



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 07:37 PM UTC

PDB ID : 1DEL / pdb_00001del
Title : DEOXYNUCLEOSIDE MONOPHOSPHATE KINASE COMPLEXED
WITH DEOXY-GMP AND AMP
Authors : Teplyakov, A.; Sebastiao, P.
Deposited on : 1996-01-09
Resolution : 2.20 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

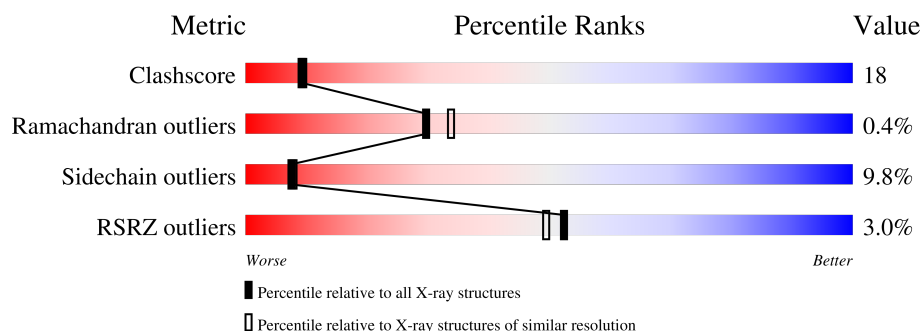
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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))
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	241	3% 61% 31% 7%
1	B	241	2% 56% 31% 7% ••

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 4061 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.

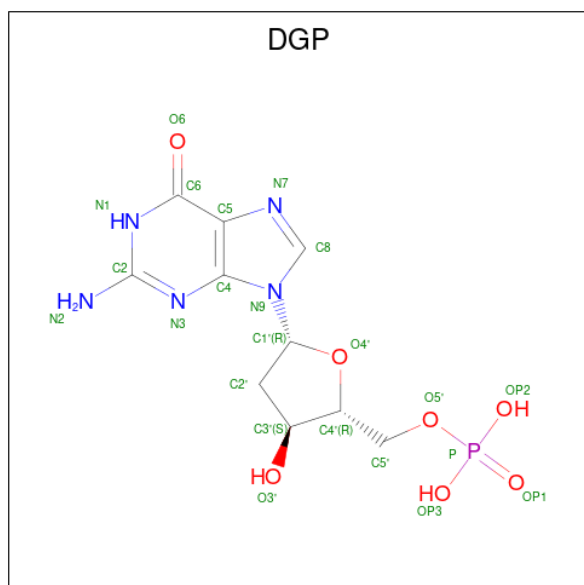
- Molecule 1 is a protein called DEOXYNUCLEOSIDE MONOPHOSPHATE KINASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	241	Total	C	N	O	S	0	0	0
			1925	1230	318	368	9			
1	B	231	Total	C	N	O	S	0	0	0
			1843	1183	302	349	9			

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

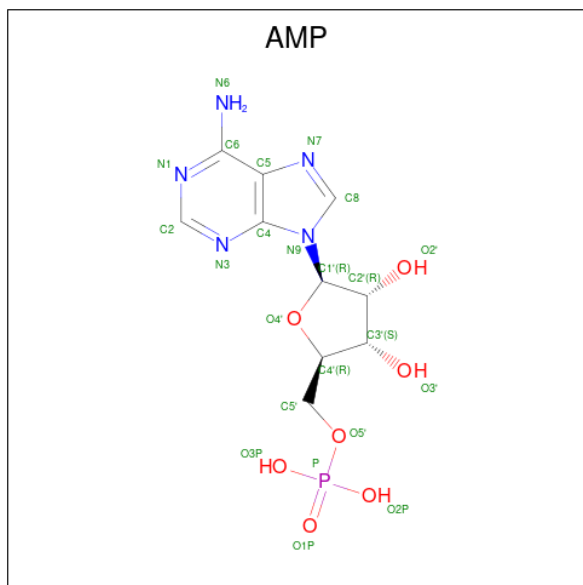
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mg	0	0
			1	1		
2	B	1	Total	Mg	0	0
			1	1		

- Molecule 3 is 2'-DEOXYGUANOSINE-5'-MONOPHOSPHATE (CCD ID: DGP) (formula: C₁₀H₁₄N₅O₇P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
3	B	1	Total	C	N	O	P	0	0
			23	10	5	7	1		

- Molecule 4 is ADENOSINE MONOPHOSPHATE (CCD ID: AMP) (formula: $C_{10}H_{14}N_5O_7P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	B	1	Total	C	N	O	P	0	0
			23	10	5	7	1		

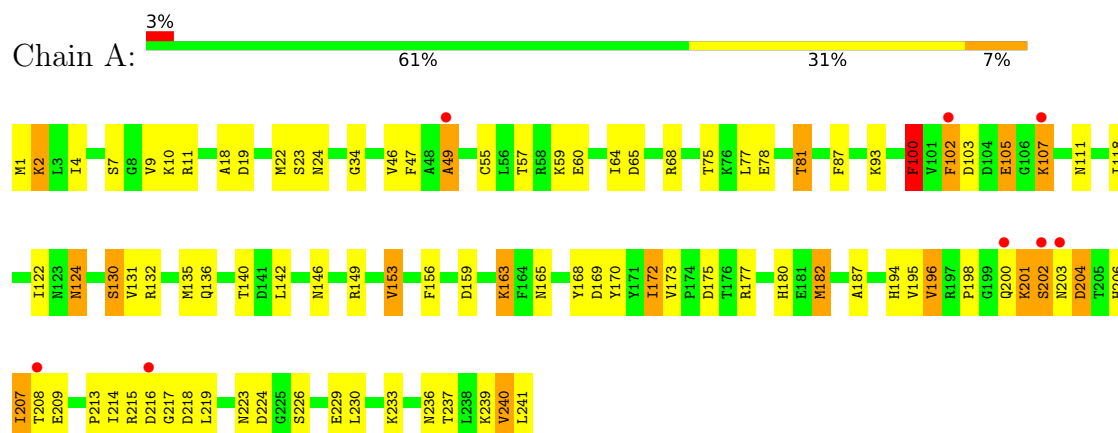
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	124	Total	O	0	0
			124	124		
5	B	98	Total	O	0	0
			98	98		

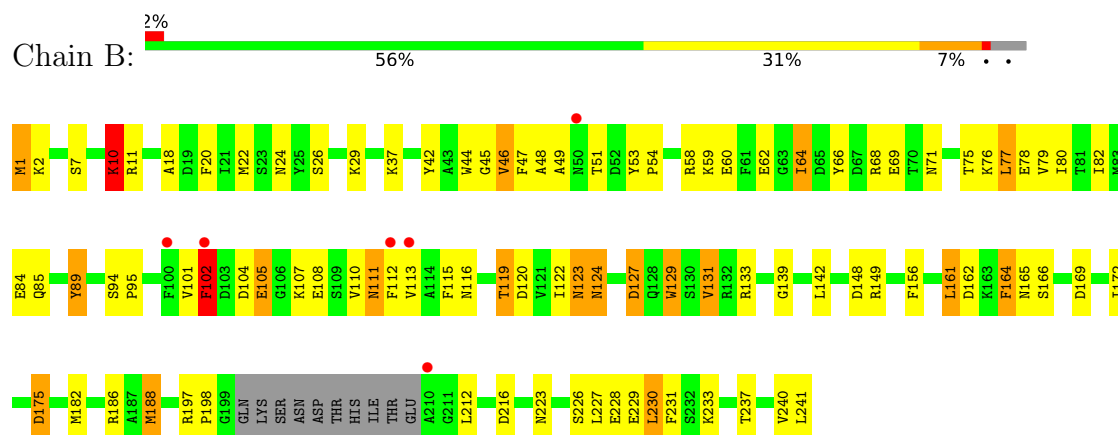
3 Residue-property plots

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: DEOXYNUCLEOSIDE MONOPHOSPHATE KINASE



• Molecule 1: DEOXYNUCLEOSIDE MONOPHOSPHATE KINASE



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	155.20Å 58.50Å 75.70Å 90.00° 108.10° 90.00°	Depositor
Resolution (Å)	10.00 – 2.20 10.00 – 2.20	Depositor EDS
% Data completeness (in resolution range)	95.9 (10.00-2.20) 95.6 (10.00-2.20)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.14 (at 2.20Å)	Xtriage
Refinement program	PROLSQ	Depositor
R, R_{free}	0.186 , (Not available) 0.183 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	31.9	Xtriage
Anisotropy	0.220	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 79.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4061	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.83% 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: DGP, MG, AMP

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	0.97	1/1963 (0.1%)	1.67	30/2655 (1.1%)
1	B	0.90	0/1879	1.63	29/2539 (1.1%)
All	All	0.94	1/3842 (0.0%)	1.65	59/5194 (1.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	173	VAL	CA-CB	-6.02	1.49	1.54

All (59) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	165	ASN	CA-CB-CG	9.49	122.09	112.60
1	A	100	PHE	CA-CB-CG	8.92	122.72	113.80
1	B	112	PHE	CA-CB-CG	-7.73	106.07	113.80
1	B	231	PHE	CA-CB-CG	-7.67	106.13	113.80
1	A	217	GLY	N-CA-C	-7.42	104.61	114.25
1	B	123	ASN	CA-CB-CG	-7.32	105.28	112.60
1	B	164	PHE	CA-CB-CG	-7.29	106.51	113.80
1	B	127	ASP	CA-CB-CG	7.12	119.72	112.60
1	B	112	PHE	N-CA-C	7.07	119.06	111.36
1	B	188	MET	CB-CA-C	-6.92	98.86	110.56
1	B	102	PHE	CA-CB-CG	6.90	120.70	113.80
1	A	24	ASN	CA-CB-CG	-6.86	105.74	112.60
1	A	196	VAL	N-CA-C	6.83	117.68	108.11
1	A	105	GLU	CA-CB-CG	6.77	127.64	114.10
1	A	19	ASP	CA-CB-CG	-6.74	105.86	112.60
1	A	172	ILE	N-CA-C	6.58	117.39	108.17
1	B	133	ARG	CD-NE-CZ	6.44	133.42	124.40
1	B	129	TRP	N-CA-C	6.42	119.29	107.99

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	175	ASP	N-CA-C	6.31	120.59	112.26
1	A	195	VAL	N-CA-CB	6.04	119.28	111.67
1	B	115	PHE	CA-CB-CG	-5.98	107.82	113.80
1	A	216	ASP	N-CA-CB	5.97	118.77	110.17
1	A	153	VAL	CA-C-N	5.92	128.21	120.28
1	A	153	VAL	C-N-CA	5.92	128.21	120.28
1	A	124	ASN	CA-CB-CG	-5.88	106.72	112.60
1	B	198	PRO	N-CA-CB	5.87	108.46	103.35
1	A	215	ARG	CA-C-O	-5.86	114.45	121.89
1	A	34	GLY	CA-C-N	5.70	125.38	119.56
1	A	34	GLY	C-N-CA	5.70	125.38	119.56
1	B	161	LEU	CA-C-N	5.69	127.91	120.28
1	B	161	LEU	C-N-CA	5.69	127.91	120.28
1	B	175	ASP	N-CA-C	5.67	119.51	112.59
1	A	213	PRO	N-CA-CB	5.66	108.60	103.34
1	A	173	VAL	N-CA-CB	5.57	117.09	111.36
1	A	159	ASP	CA-CB-CG	-5.51	107.09	112.60
1	B	64	ILE	N-CA-CB	5.46	117.79	111.90
1	A	146	ASN	CA-CB-CG	-5.45	107.15	112.60
1	B	71	ASN	CA-CB-CG	5.39	117.99	112.60
1	A	204	ASP	CA-CB-CG	5.38	117.98	112.60
1	B	212	LEU	CA-C-N	5.35	125.27	119.76
1	B	212	LEU	C-N-CA	5.35	125.27	119.76
1	A	163	LYS	N-CA-C	5.34	120.80	113.97
1	A	146	ASN	N-CA-CB	-5.30	103.57	110.88
1	A	87	PHE	N-CA-CB	5.29	117.90	110.12
1	B	166	SER	N-CA-C	5.28	119.70	113.16
1	B	148	ASP	CA-CB-CG	5.24	117.83	112.60
1	A	11	ARG	N-CA-C	5.23	118.81	111.90
1	B	230	LEU	N-CA-C	-5.16	105.57	111.14
1	A	100	PHE	N-CA-C	-5.15	101.36	109.50
1	B	89	TYR	CA-C-N	5.15	127.19	120.28
1	B	89	TYR	C-N-CA	5.15	127.19	120.28
1	B	139	GLY	N-CA-C	5.15	118.91	112.73
1	B	172	ILE	CB-CA-C	-5.14	103.46	110.96
1	B	188	MET	CA-C-O	-5.14	113.55	119.56
1	A	230	LEU	N-CA-CB	5.13	117.51	110.07
1	A	196	VAL	CA-C-O	5.12	125.77	120.39
1	A	215	ARG	N-CA-C	-5.11	102.90	110.46
1	A	169	ASP	CA-CB-CG	5.09	117.69	112.60
1	B	10	LYS	CB-CA-C	5.01	117.36	109.84

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	1925	0	1903	68	0
1	B	1843	0	1826	69	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	23	0	12	2	0
3	B	23	0	11	2	0
4	B	23	0	12	1	0
5	A	124	0	0	3	0
5	B	98	0	0	1	0
All	All	4061	0	3764	136	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 18.

All (136) 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:124:ASN:N	1:B:124:ASN:HD22	1.58	1.01
1:A:204:ASP:HB3	1:A:209:GLU:CB	1.96	0.94
1:B:104:ASP:O	1:B:107:LYS:HD2	1.69	0.93
1:B:226:SER:OG	1:B:229:GLU:HG3	1.69	0.92
1:B:104:ASP:HB3	1:B:107:LYS:HB2	1.51	0.90
1:A:204:ASP:HB3	1:A:209:GLU:HG3	1.53	0.89
1:A:204:ASP:HB3	1:A:209:GLU:CG	2.05	0.87
1:A:237:THR:O	1:A:241:LEU:HD23	1.76	0.83
1:A:226:SER:HB3	1:A:229:GLU:HG3	1.63	0.80
1:B:75:THR:HB	1:B:78:GLU:HG3	1.64	0.80
1:B:124:ASN:N	1:B:124:ASN:ND2	2.29	0.79
1:B:68:ARG:HG2	1:B:131:VAL:HG11	1.65	0.78
1:A:136:GLN:HG2	3:A:301:DGP:H2'	1.63	0.78
1:A:55:CYS:SG	5:A:348:HOH:O	2.42	0.77
1:A:57:THR:OG1	1:A:60:GLU:OE1	2.02	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1:MET:HE2	1:A:168:TYR:HB2	1.66	0.76
1:B:68:ARG:HG2	1:B:131:VAL:CG1	2.16	0.76
1:A:204:ASP:HB3	1:A:209:GLU:HB3	1.67	0.76
1:A:204:ASP:CB	1:A:209:GLU:HG3	2.16	0.74
1:A:75:THR:OG1	1:A:78:GLU:HG3	1.89	0.73
1:A:200:GLN:HG2	1:A:201:LYS:N	2.05	0.71
1:A:177:ARG:HG2	1:A:208:THR:O	1.93	0.69
1:A:194:HIS:CG	1:A:214:ILE:HD13	2.27	0.69
1:B:102:PHE:CZ	1:B:111:ASN:HA	2.29	0.68
1:A:9:VAL:HB	1:A:202:SER:OG	1.95	0.66
1:B:237:THR:O	1:B:240:VAL:HG22	1.96	0.66
1:A:204:ASP:CG	1:A:209:GLU:HG3	2.22	0.63
1:B:75:THR:HG22	1:B:77:LEU:H	1.63	0.62
1:A:102:PHE:CZ	1:A:111:ASN:HA	2.34	0.62
1:A:200:GLN:HG2	1:A:201:LYS:H	1.66	0.61
1:A:68:ARG:O	1:A:130:SER:HB2	2.01	0.60
1:A:177:ARG:HB3	1:A:208:THR:HA	1.84	0.60
1:B:69:GLU:HA	1:B:69:GLU:OE1	2.01	0.60
1:B:104:ASP:O	1:B:105:GLU:HB2	2.01	0.60
1:A:206:HIS:HB3	1:A:209:GLU:HG2	1.83	0.59
1:B:51:THR:OG1	1:B:53:TYR:HD2	1.86	0.59
1:B:53:TYR:HD1	1:B:54:PRO:HD2	1.68	0.58
1:A:100:PHE:CD2	1:A:111:ASN:HB3	2.39	0.58
1:B:122:ILE:HD11	1:B:142:LEU:HD11	1.86	0.58
1:B:75:THR:CB	1:B:78:GLU:HG3	2.33	0.58
1:B:116:ASN:HA	1:B:119:THR:OG1	2.03	0.58
1:A:10:LYS:HB2	1:A:202:SER:HB3	1.86	0.57
1:B:80:ILE:HD13	1:B:119:THR:HG22	1.86	0.57
1:B:123:ASN:C	1:B:124:ASN:HD22	2.11	0.57
1:A:102:PHE:CE2	1:A:111:ASN:HA	2.39	0.56
1:B:102:PHE:HZ	1:B:111:ASN:HA	1.71	0.56
1:A:118:ILE:O	1:A:122:ILE:HG12	2.05	0.56
1:A:4:ILE:HB	1:A:172:ILE:HD13	1.88	0.55
1:B:53:TYR:CD1	1:B:54:PRO:HD2	2.41	0.55
1:A:136:GLN:HB3	1:A:206:HIS:CE1	2.41	0.55
1:B:59:LYS:HG3	1:B:64:ILE:HD12	1.88	0.55
1:A:136:GLN:HB3	1:A:206:HIS:NE2	2.22	0.54
1:A:237:THR:O	1:A:241:LEU:CD2	2.52	0.54
1:B:101:VAL:CG1	1:B:108:GLU:HB2	2.37	0.54
1:B:68:ARG:CG	1:B:131:VAL:HG11	2.35	0.54
1:A:207:ILE:HG12	1:A:207:ILE:O	2.06	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:7:SER:HB2	1:A:182:MET:HE2	1.89	0.53
1:B:37:LYS:NZ	3:B:301:DGP:OP3	2.40	0.53
1:A:196:VAL:O	1:A:223:ASN:N	2.33	0.53
1:B:68:ARG:CG	1:B:131:VAL:CG1	2.85	0.52
1:B:120:ASP:O	1:B:124:ASN:ND2	2.43	0.52
1:A:153:VAL:O	1:A:156:PHE:HB3	2.09	0.52
1:B:226:SER:HG	1:B:229:GLU:HG3	1.74	0.52
1:A:78:GLU:O	1:A:81:THR:HB	2.10	0.52
1:A:236:ASN:HA	1:A:239:LYS:HG2	1.91	0.52
3:A:301:DGP:H5"	5:A:346:HOH:O	2.11	0.51
1:A:149:ARG:HG2	1:A:180:HIS:CD2	2.45	0.51
1:A:229:GLU:O	1:A:233:LYS:HG3	2.10	0.50
1:A:59:LYS:HG3	1:A:64:ILE:HB	1.94	0.49
1:A:64:ILE:O	1:A:64:ILE:HG22	2.13	0.49
1:A:7:SER:HB2	1:A:182:MET:CE	2.43	0.49
1:B:51:THR:OG1	1:B:53:TYR:HB2	2.12	0.49
1:B:53:TYR:CB	1:B:54:PRO:HD2	2.42	0.49
1:B:84:GLU:HG3	1:B:110:VAL:HG22	1.95	0.48
1:B:47:PHE:CE2	1:B:78:GLU:HB3	2.49	0.48
1:B:161:LEU:O	1:B:164:PHE:HB2	2.14	0.47
1:A:226:SER:HB3	1:A:229:GLU:CG	2.41	0.47
1:B:59:LYS:O	1:B:64:ILE:HB	2.13	0.47
1:A:149:ARG:CG	1:A:180:HIS:CD2	2.98	0.47
1:A:93:LYS:HA	1:A:93:LYS:HD3	1.79	0.47
1:A:187:ALA:HB2	1:B:164:PHE:HZ	1.80	0.47
1:B:47:PHE:CD1	1:B:48:ALA:N	2.83	0.47
1:A:219:LEU:HD13	1:A:237:THR:HG23	1.96	0.47
1:B:7:SER:HB2	1:B:182:MET:HE3	1.98	0.46
1:B:51:THR:OG1	1:B:53:TYR:CD2	2.67	0.46
1:B:227:LEU:O	1:B:230:LEU:HB3	2.15	0.46
1:B:1:MET:N	5:B:389:HOH:O	2.47	0.45
1:B:10:LYS:HD3	1:B:11:ARG:HE	1.82	0.45
1:B:2:LYS:HA	1:B:2:LYS:HD3	1.68	0.45
1:B:18:ALA:O	1:B:22:MET:HG2	2.17	0.45
1:A:102:PHE:HZ	1:A:111:ASN:HB3	1.82	0.45
1:B:124:ASN:HD22	1:B:124:ASN:H	1.54	0.45
1:A:194:HIS:ND1	1:A:214:ILE:HD13	2.32	0.45
1:A:46:VAL:O	1:A:49:ALA:HB2	2.17	0.44
1:B:10:LYS:NZ	1:B:11:ARG:HH21	2.15	0.44
1:B:188:MET:HE3	1:B:188:MET:HB3	1.91	0.44
1:A:2:LYS:HG3	1:A:170:TYR:CD1	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:105:GLU:C	1:A:107:LYS:H	2.25	0.44
1:A:187:ALA:CB	1:B:164:PHE:HZ	2.31	0.44
1:A:194:HIS:CG	1:A:214:ILE:CD1	3.00	0.44
1:B:78:GLU:O	1:B:82:ILE:HG13	2.18	0.44
1:A:198:PRO:HG2	1:A:224:ASP:HB3	1.99	0.44
1:B:59:LYS:HE3	1:B:59:LYS:HB2	1.55	0.43
1:A:2:LYS:HG3	1:A:170:TYR:CE1	2.53	0.43
1:B:68:ARG:CD	1:B:131:VAL:HG11	2.48	0.43
1:B:186:ARG:HH11	1:B:186:ARG:HD2	1.62	0.43
1:B:223:ASN:OD1	4:B:302:AMP:N6	2.45	0.43
1:B:44:TRP:O	1:B:45:GLY:C	2.62	0.43
1:B:68:ARG:HG2	1:B:131:VAL:CB	2.48	0.43
1:B:76:LYS:HD3	1:B:123:ASN:HA	2.01	0.43
1:A:236:ASN:HA	1:A:239:LYS:HE2	2.00	0.43
1:A:18:ALA:O	1:A:22:MET:HG2	2.19	0.43
1:B:20:PHE:O	1:B:24:ASN:ND2	2.48	0.43
1:B:104:ASP:N	1:B:107:LYS:O	2.52	0.42
1:B:175:ASP:O	3:B:301:DGP:N2	2.51	0.42
1:A:47:PHE:C	1:A:49:ALA:H	2.27	0.42
1:B:229:GLU:O	1:B:233:LYS:HG3	2.19	0.42
1:A:140:THR:HB	1:A:206:HIS:HE1	1.84	0.42
1:B:47:PHE:C	1:B:49:ALA:H	2.28	0.42
1:A:177:ARG:HG2	1:A:208:THR:C	2.45	0.42
1:A:132:ARG:HA	1:A:135:MET:HE2	2.02	0.41
1:A:102:PHE:CZ	1:A:111:ASN:CA	3.01	0.41
1:B:46:VAL:HG11	1:B:85:GLN:HG3	2.02	0.41
1:B:68:ARG:HG2	1:B:131:VAL:HB	2.02	0.41
1:B:79:VAL:HG21	1:B:129:TRP:CD2	2.55	0.41
1:A:118:ILE:HA	1:A:142:LEU:HD21	2.03	0.41
1:A:194:HIS:CD2	1:A:218:ASP:HB3	2.55	0.41
1:B:66:TYR:CD2	1:B:68:ARG:HG3	2.55	0.41
1:B:42:TYR:CD2	1:B:89:TYR:HB2	2.55	0.41
1:A:57:THR:N	1:A:60:GLU:OE1	2.48	0.41
1:A:177:ARG:NH1	5:A:346:HOH:O	2.44	0.41
1:B:26:SER:O	1:B:169:ASP:N	2.51	0.41
1:B:58:ARG:NH1	1:B:62:GLU:OE1	2.28	0.41
1:B:94:SER:HA	1:B:95:PRO:HD3	1.94	0.41
1:A:100:PHE:HD2	1:A:102:PHE:CE1	2.39	0.40
1:A:239:LYS:O	1:A:240:VAL:C	2.63	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	A	239/241 (99%)	224 (94%)	14 (6%)	1 (0%)	30	34
1	B	227/241 (94%)	219 (96%)	7 (3%)	1 (0%)	30	34
All	All	466/482 (97%)	443 (95%)	21 (4%)	2 (0%)	30	34

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	49	ALA
1	B	111	ASN

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	209/209 (100%)	189 (90%)	20 (10%)	8	8
1	B	199/209 (95%)	179 (90%)	20 (10%)	7	7
All	All	408/418 (98%)	368 (90%)	40 (10%)	7	8

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	LYS
1	A	23	SER
1	A	65	ASP
1	A	77	LEU

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Mol	Chain	Res	Type
1	A	81	THR
1	A	100	PHE
1	A	102	PHE
1	A	103	ASP
1	A	107	LYS
1	A	124	ASN
1	A	130	SER
1	A	131	VAL
1	A	163	LYS
1	A	165	ASN
1	A	182	MET
1	A	201	LYS
1	A	202	SER
1	A	203	ASN
1	A	207	ILE
1	A	240	VAL
1	B	1	MET
1	B	10	LYS
1	B	29	LYS
1	B	46	VAL
1	B	60	GLU
1	B	77	LEU
1	B	102	PHE
1	B	105	GLU
1	B	113	VAL
1	B	119	THR
1	B	124	ASN
1	B	127	ASP
1	B	131	VAL
1	B	149	ARG
1	B	156	PHE
1	B	162	ASP
1	B	197	ARG
1	B	216	ASP
1	B	228	GLU
1	B	241	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	31	GLN
1	A	165	ASN

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Mol	Chain	Res	Type
1	B	124	ASN
1	B	165	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 5 ligands modelled in this entry, 2 are monoatomic - leaving 3 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	AMP	B	302	-	25,25,25	1.29	2 (8%)	37,38,38	1.56	10 (27%)
3	DGP	A	301	-	25,25,25	0.86	0	37,38,38	1.59	6 (16%)
3	DGP	B	301	-	25,25,25	1.12	1 (4%)	37,38,38	1.74	7 (18%)

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	AMP	B	302	-	-	3/10/26/26	0/3/3/3
3	DGP	A	301	-	-	4/10/22/22	0/3/3/3
3	DGP	B	301	-	-	4/10/22/22	0/3/3/3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	302	AMP	C5-N7	-3.93	1.31	1.39
3	B	301	DGP	O4'-C4'	3.05	1.51	1.45
4	B	302	AMP	C2-N1	2.89	1.39	1.33

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	301	DGP	O4'-C1'-N9	4.98	116.70	107.86
3	B	301	DGP	C4'-O4'-C1'	-3.90	100.25	109.51
3	A	301	DGP	C1'-N9-C4	3.82	135.80	125.50
3	A	301	DGP	C1'-N9-C8	-3.67	119.69	127.91
3	A	301	DGP	C5-C4-N3	-3.63	122.61	128.39
3	A	301	DGP	O4'-C1'-N9	3.10	113.36	107.86
4	B	302	AMP	O5'-C5'-C4'	2.86	118.72	108.99
3	B	301	DGP	C1'-N9-C4	2.79	133.04	125.50
3	B	301	DGP	O5'-C5'-C4'	2.77	118.44	108.99
3	B	301	DGP	O4'-C4'-C3'	-2.76	99.36	105.65
4	B	302	AMP	O3P-P-O5'	-2.68	99.69	106.67
4	B	302	AMP	C4-N9-C8	-2.62	102.99	105.74
4	B	302	AMP	O2P-P-O5'	2.50	113.17	106.67
3	B	301	DGP	C5'-C4'-C3'	2.43	128.24	114.64
4	B	302	AMP	O3'-C3'-C4'	2.40	117.97	111.08
4	B	302	AMP	C6-C5-N7	-2.30	127.66	132.09
4	B	302	AMP	C4-C5-N7	2.29	113.20	110.58
3	A	301	DGP	C6-C5-C4	2.21	122.15	118.83
3	A	301	DGP	N2-C2-N1	-2.19	112.13	116.76
4	B	302	AMP	O3'-C3'-C2'	2.15	118.72	111.82
3	B	301	DGP	O4'-C1'-C2'	-2.13	102.27	106.25
4	B	302	AMP	O2'-C2'-C3'	2.06	118.43	111.82
4	B	302	AMP	N9-C8-N7	2.05	116.85	113.94

There are no chirality outliers.

All (11) torsion outliers are listed below:

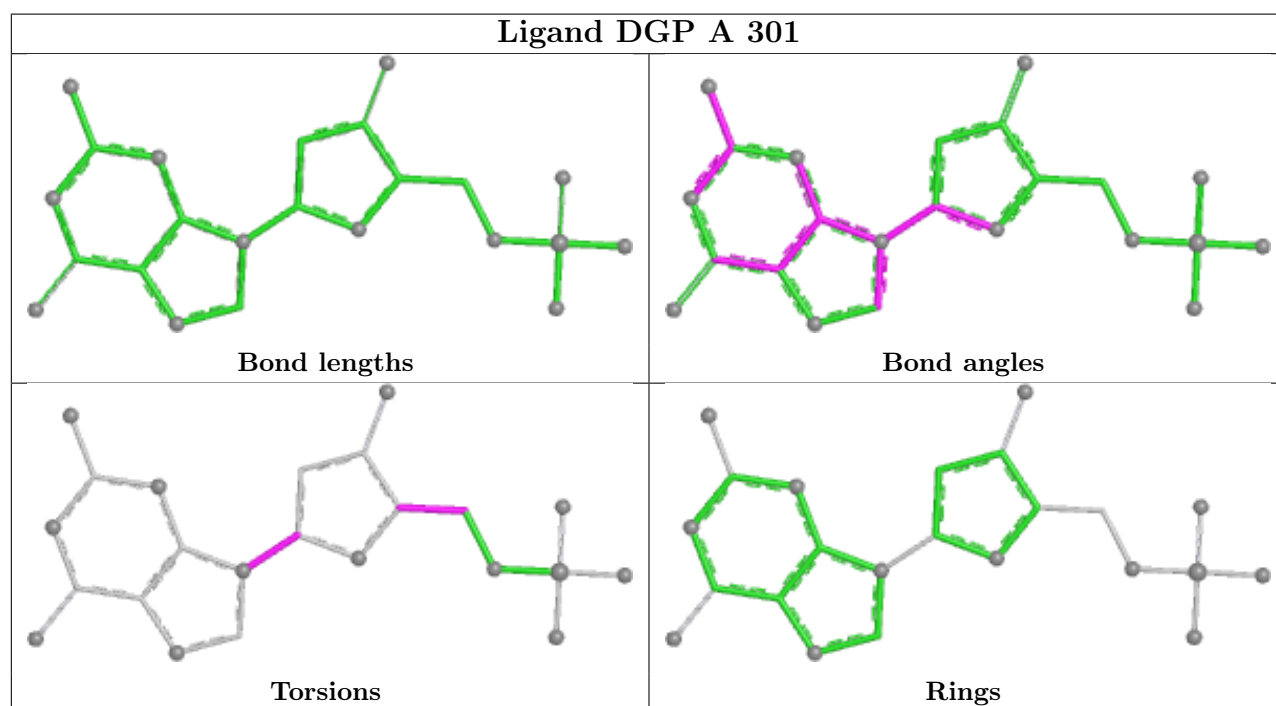
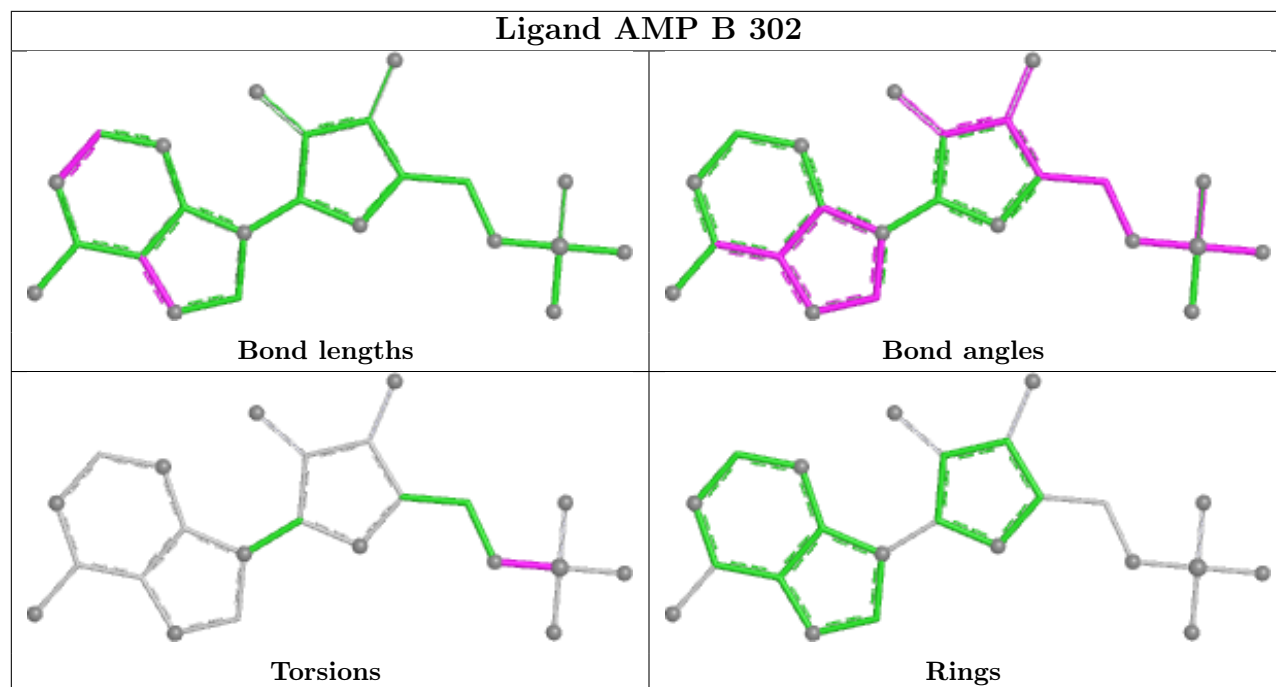
Mol	Chain	Res	Type	Atoms
3	B	301	DGP	O4'-C4'-C5'-O5'
3	B	301	DGP	O4'-C1'-N9-C8
3	B	301	DGP	O4'-C1'-N9-C4
4	B	302	AMP	C5'-O5'-P-O2P
4	B	302	AMP	C5'-O5'-P-O3P
3	A	301	DGP	C3'-C4'-C5'-O5'
3	B	301	DGP	C3'-C4'-C5'-O5'
3	A	301	DGP	O4'-C4'-C5'-O5'
4	B	302	AMP	C5'-O5'-P-O1P
3	A	301	DGP	O4'-C1'-N9-C4
3	A	301	DGP	O4'-C1'-N9-C8

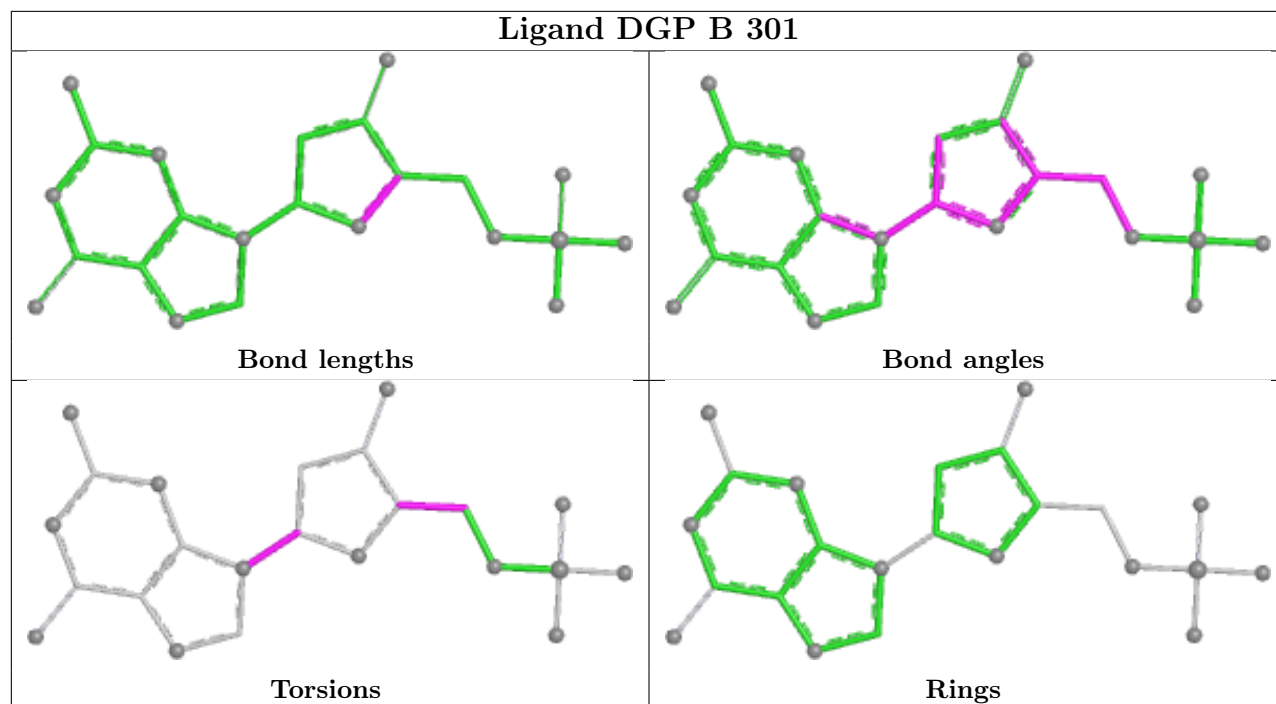
There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	302	AMP	1	0
3	A	301	DGP	2	0
3	B	301	DGP	2	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 [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	A	241/241 (100%)	-0.40	8 (3%)	49	46	12, 30, 92, 100	0
1	B	231/241 (95%)	-0.26	6 (2%)	57	54	11, 35, 86, 100	0
All	All	472/482 (97%)	-0.33	14 (2%)	52	49	11, 32, 90, 100	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	210	ALA	3.7
1	A	200	GLN	3.5
1	B	112	PHE	3.3
1	A	202	SER	3.3
1	A	216	ASP	2.9
1	B	50	ASN	2.9
1	A	203	ASN	2.7
1	A	49	ALA	2.5
1	A	208	THR	2.5
1	B	113	VAL	2.4
1	B	100	PHE	2.3
1	B	102	PHE	2.2
1	A	107	LYS	2.2
1	A	102	PHE	2.1

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](#)

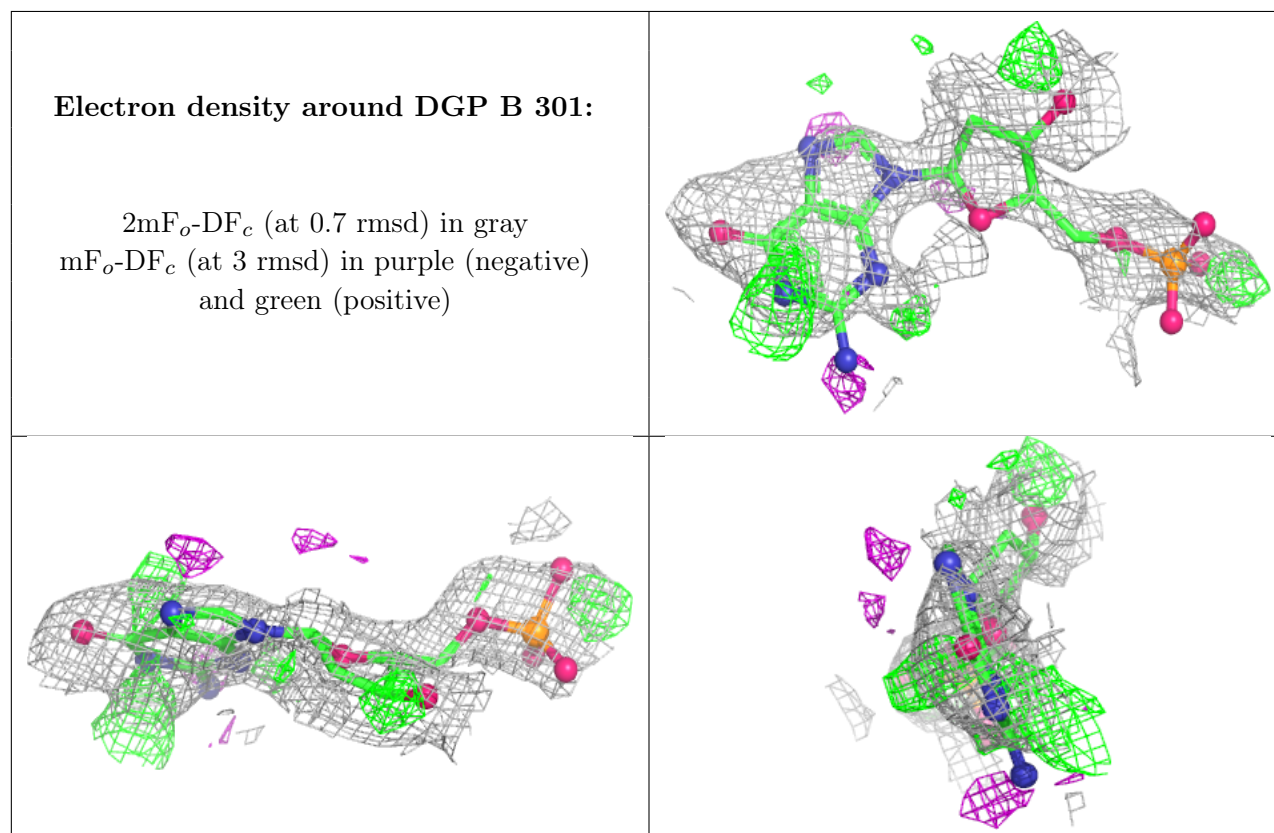
There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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.

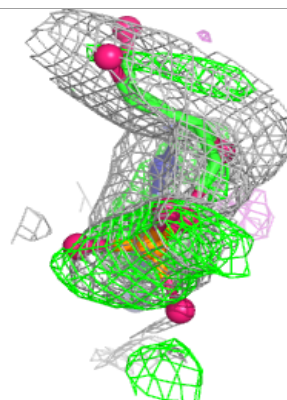
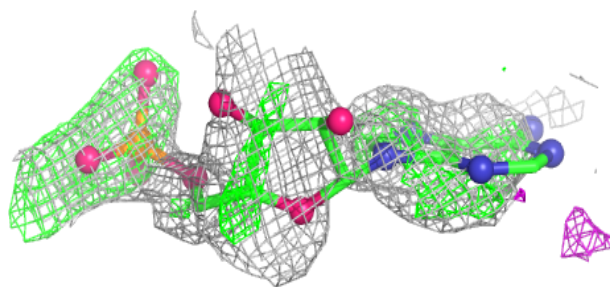
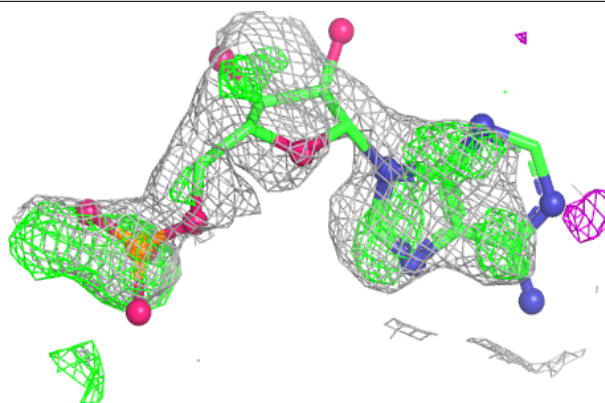
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	DGP	B	301	23/23	0.69	0.19	24,72,97,99	23
4	AMP	B	302	23/23	0.76	0.31	70,77,80,82	23
2	MG	B	300	1/1	0.85	0.07	43,43,43,43	1
3	DGP	A	301	23/23	0.89	0.10	19,69,90,99	0
2	MG	A	300	1/1	0.94	0.11	51,51,51,51	0

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.

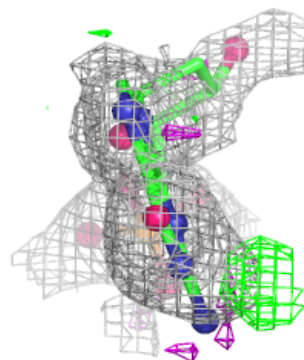
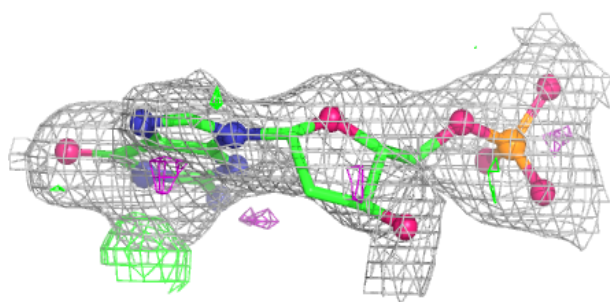
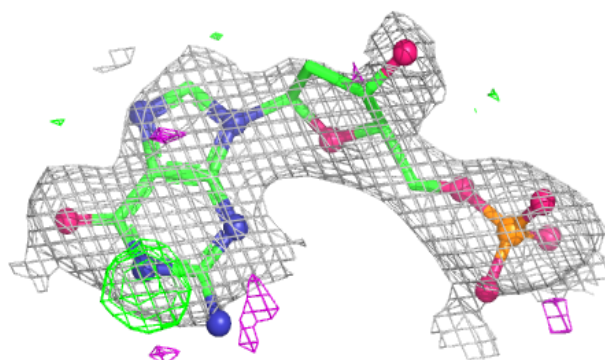


Electron density around AMP B 302:

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

**Electron density around DGP 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.