



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 12:34 PM UTC

PDB ID : 5ENS / pdb_00005ens
Title : Rhodamine bound structure of bacterial efflux pump.
Authors : Sjuts, H.; Ornik, A.R.; Pos, K.M.
Deposited on : 2015-11-09
Resolution : 2.80 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

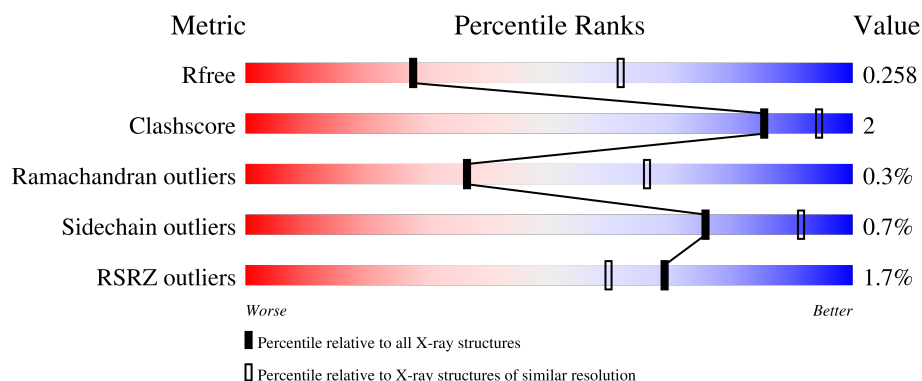
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	2% 89% 6% 5%
1	B	609	% 89% 6% 5%
1	C	609	2% 88% 6% 6%
2	D	169	% 87% 5% 8%
2	E	169	% 88% 5% 8%

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Mol	Chain	Length	Quality of chain
2	F	169	2% 90% 8%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 17010 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
			4404	2765	746	871	22			
1	C	575	Total	C	N	O	S	0	0	0
			4370	2745	742	861	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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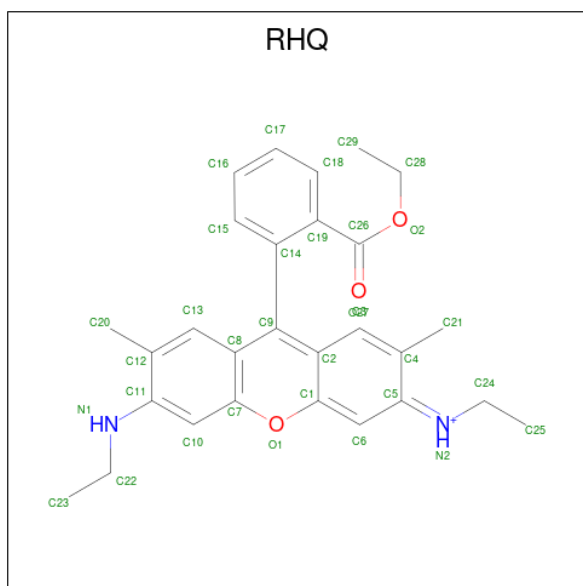
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Chain	Residue	Modelled	Actual	Comment	Reference
C	556	GLY	-	linker	UNP P31224
C	557	SER	-	linker	UNP P31224
C	558	GLY	-	linker	UNP P31224
C	559	GLY	-	linker	UNP P31224
C	560	SER	-	linker	UNP P31224

- Molecule 2 is a protein called DARPin.

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

- Molecule 3 is RHODAMINE 6G (CCD ID: RHQ) (formula: $C_{28}H_{31}N_2O_3$).



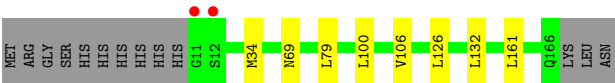
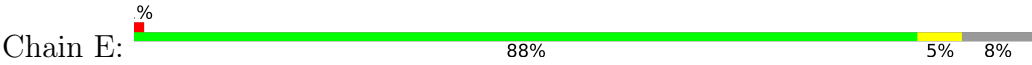
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	C	1	Total	C	N	O	0	0
			33	28	2	3		

- Molecule 4 is water.

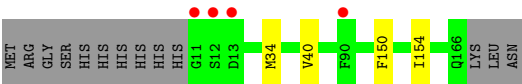
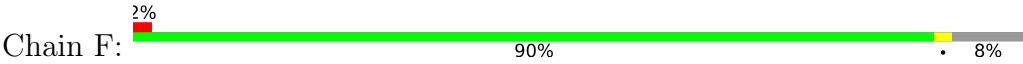
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	74	Total 74	O 74	0	0
4	B	82	Total 82	O 82	0	0
4	C	76	Total 76	O 76	0	0
4	D	6	Total 6	O 6	0	0
4	E	7	Total 7	O 7	0	0
4	F	8	Total 8	O 8	0	0



● Molecule 2: DARPin



● Molecule 2: DARPin



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	108.65Å 145.25Å 174.15Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.80 50.00 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.8 (50.00-2.80) 99.8 (50.00-2.80)	Depositor EDS
R_{merge}	0.28	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.73 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.194 , 0.258 0.198 , 0.258	Depositor DCC
R_{free} test set	3375 reflections (2.75%)	wwPDB-VP
Wilson B-factor (Å ²)	38.0	Xtriage
Anisotropy	0.419	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 28.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	17010	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.64% 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

5.1 Standard geometry

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

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	0.83	0/4494	1.00	0/6089
1	B	0.80	0/4478	0.98	0/6066
1	C	0.80	0/4444	0.97	1/6021 (0.0%)
2	D	0.78	0/1196	0.96	0/1626
2	E	0.81	0/1196	1.00	0/1626
2	F	0.79	0/1196	0.99	0/1626
All	All	0.81	0/17004	0.98	1/23054 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	70	ASN	N-CA-C	6.18	118.53	110.43

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	4419	0	4356	21	0
1	B	4404	0	4340	21	0
1	C	4370	0	4316	22	0
2	D	1177	0	1159	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	E	1177	0	1159	6	0
2	F	1177	0	1159	4	0
3	C	33	0	31	0	0
4	A	74	0	0	0	0
4	B	82	0	0	0	0
4	C	76	0	0	0	0
4	D	6	0	0	0	0
4	E	7	0	0	0	0
4	F	8	0	0	0	0
All	All	17010	0	16520	72	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 2.

All (72) 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:86:GLY:N	1:B:86:GLY:C	2.28	0.91
1:B:85:THR:HB	1:B:87:THR:HG22	1.65	0.78
1:A:733:GLN:OE1	1:A:743:ILE:HD11	1.83	0.78
1:C:56:THR:O	1:C:60:THR:HB	1.93	0.68
1:B:57:VAL:CG1	1:B:88:VAL:HG22	2.27	0.64
1:A:681:ASP:HB3	1:A:860:THR:O	1.96	0.64
1:A:39:ALA:HB1	1:A:40:PRO:CD	2.28	0.63
1:B:111:LEU:HD21	1:B:127:VAL:HG21	1.80	0.62
2:D:150:PHE:CE2	2:D:154:ILE:HD11	2.36	0.61
2:F:34:MET:HE2	2:F:34:MET:HA	1.86	0.58
2:F:34:MET:HE1	2:F:40:VAL:HG22	1.87	0.57
1:C:57:VAL:CG1	1:C:88:VAL:HG22	2.34	0.57
1:C:832:ALA:HB1	1:C:833:PRO:HD2	1.89	0.55
1:B:645:GLU:O	1:B:649:MET:HG3	2.07	0.54
1:B:213:GLN:HE22	1:B:238:THR:HG22	1.72	0.54
1:B:777:ALA:O	1:B:781:MET:HG2	2.09	0.51
1:A:620:ARG:HG2	1:A:621:GLY:N	2.25	0.51
1:B:85:THR:C	1:B:86:GLY:C	2.79	0.50
2:D:100:LEU:HD22	2:D:106:VAL:HG22	1.93	0.50
1:A:39:ALA:HB1	1:A:40:PRO:HD3	1.92	0.49
1:C:72:ILE:HD13	1:C:107:VAL:HG22	1.94	0.49
2:E:100:LEU:HD21	2:E:106:VAL:HG23	1.93	0.49
1:B:120:GLN:HE21	1:B:124:GLN:HE21	1.59	0.48
1:C:261:LEU:HD12	1:C:263:ARG:NH2	2.28	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:247:GLY:HA2	1:B:268:ILE:CD1	2.43	0.47
1:C:750:LEU:HD12	1:C:754:TRP:CD1	2.48	0.47
1:A:64:VAL:HG12	1:A:114:ALA:HB1	1.97	0.47
1:A:151:GLN:HE22	1:A:278:ILE:HG23	1.80	0.47
1:C:706:ALA:HB1	1:C:716:VAL:HG21	1.96	0.47
2:E:34:MET:HE2	2:E:69:ASN:HD22	1.79	0.47
1:A:57:VAL:CG1	1:A:88:VAL:CG2	2.94	0.46
1:A:56:THR:O	1:A:60:THR:HB	2.16	0.46
1:C:49:TYR:HE2	1:C:60:THR:HG21	1.80	0.46
1:C:188:MET:HE2	1:C:773:VAL:HG12	1.98	0.45
1:A:754:TRP:CZ2	1:A:786:ILE:HD13	2.51	0.45
1:B:213:GLN:HG3	1:C:56:THR:HG23	1.98	0.45
1:C:809:TRP:CD1	2:E:79:LEU:HD12	2.52	0.45
1:B:154:ILE:O	1:B:158:VAL:HG23	2.17	0.45
1:A:686:ASP:C	1:A:686:ASP:OD1	2.60	0.44
1:A:652:THR:HG23	1:A:664:PHE:HB2	1.99	0.44
1:C:832:ALA:HB1	1:C:833:PRO:CD	2.47	0.44
2:E:100:LEU:HD11	2:E:132:LEU:HD23	2.00	0.44
1:A:41:PRO:HB3	1:A:295:THR:HG21	2.00	0.44
1:A:151:GLN:NE2	1:A:278:ILE:HG23	2.33	0.44
1:A:754:TRP:CZ3	1:C:219:LEU:HD23	2.53	0.44
2:D:125:HIS:O	2:D:129:VAL:HG23	2.18	0.44
1:A:775:SER:HB3	1:A:780:ARG:HD3	2.00	0.44
1:C:60:THR:HG22	1:C:61:VAL:HG23	1.99	0.43
1:B:762:PHE:CZ	1:B:764:ASP:HB2	2.54	0.43
1:C:580:ALA:HB1	1:C:724:THR:HG22	2.01	0.43
1:A:741:VAL:HG12	1:A:793:ALA:HB2	2.00	0.43
2:F:150:PHE:CZ	2:F:154:ILE:HD11	2.53	0.43
1:B:101:ASP:O	1:B:105:VAL:HG23	2.19	0.43
1:B:281:PHE:O	1:B:284:GLN:HG2	2.18	0.43
1:B:684:LEU:HD11	1:B:851:LEU:HD13	2.00	0.43
1:C:242:SER:OG	1:C:245:GLU:HG3	2.19	0.42
2:D:100:LEU:CD2	2:D:106:VAL:HG22	2.48	0.42
1:C:614:GLY:HA2	1:C:621:GLY:O	2.20	0.42
1:A:51:GLY:C	1:C:215:ALA:HB1	2.45	0.42
1:B:167:SER:HB3	1:C:70:ASN:CB	2.50	0.42
2:E:34:MET:HE2	2:E:69:ASN:HB2	2.01	0.41
1:A:248:LYS:O	1:A:249:ILE:C	2.62	0.41
1:C:733:GLN:OE1	1:C:743:ILE:HD11	2.20	0.41
1:B:225:VAL:HG22	1:C:781:MET:HG3	2.02	0.41
2:D:89:HIS:CD2	2:D:119:LEU:HD22	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:160:ALA:HB1	1:A:767:ARG:HD3	2.03	0.41
2:E:126:LEU:HD23	2:E:161:LEU:HD13	2.03	0.41
1:A:144:ASN:HA	1:A:320:GLY:O	2.20	0.41
1:B:300:LEU:HD23	1:B:300:LEU:HA	1.94	0.41
2:F:150:PHE:CE2	2:F:154:ILE:HD11	2.56	0.40
1:B:309:GLU:HG3	1:B:313:MET:HE3	2.04	0.40
1:B:763:ILE:HD11	1:C:59:ASP:HB3	2.02	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 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%)	553 (96%)	20 (4%)	1 (0%)	43	72
1	B	571/609 (94%)	544 (95%)	25 (4%)	2 (0%)	30	60
1	C	569/609 (93%)	548 (96%)	19 (3%)	2 (0%)	30	60
2	D	154/169 (91%)	148 (96%)	5 (3%)	1 (1%)	21	51
2	E	154/169 (91%)	147 (96%)	7 (4%)	0	100	100
2	F	154/169 (91%)	149 (97%)	5 (3%)	0	100	100
All	All	2176/2334 (93%)	2089 (96%)	81 (4%)	6 (0%)	36	66

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	138	ASP
1	A	147	GLY
1	B	715	SER
1	C	40	PRO
1	B	634	TRP
1	C	832	ALA

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 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%)	468 (99%)	3 (1%)	78	92
1	B	470/492 (96%)	468 (100%)	2 (0%)	84	94
1	C	465/492 (94%)	457 (98%)	8 (2%)	53	83
2	D	120/132 (91%)	120 (100%)	0	100	100
2	E	120/132 (91%)	120 (100%)	0	100	100
2	F	120/132 (91%)	120 (100%)	0	100	100
All	All	1766/1872 (94%)	1753 (99%)	13 (1%)	76	91

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	741	VAL
1	A	801	PHE
1	A	814	PRO
1	B	96	SER
1	B	148	THR
1	C	60	THR
1	C	89	GLN
1	C	138	MET
1	C	615	PHE
1	C	662	MET
1	C	711	ASP
1	C	712	MET
1	C	801	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	81	ASN
1	A	109	ASN
1	A	112	GLN
1	A	144	ASN

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Mol	Chain	Res	Type
1	A	194	ASN
1	A	237	GLN
1	A	274	ASN
1	A	622	GLN
1	B	124	GLN
1	B	569	GLN
1	B	605	ASN
1	B	642	ASN
1	B	726	GLN
1	C	104	GLN
1	C	298	ASN
1	C	657	GLN
2	D	89	HIS
2	D	125	HIS
2	D	142	GLN
2	E	69	ASN
2	E	112	ASN
2	E	122	ASN
2	E	142	GLN
2	E	166	GLN
2	F	59	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	RHQ	C	901	-	35,36,36	2.31	8 (22%)	49,51,51	1.63	12 (24%)

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	RHQ	C	901	-	-	2/15/21/21	0/4/4/4

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	901	RHQ	C19-C14	6.89	1.51	1.40
3	C	901	RHQ	O2-C26	5.74	1.47	1.33
3	C	901	RHQ	C11-C12	5.28	1.51	1.40
3	C	901	RHQ	C8-C7	5.22	1.50	1.40
3	C	901	RHQ	C2-C9	3.40	1.49	1.40
3	C	901	RHQ	C2-C1	3.24	1.49	1.43
3	C	901	RHQ	C3-C4	2.33	1.40	1.35
3	C	901	RHQ	C5-C4	2.32	1.51	1.45

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	901	RHQ	C28-O2-C26	3.84	125.05	116.45
3	C	901	RHQ	O2-C26-C19	3.43	118.97	112.24
3	C	901	RHQ	O1-C7-C10	3.38	120.63	115.83
3	C	901	RHQ	C6-C1-C2	-2.87	118.89	122.86
3	C	901	RHQ	O1-C7-C8	-2.85	118.98	121.57
3	C	901	RHQ	C7-O1-C1	2.79	124.70	119.73
3	C	901	RHQ	C19-C14-C9	-2.65	119.72	123.42
3	C	901	RHQ	O2-C26-O27	-2.57	118.53	123.67
3	C	901	RHQ	C3-C2-C1	2.43	119.96	116.63
3	C	901	RHQ	O1-C1-C6	2.39	120.29	116.73
3	C	901	RHQ	C13-C12-C11	2.29	120.84	118.24
3	C	901	RHQ	C21-C4-C5	2.25	122.61	119.24

There are no chirality outliers.

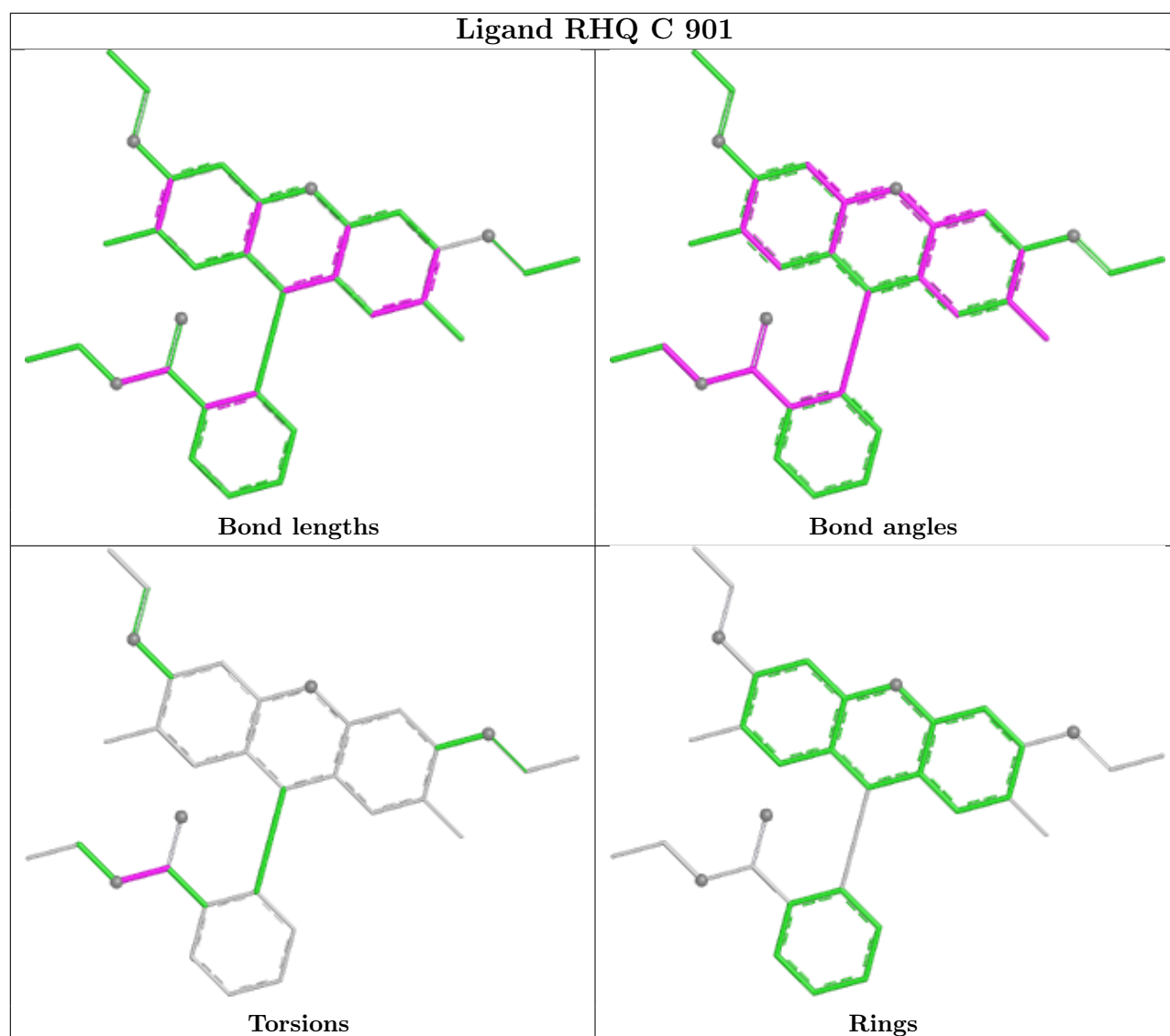
All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	901	RHQ	C19-C26-O2-C28
3	C	901	RHQ	O27-C26-O2-C28

There are no ring outliers.

No monomer is involved in short contacts.

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

6.1 Protein, DNA and RNA chains

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%)	-0.35	10 (1%) 69 60	20, 36, 68, 95	0
1	B	578/609 (94%)	-0.40	7 (1%) 76 68	16, 34, 65, 97	0
1	C	575/609 (94%)	-0.28	13 (2%) 61 51	18, 38, 78, 103	0
2	D	156/169 (92%)	-0.02	2 (1%) 75 66	31, 49, 82, 118	0
2	E	156/169 (92%)	-0.00	2 (1%) 75 66	32, 50, 89, 103	0
2	F	156/169 (92%)	0.12	4 (2%) 57 47	34, 55, 87, 137	0
All	All	2201/2334 (94%)	-0.26	38 (1%) 69 60	16, 39, 76, 137	0

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	677	ALA	4.5
2	D	11	GLY	4.3
1	B	617	PHE	3.9
1	C	677	ALA	3.8
1	A	617	PHE	3.6
1	C	39	ALA	3.4
2	E	11	GLY	3.3
1	A	668	LEU	3.3
2	E	12	SER	3.1
1	C	836	SER	3.1
1	C	133	SER	3.0
1	C	634	TRP	3.0
2	F	11	GLY	2.9
1	B	678	THR	2.9
1	C	862	MET	2.8
1	C	668	LEU	2.7
2	F	12	SER	2.7
1	A	634	TRP	2.7
1	C	618	ALA	2.6

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Mol	Chain	Res	Type	RSRZ
2	F	13	ASP	2.6
1	B	668	LEU	2.6
1	B	569	GLN	2.4
1	A	618	ALA	2.4
1	C	831	ALA	2.3
1	C	833	PRO	2.3
2	D	12	SER	2.3
1	C	134	SER	2.3
1	B	39	ALA	2.2
1	C	326	PRO	2.2
1	A	151	GLN	2.1
1	A	586	ARG	2.1
1	A	865	GLN	2.1
1	B	664	PHE	2.1
1	C	615	PHE	2.1
1	A	257	GLY	2.0
1	B	147	GLY	2.0
2	F	90	PHE	2.0
1	A	589	LYS	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.

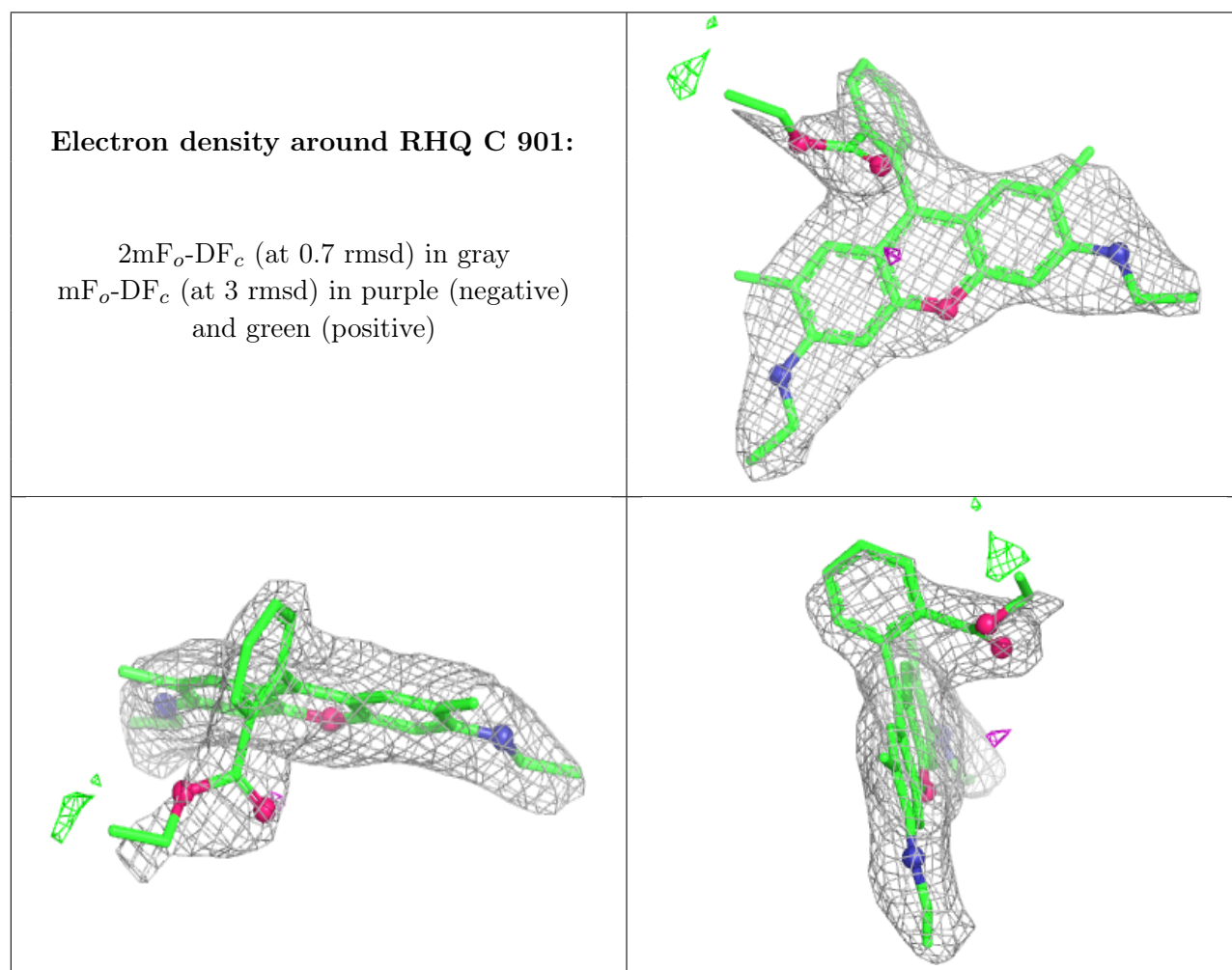
6.4 Ligands [i](#)

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	RHQ	C	901	33/33	0.88	0.16	50,54,82,87	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.



6.5 Other polymers ⓘ

There are no such residues in this entry.