



# Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 09:14 am BST

PDB ID : 5J1W  
Title : Crystal structure of human CLK1 in complex with pyrido[3,4-g]quinazoline derivative ZW31 (compound 14)  
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Deposited on : 2016-03-29  
Resolution : 2.42 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  
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

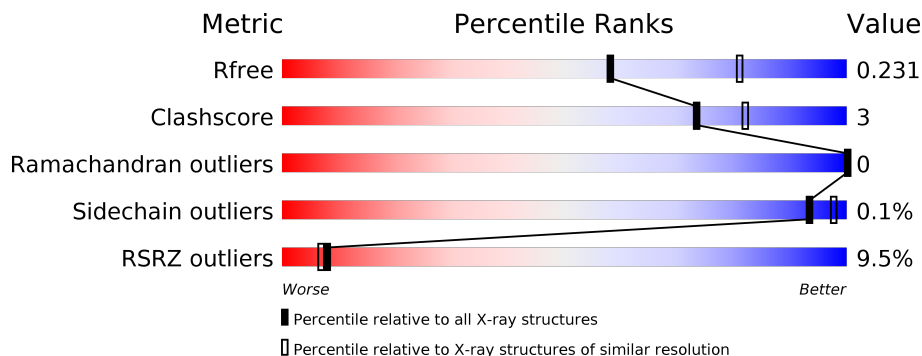
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.42 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	4647 (2.44-2.40)
Clashscore	141614	5161 (2.44-2.40)
Ramachandran outliers	138981	5073 (2.44-2.40)
Sidechain outliers	138945	5074 (2.44-2.40)
RSRZ outliers	127900	4543 (2.44-2.40)

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	339	 6% 93% 6%
1	B	339	 10% 93% 7%
1	C	339	 12% 84% 9% 7%

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
2	GOL	A	504	-	-	X	-
4	PO4	B	502	-	-	X	-

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8253 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 Dual specificity protein kinase CLK1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	338	2752	1758	477	500	17	0	1	0
1	B	337	2741	1752	473	499	17	0	1	0
1	C	316	2522	1614	428	463	17	0	0	0

There are 9 discrepancies between the modelled and reference sequences:

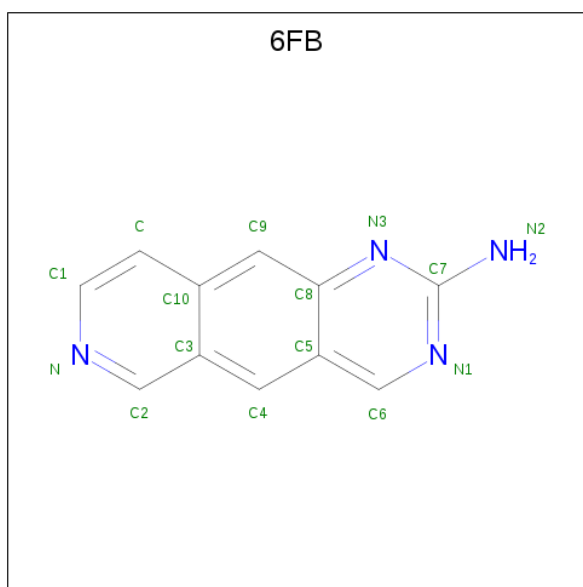
Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	SER	-	expression tag	UNP P49759
A	0	MET	-	expression tag	UNP P49759
A	432	ALA	ARG	conflict	UNP P49759
B	-1	SER	-	expression tag	UNP P49759
B	0	MET	-	expression tag	UNP P49759
B	432	ALA	ARG	conflict	UNP P49759
C	-1	SER	-	expression tag	UNP P49759
C	0	MET	-	expression tag	UNP P49759
C	432	ALA	ARG	conflict	UNP P49759

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



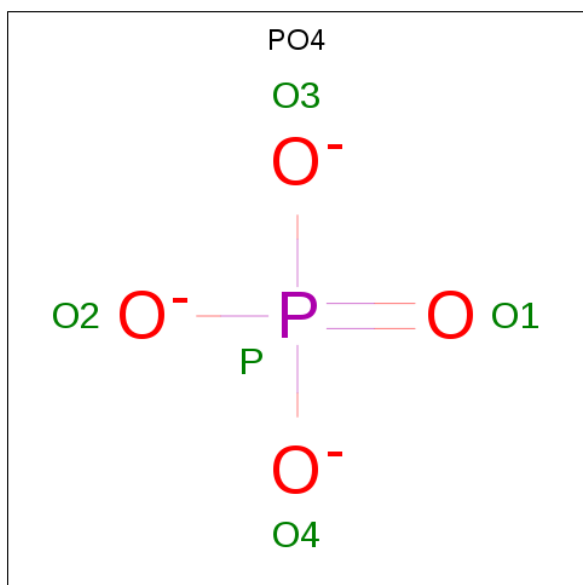
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 6 3 3	0	0
2	A	1	Total C O 6 3 3	0	0
2	A	1	Total C O 6 3 3	0	0
2	A	1	Total C O 6 3 3	0	0
2	A	1	Total C O 6 3 3	0	0
2	B	1	Total C O 6 3 3	0	0
2	C	1	Total C O 6 3 3	0	0
2	C	1	Total C O 6 3 3	0	0
2	C	1	Total C O 6 3 3	0	0

- Molecule 3 is pyrido[3,4-g]quinazolin-2-amine (three-letter code: 6FB) (formula: C<sub>11</sub>H<sub>8</sub>N<sub>4</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	N	0	0
			15	11	4		
3	B	1	Total	C	N	0	0
			15	11	4		
3	C	1	Total	C	N	0	0
			15	11	4		
3	C	1	Total	C	N	0	0
			15	11	4		

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



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	O	P	0	0
			5	4	1		
4	C	1	Total	O	P	0	0
			5	4	1		

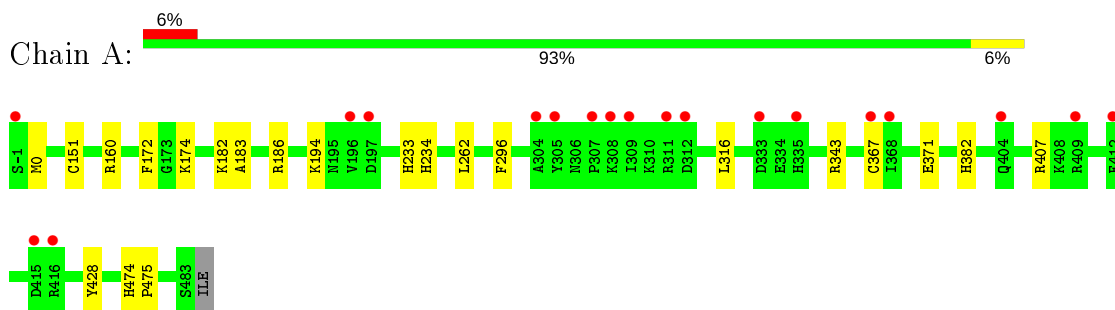
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	46	Total	O	0	0
			46	46		
5	B	32	Total	O	0	0
			32	32		
5	C	36	Total	O	0	0
			36	36		

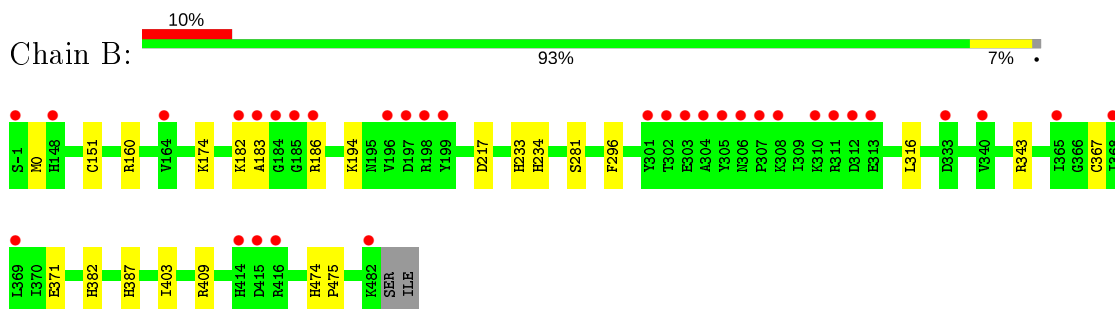
### 3 Residue-property plots [i](#)

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.

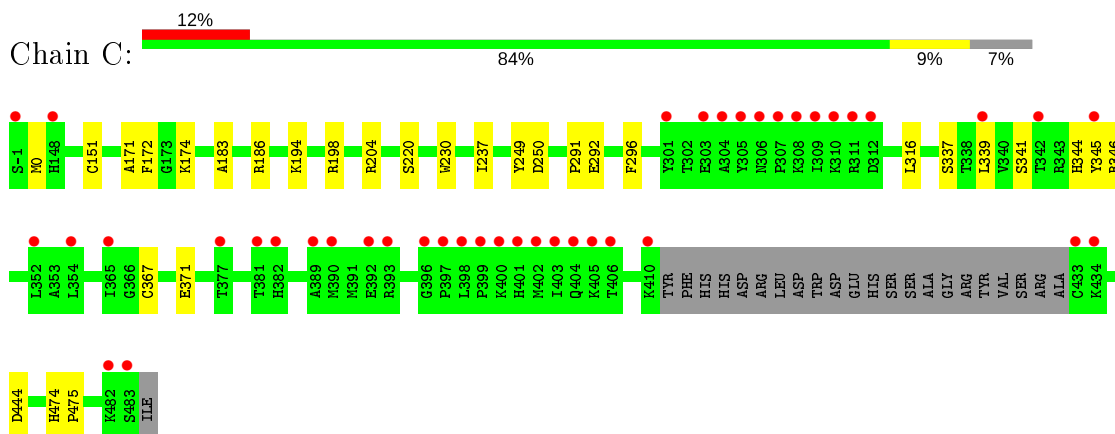
- Molecule 1: Dual specificity protein kinase CLK1



- Molecule 1: Dual specificity protein kinase CLK1



- Molecule 1: Dual specificity protein kinase CLK1





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	56.63Å 118.04Å 90.50Å 90.00° 99.55° 90.00°	Depositor
Resolution (Å)	35.59 – 2.42 35.36 – 2.42	Depositor EDS
% Data completeness (in resolution range)	99.8 (35.59-2.42) 99.8 (35.36-2.42)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.06 (at 2.42Å)	Xtrriage
Refinement program	REFMAC 5.8.0073	Depositor
R, $R_{free}$	0.189 , 0.228 0.197 , 0.231	Depositor DCC
$R_{free}$ test set	2159 reflections (4.83%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	50.4	Xtrriage
Anisotropy	0.301	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 51.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	8253	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	66.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.79% 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, 6FB

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	0.70	1/2826 (0.0%)	0.66	0/3824
1	B	0.64	0/2815	0.64	0/3810
1	C	0.70	1/2584 (0.0%)	0.66	0/3504
All	All	0.68	2/8225 (0.0%)	0.65	0/11138

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	292	GLU	CG-CD	5.34	1.59	1.51
1	A	428	TYR	CB-CG	5.12	1.59	1.51

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	2752	0	2682	15	0
1	B	2741	0	2669	14	0
1	C	2522	0	2435	22	0
2	A	30	0	40	5	0
2	B	6	0	8	0	0
2	C	18	0	24	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	15	0	0	3	0
3	B	15	0	0	0	0
3	C	30	0	0	1	0
4	B	5	0	0	2	0
4	C	5	0	0	0	0
5	A	46	0	0	1	0
5	B	32	0	0	0	0
5	C	36	0	0	2	0
All	All	8253	0	7858	52	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 3.

All (52) 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:151:CYS:SG	1:B:174:LYS:NZ	2.40	0.95
1:A:151:CYS:SG	1:A:174:LYS:NZ	2.41	0.93
1:C:151:CYS:SG	1:C:174:LYS:NZ	2.43	0.91
1:C:444:ASP:OD1	5:C:602:HOH:O	2.15	0.64
2:A:504:GOL:O3	2:A:504:GOL:O1	2.16	0.64
1:B:403:ILE:O	1:B:409:ARG:NH1	2.35	0.60
1:B:296:PHE:CZ	1:B:316:LEU:HD11	2.38	0.59
1:B:281:SER:O	1:C:230:TRP:HZ2	1.85	0.59
2:A:504:GOL:H32	3:A:506:6FB:N3	2.17	0.58
1:A:296:PHE:CZ	1:A:316:LEU:HD11	2.41	0.56
1:A:343:ARG:NH2	5:A:602:HOH:O	2.40	0.55
1:B:343:ARG:NE	1:B:382:HIS:HA	2.23	0.54
1:C:296:PHE:CZ	1:C:316:LEU:HD11	2.43	0.53
1:A:367:CYS:O	1:A:371:GLU:HG3	2.10	0.51
1:B:367:CYS:O	1:B:371:GLU:HG3	2.10	0.51
1:C:250:ASP:OD2	2:C:503:GOL:O1	2.18	0.51
1:A:343:ARG:NE	1:A:382:HIS:HA	2.26	0.50
1:B:0:MET:HE3	1:B:194:LYS:HG3	1.95	0.48
1:C:367:CYS:O	1:C:371:GLU:HG3	2.13	0.48
1:C:183:ALA:HB1	1:C:186:ARG:HD2	1.96	0.48
1:C:237:ILE:HG13	3:C:506:6FB:C2	2.43	0.48
1:C:337:SER:C	1:C:339:LEU:H	2.16	0.48
1:C:341:SER:O	1:C:346:ARG:NH2	2.47	0.48
4:B:502:PO4:O4	1:C:204:ARG:NH1	2.36	0.47
2:A:504:GOL:C3	3:A:506:6FB:N3	2.78	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:217:ASP:OD1	4:B:502:PO4:O3	2.33	0.47
1:B:183:ALA:HB1	1:B:186:ARG:HD2	1.98	0.46
1:A:474:HIS:CG	1:A:475:PRO:HD2	2.51	0.46
1:C:474:HIS:CG	1:C:475:PRO:HD2	2.51	0.46
1:A:183:ALA:HB1	1:A:186:ARG:HD2	1.99	0.45
1:B:233:HIS:O	1:B:234:HIS:C	2.55	0.45
1:A:172:PHE:CD1	1:A:172:PHE:C	2.91	0.44
1:C:220:SER:HA	5:C:609:HOH:O	2.15	0.44
1:B:281:SER:O	1:C:230:TRP:CZ2	2.68	0.44
2:A:504:GOL:O2	3:A:506:6FB:N2	2.52	0.43
1:A:262:LEU:H	2:A:501:GOL:H11	1.84	0.43
1:B:160:ARG:HD2	1:B:182:LYS:HG3	2.02	0.42
1:B:151:CYS:SG	1:B:174:LYS:HD3	2.59	0.42
1:A:151:CYS:SG	1:A:174:LYS:HD3	2.60	0.42
1:C:0:MET:HE3	1:C:194:LYS:HG3	2.01	0.42
1:A:160:ARG:HD2	1:A:182:LYS:HG3	2.02	0.42
1:C:172:PHE:C	1:C:172:PHE:CD1	2.93	0.41
1:C:344:HIS:HB2	1:C:345:TYR:CE1	2.55	0.41
1:A:0:MET:HB2	1:A:0:MET:HE3	1.98	0.41
1:C:0:MET:HE3	1:C:0:MET:HB2	1.99	0.41
1:C:151:CYS:SG	1:C:174:LYS:HD3	2.61	0.41
1:A:233:HIS:O	1:A:234:HIS:C	2.58	0.41
1:B:474:HIS:CG	1:B:475:PRO:HD2	2.55	0.41
1:C:171:ALA:O	1:C:194:LYS:NZ	2.39	0.41
1:A:407:ARG:HA	1:C:198:ARG:HD3	2.03	0.41
1:A:0:MET:HE3	1:A:194:LYS:HG3	2.03	0.40
1:C:249:TYR:HB2	1:C:291:PRO:HB2	2.03	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	337/339 (99%)	322 (96%)	15 (4%)	0	100	100
1	B	336/339 (99%)	321 (96%)	15 (4%)	0	100	100
1	C	312/339 (92%)	298 (96%)	14 (4%)	0	100	100
All	All	985/1017 (97%)	941 (96%)	44 (4%)	0	100	100

There are no Ramachandran outliers to report.

### 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	301/308 (98%)	301 (100%)	0	100	100
1	B	300/308 (97%)	299 (100%)	1 (0%)	92	97
1	C	274/308 (89%)	274 (100%)	0	100	100
All	All	875/924 (95%)	874 (100%)	1 (0%)	93	98

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	387	HIS

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

### 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](#)

15 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	GOL	C	502	-	5,5,5	0.45	0	5,5,5	0.46	0
3	6FB	C	505	-	17,17,17	0.40	0	23,24,24	0.56	1 (4%)
3	6FB	B	503	-	17,17,17	0.62	0	23,24,24	0.57	1 (4%)
3	6FB	A	506	-	17,17,17	0.43	0	23,24,24	0.52	0
3	6FB	C	506	-	17,17,17	0.22	0	23,24,24	0.51	0
2	GOL	A	503	-	5,5,5	0.59	0	5,5,5	1.18	0
2	GOL	C	503	-	5,5,5	0.37	0	5,5,5	0.78	0
2	GOL	A	502	-	5,5,5	0.79	0	5,5,5	1.01	0
4	PO4	B	502	-	4,4,4	0.73	0	6,6,6	0.83	0
4	PO4	C	504	-	4,4,4	0.65	0	6,6,6	0.60	0
2	GOL	B	501	-	5,5,5	0.47	0	5,5,5	0.23	0
2	GOL	C	501	-	5,5,5	0.34	0	5,5,5	0.80	0
2	GOL	A	501	-	5,5,5	0.41	0	5,5,5	0.86	0
2	GOL	A	504	-	5,5,5	0.16	0	5,5,5	1.13	0
2	GOL	A	505	-	5,5,5	0.43	0	5,5,5	0.24	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GOL	C	502	-	-	0/4/4/4	-
3	6FB	C	505	-	-	-	0/3/3/3
3	6FB	B	503	-	-	-	0/3/3/3
3	6FB	A	506	-	-	-	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	6FB	C	506	-	-	-	0/3/3/3
2	GOL	A	503	-	-	1/4/4/4	-
2	GOL	C	503	-	-	2/4/4/4	-
2	GOL	A	504	-	-	2/4/4/4	-
2	GOL	B	501	-	-	2/4/4/4	-
2	GOL	C	501	-	-	4/4/4/4	-
2	GOL	A	501	-	-	2/4/4/4	-
2	GOL	A	502	-	-	2/4/4/4	-
2	GOL	A	505	-	-	2/4/4/4	-

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	505	6FB	C4-C5-C6	2.09	127.35	121.95
3	B	503	6FB	C4-C5-C6	2.02	127.16	121.95

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	503	GOL	C1-C2-C3-O3
2	B	501	GOL	O1-C1-C2-O2
2	B	501	GOL	O1-C1-C2-C3
2	C	501	GOL	O1-C1-C2-O2
2	C	501	GOL	O1-C1-C2-C3
2	C	501	GOL	C1-C2-C3-O3
2	A	501	GOL	O1-C1-C2-O2
2	A	501	GOL	O1-C1-C2-C3
2	C	501	GOL	O2-C2-C3-O3
2	A	504	GOL	C1-C2-C3-O3
2	A	502	GOL	O1-C1-C2-C3
2	C	503	GOL	O2-C2-C3-O3
2	A	504	GOL	O2-C2-C3-O3
2	A	505	GOL	O2-C2-C3-O3
2	A	502	GOL	O1-C1-C2-O2
2	A	505	GOL	C1-C2-C3-O3
2	A	503	GOL	O1-C1-C2-O2

There are no ring outliers.

6 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	506	6FB	3	0
3	C	506	6FB	1	0
2	C	503	GOL	1	0
4	B	502	PO4	2	0
2	A	501	GOL	1	0
2	A	504	GOL	4	0

## 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	338/339 (99%)	0.21	19 (5%) 24 22	30, 55, 104, 134	0
1	B	337/339 (99%)	0.41	33 (9%) 7 6	40, 64, 112, 132	0
1	C	316/339 (93%)	0.65	42 (13%) 3 2	34, 62, 131, 164	0
All	All	991/1017 (97%)	0.42	94 (9%) 8 7	30, 60, 115, 164	0

All (94) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	433	CYS	7.6
1	C	304	ALA	7.4
1	C	305	TYR	7.1
1	C	309	ILE	6.7
1	A	304	ALA	6.7
1	C	404	GLN	6.6
1	C	434	LYS	6.3
1	C	382	HIS	6.3
1	C	308	LYS	6.3
1	C	307	PRO	6.2
1	C	306	ASN	5.6
1	C	403	ILE	5.5
1	B	307	PRO	5.2
1	C	406	THR	5.0
1	A	415	ASP	4.8
1	A	308	LYS	4.5
1	C	483	SER	4.3
1	B	415	ASP	4.0
1	B	482	LYS	4.0
1	C	311	ARG	4.0
1	B	-1	SER	4.0
1	C	482	LYS	4.0
1	C	402	MET	3.9

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	B	304	ALA	3.8
1	C	312	ASP	3.8
1	C	400	LYS	3.7
1	C	301	TYR	3.7
1	B	185	GLY	3.7
1	C	339	LEU	3.6
1	C	354	LEU	3.4
1	B	303	GLU	3.3
1	C	397	PRO	3.3
1	B	301	TYR	3.2
1	B	183	ALA	3.2
1	C	303	GLU	3.2
1	B	186	ARG	3.2
1	B	302	THR	3.1
1	C	148	HIS	3.1
1	C	352	LEU	3.1
1	C	392	GLU	3.0
1	C	396	GLY	3.0
1	C	381	THR	3.0
1	C	405	LYS	3.0
1	C	398	LEU	2.9
1	B	148	HIS	2.9
1	C	401	HIS	2.9
1	A	-1	SER	2.9
1	B	310	LYS	2.8
1	B	416	ARG	2.8
1	B	164	VAL	2.8
1	A	312	ASP	2.8
1	B	340	VAL	2.8
1	A	409	ARG	2.7
1	A	335	HIS	2.7
1	B	305	TYR	2.7
1	A	333	ASP	2.7
1	A	311	ARG	2.6
1	B	368	ILE	2.6
1	B	308	LYS	2.6
1	C	365	ILE	2.6
1	C	345	TYR	2.6
1	C	377	THR	2.6
1	B	414	HIS	2.5
1	C	389	ALA	2.5
1	A	412	PHE	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	367	CYS	2.4
1	B	313	GLU	2.4
1	A	307	PRO	2.4
1	A	305	TYR	2.4
1	A	309	ILE	2.4
1	C	393	ARG	2.4
1	B	182	LYS	2.4
1	B	184	GLY	2.3
1	C	310	LYS	2.3
1	C	390	MET	2.3
1	B	306	ASN	2.3
1	B	369	LEU	2.2
1	A	196	VAL	2.2
1	B	311	ARG	2.2
1	B	198	ARG	2.2
1	B	333	ASP	2.2
1	B	197	ASP	2.2
1	A	416	ARG	2.2
1	B	196	VAL	2.2
1	C	410	LYS	2.2
1	B	312	ASP	2.1
1	A	404	GLN	2.1
1	C	-1	SER	2.1
1	A	368	ILE	2.1
1	C	399	PRO	2.1
1	A	197	ASP	2.1
1	B	199	TYR	2.1
1	C	342	THR	2.0
1	B	365	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 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	GOL	A	501	6/6	0.72	0.33	59,60,67,79	0
2	GOL	A	505	6/6	0.79	0.17	67,76,86,87	0
2	GOL	A	503	6/6	0.80	0.15	63,67,70,71	0
2	GOL	A	502	6/6	0.86	0.14	67,75,80,82	0
2	GOL	C	501	6/6	0.87	0.17	78,79,81,82	0
2	GOL	B	501	6/6	0.91	0.15	64,67,70,76	0
4	PO4	C	504	5/5	0.93	0.23	76,79,96,100	0
2	GOL	C	502	6/6	0.94	0.23	71,84,89,106	0
3	6FB	C	506	15/15	0.95	0.14	48,51,68,68	0
2	GOL	C	503	6/6	0.95	0.18	47,55,66,72	0
2	GOL	A	504	6/6	0.96	0.14	52,64,65,71	0
3	6FB	B	503	15/15	0.97	0.14	27,31,44,44	0
3	6FB	A	506	15/15	0.97	0.17	26,30,36,40	0
3	6FB	C	505	15/15	0.97	0.17	25,30,32,35	0
4	PO4	B	502	5/5	0.98	0.21	60,67,70,82	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.