



# wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 16, 2017 – 11:46 AM EST

PDB ID : 5H2U  
Title : Crystal structure of PTK6 Kinase Domain complexed with Dasatinib  
Authors : Thakur, M.K.; Birudukota, S.; Swaminathan, S.; Tyagi, R.; Gosu, R.  
Deposited on : 2016-10-18  
Resolution : 2.24 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

<http://wwpdb.org/validation/2016/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 : 1.7.1 (RC1), CSD as537be (2016)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20028442  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : rb-20028442

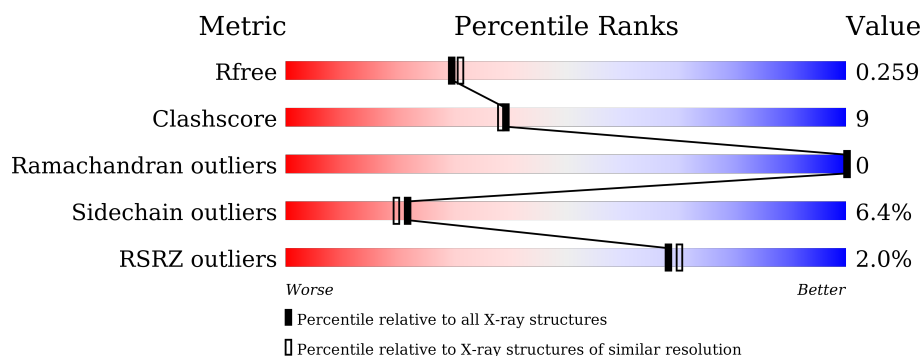
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.24 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))
$R_{free}$	91344	1611 (2.26-2.22)
Clashscore	102246	1764 (2.26-2.22)
Ramachandran outliers	100387	1724 (2.26-2.22)
Sidechain outliers	100360	1724 (2.26-2.22)
RSRZ outliers	91569	1616 (2.26-2.22)

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. 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	267	<div> <div>%</div> <div> <div></div> <div>78%</div> <div>18%</div> <div>.</div> </div> </div>
1	B	267	<div> <div></div> <div>80%</div> <div>18%</div> <div>..</div> </div>
1	C	267	<div> <div>3%</div> <div> <div></div> <div>77%</div> <div>21%</div> <div>..</div> </div> </div>
1	D	267	<div> <div>3%</div> <div> <div></div> <div>79%</div> <div>19%</div> <div>..</div> </div> </div>

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	GOL	B	504	-	-	-	X
3	GOL	C	502	-	-	-	X
3	GOL	D	501	-	-	-	X

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9541 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 Protein-tyrosine kinase 6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	266	Total	C	N	O	S	0	10	0
			2211	1413	390	390	18			
1	B	264	Total	C	N	O	S	0	12	0
			2202	1406	389	390	17			
1	C	265	Total	C	N	O	S	0	12	0
			2201	1406	387	389	19			
1	D	264	Total	C	N	O	S	0	11	0
			2190	1399	382	391	18			

There are 24 discrepancies between the modelled and reference sequences:

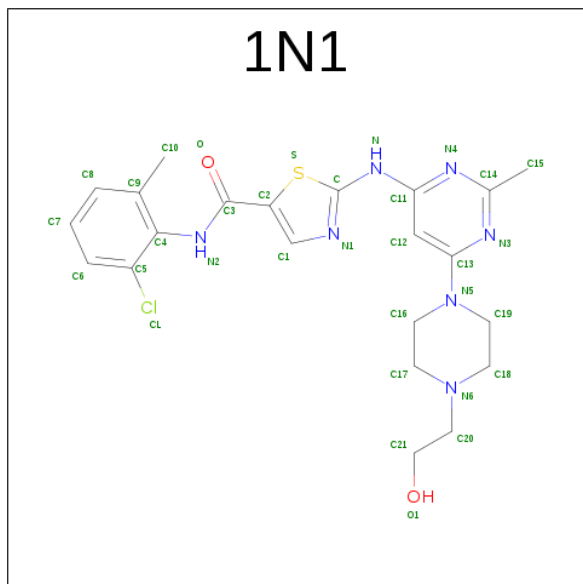
Chain	Residue	Modelled	Actual	Comment	Reference
A	184	CXM	-	expression tag	UNP Q13882
A	433	THR	CYS	engineered mutation	UNP Q13882
A	447	HIS	-	expression tag	UNP Q13882
A	448	HIS	-	expression tag	UNP Q13882
A	449	HIS	-	expression tag	UNP Q13882
A	450	HIS	-	expression tag	UNP Q13882
B	184	CXM	-	expression tag	UNP Q13882
B	433	THR	CYS	engineered mutation	UNP Q13882
B	447	HIS	-	expression tag	UNP Q13882
B	448	HIS	-	expression tag	UNP Q13882
B	449	HIS	-	expression tag	UNP Q13882
B	450	HIS	-	expression tag	UNP Q13882
C	184	CXM	-	expression tag	UNP Q13882
C	433	THR	CYS	engineered mutation	UNP Q13882
C	447	HIS	-	expression tag	UNP Q13882
C	448	HIS	-	expression tag	UNP Q13882
C	449	HIS	-	expression tag	UNP Q13882
C	450	HIS	-	expression tag	UNP Q13882
D	184	CXM	-	expression tag	UNP Q13882
D	433	THR	CYS	engineered mutation	UNP Q13882
D	447	HIS	-	expression tag	UNP Q13882

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Chain	Residue	Modelled	Actual	Comment	Reference
D	448	HIS	-	expression tag	UNP Q13882
D	449	HIS	-	expression tag	UNP Q13882
D	450	HIS	-	expression tag	UNP Q13882

- Molecule 2 is N-(2-CHLORO-6-METHYLPHENYL)-2-({6-[4-(2-HYDROXYETHYL)PIPERAZIN-1-YL]-2-METHYLPYRIMIDIN-4-YL}AMINO)-1,3-THIAZOLE-5-CARBOXAMIDE (three-letter code: 1N1) (formula: C<sub>22</sub>H<sub>26</sub>ClN<sub>7</sub>O<sub>2</sub>S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total 41	C 29	Cl 2	N 7	O 2	S 1	0	1
2	B	1	Total 41	C 29	Cl 2	N 7	O 2	S 1	0	1
2	C	1	Total 41	C 29	Cl 2	N 7	O 2	S 1	0	1
2	D	1	Total 41	C 29	Cl 2	N 7	O 2	S 1	0	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	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	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		

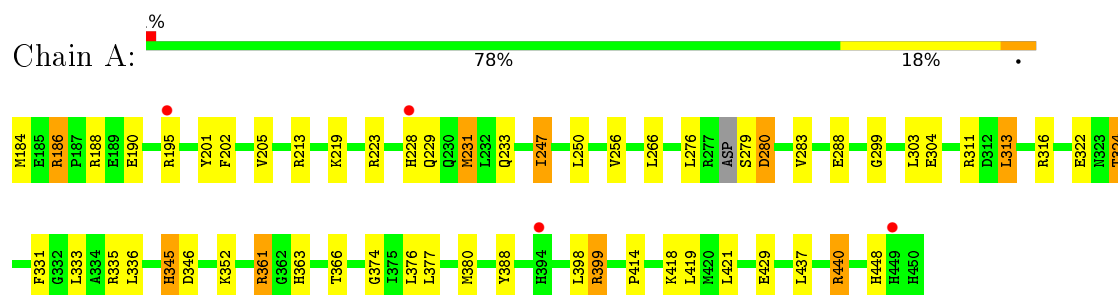
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	138	Total 138	O 138	0	0
4	B	120	Total 120	O 120	0	0
4	C	112	Total 112	O 112	0	0
4	D	131	Total 131	O 131	0	0

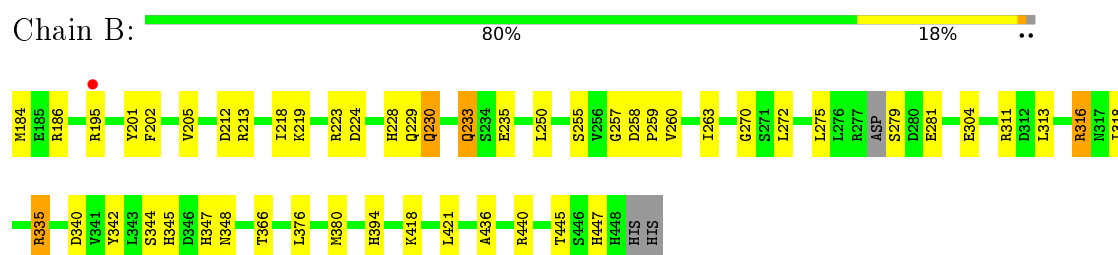
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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 ( $\text{RSRZ} > 2$ ). 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.

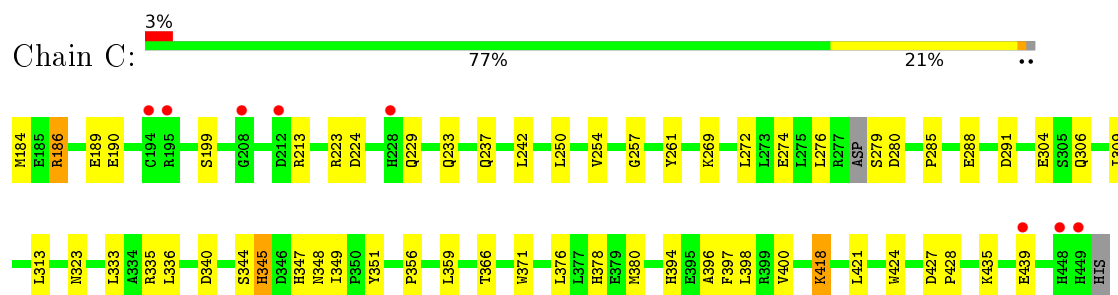
#### • Molecule 1: Protein-tyrosine kinase 6



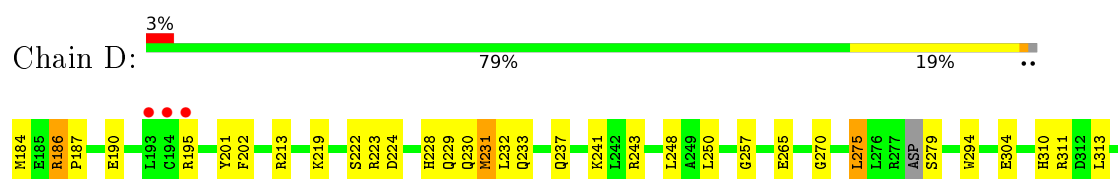
#### • Molecule 1: Protein-tyrosine kinase 6

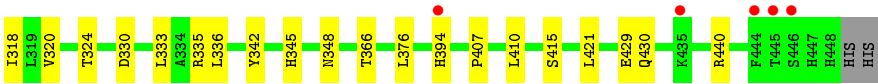


#### • Molecule 1: Protein-tyrosine kinase 6



#### • Molecule 1: Protein-tyrosine kinase 6





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	49.88Å 77.02Å 87.41Å 81.36° 75.76° 74.92°	Depositor
Resolution (Å)	84.52 – 2.24 42.18 – 2.24	Depositor EDS
% Data completeness (in resolution range)	94.5 (84.52-2.24) 92.5 (42.18-2.24)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.68 (at 2.24Å)	Xtriage
Refinement program	REFMAC	Depositor
R, $R_{free}$	0.205 , 0.260 0.204 , 0.259	Depositor DCC
$R_{free}$ test set	2789 reflections (5.34%)	DCC
Wilson B-factor (Å <sup>2</sup> )	34.8	Xtriage
Anisotropy	0.249	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 36.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	9541	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.20% 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, 1N1, CXM

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	0/2285	0.64	0/3087
1	B	0.47	0/2282	0.63	0/3085
1	C	0.45	0/2280	0.61	0/3081
1	D	0.45	0/2266	0.63	1/3063 (0.0%)
All	All	0.46	0/9113	0.63	1/12316 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	D	275	LEU	CA-CB-CG	5.90	128.87	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	184	CXM	Mainchain

## 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	2211	0	2219	41	0
1	B	2202	0	2202	37	0
1	C	2201	0	2200	45	0
1	D	2190	0	2188	36	0
2	A	41	0	14	0	0
2	B	41	0	14	0	0
2	C	41	0	14	0	0
2	D	41	0	14	0	0
3	A	18	0	24	1	0
3	B	18	0	24	5	0
3	C	18	0	24	1	0
3	D	18	0	24	0	0
4	A	138	0	0	4	0
4	B	120	0	0	2	0
4	C	112	0	0	6	0
4	D	131	0	0	3	0
All	All	9541	0	8961	155	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 9.

The worst 5 of 155 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:C:276:LEU:O	1:C:279:SER:HB3	1.55	1.05
1:A:399:ARG:HH11	1:A:399:ARG:HG2	1.18	1.04
1:C:223:ARG:NH2	1:C:229:GLN:OE1	1.94	1.00
1:B:223:ARG:NH2	1:B:229:GLN:OE1	1.96	0.99
1:D:223:ARG:HH22	1:D:229:GLN:HE22	1.14	0.96

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	272/267 (102%)	265 (97%)	7 (3%)	0	100	100
1	B	272/267 (102%)	264 (97%)	8 (3%)	0	100	100
1	C	273/267 (102%)	266 (97%)	7 (3%)	0	100	100
1	D	271/267 (102%)	265 (98%)	6 (2%)	0	100	100
All	All	1088/1068 (102%)	1060 (97%)	28 (3%)	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	244/236 (103%)	220 (90%)	24 (10%)	10	6
1	B	243/236 (103%)	230 (95%)	13 (5%)	28	28
1	C	242/236 (102%)	230 (95%)	12 (5%)	30	31
1	D	242/236 (102%)	228 (94%)	14 (6%)	25	24
All	All	971/944 (103%)	908 (94%)	63 (6%)	22	19

5 of 63 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	233	GLN
1	B	418	LYS
1	D	324	THR

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Mol	Chain	Res	Type
1	B	250	LEU
1	B	313	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 17 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	317	ASN
1	C	348	ASN
1	D	348	ASN
1	B	317	ASN
1	D	385	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 non-standard protein/DNA/RNA residues 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	$\# Z  > 2$
1	CXM	A	184	1	6,10,11	0.62	0	5,11,13	1.20	1 (20%)
1	CXM	B	184	1	6,10,11	0.55	0	5,11,13	0.99	0
1	CXM	C	184	1	6,10,11	0.63	0	5,11,13	1.08	0
1	CXM	D	184	1	6,10,11	0.69	0	5,11,13	1.05	1 (20%)

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
1	CXM	A	184	1	-	0/6/10/12	0/0/0/0
1	CXM	B	184	1	-	0/6/10/12	0/0/0/0
1	CXM	C	184	1	-	0/6/10/12	0/0/0/0
1	CXM	D	184	1	-	0/6/10/12	0/0/0/0

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	A	184	CXM	O-C-CA	-2.44	119.02	125.69
1	D	184	CXM	O-C-CA	-2.14	119.84	125.69

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	184	CXM	1	0
1	C	184	CXM	1	0
1	D	184	CXM	1	0

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

20 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	1N1	A	501[A]	-	32,36,36	0.75	0	39,50,50	1.56	8 (20%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	1N1	A	501[B]	-	32,36,36	0.78	0	39,50,50	1.61	9 (23%)
3	GOL	A	502	-	5,5,5	0.40	0	5,5,5	0.35	0
3	GOL	A	503	-	5,5,5	0.32	0	5,5,5	0.24	0
3	GOL	A	504	-	5,5,5	0.27	0	5,5,5	0.55	0
2	1N1	B	501[A]	-	32,36,36	0.73	0	39,50,50	1.51	8 (20%)
2	1N1	B	501[B]	-	32,36,36	0.75	0	39,50,50	1.54	9 (23%)
3	GOL	B	502	-	5,5,5	0.35	0	5,5,5	0.36	0
3	GOL	B	503	-	5,5,5	0.28	0	5,5,5	0.48	0
3	GOL	B	504	-	5,5,5	0.28	0	5,5,5	0.49	0
2	1N1	C	501[A]	-	32,36,36	0.78	0	39,50,50	1.87	11 (28%)
2	1N1	C	501[B]	-	32,36,36	0.80	0	39,50,50	1.80	9 (23%)
3	GOL	C	502	-	5,5,5	0.33	0	5,5,5	0.37	0
3	GOL	C	503	-	5,5,5	0.33	0	5,5,5	0.30	0
3	GOL	C	504	-	5,5,5	0.40	0	5,5,5	0.31	0
3	GOL	D	501	-	5,5,5	0.26	0	5,5,5	0.53	0
3	GOL	D	502	-	5,5,5	0.29	0	5,5,5	0.39	0
3	GOL	D	503	-	5,5,5	0.20	0	5,5,5	0.74	0
2	1N1	D	504[A]	-	32,36,36	0.79	0	39,50,50	1.45	7 (17%)
2	1N1	D	504[B]	-	32,36,36	0.81	0	39,50,50	1.53	8 (20%)

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	1N1	A	501[A]	-	-	0/13/29/29	0/4/4/4
2	1N1	A	501[B]	-	-	0/13/29/29	0/4/4/4
3	GOL	A	502	-	-	0/4/4/4	0/0/0/0
3	GOL	A	503	-	-	0/4/4/4	0/0/0/0
3	GOL	A	504	-	-	0/4/4/4	0/0/0/0
2	1N1	B	501[A]	-	-	0/13/29/29	0/4/4/4
2	1N1	B	501[B]	-	-	0/13/29/29	0/4/4/4
3	GOL	B	502	-	-	0/4/4/4	0/0/0/0
3	GOL	B	503	-	-	0/4/4/4	0/0/0/0
3	GOL	B	504	-	-	0/4/4/4	0/0/0/0
2	1N1	C	501[A]	-	-	0/13/29/29	0/4/4/4
2	1N1	C	501[B]	-	-	0/13/29/29	0/4/4/4
3	GOL	C	502	-	-	0/4/4/4	0/0/0/0
3	GOL	C	503	-	-	0/4/4/4	0/0/0/0
3	GOL	C	504	-	-	0/4/4/4	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	D	501	-	-	0/4/4/4	0/0/0/0
3	GOL	D	502	-	-	0/4/4/4	0/0/0/0
3	GOL	D	503	-	-	0/4/4/4	0/0/0/0
2	1N1	D	504[A]	-	-	0/13/29/29	0/4/4/4
2	1N1	D	504[B]	-	-	0/13/29/29	0/4/4/4

There are no bond length outliers.

The worst 5 of 69 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	501[B]	1N1	C12-C13-N5	-4.11	117.08	122.27
2	C	501[A]	1N1	C12-C13-N5	-4.11	117.08	122.27
2	C	501[A]	1N1	C5-C4-N2	-3.33	117.80	121.23
2	C	501[B]	1N1	C17-C16-N5	-3.14	104.73	110.59
2	C	501[A]	1N1	C17-C16-N5	-3.14	104.73	110.59

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	504	GOL	1	0
3	B	502	GOL	1	0
3	B	503	GOL	3	0
3	B	504	GOL	1	0
3	C	504	GOL	1	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	265/267 (99%)	0.05	4 (1%) 76 78	23, 31, 42, 55	0
1	B	263/267 (98%)	-0.02	1 (0%) 93 94	24, 31, 41, 50	0
1	C	264/267 (98%)	0.24	8 (3%) 54 57	26, 33, 44, 54	0
1	D	263/267 (98%)	0.27	8 (3%) 54 57	26, 34, 44, 52	0
All	All	1055/1068 (98%)	0.13	21 (1%) 68 70	23, 32, 44, 55	0

The worst 5 of 21 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	444	PHE	4.5
1	D	195	ARG	4.4
1	C	195[A]	ARG	3.9
1	A	449	HIS	3.4
1	C	448	HIS	3.3

### 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
1	CXM	D	184	11/12	0.92	0.15	-	41,42,46,46	0
1	CXM	B	184	11/12	0.96	0.11	-	36,37,43,43	0
1	CXM	C	184	11/12	0.89	0.19	-	42,43,44,46	0
1	CXM	A	184	11/12	0.95	0.13	-	38,39,40,41	0

### 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

### 6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	GOL	C	502	6/6	0.93	0.18	4.09	36,38,38,40	0
3	GOL	D	501	6/6	0.96	0.21	2.71	39,40,40,40	0
3	GOL	B	504	6/6	0.93	0.20	2.66	35,37,39,41	0
3	GOL	A	502	6/6	0.95	0.17	1.62	36,38,39,40	0
3	GOL	C	503	6/6	0.88	0.17	0.90	49,50,51,52	0
2	1N1	B	501[B]	33/33	0.94	0.15	0.86	30,33,41,44	8
2	1N1	B	501[A]	33/33	0.94	0.15	0.86	29,33,41,44	8
3	GOL	D	503	6/6	0.91	0.18	0.75	32,34,38,39	0
3	GOL	A	503	6/6	0.88	0.15	0.64	51,52,53,53	0
3	GOL	A	504	6/6	0.91	0.15	0.54	35,38,38,40	0
2	1N1	C	501[A]	33/33	0.92	0.15	0.51	31,36,46,48	8
3	GOL	C	504	6/6	0.90	0.16	0.44	42,43,44,46	0
2	1N1	D	504[B]	33/33	0.90	0.17	0.42	32,37,47,48	8
2	1N1	D	504[A]	33/33	0.90	0.17	0.42	32,37,47,48	8
2	1N1	A	501[B]	33/33	0.94	0.14	-0.09	27,30,39,42	8
2	1N1	A	501[A]	33/33	0.94	0.14	-0.09	27,30,39,42	8
2	1N1	C	501[B]	33/33	0.92	0.15	-0.37	30,36,46,48	8
3	GOL	B	503	6/6	0.88	0.13	-0.46	44,46,46,46	0
3	GOL	B	502	6/6	0.97	0.12	-0.48	28,30,32,32	0
3	GOL	D	502	6/6	0.93	0.14	-0.86	47,48,49,49	0

### 6.5 Other polymers ⓘ

There are no such residues in this entry.