



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 7, 2020 – 08:59 PM BST

PDB ID : 5BRP  
Title : Crystal structure of Bacillus licheniformis trehalose-6-phosphate hydrolase (TreA), mutant R201Q, in complex with PNG  
Authors : Hsiao, C.-D.; Lin, M.-G.  
Deposited on : 2015-06-01  
Resolution : 2.05 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.

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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 : 2.13.1  
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) : 2.13.1

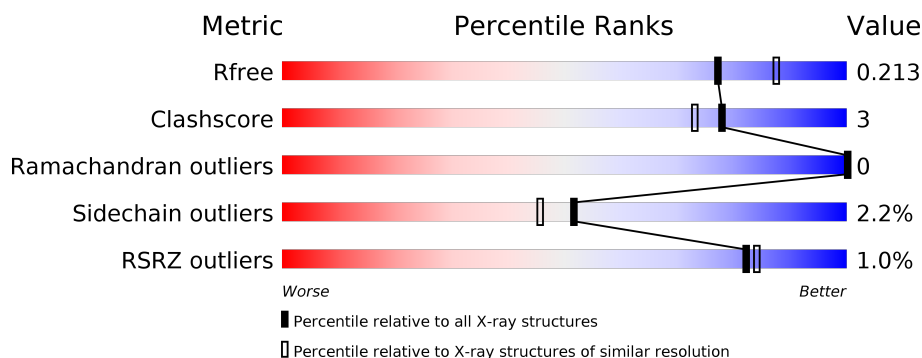
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*





The reported resolution of this entry is 2.05 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	1692 (2.04-2.04)
Clashscore	141614	1773 (2.04-2.04)
Ramachandran outliers	138981	1752 (2.04-2.04)
Sidechain outliers	138945	1752 (2.04-2.04)
RSRZ outliers	127900	1672 (2.04-2.04)

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 on 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	568	 88% 9% ••
1	B	568	 2% 92% 6% •
1	C	568	 2% 85% 13% •
1	D	568	 90% 7% •

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 20074 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 Glycoside Hydrolase Family 13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	555	Total	C	N	O	S	0	0	0
			4586	2932	755	882	17			
1	B	555	Total	C	N	O	S	0	0	0
			4586	2932	755	882	17			
1	C	555	Total	C	N	O	S	0	0	0
			4586	2932	755	882	17			
1	D	555	Total	C	N	O	S	0	0	0
			4586	2932	755	882	17			

There are 28 discrepancies between the modelled and reference sequences:

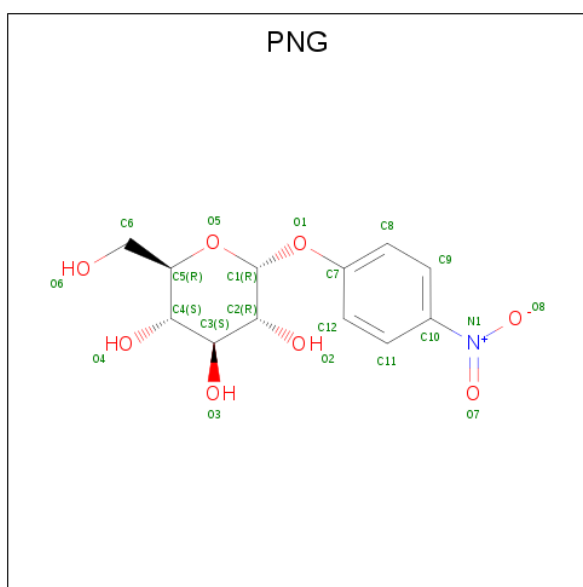
Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	HIS	-	expression tag	UNP Q65MI2
A	-4	HIS	-	expression tag	UNP Q65MI2
A	-3	HIS	-	expression tag	UNP Q65MI2
A	-2	HIS	-	expression tag	UNP Q65MI2
A	-1	HIS	-	expression tag	UNP Q65MI2
A	0	HIS	-	expression tag	UNP Q65MI2
A	201	GLN	ARG	engineered mutation	UNP Q65MI2
B	-5	HIS	-	expression tag	UNP Q65MI2
B	-4	HIS	-	expression tag	UNP Q65MI2
B	-3	HIS	-	expression tag	UNP Q65MI2
B	-2	HIS	-	expression tag	UNP Q65MI2
B	-1	HIS	-	expression tag	UNP Q65MI2
B	0	HIS	-	expression tag	UNP Q65MI2
B	201	GLN	ARG	engineered mutation	UNP Q65MI2
C	-5	HIS	-	expression tag	UNP Q65MI2
C	-4	HIS	-	expression tag	UNP Q65MI2
C	-3	HIS	-	expression tag	UNP Q65MI2
C	-2	HIS	-	expression tag	UNP Q65MI2
C	-1	HIS	-	expression tag	UNP Q65MI2
C	0	HIS	-	expression tag	UNP Q65MI2
C	201	GLN	ARG	engineered mutation	UNP Q65MI2

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-5	HIS	-	expression tag	UNP Q65MI2
D	-4	HIS	-	expression tag	UNP Q65MI2
D	-3	HIS	-	expression tag	UNP Q65MI2
D	-2	HIS	-	expression tag	UNP Q65MI2
D	-1	HIS	-	expression tag	UNP Q65MI2
D	0	HIS	-	expression tag	UNP Q65MI2
D	201	GLN	ARG	engineered mutation	UNP Q65MI2

- Molecule 2 is 4-nitrophenyl alpha-D-glucopyranoside (three-letter code: PNG) (formula:  $C_{12}H_{15}NO_8$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			21	12	1	8		
2	B	1	Total	C	N	O	0	0
			21	12	1	8		
2	C	1	Total	C	N	O	0	0
			21	12	1	8		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	1	Total 1	Mg 1	0	0
3	C	1	Total 1	Mg 1	0	0

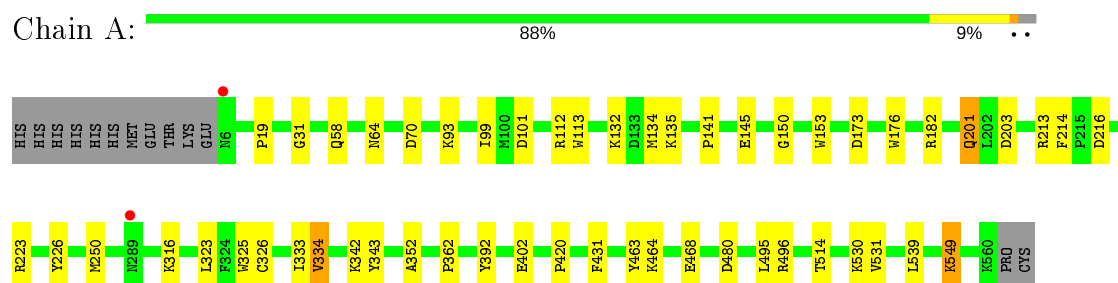
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	505	Total 505	O 505	0	0
4	B	398	Total 398	O 398	0	0
4	C	338	Total 338	O 338	0	0
4	D	422	Total 422	O 422	0	0

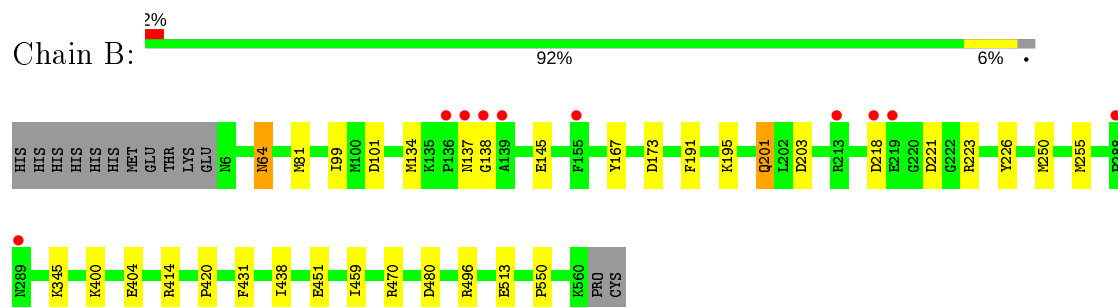
### 3 Residue-property plots [i](#)

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.

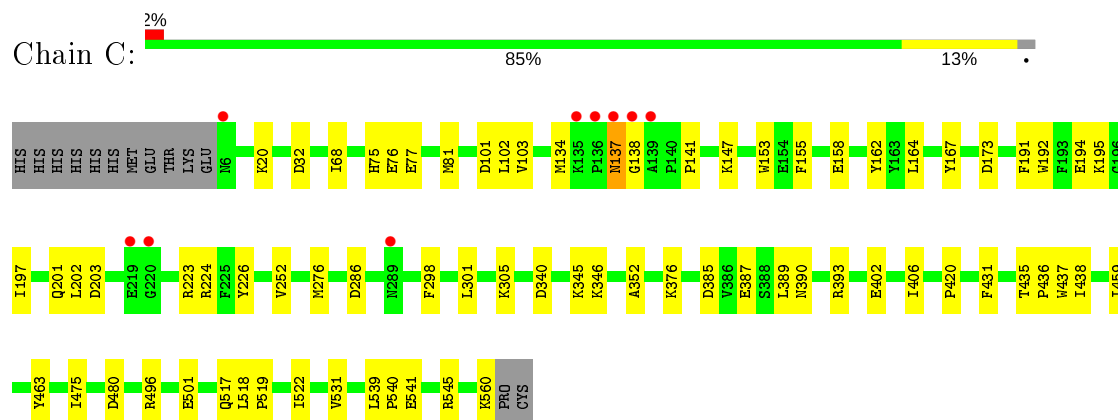
- Molecule 1: Glycoside Hydrolase Family 13



- Molecule 1: Glycoside Hydrolase Family 13

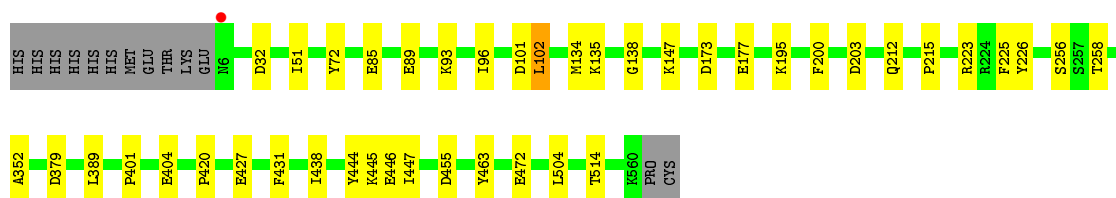
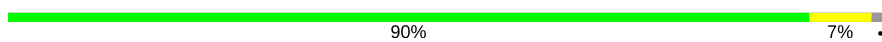


- Molecule 1: Glycoside Hydrolase Family 13



- Molecule 1: Glycoside Hydrolase Family 13

Chain D:



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	60.25Å 97.39Å 108.54Å 80.30° 88.15° 72.18°	Depositor
Resolution (Å)	30.00 – 2.05 19.55 – 2.05	Depositor EDS
% Data completeness (in resolution range)	97.5 (30.00-2.05) 49.2 (19.55-2.05)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.65 (at 2.06Å)	Xtriage
Refinement program	PHENIX 1.8.1 _1168	Depositor
R, $R_{free}$	0.171 , 0.214 0.170 , 0.213	Depositor DCC
$R_{free}$ test set	1069 reflections (1.43%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	22.3	Xtriage
Anisotropy	0.079	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 48.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.410 for h,h-k,-l	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	20074	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.51% 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: MG, PNG

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	0/4712	0.57	1/6380 (0.0%)
1	B	0.40	0/4712	0.53	0/6380
1	C	0.38	0/4712	0.53	0/6380
1	D	0.39	0/4712	0.52	0/6380
All	All	0.40	0/18848	0.54	1/25520 (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	D	0	1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	334	VAL	CB-CA-C	-6.05	99.91	111.40

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	258	THR	Peptide

## 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	4586	0	4352	31	0
1	B	4586	0	4352	20	0
1	C	4586	0	4352	35	0
1	D	4586	0	4352	19	0
2	A	21	0	15	3	0
2	B	21	0	15	3	0
2	C	21	0	15	1	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	505	0	0	1	0
4	B	398	0	0	1	0
4	C	338	0	0	1	0
4	D	422	0	0	0	0
All	All	20074	0	17453	103	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 3.

All (103) 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:C:203:ASP:OD2	2:C:601:PNG:H62	1.82	0.80
1:A:203:ASP:OD2	2:A:601:PNG:H62	1.83	0.79
1:A:326:CYS:SG	1:A:333:ILE:HG13	2.24	0.78
1:B:145:GLU:HG2	1:B:221:ASP:HB2	1.71	0.72
1:C:531:VAL:HG21	1:C:539:LEU:HD13	1.78	0.64
1:D:85:GLU:OE1	1:D:195:LYS:NZ	2.22	0.62
1:C:252:VAL:HG13	1:C:276:MET:HG3	1.80	0.61
1:A:99:ILE:HD13	1:A:250:MET:HE2	1.83	0.61
1:A:112:ARG:HH12	1:A:113:TRP:HE1	1.47	0.60
1:C:137:ASN:N	1:C:137:ASN:OD1	2.32	0.59
1:D:177:GLU:HG3	1:D:212:GLN:CD	2.23	0.58
1:B:203:ASP:OD2	2:B:601:PNG:H61	2.04	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:464:LYS:HE3	1:D:472:GLU:OE2	2.05	0.57
1:A:464:LYS:NZ	1:A:468:GLU:OE2	2.38	0.56
1:C:518:LEU:HD12	1:C:519:PRO:HD2	1.86	0.56
1:C:194:GLU:HG2	1:D:72:TYR:CZ	2.40	0.56
1:A:342:LYS:HE3	1:A:343:TYR:CZ	2.41	0.56
1:B:451:GLU:OE2	4:B:701:HOH:O	2.18	0.54
1:C:81:MET:SD	1:C:195:LYS:HE3	2.48	0.54
1:A:135:LYS:NZ	1:A:216:ASP:OD2	2.41	0.53
1:C:68:ILE:HD12	1:C:103:VAL:HB	1.91	0.53
1:D:215:PRO:HG2	1:D:225:PHE:CE1	2.43	0.52
1:C:223:ARG:HA	1:C:226:TYR:CZ	2.45	0.52
1:C:519:PRO:HG2	1:C:522:ILE:HG12	1.91	0.51
1:A:392:TYR:OH	1:A:402:GLU:OE1	2.26	0.51
1:A:549:LYS:N	1:A:549:LYS:HD2	2.27	0.50
1:D:420:PRO:HB3	1:D:431:PHE:CG	2.47	0.50
1:B:201:GLN:HG3	1:B:201:GLN:O	2.12	0.50
1:C:134:MET:SD	1:C:138:GLY:HA2	2.52	0.49
1:C:141:PRO:HG2	1:C:153:TRP:CD2	2.48	0.49
1:A:141:PRO:HG2	1:A:153:TRP:CD2	2.47	0.49
1:C:20:LYS:HE2	1:C:437:TRP:CE2	2.47	0.49
1:B:191:PHE:CZ	1:B:195:LYS:HE2	2.48	0.49
1:C:102:LEU:HD23	1:C:202:LEU:HD22	1.94	0.48
1:A:201:GLN:O	1:A:201:GLN:HG3	2.14	0.48
1:C:385:ASP:OD2	1:C:387:GLU:HB3	2.12	0.47
1:A:112:ARG:NH1	1:A:113:TRP:HE1	2.13	0.47
1:A:112:ARG:NH1	4:A:730:HOH:O	2.47	0.47
1:B:420:PRO:HB3	1:B:431:PHE:CG	2.50	0.47
1:B:513:GLU:HG3	1:B:550:PRO:HD3	1.97	0.47
1:C:352:ALA:HB2	1:C:463:TYR:CZ	2.50	0.47
1:B:400:LYS:HE3	1:B:404:GLU:HB3	1.97	0.47
1:D:102:LEU:HB2	1:D:200:PHE:CD1	2.49	0.47
1:C:20:LYS:HE2	1:C:437:TRP:CD2	2.49	0.47
1:A:112:ARG:NH1	1:A:113:TRP:NE1	2.63	0.47
1:C:192:TRP:HB3	1:C:197:ILE:HD13	1.96	0.47
1:B:420:PRO:HD3	1:B:438:ILE:HG23	1.97	0.46
1:A:64:ASN:HB2	2:A:601:PNG:O8	2.15	0.46
1:D:134:MET:SD	1:D:138:GLY:HA2	2.56	0.46
1:C:402:GLU:O	1:C:406:ILE:HG12	2.15	0.46
1:C:340:ASP:OD2	1:C:376:LYS:NZ	2.42	0.46
1:D:352:ALA:HB2	1:D:463:TYR:CZ	2.51	0.46
1:D:401:PRO:HG2	1:D:404:GLU:HG3	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:167:TYR:CE1	2:B:601:PNG:H1	2.52	0.45
1:C:420:PRO:HB3	1:C:431:PHE:CG	2.51	0.45
1:C:147:LYS:HD2	1:C:167:TYR:CE2	2.51	0.45
1:D:446:GLU:CG	1:D:447:ILE:HG13	2.47	0.45
1:B:345:LYS:HA	1:B:459:ILE:HD11	1.98	0.45
1:C:420:PRO:HD3	1:C:438:ILE:HG23	1.98	0.45
1:D:444:TYR:CZ	1:D:445:LYS:HD3	2.51	0.45
1:C:390:ASN:OD1	1:C:393:ARG:NH2	2.48	0.45
1:A:325:TRP:HB3	1:A:333:ILE:HG21	1.99	0.44
1:B:480:ASP:O	1:B:496:ARG:HA	2.16	0.44
1:B:64:ASN:HB2	2:B:601:PNG:O8	2.17	0.44
1:A:145:GLU:HG3	1:A:150:GLY:O	2.17	0.44
1:A:93:LYS:HE3	1:A:93:LYS:HB2	1.86	0.44
1:C:436:PRO:HB2	1:C:438:ILE:O	2.18	0.44
1:A:480:ASP:O	1:A:496:ARG:HA	2.17	0.44
1:B:134:MET:SD	1:B:138:GLY:HA2	2.58	0.44
1:B:414:ARG:NH1	1:B:414:ARG:HG3	2.33	0.44
1:C:75:HIS:CD2	1:C:77:GLU:HB2	2.53	0.44
1:D:446:GLU:HG2	1:D:447:ILE:HG13	2.00	0.44
1:B:218:ASP:N	1:B:218:ASP:OD1	2.51	0.43
1:A:420:PRO:HB3	1:A:431:PHE:CG	2.52	0.43
1:C:191:PHE:CZ	1:C:195:LYS:HE2	2.54	0.43
1:C:480:ASP:O	1:C:496:ARG:HA	2.18	0.43
1:C:517:GLN:OE1	1:C:545:ARG:NH1	2.52	0.43
1:A:19:PRO:O	1:A:31:GLY:HA3	2.18	0.43
1:D:223:ARG:HA	1:D:226:TYR:CZ	2.54	0.43
1:A:132:LYS:HE2	1:A:214:PHE:CD2	2.53	0.43
1:A:530:LYS:NZ	1:D:455:ASP:OD1	2.49	0.42
1:A:531:VAL:HG21	1:A:539:LEU:HD13	2.00	0.42
1:D:134:MET:HG2	1:D:135:LYS:O	2.19	0.42
1:A:134:MET:HG2	1:A:135:LYS:O	2.20	0.42
1:A:223:ARG:HA	1:A:226:TYR:CZ	2.55	0.42
1:B:64:ASN:N	1:B:64:ASN:OD1	2.52	0.42
1:C:475:ILE:HD12	1:C:501:GLU:HG2	2.01	0.42
1:C:286:ASP:HB3	1:C:298:PHE:HB3	2.02	0.41
1:A:176:TRP:CE3	1:A:182:ARG:HG2	2.55	0.41
1:D:51:ILE:HD12	1:D:96:ILE:HG21	2.02	0.41
1:B:223:ARG:HA	1:B:226:TYR:CZ	2.56	0.41
1:D:420:PRO:HD3	1:D:438:ILE:HG23	2.02	0.41
1:C:539:LEU:HA	1:C:540:PRO:HD3	1.95	0.41
1:C:345:LYS:HA	1:C:459:ILE:HD11	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:352:ALA:HB2	1:A:463:TYR:CZ	2.56	0.41
1:A:58:GLN:HB3	1:A:70:ASP:HB3	2.02	0.41
2:A:601:PNG:C12	2:A:601:PNG:H5	2.51	0.40
1:B:99:ILE:HD13	1:B:250:MET:HE2	2.03	0.40
1:C:155:PHE:HB2	1:C:162:TYR:CE2	2.56	0.40
1:A:323:LEU:HB2	1:A:362:PRO:HA	2.04	0.40
1:B:414:ARG:HH11	1:B:414:ARG:HG3	1.84	0.40
1:C:560:LYS:O	4:C:701:HOH:O	2.22	0.40
1:D:89:GLU:O	1:D:93:LYS:HG3	2.21	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	553/568 (97%)	536 (97%)	17 (3%)	0	100	100
1	B	553/568 (97%)	536 (97%)	17 (3%)	0	100	100
1	C	553/568 (97%)	536 (97%)	17 (3%)	0	100	100
1	D	553/568 (97%)	539 (98%)	14 (2%)	0	100	100
All	All	2212/2272 (97%)	2147 (97%)	65 (3%)	0	100	100

There are no Ramachandran outliers to report.

### 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	490/503 (97%)	481 (98%)	9 (2%)	59	55
1	B	490/503 (97%)	482 (98%)	8 (2%)	62	59
1	C	490/503 (97%)	475 (97%)	15 (3%)	40	33
1	D	490/503 (97%)	478 (98%)	12 (2%)	49	42
All	All	1960/2012 (97%)	1916 (98%)	44 (2%)	52	46

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	101	ASP
1	A	173	ASP
1	A	201	GLN
1	A	213	ARG
1	A	316	LYS
1	A	334	VAL
1	A	495	LEU
1	A	514	THR
1	A	549	LYS
1	B	64	ASN
1	B	81	MET
1	B	101	ASP
1	B	137	ASN
1	B	173	ASP
1	B	201	GLN
1	B	255	MET
1	B	470	ARG
1	C	32	ASP
1	C	76	GLU
1	C	101	ASP
1	C	137	ASN
1	C	158	GLU
1	C	164	LEU
1	C	173	ASP
1	C	201	GLN
1	C	224	ARG
1	C	301	LEU
1	C	305	LYS
1	C	346	LYS
1	C	389	LEU
1	C	435	THR
1	C	541	GLU
1	D	32	ASP

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Mol	Chain	Res	Type
1	D	101	ASP
1	D	102	LEU
1	D	147	LYS
1	D	173	ASP
1	D	203	ASP
1	D	256	SER
1	D	379	ASP
1	D	389	LEU
1	D	427	GLU
1	D	504	LEU
1	D	514	THR

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

Mol	Chain	Res	Type
1	D	279	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](#)

Of 7 ligands modelled in this entry, 4 are monoatomic - leaving 3 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	PNG	B	601	-	21,22,22	1.23	3 (14%)	29,31,31	1.61	4 (13%)
2	PNG	A	601	-	21,22,22	1.33	4 (19%)	29,31,31	1.78	6 (20%)
2	PNG	C	601	-	21,22,22	1.28	3 (14%)	29,31,31	1.57	3 (10%)

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	PNG	B	601	-	-	7/8/30/30	0/2/2/2
2	PNG	A	601	-	-	8/8/30/30	0/2/2/2
2	PNG	C	601	-	-	6/8/30/30	0/2/2/2

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601	PNG	O1-C7	2.73	1.43	1.38
2	C	601	PNG	O1-C7	2.64	1.43	1.38
2	A	601	PNG	O5-C5	-2.40	1.38	1.44
2	B	601	PNG	O5-C5	-2.29	1.38	1.44
2	B	601	PNG	O1-C7	2.29	1.42	1.38
2	C	601	PNG	O5-C5	-2.15	1.39	1.44
2	A	601	PNG	C4-C5	-2.13	1.48	1.53
2	C	601	PNG	O2-C2	-2.11	1.38	1.43
2	B	601	PNG	C4-C5	-2.05	1.48	1.53
2	A	601	PNG	O3-C3	-2.01	1.38	1.43

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	PNG	O1-C1-C2	6.41	116.44	107.14
2	C	601	PNG	O1-C1-C2	6.30	116.28	107.14
2	B	601	PNG	O1-C1-C2	5.23	114.73	107.14
2	B	601	PNG	O5-C1-O1	3.69	117.67	108.29
2	C	601	PNG	C1-O5-C5	3.48	120.53	113.69
2	A	601	PNG	O5-C1-O1	3.41	116.96	108.29
2	B	601	PNG	C1-O5-C5	3.13	119.82	113.69
2	B	601	PNG	C7-O1-C1	-2.96	113.45	117.79
2	C	601	PNG	O5-C1-O1	2.72	115.22	108.29
2	A	601	PNG	C1-O5-C5	2.69	118.97	113.69

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	PNG	O5-C5-C4	2.24	113.76	109.69
2	A	601	PNG	C12-C11-C10	-2.16	117.08	120.08
2	A	601	PNG	C4-C3-C2	2.10	114.48	110.82

There are no chirality outliers.

All (21) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	601	PNG	O5-C1-O1-C7
2	B	601	PNG	C9-C10-N1-O7
2	B	601	PNG	C11-C10-N1-O7
2	A	601	PNG	O5-C1-O1-C7
2	A	601	PNG	C9-C10-N1-O7
2	A	601	PNG	C11-C10-N1-O7
2	C	601	PNG	C2-C1-O1-C7
2	C	601	PNG	O5-C1-O1-C7
2	A	601	PNG	O5-C5-C6-O6
2	C	601	PNG	C4-C5-C6-O6
2	C	601	PNG	O5-C5-C6-O6
2	A	601	PNG	C4-C5-C6-O6
2	A	601	PNG	C2-C1-O1-C7
2	B	601	PNG	C12-C7-O1-C1
2	A	601	PNG	C12-C7-O1-C1
2	A	601	PNG	C8-C7-O1-C1
2	C	601	PNG	C12-C7-O1-C1
2	B	601	PNG	C8-C7-O1-C1
2	B	601	PNG	C2-C1-O1-C7
2	C	601	PNG	C8-C7-O1-C1
2	B	601	PNG	O5-C5-C6-O6

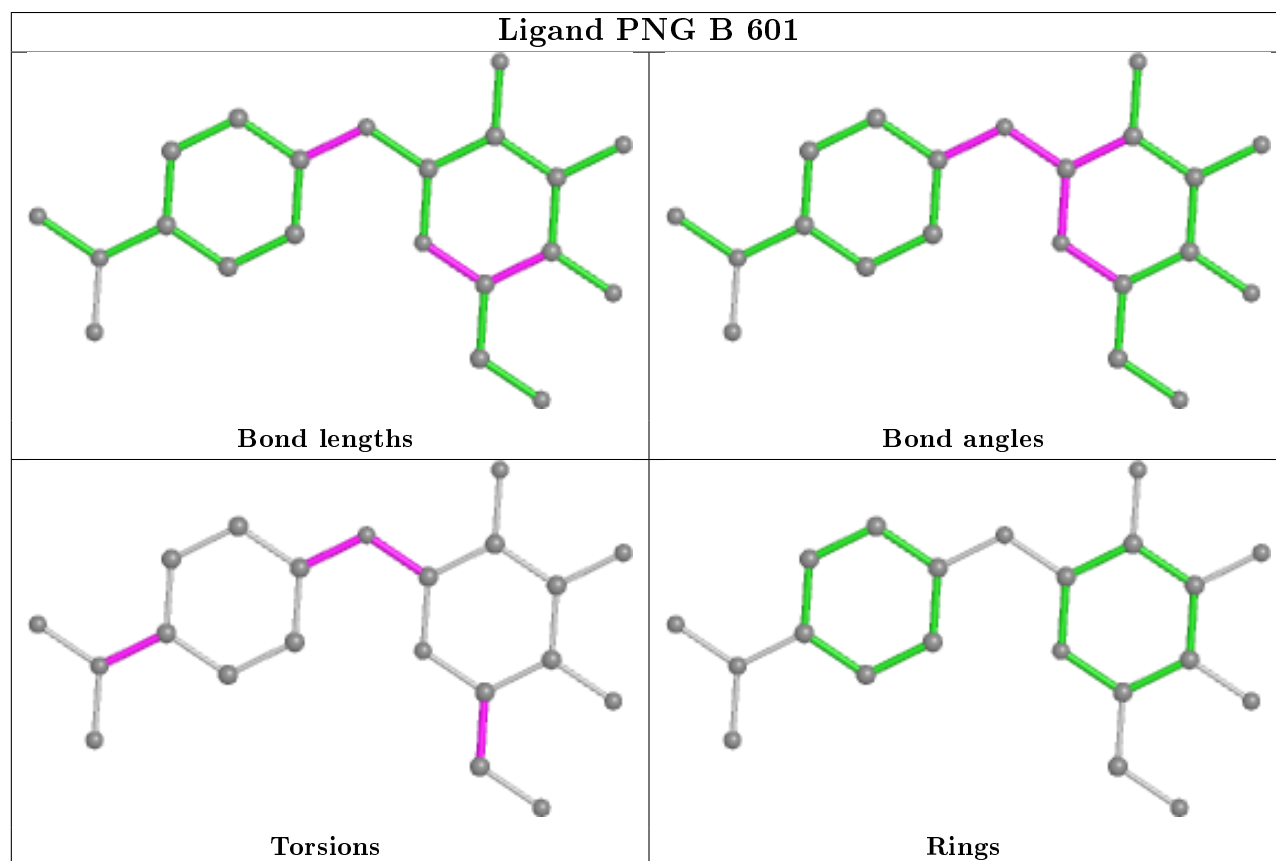
There are no ring outliers.

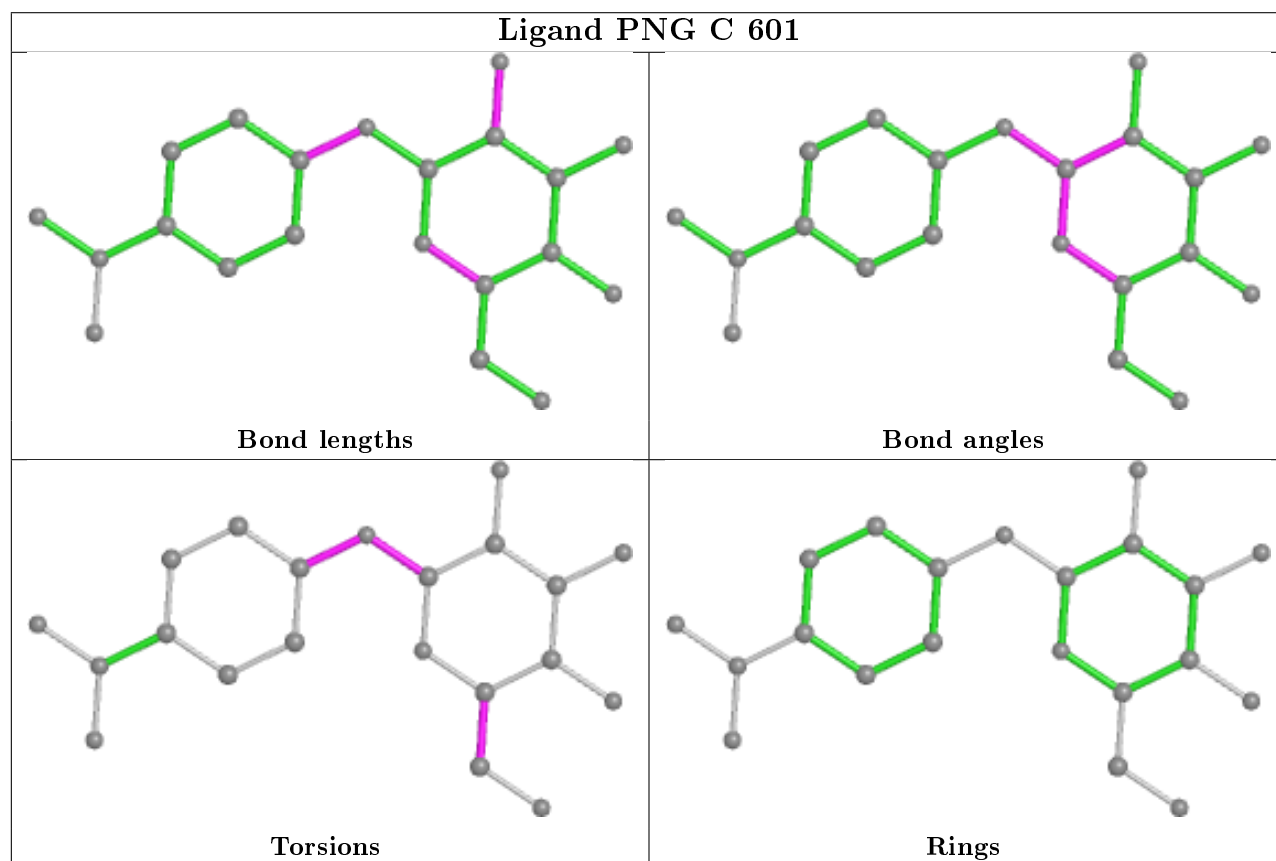
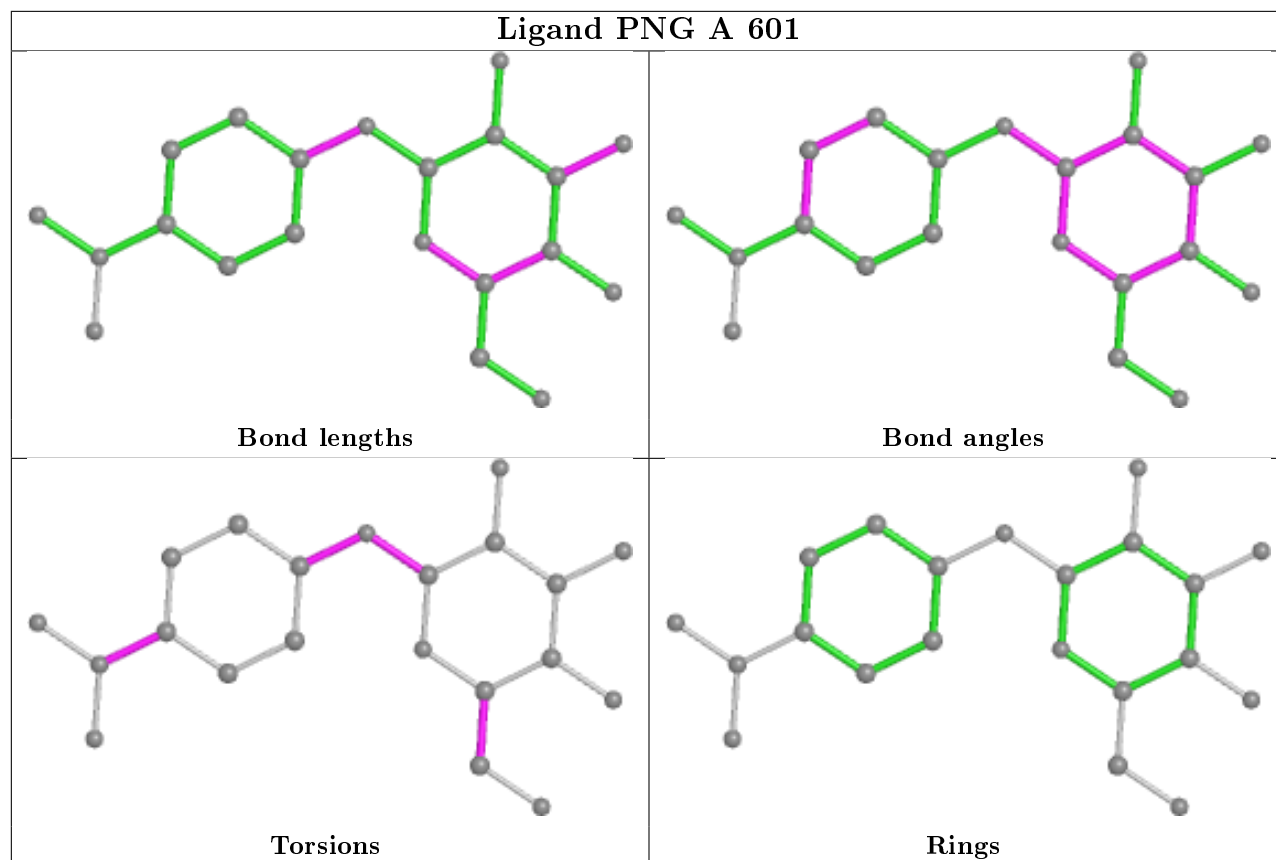
3 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	601	PNG	3	0
2	A	601	PNG	3	0
2	C	601	PNG	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

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 ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	555/568 (97%)	-0.67	2 (0%) 92 93	9, 19, 35, 55	0
1	B	555/568 (97%)	-0.54	10 (1%) 68 71	10, 23, 44, 67	0
1	C	555/568 (97%)	-0.44	9 (1%) 72 74	12, 26, 45, 64	0
1	D	555/568 (97%)	-0.59	1 (0%) 95 95	13, 23, 39, 52	0
All	All	2220/2272 (97%)	-0.56	22 (0%) 82 84	9, 23, 42, 67	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	136	PRO	4.7
1	C	220	GLY	3.6
1	A	289	ASN	3.3
1	C	136	PRO	3.3
1	C	6	ASN	3.0
1	C	137	ASN	3.0
1	D	6	ASN	2.9
1	B	219	GLU	2.8
1	C	139	ALA	2.7
1	B	138	GLY	2.7
1	C	219	GLU	2.6
1	B	288	PRO	2.6
1	B	218	ASP	2.5
1	C	138	GLY	2.5
1	B	213	ARG	2.4
1	C	289	ASN	2.3
1	B	139	ALA	2.3
1	B	137	ASN	2.2
1	B	289	ASN	2.1
1	A	6	ASN	2.1
1	C	135	LYS	2.1
1	B	155	PHE	2.0

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

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

## 6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 6.4 Ligands [i](#)

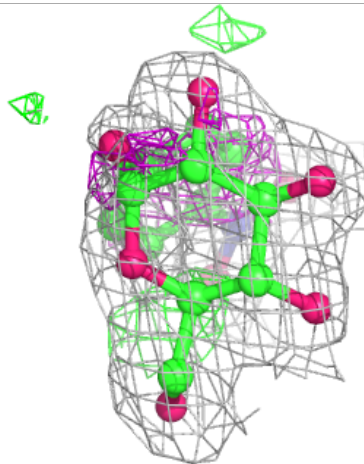
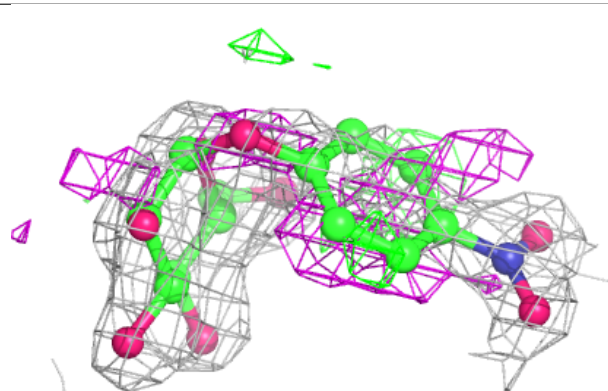
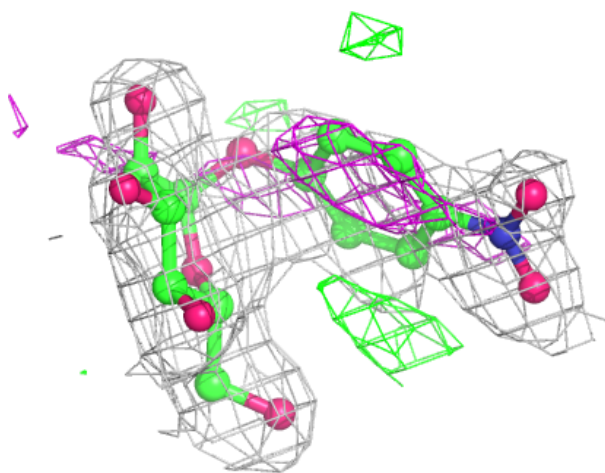
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 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
2	PNG	C	601	21/21	0.82	0.23	35,47,51,63	0
2	PNG	B	601	21/21	0.85	0.14	36,44,50,55	0
2	PNG	A	601	21/21	0.86	0.18	20,35,43,52	0
3	MG	C	602	1/1	0.97	0.04	22,22,22,22	0
3	MG	D	601	1/1	0.99	0.06	23,23,23,23	0
3	MG	A	602	1/1	0.99	0.06	18,18,18,18	0
3	MG	B	602	1/1	0.99	0.06	25,25,25,25	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.

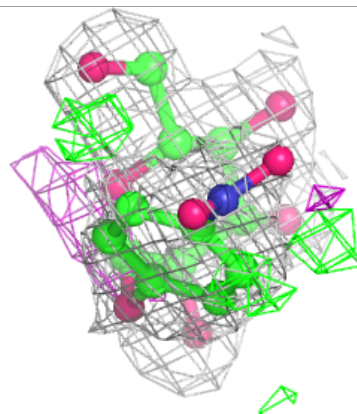
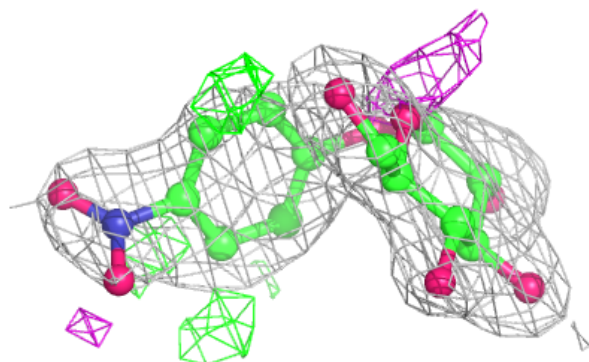
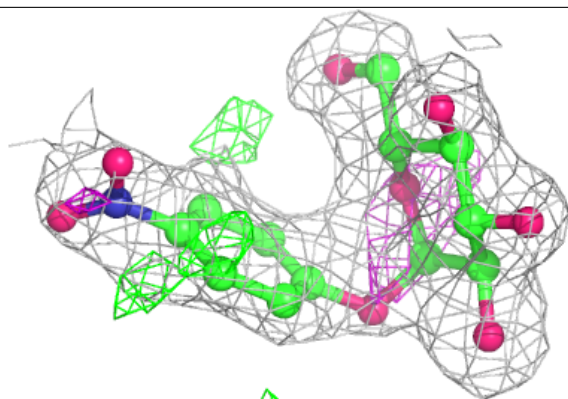
**Electron density around PNG C 601:**

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



**Electron density around PNG B 601:**

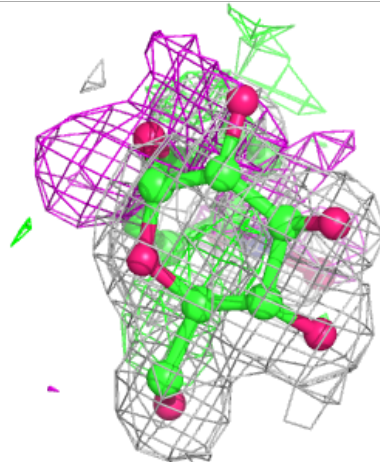
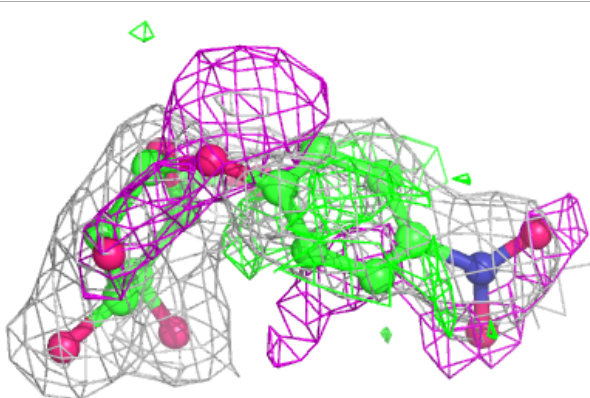
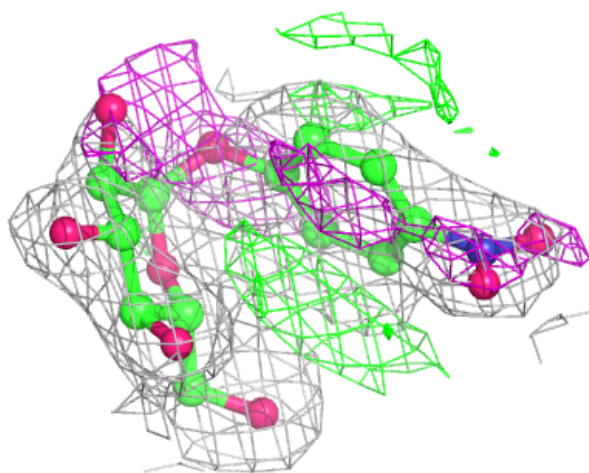
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





**Electron density around PNG A 601:**

$2mF_o-DF_c$  (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.