



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 17, 2024 – 04:52 PM EST

PDB ID : 3TDV  
Title : Structure of the GDP complex of wild-type aminoglycoside 2'-phosphotransferase-IIIa  
Authors : Smith, C.A.; Vakulenko, S.B.  
Deposited on : 2011-08-11  
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](mailto:validation@mail.wwpdb.org)

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtrriage (Phenix) : 1.13  
EDS : 2.36  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.36

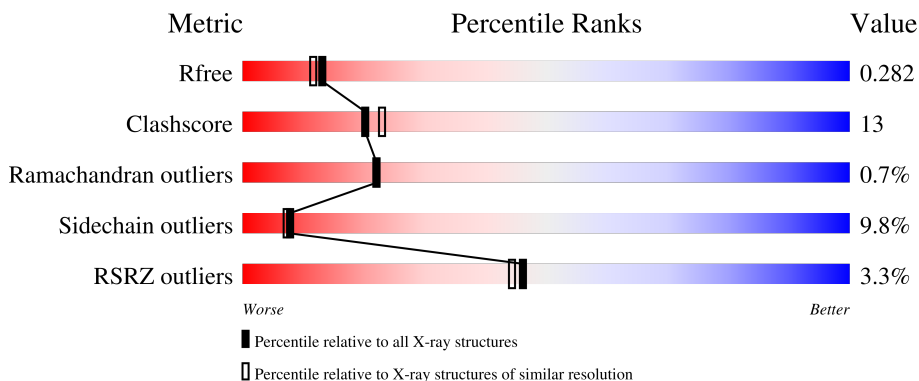
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*


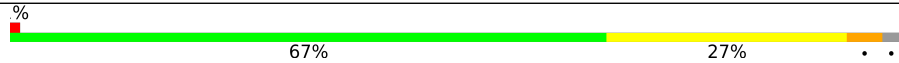
The reported resolution of this entry is 2.20 Å.

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))
$R_{free}$	130704	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	306	
1	B	306	

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 5185 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 Gentamicin resistance protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	301	Total	C	N	O	S	0	2	0
			2413	1543	397	465	8			
1	B	299	Total	C	N	O	S	0	4	0
			2419	1552	395	463	9			

- Molecule 2 is GUANOSINE-5'-DIPHOSPHATE (three-letter code: GDP) (formula: C<sub>10</sub>H<sub>15</sub>N<sub>5</sub>O<sub>11</sub>P<sub>2</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	A	1	Total	C	N	O	P	0	0
			28	10	5	11	2		
2	B	1	Total	C	N	O	P	0	0
			28	10	5	11	2		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	3	Total Mg 3 3	0	0
3	B	3	Total Mg 3 3	0	0

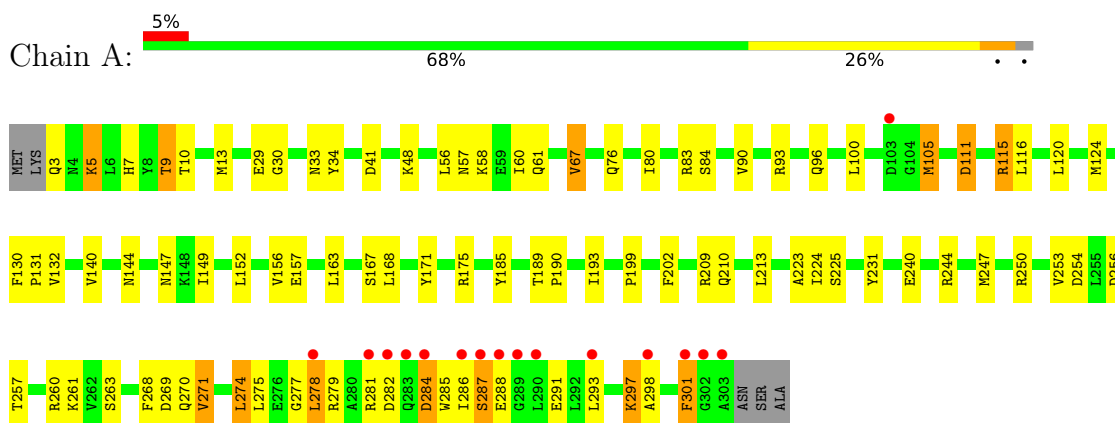
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	143	Total O 143 143	0	0
4	B	148	Total O 148 148	0	0

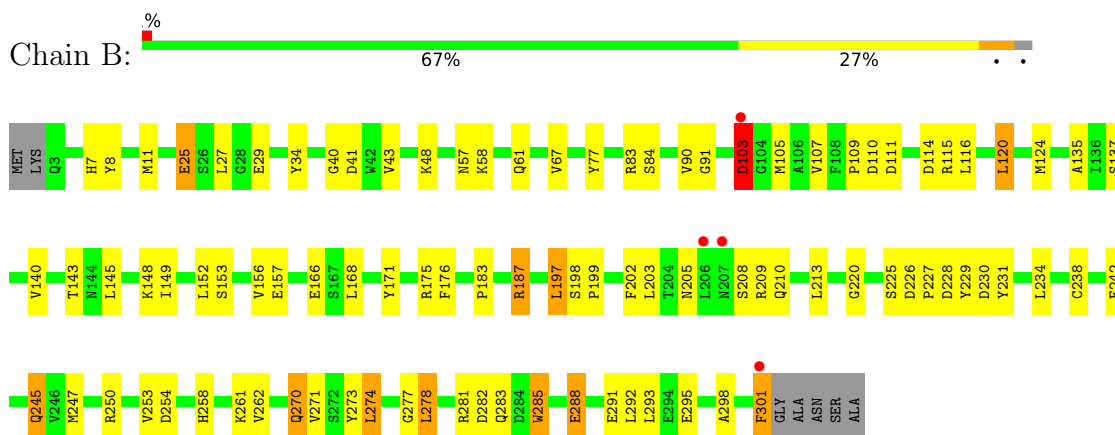
### 3 Residue-property plots [i](#)

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: Gentamicin resistance protein



- Molecule 1: Gentamicin resistance protein



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	41.68Å 54.65Å 77.16Å 89.94° 103.83° 106.84°	Depositor
Resolution (Å)	38.45 – 2.20 38.45 – 2.14	Depositor EDS
% Data completeness (in resolution range)	95.2 (38.45-2.20) 94.4 (38.45-2.14)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.11 (at 2.14Å)	Xtrriage
Refinement program	PHENIX (phenix.refine: 1.7.1_743)	Depositor
R, $R_{free}$	0.204 , 0.291 0.199 , 0.282	Depositor DCC
$R_{free}$ test set	1672 reflections (5.11%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	22.4	Xtrriage
Anisotropy	0.624	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 44.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	5185	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	29.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.22% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

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

## 5 Model quality

### 5.1 Standard geometry

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

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	0.47	0/2472	0.59	0/3357
1	B	0.47	0/2483	0.56	0/3369
All	All	0.47	0/4955	0.57	0/6726

There are no bond length outliers.

There are no bond angle outliers.

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	2413	0	2360	63	0
1	B	2419	0	2376	58	1
2	A	28	0	12	2	0
2	B	28	0	12	0	0
3	A	3	0	0	0	0
3	B	3	0	0	0	0
4	A	143	0	0	3	0
4	B	148	0	0	5	0
All	All	5185	0	4760	121	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 13.

All (121) 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:111:ASP:O	4:B:570:HOH:O	1.96	0.84
1:A:111:ASP:OD1	1:A:111:ASP:N	2.15	0.80
1:B:183:PRO:HB2	1:B:187:ARG:HH21	1.46	0.80
1:A:93:ARG:HG3	1:A:93:ARG:HH11	1.47	0.78
1:B:124:MET:HE2	1:B:250:ARG:HH11	1.48	0.77
1:A:271:VAL:HG23	1:A:293:LEU:HD11	1.68	0.74
1:A:284:ASP:OD1	1:A:287:SER:OG	2.06	0.72
1:B:225:SER:OG	1:B:226:ASP:N	2.26	0.69
1:A:76:GLN:OE1	4:A:524:HOH:O	2.11	0.67
1:B:270:GLN:CG	1:B:292:LEU:HB3	2.24	0.67
1:B:27:LEU:O	4:B:548:HOH:O	2.12	0.66
1:B:228:ASP:OD2	1:B:261:LYS:HE2	1.96	0.64
1:A:284:ASP:C	1:A:286:ILE:H	2.01	0.64
1:A:3:GLN:N	1:A:7:HIS:HD1	1.95	0.63
1:A:96:GLN:OE1	4:A:525:HOH:O	2.16	0.62
1:A:100:LEU:HD21	1:A:105:MET:HG2	1.80	0.62
1:A:124:MET:HE1	1:A:250:ARG:HH11	1.64	0.61
1:B:153:SER:HB3	1:B:176:PHE:CD1	2.36	0.60
1:B:171:TYR:O	1:B:175:ARG:HG2	2.02	0.59
1:A:240:GLU:O	1:A:244:ARG:HG3	2.03	0.58
1:B:143:THR:HB	1:B:225:SER:HB2	1.84	0.58
1:B:274:LEU:HD22	1:B:278:LEU:HD22	1.86	0.58
1:A:30:GLY:O	1:A:48:LYS:NZ	2.36	0.58
1:A:7:HIS:HE1	1:A:83:ARG:NH2	2.02	0.57
1:A:132:VAL:HG13	1:A:224:ILE:HD12	1.86	0.57
1:A:284:ASP:O	1:A:288:GLU:HB2	2.05	0.57
1:B:115:ARG:NH2	1:B:210:GLN:O	2.39	0.56
1:A:93:ARG:HG3	1:A:93:ARG:NH1	2.18	0.56
1:B:8:TYR:CE2	1:B:34:TYR:CD2	2.94	0.55
1:B:270:GLN:HG2	1:B:292:LEU:HB3	1.88	0.55
1:B:116:LEU:CD1	1:B:213:LEU:HB2	2.36	0.55
1:A:231:TYR:OH	1:A:250:ARG:HD3	2.07	0.55
1:A:282:ASP:HB3	1:A:285:TRP:HB3	1.90	0.54
1:B:205:ASN:ND2	1:B:208:SER:HB2	2.23	0.54
1:B:282:ASP:HB3	1:B:285:TRP:HB2	1.88	0.54
1:A:7:HIS:CE1	1:A:83:ARG:NH2	2.77	0.52
1:B:43:VAL:O	1:B:91:GLY:HA2	2.08	0.52
1:A:199:PRO:HA	1:A:202:PHE:CD2	2.44	0.52
1:B:124:MET:HE2	1:B:250:ARG:NH1	2.19	0.52
1:B:220:GLY:HA3	4:B:522:HOH:O	2.08	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:171:TYR:O	1:A:175:ARG:HG2	2.09	0.52
1:B:29:GLU:OE2	1:B:48:LYS:NZ	2.42	0.52
1:B:274:LEU:HD22	1:B:278:LEU:CD2	2.40	0.51
1:A:140:VAL:CG1	1:A:223:ALA:HA	2.41	0.50
1:A:9:THR:O	1:A:13:MET:HG3	2.11	0.50
1:A:288:GLU:HA	1:A:291:GLU:OE1	2.10	0.50
1:A:33:ASN:N	2:A:500:GDP:O1B	2.45	0.50
1:B:231:TYR:CZ	1:B:247:MET:HG2	2.46	0.50
1:B:103:ASP:OD2	1:B:103:ASP:N	2.45	0.50
1:B:242:PHE:O	1:B:245:GLN:HB2	2.12	0.50
1:A:244:ARG:NH2	1:A:256:ASP:OD1	2.34	0.49
1:B:231:TYR:OH	1:B:250:ARG:HD3	2.13	0.49
1:A:100:LEU:CD2	1:A:105:MET:HG2	2.43	0.49
1:A:124:MET:HE1	1:A:250:ARG:NH1	2.27	0.49
1:A:185:TYR:O	1:A:261:LYS:HE3	2.13	0.49
1:B:183:PRO:HB2	1:B:187:ARG:NH2	2.24	0.48
1:B:199:PRO:HA	1:B:202:PHE:CD2	2.48	0.48
1:A:152:LEU:O	1:A:156:VAL:HG23	2.14	0.48
1:B:120:LEU:O	1:B:124:MET:HG3	2.14	0.48
1:B:277:GLY:HA3	1:B:285:TRP:HB3	1.96	0.48
1:A:30:GLY:HA3	2:A:500:GDP:O2B	2.14	0.47
1:B:8:TYR:CE2	1:B:34:TYR:HD2	2.32	0.47
1:A:231:TYR:CZ	1:A:247:MET:HG2	2.51	0.46
1:B:152:LEU:O	1:B:156:VAL:HG23	2.15	0.46
1:B:145:LEU:O	1:B:149:ILE:HG12	2.15	0.46
1:B:137[B]:SER:OG	4:B:651:HOH:O	2.20	0.46
1:A:29:GLU:HG3	1:A:34:TYR:HD1	1.80	0.46
1:A:93:ARG:NH1	4:A:606:HOH:O	2.49	0.46
1:A:149:ILE:HD12	1:A:268:PHE:CD2	2.50	0.46
1:B:25[B]:GLU:HA	1:B:25[B]:GLU:OE1	2.14	0.46
1:A:67:VAL:HG11	1:A:76:GLN:CD	2.36	0.45
1:A:171:TYR:CG	1:A:297:LYS:HG3	2.51	0.45
1:B:135:ALA:HB1	1:B:140:VAL:HG21	1.96	0.45
1:A:284:ASP:C	1:A:286:ILE:N	2.68	0.45
1:B:273:TYR:CE2	1:B:288:GLU:HG2	2.51	0.45
1:A:9:THR:HG22	1:A:10:THR:N	2.30	0.45
1:A:277:GLY:HA3	1:A:285:TRP:O	2.17	0.45
1:B:57:ASN:O	1:B:61:GLN:HG3	2.17	0.45
1:B:253:VAL:HG23	1:B:254:ASP:CG	2.38	0.44
1:A:116:LEU:CD1	1:A:213:LEU:HB2	2.47	0.44
1:A:29:GLU:HG2	1:A:48:LYS:NZ	2.33	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:57:ASN:O	1:A:61:GLN:HG3	2.18	0.44
1:A:93:ARG:NH1	1:A:93:ARG:CG	2.80	0.44
1:A:274:LEU:HD12	1:A:293:LEU:HD12	1.99	0.44
1:B:7:HIS:O	1:B:11[A]:MET:HG3	2.17	0.44
1:B:40:GLY:HA3	4:B:542:HOH:O	2.17	0.44
1:B:153:SER:O	1:B:157:GLU:HG3	2.17	0.44
1:B:58:LYS:HE2	1:B:58:LYS:HB3	1.65	0.44
1:B:197:LEU:HD12	1:B:197:LEU:HA	1.83	0.44
1:A:144:ASN:HD21	1:A:147:ASN:HB2	1.83	0.44
1:A:56:LEU:O	1:A:60:ILE:HG13	2.18	0.43
1:A:275:LEU:O	1:A:279:ARG:HD2	2.19	0.43
1:B:258:HIS:O	1:B:262:VAL:HG23	2.19	0.43
1:A:130:PHE:HA	1:A:131:PRO:HD3	1.87	0.43
1:A:144:ASN:ND2	1:A:147:ASN:HB2	2.34	0.43
1:A:193:ILE:HD13	1:A:225:SER:HB3	2.00	0.42
1:A:58:LYS:HB3	1:A:58:LYS:HE2	1.81	0.42
1:A:254[A]:ASP:OD2	1:A:257:THR:OG1	2.26	0.42
1:B:203:LEU:HD13	1:B:203:LEU:HA	1.91	0.42
1:B:277:GLY:CA	1:B:285:TRP:HB3	2.50	0.42
1:A:278:LEU:HD12	1:A:286:ILE:CD1	2.49	0.42
1:B:120:LEU:HD13	1:B:213:LEU:HD22	2.00	0.42
1:B:226:ASP:HA	1:B:227:PRO:HD2	1.94	0.42
1:A:80:ILE:HG12	1:A:90:VAL:HG12	2.02	0.42
1:A:29:GLU:HG3	1:A:34:TYR:CD1	2.55	0.41
1:A:298:ALA:O	1:A:301:PHE:HB2	2.20	0.41
1:B:234:LEU:O	1:B:238:CYS:HB2	2.20	0.41
1:B:203:LEU:HD22	1:B:203:LEU:N	2.35	0.41
1:A:140:VAL:HG12	1:A:223:ALA:HA	2.01	0.41
1:A:260:ARG:O	1:A:263:SER:HB3	2.20	0.41
1:B:7:HIS:NE2	1:B:83:ARG:NH2	2.68	0.41
1:A:115:ARG:HA	1:A:115:ARG:HD2	1.31	0.41
1:B:77:TYR:CD2	1:B:90:VAL:HG11	2.55	0.41
1:B:120:LEU:HD12	1:B:120:LEU:HA	1.85	0.41
1:B:298:ALA:O	1:B:301[A]:PHE:HB3	2.21	0.41
1:B:271:VAL:HG23	1:B:293:LEU:HD11	2.03	0.40
1:A:5:LYS:HD3	1:A:5:LYS:HA	1.78	0.40
1:A:144:ASN:ND2	1:A:147:ASN:CB	2.84	0.40
1:A:189:THR:HA	1:A:190:PRO:HD3	1.93	0.40
1:B:115:ARG:HD2	1:B:115:ARG:HA	1.93	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)
1:B:166:GLU:OE1	1:B:209:ARG:NH2[1_545]	2.11	0.09

### 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	301/306 (98%)	286 (95%)	14 (5%)	1 (0%)	41	46
1	B	300/306 (98%)	283 (94%)	14 (5%)	3 (1%)	15	14
All	All	601/612 (98%)	569 (95%)	28 (5%)	4 (1%)	22	22

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	301	PHE
1	B	103	ASP
1	B	197	LEU
1	B	109	PRO

#### 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	266/269 (99%)	241 (91%)	25 (9%)	8	8
1	B	268/269 (100%)	239 (89%)	29 (11%)	6	6
All	All	534/538 (99%)	480 (90%)	54 (10%)	8	7

All (54) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	LYS
1	A	9	THR
1	A	41	ASP
1	A	67	VAL
1	A	84	SER
1	A	105	MET
1	A	111	ASP
1	A	115	ARG
1	A	120	LEU
1	A	157	GLU
1	A	163	LEU
1	A	167	SER
1	A	168	LEU
1	A	209	ARG
1	A	210	GLN
1	A	253	VAL
1	A	269	ASP
1	A	270	GLN
1	A	271	VAL
1	A	274	LEU
1	A	278	LEU
1	A	281	ARG
1	A	284	ASP
1	A	287	SER
1	A	297	LYS
1	B	25[A]	GLU
1	B	25[B]	GLU
1	B	41	ASP
1	B	67	VAL
1	B	84	SER
1	B	103	ASP
1	B	105	MET
1	B	107	VAL
1	B	110	ASP
1	B	114	ASP
1	B	120	LEU
1	B	148	LYS
1	B	168	LEU
1	B	187	ARG
1	B	198	SER
1	B	229	TYR
1	B	230	ASP

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Mol	Chain	Res	Type
1	B	245	GLN
1	B	270	GLN
1	B	274	LEU
1	B	278	LEU
1	B	281	ARG
1	B	283	GLN
1	B	285	TRP
1	B	288	GLU
1	B	291	GLU
1	B	295	GLU
1	B	301[A]	PHE
1	B	301[B]	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	144	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 6 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	GDP	B	500	3	24,30,30	1.05	2 (8%)	30,47,47	1.06	1 (3%)
2	GDP	A	500	3	24,30,30	1.03	2 (8%)	30,47,47	1.13	3 (10%)

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	GDP	B	500	3	-	0/12/32/32	0/3/3/3
2	GDP	A	500	3	-	0/12/32/32	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	500	GDP	O4'-C1'	2.42	1.44	1.41
2	A	500	GDP	O4'-C1'	2.26	1.44	1.41
2	B	500	GDP	C6-N1	-2.19	1.34	1.37
2	A	500	GDP	C6-N1	-2.14	1.34	1.37

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	500	GDP	PA-O3A-PB	-2.85	123.05	132.83
2	A	500	GDP	PA-O3A-PB	-2.61	123.87	132.83
2	A	500	GDP	C8-N7-C5	2.21	107.20	102.99
2	A	500	GDP	C5-C6-N1	2.17	117.79	113.95

There are no chirality outliers.

There are no torsion outliers.

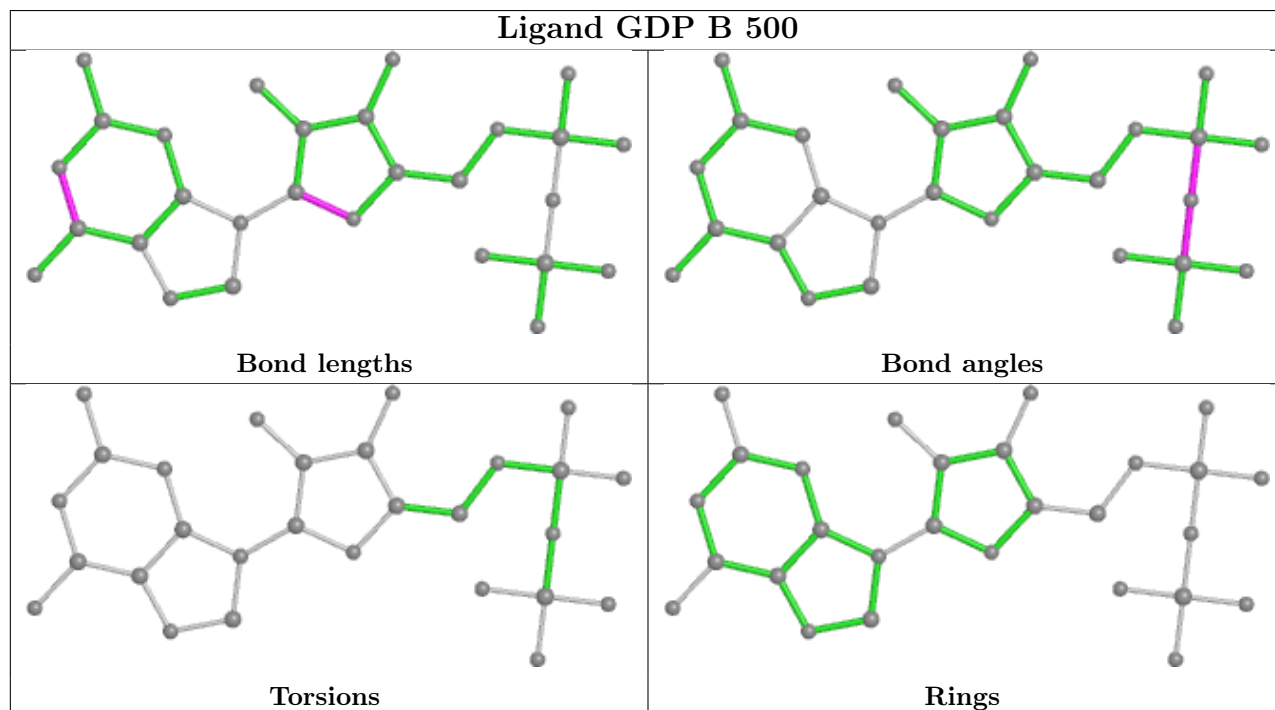
There are no ring outliers.

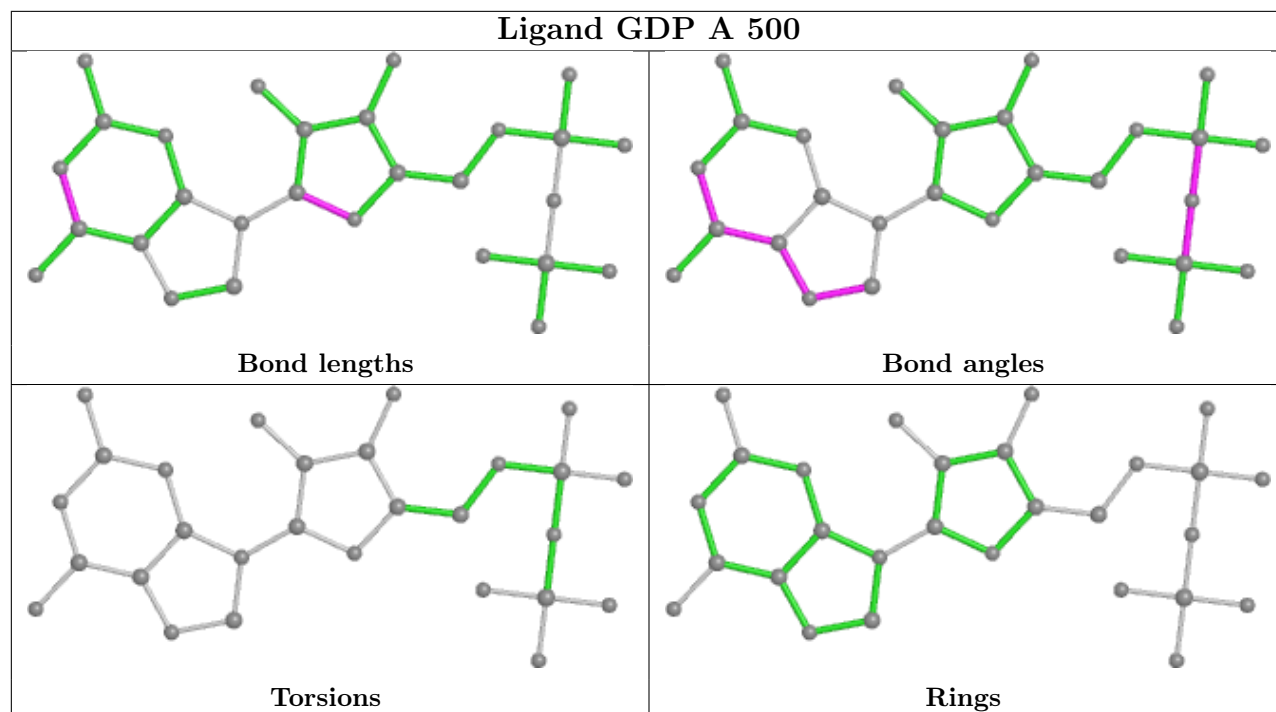
1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	500	GDP	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

### 6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	301/306 (98%)	-0.11	16 (5%) 26 25	16, 27, 60, 81	0
1	B	299/306 (97%)	-0.30	4 (1%) 77 75	16, 26, 49, 59	0
All	All	600/612 (98%)	-0.21	20 (3%) 46 44	16, 27, 53, 81	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	301	PHE	5.0
1	A	290	LEU	4.7
1	B	301[A]	PHE	3.3
1	A	288	GLU	3.3
1	B	103	ASP	3.2
1	A	281	ARG	3.0
1	A	278	LEU	2.9
1	A	293	LEU	2.8
1	B	206	LEU	2.8
1	A	103	ASP	2.5
1	A	282	ASP	2.5
1	A	287	SER	2.5
1	A	283	GLN	2.4
1	A	289	GLY	2.3
1	A	298	ALA	2.1
1	A	303	ALA	2.1
1	B	207	ASN	2.1
1	A	302	GLY	2.1
1	A	284	ASP	2.1
1	A	286	ILE	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 monosaccharides in this entry.

## 6.4 Ligands [i](#)

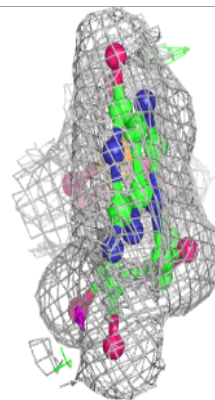
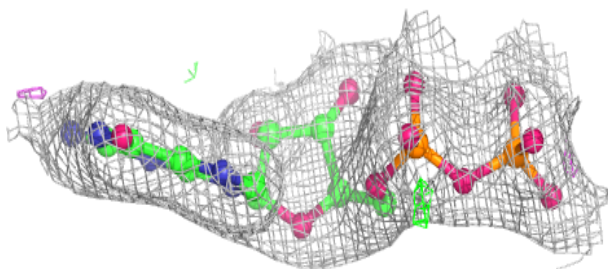
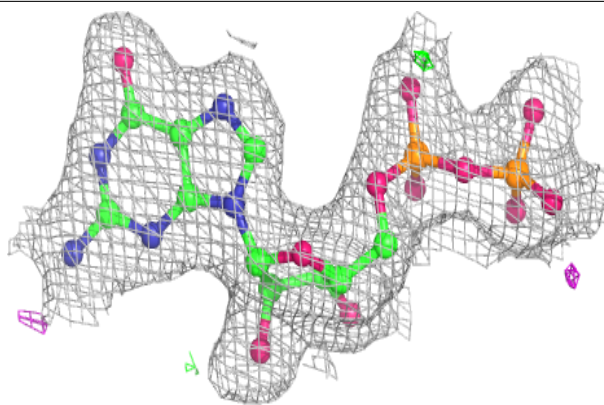
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	MG	A	503	1/1	0.89	0.07	27,27,27,27	0
3	MG	A	502	1/1	0.92	0.09	20,20,20,20	0
3	MG	B	502	1/1	0.95	0.08	16,16,16,16	0
3	MG	B	503	1/1	0.97	0.09	27,27,27,27	0
2	GDP	B	500	28/28	0.98	0.09	12,17,23,29	0
3	MG	B	501	1/1	0.98	0.03	18,18,18,18	0
3	MG	A	501	1/1	0.98	0.05	16,16,16,16	0
2	GDP	A	500	28/28	0.98	0.10	17,20,23,26	0

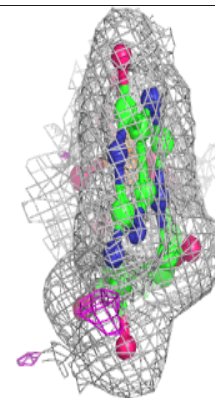
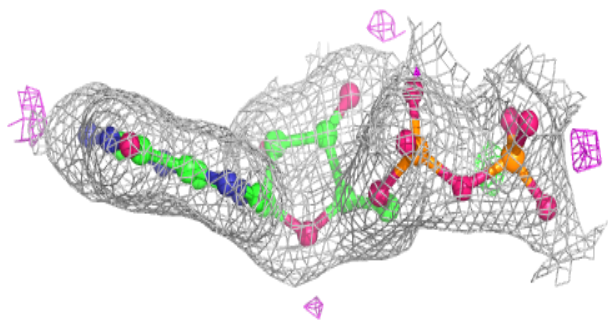
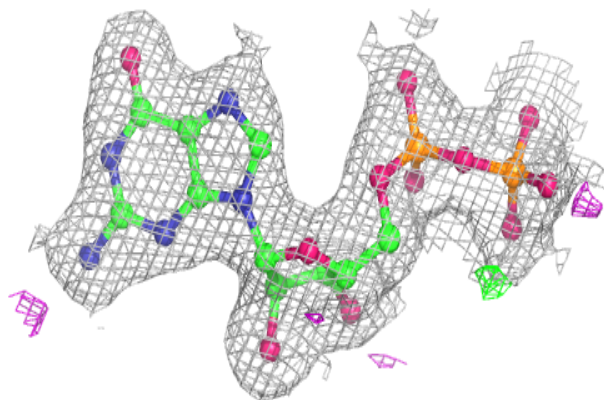
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

**Electron density around GDP B 500:**

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

**Electron density around GDP A 500:**

$2mF_o-DF_c$  (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.