



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 24, 2022 – 05:26 PM EDT

PDB ID : 8DEA  
Title : Scaffold Hopping via Ring Opening Enables Identification of Acyclic Compounds as New Complement Factor D Inhibitors  
Authors : Raman, K.; Babu, Y.S.  
Deposited on : 2022-06-20  
Resolution : 2.21 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](mailto: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>.

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

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.31.2
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.31.2

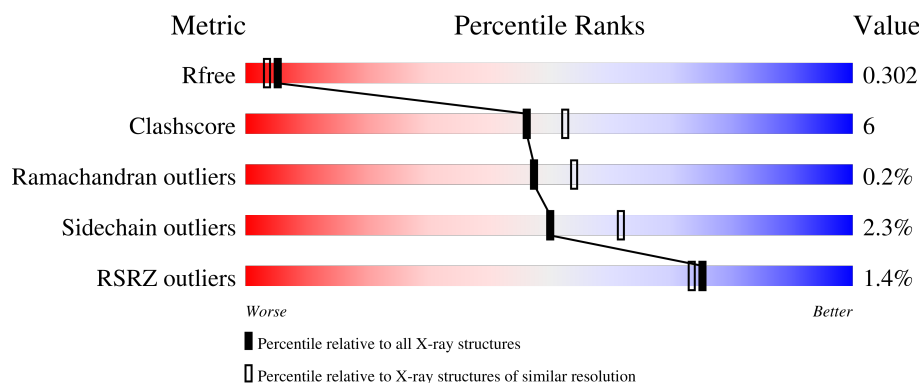
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.21 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))
$R_{free}$	130704	5912 (2.24-2.20)
Clashscore	141614	6646 (2.24-2.20)
Ramachandran outliers	138981	6543 (2.24-2.20)
Sidechain outliers	138945	6544 (2.24-2.20)
RSRZ outliers	127900	5797 (2.24-2.20)

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  $\geq 3$ , 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	230	
1	B	230	
1	C	230	
1	D	230	

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
3	GOL	C	303	-	-	-	X

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6688 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.

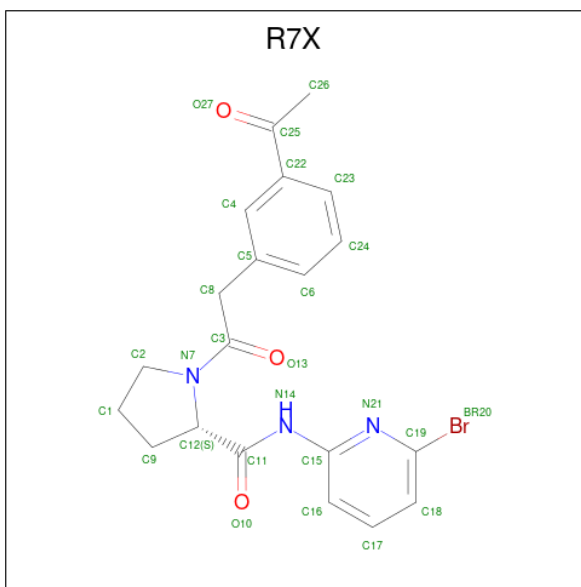
- Molecule 1 is a protein called Complement factor D.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	217	Total	C	N	O	S	0	0	0
			1621	1007	304	300	10			
1	B	217	Total	C	N	O	S	0	0	0
			1621	1007	304	300	10			
1	C	217	Total	C	N	O	S	0	0	0
			1621	1007	304	300	10			
1	D	217	Total	C	N	O	S	0	0	0
			1621	1007	304	300	10			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	229	SER	-	expression tag	UNP P00746
A	230	ALA	-	expression tag	UNP P00746
B	229	SER	-	expression tag	UNP P00746
B	230	ALA	-	expression tag	UNP P00746
C	229	SER	-	expression tag	UNP P00746
C	230	ALA	-	expression tag	UNP P00746
D	229	SER	-	expression tag	UNP P00746
D	230	ALA	-	expression tag	UNP P00746

- Molecule 2 is 1-[(3-acetylphenyl)acetyl]-N-(6-bromopyridin-2-yl)-L-prolinamide (three-letter code: R7X) (formula: C<sub>20</sub>H<sub>20</sub>BrN<sub>3</sub>O<sub>3</sub>) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	Br	C	N	O	0	0
			27	1	20	3	3		
2	B	1	Total	Br	C	N	O	0	0
			27	1	20	3	3		
2	C	1	Total	Br	C	N	O	0	0
			27	1	20	3	3		
2	D	1	Total	Br	C	N	O	0	0
			27	1	20	3	3		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).

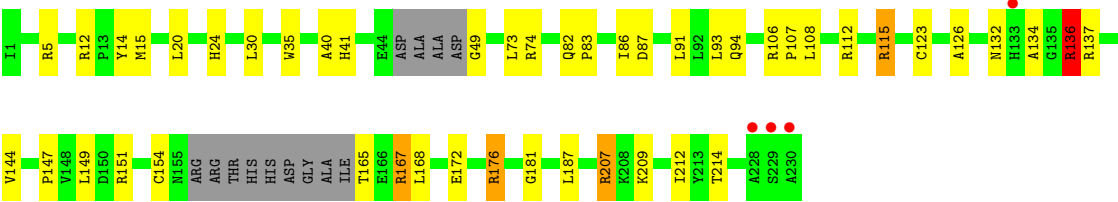


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	C	1	Total	C	O	0	0
			6	3	3		
3	C	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	21	Total	O	0	0
			21	21		
4	B	27	Total	O	0	0
			27	27		
4	C	26	Total	O	0	0
			26	26		
4	D	10	Total	O	0	0
			10	10		







## 4 Data and refinement statistics

Property	Value	Source
Space group	P 2 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	44.90Å 73.03Å 240.86Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.46 – 2.21 46.46 – 2.21	Depositor EDS
% Data completeness (in resolution range)	85.2 (46.46-2.21) 85.2 (46.46-2.21)	Depositor EDS
$R_{merge}$	0.15	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.73 (at 2.22Å)	Xtriage
Refinement program	REFMAC 5.8.0350	Depositor
R, $R_{free}$	0.225 , 0.300 0.229 , 0.302	Depositor DCC
$R_{free}$ test set	1694 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	34.0	Xtriage
Anisotropy	0.049	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 27.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	6688	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.80% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: GOL, R7X

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	0.43	0/1652	0.75	0/2246
1	B	0.41	0/1652	0.72	0/2246
1	C	0.40	0/1652	0.74	0/2246
1	D	0.39	0/1652	0.75	1/2246 (0.0%)
All	All	0.41	0/6608	0.74	1/8984 (0.0%)

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	7
1	B	0	2
1	C	0	4
1	D	0	9
All	All	0	22

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	151	ARG	NE-CZ-NH2	-5.30	117.65	120.30

There are no chirality outliers.

All (22) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	112	ARG	Sidechain

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Mol	Chain	Res	Type	Group
1	A	136	ARG	Sidechain
1	A	151	ARG	Sidechain
1	A	167	ARG	Sidechain
1	A	175	ARG	Sidechain
1	A	215	ARG	Sidechain
1	A	5	ARG	Sidechain
1	B	215	ARG	Sidechain
1	B	5	ARG	Sidechain
1	C	115	ARG	Sidechain
1	C	136	ARG	Sidechain
1	C	151	ARG	Sidechain
1	C	167	ARG	Sidechain
1	D	106	ARG	Sidechain
1	D	112	ARG	Sidechain
1	D	115	ARG	Sidechain
1	D	12	ARG	Sidechain
1	D	136	ARG	Sidechain
1	D	167	ARG	Sidechain
1	D	176	ARG	Sidechain
1	D	207	ARG	Sidechain
1	D	5	ARG	Sidechain

## 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	1621	0	1618	16	0
1	B	1621	0	1618	22	0
1	C	1621	0	1618	16	0
1	D	1621	0	1618	29	0
2	A	27	0	0	1	0
2	B	27	0	0	0	0
2	C	27	0	0	0	0
2	D	27	0	0	1	0
3	C	12	0	16	2	0
4	A	21	0	0	0	0
4	B	27	0	0	0	0
4	C	26	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	10	0	0	0	0
All	All	6688	0	6488	82	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 6.

All (82) 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:C:302:GOL:H11	4:C:416:HOH:O	1.58	1.02
1:D:165:THR:HG22	1:D:167:ARG:H	1.43	0.84
1:C:151:ARG:O	1:C:154:CYS:HB2	1.80	0.79
1:D:91:LEU:HD21	1:D:93:LEU:HD21	1.63	0.79
1:C:50:LYS:NZ	4:C:401:HOH:O	2.20	0.73
1:A:133:HIS:HE1	1:C:84:ASP:OD2	1.73	0.71
1:D:132:ASN:OD1	1:D:134:ALA:HB3	1.91	0.69
1:D:149:LEU:HD21	1:D:172:GLU:HG3	1.75	0.68
1:B:71:ASP:HB2	1:B:96:SER:OG	1.95	0.66
1:B:36:VAL:HG21	1:B:53:VAL:HG11	1.79	0.65
1:B:123:CYS:HB2	1:B:187:LEU:HD11	1.77	0.65
1:D:132:ASN:HD21	1:D:136:ARG:NH1	1.95	0.64
1:C:149:LEU:HD21	1:C:172:GLU:HG3	1.78	0.63
1:B:5:ARG:NH1	1:B:143:HIS:HE1	1.96	0.63
1:B:151:ARG:NH1	1:B:165:THR:O	2.33	0.61
1:A:151:ARG:HH21	1:A:155:ASN:HD21	1.46	0.61
1:A:133:HIS:CE1	1:C:84:ASP:OD2	2.54	0.59
1:B:165:THR:HG22	1:B:167:ARG:H	1.67	0.59
1:D:74:ARG:HH11	1:D:74:ARG:HB2	1.67	0.58
1:B:195:GLY:HA2	1:B:214:THR:O	2.04	0.58
1:B:229:SER:HA	1:D:165:THR:HG23	1.86	0.57
3:C:302:GOL:C1	4:C:416:HOH:O	2.32	0.56
1:D:132:ASN:ND2	1:D:136:ARG:NH1	2.54	0.56
1:A:5:ARG:HB2	1:A:5:ARG:NH1	2.21	0.55
1:B:5:ARG:HH11	1:B:143:HIS:HE1	1.54	0.54
1:C:108:LEU:HD12	1:C:109:PRO:HD2	1.89	0.54
1:B:169:MET:HE3	1:B:215:ARG:HA	1.91	0.52
1:D:123:CYS:HB2	1:D:187:LEU:HD11	1.91	0.52
1:D:20:LEU:HD11	1:D:49:GLY:CA	2.40	0.52
1:A:144:VAL:HG21	1:A:176:ARG:HB3	1.93	0.51
1:A:36:VAL:HG21	1:A:53:VAL:HG11	1.92	0.51
1:A:41:HIS:C	1:A:43:LEU:H	2.14	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:40:ALA:HA	1:C:91:LEU:HB2	1.92	0.51
1:C:132:ASN:OD1	1:C:134:ALA:HB3	2.11	0.50
1:D:168:LEU:HD23	1:D:214:THR:HA	1.93	0.50
1:B:71:ASP:CB	1:B:96:SER:OG	2.60	0.50
1:D:91:LEU:CD2	1:D:93:LEU:HD21	2.40	0.49
1:D:35:TRP:CH2	1:D:94:GLN:HB2	2.48	0.48
1:A:81:SER:HB3	1:A:88:HIS:O	2.13	0.48
1:A:108:LEU:HD12	1:A:109:PRO:HD2	1.96	0.48
1:D:144:VAL:HG21	1:D:176:ARG:HB3	1.95	0.48
1:B:5:ARG:HH11	1:B:143:HIS:CE1	2.31	0.47
1:A:167:ARG:HD3	1:A:219:TYR:OH	2.15	0.46
1:C:151:ARG:NH1	1:C:168:LEU:O	2.47	0.46
1:C:130:ILE:HD12	1:C:132:ASN:O	2.16	0.46
1:A:124:ASP:OD1	1:A:143:HIS:HD2	1.98	0.45
1:D:15:MET:HE1	1:D:126:ALA:HB3	1.99	0.45
1:D:147:PRO:HD2	1:D:172:GLU:HB2	1.96	0.45
1:C:15:MET:HE1	1:C:186:PRO:CG	2.47	0.44
1:D:132:ASN:ND2	1:D:136:ARG:HH11	2.15	0.44
1:D:136:ARG:NH1	1:D:136:ARG:HB2	2.32	0.44
1:B:169:MET:HE2	1:B:215:ARG:HG2	1.99	0.44
1:D:86:ILE:HD13	1:D:212:ILE:HD13	2.00	0.44
1:B:108:LEU:HD12	1:B:109:PRO:HD2	2.00	0.43
1:C:15:MET:HE1	1:C:126:ALA:HB3	2.00	0.43
1:A:15:MET:HE1	1:A:186:PRO:CG	2.49	0.43
1:D:137:ARG:HG2	2:D:301:R7X:BR20	2.73	0.43
1:C:111:GLN:HE21	1:C:113:VAL:H	1.65	0.43
1:A:147:PRO:HD2	1:A:172:GLU:HB2	1.99	0.43
1:D:24:HIS:HE1	1:D:181:GLY:O	2.01	0.43
1:C:92:LEU:HD13	1:C:227:LEU:CD1	2.49	0.43
1:B:91:LEU:HD21	1:B:93:LEU:HD21	2.02	0.42
1:A:180:LYS:HE2	2:A:301:R7X:C17	2.50	0.42
1:D:41:HIS:HE1	1:D:82:GLN:O	2.02	0.42
1:A:151:ARG:O	1:A:154:CYS:HB3	2.20	0.42
1:D:82:GLN:HE21	1:D:83:PRO:HD2	1.84	0.42
1:B:169:MET:CE	1:B:215:ARG:HG2	2.49	0.42
1:D:74:ARG:NH1	1:D:94:GLN:OE1	2.53	0.42
1:B:15:MET:CE	1:B:126:ALA:HB3	2.51	0.41
1:B:213:TYR:CD2	1:B:213:TYR:N	2.89	0.41
1:D:30:LEU:O	1:D:107:PRO:HA	2.20	0.41
1:B:15:MET:HE1	1:B:186:PRO:CG	2.50	0.41
1:B:149:LEU:HD21	1:B:172:GLU:HG3	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:116:ASP:OD1	1:B:151:ARG:NH2	2.52	0.41
1:C:43:LEU:O	1:C:44:GLU:C	2.58	0.41
1:A:136:ARG:HG3	1:A:136:ARG:HH11	1.86	0.41
1:D:132:ASN:HD21	1:D:136:ARG:HH12	1.68	0.41
1:C:24:HIS:HE1	1:C:181:GLY:O	2.03	0.41
1:D:136:ARG:HH11	1:D:136:ARG:CB	2.33	0.41
1:D:40:ALA:HA	1:D:91:LEU:HB2	2.03	0.40
1:D:14:TYR:CG	1:D:108:LEU:HB2	2.56	0.40
1:B:15:MET:HE1	1:B:126:ALA:HB3	2.03	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	211/230 (92%)	204 (97%)	6 (3%)	1 (0%)	29	30
1	B	211/230 (92%)	203 (96%)	8 (4%)	0	100	100
1	C	211/230 (92%)	200 (95%)	11 (5%)	0	100	100
1	D	211/230 (92%)	201 (95%)	9 (4%)	1 (0%)	29	30
All	All	844/920 (92%)	808 (96%)	34 (4%)	2 (0%)	47	54

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	42	CYS
1	D	73	LEU

### 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	174/183 (95%)	168 (97%)	6 (3%)	37	46
1	B	174/183 (95%)	174 (100%)	0	100	100
1	C	174/183 (95%)	170 (98%)	4 (2%)	50	62
1	D	174/183 (95%)	168 (97%)	6 (3%)	37	46
All	All	696/732 (95%)	680 (98%)	16 (2%)	50	62

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	ARG
1	A	112	ARG
1	A	151	ARG
1	A	207	ARG
1	A	209	LYS
1	A	218	SER
1	C	72	VAL
1	C	115	ARG
1	C	154	CYS
1	C	209	LYS
1	D	87	ASP
1	D	115	ARG
1	D	136	ARG
1	D	154	CYS
1	D	207	ARG
1	D	209	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	133	HIS
1	A	143	HIS
1	B	10	HIS

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Mol	Chain	Res	Type
1	B	143	HIS
1	C	21	ASN
1	C	24	HIS
1	C	88	HIS
1	C	111	GLN
1	C	143	HIS
1	D	24	HIS
1	D	41	HIS
1	D	82	GLN

### 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 monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

6 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	R7X	A	301	-	29,29,29	2.25	6 (20%)	40,40,40	2.54	8 (20%)
2	R7X	B	301	-	29,29,29	2.30	7 (24%)	40,40,40	2.05	9 (22%)
2	R7X	D	301	-	29,29,29	2.89	6 (20%)	40,40,40	2.75	8 (20%)
3	GOL	C	302	-	5,5,5	0.23	0	5,5,5	0.54	0



Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	GOL	C	303	-	5,5,5	0.29	0	5,5,5	0.25	0
2	R7X	C	301	-	29,29,29	2.05	7 (24%)	40,40,40	2.62	9 (22%)

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	R7X	A	301	-	-	1/20/30/30	0/3/3/3
2	R7X	B	301	-	-	2/20/30/30	0/3/3/3
2	R7X	D	301	-	-	0/20/30/30	0/3/3/3
3	GOL	C	302	-	-	3/4/4/4	-
3	GOL	C	303	-	-	2/4/4/4	-
2	R7X	C	301	-	-	0/20/30/30	0/3/3/3

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	301	R7X	BR20-C19	8.03	2.03	1.90
2	D	301	R7X	C8-C5	-7.50	1.39	1.51
2	A	301	R7X	C8-C5	-7.50	1.39	1.51
2	B	301	R7X	C8-C3	7.45	1.60	1.52
2	B	301	R7X	C8-C5	-6.08	1.41	1.51
2	C	301	R7X	C8-C5	-5.87	1.42	1.51
2	D	301	R7X	C22-C25	-5.65	1.39	1.49
2	D	301	R7X	C12-C11	-5.14	1.40	1.52
2	A	301	R7X	C8-C3	5.12	1.57	1.52
2	D	301	R7X	C8-C3	5.00	1.57	1.52
2	C	301	R7X	C22-C25	-4.80	1.40	1.49
2	B	301	R7X	C12-C11	-4.50	1.42	1.52
2	A	301	R7X	BR20-C19	4.42	1.97	1.90
2	D	301	R7X	C19-N21	3.91	1.36	1.32
2	C	301	R7X	C12-C11	-3.71	1.44	1.52
2	A	301	R7X	C22-C25	-3.21	1.43	1.49
2	C	301	R7X	C19-N21	3.02	1.35	1.32
2	A	301	R7X	C12-C11	-2.96	1.45	1.52
2	C	301	R7X	C12-N7	2.90	1.52	1.47
2	B	301	R7X	C22-C25	-2.75	1.44	1.49
2	B	301	R7X	C12-N7	2.69	1.52	1.47
2	C	301	R7X	C8-C3	2.63	1.55	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	301	R7X	C2-N7	2.34	1.52	1.47
2	A	301	R7X	C15-N14	-2.30	1.35	1.40
2	B	301	R7X	C15-N14	-2.23	1.35	1.40
2	B	301	R7X	C24-C6	2.20	1.43	1.38

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	301	R7X	C18-C19-N21	-10.19	117.16	125.75
2	D	301	R7X	C19-N21-C15	9.93	124.51	116.29
2	C	301	R7X	C19-N21-C15	9.73	124.34	116.29
2	C	301	R7X	C18-C19-N21	-9.36	117.86	125.75
2	A	301	R7X	C19-N21-C15	9.00	123.74	116.29
2	A	301	R7X	C18-C19-N21	-8.04	118.97	125.75
2	B	301	R7X	C19-N21-C15	7.76	122.71	116.29
2	B	301	R7X	C18-C19-N21	-5.69	120.95	125.75
2	A	301	R7X	C15-N14-C11	-5.41	119.88	127.55
2	D	301	R7X	BR20-C19-C18	5.16	123.92	118.64
2	A	301	R7X	BR20-C19-N21	4.71	119.89	115.71
2	B	301	R7X	C15-N14-C11	-4.70	120.88	127.55
2	C	301	R7X	BR20-C19-N21	4.62	119.80	115.71
2	D	301	R7X	C17-C18-C19	4.19	120.64	116.70
2	D	301	R7X	BR20-C19-N21	3.63	118.93	115.71
2	C	301	R7X	BR20-C19-C18	3.58	122.31	118.64
2	D	301	R7X	C15-N14-C11	-3.44	122.67	127.55
2	A	301	R7X	C17-C18-C19	3.22	119.72	116.70
2	A	301	R7X	C16-C15-N21	-3.07	117.98	123.16
2	C	301	R7X	C17-C18-C19	2.96	119.48	116.70
2	B	301	R7X	BR20-C19-N21	2.91	118.29	115.71
2	A	301	R7X	C2-N7-C12	-2.65	107.80	112.00
2	C	301	R7X	C16-C15-N21	-2.64	118.71	123.16
2	B	301	R7X	C9-C12-N7	2.54	106.80	103.03
2	C	301	R7X	C2-N7-C12	-2.53	107.99	112.00
2	B	301	R7X	C16-C15-N21	-2.52	118.91	123.16
2	A	301	R7X	BR20-C19-C18	2.42	121.12	118.64
2	D	301	R7X	C16-C15-N21	-2.42	119.08	123.16
2	C	301	R7X	C9-C12-N7	2.30	106.45	103.03
2	D	301	R7X	C1-C2-N7	2.30	107.29	103.25
2	B	301	R7X	C22-C4-C5	2.17	123.88	120.86
2	B	301	R7X	BR20-C19-C18	2.06	120.75	118.64
2	C	301	R7X	C26-C25-C22	-2.04	115.05	119.24
2	B	301	R7X	C2-N7-C12	-2.02	108.80	112.00

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	302	GOL	C1-C2-C3-O3
3	C	303	GOL	O1-C1-C2-C3
3	C	303	GOL	O1-C1-C2-O2
2	B	301	R7X	N21-C15-N14-C11
3	C	302	GOL	O2-C2-C3-O3
2	B	301	R7X	C16-C15-N14-C11
3	C	302	GOL	O1-C1-C2-C3
2	A	301	R7X	N21-C15-N14-C11

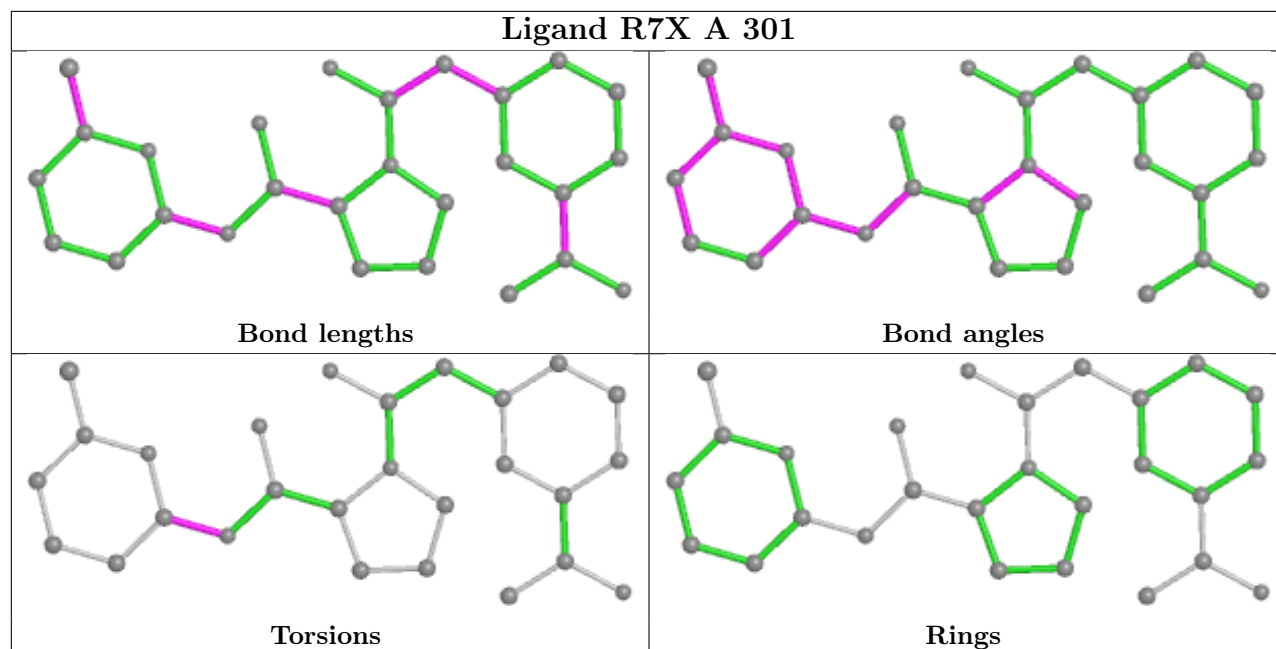
There are no ring outliers.

3 monomers are involved in 4 short contacts:

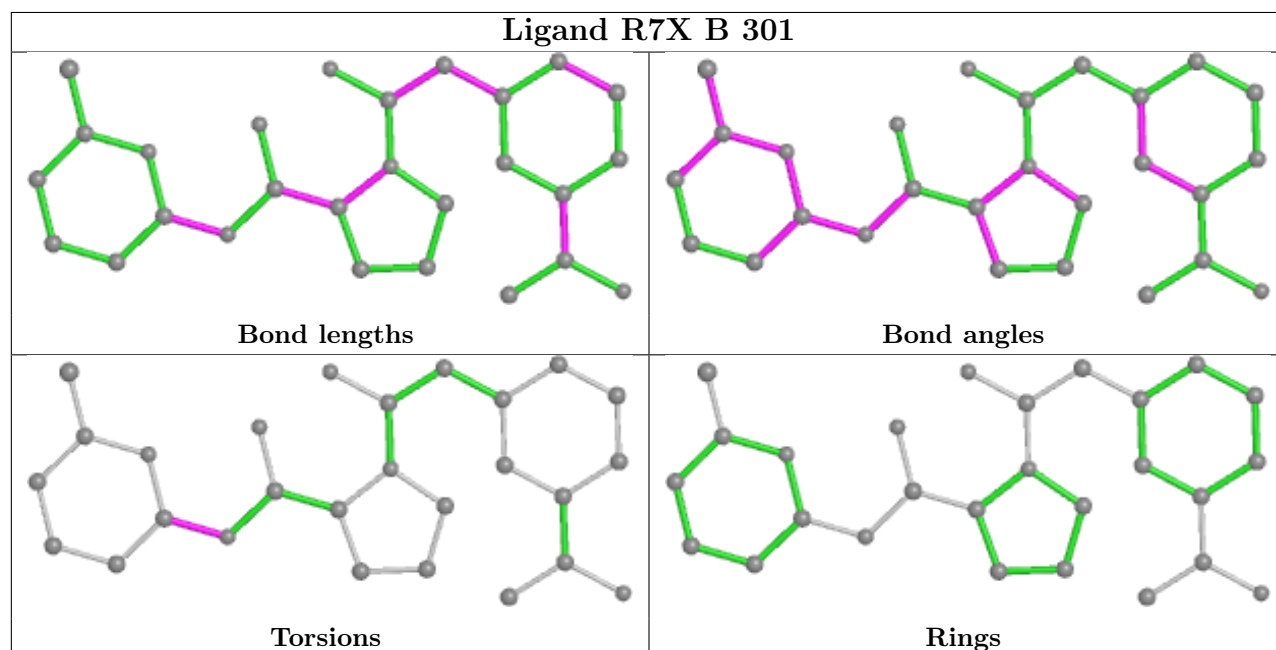
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	301	R7X	1	0
2	D	301	R7X	1	0
3	C	302	GOL	2	0

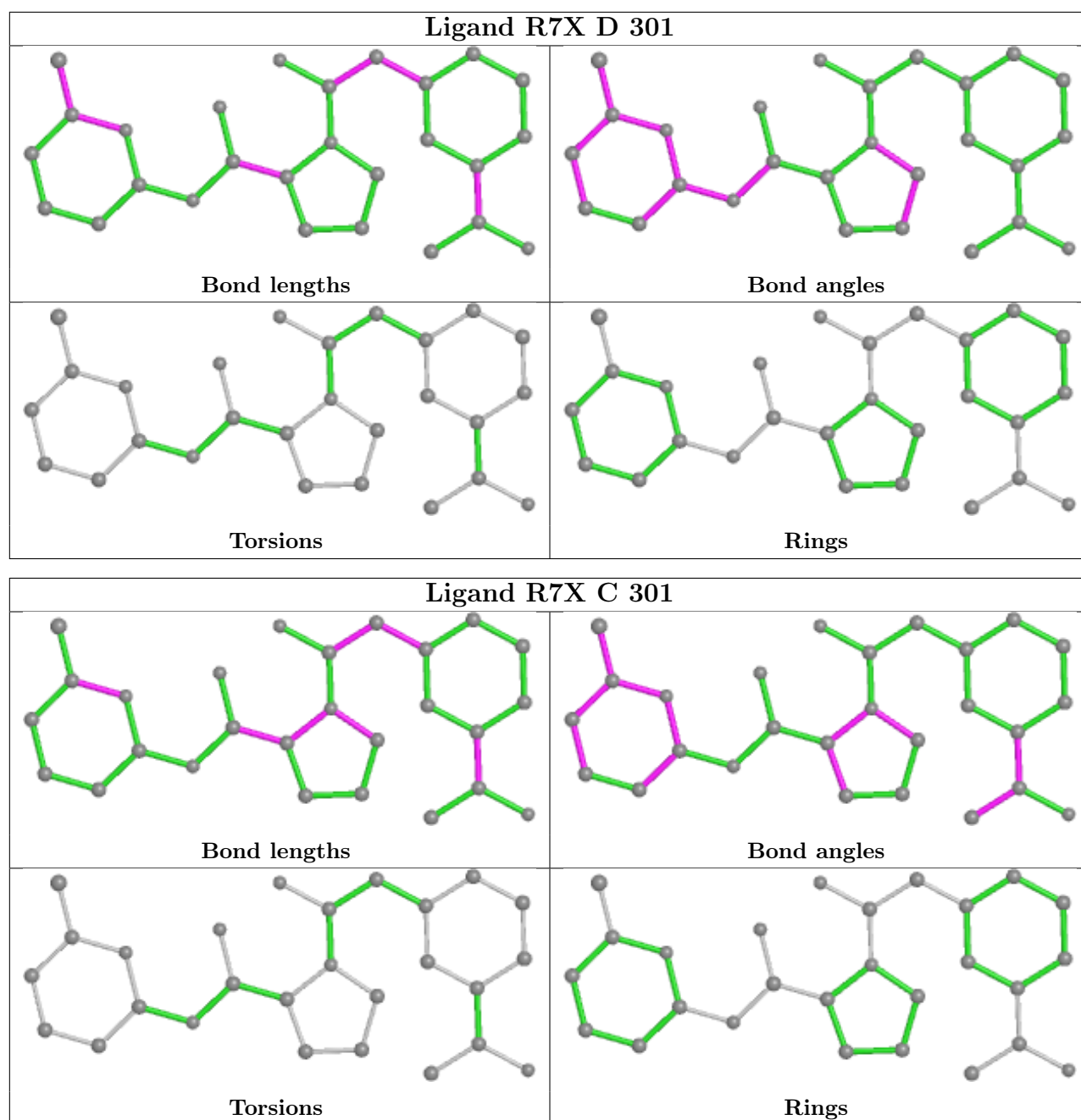
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

## Ligand R7X A 301



## Ligand R7X B 301





## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	217/230 (94%)	-0.29	3 (1%) 75 73	21, 29, 49, 101	0
1	B	217/230 (94%)	-0.26	3 (1%) 75 73	20, 32, 56, 87	0
1	C	217/230 (94%)	-0.28	2 (0%) 84 83	22, 31, 50, 93	0
1	D	217/230 (94%)	-0.07	4 (1%) 68 66	23, 35, 60, 113	0
All	All	868/920 (94%)	-0.23	12 (1%) 75 73	20, 32, 56, 113	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	230	ALA	11.0
1	B	230	ALA	10.2
1	D	229	SER	8.5
1	C	229	SER	6.9
1	A	230	ALA	6.0
1	A	229	SER	6.0
1	B	229	SER	5.0
1	C	230	ALA	4.0
1	B	49	GLY	3.1
1	D	133	HIS	2.6
1	A	49	GLY	2.4
1	D	228	ALA	2.3

### 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 monosaccharides in this entry.

## 6.4 Ligands

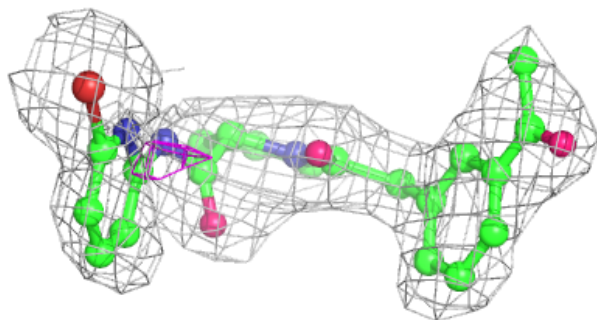
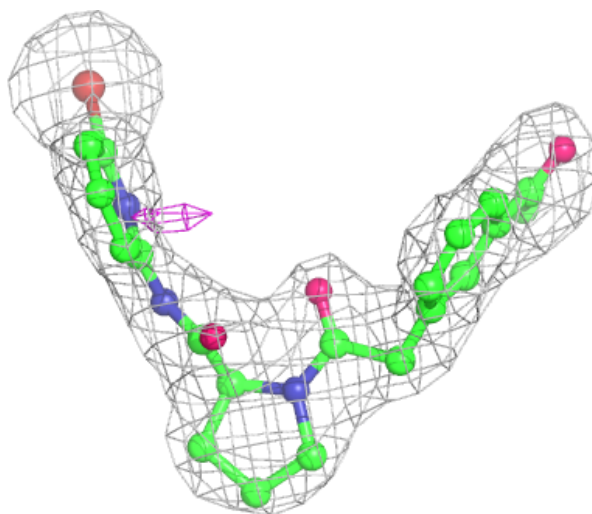
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
3	GOL	C	303	6/6	0.77	0.71	99,106,111,112	0
3	GOL	C	302	6/6	0.79	0.19	34,38,40,40	0
2	R7X	D	301	27/27	0.96	0.12	32,36,43,48	0
2	R7X	A	301	27/27	0.98	0.12	22,28,35,35	0
2	R7X	B	301	27/27	0.98	0.10	25,30,32,34	0
2	R7X	C	301	27/27	0.98	0.11	26,28,32,32	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 R7X D 301:**

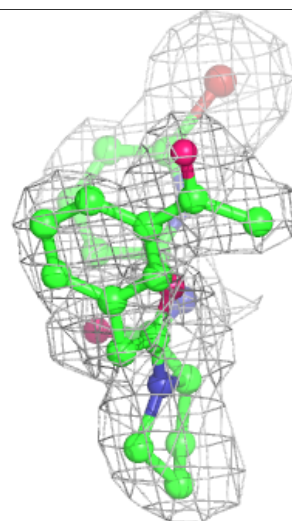
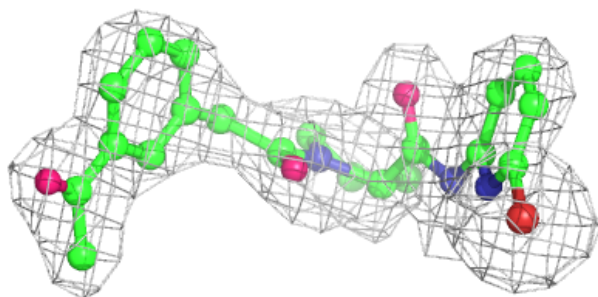
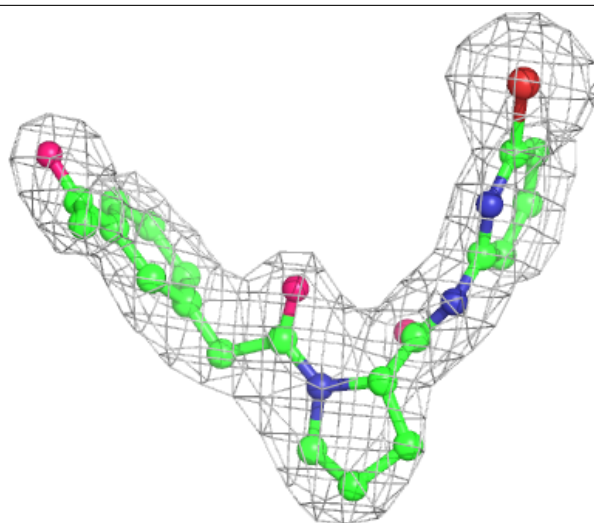
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





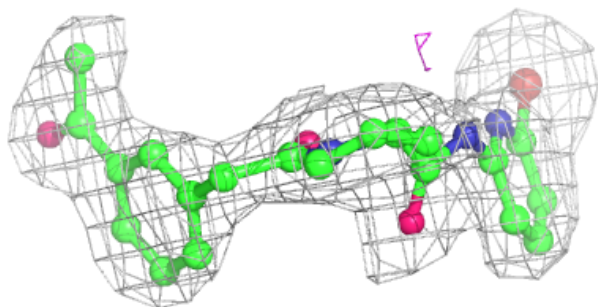
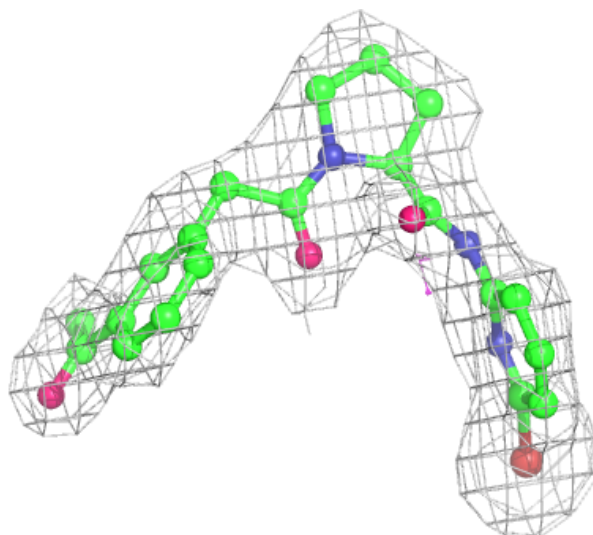
**Electron density around R7X A 301:**

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



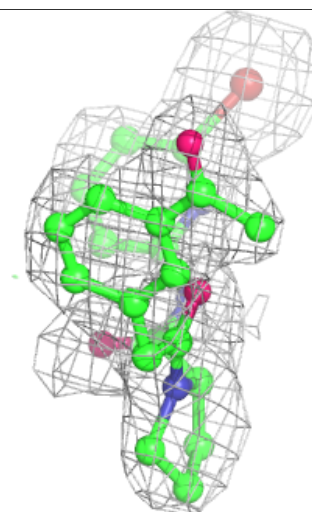
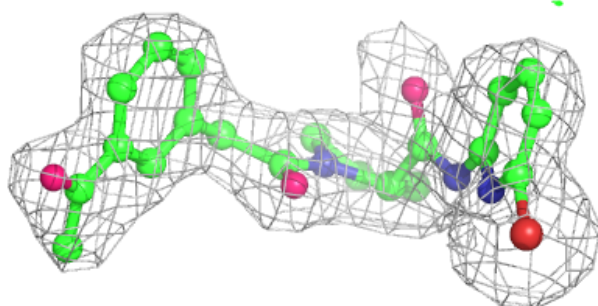
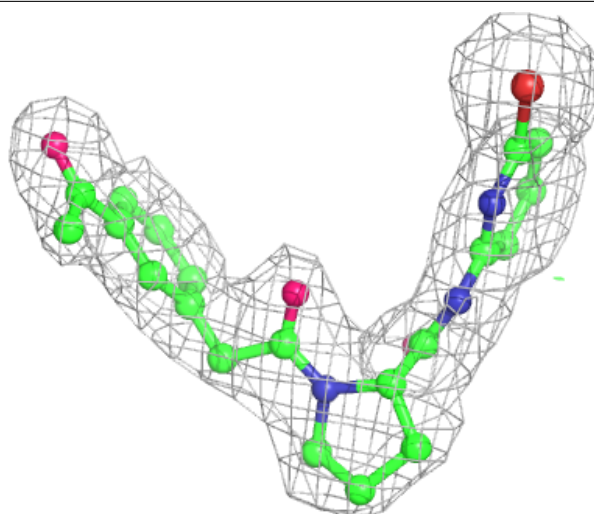
**Electron density around R7X B 301:**

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



**Electron density around R7X C 301:**

$2mF_o-DF_c$  (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.