



# Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 03:52 am BST

PDB ID : 3X08  
Title : Crystal structure of PIP4KIIBETA N203A complex with GMP  
Authors : Takeuchi, K.; Lo, Y.H.; Sumita, K.; Senda, M.; Terakawa, J.; Dimitoris, A.; Locasale, J.W.; Sasaki, M.; Yoshino, H.; Zhang, Y.; Kahoud, E.R.; Takano, T.; Yokota, T.; Emerling, B.; Asara, J.A.; Ishida, T.; Shimada, I.; Daikoku, T.; Cantley, L.C.; Senda, T.; Sasaki, A.T.  
Deposited on : 2014-10-09  
Resolution : 2.75 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.

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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)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
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) : 2.11

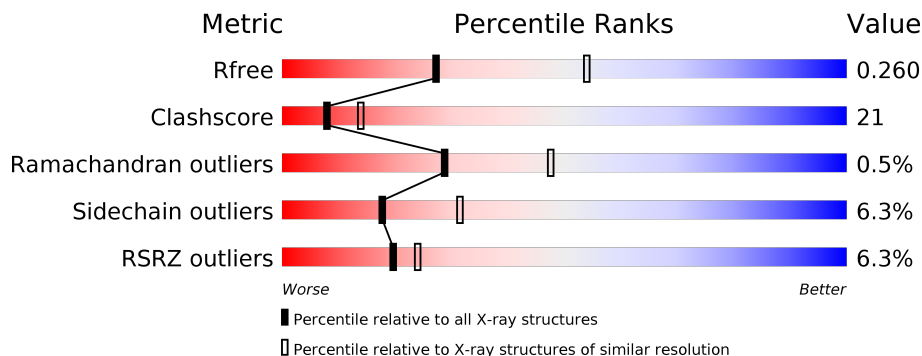
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.75 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	1235 (2.78-2.74)
Clashscore	141614	1277 (2.78-2.74)
Ramachandran outliers	138981	1257 (2.78-2.74)
Sidechain outliers	138945	1257 (2.78-2.74)
RSRZ outliers	127900	1207 (2.78-2.74)

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 on 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	393	 2% 55% 23% 19%
1	B	393	 8% 46% 28% 23%

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 5237 atoms, of which 36 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.

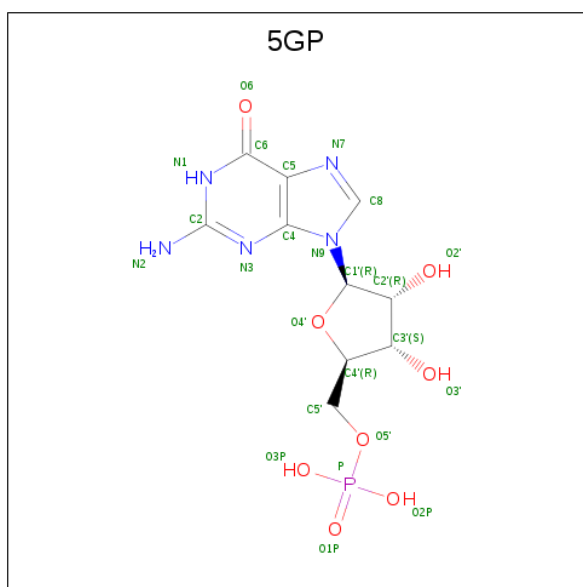
- Molecule 1 is a protein called Phosphatidylinositol 5-phosphate 4-kinase type-2 beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	319	2620	1668	446	492	14	0	0	0
1	B	303	2494	1598	425	458	13	0	0	0

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	24	GLY	-	EXPRESSION TAG	UNP P78356
A	25	PRO	-	EXPRESSION TAG	UNP P78356
A	26	ASN	-	EXPRESSION TAG	UNP P78356
A	27	CYS	-	EXPRESSION TAG	UNP P78356
A	28	ALA	-	EXPRESSION TAG	UNP P78356
A	29	PRO	-	EXPRESSION TAG	UNP P78356
A	30	GLY	-	EXPRESSION TAG	UNP P78356
A	203	ALA	ASN	ENGINEERED MUTATION	UNP P78356
B	24	GLY	-	EXPRESSION TAG	UNP P78356
B	25	PRO	-	EXPRESSION TAG	UNP P78356
B	26	ASN	-	EXPRESSION TAG	UNP P78356
B	27	CYS	-	EXPRESSION TAG	UNP P78356
B	28	ALA	-	EXPRESSION TAG	UNP P78356
B	29	PRO	-	EXPRESSION TAG	UNP P78356
B	30	GLY	-	EXPRESSION TAG	UNP P78356
B	203	ALA	ASN	ENGINEERED MUTATION	UNP P78356

- Molecule 2 is GUANOSINE-5'-MONOPHOSPHATE (three-letter code: 5GP) (formula: C<sub>10</sub>H<sub>14</sub>N<sub>5</sub>O<sub>8</sub>P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
			Total	C	H	N	O			P
2	A	1	Total	C	H	N	O	P	0	0
			36	10	12	5	8	1		
2	A	1	Total	C	H	N	O	P	0	0
			36	10	12	5	8	1		
2	A	1	Total	C	H	N	O	P	0	0
			36	10	12	5	8	1		

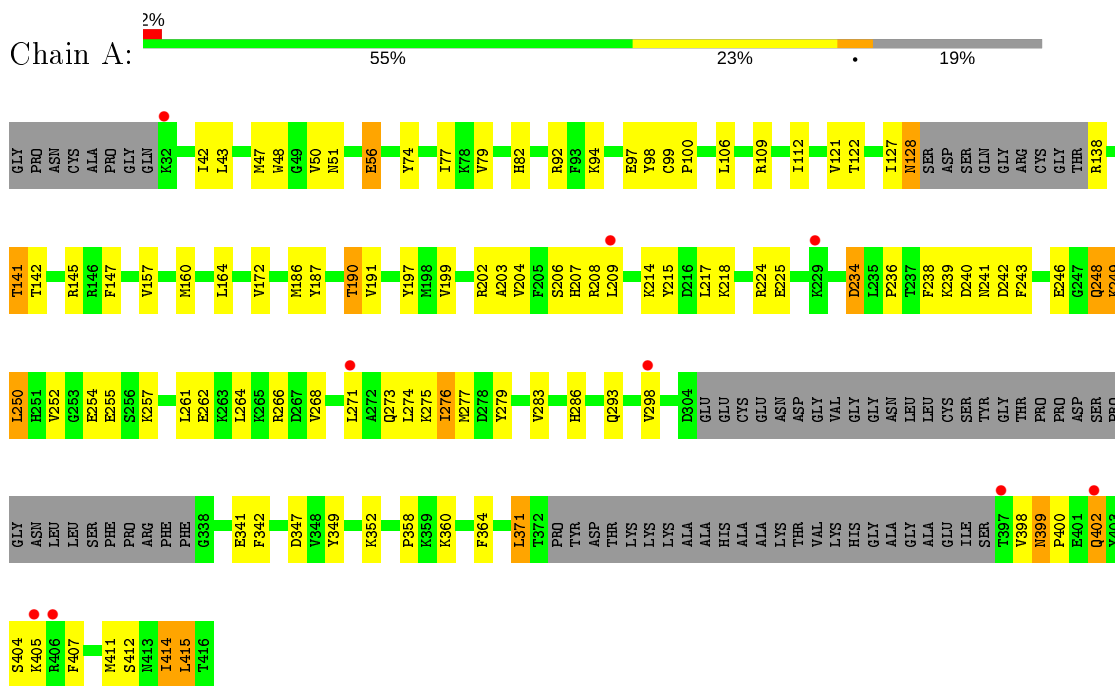
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	7	Total 7 7	0	0
3	B	8	Total 8 8	0	0

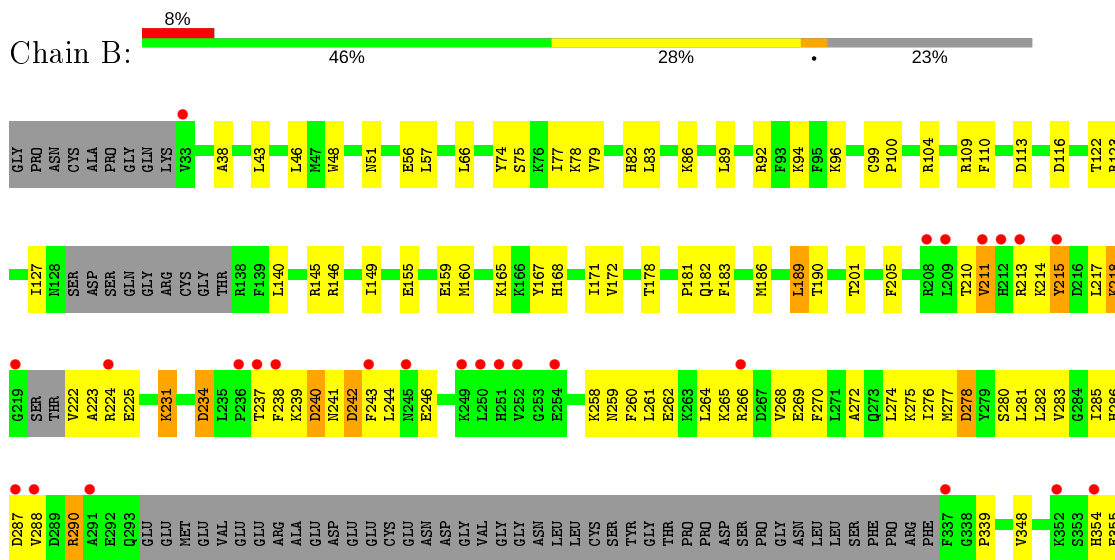
### 3 Residue-property plots [i](#)

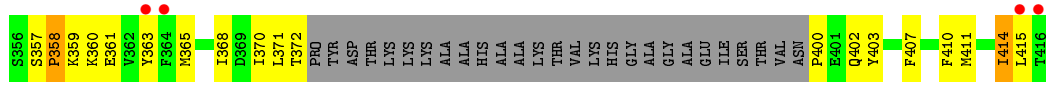
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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: Phosphatidylinositol 5-phosphate 4-kinase type-2 beta



- Molecule 1: Phosphatidylinositol 5-phosphate 4-kinase type-2 beta





## 4 Data and refinement statistics i

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	107.38Å 182.34Å 107.51Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	91.17 - 2.75 92.53 - 2.75	Depositor EDS
% Data completeness (in resolution range)	98.8 (91.17-2.75) 98.7 (92.53-2.75)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.12 (at 2.73Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE: 1.7.3_928)	Depositor
R, $R_{free}$	0.206 , 0.267 0.197 , 0.260	Depositor DCC
$R_{free}$ test set	1374 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	72.0	Xtriage
Anisotropy	0.273	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 74.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.016 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.020 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	5237	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	87.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.66% 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.49	0/2673	0.57	0/3596
1	B	0.42	0/2547	0.55	0/3424
All	All	0.46	0/5220	0.56	0/7020

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 [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2620	0	2595	102	0
1	B	2494	0	2481	127	0
2	A	72	36	36	2	0
3	A	7	0	0	0	0
3	B	8	0	0	1	0
All	All	5201	36	5112	221	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 21.

All (221) 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:213:ARG:HG2	1:B:215:TYR:CE1	1.80	1.17
1:A:77:ILE:HD11	1:B:77:ILE:HD11	1.27	1.15
1:B:213:ARG:HG2	1:B:215:TYR:HE1	0.93	1.09
1:B:78:LYS:HD2	1:B:94:LYS:HE2	1.35	1.07
1:B:215:TYR:CE2	1:B:243:PHE:HD1	1.74	1.06
1:B:213:ARG:HD3	1:B:215:TYR:OH	1.58	1.04
1:A:77:ILE:HD11	1:B:77:ILE:CD1	1.88	1.03
1:A:77:ILE:CD1	1:B:77:ILE:HD11	1.98	0.93
1:A:399:ASN:ND2	1:A:400:PRO:HD2	1.85	0.92
1:A:277:MET:HE1	1:A:398:VAL:HG23	1.51	0.92
1:B:215:TYR:CE2	1:B:243:PHE:CD1	2.60	0.89
1:A:217:LEU:HD13	1:A:414:ILE:HD11	1.55	0.87
1:A:273:GLN:C	1:A:274:LEU:HD23	1.97	0.85
1:B:217:LEU:HB2	1:B:411:MET:HE1	1.56	0.84
1:A:160:MET:HG2	1:A:371:LEU:HD21	1.60	0.83
1:B:214:LYS:C	1:B:215:TYR:HD1	1.83	0.83
1:B:215:TYR:HE2	1:B:243:PHE:HD1	1.24	0.81
1:A:42:ILE:HD13	1:A:191:VAL:HG21	1.61	0.81
1:B:214:LYS:C	1:B:215:TYR:CD1	2.56	0.79
1:A:207:HIS:CE1	1:A:208:ARG:HG3	2.19	0.78
1:B:218:LYS:N	1:B:218:LYS:CD	2.44	0.78
1:B:260:PHE:CD2	1:B:415:LEU:HD11	2.18	0.78
1:B:214:LYS:O	1:B:215:TYR:CD1	2.39	0.75
1:A:399:ASN:HD22	1:A:400:PRO:HD2	1.49	0.74
1:A:225:GLU:HG3	1:A:241:ASN:HB2	1.70	0.74
1:B:225:GLU:HA	1:B:242:ASP:OD1	1.86	0.74
1:A:141:THR:HG23	1:A:142:THR:O	1.86	0.74
1:B:371:LEU:O	1:B:372:THR:HG23	1.87	0.74
1:B:215:TYR:CD2	1:B:243:PHE:CD1	2.75	0.73
1:B:283:VAL:HG22	1:B:365:MET:HG2	1.69	0.73
1:A:217:LEU:CD1	1:A:414:ILE:HD11	2.18	0.73
1:A:224:ARG:NH1	1:A:240:ASP:OD1	2.22	0.72
1:A:42:ILE:HD13	1:A:191:VAL:CG2	2.19	0.72
1:B:277:MET:O	1:B:278:ASP:HB2	1.89	0.71
1:A:274:LEU:HD23	1:A:274:LEU:N	2.05	0.71
1:B:243:PHE:HE2	1:B:414:ILE:CG2	2.05	0.70
1:A:264:LEU:HD11	1:A:407:PHE:HE1	1.56	0.70
1:A:274:LEU:HB2	1:A:276:ILE:HG13	1.74	0.69
1:A:249:LYS:C	1:A:250:LEU:HD23	2.13	0.69
1:B:243:PHE:CE2	1:B:414:ILE:CG2	2.75	0.69
1:B:272:ALA:O	1:B:275:LYS:N	2.20	0.69
1:B:242:ASP:OD1	1:B:242:ASP:N	2.26	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:109:ARG:CZ	1:B:172:VAL:HG22	2.24	0.68
1:A:97:GLU:HG3	1:A:187:TYR:CE1	2.29	0.68
1:B:243:PHE:HE2	1:B:414:ILE:HG23	1.57	0.67
1:A:264:LEU:O	1:A:268:VAL:HG22	1.95	0.67
2:A:503:5GP:H5'1	2:A:503:5GP:H8	1.78	0.66
1:B:288:VAL:HG23	1:B:360:LYS:O	1.95	0.66
1:A:274:LEU:HB2	1:A:276:ILE:CG1	2.26	0.66
1:B:213:ARG:CG	1:B:215:TYR:CE1	2.70	0.66
1:A:218:LYS:HD2	1:A:224:ARG:NE	2.11	0.65
1:B:282:LEU:HB2	1:B:368:ILE:HD13	1.77	0.65
1:B:361:GLU:HG3	1:B:363:TYR:CZ	2.32	0.65
1:B:218:LYS:N	1:B:218:LYS:HD2	2.11	0.65
1:B:213:ARG:NH1	1:B:246:GLU:OE2	2.31	0.64
1:A:341:GLU:HG3	1:A:342:PHE:H	1.64	0.63
1:B:217:LEU:CB	1:B:411:MET:HE1	2.28	0.62
1:B:78:LYS:HD2	1:B:94:LYS:CE	2.21	0.62
1:B:361:GLU:HG3	1:B:363:TYR:CE1	2.35	0.61
1:B:43:LEU:HD21	1:B:140:LEU:HD11	1.82	0.61
1:A:186:MET:HG3	1:A:199:VAL:HG22	1.82	0.61
1:A:234:ASP:N	1:A:234:ASP:OD1	2.24	0.59
1:A:236:PRO:HB2	1:A:238:PHE:CE1	2.36	0.59
1:A:202:ARG:HG3	1:A:203:ALA:N	2.17	0.59
1:B:400:PRO:C	1:B:402:GLN:H	2.07	0.58
1:A:217:LEU:CD1	1:A:414:ILE:CD1	2.80	0.58
1:A:208:ARG:NH2	1:A:341:GLU:O	2.37	0.57
1:B:238:PHE:HD2	1:B:242:ASP:HB2	1.69	0.57
1:A:248:GLN:HG3	1:A:249:LYS:N	2.19	0.57
1:B:225:GLU:HG2	1:B:242:ASP:OD1	2.05	0.57
1:A:56:GLU:OE2	1:B:48:TRP:NE1	2.31	0.57
1:A:234:ASP:O	1:A:236:PRO:HD3	2.05	0.56
1:B:234:ASP:OD1	1:B:234:ASP:N	2.30	0.56
1:B:240:ASP:N	1:B:240:ASP:OD1	2.32	0.56
1:B:189:LEU:N	1:B:189:LEU:HD12	2.19	0.56
1:B:222:VAL:HG12	1:B:223:ALA:H	1.70	0.56
1:B:183:PHE:CE2	1:B:370:ILE:HD11	2.41	0.56
1:B:94:LYS:HB2	1:B:190:THR:HB	1.88	0.56
1:A:271:LEU:HD13	1:A:279:TYR:CD2	2.41	0.55
1:A:411:MET:O	1:A:414:ILE:HB	2.06	0.55
1:B:261:LEU:HD21	1:B:415:LEU:HD12	1.88	0.55
1:A:271:LEU:HD13	1:A:279:TYR:CE2	2.41	0.55
1:A:43:LEU:O	1:A:47:MET:HG3	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:225:GLU:O	1:A:239:LYS:HG3	2.07	0.55
1:A:127:ILE:HG13	1:A:141:THR:HG22	1.89	0.55
1:B:167:TYR:O	1:B:171:ILE:HG12	2.07	0.55
1:B:258:LYS:O	1:B:262:GLU:HG2	2.07	0.55
1:A:261:LEU:HD21	1:A:412:SER:HA	1.90	0.54
1:A:74:TYR:CZ	1:B:82:HIS:HB2	2.43	0.54
1:A:109:ARG:CZ	1:A:172:VAL:HG22	2.37	0.54
1:B:264:LEU:O	1:B:268:VAL:HG22	2.08	0.54
1:B:400:PRO:C	1:B:402:GLN:N	2.61	0.54
1:A:257:LYS:HE2	1:A:261:LEU:HD11	1.89	0.54
1:A:127:ILE:CG1	1:A:141:THR:HG22	2.38	0.53
1:A:273:GLN:O	1:A:274:LEU:HD23	2.08	0.53
1:B:243:PHE:CE2	1:B:414:ILE:HG21	2.42	0.53
1:A:99:CYS:N	1:A:100:PRO:HD3	2.24	0.53
1:B:189:LEU:N	1:B:189:LEU:CD1	2.72	0.53
1:B:210:THR:O	1:B:286:HIS:ND1	2.40	0.52
1:B:217:LEU:C	1:B:218:LYS:HD2	2.29	0.52
1:B:287:ASP:CG	1:B:359:LYS:HE2	2.30	0.52
1:A:264:LEU:HD11	1:A:407:PHE:CE1	2.42	0.51
1:B:218:LYS:N	1:B:218:LYS:HD3	2.23	0.51
1:A:48:TRP:NE1	1:B:56:GLU:OE2	2.40	0.51
1:B:260:PHE:HD2	1:B:415:LEU:HD11	1.72	0.51
1:B:400:PRO:HB2	1:B:402:GLN:HG3	1.92	0.50
1:A:79:VAL:O	1:A:92:ARG:HA	2.11	0.50
1:A:128:ASN:OD1	1:A:128:ASN:N	2.43	0.50
1:A:248:GLN:O	1:A:249:LYS:HD2	2.12	0.50
1:B:104:ARG:HD2	3:B:506:HOH:O	2.10	0.50
1:B:110:PHE:CE1	1:B:182:GLN:HG2	2.47	0.50
1:A:407:PHE:O	1:A:411:MET:HG2	2.11	0.50
1:B:285:ILE:HG23	1:B:363:TYR:CE1	2.47	0.49
1:A:145:ARG:NH2	1:A:207:HIS:O	2.45	0.49
1:A:275:LYS:O	1:A:400:PRO:HG3	2.13	0.49
1:A:293:GLN:NE2	2:A:502:5GP:C6	2.76	0.49
1:B:160:MET:HG2	1:B:186:MET:SD	2.52	0.49
1:A:277:MET:CE	1:A:398:VAL:HG23	2.33	0.49
1:B:51:ASN:HB2	1:B:122:THR:HG21	1.95	0.49
1:B:218:LYS:H	1:B:218:LYS:HD3	1.78	0.48
1:A:160:MET:CG	1:A:371:LEU:HD21	2.37	0.48
1:A:98:TYR:O	1:A:99:CYS:C	2.52	0.48
1:A:399:ASN:O	1:A:402:GLN:CG	2.62	0.48
1:B:281:LEU:HB2	1:B:407:PHE:CZ	2.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:348:VAL:O	1:B:348:VAL:CG1	2.62	0.48
1:B:46:LEU:HD21	1:B:149:ILE:HD13	1.96	0.47
1:B:259:ASN:O	1:B:262:GLU:HG2	2.15	0.47
1:B:276:ILE:HG22	1:B:277:MET:N	2.29	0.47
1:A:242:ASP:O	1:A:246:GLU:HG3	2.14	0.47
1:B:213:ARG:NH1	1:B:246:GLU:OE1	2.46	0.47
1:A:214:LYS:HA	1:A:283:VAL:O	2.15	0.47
1:A:341:GLU:HG3	1:A:342:PHE:N	2.28	0.47
1:A:50:VAL:HG21	1:A:121:VAL:HG11	1.96	0.47
1:A:74:TYR:CE2	1:B:82:HIS:CB	2.97	0.47
1:A:225:GLU:C	1:A:239:LYS:HG3	2.35	0.47
1:B:171:ILE:N	1:B:171:ILE:HD13	2.30	0.47
1:B:242:ASP:O	1:B:246:GLU:HB2	2.15	0.47
1:B:268:VAL:HG11	1:B:407:PHE:HD2	1.80	0.47
1:B:79:VAL:O	1:B:92:ARG:HA	2.14	0.47
1:B:400:PRO:HG2	1:B:403:TYR:H	1.80	0.47
1:A:157:VAL:HG21	1:A:197:TYR:CD1	2.50	0.47
1:A:207:HIS:HD2	1:A:347:ASP:OD1	1.98	0.47
1:A:82:HIS:HB2	1:B:74:TYR:CZ	2.50	0.47
1:A:207:HIS:CE1	1:A:208:ARG:CG	2.95	0.47
1:B:213:ARG:HH12	1:B:246:GLU:CD	2.19	0.47
1:B:266:ARG:HE	1:B:266:ARG:HB3	1.49	0.47
1:B:411:MET:HA	1:B:414:ILE:HD11	1.96	0.47
1:B:159:GLU:HA	1:B:159:GLU:OE2	2.15	0.46
1:A:352:LYS:HE2	1:A:360:LYS:NZ	2.31	0.46
1:B:231:LYS:O	1:B:231:LYS:HD2	2.15	0.46
1:B:38:ALA:HB1	1:B:43:LEU:HB2	1.97	0.46
1:B:182:GLN:O	1:B:201:THR:HA	2.15	0.46
1:A:112:ILE:HD13	1:A:147:PHE:HE1	1.80	0.46
1:A:250:LEU:HD23	1:A:250:LEU:N	2.29	0.46
1:A:257:LYS:HG3	1:A:415:LEU:HG	1.97	0.46
1:B:348:VAL:O	1:B:348:VAL:HG12	2.16	0.46
1:B:270:PHE:CZ	1:B:274:LEU:HD11	2.51	0.46
1:A:215:TYR:HB3	1:A:217:LEU:HD21	1.97	0.46
1:A:160:MET:HG3	1:A:186:MET:SD	2.56	0.45
1:A:145:ARG:CZ	1:A:207:HIS:O	2.64	0.45
1:A:209:LEU:HB3	1:A:286:HIS:ND1	2.30	0.45
1:A:97:GLU:HG3	1:A:187:TYR:CD1	2.51	0.45
1:B:213:ARG:NH1	1:B:246:GLU:CD	2.70	0.45
1:B:246:GLU:O	1:B:246:GLU:HG2	2.16	0.45
1:B:66:LEU:HD12	1:B:168:HIS:CE1	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:249:LYS:O	1:A:250:LEU:HD23	2.15	0.45
1:B:214:LYS:HA	1:B:283:VAL:O	2.17	0.45
1:B:358:PRO:HB2	1:B:359:LYS:H	1.56	0.45
1:B:281:LEU:HB3	1:B:411:MET:HE1	1.97	0.45
1:B:165:LYS:HB2	1:B:165:LYS:HE2	1.69	0.45
1:B:75:SER:O	1:B:96:LYS:HA	2.17	0.45
1:A:77:ILE:HD11	1:B:77:ILE:HD13	1.88	0.45
1:A:341:GLU:CG	1:A:342:PHE:H	2.30	0.45
1:B:217:LEU:CB	1:B:411:MET:CE	2.94	0.44
1:B:214:LYS:HG3	1:B:237:THR:HA	1.99	0.44
1:A:51:ASN:HB2	1:A:122:THR:HG21	1.99	0.44
1:A:209:LEU:HB3	1:A:286:HIS:CE1	2.52	0.44
1:B:48:TRP:HH2	1:B:77:ILE:HD13	1.81	0.44
1:A:207:HIS:ND1	1:A:208:ARG:HG3	2.32	0.44
1:A:262:GLU:O	1:A:266:ARG:HG2	2.18	0.44
1:B:183:PHE:HE2	1:B:370:ILE:HD11	1.81	0.44
1:B:48:TRP:HB2	1:B:89:LEU:HD21	2.00	0.44
1:A:402:GLN:HG2	1:A:402:GLN:H	1.62	0.44
1:A:204:VAL:HA	1:A:349:TYR:CD2	2.53	0.43
1:B:181:PRO:HG2	1:B:370:ILE:HG12	1.99	0.43
1:B:287:ASP:OD2	1:B:290:ARG:HB2	2.17	0.43
1:B:205:PHE:CD2	1:B:211:VAL:HG21	2.53	0.43
1:B:285:ILE:HG23	1:B:363:TYR:HE1	1.84	0.43
1:B:83:LEU:HA	1:B:86:LYS:HD2	2.00	0.43
1:A:399:ASN:O	1:A:402:GLN:HG3	2.19	0.42
1:A:106:LEU:HA	1:A:106:LEU:HD23	1.74	0.42
1:A:94:LYS:HB2	1:A:190:THR:OG1	2.19	0.42
1:B:225:GLU:HG2	1:B:241:ASN:HB2	2.00	0.42
1:A:399:ASN:O	1:A:402:GLN:HG2	2.19	0.42
1:A:215:TYR:HB2	1:A:283:VAL:HB	2.02	0.42
1:B:357:SER:HA	1:B:358:PRO:HD3	1.83	0.42
1:B:218:LYS:HD3	1:B:240:ASP:OD2	2.20	0.42
1:B:281:LEU:HB2	1:B:407:PHE:HZ	1.85	0.42
1:A:243:PHE:CE2	1:A:248:GLN:HG2	2.55	0.42
1:A:399:ASN:ND2	1:A:400:PRO:CD	2.70	0.41
1:A:112:ILE:HD13	1:A:147:PHE:CE1	2.55	0.41
1:B:265:LYS:NZ	1:B:269:GLU:OE2	2.50	0.41
1:B:244:LEU:HD12	1:B:410:PHE:HE1	1.85	0.41
1:B:259:ASN:HA	1:B:262:GLU:HG2	2.03	0.41
1:B:371:LEU:O	1:B:372:THR:CG2	2.62	0.41
1:A:206:SER:HB2	1:A:364:PHE:CE2	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:99:CYS:N	1:B:100:PRO:CD	2.83	0.41
1:A:252:VAL:HG21	1:A:415:LEU:HD12	2.02	0.41
1:B:217:LEU:HD22	1:B:240:ASP:HA	2.03	0.41
1:B:288:VAL:HG12	1:B:339:PRO:HB2	2.02	0.41
1:A:371:LEU:HD12	1:A:371:LEU:HA	1.65	0.41
1:B:57:LEU:O	1:B:104:ARG:NH2	2.54	0.41
1:B:109:ARG:NH1	1:B:172:VAL:HG22	2.36	0.41
1:A:234:ASP:C	1:A:236:PRO:HD3	2.41	0.41
1:A:160:MET:CE	1:A:164:LEU:HD13	2.50	0.40
1:A:405:LYS:HE2	1:A:405:LYS:HB3	1.79	0.40
1:B:260:PHE:HD2	1:B:415:LEU:CD1	2.34	0.40
1:B:354:HIS:CD2	1:B:355:GLU:H	2.40	0.40
1:B:113:ASP:HB3	1:B:116:ASP:OD2	2.21	0.40
1:B:288:VAL:CG1	1:B:339:PRO:HB2	2.52	0.40
1:B:224:ARG:HB3	1:B:239:LYS:HD3	2.02	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	311/393 (79%)	294 (94%)	16 (5%)	1 (0%)	41	60
1	B	293/393 (75%)	272 (93%)	19 (6%)	2 (1%)	22	39
All	All	604/786 (77%)	566 (94%)	35 (6%)	3 (0%)	29	47

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	358	PRO
1	A	358	PRO
1	B	145	ARG

### 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	295/351 (84%)	276 (94%)	19 (6%)	17	31
1	B	280/351 (80%)	263 (94%)	17 (6%)	18	33
All	All	575/702 (82%)	539 (94%)	36 (6%)	18	31

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	56	GLU
1	A	128	ASN
1	A	138	ARG
1	A	141	THR
1	A	190	THR
1	A	234	ASP
1	A	248	GLN
1	A	249	LYS
1	A	250	LEU
1	A	254	GLU
1	A	255	GLU
1	A	276	ILE
1	A	298	VAL
1	A	371	LEU
1	A	399	ASN
1	A	402	GLN
1	A	404	SER
1	A	414	ILE
1	A	415	LEU
1	B	123	ARG
1	B	127	ILE
1	B	146	ARG
1	B	155	GLU
1	B	178	THR
1	B	189	LEU
1	B	211	VAL
1	B	215	TYR

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Mol	Chain	Res	Type
1	B	218	LYS
1	B	231	LYS
1	B	234	ASP
1	B	240	ASP
1	B	242	ASP
1	B	278	ASP
1	B	280	SER
1	B	290	ARG
1	B	414	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	207	HIS
1	A	293	GLN

### 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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

3 ligands are modelled in this entry.

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	5GP	A	501	-	22,26,26	0.65	0	27,40,40	1.82	6 (22%)
2	5GP	A	503	-	22,26,26	0.81	0	27,40,40	2.07	8 (29%)
2	5GP	A	502	-	22,26,26	0.64	0	27,40,40	1.73	7 (25%)

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	5GP	A	501	-	-	2/6/26/26	0/3/3/3
2	5GP	A	503	-	-	4/6/26/26	0/3/3/3
2	5GP	A	502	-	-	3/6/26/26	0/3/3/3

There are no bond length outliers.

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	503	5GP	C2-N3-C4	5.28	121.39	115.36
2	A	503	5GP	N3-C2-N1	-5.10	120.42	127.22
2	A	501	5GP	N3-C2-N1	-4.74	120.90	127.22
2	A	501	5GP	C2-N3-C4	4.72	120.75	115.36
2	A	502	5GP	C2-N3-C4	4.14	120.09	115.36
2	A	502	5GP	N3-C2-N1	-4.10	121.75	127.22
2	A	503	5GP	N2-C2-N1	3.68	122.98	117.25
2	A	503	5GP	O3P-P-O5'	-3.36	97.80	106.73
2	A	502	5GP	O3P-P-O5'	-3.27	98.02	106.73
2	A	502	5GP	C5-C6-N1	-2.96	119.39	123.43
2	A	501	5GP	O3P-P-O5'	-2.87	99.10	106.73
2	A	502	5GP	C6-N1-C2	2.82	120.41	115.93
2	A	501	5GP	C5-C6-N1	-2.63	119.84	123.43
2	A	501	5GP	C6-N1-C2	2.62	120.09	115.93
2	A	503	5GP	C5-C6-N1	-2.38	120.18	123.43
2	A	503	5GP	C6-N1-C2	2.36	119.68	115.93
2	A	501	5GP	C4-C5-N7	-2.36	106.94	109.40
2	A	503	5GP	C4-C5-N7	-2.20	107.10	109.40
2	A	502	5GP	C4-C5-N7	-2.19	107.12	109.40
2	A	503	5GP	C3'-C2'-C1'	2.16	104.23	100.98
2	A	502	5GP	O2P-P-O1P	2.13	119.03	110.68

There are no chirality outliers.

All (9) torsion outliers are listed below:

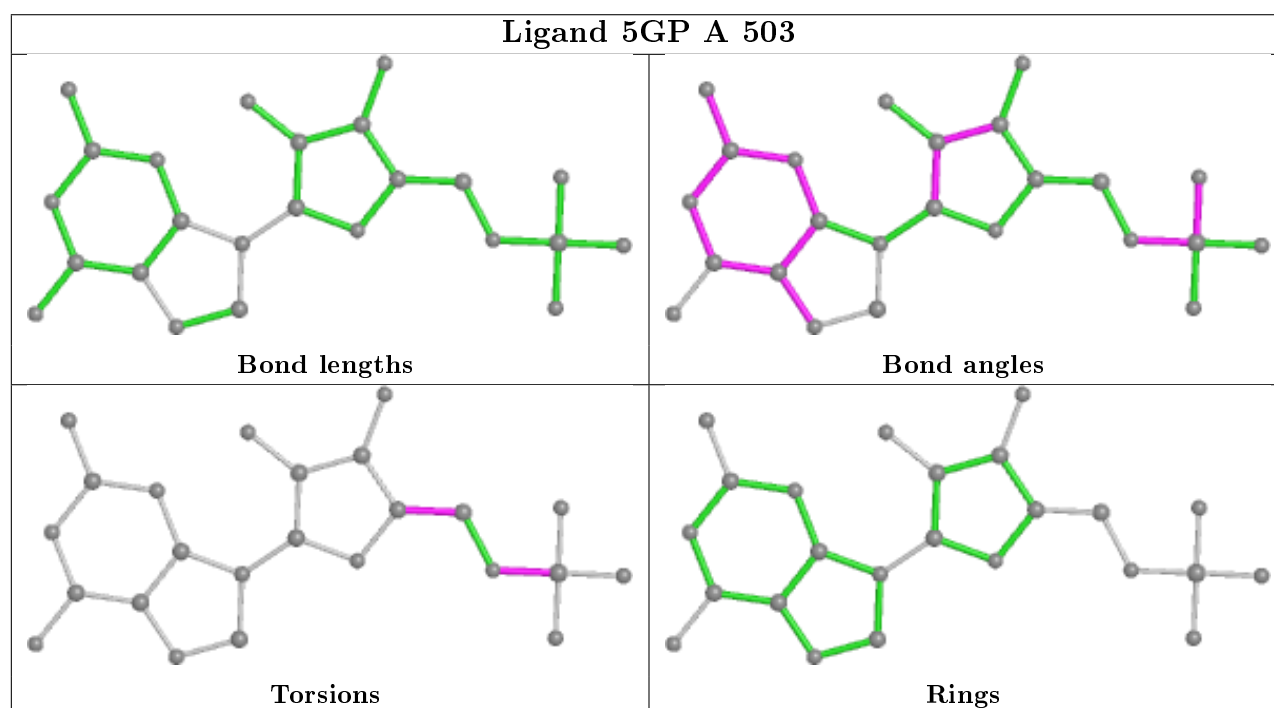
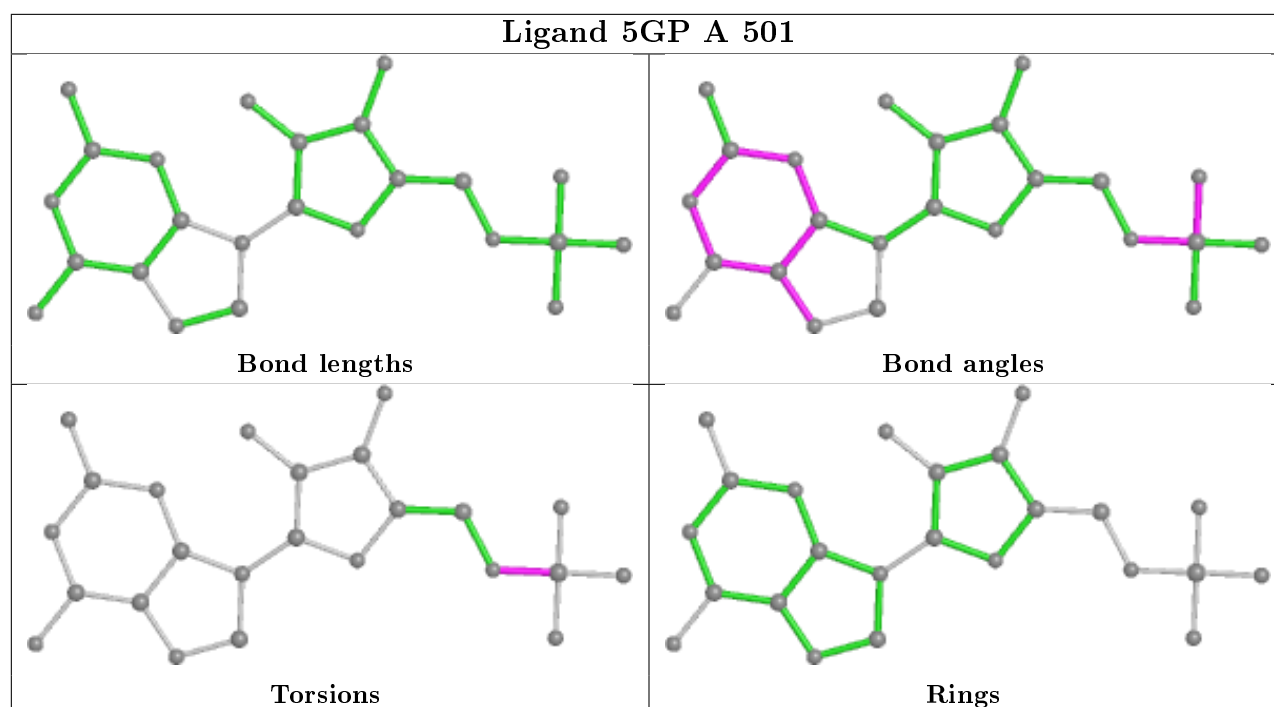
Mol	Chain	Res	Type	Atoms
2	A	501	5GP	C5'-O5'-P-O3P
2	A	503	5GP	C5'-O5'-P-O2P
2	A	503	5GP	C5'-O5'-P-O3P
2	A	502	5GP	C5'-O5'-P-O1P
2	A	502	5GP	C5'-O5'-P-O2P
2	A	502	5GP	C5'-O5'-P-O3P
2	A	501	5GP	C5'-O5'-P-O1P
2	A	503	5GP	C5'-O5'-P-O1P
2	A	503	5GP	O4'-C4'-C5'-O5'

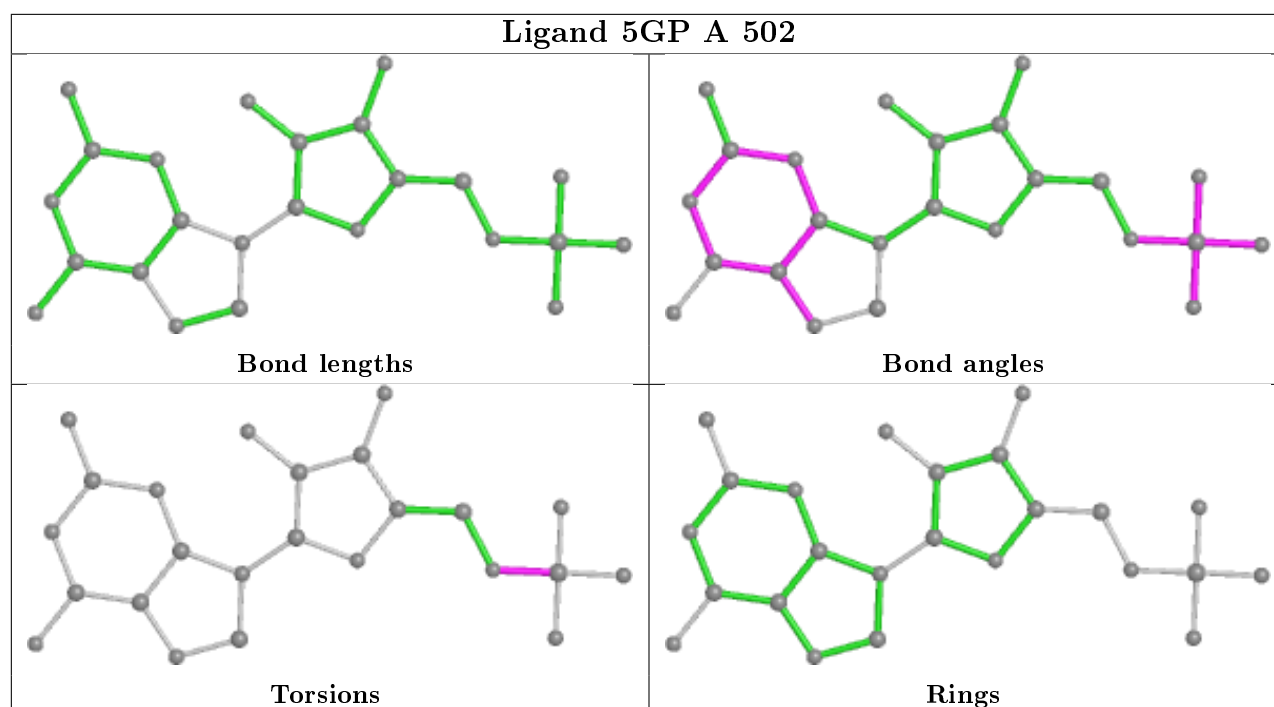
There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	503	5GP	1	0
2	A	502	5GP	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, 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	319/393 (81%)	0.70	9 (2%) 53 62	38, 76, 131, 153	0
1	B	303/393 (77%)	0.92	30 (9%) 7 8	41, 85, 157, 179	0
All	All	622/786 (79%)	0.81	39 (6%) 20 24	38, 79, 151, 179	0

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	415	LEU	8.1
1	B	291	ALA	5.2
1	B	250	LEU	4.8
1	A	406	ARG	4.1
1	B	215	TYR	3.6
1	A	397	THR	3.5
1	B	238	PHE	3.3
1	B	245	ASN	3.2
1	B	288	VAL	3.2
1	B	352	LYS	3.1
1	A	229	LYS	3.1
1	B	364	PHE	3.1
1	B	243	PHE	3.0
1	A	405	LYS	3.0
1	B	363	TYR	2.9
1	B	249	LYS	2.9
1	B	354	HIS	2.8
1	B	209	LEU	2.7
1	B	33	VAL	2.7
1	B	211	VAL	2.7
1	B	208	ARG	2.7
1	B	254	GLU	2.7
1	B	251	HIS	2.6
1	A	32	LYS	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	252	VAL	2.5
1	B	236	PRO	2.5
1	B	212	HIS	2.5
1	B	287	ASP	2.4
1	B	219	GLY	2.4
1	B	213	ARG	2.4
1	B	416	THR	2.4
1	B	224	ARG	2.4
1	A	298	VAL	2.4
1	B	266	ARG	2.3
1	A	402	GLN	2.3
1	A	271	LEU	2.2
1	A	209	LEU	2.1
1	B	337	PHE	2.1
1	B	237	THR	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 carbohydrates in this entry.

## 6.4 Ligands [i](#)

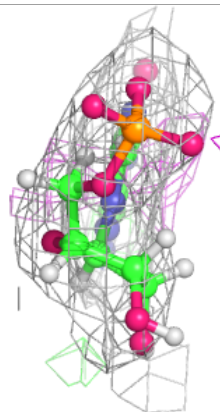
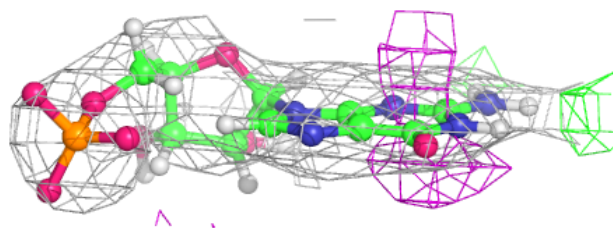
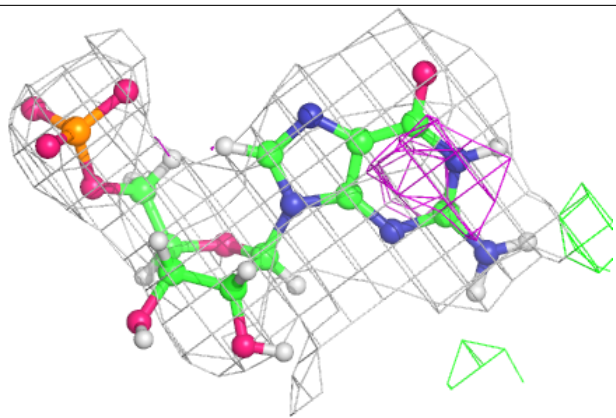
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	5GP	A	503	24/24	0.88	0.24	56,123,167,167	0
2	5GP	A	501	24/24	0.94	0.23	50,72,89,90	0
2	5GP	A	502	24/24	0.94	0.22	61,84,100,117	0

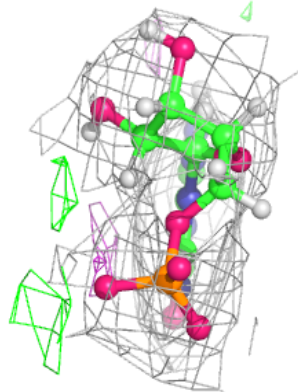
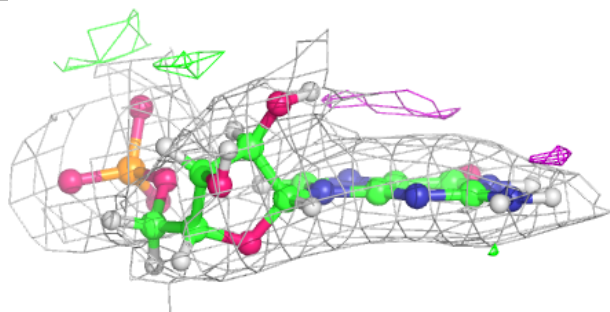
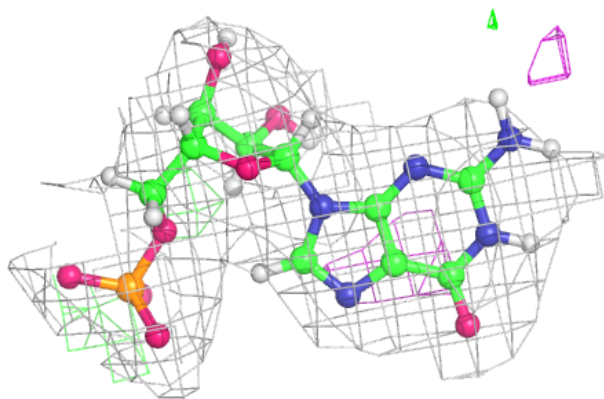
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

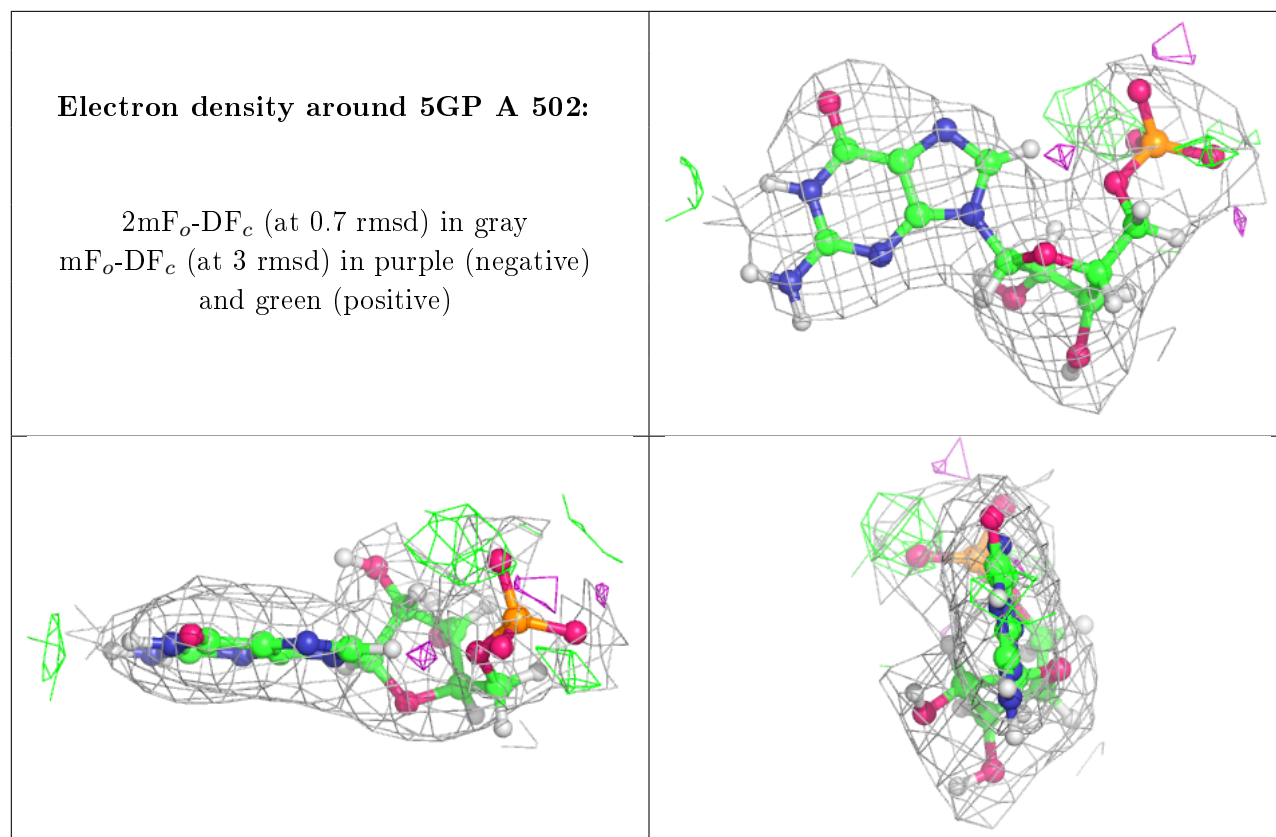
**Electron density around 5GP A 503:**

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

**Electron density around 5GP A 501:**

$2mF_o-DF_c$  (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.