



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 10:54 PM UTC

PDB ID : 5ENO / pdb_00005eno
Title : MBX2319 bound structure of bacterial efflux pump.
Authors : Sjuts, H.; Ornik, A.R.; Pos, K.M.
Deposited on : 2015-11-09
Resolution : 2.20 Å(reported)

This is a Full wwPDB X-ray Structure 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/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

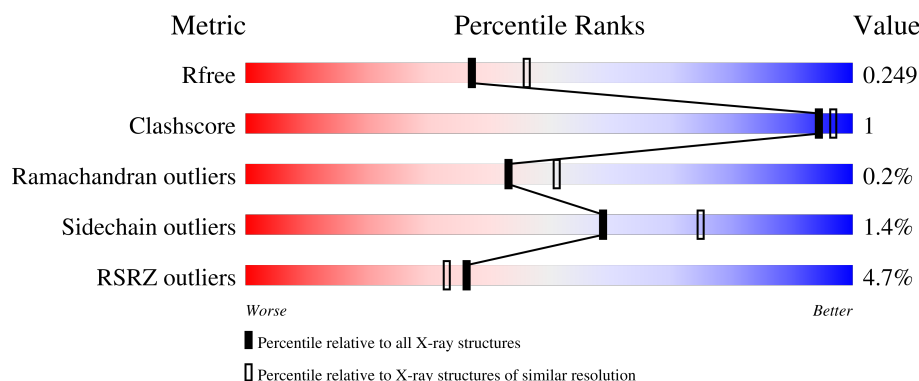
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	609	 4% 90% 5% 5%
1	B	609	 2% 90% 5% 5%
1	C	609	 5% 90% 5% 5%
2	D	169	 6% 90% 6% 8%
2	E	169	 5% 88% 5% 8%

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Mol	Chain	Length	Quality of chain
2	F	169	8% 91% 5%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 17885 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Multidrug efflux pump subunit AcrB, Multidrug efflux pump subunit AcrB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	580	Total	C	N	O	S	0	0	0
			4419	2774	749	874	22			
1	B	578	Total	C	N	O	S	0	0	0
			4405	2766	746	871	22			
1	C	577	Total	C	N	O	S	0	0	0
			4396	2761	744	869	22			

There are 27 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	552	GLY	-	linker	UNP P31224
A	553	GLY	-	linker	UNP P31224
A	554	SER	-	linker	UNP P31224
A	555	GLY	-	linker	UNP P31224
A	556	GLY	-	linker	UNP P31224
A	557	SER	-	linker	UNP P31224
A	558	GLY	-	linker	UNP P31224
A	559	GLY	-	linker	UNP P31224
A	560	SER	-	linker	UNP P31224
B	552	GLY	-	linker	UNP P31224
B	553	GLY	-	linker	UNP P31224
B	554	SER	-	linker	UNP P31224
B	555	GLY	-	linker	UNP P31224
B	556	GLY	-	linker	UNP P31224
B	557	SER	-	linker	UNP P31224
B	558	GLY	-	linker	UNP P31224
B	559	GLY	-	linker	UNP P31224
B	560	SER	-	linker	UNP P31224
C	552	GLY	-	linker	UNP P31224
C	553	GLY	-	linker	UNP P31224
C	554	SER	-	linker	UNP P31224
C	555	GLY	-	linker	UNP P31224

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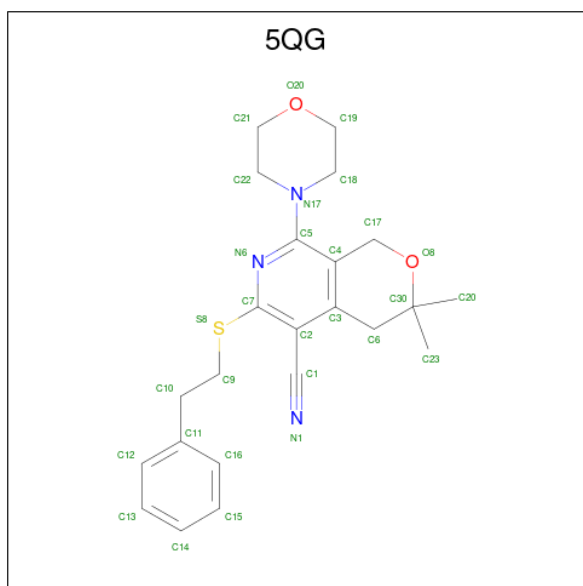
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Chain	Residue	Modelled	Actual	Comment	Reference
C	556	GLY	-	linker	UNP P31224
C	557	SER	-	linker	UNP P31224
C	558	GLY	-	linker	UNP P31224
C	559	GLY	-	linker	UNP P31224
C	560	SER	-	linker	UNP P31224

- Molecule 2 is a protein called DARPin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	162	Total	C	N	O	S	0	0	0
			1237	777	224	235	1			
2	E	156	Total	C	N	O	S	0	0	0
			1177	741	206	229	1			
2	F	162	Total	C	N	O	S	0	0	0
			1237	777	224	235	1			

- Molecule 3 is 3,3-dimethyl-8-morpholin-4-yl-6-(2-phenylethylsulfanyl)-1,4-dihydropyrano[3,4-c]pyridine-5-carbonitrile (CCD ID: 5QG) (formula: C₂₃H₂₇N₃O₂S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	C	1	Total	C	N	O	S	0	0
			29	23	3	2	1		

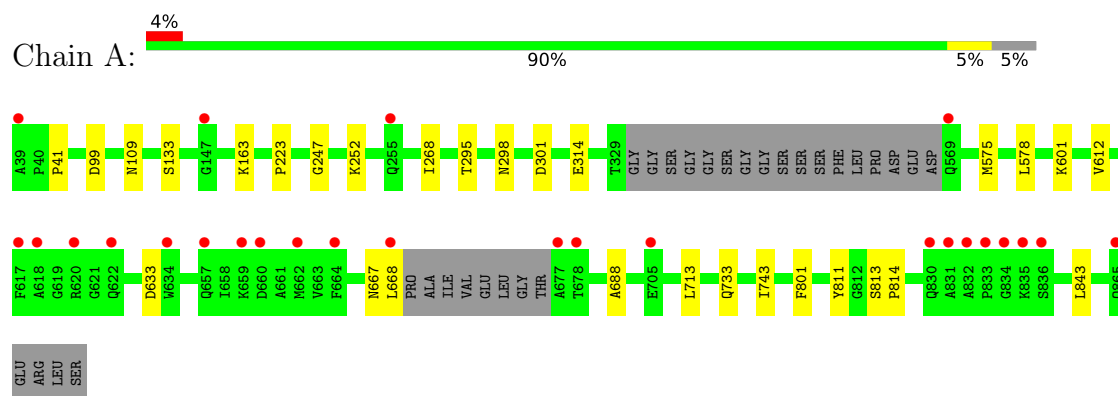
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	292	Total 292	O 292	0	0
4	B	309	Total 309	O 309	0	0
4	C	294	Total 294	O 294	0	0
4	D	36	Total 36	O 36	0	0
4	E	27	Total 27	O 27	0	0
4	F	27	Total 27	O 27	0	0

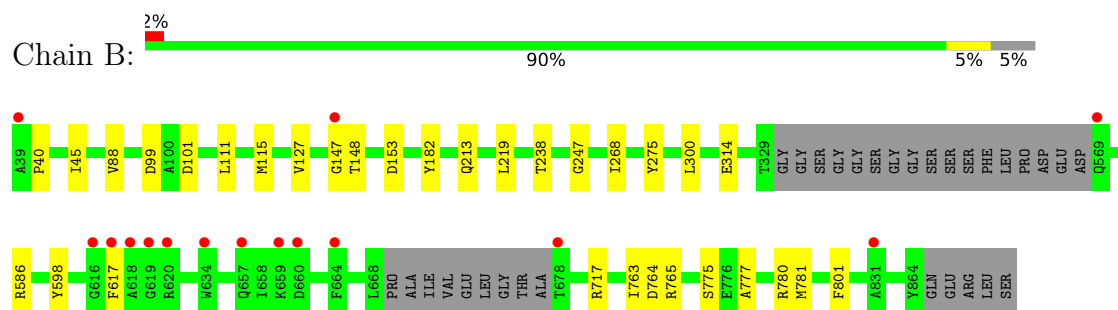
3 Residue-property plots [i](#)

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

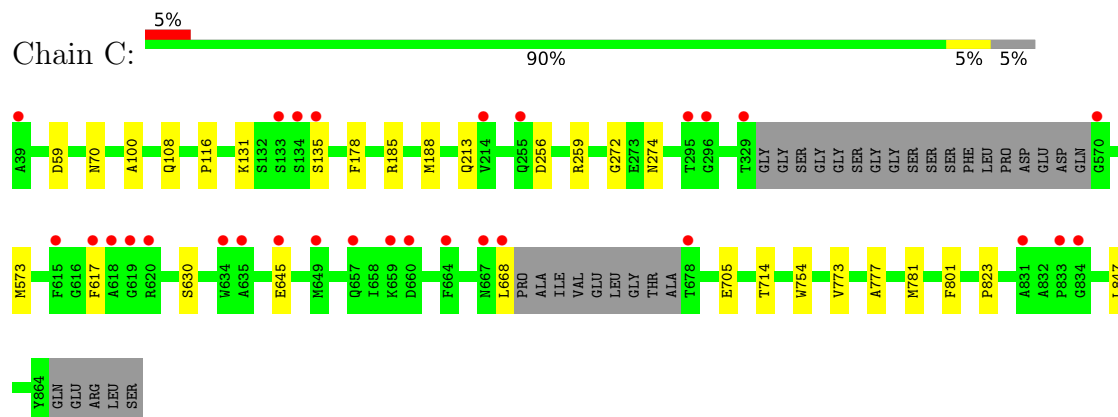
- Molecule 1: Multidrug efflux pump subunit AcrB, Multidrug efflux pump subunit AcrB



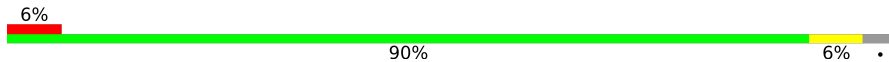
- Molecule 1: Multidrug efflux pump subunit AcrB, Multidrug efflux pump subunit AcrB



- Molecule 1: Multidrug efflux pump subunit AcrB, Multidrug efflux pump subunit AcrB



● Molecule 2: DARPin

Chain D:  6% 90% 6% .

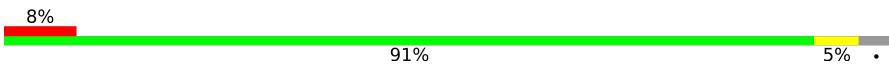


● Molecule 2: DARPin

Chain E:  5% 88% 5% 8%



● Molecule 2: DARPin

Chain F:  8% 91% 5% .



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	108.76Å 145.15Å 174.24Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.20 50.00 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.9 (50.00-2.20) 99.9 (50.00-2.20)	Depositor EDS
R_{merge}	0.31	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.68 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.199 , 0.247 0.205 , 0.249	Depositor DCC
R_{free} test set	6897 reflections (4.89%)	wwPDB-VP
Wilson B-factor (Å ²)	26.5	Xtriage
Anisotropy	0.431	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 26.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	17885	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.63% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	1.15	5/4494 (0.1%)	1.06	5/6089 (0.1%)
1	B	1.15	4/4480 (0.1%)	1.10	3/6070 (0.0%)
1	C	1.10	2/4471 (0.0%)	1.06	0/6058
2	D	1.05	0/1262	1.03	1/1716 (0.1%)
2	E	1.00	0/1196	1.07	0/1626
2	F	1.04	0/1262	1.02	0/1716
All	All	1.11	11/17165 (0.1%)	1.07	9/23275 (0.0%)

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	99	ASP	CA-CB	8.29	1.63	1.53
1	A	814	PRO	N-CA	-6.04	1.42	1.47
1	C	823	PRO	N-CA	-5.49	1.40	1.47
1	A	612	VAL	C-O	-5.39	1.18	1.24
1	B	45	ILE	C-O	-5.33	1.18	1.24
1	B	40	PRO	CA-C	5.28	1.57	1.52
1	A	688	ALA	N-CA	5.14	1.51	1.46
1	A	99	ASP	CA-C	5.12	1.59	1.52
1	A	811	TYR	C-O	-5.09	1.17	1.23
1	B	598	TYR	CA-C	5.03	1.59	1.52
1	C	213	GLN	CA-C	-5.01	1.48	1.53

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	99	ASP	CB-CG-OD2	6.53	133.42	118.40
1	A	601	LYS	N-CA-C	6.23	118.07	111.28
1	B	101	ASP	CB-CA-C	-6.10	101.06	110.81
1	A	163	LYS	N-CA-C	5.62	117.08	111.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	252	LYS	N-CA-C	5.31	116.60	108.42
1	B	717	ARG	N-CA-C	5.29	114.38	108.25
1	A	813	SER	CA-C-N	-5.13	114.83	121.04
1	A	813	SER	C-N-CA	-5.13	114.83	121.04
2	D	49	THR	N-CA-C	-5.03	104.27	110.31

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4419	0	4356	11	0
1	B	4405	0	4343	14	0
1	C	4396	0	4335	16	0
2	D	1237	0	1201	4	0
2	E	1177	0	1159	3	0
2	F	1237	0	1201	2	0
3	C	29	0	0	2	0
4	A	292	0	0	1	0
4	B	309	0	0	2	0
4	C	294	0	0	3	0
4	D	36	0	0	0	0
4	E	27	0	0	0	0
4	F	27	0	0	1	0
All	All	17885	0	16595	45	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 1.

All (45) 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:667:ASN:O	1:A:668:LEU:HG	1.92	0.70
2:F:34:MET:HE2	2:F:34:MET:HA	1.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:733:GLN:OE1	1:A:743:ILE:HD11	1.96	0.65
1:A:743:ILE:HD11	4:A:945:HOH:O	1.98	0.63
1:B:238:THR:HG21	4:B:1204:HOH:O	2.00	0.61
1:C:70:ASN:HB3	4:C:1081:HOH:O	2.02	0.60
2:D:34:MET:HA	2:D:34:MET:HE2	1.82	0.60
1:B:219:LEU:HD23	1:C:754:TRP:CZ3	2.39	0.57
2:D:153:SER:HB3	2:D:161:LEU:HD23	1.89	0.55
1:A:298:ASN:HD22	1:A:301:ASP:H	1.55	0.53
1:B:238:THR:CG2	4:B:1204:HOH:O	2.56	0.53
1:C:617:PHE:N	4:C:1005:HOH:O	2.41	0.53
1:B:115:MET:CE	1:C:116:PRO:HG2	2.40	0.51
1:B:213:GLN:HE22	1:B:238:THR:HG22	1.76	0.51
2:D:42:ALA:O	2:D:50:PRO:HD3	2.12	0.50
1:A:109:ASN:CG	1:C:108:GLN:HG2	2.36	0.49
1:B:247:GLY:HA2	1:B:268:ILE:HD13	1.93	0.49
1:A:41:PRO:HB3	1:A:295:THR:HG21	1.95	0.49
1:B:775:SER:HB3	1:B:780:ARG:HD3	1.94	0.49
2:E:51:LEU:HD11	2:E:63:VAL:HG13	1.95	0.49
1:C:705:GLU:HB3	1:C:847:LEU:HD22	1.94	0.49
1:C:714:THR:HG22	4:C:1164:HOH:O	2.13	0.48
1:B:764:ASP:OD1	1:B:765:ARG:HD3	2.14	0.47
1:C:100:ALA:HB1	1:C:131:LYS:HD2	1.97	0.47
1:C:188:MET:HE2	1:C:773:VAL:HG12	1.96	0.47
1:C:178:PHE:HB3	3:C:901:5QG:C15	2.45	0.47
1:B:153:ASP:OD1	1:B:182:TYR:OH	2.34	0.46
1:A:223:PRO:HD3	1:B:275:TYR:CD1	2.50	0.46
1:B:111:LEU:HD21	1:B:127:VAL:HG11	1.96	0.46
2:E:105:ASP:C	2:E:105:ASP:OD1	2.58	0.45
1:A:667:ASN:O	1:A:667:ASN:CG	2.57	0.45
2:F:80:GLY:O	2:F:110:ASP:HA	2.16	0.45
1:C:178:PHE:HB3	3:C:901:5QG:C14	2.48	0.44
2:E:42:ALA:O	2:E:50:PRO:HD3	2.17	0.44
1:A:578:LEU:HD12	1:A:578:LEU:N	2.32	0.43
1:B:777:ALA:O	1:B:781:MET:HG2	2.19	0.42
1:A:247:GLY:HA2	1:A:268:ILE:CD1	2.49	0.42
1:C:259:ARG:HD2	4:F:202:HOH:O	2.18	0.42
1:C:185:ARG:HD3	1:C:272:GLY:O	2.20	0.42
1:C:777:ALA:O	1:C:781:MET:HG2	2.20	0.42
1:B:763:ILE:HD11	1:C:59:ASP:HB3	2.02	0.41
1:A:713:LEU:HG	1:A:843:LEU:HD23	2.03	0.41
2:D:51:LEU:HD11	2:D:63:VAL:HG13	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:573:MET:HE1	1:C:668:LEU:HD22	2.02	0.41
1:B:247:GLY:HA2	1:B:268:ILE:CD1	2.51	0.41

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	574/609 (94%)	555 (97%)	18 (3%)	1 (0%)	43	51
1	B	572/609 (94%)	554 (97%)	16 (3%)	2 (0%)	36	42
1	C	571/609 (94%)	556 (97%)	15 (3%)	0	100	100
2	D	160/169 (95%)	156 (98%)	3 (2%)	1 (1%)	21	23
2	E	154/169 (91%)	151 (98%)	3 (2%)	0	100	100
2	F	160/169 (95%)	156 (98%)	3 (2%)	1 (1%)	21	23
All	All	2191/2334 (94%)	2128 (97%)	58 (3%)	5 (0%)	43	51

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	633	ASP
1	B	147	GLY
2	F	9	HIS
2	D	12	SER
1	B	617	PHE

5.3.2 Protein sidechains [i](#)

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

resolution.

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	471/492 (96%)	467 (99%)	4 (1%)	73	85
1	B	470/492 (96%)	464 (99%)	6 (1%)	61	76
1	C	469/492 (95%)	463 (99%)	6 (1%)	61	76
2	D	126/132 (96%)	125 (99%)	1 (1%)	73	85
2	E	120/132 (91%)	117 (98%)	3 (2%)	42	56
2	F	126/132 (96%)	121 (96%)	5 (4%)	28	38
All	All	1782/1872 (95%)	1757 (99%)	25 (1%)	59	75

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	133	SER
1	A	314	GLU
1	A	575	MET
1	A	801	PHE
1	B	88	VAL
1	B	148	THR
1	B	300	LEU
1	B	314	GLU
1	B	586	ARG
1	B	801	PHE
1	C	135	SER
1	C	256	ASP
1	C	274	ASN
1	C	630	SER
1	C	645	GLU
1	C	801	PHE
2	D	101	LYS
2	E	97	GLU
2	E	134	LYS
2	E	139	VAL
2	F	20	GLU
2	F	27	ASP
2	F	139	VAL
2	F	163	GLU
2	F	166	GLN

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	68	ASN
1	A	194	ASN
1	A	284	GLN
1	A	604	ASN
1	A	687	GLN
1	A	709	HIS
1	A	737	GLN
1	B	58	GLN
1	B	68	ASN
1	B	109	ASN
1	B	112	GLN
1	B	151	GLN
1	B	229	GLN
1	B	231	ASN
1	B	604	ASN
1	B	622	GLN
1	B	697	GLN
1	C	58	GLN
1	C	81	ASN
1	C	89	GLN
1	C	104	GLN
1	C	108	GLN
1	C	125	GLN
1	C	254	ASN
1	C	284	GLN
1	C	604	ASN
2	D	5	HIS
2	D	142	GLN
2	E	112	ASN
2	E	125	HIS
2	F	92	HIS
2	F	112	ASN
2	F	140	ASN

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](#)

1 ligand is 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	5QG	C	901	-	32,32,32	2.13	5 (15%)	39,45,45	2.39	9 (23%)

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	5QG	C	901	-	-	0/11/31/31	0/4/4/4

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	901	5QG	C2-C1	-8.17	1.30	1.44
3	C	901	5QG	O8-C30	-5.60	1.41	1.45
3	C	901	5QG	C7-S8	-2.99	1.72	1.76
3	C	901	5QG	C5-N17	2.84	1.44	1.37
3	C	901	5QG	C7-N6	2.29	1.36	1.33

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	901	5QG	C9-S8-C7	6.04	110.87	101.71
3	C	901	5QG	C17-O8-C30	5.54	117.66	114.31
3	C	901	5QG	C2-C7-S8	5.40	123.84	118.06
3	C	901	5QG	C2-C7-N6	-4.78	119.07	122.89
3	C	901	5QG	O8-C30-C6	4.71	112.11	107.80
3	C	901	5QG	C4-C5-N6	-4.10	116.35	122.60
3	C	901	5QG	C5-N6-C7	3.67	125.88	116.24
3	C	901	5QG	C17-C4-C3	-3.14	117.39	120.64
3	C	901	5QG	C3-C2-C1	2.77	123.18	119.57

There are no chirality outliers.

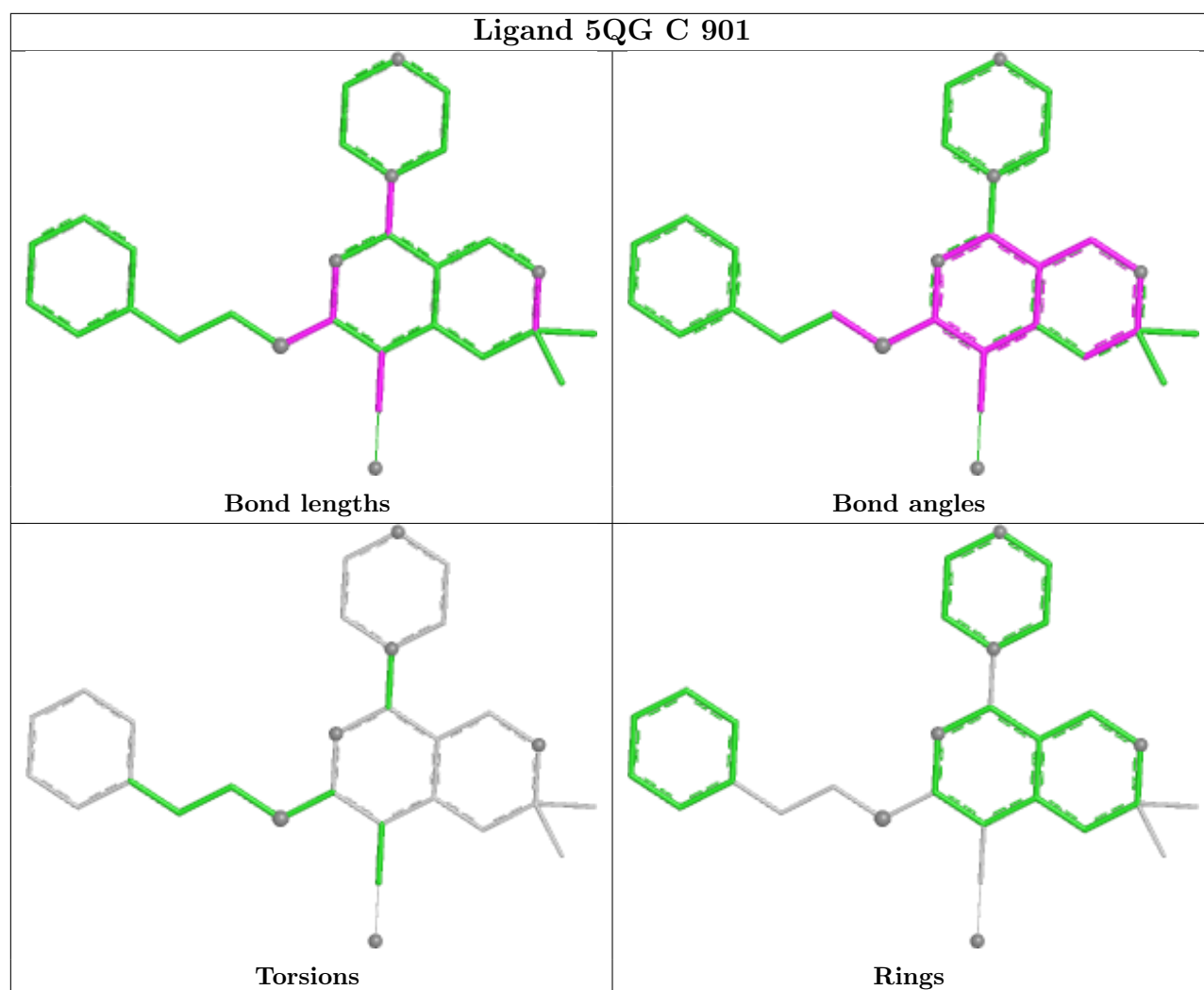
There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	901	5QG	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 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å ²)	Q<0.9
1	A	580/609 (95%)	-0.07	26 (4%)	38	35	14, 26, 61, 94	0
1	B	578/609 (94%)	-0.23	15 (2%)	57	54	14, 24, 53, 128	0
1	C	577/609 (94%)	-0.04	29 (5%)	34	31	12, 27, 65, 88	0
2	D	162/169 (95%)	0.22	10 (6%)	26	23	21, 35, 67, 113	0
2	E	156/169 (92%)	0.39	9 (5%)	29	25	22, 36, 71, 109	0
2	F	162/169 (95%)	0.69	14 (8%)	16	14	23, 40, 88, 131	0
All	All	2215/2334 (94%)	0.00	103 (4%)	36	33	12, 28, 64, 131	0

All (103) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	617	PHE	10.8
1	A	668	LEU	7.9
1	C	678	THR	6.3
2	F	8	HIS	6.1
2	F	11	GLY	6.1
1	A	677	ALA	6.0
1	C	668	LEU	6.0
1	A	634	TRP	6.0
2	E	11	GLY	5.8
1	A	617	PHE	5.8
1	C	618	ALA	5.3
1	C	664	PHE	5.1
2	E	12	SER	5.0
1	B	678	THR	4.9
2	F	10	HIS	4.8
1	A	39	ALA	4.6
1	B	618	ALA	4.6
2	F	9	HIS	4.6
1	B	619	GLY	4.4

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Mol	Chain	Res	Type	RSRZ
1	A	865	GLN	4.4
2	D	8	HIS	4.3
1	A	618	ALA	4.2
1	C	634	TRP	4.0
1	A	831	ALA	4.0
2	D	10	HIS	3.9
2	D	11	GLY	3.9
2	F	5	HIS	3.9
2	D	9	HIS	3.9
1	C	39	ALA	3.7
1	A	664	PHE	3.6
1	C	133	SER	3.6
1	C	617	PHE	3.6
1	A	678	THR	3.5
2	F	166	GLN	3.5
1	C	619	GLY	3.5
2	F	12	SER	3.4
1	B	657	GLN	3.3
2	F	6	HIS	3.2
2	E	166	GLN	3.2
2	D	12	SER	3.2
1	C	620	ARG	3.1
2	E	134	LYS	3.1
2	F	13	ASP	3.1
1	C	329	THR	3.0
2	E	76	TYR	3.0
2	E	165	LEU	3.0
1	A	660	ASP	3.0
1	A	662	MET	3.0
1	C	135	SER	2.9
1	A	147	GLY	2.9
1	B	569	GLN	2.8
1	B	39	ALA	2.8
1	C	833	PRO	2.8
1	B	634	TRP	2.7
1	C	296	GLY	2.7
1	B	660	ASP	2.7
1	C	255	GLN	2.7
1	A	659	LYS	2.7
1	B	147	GLY	2.7
2	E	34	MET	2.6
1	C	657	GLN	2.6

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Mol	Chain	Res	Type	RSRZ
1	C	834	GLY	2.6
1	C	134	SER	2.6
2	F	164	ILE	2.6
1	A	705	GLU	2.5
1	C	635	ALA	2.5
1	C	660	ASP	2.5
2	E	68	LYS	2.5
1	C	615	PHE	2.5
1	C	214	VAL	2.5
2	D	166	GLN	2.5
1	A	569	GLN	2.4
1	A	830	GLN	2.4
1	A	835	LYS	2.4
1	A	622	GLN	2.4
1	B	620	ARG	2.4
1	A	836	SER	2.4
1	B	664	PHE	2.4
2	E	13	ASP	2.3
1	A	255	GLN	2.3
1	A	657	GLN	2.3
1	A	834	GLY	2.3
1	C	570	GLY	2.3
1	C	649	MET	2.3
2	D	163	GLU	2.3
2	F	40	VAL	2.3
1	A	620	ARG	2.3
2	D	13	ASP	2.2
2	F	31	ARG	2.2
2	D	5	HIS	2.2
1	B	831	ALA	2.2
1	B	616	GLY	2.2
1	B	659	LYS	2.2
1	C	659	LYS	2.2
2	F	7	HIS	2.1
2	D	31	ARG	2.1
1	A	832	ALA	2.1
2	F	14	LEU	2.1
1	C	645	GLU	2.1
1	C	667	ASN	2.1
1	C	295	THR	2.0
1	C	831	ALA	2.0
1	A	833	PRO	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

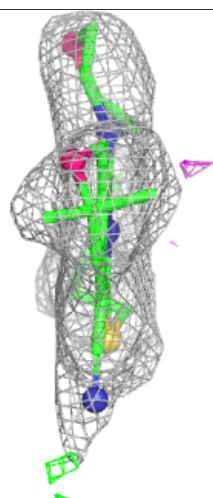
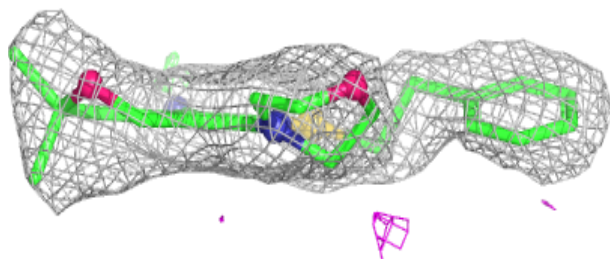
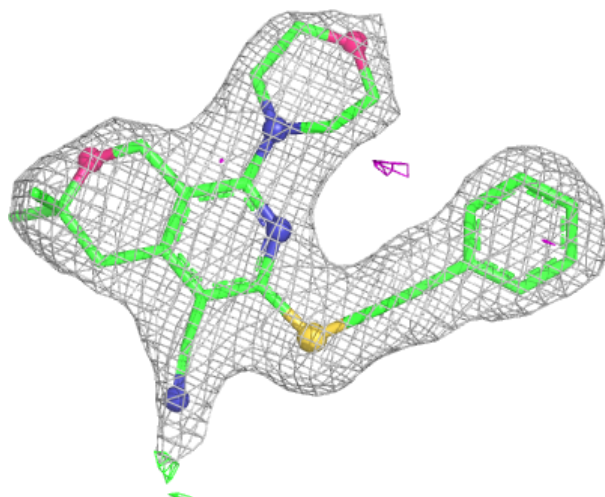
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	5QG	C	901	29/29	0.90	0.13	39,51,63,63	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around 5QG C 901:

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



6.5 Other polymers [i](#)

There are no such residues in this entry.