



# wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 22, 2025 – 05:06 PM EDT

PDB ID : 4RPU  
Title : Crystal Structure of Human Presequence Protease in Complex with Inhibitor MitoBloCK-60  
Authors : Mo, S.M.; Liang, W.G.; King, J.V.; Wijaya, J.; Koehler, C.M.; Tang, W.J.  
Deposited on : 2014-10-31  
Resolution : 2.27 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](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 : 2022.3.0, CSD as543be (2022)  
Xtriage (Phenix) : 1.21  
EDS : 3.0  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)  
CCP4 : 9.0.004 (Gargrove)  
Density-Fitness : 1.0.11  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.41.4

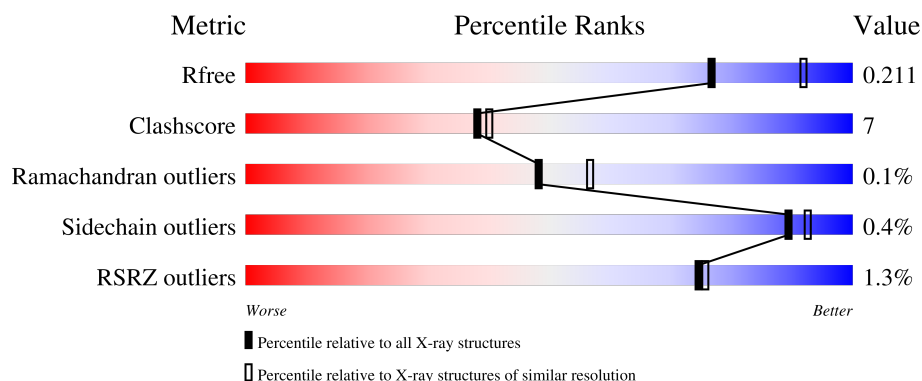
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.27 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))
$R_{free}$	164625	1763 (2.26-2.26)
Clashscore	180529	1919 (2.26-2.26)
Ramachandran outliers	177936	1884 (2.26-2.26)
Sidechain outliers	177891	1885 (2.26-2.26)
RSRZ outliers	164620	1763 (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  $\geq 3$ , 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	1014	<div> <div>%</div> <div> <div></div> <div>86%</div> <div>11%</div> <div>••</div> </div> </div>
2	B	1014	<div> <div>%</div> <div> <div></div> <div>82%</div> <div>15%</div> <div>•</div> </div> </div>

## 2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 17042 atoms, of which 100 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 Presequence protease, mitochondrial.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	As	C	N	O	S			
1	A	988	7985	1	5101	1350	1493	40	0	5	0

There are 13 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	24	MET	-	expression tag	UNP Q5JRX3
A	25	HIS	-	expression tag	UNP Q5JRX3
A	26	HIS	-	expression tag	UNP Q5JRX3
A	27	HIS	-	expression tag	UNP Q5JRX3
A	28	HIS	-	expression tag	UNP Q5JRX3
A	29	HIS	-	expression tag	UNP Q5JRX3
A	30	HIS	-	expression tag	UNP Q5JRX3
A	31	ALA	-	expression tag	UNP Q5JRX3
A	32	ALA	-	expression tag	UNP Q5JRX3
A	107	GLN	GLU	engineered mutation	UNP Q5JRX3
A	328	VAL	ILE	SEE REMARK 999	UNP Q5JRX3
A	397	VAL	ALA	SEE REMARK 999	UNP Q5JRX3
A	1037	ARG	GLN	SEE REMARK 999	UNP Q5JRX3

- Molecule 2 is a protein called Presequence protease, mitochondrial.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	As	C	N	O	S			
2	B	979	7878	1	5036	1327	1474	40	0	1	0

There are 13 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	24	MET	-	expression tag	UNP Q5JRX3
B	25	HIS	-	expression tag	UNP Q5JRX3
B	26	HIS	-	expression tag	UNP Q5JRX3

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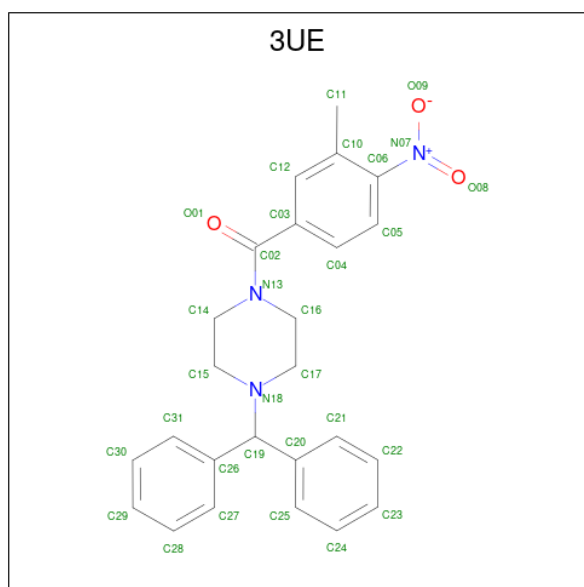
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Chain	Residue	Modelled	Actual	Comment	Reference
B	27	HIS	-	expression tag	UNP Q5JRX3
B	28	HIS	-	expression tag	UNP Q5JRX3
B	29	HIS	-	expression tag	UNP Q5JRX3
B	30	HIS	-	expression tag	UNP Q5JRX3
B	31	ALA	-	expression tag	UNP Q5JRX3
B	32	ALA	-	expression tag	UNP Q5JRX3
B	107	GLN	GLU	engineered mutation	UNP Q5JRX3
B	328	VAL	ILE	SEE REMARK 999	UNP Q5JRX3
B	397	VAL	ALA	SEE REMARK 999	UNP Q5JRX3
B	1037	ARG	GLN	SEE REMARK 999	UNP Q5JRX3

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Zn 1 1	0	0
3	B	1	Total Zn 1 1	0	0

- Molecule 4 is [4-(diphenylmethyl)piperazin-1-yl](3-methyl-4-nitrophenyl)methanone (three-letter code: 3UE) (formula: C<sub>25</sub>H<sub>25</sub>N<sub>3</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C H N O 56 25 25 3 3	0	0

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	H	N	O	0	0
			56	25	25	3	3		
4	B	1	Total	C	H	N	O	0	0
			56	25	25	3	3		
4	B	1	Total	C	H	N	O	0	0
			56	25	25	3	3		

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



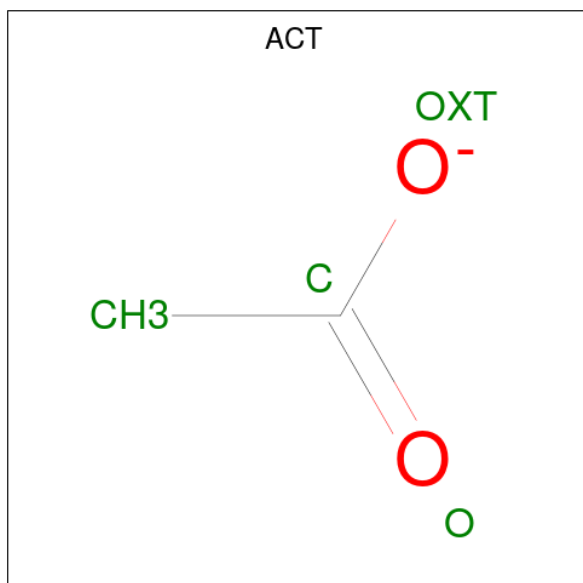
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			6	3	3		
5	A	1	Total	C	O	0	0
			6	3	3		
5	A	1	Total	C	O	0	0
			6	3	3		
5	A	1	Total	C	O	0	0
			6	3	3		
5	A	1	Total	C	O	0	0
			6	3	3		
5	A	1	Total	C	O	0	0
			6	3	3		
5	A	1	Total	C	O	0	0
			6	3	3		

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

- Molecule 6 is ACETATE ION (three-letter code: ACT) (formula:  $C_2H_3O_2$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			4	2	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			4	2	2		
6	A	1	Total	C	O	0	0
			4	2	2		
6	B	1	Total	C	O	0	0
			4	2	2		
6	B	1	Total	C	O	0	0
			4	2	2		
6	B	1	Total	C	O	0	0
			4	2	2		
6	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 7 is CALCIUM ION (three-letter code: CA) (formula: Ca).

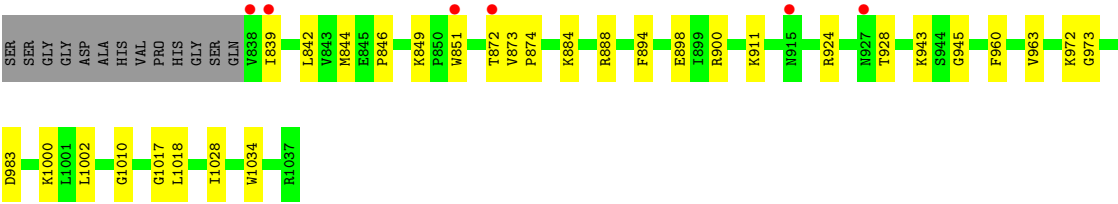
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	B	2	Total	Ca	0	0
			2	2		

- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	420	Total	O	0	0
			420	420		
8	B	391	Total	O	0	0
			391	391		







## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	245.62Å 85.49Å 158.21Å 90.00° 127.53° 90.00°	Depositor
Resolution (Å)	44.85 – 2.27 44.85 – 2.27	Depositor EDS
% Data completeness (in resolution range)	98.5 (44.85-2.27) 93.2 (44.85-2.27)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.86 (at 2.27Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, $R_{free}$	0.175 , 0.208 0.181 , 0.211	Depositor DCC
$R_{free}$ test set	1992 reflections (1.67%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	33.3	Xtriage
Anisotropy	0.393	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 34.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.015 for -h-2*k,-k,l	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	17042	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.05% 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 3UE, ACT, CA, GOL, ZN, CAS, MLZ, MLY

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	0.37	0/7776	0.54	4/10565 (0.0%)
2	B	0.35	0/7640	0.51	1/10365 (0.0%)
All	All	0.36	0/15416	0.53	5/20930 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	283	GLU	CB-CA-C	5.97	122.33	110.40
1	A	894	PHE	N-CA-C	5.81	126.69	111.00
1	A	228	ASP	CB-CA-C	-5.75	98.91	110.40
2	B	928	THR	N-CA-C	5.34	125.42	111.00
1	A	894	PHE	N-CA-CB	-5.29	101.08	110.60

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	7985	0	7877	97	0
2	B	7878	0	7775	112	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	62	50	50	4	0
4	B	62	50	50	2	0
5	A	66	0	88	7	0
5	B	42	0	56	11	0
6	A	12	0	9	0	0
6	B	20	0	15	2	0
7	B	2	0	0	0	0
8	A	420	0	0	6	0
8	B	391	0	0	7	0
All	All	16942	100	15920	217	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 7.

The worst 5 of 217 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:822:VAL:HB	1:A:838:VAL:HG13	1.40	1.00
2:B:380:TYR:HB3	5:B:1105:GOL:H31	1.57	0.87
1:A:248:GLU:OE2	1:A:251:MLY:HH12	1.74	0.86
1:A:326:THR:HG21	1:A:401:ILE:HD11	1.57	0.85
2:B:398:GLU:HG3	2:B:505:ARG:HH21	1.46	0.81

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	950/1014 (94%)	937 (99%)	12 (1%)	1 (0%)	48 57
2	B	931/1014 (92%)	919 (99%)	11 (1%)	1 (0%)	48 57
All	All	1881/2028 (93%)	1856 (99%)	23 (1%)	2 (0%)	48 57

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	147	PRO
1	A	147	PRO

### 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	848/864 (98%)	842 (99%)	6 (1%)	81	88
2	B	834/862 (97%)	832 (100%)	2 (0%)	92	95
All	All	1682/1726 (98%)	1674 (100%)	8 (0%)	89	91

5 of 8 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	818	VAL
2	B	107	GLN
1	A	777[B]	ASN
1	A	777[A]	ASN
1	A	811	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	915	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](#)

76 non-standard protein/DNA/RNA residues 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	MLZ	B	1000	2	8,9,10	1.12	0	4,9,11	0.97	0
1	MLY	A	278	1	9,10,11	0.52	0	6,11,13	0.64	0
1	MLZ	A	207	1	8,9,10	1.07	0	4,9,11	0.78	0
2	MLZ	B	884	2	8,9,10	1.12	1 (12%)	4,9,11	0.99	0
1	MLZ	A	937	1	8,9,10	1.10	0	4,9,11	0.77	0
1	MLY	A	154	1	9,10,11	0.54	0	6,11,13	0.83	0
1	MLY	A	488	1	9,10,11	0.52	0	6,11,13	0.59	0
1	MLY	A	750	1	9,10,11	0.47	0	6,11,13	0.88	0
1	MLZ	A	769	1	8,9,10	1.11	0	4,9,11	0.81	0
1	MLY	A	794	1	9,10,11	0.56	0	6,11,13	0.53	0
2	MLZ	B	494	2	8,9,10	1.09	0	4,9,11	0.77	0
1	MLZ	A	1000	1	8,9,10	1.16	1 (12%)	4,9,11	0.86	0
2	MLY	B	287	2	9,10,11	0.54	0	6,11,13	0.87	0
1	MLY	A	759	1	9,10,11	0.50	0	6,11,13	0.72	0
2	MLY	B	540	2	9,10,11	0.46	0	6,11,13	0.70	0
2	CAS	B	112	2	5,8,9	1.39	1 (20%)	1,9,11	0.55	0
2	MLZ	B	854	2	8,9,10	1.10	0	4,9,11	0.84	0
2	MLY	B	972	2	9,10,11	0.59	0	6,11,13	0.79	0
2	MLY	B	986	2	9,10,11	0.49	0	6,11,13	0.71	0
2	MLY	B	116	2	9,10,11	0.52	0	6,11,13	0.90	0
1	MLY	A	290	1	9,10,11	0.60	0	6,11,13	0.74	0
1	MLY	A	700	1	9,10,11	0.53	0	6,11,13	1.11	0
1	MLZ	A	884	1	8,9,10	1.13	0	4,9,11	0.85	0
2	MLY	B	911	2	9,10,11	0.55	0	6,11,13	0.90	0
1	MLY	A	704	1	9,10,11	0.64	0	6,11,13	0.77	0
2	MLZ	B	490	2	8,9,10	1.20	0	4,9,11	1.01	0
2	MLZ	B	125	2	8,9,10	1.06	0	4,9,11	0.63	0
1	MLZ	A	550	1	8,9,10	1.14	0	4,9,11	0.99	0
2	MLZ	B	769	2	8,9,10	1.13	0	4,9,11	0.95	0
2	MLZ	B	764	2	8,9,10	1.14	0	4,9,11	0.92	0
2	MLY	B	525	2	9,10,11	0.60	0	6,11,13	0.69	0
1	MLY	A	642	1	9,10,11	0.45	0	6,11,13	0.98	0
1	MLY	A	946	1	9,10,11	0.52	0	6,11,13	0.87	0
2	MLZ	B	759	2	8,9,10	1.05	0	4,9,11	0.86	0
2	MLY	B	290	2	9,10,11	0.60	0	6,11,13	0.77	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
1	MLY	A	116	1	9,10,11	0.45	0	6,11,13	0.99	0
2	MLZ	B	199	2	8,9,10	1.10	0	4,9,11	0.98	0
2	MLY	B	363	2	9,10,11	0.69	0	6,11,13	0.64	0
1	MLY	A	911	1	9,10,11	0.66	0	6,11,13	0.94	0
1	MLY	A	251	1	9,10,11	0.61	0	6,11,13	0.90	0
2	MLY	B	431	2	9,10,11	0.51	0	6,11,13	1.16	0
2	MLZ	B	66	2	8,9,10	0.96	0	4,9,11	0.57	0
1	MLY	A	854	1	9,10,11	0.61	0	6,11,13	0.65	0
2	MLZ	B	956	2	8,9,10	1.04	0	4,9,11	0.66	0
1	MLZ	A	66	1	8,9,10	1.05	0	4,9,11	0.81	0
1	MLY	A	287	1	9,10,11	0.45	0	6,11,13	0.63	0
2	MLZ	B	937	2	8,9,10	1.02	0	4,9,11	0.67	0
2	MLY	B	946	2	9,10,11	0.46	0	6,11,13	1.00	0
2	MLY	B	513	2	9,10,11	0.52	0	6,11,13	0.78	0
1	MLY	A	363	1	9,10,11	0.58	0	6,11,13	0.77	0
1	CAS	A	112	1	5,8,9	1.36	1 (20%)	1,9,11	1.17	0
2	MLY	B	466	2	9,10,11	0.55	0	6,11,13	0.89	0
2	MLY	B	41	2	9,10,11	0.52	0	6,11,13	0.95	0
1	MLY	A	431	1	9,10,11	0.59	0	6,11,13	1.00	0
2	MLZ	B	902	2	8,9,10	1.05	0	4,9,11	0.88	0
2	MLY	B	499	2	9,10,11	0.53	0	6,11,13	0.81	0
1	MLY	A	521	1	9,10,11	0.59	0	6,11,13	0.84	0
2	MLY	B	550	2	9,10,11	0.55	0	6,11,13	1.01	0
2	MLY	B	642	2	9,10,11	0.48	0	6,11,13	0.82	0
1	MLZ	A	972	1	8,9,10	1.07	0	4,9,11	0.87	0
2	MLZ	B	324	2	8,9,10	1.15	0	4,9,11	0.66	0
2	MLZ	B	624	2	8,9,10	1.06	0	4,9,11	0.93	0
2	MLY	B	750	2	9,10,11	0.53	0	6,11,13	0.76	0
1	MLY	A	540	1	9,10,11	0.55	0	6,11,13	0.73	0
1	MLZ	A	943	1	8,9,10	1.13	0	4,9,11	0.82	0
2	MLY	B	943	2	9,10,11	0.62	0	6,11,13	0.77	0
2	MLY	B	488	2	9,10,11	0.49	0	6,11,13	0.84	0
1	MLZ	A	437	1	8,9,10	1.07	0	4,9,11	0.74	0
2	MLZ	B	1013	2	8,9,10	1.18	0	4,9,11	0.66	0
1	MLY	A	466	1	9,10,11	0.48	0	6,11,13	0.84	0
1	MLY	A	513	1	9,10,11	0.57	0	6,11,13	0.79	0
1	MLZ	A	525	1	8,9,10	1.17	0	4,9,11	1.04	0
1	MLY	A	956	1	9,10,11	0.54	0	6,11,13	1.05	0
1	MLY	A	902	1	9,10,11	0.56	0	6,11,13	0.56	0
1	MLY	A	764	1	9,10,11	0.48	0	6,11,13	0.84	0
2	MLY	B	278	2	9,10,11	0.52	0	6,11,13	0.74	0

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	MLZ	B	1000	2	-	3/7/8/10	-
1	MLY	A	278	1	-	0/8/9/11	-
1	MLZ	A	207	1	-	2/7/8/10	-
2	MLZ	B	884	2	-	0/7/8/10	-
1	MLZ	A	937	1	-	3/7/8/10	-
1	MLY	A	154	1	-	0/8/9/11	-
1	MLY	A	488	1	-	2/8/9/11	-
1	MLY	A	750	1	-	2/8/9/11	-
1	MLZ	A	769	1	-	2/7/8/10	-
1	MLY	A	794	1	-	3/8/9/11	-
2	MLZ	B	494	2	-	0/7/8/10	-
1	MLZ	A	1000	1	-	3/7/8/10	-
2	MLY	B	287	2	-	0/8/9/11	-
1	MLY	A	759	1	-	3/8/9/11	-
2	MLY	B	540	2	-	2/8/9/11	-
2	CAS	B	112	2	-	0/0/7/9	-
2	MLZ	B	854	2	-	1/7/8/10	-
2	MLY	B	972	2	-	4/8/9/11	-
2	MLY	B	986	2	-	1/8/9/11	-
2	MLY	B	116	2	-	3/8/9/11	-
1	MLY	A	290	1	-	1/8/9/11	-
1	MLY	A	700	1	-	2/8/9/11	-
1	MLZ	A	884	1	-	1/7/8/10	-
2	MLY	B	911	2	-	0/8/9/11	-
1	MLY	A	704	1	-	0/8/9/11	-
2	MLZ	B	490	2	-	0/7/8/10	-
2	MLZ	B	125	2	-	1/7/8/10	-
1	MLZ	A	550	1	-	2/7/8/10	-
2	MLZ	B	769	2	-	2/7/8/10	-
2	MLZ	B	764	2	-	3/7/8/10	-
2	MLY	B	525	2	-	2/8/9/11	-
1	MLY	A	642	1	-	0/8/9/11	-
1	MLY	A	946	1	-	0/8/9/11	-
2	MLZ	B	759	2	-	1/7/8/10	-
2	MLY	B	290	2	-	0/8/9/11	-
1	MLY	A	116	1	-	0/8/9/11	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	MLZ	B	199	2	-	2/7/8/10	-
2	MLY	B	363	2	-	0/8/9/11	-
1	MLY	A	911	1	-	0/8/9/11	-
1	MLY	A	251	1	-	0/8/9/11	-
2	MLY	B	431	2	-	1/8/9/11	-
2	MLZ	B	66	2	-	2/7/8/10	-
1	MLY	A	854	1	-	0/8/9/11	-
2	MLZ	B	956	2	-	1/7/8/10	-
1	MLZ	A	66	1	-	3/7/8/10	-
1	MLY	A	287	1	-	0/8/9/11	-
2	MLZ	B	937	2	-	2/7/8/10	-
2	MLY	B	946	2	-	0/8/9/11	-
2	MLY	B	513	2	-	2/8/9/11	-
1	MLY	A	363	1	-	0/8/9/11	-
1	CAS	A	112	1	-	0/0/7/9	-
2	MLY	B	466	2	-	3/8/9/11	-
2	MLY	B	41	2	-	0/8/9/11	-
1	MLY	A	431	1	-	1/8/9/11	-
2	MLZ	B	902	2	-	0/7/8/10	-
2	MLY	B	499	2	-	1/8/9/11	-
1	MLY	A	521	1	-	2/8/9/11	-
2	MLY	B	550	2	-	0/8/9/11	-
2	MLY	B	642	2	-	0/8/9/11	-
1	MLZ	A	972	1	-	2/7/8/10	-
2	MLZ	B	324	2	-	4/7/8/10	-
2	MLZ	B	624	2	-	0/7/8/10	-
2	MLY	B	750	2	-	1/8/9/11	-
1	MLY	A	540	1	-	2/8/9/11	-
1	MLZ	A	943	1	-	2/7/8/10	-
2	MLY	B	943	2	-	3/8/9/11	-
2	MLY	B	488	2	-	0/8/9/11	-
1	MLZ	A	437	1	-	4/7/8/10	-
2	MLZ	B	1013	2	-	5/7/8/10	-
1	MLY	A	466	1	-	2/8/9/11	-
1	MLY	A	513	1	-	1/8/9/11	-
1	MLZ	A	525	1	-	2/7/8/10	-
1	MLY	A	956	1	-	1/8/9/11	-
1	MLY	A	902	1	-	2/8/9/11	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	MLY	A	764	1	-	0/8/9/11	-
2	MLY	B	278	2	-	2/8/9/11	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1000	MLZ	CA-N	-2.17	1.42	1.48
1	A	112	CAS	O-C	2.12	1.28	1.20
2	B	112	CAS	O-C	2.07	1.27	1.20
2	B	884	MLZ	CA-N	-2.01	1.42	1.48

There are no bond angle outliers.

There are no chirality outliers.

5 of 102 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	437	MLZ	O-C-CA-CB
1	A	437	MLZ	CD-CE-NZ-CM
1	A	769	MLZ	C-CA-CB-CG
1	A	1000	MLZ	C-CA-CB-CG
1	A	1000	MLZ	CD-CE-NZ-CM

There are no ring outliers.

34 monomers are involved in 57 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1000	MLZ	1	0
1	A	488	MLY	2	0
1	A	750	MLY	1	0
1	A	794	MLY	2	0
1	A	1000	MLZ	1	0
1	A	759	MLY	5	0
2	B	540	MLY	2	0
2	B	112	CAS	1	0
2	B	972	MLY	1	0
1	A	700	MLY	2	0
2	B	911	MLY	1	0
1	A	704	MLY	1	0
2	B	125	MLZ	1	0
2	B	764	MLZ	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	946	MLY	3	0
2	B	759	MLZ	3	0
2	B	290	MLY	2	0
2	B	199	MLZ	2	0
1	A	251	MLY	2	0
2	B	66	MLZ	1	0
1	A	854	MLY	1	0
1	A	287	MLY	2	0
2	B	513	MLY	2	0
1	A	112	CAS	1	0
2	B	466	MLY	3	0
2	B	41	MLY	1	0
2	B	499	MLY	2	0
2	B	550	MLY	1	0
2	B	324	MLZ	2	0
1	A	540	MLY	2	0
2	B	943	MLY	1	0
1	A	437	MLZ	1	0
1	A	902	MLY	2	0
2	B	278	MLY	1	0

## 5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 34 ligands modelled in this entry, 4 are monoatomic - leaving 30 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	$\# Z  > 2$
4	3UE	B	1103	-	34,34,34	4.06	14 (41%)	44,47,47	1.45	7 (15%)
5	GOL	B	1105	-	5,5,5	0.47	0	5,5,5	0.63	0
5	GOL	A	1108	-	5,5,5	0.44	0	5,5,5	0.34	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
5	GOL	A	1106	-	5,5,5	0.37	0	5,5,5	0.32	0
5	GOL	B	1104	-	5,5,5	0.42	0	5,5,5	0.38	0
6	ACT	B	1113	-	3,3,3	0.77	0	3,3,3	1.35	0
6	ACT	B	1112	-	3,3,3	0.80	0	3,3,3	1.35	0
5	GOL	A	1112	-	5,5,5	0.33	0	5,5,5	0.35	0
5	GOL	A	1114	-	5,5,5	0.42	0	5,5,5	0.27	0
5	GOL	B	1109	-	5,5,5	0.40	0	5,5,5	0.31	0
5	GOL	A	1104	-	5,5,5	0.32	0	5,5,5	0.58	0
6	ACT	B	1111	-	3,3,3	0.93	0	3,3,3	1.12	0
5	GOL	A	1105	-	5,5,5	0.38	0	5,5,5	0.35	0
5	GOL	A	1111	-	5,5,5	0.36	0	5,5,5	0.39	0
5	GOL	B	1106	-	5,5,5	0.35	0	5,5,5	0.45	0
4	3UE	A	1102	-	34,34,34	4.03	12 (35%)	44,47,47	1.19	5 (11%)
4	3UE	A	1103	-	34,34,34	4.05	15 (44%)	44,47,47	1.55	9 (20%)
5	GOL	B	1107	-	5,5,5	0.29	0	5,5,5	0.45	0
6	ACT	A	1116	-	3,3,3	0.82	0	3,3,3	1.38	0
5	GOL	A	1107	-	5,5,5	0.36	0	5,5,5	0.29	0
5	GOL	A	1110	-	5,5,5	0.36	0	5,5,5	0.33	0
5	GOL	A	1113	-	5,5,5	0.38	0	5,5,5	0.36	0
6	ACT	A	1115	-	3,3,3	0.86	0	3,3,3	1.39	0
6	ACT	B	1114	-	3,3,3	0.83	0	3,3,3	1.50	0
5	GOL	B	1108	-	5,5,5	0.39	0	5,5,5	0.44	0
6	ACT	A	1117	3	3,3,3	0.83	0	3,3,3	1.19	0
4	3UE	B	1102	-	34,34,34	3.99	14 (41%)	44,47,47	1.29	8 (18%)
5	GOL	A	1109	-	5,5,5	0.36	0	5,5,5	0.57	0
5	GOL	B	1110	-	5,5,5	0.40	0	5,5,5	0.33	0
6	ACT	B	1115	3	3,3,3	0.77	0	3,3,3	1.39	0

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
4	3UE	B	1103	-	-	0/22/34/34	0/4/4/4
5	GOL	B	1105	-	-	2/4/4/4	-
5	GOL	A	1108	-	-	2/4/4/4	-
5	GOL	A	1106	-	-	4/4/4/4	-
5	GOL	B	1104	-	-	2/4/4/4	-
5	GOL	A	1112	-	-	2/4/4/4	-
5	GOL	A	1114	-	-	4/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	GOL	B	1109	-	-	0/4/4/4	-
5	GOL	A	1104	-	-	0/4/4/4	-
5	GOL	A	1105	-	-	2/4/4/4	-
5	GOL	A	1111	-	-	3/4/4/4	-
5	GOL	B	1106	-	-	1/4/4/4	-
4	3UE	A	1102	-	-	0/22/34/34	0/4/4/4
4	3UE	A	1103	-	-	0/22/34/34	0/4/4/4
5	GOL	B	1107	-	-	0/4/4/4	-
5	GOL	A	1107	-	-	2/4/4/4	-
5	GOL	A	1110	-	-	2/4/4/4	-
5	GOL	A	1113	-	-	4/4/4/4	-
5	GOL	B	1108	-	-	2/4/4/4	-
4	3UE	B	1102	-	-	0/22/34/34	0/4/4/4
5	GOL	A	1109	-	-	2/4/4/4	-
5	GOL	B	1110	-	-	0/4/4/4	-

The worst 5 of 55 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	1103	3UE	C19-N18	-12.76	1.27	1.48
4	A	1103	3UE	C19-N18	-12.41	1.27	1.48
4	A	1102	3UE	C19-N18	-12.26	1.28	1.48
4	B	1102	3UE	C19-N18	-12.22	1.28	1.48
4	B	1103	3UE	O08-N07	10.82	1.41	1.22

The worst 5 of 29 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1103	3UE	C16-N13-C14	5.69	124.29	112.68
4	B	1103	3UE	C16-N13-C14	4.60	122.06	112.68
4	B	1103	3UE	C03-C02-N13	3.56	123.11	118.66
4	A	1103	3UE	C03-C02-N13	3.36	122.86	118.66
4	A	1102	3UE	C14-C15-N18	3.27	116.36	110.61

There are no chirality outliers.

5 of 34 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	1105	GOL	O1-C1-C2-O2
5	A	1105	GOL	O1-C1-C2-C3

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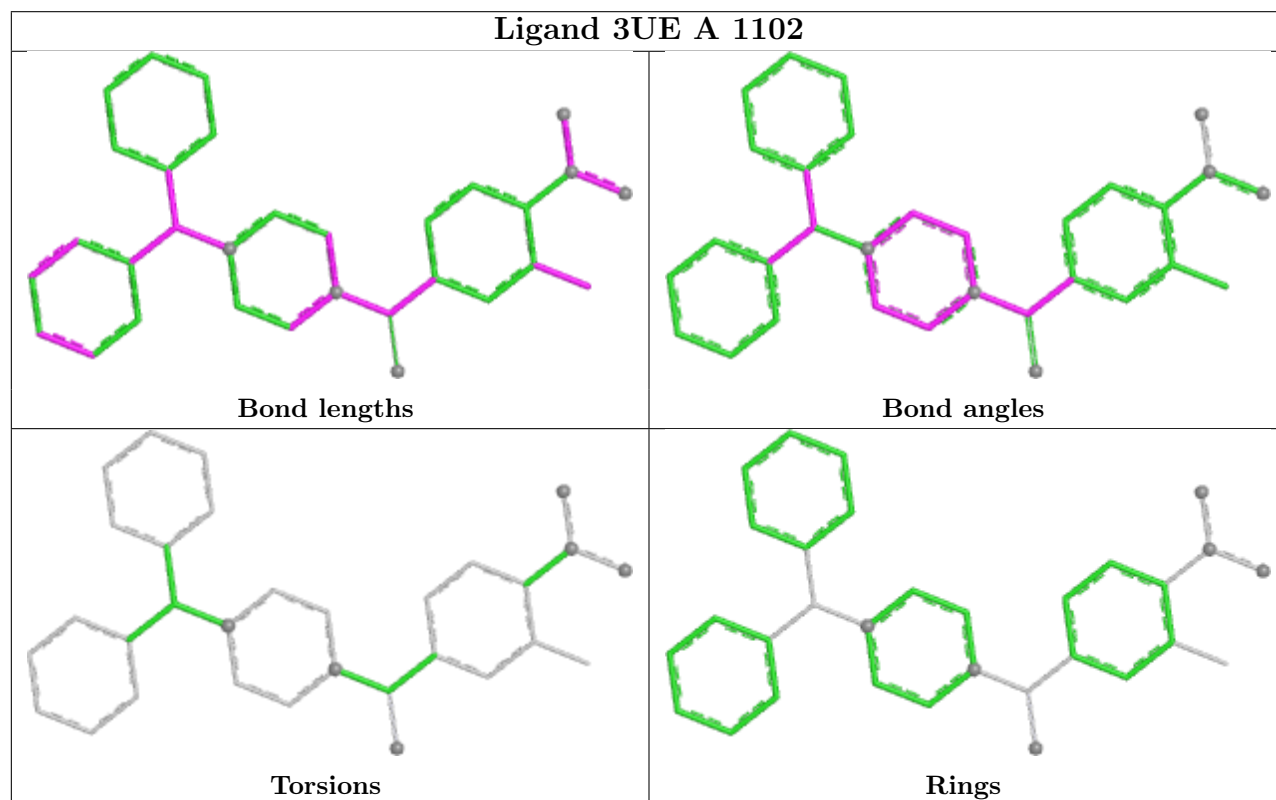
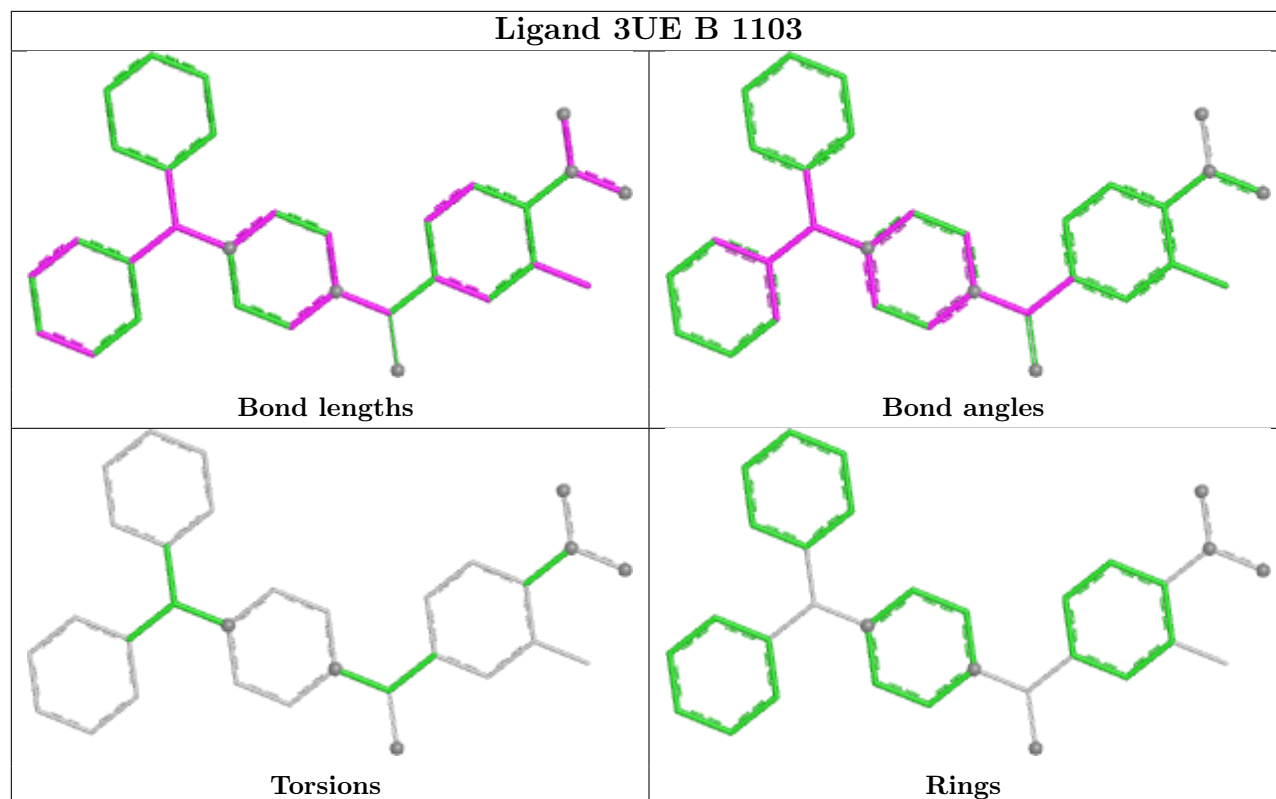
Mol	Chain	Res	Type	Atoms
5	A	1106	GOL	O1-C1-C2-C3
5	A	1106	GOL	C1-C2-C3-O3
5	A	1108	GOL	O1-C1-C2-C3

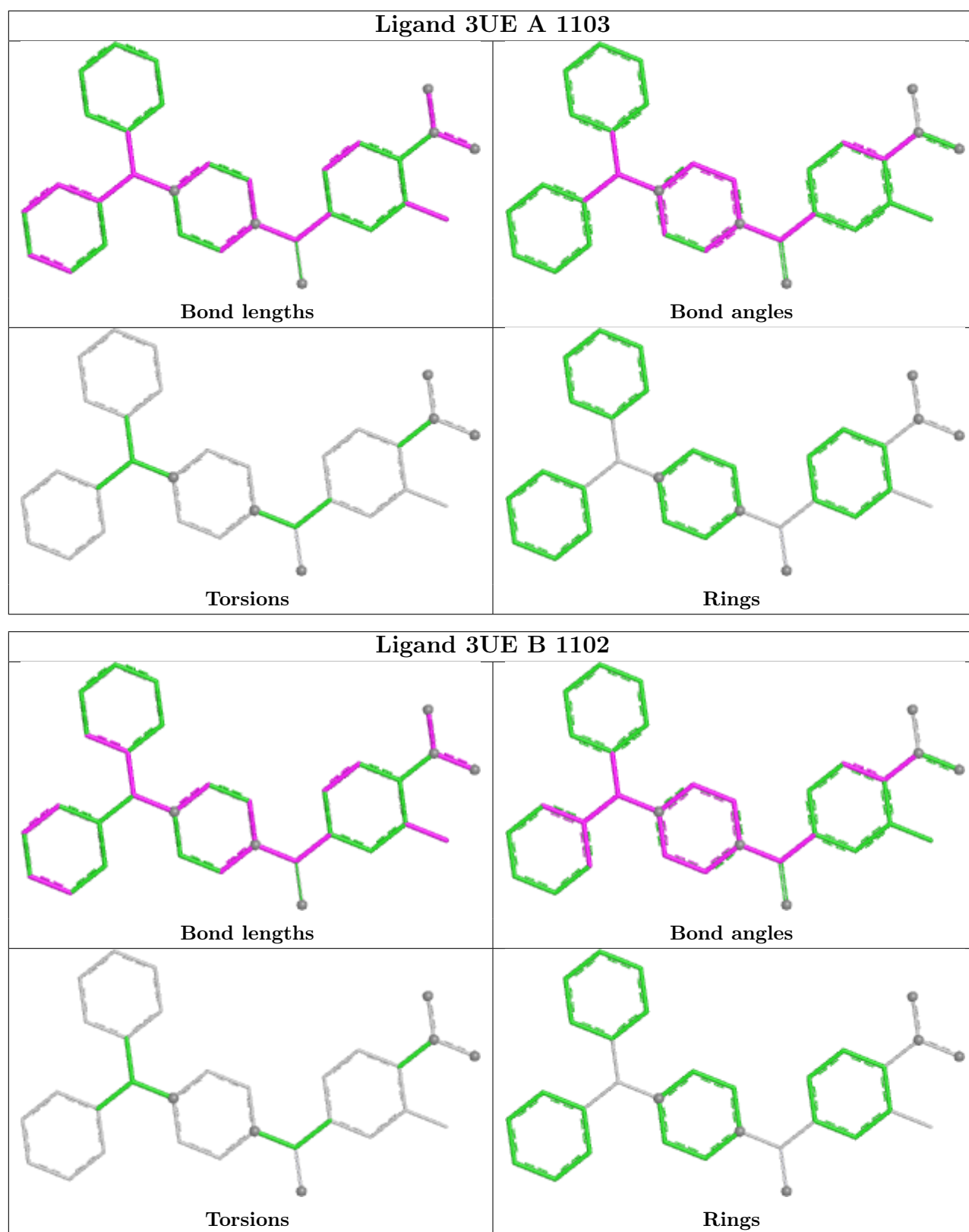
There are no ring outliers.

16 monomers are involved in 26 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	1103	3UE	1	0
5	B	1105	GOL	1	0
5	A	1106	GOL	2	0
5	B	1104	GOL	1	0
6	B	1113	ACT	1	0
5	A	1112	GOL	1	0
5	A	1114	GOL	1	0
5	B	1109	GOL	3	0
5	A	1111	GOL	2	0
5	B	1106	GOL	3	0
4	A	1102	3UE	2	0
4	A	1103	3UE	2	0
5	B	1107	GOL	3	0
5	A	1110	GOL	1	0
6	B	1114	ACT	1	0
4	B	1102	3UE	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 ⓘ

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	951/1014 (93%)	-0.15	10 (1%) 77 79	15, 32, 51, 82	5 (0%)
2	B	940/1014 (92%)	-0.08	15 (1%) 70 71	16, 34, 54, 94	1 (0%)
All	All	1891/2028 (93%)	-0.11	25 (1%) 74 76	15, 33, 53, 94	6 (0%)

The worst 5 of 25 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	915	ASN	4.3
2	B	851	TRP	4.1
1	A	927	ASN	3.8
2	B	318	PHE	3.7
1	A	838	VAL	3.7

### 6.2 Non-standard residues in protein, DNA, RNA chains [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, 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
2	MLZ	B	324	10/11	0.77	0.23	46,61,69,70	0
1	MLY	A	521	11/12	0.81	0.17	42,46,60,63	0
1	MLY	A	794	11/12	0.84	0.17	37,51,66,67	0
1	MLY	A	700	11/12	0.85	0.17	31,32,61,61	0
2	MLY	B	499	11/12	0.85	0.18	33,41,65,65	0
2	MLZ	B	1013	10/11	0.85	0.14	40,45,54,58	0
1	MLY	A	750	11/12	0.86	0.16	31,39,54,57	0
2	MLY	B	513	11/12	0.88	0.12	34,39,47,49	0
2	MLY	B	972	11/12	0.88	0.16	32,34,56,59	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	MLZ	B	494	10/11	0.88	0.12	33,40,54,57	0
1	MLY	A	704	11/12	0.89	0.13	30,31,51,53	0
2	MLY	B	750	11/12	0.90	0.12	35,41,47,53	0
2	MLY	B	41	11/12	0.90	0.12	34,35,46,49	0
2	MLY	B	986	11/12	0.90	0.14	32,42,54,56	0
2	MLY	B	290	11/12	0.90	0.13	33,37,53,54	0
1	MLY	A	363	11/12	0.91	0.12	30,32,52,56	0
2	MLZ	B	769	10/11	0.91	0.10	38,39,42,46	0
2	MLZ	B	902	10/11	0.91	0.11	28,33,39,42	0
2	MLY	B	116	11/12	0.91	0.13	28,39,47,54	0
2	MLY	B	287	11/12	0.91	0.11	31,32,41,41	0
1	MLY	A	290	11/12	0.91	0.11	34,37,48,50	0
2	MLZ	B	66	10/11	0.92	0.11	31,31,47,50	0
1	MLY	A	251	11/12	0.92	0.10	31,31,45,46	0
2	MLZ	B	125	10/11	0.92	0.10	27,27,39,41	0
2	MLY	B	550	11/12	0.92	0.11	27,32,38,44	0
2	MLY	B	278	11/12	0.92	0.12	31,35,50,53	0
2	MLZ	B	759	10/11	0.92	0.10	39,45,47,50	0
1	MLY	A	431	11/12	0.92	0.12	28,31,52,53	0
1	MLY	A	911	11/12	0.92	0.11	27,27,38,46	0
2	MLY	B	911	11/12	0.92	0.11	28,29,41,45	0
1	MLY	A	466	11/12	0.92	0.11	29,33,44,46	0
2	MLY	B	363	11/12	0.92	0.12	33,34,51,55	0
2	MLZ	B	490	10/11	0.92	0.13	32,38,59,60	0
1	MLY	A	956	11/12	0.93	0.12	27,28,53,55	0
2	MLZ	B	199	10/11	0.93	0.11	29,38,49,50	0
1	MLZ	A	972	10/11	0.93	0.11	29,33,49,51	0
2	MLZ	B	884	10/11	0.93	0.10	28,32,43,45	0
1	MLZ	A	525	10/11	0.93	0.09	37,38,44,45	0
1	MLY	A	759	11/12	0.93	0.10	35,35,41,44	0
2	MLZ	B	937	10/11	0.93	0.10	28,33,42,47	0
2	MLY	B	943	11/12	0.93	0.10	28,31,48,52	0
2	MLZ	B	956	10/11	0.93	0.09	28,31,42,42	0
2	MLY	B	525	11/12	0.93	0.09	32,34,42,43	0
2	MLY	B	540	11/12	0.93	0.09	28,29,34,38	0
1	MLY	A	946	11/12	0.93	0.10	26,34,48,48	0
1	MLZ	A	66	10/11	0.94	0.09	31,32,42,45	0
2	MLZ	B	624	10/11	0.94	0.08	33,34,39,42	0
1	MLY	A	540	11/12	0.94	0.08	30,31,33,35	0
1	MLY	A	642	11/12	0.94	0.09	29,31,35,40	0
2	MLZ	B	764	10/11	0.94	0.08	39,43,45,47	0
1	MLZ	A	437	10/11	0.94	0.09	28,29,47,48	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	MLZ	B	854	10/11	0.94	0.08	32,33,36,40	0
2	MLZ	B	1000	10/11	0.94	0.11	31,36,54,56	0
1	MLY	A	902	11/12	0.94	0.10	28,29,37,39	0
1	MLY	A	116	11/12	0.95	0.08	29,29,30,37	0
1	MLY	A	764	11/12	0.95	0.08	34,34,40,46	0
2	MLY	B	431	11/12	0.95	0.10	31,32,48,56	0
2	MLY	B	466	11/12	0.95	0.08	30,31,34,40	0
1	MLZ	A	1000	10/11	0.95	0.08	27,31,44,44	0
1	MLZ	A	769	10/11	0.95	0.08	33,34,39,43	0
1	MLY	A	488	11/12	0.95	0.07	30,30,32,35	0
1	MLY	A	854	11/12	0.95	0.08	31,31,38,40	0
1	MLZ	A	884	10/11	0.95	0.08	27,27,29,31	0
2	MLY	B	946	11/12	0.95	0.10	27,35,47,47	0
1	MLY	A	513	11/12	0.95	0.07	33,36,44,51	0
1	MLY	A	278	11/12	0.95	0.07	31,32,35,36	0
1	MLZ	A	943	10/11	0.95	0.09	27,35,38,42	0
2	MLY	B	642	11/12	0.95	0.09	33,34,39,41	0
1	MLZ	A	207	10/11	0.95	0.10	29,29,46,53	0
1	MLY	A	154	11/12	0.96	0.09	28,30,43,43	0
1	MLY	A	287	11/12	0.96	0.07	31,32,34,38	0
1	MLZ	A	937	10/11	0.96	0.07	28,29,43,44	0
2	CAS	B	112	9/10	0.97	0.10	27,28,60,92	0
1	CAS	A	112	9/10	0.97	0.10	28,32,68,93	0
2	MLY	B	488	11/12	0.97	0.07	31,32,37,40	0
1	MLZ	A	550	10/11	0.97	0.06	27,30,37,41	0

## 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, 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( $\text{\AA}^2$ )	Q<0.9
6	ACT	A	1115	4/4	0.67	0.21	56,63,63,67	0
5	GOL	B	1108	6/6	0.77	0.24	69,70,74,75	0
6	ACT	B	1112	4/4	0.77	0.17	43,47,52,61	0

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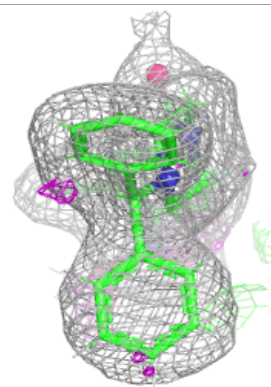
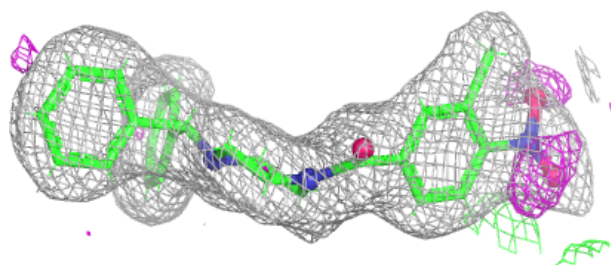
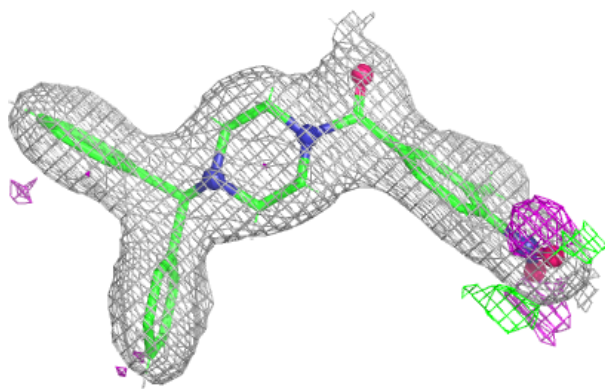
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
5	GOL	B	1107	6/6	0.80	0.20	53,60,68,75	0
5	GOL	A	1109	6/6	0.80	0.23	54,62,73,77	0
6	ACT	B	1114	4/4	0.80	0.17	41,51,52,59	0
5	GOL	A	1110	6/6	0.82	0.15	51,66,68,69	0
5	GOL	A	1106	6/6	0.82	0.20	58,61,65,65	0
5	GOL	A	1105	6/6	0.82	0.18	50,51,65,72	0
5	GOL	A	1108	6/6	0.84	0.17	51,62,70,72	0
6	ACT	B	1111	4/4	0.84	0.14	41,44,52,53	0
5	GOL	A	1107	6/6	0.85	0.14	48,59,68,72	0
6	ACT	B	1113	4/4	0.85	0.14	43,47,49,49	0
5	GOL	A	1112	6/6	0.85	0.18	32,46,57,63	0
5	GOL	A	1111	6/6	0.86	0.18	35,50,60,67	0
5	GOL	B	1105	6/6	0.87	0.14	39,41,49,51	0
5	GOL	A	1114	6/6	0.87	0.14	54,60,72,80	0
5	GOL	B	1106	6/6	0.88	0.16	51,57,58,61	0
5	GOL	A	1113	6/6	0.88	0.14	54,58,61,62	0
5	GOL	A	1104	6/6	0.90	0.16	38,41,46,48	0
5	GOL	B	1109	6/6	0.91	0.12	28,43,48,52	0
5	GOL	B	1104	6/6	0.91	0.12	41,48,52,54	0
6	ACT	B	1115	4/4	0.92	0.10	28,42,43,45	0
6	ACT	A	1116	4/4	0.93	0.10	33,34,39,39	0
5	GOL	B	1110	6/6	0.93	0.10	35,41,46,54	0
4	3UE	B	1102	31/31	0.93	0.08	29,31,36,49	0
6	ACT	A	1117	4/4	0.94	0.11	28,37,44,46	0
4	3UE	A	1103	31/31	0.94	0.07	28,32,39,58	0
4	3UE	B	1103	31/31	0.94	0.07	29,35,40,56	0
4	3UE	A	1102	31/31	0.95	0.07	27,33,45,64	0
7	CA	B	1116	1/1	0.97	0.08	41,41,41,41	0
7	CA	B	1117	1/1	0.98	0.14	54,54,54,54	0
3	ZN	A	1101	1/1	1.00	0.02	38,38,38,38	0
3	ZN	B	1101	1/1	1.00	0.03	42,42,42,42	0

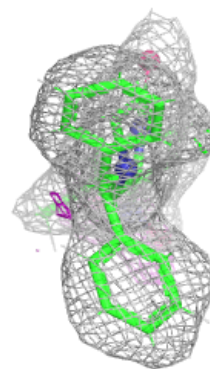
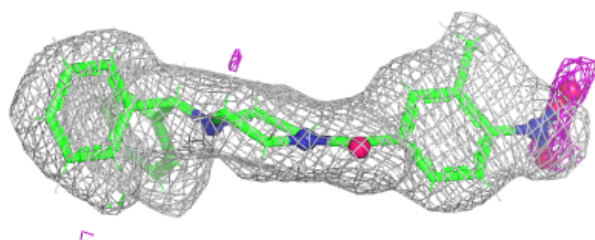
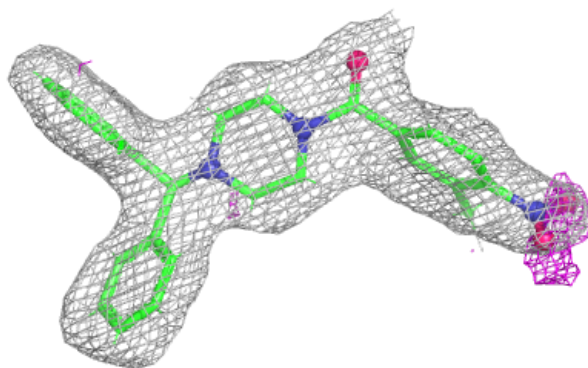
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

**Electron density around 3UE B 1102:**

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

**Electron density around 3UE A 1103:**

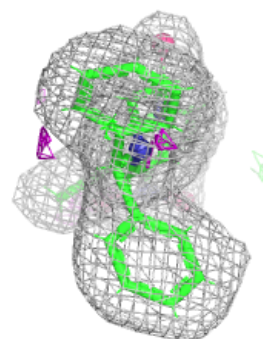
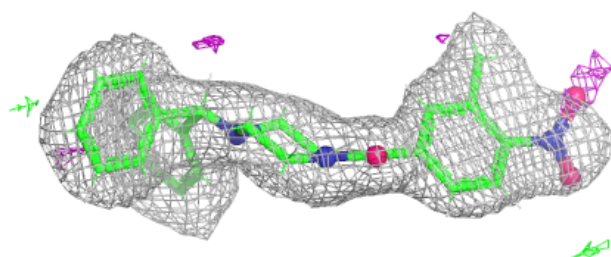
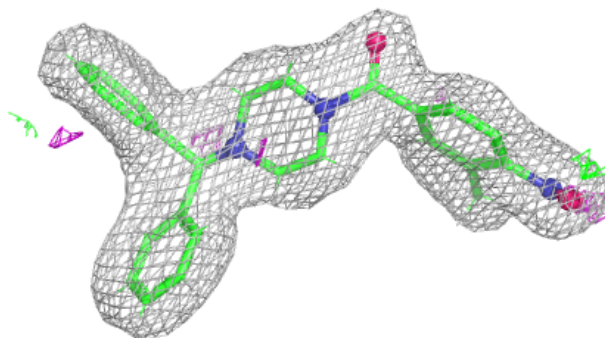
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



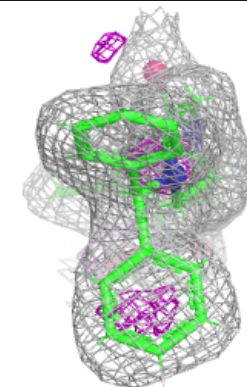
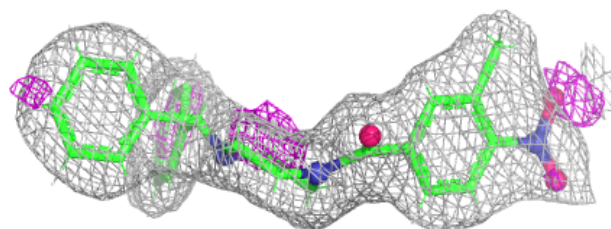
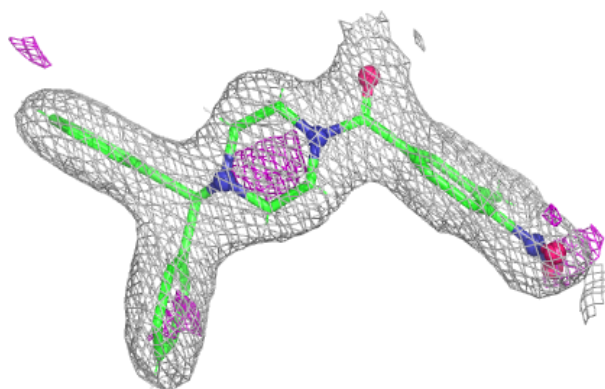


**Electron density around 3UE B 1103:**

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

**Electron density around 3UE A 1102:**

$2mF_o-DF_c$  (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.