



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 09:14 AM UTC

PDB ID : 6XJT / pdb_00006xjt
Title : Crystal Structure of KPT-8602 bound to CRM1 (537-DLTVK-541 to GLCEQ)
Authors : Baumhardt, J.M.; Chook, Y.M.
Deposited on : 2020-06-24
Resolution : 2.41 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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

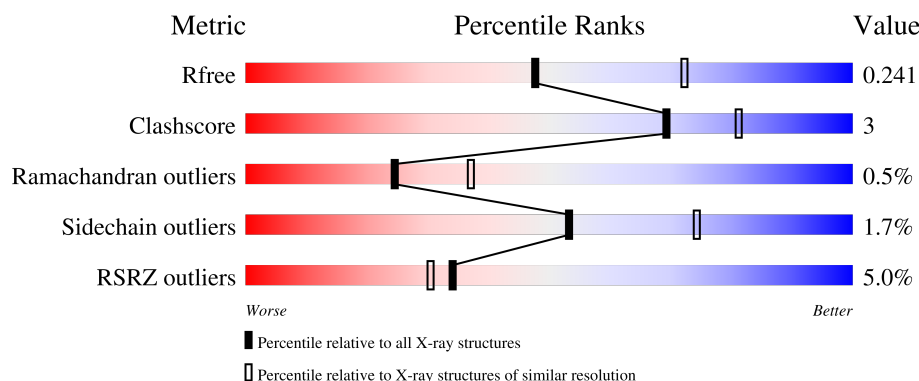
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.41 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}	180053	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	216	
2	B	140	
3	C	1024	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 22274 atoms, of which 11178 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 GTP-binding nuclear protein Ran.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	208	Total	C	H	N	O	S	0	5	0
			3391	1093	1697	291	304	6			

- Molecule 2 is a protein called Ran-specific GTPase-activating protein 1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	123	Total	C	H	N	O	S	0	3	0
			2060	652	1031	181	191	5			

- Molecule 3 is a protein called Exportin-1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	C	1011	Total	C	H	N	O	S	0	42	0
			16736	5336	8426	1367	1562	45			

There are 45 discrepancies between the modelled and reference sequences:

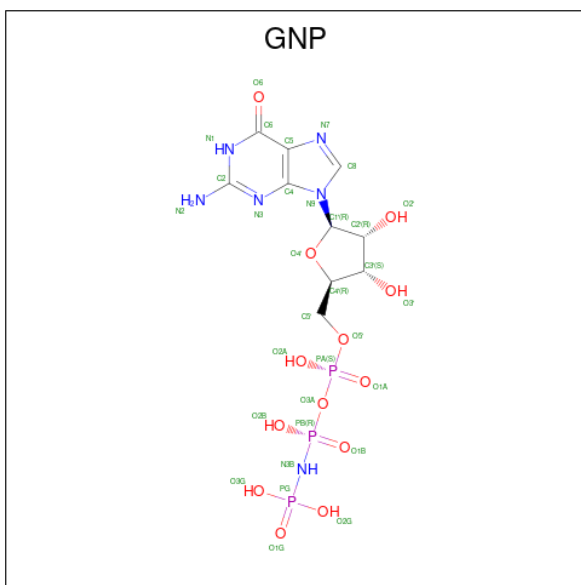
Chain	Residue	Modelled	Actual	Comment	Reference
C	-2	GLY	-	expression tag	UNP P30822
C	-1	GLY	-	expression tag	UNP P30822
C	0	SER	-	expression tag	UNP P30822
C	?	-	VAL	deletion	UNP P30822
C	?	-	GLN	deletion	UNP P30822
C	?	-	ARG	deletion	UNP P30822
C	?	-	LEU	deletion	UNP P30822
C	?	-	PRO	deletion	UNP P30822
C	?	-	ALA	deletion	UNP P30822
C	?	-	THR	deletion	UNP P30822
C	?	-	GLU	deletion	UNP P30822
C	?	-	MET	deletion	UNP P30822
C	?	-	SER	deletion	UNP P30822
C	?	-	PRO	deletion	UNP P30822

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Chain	Residue	Modelled	Actual	Comment	Reference
C	?	-	LEU	deletion	UNP P30822
C	?	-	ILE	deletion	UNP P30822
C	?	-	GLN	deletion	UNP P30822
C	?	-	LEU	deletion	UNP P30822
C	?	-	SER	deletion	UNP P30822
C	?	-	VAL	deletion	UNP P30822
C	?	-	GLY	deletion	UNP P30822
C	?	-	SER	deletion	UNP P30822
C	?	-	GLN	deletion	UNP P30822
C	?	-	ALA	deletion	UNP P30822
C	?	-	ILE	deletion	UNP P30822
C	?	-	SER	deletion	UNP P30822
C	?	-	THR	deletion	UNP P30822
C	?	-	GLY	deletion	UNP P30822
C	?	-	SER	deletion	UNP P30822
C	?	-	GLY	deletion	UNP P30822
C	?	-	ALA	deletion	UNP P30822
C	?	-	LEU	deletion	UNP P30822
C	?	-	ASN	deletion	UNP P30822
C	?	-	PRO	deletion	UNP P30822
C	?	-	GLU	deletion	UNP P30822
C	?	-	TYR	deletion	UNP P30822
C	?	-	MET	deletion	UNP P30822
C	?	-	LYS	deletion	UNP P30822
C	?	-	ARG	deletion	UNP P30822
C	?	-	PHE	deletion	UNP P30822
C	537	GLY	ASP	engineered mutation	UNP P30822
C	539	CYS	THR	engineered mutation	UNP P30822
C	540	GLU	VAL	engineered mutation	UNP P30822
C	541	GLN	LYS	engineered mutation	UNP P30822
C	1022	CYS	TYR	conflict	UNP P30822

- Molecule 4 is PHOSPHOAMINOPHOSPHONIC ACID-GUANYLATE ESTER (CCD ID: GNP) (formula: $C_{10}H_{17}N_6O_{13}P_3$).

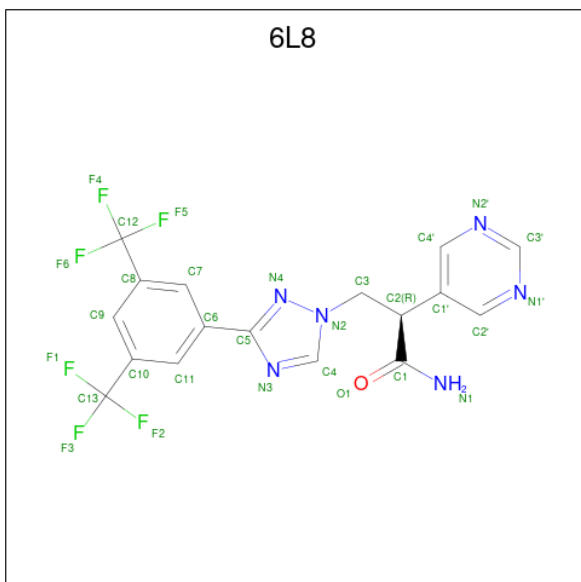


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	H	N	O	P	
			44	10	12	6	13	3	

- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Mg	0	0
			1	1		

- Molecule 6 is (2R)-3-{3-[3,5-bis(trifluoromethyl)phenyl]-1H-1,2,4-triazol-1-yl}-2-(pyrimidin-5-yl)propanamide (CCD ID: 6L8) (formula: C₁₇H₁₂F₆N₆O) (labeled as "Ligand of Interest" by depositor).

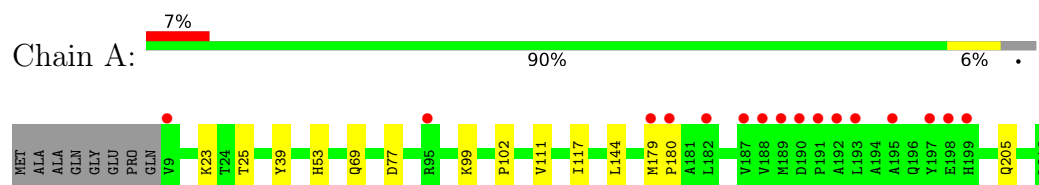


Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
6	C	1	Total	C	F	H	N	O	0	0
			42	17	6	12	6	1		

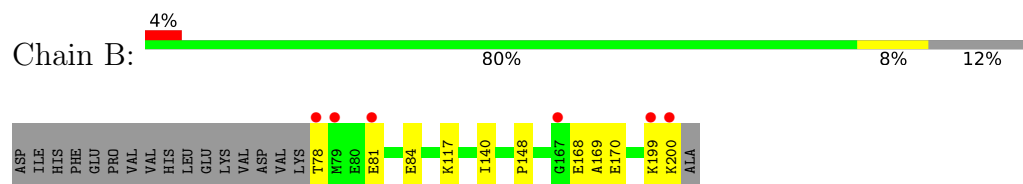
3 Residue-property plots [i](#)

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.

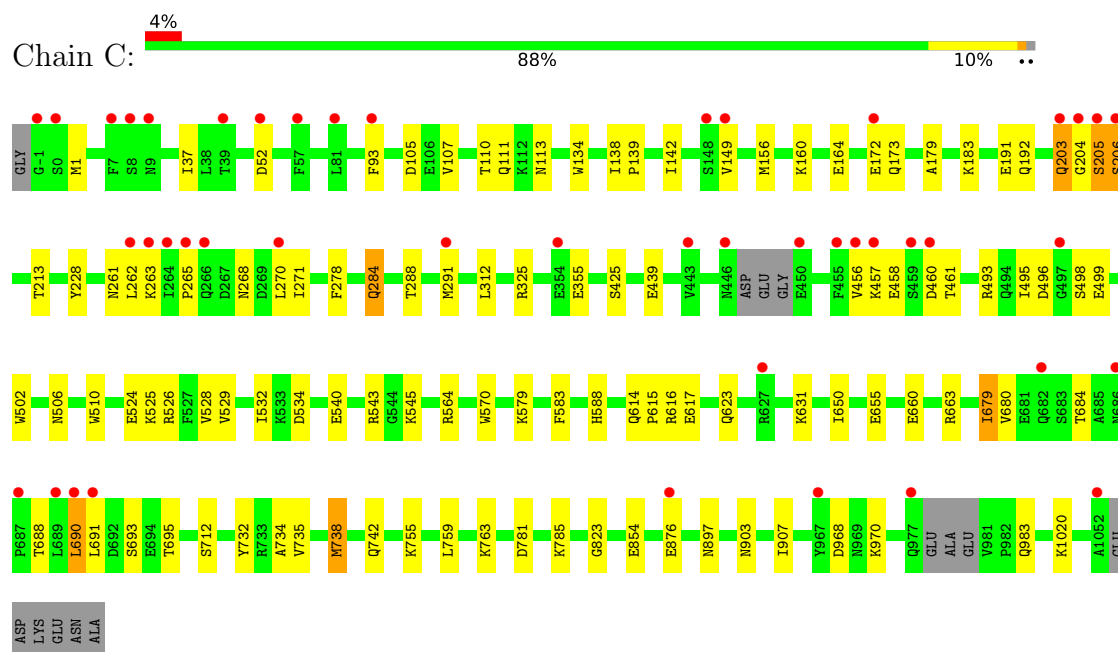
- Molecule 1: GTP-binding nuclear protein Ran



- Molecule 2: Ran-specific GTPase-activating protein 1



- Molecule 3: Exportin-1



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	106.67Å 106.67Å 306.07Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.55 – 2.41 45.55 – 2.41	Depositor EDS
% Data completeness (in resolution range)	97.2 (45.55-2.41) 97.2 (45.55-2.41)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.91 (at 2.39Å)	Xtriage
Refinement program	PHENIX 1.11.1_2575	Depositor
R, R_{free}	0.194 , 0.246 0.191 , 0.241	Depositor DCC
R_{free} test set	2000 reflections (2.96%)	wwPDB-VP
Wilson B-factor (Å ²)	35.1	Xtriage
Anisotropy	0.026	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 39.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	22274	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.54% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MG, GNP, 6L8

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	0.67	0/1754	0.74	0/2379
2	B	0.52	0/1059	0.63	0/1413
3	C	0.62	0/8608	0.71	0/11659
All	All	0.62	0/11421	0.71	0/15451

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
3	C	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	C	203	GLN	Peptide

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	1694	1697	1676	13	0
2	B	1029	1031	1021	9	0
3	C	8310	8426	8258	60	1
4	A	32	12	12	3	0
5	A	1	0	0	0	0
6	C	30	12	0	0	0
All	All	11096	11178	10967	77	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 3.

All (77) 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)
3:C:691:LEU:O	3:C:755:LYS:NZ	2.07	0.87
3:C:460:ASP:OD1	3:C:460:ASP:N	2.17	0.74
3:C:439:GLU:OE1	3:C:457:LYS:NZ	2.16	0.67
3:C:205[A]:SER:OG	3:C:206[A]:SER:N	2.22	0.64
3:C:690:LEU:HD13	3:C:738:MET:HE1	1.84	0.60
3:C:525:LYS:O	3:C:529:VAL:HG23	2.01	0.59
3:C:680:VAL:O	3:C:684:THR:HG23	2.05	0.57
3:C:263:LYS:HD2	3:C:325:ARG:HD2	1.87	0.56
3:C:528:VAL:O	3:C:532:ILE:HG12	2.07	0.55
1:A:23:LYS:NZ	4:A:301:GNP:O2G	2.32	0.54
3:C:268:ASN:HB3	3:C:271:ILE:HD12	1.90	0.54
3:C:496:ASP:C	3:C:498:SER:H	2.15	0.53
1:A:180:PRO:HA	2:B:78:THR:O	2.08	0.53
1:A:111:VAL:O	3:C:113[B]:ASN:ND2	2.43	0.53
3:C:52:ASP:HB2	3:C:93:PHE:CE2	2.44	0.52
3:C:781:ASP:O	3:C:785:LYS:HE2	2.08	0.52
3:C:179:ALA:O	3:C:183:LYS:HG3	2.10	0.52
3:C:570:TRP:CE2	3:C:614:GLN:HG3	2.45	0.51
1:A:111:VAL:HA	3:C:113[B]:ASN:OD1	2.11	0.50
3:C:614:GLN:HB2	3:C:617:GLU:HG3	1.94	0.49
3:C:579:LYS:HE3	3:C:583:PHE:CE1	2.48	0.49
2:B:199:LYS:O	2:B:200:LYS:HG3	2.12	0.49
3:C:142:ILE:CD1	3:C:192[B]:GLN:HB3	2.44	0.48
3:C:688:THR:O	3:C:691:LEU:HG	2.14	0.48
3:C:160:LYS:O	3:C:164:GLU:HG3	2.14	0.48
1:A:77:ASP:OD1	1:A:77:ASP:N	2.47	0.48
3:C:545:LYS:HE3	3:C:588:HIS:HB2	1.94	0.48
3:C:506:ASN:HB3	3:C:510:TRP:CZ2	2.49	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:897:ASN:OD1	3:C:897:ASN:C	2.56	0.48
3:C:735[B]:VAL:HG13	3:C:759:LEU:HB3	1.96	0.47
3:C:496:ASP:C	3:C:498:SER:N	2.72	0.47
3:C:655:GLU:HG2	3:C:660:GLU:HG3	1.95	0.47
3:C:156:MET:HE3	3:C:213:THR:HG23	1.97	0.47
3:C:205[A]:SER:HG	3:C:206[A]:SER:H	1.59	0.47
3:C:205[B]:SER:O	3:C:206[B]:SER:OG	2.23	0.47
3:C:495:ILE:HD11	3:C:534:ASP:HB3	1.97	0.47
1:A:117:ILE:HB	1:A:144:LEU:HD22	1.96	0.46
3:C:261:ASN:O	3:C:325:ARG:NH2	2.44	0.46
1:A:39:TYR:CD1	1:A:39:TYR:C	2.91	0.46
3:C:524:GLU:CD	3:C:564:ARG:HH21	2.23	0.46
3:C:903:ASN:O	3:C:907:ILE:HG13	2.17	0.45
1:A:205:GLN:HG2	2:B:148:PRO:O	2.17	0.45
3:C:679:ILE:HD11	3:C:695:THR:HG23	1.99	0.45
3:C:854:GLU:H	3:C:854:GLU:CD	2.24	0.45
1:A:53:HIS:CD2	1:A:179:MET:HE3	2.52	0.45
3:C:1:MET:CB	3:C:37:ILE:HD13	2.47	0.44
1:A:25:THR:OG1	4:A:301:GNP:O2A	2.30	0.44
2:B:78:THR:HG21	2:B:81:GLU:OE1	2.17	0.44
3:C:138:ILE:HB	3:C:139:PRO:HD3	2.00	0.44
2:B:168:GLU:O	2:B:170:GLU:N	2.51	0.43
3:C:734:ALA:O	3:C:738:MET:HE3	2.19	0.43
3:C:278:PHE:CD1	3:C:278:PHE:C	2.96	0.43
3:C:191:GLU:OE1	3:C:228:TYR:OH	2.32	0.43
3:C:156:MET:HE3	3:C:213:THR:OG1	2.19	0.43
2:B:84:GLU:OE1	2:B:117:LYS:NZ	2.48	0.43
3:C:650:ILE:HG13	3:C:712:SER:HB2	2.00	0.43
3:C:284:GLN:O	3:C:288:THR:HG23	2.20	0.42
1:A:99:LYS:O	1:A:102:PRO:HD2	2.19	0.42
3:C:268:ASN:OD1	3:C:268:ASN:C	2.62	0.42
3:C:732:TYR:CZ	3:C:763:LYS:HE3	2.55	0.42
1:A:23:LYS:NZ	4:A:301:GNP:O2B	2.44	0.42
3:C:291[B]:MET:HE1	3:C:312:LEU:HD21	2.01	0.42
3:C:204[B]:GLY:O	3:C:205[B]:SER:OG	2.32	0.41
3:C:493:ARG:HE	3:C:499:GLU:CD	2.28	0.41
3:C:615:PRO:O	3:C:616:ARG:HB2	2.20	0.41
3:C:968:ASP:HB2	3:C:970:LYS:HG3	2.02	0.41
1:A:180:PRO:HA	2:B:78:THR:C	2.46	0.41
3:C:107:VAL:HG13	3:C:111:GLN:HG3	2.03	0.41
3:C:876:GLU:CD	3:C:876:GLU:C	2.89	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:983:GLN:CD	3:C:983:GLN:C	2.88	0.41
2:B:140:ILE:HG22	2:B:140:ILE:O	2.21	0.41
3:C:172:GLU:HG2	3:C:173:GLN:HG3	2.03	0.41
2:B:117:LYS:HA	2:B:117:LYS:HE2	2.02	0.40
3:C:268:ASN:OD1	3:C:270:LEU:N	2.50	0.40
3:C:691:LEU:O	3:C:755:LYS:CE	2.69	0.40

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:660:GLU:OE2	3:C:663:ARG:HH22[7_555]	1.48	0.12

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	211/216 (98%)	203 (96%)	8 (4%)	0	100	100
2	B	124/140 (89%)	116 (94%)	7 (6%)	1 (1%)	16	25
3	C	1047/1024 (102%)	1006 (96%)	33 (3%)	8 (1%)	16	25
All	All	1382/1380 (100%)	1325 (96%)	48 (4%)	9 (1%)	24	28

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	205[A]	SER
3	C	205[B]	SER
3	C	206[A]	SER
3	C	206[B]	SER
3	C	458	GLU
2	B	169	ALA

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Mol	Chain	Res	Type
3	C	355	GLU
3	C	134	TRP
3	C	265	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	185/185 (100%)	183 (99%)	2 (1%)	65	82
2	B	108/121 (89%)	108 (100%)	0	100	100
3	C	960/933 (103%)	940 (98%)	20 (2%)	47	69
All	All	1253/1239 (101%)	1231 (98%)	22 (2%)	53	73

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

Mol	Chain	Res	Type
1	A	69[A]	GLN
1	A	69[B]	GLN
3	C	105	ASP
3	C	110	THR
3	C	149	VAL
3	C	203	GLN
3	C	262	LEU
3	C	284	GLN
3	C	425	SER
3	C	456	VAL
3	C	461	THR
3	C	502	TRP
3	C	526	ARG
3	C	540	GLU
3	C	543	ARG
3	C	623	GLN
3	C	631	LYS
3	C	679	ILE
3	C	690	LEU

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Mol	Chain	Res	Type
3	C	693	SER
3	C	738	MET
3	C	742	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	62	ASN
1	A	103	ASN
1	A	105	HIS
1	A	143	ASN
1	A	205	GLN
3	C	31	GLN
3	C	33	GLN
3	C	35	GLN
3	C	92	ASN
3	C	313	GLN
3	C	463	GLN
3	C	609	HIS
3	C	642	HIS

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 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
6	6L8	C	1101	3	32,32,32	1.51	3 (9%)	43,48,48	2.22	12 (27%)
4	GNP	A	301	5	34,34,34	3.72	16 (47%)	47,54,54	1.66	8 (17%)

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
6	6L8	C	1101	3	-	4/28/28/28	0/3/3/3
4	GNP	A	301	5	-	5/18/38/38	0/3/3/3

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	301	GNP	C2-N2	8.81	1.54	1.34
4	A	301	GNP	O4'-C1'	8.54	1.61	1.42
4	A	301	GNP	PA-O3A	7.47	1.67	1.59
4	A	301	GNP	PB-O3A	7.08	1.67	1.59
4	A	301	GNP	C2'-C1'	-7.02	1.31	1.53
4	A	301	GNP	PB-O1B	5.83	1.55	1.46
4	A	301	GNP	O4'-C4'	-5.39	1.33	1.45
6	C	1101	6L8	C1-N1	5.25	1.45	1.32
4	A	301	GNP	PG-O1G	4.12	1.52	1.46
4	A	301	GNP	O2'-C2'	3.93	1.52	1.43
6	C	1101	6L8	C6-C5	3.61	1.53	1.47
4	A	301	GNP	C6-N1	-3.21	1.32	1.38
4	A	301	GNP	O3'-C3'	-2.88	1.35	1.43
4	A	301	GNP	PB-N3B	2.71	1.70	1.63
6	C	1101	6L8	O1-C1	-2.69	1.19	1.23
4	A	301	GNP	O6-C6	-2.61	1.18	1.23
4	A	301	GNP	PA-O5'	2.50	1.69	1.59
4	A	301	GNP	PG-N3B	2.46	1.69	1.63
4	A	301	GNP	PB-O2B	-2.03	1.51	1.56

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	C	1101	6L8	N2-C4-N3	-6.54	107.33	110.82
6	C	1101	6L8	C4-N3-C5	5.46	107.41	102.71
4	A	301	GNP	C5-C4-N3	-5.36	119.86	128.39
6	C	1101	6L8	C4'-N2'-C3'	5.01	121.72	115.80
4	A	301	GNP	C2-N3-C4	4.51	120.08	112.30
6	C	1101	6L8	C2'-N1'-C3'	4.14	120.68	115.80
6	C	1101	6L8	C1'-C4'-N2'	-3.21	119.03	124.05
6	C	1101	6L8	N4-C5-N3	-3.21	108.10	114.62
6	C	1101	6L8	N1'-C3'-N2'	-3.13	121.26	126.73
4	A	301	GNP	N9-C4-N3	3.05	132.05	125.95
4	A	301	GNP	N9-C8-N7	-2.84	108.13	113.40
4	A	301	GNP	C2-N1-C6	-2.77	120.08	125.11
6	C	1101	6L8	C1'-C2'-N1'	-2.62	119.95	124.05
4	A	301	GNP	C5-C6-N1	2.62	119.93	113.25
4	A	301	GNP	O6-C6-C5	-2.48	120.00	126.53
6	C	1101	6L8	F4-C12-C8	-2.39	107.78	112.90
6	C	1101	6L8	C2-C1-N1	2.30	119.15	116.75
6	C	1101	6L8	C2'-C1'-C4'	2.27	117.35	115.26
4	A	301	GNP	C8-N7-C5	2.06	107.93	104.26
6	C	1101	6L8	C3-N2-N4	2.05	125.45	120.75

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	301	GNP	PG-N3B-PB-O1B
4	A	301	GNP	PG-N3B-PB-O3A
6	C	1101	6L8	C2'-C1'-C2-C1
6	C	1101	6L8	C4'-C1'-C2-C1
6	C	1101	6L8	O1-C1-C2-C1'
4	A	301	GNP	C5'-O5'-PA-O3A
4	A	301	GNP	C5'-O5'-PA-O1A
4	A	301	GNP	C5'-O5'-PA-O2A
6	C	1101	6L8	N1-C1-C2-C1'

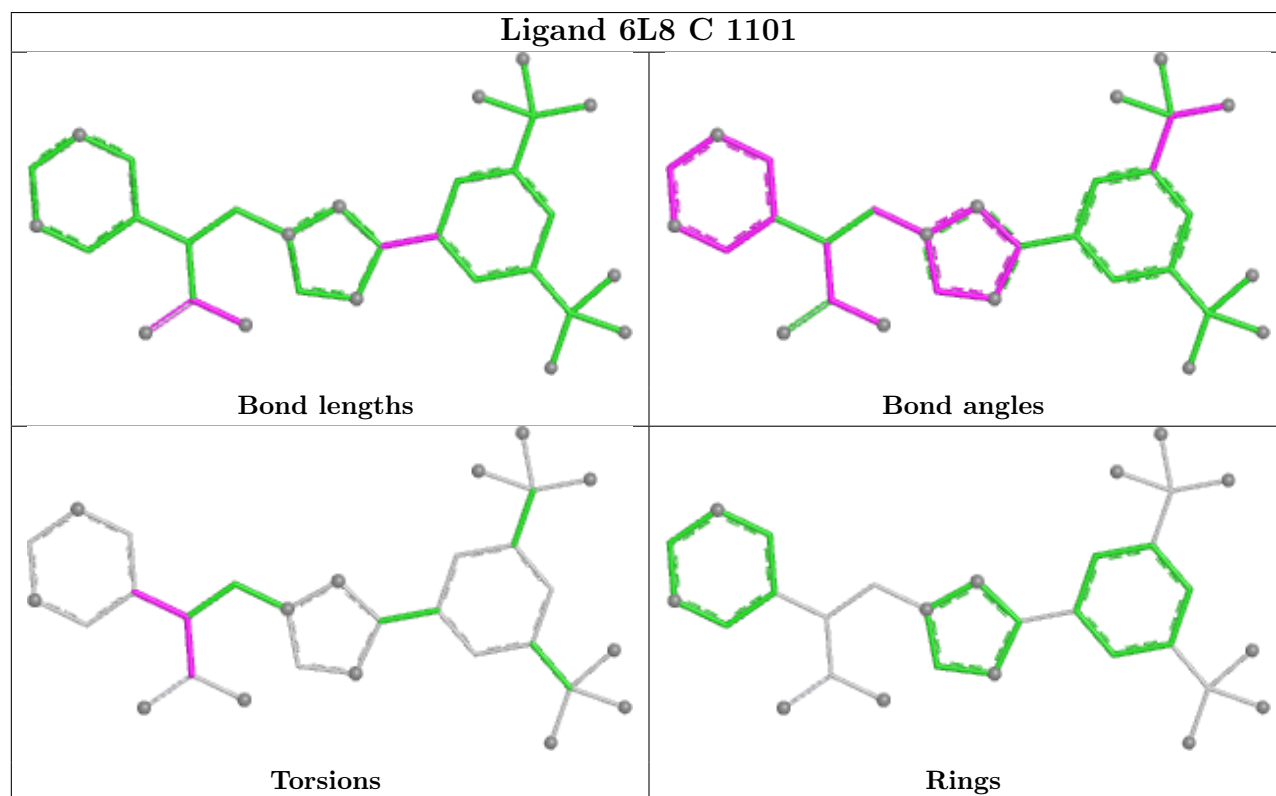
There are no ring outliers.

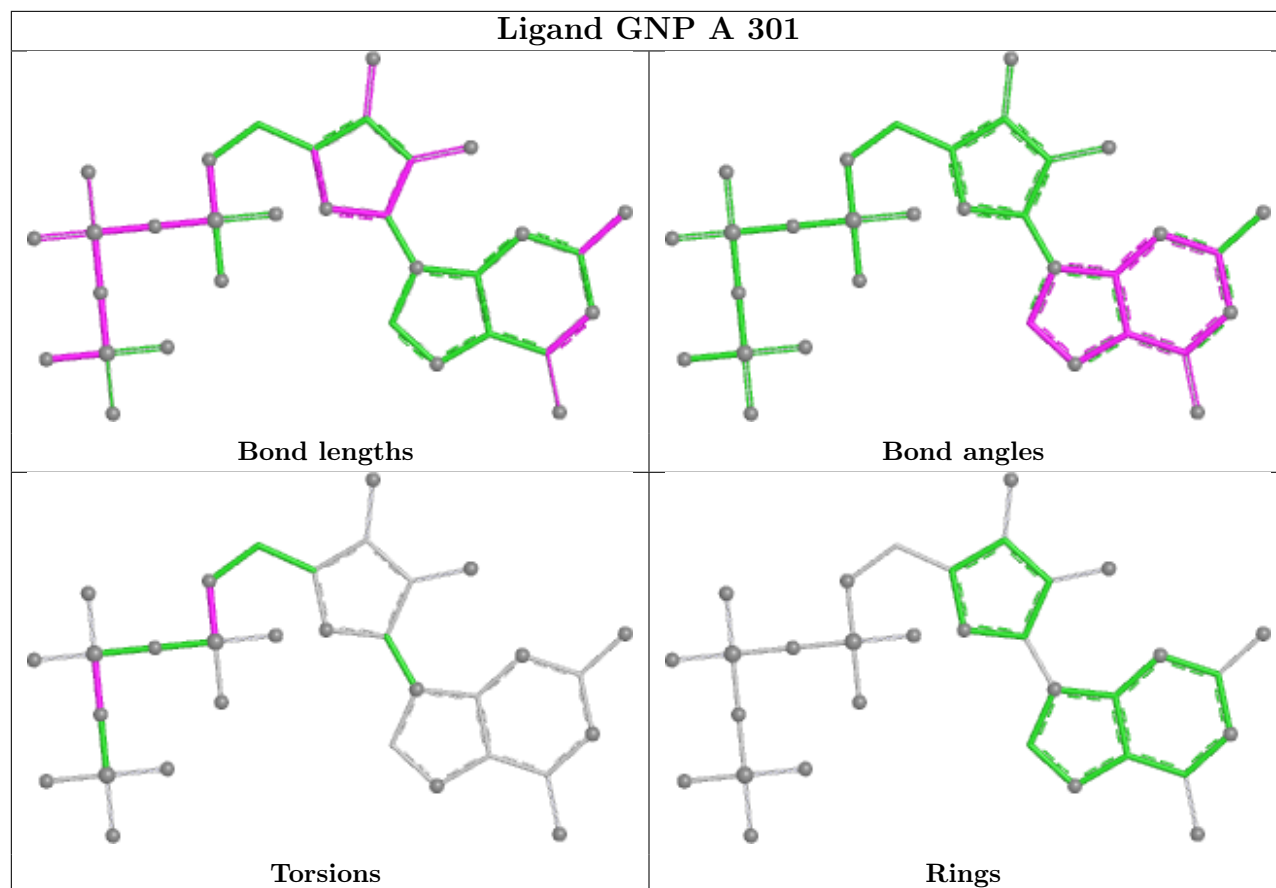
1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	301	GNP	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths,

bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

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, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	208/216 (96%)	-0.06	16 (7%) 19 16	18, 39, 111, 159	3 (1%)
2	B	123/140 (87%)	0.29	6 (4%) 35 31	19, 60, 93, 155	2 (1%)
3	C	1011/1024 (98%)	-0.08	45 (4%) 38 34	11, 43, 81, 143	25 (2%)
All	All	1342/1380 (97%)	-0.05	67 (4%) 34 30	11, 44, 86, 159	30 (2%)

All (67) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	C	205[A]	SER	9.9
3	C	446	ASN	8.8
1	A	191	PRO	6.3
3	C	457	LYS	5.9
3	C	206[A]	SER	5.8
1	A	193	LEU	5.2
3	C	455	PHE	5.1
2	B	79	MET	5.0
1	A	188	VAL	4.9
3	C	691	LEU	4.8
3	C	689	LEU	4.7
3	C	456	VAL	4.5
3	C	-1	GLY	4.4
2	B	78	THR	4.2
1	A	195	ALA	4.2
1	A	197	TYR	4.2
3	C	39[A]	THR	4.2
3	C	57	PHE	4.2
3	C	687	PRO	4.1
1	A	189	MET	4.0
3	C	264	ILE	3.9
3	C	459	SER	3.9
3	C	690	LEU	3.7

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Mol	Chain	Res	Type	RSRZ
3	C	81	LEU	3.6
3	C	977	GLN	3.5
3	C	204[A]	GLY	3.4
1	A	190	ASP	3.3
1	A	198	GLU	3.3
3	C	497	GLY	3.2
2	B	199	LYS	3.1
1	A	187	VAL	3.1
2	B	200	LYS	3.0
1	A	179	MET	2.9
3	C	266	GLN	2.9
3	C	149	VAL	2.8
2	B	81	GLU	2.8
3	C	627	ARG	2.8
3	C	52	ASP	2.8
3	C	93	PHE	2.8
1	A	9	VAL	2.7
3	C	450	GLU	2.6
3	C	686	ASN	2.6
3	C	203	GLN	2.6
3	C	265	PRO	2.5
3	C	9	ASN	2.5
3	C	270	LEU	2.5
3	C	443	VAL	2.4
3	C	876	GLU	2.4
3	C	460	ASP	2.4
3	C	967	TYR	2.4
1	A	182	LEU	2.4
3	C	1052	ALA	2.4
1	A	199	HIS	2.3
3	C	263	LYS	2.3
1	A	95	ARG	2.3
3	C	8	SER	2.3
3	C	172	GLU	2.3
2	B	167	GLY	2.3
3	C	354	GLU	2.2
1	A	192	ALA	2.2
3	C	291[A]	MET	2.2
3	C	682	GLN	2.2
3	C	148[A]	SER	2.1
3	C	7	PHE	2.1
1	A	180	PRO	2.1

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Mol	Chain	Res	Type	RSRZ
3	C	262	LEU	2.0
3	C	0	SER	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

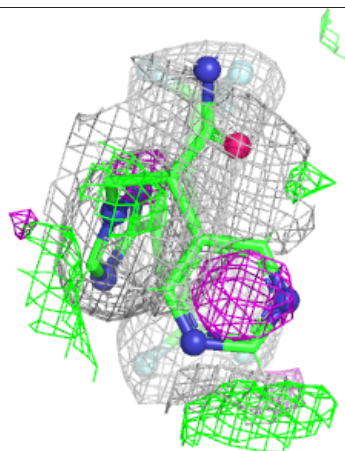
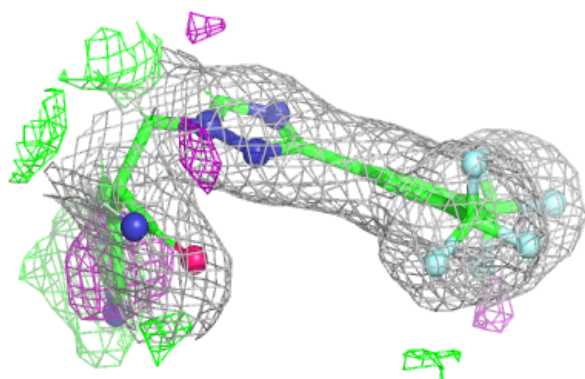
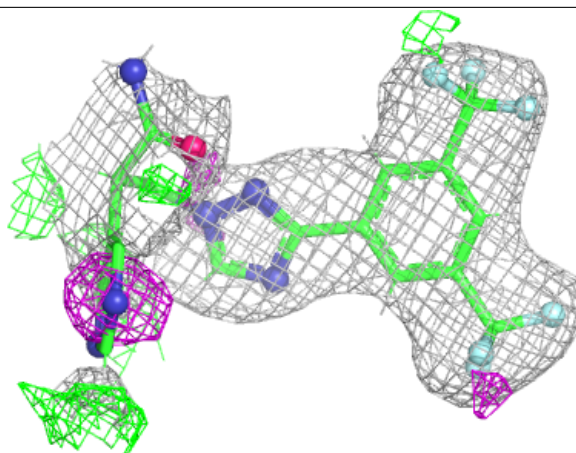
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	MG	A	302	1/1	0.79	0.12	25,25,25,25	0
6	6L8	C	1101	30/30	0.80	0.15	43,57,90,98	0
4	GNP	A	301	32/32	0.91	0.10	25,29,37,39	0

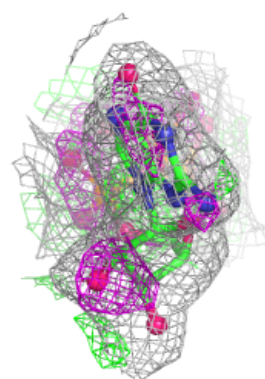
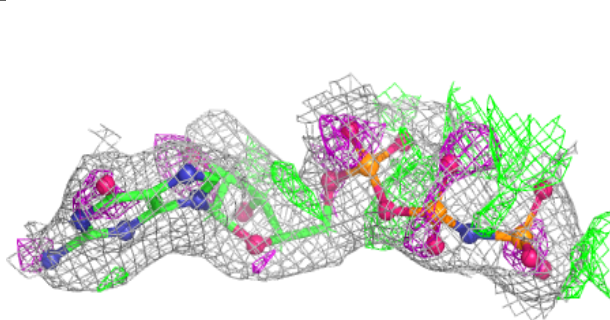
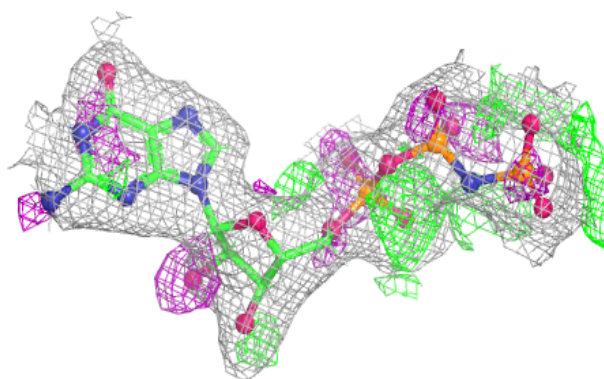
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around 6L8 C 1101:

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

**Electron density around GNP A 301:**

$2mF_o-DF_c$ (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.