



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 14, 2020 – 04:30 am BST

PDB ID : 1LVW
Title : Crystal structure of glucose-1-phosphate thymidyltransferase, RmlA, complex with dTDP
Authors : Dong, A.; Christendat, D.; Pai, E.F.; Northeast Structural Genomics Consortium (NESG)
Deposited on : 2002-05-29
Resolution : 1.70 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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

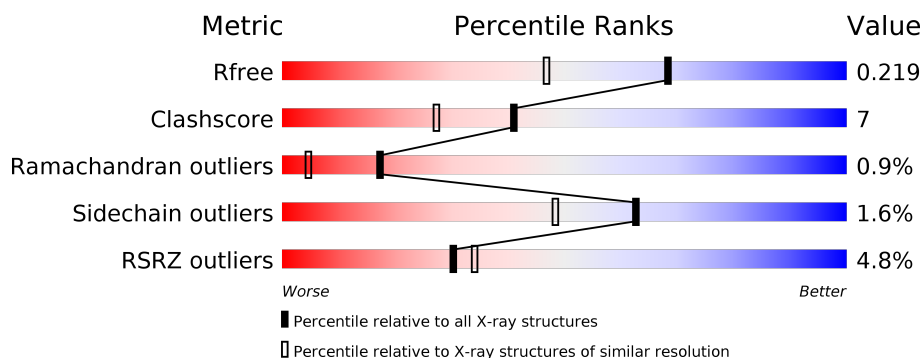
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	4298 (1.70-1.70)
Clashscore	141614	4695 (1.70-1.70)
Ramachandran outliers	138981	4610 (1.70-1.70)
Sidechain outliers	138945	4610 (1.70-1.70)
RSRZ outliers	127900	4222 (1.70-1.70)

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 ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	295	<div> <div>3%</div> <div> <div></div> <div>88%</div> <div>11%</div> <div>.</div> </div> </div>
1	B	295	<div> <div>2%</div> <div> <div></div> <div>86%</div> <div>12%</div> <div>..</div> </div> </div>
1	C	295	<div> <div>4%</div> <div> <div></div> <div>82%</div> <div>15%</div> <div>..</div> </div> </div>
1	D	295	<div> <div>10%</div> <div> <div></div> <div>80%</div> <div>19%</div> <div>.</div> </div> </div>

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 10472 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	295	Total	C	N	O	S	0	7	0
			2392	1529	408	447	8			
1	B	292	Total	C	N	O	S	0	7	0
			2376	1521	402	444	9			
1	C	292	Total	C	N	O	S	0	4	0
			2341	1503	397	433	8			
1	D	291	Total	C	N	O	S	0	4	0
			2339	1499	396	436	8			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	CLONING ARTIFACT	UNP O27819
A	-1	ALA	-	CLONING ARTIFACT	UNP O27819
A	0	HIS	-	CLONING ARTIFACT	UNP O27819
B	-2	GLY	-	CLONING ARTIFACT	UNP O27819
B	-1	ALA	-	CLONING ARTIFACT	UNP O27819
B	0	HIS	-	CLONING ARTIFACT	UNP O27819
C	-2	GLY	-	CLONING ARTIFACT	UNP O27819
C	-1	ALA	-	CLONING ARTIFACT	UNP O27819
C	0	HIS	-	CLONING ARTIFACT	UNP O27819
D	-2	GLY	-	CLONING ARTIFACT	UNP O27819
D	-1	ALA	-	CLONING ARTIFACT	UNP O27819
D	0	HIS	-	CLONING ARTIFACT	UNP O27819

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Cl	0	0
			1	1		
2	A	1	Total	Cl	0	0
			1	1		

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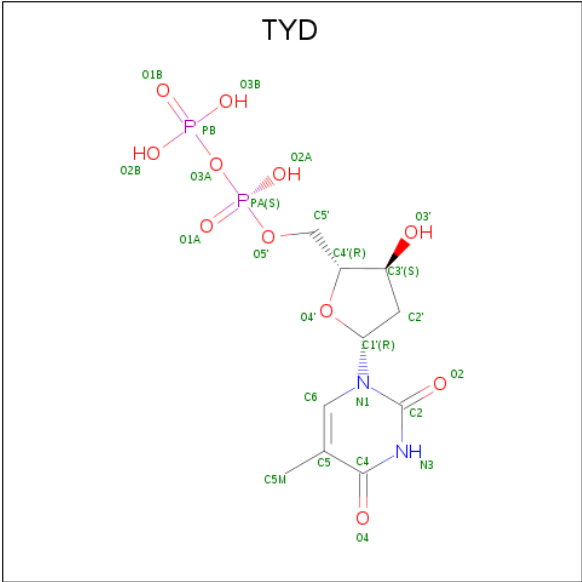
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	D	1	Total	Cl	0	0
			1	1		
2	C	2	Total	Cl	0	0
			2	2		

- Molecule 3 is SULFATE ION (three-letter code: SO₄) (formula: O₄S).



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

- Molecule 4 is THYMIDINE-5'-DIPHOSPHATE (three-letter code: TYD) (formula: C₁₀H₁₆N₂O₁₁P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			25	10	2	11	2		
4	A	1	Total	C	N	O	P	0	0
			25	10	2	11	2		
4	B	1	Total	C	N	O	P	0	0
			25	10	2	11	2		
4	B	1	Total	C	N	O	P	0	0
			25	10	2	11	2		
4	C	1	Total	C	N	O	P	0	0
			25	10	2	11	2		
4	C	1	Total	C	N	O	P	0	0
			25	10	2	11	2		
4	D	1	Total	C	N	O	P	0	0
			25	10	2	11	2		
4	D	1	Total	C	N	O	P	0	0
			25	10	2	11	2		

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



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

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	221	Total	O	0	0
			221	221		

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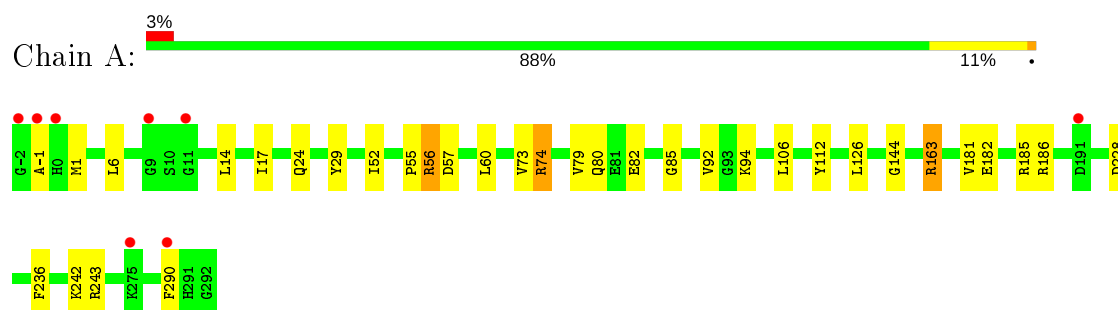
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	218	Total 218	O 218	0	0
6	C	175	Total 175	O 175	0	0
6	D	129	Total 129	O 129	0	0

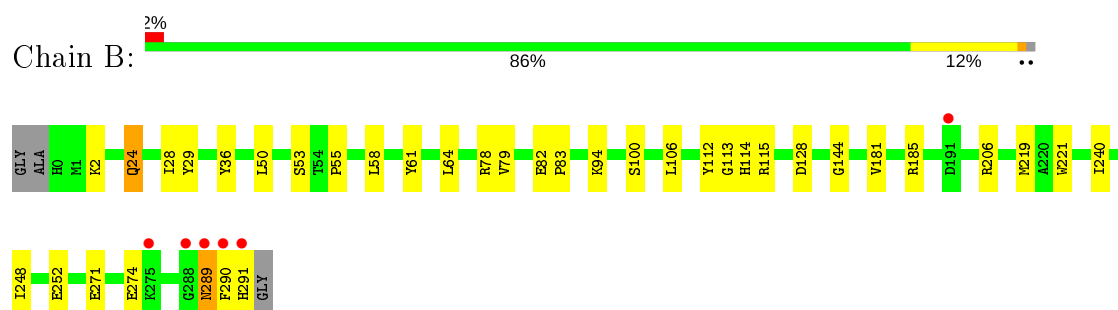
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

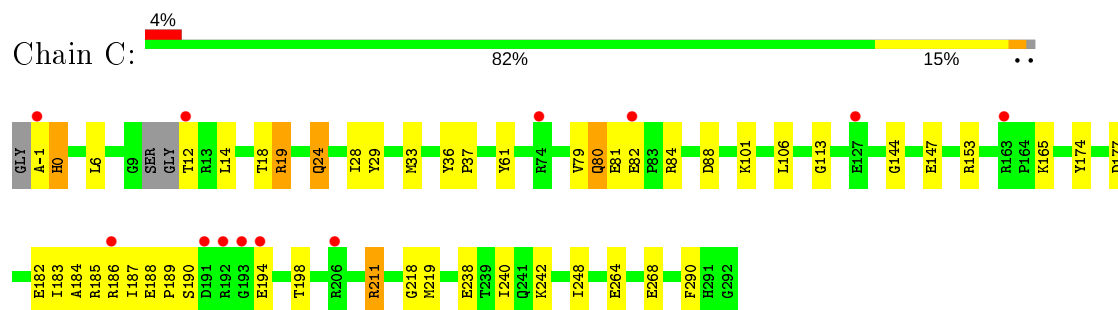
- Molecule 1: glucose-1-phosphate thymidyltransferase



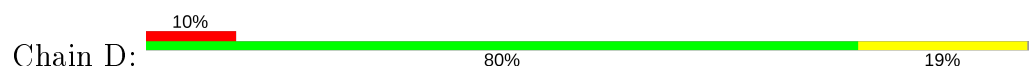
- Molecule 1: glucose-1-phosphate thymidyltransferase

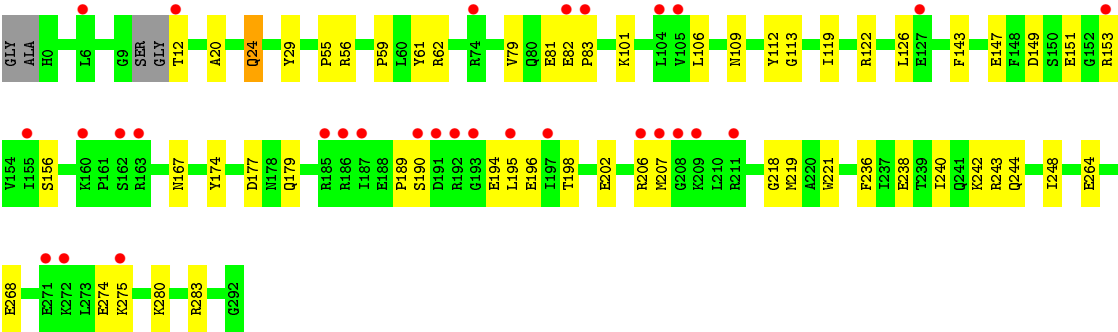


- Molecule 1: glucose-1-phosphate thymidyltransferase



- Molecule 1: glucose-1-phosphate thymidyltransferase





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	112.68Å 115.88Å 116.91Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	31.52 – 1.70 31.52 – 1.70	Depositor EDS
% Data completeness (in resolution range)	95.9 (31.52-1.70) 95.6 (31.52-1.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.34 (at 1.70Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.187 , 0.215 0.193 , 0.219	Depositor DCC
R_{free} test set	8072 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	15.4	Xtriage
Anisotropy	0.096	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 51.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.015 for -h,l,k 0.016 for -l,-k,-h 0.015 for k,h,-l 0.003 for k,l,h 0.003 for l,h,k	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	10472	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.09% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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: TYD, GOL, SO4, CL

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	0.58	0/2442	0.77	1/3299 (0.0%)
1	B	0.59	0/2425	0.75	1/3275 (0.0%)
1	C	0.56	0/2389	0.73	2/3228 (0.1%)
1	D	0.51	0/2387	0.71	0/3224
All	All	0.56	0/9643	0.74	4/13026 (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	B	144	GLY	N-CA-C	-5.97	98.18	113.10
1	C	144	GLY	N-CA-C	-5.31	99.83	113.10
1	C	80	GLN	N-CA-C	-5.12	97.17	111.00
1	A	144	GLY	N-CA-C	-5.04	100.50	113.10

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	2392	0	2367	32	0
1	B	2376	0	2359	35	0
1	C	2341	0	2324	38	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	2339	0	2311	42	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	2	0	0	1	0
2	D	1	0	0	1	0
3	A	5	0	0	0	0
3	D	5	0	0	0	0
4	A	50	0	26	4	0
4	B	50	0	26	2	0
4	C	50	0	26	2	0
4	D	50	0	26	3	0
5	A	18	0	24	0	0
5	B	18	0	24	1	0
5	C	18	0	24	0	0
5	D	12	0	16	2	0
6	A	221	0	0	0	0
6	B	218	0	0	2	0
6	C	175	0	0	1	0
6	D	129	0	0	2	0
All	All	10472	0	9553	137	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 7.

All (137) 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:82:GLU:HG3	1:B:83:PRO:HD2	1.31	1.07
1:B:82:GLU:HG3	1:B:83:PRO:CD	2.12	0.80
1:C:19:ARG:HH11	1:C:19:ARG:CG	1.93	0.80
1:C:12:THR:HG22	1:C:14:LEU:H	1.45	0.80
1:B:55:PRO:HG3	1:B:79:VAL:HG11	1.65	0.77
1:D:280:LYS:HE3	1:D:283:ARG:HH22	1.52	0.73
1:A:236:PHE:HE1	1:C:240:ILE:HD11	1.54	0.73
1:B:94:LYS:HE3	1:B:185:ARG:NH2	2.03	0.73
1:A:182:GLU:HG3	1:A:186:ARG:NH1	2.06	0.71
1:B:94:LYS:HE3	1:B:185:ARG:HH22	1.56	0.70
1:C:81:GLU:HB2	1:C:84:ARG:HH21	1.56	0.70
1:A:242:LYS:CE	1:A:243:ARG:HH12	2.06	0.69
1:B:78:ARG:HH11	1:B:78:ARG:HB3	1.58	0.69
1:B:291:HIS:HB2	1:D:167:ASN:OD1	1.93	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:78:ARG:NH1	1:B:78:ARG:HB3	2.09	0.67
1:A:182:GLU:HG3	1:A:186:ARG:CZ	2.24	0.67
1:C:198:THR:HG22	2:C:4003:CL:CL	2.32	0.67
1:A:163:ARG:HH11	1:A:163:ARG:HB3	1.61	0.65
1:D:280:LYS:HG3	1:D:283:ARG:NH2	2.11	0.64
1:C:190:SER:OG	1:C:194:GLU:HG2	1.98	0.64
1:C:153:ARG:HB2	1:C:211:ARG:NH2	2.13	0.63
1:B:113:GLY:HA3	1:B:219[A]:MET:HE2	1.80	0.61
1:D:20:ALA:C	1:D:56:ARG:HH12	2.03	0.61
1:C:147:GLU:OE1	1:C:165:LYS:HD3	2.02	0.60
1:C:19:ARG:HH11	1:C:19:ARG:HG3	1.65	0.60
1:A:242:LYS:HE2	1:A:243:ARG:HH12	1.66	0.60
1:B:55:PRO:HG3	1:B:79:VAL:CG1	2.31	0.60
5:B:2011:GOL:H11	1:D:221:TRP:O	2.01	0.59
1:B:113:GLY:HA3	1:B:219[A]:MET:CE	2.32	0.59
1:C:240:ILE:HG21	1:C:248[B]:ILE:HD11	1.83	0.59
1:D:143:PHE:HE1	5:D:2006:GOL:H32	1.68	0.59
1:C:19:ARG:HH11	1:C:19:ARG:HG2	1.66	0.58
1:B:94:LYS:HG3	1:B:181:VAL:HG11	1.86	0.58
1:C:-1:ALA:O	1:C:0:HIS:O	2.20	0.58
1:A:56:ARG:HH12	1:A:60:LEU:HD11	1.67	0.58
1:D:280:LYS:HE3	1:D:283:ARG:NH2	2.18	0.58
1:B:113:GLY:HA3	1:B:219[B]:MET:SD	2.44	0.58
1:B:206:ARG:NE	6:B:1671:HOH:O	2.36	0.58
1:A:106:LEU:HB3	4:A:3001:TYD:H4'	1.86	0.57
1:C:101:LYS:HG2	1:C:177:ASP:HA	1.86	0.57
1:B:252:GLU:HG2	6:B:1430:HOH:O	2.04	0.57
1:D:198:THR:HG22	2:D:4004:CL:CL	2.42	0.57
1:D:238[A]:GLU:OE2	1:D:242:LYS:HE3	2.06	0.56
1:D:190:SER:HB3	1:D:196:GLU:OE2	2.04	0.56
1:D:147:GLU:HB3	1:D:156:SER:HB3	1.86	0.56
1:A:74:ARG:N	1:A:74:ARG:HD3	2.20	0.55
1:D:112:TYR:CE2	1:D:240:ILE:HD13	2.42	0.55
1:A:73:VAL:C	1:A:74:ARG:HD3	2.27	0.55
1:D:143:PHE:CE1	5:D:2006:GOL:H32	2.42	0.54
1:A:185:ARG:HB3	1:A:185:ARG:NH1	2.22	0.54
1:D:149:ASP:OD2	1:D:153:ARG:HB3	2.08	0.54
1:C:12:THR:HG21	1:C:18:THR:OG1	2.08	0.54
1:B:64:LEU:HA	1:C:19:ARG:HD3	1.90	0.54
1:C:84:ARG:HD2	1:C:88:ASP:OD2	2.09	0.53
1:B:112:TYR:OH	1:D:218:GLY:HA2	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:264:GLU:O	1:C:268:GLU:HG3	2.10	0.52
1:D:12:THR:N	6:D:1133:HOH:O	2.42	0.52
1:A:56:ARG:NH2	1:A:57:ASP:OD1	2.42	0.52
1:D:106:LEU:HB3	4:D:3007:TYD:H4'	1.92	0.52
1:B:24:GLN:HB2	1:B:61:TYR:OH	2.09	0.51
1:C:113:GLY:HA3	1:C:219:MET:SD	2.51	0.51
1:D:206:ARG:O	1:D:206:ARG:HG3	2.11	0.51
1:A:112:TYR:OH	1:C:218:GLY:HA2	2.11	0.51
1:D:113:GLY:HA3	1:D:219:MET:SD	2.51	0.50
1:A:55:PRO:HG3	1:A:79:VAL:HG11	1.93	0.50
1:C:182:GLU:OE2	1:C:185:ARG:NE	2.45	0.50
1:D:244:GLN:HB3	6:D:1059:HOH:O	2.10	0.50
1:D:240:ILE:HG21	1:D:248[A]:ILE:HD11	1.93	0.49
1:B:94:LYS:CE	1:B:185:ARG:NH2	2.73	0.49
1:C:19:ARG:HG3	1:C:19:ARG:NH1	2.27	0.49
1:B:106:LEU:HD13	4:B:3003:TYD:H5'1	1.94	0.49
1:C:81:GLU:H	1:C:84:ARG:NH2	2.10	0.49
1:A:182:GLU:CG	1:A:186:ARG:CZ	2.91	0.48
1:A:55:PRO:HG3	1:A:79:VAL:CG1	2.43	0.48
1:D:119:ILE:HG22	1:D:122:ARG:NH2	2.28	0.48
1:B:289:ASN:O	1:B:290:PHE:HB2	2.13	0.48
1:A:80:GLN:HG2	1:A:82:GLU:O	2.14	0.48
1:B:53[A]:SER:OG	1:B:58:LEU:CD1	2.62	0.48
1:B:106:LEU:HB3	4:B:3003:TYD:H4'	1.96	0.47
1:D:106:LEU:HD13	4:D:3007:TYD:H5'1	1.95	0.47
1:A:182:GLU:CG	1:A:186:ARG:NH2	2.77	0.47
1:A:242:LYS:HE2	1:A:243:ARG:HH22	1.79	0.47
1:B:271:GLU:O	1:B:274:GLU:HG2	2.15	0.47
1:B:221:TRP:O	1:D:243[B]:ARG:HG2	2.15	0.46
1:D:59:PRO:HA	1:D:62:ARG:HD2	1.98	0.46
1:C:211:ARG:HD3	6:C:1477:HOH:O	2.16	0.46
1:D:236:PHE:CZ	1:D:240:ILE:HD11	2.50	0.46
1:C:12:THR:HG22	1:C:14:LEU:N	2.23	0.45
1:C:79[A]:VAL:CG1	1:C:80:GLN:N	2.79	0.45
1:D:198:THR:O	1:D:202:GLU:HG2	2.17	0.45
1:C:6:LEU:HG	4:C:3005:TYD:C2	2.47	0.45
1:A:85:GLY:HA2	4:A:3001:TYD:H52	1.98	0.45
1:C:106:LEU:HB3	4:C:3005:TYD:H4'	1.99	0.45
1:C:238:GLU:OE2	1:C:242:LYS:HE3	2.16	0.44
1:A:163:ARG:CB	1:A:163:ARG:HH11	2.28	0.44
1:B:112:TYR:CZ	1:D:218:GLY:HA2	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:82:GLU:O	1:C:84:ARG:HG3	2.18	0.44
1:A:228[B]:ASP:OD1	1:D:238[B]:GLU:OE2	2.34	0.44
1:C:183:ILE:O	1:C:187:ILE:HG23	2.16	0.44
1:B:78:ARG:CB	1:B:78:ARG:HH11	2.29	0.44
1:D:106:LEU:HD13	4:D:3007:TYD:C5'	2.47	0.44
1:B:114:HIS:CD2	1:B:115:ARG:HG3	2.53	0.43
1:D:24:GLN:HB2	1:D:61:TYR:OH	2.18	0.43
1:C:24:GLN:HB2	1:C:61:TYR:OH	2.19	0.43
1:D:55:PRO:HG3	1:D:79:VAL:HG21	2.00	0.43
1:A:112:TYR:CZ	1:C:218:GLY:HA2	2.54	0.43
1:B:53[A]:SER:OG	1:B:58:LEU:HD13	2.18	0.43
1:D:177:ASP:OD1	1:D:179:GLN:HB2	2.18	0.43
1:B:240:ILE:HG21	1:B:248:ILE:HD11	2.02	0.42
1:D:81:GLU:C	1:D:82:GLU:HG3	2.39	0.42
1:D:81:GLU:O	1:D:82:GLU:HG3	2.20	0.42
1:A:106:LEU:HD13	4:A:3001:TYD:C5'	2.50	0.42
1:A:14:LEU:O	1:A:17:ILE:HG12	2.20	0.42
1:B:50:LEU:HD11	1:B:78:ARG:HG2	2.01	0.42
1:D:202:GLU:O	1:D:206:ARG:HG2	2.20	0.41
1:C:184:ALA:O	1:C:187:ILE:HG12	2.20	0.41
1:C:33:MET:O	1:C:37:PRO:HD2	2.20	0.41
1:A:163:ARG:HH11	1:A:163:ARG:CG	2.33	0.41
1:A:56:ARG:HH12	1:A:60:LEU:CD1	2.32	0.41
1:B:2[A]:LYS:HE3	1:B:100:SER:OG	2.19	0.41
1:D:149:ASP:OD1	1:D:151:GLU:N	2.50	0.41
1:A:1:MET:CE	1:A:126[A]:LEU:HD21	2.50	0.41
1:B:64:LEU:HD12	1:C:19:ARG:HD3	2.03	0.41
1:D:274:GLU:O	1:D:275:LYS:CB	2.68	0.41
1:A:52:ILE:HD13	1:A:92[B]:VAL:CG2	2.51	0.41
1:A:6:LEU:HG	4:A:3001:TYD:O2	2.21	0.41
1:C:182:GLU:HG3	1:C:186:ARG:CZ	2.51	0.41
1:D:101:LYS:HG2	1:D:177:ASP:HA	2.03	0.41
1:B:28:ILE:HB	1:B:36:TYR:CE1	2.57	0.40
1:C:188:GLU:HA	1:C:189:PRO:HD3	1.94	0.40
1:D:190:SER:OG	1:D:194:GLU:HB2	2.21	0.40
1:A:94:LYS:HG3	1:A:181:VAL:HG11	2.03	0.40
1:B:289:ASN:HD22	1:B:289:ASN:HA	1.61	0.40
1:A:185:ARG:HH11	1:A:185:ARG:CB	2.35	0.40
1:C:28:ILE:HB	1:C:36:TYR:CE1	2.56	0.40
1:D:264:GLU:OE2	1:D:268:GLU:HG3	2.22	0.40
1:D:82:GLU:HA	1:D:83:PRO:HD3	1.91	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	300/295 (102%)	294 (98%)	3 (1%)	3 (1%)	15	4
1	B	297/295 (101%)	290 (98%)	6 (2%)	1 (0%)	41	24
1	C	292/295 (99%)	283 (97%)	6 (2%)	3 (1%)	15	4
1	D	291/295 (99%)	281 (97%)	7 (2%)	3 (1%)	15	4
All	All	1180/1180 (100%)	1148 (97%)	22 (2%)	10 (1%)	17	6

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	0	HIS
1	D	207	MET
1	A	-1	ALA
1	D	29	TYR
1	A	290	PHE
1	B	29	TYR
1	C	29	TYR
1	A	29	TYR
1	C	290	PHE
1	D	189	PRO

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	252/248 (102%)	248 (98%)	4 (2%)	62	48
1	B	252/248 (102%)	249 (99%)	3 (1%)	71	59
1	C	245/248 (99%)	241 (98%)	4 (2%)	62	48
1	D	245/248 (99%)	240 (98%)	5 (2%)	55	38
All	All	994/992 (100%)	978 (98%)	16 (2%)	62	48

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	GLN
1	A	56	ARG
1	A	74	ARG
1	A	163	ARG
1	B	24	GLN
1	B	128	ASP
1	B	289	ASN
1	C	19	ARG
1	C	24	GLN
1	C	174	TYR
1	C	211	ARG
1	D	24	GLN
1	D	109	ASN
1	D	126	LEU
1	D	174	TYR
1	D	195	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	289	ASN
1	D	201	ASN

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 [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 26 ligands modelled in this entry, 5 are monoatomic - leaving 21 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
4	TYD	B	3004	-	23,26,26	2.27	8 (34%)	33,40,40	1.77	3 (9%)
4	TYD	C	3006	-	23,26,26	2.38	8 (34%)	33,40,40	1.80	5 (15%)
5	GOL	B	2011	-	5,5,5	0.16	0	5,5,5	0.23	0
4	TYD	A	3002	-	23,26,26	2.17	6 (26%)	33,40,40	1.72	3 (9%)
5	GOL	A	2007	-	5,5,5	0.30	0	5,5,5	0.29	0
5	GOL	D	2004	-	5,5,5	0.26	0	5,5,5	0.21	0
5	GOL	B	2002	-	5,5,5	0.37	0	5,5,5	0.20	0
4	TYD	B	3003	-	23,26,26	2.22	5 (21%)	33,40,40	1.89	4 (12%)
4	TYD	C	3005	-	23,26,26	2.12	8 (34%)	33,40,40	1.86	4 (12%)
3	SO4	D	5002	-	4,4,4	0.39	0	6,6,6	0.32	0
4	TYD	D	3008	-	23,26,26	2.16	7 (30%)	33,40,40	1.74	3 (9%)
5	GOL	C	2001	-	5,5,5	0.29	0	5,5,5	0.28	0
5	GOL	B	2008	-	5,5,5	0.43	0	5,5,5	0.22	0
5	GOL	D	2006	-	5,5,5	0.32	0	5,5,5	0.40	0
5	GOL	A	2005	-	5,5,5	0.36	0	5,5,5	0.21	0
5	GOL	C	2003	-	5,5,5	0.35	0	5,5,5	0.26	0
5	GOL	A	2009	-	5,5,5	0.36	0	5,5,5	0.31	0
5	GOL	C	2010	-	5,5,5	0.21	0	5,5,5	0.28	0
4	TYD	D	3007	-	23,26,26	2.31	9 (39%)	33,40,40	1.82	4 (12%)
3	SO4	A	5001	-	4,4,4	0.30	0	6,6,6	0.11	0
4	TYD	A	3001	-	23,26,26	2.08	8 (34%)	33,40,40	1.75	5 (15%)

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
4	TYD	B	3004	-	-	2/16/28/28	0/2/2/2
4	TYD	C	3006	-	-	0/16/28/28	0/2/2/2
5	GOL	B	2011	-	-	0/4/4/4	-
4	TYD	A	3002	-	-	1/16/28/28	0/2/2/2
5	GOL	A	2007	-	-	0/4/4/4	-
5	GOL	D	2004	-	-	0/4/4/4	-
5	GOL	B	2002	-	-	0/4/4/4	-
4	TYD	B	3003	-	-	1/16/28/28	0/2/2/2
4	TYD	C	3005	-	-	3/16/28/28	0/2/2/2
4	TYD	D	3008	-	-	0/16/28/28	0/2/2/2
5	GOL	C	2001	-	-	0/4/4/4	-
5	GOL	B	2008	-	-	0/4/4/4	-
5	GOL	D	2006	-	-	0/4/4/4	-
5	GOL	A	2005	-	-	0/4/4/4	-
5	GOL	C	2003	-	-	0/4/4/4	-
5	GOL	A	2009	-	-	2/4/4/4	-
4	TYD	D	3007	-	-	1/16/28/28	0/2/2/2
5	GOL	C	2010	-	-	0/4/4/4	-
4	TYD	A	3001	-	-	1/16/28/28	0/2/2/2

All (59) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	3005	TYD	C6-C5	-6.00	1.33	1.51
4	A	3002	TYD	C6-C5	-5.99	1.33	1.51
4	C	3006	TYD	C6-C5	-5.97	1.33	1.51
4	D	3007	TYD	C6-C5	-5.94	1.33	1.51
4	B	3003	TYD	C6-C5	-5.93	1.33	1.51
4	D	3008	TYD	C6-C5	-5.84	1.33	1.51
4	B	3004	TYD	C6-C5	-5.68	1.34	1.51
4	A	3001	TYD	C6-C5	-5.64	1.34	1.51
4	B	3003	TYD	C1'-N1	5.18	1.52	1.45
4	D	3007	TYD	C1'-N1	5.03	1.52	1.45
4	C	3006	TYD	C6-N1	-4.87	1.40	1.46
4	C	3006	TYD	C1'-N1	4.79	1.52	1.45
4	B	3004	TYD	C6-N1	-4.71	1.40	1.46
4	D	3008	TYD	C1'-N1	4.65	1.51	1.45
4	B	3003	TYD	C6-N1	-4.62	1.41	1.46
4	B	3004	TYD	C1'-N1	4.61	1.51	1.45
4	A	3001	TYD	C1'-N1	4.28	1.51	1.45
4	D	3007	TYD	C6-N1	-4.24	1.41	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	3002	TYD	C6-N1	-4.21	1.41	1.46
4	C	3005	TYD	C1'-N1	4.16	1.51	1.45
4	D	3008	TYD	C6-N1	-4.04	1.41	1.46
4	A	3001	TYD	C6-N1	-3.93	1.41	1.46
4	C	3005	TYD	C6-N1	-3.85	1.42	1.46
4	A	3002	TYD	C1'-N1	3.75	1.50	1.45
4	C	3006	TYD	C2-N1	3.64	1.40	1.35
4	A	3002	TYD	C2-N1	3.16	1.40	1.35
4	B	3004	TYD	C2-N1	3.14	1.40	1.35
4	D	3007	TYD	C2-N1	3.13	1.40	1.35
4	D	3008	TYD	PB-O1B	-2.72	1.41	1.50
4	A	3002	TYD	PB-O1B	-2.70	1.41	1.50
4	B	3004	TYD	PB-O2B	-2.69	1.44	1.54
4	B	3004	TYD	PB-O1B	-2.61	1.42	1.50
4	A	3002	TYD	PB-O2B	-2.59	1.44	1.54
4	C	3006	TYD	PB-O2B	-2.54	1.45	1.54
4	C	3006	TYD	PB-O1B	-2.52	1.42	1.50
4	D	3007	TYD	PB-O3B	2.51	1.64	1.54
4	D	3007	TYD	PB-O1B	-2.46	1.42	1.50
4	C	3005	TYD	C2-N1	2.46	1.39	1.35
4	C	3006	TYD	C4-N3	2.45	1.41	1.37
4	B	3004	TYD	PA-O2A	-2.42	1.44	1.55
4	D	3008	TYD	PA-O2A	-2.38	1.44	1.55
4	A	3001	TYD	PB-O3B	2.36	1.64	1.54
4	D	3008	TYD	PB-O2B	-2.32	1.45	1.54
4	C	3006	TYD	PA-O2A	-2.29	1.44	1.55
4	D	3007	TYD	O4-C4	-2.28	1.18	1.23
4	D	3008	TYD	C2-N1	2.25	1.38	1.35
4	A	3001	TYD	PB-O1B	-2.22	1.43	1.50
4	B	3003	TYD	PB-O3B	2.22	1.63	1.54
4	A	3001	TYD	C2-N1	2.19	1.38	1.35
4	A	3001	TYD	PA-O2A	-2.18	1.45	1.55
4	C	3005	TYD	PB-O3B	2.18	1.63	1.54
4	C	3005	TYD	PB-O2B	-2.17	1.46	1.54
4	B	3004	TYD	PB-O3B	2.17	1.63	1.54
4	D	3007	TYD	PB-O2B	-2.16	1.46	1.54
4	A	3001	TYD	PB-O2B	-2.14	1.46	1.54
4	B	3003	TYD	PA-O2A	-2.10	1.45	1.55
4	D	3007	TYD	PA-O2A	-2.06	1.45	1.55
4	C	3005	TYD	O4-C4	-2.05	1.19	1.23
4	C	3005	TYD	PA-O2A	-2.04	1.45	1.55

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	3006	TYD	C5-C6-N1	6.92	124.74	111.11
4	C	3005	TYD	C5-C6-N1	6.77	124.45	111.11
4	B	3003	TYD	C5-C6-N1	6.75	124.41	111.11
4	B	3004	TYD	C5-C6-N1	6.73	124.37	111.11
4	A	3002	TYD	C5-C6-N1	6.70	124.30	111.11
4	D	3007	TYD	C5-C6-N1	6.69	124.30	111.11
4	A	3001	TYD	C5-C6-N1	6.61	124.12	111.11
4	D	3008	TYD	C5-C6-N1	6.52	123.94	111.11
4	B	3003	TYD	C5M-C5-C6	5.19	123.44	112.34
4	C	3005	TYD	C5M-C5-C6	5.11	123.25	112.34
4	D	3007	TYD	C5M-C5-C6	5.10	123.24	112.34
4	D	3008	TYD	C5M-C5-C6	5.04	123.11	112.34
4	A	3001	TYD	C5M-C5-C6	4.94	122.88	112.34
4	C	3006	TYD	C5M-C5-C6	4.93	122.86	112.34
4	B	3004	TYD	C5M-C5-C6	4.86	122.73	112.34
4	A	3002	TYD	C5M-C5-C6	4.26	121.44	112.34
4	C	3005	TYD	N3-C2-N1	-3.54	112.91	116.65
4	B	3003	TYD	N3-C2-N1	-3.45	113.01	116.65
4	A	3002	TYD	N3-C2-N1	-3.21	113.26	116.65
4	D	3007	TYD	N3-C2-N1	-3.20	113.27	116.65
4	C	3006	TYD	N3-C2-N1	-3.17	113.30	116.65
4	B	3004	TYD	N3-C2-N1	-3.13	113.34	116.65
4	D	3008	TYD	N3-C2-N1	-2.91	113.57	116.65
4	A	3001	TYD	N3-C2-N1	-2.62	113.88	116.65
4	D	3007	TYD	O2B-PB-O3A	2.47	112.93	104.64
4	B	3003	TYD	O4'-C1'-N1	-2.37	105.39	108.41
4	C	3006	TYD	O4'-C1'-N1	-2.30	105.48	108.41
4	A	3001	TYD	O5'-C5'-C4'	2.17	116.46	108.99
4	C	3005	TYD	O2B-PB-O3A	2.12	111.74	104.64
4	A	3001	TYD	O2B-PB-O3A	2.06	111.55	104.64
4	C	3006	TYD	O2B-PB-O3A	2.04	111.49	104.64

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	2009	GOL	C1-C2-C3-O3
4	C	3005	TYD	O4'-C4'-C5'-O5'
5	A	2009	GOL	O2-C2-C3-O3
4	C	3005	TYD	PB-O3A-PA-O5'
4	B	3003	TYD	PB-O3A-PA-O5'
4	A	3001	TYD	PB-O3A-PA-O5'
4	B	3004	TYD	PB-O3A-PA-O1A

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Mol	Chain	Res	Type	Atoms
4	C	3005	TYD	C3'-C4'-C5'-O5'
4	D	3007	TYD	PA-O3A-PB-O1B
4	A	3002	TYD	C5'-O5'-PA-O3A
4	B	3004	TYD	PB-O3A-PA-O2A

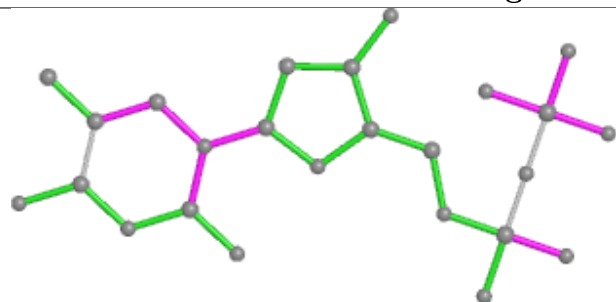
There are no ring outliers.

6 monomers are involved in 14 short contacts:

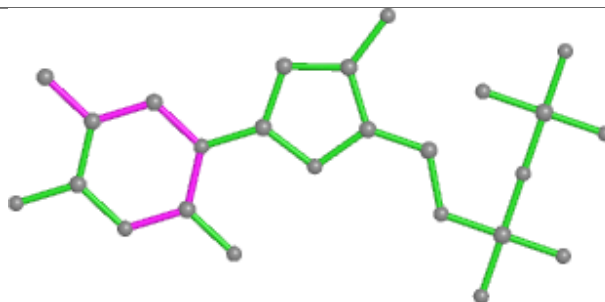
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	2011	GOL	1	0
4	B	3003	TYD	2	0
4	C	3005	TYD	2	0
5	D	2006	GOL	2	0
4	D	3007	TYD	3	0
4	A	3001	TYD	4	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

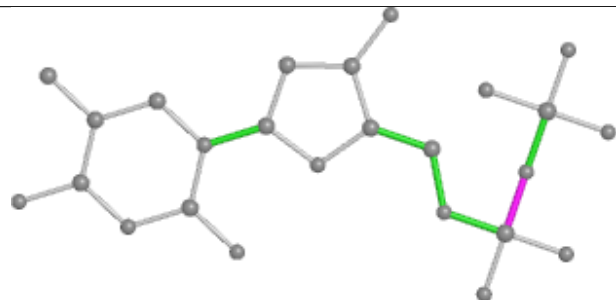
Ligand TYD B 3004



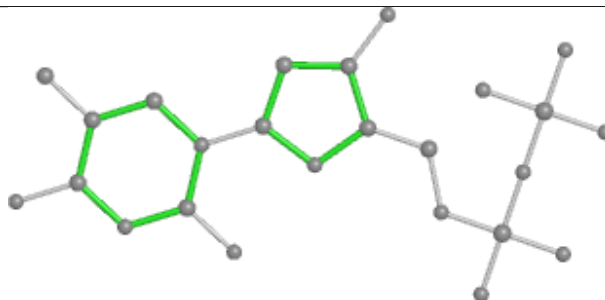
Bond lengths



Bond angles

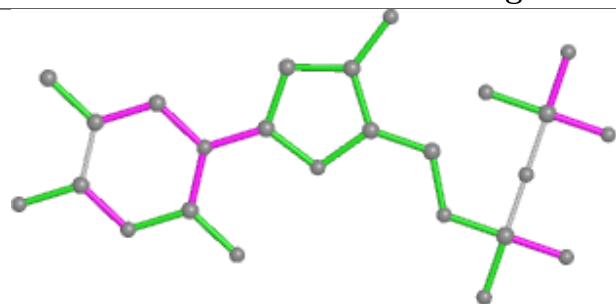


Torsions

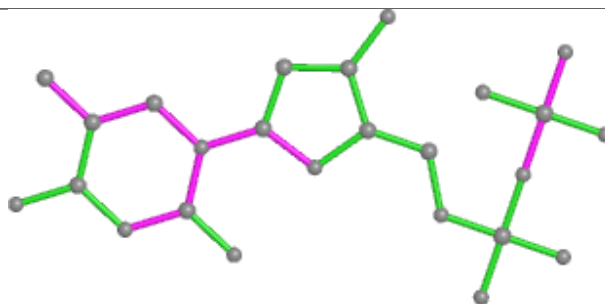


Rings

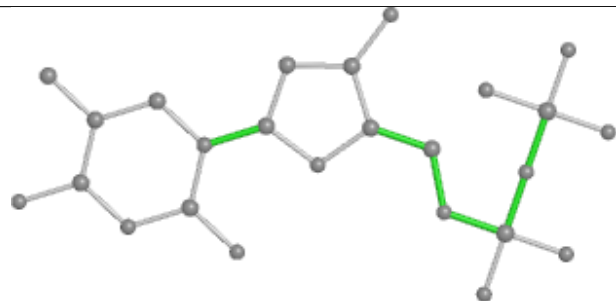
Ligand TYD C 3006



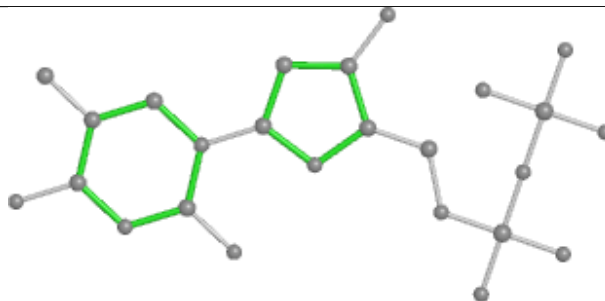
Bond lengths



Bond angles

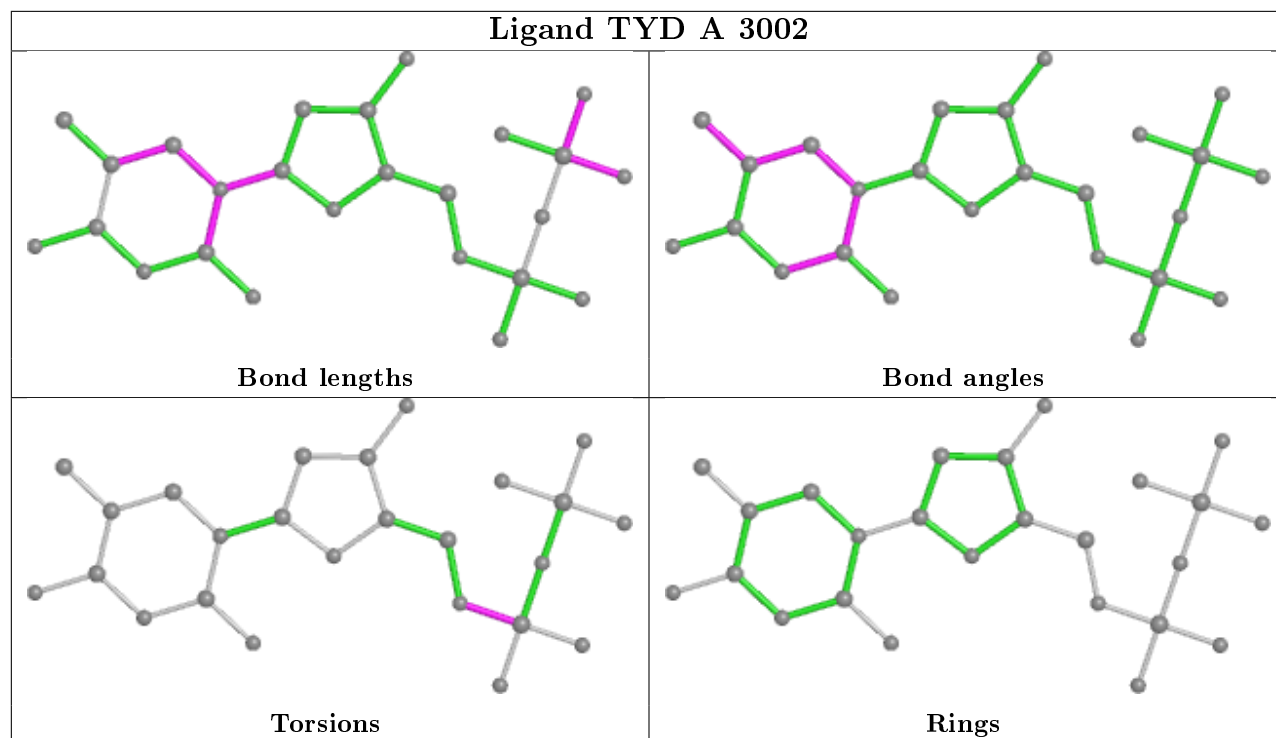


Torsions

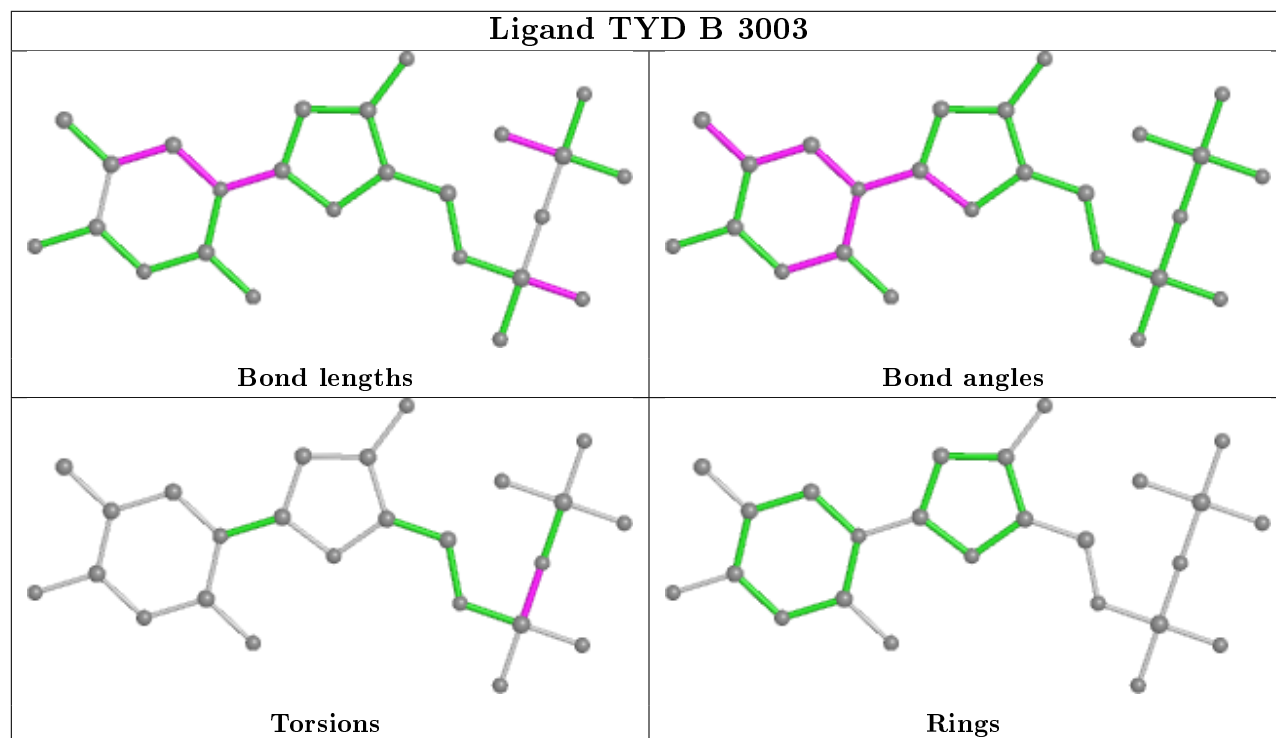


Rings

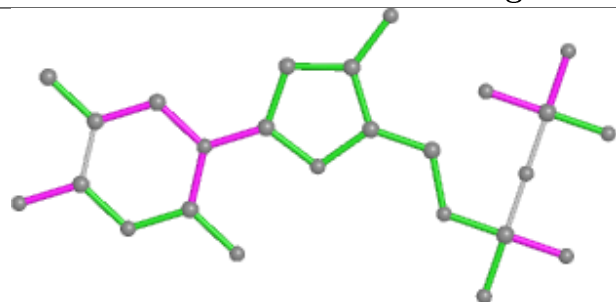
Ligand TYD A 3002



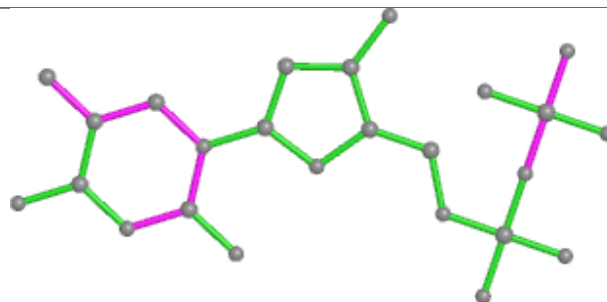
Ligand TYD B 3003



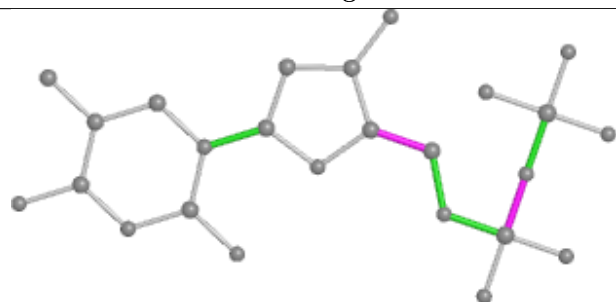
Ligand TYD C 3005



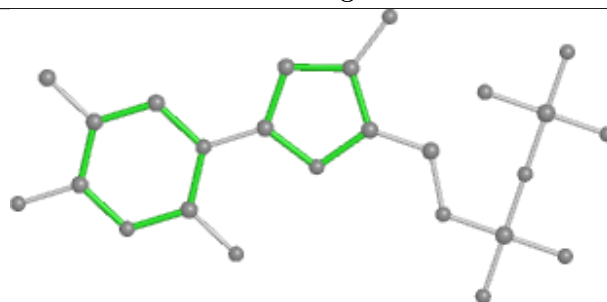
Bond lengths



Bond angles

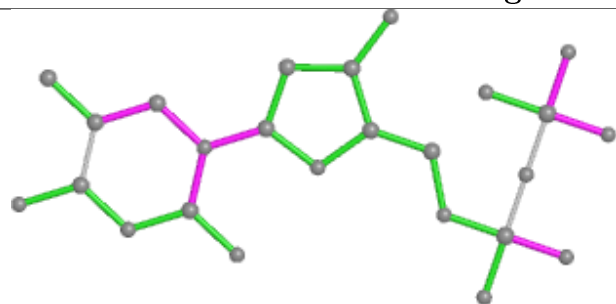


Torsions

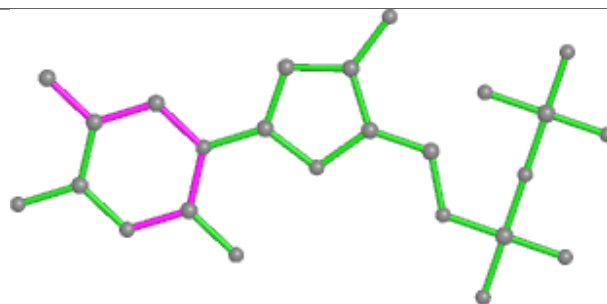


Rings

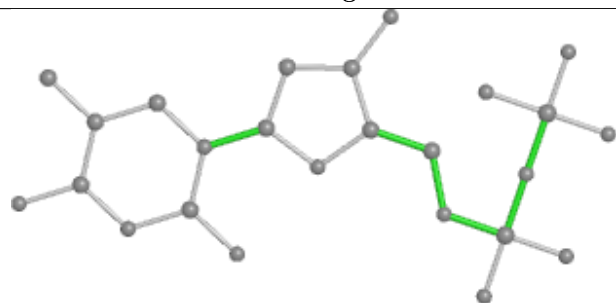
Ligand TYD D 3008



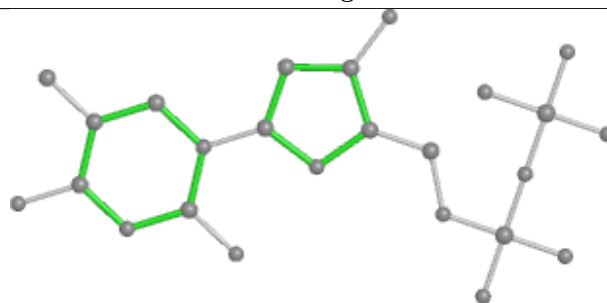
Bond lengths



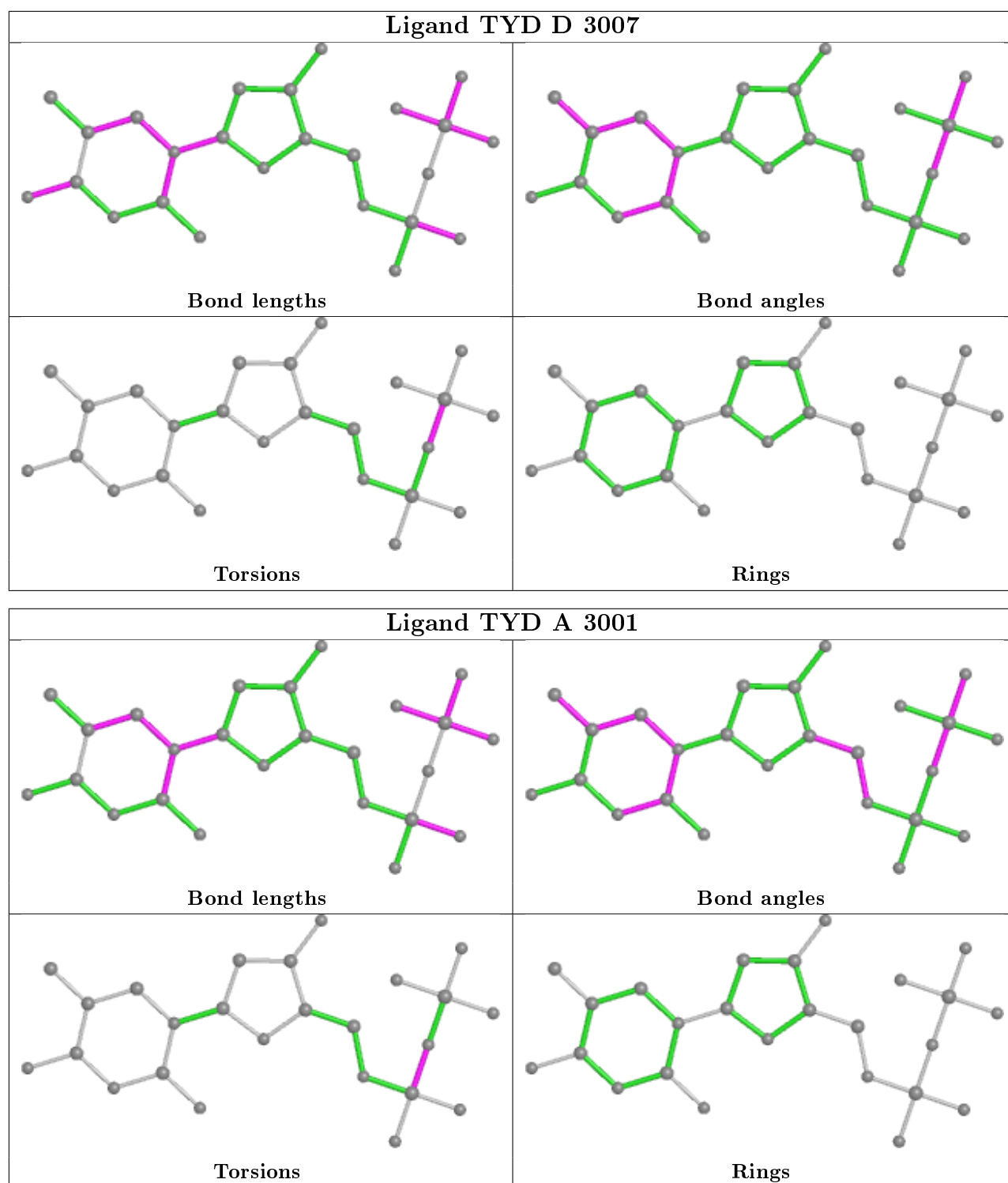
Bond angles



Torsions



Rings



5.7 Other polymers [i](#)

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 ⓘ

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, 95th 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 ²)	Q<0.9
1	A	295/295 (100%)	-0.10	8 (2%) 54 58	9, 15, 34, 50	0
1	B	292/295 (98%)	-0.05	6 (2%) 63 67	8, 16, 34, 65	0
1	C	292/295 (98%)	0.15	12 (4%) 37 41	9, 19, 40, 59	0
1	D	291/295 (98%)	0.47	30 (10%) 6 7	9, 24, 47, 60	0
All	All	1170/1180 (99%)	0.12	56 (4%) 30 34	8, 18, 40, 65	0

All (56) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	290	PHE	10.8
1	A	-2	GLY	10.6
1	B	291	HIS	7.4
1	C	-1	ALA	7.3
1	A	-1	ALA	7.0
1	B	289	ASN	6.9
1	C	192	ARG	6.5
1	D	192	ARG	6.0
1	C	191	ASP	5.5
1	C	193	GLY	4.3
1	D	193	GLY	4.3
1	D	191	ASP	4.0
1	D	163	ARG	3.7
1	A	11	GLY	3.3
1	D	186	ARG	3.3
1	A	0	HIS	3.3
1	D	206	ARG	3.2
1	B	191	ASP	3.2
1	C	206	ARG	3.2
1	D	271	GLU	3.1
1	D	105	VAL	3.1

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Mol	Chain	Res	Type	RSRZ
1	C	12	THR	2.9
1	D	12	THR	2.9
1	D	160	LYS	2.8
1	A	9	GLY	2.8
1	D	187	ILE	2.8
1	B	288	GLY	2.8
1	C	194	GLU	2.7
1	D	127	GLU	2.7
1	D	185	ARG	2.7
1	D	83	PRO	2.7
1	C	127	GLU	2.6
1	C	163	ARG	2.6
1	D	155	ILE	2.5
1	D	74	ARG	2.4
1	D	211	ARG	2.4
1	C	186	ARG	2.4
1	B	275	LYS	2.4
1	A	275	LYS	2.3
1	A	191	ASP	2.3
1	C	74	ARG	2.3
1	D	209	LYS	2.3
1	D	82	GLU	2.3
1	D	207	MET	2.3
1	A	290	PHE	2.3
1	D	104	LEU	2.2
1	D	6	LEU	2.2
1	D	195	LEU	2.2
1	D	197	ILE	2.2
1	D	190	SER	2.1
1	D	208	GLY	2.1
1	D	153	ARG	2.1
1	D	275	LYS	2.1
1	C	82	GLU	2.1
1	D	162	SER	2.1
1	D	272	LYS	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 ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

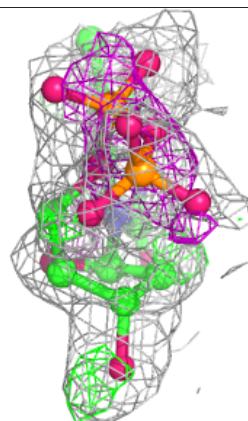
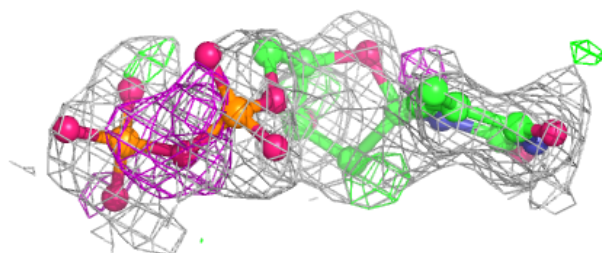
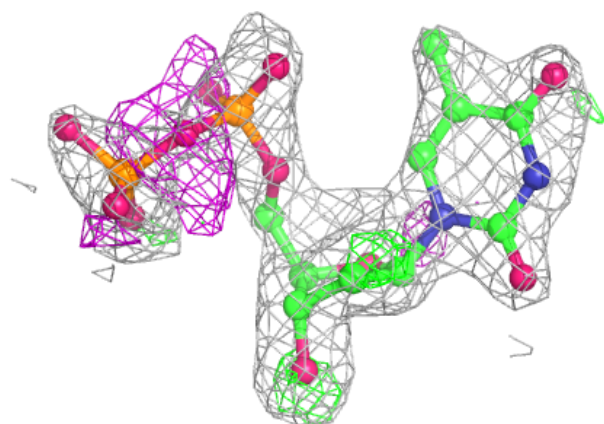
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	TYD	D	3007	25/25	0.65	0.22	36,48,57,57	0
5	GOL	B	2008	6/6	0.76	0.20	28,37,41,42	0
4	TYD	A	3001	25/25	0.83	0.15	21,30,54,55	0
4	TYD	C	3005	25/25	0.84	0.15	26,34,47,48	0
4	TYD	B	3003	25/25	0.85	0.13	18,25,51,53	0
5	GOL	A	2007	6/6	0.86	0.17	24,31,33,38	0
5	GOL	C	2003	6/6	0.87	0.14	17,24,28,32	0
5	GOL	B	2011	6/6	0.88	0.14	23,28,30,34	0
5	GOL	A	2009	6/6	0.88	0.14	22,31,33,35	0
5	GOL	D	2004	6/6	0.91	0.10	19,26,30,35	0
2	CL	D	4004	1/1	0.91	0.07	44,44,44,44	0
5	GOL	D	2006	6/6	0.91	0.15	20,29,31,32	0
3	SO4	A	5001	5/5	0.92	0.23	52,53,53,54	0
5	GOL	C	2010	6/6	0.92	0.10	18,20,22,23	0
5	GOL	A	2005	6/6	0.95	0.09	13,18,21,24	0
5	GOL	B	2002	6/6	0.95	0.11	15,19,22,27	0
3	SO4	D	5002	5/5	0.96	0.12	29,30,31,32	0
5	GOL	C	2001	6/6	0.97	0.07	17,22,24,27	0
4	TYD	D	3008	25/25	0.97	0.06	13,16,27,31	0
4	TYD	B	3004	25/25	0.97	0.08	11,13,32,34	0
2	CL	C	4003	1/1	0.98	0.05	26,26,26,26	0
4	TYD	A	3002	25/25	0.98	0.08	9,12,25,26	0
4	TYD	C	3006	25/25	0.98	0.06	10,12,27,30	0
2	CL	C	4005	1/1	0.99	0.04	24,24,24,24	0
2	CL	A	4001	1/1	1.00	0.04	17,17,17,17	0
2	CL	B	4002	1/1	1.00	0.04	16,16,16,16	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

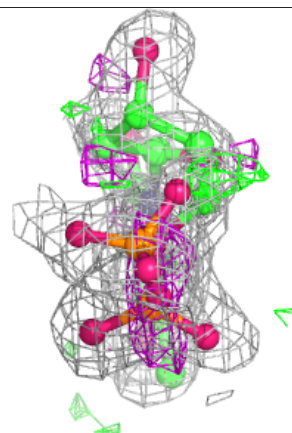
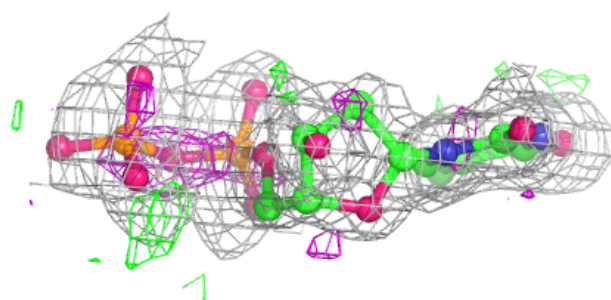
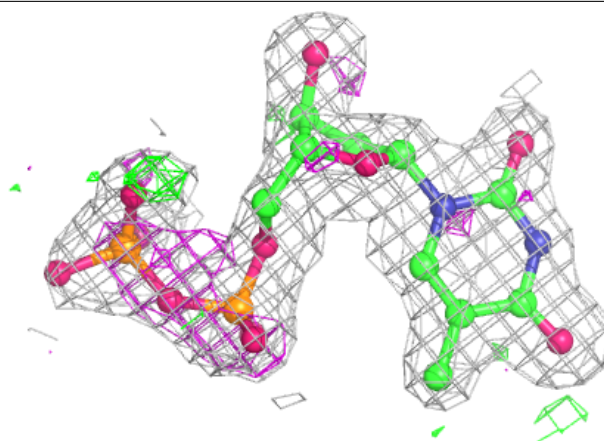
Electron density around TYD D 3007:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

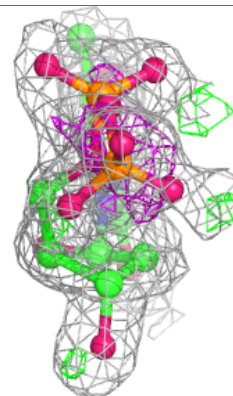
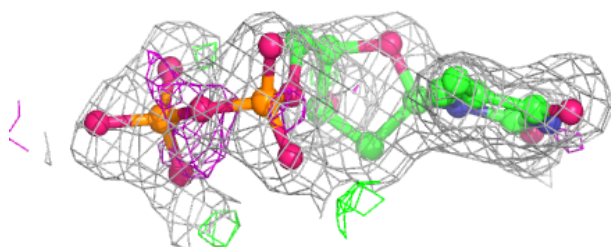
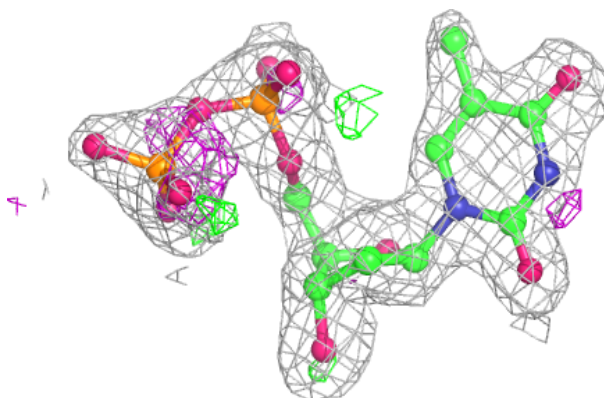


Electron density around TYD A 3001:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

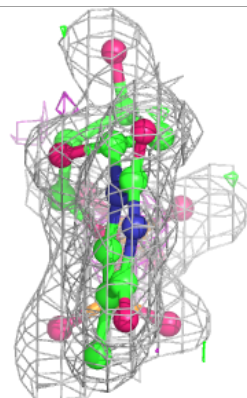
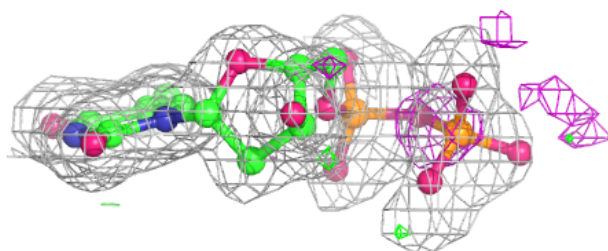
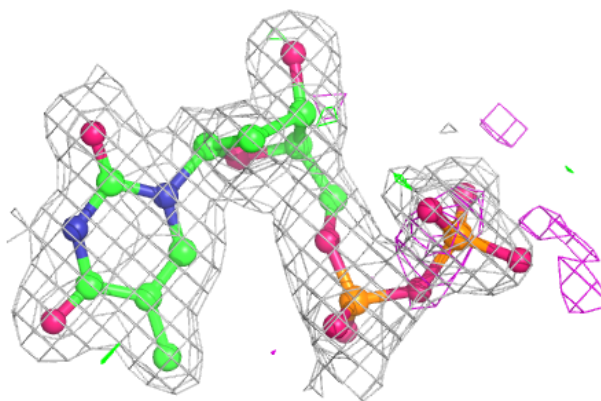
**Electron density around TYD C 3005:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



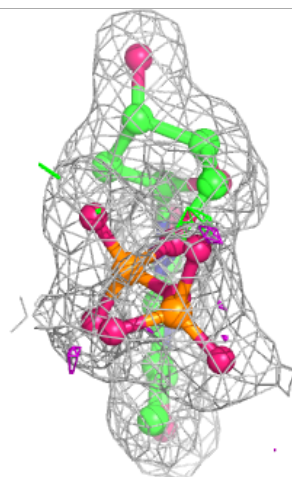
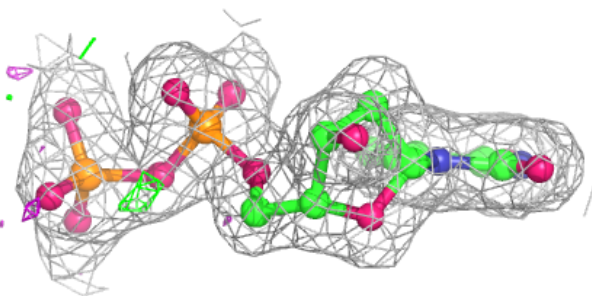
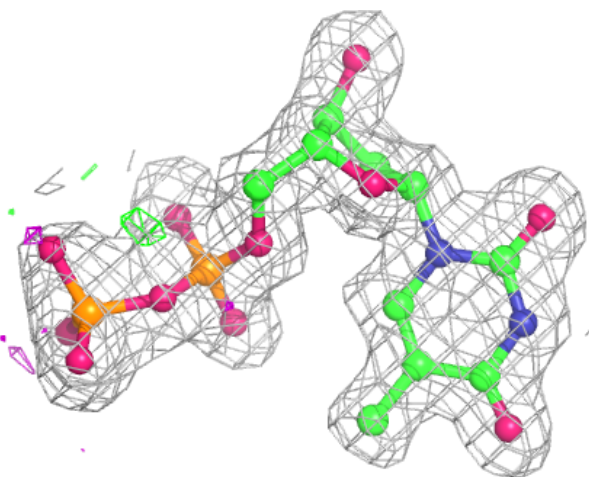
Electron density around TYD B 3003:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



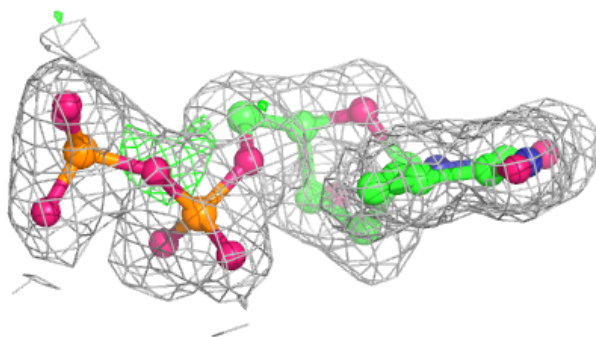
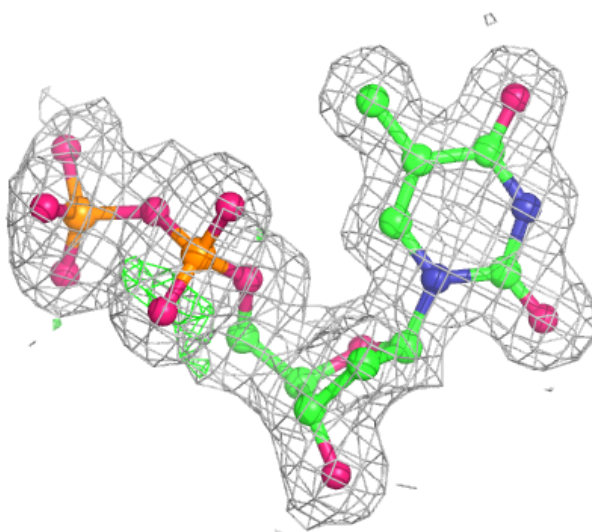
Electron density around TYD D 3008:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



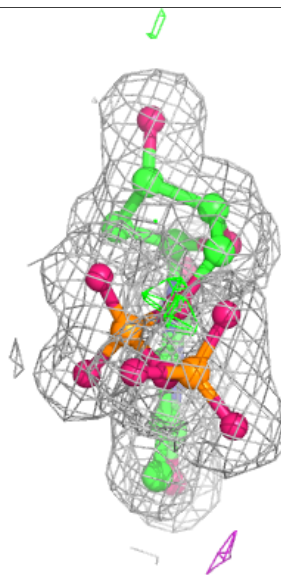
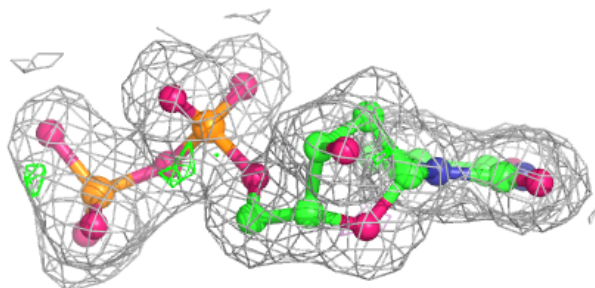
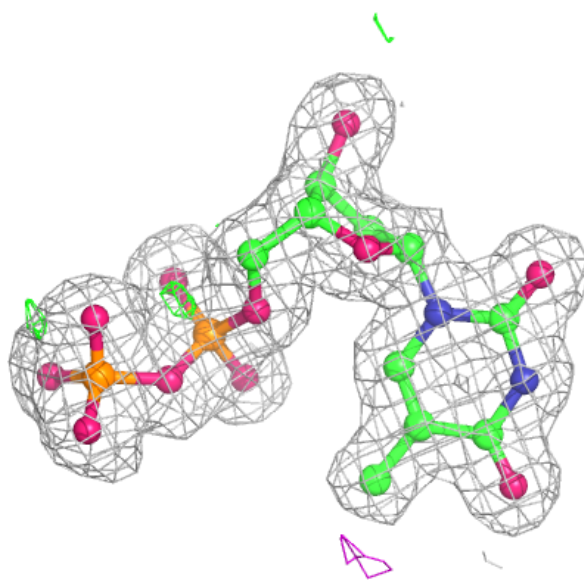
Electron density around TYD B 3004:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



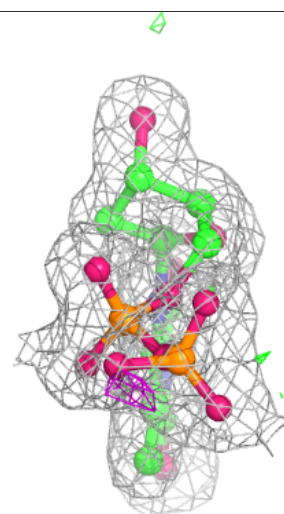
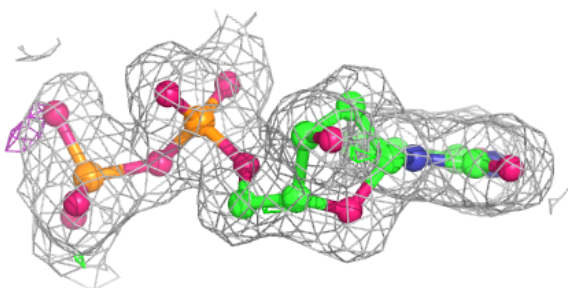
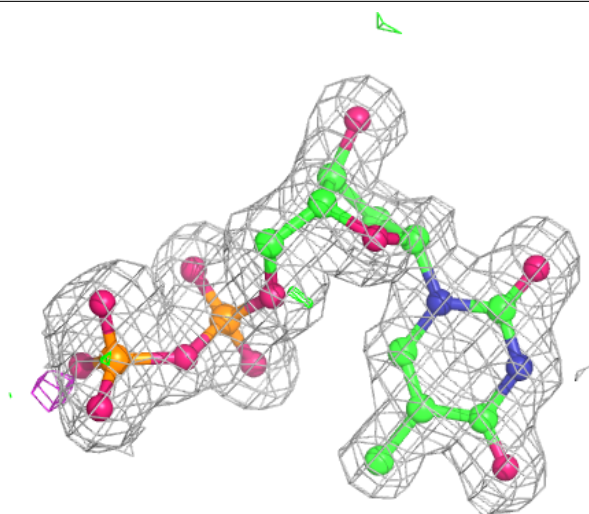
Electron density around TYD A 3002:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around TYD C 3006:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.