



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 07:58 pm GMT

PDB ID : 5IDT  
Title : Crystal Structure of a Glucose-1-phosphate Thymidyltransferase from  
Burkholderia vietnamiensis with bound Thymidine  
Authors : Seattle Structural Genomics Center for Infectious Disease (SSGCID)  
Deposited on : 2016-02-24  
Resolution : 2.35 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 : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : trunk28620  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
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) : recalc28949

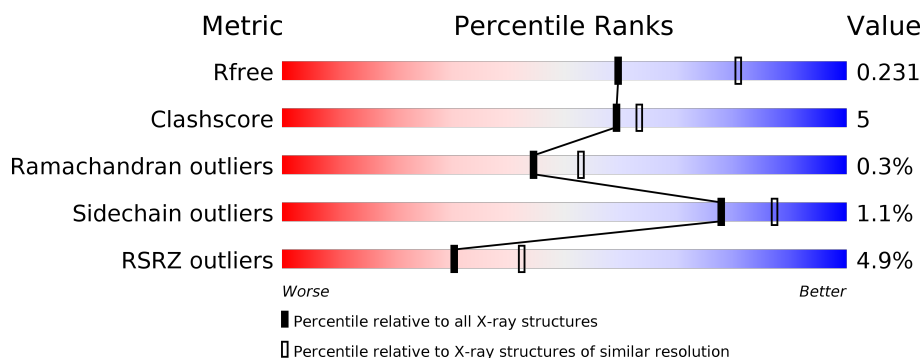
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.35 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))
$R_{free}$	100719	1522 (2.38-2.34)
Clashscore	112137	1626 (2.38-2.34)
Ramachandran outliers	110173	1605 (2.38-2.34)
Sidechain outliers	110143	1606 (2.38-2.34)
RSRZ outliers	101464	1528 (2.38-2.34)

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	302	<div> <div>3%</div> <div> <div></div> <div>84%</div> <div>11%</div> <div>• •</div> </div> </div>
1	B	302	<div> <div>8%</div> <div> <div></div> <div>82%</div> <div>14%</div> <div>•</div> </div> </div>
1	C	302	<div> <div>6%</div> <div> <div></div> <div>82%</div> <div>11%</div> <div>• 5%</div> </div> </div>
1	D	302	<div> <div>2%</div> <div> <div></div> <div>81%</div> <div>15%</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
2	THM	D	301	-	-	-	X
3	EDO	B	301	-	-	-	X
3	EDO	D	302	-	-	-	X

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8678 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 thymidyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	289	Total	C	N	O	S	0	1	0
			2157	1377	361	411	8			
1	B	289	Total	C	N	O	S	0	1	0
			2068	1329	344	386	9			
1	C	286	Total	C	N	O	S	0	1	0
			2108	1351	353	396	8			
1	D	289	Total	C	N	O	S	0	0	0
			2133	1367	361	397	8			

There are 32 discrepancies between the modelled and reference sequences:

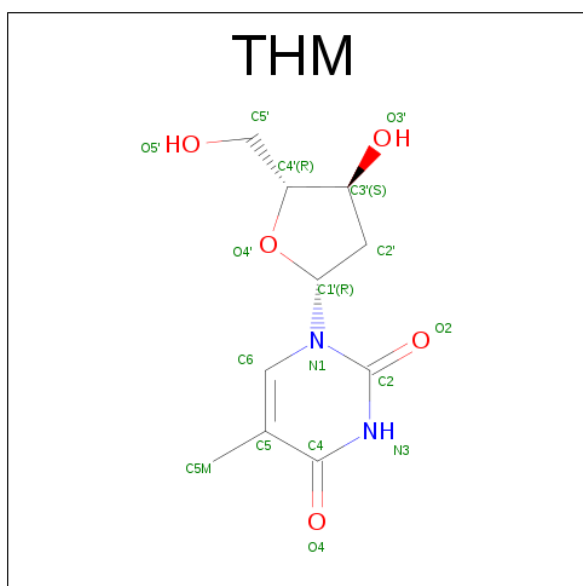
Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	MET	-	initiating methionine	UNP A4JIV2
A	-6	ALA	-	expression tag	UNP A4JIV2
A	-5	HIS	-	expression tag	UNP A4JIV2
A	-4	HIS	-	expression tag	UNP A4JIV2
A	-3	HIS	-	expression tag	UNP A4JIV2
A	-2	HIS	-	expression tag	UNP A4JIV2
A	-1	HIS	-	expression tag	UNP A4JIV2
A	0	HIS	-	expression tag	UNP A4JIV2
B	-7	MET	-	initiating methionine	UNP A4JIV2
B	-6	ALA	-	expression tag	UNP A4JIV2
B	-5	HIS	-	expression tag	UNP A4JIV2
B	-4	HIS	-	expression tag	UNP A4JIV2
B	-3	HIS	-	expression tag	UNP A4JIV2
B	-2	HIS	-	expression tag	UNP A4JIV2
B	-1	HIS	-	expression tag	UNP A4JIV2
B	0	HIS	-	expression tag	UNP A4JIV2
C	-7	MET	-	initiating methionine	UNP A4JIV2
C	-6	ALA	-	expression tag	UNP A4JIV2
C	-5	HIS	-	expression tag	UNP A4JIV2
C	-4	HIS	-	expression tag	UNP A4JIV2
C	-3	HIS	-	expression tag	UNP A4JIV2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-2	HIS	-	expression tag	UNP A4JIV2
C	-1	HIS	-	expression tag	UNP A4JIV2
C	0	HIS	-	expression tag	UNP A4JIV2
D	-7	MET	-	initiating methionine	UNP A4JIV2
D	-6	ALA	-	expression tag	UNP A4JIV2
D	-5	HIS	-	expression tag	UNP A4JIV2
D	-4	HIS	-	expression tag	UNP A4JIV2
D	-3	HIS	-	expression tag	UNP A4JIV2
D	-2	HIS	-	expression tag	UNP A4JIV2
D	-1	HIS	-	expression tag	UNP A4JIV2
D	0	HIS	-	expression tag	UNP A4JIV2

- Molecule 2 is THYMIDINE (three-letter code: THM) (formula:  $C_{10}H_{14}N_2O_5$ ).



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

- Molecule 3 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula:  $C_2H_6O_2$ ).



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

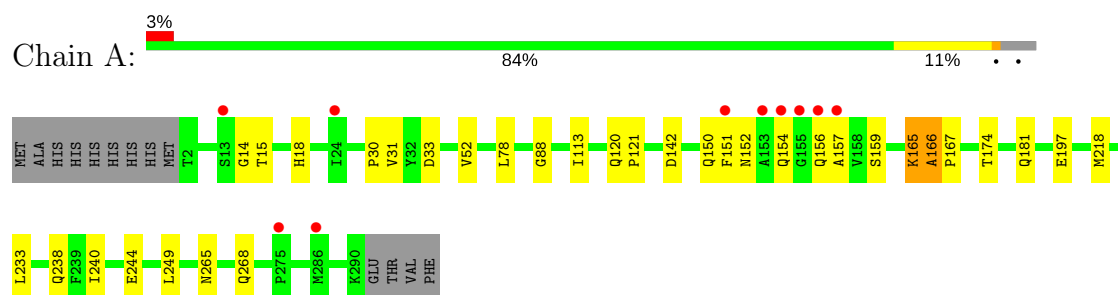
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	47	Total	O	0	0
			47	47		
4	B	28	Total	O	0	0
			28	28		
4	C	45	Total	O	0	0
			45	45		
4	D	38	Total	O	0	0
			38	38		

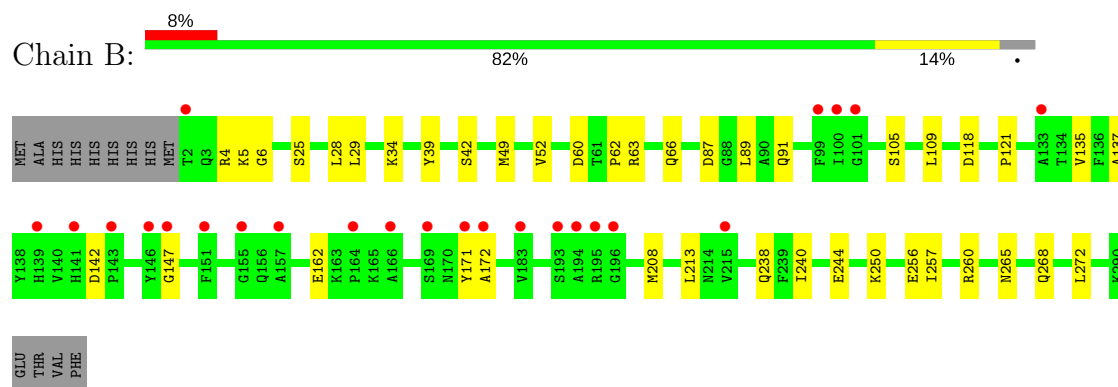
### 3 Residue-property plots [i](#)

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.

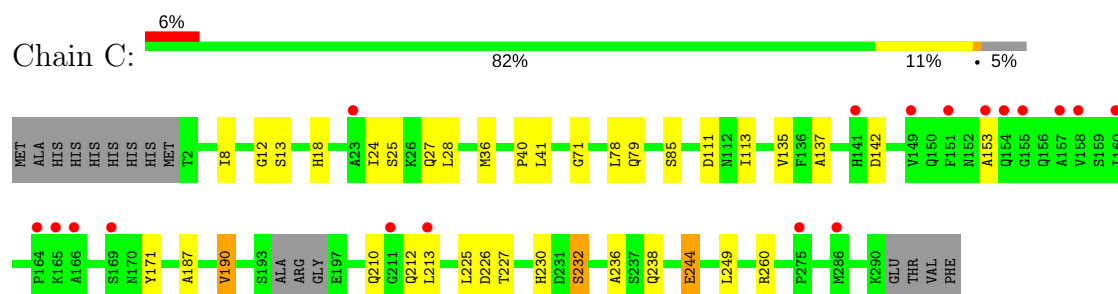
- Molecule 1: Glucose-1-phosphate thymidyltransferase



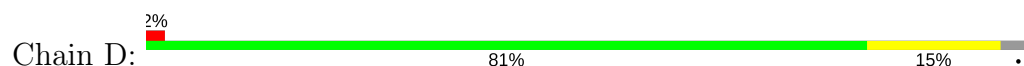
- Molecule 1: Glucose-1-phosphate thymidyltransferase

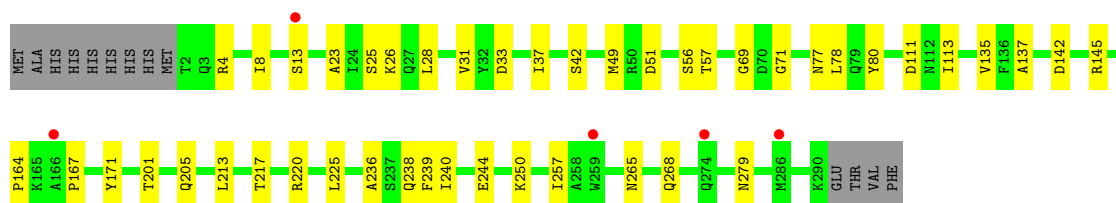


- Molecule 1: Glucose-1-phosphate thymidyltransferase



- Molecule 1: Glucose-1-phosphate thymidyltransferase







## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	112.11Å 114.86Å 95.78Å 90.00° 90.58° 90.00°	Depositor
Resolution (Å)	48.16 – 2.35 48.16 – 2.35	Depositor EDS
% Data completeness (in resolution range)	99.2 (48.16-2.35) 99.2 (48.16-2.35)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.98 (at 2.34Å)	Xtriage
Refinement program	PHENIX (dev_2271: ???)	Depositor
R, $R_{free}$	0.176 , 0.231 0.173 , 0.231	Depositor DCC
$R_{free}$ test set	1863 reflections (3.72%)	DCC
Wilson B-factor (Å <sup>2</sup> )	44.4	Xtriage
Anisotropy	0.163	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 44.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.016 for k,h,-l 0.011 for -k,-h,-l 0.016 for -h,-k,l	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	8678	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.45% 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: THM, EDO

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	0.45	0/2211	0.62	1/3017 (0.0%)
1	B	0.39	0/2121	0.55	0/2907
1	C	0.47	2/2160 (0.1%)	0.61	1/2953 (0.0%)
1	D	0.42	0/2183	0.58	1/2984 (0.0%)
All	All	0.44	2/8675 (0.0%)	0.59	3/11861 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	190	VAL	C-N	6.46	1.49	1.34
1	C	244	GLU	CG-CD	5.05	1.59	1.51

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	165	LYS	C-N-CA	7.60	140.69	121.70
1	C	190	VAL	C-N-CA	-6.09	106.47	121.70
1	D	13	SER	C-N-CA	-5.05	111.69	122.30

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	2157	0	2043	25	0
1	B	2068	0	1882	25	0
1	C	2108	0	1973	20	0
1	D	2133	0	2023	26	0
2	A	17	0	14	0	0
2	D	17	0	14	1	0
3	A	12	0	18	1	0
3	B	4	0	6	0	0
3	D	4	0	6	0	0
4	A	47	0	0	1	0
4	B	28	0	0	0	0
4	C	45	0	0	0	0
4	D	38	0	0	1	0
All	All	8678	0	7979	89	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 5.

All (89) 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:137:ALA:HB1	1:B:171:TYR:HB3	1.68	0.74
1:C:226:ASP:O	1:C:232:SER:OG	2.09	0.70
1:A:152:ASN:N	1:A:156:GLN:O	2.21	0.69
1:C:210:GLN:NE2	1:C:212:GLN:OE1	2.28	0.67
1:A:165:LYS:N	1:A:166:ALA:HB3	2.09	0.67
1:A:166:ALA:N	1:A:167:PRO:HD3	2.10	0.67
1:C:137:ALA:HB1	1:C:171:TYR:HB3	1.78	0.66
1:A:14:GLY:O	4:A:401:HOH:O	2.17	0.59
1:A:165:LYS:H	1:A:166:ALA:HB3	1.67	0.58
1:C:244:GLU:HG2	1:C:249:LEU:O	2.04	0.57
1:B:208:MET:HE2	1:B:213:LEU:HD23	1.85	0.57
1:B:28:LEU:HD11	1:B:63:ARG:HB3	1.86	0.56
1:A:30:PRO:HG3	1:B:29:LEU:HD22	1.86	0.56
1:D:26:LYS:NZ	4:D:403:HOH:O	2.36	0.56
1:D:240:ILE:O	1:D:244:GLU:HG3	2.07	0.55
1:A:174:THR:HG21	1:A:218:MET:HE1	1.90	0.54
1:D:201:THR:O	1:D:205:GLN:HG2	2.08	0.54
1:B:147:GLY:HA2	1:B:162:GLU:HA	1.89	0.53
1:B:268:GLN:O	1:B:272:LEU:HD13	2.08	0.53
1:D:137:ALA:HB1	1:D:171:TYR:HB3	1.89	0.53
1:B:256:GLU:OE1	1:B:260:ARG:NH1	2.42	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:225:LEU:HD13	1:C:236:ALA:HA	1.92	0.51
1:D:135:VAL:HG13	1:D:213:LEU:HD11	1.93	0.51
1:D:111:ASP:OD1	2:D:301:THM:H5'1	2.11	0.50
1:D:164:PRO:HG2	1:D:167:PRO:HB3	1.92	0.50
1:C:12:GLY:H	1:C:27:GLN:HE22	1.60	0.49
1:A:244:GLU:HG2	1:A:249:LEU:O	2.13	0.49
1:B:42:SER:HB2	1:B:257:ILE:CD1	2.43	0.49
1:B:63:ARG:NH1	1:C:79[B]:GLN:OE1	2.43	0.48
1:B:265:ASN:OD1	1:B:268:GLN:HG3	2.14	0.48
1:C:230:HIS:CE1	1:D:33:ASP:HB2	2.48	0.48
1:D:142:ASP:OD1	1:D:142:ASP:N	2.47	0.48
1:B:6:GLY:O	1:B:52:VAL:HA	2.13	0.47
1:A:151:PHE:HA	1:A:157:ALA:HA	1.95	0.47
1:D:250:LYS:HA	1:D:250:LYS:HD2	1.71	0.47
1:B:62:PRO:O	1:B:66:GLN:HG3	2.14	0.47
1:A:233:LEU:HA	1:A:233:LEU:HD12	1.56	0.47
1:C:28:LEU:HD22	1:D:23:ALA:HB1	1.97	0.47
1:D:142:ASP:O	1:D:145:ARG:HG3	2.14	0.47
1:D:56:SER:OG	1:D:57:THR:N	2.46	0.47
1:C:36:MET:O	1:C:40:PRO:HD2	2.15	0.47
1:D:71:GLY:HA3	1:D:78:LEU:HG	1.97	0.46
1:B:135:VAL:HG13	1:B:213:LEU:HD11	1.97	0.46
1:A:166:ALA:H	1:A:167:PRO:HD3	1.80	0.46
1:A:181:GLN:OE1	1:A:181:GLN:N	2.49	0.46
1:A:88:GLY:HA2	1:A:197:GLU:OE2	2.15	0.46
1:D:69:GLY:O	1:D:80:TYR:OH	2.14	0.46
1:C:113:ILE:HB	1:C:225:LEU:HB2	1.98	0.46
1:B:87:ASP:HB2	1:B:91:GLN:OE1	2.15	0.45
1:B:240:ILE:O	1:B:244:GLU:HG3	2.15	0.45
1:A:238:GLN:OE1	1:B:238:GLN:NE2	2.49	0.45
1:C:238:GLN:NE2	1:D:238:GLN:OE1	2.46	0.45
1:A:152:ASN:HA	3:A:303:EDO:O2	2.17	0.45
1:D:265:ASN:OD1	1:D:268:GLN:HG3	2.16	0.44
1:C:142:ASP:N	1:C:142:ASP:OD1	2.50	0.44
1:C:111:ASP:HB2	1:C:227:THR:OG1	2.18	0.44
1:A:15:THR:HG22	1:A:18:HIS:CD2	2.52	0.44
1:C:135:VAL:HG13	1:C:213:LEU:HD11	1.99	0.43
1:A:33:ASP:N	1:A:33:ASP:OD1	2.50	0.43
1:C:71:GLY:HA3	1:C:78:LEU:HG	2.00	0.43
1:D:31:VAL:HG21	1:D:113:ILE:HD11	2.00	0.43
1:D:8:ILE:HG21	1:D:37:ILE:HD11	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:42:SER:HB2	1:D:257:ILE:CD1	2.49	0.43
1:B:4:ARG:O	1:B:49:MET:HG2	2.19	0.43
1:B:5:LYS:O	1:B:105:SER:HA	2.19	0.43
1:D:225:LEU:HD11	1:D:239:PHE:CD2	2.54	0.43
1:B:34:LYS:HG3	1:B:39:TYR:CE1	2.53	0.42
1:A:142:ASP:N	1:A:142:ASP:OD1	2.52	0.42
1:A:244:GLU:HA	1:A:249:LEU:O	2.19	0.42
1:A:31:VAL:HG21	1:A:113:ILE:HD11	2.01	0.42
1:B:89:LEU:HD21	1:B:109:LEU:HD11	2.02	0.42
1:A:265:ASN:OD1	1:A:268:GLN:HG3	2.20	0.42
1:B:250:LYS:HD2	1:B:250:LYS:HA	1.84	0.41
1:C:24:ILE:HD12	1:C:24:ILE:HA	1.94	0.41
1:A:120:GLN:HB3	1:A:121:PRO:HD3	2.02	0.41
1:B:118:ASP:C	1:B:121:PRO:HD2	2.41	0.41
1:B:137:ALA:HA	1:B:172:ALA:O	2.20	0.41
1:A:152:ASN:C	1:A:154:GLN:H	2.23	0.41
1:A:240:ILE:O	1:A:244:GLU:HG3	2.19	0.41
1:C:18:HIS:O	1:D:279:ASN:ND2	2.51	0.41
1:D:28:LEU:HD23	1:D:28:LEU:HA	1.94	0.41
1:B:25:SER:HB2	1:B:60:ASP:OD1	2.21	0.41
1:D:4:ARG:O	1:D:49:MET:HG2	2.21	0.41
1:C:8:ILE:HD11	1:C:41:LEU:HA	2.02	0.41
1:D:51:ASP:OD1	1:D:77:ASN:HB3	2.21	0.41
1:A:52:VAL:HB	1:A:78:LEU:CD2	2.51	0.40
1:B:142:ASP:OD1	1:B:142:ASP:N	2.54	0.40
1:C:187:ALA:O	1:C:190:VAL:HG22	2.21	0.40
1:D:225:LEU:HD13	1:D:236:ALA:HA	2.02	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	288/302 (95%)	281 (98%)	6 (2%)	1 (0%)	44	53
1	B	288/302 (95%)	285 (99%)	3 (1%)	0	100	100
1	C	283/302 (94%)	275 (97%)	6 (2%)	2 (1%)	25	27
1	D	287/302 (95%)	282 (98%)	5 (2%)	0	100	100
All	All	1146/1208 (95%)	1123 (98%)	20 (2%)	3 (0%)	44	53

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	166	ALA
1	C	13	SER
1	C	153	ALA

### 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	212/246 (86%)	210 (99%)	2 (1%)	82	90
1	B	184/246 (75%)	184 (100%)	0	100	100
1	C	201/246 (82%)	197 (98%)	4 (2%)	60	73
1	D	205/246 (83%)	202 (98%)	3 (2%)	70	81
All	All	802/984 (82%)	793 (99%)	9 (1%)	78	87

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	150	GLN
1	A	159	SER
1	C	25	SER
1	C	85	SER
1	C	232	SER
1	C	260	ARG
1	D	25	SER
1	D	217	THR

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Mol	Chain	Res	Type
1	D	220	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	91	GLN

### 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 ⓘ

7 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	THM	A	301	-	13,18,18	4.38	7 (53%)	17,26,26	2.29	3 (17%)
3	EDO	A	302	-	3,3,3	0.34	0	2,2,2	0.61	0
3	EDO	A	303	-	3,3,3	0.51	0	2,2,2	0.32	0
3	EDO	A	304	-	3,3,3	0.49	0	2,2,2	0.62	0
3	EDO	B	301	-	3,3,3	0.54	0	2,2,2	0.28	0
2	THM	D	301	-	13,18,18	4.50	8 (61%)	17,26,26	2.34	5 (29%)
3	EDO	D	302	-	3,3,3	0.50	0	2,2,2	0.19	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	THM	A	301	-	-	0/2/18/18	0/2/2/2
3	EDO	A	302	-	-	0/1/1/1	0/0/0/0
3	EDO	A	303	-	-	0/1/1/1	0/0/0/0
3	EDO	A	304	-	-	0/1/1/1	0/0/0/0
3	EDO	B	301	-	-	0/1/1/1	0/0/0/0
2	THM	D	301	-	-	0/2/18/18	0/2/2/2
3	EDO	D	302	-	-	0/1/1/1	0/0/0/0

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	301	THM	C3'-C4'	-7.89	1.31	1.53
2	A	301	THM	C3'-C4'	-7.64	1.31	1.53
2	D	301	THM	O4'-C1'	-5.08	1.30	1.42
2	A	301	THM	O4'-C1'	-4.69	1.31	1.42
2	D	301	THM	O4-C4	-2.19	1.19	1.24
2	D	301	THM	O3'-C3'	2.47	1.48	1.43
2	A	301	THM	O3'-C3'	2.85	1.49	1.43
2	D	301	THM	C6-C5	3.44	1.49	1.40
2	A	301	THM	C6-C5	3.84	1.50	1.40
2	A	301	THM	C2-N3	5.70	1.49	1.38
2	D	301	THM	C2-N3	6.06	1.50	1.38
2	A	301	THM	C4-N3	6.68	1.45	1.33
2	D	301	THM	C4-N3	7.12	1.45	1.33
2	D	301	THM	O4'-C4'	7.87	1.62	1.45
2	A	301	THM	O4'-C4'	7.88	1.62	1.45

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301	THM	C5-C4-N3	-5.63	119.04	125.24
2	D	301	THM	C5-C4-N3	-4.99	119.74	125.24
2	D	301	THM	O4'-C1'-C2'	-2.77	100.94	106.25
2	D	301	THM	C4'-O4'-C1'	-2.65	102.93	109.42
2	A	301	THM	C2'-C3'-C4'	2.40	107.84	102.73
2	D	301	THM	O4'-C1'-N1	3.87	114.30	107.78
2	D	301	THM	C4-N3-C2	5.54	120.00	115.16
2	A	301	THM	C4-N3-C2	6.05	120.45	115.16



There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	303	EDO	1	0
2	D	301	THM	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	289/302 (95%)	0.22	10 (3%) 44 57	34, 51, 93, 134	0
1	B	289/302 (95%)	0.41	24 (8%) 12 17	39, 71, 117, 140	0
1	C	286/302 (94%)	0.41	18 (6%) 21 29	34, 57, 100, 133	0
1	D	289/302 (95%)	0.10	5 (1%) 70 80	36, 60, 88, 153	0
All	All	1153/1208 (95%)	0.28	57 (4%) 30 42	34, 59, 105, 153	0

All (57) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	151	PHE	6.2
1	A	155	GLY	4.5
1	B	171	TYR	4.1
1	A	153	ALA	4.0
1	C	286	MET	4.0
1	B	194	ALA	3.9
1	C	153	ALA	3.9
1	B	193	SER	3.8
1	D	13	SER	3.8
1	C	166	ALA	3.7
1	C	213	LEU	3.5
1	C	151	PHE	3.5
1	B	164	PRO	3.5
1	B	133	ALA	3.5
1	B	2	THR	3.5
1	B	196	GLY	3.5
1	B	139	HIS	3.4
1	B	141	HIS	3.4
1	A	156	GLN	3.3
1	B	146	TYR	3.3
1	C	275	PRO	3.3

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Mol	Chain	Res	Type	RSRZ
1	B	169	SER	3.3
1	B	166	ALA	3.2
1	B	151	PHE	3.2
1	D	274	GLN	3.1
1	B	155	GLY	3.1
1	A	157	ALA	3.1
1	A	275	PRO	3.0
1	A	24	ILE	3.0
1	C	164	PRO	3.0
1	B	143	PRO	2.8
1	C	155	GLY	2.7
1	B	195	ARG	2.7
1	B	100	ILE	2.7
1	B	147	GLY	2.7
1	A	286	MET	2.5
1	B	157	ALA	2.4
1	D	166	ALA	2.4
1	C	157	ALA	2.4
1	B	215	VAL	2.4
1	C	165	LYS	2.4
1	B	101	GLY	2.3
1	C	141	HIS	2.3
1	C	154	GLN	2.3
1	C	149	VAL	2.3
1	C	158	VAL	2.2
1	B	99	PHE	2.2
1	B	172	ALA	2.2
1	D	286	MET	2.2
1	C	160	ILE	2.2
1	A	154	GLN	2.1
1	D	259	TRP	2.1
1	B	183	VAL	2.1
1	A	13	SER	2.1
1	C	23	ALA	2.1
1	C	169	SER	2.0
1	C	211	GLY	2.0

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

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. 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
2	THM	D	301	17/17	0.82	0.30	5.97	80,93,99,103	0
3	EDO	D	302	4/4	0.94	0.20	3.69	57,60,63,64	0
3	EDO	B	301	4/4	0.79	0.21	3.64	64,74,79,81	0
2	THM	A	301	17/17	0.85	0.22	1.33	59,81,88,94	0
3	EDO	A	304	4/4	0.85	0.17	-	67,69,70,78	0
3	EDO	A	303	4/4	0.87	0.16	-	77,79,83,84	0
3	EDO	A	302	4/4	0.89	0.22	-	54,57,62,67	0

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

There are no such residues in this entry.