



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 06:48 PM GMT

PDB ID : 1CJB  
Title : MALARIAL PURINE PHOSPHORIBOSYLTRANSFERASE  
Authors : Shi, W.; Li, C.M.; Tyler, P.C.; Furneaux, R.H.; Cahill, S.M.; Girvin, M.E.;  
Grubmeyer, C.; Schramm, V.L.; Almo, S.C.  
Deposited on : 1999-04-08  
Resolution : 2.00 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  
<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 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

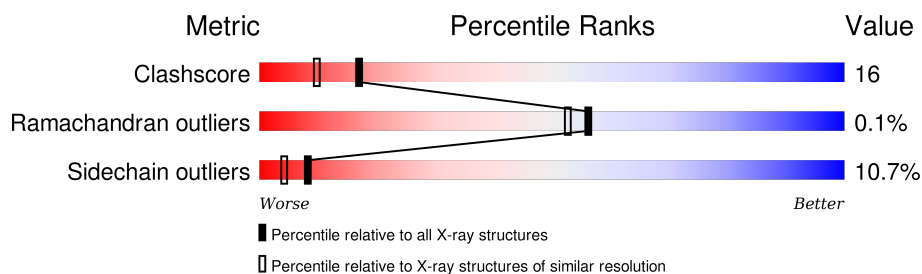
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.00 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))
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)

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. 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	231	
1	B	231	
1	C	231	
1	D	231	

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 7876 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.

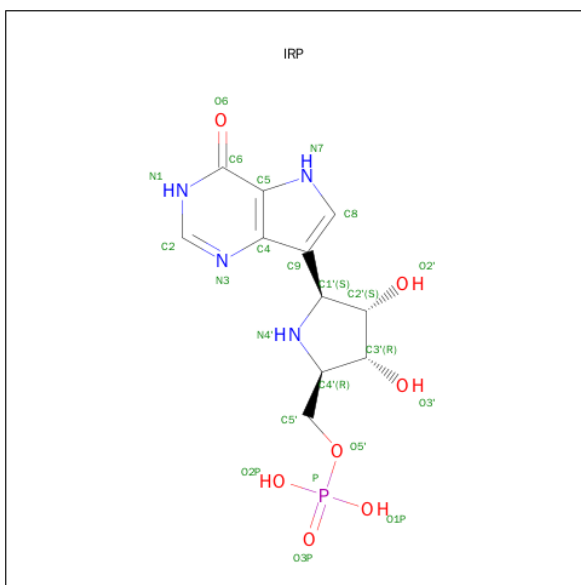
- Molecule 1 is a protein called PROTEIN (HYPOXANTHINE-GUANINE PHOSPHORIBOSYLTRANSFERASE).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	228	Total	C	N	O	S	0	0	1
			1828	1186	301	333	8			
1	B	228	Total	C	N	O	S	0	0	1
			1828	1186	301	333	8			
1	C	230	Total	C	N	O	S	0	0	0
			1849	1199	303	339	8			
1	D	228	Total	C	N	O	S	0	0	1
			1828	1186	301	333	8			

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

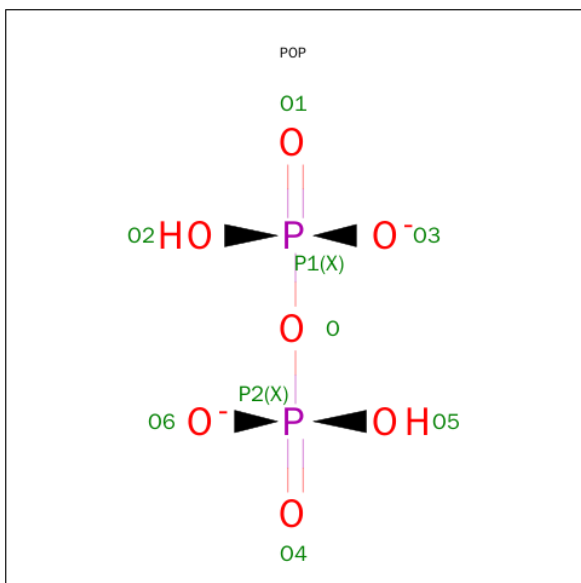
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	2	Total	Mg	0	0
			2	2		
2	A	2	Total	Mg	0	0
			2	2		
2	D	2	Total	Mg	0	0
			2	2		
2	C	2	Total	Mg	0	0
			2	2		

- Molecule 3 is (1S)-1(9-DEAZAHYPOXANTHIN-9YL)1,4-DIDEOXY-1,4-IMINO-D-RIBITOL-5-PHOSPHATE (three-letter code: IRP) (formula: C<sub>11</sub>H<sub>15</sub>N<sub>4</sub>O<sub>7</sub>P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			23	11	4	7	1		
3	B	1	Total	C	N	O	P	0	0
			23	11	4	7	1		
3	C	1	Total	C	N	O	P	0	0
			23	11	4	7	1		
3	D	1	Total	C	N	O	P	0	0
			23	11	4	7	1		

- Molecule 4 is PYROPHOSPHATE 2- (three-letter code: POP) (formula:  $\text{H}_2\text{O}_7\text{P}_2$ ).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total O P 9 7 2	0	0
4	B	1	Total O P 9 7 2	0	0
4	C	1	Total O P 9 7 2	0	0
4	D	1	Total O P 9 7 2	0	0

- Molecule 5 is water.

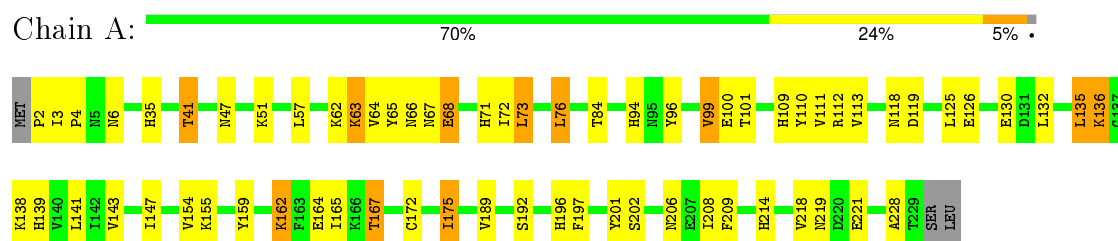
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	118	Total O 118 118	0	0
5	B	128	Total O 128 128	0	0
5	C	92	Total O 92 92	0	0
5	D	69	Total O 69 69	0	0

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

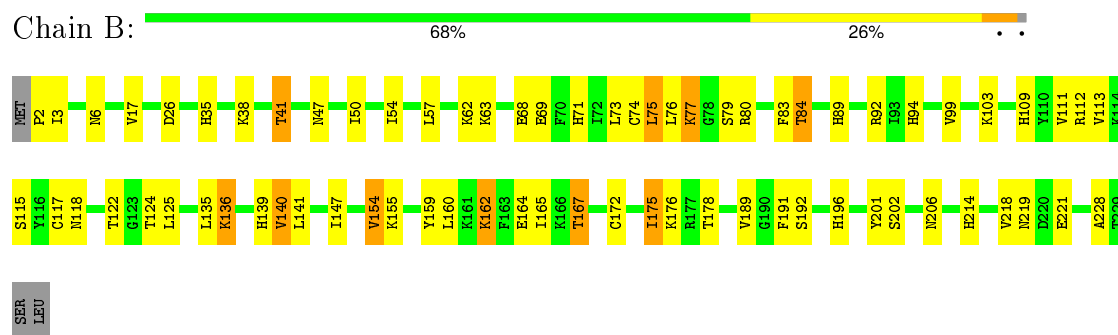
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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 ( $RSRZ > 2$ ). 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.

Note EDS was not executed.

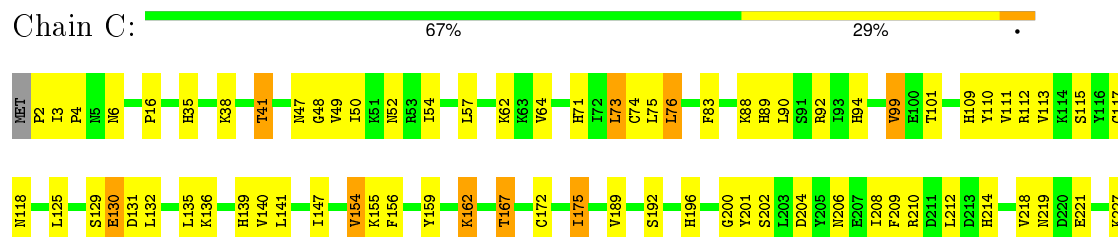
- Molecule 1: PROTEIN (HYPOXANTHINE-GUANINE PHOSPHORIBOSYLTRANSFERASE)



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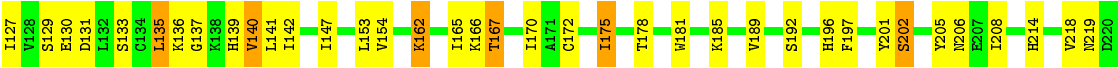


- Molecule 1: PROTEIN (HYPOXANTHINE-GUANINE PHOSPHORIBOSYLTRANSFERASE)





● Molecule 1: PROTEIN (HYPOXANTHINE-GUANINE PHOSPHORIBOSYLTRANSFERASE )



## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	105.49 Å   110.39 Å   173.70 Å 90.00°   90.00°   90.00°	Depositor
Resolution (Å)	20.00 – 2.00	Depositor
% Data completeness (in resolution range)	98.0 (20.00-2.00)	Depositor
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.05	Depositor
Refinement program	X-PLOR 3.8	Depositor
R, $R_{free}$	0.196 , 0.239	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7876	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	30.0	wwPDB-VP



## 5 Model quality

### 5.1 Standard geometry

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

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	0.37	0/1872	0.63	1/2529 (0.0%)
1	B	0.39	0/1872	0.64	1/2529 (0.0%)
1	C	0.35	0/1893	0.61	1/2556 (0.0%)
1	D	0.38	0/1872	0.61	1/2529 (0.0%)
All	All	0.37	0/7509	0.62	4/10143 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	2	PRO	N-CA-CB	5.30	109.66	103.30
1	A	2	PRO	N-CA-CB	5.18	109.51	103.30
1	C	2	PRO	N-CA-CB	5.06	109.37	103.30
1	B	2	PRO	N-CA-CB	5.05	109.36	103.30

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1828	0	1822	61	0
1	B	1828	0	1822	57	0
1	C	1849	0	1845	63	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	1828	0	1822	73	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
3	A	23	0	12	0	0
3	B	23	0	11	0	0
3	C	23	0	12	0	0
3	D	23	0	12	0	0
4	A	9	0	0	0	0
4	B	9	0	0	0	0
4	C	9	0	0	0	0
4	D	9	0	0	0	0
5	A	118	0	0	6	0
5	B	128	0	0	7	0
5	C	92	0	0	4	0
5	D	69	0	0	5	0
All	All	7876	0	7358	241	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 16.

All (241) 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:118:ASN:HD21	1:C:228:ALA:H	1.10	0.99
1:C:47:ASN:H	1:C:214:HIS:HD2	1.07	0.96
1:D:47:ASN:H	1:D:214:HIS:HD2	1.10	0.94
1:C:48:GLY:H	1:D:52:ASN:HD21	1.13	0.94
1:B:118:ASN:HD21	1:B:228:ALA:H	1.09	0.94
1:C:52:ASN:HD21	1:D:48:GLY:H	1.15	0.92
1:A:47:ASN:H	1:A:214:HIS:HD2	1.06	0.92
1:C:139:HIS:HD2	1:C:167:THR:HB	1.33	0.90
1:D:118:ASN:HD21	1:D:228:ALA:H	1.15	0.90
1:B:113:VAL:HG13	1:B:125:LEU:HD11	1.51	0.89
1:D:118:ASN:HD22	1:D:206:ASN:HD21	1.20	0.88
1:A:41:THR:HG22	1:A:219:ASN:HB3	1.56	0.88
1:B:47:ASN:H	1:B:214:HIS:HD2	1.23	0.87
1:D:111:VAL:HG13	1:D:129:SER:HB3	1.56	0.87
1:C:41:THR:HG22	1:C:219:ASN:HB3	1.59	0.84
1:D:47:ASN:H	1:D:214:HIS:CD2	1.95	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:41:THR:HG22	1:B:219:ASN:HB3	1.60	0.83
1:A:118:ASN:HD22	1:A:206:ASN:HD21	1.26	0.82
1:B:113:VAL:CG1	1:B:125:LEU:HD11	2.10	0.82
1:B:139:HIS:HD2	1:B:167:THR:HB	1.44	0.81
1:C:118:ASN:HD22	1:C:206:ASN:HD21	1.27	0.81
1:A:118:ASN:HD21	1:A:228:ALA:H	1.24	0.81
1:C:47:ASN:H	1:C:214:HIS:CD2	1.98	0.79
1:A:139:HIS:HD2	1:A:167:THR:HB	1.48	0.78
1:C:113:VAL:CG1	1:C:125:LEU:HD11	2.15	0.77
1:C:52:ASN:ND2	1:D:48:GLY:H	1.81	0.76
1:C:48:GLY:H	1:D:52:ASN:ND2	1.84	0.76
1:A:118:ASN:ND2	1:A:206:ASN:HD21	1.83	0.76
1:B:38:LYS:HE3	5:B:804:HOH:O	1.87	0.74
1:D:118:ASN:ND2	1:D:206:ASN:HD21	1.85	0.74
1:A:47:ASN:H	1:A:214:HIS:CD2	1.99	0.73
1:C:62:LYS:NZ	1:C:94:HIS:HD2	1.86	0.72
1:C:49:VAL:HG22	1:D:49:VAL:HG23	1.70	0.72
1:A:71:HIS:HD2	1:A:109:HIS:HE1	1.38	0.71
1:C:167:THR:HG22	5:C:519:HOH:O	1.91	0.70
1:A:196:HIS:HE1	5:A:775:HOH:O	1.73	0.70
1:D:139:HIS:HA	1:D:167:THR:HG22	1.74	0.70
1:D:111:VAL:HG22	1:D:129:SER:HB2	1.74	0.69
1:C:71:HIS:CD2	1:C:109:HIS:HE1	2.09	0.69
1:D:71:HIS:HD2	1:D:109:HIS:HE1	1.41	0.69
1:B:118:ASN:ND2	1:B:206:ASN:HD21	1.91	0.68
1:C:118:ASN:ND2	1:C:228:ALA:H	1.88	0.68
1:B:118:ASN:HD22	1:B:206:ASN:HD21	1.42	0.67
1:D:41:THR:HG22	1:D:219:ASN:HB3	1.77	0.67
1:D:71:HIS:CD2	1:D:109:HIS:HE1	2.12	0.66
1:B:71:HIS:HD2	1:B:109:HIS:HE1	1.41	0.66
1:C:201:TYR:O	1:C:202:SER:HB2	1.95	0.66
1:C:71:HIS:HD2	1:C:109:HIS:HE1	1.41	0.66
1:A:71:HIS:CD2	1:A:109:HIS:HE1	2.14	0.66
1:C:113:VAL:HG13	1:C:125:LEU:HD11	1.78	0.66
1:C:52:ASN:HD21	1:D:48:GLY:N	1.92	0.65
1:C:38:LYS:HE3	5:C:587:HOH:O	1.96	0.65
1:C:49:VAL:HG22	1:D:49:VAL:CG2	2.27	0.65
1:C:147:ILE:HB	1:C:175:ILE:HG12	1.78	0.65
1:C:118:ASN:ND2	1:C:206:ASN:HD21	1.94	0.65
1:A:41:THR:HG21	5:A:901:HOH:O	1.97	0.64
1:A:3:ILE:HG21	1:A:167:THR:HG21	1.80	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:139:HIS:CD2	1:B:167:THR:HB	2.30	0.63
1:B:201:TYR:O	1:B:202:SER:HB2	1.98	0.62
1:B:62:LYS:NZ	1:B:94:HIS:HD2	1.97	0.62
1:A:62:LYS:NZ	1:A:94:HIS:HD2	1.96	0.62
1:D:175:ILE:HG22	1:D:192:SER:HA	1.82	0.61
1:A:136:LYS:HG3	1:A:164:GLU:HB2	1.81	0.61
1:B:47:ASN:H	1:B:214:HIS:CD2	2.11	0.61
1:C:62:LYS:HZ3	1:C:94:HIS:HD2	1.46	0.61
1:B:71:HIS:CD2	1:B:109:HIS:HE1	2.18	0.60
1:D:62:LYS:NZ	1:D:94:HIS:HD2	1.99	0.60
1:D:201:TYR:O	1:D:202:SER:HB2	2.01	0.60
1:A:147:ILE:HG13	1:A:172:CYS:SG	2.42	0.60
1:B:75:LEU:O	1:B:79:SER:HB2	2.01	0.59
1:A:41:THR:HG22	1:A:219:ASN:CB	2.32	0.59
1:C:73:LEU:HD21	1:C:111:VAL:HG23	1.85	0.59
1:B:160:LEU:HD22	1:B:165:ILE:CD1	2.33	0.58
1:B:139:HIS:HD2	1:B:167:THR:CB	2.15	0.58
1:C:48:GLY:N	1:D:52:ASN:HD21	1.94	0.58
1:C:74:CYS:HB2	1:C:83:PHE:CD1	2.38	0.58
1:D:181:TRP:HH2	1:D:185:LYS:HE3	1.68	0.58
1:B:118:ASN:HD21	1:B:228:ALA:N	1.92	0.58
1:B:167:THR:CG2	5:B:506:HOH:O	2.52	0.57
1:A:167:THR:CG2	5:A:525:HOH:O	2.51	0.57
1:A:35:HIS:CD2	1:A:35:HIS:H	2.22	0.57
1:D:47:ASN:N	1:D:214:HIS:HD2	1.91	0.57
1:B:35:HIS:CD2	1:B:35:HIS:H	2.23	0.56
1:B:175:ILE:HG22	1:B:192:SER:HA	1.87	0.56
1:C:16:PRO:HB3	1:C:189:VAL:HG13	1.87	0.56
1:B:69:GLU:OE2	1:B:71:HIS:HE1	1.87	0.56
1:A:147:ILE:HB	1:A:175:ILE:HG12	1.87	0.56
1:D:118:ASN:HD21	1:D:228:ALA:N	1.96	0.55
1:A:113:VAL:CG1	1:A:125:LEU:HD11	2.37	0.55
1:A:73:LEU:HD21	1:A:111:VAL:HG23	1.89	0.55
1:B:136:LYS:HE3	1:B:164:GLU:HB2	1.89	0.55
1:C:147:ILE:HG13	1:C:172:CYS:SG	2.47	0.55
1:D:116:TYR:CE1	1:D:121:SER:HB2	2.41	0.55
1:A:100:GLU:OE2	1:B:38:LYS:HE2	2.07	0.54
1:B:122:THR:OG1	1:B:124:THR:HG22	2.08	0.54
1:B:176:LYS:HE2	1:B:196:HIS:O	2.08	0.54
1:A:63:LYS:HE3	5:A:710:HOH:O	2.07	0.54
1:D:147:ILE:HG12	1:D:153:LEU:HD12	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:139:HIS:HD2	1:D:167:THR:CG2	2.21	0.53
1:A:139:HIS:HD2	1:A:167:THR:CB	2.20	0.53
1:A:3:ILE:CG2	1:A:167:THR:HG21	2.38	0.53
1:B:147:ILE:HB	1:B:175:ILE:HG12	1.89	0.53
1:C:196:HIS:HE1	5:C:630:HOH:O	1.91	0.53
1:C:139:HIS:CD2	1:C:167:THR:HB	2.26	0.53
1:B:89:HIS:HD2	1:B:92:ARG:HH21	1.57	0.52
1:D:139:HIS:HD2	1:D:167:THR:HB	1.75	0.52
1:D:47:ASN:O	1:D:51:LYS:HG2	2.10	0.52
1:D:62:LYS:HZ1	1:D:94:HIS:CD2	2.28	0.52
1:C:88:LYS:HE3	1:C:92:ARG:HH12	1.74	0.52
1:C:172:CYS:O	1:C:189:VAL:HA	2.10	0.51
1:B:3:ILE:HG21	1:B:167:THR:HG21	1.91	0.51
1:A:84:THR:HG21	1:B:84:THR:HG23	1.91	0.51
1:A:84:THR:HG21	1:B:84:THR:CG2	2.40	0.51
1:D:35:HIS:CD2	1:D:35:HIS:H	2.27	0.51
1:D:62:LYS:HZ1	1:D:94:HIS:HD2	1.59	0.51
1:A:139:HIS:CD2	1:A:167:THR:HB	2.37	0.51
1:C:41:THR:HG22	1:C:219:ASN:CB	2.37	0.51
1:A:201:TYR:O	1:A:202:SER:HB2	2.11	0.51
1:D:113:VAL:CG1	1:D:125:LEU:HD11	2.41	0.51
1:D:47:ASN:HD21	1:D:51:LYS:HZ2	1.58	0.50
1:B:68:GLU:HG2	5:B:733:HOH:O	2.10	0.50
1:D:71:HIS:HD2	1:D:109:HIS:CE1	2.27	0.50
1:D:99:VAL:HG13	1:D:100:GLU:N	2.26	0.50
1:A:62:LYS:HZ2	1:A:94:HIS:HD2	1.58	0.50
1:D:137:GLY:C	1:D:166:LYS:HB2	2.32	0.50
1:D:73:LEU:CD2	1:D:111:VAL:HG23	2.42	0.50
1:B:154:VAL:HG12	1:B:155:LYS:N	2.26	0.50
1:B:73:LEU:HD21	1:B:111:VAL:CG2	2.42	0.49
1:B:74:CYS:HB2	1:B:83:PHE:CE1	2.47	0.49
1:B:62:LYS:HZ3	1:B:94:HIS:HD2	1.58	0.49
1:D:115:SER:HB3	1:D:125:LEU:HD12	1.94	0.49
1:C:50:ILE:O	1:C:54:ILE:HG13	2.13	0.49
1:C:74:CYS:HB2	1:C:83:PHE:CE1	2.49	0.48
1:A:72:ILE:HG23	1:A:143:VAL:CG2	2.44	0.48
1:D:172:CYS:O	1:D:189:VAL:HA	2.14	0.48
1:B:26:ASP:OD2	5:B:903:HOH:O	2.20	0.48
1:D:139:HIS:CD2	1:D:167:THR:CG2	2.97	0.48
1:D:196:HIS:HE1	5:D:884:HOH:O	1.96	0.48
1:A:154:VAL:HG13	1:A:155:LYS:N	2.29	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:175:ILE:HG22	1:C:192:SER:HA	1.96	0.47
1:A:35:HIS:CE1	5:B:674:HOH:O	2.67	0.47
1:A:99:VAL:HG13	1:A:100:GLU:N	2.29	0.47
1:C:99:VAL:HG13	1:C:101:THR:H	1.79	0.47
1:D:140:VAL:HG13	1:D:165:ILE:HG21	1.96	0.47
1:A:72:ILE:HG23	1:A:143:VAL:HG23	1.96	0.47
1:D:139:HIS:CA	1:D:167:THR:HG22	2.43	0.47
1:C:89:HIS:HD2	1:C:92:ARG:NH2	2.13	0.47
1:A:66:ASN:HB3	1:A:67:ASN:H	1.54	0.47
1:D:122:THR:OG1	1:D:124:THR:HG22	2.15	0.47
1:C:208:ILE:O	1:C:209:PHE:HB2	2.15	0.47
1:C:62:LYS:NZ	1:C:94:HIS:CD2	2.76	0.47
1:C:3:ILE:CG2	1:C:167:THR:HG21	2.45	0.46
1:A:110:TYR:OH	1:B:80:ARG:HG3	2.15	0.46
1:D:167:THR:HG23	5:D:571:HOH:O	2.15	0.46
1:C:49:VAL:CG2	1:D:49:VAL:CG2	2.93	0.46
1:C:35:HIS:H	1:C:35:HIS:CD2	2.31	0.46
1:A:113:VAL:HG12	1:A:125:LEU:HD11	1.97	0.46
1:A:135:LEU:HB3	1:A:165:ILE:HG12	1.96	0.46
1:D:208:ILE:HG13	1:D:226:TYR:CD1	2.50	0.46
1:A:175:ILE:HG22	1:A:192:SER:HA	1.97	0.46
1:A:172:CYS:O	1:A:189:VAL:HA	2.16	0.46
1:A:73:LEU:CD2	1:A:111:VAL:HG23	2.44	0.46
1:B:89:HIS:CD2	1:B:92:ARG:HH21	2.33	0.46
1:D:113:VAL:HG22	1:D:127:ILE:HG12	1.97	0.46
1:A:76:LEU:HD22	1:A:110:TYR:HB3	1.98	0.46
1:A:63:LYS:HE2	5:A:510:HOH:O	2.16	0.46
1:C:159:TYR:O	1:C:162:LYS:HG3	2.15	0.46
1:B:159:TYR:O	1:B:162:LYS:HG3	2.15	0.46
1:D:76:LEU:HA	1:D:77:LYS:HA	1.59	0.45
1:A:136:LYS:HE3	1:A:164:GLU:HG3	1.98	0.45
1:B:89:HIS:CD2	1:B:92:ARG:NH2	2.84	0.45
1:A:154:VAL:CG1	1:A:155:LYS:N	2.79	0.45
1:A:119:ASP:HA	1:A:197:PHE:CE1	2.52	0.45
1:D:135:LEU:HB3	1:D:165:ILE:HG12	1.99	0.45
1:C:115:SER:HB3	1:C:125:LEU:HD12	1.99	0.45
1:D:140:VAL:HG13	1:D:165:ILE:CG2	2.47	0.45
1:C:154:VAL:HG12	1:C:155:LYS:N	2.31	0.45
1:D:63:LYS:HE2	5:D:589:HOH:O	2.16	0.44
1:C:117:CYS:O	1:C:118:ASN:HB2	2.18	0.44
1:B:117:CYS:O	1:B:118:ASN:HB2	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:117:CYS:O	1:D:118:ASN:HB2	2.18	0.44
1:A:118:ASN:HD22	1:A:206:ASN:ND2	2.06	0.44
1:A:136:LYS:HE3	1:A:164:GLU:HB2	1.99	0.44
1:B:140:VAL:HG13	1:B:165:ILE:CG2	2.47	0.44
5:A:619:HOH:O	1:B:77:LYS:HB2	2.17	0.44
1:B:69:GLU:OE2	1:B:71:HIS:CE1	2.69	0.44
1:C:3:ILE:HG22	1:C:167:THR:HG21	1.99	0.44
1:B:50:ILE:O	1:B:54:ILE:HG13	2.18	0.44
1:D:162:LYS:HG2	1:D:162:LYS:H	1.39	0.44
1:A:4:PRO:HD3	1:A:64:VAL:HG11	2.00	0.44
1:A:47:ASN:HD21	1:A:51:LYS:NZ	2.16	0.44
1:D:76:LEU:HD22	1:D:110:TYR:HB3	1.99	0.43
1:D:3:ILE:HG21	1:D:167:THR:HG21	1.99	0.43
1:D:167:THR:CG2	5:D:571:HOH:O	2.65	0.43
1:D:72:ILE:CG2	1:D:83:PHE:HE1	2.32	0.43
1:C:130:GLU:O	1:C:132:LEU:HG	2.19	0.43
1:C:200:GLY:HA2	1:C:212:LEU:HB3	2.01	0.43
1:A:159:TYR:O	1:A:162:LYS:HG3	2.18	0.43
1:C:113:VAL:HG13	1:C:125:LEU:CD1	2.47	0.43
1:C:167:THR:CG2	5:C:519:HOH:O	2.59	0.42
1:B:167:THR:HG22	5:B:506:HOH:O	2.18	0.42
1:B:76:LEU:HA	1:B:77:LYS:HA	1.70	0.42
1:A:51:LYS:HZ2	1:A:51:LYS:HG3	1.63	0.42
1:B:136:LYS:HE2	1:B:164:GLU:OE2	2.19	0.42
1:D:111:VAL:HG22	1:D:129:SER:CB	2.46	0.42
1:A:109:HIS:HD2	1:A:130:GLU:OE2	2.02	0.42
1:D:139:HIS:HD2	1:D:167:THR:CB	2.32	0.42
1:C:130:GLU:O	1:C:132:LEU:N	2.52	0.42
1:B:196:HIS:HE1	5:B:683:HOH:O	2.01	0.42
1:C:162:LYS:H	1:C:162:LYS:HG2	1.56	0.42
1:D:175:ILE:CG2	1:D:175:ILE:O	2.68	0.42
1:B:17:VAL:HB	1:B:191:PHE:CD2	2.54	0.42
1:C:4:PRO:HD3	1:C:64:VAL:HG11	2.02	0.41
1:B:3:ILE:CG2	1:B:167:THR:HG21	2.50	0.41
1:A:99:VAL:HG13	1:A:101:THR:H	1.84	0.41
1:B:3:ILE:HD11	1:B:68:GLU:HG2	2.02	0.41
1:D:119:ASP:HA	1:D:197:PHE:CE1	2.55	0.41
1:D:73:LEU:HD21	1:D:111:VAL:HG23	2.02	0.41
1:D:41:THR:HG22	1:D:219:ASN:CB	2.47	0.41
1:C:76:LEU:HD22	1:C:110:TYR:HB3	2.03	0.41
1:A:208:ILE:O	1:A:209:PHE:HB2	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:47:ASN:ND2	5:D:888:HOH:O	2.44	0.41
1:A:71:HIS:CE1	1:A:138:LYS:HD2	2.55	0.41
1:D:147:ILE:CG1	1:D:153:LEU:HD12	2.50	0.41
1:B:115:SER:HB3	1:B:125:LEU:HD12	2.03	0.41
1:C:111:VAL:HG11	1:C:156:PHE:CE1	2.55	0.41
1:A:99:VAL:CG1	1:A:101:THR:H	2.34	0.41
1:C:89:HIS:CD2	1:C:92:ARG:NH2	2.89	0.41
1:C:204:ASP:HB3	1:C:210:ARG:HG2	2.03	0.41
1:B:74:CYS:HB2	1:B:83:PHE:CD1	2.56	0.40
1:D:99:VAL:CG1	1:D:100:GLU:N	2.84	0.40
1:A:65:TYR:O	1:A:68:GLU:HB2	2.21	0.40
1:A:139:HIS:HD2	1:A:167:THR:CG2	2.34	0.40
1:D:205:TYR:CE2	1:D:223:LYS:HA	2.56	0.40
1:D:131:ASP:OD1	1:D:133:SER:HB3	2.20	0.40
1:D:142:ILE:HD12	1:D:170:ILE:HD13	2.02	0.40
1:C:111:VAL:HG12	1:C:113:VAL:HG23	2.03	0.40
1:D:99:VAL:CG1	1:D:101:THR:H	2.35	0.40
1:A:132:LEU:HD12	1:A:159:TYR:HE2	1.85	0.40
1:C:90:LEU:HD23	1:C:90:LEU:HA	1.96	0.40
1:B:172:CYS:O	1:B:189:VAL:HA	2.22	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	226/231 (98%)	217 (96%)	9 (4%)	0	100	100
1	B	226/231 (98%)	218 (96%)	8 (4%)	0	100	100
1	C	228/231 (99%)	219 (96%)	8 (4%)	1 (0%)	39	33
1	D	226/231 (98%)	216 (96%)	10 (4%)	0	100	100
All	All	906/924 (98%)	870 (96%)	35 (4%)	1 (0%)	56	53



All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	131	ASP

### 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	202/207 (98%)	183 (91%)	19 (9%)	11	6
1	B	202/207 (98%)	181 (90%)	21 (10%)	9	4
1	C	205/207 (99%)	183 (89%)	22 (11%)	8	4
1	D	202/207 (98%)	177 (88%)	25 (12%)	6	3
All	All	811/828 (98%)	724 (89%)	87 (11%)	8	4

All (87) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	ASN
1	A	41	THR
1	A	57	LEU
1	A	63	LYS
1	A	68	GLU
1	A	73	LEU
1	A	76	LEU
1	A	96	TYR
1	A	99	VAL
1	A	112	ARG
1	A	126	GLU
1	A	135	LEU
1	A	136	LYS
1	A	141	LEU
1	A	162	LYS
1	A	167	THR
1	A	175	ILE
1	A	218	VAL
1	A	221	GLU

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Mol	Chain	Res	Type
1	B	6	ASN
1	B	41	THR
1	B	57	LEU
1	B	63	LYS
1	B	75	LEU
1	B	77	LYS
1	B	84	THR
1	B	99	VAL
1	B	103	LYS
1	B	112	ARG
1	B	135	LEU
1	B	136	LYS
1	B	140	VAL
1	B	141	LEU
1	B	154	VAL
1	B	162	LYS
1	B	167	THR
1	B	175	ILE
1	B	178	THR
1	B	218	VAL
1	B	221	GLU
1	C	6	ASN
1	C	41	THR
1	C	57	LEU
1	C	73	LEU
1	C	75	LEU
1	C	76	LEU
1	C	99	VAL
1	C	112	ARG
1	C	129	SER
1	C	130	GLU
1	C	135	LEU
1	C	136	LYS
1	C	140	VAL
1	C	141	LEU
1	C	154	VAL
1	C	162	LYS
1	C	167	THR
1	C	175	ILE
1	C	218	VAL
1	C	221	GLU
1	C	227	LYS

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Mol	Chain	Res	Type
1	C	230	SER
1	D	6	ASN
1	D	41	THR
1	D	57	LEU
1	D	69	GLU
1	D	73	LEU
1	D	75	LEU
1	D	76	LEU
1	D	77	LYS
1	D	96	TYR
1	D	99	VAL
1	D	112	ARG
1	D	124	THR
1	D	130	GLU
1	D	135	LEU
1	D	136	LYS
1	D	140	VAL
1	D	141	LEU
1	D	154	VAL
1	D	162	LYS
1	D	167	THR
1	D	175	ILE
1	D	178	THR
1	D	202	SER
1	D	218	VAL
1	D	221	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (48) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	6	ASN
1	A	20	ASN
1	A	35	HIS
1	A	47	ASN
1	A	71	HIS
1	A	89	HIS
1	A	94	HIS
1	A	109	HIS
1	A	118	ASN
1	A	139	HIS
1	A	196	HIS
1	A	214	HIS

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Mol	Chain	Res	Type
1	B	6	ASN
1	B	20	ASN
1	B	35	HIS
1	B	47	ASN
1	B	71	HIS
1	B	89	HIS
1	B	94	HIS
1	B	109	HIS
1	B	118	ASN
1	B	139	HIS
1	B	196	HIS
1	B	214	HIS
1	C	6	ASN
1	C	20	ASN
1	C	35	HIS
1	C	47	ASN
1	C	52	ASN
1	C	71	HIS
1	C	89	HIS
1	C	94	HIS
1	C	109	HIS
1	C	118	ASN
1	C	139	HIS
1	C	196	HIS
1	C	214	HIS
1	D	6	ASN
1	D	20	ASN
1	D	35	HIS
1	D	47	ASN
1	D	52	ASN
1	D	71	HIS
1	D	94	HIS
1	D	109	HIS
1	D	118	ASN
1	D	139	HIS
1	D	214	HIS

### 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 carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 16 ligands modelled in this entry, 8 are monoatomic - leaving 8 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
3	IRP	A	300	2	24,25,25	2.27	8 (33%)	18,38,38	2.93	6 (33%)
4	POP	A	400	2	8,8,8	1.66	2 (25%)	13,13,13	0.76	0
3	IRP	B	300	2	24,25,25	2.47	8 (33%)	18,38,38	3.01	6 (33%)
4	POP	B	400	2	8,8,8	1.62	3 (37%)	13,13,13	0.78	0
3	IRP	C	300	2	24,25,25	2.30	8 (33%)	18,38,38	2.96	6 (33%)
4	POP	C	400	2	8,8,8	1.81	3 (37%)	13,13,13	0.70	0
3	IRP	D	300	2	24,25,25	2.38	10 (41%)	18,38,38	2.98	6 (33%)
4	POP	D	400	2	8,8,8	1.57	1 (12%)	13,13,13	0.71	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
3	IRP	A	300	2	-	0/6/26/26	0/3/3/3
4	POP	A	400	2	-	0/6/6/6	0/0/0/0
3	IRP	B	300	2	-	0/6/26/26	0/3/3/3
4	POP	B	400	2	-	0/6/6/6	0/0/0/0
3	IRP	C	300	2	-	0/6/26/26	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	POP	C	400	2	-	0/6/6/6	0/0/0/0
3	IRP	D	300	2	-	0/6/26/26	0/3/3/3
4	POP	D	400	2	-	0/6/6/6	0/0/0/0

All (43) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	300	IRP	C3'-C4'	-6.47	1.48	1.53
3	D	300	IRP	C3'-C4'	-4.90	1.49	1.53
3	C	300	IRP	C1'-N4'	-4.64	1.41	1.47
3	A	300	IRP	C3'-C4'	-4.33	1.50	1.53
3	D	300	IRP	C1'-N4'	-4.27	1.42	1.47
3	A	300	IRP	C1'-N4'	-4.26	1.42	1.47
3	C	300	IRP	C3'-C4'	-4.20	1.50	1.53
3	B	300	IRP	C1'-N4'	-3.90	1.42	1.47
3	C	300	IRP	C4'-N4'	-3.55	1.42	1.48
3	B	300	IRP	C8-N7	-3.26	1.30	1.36
3	D	300	IRP	C4'-N4'	-3.26	1.43	1.48
3	A	300	IRP	C8-N7	-3.24	1.30	1.36
3	A	300	IRP	C4'-N4'	-3.20	1.43	1.48
3	D	300	IRP	C8-N7	-3.10	1.30	1.36
3	B	300	IRP	C4'-N4'	-2.90	1.43	1.48
3	C	300	IRP	C8-N7	-2.76	1.31	1.36
4	C	400	POP	P1-O	-2.57	1.55	1.61
3	C	300	IRP	P-O2P	-2.47	1.45	1.54
3	D	300	IRP	P-O2P	-2.28	1.46	1.54
3	B	300	IRP	P-O2P	-2.26	1.46	1.54
4	A	400	POP	P1-O	-2.21	1.56	1.61
3	A	300	IRP	P-O2P	-2.18	1.46	1.54
4	B	400	POP	P2-O5	-2.17	1.46	1.55
4	B	400	POP	P1-O	-2.16	1.56	1.61
4	C	400	POP	P2-O5	-2.08	1.46	1.55
3	D	300	IRP	P-O1P	-2.02	1.47	1.54
3	D	300	IRP	C4-N3	2.20	1.40	1.37
4	B	400	POP	P2-O6	2.51	1.58	1.50
4	D	400	POP	P2-O6	2.70	1.59	1.50
4	A	400	POP	P2-O6	2.71	1.59	1.50
4	C	400	POP	P2-O6	2.87	1.60	1.50
3	B	300	IRP	C2-N3	2.88	1.37	1.32
3	A	300	IRP	C2-N3	2.89	1.37	1.32
3	D	300	IRP	C2-N3	2.94	1.37	1.32
3	C	300	IRP	C2-N3	2.95	1.37	1.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	300	IRP	C2-N1	3.75	1.41	1.33
3	C	300	IRP	C2-N1	3.90	1.41	1.33
3	D	300	IRP	C2-N1	3.96	1.41	1.33
3	D	300	IRP	C6-N1	4.19	1.40	1.33
3	B	300	IRP	C2-N1	4.23	1.41	1.33
3	C	300	IRP	C6-N1	4.24	1.41	1.33
3	A	300	IRP	C6-N1	4.40	1.41	1.33
3	B	300	IRP	C6-N1	4.53	1.41	1.33

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	300	IRP	N3-C2-N1	-8.66	122.26	128.89
3	D	300	IRP	N3-C2-N1	-8.66	122.27	128.89
3	C	300	IRP	N3-C2-N1	-8.64	122.28	128.89
3	A	300	IRP	N3-C2-N1	-8.55	122.34	128.89
3	C	300	IRP	C9-C1'-N4'	-4.51	106.66	113.38
3	B	300	IRP	C9-C1'-N4'	-4.44	106.78	113.38
3	D	300	IRP	C9-C1'-N4'	-4.18	107.16	113.38
3	A	300	IRP	C9-C1'-N4'	-4.12	107.24	113.38
3	B	300	IRP	C5'-C4'-C3'	-3.09	108.78	114.11
3	D	300	IRP	C5'-C4'-C3'	-2.52	109.77	114.11
3	A	300	IRP	C5'-C4'-C3'	-2.51	109.79	114.11
3	A	300	IRP	O5'-P-O3P	-2.23	101.48	107.14
3	C	300	IRP	C5'-C4'-C3'	-2.11	110.47	114.11
3	D	300	IRP	O5'-P-O3P	-2.11	101.77	107.14
3	C	300	IRP	O5'-P-O3P	-2.09	101.82	107.14
3	B	300	IRP	O5'-P-O3P	-2.04	101.96	107.14
3	C	300	IRP	C9-C8-N7	3.41	115.31	108.79
3	A	300	IRP	C9-C8-N7	3.43	115.33	108.79
3	D	300	IRP	C9-C8-N7	3.56	115.59	108.79
3	B	300	IRP	C9-C8-N7	3.71	115.86	108.79
3	B	300	IRP	C2-N1-C6	5.66	124.62	116.04
3	A	300	IRP	C2-N1-C6	5.69	124.66	116.04
3	C	300	IRP	C2-N1-C6	5.75	124.75	116.04
3	D	300	IRP	C2-N1-C6	5.79	124.82	116.04

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.7 Other polymers

There are no such residues in this entry.

## 5.8 Polymer linkage issues

There are no chain breaks in this entry.



## 6 Fit of model and data [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

### 6.3 Carbohydrates [i](#)

EDS was not executed - this section will therefore be empty.

### 6.4 Ligands [i](#)

EDS was not executed - this section will therefore be empty.

### 6.5 Other polymers [i](#)

EDS was not executed - this section will therefore be empty.