



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 8, 2020 – 07:07 PM BST

PDB ID : 4RGG  
Title : Structure of the lactococcal phage 1358 receptor binding protein in complex with GlcNAc-1P  
Authors : Spinelli, S.; Moineau, S.; Cambillau, C.  
Deposited on : 2014-09-30  
Resolution : 2.15 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

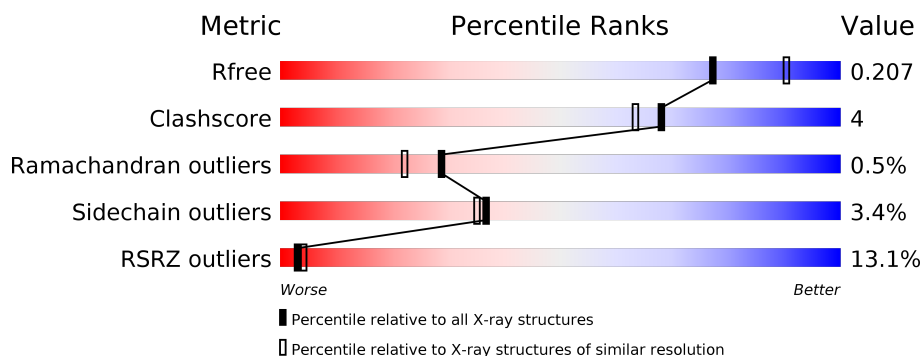
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.15 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	1479 (2.16-2.16)
Clashscore	141614	1585 (2.16-2.16)
Ramachandran outliers	138981	1560 (2.16-2.16)
Sidechain outliers	138945	1559 (2.16-2.16)
RSRZ outliers	127900	1456 (2.16-2.16)

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	393	<div> <div>15%</div> <div>88%</div> <div>9%</div> <div>..</div> </div>
1	B	393	<div> <div>11%</div> <div>87%</div> <div>9%</div> <div>..</div> </div>

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 6596 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.

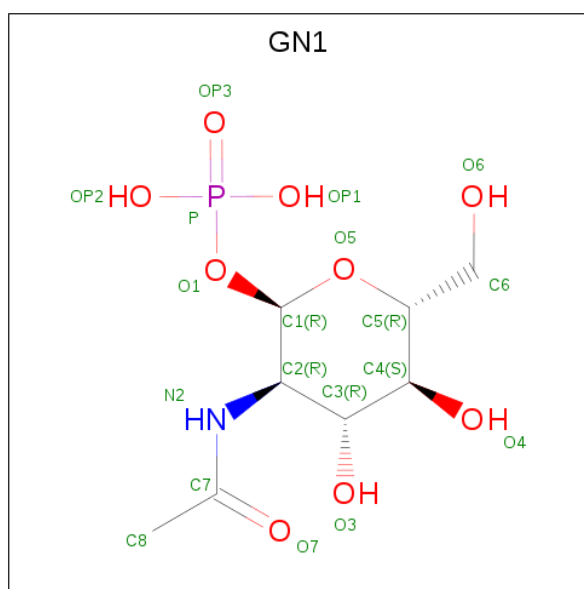
- Molecule 1 is a protein called Putative phage structural protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	390	Total	C	N	O	S	0	0	0
			3028	1902	522	595	9			
1	B	390	Total	C	N	O	S	0	0	0
			3028	1902	522	595	9			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	expression tag	UNP D3W0F1
B	1	GLY	-	expression tag	UNP D3W0F1

- Molecule 2 is 2-acetamido-2-deoxy-1-O-phosphono-alpha-D-glucopyranose (three-letter code: GN1) (formula: C<sub>8</sub>H<sub>16</sub>NO<sub>9</sub>P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			19	8	1	9	1		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	N	O	P	0	0
			19	8	1	9	1		

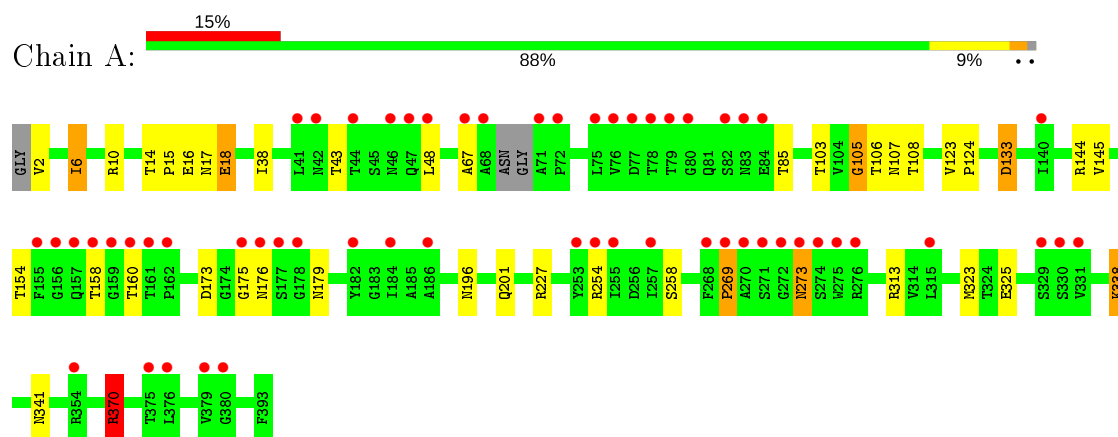
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	255	Total	O	0	0
			255	255		
3	B	247	Total	O	0	0
			247	247		

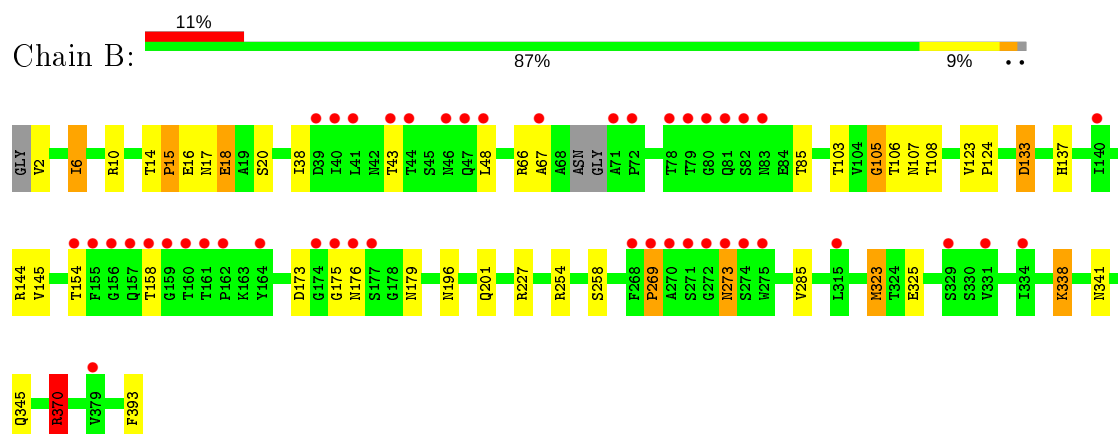
### 3 Residue-property plots [i](#)

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: Putative phage structural protein



- Molecule 1: Putative phage structural protein



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 3	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	165.91Å 165.91Å 165.91Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.89 – 2.15 46.02 – 2.15	Depositor EDS
% Data completeness (in resolution range)	100.0 (47.89-2.15) 100.0 (46.02-2.15)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.20 (at 2.16Å)	Xtriage
Refinement program	REFMAC 5.8.0069	Depositor
R, $R_{free}$	0.180 , 0.204 0.187 , 0.207	Depositor DCC
$R_{free}$ test set	3914 reflections (4.74%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	38.2	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 42.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.039 for l,-k,h	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	6596	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.09% 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: GN1

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	0.98	4/3094 (0.1%)	1.06	15/4216 (0.4%)
1	B	0.99	3/3094 (0.1%)	1.08	13/4216 (0.3%)
All	All	0.98	7/6188 (0.1%)	1.07	28/8432 (0.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	5
1	B	0	6
All	All	0	11

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	258	SER	CB-OG	-6.82	1.33	1.42
1	B	370	ARG	CD-NE	-6.69	1.35	1.46
1	A	338	LYS	CE-NZ	6.54	1.65	1.49
1	A	258	SER	CB-OG	-6.35	1.33	1.42
1	B	338	LYS	CD-CE	6.07	1.66	1.51
1	A	338	LYS	CD-CE	5.79	1.65	1.51
1	A	18	GLU	CG-CD	5.27	1.59	1.51

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	370	ARG	NE-CZ-NH2	-22.42	109.09	120.30
1	A	370	ARG	NE-CZ-NH2	-19.39	110.61	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	370	ARG	NE-CZ-NH1	18.80	129.70	120.30
1	A	370	ARG	NE-CZ-NH1	15.60	128.10	120.30
1	A	227	ARG	NE-CZ-NH2	-11.62	114.49	120.30
1	B	227	ARG	NE-CZ-NH2	-9.68	115.46	120.30
1	B	254	ARG	NE-CZ-NH2	-8.77	115.92	120.30
1	B	10	ARG	NE-CZ-NH2	-8.19	116.21	120.30
1	A	254	ARG	NE-CZ-NH2	-7.88	116.36	120.30
1	A	338	LYS	CG-CD-CE	7.63	134.79	111.90
1	B	370	ARG	CD-NE-CZ	7.60	134.25	123.60
1	A	370	ARG	CD-NE-CZ	7.48	134.07	123.60
1	A	10	ARG	NE-CZ-NH2	-7.20	116.70	120.30
1	B	338	LYS	CG-CD-CE	7.03	133.00	111.90
1	B	18	GLU	OE1-CD-OE2	-6.79	115.16	123.30
1	A	133	ASP	CB-CG-OD2	-6.38	112.56	118.30
1	A	227	ARG	NE-CZ-NH1	6.24	123.42	120.30
1	B	227	ARG	CG-CD-NE	-6.19	98.80	111.80
1	A	227	ARG	CG-CD-NE	-6.18	98.82	111.80
1	B	133	ASP	CB-CG-OD2	-5.91	112.98	118.30
1	A	370	ARG	CG-CD-NE	-5.83	99.55	111.80
1	A	133	ASP	CB-CG-OD1	5.68	123.41	118.30
1	B	323	MET	CA-CB-CG	5.58	122.79	113.30
1	A	254	ARG	NE-CZ-NH1	5.57	123.09	120.30
1	A	10	ARG	CG-CD-NE	-5.43	100.39	111.80
1	A	313	ARG	NE-CZ-NH2	-5.43	117.59	120.30
1	B	370	ARG	CG-CD-NE	-5.39	100.47	111.80
1	B	10	ARG	CG-CD-NE	-5.36	100.54	111.80

There are no chirality outliers.

All (11) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	105	GLY	Peptide
1	A	173	ASP	Peptide
1	A	175	GLY	Peptide
1	A	6	ILE	Peptide
1	A	67	ALA	Peptide
1	B	105	GLY	Peptide
1	B	173	ASP	Peptide
1	B	175	GLY	Peptide
1	B	370	ARG	Sidechain
1	B	6	ILE	Peptide
1	B	67	ALA	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	3028	0	2893	19	2
1	B	3028	0	2893	26	3
2	A	19	0	0	0	0
2	B	19	0	0	0	0
3	A	255	0	0	2	1
3	B	247	0	0	6	1
All	All	6596	0	5786	45	7

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 4.

All (45) 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:133:ASP:OD1	1:A:338:LYS:NZ	1.93	1.01
1:B:133:ASP:OD1	1:B:338:LYS:NZ	1.98	0.97
1:B:137:HIS:ND1	3:B:746:HOH:O	2.07	0.88
1:A:38:ILE:HD11	1:A:48:LEU:HD11	1.59	0.85
1:A:106:THR:HG22	1:A:107:ASN:N	1.92	0.83
1:B:196:ASN:H	1:B:201:GLN:HE22	1.27	0.82
1:B:106:THR:HG22	1:B:107:ASN:N	1.95	0.81
1:B:38:ILE:HD11	1:B:48:LEU:HD11	1.61	0.81
1:A:196:ASN:H	1:A:201:GLN:HE22	1.30	0.79
1:A:85:THR:OG1	1:A:154:THR:HG22	1.84	0.78
1:A:106:THR:HG22	1:A:107:ASN:H	1.49	0.77
1:B:106:THR:HG22	1:B:107:ASN:H	1.51	0.74
1:B:85:THR:OG1	1:B:154:THR:HG22	1.87	0.74
1:B:20:SER:OG	3:B:746:HOH:O	2.06	0.73
1:B:338:LYS:NZ	3:B:744:HOH:O	2.13	0.73
1:A:325:GLU:OE2	1:A:370:ARG:HD3	1.93	0.67
1:B:106:THR:HG21	1:B:108:THR:HG23	1.77	0.66
1:B:106:THR:CG2	1:B:108:THR:HG23	2.26	0.66
1:B:325:GLU:OE2	1:B:370:ARG:HD3	1.98	0.63
1:A:106:THR:CG2	1:A:108:THR:HG23	2.28	0.63
1:B:393:PHE:OXT	3:B:747:HOH:O	2.16	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:106:THR:HG21	1:A:108:THR:HG23	1.80	0.61
1:A:176:ASN:ND2	3:A:702:HOH:O	2.37	0.57
1:A:106:THR:CG2	1:A:107:ASN:N	2.64	0.54
1:B:106:THR:CG2	1:B:107:ASN:N	2.66	0.54
1:B:66:ARG:HD2	3:B:666:HOH:O	2.07	0.53
1:B:196:ASN:N	1:B:201:GLN:HE22	2.03	0.52
1:B:176:ASN:ND2	3:B:699:HOH:O	2.43	0.51
1:A:160:THR:HG21	3:A:639:HOH:O	2.12	0.50
1:A:43:THR:HG23	1:A:48:LEU:HA	1.96	0.48
1:B:43:THR:HG23	1:B:48:LEU:HA	1.95	0.47
1:A:106:THR:CG2	1:A:107:ASN:H	2.24	0.47
1:A:123:VAL:HB	1:A:124:PRO:HD3	1.97	0.46
1:A:14:THR:H	1:A:17:ASN:HD22	1.64	0.46
1:B:106:THR:CG2	1:B:107:ASN:H	2.26	0.46
1:A:103:THR:HG23	1:A:106:THR:H	1.80	0.46
1:B:105:GLY:N	1:B:106:THR:O	2.49	0.45
1:B:103:THR:HG23	1:B:106:THR:H	1.80	0.45
1:B:269:PRO:HA	1:B:273:ASN:HD21	1.82	0.44
1:A:105:GLY:N	1:A:106:THR:O	2.50	0.43
1:B:14:THR:H	1:B:17:ASN:HD22	1.67	0.43
1:A:269:PRO:HA	1:A:273:ASN:HD21	1.83	0.43
1:B:123:VAL:HB	1:B:124:PRO:HD3	2.03	0.41
1:B:14:THR:HA	1:B:15:PRO:HD2	1.88	0.41
1:B:285:VAL:HG22	1:B:345:GLN:HB2	2.04	0.40

All (7) 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:A:18:GLU:OE1	1:A:18:GLU:OE2[12_455]	1.30	0.90
1:B:18:GLU:OE1	1:B:18:GLU:OE2[5_555]	1.72	0.48
3:B:714:HOH:O	3:B:731:HOH:O[5_555]	1.87	0.33
1:B:18:GLU:OE2	1:B:18:GLU:OE2[5_555]	1.98	0.22
1:A:18:GLU:CD	1:A:18:GLU:OE2[12_455]	2.04	0.16
1:B:18:GLU:CD	1:B:18:GLU:OE2[5_555]	2.10	0.10
3:A:680:HOH:O	3:A:680:HOH:O[6_445]	2.11	0.09

## 5.3 Torsion angles

### 5.3.1 Protein backbone

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	386/393 (98%)	367 (95%)	17 (4%)	2 (0%)	29	22
1	B	386/393 (98%)	369 (96%)	15 (4%)	2 (0%)	29	22
All	All	772/786 (98%)	736 (95%)	32 (4%)	4 (0%)	29	22

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	269	PRO
1	B	15	PRO
1	A	15	PRO
1	B	269	PRO

### 5.3.2 Protein sidechains

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	327/332 (98%)	316 (97%)	11 (3%)	37	35
1	B	327/332 (98%)	316 (97%)	11 (3%)	37	35
All	All	654/664 (98%)	632 (97%)	22 (3%)	37	35

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	VAL
1	A	6	ILE

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Mol	Chain	Res	Type
1	A	16	GLU
1	A	144	ARG
1	A	145	VAL
1	A	158	THR
1	A	179	ASN
1	A	273	ASN
1	A	323	MET
1	A	341	ASN
1	A	370	ARG
1	B	2	VAL
1	B	6	ILE
1	B	16	GLU
1	B	144	ARG
1	B	145	VAL
1	B	158	THR
1	B	179	ASN
1	B	273	ASN
1	B	323	MET
1	B	341	ASN
1	B	370	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	17	ASN
1	A	28	ASN
1	A	81	GLN
1	A	157	GLN
1	A	179	ASN
1	A	196	ASN
1	A	201	GLN
1	A	222	ASN
1	A	273	ASN
1	A	294	ASN
1	A	371	ASN
1	B	17	ASN
1	B	28	ASN
1	B	52	HIS
1	B	81	GLN
1	B	157	GLN
1	B	179	ASN
1	B	196	ASN

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Mol	Chain	Res	Type
1	B	201	GLN
1	B	222	ASN
1	B	273	ASN
1	B	294	ASN
1	B	371	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](#)

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
2	GN1	B	401	-	18,19,19	1.46	2 (11%)	27,28,28	1.18	3 (11%)
2	GN1	A	401	-	18,19,19	1.26	3 (16%)	27,28,28	1.53	8 (29%)

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	GN1	B	401	-	-	1/10/31/31	0/1/1/1
2	GN1	A	401	-	-	1/10/31/31	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	GN1	C2-N2	3.00	1.50	1.45
2	A	401	GN1	P-O1	2.98	1.64	1.59
2	B	401	GN1	P-O1	2.58	1.64	1.59
2	A	401	GN1	P-OP3	2.30	1.57	1.50
2	A	401	GN1	C3-C2	2.08	1.57	1.53

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	GN1	OP2-P-OP1	3.18	119.78	107.64
2	A	401	GN1	C1-C2-N2	-2.84	106.12	111.00
2	B	401	GN1	C3-C4-C5	2.64	114.95	110.24
2	A	401	GN1	O4-C4-C5	-2.62	102.79	109.30
2	B	401	GN1	O4-C4-C3	-2.51	104.55	110.35
2	B	401	GN1	OP2-P-OP3	2.50	120.48	110.68
2	A	401	GN1	C3-C2-N2	-2.49	105.92	110.62
2	A	401	GN1	OP2-P-O1	-2.26	95.86	105.99
2	A	401	GN1	O3-C3-C4	2.21	115.46	110.35
2	A	401	GN1	O3-C3-C2	2.09	113.89	109.66
2	A	401	GN1	O5-C5-C4	2.03	113.37	109.69

There are no chirality outliers.

All (2) torsion outliers are listed below:

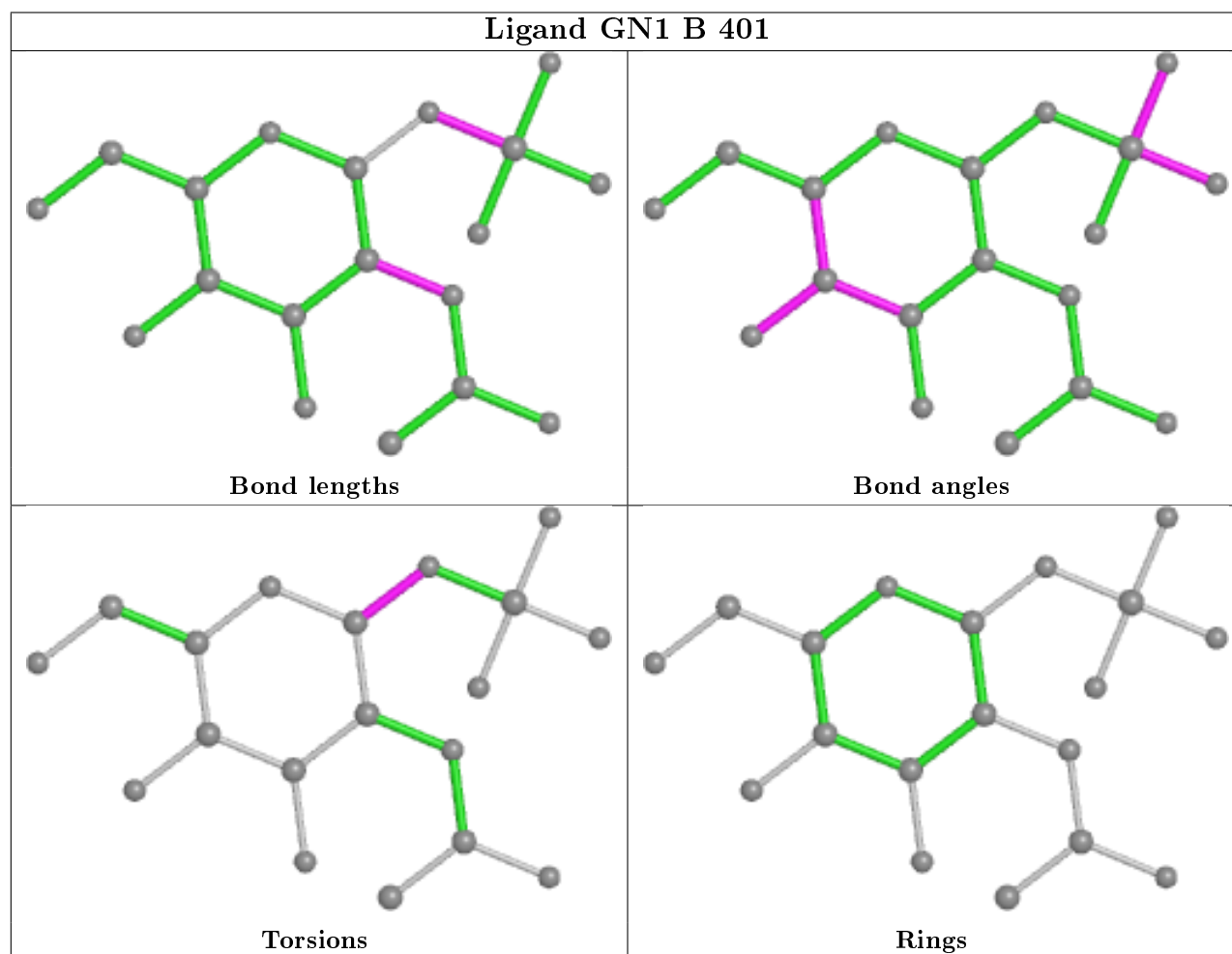
Mol	Chain	Res	Type	Atoms
2	B	401	GN1	O5-C1-O1-P
2	A	401	GN1	O5-C1-O1-P

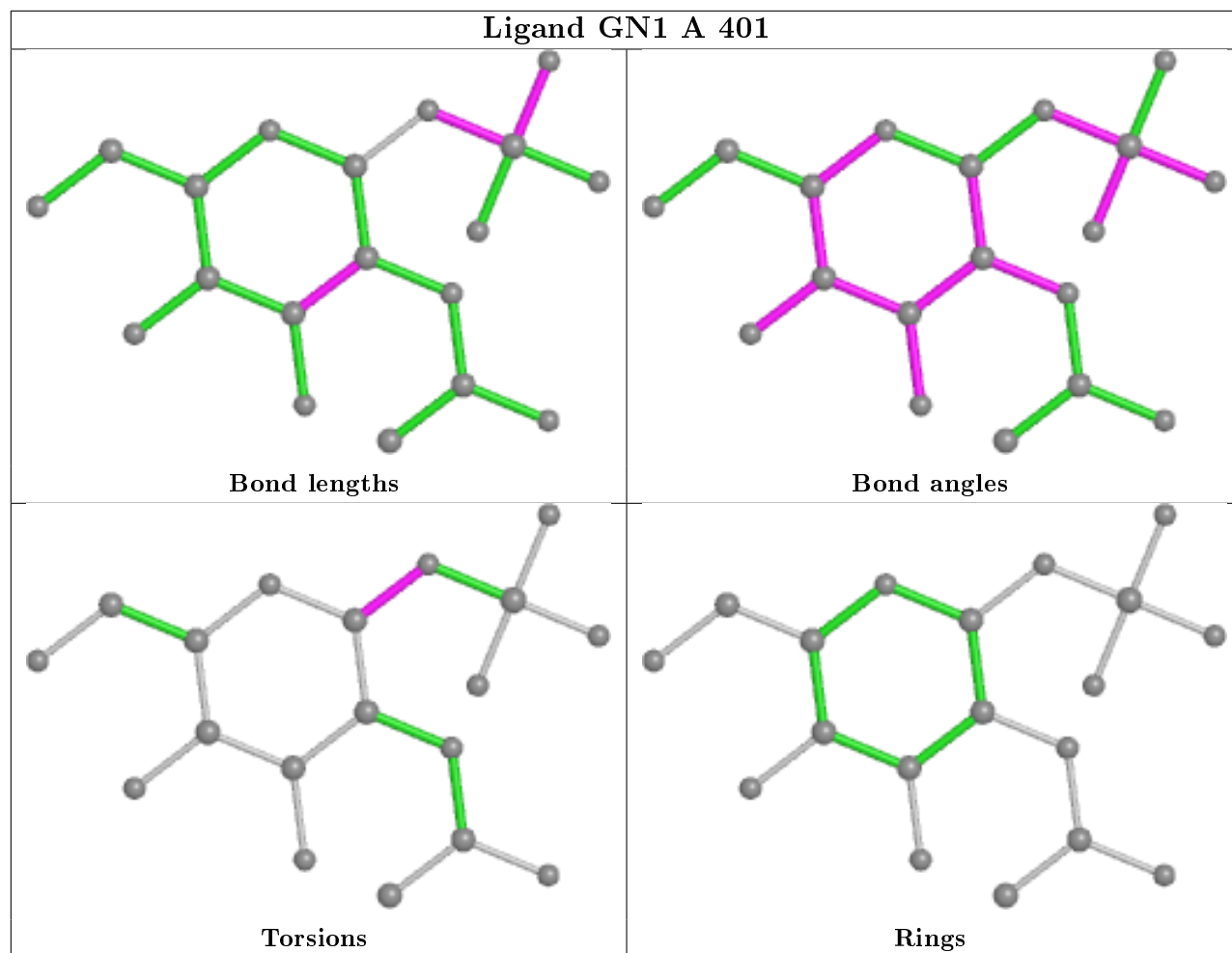
There are no ring outliers.

No monomer is involved in short contacts.

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 ⓘ

### 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	390/393 (99%)	0.56	57 (14%) 2 3	23, 40, 91, 129	0
1	B	390/393 (99%)	0.35	45 (11%) 4 6	23, 40, 88, 125	0
All	All	780/786 (99%)	0.45	102 (13%) 3 4	23, 40, 89, 129	0

All (102) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	275	TRP	9.6
1	B	270	ALA	6.3
1	B	71	ALA	6.1
1	A	270	ALA	5.6
1	A	271	SER	5.6
1	A	140	ILE	5.5
1	B	275	TRP	5.3
1	A	160	THR	4.8
1	B	160	THR	4.8
1	A	156	GLY	4.8
1	A	276	ARG	4.6
1	A	68	ALA	4.5
1	B	155	PHE	4.4
1	A	44	THR	4.4
1	A	272	GLY	4.3
1	A	46	ASN	4.2
1	A	268	PHE	4.1
1	B	271	SER	4.1
1	B	174	GLY	4.0
1	B	159	GLY	4.0
1	A	379	VAL	3.9
1	A	41	LEU	3.9
1	A	155	PHE	3.8
1	B	47	GLN	3.8

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Mol	Chain	Res	Type	RSRZ
1	B	80	GLY	3.8
1	B	161	THR	3.8
1	A	67	ALA	3.8
1	B	79	THR	3.7
1	A	71	ALA	3.7
1	A	331	VAL	3.7
1	B	140	ILE	3.7
1	B	157	GLN	3.6
1	B	156	GLY	3.6
1	B	48	LEU	3.5
1	B	268	PHE	3.5
1	A	83	ASN	3.5
1	B	158	THR	3.5
1	A	273	ASN	3.5
1	B	40	ILE	3.5
1	B	177	SER	3.4
1	A	177	SER	3.4
1	A	158	THR	3.4
1	B	46	ASN	3.4
1	A	79	THR	3.4
1	A	75	LEU	3.3
1	B	44	THR	3.3
1	A	329	SER	3.3
1	B	78	THR	3.2
1	B	43	THR	3.2
1	B	83	ASN	3.2
1	A	47	GLN	3.2
1	A	255	ILE	3.1
1	A	80	GLY	3.1
1	A	161	THR	3.1
1	B	272	GLY	3.0
1	A	157	GLN	2.9
1	B	176	ASN	2.9
1	A	375	THR	2.8
1	B	175	GLY	2.8
1	B	273	ASN	2.8
1	B	331	VAL	2.7
1	B	162	PRO	2.7
1	B	379	VAL	2.7
1	A	330	SER	2.7
1	B	72	PRO	2.6
1	A	354	ARG	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	376	LEU	2.6
1	B	269	PRO	2.6
1	A	182	TYR	2.5
1	B	39	ASP	2.5
1	A	274	SER	2.5
1	A	48	LEU	2.5
1	A	78	THR	2.5
1	A	162	PRO	2.5
1	A	178	GLY	2.4
1	A	380	GLY	2.4
1	A	175	GLY	2.4
1	A	254	ARG	2.4
1	A	184	ILE	2.4
1	A	257	ILE	2.4
1	A	186	ALA	2.4
1	A	84	GLU	2.4
1	B	81	GLN	2.4
1	B	334	ILE	2.3
1	B	41	LEU	2.3
1	A	42	ASN	2.3
1	B	329	SER	2.3
1	A	176	ASN	2.2
1	A	77	ASP	2.2
1	B	154	THR	2.2
1	A	269	PRO	2.2
1	A	315	LEU	2.2
1	B	315	LEU	2.2
1	B	67	ALA	2.2
1	A	76	VAL	2.1
1	B	164	TYR	2.1
1	B	274	SER	2.1
1	A	72	PRO	2.1
1	A	159	GLY	2.1
1	A	253	TYR	2.1
1	A	82	SER	2.0
1	B	82	SER	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

### 6.3 Carbohydrates

There are no monosaccharides in this entry.

### 6.4 Ligands

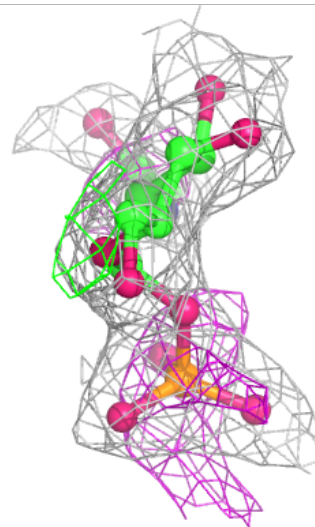
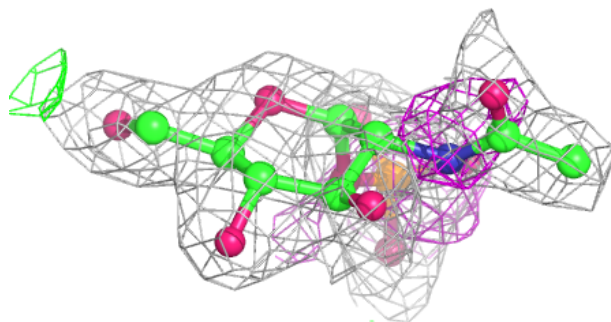
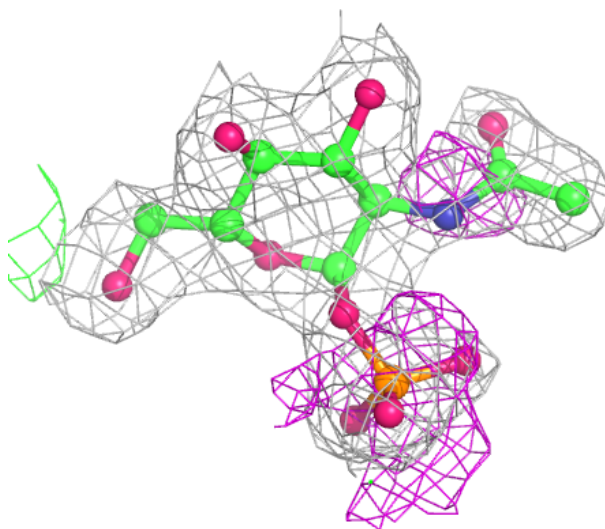
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

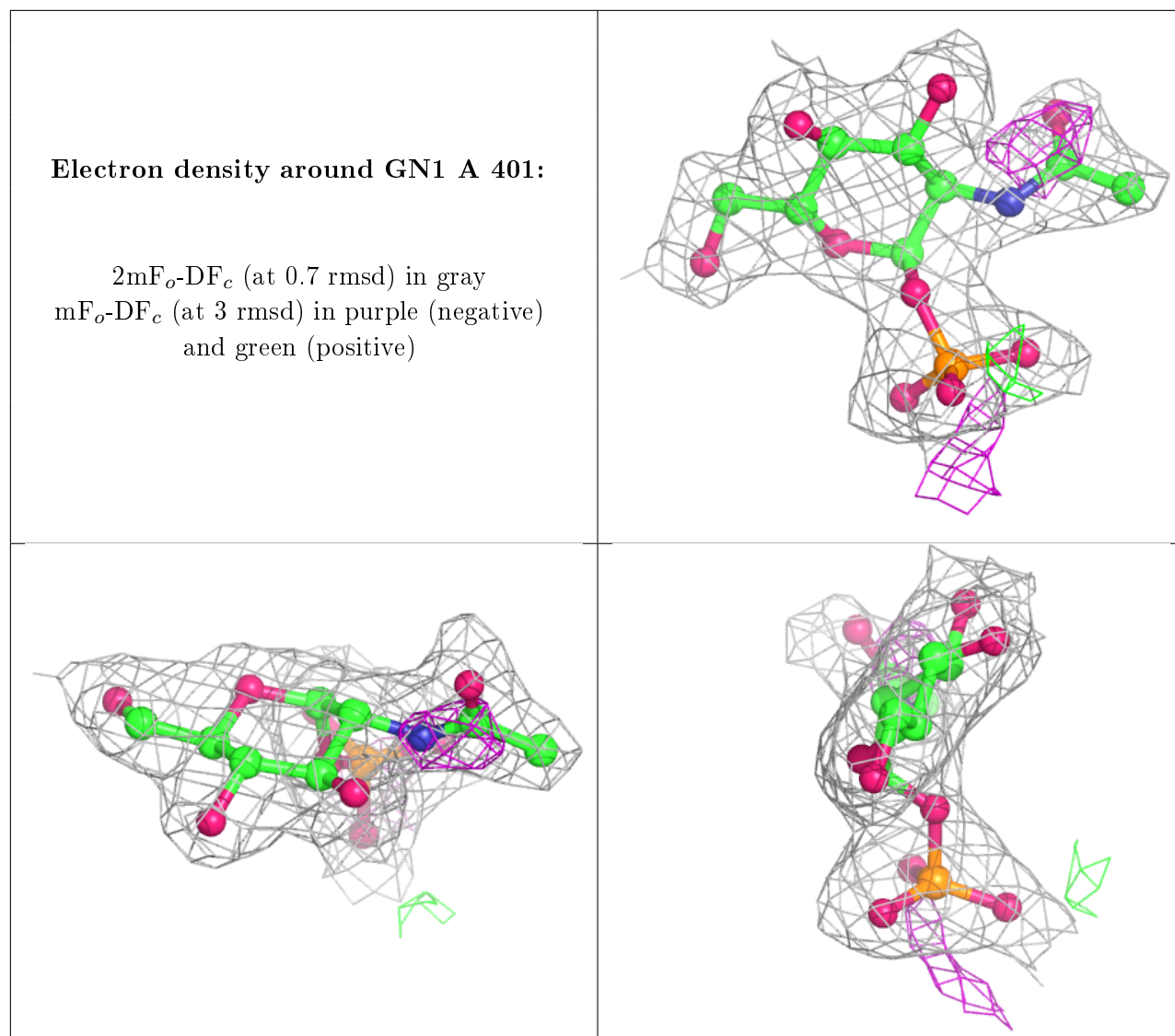
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	GN1	B	401	19/19	0.85	0.25	37,52,62,63	0
2	GN1	A	401	19/19	0.93	0.19	28,41,54,55	0

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 GN1 B 401:**

$2mF_o-DF_c$  (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.