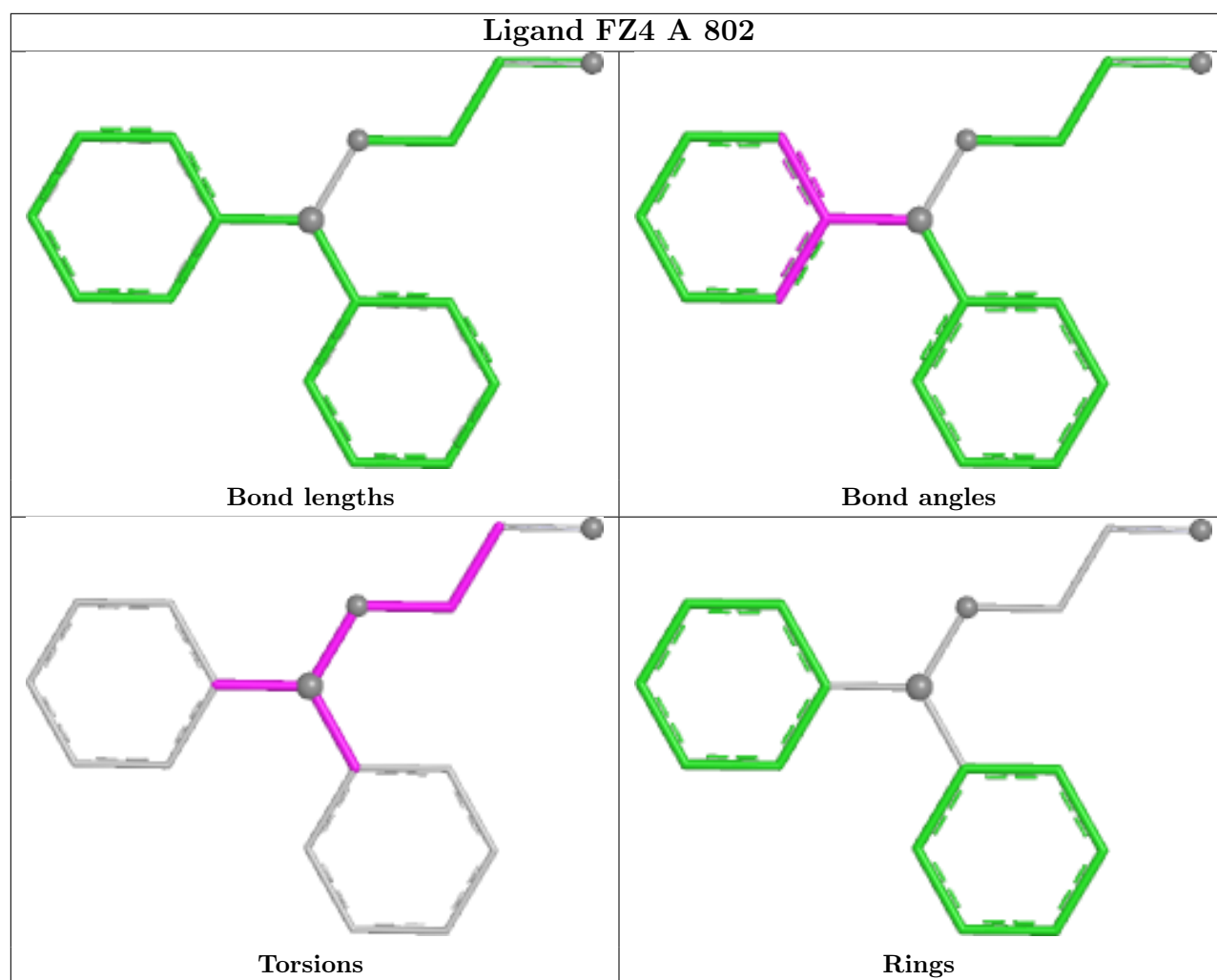
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	802	FZ4	1	0
2	A	801	FZ4	1	0

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









5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Map visualisation

This section contains visualisations of the EMDB entry EMD-24109. These allow visual inspection of the internal detail of the map and identification of artifacts.

No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections

This section was not generated.

6.2 Central slices

This section was not generated.

6.3 Largest variance slices

This section was not generated.

6.4 Orthogonal standard-deviation projections (False-color)

This section was not generated.

6.5 Orthogonal surface views

This section was not generated.

6.6 Mask visualisation

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

7 Map analysis ⓘ

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

7.1 Map-value distribution ⓘ

This section was not generated.

7.2 Volume estimate versus contour level ⓘ

This section was not generated.

7.3 Rotationally averaged power spectrum ⓘ

This section was not generated. The rotationally averaged power spectrum had issues being displayed.

8 Fourier-Shell correlation

This section was not generated. No FSC curve or half-maps provided.

9 Map-model fit

This section was not generated.