



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:46 PM GMT

PDB ID : 1H5R  
Title : THYMIDYLYLTRANSFERASE COMPLEXED WITH THIMIDINE AND  
GLUCOSE-1-PHOSPATE  
Authors : Rosano, C.; Zuccotti, S.; Bolognesi, M.  
Deposited on : 2001-05-25  
Resolution : 1.90 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

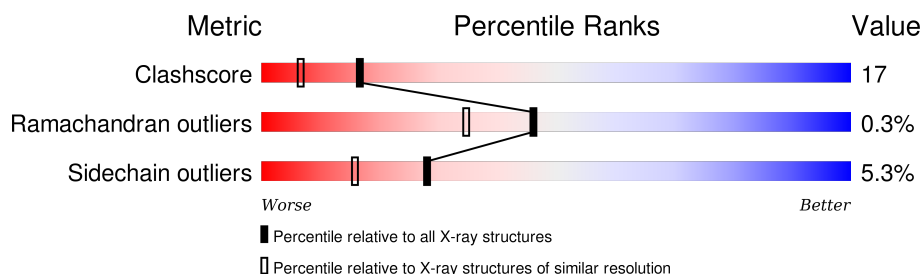
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.90 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	5398 (1.90-1.90)
Ramachandran outliers	100387	5338 (1.90-1.90)
Sidechain outliers	100360	5339 (1.90-1.90)

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	293	 78% 18% ...
1	C	293	 76% 20% ...
1	D	293	 73% 24% ..
2	B	293	 70% 25% ...

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	THM	B	1292	-	-	X	-
4	G1P	A	1294	-	-	X	-
4	G1P	D	1294	X	-	-	-
5	SO4	C	1295	-	-	X	-

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 9791 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 GLUCOSE-1-PHOSPHATE THYMIDYLYLTRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	290	Total	C	N	O	S	8	0	0
			2277	1459	378	429	11			
1	C	290	Total	C	N	O	S	9	0	0
			2273	1457	378	427	11			
1	D	290	Total	C	N	O	S	9	0	0
			2273	1457	378	427	11			

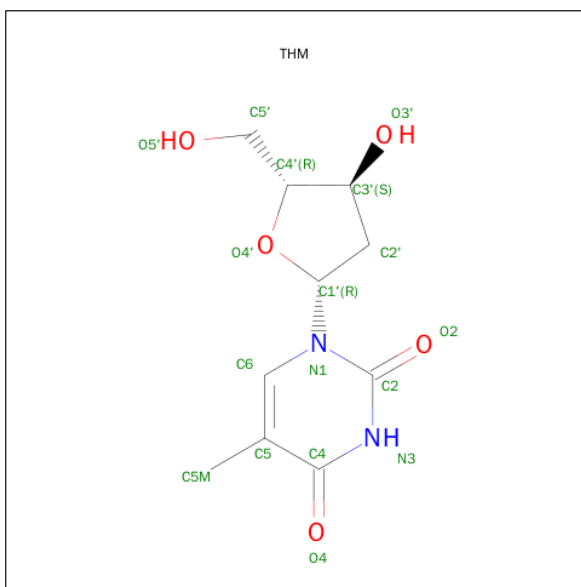
- Molecule 2 is a protein called GLUCOSE-1-PHOSPHATE THYMIDYLYLTRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	290	Total	C	N	O	S	9	0	0
			2270	1455	378	428	9			

There are 2 discrepancies between the modelled and reference sequences:

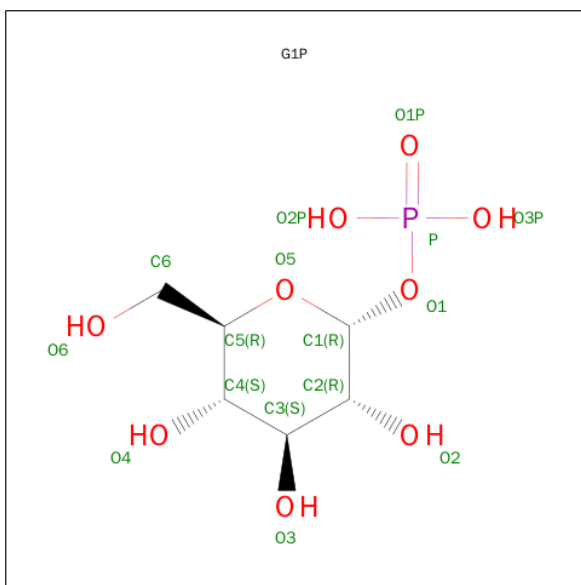
Chain	Residue	Modelled	Actual	Comment	Reference
B	217	LEU	MET	CONFLICT	UNP P37744
B	288	GLN	MET	CONFLICT	UNP P37744

- Molecule 3 is THYMIDINE (three-letter code: THM) (formula: C<sub>10</sub>H<sub>14</sub>N<sub>2</sub>O<sub>5</sub>).



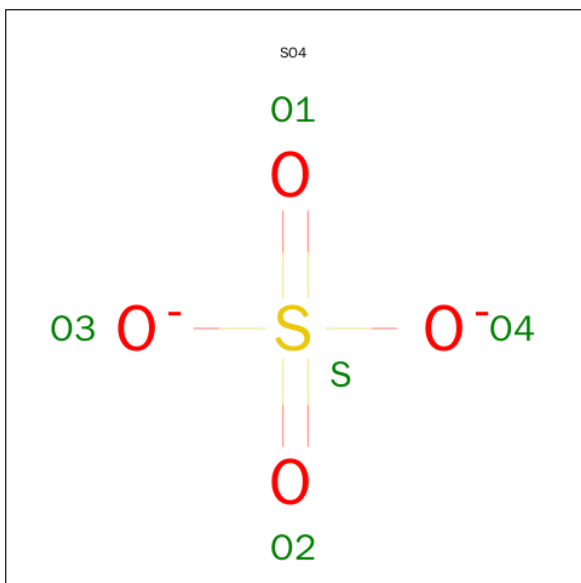
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			17	10	2	5		
3	A	1	Total	C	N	O	0	0
			17	10	2	5		
3	B	1	Total	C	N	O	0	0
			17	10	2	5		
3	B	1	Total	C	N	O	0	0
			17	10	2	5		
3	C	1	Total	C	N	O	0	0
			17	10	2	5		
3	C	1	Total	C	N	O	0	0
			17	10	2	5		
3	D	1	Total	C	N	O	0	0
			17	10	2	5		
3	D	1	Total	C	N	O	0	0
			17	10	2	5		

- Molecule 4 is SUGAR (ALPHA-D-GLUCOSE-1-PHOSPHATE) (three-letter code: G1P) (formula: C<sub>6</sub>H<sub>13</sub>O<sub>9</sub>P).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	O	P	0	0
			16	6	9	1		
4	B	1	Total	C	O	P	0	0
			16	6	9	1		
4	C	1	Total	C	O	P	0	0
			16	6	9	1		
4	D	1	Total	C	O	P	1	0
			16	6	9	1		

- Molecule 5 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total O S 5 4 1	0	0
5	A	1	Total O S 5 4 1	0	0
5	B	1	Total O S 5 4 1	0	0
5	C	1	Total O S 5 4 1	0	0
5	D	1	Total O S 5 4 1	0	0

- Molecule 6 is water.

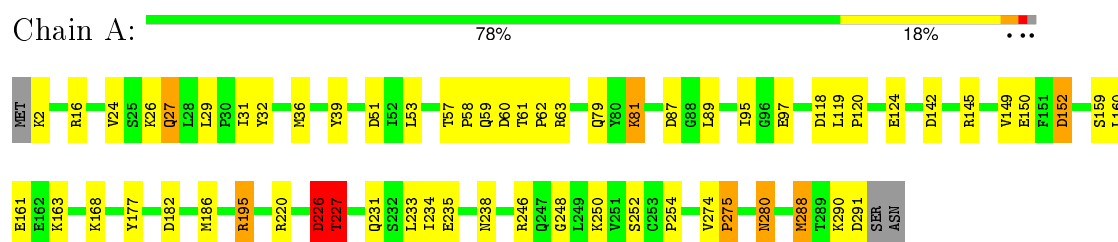
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	134	Total O 134 134	0	0
6	B	97	Total O 97 97	0	0
6	C	108	Total O 108 108	0	0
6	D	134	Total O 134 134	0	0

### 3 Residue-property plots

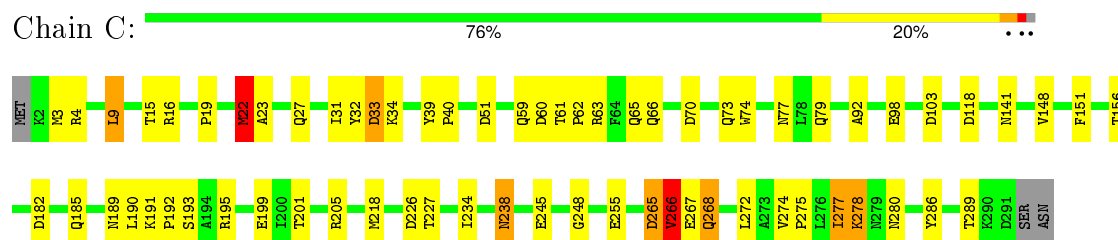
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according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

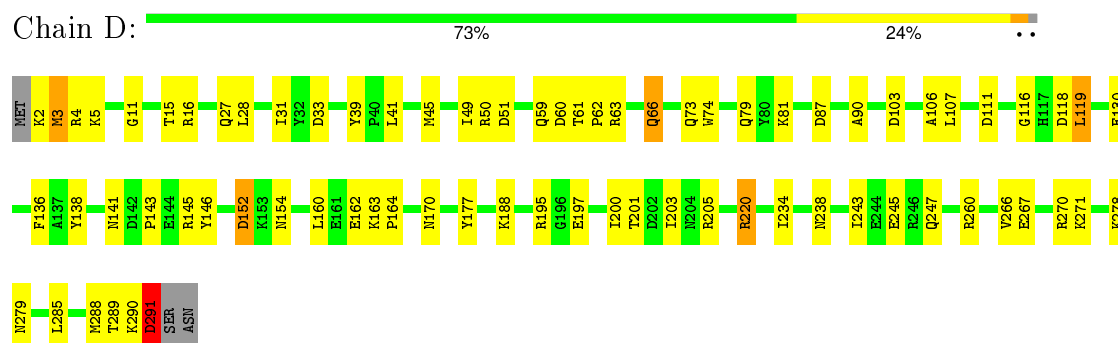
#### • Molecule 1: GLUCOSE-1-PHOSPHATE THYMIDYLTRANSFERASE



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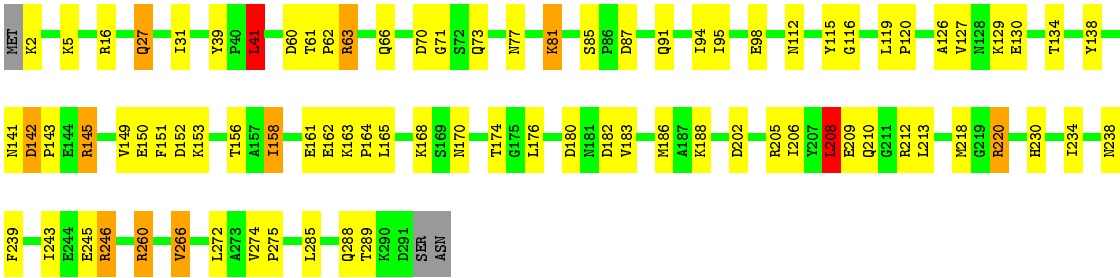
#### • Molecule 1: GLUCOSE-1-PHOSPHATE THYMIDYLTRANSFERASE



#### • Molecule 2: GLUCOSE-1-PHOSPHATE THYMIDYLTRANSFERASE







## 4 Data and refinement statistics

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

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	72.81 Å   118.96 Å   80.59 Å 90.00°   112.60°   90.00°	Depositor
Resolution (Å)	12.00 – 1.90	Depositor
% Data completeness (in resolution range)	85.4 (12.00-1.90)	Depositor
$R_{merge}$	0.03	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	REFMAC	Depositor
R, $R_{free}$	0.173 , 0.234	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	9791	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	28.0	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: THM, SO4, G1P

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	0.97	3/2324 (0.1%)	1.19	14/3145 (0.4%)
1	C	0.93	2/2320 (0.1%)	1.05	13/3139 (0.4%)
1	D	0.81	2/2320 (0.1%)	1.02	10/3140 (0.3%)
2	B	0.79	1/2317 (0.0%)	1.00	10/3138 (0.3%)
All	All	0.88	8/9281 (0.1%)	1.07	47/12562 (0.4%)

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	A	0	1
2	B	1	1
All	All	1	2

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	16	ARG	CG-CD	-21.04	0.99	1.51
1	A	152	ASP	CA-CB	20.32	1.98	1.53
1	A	16	ARG	CD-NE	-12.85	1.24	1.46
2	B	16	ARG	CG-CD	8.40	1.73	1.51
1	D	16	ARG	CG-CD	-8.18	1.31	1.51
1	C	218	MET	SD-CE	-6.59	1.41	1.77
1	D	138	TYR	CD1-CE1	5.30	1.47	1.39
1	A	177	TYR	CD2-CE2	5.15	1.47	1.39

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	16	ARG	CD-NE-CZ	-24.89	88.75	123.60
1	A	16	ARG	CG-CD-NE	16.33	146.09	111.80
1	A	152	ASP	CB-CA-C	-12.27	85.85	110.40
1	A	152	ASP	N-CA-CB	-11.60	89.72	110.60
2	B	152	ASP	N-CA-CB	10.67	129.81	110.60
1	C	16	ARG	CB-CG-CD	9.18	135.47	111.60
2	B	152	ASP	CB-CG-OD2	7.83	125.35	118.30
1	C	63	ARG	NE-CZ-NH1	7.18	123.89	120.30
2	B	60	ASP	CB-CG-OD2	7.14	124.73	118.30
1	C	103	ASP	CB-CG-OD2	7.12	124.71	118.30
1	C	22	MET	CG-SD-CE	-6.83	89.26	100.20
1	D	145	ARG	NE-CZ-NH2	-6.74	116.93	120.30
2	B	246	ARG	NE-CZ-NH1	6.50	123.55	120.30
1	D	145	ARG	NE-CZ-NH1	6.43	123.51	120.30
1	A	195	ARG	NE-CZ-NH2	-6.42	117.09	120.30
1	A	220	ARG	NE-CZ-NH2	-6.27	117.17	120.30
1	A	118	ASP	CB-CG-OD2	6.17	123.86	118.30
1	C	118	ASP	CB-CG-OD2	6.09	123.78	118.30
1	D	60	ASP	CB-CG-OD2	5.98	123.68	118.30
1	A	51	ASP	CB-CG-OD2	5.89	123.60	118.30
1	C	218	MET	CG-SD-CE	-5.88	90.79	100.20
1	D	33	ASP	CB-CG-OD2	5.79	123.51	118.30
2	B	213	LEU	CB-CG-CD1	-5.71	101.29	111.00
1	C	148	VAL	CG1-CB-CG2	5.71	120.03	110.90
1	C	265	ASP	CB-CG-OD2	5.70	123.43	118.30
1	A	87	ASP	CB-CG-OD2	5.67	123.41	118.30
1	D	291	ASP	CB-CG-OD2	5.66	123.39	118.30
1	A	220	ARG	NE-CZ-NH1	5.65	123.13	120.30
1	D	118	ASP	CB-CG-OD2	5.64	123.37	118.30
2	B	208	LEU	CB-CG-CD2	5.62	120.56	111.00
2	B	16	ARG	CB-CG-CD	5.61	126.17	111.60
1	A	227	THR	N-CA-CB	-5.60	99.67	110.30
1	C	70	ASP	CB-CG-OD1	5.59	123.33	118.30
1	A	60	ASP	CB-CG-OD2	5.57	123.32	118.30
2	B	246	ARG	NE-CZ-NH2	-5.53	117.53	120.30
1	D	103	ASP	CB-CG-OD2	5.47	123.22	118.30
2	B	142	ASP	CB-CG-OD2	5.43	123.19	118.30
1	D	111	ASP	CB-CG-OD2	5.40	123.16	118.30
1	A	226	ASP	N-CA-CB	5.35	120.22	110.60
1	C	33	ASP	CB-CG-OD2	5.34	123.11	118.30
1	A	152	ASP	CB-CG-OD2	5.32	123.08	118.30
2	B	41	LEU	CB-CG-CD1	5.24	119.91	111.00
1	C	148	VAL	CA-CB-CG1	5.23	118.74	110.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	87	ASP	CB-CG-OD2	5.21	122.99	118.30
1	C	60	ASP	CB-CG-OD2	5.08	122.87	118.30
1	C	182	ASP	CB-CG-OD2	5.05	122.84	118.30
1	D	152	ASP	N-CA-CB	5.01	119.61	110.60

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	B	174	THR	CB

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	195	ARG	Sidechain
2	B	208	LEU	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	2277	0	2284	68	0
1	C	2273	0	2281	68	0
1	D	2273	0	2280	92	0
2	B	2270	0	2270	87	0
3	A	34	0	28	0	0
3	B	34	0	28	9	0
3	C	34	0	28	5	0
3	D	34	0	28	5	0
4	A	16	0	11	13	0
4	B	16	0	11	1	0
4	C	16	0	11	0	0
4	D	16	0	10	2	0
5	A	10	0	0	0	0
5	B	5	0	0	0	0
5	C	5	0	0	4	0
5	D	5	0	0	0	0
6	A	134	0	0	16	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	B	97	0	0	12	0
6	C	108	0	0	5	0
6	D	134	0	0	25	0
All	All	9791	0	9270	306	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 17.

All (306) 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)
2:B:245:GLU:O	1:C:141:ASN:ND2	1.65	1.30
4:D:1294:G1P:P	4:D:1294:G1P:C1	2.18	1.30
2:B:142:ASP:OD2	2:B:145:ARG:NH1	1.61	1.30
2:B:161:GLU:CD	6:B:2055:HOH:O	1.73	1.25
2:B:161:GLU:OE1	6:B:2055:HOH:O	1.54	1.19
1:A:186:MET:HE3	6:A:2092:HOH:O	1.38	1.19
2:B:260:ARG:HH11	2:B:260:ARG:HG2	1.06	1.17
2:B:115:TYR:HA	3:B:1292:THM:O3'	1.45	1.16
1:A:226:ASP:OD2	6:A:2103:HOH:O	1.60	1.16
1:A:26:LYS:CE	1:A:227:THR:HG22	1.74	1.16
1:D:270:ARG:NH2	1:D:290:LYS:HG3	1.59	1.14
1:C:3:MET:CE	1:C:51:ASP:HB2	1.78	1.12
1:A:26:LYS:CE	1:A:227:THR:CG2	2.26	1.11
1:A:226:ASP:CG	6:A:2103:HOH:O	1.86	1.10
1:D:290:LYS:CD	6:D:2129:HOH:O	1.96	1.09
3:B:1292:THM:H6	3:B:1292:THM:H5'2	1.19	1.09
2:B:205:ARG:O	2:B:209:GLU:HG3	1.53	1.08
1:D:290:LYS:HG2	6:D:2129:HOH:O	1.53	1.07
1:C:191:LYS:HD3	1:C:192:PRO:CD	1.85	1.07
1:A:182:ASP:CB	6:A:2092:HOH:O	2.00	1.06
1:A:97:GLU:OE2	6:A:2038:HOH:O	1.71	1.05
1:D:290:LYS:CG	6:D:2129:HOH:O	2.02	1.03
1:C:3:MET:HE3	1:C:51:ASP:CB	1.88	1.03
1:A:26:LYS:HE3	1:A:227:THR:HG22	1.38	1.02
1:A:26:LYS:HE3	1:A:227:THR:CG2	1.88	1.02
1:D:119:LEU:HB2	3:D:1292:THM:H5'1	1.41	1.01
3:B:1292:THM:C6	3:B:1292:THM:H5'2	1.91	1.00
1:D:289:THR:O	1:D:290:LYS:HG2	1.61	1.00
1:A:26:LYS:HE2	1:A:227:THR:HG21	1.41	1.00
4:A:1294:G1P:H4	6:A:2130:HOH:O	1.60	0.99

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:195:ARG:HD3	1:C:199:GLU:OE2	1.63	0.97
1:C:3:MET:HE3	1:C:51:ASP:HB2	0.98	0.96
3:C:1292:THM:H5'2	5:C:1295:SO4:O2	1.62	0.96
1:C:191:LYS:HD3	1:C:192:PRO:HD3	1.46	0.96
2:B:142:ASP:CG	2:B:145:ARG:NH1	2.19	0.95
2:B:145:ARG:HH11	2:B:145:ARG:HG3	1.33	0.94
2:B:260:ARG:CG	2:B:260:ARG:HH11	1.81	0.94
2:B:63:ARG:CG	2:B:63:ARG:HH11	1.81	0.93
2:B:73:GLN:N	2:B:73:GLN:OE1	2.00	0.93
1:A:26:LYS:HE2	1:A:227:THR:CG2	1.96	0.93
2:B:115:TYR:HA	3:B:1292:THM:HO3'	1.26	0.92
1:D:220:ARG:HG3	6:D:2099:HOH:O	1.67	0.91
2:B:63:ARG:HG3	2:B:63:ARG:HH11	1.34	0.91
1:A:182:ASP:HB2	6:A:2092:HOH:O	1.63	0.90
1:A:227:THR:HG23	1:A:233:LEU:HD21	1.52	0.90
1:D:188:LYS:O	6:D:2082:HOH:O	1.89	0.90
3:B:1292:THM:C5'	3:B:1292:THM:H6	2.02	0.89
1:C:191:LYS:HD3	1:C:192:PRO:HD2	1.52	0.89
1:D:2:LYS:O	1:D:51:ASP:HB2	1.73	0.89
2:B:285:LEU:O	2:B:288:GLN:HG2	1.72	0.88
1:A:227:THR:O	1:A:227:THR:HG22	1.71	0.87
1:A:227:THR:O	1:A:227:THR:CG2	2.20	0.87
1:D:245:GLU:HG3	6:D:2021:HOH:O	1.73	0.87
1:D:266:VAL:HG12	1:D:289:THR:HG23	1.57	0.86
4:A:1294:G1P:P	4:A:1294:G1P:O2	2.32	0.86
1:A:226:ASP:CB	6:A:2103:HOH:O	2.23	0.85
1:D:270:ARG:HH21	1:D:290:LYS:HG3	1.38	0.85
2:B:71:GLY:HA2	2:B:73:GLN:HE22	1.41	0.83
1:D:160:LEU:O	1:D:205:ARG:NH1	2.13	0.82
1:A:182:ASP:HB3	6:A:2092:HOH:O	1.66	0.82
1:D:290:LYS:HD3	6:D:2129:HOH:O	1.68	0.80
1:D:289:THR:O	1:D:290:LYS:CG	2.30	0.79
1:D:290:LYS:CE	6:D:2122:HOH:O	2.30	0.79
1:A:161:GLU:OE1	1:A:168:LYS:HE2	1.82	0.78
1:C:185:GLN:HE21	1:C:189:ASN:HD21	1.30	0.77
2:B:77:ASN:CG	6:B:2025:HOH:O	2.21	0.77
1:C:3:MET:CE	1:C:51:ASP:CB	2.53	0.76
2:B:220:ARG:HG3	6:B:2073:HOH:O	1.85	0.76
2:B:77:ASN:ND2	6:B:2025:HOH:O	2.19	0.75
2:B:145:ARG:HG3	2:B:145:ARG:NH1	2.01	0.74
3:C:1292:THM:H5'2	5:C:1295:SO4:S	2.27	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:270:ARG:HH22	1:D:290:LYS:HG3	1.51	0.73
1:A:89:LEU:HD11	4:A:1294:G1P:H1	1.68	0.73
1:D:290:LYS:HE2	6:D:2122:HOH:O	1.86	0.73
2:B:260:ARG:HG2	2:B:260:ARG:NH1	1.89	0.73
2:B:81:LYS:HD3	6:B:2030:HOH:O	1.88	0.73
1:D:119:LEU:CB	3:D:1292:THM:H5'1	2.19	0.73
1:A:124:GLU:HG3	6:A:2023:HOH:O	1.89	0.73
1:A:26:LYS:NZ	1:A:227:THR:HG22	2.05	0.72
2:B:27:GLN:H	2:B:27:GLN:HE21	1.36	0.72
1:D:238:ASN:HB3	6:D:2108:HOH:O	1.89	0.72
1:C:195:ARG:CD	1:C:199:GLU:OE2	2.38	0.71
1:C:15:THR:HG21	1:D:278:LYS:HE2	1.70	0.71
2:B:73:GLN:H	2:B:73:GLN:CD	1.93	0.70
1:D:4:ARG:HD3	1:D:5:LYS:H	1.56	0.70
1:D:289:THR:HG22	1:D:289:THR:O	1.92	0.69
1:C:3:MET:HE1	1:C:77:ASN:HB3	1.75	0.69
2:B:27:GLN:H	2:B:27:GLN:NE2	1.90	0.69
1:D:260:ARG:HG2	1:D:260:ARG:HH11	1.58	0.69
1:A:26:LYS:CE	1:A:227:THR:HG21	2.05	0.68
2:B:180:ASP:OD1	2:B:182:ASP:HB2	1.93	0.68
2:B:127:VAL:O	6:B:2044:HOH:O	2.12	0.68
2:B:71:GLY:HA2	2:B:73:GLN:NE2	2.08	0.68
1:D:2:LYS:HA	1:D:50:ARG:HB2	1.75	0.67
2:B:174:THR:HG22	6:B:2061:HOH:O	1.95	0.67
2:B:94:ILE:HD12	2:B:188:LYS:HG3	1.77	0.67
1:D:152:ASP:O	1:D:154:ASN:O	2.14	0.66
1:A:290:LYS:O	1:A:291:ASP:CB	2.44	0.66
1:D:290:LYS:HE3	6:D:2122:HOH:O	1.92	0.65
1:C:192:PRO:HD2	6:C:2067:HOH:O	1.96	0.65
1:C:59:GLN:OE1	1:D:63:ARG:NH1	2.28	0.65
1:A:163:LYS:HE3	4:A:1294:G1P:O1P	1.97	0.65
1:D:90:ALA:HB1	1:D:203:ILE:HD11	1.78	0.65
1:A:163:LYS:CE	4:A:1294:G1P:O1P	2.45	0.65
1:D:116:GLY:O	3:D:1292:THM:H4'	1.97	0.64
2:B:126:ALA:HB2	2:B:134:THR:HG21	1.80	0.64
1:A:234:ILE:HG21	2:B:238:ASN:OD1	1.98	0.64
1:C:238:ASN:ND2	1:D:234:ILE:HG21	2.13	0.63
2:B:274:VAL:HB	2:B:275:PRO:HD3	1.80	0.63
2:B:94:ILE:HD12	2:B:188:LYS:CG	2.28	0.63
1:C:98:GLU:HG2	6:C:2031:HOH:O	1.97	0.63
1:A:238:ASN:OD1	2:B:234:ILE:HG21	1.98	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:162:GLU:OE2	4:B:1294:G1P:O2	2.17	0.62
2:B:174:THR:HG23	2:B:176:LEU:H	1.64	0.62
2:B:116:GLY:H	3:B:1292:THM:H4'	1.63	0.62
2:B:119:LEU:HB3	3:B:1292:THM:H5'1	1.79	0.62
3:C:1292:THM:C5'	5:C:1295:SO4:O2	2.45	0.62
1:D:90:ALA:CB	1:D:203:ILE:HD11	2.29	0.61
2:B:260:ARG:CG	2:B:260:ARG:NH1	2.48	0.61
1:A:274:VAL:HB	1:A:275:PRO:HD3	1.82	0.61
1:A:36:MET:HE1	1:A:227:THR:HG21	1.83	0.61
1:D:270:ARG:NH2	1:D:290:LYS:CG	2.50	0.61
1:C:9:LEU:CD2	1:C:92:ALA:HB2	2.30	0.61
1:D:4:ARG:HD2	1:D:49:ILE:HG12	1.82	0.61
1:D:238:ASN:CB	6:D:2108:HOH:O	2.47	0.61
2:B:2:LYS:HD2	6:B:2039:HOH:O	2.01	0.61
1:C:185:GLN:NE2	1:C:189:ASN:HD21	1.98	0.60
3:C:1292:THM:H5'1	3:C:1292:THM:H6	1.82	0.60
1:C:278:LYS:HE2	1:D:15:THR:CG2	2.32	0.60
2:B:138:TYR:HD1	2:B:218:MET:HG2	1.66	0.60
1:D:266:VAL:CG1	1:D:289:THR:HG23	2.30	0.59
1:A:89:LEU:CD1	4:A:1294:G1P:H2	2.31	0.59
1:A:290:LYS:O	1:A:291:ASP:HB2	2.02	0.59
1:C:15:THR:CG2	1:D:278:LYS:HE2	2.32	0.59
1:D:119:LEU:HB2	3:D:1292:THM:C5'	2.27	0.59
1:C:227:THR:N	6:C:2088:HOH:O	1.96	0.59
1:D:271:LYS:NZ	6:D:2124:HOH:O	2.34	0.59
1:C:185:GLN:HE21	1:C:189:ASN:ND2	1.99	0.59
1:D:200:ILE:HA	1:D:203:ILE:HD12	1.85	0.58
1:C:265:ASP:O	1:C:265:ASP:OD1	2.21	0.58
1:D:267:GLU:HG2	6:D:2054:HOH:O	2.02	0.58
1:D:270:ARG:HH22	1:D:290:LYS:HE3	1.67	0.58
1:A:89:LEU:HD11	4:A:1294:G1P:H2	1.85	0.58
1:A:159:SER:OG	6:A:2066:HOH:O	2.09	0.58
1:A:238:ASN:OD1	2:B:234:ILE:CG2	2.53	0.57
1:C:266:VAL:CG1	1:C:289:THR:HB	2.33	0.57
1:D:260:ARG:HH11	1:D:260:ARG:CG	2.18	0.56
1:C:266:VAL:HG23	6:C:2102:HOH:O	2.04	0.56
1:A:27:GLN:H	1:A:27:GLN:HE21	1.52	0.56
1:A:231:GLN:O	1:A:235:GLU:HG3	2.04	0.56
3:C:1292:THM:H5'2	5:C:1295:SO4:O1	2.04	0.56
1:C:226:ASP:HA	6:C:2088:HOH:O	2.05	0.56
1:A:227:THR:CG2	1:A:233:LEU:HD21	2.31	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:107:LEU:HD23	1:D:200:ILE:HD12	1.88	0.56
1:A:27:GLN:H	1:A:27:GLN:NE2	2.03	0.56
2:B:150:GLU:CD	2:B:168:LYS:HD3	2.25	0.56
1:A:161:GLU:HG3	6:A:2066:HOH:O	2.05	0.55
1:C:278:LYS:HE2	1:D:15:THR:HG23	1.88	0.55
2:B:183:VAL:HA	2:B:186:MET:HE3	1.87	0.55
2:B:150:GLU:OE1	2:B:168:LYS:HD3	2.07	0.55
1:D:143:PRO:HG3	1:D:170:ASN:HA	1.88	0.55
1:A:142:ASP:OD2	1:A:145:ARG:HD3	2.06	0.55
1:A:250:LYS:NZ	1:A:280:ASN:HD21	2.04	0.55
1:C:61:THR:HB	1:C:62:PRO:HD3	1.88	0.55
2:B:73:GLN:N	2:B:73:GLN:CD	2.57	0.54
1:A:26:LYS:HZ1	1:A:227:THR:HG22	1.72	0.54
1:C:193:SER:HB3	1:C:195:ARG:HG2	1.88	0.54
1:C:234:ILE:HG21	1:D:238:ASN:OD1	2.07	0.54
1:C:39:TYR:N	1:C:40:PRO:HD2	2.23	0.54
1:C:9:LEU:HD22	1:C:92:ALA:HB2	1.90	0.53
1:A:61:THR:N	1:A:62:PRO:CD	2.71	0.53
1:C:277:ILE:HD11	1:C:286:TYR:HB2	1.90	0.53
1:A:2:LYS:HB2	6:A:2001:HOH:O	2.08	0.53
4:A:1294:G1P:O1P	4:A:1294:G1P:O2	2.25	0.53
1:D:270:ARG:HH22	1:D:290:LYS:CG	2.20	0.53
2:B:116:GLY:O	3:B:1292:THM:H4'	2.09	0.53
1:D:3:MET:HE3	6:D:2028:HOH:O	2.08	0.53
4:A:1294:G1P:P	4:A:1294:G1P:HO2	2.31	0.52
2:B:77:ASN:OD1	6:B:2025:HOH:O	2.18	0.52
1:D:73:GLN:HG2	1:D:74:TRP:CD1	2.45	0.52
1:D:195:ARG:NH2	1:D:197:GLU:OE1	2.43	0.52
1:C:9:LEU:HD21	1:C:92:ALA:HB2	1.91	0.51
4:A:1294:G1P:O2	4:A:1294:G1P:O2P	2.27	0.51
2:B:138:TYR:CD1	2:B:218:MET:HG2	2.45	0.51
1:D:289:THR:O	1:D:289:THR:CG2	2.58	0.51
1:A:89:LEU:HD11	4:A:1294:G1P:C1	2.36	0.51
1:A:119:LEU:N	1:A:120:PRO:CD	2.72	0.51
1:A:234:ILE:CG2	2:B:238:ASN:OD1	2.59	0.51
1:A:81:LYS:HD2	1:A:95:ILE:HG23	1.93	0.50
1:C:277:ILE:CD1	1:C:286:TYR:HB2	2.42	0.50
1:A:288:MET:O	1:A:288:MET:HG3	2.09	0.50
2:B:143:PRO:HG3	2:B:170:ASN:HD22	1.76	0.50
1:D:245:GLU:CG	6:D:2021:HOH:O	2.46	0.50
2:B:202:ASP:OD1	2:B:205:ARG:NH2	2.44	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:238:ASN:ND2	1:D:234:ILE:CG2	2.75	0.50
1:D:146:TYR:HD2	4:D:1294:G1P:H5	1.77	0.50
1:D:290:LYS:HB3	1:D:290:LYS:NZ	2.26	0.50
2:B:71:GLY:CA	2:B:73:GLN:HE22	2.19	0.49
1:C:265:ASP:OD1	1:C:265:ASP:C	2.51	0.49
1:D:163:LYS:N	1:D:164:PRO:CD	2.75	0.49
2:B:141:ASN:HB3	1:C:248:GLY:HA3	1.94	0.49
1:D:266:VAL:CG2	6:D:2120:HOH:O	2.61	0.49
1:D:285:LEU:O	1:D:288:MET:HB2	2.12	0.49
1:D:4:ARG:NH2	6:D:2003:HOH:O	2.45	0.49
1:A:31:ILE:HB	1:A:39:TYR:CE1	2.48	0.49
1:C:274:VAL:N	1:C:275:PRO:CD	2.75	0.49
2:B:119:LEU:HB3	2:B:120:PRO:HD3	1.95	0.49
1:D:266:VAL:HG23	6:D:2120:HOH:O	2.12	0.49
2:B:70:ASP:OD1	2:B:70:ASP:C	2.51	0.49
1:D:107:LEU:HD23	1:D:200:ILE:CD1	2.42	0.48
1:A:226:ASP:CA	6:A:2103:HOH:O	2.60	0.48
1:D:267:GLU:CG	6:D:2054:HOH:O	2.58	0.48
2:B:94:ILE:HD11	2:B:188:LYS:N	2.29	0.48
1:C:22:MET:SD	6:D:2025:HOH:O	2.60	0.48
1:C:31:ILE:HB	1:C:39:TYR:CE1	2.48	0.48
1:C:23:ALA:HB1	1:D:28:LEU:HD22	1.95	0.48
2:B:61:THR:N	2:B:62:PRO:CD	2.77	0.48
1:C:277:ILE:HD11	1:C:286:TYR:CB	2.44	0.47
2:B:63:ARG:HG3	2:B:63:ARG:NH1	2.12	0.47
1:D:260:ARG:CG	1:D:260:ARG:NH1	2.77	0.47
1:A:53:LEU:HD11	1:A:81:LYS:HB2	1.96	0.47
2:B:158:ILE:HG23	2:B:158:ILE:O	2.14	0.47
2:B:94:ILE:CD1	2:B:188:LYS:HG3	2.42	0.47
1:C:33:ASP:C	1:C:34:LYS:HG3	2.34	0.47
1:D:177:TYR:CZ	1:D:200:ILE:HD13	2.49	0.47
2:B:85:SER:OG	2:B:87:ASP:OD2	2.22	0.47
1:C:268:GLN:O	1:C:272:LEU:HG	2.15	0.47
1:A:142:ASP:N	1:A:142:ASP:OD1	2.48	0.47
2:B:151:PHE:HA	2:B:156:THR:O	2.15	0.46
1:D:290:LYS:HA	1:D:291:ASP:HA	1.65	0.46
1:C:19:PRO:HA	1:C:22:MET:HG3	1.96	0.46
6:A:2016:HOH:O	2:B:230:HIS:CE1	2.68	0.46
1:D:61:THR:HB	1:D:62:PRO:HD3	1.96	0.46
1:C:278:LYS:HE2	1:D:15:THR:HG21	1.98	0.46
1:C:265:ASP:O	1:C:266:VAL:C	2.54	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:205:ARG:HD3	6:D:2090:HOH:O	2.14	0.46
1:A:142:ASP:OD2	1:A:145:ARG:CD	2.63	0.46
2:B:205:ARG:O	2:B:209:GLU:CG	2.45	0.46
1:A:150:GLU:CD	1:A:168:LYS:HD3	2.36	0.46
1:C:151:PHE:HA	1:C:156:THR:O	2.15	0.46
1:A:288:MET:HB3	1:A:288:MET:HE3	1.57	0.45
2:B:266:VAL:HG22	2:B:289:THR:HB	1.99	0.45
1:D:41:LEU:HD13	1:D:41:LEU:C	2.37	0.45
1:A:288:MET:O	1:A:288:MET:CG	2.60	0.45
2:B:158:ILE:CG2	2:B:158:ILE:O	2.63	0.45
1:A:63:ARG:HH11	1:A:63:ARG:HG3	1.80	0.45
1:C:265:ASP:O	1:C:268:GLN:N	2.47	0.45
2:B:141:ASN:HB3	1:C:248:GLY:CA	2.47	0.45
1:A:227:THR:HG23	1:A:233:LEU:CD2	2.35	0.45
1:C:59:GLN:HE22	1:D:66:GLN:NE2	2.15	0.45
1:C:34:LYS:HE2	1:C:34:LYS:HB2	1.57	0.45
2:B:63:ARG:HG2	2:B:63:ARG:HH11	1.75	0.44
1:A:248:GLY:HA3	1:D:141:ASN:HB3	1.99	0.44
1:D:267:GLU:HG3	6:D:2120:HOH:O	2.17	0.44
2:B:206:ILE:O	2:B:210:GLN:HG3	2.17	0.44
1:D:31:ILE:HB	1:D:39:TYR:CE1	2.52	0.44
2:B:119:LEU:CD2	6:B:2016:HOH:O	2.66	0.44
2:B:182:ASP:OD2	2:B:212:ARG:NH2	2.47	0.44
1:C:19:PRO:CA	1:C:22:MET:HG3	2.48	0.44
1:A:24:VAL:HG11	1:A:29:LEU:HD23	1.99	0.44
2:B:119:LEU:CB	3:B:1292:THM:H5'1	2.48	0.44
1:D:4:ARG:HD3	1:D:5:LYS:N	2.28	0.44
1:D:41:LEU:CD1	1:D:45:MET:SD	3.06	0.44
2:B:31:ILE:HB	2:B:39:TYR:CE1	2.53	0.44
1:A:89:LEU:HD12	4:A:1294:G1P:H2	1.99	0.43
2:B:239:PHE:CE2	2:B:243:ILE:HD11	2.54	0.43
1:D:270:ARG:NE	1:D:289:THR:HG21	2.33	0.43
1:C:39:TYR:N	1:C:40:PRO:CD	2.81	0.43
1:C:22:MET:HB3	1:C:22:MET:HE2	1.77	0.43
1:D:11:GLY:O	3:D:1293:THM:H2'1	2.19	0.43
1:A:119:LEU:HB3	1:A:120:PRO:HD3	1.99	0.43
1:D:41:LEU:HD13	1:D:45:MET:SD	2.59	0.43
1:C:74:TRP:NE1	1:C:255:GLU:HG3	2.33	0.43
1:D:238:ASN:CG	6:D:2108:HOH:O	2.56	0.43
2:B:63:ARG:CG	2:B:63:ARG:NH1	2.53	0.43
2:B:27:GLN:N	2:B:27:GLN:NE2	2.61	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:136:PHE:N	6:D:2069:HOH:O	2.16	0.43
1:C:62:PRO:HA	1:C:65:GLN:OE1	2.19	0.42
2:B:141:ASN:ND2	1:C:245:GLU:O	2.52	0.42
2:B:91:GLN:HG2	2:B:95:ILE:CD1	2.49	0.42
2:B:5:LYS:NZ	6:B:2003:HOH:O	2.41	0.42
2:B:41:LEU:HD22	2:B:41:LEU:O	2.19	0.42
1:D:3:MET:CE	1:D:3:MET:H	2.33	0.42
1:C:265:ASP:O	1:C:267:GLU:N	2.53	0.41
1:D:243:ILE:O	1:D:247:GLN:HG3	2.19	0.41
1:C:191:LYS:CD	1:C:192:PRO:HD2	2.37	0.41
1:D:4:ARG:HH12	1:D:106:ALA:N	2.19	0.41
1:C:266:VAL:HG12	1:C:289:THR:HB	2.00	0.41
1:C:201:THR:O	1:C:205:ARG:HG3	2.20	0.41
2:B:73:GLN:HB2	2:B:272:LEU:HD21	2.02	0.41
1:D:61:THR:N	1:D:62:PRO:CD	2.83	0.41
1:A:59:GLN:HG3	6:A:2026:HOH:O	2.20	0.41
2:B:163:LYS:N	2:B:164:PRO:CD	2.83	0.41
1:C:73:GLN:HG2	1:C:74:TRP:CD1	2.54	0.41
1:A:57:THR:HB	1:A:58:PRO:HD2	2.03	0.41
2:B:163:LYS:N	2:B:164:PRO:HD3	2.35	0.41
1:D:279:ASN:OD1	1:D:279:ASN:C	2.58	0.41
2:B:94:ILE:HD12	2:B:188:LYS:HG2	2.01	0.41
1:C:19:PRO:HA	1:C:22:MET:CG	2.51	0.41
1:A:89:LEU:HD11	4:A:1294:G1P:C2	2.51	0.41
1:C:238:ASN:HD22	1:D:234:ILE:CG2	2.33	0.41
2:B:63:ARG:HD3	2:B:63:ARG:HA	1.80	0.40
1:A:248:GLY:CA	1:D:141:ASN:HB3	2.51	0.40
1:C:66:GLN:HE22	1:D:59:GLN:HE22	1.68	0.40
1:A:252:SER:O	1:A:254:PRO:HD3	2.21	0.40
1:D:162:GLU:OE1	1:D:201:THR:OG1	2.40	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

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	288/293 (98%)	286 (99%)	1 (0%)	1 (0%)	46	35
1	C	288/293 (98%)	282 (98%)	4 (1%)	2 (1%)	26	14
1	D	288/293 (98%)	284 (99%)	4 (1%)	0	100	100
2	B	288/293 (98%)	280 (97%)	8 (3%)	0	100	100
All	All	1152/1172 (98%)	1132 (98%)	17 (2%)	3 (0%)	46	35

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	32	TYR
1	C	32	TYR
1	C	266	VAL

### 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	244/247 (99%)	232 (95%)	12 (5%)	31	18
1	C	243/247 (98%)	231 (95%)	12 (5%)	31	18
1	D	243/247 (98%)	234 (96%)	9 (4%)	41	29
2	B	242/247 (98%)	223 (92%)	19 (8%)	15	6
All	All	972/988 (98%)	920 (95%)	52 (5%)	28	16

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	GLN
1	A	79	GLN
1	A	81	LYS
1	A	149	VAL
1	A	152	ASP

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Mol	Chain	Res	Type
1	A	160	LEU
1	A	226	ASP
1	A	227	THR
1	A	246	ARG
1	A	275	PRO
1	A	280	ASN
1	A	288	MET
2	B	27	GLN
2	B	41	LEU
2	B	63	ARG
2	B	66	GLN
2	B	81	LYS
2	B	98	GLU
2	B	112	ASN
2	B	129	LYS
2	B	130	GLU
2	B	145	ARG
2	B	149	VAL
2	B	153	LYS
2	B	158	ILE
2	B	165	LEU
2	B	208	LEU
2	B	220	ARG
2	B	246	ARG
2	B	260	ARG
2	B	266	VAL
1	C	4	ARG
1	C	9	LEU
1	C	22	MET
1	C	27	GLN
1	C	79	GLN
1	C	190	LEU
1	C	238	ASN
1	C	266	VAL
1	C	268	GLN
1	C	277	ILE
1	C	278	LYS
1	C	280	ASN
1	D	3	MET
1	D	27	GLN
1	D	66	GLN
1	D	79	GLN

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Mol	Chain	Res	Type
1	D	81	LYS
1	D	119	LEU
1	D	130	GLU
1	D	220	ARG
1	D	291	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	27	GLN
1	A	59	GLN
1	A	66	GLN
1	A	128	ASN
1	A	280	ASN
2	B	27	GLN
2	B	170	ASN
2	B	230	HIS
1	C	79	GLN
1	C	185	GLN
1	C	230	HIS
1	C	238	ASN
1	C	268	GLN
1	D	27	GLN
1	D	59	GLN
1	D	66	GLN
1	D	77	ASN
1	D	230	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

17 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
3	THM	A	1292	-	13,18,18	1.05	1 (7%)	16,26,26	2.44	4 (25%)
3	THM	A	1293	-	13,18,18	0.89	1 (7%)	16,26,26	1.72	5 (31%)
4	G1P	A	1294	-	15,16,16	1.22	2 (13%)	21,24,24	3.99	9 (42%)
5	SO4	A	1295	-	4,4,4	1.00	0	6,6,6	2.35	1 (16%)
5	SO4	A	1296	-	4,4,4	0.32	0	6,6,6	0.70	0
3	THM	B	1292	-	13,18,18	1.45	1 (7%)	16,26,26	2.35	5 (31%)
3	THM	B	1293	-	13,18,18	1.13	1 (7%)	16,26,26	2.02	5 (31%)
4	G1P	B	1294	-	15,16,16	0.82	1 (6%)	21,24,24	1.22	1 (4%)
5	SO4	B	1295	-	4,4,4	0.72	0	6,6,6	0.18	0
3	THM	C	1292	-	13,18,18	1.23	1 (7%)	16,26,26	2.73	3 (18%)
3	THM	C	1293	-	13,18,18	1.27	1 (7%)	16,26,26	2.16	5 (31%)
4	G1P	C	1294	-	15,16,16	0.77	1 (6%)	21,24,24	1.53	3 (14%)
5	SO4	C	1295	-	4,4,4	0.65	0	6,6,6	0.48	0
3	THM	D	1292	-	13,18,18	1.22	1 (7%)	16,26,26	2.74	5 (31%)
3	THM	D	1293	-	13,18,18	1.33	2 (15%)	16,26,26	2.72	6 (37%)
4	G1P	D	1294	-	15,16,16	5.08	2 (13%)	21,24,24	5.61	6 (28%)
5	SO4	D	1295	-	4,4,4	0.76	0	6,6,6	0.47	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	THM	A	1292	-	-	0/2/18/18	0/2/2/2
3	THM	A	1293	-	-	0/2/18/18	0/2/2/2
4	G1P	A	1294	-	-	0/7/27/27	0/1/1/1
5	SO4	A	1295	-	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	SO4	A	1296	-	-	0/0/0/0	0/0/0/0
3	THM	B	1292	-	-	0/2/18/18	0/2/2/2
3	THM	B	1293	-	-	0/2/18/18	0/2/2/2
4	G1P	B	1294	-	-	0/7/27/27	0/1/1/1
5	SO4	B	1295	-	-	0/0/0/0	0/0/0/0
3	THM	C	1292	-	-	0/2/18/18	0/2/2/2
3	THM	C	1293	-	-	0/2/18/18	0/2/2/2
4	G1P	C	1294	-	-	0/7/27/27	0/1/1/1
5	SO4	C	1295	-	-	0/0/0/0	0/0/0/0
3	THM	D	1292	-	-	0/2/18/18	0/2/2/2
3	THM	D	1293	-	-	0/2/18/18	0/2/2/2
4	G1P	D	1294	-	1/1/6/6	0/7/27/27	0/1/1/1
5	SO4	D	1295	-	-	0/0/0/0	0/0/0/0

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	1294	G1P	P-O1	-19.40	1.01	1.60
4	B	1294	G1P	P-O1	-2.64	1.52	1.60
4	C	1294	G1P	P-O1	-2.17	1.53	1.60
3	A	1293	THM	O4-C4	2.02	1.29	1.24
4	D	1294	G1P	O4-C4	2.09	1.48	1.43
3	D	1293	THM	O3'-C3'	2.23	1.48	1.43
4	A	1294	G1P	C4-C3	2.44	1.58	1.52
4	A	1294	G1P	O4-C4	2.78	1.49	1.43
3	A	1292	THM	O4-C4	3.22	1.32	1.24
3	D	1293	THM	O4-C4	3.47	1.33	1.24
3	C	1292	THM	O4-C4	3.68	1.33	1.24
3	B	1293	THM	O4-C4	3.84	1.33	1.24
3	C	1293	THM	O4-C4	3.87	1.34	1.24
3	D	1292	THM	O4-C4	3.99	1.34	1.24
3	B	1292	THM	O4-C4	4.26	1.34	1.24

All (58) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1294	G1P	O3-C3-C2	-9.48	88.99	110.34
3	D	1292	THM	C5-C4-N3	-5.99	118.46	125.14
3	C	1292	THM	C5-C4-N3	-5.87	118.61	125.14
3	D	1293	THM	C5-C4-N3	-5.55	118.96	125.14
3	B	1292	THM	C5-C4-N3	-5.36	119.17	125.14
3	A	1292	THM	C5-C4-N3	-4.99	119.58	125.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1292	THM	C5'-C4'-C3'	-4.68	102.34	114.80
3	C	1293	THM	C5-C4-N3	-4.56	120.06	125.14
3	B	1293	THM	C5-C4-N3	-4.50	120.13	125.14
3	D	1293	THM	O4'-C4'-C5'	-3.31	102.00	109.17
4	D	1294	G1P	O5-C1-C2	-3.15	103.80	110.28
4	A	1294	G1P	O6-C6-C5	-3.14	100.97	111.33
3	D	1293	THM	C5'-C4'-C3'	-3.07	106.61	114.80
3	D	1293	THM	C2'-C3'-C4'	-3.02	96.51	102.77
3	C	1293	THM	O4'-C1'-N1	-2.92	102.67	107.72
3	D	1292	THM	C5'-C4'-C3'	-2.78	107.39	114.80
3	A	1293	THM	C5-C4-N3	-2.49	122.36	125.14
3	B	1292	THM	O4'-C4'-C3'	-2.37	99.70	105.67
3	A	1293	THM	O3'-C3'-C2'	-2.24	103.33	110.74
3	B	1293	THM	O4'-C1'-C2'	-2.14	102.01	106.27
3	C	1293	THM	O5'-C5'-C4'	-2.02	104.65	111.33
3	A	1293	THM	C4'-O4'-C1'	2.00	114.52	109.47
3	B	1293	THM	C4'-O4'-C1'	2.04	114.62	109.47
3	C	1293	THM	C4'-O4'-C1'	2.20	115.02	109.47
4	D	1294	G1P	O1-C1-C2	2.26	112.62	108.39
3	B	1293	THM	O4'-C4'-C5'	2.41	114.39	109.17
3	B	1292	THM	O4'-C1'-N1	2.41	111.89	107.72
4	D	1294	G1P	C6-C5-C4	2.45	119.06	113.02
3	A	1293	THM	C2'-C3'-C4'	2.68	108.32	102.77
3	A	1292	THM	C4'-O4'-C1'	2.72	116.33	109.47
3	A	1293	THM	C4-N3-C2	2.72	117.60	115.25
4	C	1294	G1P	O2-C2-C1	2.73	116.00	110.02
4	A	1294	G1P	C4-C3-C2	2.87	116.15	110.79
4	C	1294	G1P	O4-C4-C5	3.37	118.18	109.24
4	D	1294	G1P	C1-O5-C5	3.42	120.38	113.75
3	B	1292	THM	O5'-C5'-C4'	3.80	123.88	111.33
3	D	1293	THM	C2'-C1'-N1	3.89	123.61	114.16
3	D	1292	THM	C2'-C1'-N1	4.05	124.01	114.16
4	B	1294	G1P	C1-O5-C5	4.16	121.83	113.75
4	C	1294	G1P	C1-O5-C5	4.21	121.91	113.75
3	D	1292	THM	O4'-C1'-N1	4.34	115.22	107.72
3	C	1292	THM	O4'-C1'-N1	4.52	115.54	107.72
3	C	1293	THM	C4-N3-C2	4.58	119.21	115.25
3	B	1293	THM	C4-N3-C2	4.70	119.31	115.25
3	B	1292	THM	C4-N3-C2	4.74	119.34	115.25
4	D	1294	G1P	O4-C4-C5	5.12	122.80	109.24
5	A	1295	SO4	O2-S-O1	5.19	125.95	109.50
3	A	1292	THM	C4-N3-C2	5.22	119.76	115.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	1292	THM	C4-N3-C2	5.50	120.00	115.25
4	A	1294	G1P	C1-C2-C3	5.57	120.96	109.97
4	A	1294	G1P	O4-C4-C3	5.73	123.25	110.34
3	D	1293	THM	C4-N3-C2	5.74	120.21	115.25
4	A	1294	G1P	C1-O5-C5	5.80	125.00	113.75
4	A	1294	G1P	O5-C5-C4	6.18	121.29	109.68
4	A	1294	G1P	O3-C3-C4	6.24	124.38	110.34
4	A	1294	G1P	O5-C1-O1	6.53	119.97	111.36
3	C	1292	THM	C4-N3-C2	7.57	121.79	115.25
4	D	1294	G1P	O5-C1-O1	24.32	143.44	111.36

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
4	D	1294	G1P	C1

There are no torsion outliers.

There are no ring outliers.

8 monomers are involved in 35 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1294	G1P	13	0
3	B	1292	THM	9	0
4	B	1294	G1P	1	0
3	C	1292	THM	5	0
5	C	1295	SO4	4	0
3	D	1292	THM	4	0
3	D	1293	THM	1	0
4	D	1294	G1P	2	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 ⓘ

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

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

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

### 6.3 Carbohydrates ⓘ

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

### 6.4 Ligands ⓘ

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

### 6.5 Other polymers ⓘ

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