



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 20, 2024 – 12:28 pm GMT

PDB ID : 8PXT
Title : Targeting extended blood antigens by Akkermansia muciniphila enzymes unveils a missing link for generating universal donor blood
Authors : Jensen, M.; Abou Hachem, M.; Morth, J.P.
Deposited on : 2023-07-24
Resolution : 2.25 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 ⓘ 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 : 4.02b-467
Mogul : 1.8.4, 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

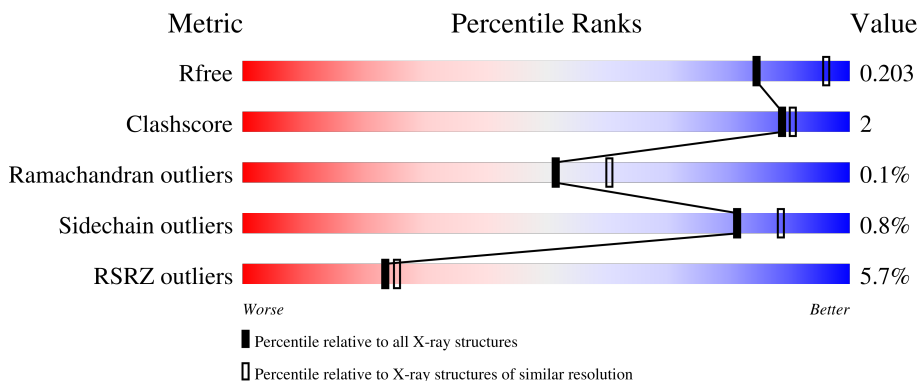
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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	1377 (2.26-2.26)
Clashscore	141614	1487 (2.26-2.26)
Ramachandran outliers	138981	1449 (2.26-2.26)
Sidechain outliers	138945	1450 (2.26-2.26)
RSRZ outliers	127900	1356 (2.26-2.26)

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	655	 5% 92% 5% •
1	B	655	 10% 90% 7% •
1	C	655	 2% 93% • •

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 30647 atoms, of which 14980 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 Beta-N-acetylhexosaminidase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	636	9980	3194	4962	879	928	17	0	0	0
1	B	635	9974	3192	4960	878	927	17	0	0	0
1	C	634	9957	3186	4953	875	926	17	0	0	0

There are 30 discrepancies between the modelled and reference sequences:

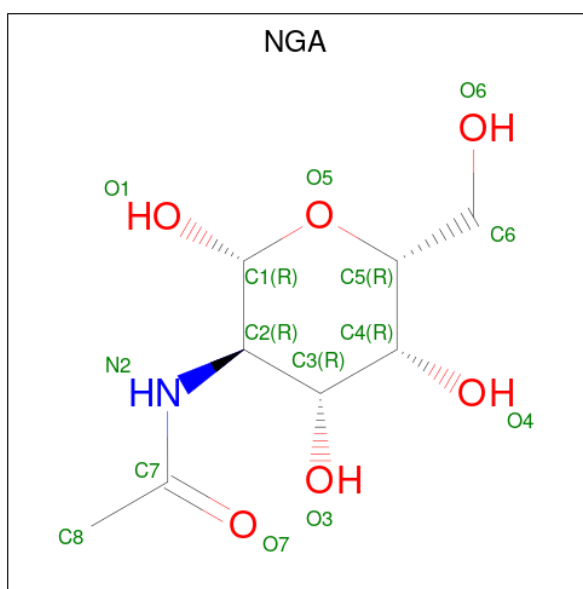
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP B2UN02
A	321	ALA	GLU	conflict	UNP B2UN02
A	648	LEU	-	expression tag	UNP B2UN02
A	649	GLU	-	expression tag	UNP B2UN02
A	650	HIS	-	expression tag	UNP B2UN02
A	651	HIS	-	expression tag	UNP B2UN02
A	652	HIS	-	expression tag	UNP B2UN02
A	653	HIS	-	expression tag	UNP B2UN02
A	654	HIS	-	expression tag	UNP B2UN02
A	655	HIS	-	expression tag	UNP B2UN02
B	1	MET	-	initiating methionine	UNP B2UN02
B	321	ALA	GLU	conflict	UNP B2UN02
B	648	LEU	-	expression tag	UNP B2UN02
B	649	GLU	-	expression tag	UNP B2UN02
B	650	HIS	-	expression tag	UNP B2UN02
B	651	HIS	-	expression tag	UNP B2UN02
B	652	HIS	-	expression tag	UNP B2UN02
B	653	HIS	-	expression tag	UNP B2UN02
B	654	HIS	-	expression tag	UNP B2UN02
B	655	HIS	-	expression tag	UNP B2UN02
C	1	MET	-	initiating methionine	UNP B2UN02
C	321	ALA	GLU	conflict	UNP B2UN02
C	648	LEU	-	expression tag	UNP B2UN02

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Chain	Residue	Modelled	Actual	Comment	Reference
C	649	GLU	-	expression tag	UNP B2UN02
C	650	HIS	-	expression tag	UNP B2UN02
C	651	HIS	-	expression tag	UNP B2UN02
C	652	HIS	-	expression tag	UNP B2UN02
C	653	HIS	-	expression tag	UNP B2UN02
C	654	HIS	-	expression tag	UNP B2UN02
C	655	HIS	-	expression tag	UNP B2UN02

- Molecule 2 is 2-acetamido-2-deoxy-beta-D-galactopyranose (three-letter code: NGA) (formula: C₈H₁₅NO₆) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
			Total	C	H	N			O
2	A	1	Total	C	H	N	O	0	0
			30	8	15	1	6		
2	A	1	Total	C	H	N	O	0	0
			30	8	15	1	6		
2	A	1	Total	C	H	N	O	0	0
			30	8	15	1	6		
2	B	1	Total	C	H	N	O	0	0
			30	8	15	1	6		
2	B	1	Total	C	H	N	O	0	0
			30	8	15	1	6		
2	C	1	Total	C	H	N	O	0	0
			30	8	15	1	6		
2	C	1	Total	C	H	N	O	0	0
			30	8	15	1	6		

- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total 2	Na 2	0	0
3	C	2	Total 2	Na 2	0	0

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	1	Total 1	Cl 1	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	162	Total 162	O 162	0	0
5	B	167	Total 167	O 167	0	0
5	C	192	Total 192	O 192	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	92.23Å 169.06Å 171.97Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	62.32 – 2.25 62.32 – 2.25	Depositor EDS
% Data completeness (in resolution range)	100.0 (62.32-2.25) 100.0 (62.32-2.25)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.61 (at 2.25Å)	Xtrriage
Refinement program	PHENIX 1.19.2_4158, PHENIX 1.19.2_4158	Depositor
R, R_{free}	0.180 , 0.210 0.174 , 0.203	Depositor DCC
R_{free} test set	6408 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	47.0	Xtrriage
Anisotropy	0.352	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.43 , 45.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.000 for -h,l,k	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	30647	wwPDB-VP
Average B, all atoms (Å ²)	63.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.55% 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: NGA, NA, CL

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	0.27	0/5145	0.53	0/6965
1	B	0.26	0/5141	0.52	0/6960
1	C	0.27	0/5130	0.52	0/6945
All	All	0.27	0/15416	0.52	0/20870

There are no bond length outliers.

There are no bond angle outliers.

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	5018	4962	4961	18	0
1	B	5014	4960	4959	27	1
1	C	5004	4953	4951	15	1
2	A	45	45	43	0	0
2	B	30	30	28	0	0
2	C	30	30	28	2	0
3	A	2	0	0	0	0
3	C	2	0	0	0	0
4	C	1	0	0	0	0
5	A	162	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	B	167	0	0	0	0
5	C	192	0	0	3	0
All	All	15667	14980	14970	58	1

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.

The worst 5 of 58 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:186:LEU:CD1	1:B:194:ILE:HD11	2.20	0.70
1:B:186:LEU:HD13	1:B:194:ILE:HD11	1.75	0.69
1:C:120:VAL:HG23	5:C:975:HOH:O	1.93	0.67
1:B:29:ILE:HD11	1:B:72:VAL:HG22	1.77	0.65
1:A:566:HIS:ND1	1:A:646:THR:HG22	2.13	0.63

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:333:GLU:OE2	1:C:590:SER:OG[3_645]	2.18	0.02

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	632/655 (96%)	613 (97%)	19 (3%)	0	100 100
1	B	631/655 (96%)	613 (97%)	18 (3%)	0	100 100
1	C	630/655 (96%)	614 (98%)	15 (2%)	1 (0%)	47 55
All	All	1893/1965 (96%)	1840 (97%)	52 (3%)	1 (0%)	51 60

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	3	GLU

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	527/541 (97%)	522 (99%)	5 (1%)	78	86
1	B	527/541 (97%)	523 (99%)	4 (1%)	81	88
1	C	526/541 (97%)	523 (99%)	3 (1%)	86	91
All	All	1580/1623 (97%)	1568 (99%)	12 (1%)	81	88

5 of 12 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	184	TRP
1	B	255	GLU
1	C	372	ASN
1	C	169	GLN
1	A	535	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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](#)

Of 12 ligands modelled in this entry, 5 are monoatomic - leaving 7 for Mogul analysis.

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	NGA	A	702	-	15,15,15	1.82	4 (26%)	21,21,21	1.61	4 (19%)
2	NGA	A	701	-	15,15,15	1.76	4 (26%)	21,21,21	1.26	2 (9%)
2	NGA	A	703	-	15,15,15	1.89	4 (26%)	21,21,21	1.40	3 (14%)
2	NGA	B	701	-	15,15,15	1.78	4 (26%)	21,21,21	0.98	1 (4%)
2	NGA	C	701	-	15,15,15	1.82	4 (26%)	21,21,21	1.24	3 (14%)
2	NGA	B	702	-	15,15,15	1.72	4 (26%)	21,21,21	1.16	3 (14%)
2	NGA	C	702	-	15,15,15	1.84	4 (26%)	21,21,21	1.24	4 (19%)

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	NGA	A	702	-	-	2/6/26/26	0/1/1/1
2	NGA	A	701	-	-	2/6/26/26	0/1/1/1
2	NGA	A	703	-	-	1/6/26/26	0/1/1/1
2	NGA	B	701	-	-	2/6/26/26	0/1/1/1
2	NGA	C	701	-	-	1/6/26/26	0/1/1/1
2	NGA	B	702	-	-	2/6/26/26	0/1/1/1
2	NGA	C	702	-	-	2/6/26/26	0/1/1/1

The worst 5 of 28 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	701	NGA	C7-N2	3.77	1.47	1.34
2	C	702	NGA	C7-N2	3.77	1.47	1.34
2	A	703	NGA	C7-N2	3.69	1.47	1.34
2	B	701	NGA	C7-N2	3.69	1.47	1.34
2	A	702	NGA	C7-N2	3.65	1.46	1.34

The worst 5 of 20 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	702	NGA	C3-C4-C5	4.54	118.33	110.24
2	A	703	NGA	O5-C1-C2	4.07	113.61	109.52
2	A	701	NGA	O5-C1-C2	3.81	113.34	109.52
2	A	702	NGA	C4-C3-C2	3.26	115.11	110.34
2	A	702	NGA	O5-C5-C4	2.72	114.63	109.69

There are no chirality outliers.

5 of 12 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	703	NGA	C3-C2-N2-C7
2	C	702	NGA	C1-C2-N2-C7
2	B	701	NGA	C4-C5-C6-O6
2	B	701	NGA	O5-C5-C6-O6
2	C	701	NGA	O5-C5-C6-O6

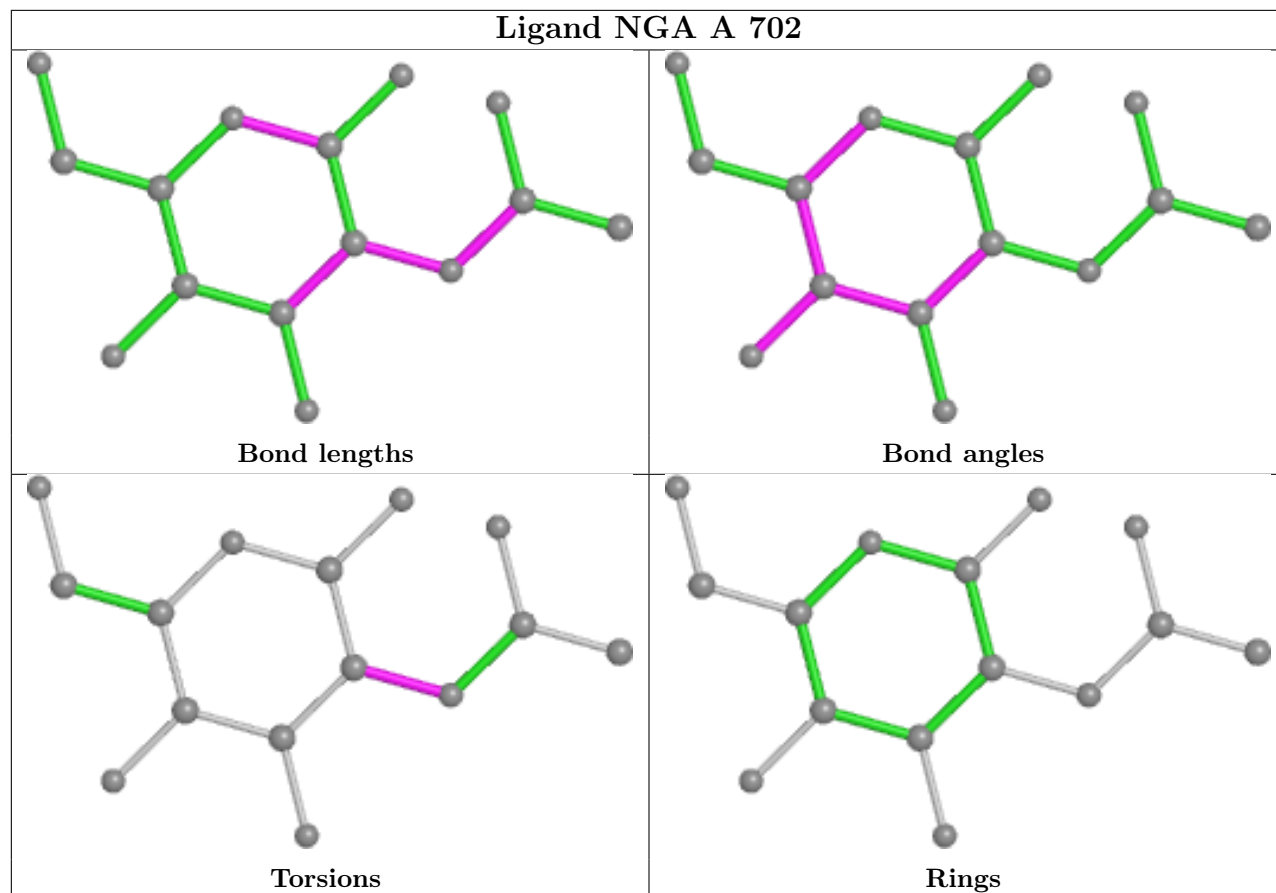
There are no ring outliers.

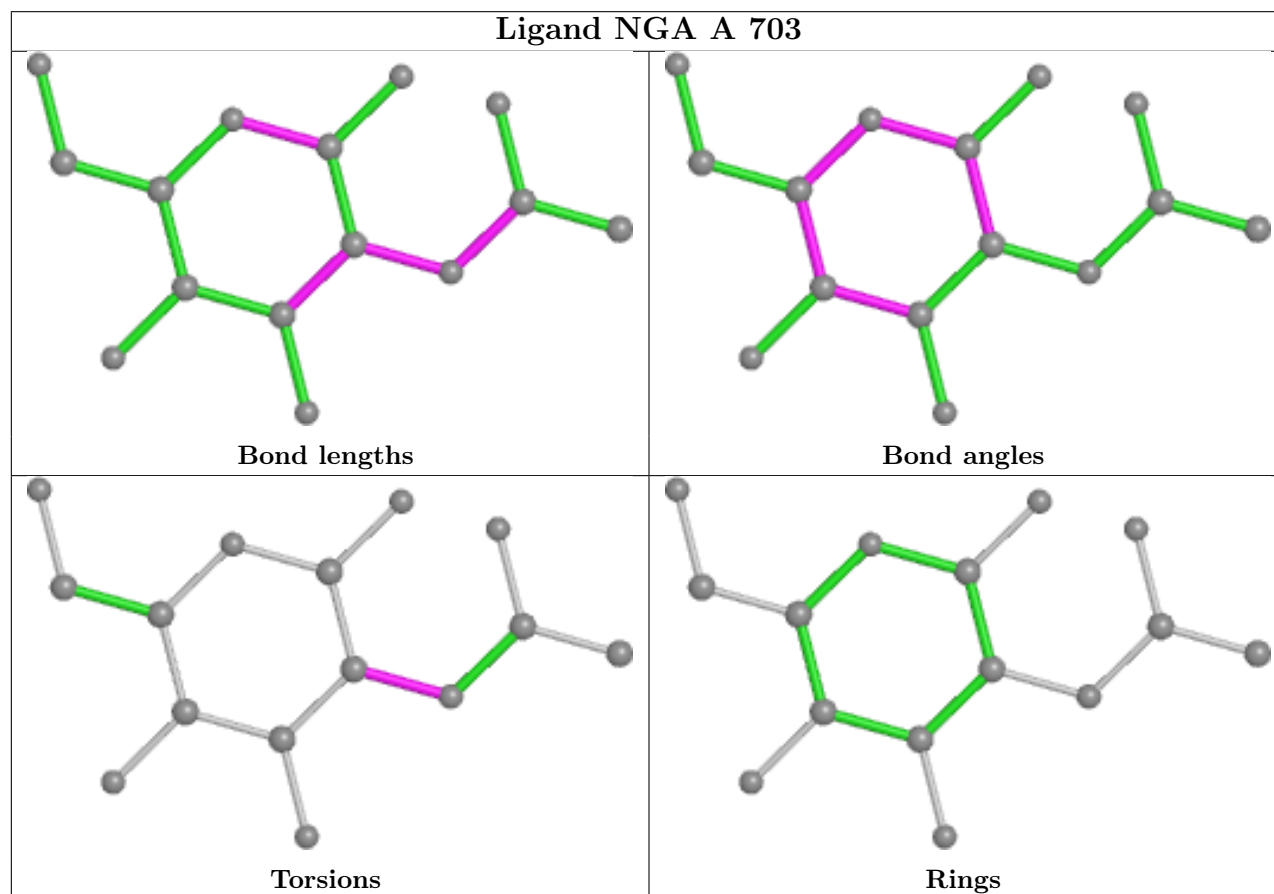
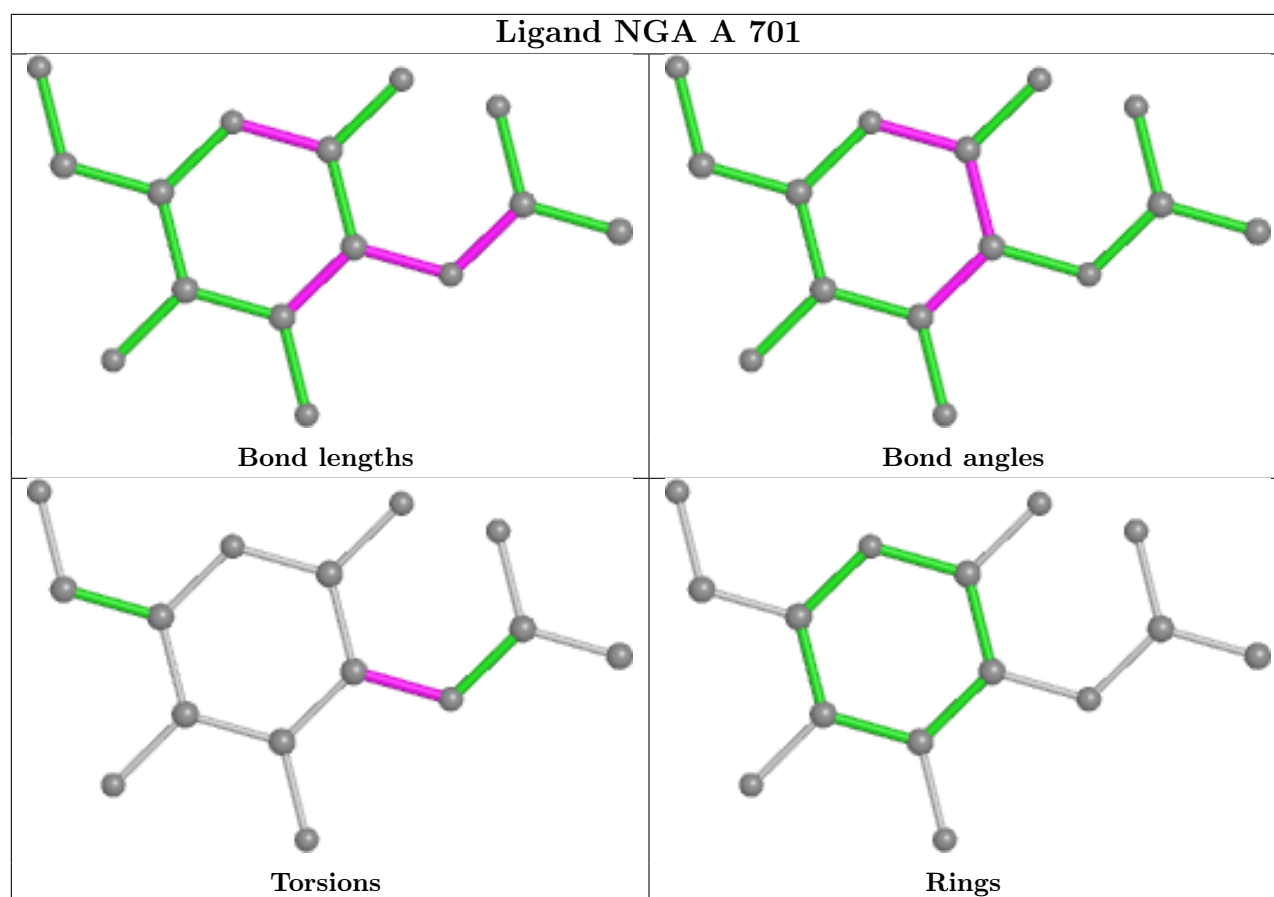
2 monomers are involved in 2 short contacts:

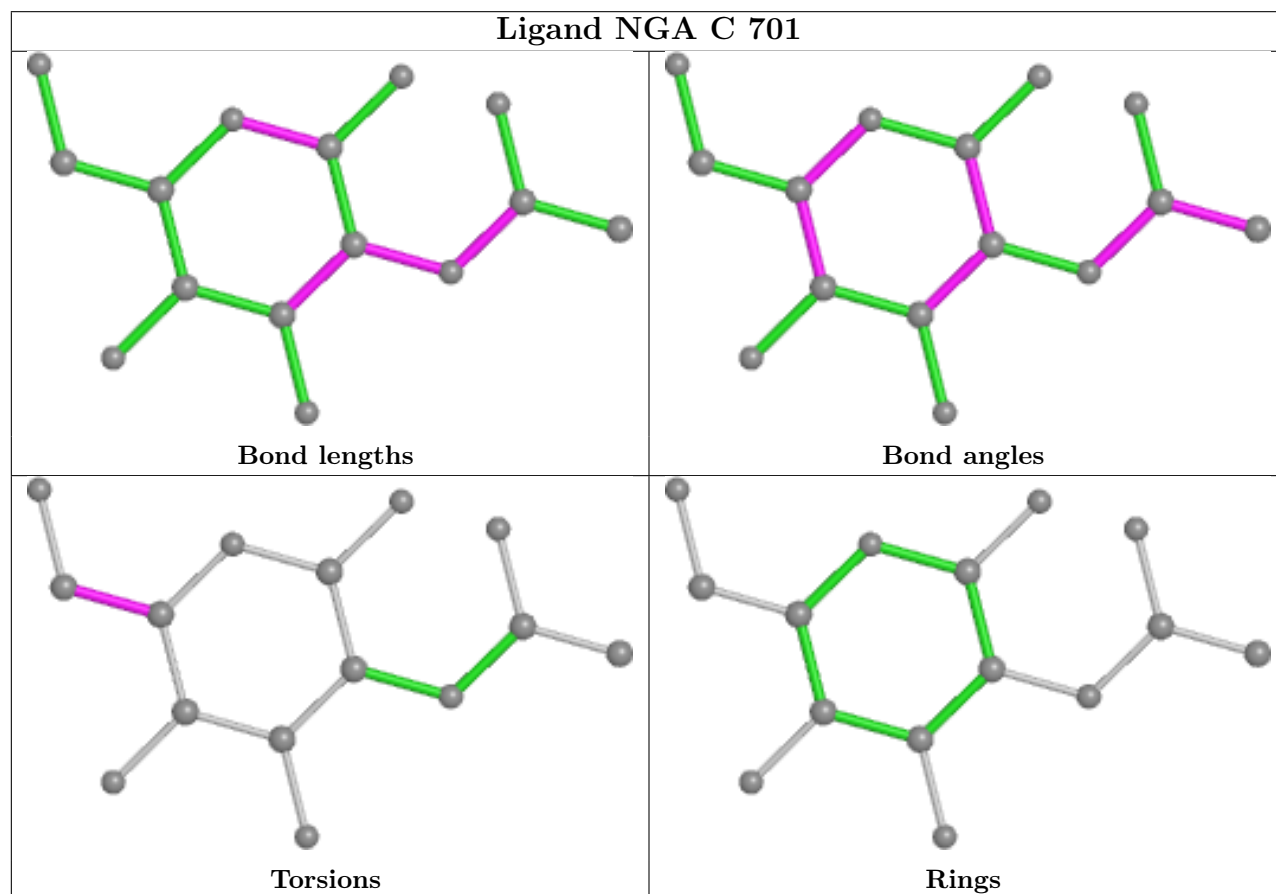
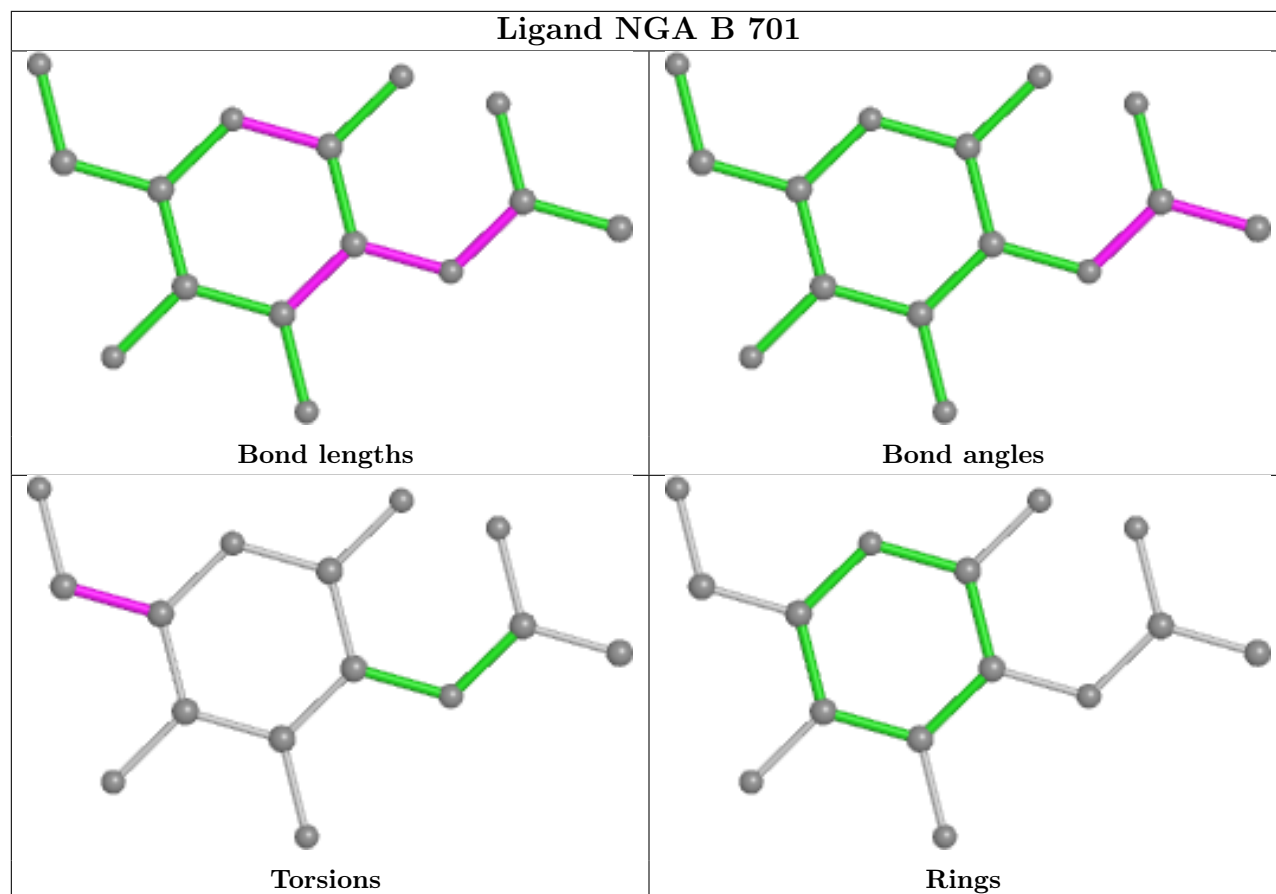
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	701	NGA	1	0
2	C	702	NGA	1	0

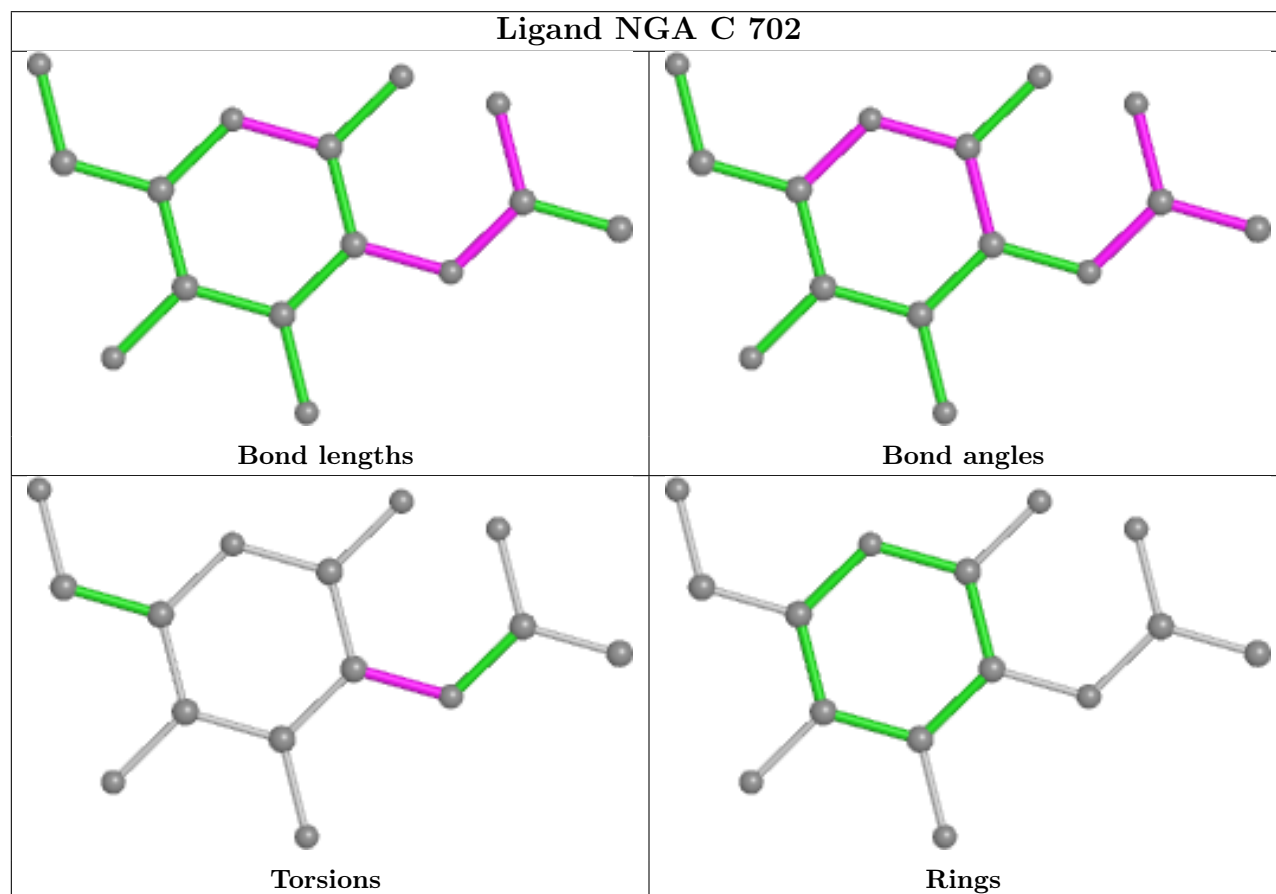
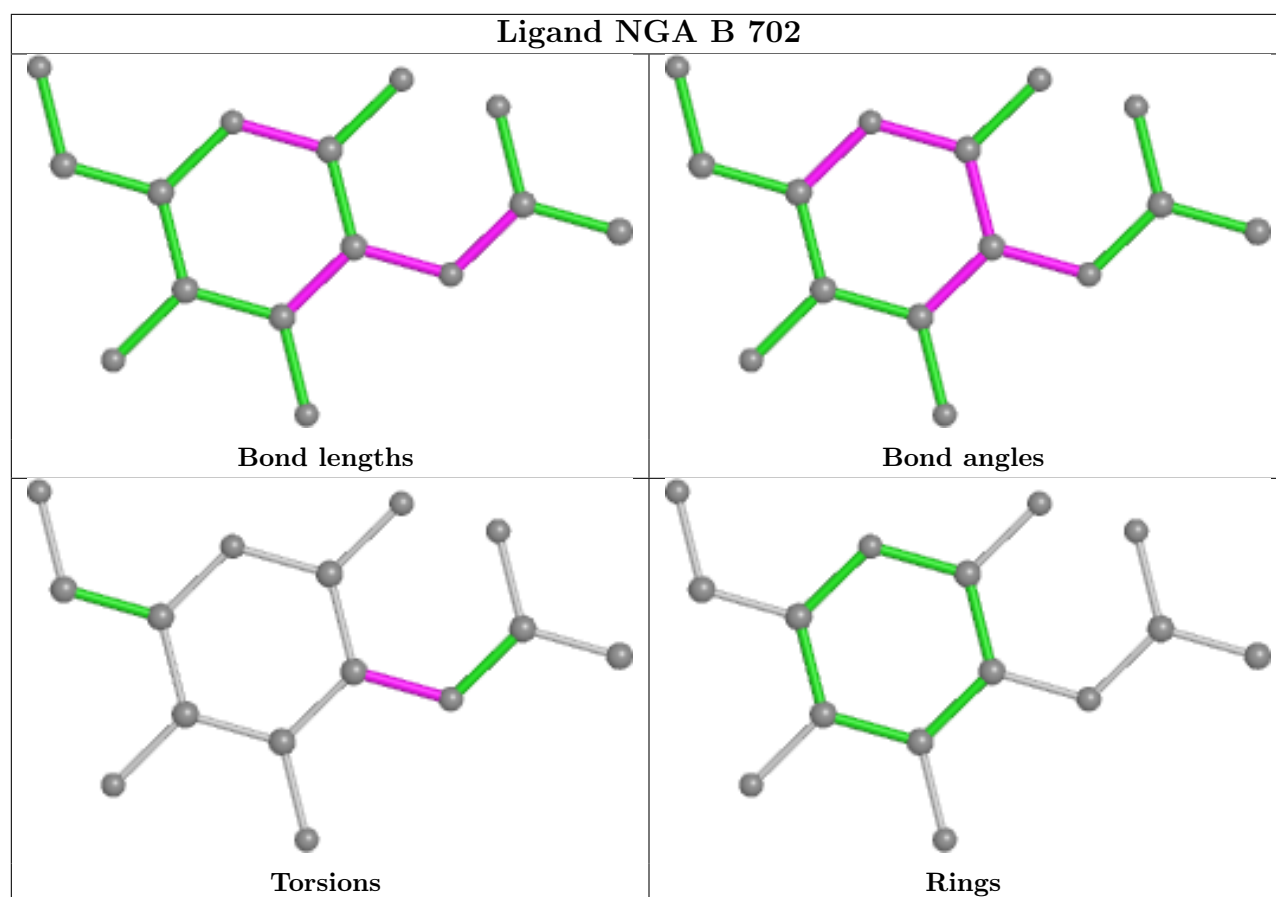
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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	636/655 (97%)	0.56	35 (5%) 25 27	36, 54, 91, 153	0
1	B	635/655 (96%)	0.62	64 (10%) 7 7	38, 57, 104, 156	0
1	C	634/655 (96%)	0.30	10 (1%) 72 74	34, 49, 78, 126	0
All	All	1905/1965 (96%)	0.50	109 (5%) 23 25	34, 53, 96, 156	0

The worst 5 of 109 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	66	LYS	8.5
1	A	64	ALA	8.3
1	A	389	ALA	7.6
1	A	388	GLY	6.3
1	B	72	VAL	6.2

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](#)

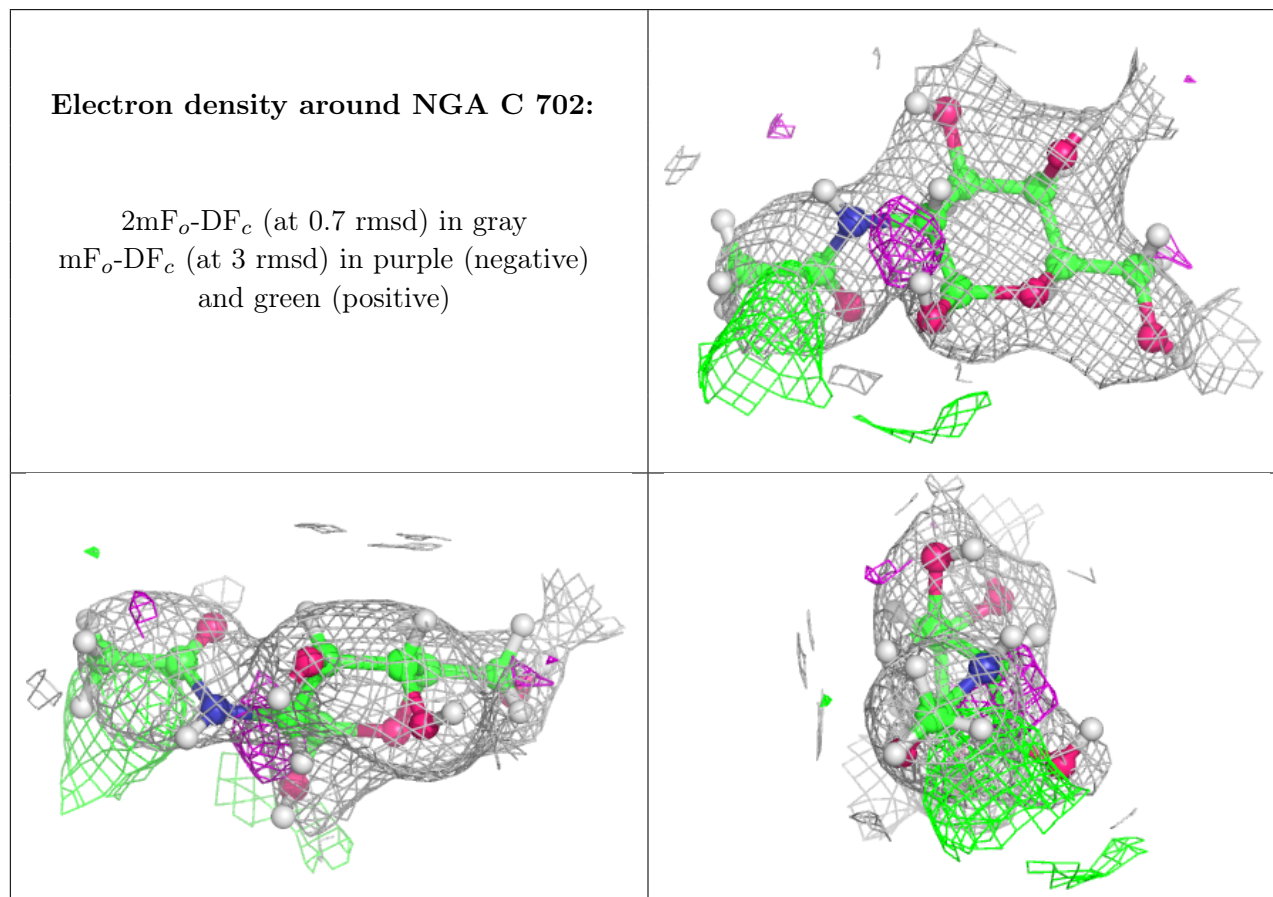
There are no monosaccharides in this entry.

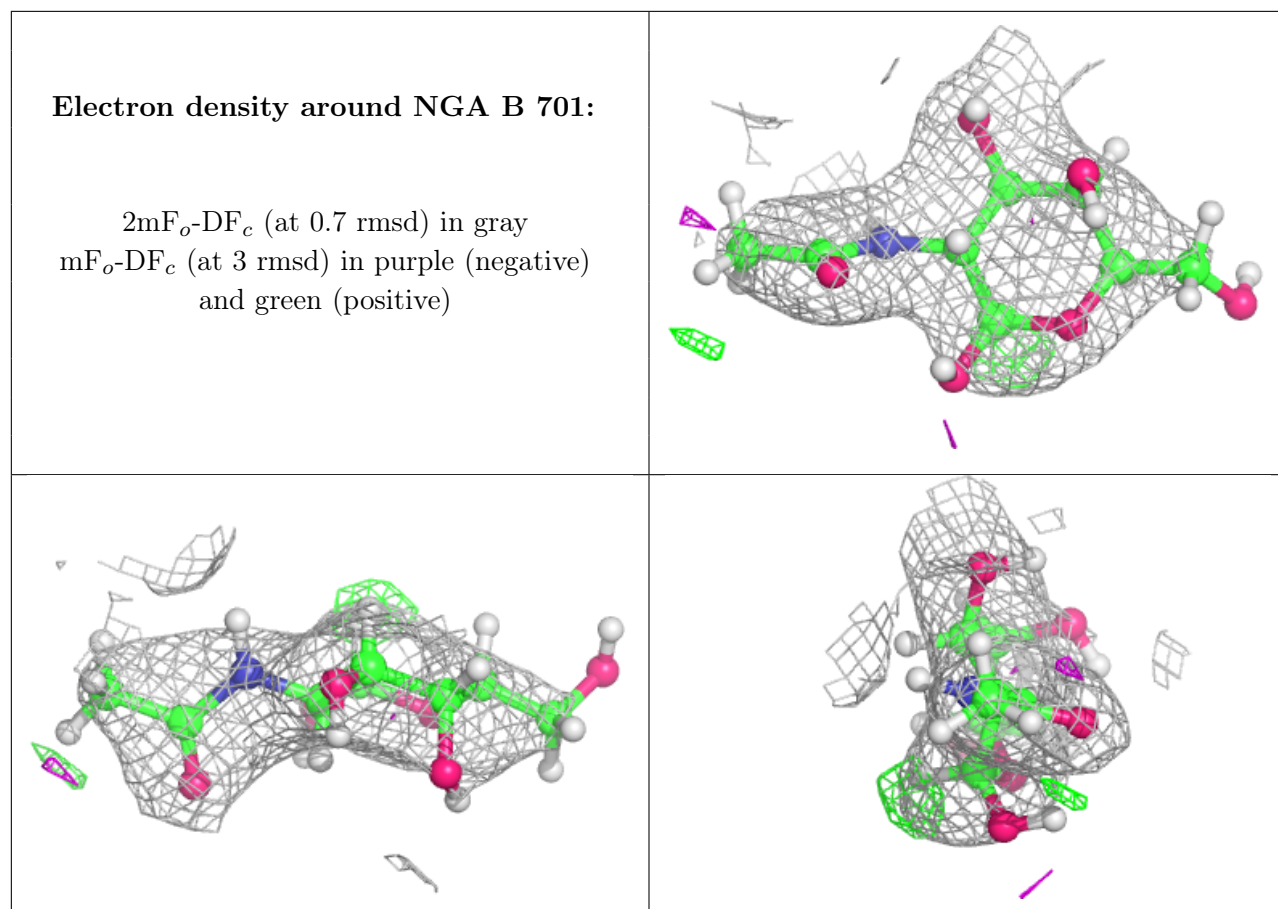
6.4 Ligands [i](#)

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(\AA^2)	Q<0.9
3	NA	A	704	1/1	0.80	0.25	54,54,54,54	0
4	CL	C	705	1/1	0.83	0.09	85,85,85,85	0
2	NGA	C	702	15/15	0.86	0.18	46,57,69,80	0
3	NA	C	703	1/1	0.87	0.32	59,59,59,59	0
2	NGA	B	701	15/15	0.90	0.18	61,76,96,99	0
2	NGA	A	703	15/15	0.91	0.19	55,78,100,111	0
3	NA	C	704	1/1	0.91	0.12	55,55,55,55	0
2	NGA	C	701	15/15	0.91	0.26	70,85,104,106	0
2	NGA	A	702	15/15	0.93	0.16	54,67,86,96	0
3	NA	A	705	1/1	0.94	0.41	72,72,72,72	0
2	NGA	B	702	15/15	0.97	0.16	39,46,54,56	0
2	NGA	A	701	15/15	0.97	0.17	38,47,55,56	0

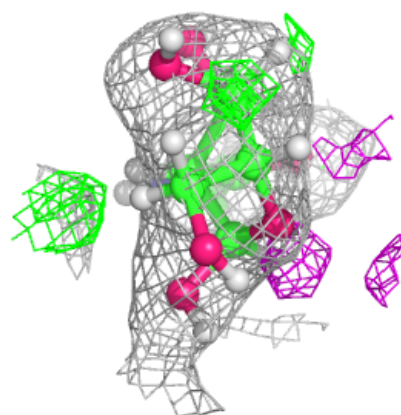
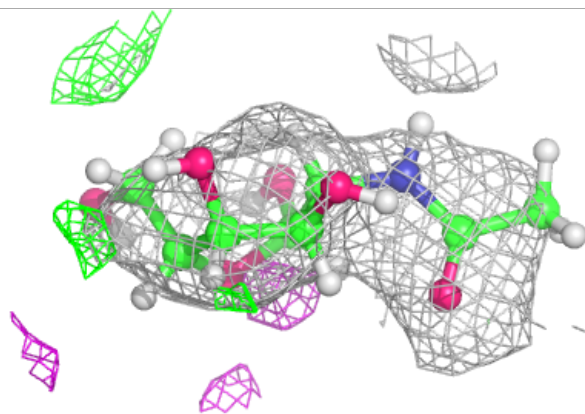
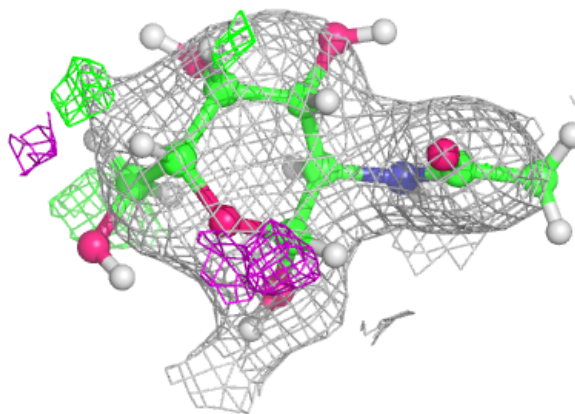
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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





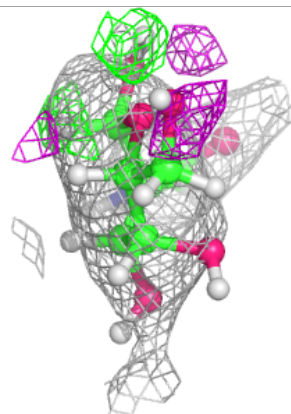
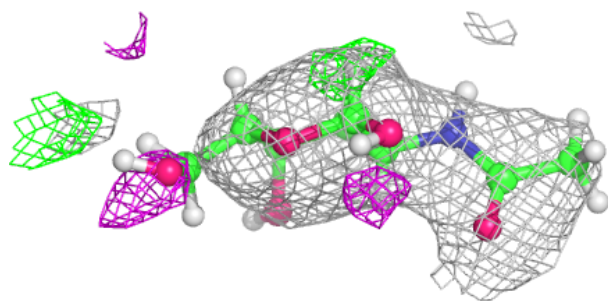
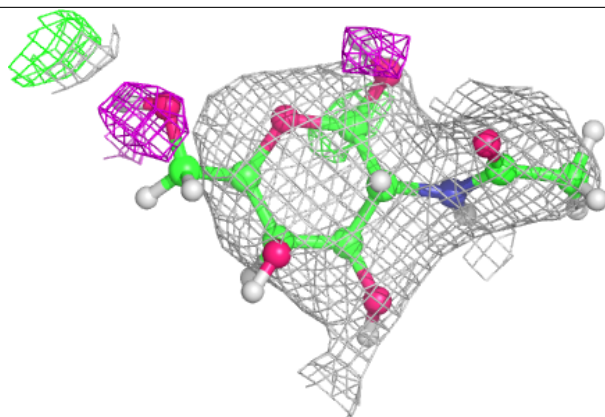
Electron density around NGA A 703:

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

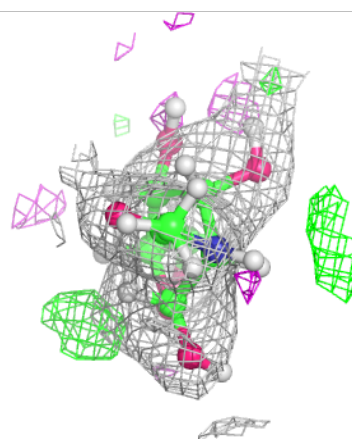
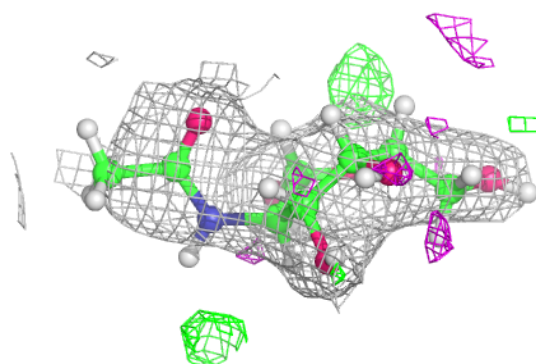
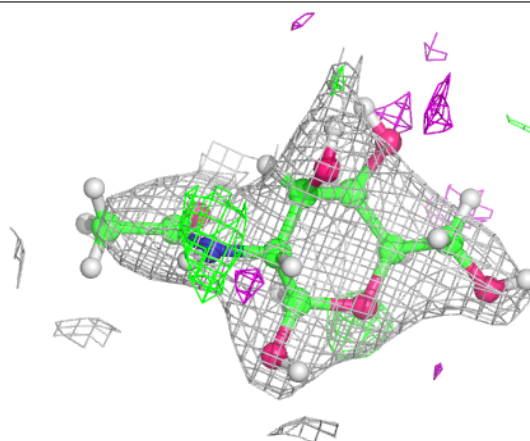


Electron density around NGA C 701:

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

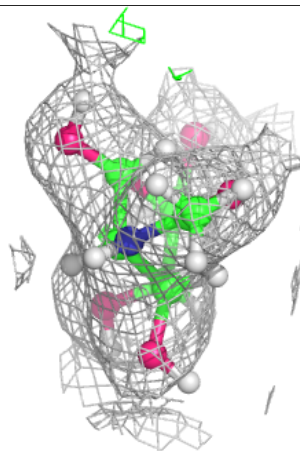
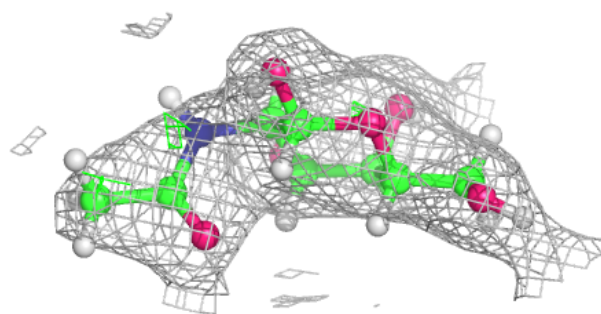
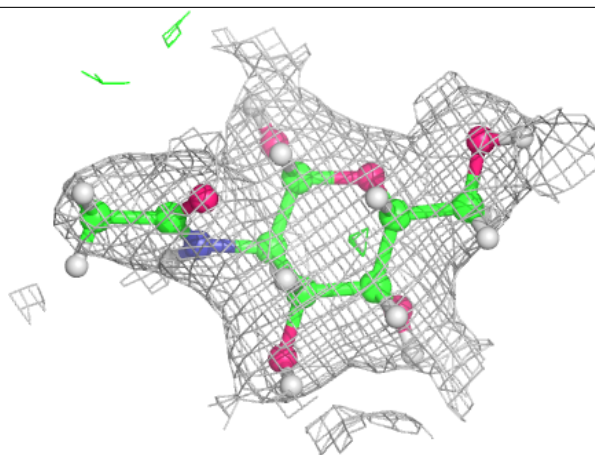
**Electron density around NGA A 702:**

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

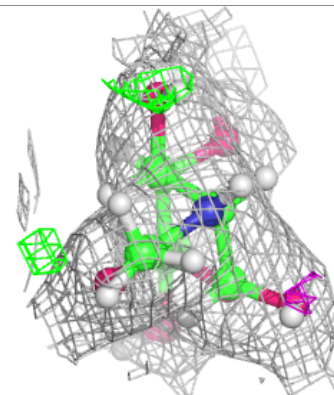
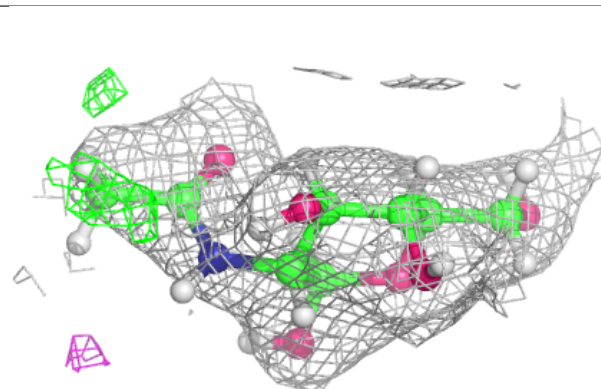
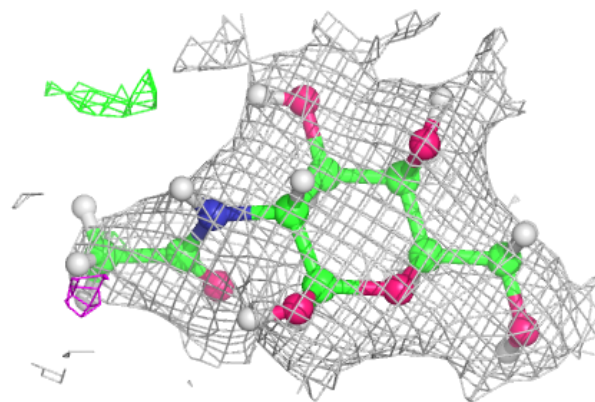


Electron density around NGA B 702:

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

**Electron density around NGA A 701:**

$2mF_o-DF_c$ (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.