



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 11:38 PM UTC

PDB ID : 7P97 / pdb_00007p97
Title : Structure of 3-phospho-D-glycerate guanylyltransferase with product 3-GPPG bound
Authors : Palm, G.J.; Berndt, L.; Lammers, M.
Deposited on : 2021-07-26
Resolution : 2.35 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

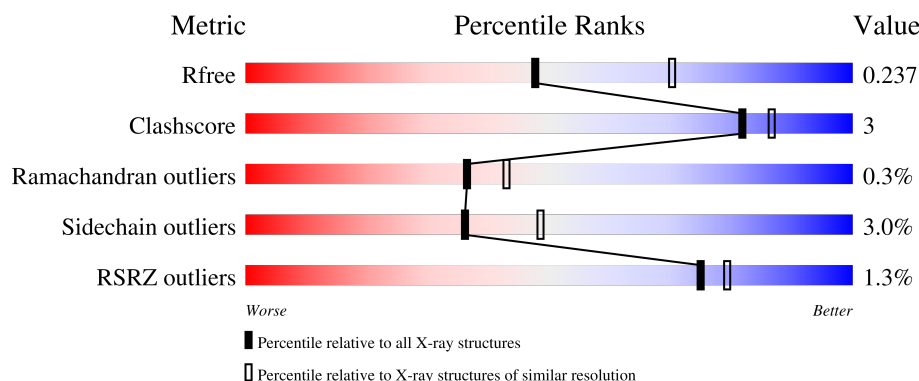
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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	1596 (2.36-2.36)
Clashscore	190562	1663 (2.36-2.36)
Ramachandran outliers	187476	1646 (2.36-2.36)
Sidechain outliers	187428	1646 (2.36-2.36)
RSRZ outliers	180081	1598 (2.36-2.36)

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	AAA	243	% 74% 9% • 16%
1	BBB	243	% 77% • 20%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6206 atoms, of which 2987 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 3-phospho-D-glycerate guanylyltransferase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	AAA	203	Total	C	H	N	O	S	35	1	0
			3022	946	1514	277	273	12			
1	BBB	195	Total	C	H	N	O	S	34	1	0
			2884	907	1441	262	263	11			

There are 68 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
AAA	-26	MET	-	initiating methionine	UNP E5ASS2
AAA	-25	GLY	-	expression tag	UNP E5ASS2
AAA	-24	SER	-	expression tag	UNP E5ASS2
AAA	-23	SER	-	expression tag	UNP E5ASS2
AAA	-22	HIS	-	expression tag	UNP E5ASS2
AAA	-21	HIS	-	expression tag	UNP E5ASS2
AAA	-20	HIS	-	expression tag	UNP E5ASS2
AAA	-19	HIS	-	expression tag	UNP E5ASS2
AAA	-18	HIS	-	expression tag	UNP E5ASS2
AAA	-17	HIS	-	expression tag	UNP E5ASS2
AAA	-16	SER	-	expression tag	UNP E5ASS2
AAA	-15	GLN	-	expression tag	UNP E5ASS2
AAA	-14	ASP	-	expression tag	UNP E5ASS2
AAA	-13	PRO	-	expression tag	UNP E5ASS2
AAA	-12	ASN	-	expression tag	UNP E5ASS2
AAA	-11	SER	-	expression tag	UNP E5ASS2
AAA	-10	SER	-	expression tag	UNP E5ASS2
AAA	-9	SER	-	expression tag	UNP E5ASS2
AAA	-8	ALA	-	expression tag	UNP E5ASS2
AAA	-7	ARG	-	expression tag	UNP E5ASS2
AAA	-6	LEU	-	expression tag	UNP E5ASS2
AAA	-5	GLN	-	expression tag	UNP E5ASS2
AAA	-4	VAL	-	expression tag	UNP E5ASS2
AAA	-3	ASP	-	expression tag	UNP E5ASS2
AAA	-2	LYS	-	expression tag	UNP E5ASS2

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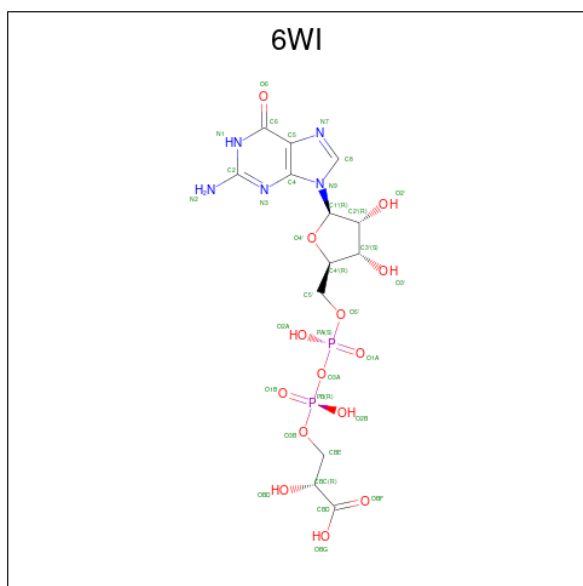
Chain	Residue	Modelled	Actual	Comment	Reference
AAA	-1	LEU	-	expression tag	UNP E5ASS2
AAA	0	ALA	-	expression tag	UNP E5ASS2
AAA	30	ALA	SER	variant	UNP E5ASS2
AAA	122	GLU	GLY	variant	UNP E5ASS2
AAA	175	ARG	CYS	variant	UNP E5ASS2
AAA	182	VAL	ILE	variant	UNP E5ASS2
AAA	205	VAL	ILE	variant	UNP E5ASS2
AAA	206	ARG	LYS	variant	UNP E5ASS2
AAA	216	ASN	-	expression tag	UNP E5ASS2
BBB	-26	MET	-	initiating methionine	UNP E5ASS2
BBB	-25	GLY	-	expression tag	UNP E5ASS2
BBB	-24	SER	-	expression tag	UNP E5ASS2
BBB	-23	SER	-	expression tag	UNP E5ASS2
BBB	-22	HIS	-	expression tag	UNP E5ASS2
BBB	-21	HIS	-	expression tag	UNP E5ASS2
BBB	-20	HIS	-	expression tag	UNP E5ASS2
BBB	-19	HIS	-	expression tag	UNP E5ASS2
BBB	-18	HIS	-	expression tag	UNP E5ASS2
BBB	-17	HIS	-	expression tag	UNP E5ASS2
BBB	-16	SER	-	expression tag	UNP E5ASS2
BBB	-15	GLN	-	expression tag	UNP E5ASS2
BBB	-14	ASP	-	expression tag	UNP E5ASS2
BBB	-13	PRO	-	expression tag	UNP E5ASS2
BBB	-12	ASN	-	expression tag	UNP E5ASS2
BBB	-11	SER	-	expression tag	UNP E5ASS2
BBB	-10	SER	-	expression tag	UNP E5ASS2
BBB	-9	SER	-	expression tag	UNP E5ASS2
BBB	-8	ALA	-	expression tag	UNP E5ASS2
BBB	-7	ARG	-	expression tag	UNP E5ASS2
BBB	-6	LEU	-	expression tag	UNP E5ASS2
BBB	-5	GLN	-	expression tag	UNP E5ASS2
BBB	-4	VAL	-	expression tag	UNP E5ASS2
BBB	-3	ASP	-	expression tag	UNP E5ASS2
BBB	-2	LYS	-	expression tag	UNP E5ASS2
BBB	-1	LEU	-	expression tag	UNP E5ASS2
BBB	0	ALA	-	expression tag	UNP E5ASS2
BBB	30	ALA	SER	variant	UNP E5ASS2
BBB	122	GLU	GLY	variant	UNP E5ASS2
BBB	175	ARG	CYS	variant	UNP E5ASS2
BBB	182	VAL	ILE	variant	UNP E5ASS2
BBB	205	VAL	ILE	variant	UNP E5ASS2
BBB	206	ARG	LYS	variant	UNP E5ASS2

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Chain	Residue	Modelled	Actual	Comment	Reference
BBB	216	ASN	-	expression tag	UNP E5ASS2

- Molecule 2 is 3-(guanosine-5'-diphospho)-D-glycerate (CCD ID: 6WI) (formula: $C_{13}H_{19}N_5O_{14}P_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	AAA	1	Total	C	H	N	O	P	3	0
			50	13	16	5	14	2		
2	BBB	1	Total	C	H	N	O	P	3	0
			50	13	16	5	14	2		

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	AAA	2	Total	Mg	0	0
			2	2		
3	BBB	2	Total	Mg	0	0
			2	2		

- Molecule 4 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

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

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	BBB	2	Total	Cl	0	0
			2	2		

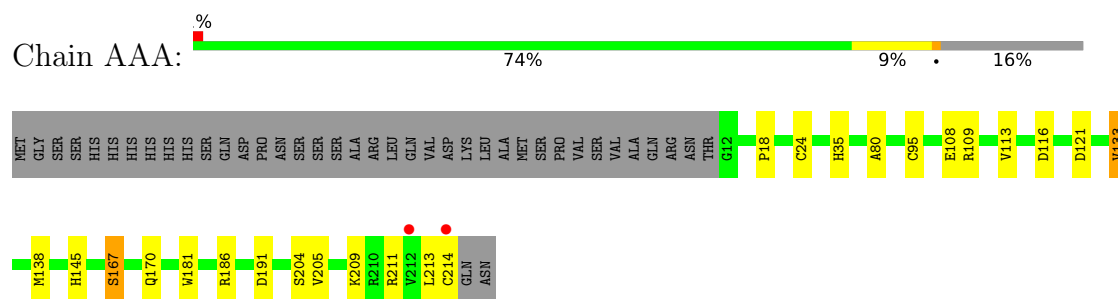
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	AAA	99	Total	O	0	0
			99	99		
5	BBB	94	Total	O	0	0
			94	94		

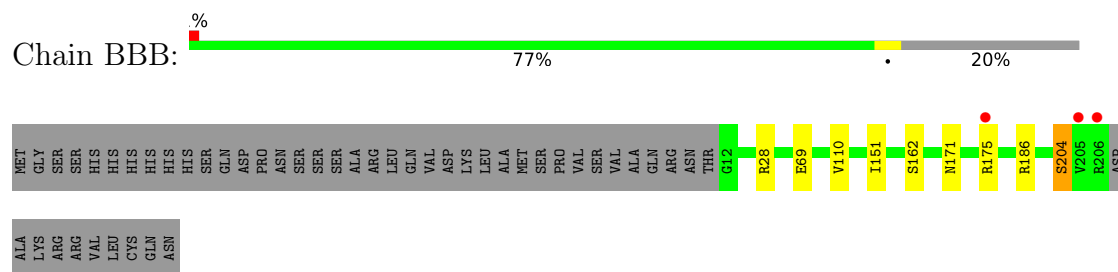
3 Residue-property plots [i](#)

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: 3-phospho-D-glycerate guanylyltransferase



- Molecule 1: 3-phospho-D-glycerate guanylyltransferase



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	57.09Å 57.09Å 228.97Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.49 – 2.35 49.49 – 2.35	Depositor EDS
% Data completeness (in resolution range)	99.8 (49.49-2.35) 99.9 (49.49-2.35)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.47	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.58 (at 2.34Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.186 , 0.238 0.187 , 0.237	Depositor DCC
R_{free} test set	878 reflections (4.61%)	wwPDB-VP
Wilson B-factor (Å ²)	24.2	Xtriage
Anisotropy	0.181	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.42 , 43.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.038 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6206	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.31% 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: 6WI, MG, CL

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	AAA	1.21	3/1544 (0.2%)	1.54	1/2099 (0.0%)
1	BBB	1.16	0/1479	1.58	1/2013 (0.0%)
All	All	1.19	3/3023 (0.1%)	1.56	2/4112 (0.0%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	AAA	145	HIS	CE1-NE2	5.72	1.38	1.32
1	AAA	80	ALA	C-O	-5.23	1.17	1.23
1	AAA	113	VAL	N-CA	5.09	1.49	1.45

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	BBB	69	GLU	CB-CG-CD	6.15	123.06	112.60
1	AAA	121	ASP	CA-CB-CG	5.30	117.90	112.60

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	AAA	1508	1514	1509	9	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	BBB	1443	1441	1436	6	0
2	AAA	34	16	0	0	0
2	BBB	34	16	0	0	0
3	AAA	2	0	0	0	0
3	BBB	2	0	0	0	0
4	AAA	1	0	0	0	0
4	BBB	2	0	0	0	0
5	AAA	99	0	0	2	0
5	BBB	94	0	0	1	0
All	All	3219	2987	2945	15	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 (15) 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:BBB:186:ARG:HH21	1:BBB:204:SER:HB2	1.61	0.65
1:BBB:171:ASN:HB3	1:BBB:175:ARG:NH1	2.13	0.64
1:BBB:171:ASN:HB3	1:BBB:175:ARG:HH12	1.74	0.52
1:AAA:181:TRP:HB3	5:AAA:483:HOH:O	2.10	0.50
1:BBB:28:ARG:HB2	5:BBB:433:HOH:O	2.12	0.50
1:AAA:109:ARG:HD2	1:AAA:133:VAL:HG12	1.93	0.50
1:BBB:186:ARG:HH11	1:BBB:186:ARG:HG3	1.78	0.49
1:AAA:213:LEU:O	1:AAA:214:CYS:SG	2.71	0.48
1:AAA:18:PRO:HG3	1:AAA:95:CYS:SG	2.55	0.47
1:AAA:209:LYS:HA	1:AAA:213:LEU:HD12	1.97	0.46
1:AAA:116:ASP:O	1:AAA:191:ASP:HA	2.19	0.42
1:AAA:167:SER:HB2	5:AAA:417:HOH:O	2.19	0.42
1:AAA:186:ARG:NH2	1:AAA:204:SER:HB3	2.35	0.41
1:BBB:110:VAL:O	1:BBB:151:ILE:HA	2.21	0.41
1:AAA:138:MET:HE3	1:AAA:138:MET:HB2	1.93	0.40

There are no symmetry-related clashes.

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	AAA	202/243 (83%)	198 (98%)	4 (2%)	0	100	100
1	BBB	194/243 (80%)	189 (97%)	4 (2%)	1 (0%)	24	27
All	All	396/486 (82%)	387 (98%)	8 (2%)	1 (0%)	36	43

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	BBB	162	SER

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	AAA	153/188 (81%)	144 (94%)	9 (6%)	18	22
1	BBB	146/188 (78%)	145 (99%)	1 (1%)	76	86
All	All	299/376 (80%)	289 (97%)	10 (3%)	36	44

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	AAA	24[A]	CYS
1	AAA	24[B]	CYS
1	AAA	35	HIS
1	AAA	108	GLU
1	AAA	133	VAL

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Mol	Chain	Res	Type
1	AAA	167	SER
1	AAA	170	GLN
1	AAA	205	VAL
1	AAA	211	ARG
1	BBB	204	SER

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

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 7 are 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	$\# Z > 2$
2	6WI	BBB	301	3	35,36,36	1.84	7 (20%)	50,55,55	1.85	11 (22%)
2	6WI	AAA	301	3	35,36,36	1.18	5 (14%)	50,55,55	1.47	8 (16%)

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	6WI	BBB	301	3	-	2/26/42/42	0/3/3/3
2	6WI	AAA	301	3	-	2/26/42/42	0/3/3/3

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	BBB	301	6WI	PB-O3A	4.77	1.64	1.59
2	BBB	301	6WI	CBC-CBD	4.43	1.58	1.52
2	BBB	301	6WI	PA-O3A	3.81	1.63	1.59
2	BBB	301	6WI	C6-N1	3.39	1.45	1.38
2	AAA	301	6WI	C5-C6	-2.80	1.34	1.44
2	BBB	301	6WI	PA-O5'	2.55	1.69	1.59
2	AAA	301	6WI	O3'-C3'	2.44	1.49	1.43
2	AAA	301	6WI	OBF-CBD	2.34	1.29	1.22
2	BBB	301	6WI	PB-O1B	2.27	1.58	1.50
2	BBB	301	6WI	C5-N7	-2.23	1.34	1.39
2	AAA	301	6WI	PA-O3A	2.23	1.61	1.59
2	AAA	301	6WI	PA-O5'	2.06	1.67	1.59

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	BBB	301	6WI	C2-N3-C4	5.79	122.27	112.30
2	BBB	301	6WI	C5-C4-N3	-5.26	120.01	128.39
2	BBB	301	6WI	C2-N1-C6	-4.49	116.96	125.11
2	AAA	301	6WI	C2-N3-C4	4.02	119.22	112.30
2	AAA	301	6WI	C5-C4-N3	-3.71	122.49	128.39
2	BBB	301	6WI	O4'-C4'-C5'	-3.19	99.12	109.33
2	BBB	301	6WI	N9-C4-N3	2.84	131.63	125.95
2	AAA	301	6WI	OBF-CBD-CBC	-2.74	117.12	122.60
2	BBB	301	6WI	C4-C5-N7	-2.68	106.43	110.67
2	BBB	301	6WI	N9-C8-N7	-2.64	108.50	113.40
2	AAA	301	6WI	O2B-PB-O3A	2.56	114.18	107.27
2	BBB	301	6WI	OBF-CBD-CBC	-2.55	117.50	122.60
2	BBB	301	6WI	OBG-CBD-CBC	2.47	117.96	112.74
2	BBB	301	6WI	C8-N7-C5	2.46	108.64	104.26
2	BBB	301	6WI	C5-C6-N1	2.45	119.48	113.25
2	AAA	301	6WI	C5'-C4'-C3'	-2.30	106.92	115.21
2	AAA	301	6WI	O3A-PA-O1A	2.29	117.58	110.70
2	AAA	301	6WI	N9-C4-N3	2.10	130.15	125.95

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	AAA	301	6WI	N1-C2-N3	-2.10	119.48	123.32

There are no chirality outliers.

All (4) torsion outliers are listed below:

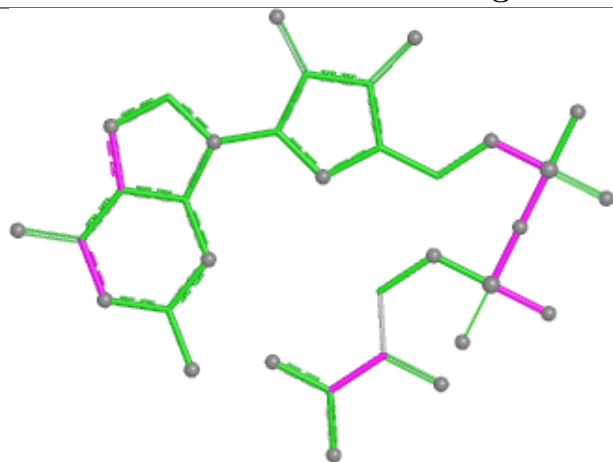
Mol	Chain	Res	Type	Atoms
2	AAA	301	6WI	CBE-O3B-PB-O1B
2	AAA	301	6WI	CBE-O3B-PB-O3A
2	BBB	301	6WI	CBE-O3B-PB-O1B
2	BBB	301	6WI	CBE-O3B-PB-O3A

There are no ring outliers.

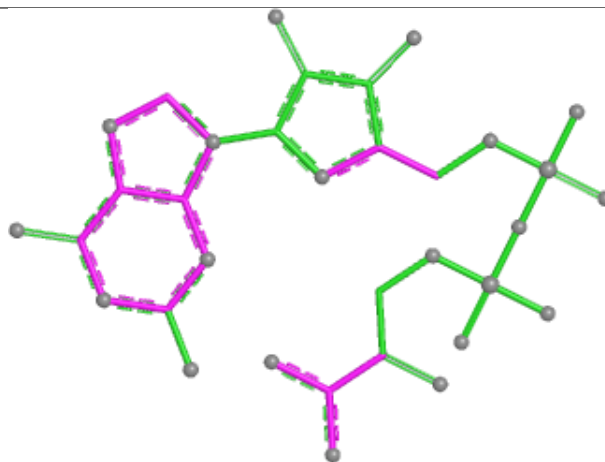
No monomer is involved in short contacts.

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.

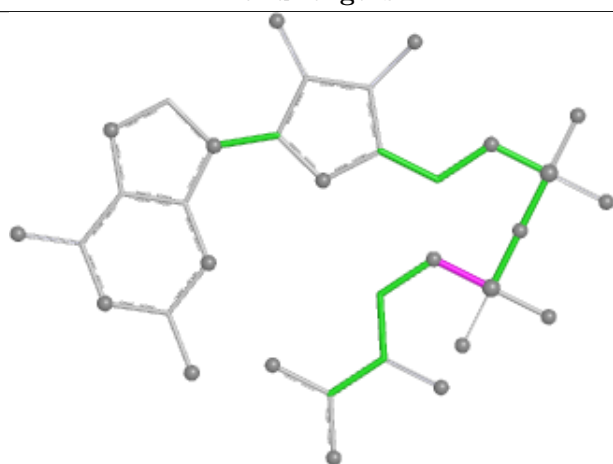
Ligand 6WI BBB 301



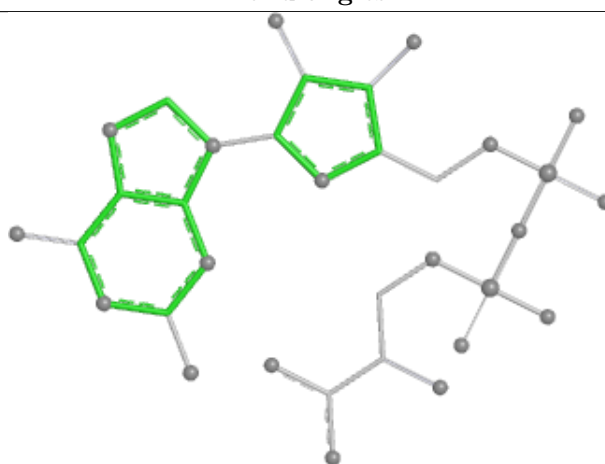
Bond lengths



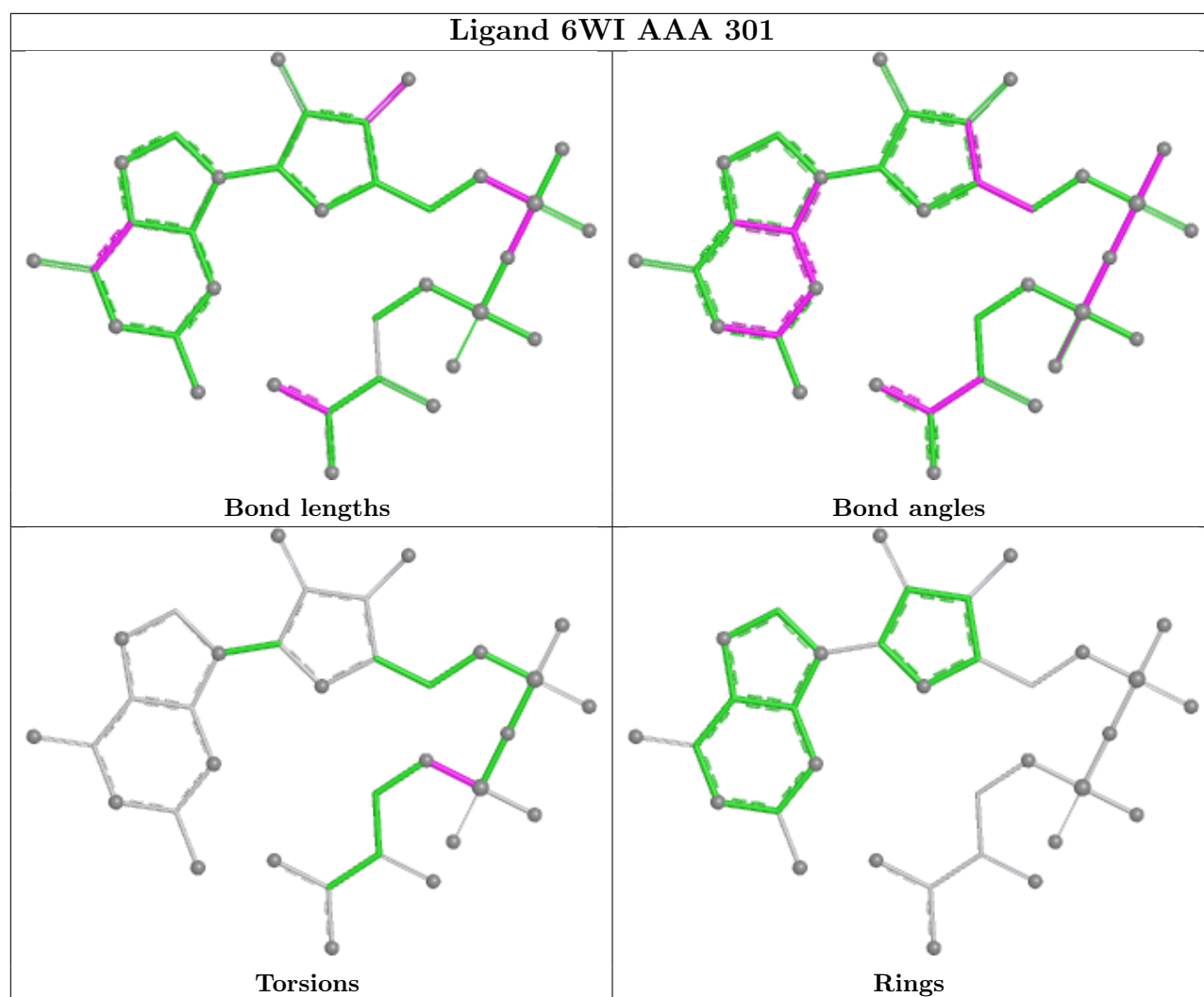
Bond angles



Torsions



Rings



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	AAA	203/243 (83%)	-0.40	2 (0%) 79 83	14, 25, 41, 75	1 (0%)
1	BBB	195/243 (80%)	-0.37	3 (1%) 72 77	16, 26, 45, 75	1 (0%)
All	All	398/486 (81%)	-0.38	5 (1%) 75 79	14, 26, 45, 75	2 (0%)

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	AAA	214	CYS	3.4
1	BBB	206	ARG	2.3
1	AAA	212	VAL	2.3
1	BBB	205	VAL	2.2
1	BBB	175	ARG	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](#)

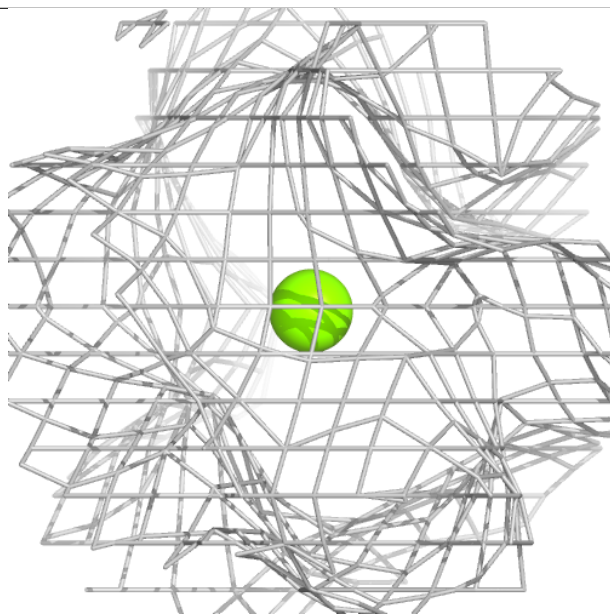
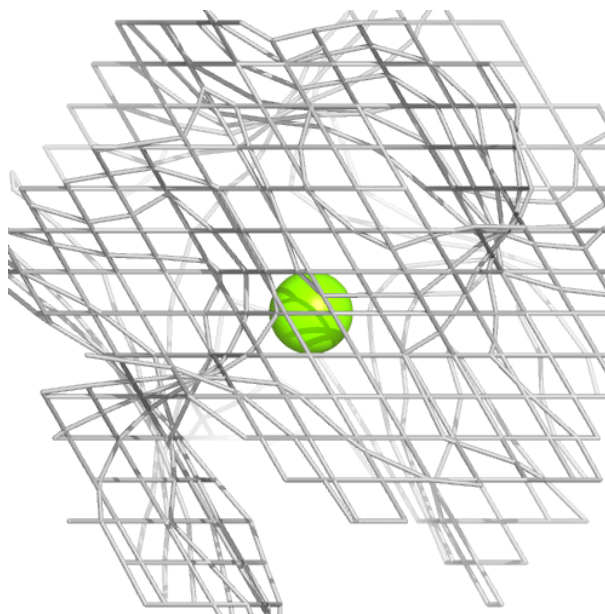
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists 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
4	CL	BBB	305	1/1	0.84	0.16	45,45,45,45	1
4	CL	BBB	304	1/1	0.97	0.05	28,28,28,28	0
3	MG	AAA	302	1/1	0.98	0.03	26,26,26,26	0
3	MG	AAA	303	1/1	0.99	0.03	22,22,22,22	0
3	MG	BBB	303	1/1	0.99	0.03	23,23,23,23	0
4	CL	AAA	304	1/1	0.99	0.03	26,26,26,26	0
2	6WI	BBB	301	34/34	0.99	0.04	15,18,21,22	3
2	6WI	AAA	301	34/34	0.99	0.04	13,15,18,19	3
3	MG	BBB	302	1/1	1.00	0.01	19,19,19,19	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.

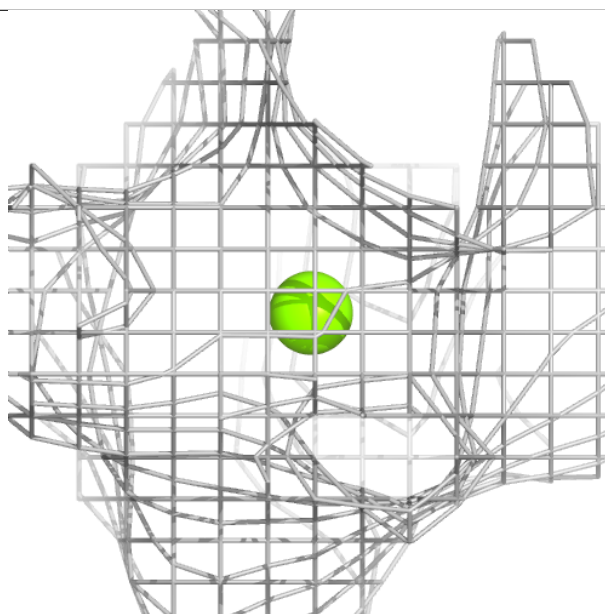
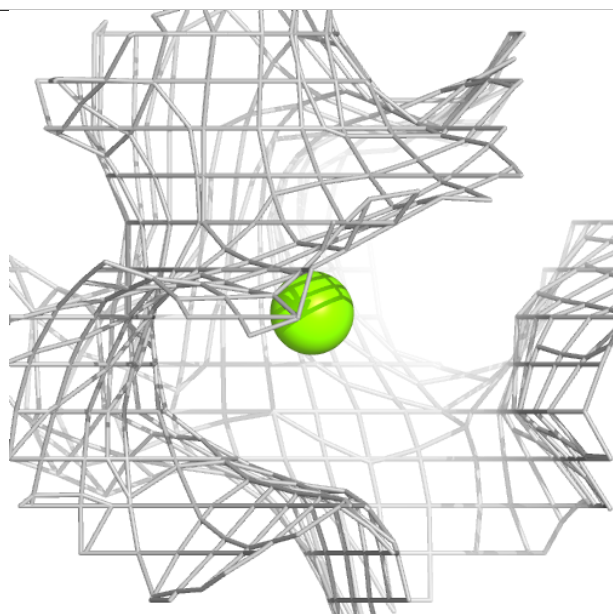
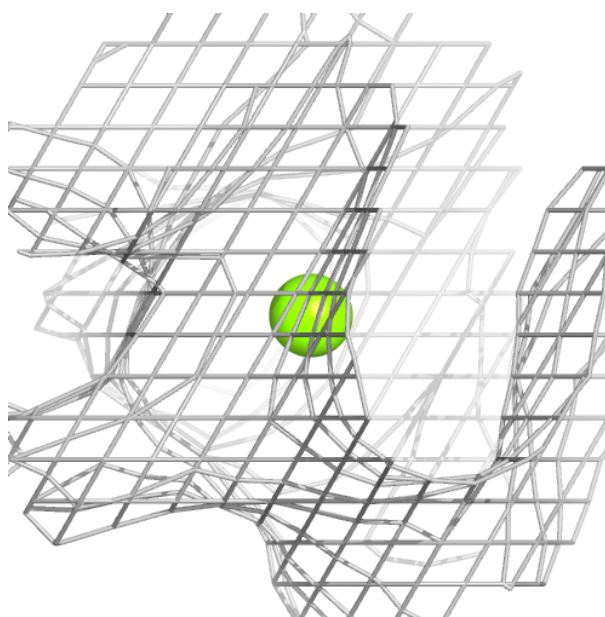
Electron density around MG AAA 302:

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



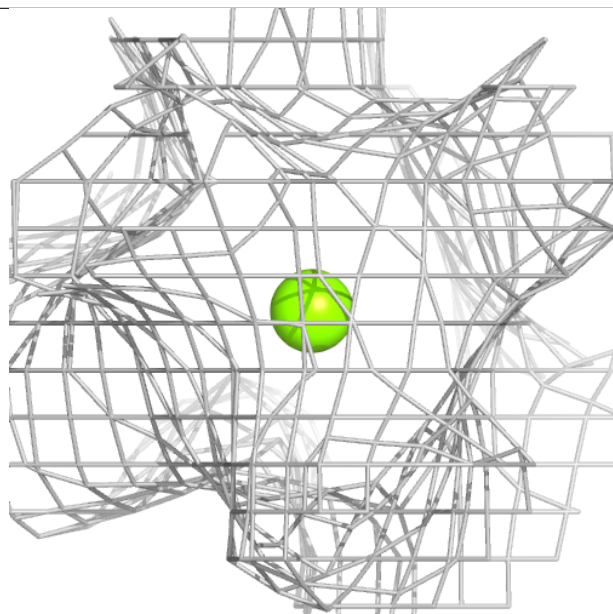
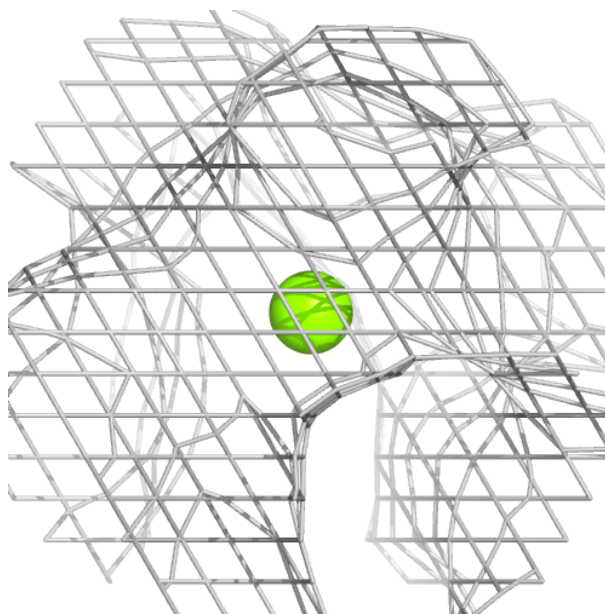
Electron density around MG AAA 303:

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



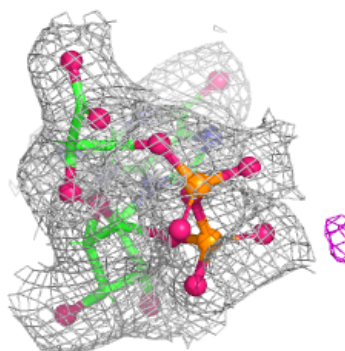
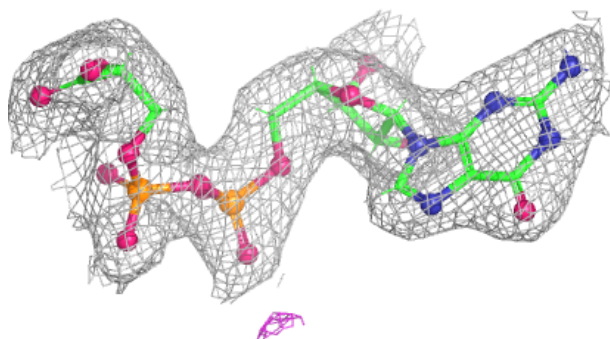
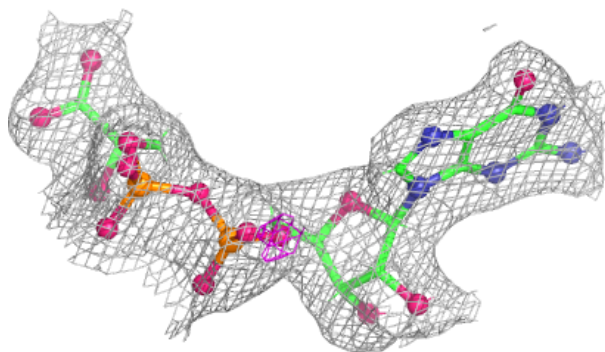
Electron density around MG BBB 303:

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

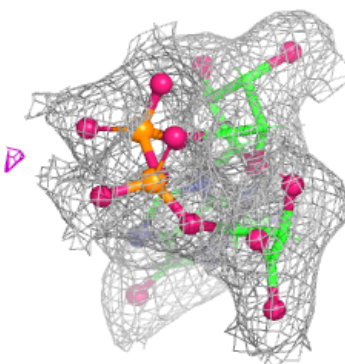
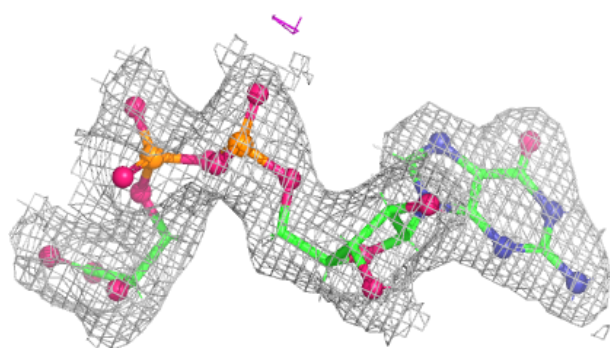
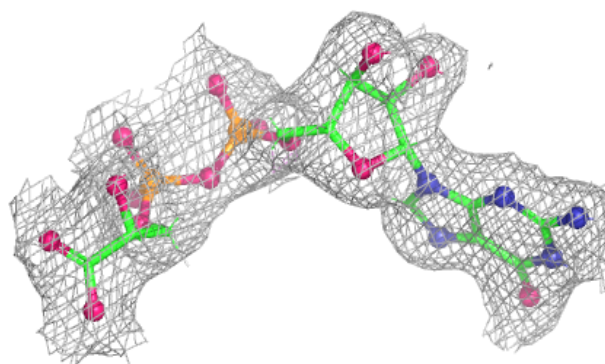


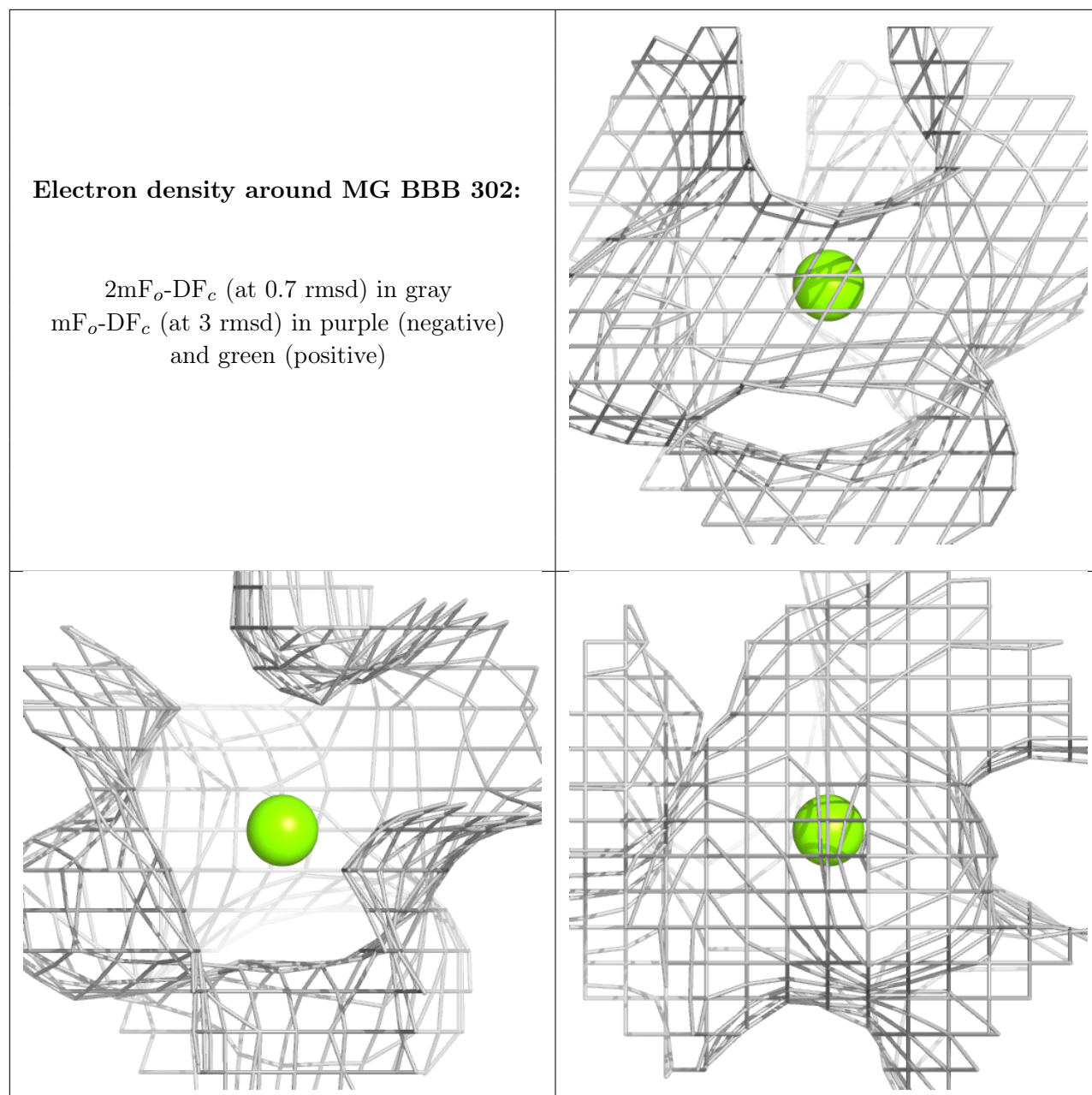
Electron density around 6WI BBB 301:

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

**Electron density around 6WI AAA 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 ⓘ

There are no such residues in this entry.