



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 24, 2023 – 10:32 AM EDT

PDB ID : 5TXN
Title : STRUCTURE OF Q151M MUTANT HIV-1 REVERSE TRANSCRIPTASE (RT) TERNARY COMPLEX WITH A DOUBLE STRANDED DNA AND AN INCOMING DATP
Authors : Das, K.; Martinez, S.M.; Arnold, E.
Deposited on : 2016-11-17
Resolution : 2.55 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtrriage (Phenix) : 1.13
EDS : 2.35.1
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) : 2.35.1

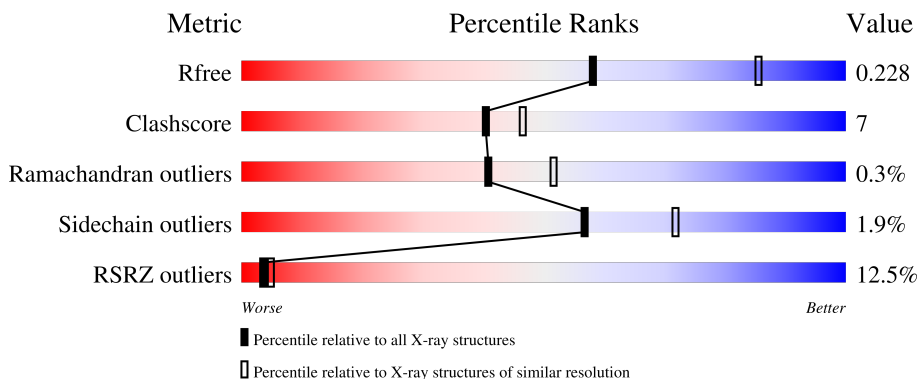
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.55 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	1284 (2.56-2.52)
Clashscore	141614	1332 (2.56-2.52)
Ramachandran outliers	138981	1315 (2.56-2.52)
Sidechain outliers	138945	1315 (2.56-2.52)
RSRZ outliers	127900	1272 (2.56-2.52)

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	556	 15% 82% 17%
1	C	556	 15% 80% 20%
2	B	428	 7% 82% 14%
2	D	428	 11% 82% 14%

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Mol	Chain	Length	Quality of chain
3	E	27	<p>15% 37% 52% 11%</p>
3	T	27	<p>22% 59% 26% 11%</p>
4	F	21	<p>5% 43% 48% 5% 5%</p>
4	P	21	<p>10% 38% 48% 10% 5%</p>
5	G	2	<p>100%</p>
5	H	2	<p>50% 50%</p>

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 17923 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 HIV-1 REVERSE TRANSCRIPTASE P66 SUBUNIT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	555	4510	2920	750	831	9	0	0	0
1	C	555	4510	2920	750	831	9	0	0	0

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MET	-	initiating methionine	UNP P03366
A	0	VAL	-	expression tag	UNP P03366
A	151	MET	GLN	engineered mutation	UNP P03366
A	258	CYS	GLN	engineered mutation	UNP P03366
A	280	SER	CYS	engineered mutation	UNP P03366
A	498	ASN	ASP	engineered mutation	UNP P03366
C	-1	MET	-	initiating methionine	UNP P03366
C	0	VAL	-	expression tag	UNP P03366
C	151	MET	GLN	engineered mutation	UNP P03366
C	258	CYS	GLN	engineered mutation	UNP P03366
C	280	SER	CYS	engineered mutation	UNP P03366
C	498	ASN	ASP	engineered mutation	UNP P03366

- Molecule 2 is a protein called HIV-1 REVERSE TRANSCRIPTASE P51 SUBUNIT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	412	3400	2212	563	619	6	0	0	0
2	D	412	3400	2212	563	619	6	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	280	SER	CYS	engineered mutation	UNP P03366
D	280	SER	CYS	engineered mutation	UNP P03366

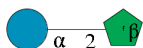
- Molecule 3 is a DNA chain called DNA (5'-D(*TP*GP*GP*TP*CP*GP*GP*CP*GP*CP*CP*GP*AP*AP*CP*AP*GP*GP*GP*AP*CP*TP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
3	T	24	494	233	97	141	23	0	0	0
3	E	24	494	233	97	141	23	0	0	0

- Molecule 4 is a DNA chain called DNA (5'-D(*CP*AP*GP*TP*CP*CP*CP*TP*GP*TP*TP*CP*GP*GP*(MRG)P*CP*GP*CP*CP*G)-3').

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	P	S			
4	P	20	407	195	72	120	19	1	0	0	0
4	F	20	407	195	72	120	19	1	0	0	0

- Molecule 5 is an oligosaccharide called beta-D-fructofuranose-(2-1)-alpha-D-glucopyranose.



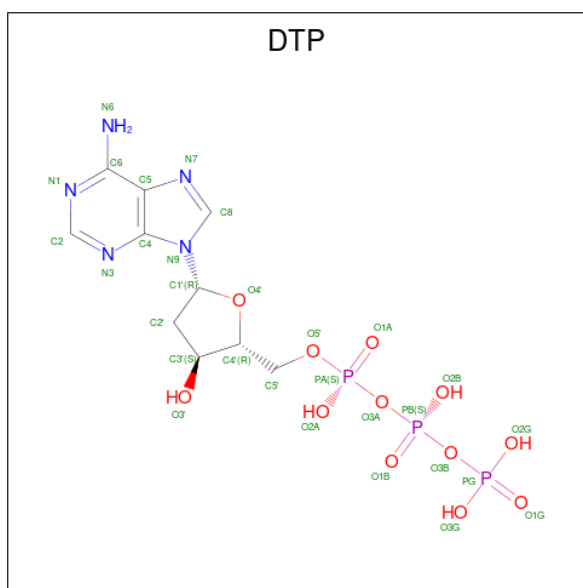
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
			Total	C	O			
5	G	2	23	12	11	0	0	0
5	H	2	23	12	11	0	0	0

- Molecule 6 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

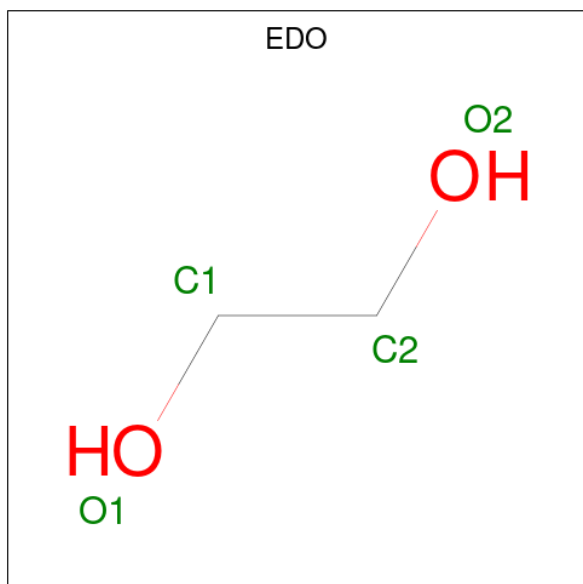
- Molecule 7 is 2'-DEOXYADENOSINE 5'-TRIPHOSPHATE (three-letter code: DTP) (for-

mula: C₁₀H₁₆N₅O₁₂P₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
7	A	1	Total	C	N	O	P	0	0
			30	10	5	12	3		
7	C	1	Total	C	N	O	P	0	0
			30	10	5	12	3		

- Molecule 8 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



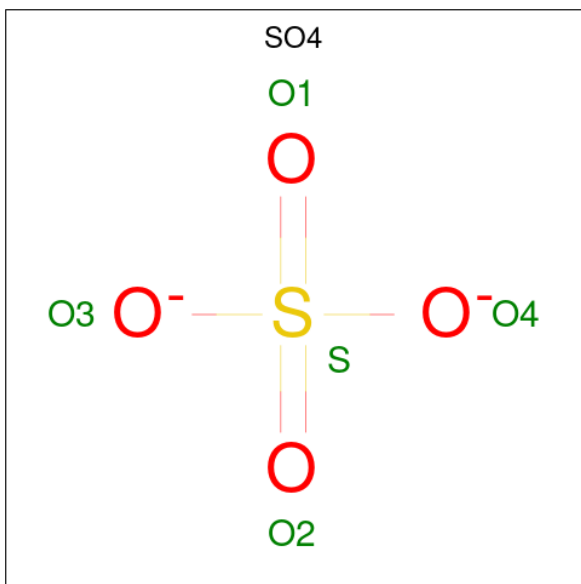
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	
8	B	1	Total	C	O	0	0
			4	2	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
8	D	1	4	2	2	0	0

- Molecule 9 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
10	A	50	50	50	0	0
10	B	42	42	42	0	0
10	T	1	1	1	0	0
10	P	1	1	1	0	0
10	C	48	48	48	0	0
10	D	33	33	33	0	0
10	E	1	1	1	0	0

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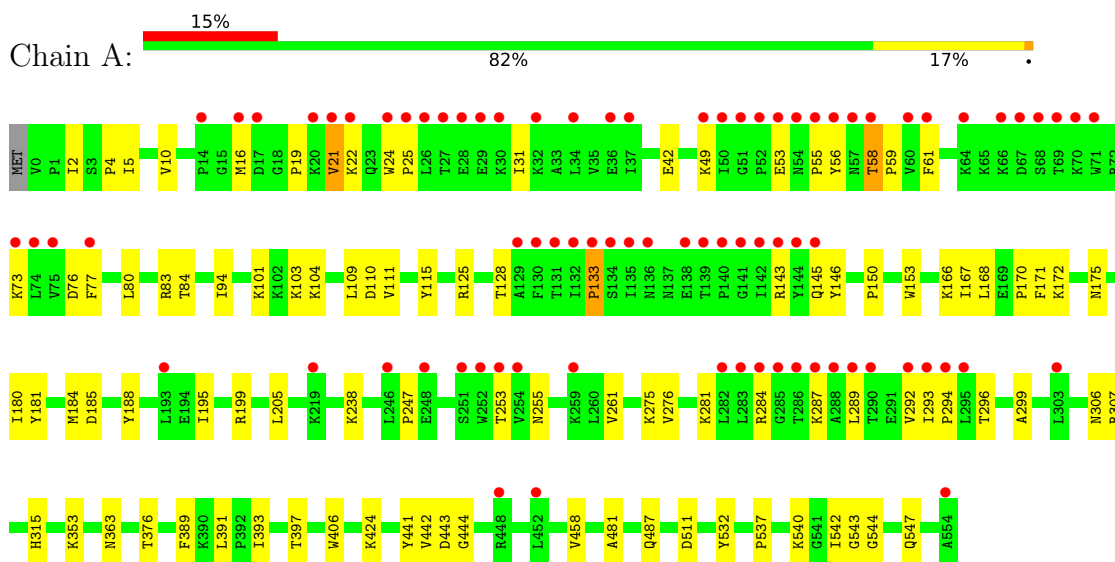
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	F	2	Total	O	0	0
			2	2		

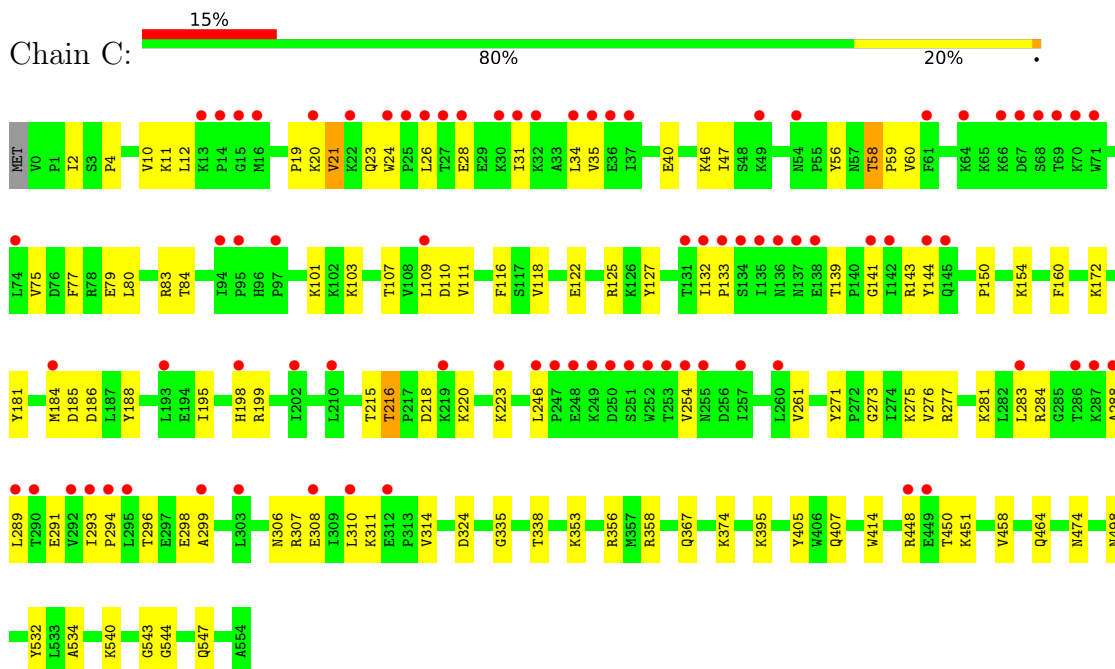
3 Residue-property plots [i](#)

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.

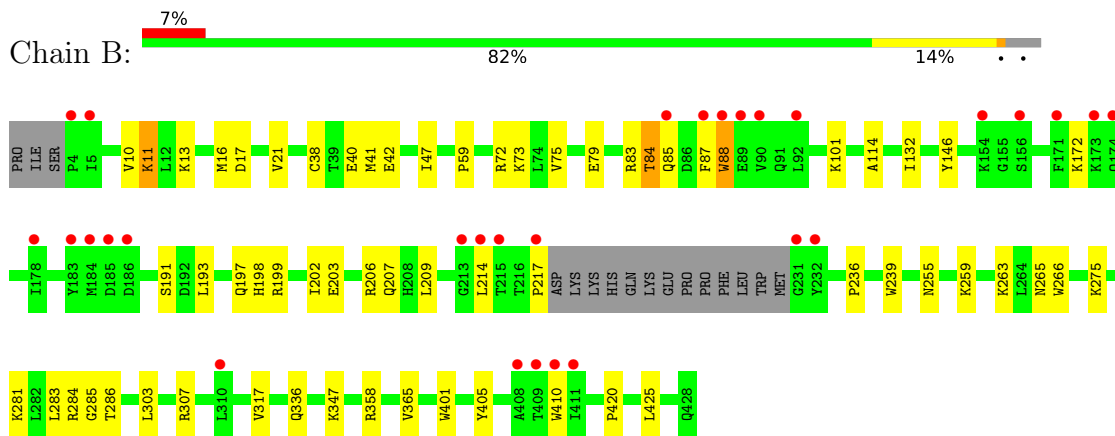
- Molecule 1: HIV-1 REVERSE TRANSCRIPTASE P66 SUBUNIT



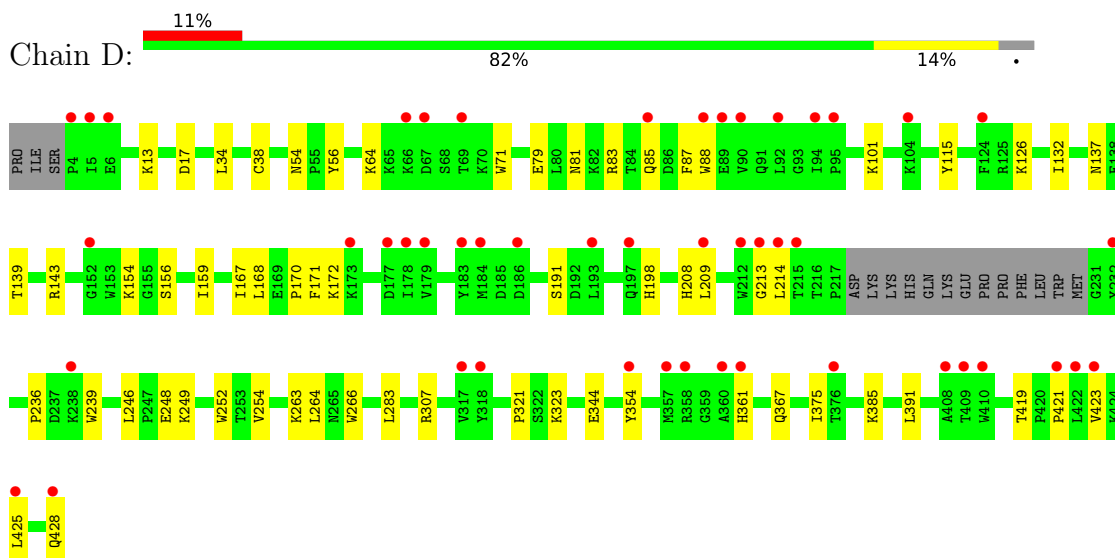
- Molecule 1: HIV-1 REVERSE TRANSCRIPTASE P66 SUBUNIT



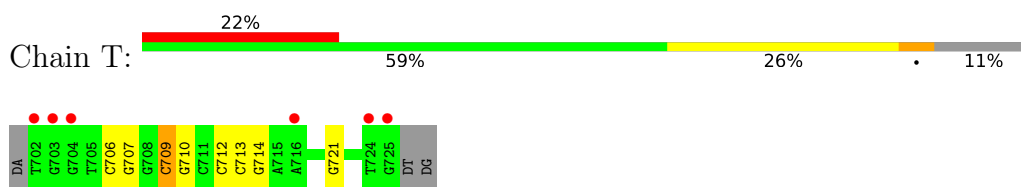
- Molecule 2: HIV-1 REVERSE TRANSCRIPTASE P51 SUBUNIT



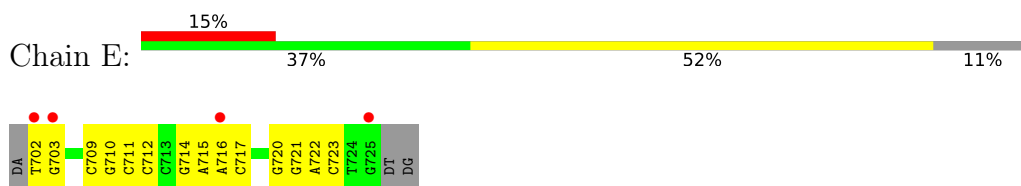
- Molecule 2: HIV-1 REVERSE TRANSCRIPTASE P51 SUBUNIT



- Molecule 3: DNA (5'-D(*TP*GP*GP*TP*CP*GP*GP*CP*GP*CP*CP*CP*GP*AP*AP*CP*AP*GP*GP*GP*AP*CP*TP*G)-3')

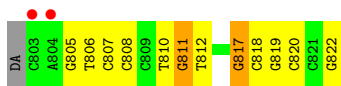


- Molecule 3: DNA (5'-D(*TP*GP*GP*TP*CP*GP*GP*CP*GP*CP*CP*CP*GP*AP*AP*CP*AP*GP*GP*GP*AP*CP*TP*G)-3')



- Molecule 4: DNA (5'-D(*CP*AP*GP*TP*CP*CP*CP*TP*GP*TP*TP*CP*GP*GP*(MRG)P*CP*GP*CP*CP*G)-3')

Chain P: 



- Molecule 4: DNA (5'-D(*CP*AP*GP*TP*CP*CP*CP*TP*GP*TP*TP*CP*GP*GP*(MRG)P*CP*GP*CP*CP*G)-3')

Chain F: 



- Molecule 5: beta-D-fructofuranose-(2-1)-alpha-D-glucopyranose

Chain G: 



- Molecule 5: beta-D-fructofuranose-(2-1)-alpha-D-glucopyranose

Chain H: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	90.04Å 133.63Å 139.21Å 90.00° 97.88° 90.00°	Depositor
Resolution (Å)	33.38 – 2.55 36.37 – 2.53	Depositor EDS
% Data completeness (in resolution range)	98.2 (33.38-2.55) 97.4 (36.37-2.53)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.86 (at 2.54Å)	Xtrriage
Refinement program	PHENIX (1.10_2155: ???)	Depositor
R, R_{free}	0.196 , 0.226 0.197 , 0.228	Depositor DCC
R_{free} test set	3101 reflections (2.92%)	wwPDB-VP
Wilson B-factor (Å ²)	57.2	Xtrriage
Anisotropy	0.201	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 63.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	17923	wwPDB-VP
Average B, all atoms (Å ²)	89.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.79% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: FRU, GLC, MRG, EDO, SO4, DTP, MG

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	0.25	0/4628	0.45	1/6288 (0.0%)
1	C	0.27	1/4628 (0.0%)	0.43	1/6288 (0.0%)
2	B	0.27	1/3497 (0.0%)	0.45	1/4751 (0.0%)
2	D	0.28	2/3497 (0.1%)	0.44	1/4751 (0.0%)
3	E	0.52	0/555	0.85	0/856
3	T	0.51	0/555	0.87	1/856 (0.1%)
4	F	0.54	0/424	0.94	0/649
4	P	0.61	1/424 (0.2%)	1.03	2/649 (0.3%)
All	All	0.31	5/18208 (0.0%)	0.52	7/25088 (0.0%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	172	LYS	CE-NZ	-6.31	1.33	1.49
2	D	172	LYS	CE-NZ	-6.27	1.33	1.49
1	C	172	LYS	CE-NZ	-6.27	1.33	1.49
2	D	391	LEU	C-N	5.66	1.45	1.34
4	P	811	DG	C4'-O4'	5.28	1.50	1.45

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	P	820	DC	O4'-C4'-C3'	-9.46	100.32	106.00
3	T	709	DC	O4'-C4'-C3'	-6.50	101.90	104.50
1	A	73	LYS	CB-CA-C	-5.77	98.86	110.40
2	B	172	LYS	CD-CE-NZ	5.37	124.04	111.70
2	D	172	LYS	CD-CE-NZ	5.36	124.03	111.70
1	C	172	LYS	CD-CE-NZ	5.28	123.84	111.70
4	P	805	DG	C5'-C4'-O4'	5.00	118.80	109.30

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	4510	0	4571	67	0
1	C	4510	0	4571	61	0
2	B	3400	0	3433	46	0
2	D	3400	0	3433	41	0
3	E	494	0	269	9	0
3	T	494	0	269	9	0
4	F	407	0	229	13	0
4	P	407	0	229	11	0
5	G	23	0	21	0	0
5	H	23	0	21	0	0
6	A	2	0	0	0	0
6	C	2	0	0	0	0
7	A	30	0	12	2	0
7	C	30	0	12	0	0
8	B	4	0	6	0	0
8	D	4	0	6	3	0
9	D	5	0	0	0	0
10	A	50	0	0	2	0
10	B	42	0	0	2	0
10	C	48	0	0	2	0
10	D	33	0	0	4	0
10	E	1	0	0	0	0
10	F	2	0	0	0	0
10	P	1	0	0	1	0
10	T	1	0	0	0	0
All	All	17923	0	17082	245	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 (245) 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)
4:P:817:MRG:H231	4:P:817:MRG:N3	1.90	0.87
1:A:167:ILE:O	1:A:170:PRO:HD2	1.79	0.82
1:A:542:ILE:HG23	2:B:283:LEU:HD13	1.65	0.77
2:D:263:LYS:HE3	2:D:425:LEU:HA	1.69	0.75
1:A:171:PHE:CD2	1:A:205:LEU:HD13	2.24	0.71
1:A:543:GLY:O	1:A:547:GLN:NE2	2.23	0.71
1:A:111:VAL:HB	1:A:185:ASP:HB2	1.74	0.69
3:E:722:DA:H2''	3:E:723:DC:H5''	1.75	0.69
4:F:817:MRG:N3	4:F:817:MRG:C22	2.56	0.68
2:B:114:ALA:HB2	2:B:214:LEU:HD22	1.76	0.67
1:C:448:ARG:O	1:C:451:LYS:NZ	2.28	0.67
1:C:261:VAL:HG13	1:C:276:VAL:HG11	1.76	0.67
2:B:199:ARG:HA	2:B:202:ILE:HD12	1.77	0.67
2:D:56:TYR:HE2	2:D:126:LYS:HE2	1.60	0.66
2:D:421:PRO:HD3	10:D:610:HOH:O	1.95	0.66
1:C:288:ALA:HB3	1:C:291:GLU:HB2	1.77	0.66
4:F:817:MRG:N3	4:F:817:MRG:H221	2.10	0.66
1:C:543:GLY:O	1:C:547:GLN:NE2	2.28	0.65
1:A:253:THR:HA	1:A:292:VAL:HA	1.78	0.65
1:A:296:THR:HG23	1:A:299:ALA:H	1.61	0.65
1:C:276:VAL:HG23	1:C:353:LYS:HE2	1.79	0.65
1:C:498:ASN:ND2	10:C:702:HOH:O	2.30	0.64
1:C:60:VAL:HG22	1:C:75:VAL:HG13	1.79	0.64
1:A:542:ILE:HG23	2:B:283:LEU:CD1	2.28	0.64
1:C:139:THR:HG22	1:C:141:GLY:H	1.63	0.64
1:A:172:LYS:HE2	1:A:180:ILE:HB	1.80	0.63
1:A:542:ILE:CG2	2:B:283:LEU:HD13	2.27	0.63
2:B:193:LEU:HD13	2:B:197:GLN:HB3	1.80	0.63
1:C:246:LEU:HD11	1:C:310:LEU:HD22	1.80	0.63
2:B:17:ASP:O	2:B:83:ARG:NH1	2.32	0.62
2:D:191:SER:OG	2:D:198:HIS:ND1	2.25	0.62
2:B:365:VAL:HG11	2:B:401:TRP:HB2	1.82	0.61
1:A:487:GLN:NE2	10:A:705:HOH:O	2.33	0.61
2:D:137:ASN:ND2	10:D:606:HOH:O	2.34	0.61
1:C:19:PRO:HG3	1:C:80:LEU:HB2	1.83	0.61
1:C:353:LYS:O	1:C:374:LYS:NZ	2.34	0.60
2:D:246:LEU:HD11	2:D:264:LEU:HD21	1.84	0.60
1:C:395:LYS:HE3	10:C:727:HOH:O	2.00	0.60
1:A:128:THR:HG21	1:A:146:TYR:HB2	1.84	0.60
1:A:261:VAL:HG13	1:A:276:VAL:HG11	1.83	0.59
1:A:110:ASP:OD1	1:A:111:VAL:N	2.36	0.59
1:A:24:TRP:HE3	1:A:25:PRO:HD2	1.66	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:72:ARG:NH1	10:B:601:HOH:O	2.21	0.58
1:C:107:THR:OG1	1:C:198:HIS:NE2	2.34	0.58
3:T:713:DC:H2'	3:T:714:DG:C8	2.38	0.58
4:F:807:DC:H2''	4:F:808:DC:H5''	1.86	0.58
3:E:702:DT:H3'	3:E:703:DG:H8	1.69	0.57
3:E:720:DG:H2''	3:E:721:DG:C8	2.39	0.57
1:A:58:THR:HG21	1:A:77:PHE:HE1	1.70	0.56
1:C:216:THR:O	1:C:216:THR:OG1	2.16	0.56
2:B:209:LEU:HD22	2:B:214:LEU:HD23	1.88	0.55
2:B:203:GLU:O	2:B:207:GLN:HG2	2.07	0.55
1:A:171:PHE:CE2	1:A:205:LEU:HD13	2.41	0.55
1:C:111:VAL:HB	1:C:185:ASP:HB2	1.88	0.55
2:B:255:ASN:ND2	10:B:603:HOH:O	2.32	0.55
2:B:13:LYS:HD2	2:B:85:GLN:HB3	1.88	0.54
2:D:87:PHE:HZ	2:D:159:ILE:HG13	1.73	0.54
1:A:56:TYR:O	1:A:143:ARG:NH2	2.41	0.54
1:C:56:TYR:O	1:C:143:ARG:NH2	2.41	0.54
1:A:389:PHE:HB3	1:A:391:LEU:HD13	1.90	0.54
2:D:209:LEU:HD22	2:D:214:LEU:HD12	1.90	0.54
4:F:804:DA:H2'	4:F:805:DG:C8	2.43	0.54
2:D:168:LEU:O	2:D:171:PHE:N	2.40	0.53
1:C:181:TYR:HB2	1:C:188:TYR:HB3	1.89	0.53
2:D:64:LYS:HE3	2:D:71:TRP:CE2	2.44	0.53
2:B:266:TRP:HD1	2:B:425:LEU:HD13	1.74	0.53
2:B:79:GLU:HG3	2:B:83:ARG:HE	1.74	0.53
1:C:254:VAL:HG13	1:C:283:LEU:HD22	1.91	0.53
2:D:249:LYS:HB2	2:D:252:TRP:CE2	2.44	0.53
1:A:391:LEU:HB3	1:A:393:ILE:HG22	1.91	0.52
1:C:335:GLY:HA2	1:C:367:GLN:HE22	1.74	0.52
1:C:543:GLY:HA3	10:D:623:HOH:O	2.09	0.52
4:F:817:MRG:H2'	4:F:818:DC:C6	2.44	0.52
1:C:122:GLU:HA	1:C:125:ARG:HD2	1.92	0.52
2:B:317:VAL:HG12	2:B:347:LYS:HD3	1.92	0.52
2:D:248:GLU:HA	8:D:503:EDO:O2	2.09	0.52
1:A:166:LYS:O	1:A:170:PRO:HD3	2.10	0.52
4:F:814:DC:H2''	4:F:815:DG:C8	2.45	0.52
1:A:238:LYS:HD2	1:A:315:HIS:CG	2.45	0.52
1:C:395:LYS:NZ	1:C:414:TRP:O	2.43	0.51
1:A:80:LEU:O	1:A:84:THR:OG1	2.24	0.51
1:A:31:ILE:HG23	1:A:133:PRO:HG2	1.92	0.51
3:E:702:DT:H3'	3:E:703:DG:C8	2.44	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:F:807:DC:H4'	4:F:808:DC:OP1	2.10	0.51
2:B:101:LYS:O	2:B:236:PRO:HB2	2.11	0.51
4:P:807:DC:H4'	4:P:808:DC:OP1	2.10	0.51
1:C:458:VAL:HG13	1:C:464:GLN:HG2	1.92	0.51
2:D:54:ASN:HB3	2:D:143:ARG:HH21	1.77	0.50
2:D:254:VAL:HG13	2:D:283:LEU:HD22	1.93	0.50
2:B:317:VAL:HG12	2:B:347:LYS:HB3	1.94	0.50
2:D:85:GLN:HA	2:D:88:TRP:CE2	2.46	0.50
2:D:213:GLY:O	2:D:214:LEU:HD23	2.11	0.50
4:F:818:DC:H2'	4:F:819:DG:C8	2.47	0.50
1:C:281:LYS:O	1:C:284:ARG:HG2	2.11	0.50
1:C:195:ILE:HD11	1:C:199:ARG:HH21	1.76	0.50
2:D:167:ILE:HD13	2:D:214:LEU:HD11	1.93	0.50
1:A:125:ARG:O	1:A:128:THR:OG1	2.23	0.50
2:B:191:SER:OG	2:B:198:HIS:ND1	2.35	0.50
2:B:38:CYS:SG	2:B:132:ILE:HD11	2.51	0.49
1:C:356:ARG:HG2	1:C:358:ARG:HG3	1.94	0.49
2:D:249:LYS:HB2	2:D:252:TRP:CZ2	2.47	0.49
2:B:266:TRP:NE1	2:B:425:LEU:HD22	2.28	0.49
1:C:273:GLY:HA2	1:C:338:THR:HG21	1.94	0.49
1:A:397:THR:HG21	1:A:424:LYS:HA	1.95	0.49
1:A:94:ILE:O	10:A:701:HOH:O	2.20	0.49
1:A:184:MET:HG2	4:P:822:DG:H2''	1.94	0.49
1:A:543:GLY:HA2	2:B:283:LEU:O	2.12	0.49
1:C:24:TRP:HB2	1:C:26:LEU:HG	1.94	0.49
2:D:38:CYS:SG	2:D:132:ILE:HD11	2.52	0.49
1:C:184:MET:HG2	4:F:822:DG:H2''	1.94	0.49
1:C:543:GLY:HA2	2:D:283:LEU:O	2.13	0.49
3:T:713:DC:H2''	3:T:714:DG:H5'	1.95	0.49
2:D:421:PRO:HB2	2:D:423:VAL:HG23	1.94	0.49
1:A:19:PRO:HD3	1:A:80:LEU:HD13	1.95	0.49
3:E:716:DA:H2''	3:E:717:DC:OP2	2.13	0.49
4:F:805:DG:H4'	4:F:806:DT:OP1	2.13	0.49
2:B:236:PRO:HA	2:B:239:TRP:CD2	2.48	0.48
1:C:58:THR:HG21	1:C:77:PHE:CE1	2.48	0.48
1:A:276:VAL:HG23	1:A:353:LYS:HE2	1.95	0.48
1:C:47:ILE:HG12	1:C:144:TYR:HB3	1.95	0.48
1:C:296:THR:HG23	1:C:299:ALA:H	1.76	0.48
4:P:810:DT:H2''	4:P:811:DG:H5'	1.94	0.48
2:D:323:LYS:NZ	2:D:344:GLU:OE2	2.31	0.48
2:D:79:GLU:HG3	2:D:83:ARG:HE	1.79	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:10:VAL:HG11	1:A:153:TRP:CZ2	2.49	0.48
4:P:818:DC:H2'	4:P:819:DG:C8	2.49	0.48
1:A:443:ASP:OD1	1:A:444:GLY:N	2.47	0.47
2:B:84:THR:HG22	2:B:87:PHE:HB3	1.95	0.47
3:T:706:DC:H2'	3:T:707:DG:C8	2.49	0.47
2:D:354:TYR:HE2	2:D:375:ILE:HG13	1.79	0.47
8:D:503:EDO:O1	10:D:601:HOH:O	2.20	0.47
2:B:266:TRP:HE1	2:B:425:LEU:HD22	1.80	0.47
1:C:58:THR:HG21	1:C:77:PHE:HE1	1.79	0.47
1:C:307:ARG:O	1:C:311:LYS:HG3	2.14	0.47
2:D:13:LYS:HD2	2:D:85:GLN:HB3	1.97	0.47
2:D:423:VAL:HG12	2:D:423:VAL:O	2.14	0.47
4:F:806:DT:H2'	4:F:807:DC:C6	2.50	0.47
2:B:47:ILE:HG22	2:B:146:TYR:HA	1.96	0.47
1:C:308:GLU:HA	1:C:311:LYS:HE2	1.97	0.47
3:T:709:DC:H2'	3:T:710:DG:C8	2.50	0.47
2:D:323:LYS:O	2:D:385:LYS:NZ	2.48	0.47
1:A:458:VAL:HG12	2:B:286:THR:HG21	1.97	0.47
1:C:79:GLU:OE1	1:C:83:ARG:NH2	2.48	0.46
3:E:709:DC:H2'	3:E:710:DG:H8	1.80	0.46
1:A:115:TYR:CD2	7:A:603:DTP:H2'1	2.51	0.46
1:C:199:ARG:NH1	1:C:223:LYS:HD2	2.30	0.46
1:C:544:GLY:HA2	1:C:547:GLN:HE21	1.81	0.46
2:B:358:ARG:NH2	2:B:405:TYR:O	2.44	0.46
1:A:406:TRP:CZ2	2:B:420:PRO:HB3	2.49	0.46
1:C:21:VAL:HG23	1:C:59:PRO:HD3	1.97	0.46
1:C:405:TYR:CE2	1:C:407:GLN:HB2	2.51	0.46
4:P:807:DC:H2''	4:P:808:DC:H5''	1.98	0.45
1:A:441:TYR:CG	1:A:544:GLY:HA3	2.51	0.45
1:C:31:ILE:O	1:C:35:VAL:HG23	2.17	0.45
1:A:247:PRO:O	1:A:307:ARG:NH2	2.50	0.45
4:P:817:MRG:H2'	4:P:818:DC:C6	2.51	0.45
1:C:275:LYS:O	1:C:306:ASN:ND2	2.43	0.45
1:C:395:LYS:HD2	1:C:414:TRP:CH2	2.52	0.45
1:C:46:LYS:HD3	1:C:116:PHE:HB3	1.97	0.45
2:D:428:GLN:H	2:D:428:GLN:CD	2.19	0.45
1:A:10:VAL:HG11	1:A:153:TRP:CH2	2.51	0.45
1:A:21:VAL:HG23	1:A:59:PRO:HD3	1.98	0.45
1:C:31:ILE:HD13	1:C:133:PRO:HB2	1.98	0.45
1:A:22:LYS:HA	1:A:22:LYS:HD3	1.77	0.45
1:A:171:PHE:CE1	1:A:175:ASN:OD1	2.70	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:181:TYR:HB2	1:A:188:TYR:HB3	1.98	0.44
4:P:818:DC:H2'	4:P:819:DG:H8	1.81	0.44
1:C:199:ARG:HH11	1:C:223:LYS:HD2	1.81	0.44
4:F:818:DC:H2'	4:F:819:DG:H8	1.82	0.44
3:T:721:DG:H5'	3:T:721:DG:H8	1.82	0.44
1:A:376:THR:HG21	2:B:401:TRP:CH2	2.53	0.44
4:P:812:DT:OP1	10:P:901:HOH:O	2.20	0.44
1:C:103:LYS:HA	1:C:103:LYS:HD3	1.76	0.44
1:A:103:LYS:HD3	1:A:103:LYS:HA	1.84	0.44
1:A:255:ASN:HB2	1:A:289:LEU:HG	2.00	0.44
2:B:114:ALA:N	2:B:214:LEU:HD13	2.33	0.44
1:A:21:VAL:HG23	1:A:58:THR:HA	1.99	0.43
1:A:61:PHE:HE2	1:A:76:ASP:HB2	1.83	0.43
1:A:363:ASN:HA	1:A:511:ASP:OD1	2.18	0.43
2:B:206:ARG:NH2	2:B:217:PRO:O	2.49	0.43
7:A:603:DTP:O2B	7:A:603:DTP:O1A	2.36	0.43
3:T:712:DC:H2''	3:T:713:DC:H6	1.83	0.43
1:C:110:ASP:HB3	1:C:220:LYS:HB3	1.99	0.43
1:A:281:LYS:O	1:A:284:ARG:HG2	2.19	0.43
2:D:266:TRP:CE3	2:D:425:LEU:HG	2.53	0.43
2:D:307:ARG:HH22	8:D:503:EDO:H12	1.83	0.43
2:D:321:PRO:O	2:D:385:LYS:NZ	2.38	0.43
1:A:42:GLU:OE1	1:A:49:LYS:HE3	2.19	0.43
1:A:287:LYS:H	1:A:287:LYS:HG2	1.56	0.43
1:A:195:ILE:O	1:A:199:ARG:HG3	2.19	0.43
1:C:35:VAL:HG22	1:C:132:ILE:HG21	2.00	0.43
1:C:118:VAL:HG21	1:C:160:PHE:HD1	1.84	0.43
2:D:81:ASN:HB3	2:D:154:LYS:HD2	2.01	0.43
2:B:21:VAL:HB	2:B:59:PRO:HD3	2.00	0.43
1:C:254:VAL:HG12	1:C:289:LEU:HD12	2.00	0.43
1:A:168:LEU:HD23	1:A:168:LEU:HA	1.81	0.43
2:B:266:TRP:CD1	2:B:425:LEU:HD22	2.54	0.43
2:B:42:GLU:HA	2:B:47:ILE:O	2.19	0.43
3:T:721:DG:H5'	3:T:721:DG:C8	2.54	0.43
1:A:275:LYS:O	1:A:306:ASN:ND2	2.43	0.42
1:C:12:LEU:HD11	1:C:127:TYR:CZ	2.54	0.42
1:A:532:TYR:OH	2:B:259:LYS:NZ	2.48	0.42
1:C:293:ILE:HG13	1:C:294:PRO:HD2	2.02	0.42
1:C:532:TYR:CE2	1:C:534:ALA:HB2	2.55	0.42
2:D:115:TYR:CD2	2:D:156:SER:HB3	2.54	0.42
1:A:293:ILE:HG13	1:A:294:PRO:HD2	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:442:VAL:HB	1:A:481:ALA:HB1	2.01	0.42
2:B:41:MET:HB3	2:B:47:ILE:HG12	2.01	0.42
1:C:20:LYS:HG2	1:C:56:TYR:HA	2.02	0.42
1:A:53:GLU:O	1:A:55:PRO:HD3	2.19	0.42
2:B:85:GLN:HA	2:B:88:TRP:CE2	2.54	0.42
2:B:303:LEU:HD21	2:B:307:ARG:HH22	1.85	0.42
3:E:714:DG:H2'	3:E:715:DA:C8	2.55	0.42
1:A:184:MET:HE3	1:A:184:MET:HA	2.01	0.42
3:T:709:DC:H2'	3:T:710:DG:H8	1.84	0.42
1:C:10:VAL:HG12	1:C:11:LYS:H	1.85	0.42
3:T:712:DC:O2	4:P:817:MRG:H212	2.20	0.42
2:B:263:LYS:HE3	2:B:425:LEU:HA	2.02	0.42
2:D:170:PRO:HB2	2:D:208:HIS:CE1	2.55	0.42
3:E:711:DC:H2'	3:E:712:DC:C6	2.55	0.42
1:A:145:GLN:HG3	1:A:145:GLN:O	2.19	0.41
2:B:281:LYS:O	2:B:284:ARG:HG3	2.20	0.41
4:P:806:DT:H2'	4:P:807:DC:C6	2.55	0.41
2:D:361:HIS:H	2:D:367:GLN:NE2	2.18	0.41
4:F:807:DC:H2''	4:F:808:DC:C5'	2.50	0.41
1:C:271:TYR:CE2	1:C:310:LEU:HA	2.55	0.41
1:C:84:THR:HB	1:C:154:LYS:HE2	2.03	0.41
2:B:73:LYS:HZ3	2:B:75:VAL:HG22	1.85	0.41
2:B:275:LYS:HA	2:B:275:LYS:HD3	1.80	0.41
3:E:709:DC:H2'	3:E:710:DG:C8	2.54	0.41
1:A:543:GLY:HA2	2:B:285:GLY:O	2.20	0.41
2:D:34:LEU:HD23	2:D:34:LEU:HA	1.91	0.41
2:D:101:LYS:O	2:D:236:PRO:HB2	2.21	0.41
2:D:419:THR:O	2:D:421:PRO:HD3	2.21	0.41
1:C:540:LYS:HD3	1:C:540:LYS:HA	1.82	0.41
1:A:537:PRO:HG2	1:A:540:LYS:HG3	2.03	0.40
1:C:271:TYR:CE2	1:C:310:LEU:HG	2.55	0.40
2:D:17:ASP:O	2:D:83:ARG:HD3	2.21	0.40
1:A:540:LYS:HD3	1:A:540:LYS:HA	1.79	0.40
2:B:10:VAL:O	2:B:11:LYS:HD2	2.21	0.40
2:D:87:PHE:CZ	2:D:159:ILE:HG13	2.54	0.40
1:A:16:MET:HG2	1:A:83:ARG:HD3	2.03	0.40
1:A:540:LYS:HE3	2:B:265:ASN:OD1	2.22	0.40
2:D:236:PRO:HA	2:D:239:TRP:CE2	2.57	0.40
1:A:104:LYS:HE3	1:A:104:LYS:HB2	1.86	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	553/556 (100%)	533 (96%)	17 (3%)	3 (0%)	29	40
1	C	553/556 (100%)	534 (97%)	17 (3%)	2 (0%)	34	46
2	B	408/428 (95%)	399 (98%)	9 (2%)	0	100	100
2	D	408/428 (95%)	394 (97%)	14 (3%)	0	100	100
All	All	1922/1968 (98%)	1860 (97%)	57 (3%)	5 (0%)	41	51

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	4	PRO
1	C	4	PRO
1	A	150	PRO
1	A	133	PRO
1	C	150	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	495/497 (100%)	489 (99%)	6 (1%)	71	81
1	C	495/497 (100%)	476 (96%)	19 (4%)	33	45
2	B	374/390 (96%)	367 (98%)	7 (2%)	57	72
2	D	374/390 (96%)	373 (100%)	1 (0%)	92	96
All	All	1738/1774 (98%)	1705 (98%)	33 (2%)	57	72

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	ILE
1	A	5	ILE
1	A	21	VAL
1	A	58	THR
1	A	101	LYS
1	A	109	LEU
2	B	11	LYS
2	B	16	MET
2	B	40	GLU
2	B	84	THR
2	B	88	TRP
2	B	336	GLN
2	B	410	TRP
1	C	2	ILE
1	C	21	VAL
1	C	23	GLN
1	C	28	GLU
1	C	34	LEU
1	C	40	GLU
1	C	58	THR
1	C	101	LYS
1	C	109	LEU
1	C	186	ASP
1	C	215	THR
1	C	216	THR
1	C	218	ASP
1	C	277	ARG
1	C	298	GLU
1	C	314	VAL
1	C	324	ASP
1	C	450	THR
1	C	474	ASN
2	D	139	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	174	GLN
1	A	182	GLN
1	C	367	GLN
1	C	498	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

2 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
4	MRG	P	817	1,3,4	21,28,29	2.47	8 (38%)	19,39,42	1.61	5 (26%)
4	MRG	F	817	1,3,4	21,28,29	2.44	8 (38%)	19,39,42	1.57	5 (26%)

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	MRG	P	817	1,3,4	-	3/8/26/27	0/3/3/3
4	MRG	F	817	1,3,4	-	2/8/26/27	0/3/3/3

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	P	817	MRG	C2-N2	6.98	1.49	1.34
4	F	817	MRG	C2-N2	6.81	1.49	1.34
4	F	817	MRG	C2-N1	4.36	1.43	1.36
4	P	817	MRG	C2-N1	4.32	1.43	1.36
4	P	817	MRG	C4-N3	3.41	1.45	1.37
4	F	817	MRG	C4-N3	3.31	1.45	1.37
4	P	817	MRG	C2'-C3'	-3.28	1.44	1.52
4	F	817	MRG	C2'-C3'	-3.25	1.44	1.52
4	F	817	MRG	C6-N1	2.96	1.42	1.37
4	P	817	MRG	C6-N1	2.85	1.42	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	F	817	MRG	O4'-C4'	-2.65	1.39	1.45
4	F	817	MRG	O3'-C3'	-2.61	1.37	1.43
4	P	817	MRG	O3'-C3'	-2.58	1.37	1.43
4	P	817	MRG	O4'-C4'	-2.57	1.39	1.45
4	P	817	MRG	O5'-C5'	-2.40	1.38	1.44
4	F	817	MRG	O5'-C5'	-2.33	1.39	1.44

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	P	817	MRG	C5-C6-N1	3.48	120.09	113.95
4	F	817	MRG	C5-C6-N1	3.40	119.95	113.95
4	P	817	MRG	C22-C21-N2	3.04	120.89	112.21
4	P	817	MRG	O5'-C5'-C4'	2.77	118.43	108.99
4	F	817	MRG	O5'-C5'-C4'	2.76	118.39	108.99
4	F	817	MRG	C8-N7-C5	2.65	108.04	102.99
4	P	817	MRG	C8-N7-C5	2.61	107.96	102.99
4	F	817	MRG	O6-C6-C5	-2.26	119.95	124.37
4	F	817	MRG	C22-C21-N2	2.24	118.62	112.21
4	P	817	MRG	O6-C6-C5	-2.23	120.02	124.37

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	P	817	MRG	N3-C2-N2-C21
4	P	817	MRG	N1-C2-N2-C21
4	P	817	MRG	C21-C22-C23-S24
4	F	817	MRG	C22-C21-N2-C2
4	F	817	MRG	C21-C22-C23-S24

There are no ring outliers.

2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	P	817	MRG	3	0
4	F	817	MRG	3	0

5.5 Carbohydrates i

4 monosaccharides 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
5	GLC	G	1	5	11,11,12	0.30	0	15,15,17	0.78	0
5	FRU	G	2	5	11,12,12	0.47	0	10,18,18	0.55	0
5	GLC	H	1	5	11,11,12	0.21	0	15,15,17	1.18	2 (13%)
5	FRU	H	2	5	11,12,12	0.44	0	10,18,18	0.64	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. '2' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	GLC	G	1	5	-	0/2/19/22	0/1/1/1
5	FRU	G	2	5	-	3/5/24/24	0/1/1/1
5	GLC	H	1	5	-	1/2/19/22	0/1/1/1
5	FRU	H	2	5	-	3/5/24/24	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	H	1	GLC	C1-O5-C5	2.27	115.26	112.19
5	H	1	GLC	C6-C5-C4	-2.16	107.94	113.00

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	G	2	FRU	O1-C1-C2-C3
5	G	2	FRU	O1-C1-C2-O2

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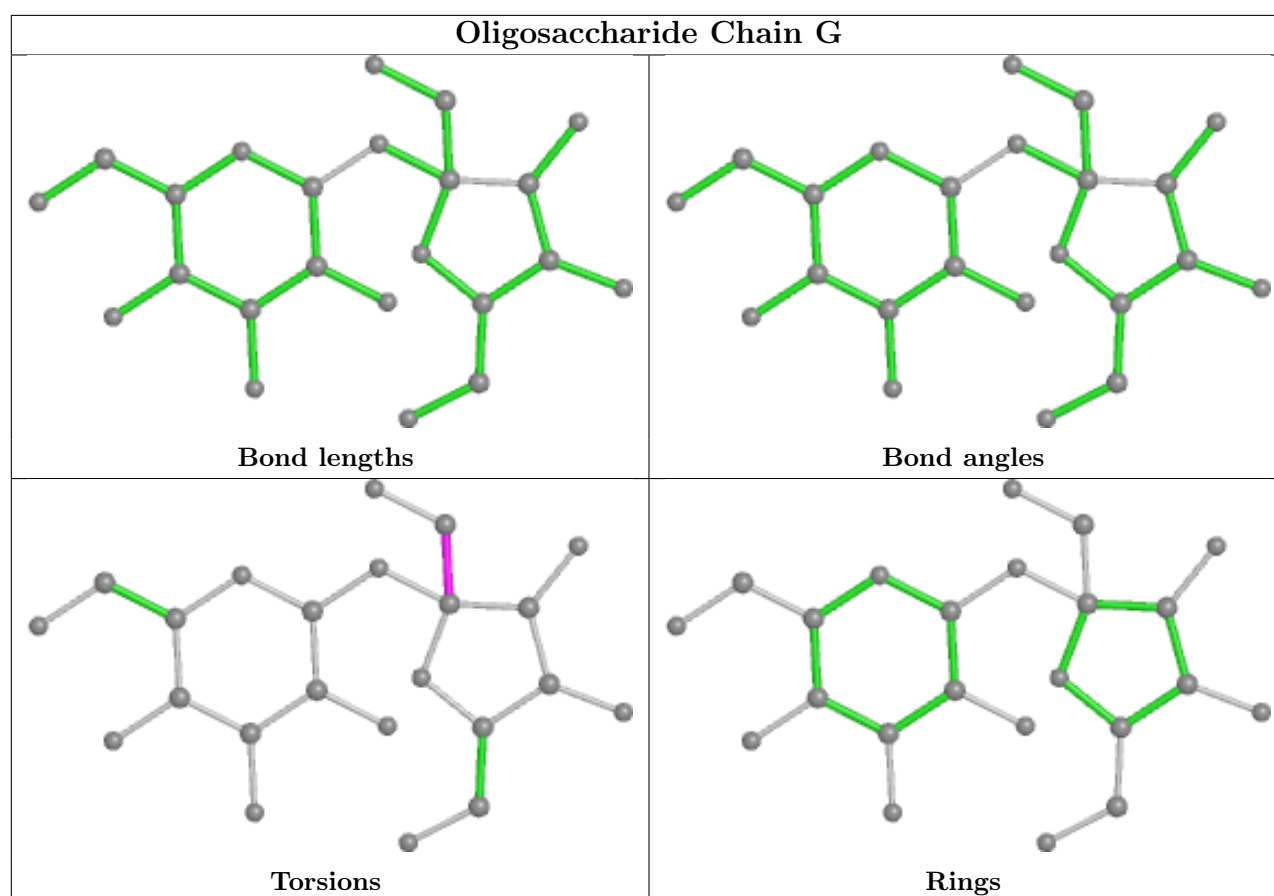
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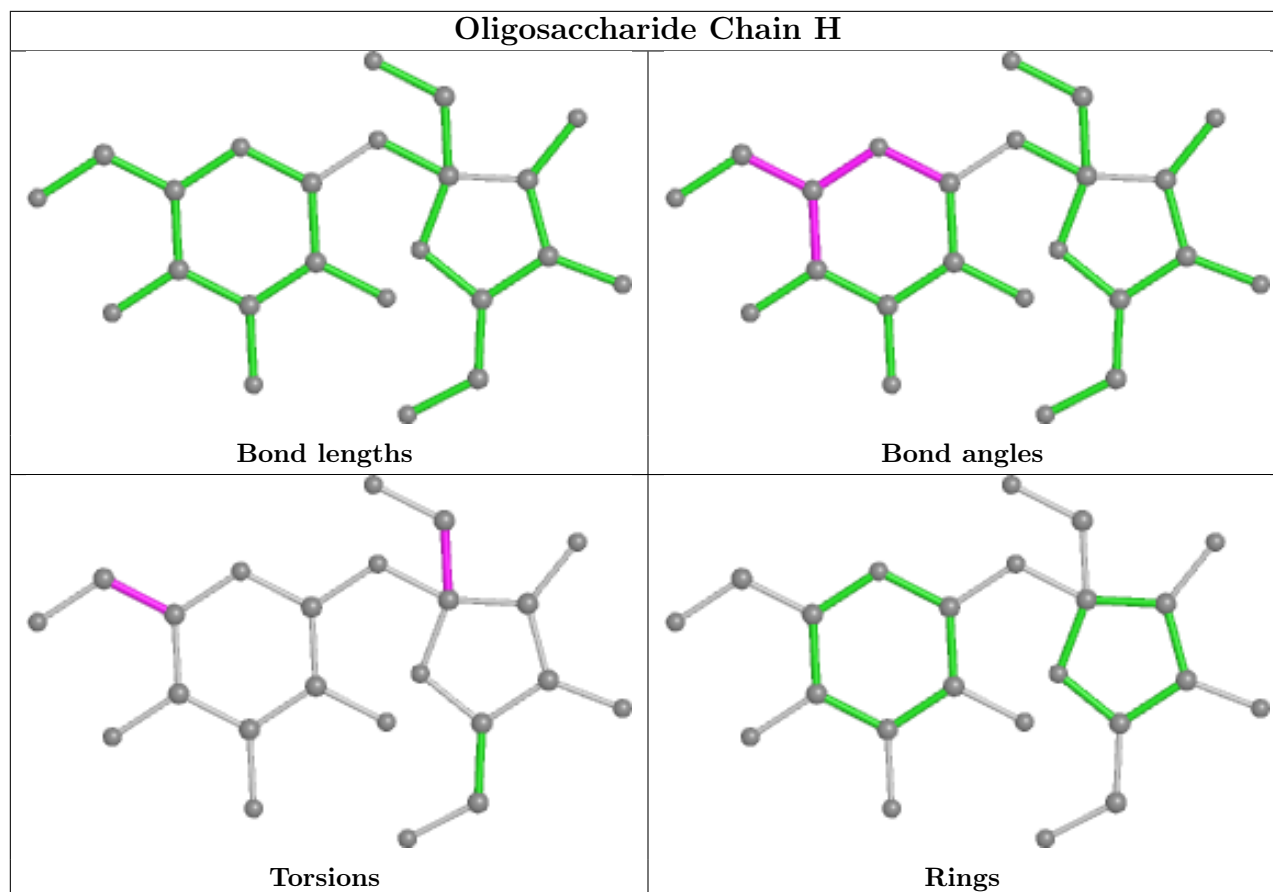
Mol	Chain	Res	Type	Atoms
5	H	2	FRU	O1-C1-C2-O2
5	G	2	FRU	O1-C1-C2-O5
5	H	2	FRU	O1-C1-C2-C3
5	H	1	GLC	C4-C5-C6-O6
5	H	2	FRU	O1-C1-C2-O5

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.





5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 4 are monoatomic - leaving 5 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	$\# Z > 2$
9	SO4	D	502	-	4,4,4	0.14	0	6,6,6	0.05	0
8	EDO	D	503	-	3,3,3	0.45	0	2,2,2	0.21	0
8	EDO	B	502	-	3,3,3	0.37	0	2,2,2	0.43	0
7	DTP	C	601	6	26,32,32	1.50	6 (23%)	30,50,50	1.51	5 (16%)
7	DTP	A	603	6	26,32,32	1.50	6 (23%)	30,50,50	1.51	4 (13%)

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
8	EDO	B	502	-	-	0/1/1/1	-
7	DTP	C	601	6	-	1/18/34/34	0/3/3/3
8	EDO	D	503	-	-	1/1/1/1	-
7	DTP	A	603	6	-	3/18/34/34	0/3/3/3

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	C	601	DTP	C6-N6	3.53	1.46	1.34
7	A	603	DTP	C6-N6	3.52	1.46	1.34
7	A	603	DTP	C5'-C4'	-3.13	1.41	1.51
7	C	601	DTP	C5'-C4'	-3.12	1.41	1.51
7	A	603	DTP	C2'-C3'	-2.61	1.46	1.52
7	C	601	DTP	C2'-C3'	-2.60	1.46	1.52
7	C	601	DTP	O5'-C5'	-2.53	1.35	1.44
7	A	603	DTP	O5'-C5'	-2.53	1.35	1.44
7	C	601	DTP	O3'-C3'	-2.36	1.38	1.43
7	A	603	DTP	O3'-C3'	-2.34	1.38	1.43
7	C	601	DTP	C3'-C4'	-2.19	1.47	1.53
7	A	603	DTP	C3'-C4'	-2.18	1.47	1.53

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	C	601	DTP	N3-C2-N1	-4.38	121.83	128.68
7	A	603	DTP	N3-C2-N1	-4.33	121.92	128.68
7	C	601	DTP	PB-O3B-PG	-3.81	119.74	132.83
7	A	603	DTP	PB-O3B-PG	-3.70	120.12	132.83
7	A	603	DTP	PA-O3A-PB	-3.55	120.65	132.83
7	C	601	DTP	PA-O3A-PB	-3.35	121.32	132.83
7	A	603	DTP	O5'-C5'-C4'	2.30	116.92	108.99
7	C	601	DTP	O5'-C5'-C4'	2.30	116.91	108.99
7	C	601	DTP	C4-C5-N7	-2.11	107.20	109.40

There are no chirality outliers.

All (5) torsion outliers are listed below:

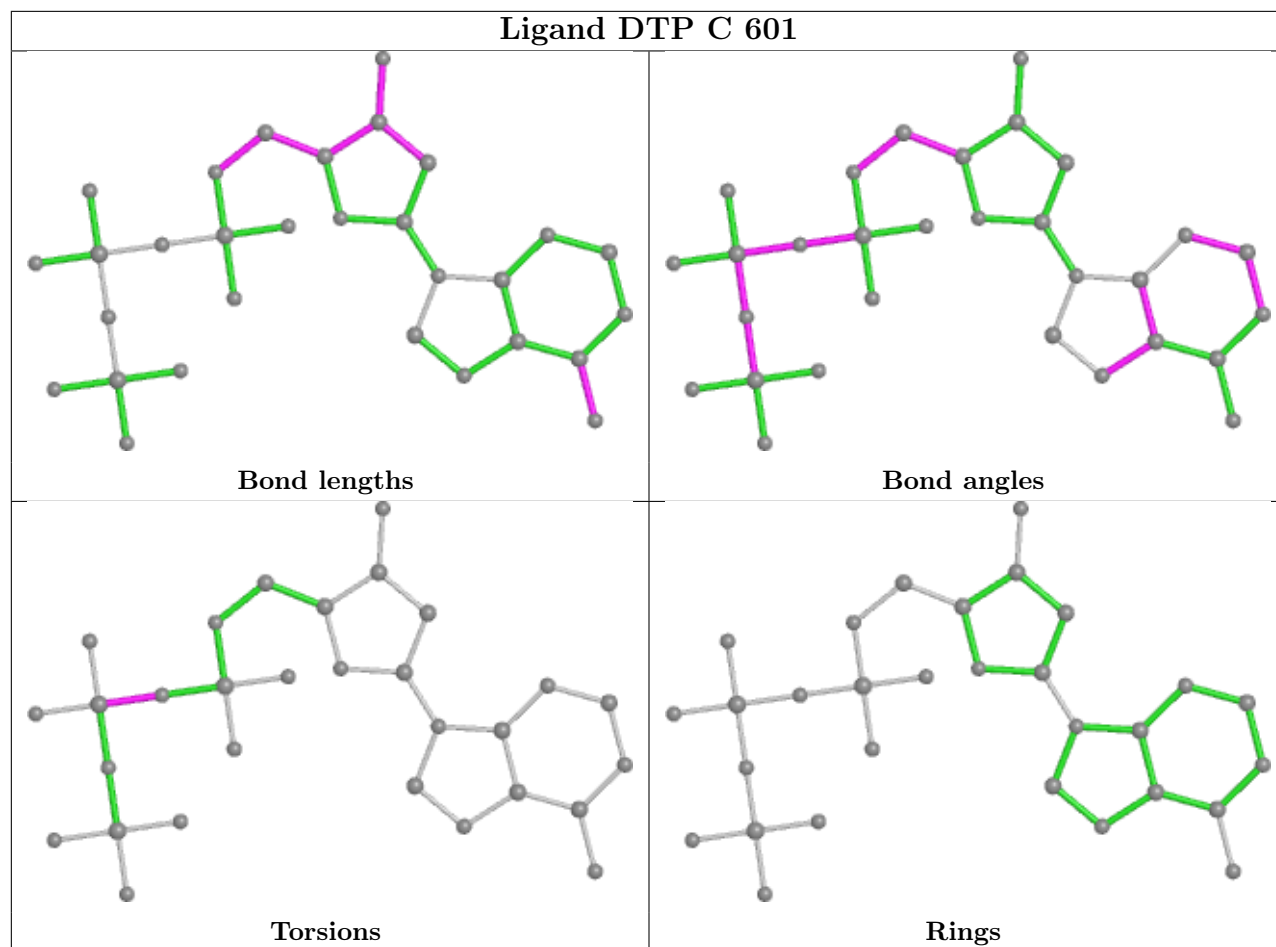
Mol	Chain	Res	Type	Atoms
7	C	601	DTP	PA-O3A-PB-O2B
8	D	503	EDO	O1-C1-C2-O2
7	A	603	DTP	PA-O3A-PB-O2B
7	A	603	DTP	PG-O3B-PB-O1B
7	A	603	DTP	PA-O3A-PB-O1B

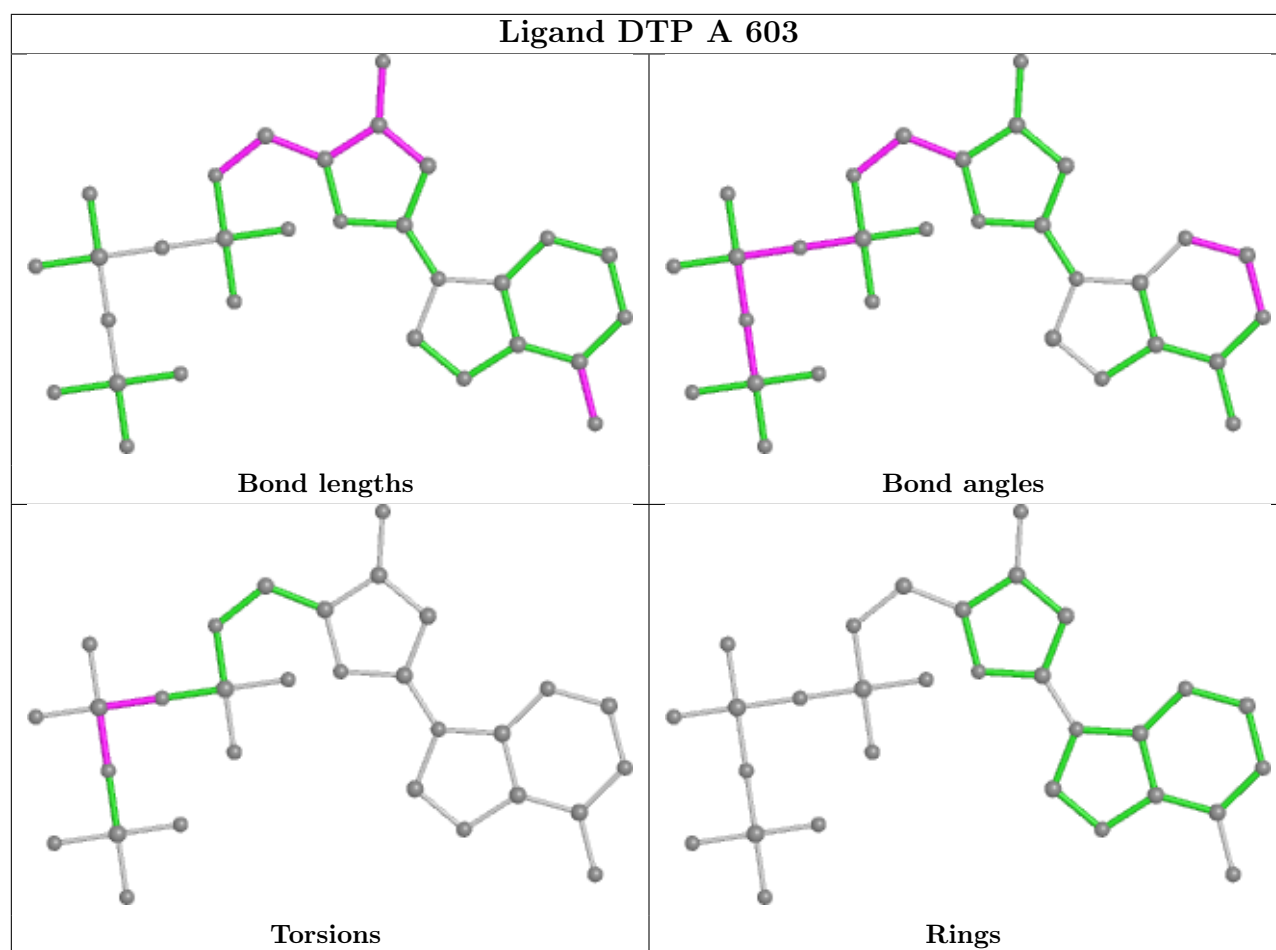
There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	D	503	EDO	3	0
7	A	603	DTP	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	555/556 (99%)	0.73	82 (14%) 2 2	33, 84, 169, 230	0
1	C	555/556 (99%)	0.71	81 (14%) 2 3	36, 91, 160, 232	0
2	B	412/428 (96%)	0.42	29 (7%) 16 19	35, 70, 124, 162	0
2	D	412/428 (96%)	0.62	48 (11%) 4 6	37, 82, 142, 174	0
3	E	24/27 (88%)	1.01	4 (16%) 1 1	76, 116, 193, 231	0
3	T	24/27 (88%)	1.18	6 (25%) 0 0	76, 110, 203, 249	0
4	F	19/21 (90%)	0.55	1 (5%) 26 31	62, 99, 182, 186	0
4	P	19/21 (90%)	0.58	2 (10%) 6 8	69, 100, 158, 180	0
All	All	2020/2064 (97%)	0.65	253 (12%) 3 5	33, 81, 156, 249	0

All (253) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	214	LEU	14.6
3	T	702	DT	10.6
2	B	4	PRO	10.4
2	D	214	LEU	10.3
1	A	135	ILE	9.8
1	C	252	TRP	9.4
2	D	4	PRO	9.2
2	B	5	ILE	8.8
2	D	360	ALA	8.8
1	A	133	PRO	7.9
1	A	286	THR	7.8
1	A	131	THR	7.7
1	A	67	ASP	7.7
1	A	141	GLY	7.3
1	A	288	ALA	7.1
1	C	135	ILE	7.0

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Mol	Chain	Res	Type	RSRZ
1	C	246	LEU	6.7
1	C	292	VAL	6.3
1	A	140	PRO	6.1
1	C	142	ILE	6.0
1	A	138	GLU	5.9
1	A	142	ILE	5.8
1	A	285	GLY	5.7
1	A	293	ILE	5.6
1	A	254	VAL	5.6
1	C	26	LEU	5.6
1	C	67	ASP	5.6
1	A	50	ILE	5.5
1	A	52	PRO	5.5
3	E	702	DT	5.4
1	C	133	PRO	5.4
4	P	803	DC	5.3
2	B	89	GLU	5.3
1	C	293	ILE	5.3
2	D	358	ARG	5.3
2	D	90	VAL	5.2
1	A	71	TRP	5.2
1	C	295	LEU	5.1
1	A	54	ASN	5.1
1	C	24	TRP	5.1
1	A	26	LEU	5.1
1	A	30	LYS	5.0
1	C	251	SER	5.0
2	D	361	HIS	5.0
1	A	292	VAL	5.0
1	C	247	PRO	4.8
1	A	287	LYS	4.8
3	T	703	DG	4.7
2	D	318	TYR	4.7
1	C	303	LEU	4.7
1	C	138	GLU	4.6
1	C	249	LYS	4.6
1	C	37	ILE	4.6
2	D	85	GLN	4.6
1	C	287	LYS	4.5
3	T	725	DG	4.5
1	A	289	LEU	4.4
1	A	139	THR	4.4

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Mol	Chain	Res	Type	RSRZ
1	A	66	LYS	4.4
1	C	32	LYS	4.4
2	D	89	GLU	4.3
2	D	422	LEU	4.3
3	E	703	DG	4.3
2	D	428	GLN	4.2
2	B	215	THR	4.2
1	C	288	ALA	4.1
1	A	24	TRP	4.1
1	C	253	THR	4.1
1	C	74	LEU	4.1
1	C	71	TRP	4.1
1	A	284	ARG	4.1
3	E	725	DG	4.1
1	C	132	ILE	4.0
1	C	289	LEU	4.0
1	A	27	THR	4.0
1	C	294	PRO	4.0
1	A	16	MET	4.0
1	A	294	PRO	4.0
2	D	95	PRO	3.9
1	C	134	SER	3.9
2	D	88	TRP	3.9
2	B	217	PRO	3.9
2	D	67	ASP	3.9
2	B	184	MET	3.8
2	D	212	TRP	3.7
1	C	31	ILE	3.7
1	C	28	GLU	3.7
1	C	136	ASN	3.7
1	C	299	ALA	3.7
1	A	295	LEU	3.6
1	A	69	THR	3.6
1	A	37	ILE	3.6
1	A	32	LYS	3.6
1	A	55	PRO	3.6
1	A	56	TYR	3.6
1	A	143	ARG	3.6
1	C	286	THR	3.6
2	D	409	THR	3.6
1	A	132	ILE	3.6
1	C	15	GLY	3.5

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Mol	Chain	Res	Type	RSRZ
1	A	60	VAL	3.5
1	C	250	ASP	3.5
2	B	85	GLN	3.5
1	A	73	LYS	3.5
2	D	184	MET	3.5
2	D	317	VAL	3.5
1	A	130	PHE	3.4
1	C	35	VAL	3.4
1	A	29	GLU	3.4
2	D	232	TYR	3.4
1	A	252	TRP	3.4
2	D	215	THR	3.4
2	D	173	LYS	3.4
1	C	68	SER	3.3
1	C	131	THR	3.3
1	A	61	PHE	3.2
1	C	54	ASN	3.2
2	D	425	LEU	3.2
1	A	134	SER	3.2
2	D	193	LEU	3.2
1	C	14	PRO	3.1
1	A	282	LEU	3.1
2	D	6	GLU	3.1
1	C	137	ASN	3.1
1	A	246	LEU	3.0
1	C	34	LEU	3.0
1	A	136	ASN	3.0
2	D	5	ILE	3.0
1	A	20	LYS	3.0
1	A	70	LYS	3.0
1	A	34	LEU	3.0
1	C	25	PRO	3.0
2	D	197	GLN	3.0
1	A	21	VAL	2.9
1	C	70	LYS	2.9
2	B	410	TRP	2.9
2	D	177	ASP	2.9
2	B	92	LEU	2.9
1	C	141	GLY	2.9
1	C	61	PHE	2.9
1	C	69	THR	2.8
1	A	129	ALA	2.8

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Mol	Chain	Res	Type	RSRZ
2	B	232	TYR	2.8
1	A	259	LYS	2.8
2	B	90	VAL	2.8
2	D	410	TRP	2.8
2	B	186	ASP	2.8
2	B	174	GLN	2.7
2	D	209	LEU	2.7
1	C	95	PRO	2.7
2	D	423	VAL	2.7
1	C	260	LEU	2.7
2	D	92	LEU	2.7
1	C	66	LYS	2.7
1	C	144	TYR	2.7
2	D	357	MET	2.7
1	C	290	THR	2.7
2	D	238	LYS	2.7
1	C	27	THR	2.7
2	B	178	ILE	2.7
1	A	51	GLY	2.7
1	C	20	LYS	2.6
2	B	409	THR	2.6
2	B	213	GLY	2.6
1	C	223	LYS	2.6
1	A	144	TYR	2.6
1	C	255	ASN	2.6
1	C	16	MET	2.6
1	A	74	LEU	2.6
1	A	17	ASP	2.6
1	A	25	PRO	2.6
1	C	219	LYS	2.6
1	C	308	GLU	2.6
1	A	283	LEU	2.5
2	D	178	ILE	2.5
1	C	94	ILE	2.5
1	C	30	LYS	2.5
2	B	88	TRP	2.5
1	C	36	GLU	2.5
1	C	202	ILE	2.5
2	D	66	LYS	2.5
1	C	448	ARG	2.4
4	F	803	DC	2.4
1	C	312	GLU	2.4

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Mol	Chain	Res	Type	RSRZ
2	B	171	PHE	2.4
1	C	109	LEU	2.4
1	A	64	LYS	2.4
1	A	290	THR	2.4
2	D	186	ASP	2.4
2	D	152	GLY	2.4
2	D	213	GLY	2.4
1	C	254	VAL	2.3
3	E	716	DA	2.3
2	B	173	LYS	2.3
2	B	231	GLY	2.3
2	D	183	TYR	2.3
1	A	448	ARG	2.3
2	D	124	PHE	2.3
1	A	554	ALA	2.3
2	D	69	THR	2.3
1	C	145	GLN	2.3
3	T	724	DT	2.3
2	B	411	ILE	2.3
1	A	68	SER	2.3
1	A	28	GLU	2.3
2	D	376	THR	2.3
1	A	75	VAL	2.2
2	B	87	PHE	2.2
1	A	452	LEU	2.2
1	C	449	GLU	2.2
2	D	421	PRO	2.2
1	C	13	LYS	2.2
1	C	257	ILE	2.2
1	C	193	LEU	2.2
2	B	154	LYS	2.2
2	D	94	ILE	2.2
1	A	248	GLU	2.2
1	C	248	GLU	2.2
1	C	310	LEU	2.2
2	B	310	LEU	2.2
2	B	156	SER	2.2
1	C	184	MET	2.2
1	C	198	HIS	2.2
1	C	97	PRO	2.2
1	A	251	SER	2.2
1	A	58	THR	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	14	PRO	2.1
1	C	22	LYS	2.1
1	A	303	LEU	2.1
1	A	145	GLN	2.1
1	A	36	GLU	2.1
2	D	104	LYS	2.1
2	B	408	ALA	2.1
1	A	77	PHE	2.1
1	A	193	LEU	2.1
1	A	57	ASN	2.1
1	C	283	LEU	2.1
1	C	49	LYS	2.1
2	D	408	ALA	2.1
2	B	183	TYR	2.1
2	B	185	ASP	2.1
1	A	219	LYS	2.1
1	A	253	THR	2.0
2	D	179	VAL	2.0
3	T	704	DG	2.0
1	A	22	LYS	2.0
1	A	49	LYS	2.0
2	D	354	TYR	2.0
3	T	716	DA	2.0
4	P	804	DA	2.0
1	C	64	LYS	2.0
1	C	210	LEU	2.0
1	A	53	GLU	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
4	MRG	F	817	26/27	0.85	0.16	105,119,149,157	0
4	MRG	P	817	26/27	0.90	0.14	90,114,145,155	0

6.3 Carbohydrates

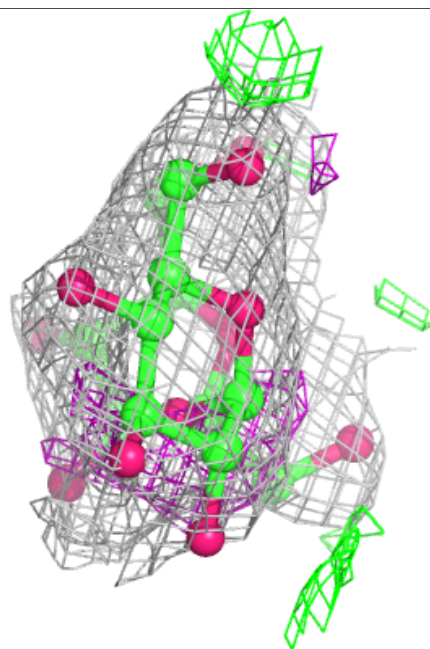
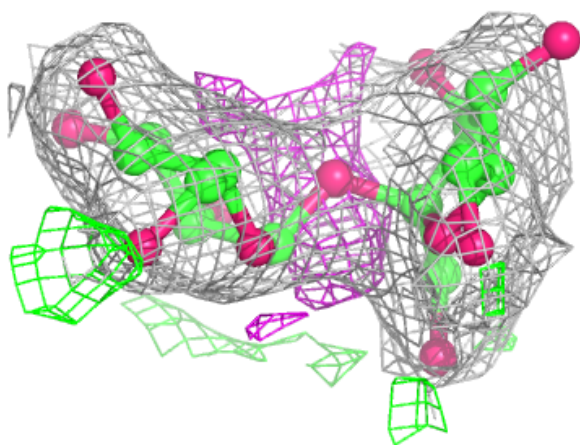
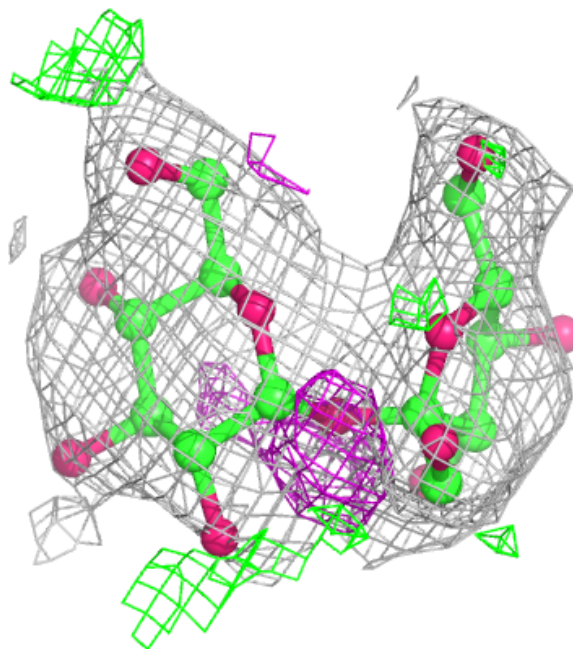
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

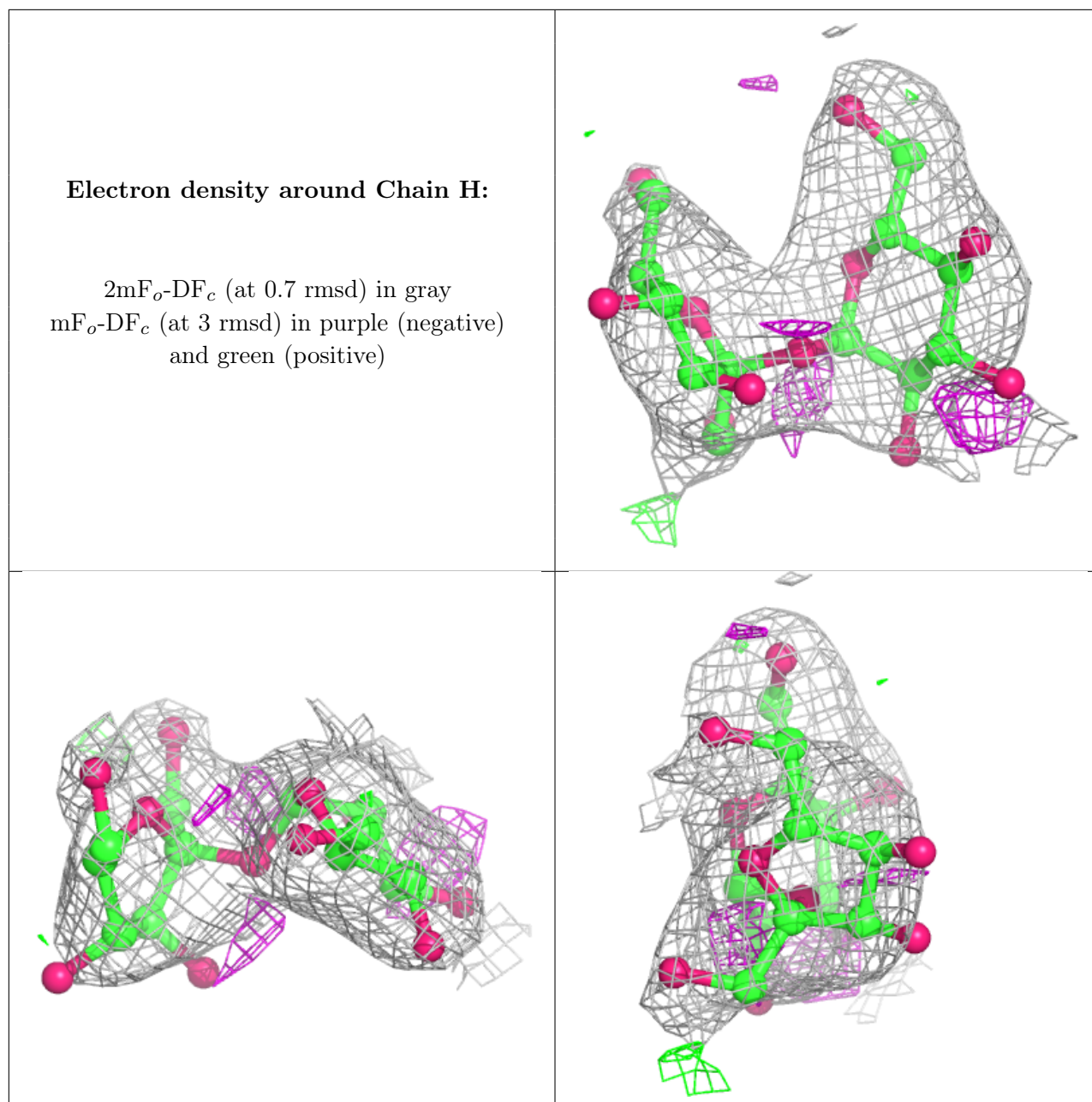
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	GLC	H	1	11/12	0.79	0.21	79,102,114,114	0
5	FRU	G	2	12/12	0.84	0.18	102,114,119,121	0
5	FRU	H	2	12/12	0.87	0.20	95,115,120,123	0
5	GLC	G	1	11/12	0.89	0.18	69,91,106,110	0

The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain G:

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





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.

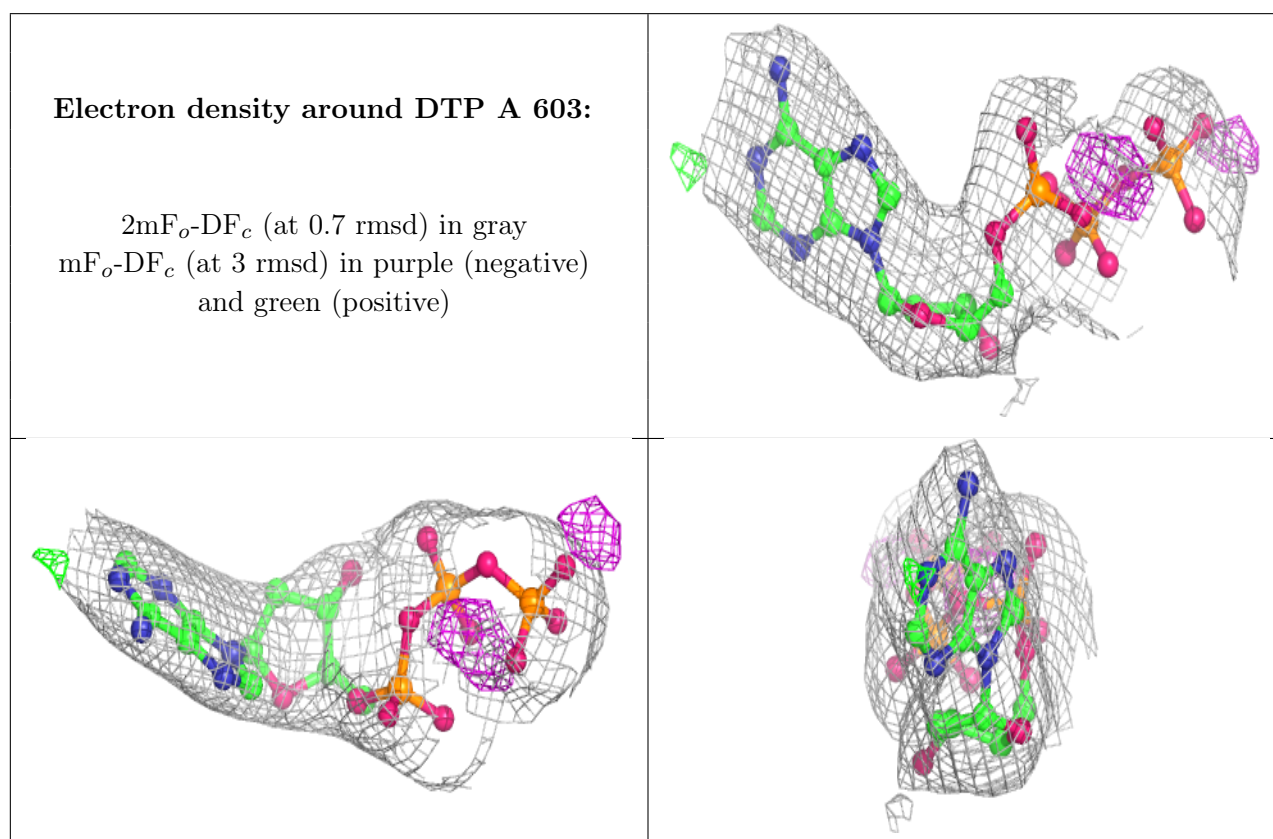
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	MG	A	602	1/1	0.55	0.39	85,85,85,85	0
9	SO4	D	502	5/5	0.71	0.31	160,162,164,165	0
6	MG	A	601	1/1	0.84	0.04	95,95,95,95	0

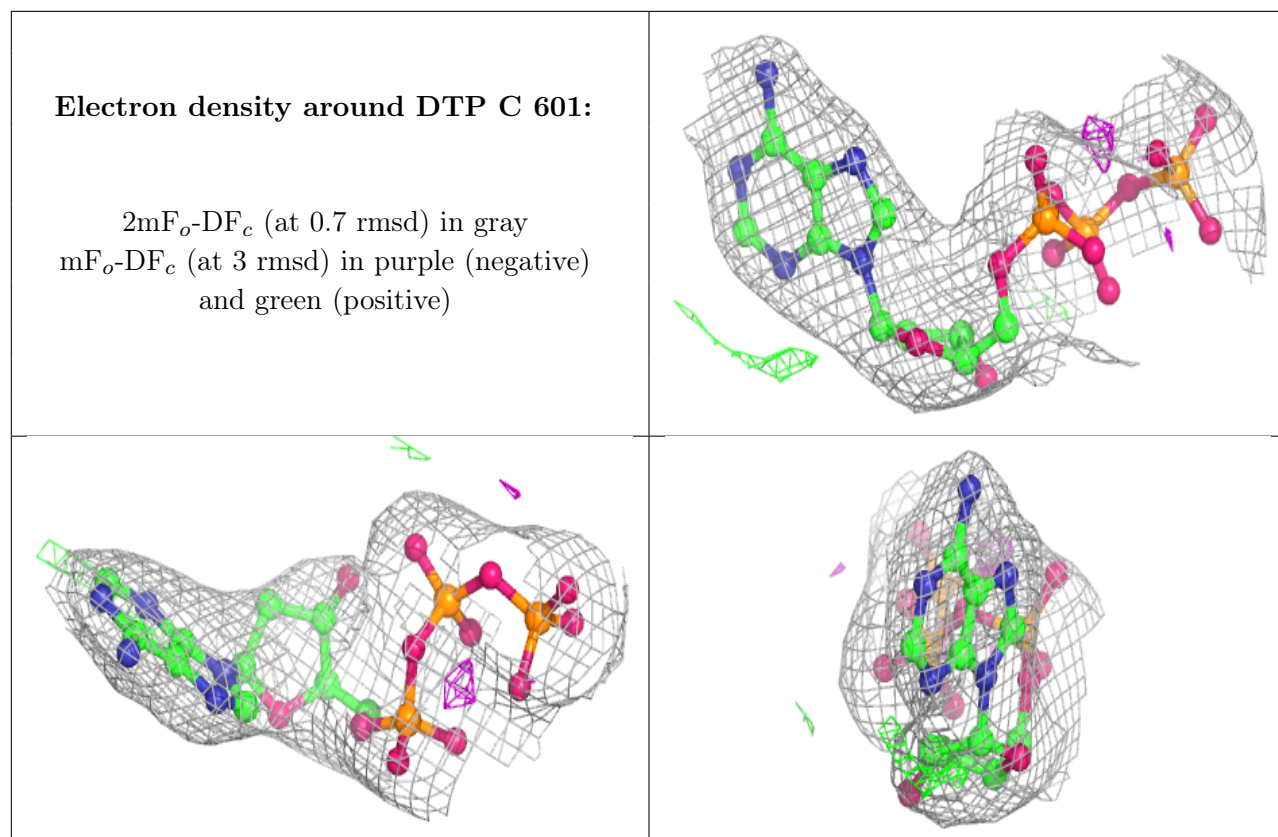
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
8	EDO	D	503	4/4	0.88	0.25	70,74,74,77	0
6	MG	C	602	1/1	0.90	0.07	80,80,80,80	0
6	MG	C	603	1/1	0.92	0.25	96,96,96,96	0
7	DTP	A	603	30/30	0.92	0.13	73,104,136,139	0
8	EDO	B	502	4/4	0.94	0.15	59,66,71,79	0
7	DTP	C	601	30/30	0.95	0.15	76,97,131,138	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.





6.5 Other polymers [i](#)

There are no such residues in this entry.