wwPDB X-ray Structure Validation Summary Report (i)

Mar 5, 2024 – 06:17 AM EST

PDB ID	:	1ZGS
Title	:	Parkia platycephala seed lectin in complex with 5-bromo-4-chloro-3-indolyl-a-D-mannose
Authors	:	Gallego del Sol, F.; Cavada, B.S.; Calvete, J.J.
Deposited on	:	2005-04-22
Resolution	:	2.50 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 (i) 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 (i)) were used in the production of this report:

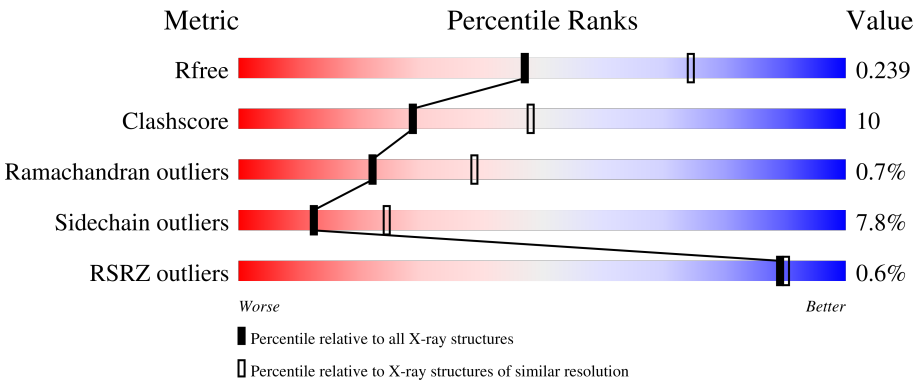
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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 <=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	447	75% 21% • ••
1	B	447	% 78% 17% • ••

2 Entry composition (i)

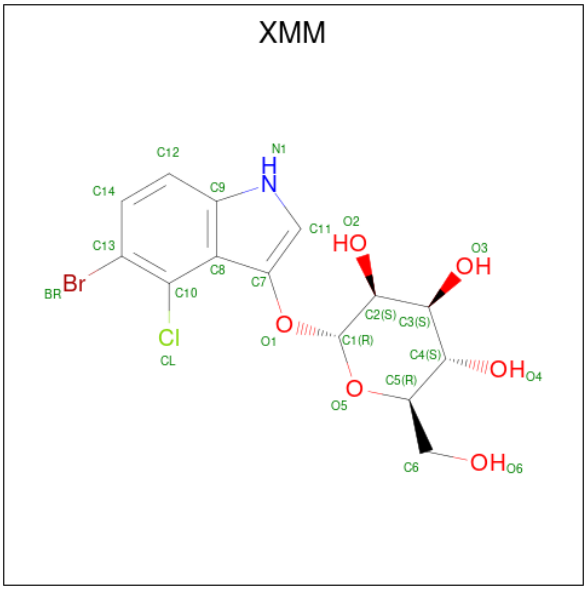
There are 3 unique types of molecules in this entry. The entry contains 7009 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Mannose/glucose-specific lectin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	442	3334	2142	543	646	3	36	0	0
1	B	442	3334	2142	543	646	3	30	0	0

- Molecule 2 is 5-bromo-4-chloro-1H-indol-3-yl alpha-D-mannopyranoside (three-letter code: XMM) (formula: $C_{14}H_{15}BrClNO_6$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
			Total	Br	C	Cl	N	O		
2	A	1	23	1	14	1	1	6	0	0
2	A	1	23	1	14	1	1	6	0	0
2	A	1	23	1	14	1	1	6	0	0
2	B	1	23	1	14	1	1	6	0	0

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Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
			Total	Br	C	Cl	N	O		
2	B	1	23	1	14	1	1	6	0	0
			Total	Br	C	Cl	N	O		
2	B	1	23	1	14	1	1	6	0	0

- Molecule 3 is water.

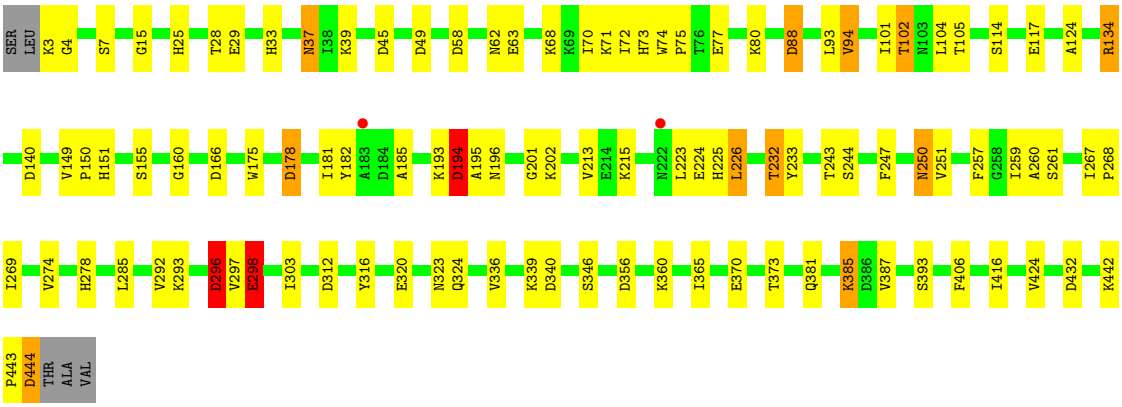
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	103	103	103	0	0
			Total	O		
3	B	100	100	100	0	0

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.

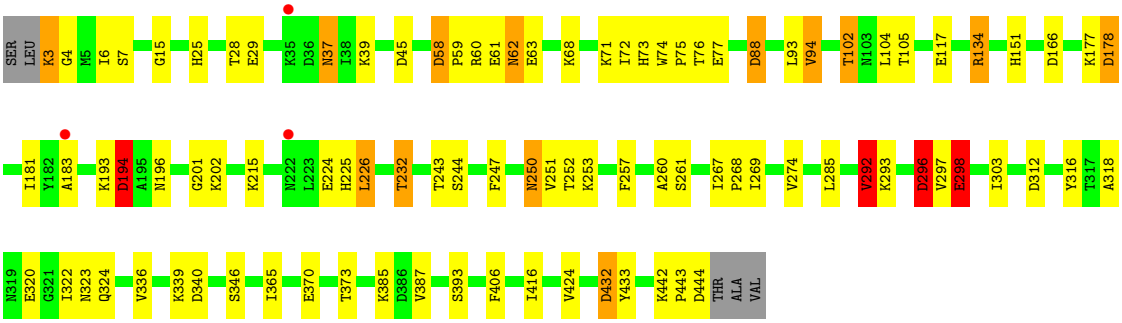
- Molecule 1: Mannose/glucose-specific lectin

Chain A:



- Molecule 1: Mannose/glucose-specific lectin

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	80.24Å 114.18Å 80.30Å 90.00° 119.88° 90.00°	Depositor
Resolution (Å)	30.00 – 2.50 69.63 – 2.50	Depositor EDS
% Data completeness (in resolution range)	100.0 (30.00-2.50) 98.1 (69.63-2.50)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	0.04	Depositor
$< I/\sigma(I) >$ [1]	9.31 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.2.0003	Depositor
R, R_{free}	0.208 , 0.244 0.212 , 0.239	Depositor DCC
R_{free} test set	3932 reflections (9.05%)	wwPDB-VP
Wilson B-factor (Å^2)	29.1	Xtriage
Anisotropy	0.165	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 2.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	0.017 for -h-l,k,h 0.017 for l,k,-h-l 0.018 for h,-k,-h-l 0.019 for -h-l,-k,l 0.487 for l,-k,h	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	7009	wwPDB-VP
Average B, all atoms (Å^2)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.15% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: XMM

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	1.27	6/3428 (0.2%)	0.95	18/4650 (0.4%)
1	B	1.27	5/3428 (0.1%)	0.97	18/4650 (0.4%)
All	All	1.27	11/6856 (0.2%)	0.96	36/9300 (0.4%)

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	3
1	B	0	3
All	All	0	6

The worst 5 of 11 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	298	GLU	C-N	-56.91	0.30	1.33
1	B	298	GLU	C-N	-55.75	0.32	1.33
1	B	296	ASP	C-N	-18.21	0.92	1.34
1	A	296	ASP	C-N	-18.19	0.92	1.34
1	B	292	VAL	CB-CG2	-7.95	1.36	1.52

The worst 5 of 36 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	3	LYS	O-C-N	-11.56	103.55	123.20
1	B	3	LYS	O-C-N	-11.53	103.60	123.20
1	B	3	LYS	CA-C-N	11.20	138.59	116.20
1	A	3	LYS	CA-C-N	11.08	138.37	116.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	3	LYS	C-N-CA	10.37	144.07	122.30

There are no chirality outliers.

5 of 6 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	201	GLY	Peptide
1	A	296	ASP	Mainchain
1	A	298	GLU	Mainchain
1	B	201	GLY	Peptide
1	B	296	ASP	Mainchain

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	3334	0	3200	69	0
1	B	3334	0	3200	59	0
2	A	69	0	42	0	0
2	B	69	0	42	2	0
3	A	103	0	0	4	0
3	B	100	0	0	1	0
All	All	7009	0	6484	128	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 10.

The worst 5 of 128 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:37:ASN:HD21	1:B:39:LYS:HZ3	1.13	0.92
1:B:37:ASN:HD21	1:B:39:LYS:NZ	1.70	0.88
1:A:37:ASN:HD21	1:A:39:LYS:NZ	1.73	0.86
1:A:37:ASN:HD21	1:A:39:LYS:HZ3	1.26	0.80
1:A:102:THR:CG2	1:A:104:LEU:H	1.95	0.78

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	440/447 (98%)	416 (94%)	22 (5%)	2 (0%)	29	48
1	B	440/447 (98%)	418 (95%)	18 (4%)	4 (1%)	17	31
All	All	880/894 (98%)	834 (95%)	40 (4%)	6 (1%)	22	39

5 of 6 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	298	GLU
1	B	298	GLU
1	B	4	GLY
1	B	62	ASN
1	B	194	ASP

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	353/358 (99%)	326 (92%)	27 (8%)	13	25
1	B	353/358 (99%)	325 (92%)	28 (8%)	12	24
All	All	706/716 (99%)	651 (92%)	55 (8%)	12	24

5 of 55 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	37	ASN
1	B	94	VAL
1	B	416	ILE
1	B	336	VAL
1	B	59	PRO

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 22 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	37	ASN
1	B	151	HIS
1	B	73	HIS
1	B	188	ASN
1	A	188	ASN

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	XMM	A	450	-	24,25,25	5.40	7 (29%)	26,37,37	6.55	13 (50%)
2	XMM	A	451	-	24,25,25	5.35	7 (29%)	26,37,37	6.12	12 (46%)
2	XMM	A	452	-	24,25,25	5.05	7 (29%)	26,37,37	5.93	11 (42%)
2	XMM	B	453	-	24,25,25	5.08	8 (33%)	26,37,37	6.17	12 (46%)
2	XMM	B	455	-	24,25,25	5.45	7 (29%)	26,37,37	6.48	12 (46%)
2	XMM	B	454	-	24,25,25	5.31	7 (29%)	26,37,37	6.06	8 (30%)

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	XMM	A	450	-	-	0/4/26/26	0/3/3/3
2	XMM	A	451	-	-	0/4/26/26	0/3/3/3
2	XMM	A	452	-	-	4/4/26/26	0/3/3/3
2	XMM	B	453	-	-	4/4/26/26	0/3/3/3
2	XMM	B	455	-	-	2/4/26/26	0/3/3/3
2	XMM	B	454	-	-	0/4/26/26	0/3/3/3

The worst 5 of 43 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	455	XMM	BR-C13	-19.63	1.45	1.89
2	A	450	XMM	BR-C13	-19.49	1.45	1.89
2	A	451	XMM	BR-C13	-19.21	1.46	1.89
2	B	454	XMM	BR-C13	-18.98	1.46	1.89
2	B	453	XMM	BR-C13	-18.15	1.48	1.89

The worst 5 of 68 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	450	XMM	C12-C14-C13	26.63	153.70	119.91
2	B	454	XMM	C12-C14-C13	26.55	153.59	119.91
2	B	455	XMM	C12-C14-C13	26.46	153.47	119.91
2	A	451	XMM	C12-C14-C13	26.29	153.26	119.91
2	B	453	XMM	C12-C14-C13	26.16	153.10	119.91

There are no chirality outliers.

5 of 10 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	455	XMM	O5-C5-C6-O6
2	B	455	XMM	C4-C5-C6-O6
2	B	453	XMM	O5-C5-C6-O6
2	B	453	XMM	C4-C5-C6-O6
2	A	452	XMM	C2-C1-O1-C7

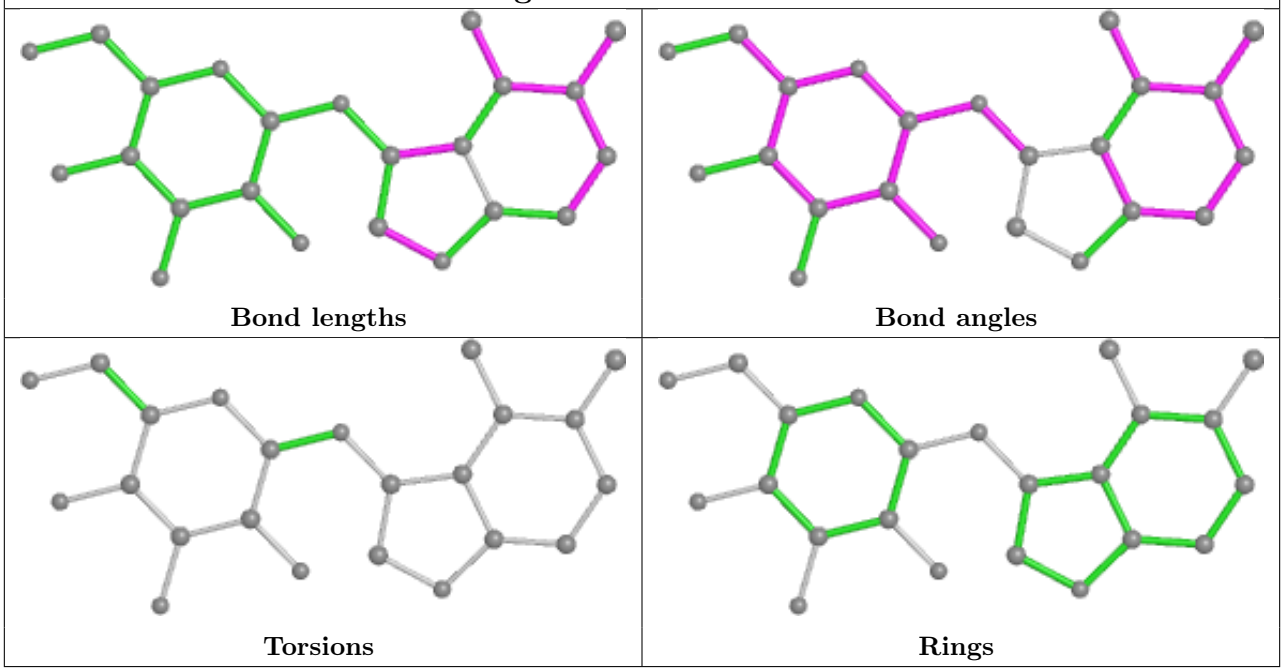
There are no ring outliers.

1 monomer is involved in 2 short contacts:

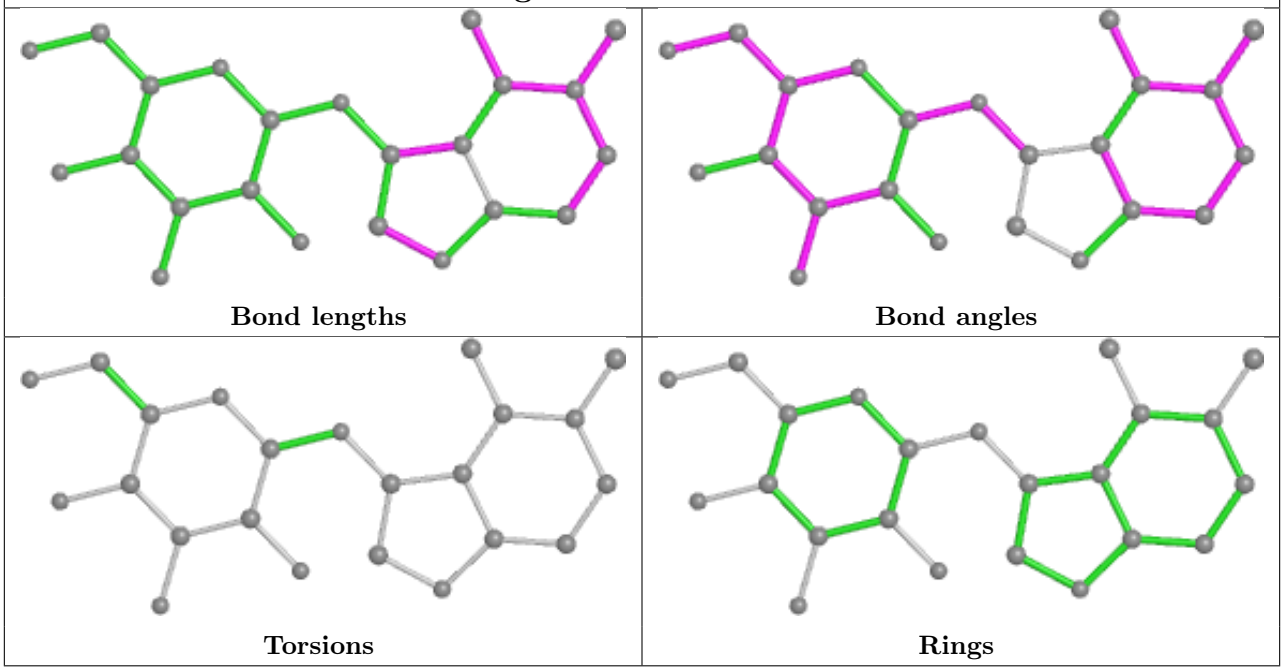
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	455	XMM	2	0

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 then 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.

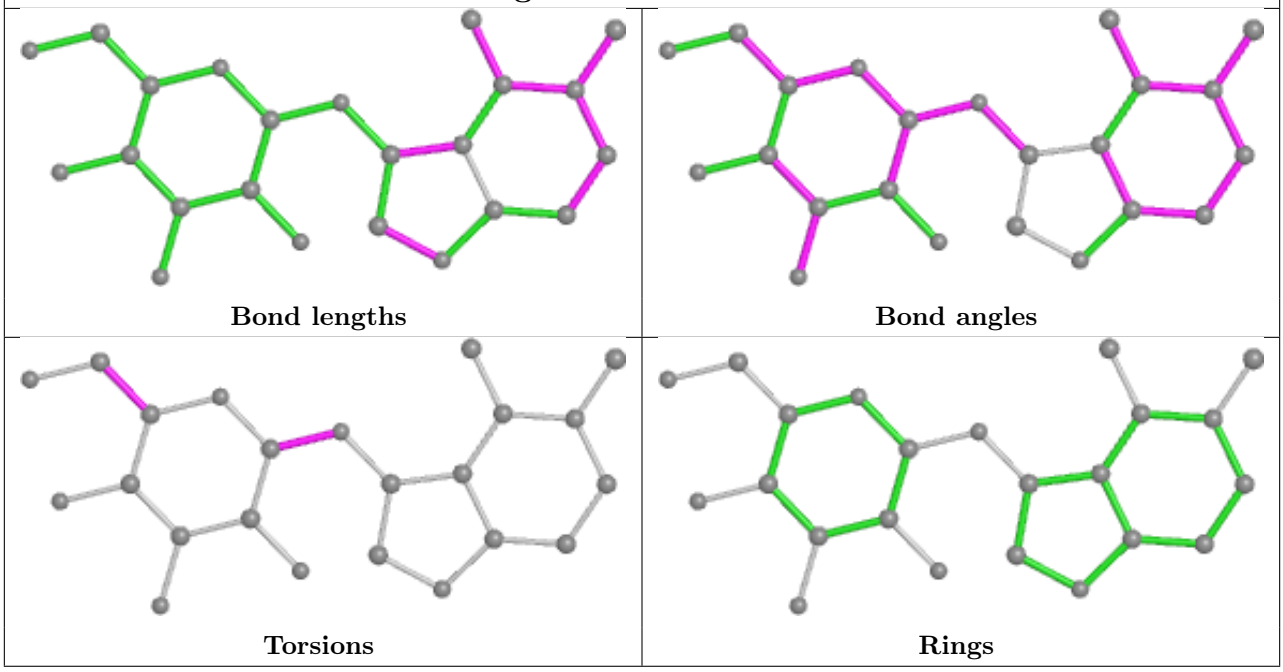
Ligand XMM A 450



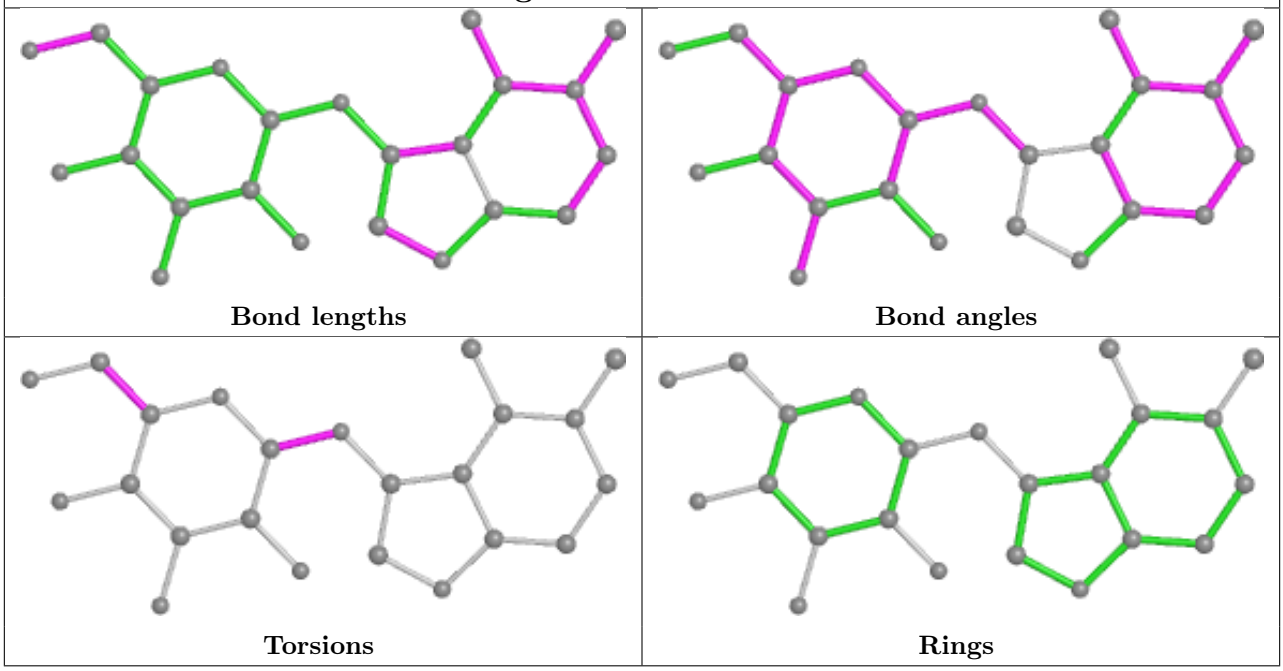
Ligand XMM A 451



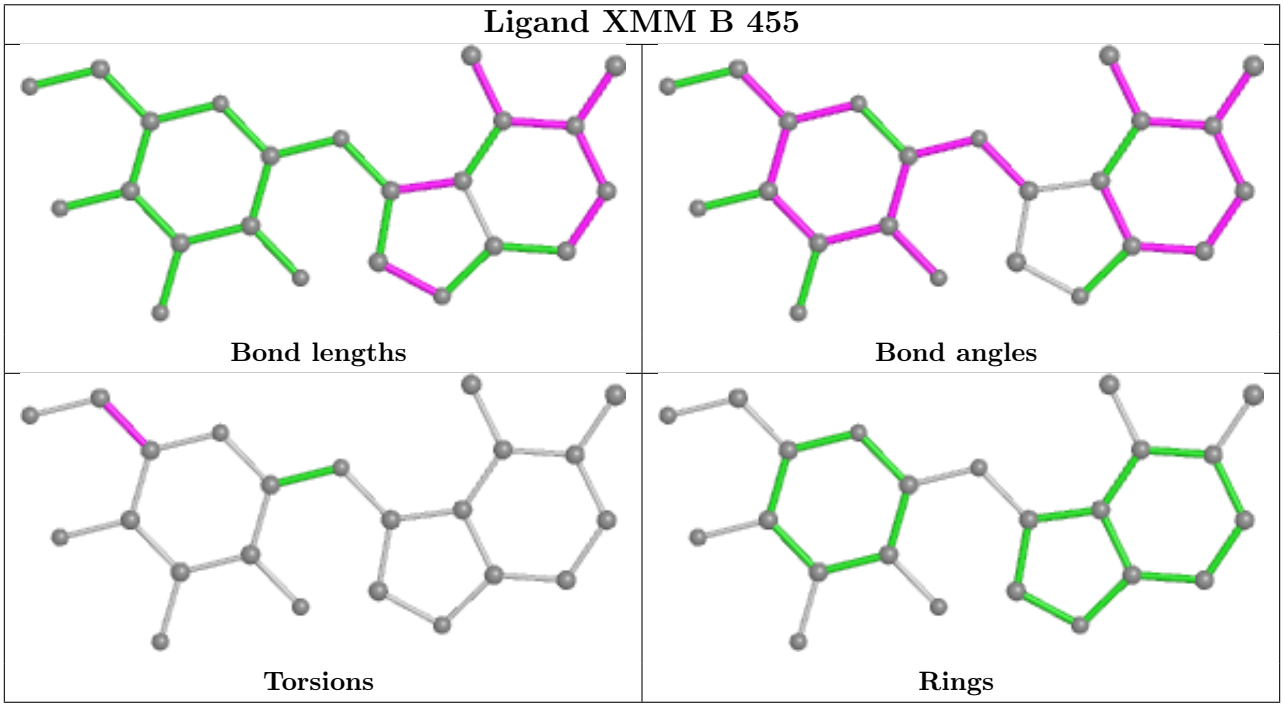
Ligand XMM A 452

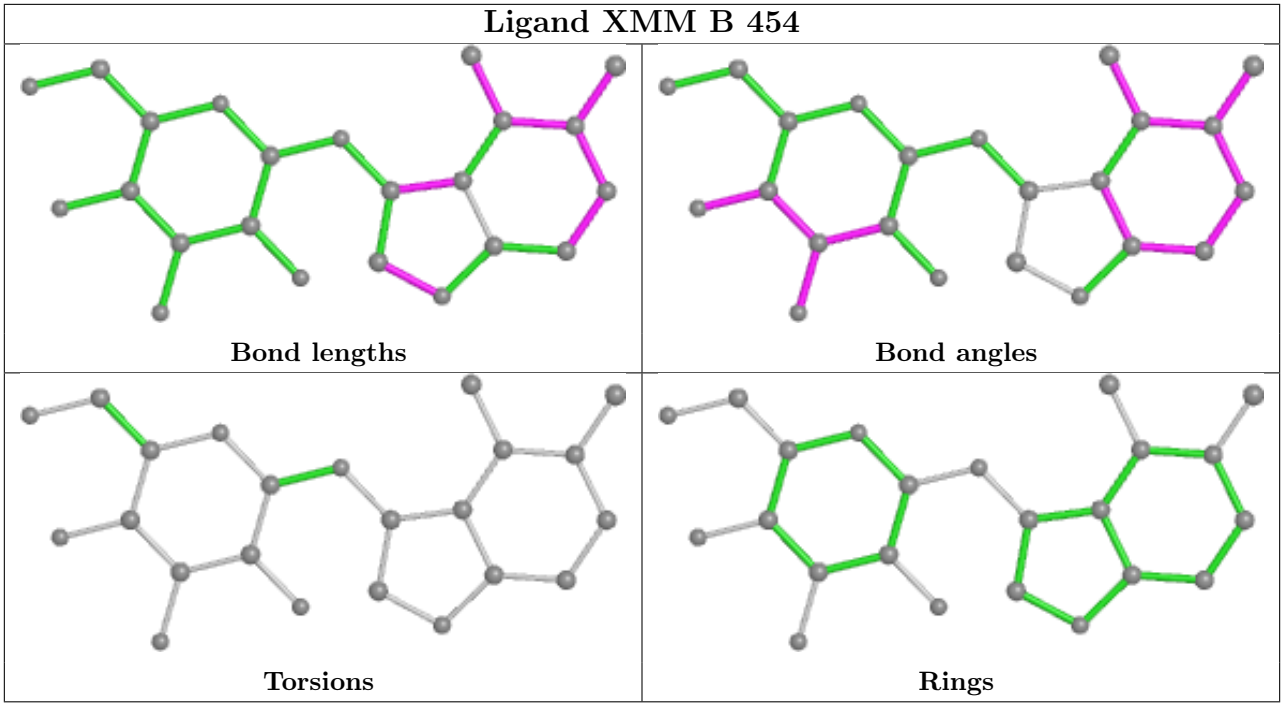


Ligand XMM B 453



Ligand XMM B 455





5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	3
1	B	2

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	3:LYS	C	4:GLY	N	1.19
1	A	296:ASP	C	297:VAL	N	0.92
1	B	296:ASP	C	297:VAL	N	0.92
1	B	298:GLU	C	299:GLY	N	0.32
1	A	298:GLU	C	299:GLY	N	0.30

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	439/447 (98%)	-0.30	2 (0%)	91	91	9, 24, 41, 54	4 (0%)
1	B	439/447 (98%)	-0.30	3 (0%)	87	89	9, 24, 41, 54	2 (0%)
All	All	878/894 (98%)	-0.30	5 (0%)	89	90	9, 24, 41, 54	6 (0%)

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	222	ASN	2.8
1	A	222	ASN	2.3
1	A	183	ALA	2.1
1	B	35	LYS	2.1
1	B	183	ALA	2.1

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 (i)

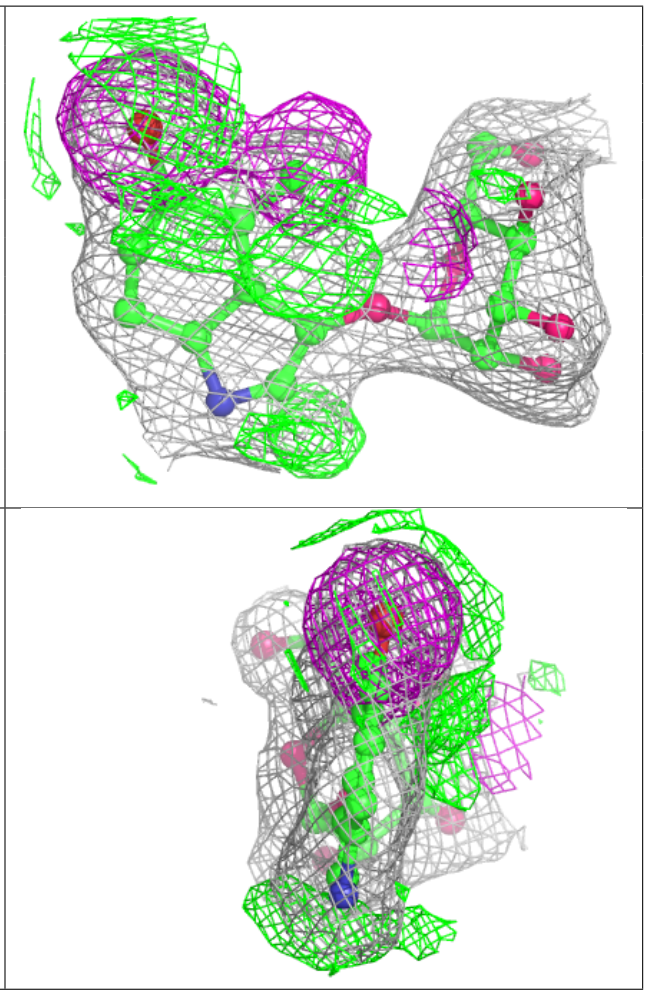
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	XMM	A	452	23/23	0.83	0.17	11,20,25,29	0
2	XMM	B	453	23/23	0.83	0.17	12,21,27,32	0
2	XMM	A	451	23/23	0.87	0.17	8,18,26,36	0
2	XMM	B	454	23/23	0.87	0.16	9,18,24,28	0
2	XMM	B	455	23/23	0.87	0.16	11,22,30,35	0
2	XMM	A	450	23/23	0.90	0.15	10,21,32,38	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.

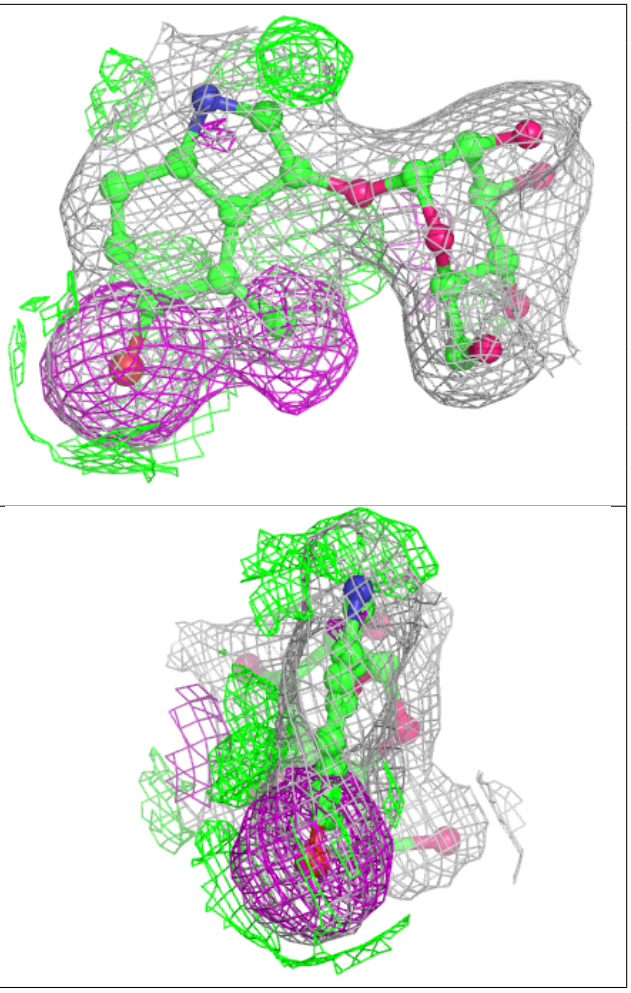
Electron density around XMM A 452:

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



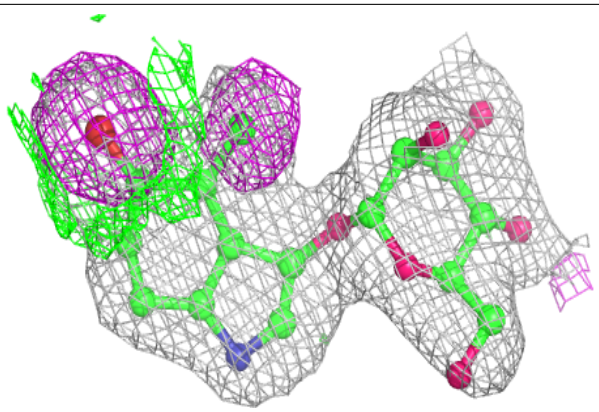
Electron density around XMM B 453:

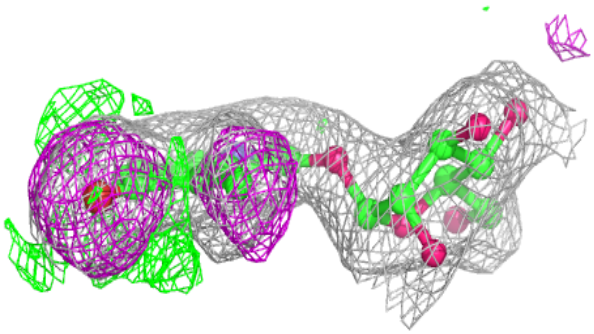
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

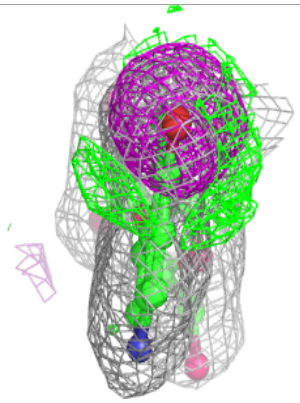


Electron density around XMM A 451:

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

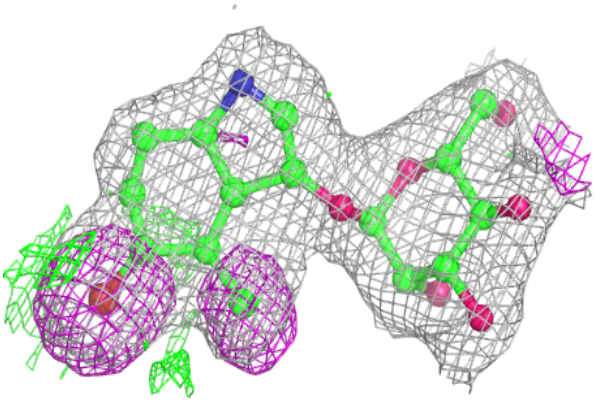


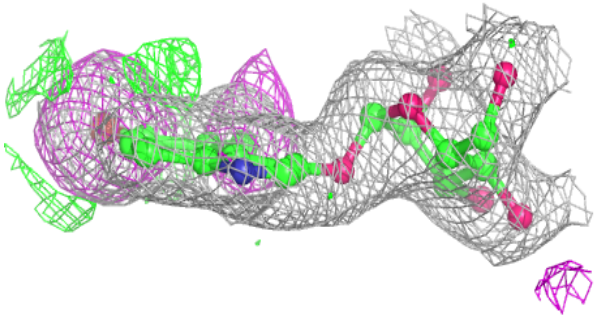


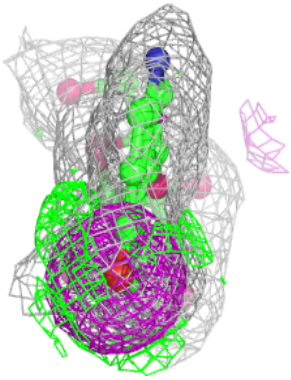


Electron density around XMM B 454:

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

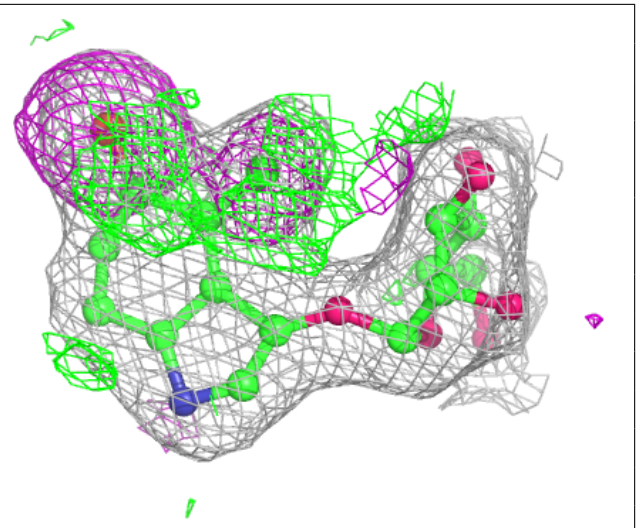


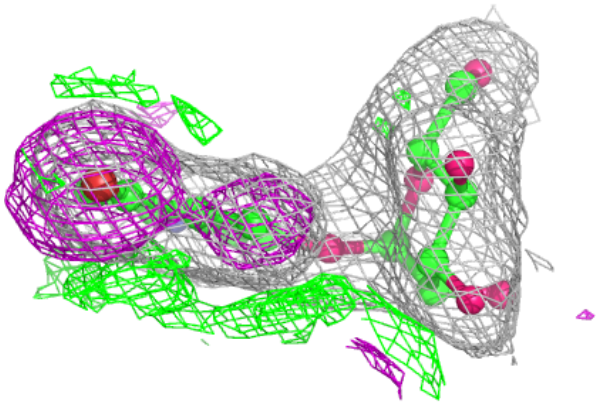


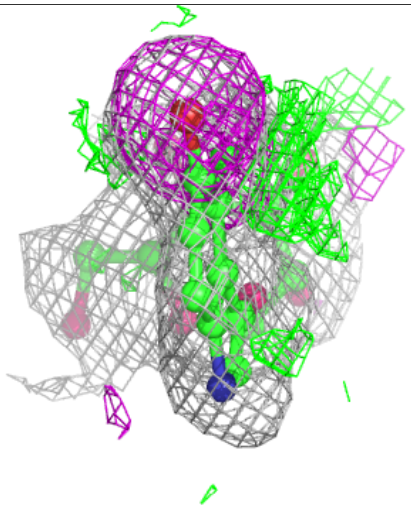


Electron density around XMM B 455:

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

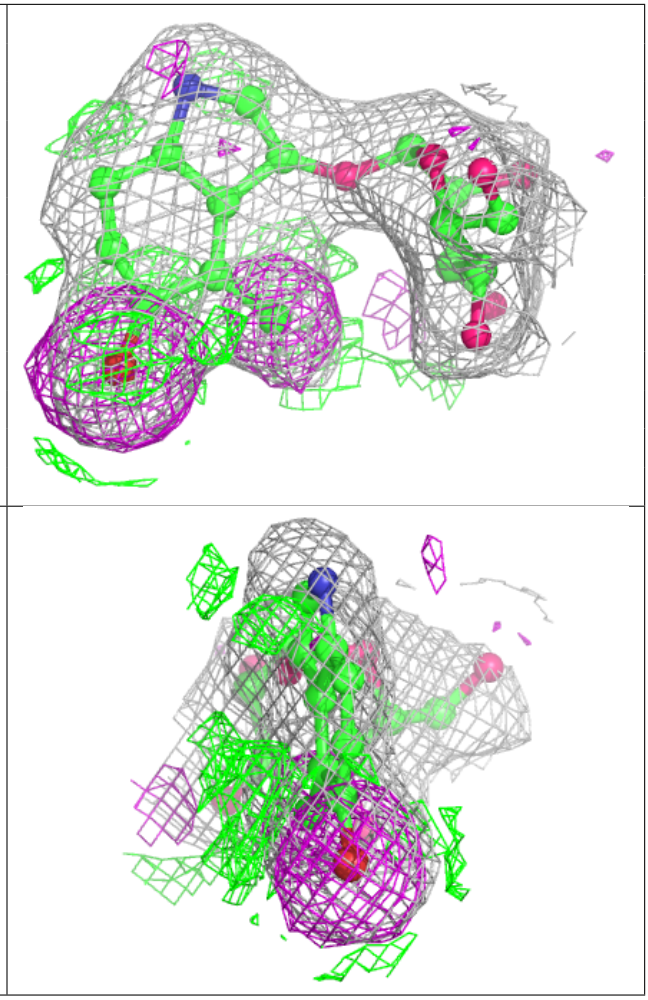






Electron density around XMM A 450:

$2mF_o$-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.