



# wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 17, 2020 – 12:23 PM BST

PDB ID : 6PWZ  
Title : Crystal structure of human uridine-cytidine kinase 2 complexed with 2'-azidocytidine  
Authors : Cuthbert, B.J.; Goulding, C.W.  
Deposited on : 2019-07-24  
Resolution : 3.00 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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	:	2.13.1
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)	:	2.13.1

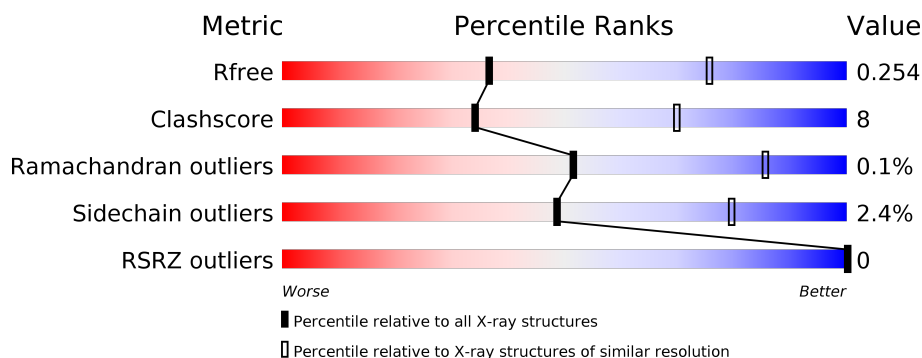
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.00 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	2092 (3.00-3.00)
Clashscore	141614	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)
RSRZ outliers	127900	1990 (3.00-3.00)

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	250	
1	B	250	
1	C	250	
1	D	250	
1	E	250	
1	F	250	

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Mol	Chain	Length	Quality of chain
1	G	250	
1	H	250	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	GOL	C	304	-	-	-	X
3	GOL	H	302	-	-	-	X

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 13675 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 Uridine-cytidine kinase 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	213	Total	C	N	O	S	0	0	0
			1616	1036	267	310	3			
1	B	215	Total	C	N	O	S	0	0	0
			1646	1059	271	313	3			
1	C	212	Total	C	N	O	S	0	0	0
			1659	1066	274	316	3			
1	D	212	Total	C	N	O	S	0	1	0
			1657	1063	281	310	3			
1	E	213	Total	C	N	O	S	0	0	0
			1671	1070	284	314	3			
1	F	213	Total	C	N	O	S	0	0	0
			1657	1066	275	313	3			
1	G	213	Total	C	N	O	S	0	0	0
			1615	1036	275	301	3			
1	H	205	Total	C	N	O	S	0	0	0
			1555	996	263	293	3			

- Molecule 2 is PHOSPHATE ION (three-letter code: PO4) (formula: O<sub>4</sub>P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		
2	C	1	Total	O	P	0	0
			5	4	1		
2	D	1	Total	O	P	0	0
			5	4	1		
2	E	1	Total	O	P	0	0
			5	4	1		
2	F	1	Total	O	P	0	0
			5	4	1		
2	G	1	Total	O	P	0	0
			5	4	1		
2	H	1	Total	O	P	0	0
			5	4	1		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



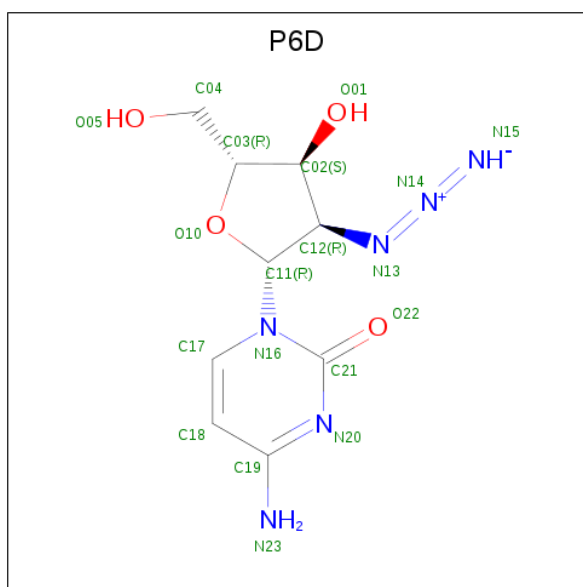
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	C	1	Total	C	O	0	0
			6	3	3		
3	C	1	Total	C	O	0	0
			6	3	3		
3	C	1	Total	C	O	0	0
			6	3	3		
3	C	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	D	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		
3	E	1	Total	C	O	0	0
			6	3	3		
3	E	1	Total	C	O	0	0
			6	3	3		
3	E	1	Total	C	O	0	0
			6	3	3		
3	E	1	Total	C	O	0	0
			6	3	3		
3	E	1	Total	C	O	0	0
			6	3	3		
3	F	1	Total	C	O	0	0
			6	3	3		
3	F	1	Total	C	O	0	0
			6	3	3		
3	F	1	Total	C	O	0	0
			6	3	3		
3	G	1	Total	C	O	0	0
			6	3	3		
3	H	1	Total	C	O	0	0
			6	3	3		
3	H	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is 2'-azidocytidine (three-letter code: P6D) (formula: C<sub>9</sub>H<sub>13</sub>N<sub>6</sub>O<sub>4</sub>) (labeled as "Ligand of Interest" by author).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	D	1	Total	C	N	O	0	0
			19	9	6	4		
4	E	1	Total	C	N	O	0	0
			19	9	6	4		
4	H	1	Total	C	N	O	0	0
			19	9	6	4		

- Molecule 5 is water.

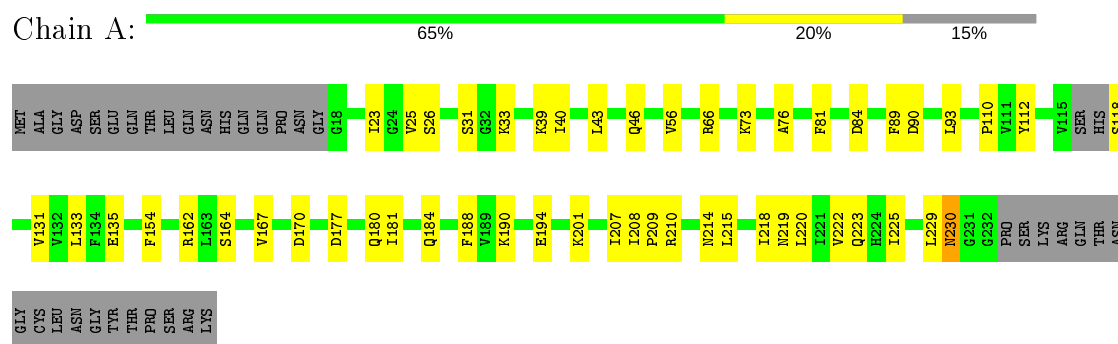
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	49	Total	O	0	0
			49	49		
5	B	42	Total	O	0	0
			42	42		
5	C	36	Total	O	0	0
			36	36		
5	D	38	Total	O	0	0
			38	38		
5	E	41	Total	O	0	0
			41	41		
5	F	46	Total	O	0	0
			46	46		
5	G	35	Total	O	0	0
			35	35		
5	H	35	Total	O	0	0
			35	35		



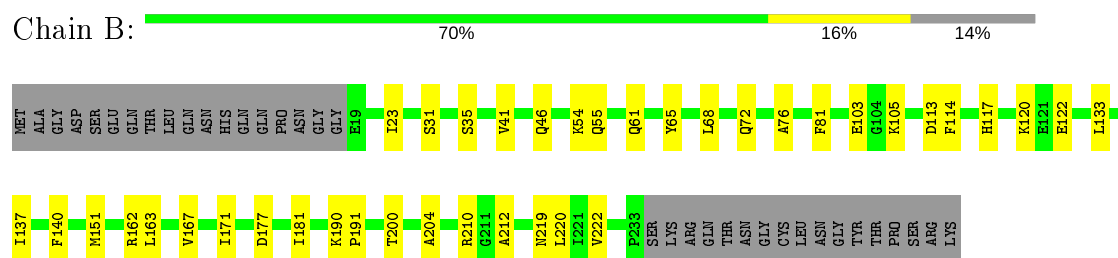
### 3 Residue-property plots

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.

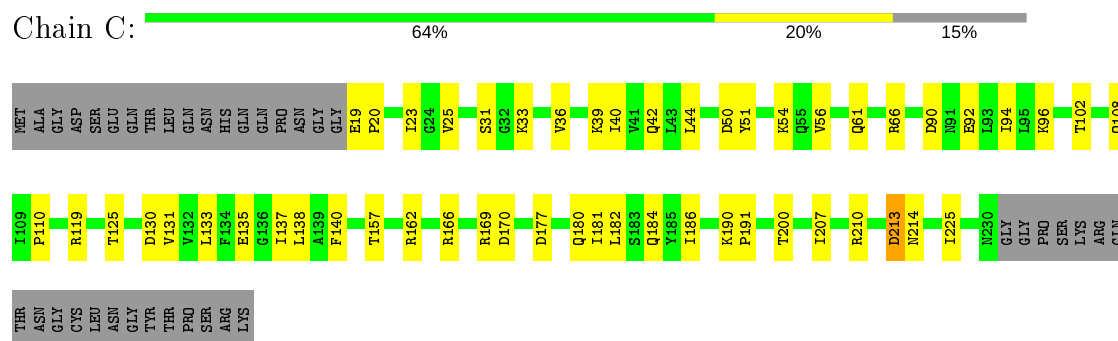
#### • Molecule 1: Uridine-cytidine kinase 2



#### • Molecule 1: Uridine-cytidine kinase 2

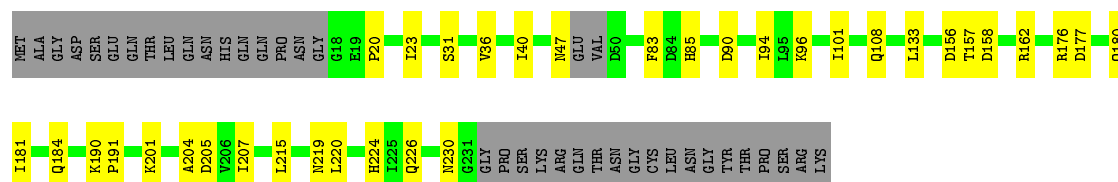


#### • Molecule 1: Uridine-cytidine kinase 2



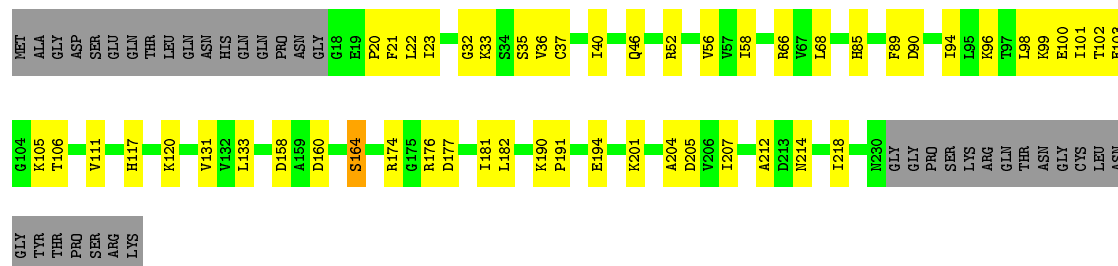
#### • Molecule 1: Uridine-cytidine kinase 2





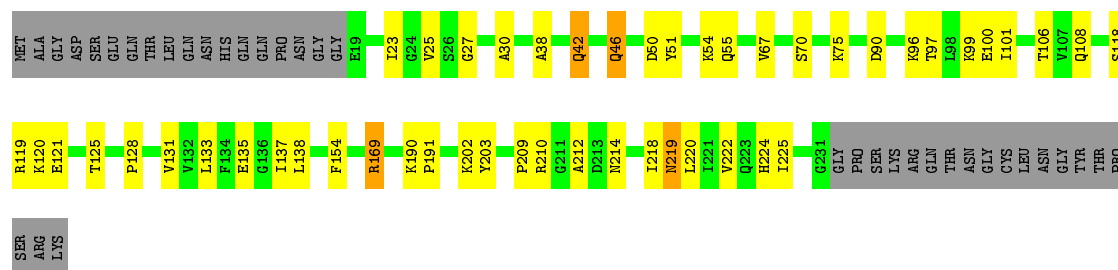
- Molecule 1: Uridine-cytidine kinase 2

Chain E: 64% 20% 15%



- Molecule 1: Uridine-cytidine kinase 2

Chain F: 66% 18% 15%



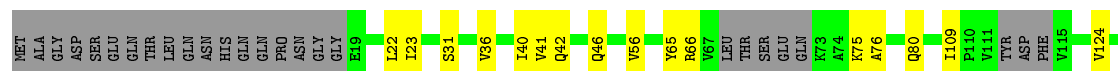
- Molecule 1: Uridine-cytidine kinase 2

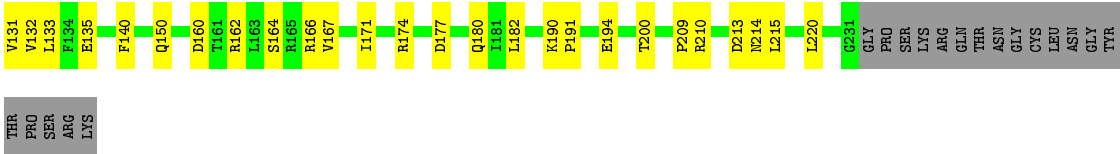
Chain G: 65% 20% 15%



- Molecule 1: Uridine-cytidine kinase 2

Chain H: 65% 17% 18%





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	93.73Å 84.74Å 153.58Å 90.00° 95.36° 90.00°	Depositor
Resolution (Å)	83.20 – 3.00 93.32 – 3.00	Depositor EDS
% Data completeness (in resolution range)	99.8 (83.20-3.00) 99.8 (93.32-3.00)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.65 (at 3.01Å)	Xtriage
Refinement program	PHENIX 1.16_3549	Depositor
R, $R_{free}$	0.197 , 0.254 0.197 , 0.254	Depositor DCC
$R_{free}$ test set	1119 reflections (2.32%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	69.6	Xtriage
Anisotropy	0.207	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 74.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	13675	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.61% 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: GOL, PO4, P6D

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	0.27	0/1640	0.46	0/2226
1	B	0.27	0/1674	0.45	0/2273
1	C	0.25	0/1688	0.42	0/2289
1	D	0.26	0/1662	0.44	0/2248
1	E	0.27	0/1699	0.45	0/2300
1	F	0.26	0/1686	0.42	0/2287
1	G	0.26	0/1643	0.45	0/2233
1	H	0.26	0/1578	0.47	0/2142
All	All	0.26	0/13270	0.45	0/17998

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	1616	0	1588	35	0
1	B	1646	0	1622	23	0
1	C	1659	0	1648	31	0
1	D	1657	0	1652	30	1
1	E	1671	0	1675	33	2

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	F	1657	0	1650	32	1
1	G	1615	0	1563	31	0
1	H	1555	0	1518	24	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
2	C	5	0	0	0	0
2	D	5	0	0	0	0
2	E	5	0	0	1	0
2	F	5	0	0	0	0
2	G	5	0	0	0	0
2	H	5	0	0	0	0
3	A	36	0	48	4	0
3	B	18	0	24	1	0
3	C	24	0	32	2	0
3	D	36	0	48	8	0
3	E	30	0	40	2	0
3	F	18	0	24	2	0
3	G	6	0	8	0	0
3	H	12	0	16	0	0
4	D	19	0	0	1	0
4	E	19	0	0	1	0
4	H	19	0	0	2	0
5	A	49	0	0	0	0
5	B	42	0	0	1	0
5	C	36	0	0	2	0
5	D	38	0	0	0	0
5	E	41	0	0	1	0
5	F	46	0	0	0	0
5	G	35	0	0	1	0
5	H	35	0	0	0	0
All	All	13675	0	13156	222	2

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 8.

The worst 5 of 222 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)
4:H:304:P6D:C03	4:H:304:P6D:O10	1.65	1.40
4:D:308:P6D:C03	4:D:308:P6D:O10	1.64	1.35
1:G:56:VAL:HG12	1:G:131:VAL:HB	1.56	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:46:GLN:HG3	1:A:56:VAL:HG22	1.61	0.82
1:A:40:ILE:HG13	1:A:222:VAL:HG22	1.63	0.80

All (2) 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:E:106:THR:OG1	1:F:50:ASP:OD2[2_656]	2.04	0.16
1:D:180:GLN:NE2	1:E:177:ASP:OD1[1_545]	2.19	0.01

## 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	209/250 (84%)	204 (98%)	4 (2%)	1 (0%)	29	68
1	B	213/250 (85%)	208 (98%)	5 (2%)	0	100	100
1	C	210/250 (84%)	208 (99%)	2 (1%)	0	100	100
1	D	209/250 (84%)	204 (98%)	5 (2%)	0	100	100
1	E	211/250 (84%)	206 (98%)	5 (2%)	0	100	100
1	F	211/250 (84%)	207 (98%)	4 (2%)	0	100	100
1	G	211/250 (84%)	199 (94%)	12 (6%)	0	100	100
1	H	199/250 (80%)	193 (97%)	6 (3%)	0	100	100
All	All	1673/2000 (84%)	1629 (97%)	43 (3%)	1 (0%)	51	85

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	112	TYR

### 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	171/220 (78%)	167 (98%)	4 (2%)	50	80
1	B	175/220 (80%)	174 (99%)	1 (1%)	86	95
1	C	181/220 (82%)	176 (97%)	5 (3%)	43	77
1	D	177/220 (80%)	177 (100%)	0	100	100
1	E	182/220 (83%)	180 (99%)	2 (1%)	73	90
1	F	180/220 (82%)	174 (97%)	6 (3%)	38	73
1	G	165/220 (75%)	157 (95%)	8 (5%)	25	62
1	H	162/220 (74%)	155 (96%)	7 (4%)	29	66
All	All	1393/1760 (79%)	1360 (98%)	33 (2%)	49	79

5 of 33 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	F	70	SER
1	G	42	GLN
1	H	135	GLU
1	F	75	LYS
1	F	169	ARG

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

### 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 monosaccharides in this entry.

## 5.6 Ligand geometry

41 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	PO4	E	301	-	4,4,4	0.97	0	6,6,6	0.47	0
2	PO4	C	301	-	4,4,4	0.89	0	6,6,6	0.41	0
3	GOL	C	305	-	5,5,5	0.88	0	5,5,5	1.00	0
3	GOL	A	302	-	5,5,5	0.94	0	5,5,5	0.93	0
3	GOL	A	306	-	5,5,5	0.93	0	5,5,5	0.94	0
3	GOL	B	303	-	5,5,5	0.88	0	5,5,5	1.01	0
3	GOL	B	304	-	5,5,5	0.91	0	5,5,5	0.99	0
3	GOL	E	304	-	5,5,5	0.93	0	5,5,5	0.96	0
3	GOL	H	303	-	5,5,5	0.88	0	5,5,5	0.99	0
2	PO4	H	301	-	4,4,4	0.94	0	6,6,6	0.42	0
3	GOL	F	302	-	5,5,5	0.91	0	5,5,5	1.01	0
2	PO4	B	301	-	4,4,4	0.92	0	6,6,6	0.57	0
4	P6D	D	308	-	15,20,20	5.32	11 (73%)	14,28,28	2.02	4 (28%)
3	GOL	E	306	-	5,5,5	0.91	0	5,5,5	0.99	0
3	GOL	F	304	-	5,5,5	0.94	0	5,5,5	0.92	0
3	GOL	E	302	-	5,5,5	0.94	0	5,5,5	0.96	0
3	GOL	C	303	-	5,5,5	0.90	0	5,5,5	1.01	0
2	PO4	G	301	-	4,4,4	0.90	0	6,6,6	0.49	0
3	GOL	C	304	-	5,5,5	0.93	0	5,5,5	0.92	0
2	PO4	A	301	-	4,4,4	0.92	0	6,6,6	0.55	0
3	GOL	D	302	-	5,5,5	0.90	0	5,5,5	0.99	0
4	P6D	H	304	-	15,20,20	5.55	11 (73%)	14,28,28	1.87	4 (28%)
2	PO4	F	301	-	4,4,4	0.96	0	6,6,6	0.39	0
3	GOL	G	302	-	5,5,5	0.92	0	5,5,5	0.95	0
3	GOL	D	307	-	5,5,5	0.89	0	5,5,5	0.96	0
3	GOL	F	303	-	5,5,5	0.88	0	5,5,5	0.99	0
3	GOL	A	303	-	5,5,5	0.91	0	5,5,5	0.94	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	GOL	D	303	-	5,5,5	0.96	0	5,5,5	0.91	0
3	GOL	A	307	-	5,5,5	0.83	0	5,5,5	1.00	0
3	GOL	A	305	-	5,5,5	1.01	0	5,5,5	0.82	0
3	GOL	E	305	-	5,5,5	0.88	0	5,5,5	0.99	0
3	GOL	H	302	-	5,5,5	1.06	0	5,5,5	0.86	0
2	PO4	D	301	-	4,4,4	0.93	0	6,6,6	0.43	0
3	GOL	C	302	-	5,5,5	0.87	0	5,5,5	1.02	0
3	GOL	D	304	-	5,5,5	0.92	0	5,5,5	1.00	0
3	GOL	D	305	-	5,5,5	0.96	0	5,5,5	0.91	0
3	GOL	E	303	-	5,5,5	0.93	0	5,5,5	0.97	0
3	GOL	B	302	-	5,5,5	0.87	0	5,5,5	1.01	0
3	GOL	A	304	-	5,5,5	0.90	0	5,5,5	1.08	0
3	GOL	D	306	-	5,5,5	0.91	0	5,5,5	0.97	0
4	P6D	E	307	-	15,20,20	5.54	12 (80%)	14,28,28	2.20	6 (42%)

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
3	GOL	C	305	-	-	0/4/4/4	-
3	GOL	A	302	-	-	1/4/4/4	-
3	GOL	A	306	-	-	0/4/4/4	-
3	GOL	B	303	-	-	1/4/4/4	-
3	GOL	B	304	-	-	2/4/4/4	-
3	GOL	E	304	-	-	1/4/4/4	-
3	GOL	H	303	-	-	2/4/4/4	-
3	GOL	F	302	-	-	0/4/4/4	-
4	P6D	D	308	-	-	3/7/25/25	0/2/2/2
3	GOL	E	306	-	-	2/4/4/4	-
3	GOL	F	304	-	-	4/4/4/4	-
3	GOL	E	302	-	-	2/4/4/4	-
3	GOL	C	303	-	-	0/4/4/4	-
3	GOL	C	304	-	-	0/4/4/4	-
3	GOL	D	305	-	-	0/4/4/4	-
3	GOL	D	302	-	-	2/4/4/4	-
4	P6D	H	304	-	-	5/7/25/25	0/2/2/2
3	GOL	G	302	-	-	1/4/4/4	-
3	GOL	D	307	-	-	0/4/4/4	-
3	GOL	F	303	-	-	0/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	A	303	-	-	2/4/4/4	-
3	GOL	D	303	-	-	4/4/4/4	-
3	GOL	A	307	-	-	0/4/4/4	-
3	GOL	E	305	-	-	0/4/4/4	-
3	GOL	H	302	-	-	1/4/4/4	-
3	GOL	A	305	-	-	2/4/4/4	-
3	GOL	C	302	-	-	0/4/4/4	-
3	GOL	D	304	-	-	2/4/4/4	-
3	GOL	E	303	-	-	2/4/4/4	-
3	GOL	B	302	-	-	0/4/4/4	-
3	GOL	A	304	-	-	2/4/4/4	-
3	GOL	D	306	-	-	0/4/4/4	-
4	P6D	E	307	-	-	4/7/25/25	0/2/2/2

The worst 5 of 34 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	H	304	P6D	C17-N16	10.35	1.48	1.35
4	E	307	P6D	C17-N16	9.26	1.47	1.35
4	H	304	P6D	O10-C03	8.99	1.65	1.45
4	E	307	P6D	C12-C11	8.90	1.67	1.54
4	D	308	P6D	C17-N16	8.75	1.46	1.35

The worst 5 of 14 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	308	P6D	C21-N20-C19	4.50	120.90	116.34
4	H	304	P6D	C12-N13-N14	3.64	123.80	115.54
4	E	307	P6D	N23-C19-N20	3.63	122.23	116.49
4	E	307	P6D	C21-N20-C19	3.62	120.01	116.34
4	D	308	P6D	C17-N16-C21	-3.29	115.97	121.20

There are no chirality outliers.

5 of 45 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	H	303	GOL	O1-C1-C2-C3
4	D	308	P6D	O10-C03-C04-O05
3	E	306	GOL	O1-C1-C2-C3
3	F	304	GOL	O2-C2-C3-O3

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Mol	Chain	Res	Type	Atoms
3	E	302	GOL	O1-C1-C2-C3

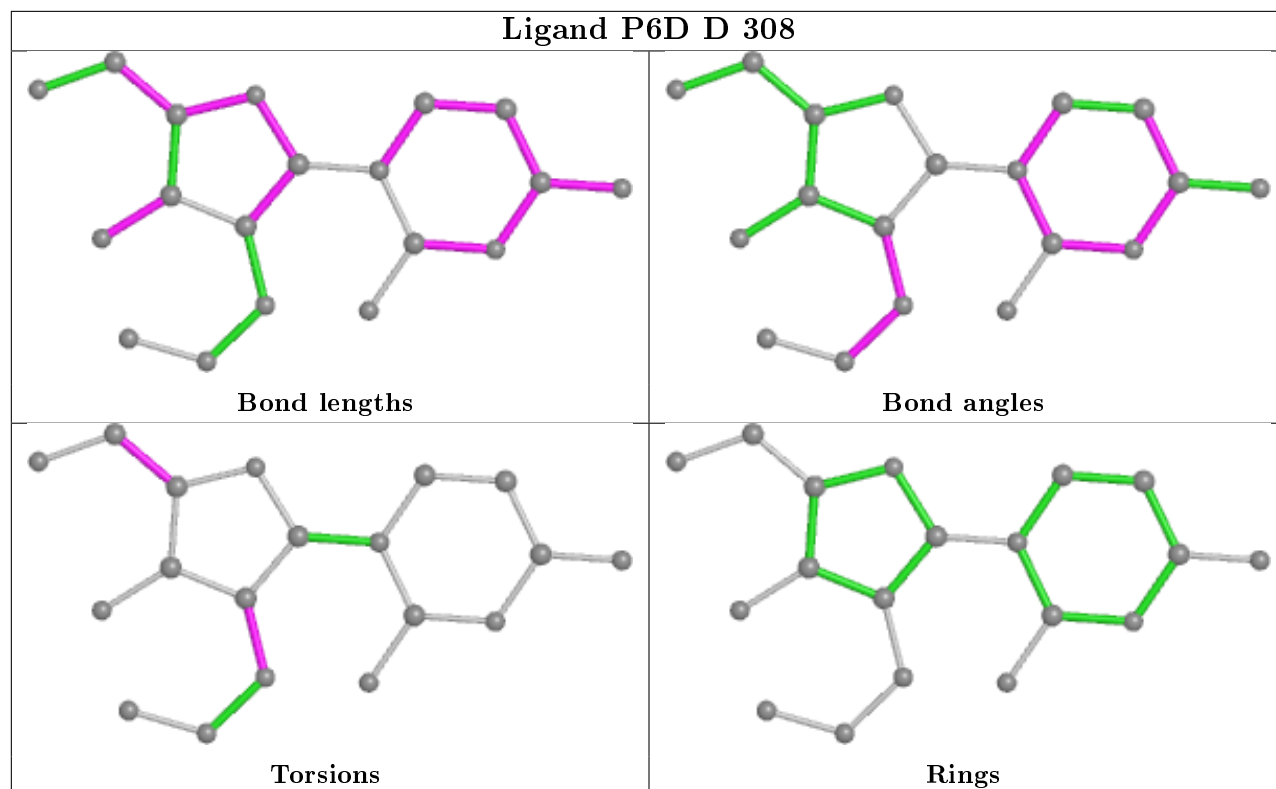
There are no ring outliers.

15 monomers are involved in 24 short contacts:

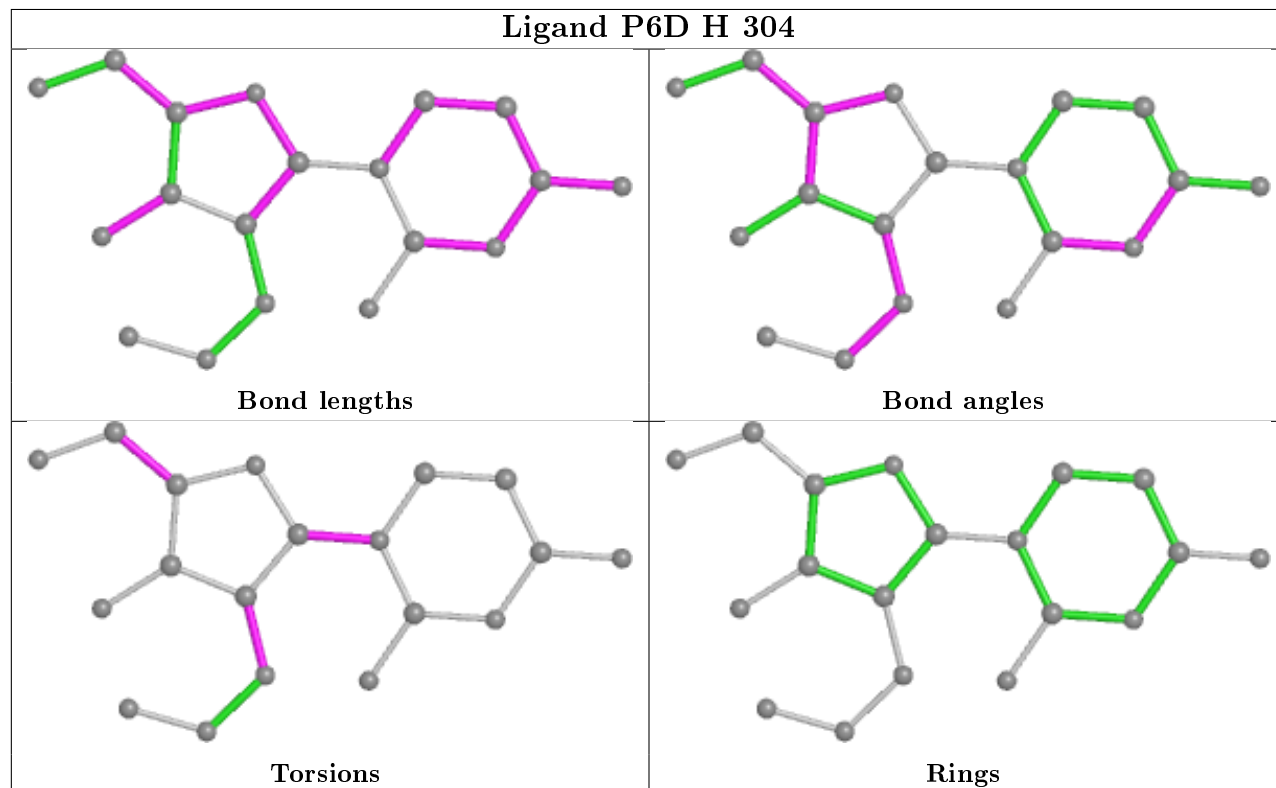
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	301	PO4	1	0
3	C	305	GOL	2	0
3	A	306	GOL	2	0
3	B	304	GOL	1	0
3	F	302	GOL	2	0
4	D	308	P6D	1	0
3	E	306	GOL	1	0
3	D	302	GOL	1	0
4	H	304	P6D	2	0
3	A	307	GOL	2	0
3	D	304	GOL	3	0
3	D	305	GOL	3	0
3	E	303	GOL	1	0
3	D	306	GOL	1	0
4	E	307	P6D	1	0

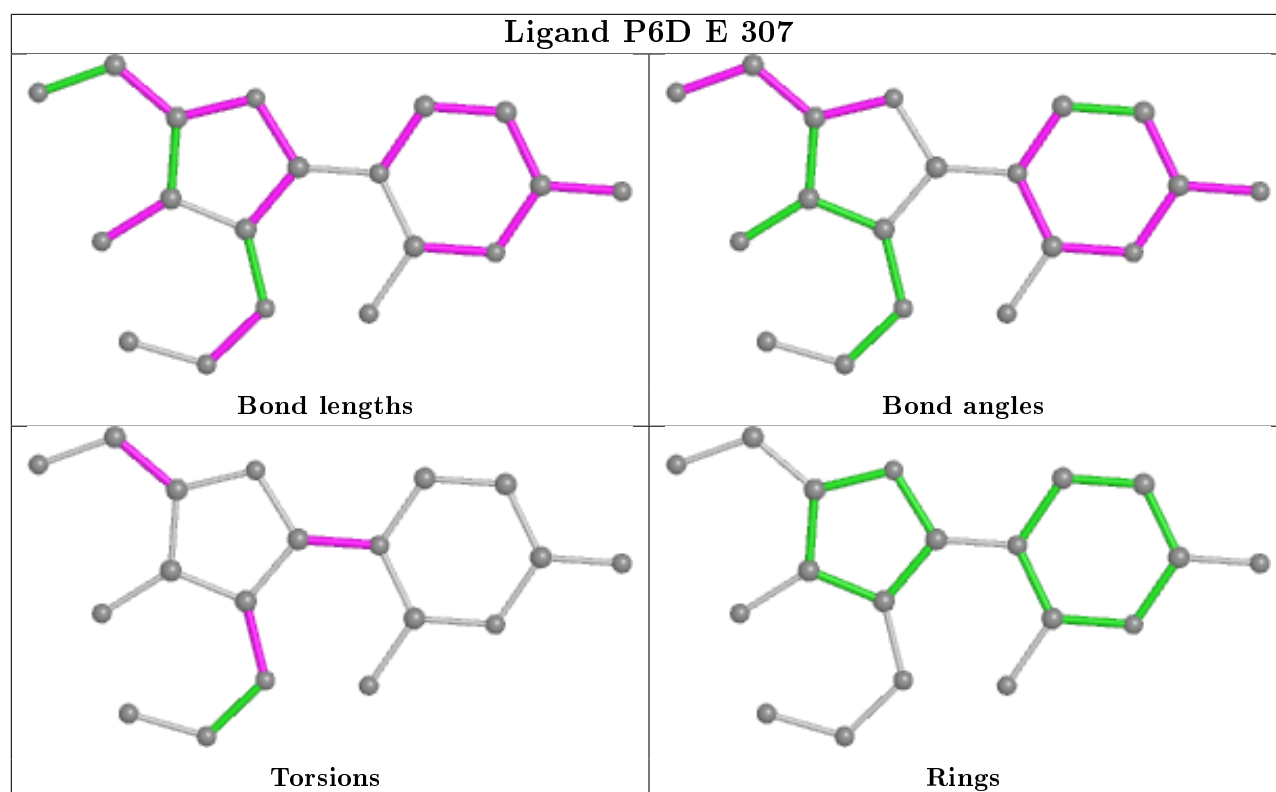
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

## Ligand P6D D 308



## Ligand P6D H 304





## 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, 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	213/250 (85%)	-0.09	0 100 100	31, 54, 107, 124	0
1	B	215/250 (86%)	-0.08	0 100 100	31, 55, 108, 128	0
1	C	212/250 (84%)	-0.07	0 100 100	29, 62, 92, 100	0
1	D	212/250 (84%)	-0.21	0 100 100	36, 55, 82, 110	0
1	E	213/250 (85%)	-0.15	0 100 100	36, 55, 88, 115	0
1	F	213/250 (85%)	-0.06	0 100 100	36, 64, 91, 117	0
1	G	213/250 (85%)	-0.14	0 100 100	34, 65, 130, 148	0
1	H	205/250 (82%)	-0.13	0 100 100	31, 67, 110, 122	0
All	All	1696/2000 (84%)	-0.12	0 100 100	29, 58, 105, 148	0

There are no RSRZ outliers to report.

### 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	GOL	H	302	6/6	0.64	0.46	36,57,61,67	6
3	GOL	C	304	6/6	0.67	0.57	55,68,78,78	0
4	P6D	H	304	19/19	0.71	0.33	72,82,95,96	0
3	GOL	F	304	6/6	0.72	0.25	81,84,93,101	0
4	P6D	E	307	19/19	0.74	0.31	41,71,96,107	19
3	GOL	F	303	6/6	0.75	0.40	62,66,72,75	6
4	P6D	D	308	19/19	0.75	0.31	56,72,96,105	19
3	GOL	F	302	6/6	0.75	0.24	73,73,92,110	0
3	GOL	E	306	6/6	0.83	0.25	74,78,87,100	0
3	GOL	B	304	6/6	0.84	0.43	69,77,87,95	0
3	GOL	A	303	6/6	0.84	0.33	57,60,73,77	0
3	GOL	A	306	6/6	0.85	0.32	48,67,79,88	0
3	GOL	B	303	6/6	0.86	0.29	65,71,76,78	0
3	GOL	E	305	6/6	0.87	0.30	64,77,82,84	0
3	GOL	A	302	6/6	0.87	0.25	63,69,72,73	0
3	GOL	G	302	6/6	0.87	0.27	69,81,87,90	0
3	GOL	A	307	6/6	0.88	0.27	40,52,67,71	0
3	GOL	D	303	6/6	0.88	0.33	53,68,81,83	0
3	GOL	E	302	6/6	0.89	0.24	51,65,75,79	0
3	GOL	E	304	6/6	0.89	0.48	49,54,69,70	6
3	GOL	D	302	6/6	0.89	0.26	42,52,59,62	0
3	GOL	H	303	6/6	0.89	0.17	79,81,93,97	0
3	GOL	E	303	6/6	0.90	0.35	57,84,87,92	0
3	GOL	C	305	6/6	0.90	0.20	62,77,92,101	0
3	GOL	D	307	6/6	0.91	0.36	71,78,89,93	0
3	GOL	A	305	6/6	0.91	0.21	49,70,79,86	0
3	GOL	B	302	6/6	0.92	0.27	51,73,75,77	0
3	GOL	D	304	6/6	0.93	0.42	59,77,85,92	0
3	GOL	D	306	6/6	0.93	0.26	55,61,66,71	0
3	GOL	C	303	6/6	0.93	0.33	81,93,97,118	0
3	GOL	A	304	6/6	0.95	0.25	49,50,52,54	0
3	GOL	C	302	6/6	0.95	0.22	49,56,59,63	0
3	GOL	D	305	6/6	0.95	0.17	60,68,82,83	0
2	PO4	D	301	5/5	0.96	0.21	41,54,59,67	0
2	PO4	G	301	5/5	0.97	0.16	41,51,55,56	0
2	PO4	H	301	5/5	0.98	0.18	40,41,47,51	0
2	PO4	B	301	5/5	0.99	0.19	35,38,41,49	0
2	PO4	F	301	5/5	0.99	0.17	42,44,52,66	0
2	PO4	C	301	5/5	0.99	0.16	26,37,57,64	0
2	PO4	A	301	5/5	0.99	0.17	34,41,46,49	0
2	PO4	E	301	5/5	0.99	0.17	51,53,61,65	0

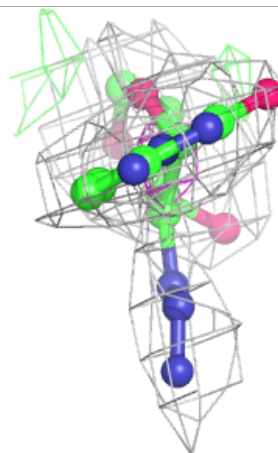
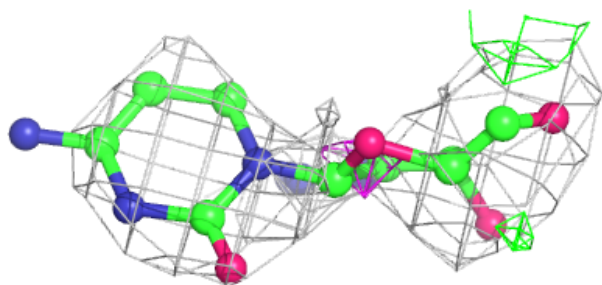
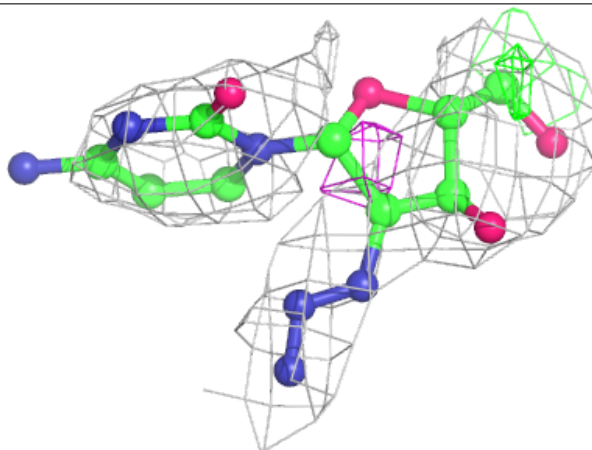
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers



as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

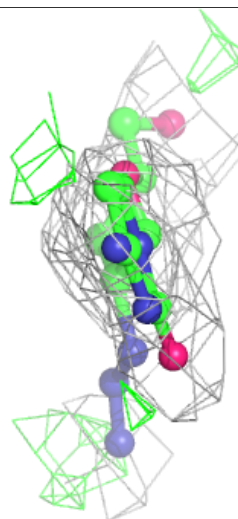
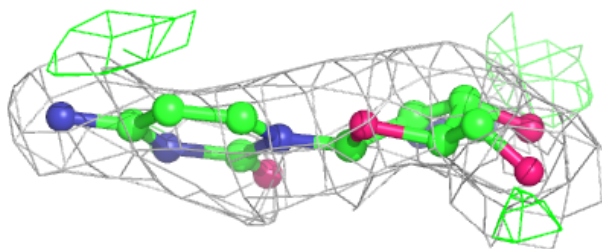
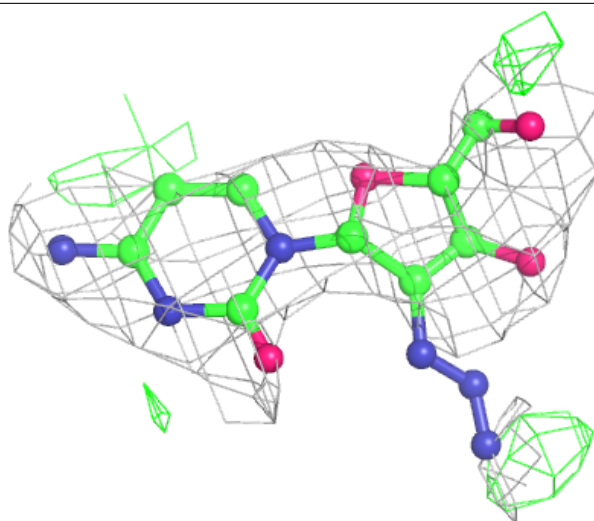
**Electron density around P6D H 304:**

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



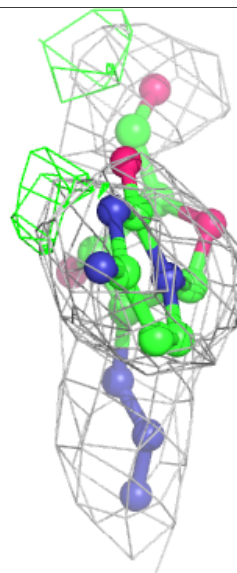
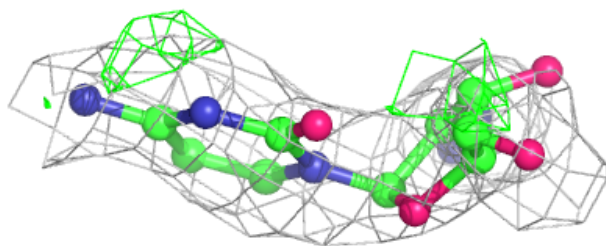
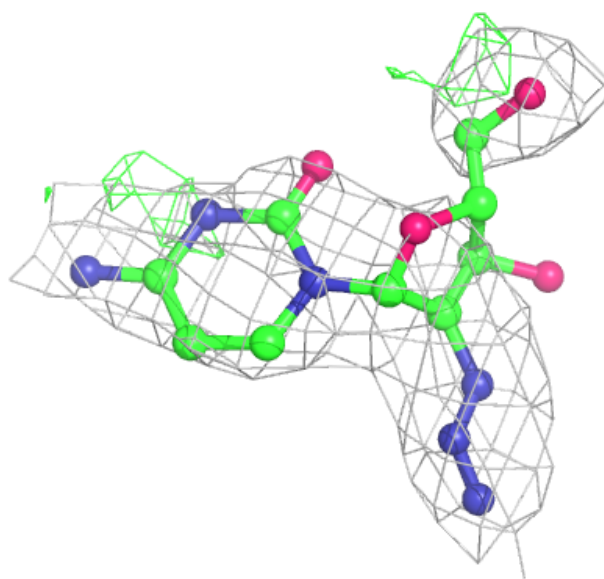
**Electron density around P6D E 307:**

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



**Electron density around P6D D 308:**

$2mF_o-DF_c$  (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.