



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 09:53 AM UTC

PDB ID : 1RRV / pdb_00001rrv
Title : X-ray crystal structure of TDP-vancosaminyltransferase GtfD as a complex with TDP and the natural substrate, desvancosaminyl vancomycin.
Authors : Mulichak, A.M.; Lu, W.; Losey, H.C.; Walsh, C.T.; Garavito, R.M.
Deposited on : 2003-12-09
Resolution : 2.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

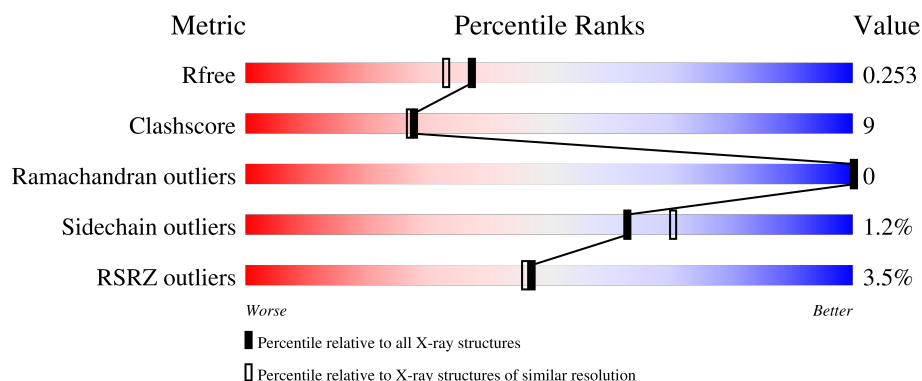
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of 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	416	2% 81% 15% .
1	B	416	4% 75% 20% . .
2	C	7	86% 14%
2	D	7	86% 14%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 6567 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 GLYCOSYLTRANSFERASE GTFD.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	401	Total	C	N	O	S	0	0	0
			2915	1848	519	536	12			
1	B	400	Total	C	N	O	S	0	0	0
			2900	1838	515	535	12			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	195	LEU	CYS	cloning artifact	UNP Q9AFC7
A	196	LEU	TRP	cloning artifact	UNP Q9AFC7
A	197	ALA	PRO	cloning artifact	UNP Q9AFC7
A	409	LEU	-	cloning artifact	UNP Q9AFC7
A	410	GLU	-	cloning artifact	UNP Q9AFC7
A	411	HIS	-	expression tag	UNP Q9AFC7
A	412	HIS	-	expression tag	UNP Q9AFC7
A	413	HIS	-	expression tag	UNP Q9AFC7
A	414	HIS	-	expression tag	UNP Q9AFC7
A	415	HIS	-	expression tag	UNP Q9AFC7
A	416	HIS	-	expression tag	UNP Q9AFC7
B	195	LEU	CYS	cloning artifact	UNP Q9AFC7
B	196	LEU	TRP	cloning artifact	UNP Q9AFC7
B	197	ALA	PRO	cloning artifact	UNP Q9AFC7
B	409	LEU	-	cloning artifact	UNP Q9AFC7
B	410	GLU	-	cloning artifact	UNP Q9AFC7
B	411	HIS	-	expression tag	UNP Q9AFC7
B	412	HIS	-	expression tag	UNP Q9AFC7
B	413	HIS	-	expression tag	UNP Q9AFC7
B	414	HIS	-	expression tag	UNP Q9AFC7
B	415	HIS	-	expression tag	UNP Q9AFC7
B	416	HIS	-	expression tag	UNP Q9AFC7

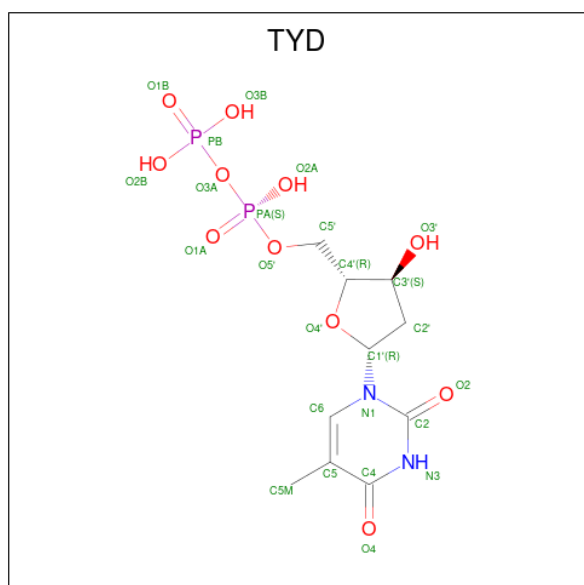
- Molecule 2 is a protein called DESVANCOSAMINYL VANCOMYCIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	7	Total 80	C 53	Cl 2	N 8	O 17	0	0	0
2	D	7	Total 80	C 53	Cl 2	N 8	O 17	0	0	0

- Molecule 3 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	K	0	0
			1	1		
3	B	1	Total	K	0	0
			1	1		

- Molecule 4 is THYMIDINE-5'-DIPHOSPHATE (CCD ID: TYD) (formula: C₁₀H₁₆N₂O₁₁P₂).



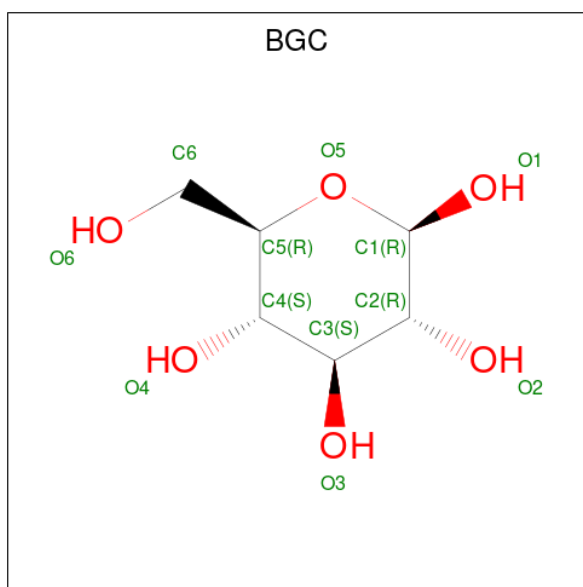
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total 25	C 10	N 2	O 11	P 2	0	0
4	B	1	Total 25	C 10	N 2	O 11	P 2	0	0

- Molecule 5 is GLYCEROL (CCD ID: 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	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 6 is beta-D-glucopyranose (CCD ID: BGC) (formula: $C_6H_{12}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	C	1	Total	C	O	0	0
			11	6	5		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	D	1	Total	C	O	0	0
			11	6	5		

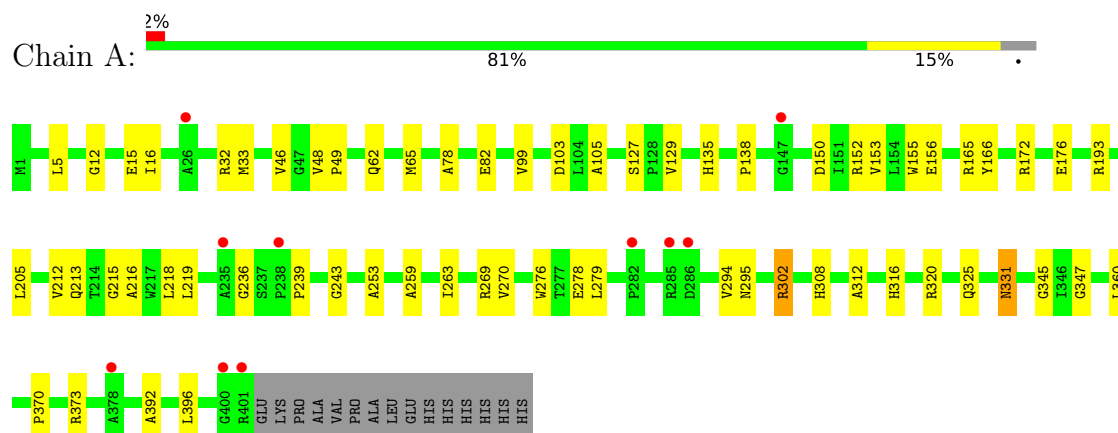
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	303	Total	O	0	0
			303	303		
7	B	177	Total	O	0	0
			177	177		
7	C	13	Total	O	0	0
			13	13		
7	D	7	Total	O	0	0
			7	7		

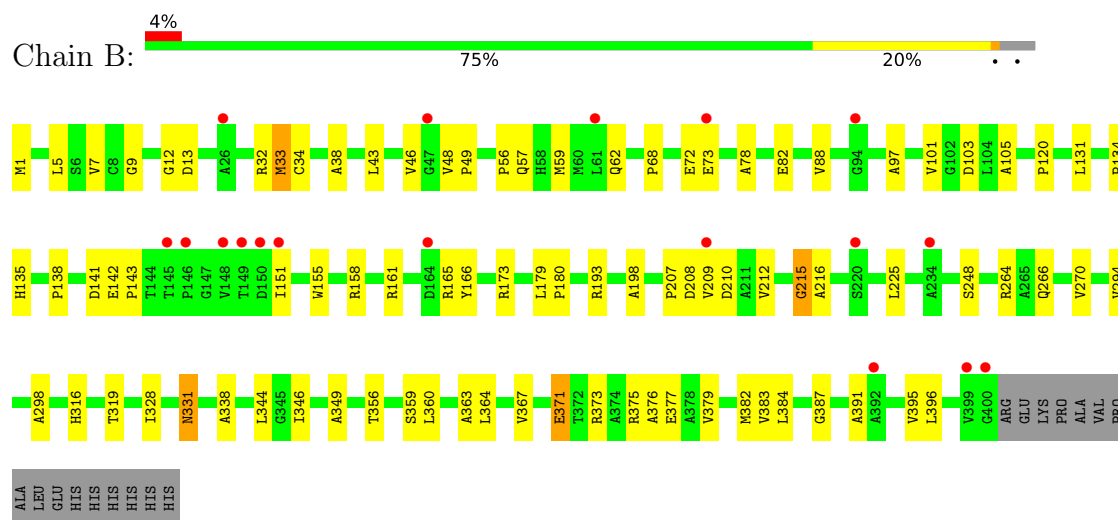
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

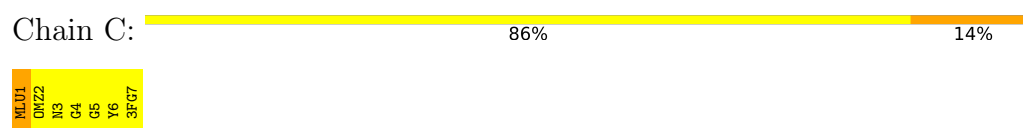
• Molecule 1: GLYCOSYLTRANSFERASE GTFD



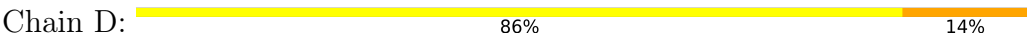
• Molecule 1: GLYCOSYLTRANSFERASE GTFD



• Molecule 2: DESVANCOSAMINYL VANCOMYCIN



• Molecule 2: DESVANCOSAMINYL VANCOMYCIN



HEL1
CR22
N3
C4
C5
Y6
3R67

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	50.67Å 64.12Å 144.08Å 90.00° 91.73° 90.00°	Depositor
Resolution (Å)	30.00 – 2.00 30.00 – 2.08	Depositor EDS
% Data completeness (in resolution range)	(Not available) (30.00-2.00) 99.6 (30.00-2.08)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.14 (at 2.08Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.210 , 0.252 0.212 , 0.253	Depositor DCC
R_{free} test set	2678 reflections (4.82%)	wwPDB-VP
Wilson B-factor (Å ²)	31.9	Xtriage
Anisotropy	0.477	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 47.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.029 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6567	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.76% 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: GHP, K, 3FG, GOL, MLU, OMY, TYD, BGC, OMZ

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	0.43	0/2985	0.93	5/4093 (0.1%)
1	B	0.38	0/2970	0.92	7/4075 (0.2%)
2	C	1.02	0/7	1.84	0/8
2	D	1.14	0/7	1.90	0/8
All	All	0.41	0/5969	0.93	12/8184 (0.1%)

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	215	GLY	N-CA-C	-9.33	102.53	112.04
1	B	215	GLY	N-CA-C	-6.37	105.90	111.95
1	B	88	VAL	CB-CA-C	-6.12	107.88	113.70
1	A	312	ALA	N-CA-C	5.99	118.30	111.11
1	B	328	ILE	CA-C-N	5.47	125.42	119.78
1	B	328	ILE	C-N-CA	5.47	125.42	119.78
1	A	135	HIS	N-CA-C	5.39	119.03	112.23
1	A	278	GLU	N-CA-C	5.30	117.65	111.02
1	B	33	MET	N-CA-C	5.17	116.85	108.26
1	B	34	CYS	N-CA-C	-5.16	100.76	108.96
1	B	270	VAL	N-CA-C	5.16	115.90	108.48
1	A	33	MET	N-CA-C	5.05	116.62	108.34

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	2915	0	2891	50	0
1	B	2900	0	2867	58	0
2	C	80	0	44	1	0
2	D	80	0	44	3	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	25	0	13	3	0
4	B	25	0	13	2	0
5	A	12	0	14	2	0
5	B	6	0	6	0	0
6	C	11	0	10	0	0
6	D	11	0	10	0	0
7	A	303	0	0	9	0
7	B	177	0	0	2	0
7	C	13	0	0	0	0
7	D	7	0	0	0	0
All	All	6567	0	5912	108	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 9.

All (108) 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:379:VAL:HA	1:B:382:MET:HE2	1.56	0.85
1:A:205:LEU:H	1:A:213:GLN:HE22	1.28	0.80
1:B:12:GLY:HA2	4:B:1004:TYD:H5'1	1.70	0.73
1:A:345:GLY:HA2	7:A:2274:HOH:O	1.91	0.70
1:B:166:TYR:OH	2:D:1:MLU:HCN3	1.93	0.69
1:B:319:THR:HG22	1:B:383:VAL:HG21	1.77	0.66
1:B:193:ARG:HD2	1:B:212:VAL:HG23	1.79	0.65
1:B:373:ARG:O	1:B:377:GLU:HG3	1.97	0.64
1:B:12:GLY:HA2	4:B:1004:TYD:C5'	2.28	0.64
1:B:346:ILE:HG22	1:B:376:ALA:HA	1.78	0.64
1:B:46:VAL:HG23	1:B:48:VAL:HG23	1.81	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:166:TYR:OH	2:C:1:MLU:HCN3	1.99	0.61
1:A:12:GLY:O	1:A:16:ILE:HG12	2.01	0.61
1:B:135:HIS:NE2	1:B:207:PRO:HG2	2.15	0.61
1:B:173:ARG:HD2	7:B:2067:HOH:O	2.00	0.60
1:A:276:TRP:HZ3	5:A:1005:GOL:H12	1.67	0.60
1:A:78:ALA:O	1:A:82:GLU:HG3	2.02	0.59
1:B:216:ALA:HB2	1:B:316:HIS:CD2	2.37	0.59
1:B:32:ARG:HA	1:B:49:PRO:HG2	1.85	0.59
1:A:12:GLY:HA2	4:A:1003:TYD:H5'1	1.85	0.59
1:A:12:GLY:HA2	4:A:1003:TYD:C5'	2.33	0.58
1:A:129:VAL:HG22	7:A:2120:HOH:O	2.03	0.58
1:B:379:VAL:HA	1:B:382:MET:CE	2.30	0.57
1:B:384:LEU:HD21	7:B:2144:HOH:O	2.05	0.56
1:A:127:SER:HB2	7:A:2120:HOH:O	2.05	0.56
1:A:331:ASN:HD22	1:A:331:ASN:C	2.14	0.55
1:A:46:VAL:HG23	1:A:48:VAL:HG23	1.87	0.55
1:A:219:LEU:C	1:A:219:LEU:HD12	2.32	0.55
1:A:32:ARG:HA	1:A:49:PRO:HG2	1.87	0.55
1:A:259:ALA:HA	1:A:360:LEU:CD2	2.38	0.54
1:A:150:ASP:OD2	1:A:153:VAL:HG23	2.08	0.54
1:A:259:ALA:HA	1:A:360:LEU:HD21	1.91	0.53
1:B:266:GLN:HG3	1:B:364:LEU:HD23	1.92	0.52
1:A:172:ARG:O	1:A:176:GLU:HG3	2.09	0.52
1:B:1:MET:HE1	1:B:396:LEU:CD2	2.39	0.52
1:B:43:LEU:HD22	1:B:48:VAL:HB	1.92	0.52
1:A:152:ARG:O	1:A:156:GLU:HG3	2.11	0.50
1:A:205:LEU:H	1:A:213:GLN:NE2	2.03	0.50
1:B:135:HIS:CD2	1:B:207:PRO:HG2	2.46	0.50
1:A:138:PRO:HG3	1:A:155:TRP:CD2	2.46	0.50
1:B:338:ALA:O	1:B:349:ALA:HB2	2.12	0.50
1:B:56:PRO:HG2	1:B:59:MET:HG2	1.94	0.50
1:B:165:ARG:HD3	2:D:1:MLU:HCN1	1.94	0.50
1:B:193:ARG:NH1	1:B:210:ASP:OD2	2.45	0.49
1:A:16:ILE:HD11	7:A:2256:HOH:O	2.12	0.49
1:B:207:PRO:O	1:B:208:ASP:HB2	2.12	0.49
1:A:193:ARG:HD2	1:A:212:VAL:HG23	1.95	0.48
1:B:331:ASN:HD22	1:B:331:ASN:C	2.21	0.48
1:A:392:ALA:O	1:A:396:LEU:HG	2.14	0.48
1:A:236:GLY:HA3	1:A:269:ARG:CZ	2.44	0.48
1:A:219:LEU:HD12	1:A:219:LEU:O	2.14	0.47
1:B:383:VAL:HG23	1:B:383:VAL:O	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:302:ARG:HB3	1:A:302:ARG:HH11	1.79	0.47
1:B:68:PRO:HG2	1:B:73:GLU:HG2	1.97	0.47
1:B:62:GLN:HG2	1:B:248:SER:HB2	1.95	0.47
1:B:225:LEU:HD21	1:B:298:ALA:HB3	1.97	0.47
1:A:62:GLN:HB2	1:A:65:MET:HE3	1.97	0.47
1:B:225:LEU:HD22	1:B:294:VAL:CG1	2.45	0.47
1:A:103:ASP:OD1	1:A:105:ALA:HB3	2.14	0.47
1:A:263:ILE:HG21	1:A:270:VAL:HG22	1.96	0.47
1:B:193:ARG:HD2	1:B:212:VAL:CG2	2.46	0.46
1:B:141:ASP:OD1	1:B:142:GLU:HG2	2.15	0.46
1:A:243:GLY:O	1:A:308:HIS:HA	2.16	0.46
1:A:138:PRO:HG3	1:A:155:TRP:CE2	2.51	0.45
1:A:259:ALA:O	1:A:263:ILE:HG13	2.15	0.45
1:A:325:GLN:O	1:A:347:GLY:HA3	2.16	0.45
1:B:78:ALA:O	1:B:82:GLU:HG3	2.16	0.45
1:B:143:PRO:HD3	2:D:7:3FG:O	2.17	0.45
1:B:5:LEU:O	1:B:33:MET:HA	2.17	0.45
1:A:218:LEU:HD23	7:A:2208:HOH:O	2.17	0.45
1:A:370:PRO:HG2	7:A:2290:HOH:O	2.18	0.44
1:B:68:PRO:HB2	1:B:72:GLU:HB3	1.99	0.44
1:B:356:THR:H	1:B:359:SER:HG	1.63	0.44
1:A:216:ALA:HB2	1:A:316:HIS:CD2	2.52	0.44
1:B:9:GLY:HA3	1:B:13:ASP:OD2	2.18	0.44
1:B:131:LEU:HD13	1:B:209:VAL:HG11	2.00	0.44
1:A:239:PRO:HB3	1:A:269:ARG:HB2	2.00	0.44
1:A:294:VAL:CG1	1:A:295:ASN:N	2.81	0.44
1:A:15:GLU:HG3	7:A:2208:HOH:O	2.19	0.43
7:A:2288:HOH:O	1:B:264:ARG:HB3	2.19	0.43
1:B:371:GLU:H	1:B:371:GLU:CD	2.25	0.43
1:B:138:PRO:HG3	1:B:155:TRP:CE2	2.54	0.43
1:A:253:ALA:HA	1:A:279:LEU:HD22	2.00	0.43
1:A:320:ARG:HD3	7:A:2253:HOH:O	2.18	0.43
1:B:344:LEU:O	1:B:375:ARG:NH1	2.52	0.43
1:B:193:ARG:HH11	1:B:193:ARG:HG3	1.83	0.43
1:A:5:LEU:HD23	1:A:99:VAL:HB	2.01	0.43
1:B:179:LEU:HA	1:B:180:PRO:HD3	1.88	0.42
1:B:198:ALA:O	1:B:215:GLY:HA2	2.19	0.42
1:B:103:ASP:OD1	1:B:105:ALA:HB3	2.20	0.42
1:B:158:ARG:HA	1:B:161:ARG:NH2	2.34	0.42
1:A:316:HIS:O	1:A:320:ARG:HG2	2.19	0.42
1:A:12:GLY:HA2	4:A:1003:TYD:H5'2	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:331:ASN:C	1:A:331:ASN:ND2	2.77	0.42
1:B:97:ALA:CB	1:B:120:PRO:HG2	2.49	0.41
1:B:391:ALA:O	1:B:395:VAL:HG23	2.19	0.41
1:B:363:ALA:O	1:B:367:VAL:HG23	2.20	0.41
1:B:138:PRO:HG3	1:B:155:TRP:CD2	2.55	0.41
1:B:215:GLY:O	1:B:387:GLY:HA3	2.21	0.41
1:B:360:LEU:HD12	1:B:360:LEU:HA	1.86	0.41
1:A:294:VAL:HG12	1:A:295:ASN:N	2.35	0.41
1:B:7:VAL:HG22	1:B:101:VAL:HB	2.01	0.41
1:B:38:ALA:HB2	1:B:57:GLN:HB2	2.03	0.41
1:B:375:ARG:HA	1:B:375:ARG:HD2	1.85	0.41
1:A:165:ARG:HH11	1:A:165:ARG:HG3	1.86	0.41
1:A:276:TRP:CZ3	5:A:1005:GOL:H12	2.52	0.40
1:A:219:LEU:C	1:A:219:LEU:CD1	2.94	0.40
1:B:134:PRO:O	1:B:151:ILE:HD12	2.22	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	399/416 (96%)	390 (98%)	9 (2%)	0	100	100
1	B	398/416 (96%)	384 (96%)	14 (4%)	0	100	100
2	C	1/7 (14%)	1 (100%)	0	0	100	100
2	D	1/7 (14%)	1 (100%)	0	0	100	100
All	All	799/846 (94%)	776 (97%)	23 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	289/314 (92%)	286 (99%)	3 (1%)	68	75
1	B	287/314 (91%)	285 (99%)	2 (1%)	76	82
2	C	1/1 (100%)	0	1 (100%)	0	0
2	D	1/1 (100%)	0	1 (100%)	0	0
All	All	578/630 (92%)	571 (99%)	7 (1%)	63	70

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	302	ARG
1	A	331	ASN
1	A	373	ARG
1	B	331	ASN
1	B	371	GLU
2	C	3	ASN
2	D	3	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	30	GLN
1	A	213	GLN
1	A	266	GLN
1	A	331	ASN
1	B	266	GLN
1	B	331	ASN
2	C	3	ASN
2	D	3	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

12 non-standard protein/DNA/RNA residues are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
2	OMY	D	6	2	12,14,15	2.33	5 (41%)	17,19,21	1.78	2 (11%)
2	OMZ	C	2	2	12,14,15	1.54	4 (33%)	17,19,21	0.95	1 (5%)
2	3FG	C	7	2	12,13,13	2.72	7 (58%)	14,18,18	1.24	1 (7%)
2	MLU	C	1	2	7,8,9	1.01	1 (14%)	7,9,11	1.18	1 (14%)
2	GHP	D	4	2,6	10,11,12	1.84	4 (40%)	11,14,16	1.16	1 (9%)
2	GHP	C	4	2,6	10,11,12	1.74	2 (20%)	11,14,16	1.22	2 (18%)
2	MLU	D	1	2	7,8,9	0.93	0	7,9,11	1.05	0
2	OMZ	D	2	2	12,14,15	1.76	4 (33%)	17,19,21	1.00	1 (5%)
2	OMY	C	6	2	12,14,15	1.99	3 (25%)	17,19,21	1.73	2 (11%)
2	3FG	D	7	2	12,13,13	3.12	8 (66%)	14,18,18	1.31	1 (7%)
2	GHP	C	5	2	10,11,12	1.80	3 (30%)	11,14,16	1.91	5 (45%)
2	GHP	D	5	2	10,11,12	2.09	2 (20%)	11,14,16	1.91	5 (45%)

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	OMY	D	6	2	-	1/9/10/12	0/1/1/1
2	OMZ	C	2	2	-	1/9/10/12	0/1/1/1
2	3FG	C	7	2	-	5/8/8/8	0/1/1/1
2	MLU	C	1	2	-	3/5/8/10	-
2	GHP	D	4	2,6	-	0/4/6/8	0/1/1/1
2	GHP	C	4	2,6	-	0/4/6/8	0/1/1/1
2	MLU	D	1	2	-	3/5/8/10	-
2	OMZ	D	2	2	-	1/9/10/12	0/1/1/1
2	OMY	C	6	2	-	1/9/10/12	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	3FG	D	7	2	-	5/8/8/8	0/1/1/1
2	GHP	C	5	2	-	0/4/6/8	0/1/1/1
2	GHP	D	5	2	-	0/4/6/8	0/1/1/1

All (43) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	7	3FG	CG1-CB	5.93	1.47	1.39
2	C	7	3FG	CG1-CB	5.39	1.47	1.39
2	D	6	OMY	CZ-CE1	5.18	1.44	1.39
2	D	7	3FG	CB-CA	5.15	1.58	1.52
2	D	5	GHP	CA-C	4.78	1.59	1.51
2	C	7	3FG	CB-CA	4.36	1.57	1.52
2	C	6	OMY	CZ-CE1	4.36	1.43	1.39
2	C	4	GHP	C3-C2	3.77	1.44	1.38
2	C	5	GHP	CA-C	3.73	1.57	1.51
2	D	7	3FG	CZ-CD1	3.29	1.44	1.39
2	D	4	GHP	C3-C2	3.25	1.44	1.38
2	C	7	3FG	OXT-C	-3.14	1.20	1.30
2	D	7	3FG	CG1-CD1	3.10	1.43	1.39
2	D	7	3FG	OXT-C	-3.09	1.20	1.30
2	D	7	3FG	CG2-CD2	3.09	1.43	1.39
2	D	2	OMZ	OH-CZ	3.05	1.42	1.36
2	D	2	OMZ	CE2-CZ	3.02	1.44	1.39
2	D	6	OMY	CD2-CG	2.90	1.43	1.39
2	C	7	3FG	CZ-CD1	2.76	1.43	1.39
2	C	5	GHP	C3-C4	2.73	1.44	1.39
2	C	2	OMZ	CD1-CG	2.72	1.43	1.39
2	D	4	GHP	C5-C4	2.72	1.44	1.39
2	D	6	OMY	OCZ-CZ	2.71	1.41	1.36
2	D	2	OMZ	CD2-CG	2.69	1.43	1.39
2	C	6	OMY	CD2-CG	2.62	1.43	1.39
2	D	5	GHP	C3-C4	2.61	1.43	1.39
2	D	2	OMZ	CD1-CG	2.56	1.42	1.39
2	C	2	OMZ	CE2-CZ	2.55	1.44	1.39
2	D	4	GHP	O4-C4	2.46	1.42	1.37
2	D	7	3FG	CG2-CB	2.45	1.42	1.39
2	C	2	OMZ	CD2-CG	2.44	1.43	1.39
2	C	6	OMY	ODE-CB	-2.39	1.38	1.42
2	C	7	3FG	CG2-CD2	2.38	1.42	1.39
2	C	1	MLU	CA-N	2.37	1.51	1.47
2	C	7	3FG	CG1-CD1	2.33	1.42	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	2	OMZ	OH-CZ	2.30	1.41	1.36
2	D	6	OMY	CD1-CG	2.26	1.42	1.39
2	C	4	GHP	C1-CA	2.13	1.54	1.52
2	D	6	OMY	CD1-CE1	2.10	1.42	1.38
2	C	5	GHP	C6-C1	2.05	1.42	1.39
2	D	4	GHP	C1-CA	2.02	1.54	1.52
2	D	7	3FG	CA-C	2.01	1.59	1.54
2	C	7	3FG	CA-C	2.00	1.59	1.54

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	6	OMY	CG-CB-CA	-6.23	103.40	111.58
2	C	6	OMY	CG-CB-CA	-6.20	103.44	111.58
2	C	5	GHP	C2-C1-CA	-3.30	115.51	120.64
2	D	5	GHP	C2-C1-CA	-2.82	116.25	120.64
2	D	5	GHP	C2-C3-C4	-2.63	117.10	119.88
2	D	5	GHP	C1-CA-C	-2.55	104.93	112.09
2	D	5	GHP	C5-C6-C1	-2.53	118.66	121.18
2	C	5	GHP	C5-C6-C1	-2.51	118.68	121.18
2	C	5	GHP	C6-C1-CA	2.49	124.52	120.64
2	D	4	GHP	C5-C6-C1	2.41	123.59	121.18
2	C	7	3FG	CB-CA-C	2.41	115.32	110.10
2	C	4	GHP	C5-C6-C1	2.38	123.56	121.18
2	D	2	OMZ	O-C-CA	-2.38	118.66	124.77
2	C	2	OMZ	O-C-CA	-2.36	118.70	124.77
2	D	6	OMY	OCZ-CZ-CE1	2.32	124.84	119.25
2	D	7	3FG	CB-CA-C	2.29	115.05	110.10
2	C	5	GHP	C2-C3-C4	-2.22	117.53	119.88
2	C	4	GHP	C2-C1-CA	2.14	123.97	120.64
2	C	5	GHP	C1-CA-C	-2.06	106.32	112.09
2	D	5	GHP	C6-C1-CA	2.05	123.83	120.64
2	C	6	OMY	OCZ-CZ-CE1	2.03	124.14	119.25
2	C	1	MLU	CB-CA-C	-2.00	107.91	110.99

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	6	OMY	O-C-CA-CB
2	D	6	OMY	O-C-CA-CB
2	C	7	3FG	C-CA-CB-CG1

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Mol	Chain	Res	Type	Atoms
2	C	7	3FG	C-CA-CB-CG2
2	C	7	3FG	OXT-C-CA-N
2	D	7	3FG	C-CA-CB-CG1
2	D	7	3FG	C-CA-CB-CG2
2	D	7	3FG	OXT-C-CA-N
2	D	1	MLU	N-CA-CB-CG
2	C	1	MLU	CA-CB-CG-CD1
2	C	1	MLU	CA-CB-CG-CD2
2	D	1	MLU	CA-CB-CG-CD1
2	D	7	3FG	O-C-CA-CB
2	D	7	3FG	OXT-C-CA-CB
2	C	1	MLU	N-CA-CB-CG
2	D	1	MLU	CA-CB-CG-CD2
2	C	7	3FG	O-C-CA-CB
2	C	7	3FG	OXT-C-CA-CB
2	C	2	OMZ	O-C-CA-CB
2	D	2	OMZ	O-C-CA-CB

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	1	MLU	1	0
2	D	1	MLU	2	0
2	D	7	3FG	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 2 are monoatomic - leaving 7 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
6	BGC	C	8	2	11,11,12	1.16	0	15,15,17	1.65	2 (13%)
6	BGC	D	8	2	11,11,12	1.17	0	15,15,17	1.66	2 (13%)
4	TYD	A	1003	-	25,26,26	0.87	1 (4%)	38,40,40	0.66	0
5	GOL	A	1005	-	5,5,5	0.30	0	5,5,5	0.18	0
4	TYD	B	1004	-	25,26,26	0.83	1 (4%)	38,40,40	0.63	0
5	GOL	A	1006	3	5,5,5	0.37	0	5,5,5	0.32	0
5	GOL	B	1007	3	5,5,5	0.30	0	5,5,5	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
6	BGC	C	8	2	-	2/2/19/22	0/1/1/1
6	BGC	D	8	2	-	0/2/19/22	0/1/1/1
4	TYD	A	1003	-	-	0/16/28/28	0/2/2/2
5	GOL	A	1005	-	-	0/4/4/4	-
4	TYD	B	1004	-	-	0/16/28/28	0/2/2/2
5	GOL	A	1006	3	-	0/4/4/4	-
5	GOL	B	1007	3	-	0/4/4/4	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1003	TYD	PA-O3A	3.54	1.63	1.59
4	B	1004	TYD	PA-O3A	3.39	1.63	1.59

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	D	8	BGC	O6-C6-C5	4.95	128.18	111.33
6	C	8	BGC	O6-C6-C5	4.79	127.64	111.33
6	C	8	BGC	C2-C3-C4	-2.28	106.85	110.86
6	D	8	BGC	C2-C3-C4	-2.14	107.10	110.86

There are no chirality outliers.

All (2) torsion outliers are listed below:

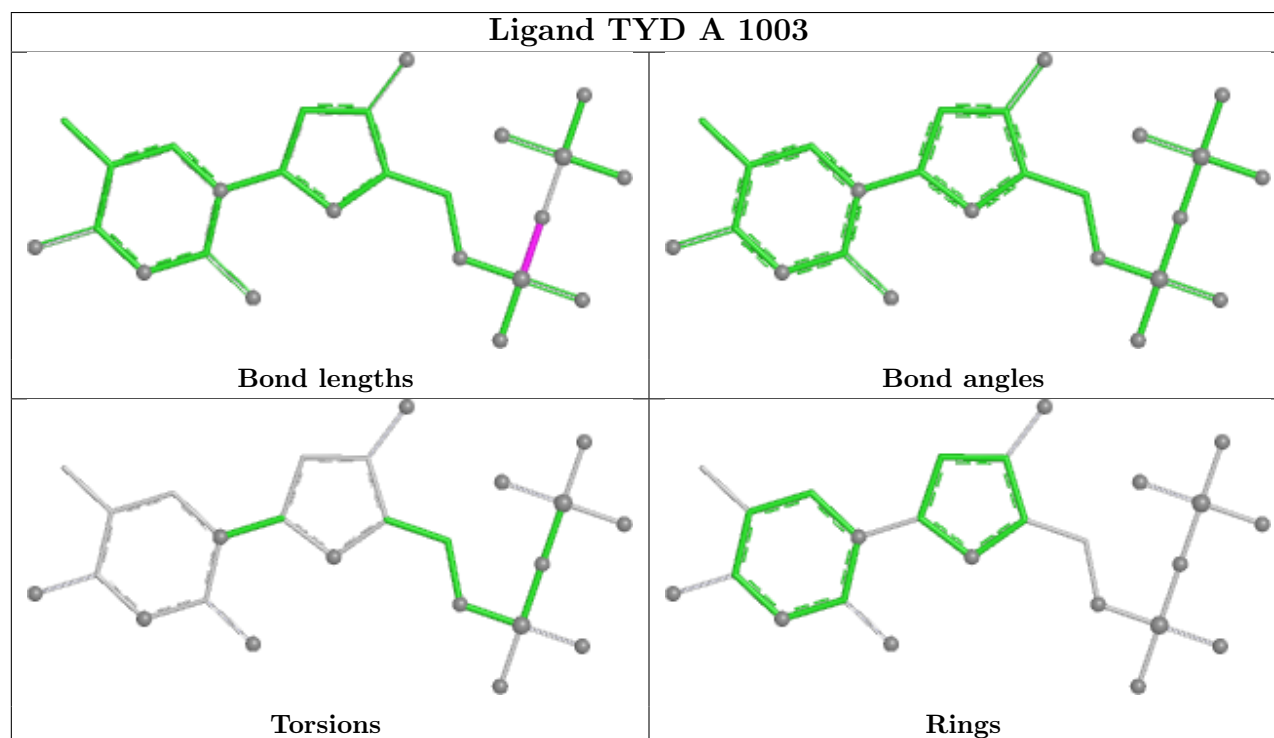
Mol	Chain	Res	Type	Atoms
6	C	8	BGC	C4-C5-C6-O6
6	C	8	BGC	O5-C5-C6-O6

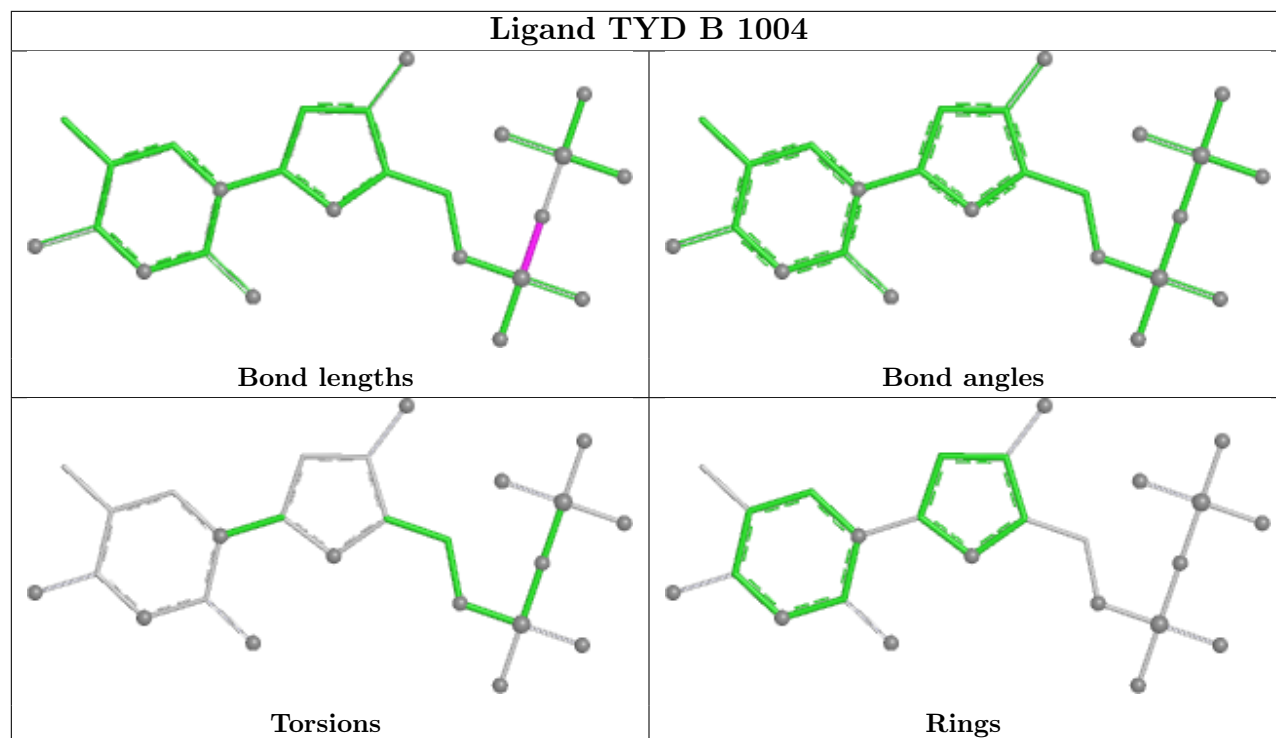
There are no ring outliers.

3 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1003	TYD	3	0
5	A	1005	GOL	2	0
4	B	1004	TYD	2	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.





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, 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	401/416 (96%)	0.18	10 (2%) 58 58	18, 31, 52, 60	0
1	B	400/416 (96%)	0.67	18 (4%) 38 37	29, 40, 58, 70	0
2	C	1/7 (14%)	-0.76	0 100 100	24, 24, 24, 24	0
2	D	1/7 (14%)	1.18	0 100 100	37, 37, 37, 37	0
All	All	803/846 (94%)	0.42	28 (3%) 47 46	18, 37, 55, 70	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	401	ARG	3.5
1	A	400	GLY	3.5
1	B	400	GLY	3.4
1	B	151	ILE	2.7
1	A	378	ALA	2.7
1	B	73	GLU	2.5
1	B	399	VAL	2.4
1	A	235	ALA	2.4
1	B	26	ALA	2.4
1	B	150	ASP	2.4
1	B	149	THR	2.4
1	B	220	SER	2.3
1	B	94	GLY	2.3
1	B	145	THR	2.3
1	B	47	GLY	2.3
1	B	209	VAL	2.3
1	B	146	PRO	2.3
1	A	285	ARG	2.2
1	A	286	ASP	2.2
1	B	234	ALA	2.2
1	A	238	PRO	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	147	GLY	2.2
1	A	282	PRO	2.1
1	B	164	ASP	2.1
1	B	148	VAL	2.1
1	B	61	LEU	2.1
1	A	26	ALA	2.1
1	B	392	ALA	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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. 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(Å ²)	Q<0.9
2	OMZ	D	2	14/15	0.58	0.16	35,37,39,40	0
2	OMY	D	6	14/15	0.60	0.16	31,34,40,40	0
2	OMY	C	6	14/15	0.61	0.14	18,23,26,30	0
2	OMZ	C	2	14/15	0.73	0.12	18,22,26,27	0
2	3FG	D	7	13/13	0.79	0.13	42,46,53,54	0
2	MLU	D	1	9/10	0.84	0.16	44,46,51,53	0
2	3FG	C	7	13/13	0.88	0.11	27,29,41,41	0
2	GHP	D	4	11/12	0.90	0.10	31,33,36,38	0
2	GHP	D	5	11/12	0.90	0.12	37,41,43,44	0
2	MLU	C	1	9/10	0.90	0.12	24,38,40,40	0
2	GHP	C	5	11/12	0.94	0.08	17,23,27,27	0
2	GHP	C	4	11/12	0.96	0.07	19,22,25,27	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

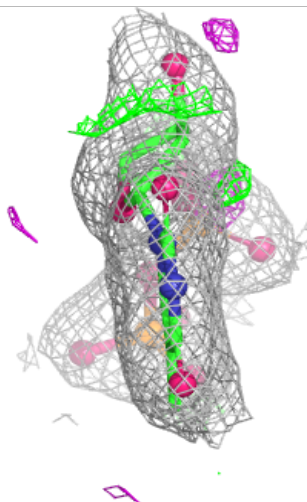
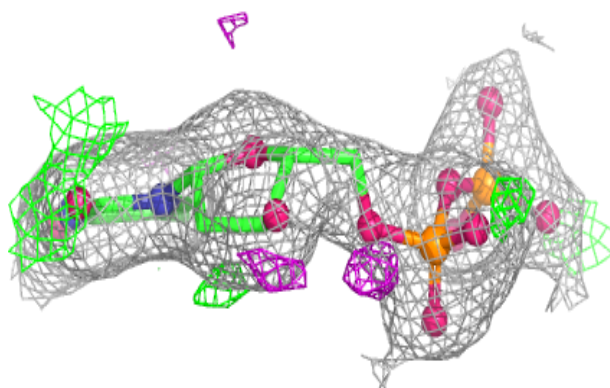
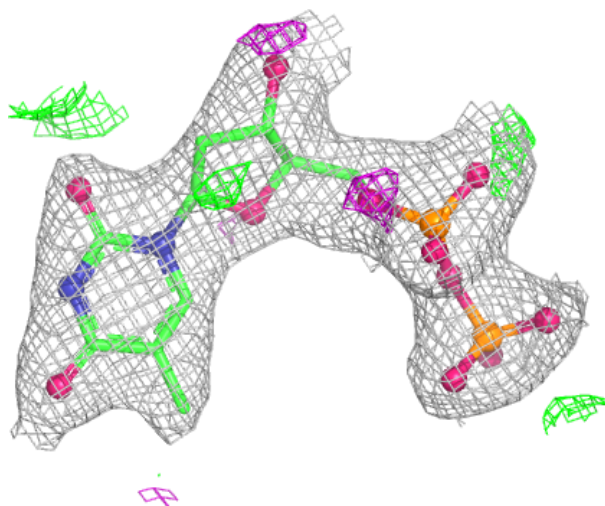
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	GOL	B	1007	6/6	0.85	0.16	37,44,46,50	0
6	BGC	C	8	11/12	0.87	0.11	30,32,34,39	0
6	BGC	D	8	11/12	0.88	0.09	41,44,46,51	0
5	GOL	A	1006	6/6	0.89	0.12	32,43,45,49	0
4	TYD	B	1004	25/25	0.95	0.09	31,37,38,40	0
5	GOL	A	1005	6/6	0.95	0.08	36,38,38,40	0
4	TYD	A	1003	25/25	0.96	0.09	21,32,35,37	0
3	K	B	417	1/1	0.97	0.08	38,38,38,38	0
3	K	A	417	1/1	0.99	0.07	25,25,25,25	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.

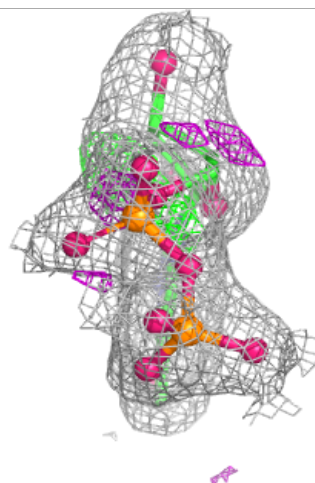
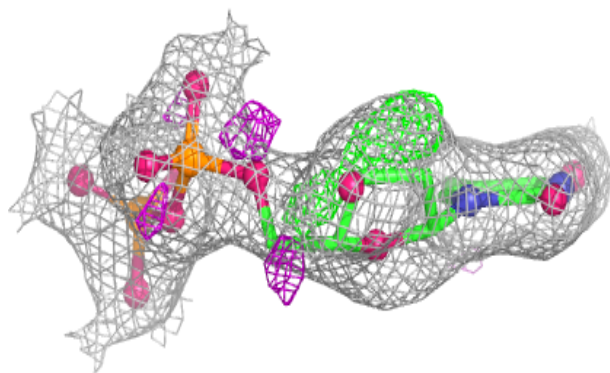
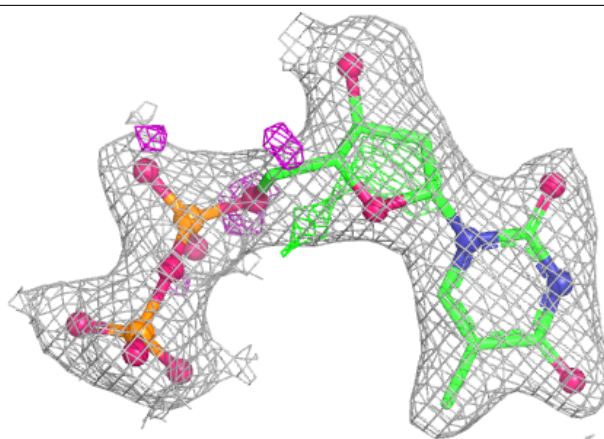
Electron density around TYD B 1004:

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 1003:

$2mF_o-DF_c$ (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.