



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 20, 2026 – 11:29 am BST

PDB ID : 2J41 / pdb_00002j41
Title : Crystal structure of Staphylococcus aureus guanylate monophosphate kinase
Authors : El Omari, K.; Dhaliwal, B.; Lockyer, M.; Charles, I.; Hawkins, A.R.; Stammers, D.K.
Deposited on : 2006-08-24
Resolution : 1.90 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	:	1.8.4, CSD as541be (2020)
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

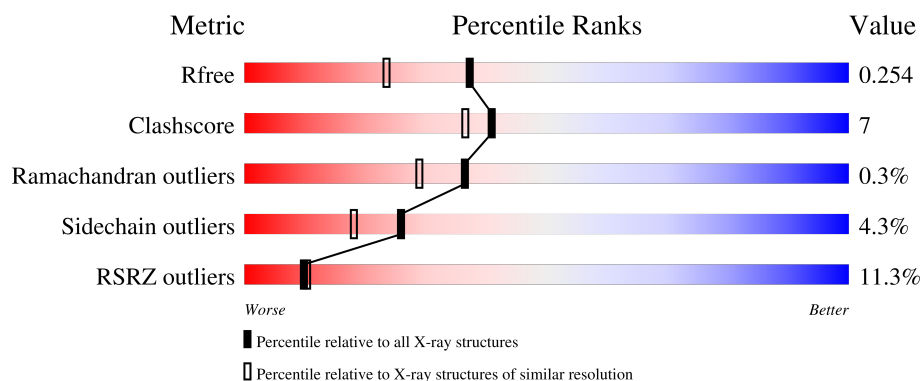
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	207	7% 65% 14% • 19%
1	B	207	11% 74% 8% • 16%
1	C	207	12% 74% 8% • 16%
1	D	207	8% 77% 7% • 14%

2 Entry composition [i](#)

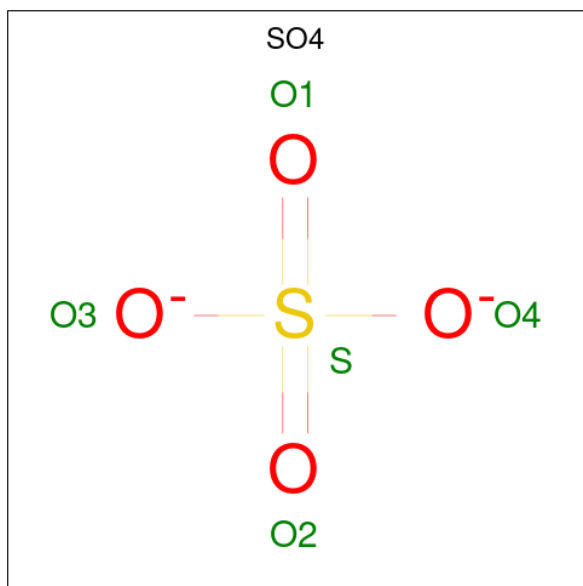
There are 5 unique types of molecules in this entry. The entry contains 6201 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 GUANYLATE KINASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	167	Total	C	N	O	S	0	1	0
			1363	873	221	263	6			
1	B	173	Total	C	N	O	S	0	0	0
			1406	900	230	269	7			
1	C	174	Total	C	N	O	S	0	0	0
			1412	898	236	272	6			
1	D	178	Total	C	N	O	S	0	0	0
			1449	926	240	276	7			

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

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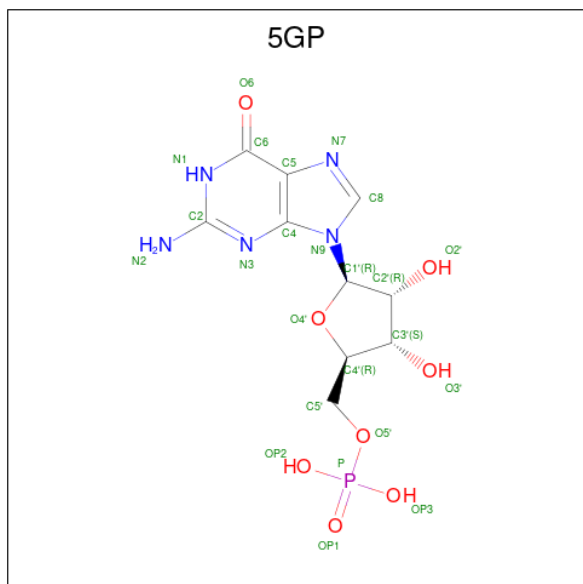
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is POTASSIUM ION (CCD ID: K) (formula: K).

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

- Molecule 4 is GUANOSINE-5'-MONOPHOSPHATE (CCD ID: 5GP) (formula: C₁₀H₁₄N₅O₈P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	B	1	Total	C	N	O	P	0	0
			24	10	5	8	1		
4	C	1	Total	C	N	O	P	0	0
			24	10	5	8	1		
4	D	1	Total	C	N	O	P	0	0
			24	10	5	8	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	95	Total 95	O 95	0	0
5	B	111	Total 111	O 111	0	0
5	C	130	Total 130	O 130	0	0
5	D	135	Total 135	O 135	0	0

- Molecule 1: GUANYLATE KINASE





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	70.01Å 93.99Å 83.94Å 90.00° 110.59° 90.00°	Depositor
Resolution (Å)	29.40 – 1.90 29.40 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.1 (29.40-1.90) 99.1 (29.40-1.90)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.00 (at 1.91Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.207 , 0.239 0.228 , 0.254	Depositor DCC
R_{free} test set	3991 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	28.8	Xtriage
Anisotropy	0.123	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 45.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	6201	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.88% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, K, 5GP

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	0.61	0/1392	0.83	1/1873 (0.1%)
1	B	0.72	1/1432 (0.1%)	0.81	0/1926
1	C	0.77	2/1437 (0.1%)	0.86	3/1932 (0.2%)
1	D	0.65	0/1475	0.83	1/1981 (0.1%)
All	All	0.69	3/5736 (0.1%)	0.83	5/7712 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	161	MET	SD-CE	15.68	2.18	1.79
1	B	161	MET	SD-CE	10.03	2.04	1.79
1	C	161	MET	CG-SD	6.18	1.96	1.80

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	59	THR	N-CA-CB	-6.36	101.02	110.17
1	C	161	MET	CG-SD-CE	-6.15	87.36	100.90
1	C	74	GLU	N-CA-CB	-5.59	101.12	110.57
1	C	74	GLU	CA-CB-CG	5.23	124.56	114.10
1	D	74	GLU	N-CA-CB	-5.16	101.85	110.57

There are no chirality outliers.

There are no planarity outliers.

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	1363	0	1341	26	0
1	B	1406	0	1388	19	0
1	C	1412	0	1391	20	0
1	D	1449	0	1438	12	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
2	C	5	0	0	0	0
2	D	10	0	0	1	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	B	24	0	12	2	0
4	C	24	0	12	0	0
4	D	24	0	12	0	0
5	A	95	0	0	4	0
5	B	111	0	0	1	0
5	C	130	0	0	5	0
5	D	135	0	0	4	0
All	All	6201	0	5594	74	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 7.

All (74) 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:B:161:MET:SD	1:B:161:MET:CE	2.04	1.44
1:C:161:MET:CE	1:C:161:MET:SD	2.18	1.32
1:B:199:ARG:HH21	1:B:199:ARG:HG3	1.15	1.10
1:B:161:MET:HE3	1:B:167:VAL:HG21	1.51	0.92
1:B:199:ARG:HH21	1:B:199:ARG:CG	1.89	0.84
1:A:59:THR:HG22	1:A:62:ALA:H	1.43	0.82
1:B:199:ARG:HG3	1:B:199:ARG:NH2	1.89	0.81
1:A:23:ARG:HD2	1:A:36[A]:TYR:CE1	2.23	0.74
1:C:181:GLN:HG3	5:C:2114:HOH:O	1.89	0.73
1:A:106:VAL:HG12	1:A:160:MET:HG2	1.72	0.70
1:C:161:MET:CE	1:C:161:MET:HB3	2.22	0.69
1:C:161:MET:HB3	1:C:161:MET:HE2	1.75	0.69
2:D:1207:SO4:O2	5:D:2127:HOH:O	2.10	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:B:1208:5GP:O3'	5:B:2111:HOH:O	2.12	0.68
1:C:161:MET:HE3	1:D:191:ARG:HG2	1.74	0.68
1:A:59:THR:CG2	1:A:62:ALA:H	2.07	0.67
1:C:191:ARG:HD3	1:D:161:MET:O	1.95	0.67
1:C:25:ARG:HG3	1:C:177:LYS:HD3	1.78	0.65
1:A:43:ARG:NH2	1:A:49:GLU:OE1	2.29	0.65
1:A:106:VAL:HG23	5:A:2062:HOH:O	1.97	0.63
1:D:60:ARG:HD2	5:D:2049:HOH:O	1.96	0.63
1:A:43:ARG:NH2	1:A:46:ARG:HG3	2.14	0.62
1:A:191:ARG:O	1:A:195:GLU:HB2	1.99	0.62
1:A:43:ARG:HH22	1:A:49:GLU:CD	2.11	0.58
1:C:113:ARG:HD2	5:C:2104:HOH:O	2.03	0.58
1:B:113:ARG:NH1	1:B:163:LEU:O	2.38	0.56
1:A:40:MET:HG2	1:A:56:PHE:HB2	1.88	0.56
1:C:44:GLN:H	1:C:44:GLN:CD	2.11	0.56
1:C:181:GLN:CG	5:C:2114:HOH:O	2.47	0.56
1:B:25:ARG:HG3	1:B:177:LYS:HD3	1.88	0.56
1:D:14:PRO:HD3	1:D:123:PHE:CZ	2.42	0.55
1:C:161:MET:CE	1:C:161:MET:CB	2.85	0.54
1:D:153:GLU:HB3	1:D:155:ARG:HH21	1.72	0.54
1:B:172:GLU:OE1	1:C:25:ARG:NH1	2.39	0.54
1:B:199:ARG:HD3	1:B:203:LEU:HD22	1.89	0.53
1:D:187:GLU:OE2	1:D:190:LYS:HE3	2.09	0.53
1:A:23:ARG:HD2	1:A:36[A]:TYR:HE1	1.74	0.53
1:C:161:MET:CE	1:C:161:MET:CG	2.86	0.53
1:A:157:GLU:C	1:A:159:GLU:H	2.16	0.52
1:A:158:VAL:HG12	1:A:161:MET:HE3	1.91	0.52
1:A:106:VAL:CG1	1:A:160:MET:HG2	2.40	0.52
1:A:128:SER:HA	5:A:2073:HOH:O	2.10	0.52
1:A:162:ASN:HB3	1:B:195:GLU:OE2	2.10	0.51
1:B:161:MET:CE	1:B:161:MET:HB3	2.40	0.51
1:D:47:GLU:HG3	5:D:2039:HOH:O	2.11	0.50
1:C:128:SER:HB3	5:C:2101:HOH:O	2.12	0.50
1:B:161:MET:CE	1:B:161:MET:CB	2.90	0.49
1:D:199:ARG:O	1:D:202:ILE:HD13	2.13	0.49
1:A:25:ARG:HG3	1:A:177:LYS:HD3	1.93	0.48
1:A:75:TYR:H	1:A:111:GLN:HE22	1.61	0.48
1:A:36[A]:TYR:CD1	5:A:2058:HOH:O	2.56	0.48
1:D:198:TYR:O	1:D:202:ILE:HG23	2.17	0.44
1:C:14:PRO:HD3	1:C:123:PHE:CZ	2.52	0.44
1:C:150:ARG:HA	5:C:2102:HOH:O	2.16	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:113:ARG:HD2	1:B:163:LEU:HD13	1.99	0.44
1:C:126:PRO:C	1:C:128:SER:H	2.25	0.43
1:B:161:MET:HB3	1:B:161:MET:HE2	2.00	0.43
1:A:190:LYS:HG3	1:A:192:GLU:OE1	2.18	0.43
1:A:158:VAL:HG12	1:A:158:VAL:O	2.19	0.42
1:D:120:LEU:HD22	1:D:187:GLU:HG3	2.00	0.42
1:B:25:ARG:NH2	1:C:172:GLU:OE2	2.50	0.42
1:A:111:GLN:HG3	5:A:2063:HOH:O	2.19	0.42
1:B:119:ALA:HB3	1:B:121:PHE:CE2	2.53	0.42
1:D:155:ARG:CZ	1:D:155:ARG:H	2.33	0.42
1:A:59:THR:HG22	1:A:62:ALA:CB	2.50	0.41
1:B:192:GLU:H	1:B:192:GLU:CD	2.28	0.41
1:A:43:ARG:HH21	1:A:46:ARG:HG3	1.85	0.41
1:B:199:ARG:O	1:B:202:ILE:HG22	2.20	0.41
1:A:8:LEU:HD11	1:A:102:LEU:HG	2.02	0.41
1:C:160:MET:HE3	1:C:160:MET:HB2	1.93	0.41
1:D:187:GLU:HG2	5:D:2104:HOH:O	2.20	0.41
1:A:26:ILE:HD13	1:A:101:PHE:CZ	2.56	0.40
1:C:161:MET:HE2	1:C:161:MET:CB	2.45	0.40
1:B:104:ILE:HG22	4:B:1208:5GP:C6	2.52	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	164/207 (79%)	161 (98%)	2 (1%)	1 (1%)	21	13
1	B	169/207 (82%)	167 (99%)	2 (1%)	0	100	100
1	C	170/207 (82%)	167 (98%)	2 (1%)	1 (1%)	21	13
1	D	174/207 (84%)	172 (99%)	2 (1%)	0	100	100
All	All	677/828 (82%)	667 (98%)	8 (1%)	2 (0%)	36	29

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	127	PRO
1	A	158	VAL

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	148/183 (81%)	143 (97%)	5 (3%)	32	25
1	B	152/183 (83%)	146 (96%)	6 (4%)	28	21
1	C	153/183 (84%)	146 (95%)	7 (5%)	24	16
1	D	156/183 (85%)	148 (95%)	8 (5%)	21	13
All	All	609/732 (83%)	583 (96%)	26 (4%)	26	18

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	43	ARG
1	A	44	GLN
1	A	59	THR
1	A	70	ASP
1	A	79	VAL
1	B	34	TYR
1	B	105	GLU
1	B	199	ARG
1	B	200	LYS
1	B	202	ILE
1	B	204	GLU
1	C	34	TYR
1	C	44	GLN
1	C	69	ASP
1	C	74	GLU
1	C	155	ARG
1	C	185	GLU
1	C	191	ARG

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Mol	Chain	Res	Type
1	D	24	LYS
1	D	34	TYR
1	D	60	ARG
1	D	74	GLU
1	D	105	GLU
1	D	155	ARG
1	D	156	LYS
1	D	202	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	111	GLN
1	C	98	HIS
1	D	81	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 11 ligands modelled in this entry, 3 are monoatomic - leaving 8 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	SO4	C	1198	-	4,4,4	0.25	0	6,6,6	0.49	0
2	SO4	D	1208	-	4,4,4	0.16	0	6,6,6	0.42	0
2	SO4	A	1198	-	4,4,4	0.16	0	6,6,6	0.42	0
4	5GP	D	1210	3	26,26,26	1.14	2 (7%)	40,40,40	1.75	9 (22%)
2	SO4	B	1206	-	4,4,4	0.23	0	6,6,6	0.39	0
4	5GP	B	1208	3	26,26,26	1.23	4 (15%)	40,40,40	1.62	7 (17%)
4	5GP	C	1200	3	26,26,26	1.02	1 (3%)	40,40,40	1.71	9 (22%)
2	SO4	D	1207	-	4,4,4	0.24	0	6,6,6	0.73	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	5GP	B	1208	3	-	0/10/26/26	0/3/3/3
4	5GP	C	1200	3	-	0/10/26/26	0/3/3/3
4	5GP	D	1210	3	-	0/10/26/26	0/3/3/3

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	1210	5GP	C6-N1	-3.18	1.32	1.38
4	B	1208	5GP	C5-N7	-2.70	1.33	1.39
4	B	1208	5GP	C6-N1	-2.65	1.33	1.38
4	D	1210	5GP	C5-N7	-2.64	1.34	1.39
4	C	1200	5GP	C6-N1	-2.23	1.34	1.38
4	B	1208	5GP	C4-N9	-2.17	1.32	1.38
4	B	1208	5GP	O6-C6	2.15	1.27	1.23

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	1210	5GP	C5-C4-N3	-5.36	119.77	128.46
4	B	1208	5GP	C5-C4-N3	-5.03	120.30	128.46
4	C	1200	5GP	C5-C4-N3	-4.76	120.73	128.46
4	C	1200	5GP	C2-N3-C4	4.17	119.73	112.30
4	D	1210	5GP	C2-N3-C4	3.98	119.39	112.30
4	B	1208	5GP	C2-N3-C4	3.65	118.79	112.30
4	C	1200	5GP	N9-C4-N3	3.41	132.78	125.94
4	D	1210	5GP	N9-C4-N3	3.11	132.18	125.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	1210	5GP	N9-C8-N7	-3.10	107.56	113.39
4	B	1208	5GP	N9-C4-N3	3.06	132.07	125.94
4	C	1200	5GP	N9-C8-N7	-2.82	108.08	113.39
4	B	1208	5GP	N9-C8-N7	-2.51	108.67	113.39
4	D	1210	5GP	C4-C5-N7	-2.37	106.97	110.72
4	C	1200	5GP	C5-C6-N1	2.34	119.14	113.19
4	B	1208	5GP	C4-C5-N7	-2.33	107.04	110.72
4	C	1200	5GP	O3'-C3'-C4'	-2.32	104.35	111.05
4	D	1210	5GP	C8-N7-C5	2.31	108.42	104.24
4	D	1210	5GP	O3'-C3'-C4'	-2.26	104.52	111.05
4	C	1200	5GP	C8-N9-C4	2.21	110.25	106.05
4	D	1210	5GP	C5-C6-N1	2.18	118.72	113.19
4	C	1200	5GP	O4'-C1'-N9	2.14	113.27	108.36
4	C	1200	5GP	C6-C5-N7	2.13	134.21	130.25
4	B	1208	5GP	O3'-C3'-C4'	-2.08	105.04	111.05
4	D	1210	5GP	O6-C6-C5	-2.01	121.26	126.60
4	B	1208	5GP	O6-C6-C5	-2.00	121.29	126.60

There are no chirality outliers.

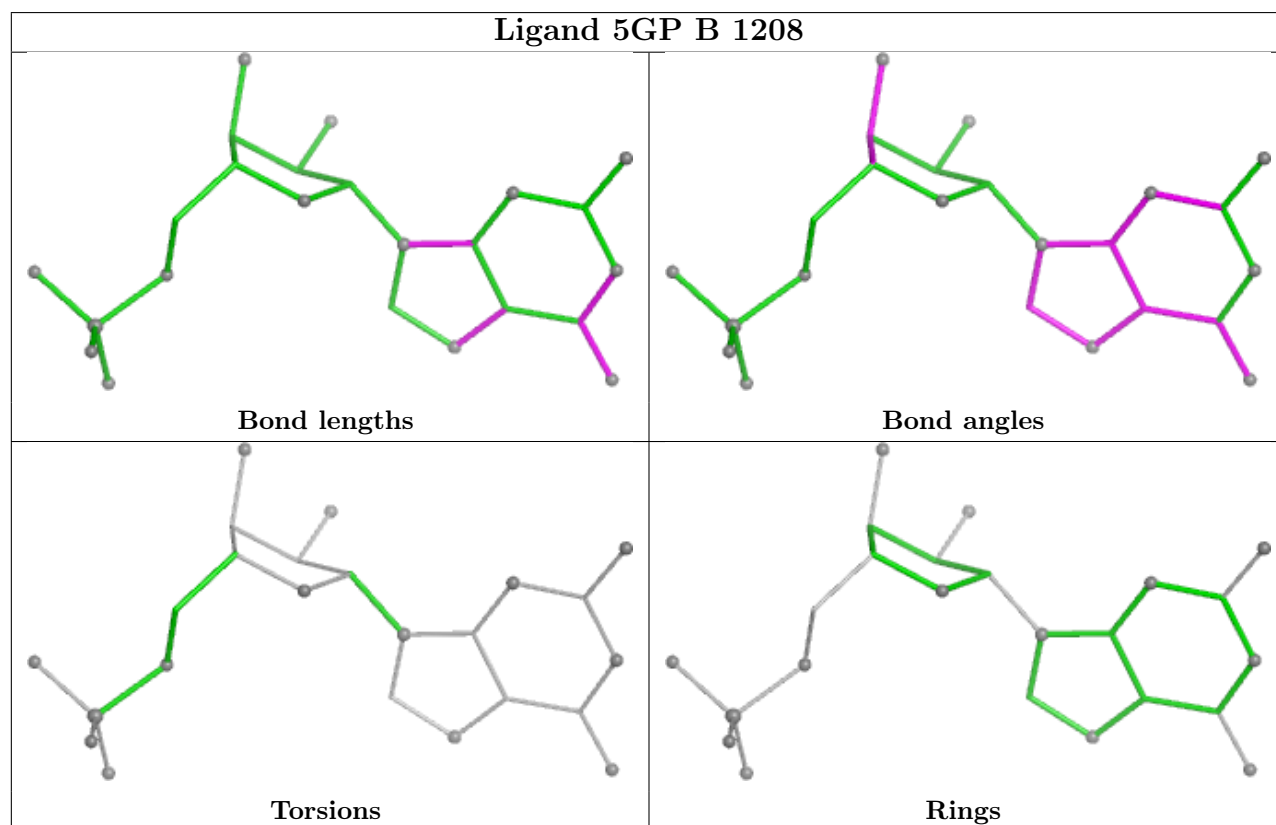
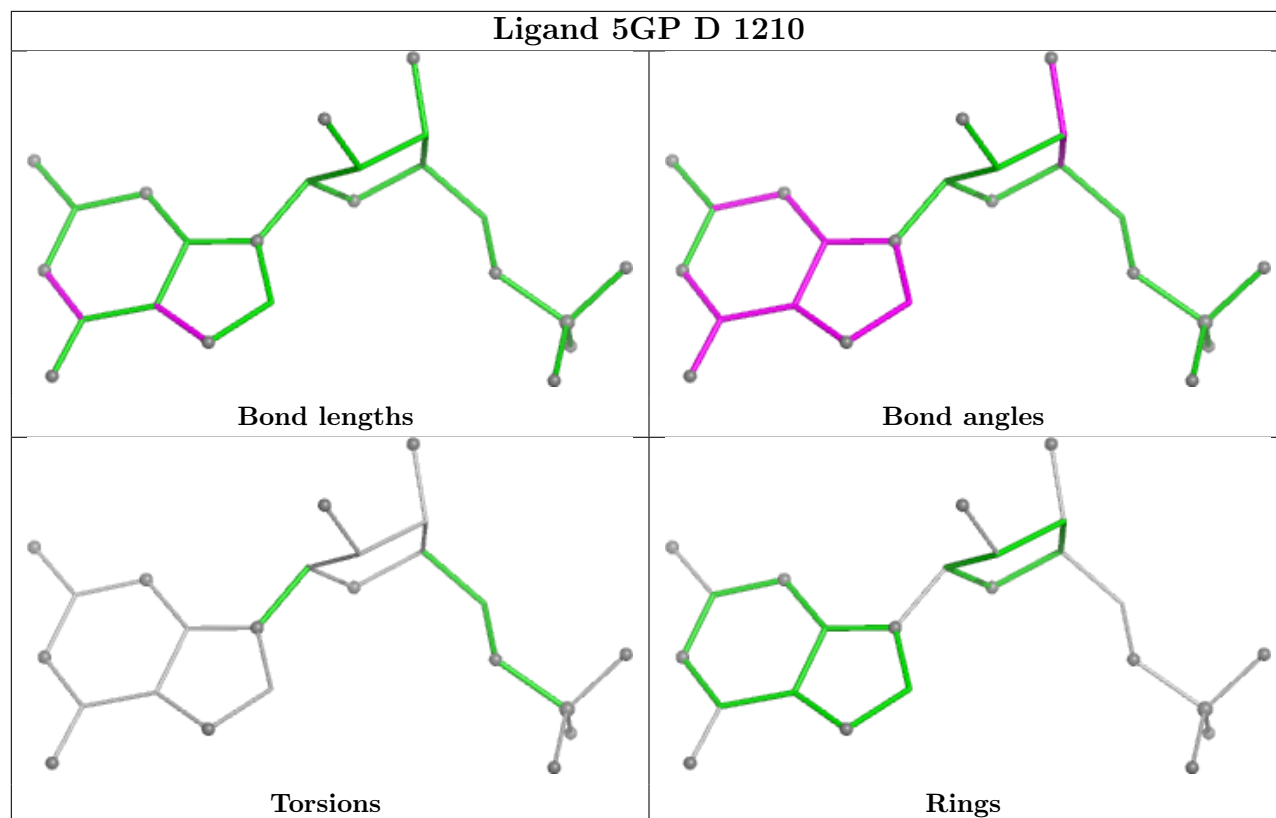
There are no torsion outliers.

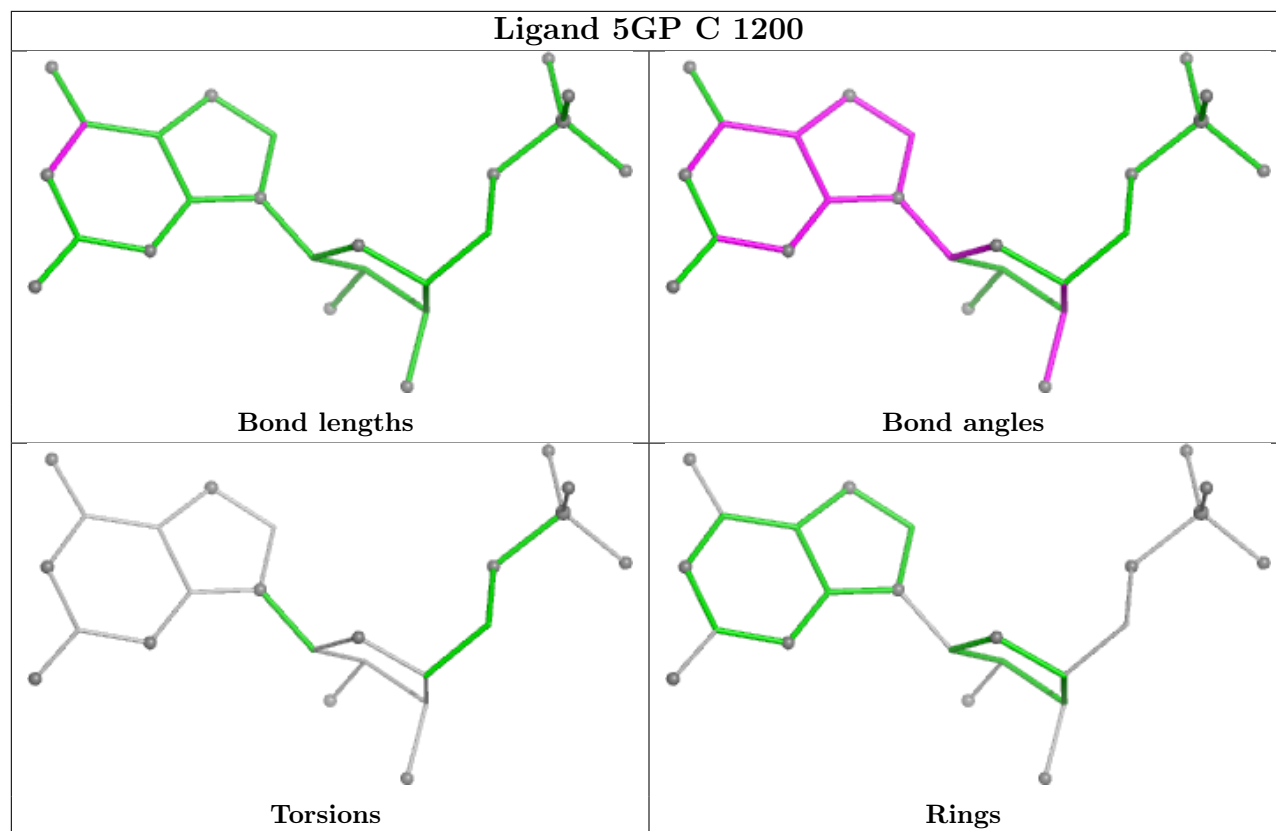
There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	1208	5GP	2	0
2	D	1207	SO4	1	0

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





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [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	167/207 (80%)	0.85	14 (8%) 17 18	15, 24, 35, 41	1 (0%)
1	B	173/207 (83%)	0.88	22 (12%) 8 8	14, 22, 44, 49	0
1	C	174/207 (84%)	0.87	25 (14%) 6 6	14, 21, 44, 56	0
1	D	178/207 (85%)	0.86	17 (9%) 13 14	14, 21, 45, 55	0
All	All	692/828 (83%)	0.87	78 (11%) 10 10	14, 22, 41, 56	1 (0%)

All (78) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	155	ARG	6.0
1	B	205	ALA	5.6
1	C	149	SER	4.9
1	B	159	GLU	4.8
1	D	156	LYS	4.7
1	A	158	VAL	4.7
1	D	162	ASN	4.6
1	C	15	SER	4.6
1	C	151	ILE	4.5
1	A	78	TYR	4.4
1	B	158	VAL	4.4
1	D	128	SER	4.3
1	C	150	ARG	4.3
1	D	154	ALA	4.2
1	C	158	VAL	4.1
1	A	196	ALA	4.1
1	C	155	ARG	4.1
1	C	152	ASN	4.0
1	B	198	TYR	4.0
1	B	201	MET	3.9
1	B	47	GLU	3.8

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Mol	Chain	Res	Type	RSRZ
1	B	161	MET	3.7
1	D	204	GLU	3.7
1	C	161	MET	3.6
1	D	157	GLU	3.5
1	B	163	LEU	3.5
1	B	202	ILE	3.4
1	D	160	MET	3.3
1	A	159	GLU	3.3
1	C	148	GLN	3.3
1	D	202	ILE	3.2
1	A	36[A]	TYR	3.1
1	D	203	LEU	3.1
1	C	153	GLU	3.1
1	C	154	ALA	3.0
1	A	160	MET	3.0
1	C	47	GLU	3.0
1	A	156	LYS	2.9
1	B	203	LEU	2.9
1	D	158	VAL	2.9
1	C	44	GLN	2.8
1	B	160	MET	2.7
1	C	159	GLU	2.7
1	C	197	LYS	2.7
1	B	204	GLU	2.6
1	C	128	SER	2.6
1	D	206	LYS	2.6
1	A	76	ALA	2.6
1	C	196	ALA	2.6
1	D	196	ALA	2.6
1	B	45	MET	2.6
1	B	196	ALA	2.6
1	B	15	SER	2.6
1	B	78	TYR	2.5
1	B	157	GLU	2.5
1	C	163	LEU	2.5
1	A	194	VAL	2.5
1	B	50	VAL	2.5
1	A	29	ASP	2.5
1	B	48	GLY	2.4
1	C	14	PRO	2.4
1	A	157	GLU	2.4
1	D	201	MET	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	15	SER	2.3
1	C	95	ASP	2.3
1	D	95	ASP	2.3
1	B	162	ASN	2.3
1	C	182	CYS	2.3
1	C	156	LYS	2.3
1	C	162	ASN	2.2
1	A	17	VAL	2.2
1	D	163	LEU	2.2
1	C	48	GLY	2.1
1	A	47	GLU	2.1
1	B	16	GLY	2.1
1	D	68	LYS	2.0
1	B	182	CYS	2.0
1	C	29	ASP	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.

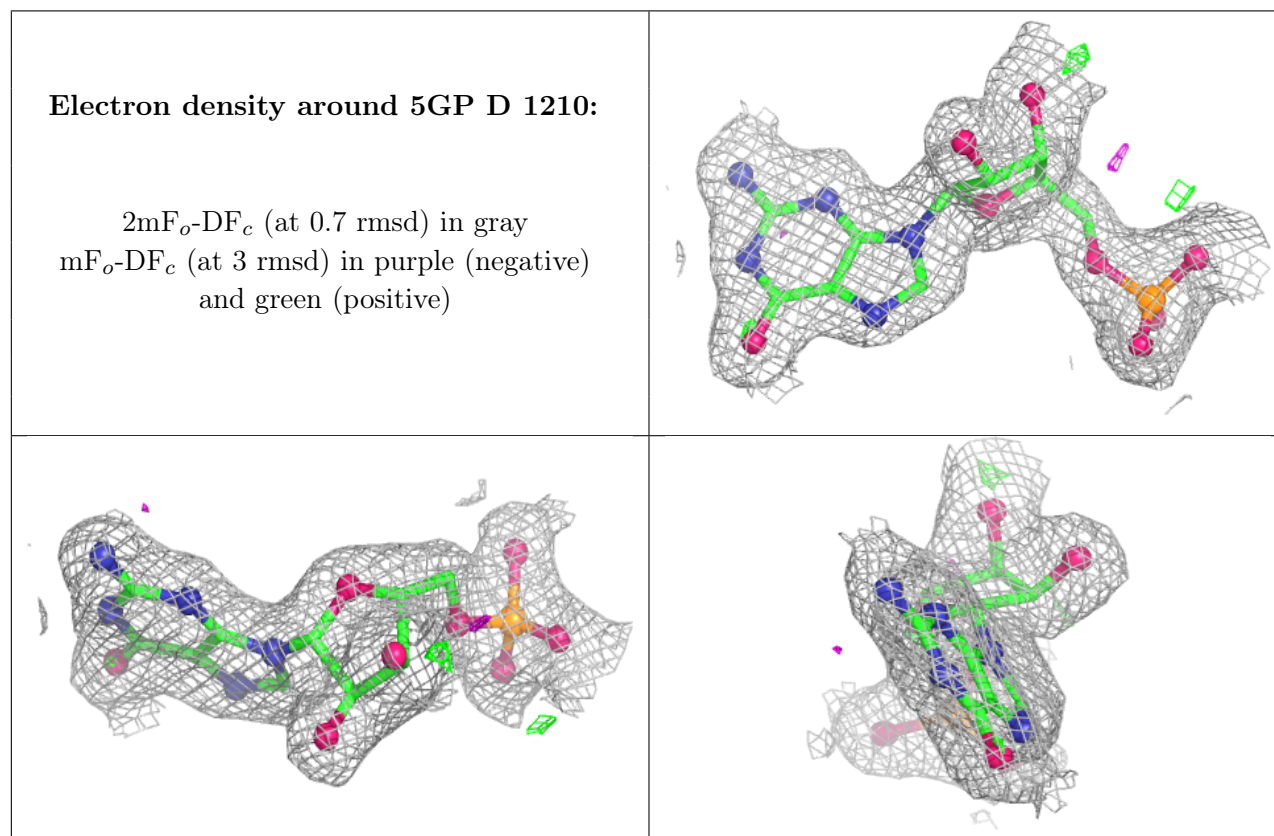
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	SO4	D	1207	5/5	0.93	0.12	31,32,38,39	0
2	SO4	B	1206	5/5	0.94	0.12	32,34,38,39	0
2	SO4	C	1198	5/5	0.95	0.11	30,32,37,37	0
4	5GP	D	1210	24/24	0.95	0.09	18,21,28,29	0
3	K	D	1209	1/1	0.96	0.09	31,31,31,31	0
4	5GP	B	1208	24/24	0.96	0.09	18,23,32,32	0
4	5GP	C	1200	24/24	0.96	0.08	16,20,25,28	0
2	SO4	A	1198	5/5	0.96	0.09	39,40,41,42	0

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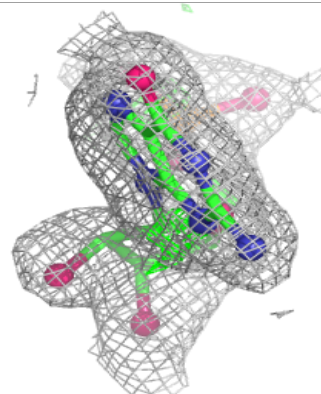
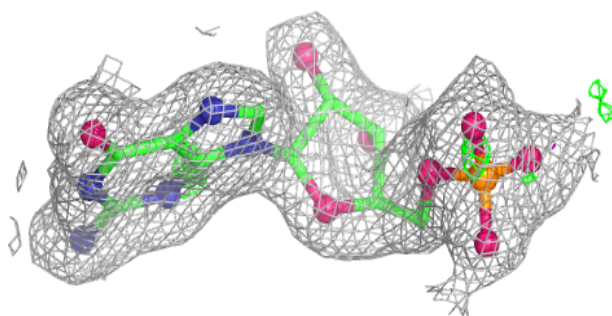
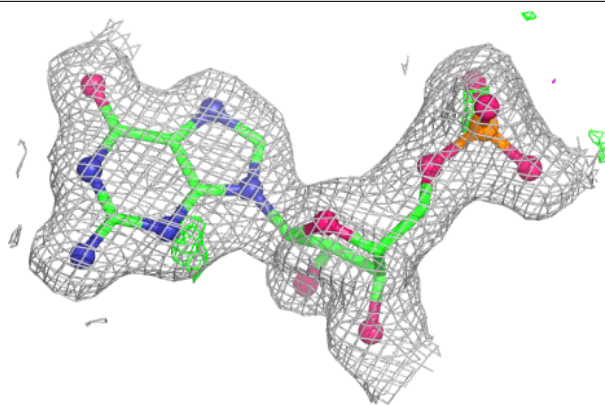
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	SO4	D	1208	5/5	0.97	0.07	31,31,33,34	0
3	K	B	1207	1/1	0.97	0.07	31,31,31,31	0
3	K	C	1199	1/1	0.98	0.06	33,33,33,33	0

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.

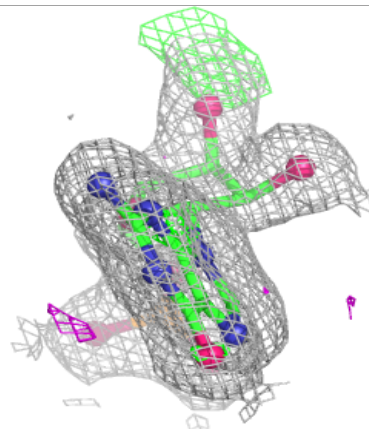
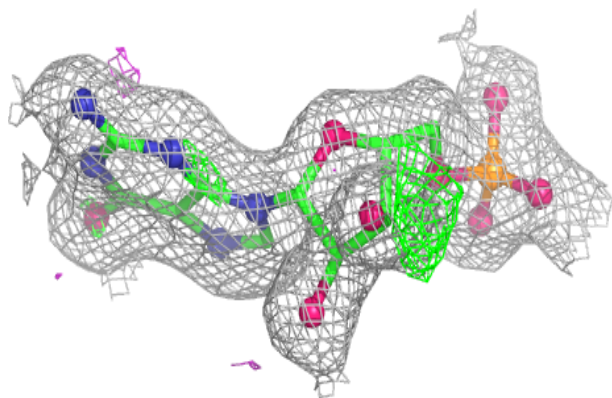
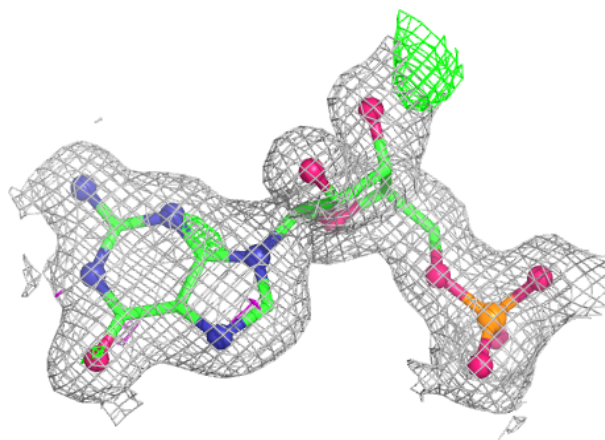


Electron density around 5GP B 1208:

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

**Electron density around 5GP C 1200:**

$2mF_o-DF_c$ (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.