



wwPDB EM Validation Summary Report ⓘ

Mar 17, 2026 – 04:00 PM UTC

PDB ID : 5NG5 / pdb_00005ng5
EMDB ID : EMD-3636
Title : multi-drug efflux; membrane transport; RND superfamily; Drug resistance
Authors : Wang, Z.; Fan, G.; Hryc, C.F.; Blaza, J.N.; Serysheva, I.I.; Schmid, M.F.;
Chiu, W.; Luisi, B.F.; Du, D.
Deposited on : 2017-03-16
Resolution : 6.50 Å(reported)

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<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>
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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

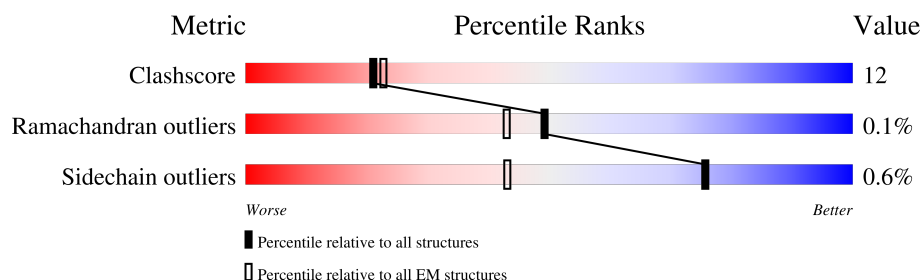
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 6.50 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	373	68% 22% • 9%
1	B	373	69% 22% • 9%
1	D	373	60% 28% • 9%
1	E	373	64% 24% • 9%
1	G	373	61% 29% • 9%
1	H	373	66% 24% • 9%
2	C	493	65% 22% 13%
2	F	493	66% 20% 13%
2	I	493	63% 24% 13%

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Mol	Chain	Length	Quality of chain
3	J	1049	
3	K	1049	
3	L	1049	
4	M	54	
4	N	54	
4	O	54	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	5QF	J	1101	-	-	X	-
5	5QF	K	1101	-	-	X	-
5	5QF	L	1101	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 49920 atoms, of which 114 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 Multidrug efflux pump subunit AcrA.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	E	340	Total	C	N	O	S	0	0
			2553	1591	451	506	5		
1	D	340	Total	C	N	O	S	0	0
			2553	1591	451	506	5		
1	G	340	Total	C	N	O	S	0	0
			2553	1591	451	506	5		
1	H	340	Total	C	N	O	S	0	0
			2553	1591	451	506	5		
1	A	340	Total	C	N	O	S	0	0
			2553	1591	451	506	5		
1	B	340	Total	C	N	O	S	0	0
			2553	1591	451	506	5		

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	223	MET	PHE	conflict	UNP P0AE06
E	224	MET	LEU	conflict	UNP P0AE06
E	287	MET	LEU	conflict	UNP P0AE06
E	288	MET	LEU	conflict	UNP P0AE06
D	223	MET	PHE	conflict	UNP P0AE06
D	224	MET	LEU	conflict	UNP P0AE06
D	287	MET	LEU	conflict	UNP P0AE06
D	288	MET	LEU	conflict	UNP P0AE06
G	223	MET	PHE	conflict	UNP P0AE06
G	224	MET	LEU	conflict	UNP P0AE06
G	287	MET	LEU	conflict	UNP P0AE06
G	288	MET	LEU	conflict	UNP P0AE06
H	223	MET	PHE	conflict	UNP P0AE06
H	224	MET	LEU	conflict	UNP P0AE06
H	287	MET	LEU	conflict	UNP P0AE06
H	288	MET	LEU	conflict	UNP P0AE06
A	223	MET	PHE	conflict	UNP P0AE06
A	224	MET	LEU	conflict	UNP P0AE06

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Chain	Residue	Modelled	Actual	Comment	Reference
A	287	MET	LEU	conflict	UNP P0AE06
A	288	MET	LEU	conflict	UNP P0AE06
B	223	MET	PHE	conflict	UNP P0AE06
B	224	MET	LEU	conflict	UNP P0AE06
B	287	MET	LEU	conflict	UNP P0AE06
B	288	MET	LEU	conflict	UNP P0AE06

- Molecule 2 is a protein called Outer membrane protein TolC.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	F	428	Total	C	N	O	S	0	0
			3304	2037	586	676	5		
2	I	428	Total	C	N	O	S	0	0
			3304	2037	586	676	5		
2	C	428	Total	C	N	O	S	0	0
			3304	2037	586	676	5		

- Molecule 3 is a protein called Multidrug efflux pump subunit AcrB.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	K	1037	Total	C	N	O	S	0	0
			7819	5032	1290	1453	44		
3	L	1037	Total	C	N	O	S	0	0
			7819	5032	1290	1453	44		
3	J	1033	Total	C	N	O	S	0	0
			7849	5052	1295	1458	44		

- Molecule 4 is a protein called Multidrug efflux pump accessory protein AcrZ.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	M	46	Total	C	N	O	S	0	0
			326	219	48	56	3		
4	N	46	Total	C	N	O	S	0	0
			326	219	48	56	3		
4	O	46	Total	C	N	O	S	0	0
			326	219	48	56	3		

There are 15 discrepancies between the modelled and reference sequences:

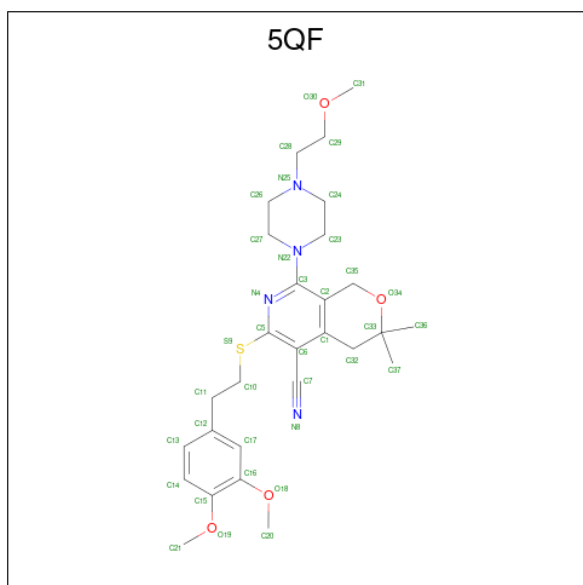
Chain	Residue	Modelled	Actual	Comment	Reference
M	50	HIS	-	expression tag	UNP P0AAW9
M	51	HIS	-	expression tag	UNP P0AAW9

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Chain	Residue	Modelled	Actual	Comment	Reference
M	52	HIS	-	expression tag	UNP P0AAW9
M	53	HIS	-	expression tag	UNP P0AAW9
M	54	HIS	-	expression tag	UNP P0AAW9
N	50	HIS	-	expression tag	UNP P0AAW9
N	51	HIS	-	expression tag	UNP P0AAW9
N	52	HIS	-	expression tag	UNP P0AAW9
N	53	HIS	-	expression tag	UNP P0AAW9
N	54	HIS	-	expression tag	UNP P0AAW9
O	50	HIS	-	expression tag	UNP P0AAW9
O	51	HIS	-	expression tag	UNP P0AAW9
O	52	HIS	-	expression tag	UNP P0AAW9
O	53	HIS	-	expression tag	UNP P0AAW9
O	54	HIS	-	expression tag	UNP P0AAW9

- Molecule 5 is 6-[2-(3,4-dimethoxyphenyl)ethylsulfanyl]-8-[4-(2-methoxyethyl)piperazin-1-yl]-3,3-dimethyl-1,4-dihydropyrano[3,4-c]pyridine-5-carbonitrile (CCD ID: 5QF) (formula: $C_{28}H_{38}N_4O_4S$).

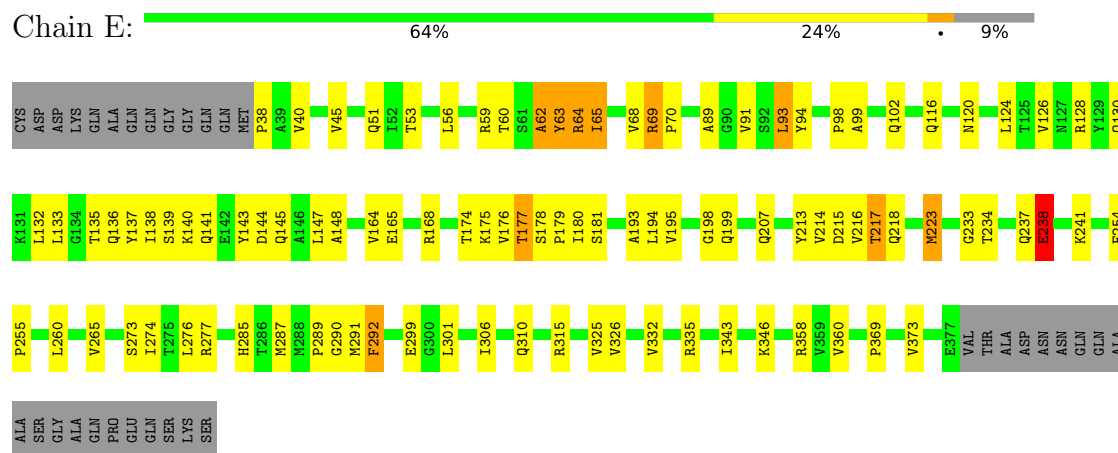


Mol	Chain	Residues	Atoms						AltConf
5	K	1	Total	C	H	N	O	S	0
			75	28	38	4	4	1	
5	L	1	Total	C	H	N	O	S	0
			75	28	38	4	4	1	
5	J	1	Total	C	H	N	O	S	0
			75	28	38	4	4	1	

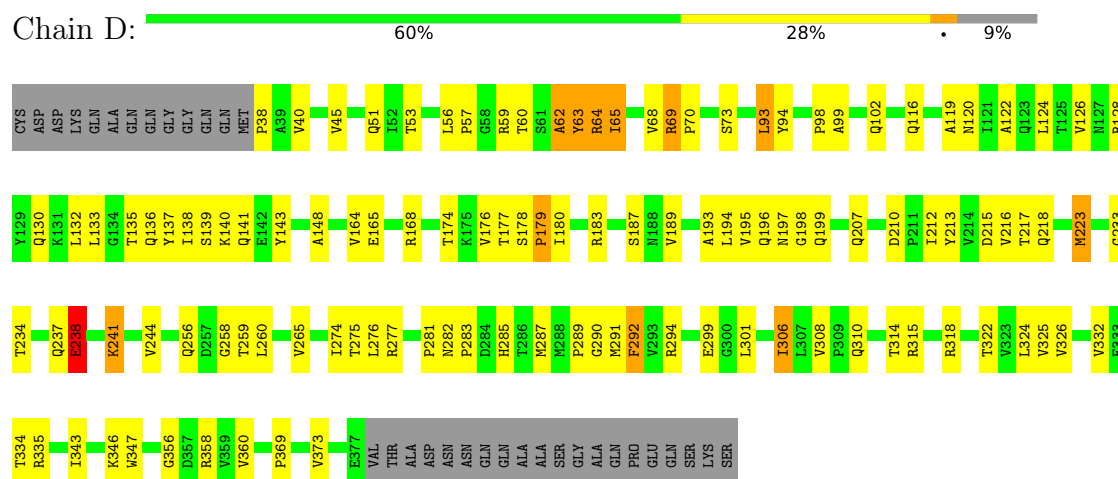
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

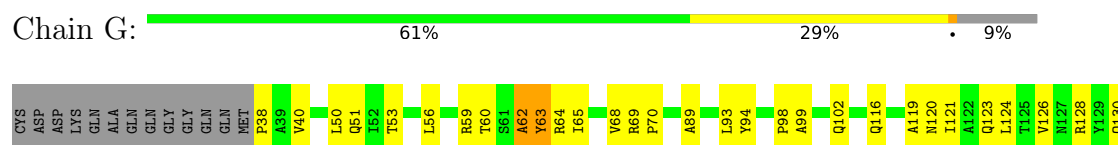
- Molecule 1: Multidrug efflux pump subunit AcrA

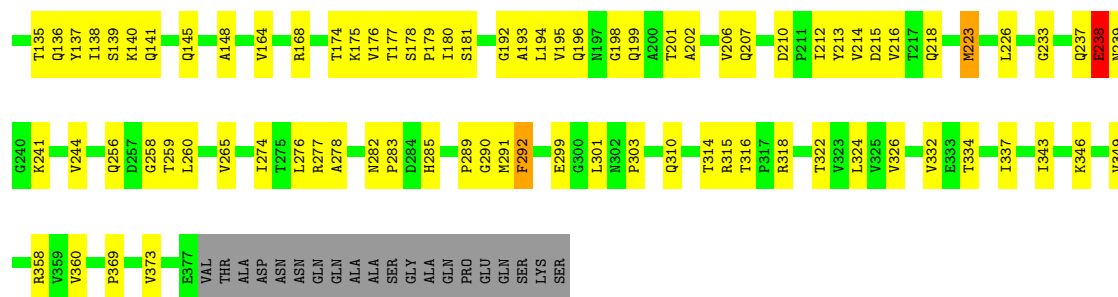


- Molecule 1: Multidrug efflux pump subunit AcrA



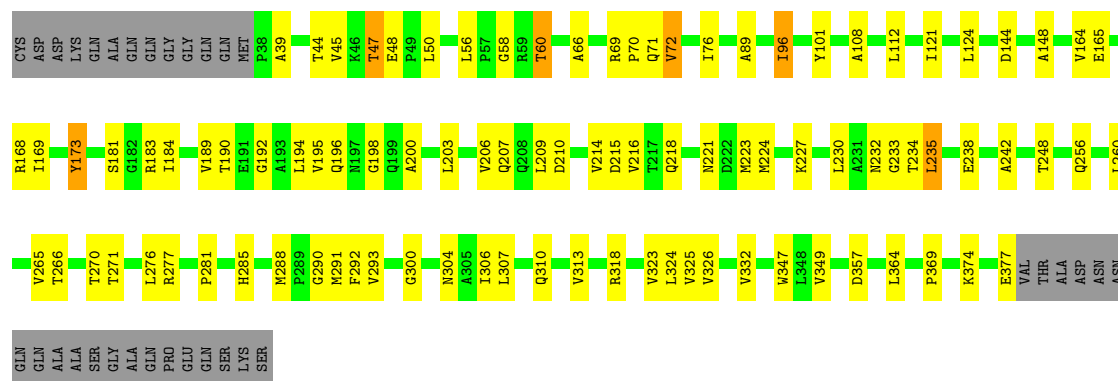
- Molecule 1: Multidrug efflux pump subunit AcrA





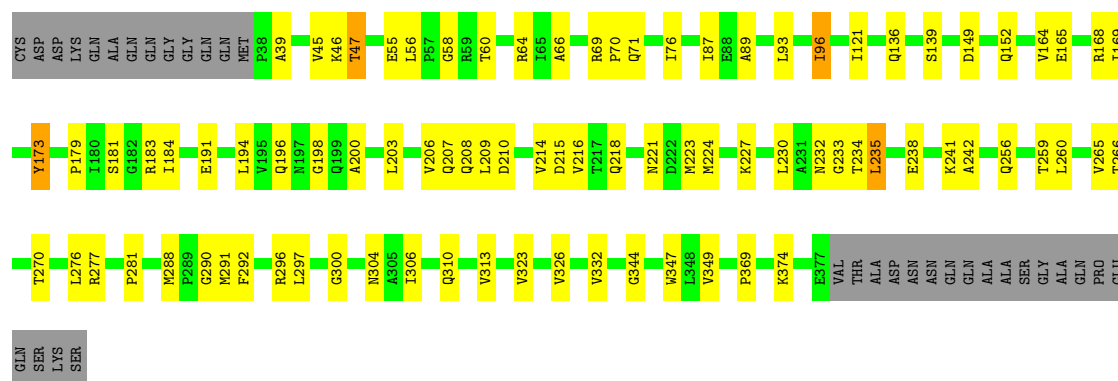
• Molecule 1: Multidrug efflux pump subunit AcrA

Chain H: 66% 24% 9%



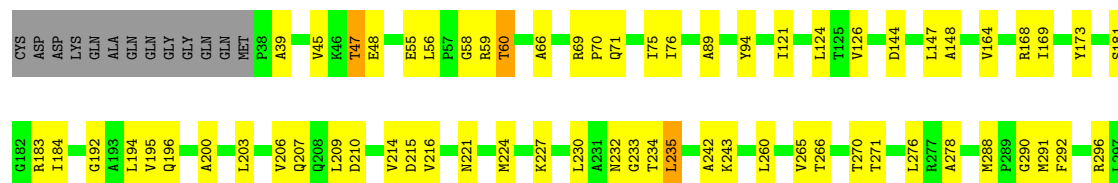
• Molecule 1: Multidrug efflux pump subunit AcrA

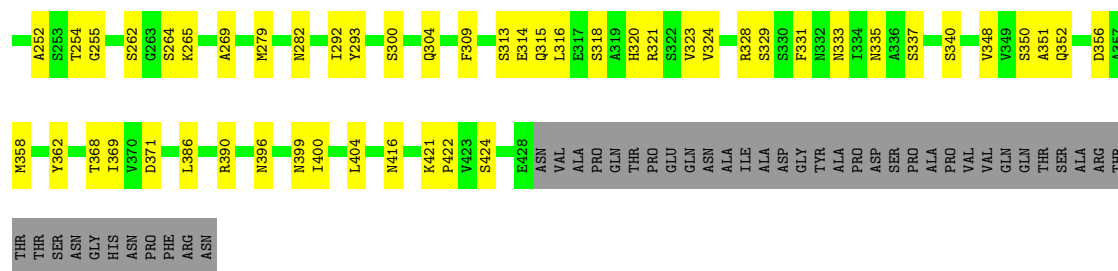
Chain A: 68% 22% 9%



• Molecule 1: Multidrug efflux pump subunit AcrA

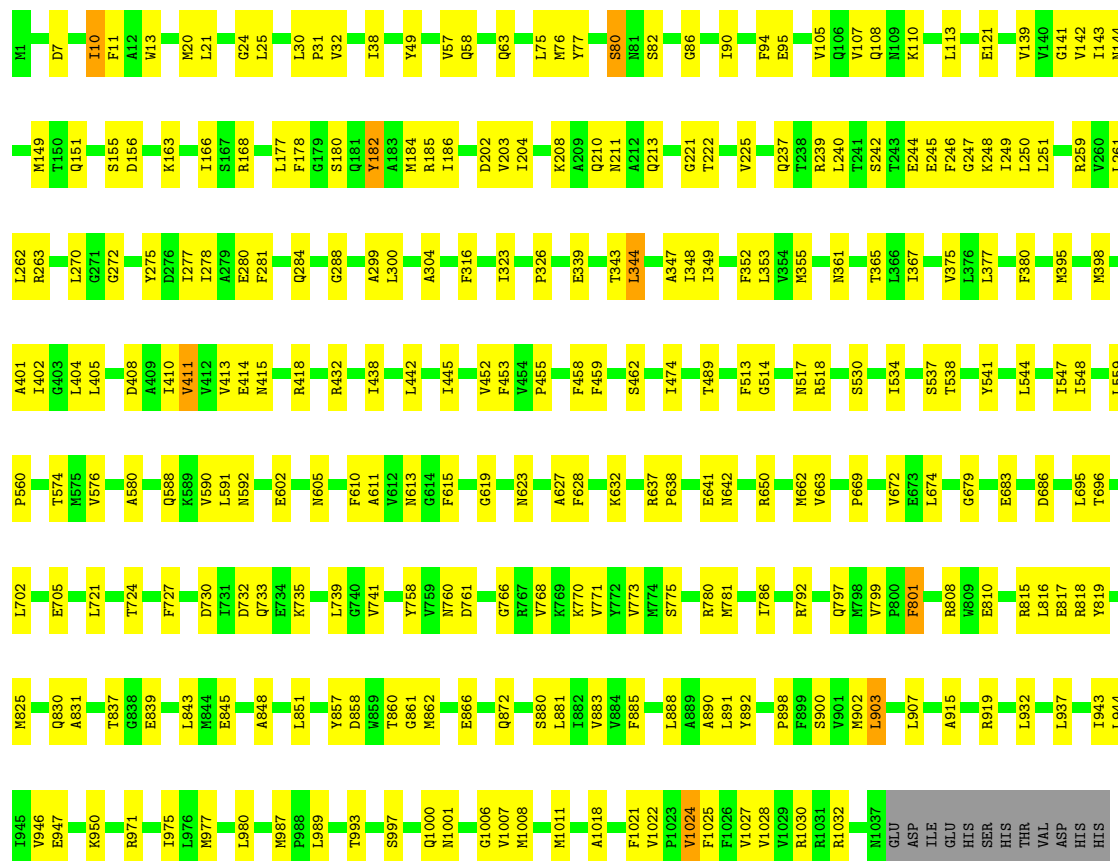
Chain B: 69% 22% 9%





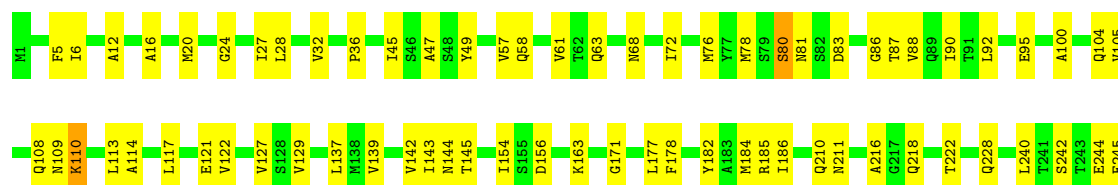
• Molecule 3: Multidrug efflux pump subunit AcrB

Chain K: 73% 25% ..



• Molecule 3: Multidrug efflux pump subunit AcrB

Chain L: 73% 25% .



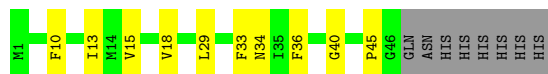


Response	Percentage
Doing a good job	71%
Doing a bad job	27%



- Molecule 4: Multidrug efflux pump accessory protein AcrZ

Chain M:  67% 19% 15%



- Molecule 4: Multidrug efflux pump accessory protein AcrZ

Chain N:  74% 11% 15%



- Molecule 4: Multidrug efflux pump accessory protein AcrZ

Chain O:  70% 15% 15%



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	13544	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING ONLY	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	1	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 5QF

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	0.55	0/2585	0.84	6/3512 (0.2%)
1	B	0.56	0/2585	0.85	3/3512 (0.1%)
1	D	0.55	2/2585 (0.1%)	0.92	5/3512 (0.1%)
1	E	0.58	2/2585 (0.1%)	0.91	5/3512 (0.1%)
1	G	0.56	1/2585 (0.0%)	0.94	4/3512 (0.1%)
1	H	0.56	0/2585	0.85	5/3512 (0.1%)
2	C	0.41	0/3345	0.65	2/4544 (0.0%)
2	F	0.41	0/3345	0.65	1/4544 (0.0%)
2	I	0.41	0/3345	0.65	4/4544 (0.1%)
3	J	0.46	0/7999	0.74	5/10863 (0.0%)
3	K	0.46	0/7968	0.75	4/10826 (0.0%)
3	L	0.46	0/7968	0.75	5/10826 (0.0%)
4	M	0.27	0/330	0.68	0/448
4	N	0.30	0/330	0.76	0/448
4	O	0.32	0/330	0.68	0/448
All	All	0.48	5/50470 (0.0%)	0.77	49/68563 (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	6
1	B	0	4
1	D	0	10
1	E	0	10
1	G	0	6
1	H	0	5
3	J	0	9
3	K	0	4

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Mol	Chain	#Chirality outliers	#Planarity outliers
3	L	0	3
All	All	0	57

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	68	VAL	C-N	-8.48	1.16	1.33
1	E	69	ARG	C-N	-5.87	1.27	1.33
1	D	69	ARG	C-N	-5.65	1.27	1.33
1	G	68	VAL	C-N	-5.45	1.22	1.33
1	D	68	VAL	C-N	-5.14	1.23	1.33

The worst 5 of 49 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	63	TYR	CA-C-N	9.28	138.41	121.70
1	E	63	TYR	C-N-CA	9.28	138.41	121.70
1	D	63	TYR	CA-C-N	8.94	137.79	121.70
1	D	63	TYR	C-N-CA	8.94	137.79	121.70
1	G	63	TYR	CA-C-N	8.81	137.56	121.70

There are no chirality outliers.

5 of 57 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	E	135	THR	Peptide
1	E	174	THR	Peptide
1	E	177	THR	Peptide
1	E	62	ALA	Peptide
1	E	93	LEU	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	2553	0	2610	61	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	2553	0	2610	58	0
1	D	2553	0	2610	84	0
1	E	2553	0	2610	68	0
1	G	2553	0	2610	78	0
1	H	2553	0	2610	65	0
2	C	3304	0	3254	80	0
2	F	3304	0	3254	78	0
2	I	3304	0	3254	89	0
3	J	7849	0	8001	233	0
3	K	7819	0	7919	212	0
3	L	7819	0	7919	208	0
4	M	326	0	346	8	0
4	N	326	0	346	4	0
4	O	326	0	346	6	0
5	J	37	38	0	49	0
5	K	37	38	0	43	0
5	L	37	38	0	42	0
All	All	49806	114	50299	1188	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 12.

The worst 5 of 1188 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)
3:J:628:PHE:HE2	5:J:1101:5QF:C5	1.07	1.64
3:J:628:PHE:CE2	5:J:1101:5QF:C5	1.84	1.59
3:K:277:ILE:HG21	5:K:1101:5QF:C16	1.31	1.57
3:L:178:PHE:HB3	5:L:1101:5QF:C15	1.39	1.49
3:J:628:PHE:HE2	5:J:1101:5QF:N4	1.18	1.39

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	338/373 (91%)	299 (88%)	39 (12%)	0	100	100
1	B	338/373 (91%)	300 (89%)	38 (11%)	0	100	100
1	D	338/373 (91%)	288 (85%)	48 (14%)	2 (1%)	21	59
1	E	338/373 (91%)	286 (85%)	50 (15%)	2 (1%)	21	59
1	G	338/373 (91%)	283 (84%)	53 (16%)	2 (1%)	21	59
1	H	338/373 (91%)	297 (88%)	41 (12%)	0	100	100
2	C	426/493 (86%)	407 (96%)	19 (4%)	0	100	100
2	F	426/493 (86%)	407 (96%)	19 (4%)	0	100	100
2	I	426/493 (86%)	408 (96%)	18 (4%)	0	100	100
3	J	1031/1049 (98%)	937 (91%)	93 (9%)	1 (0%)	48	83
3	K	1035/1049 (99%)	925 (89%)	110 (11%)	0	100	100
3	L	1035/1049 (99%)	935 (90%)	100 (10%)	0	100	100
4	M	44/54 (82%)	39 (89%)	5 (11%)	0	100	100
4	N	44/54 (82%)	40 (91%)	4 (9%)	0	100	100
4	O	44/54 (82%)	41 (93%)	3 (7%)	0	100	100
All	All	6539/7026 (93%)	5892 (90%)	640 (10%)	7 (0%)	49	83

5 of 7 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	140	LYS
1	D	140	LYS
1	G	140	LYS
1	D	179	PRO
3	J	257	GLY

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	274/299 (92%)	272 (99%)	2 (1%)	76	81
1	B	274/299 (92%)	272 (99%)	2 (1%)	76	81
1	D	274/299 (92%)	271 (99%)	3 (1%)	65	76
1	E	274/299 (92%)	271 (99%)	3 (1%)	65	76
1	G	274/299 (92%)	270 (98%)	4 (2%)	57	72
1	H	274/299 (92%)	270 (98%)	4 (2%)	57	72
2	C	358/412 (87%)	358 (100%)	0	100	100
2	F	358/412 (87%)	358 (100%)	0	100	100
2	I	358/412 (87%)	358 (100%)	0	100	100
3	J	839/855 (98%)	835 (100%)	4 (0%)	81	83
3	K	826/855 (97%)	820 (99%)	6 (1%)	76	81
3	L	826/855 (97%)	822 (100%)	4 (0%)	81	83
4	M	34/46 (74%)	34 (100%)	0	100	100
4	N	34/46 (74%)	34 (100%)	0	100	100
4	O	34/46 (74%)	34 (100%)	0	100	100
All	All	5311/5733 (93%)	5279 (99%)	32 (1%)	76	83

5 of 32 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	J	253	VAL
3	J	256	ASP
1	H	72	VAL
1	H	60	THR
3	J	695	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 87 such sidechains are listed below:

Mol	Chain	Res	Type
2	C	332	ASN
3	K	726	GLN
2	C	392	ASN
3	K	191	ASN
3	L	254	ASN

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 ⓘ

3 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
5	5QF	L	1101	-	40,40,40	2.01	12 (30%)	50,56,56	2.66	17 (34%)
5	5QF	K	1101	-	40,40,40	2.05	9 (22%)	50,56,56	2.38	16 (32%)
5	5QF	J	1101	-	40,40,40	2.11	11 (27%)	50,56,56	2.27	13 (26%)

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
5	5QF	L	1101	-	-	11/19/41/41	0/4/4/4
5	5QF	K	1101	-	-	11/19/41/41	0/4/4/4
5	5QF	J	1101	-	-	10/19/41/41	0/4/4/4

The worst 5 of 32 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	K	1101	5QF	C3-N22	4.88	1.49	1.37
5	K	1101	5QF	C5-S9	4.83	1.83	1.76
5	K	1101	5QF	C35-C2	4.79	1.57	1.51
5	K	1101	5QF	C6-C7	4.59	1.51	1.44
5	J	1101	5QF	C3-N22	4.55	1.48	1.37

The worst 5 of 46 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	L	1101	5QF	O34-C33-C32	9.53	116.53	107.80
5	K	1101	5QF	C10-S9-C5	7.36	112.87	101.71
5	L	1101	5QF	C10-S9-C5	7.15	112.55	101.71
5	J	1101	5QF	O18-C16-C15	6.78	124.60	115.40
5	K	1101	5QF	O34-C33-C32	6.75	113.98	107.80

There are no chirality outliers.

5 of 32 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	L	1101	5QF	C11-C10-S9-C5
5	J	1101	5QF	S9-C10-C11-C12
5	L	1101	5QF	C16-C15-O19-C21
5	K	1101	5QF	C15-C16-O18-C20
5	J	1101	5QF	C15-C16-O18-C20

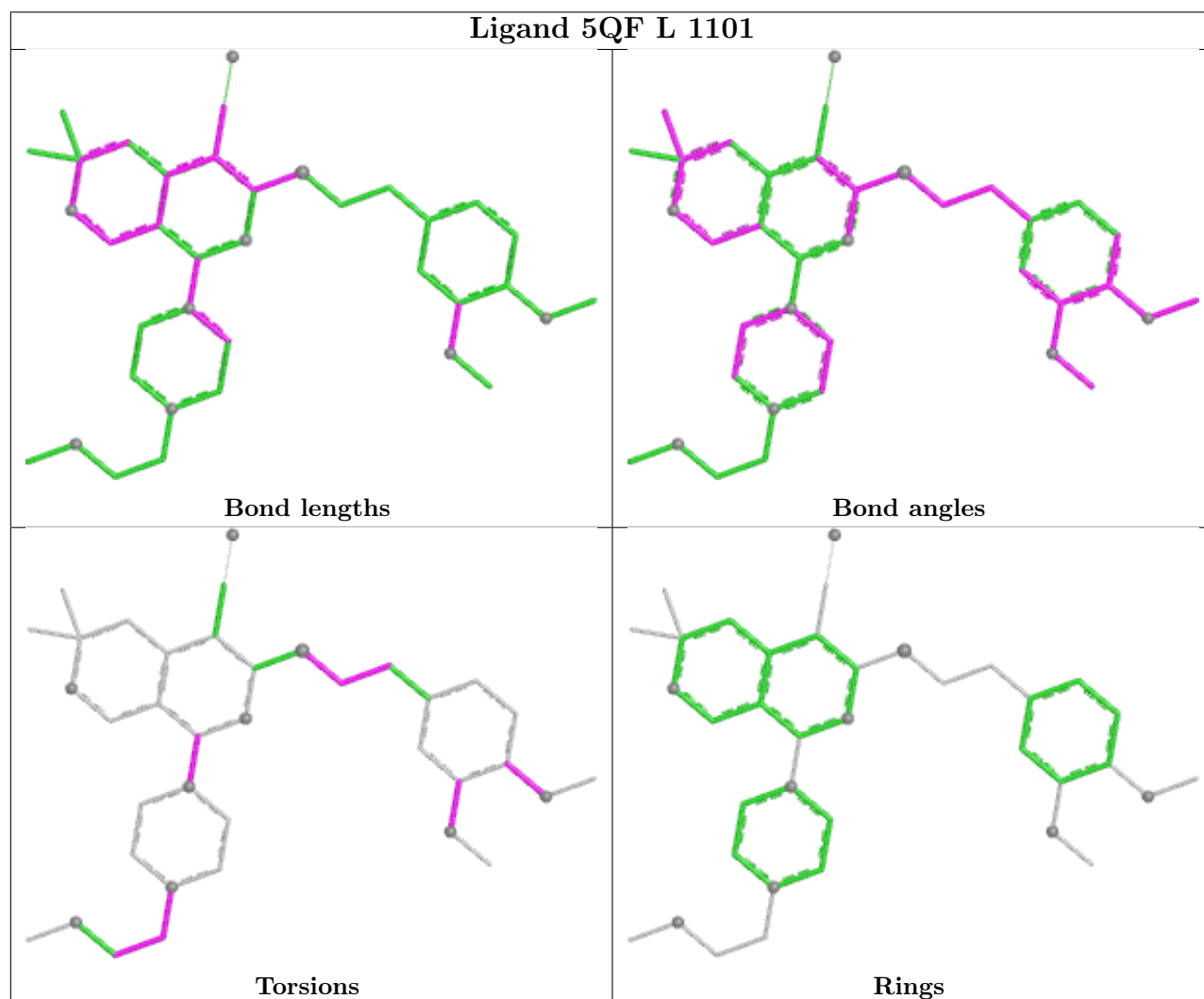
There are no ring outliers.

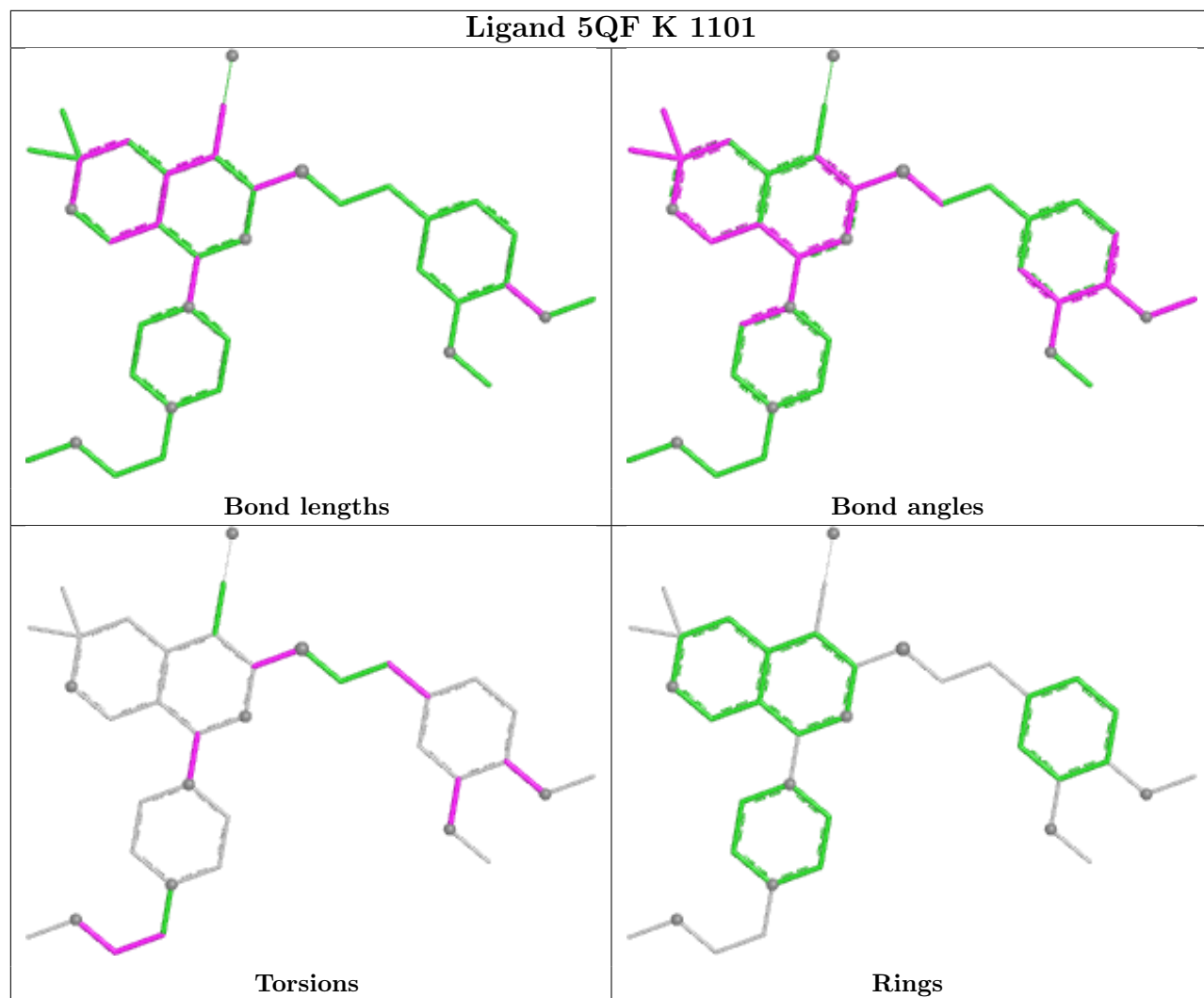
3 monomers are involved in 134 short contacts:

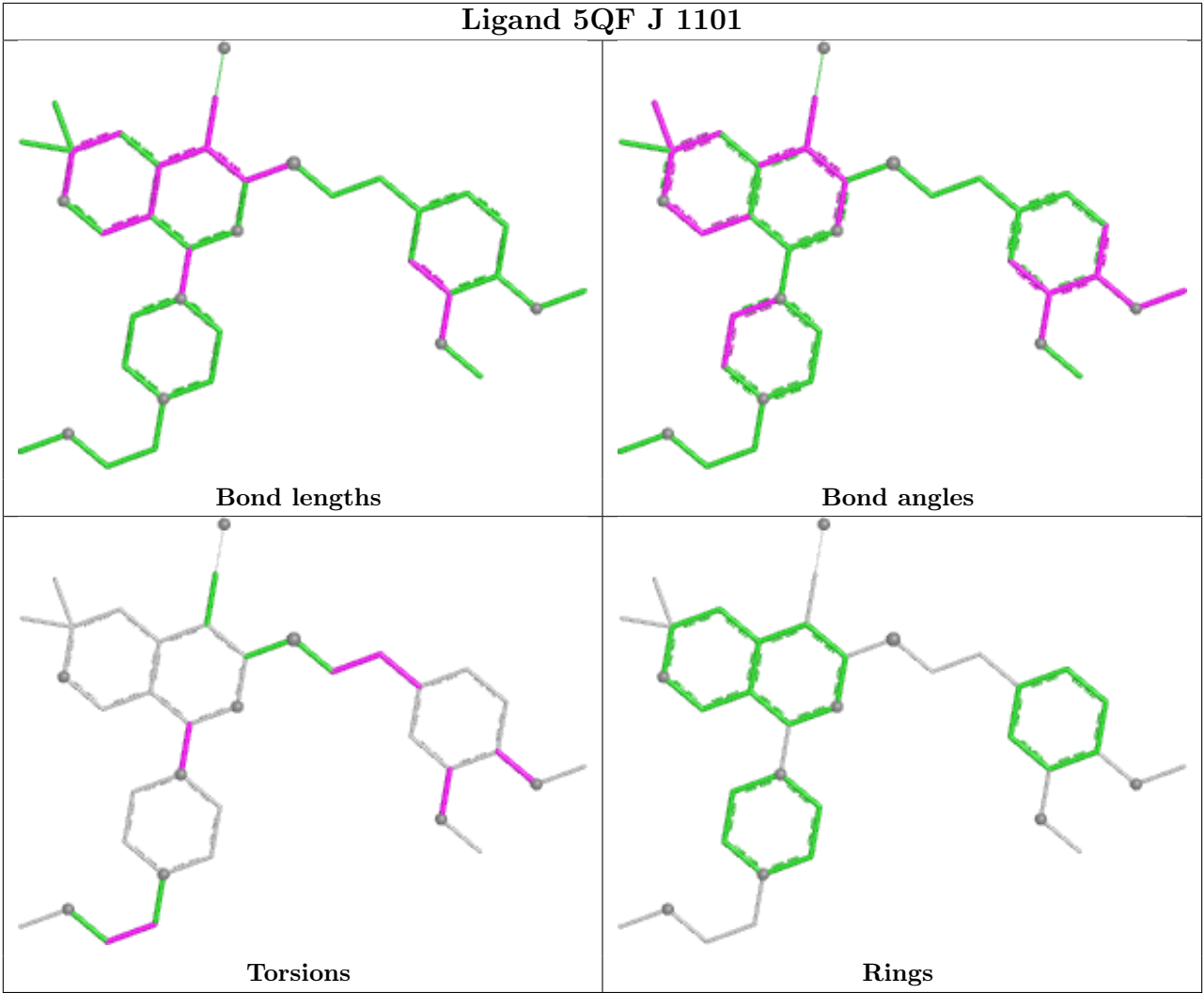
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	L	1101	5QF	42	0
5	K	1101	5QF	43	0
5	J	1101	5QF	49	0

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 ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	E	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	E	68:VAL	C	69:ARG	N	1.16

6 Map visualisation

This section contains visualisations of the EMDB entry EMD-3636. 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.