



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:19 AM UTC

PDB ID : 5TXM / pdb_00005txm
Title : STRUCTURE OF HIV-1 REVERSE TRANSCRIPTASE (RT) TERNARY COMPLEX WITH A DOUBLE STRANDED DNA AND AN INCOMING DDATP
Authors : Das, K.; Martinez, S.M.; Arnold, E.
Deposited on : 2016-11-17
Resolution : 2.70 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 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

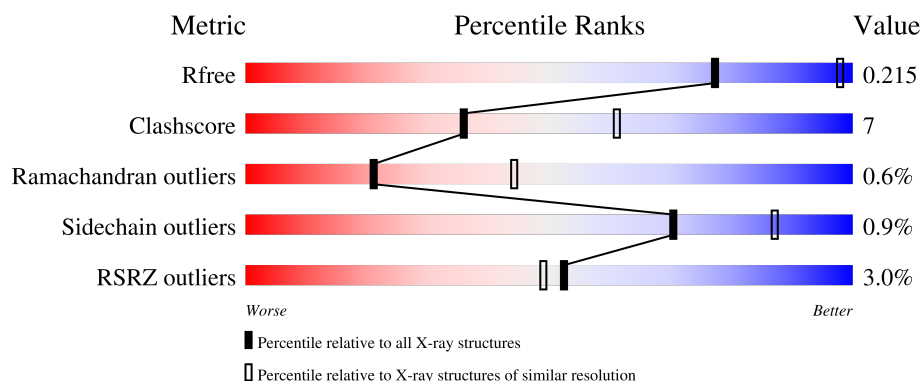
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

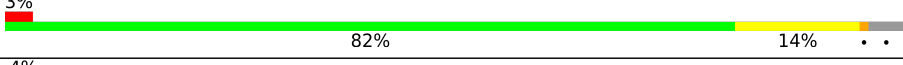
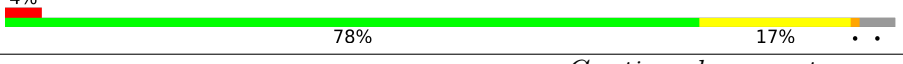
The reported resolution of this entry is 2.70 Å.

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}	180053	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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 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	557	
1	C	557	
2	B	428	
2	D	428	

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Mol	Chain	Length	Quality of chain
3	E	27	
3	T	27	
4	F	21	
4	P	21	
5	G	2	
5	H	2	

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 17950 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
1	A	556	Total	C	N	O	S	0	0	0
			4516	2923	752	833	8			
1	C	556	Total	C	N	O	S	0	0	0
			4516	2923	752	833	8			

There are 10 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	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	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
2	B	411	Total	C	N	O	S	0	0	0
			3392	2207	562	616	7			
2	D	412	Total	C	N	O	S	0	0	0
			3400	2212	563	619	6			

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

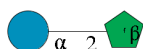
- Molecule 3 is a DNA chain called DNA (5'-D(*TP*GP*GP*TP*CP*GP*GP*CP*GP*CP*CP*GP*AP*AP*CP*AP*GP*GP*GP*AP*CP*TP*G)-3').

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

- 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
4	P	20	Total	C	N	O	P	S	0	0	0
			407	195	72	120	19	1			
4	F	20	Total	C	N	O	P	S	0	0	0
			407	195	72	120	19	1			

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



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

- Molecule 6 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

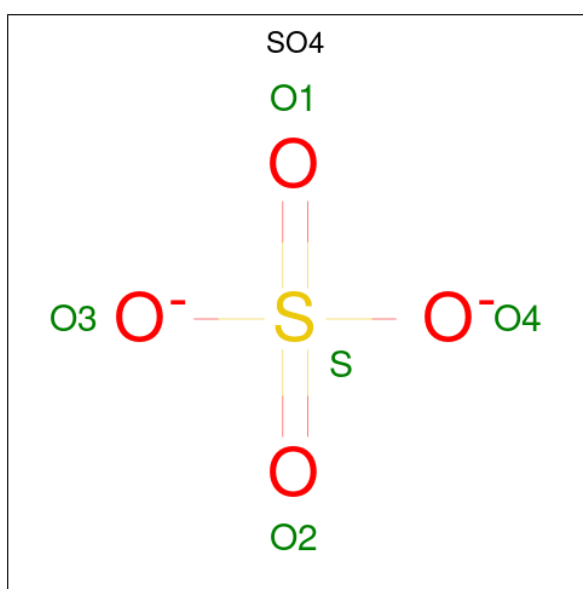
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	2	Total	Mg	0	0
			2	2		
6	C	2	Total	Mg	0	0
			2	2		

- Molecule 7 is 2',3'-dideoxyadenosine triphosphate (CCD ID: DDS) (formula: C₁₀H₁₆N₅O₁₁P₃).

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

- Molecule 9 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	52	Total	O	0	0
			52	52		
10	B	48	Total	O	0	0
			48	48		
10	T	2	Total	O	0	0
			2	2		
10	P	1	Total	O	0	0
			1	1		

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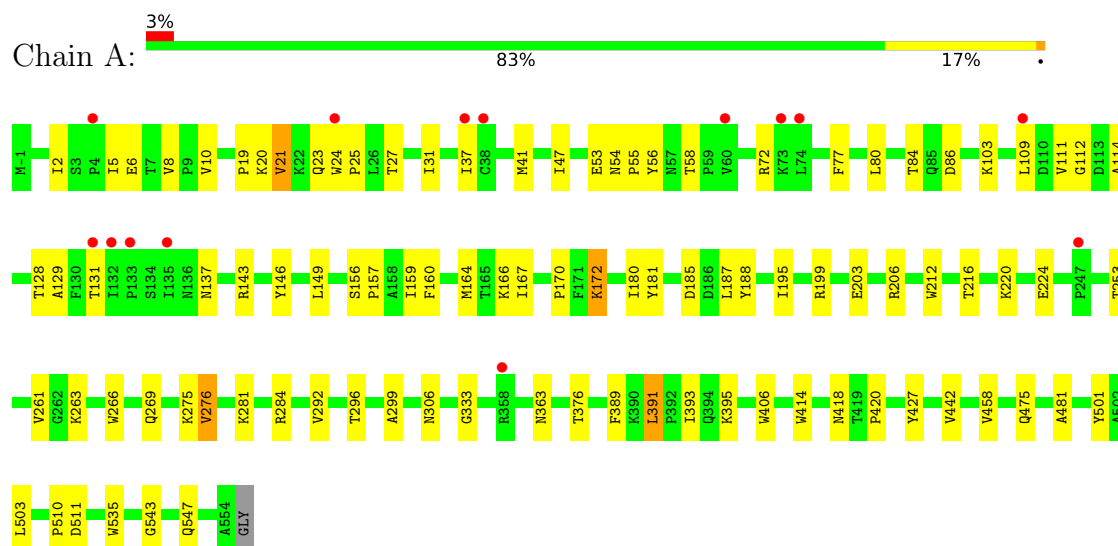
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	C	50	Total	O	0	0
			50	50		
10	D	34	Total	O	0	0
			34	34		

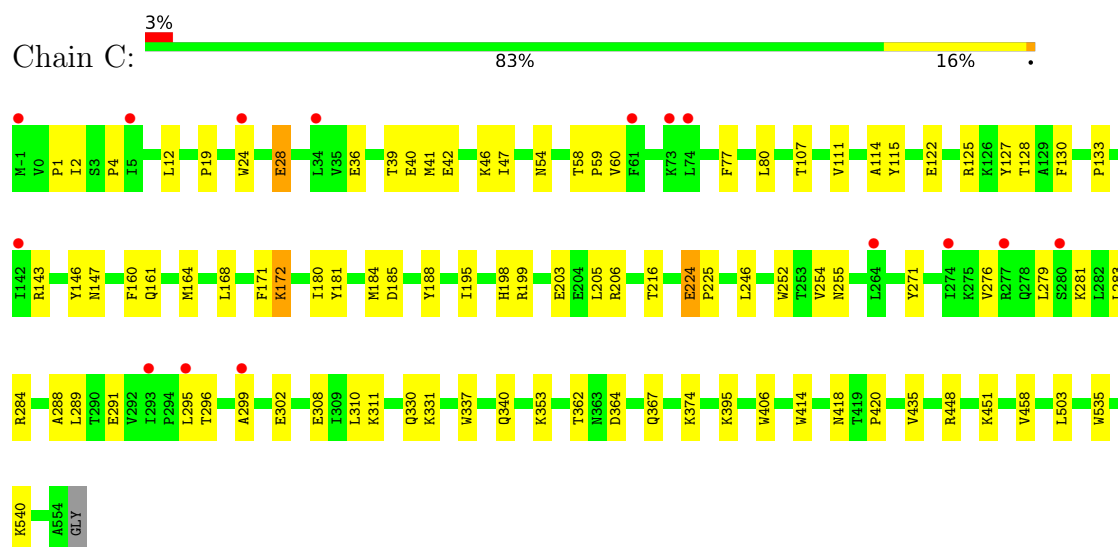
3 Residue-property plots [i](#)

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.

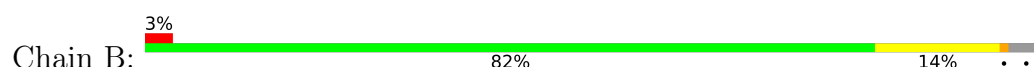
- Molecule 1: HIV-1 Reverse Transcriptase P66 subunit

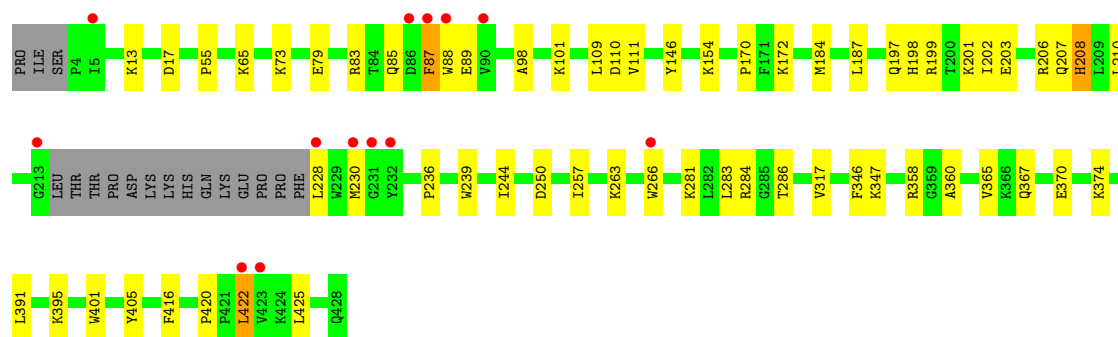


- Molecule 1: HIV-1 Reverse Transcriptase P66 subunit

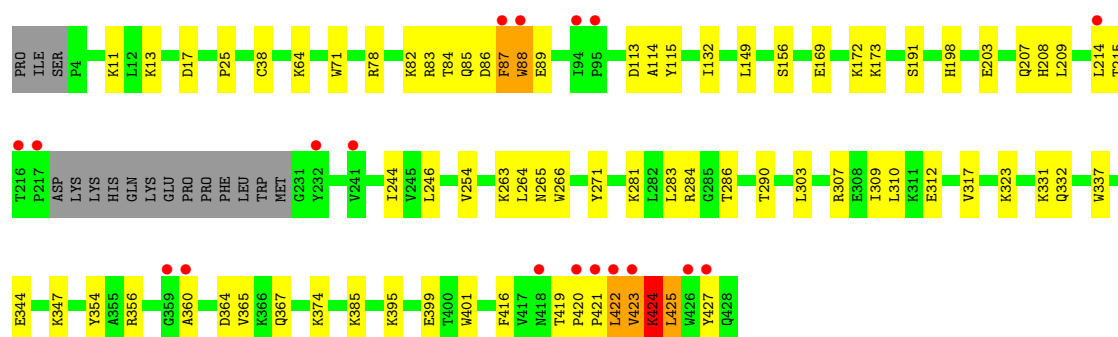
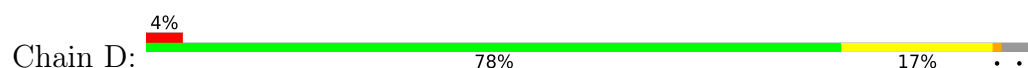


- Molecule 2: HIV-1 Reverse Transcriptase P51 subunit





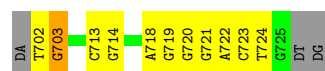
● Molecule 2: HIV-1 Reverse Transcriptase P51 subunit



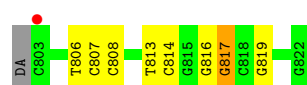
● Molecule 3: DNA (5'-D(*TP*GP*GP*TP*CP*GP*GP*CP*GP*CP*CP*CP*GP*AP*AP*CP*AP*GP*GP*GP*AP*CP*TP*G)-3')



● 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')



- 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:  48% 43% 5% 5%

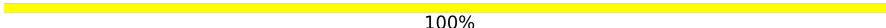


- Molecule 5: beta-D-fructofuranose-(2-1)-alpha-D-glucopyranose

Chain G:  50% 50%



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

Chain H:  100%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	90.34Å 133.95Å 139.40Å 90.00° 97.71° 90.00°	Depositor
Resolution (Å)	46.05 – 2.70 46.05 – 2.70	Depositor EDS
% Data completeness (in resolution range)	97.4 (46.05-2.70) 97.3 (46.05-2.70)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.79 (at 2.61Å)	Xtriage
Refinement program	PHENIX (1.10_2155: ???)	Depositor
R, R_{free}	0.184 , 0.223 (Not available) , 0.215	Depositor DCC
R_{free} test set	2864 reflections (2.84%)	wwPDB-VP
Wilson B-factor (Å ²)	61.9	Xtriage
Anisotropy	0.160	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 76.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	17950	wwPDB-VP
Average B, all atoms (Å ²)	94.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.69% 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: DDS, GLC, SO4, MRG, FRU, EDO, 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	0.26	1/4634 (0.0%)	0.45	2/6297 (0.0%)
1	C	0.25	1/4634 (0.0%)	0.46	2/6297 (0.0%)
2	B	0.30	2/3488 (0.1%)	0.47	0/4736
2	D	0.29	1/3497 (0.0%)	0.63	3/4751 (0.1%)
3	E	0.40	0/555	0.70	1/856 (0.1%)
3	T	0.24	0/555	0.42	1/856 (0.1%)
4	F	0.35	0/424	0.70	0/649
4	P	0.28	0/424	0.48	0/649
All	All	0.28	5/18211 (0.0%)	0.51	9/25091 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
2	D	0	2

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	391	LEU	C-N	7.67	1.51	1.33
1	A	172	LYS	CE-NZ	-5.37	1.33	1.49
1	C	172	LYS	CE-NZ	-5.35	1.33	1.49
2	D	172	LYS	CE-NZ	-5.31	1.33	1.49
2	B	172	LYS	CE-NZ	-5.27	1.33	1.49

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	89	GLU	N-CA-C	9.26	121.14	111.14
1	C	224	GLU	CA-C-N	8.24	125.59	119.66
1	C	224	GLU	C-N-CA	8.24	125.59	119.66
2	D	424	LYS	CA-C-N	6.92	129.85	120.44
2	D	424	LYS	C-N-CA	6.92	129.85	120.44
1	A	224	GLU	CA-C-N	5.55	123.66	119.66
1	A	224	GLU	C-N-CA	5.55	123.66	119.66
3	E	703	DG	N9-C1'-C2'	5.17	121.25	113.50
3	T	722	DA	C5'-C4'-O4'	5.09	117.03	109.40

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	D	424	LYS	Peptide
2	D	88	TRP	Mainchain

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4516	0	4572	61	0
1	C	4516	0	4572	58	0
2	B	3392	0	3423	45	0
2	D	3400	0	3433	55	0
3	E	494	0	269	13	0
3	T	494	0	269	5	0
4	F	407	0	229	17	0
4	P	407	0	229	6	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	29	0	12	4	0
7	C	29	0	12	2	0
8	A	4	0	6	1	0
8	B	8	0	12	1	0
8	D	8	0	12	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	T	4	0	6	0	0
9	B	5	0	0	0	0
10	A	52	0	0	0	0
10	B	48	0	0	0	0
10	C	50	0	0	0	0
10	D	34	0	0	0	0
10	P	1	0	0	0	0
10	T	2	0	0	0	0
All	All	17950	0	17098	253	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 (253) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:419:THR:O	2:D:421:PRO:HD3	1.67	0.94
1:A:172:LYS:HE2	1:A:180:ILE:HB	1.49	0.92
2:B:109:LEU:HD23	2:B:230:MET:HB2	1.57	0.86
3:E:722:DA:H2''	3:E:723:DC:H5''	1.59	0.85
2:D:266:TRP:CZ3	2:D:425:LEU:HG	2.13	0.83
2:D:209:LEU:HD22	2:D:214:LEU:HD23	1.59	0.82
4:F:804:DA:H2'	4:F:805:DG:C8	2.18	0.78
3:E:719:DG:H2''	3:E:720:DG:C8	2.19	0.78
1:C:448:ARG:O	1:C:451:LYS:NZ	2.16	0.78
1:A:543:GLY:O	1:A:547:GLN:NE2	2.17	0.76
2:D:244:ILE:HG12	2:D:425:LEU:HD11	1.66	0.76
2:B:266:TRP:CD1	2:B:425:LEU:HD22	2.22	0.74
2:D:421:PRO:HG2	2:D:423:VAL:HB	1.70	0.73
1:A:195:ILE:HD11	1:A:199:ARG:HH21	1.54	0.73
2:B:110:ASP:H	2:B:230:MET:HG3	1.53	0.73
1:C:203:GLU:OE2	1:C:206:ARG:NH1	2.22	0.72
1:A:128:THR:HG21	1:A:146:TYR:HB2	1.70	0.72
2:B:79:GLU:OE2	2:B:83:ARG:NH2	2.23	0.70
2:D:323:LYS:NZ	2:D:344:GLU:OE2	2.26	0.68
7:A:603:DDS:H5'	7:A:603:DDS:H8	1.77	0.67
2:B:365:VAL:HG11	2:B:401:TRP:HB2	1.78	0.67
2:B:206:ARG:NH2	2:B:230:MET:O	2.28	0.66
3:E:718:DA:H2''	3:E:719:DG:C8	2.31	0.66
4:F:805:DG:H4'	4:F:806:DT:OP1	1.95	0.66
3:E:723:DC:H2'	3:E:724:DT:H72	1.79	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:C:601:DDS:H5'	7:C:601:DDS:H8	1.78	0.65
2:B:73:LYS:NZ	2:B:146:TYR:OH	2.29	0.65
2:D:254:VAL:HG13	2:D:283:LEU:HD22	1.80	0.64
1:A:296:THR:HG23	1:A:299:ALA:H	1.63	0.64
1:A:56:TYR:O	1:A:143:ARG:NH2	2.31	0.63
3:E:718:DA:H2''	3:E:719:DG:H8	1.62	0.63
1:A:164:MET:HE2	1:A:187:LEU:HD11	1.80	0.63
1:C:54:ASN:HB3	1:C:143:ARG:HH12	1.64	0.62
2:D:263:LYS:HE3	2:D:425:LEU:HA	1.80	0.62
1:A:54:ASN:HB3	1:A:143:ARG:HH12	1.65	0.61
2:D:246:LEU:HD11	2:D:264:LEU:HD21	1.83	0.61
2:D:244:ILE:HD12	2:D:271:TYR:HE2	1.67	0.60
3:E:702:DT:H3'	3:E:703:DG:H8	1.67	0.60
1:A:261:VAL:HG13	1:A:276:VAL:HG11	1.82	0.59
1:A:253:THR:HA	1:A:292:VAL:HA	1.84	0.59
2:D:421:PRO:HD2	2:D:423:VAL:HG12	1.83	0.59
2:B:346:PHE:H	8:B:504:EDO:H11	1.68	0.59
1:C:246:LEU:HD11	1:C:310:LEU:HD22	1.85	0.58
1:C:503:LEU:HD22	1:C:535:TRP:HB2	1.85	0.58
7:A:603:DDS:H5'	7:A:603:DDS:C8	2.35	0.57
2:B:199:ARG:HA	2:B:202:ILE:HD12	1.85	0.57
2:B:370:GLU:HG3	2:B:374:LYS:HD2	1.86	0.57
2:B:203:GLU:O	2:B:207:GLN:HG2	2.04	0.56
1:C:418:ASN:O	1:C:420:PRO:HD3	2.05	0.56
1:A:503:LEU:HD22	1:A:535:TRP:HB2	1.88	0.56
3:E:721:DG:C8	3:E:721:DG:H5'	2.40	0.56
4:F:817:MRG:N3	4:F:817:MRG:H221	2.21	0.56
1:A:37:ILE:HG22	1:A:41:MET:HE2	1.88	0.56
1:C:58:THR:HG21	1:C:77:PHE:HE1	1.71	0.55
3:T:713:DC:H2'	3:T:714:DG:C8	2.41	0.55
1:C:58:THR:HG21	1:C:77:PHE:CE1	2.42	0.55
2:B:422:LEU:H	2:B:422:LEU:HD22	1.71	0.55
4:F:816:DG:H2'	4:F:817:MRG:C8	2.36	0.55
1:A:114:ALA:HB1	1:A:160:PHE:CZ	2.42	0.55
1:A:19:PRO:HD3	1:A:80:LEU:HD13	1.88	0.55
1:C:395:LYS:HD2	1:C:414:TRP:CH2	2.41	0.54
2:D:309:ILE:O	2:D:312:GLU:HG2	2.08	0.54
1:C:296:THR:HG23	1:C:299:ALA:H	1.72	0.54
1:A:111:VAL:HB	1:A:185:ASP:HB2	1.89	0.54
1:C:279:LEU:N	1:C:302:GLU:OE1	2.40	0.54
1:A:172:LYS:CE	1:A:180:ILE:HB	2.30	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:P:807:DC:H2''	4:P:808:DC:H5''	1.90	0.54
1:C:12:LEU:HD11	1:C:127:TYR:CZ	2.43	0.54
2:D:85:GLN:HA	2:D:88:TRP:CE2	2.44	0.53
3:E:713:DC:H2''	3:E:714:DG:H5'	1.89	0.53
3:T:713:DC:H2''	3:T:714:DG:H5'	1.91	0.53
1:C:164:MET:HE2	1:C:168:LEU:HG	1.92	0.52
2:B:197:GLN:O	2:B:201:LYS:HG2	2.10	0.52
1:C:125:ARG:HE	1:C:147:ASN:HA	1.74	0.52
4:F:817:MRG:H2'	4:F:818:DC:C6	2.44	0.52
1:A:389:PHE:HB3	1:A:391:LEU:HD13	1.92	0.52
2:D:86:ASP:OD2	2:D:87:PHE:N	2.42	0.52
2:D:114:ALA:HB2	2:D:214:LEU:HD13	1.91	0.52
2:B:85:GLN:HA	2:B:88:TRP:NE1	2.25	0.52
1:A:333:GLY:HA2	8:A:604:EDO:H21	1.92	0.52
3:E:713:DC:H2'	3:E:714:DG:C8	2.45	0.52
1:C:107:THR:OG1	1:C:198:HIS:NE2	2.38	0.52
2:D:203:GLU:O	2:D:207:GLN:HG2	2.10	0.52
1:C:254:VAL:HG12	1:C:289:LEU:HD12	1.91	0.52
1:C:288:ALA:HB3	1:C:291:GLU:HB2	1.91	0.52
2:D:191:SER:OG	2:D:198:HIS:ND1	2.30	0.52
1:C:254:VAL:HG13	1:C:283:LEU:HD22	1.91	0.51
2:D:244:ILE:HG12	2:D:425:LEU:CD1	2.39	0.51
2:D:266:TRP:HZ3	2:D:425:LEU:HG	1.72	0.51
1:C:60:VAL:HG21	1:C:130:PHE:CD1	2.46	0.51
1:A:167:ILE:O	1:A:170:PRO:HD2	2.09	0.51
1:C:36:GLU:O	1:C:40:GLU:HG2	2.10	0.51
1:C:308:GLU:HA	1:C:311:LYS:HE2	1.91	0.51
1:C:181:TYR:HB2	1:C:188:TYR:HB3	1.92	0.51
2:B:170:PRO:HB2	2:B:208:HIS:CE1	2.46	0.51
1:C:122:GLU:HA	1:C:125:ARG:HH11	1.76	0.51
2:B:257:ILE:HB	2:B:283:LEU:HD21	1.93	0.51
1:C:12:LEU:H	1:C:12:LEU:HD12	1.75	0.51
1:A:109:LEU:HD13	1:A:216:THR:HG21	1.93	0.50
1:C:128:THR:HG21	1:C:146:TYR:HB2	1.93	0.50
1:C:206:ARG:NE	1:C:216:THR:OG1	2.38	0.50
4:F:807:DC:H4'	4:F:808:DC:OP1	2.11	0.50
1:C:125:ARG:O	1:C:128:THR:OG1	2.27	0.50
1:C:362:THR:HG21	1:C:367:GLN:HE21	1.77	0.50
2:D:114:ALA:HB2	2:D:214:LEU:HD22	1.93	0.50
3:E:721:DG:H5'	3:E:721:DG:H8	1.75	0.50
1:A:23:GLN:HG3	1:A:131:THR:HG23	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:47:ILE:HD12	1:A:146:TYR:HA	1.94	0.50
4:F:818:DC:H2'	4:F:819:DG:C8	2.47	0.50
1:C:115:TYR:CD1	7:C:601:DDS:H3'A	2.48	0.49
2:D:113:ASP:OD2	2:D:215:THR:OG1	2.30	0.49
2:D:332:GLN:CD	2:D:427:TYR:HB2	2.36	0.49
2:D:13:LYS:HD2	2:D:85:GLN:HB3	1.94	0.49
4:F:807:DC:H2''	4:F:808:DC:H5''	1.93	0.49
2:D:17:ASP:O	2:D:83:ARG:HD3	2.12	0.49
4:F:804:DA:H2''	4:F:805:DG:C5'	2.43	0.49
1:C:435:VAL:HG22	2:D:290:THR:HG21	1.95	0.49
2:D:331:LYS:HB2	2:D:337:TRP:CZ3	2.48	0.49
4:F:803:DC:H2''	4:F:804:DA:N7	2.28	0.49
2:B:244:ILE:HD13	2:B:425:LEU:HD11	1.95	0.49
1:C:195:ILE:HD11	1:C:199:ARG:NH2	2.28	0.49
1:C:271:TYR:HE2	1:C:310:LEU:HG	1.78	0.49
1:C:281:LYS:O	1:C:284:ARG:HG2	2.13	0.48
1:A:458:VAL:HG12	2:B:286:THR:HG21	1.95	0.48
2:B:85:GLN:HA	2:B:88:TRP:CE2	2.48	0.48
1:C:39:THR:O	1:C:42:GLU:HG2	2.13	0.48
2:D:365:VAL:HG11	2:D:401:TRP:HB2	1.95	0.48
1:C:331:LYS:HE3	1:C:364:ASP:OD2	2.14	0.48
1:A:5:ILE:HG22	1:A:212:TRP:HD1	1.77	0.48
2:D:354:TYR:HD1	2:D:374:LYS:HE2	1.79	0.48
2:B:198:HIS:CD2	2:B:202:ILE:HD11	2.49	0.47
2:D:244:ILE:HB	2:D:310:LEU:HD22	1.96	0.47
2:B:13:LYS:HD2	2:B:85:GLN:HB3	1.96	0.47
2:B:266:TRP:NE1	2:B:425:LEU:HD22	2.29	0.47
1:A:203:GLU:OE2	1:A:206:ARG:NH1	2.48	0.47
1:C:331:LYS:HB2	1:C:337:TRP:CZ3	2.48	0.47
1:A:475:GLN:HB3	1:A:501:TYR:CE2	2.49	0.47
1:A:442:VAL:HB	1:A:481:ALA:HB1	1.97	0.47
4:F:818:DC:H2'	4:F:819:DG:H8	1.80	0.47
4:F:804:DA:H2''	4:F:805:DG:O5'	2.14	0.46
1:A:58:THR:HG22	1:A:129:ALA:O	2.16	0.46
1:A:114:ALA:HB1	1:A:160:PHE:CE2	2.50	0.46
3:E:719:DG:H2''	3:E:720:DG:H8	1.72	0.46
2:B:317:VAL:HG12	2:B:347:LYS:HB3	1.98	0.46
1:C:41:MET:HB3	1:C:46:LYS:HB2	1.97	0.46
2:D:360:ALA:HA	2:D:367:GLN:NE2	2.31	0.46
4:F:816:DG:H2'	4:F:817:MRG:H8	1.98	0.46
2:D:244:ILE:CG1	2:D:425:LEU:HD11	2.42	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:F:803:DC:H2''	4:F:804:DA:C8	2.51	0.46
1:A:72:ARG:HH22	7:A:603:DDS:PA	2.39	0.46
1:C:114:ALA:HB1	1:C:160:PHE:CZ	2.50	0.46
3:E:702:DT:H3'	3:E:703:DG:C8	2.48	0.46
3:T:721:DG:H5'	3:T:721:DG:C8	2.51	0.46
2:D:78:ARG:O	2:D:82:LYS:HG3	2.16	0.46
1:C:111:VAL:HB	1:C:185:ASP:HB2	1.96	0.45
1:A:24:TRP:HE3	1:A:25:PRO:HD2	1.81	0.45
1:A:181:TYR:HB2	1:A:188:TYR:HB3	1.98	0.45
1:C:353:LYS:O	1:C:374:LYS:NZ	2.49	0.45
2:B:111:VAL:HG11	2:B:187:LEU:HD22	1.98	0.45
1:A:391:LEU:HB3	1:A:393:ILE:HG22	1.99	0.45
4:P:807:DC:H4'	4:P:808:DC:OP1	2.17	0.45
2:D:38:CYS:SG	2:D:132:ILE:HD11	2.57	0.45
1:A:263:LYS:HD2	4:P:819:DG:H4'	1.99	0.45
1:A:281:LYS:O	1:A:284:ARG:HG2	2.17	0.45
2:B:98:ALA:O	2:B:101:LYS:NZ	2.47	0.45
1:C:28:GLU:H	1:C:28:GLU:HG2	1.55	0.45
2:B:198:HIS:NE2	2:B:202:ILE:HD11	2.31	0.45
3:T:706:DC:H2'	3:T:707:DG:C8	2.51	0.45
1:C:19:PRO:HG3	1:C:80:LEU:HB2	1.98	0.45
1:A:195:ILE:HG12	1:A:199:ARG:HE	1.81	0.45
1:A:275:LYS:O	1:A:306:ASN:ND2	2.48	0.45
2:B:358:ARG:NH2	2:B:405:TYR:O	2.50	0.45
2:D:422:LEU:C	2:D:424:LYS:H	2.24	0.45
1:A:395:LYS:HD2	1:A:414:TRP:CH2	2.51	0.45
1:C:255:ASN:HB2	1:C:289:LEU:HG	1.98	0.45
1:C:448:ARG:NH1	4:F:806:DT:O2	2.50	0.45
1:A:56:TYR:HB2	1:A:129:ALA:HB3	1.99	0.45
2:B:184:MET:HE3	2:B:184:MET:HB3	1.64	0.44
1:A:20:LYS:O	1:A:21:VAL:HG23	2.17	0.44
1:A:112:GLY:HA2	7:A:603:DDS:O2G	2.17	0.44
2:B:236:PRO:HA	2:B:239:TRP:CD2	2.52	0.44
2:D:356:ARG:HD2	2:D:360:ALA:HB1	1.99	0.44
4:P:806:DT:H2'	4:P:807:DC:C6	2.52	0.44
2:D:85:GLN:HA	2:D:88:TRP:NE1	2.33	0.44
2:D:395:LYS:HG3	2:D:416:PHE:CE2	2.53	0.44
1:A:376:THR:HG21	2:B:401:TRP:CH2	2.53	0.43
2:B:17:ASP:O	2:B:83:ARG:HD3	2.18	0.43
1:C:252:TRP:NE1	1:C:295:LEU:HD11	2.33	0.43
2:D:395:LYS:O	2:D:399:GLU:HG3	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:540:LYS:HE3	2:D:265:ASN:OD1	2.18	0.43
2:B:101:LYS:O	2:B:236:PRO:HB2	2.18	0.43
1:A:137:ASN:OD1	1:A:137:ASN:N	2.43	0.43
3:T:721:DG:H5'	3:T:721:DG:H8	1.84	0.43
2:D:84:THR:HG22	2:D:84:THR:O	2.19	0.43
1:A:266:TRP:O	1:A:269:GLN:HG2	2.19	0.43
2:B:88:TRP:CE3	2:B:89:GLU:HG3	2.53	0.43
2:B:395:LYS:HG3	2:B:416:PHE:CE2	2.54	0.43
1:C:308:GLU:HA	1:C:311:LYS:HG3	2.00	0.43
1:C:330:GLN:NE2	1:C:340:GLN:OE1	2.38	0.43
2:D:281:LYS:HG2	2:D:284:ARG:HE	1.84	0.43
4:F:804:DA:H2''	4:F:805:DG:H5'	1.99	0.43
2:D:303:LEU:HG	2:D:307:ARG:NH1	2.34	0.42
1:C:458:VAL:HG12	2:D:286:THR:HG21	2.00	0.42
2:D:323:LYS:O	2:D:385:LYS:NZ	2.51	0.42
2:B:250:ASP:OD1	2:B:250:ASP:N	2.51	0.42
1:C:540:LYS:HD3	1:C:540:LYS:HA	1.74	0.42
1:A:80:LEU:O	1:A:84:THR:OG1	2.30	0.42
1:A:58:THR:HG21	1:A:77:PHE:CE1	2.55	0.42
1:C:406:TRP:CD2	2:D:420:PRO:HB3	2.55	0.42
2:D:25:PRO:HG3	8:D:503:EDO:H11	2.00	0.42
1:A:149:LEU:HD11	1:A:159:ILE:HG22	2.02	0.42
1:C:172:LYS:HG2	1:C:180:ILE:HD12	2.02	0.42
1:C:24:TRP:CD1	1:C:59:PRO:HB3	2.55	0.41
1:A:27:THR:O	1:A:31:ILE:HG13	2.19	0.41
1:C:47:ILE:HD12	1:C:146:TYR:HA	2.03	0.41
2:D:331:LYS:NZ	2:D:364:ASP:OD2	2.52	0.41
1:A:8:VAL:O	1:A:10:VAL:HG23	2.20	0.41
1:A:363:ASN:HA	1:A:511:ASP:OD1	2.20	0.41
1:A:406:TRP:CZ2	2:B:420:PRO:HG3	2.55	0.41
1:C:224:GLU:HA	1:C:225:PRO:HD3	1.84	0.41
2:D:115:TYR:CD2	2:D:156:SER:HB3	2.55	0.41
2:D:169:GLU:O	2:D:173:LYS:HG2	2.21	0.41
2:B:210:LEU:HD12	2:B:228:LEU:HD23	2.03	0.41
1:C:252:TRP:CD1	1:C:295:LEU:HD21	2.55	0.41
2:D:115:TYR:O	2:D:149:LEU:HB2	2.21	0.41
1:A:103:LYS:HD3	1:A:103:LYS:HA	1.84	0.41
1:A:418:ASN:O	1:A:420:PRO:HD3	2.20	0.41
1:A:427:TYR:OH	1:A:510:PRO:HD2	2.21	0.41
2:B:87:PHE:CD2	2:B:87:PHE:C	2.99	0.41
2:B:263:LYS:HE3	2:B:425:LEU:HA	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:281:LYS:HG2	2:B:284:ARG:HH21	1.85	0.41
4:P:813:DT:H2''	4:P:814:DC:C6	2.56	0.41
2:D:423:VAL:HG22	2:D:423:VAL:O	2.21	0.41
1:A:6:GLU:OE1	1:A:166:LYS:NZ	2.54	0.41
1:A:220:LYS:HD3	1:A:220:LYS:HA	1.89	0.41
2:B:230:MET:HE3	2:B:230:MET:HB3	1.89	0.41
1:C:171:PHE:CD2	1:C:205:LEU:HD13	2.56	0.41
3:E:702:DT:H2'	3:E:703:DG:O4'	2.21	0.41
1:A:53:GLU:O	1:A:55:PRO:HD3	2.21	0.40
2:B:360:ALA:HA	2:B:367:GLN:CD	2.46	0.40
2:D:423:VAL:C	2:D:425:LEU:H	2.29	0.40
1:A:156:SER:HB2	1:A:157:PRO:HD3	2.03	0.40
1:A:206:ARG:NH2	1:A:216:THR:O	2.54	0.40
2:B:65:LYS:NZ	2:B:230:MET:HE1	2.36	0.40
4:P:816:DG:H2'	4:P:817:MRG:H8	2.04	0.40
2:D:64:LYS:HE3	2:D:71:TRP:CE2	2.56	0.40
2:D:317:VAL:HG12	2:D:347:LYS:HB3	2.03	0.40
4:F:806:DT:H2'	4:F:807:DC:C6	2.57	0.40
2:B:88:TRP:CE3	2:B:154:LYS:HD2	2.55	0.40
2:D:425:LEU:HD13	2:D:425:LEU:C	2.47	0.40
1:A:58:THR:HG21	1:A:77:PHE:HE1	1.86	0.40
1:A:86:ASP:O	2:B:55:PRO:HB3	2.21	0.40
1:C:161:GLN:HG3	1:C:184:MET:HE3	2.03	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	116/557 (21%)	114 (98%)	2 (2%)	0	100	100
1	C	255/557 (46%)	241 (94%)	11 (4%)	3 (1%)	10	27

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	D	284/428 (66%)	276 (97%)	7 (2%)	1 (0%)	30	54
All	All	655/1542 (42%)	631 (96%)	20 (3%)	4 (1%)	21	44

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	4	PRO
1	C	1	PRO
2	D	423	VAL
1	C	133	PRO

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	495/497 (100%)	491 (99%)	4 (1%)	73	88
1	C	495/497 (100%)	492 (99%)	3 (1%)	78	91
2	B	372/390 (95%)	369 (99%)	3 (1%)	73	88
2	D	374/390 (96%)	369 (99%)	5 (1%)	61	83
All	All	1736/1774 (98%)	1721 (99%)	15 (1%)	70	87

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	ILE
1	A	21	VAL
1	A	276	VAL
1	A	391	LEU
2	B	87	PHE
2	B	208	HIS
2	B	422	LEU
1	C	2	ILE
1	C	28	GLU
1	C	276	VAL

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Mol	Chain	Res	Type
2	D	11	LYS
2	D	87	PHE
2	D	208	HIS
2	D	422	LEU
2	D	425	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	145	GLN
1	A	547	GLN
2	B	255	ASN
2	B	269	GLN
2	B	428	GLN
1	C	255	ASN
1	C	367	GLN
1	C	407	GLN
1	C	547	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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	25,28,29	2.72	7 (28%)	32,39,42	2.39	13 (40%)
4	MRG	F	817	1,3,4	25,28,29	2.68	7 (28%)	32,39,42	2.38	12 (37%)

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	-	2/12/26/27	0/3/3/3
4	MRG	F	817	1,3,4	-	2/12/26/27	0/3/3/3

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	P	817	MRG	C2-N2	7.18	1.49	1.34
4	F	817	MRG	C2-N2	7.11	1.49	1.34
4	P	817	MRG	C2-N3	6.74	1.44	1.32
4	F	817	MRG	C2-N3	6.55	1.44	1.32
4	P	817	MRG	C4-N3	4.96	1.45	1.34
4	F	817	MRG	C4-N3	4.76	1.45	1.34
4	F	817	MRG	C2-N1	4.51	1.43	1.36
4	P	817	MRG	C2-N1	4.30	1.43	1.36
4	F	817	MRG	C2'-C3'	-3.37	1.44	1.52
4	P	817	MRG	C2'-C3'	-3.33	1.44	1.52
4	P	817	MRG	O3'-C3'	-2.64	1.37	1.43
4	F	817	MRG	O4'-C4'	-2.48	1.39	1.45
4	P	817	MRG	O4'-C4'	-2.46	1.39	1.45
4	F	817	MRG	O3'-C3'	-2.31	1.38	1.43

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	P	817	MRG	C2-N3-C4	7.27	121.10	112.00
4	F	817	MRG	C2-N3-C4	7.26	121.08	112.00
4	P	817	MRG	C5-C4-N3	-5.53	119.59	128.39
4	F	817	MRG	C5-C4-N3	-5.20	120.11	128.39
4	F	817	MRG	C22-C21-N2	4.04	123.52	112.20
4	F	817	MRG	O4'-C1'-N9	3.91	114.80	107.86
4	P	817	MRG	O4'-C1'-N9	3.79	114.59	107.86
4	P	817	MRG	C22-C21-N2	3.46	121.91	112.20
4	F	817	MRG	C2-N1-C6	-3.43	120.40	124.55
4	P	817	MRG	N9-C4-N3	3.42	132.80	125.95
4	P	817	MRG	C2-N1-C6	-3.26	120.61	124.55
4	P	817	MRG	N1-C2-N3	-2.83	118.91	123.68
4	F	817	MRG	N1-C2-N3	-2.80	118.96	123.68
4	F	817	MRG	N9-C4-N3	2.72	131.40	125.95
4	F	817	MRG	C5-C6-N1	2.63	119.94	113.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	P	817	MRG	C5-C6-N1	2.55	119.73	113.25
4	P	817	MRG	O6-C6-C5	-2.54	119.83	126.53
4	F	817	MRG	C4-C5-N7	-2.47	106.75	110.67
4	F	817	MRG	O6-C6-C5	-2.40	120.19	126.53
4	F	817	MRG	C8-N7-C5	2.09	107.98	104.26
4	P	817	MRG	C4-C5-N7	-2.08	107.37	110.67
4	P	817	MRG	N9-C8-N7	-2.07	109.56	113.40
4	F	817	MRG	N9-C8-N7	-2.07	109.57	113.40
4	P	817	MRG	C8-N7-C5	2.06	107.94	104.26
4	P	817	MRG	N2-C2-N3	2.06	123.03	120.19

There are no chirality outliers.

All (4) torsion outliers are listed below:

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

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	P	817	MRG	1	0
4	F	817	MRG	4	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	0.85	0	15,15,17	0.84	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	FRU	G	2	5	11,12,12	2.40	5 (45%)	10,18,18	5.25	3 (30%)
5	GLC	H	1	5	11,11,12	0.90	1 (9%)	15,15,17	1.18	1 (6%)
5	FRU	H	2	5	11,12,12	3.06	4 (36%)	10,18,18	5.37	5 (50%)

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
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	-	0/2/19/22	0/1/1/1
5	FRU	H	2	5	-	2/5/24/24	0/1/1/1

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	H	2	FRU	C4-C5	-7.68	1.33	1.53
5	G	2	FRU	O5-C2	5.23	1.51	1.43
5	H	2	FRU	C4-C3	4.18	1.69	1.53
5	H	2	FRU	O2-C2	3.41	1.46	1.40
5	G	2	FRU	O5-C5	-3.27	1.36	1.43
5	G	2	FRU	C4-C5	2.88	1.60	1.53
5	H	2	FRU	O5-C2	-2.85	1.39	1.43
5	G	2	FRU	O2-C2	2.45	1.45	1.40
5	G	2	FRU	O3-C3	2.10	1.46	1.42
5	H	1	GLC	O5-C1	2.08	1.47	1.43

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	G	2	FRU	O2-C2-O5	-13.75	82.93	109.33
5	H	2	FRU	C5-C4-C3	-9.52	72.54	102.07
5	H	2	FRU	O5-C5-C6	-8.79	84.91	108.89
5	H	2	FRU	C6-C5-C4	-8.27	95.57	115.10
5	G	2	FRU	O5-C5-C4	-7.92	85.43	105.42
5	H	2	FRU	O5-C5-C4	6.24	121.18	105.42
5	G	2	FRU	O3-C3-C4	3.76	126.51	113.25
5	H	2	FRU	O4-C4-C5	-2.45	104.04	111.08
5	H	1	GLC	C6-C5-C4	-2.20	107.62	113.02

There are no chirality outliers.

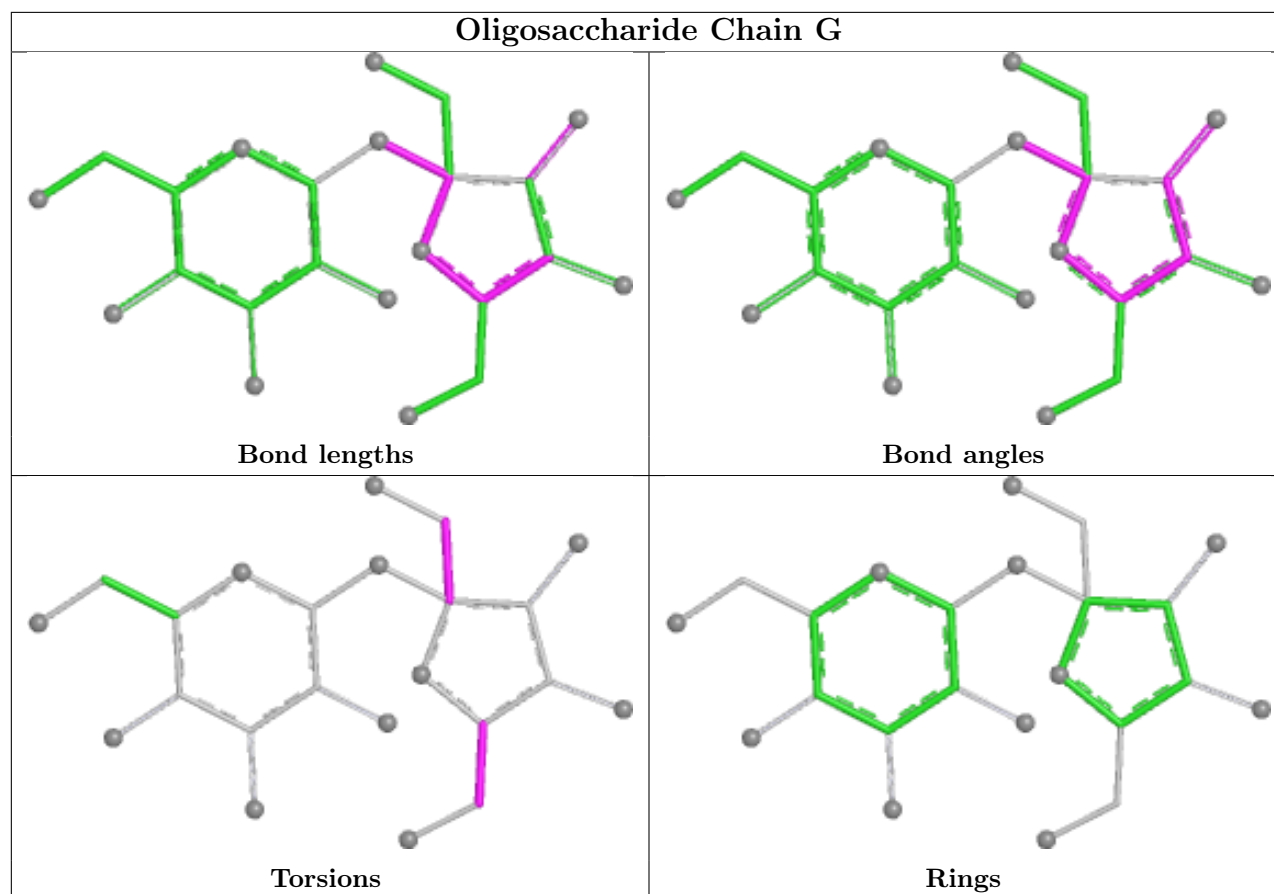
All (5) torsion outliers are listed below:

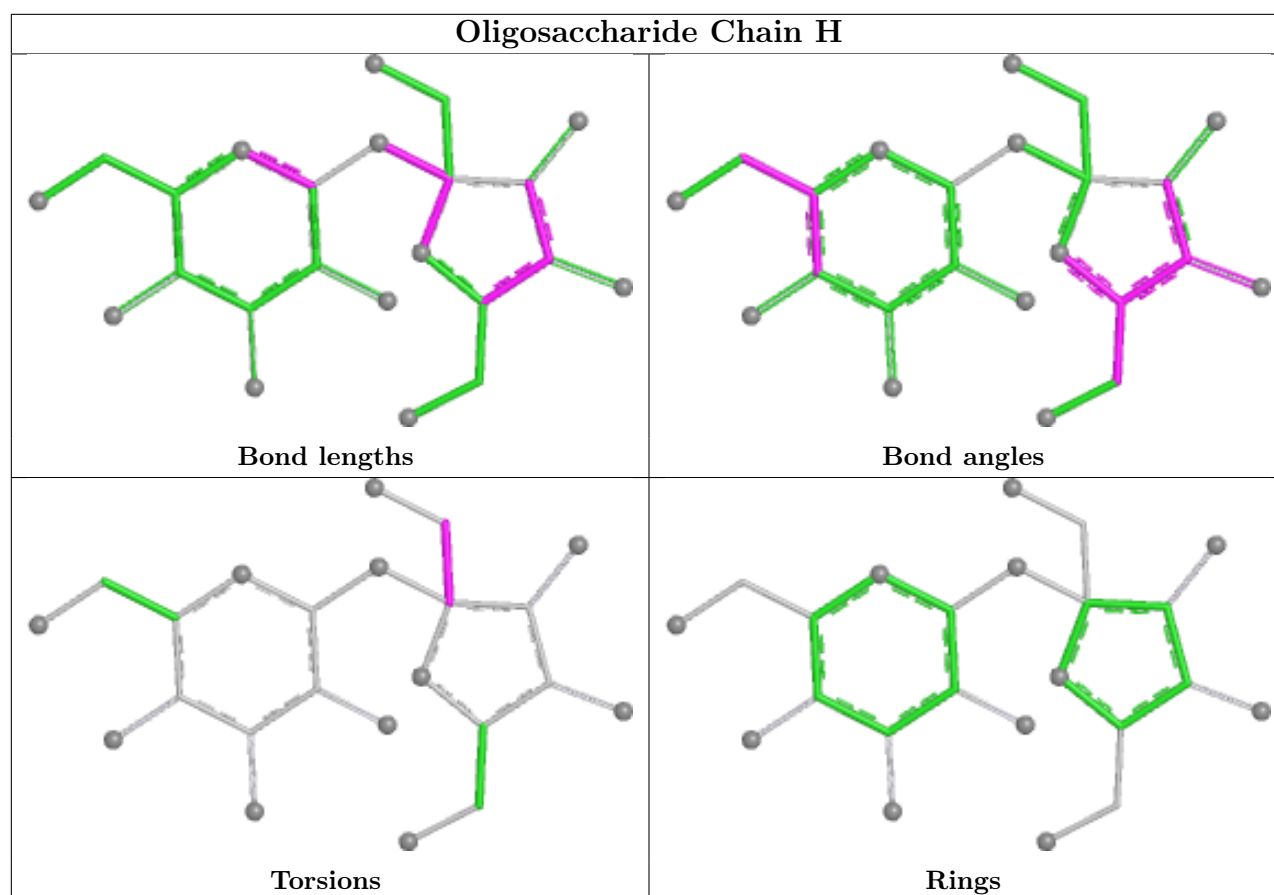
Mol	Chain	Res	Type	Atoms
5	H	2	FRU	O1-C1-C2-O2
5	G	2	FRU	O5-C5-C6-O6
5	G	2	FRU	O1-C1-C2-C3
5	H	2	FRU	O1-C1-C2-C3
5	G	2	FRU	O1-C1-C2-O2

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 13 ligands modelled in this entry, 4 are monoatomic - leaving 9 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	$\# Z > 2$
8	EDO	A	604	-	3,3,3	0.43	0	2,2,2	0.32	0
8	EDO	B	503	-	3,3,3	0.44	0	2,2,2	0.42	0
8	EDO	B	504	-	3,3,3	0.51	0	2,2,2	0.32	0
8	EDO	D	502	-	3,3,3	0.44	0	2,2,2	0.41	0
7	DDS	C	601	6	30,31,31	1.72	5 (16%)	42,48,48	1.97	10 (23%)
8	EDO	D	503	-	3,3,3	0.49	0	2,2,2	0.40	0
7	DDS	A	603	6	30,31,31	1.71	5 (16%)	42,48,48	1.95	11 (26%)
8	EDO	T	801	-	3,3,3	0.45	0	2,2,2	0.30	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
9	SO4	B	502	-	4,4,4	0.20	0	6,6,6	0.16	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
8	EDO	A	604	-	-	0/1/1/1	-
8	EDO	B	503	-	-	1/1/1/1	-
8	EDO	B	504	-	-	0/1/1/1	-
8	EDO	D	502	-	-	1/1/1/1	-
7	DDS	C	601	6	-	5/22/31/31	0/3/3/3
8	EDO	D	503	-	-	1/1/1/1	-
7	DDS	A	603	6	-	5/22/31/31	0/3/3/3
8	EDO	T	801	-	-	0/1/1/1	-

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	A	603	DDS	C6-N6	5.07	1.47	1.34
7	C	601	DDS	C6-N6	5.06	1.47	1.34
7	C	601	DDS	PB-O3A	-4.48	1.54	1.59
7	A	603	DDS	PB-O3A	-4.05	1.55	1.59
7	C	601	DDS	C5'-C4'	-2.98	1.41	1.50
7	A	603	DDS	C5'-C4'	-2.84	1.41	1.50
7	A	603	DDS	C3'-C2'	-2.68	1.46	1.54
7	C	601	DDS	C3'-C2'	-2.67	1.46	1.54
7	C	601	DDS	O5'-C5'	-2.66	1.34	1.44
7	A	603	DDS	O5'-C5'	-2.52	1.35	1.44

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	A	603	DDS	C5-C4-N3	-5.40	119.28	126.72
7	C	601	DDS	C5-C4-N3	-5.28	119.44	126.72
7	A	603	DDS	N3-C2-N1	-4.24	122.16	128.58
7	C	601	DDS	N3-C2-N1	-4.09	122.39	128.58
7	C	601	DDS	N3-C4-N9	4.03	134.02	127.17
7	A	603	DDS	N3-C4-N9	4.01	133.99	127.17
7	C	601	DDS	O4'-C1'-N9	3.77	114.73	107.96
7	A	603	DDS	N9-C8-N7	-3.66	108.75	113.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	C	601	DDS	N9-C8-N7	-3.54	108.92	113.94
7	A	603	DDS	O4'-C1'-N9	3.51	114.27	107.96
7	A	603	DDS	C2-N3-C4	3.47	120.31	111.83
7	C	601	DDS	C2-N3-C4	3.46	120.29	111.83
7	C	601	DDS	C3'-C2'-C1'	3.29	106.66	102.87
7	A	603	DDS	C5-N7-C8	3.21	108.50	103.45
7	C	601	DDS	C4-N9-C8	3.20	109.10	105.74
7	A	603	DDS	C4-C5-N7	-3.00	107.15	110.58
7	A	603	DDS	C4-N9-C8	2.98	108.87	105.74
7	C	601	DDS	C5-N7-C8	2.95	108.08	103.45
7	C	601	DDS	C4-C5-N7	-2.82	107.36	110.58
7	A	603	DDS	C3'-C2'-C1'	2.29	105.51	102.87
7	A	603	DDS	C4'-O4'-C1'	-2.10	107.82	109.81

There are no chirality outliers.

All (13) torsion outliers are listed below:

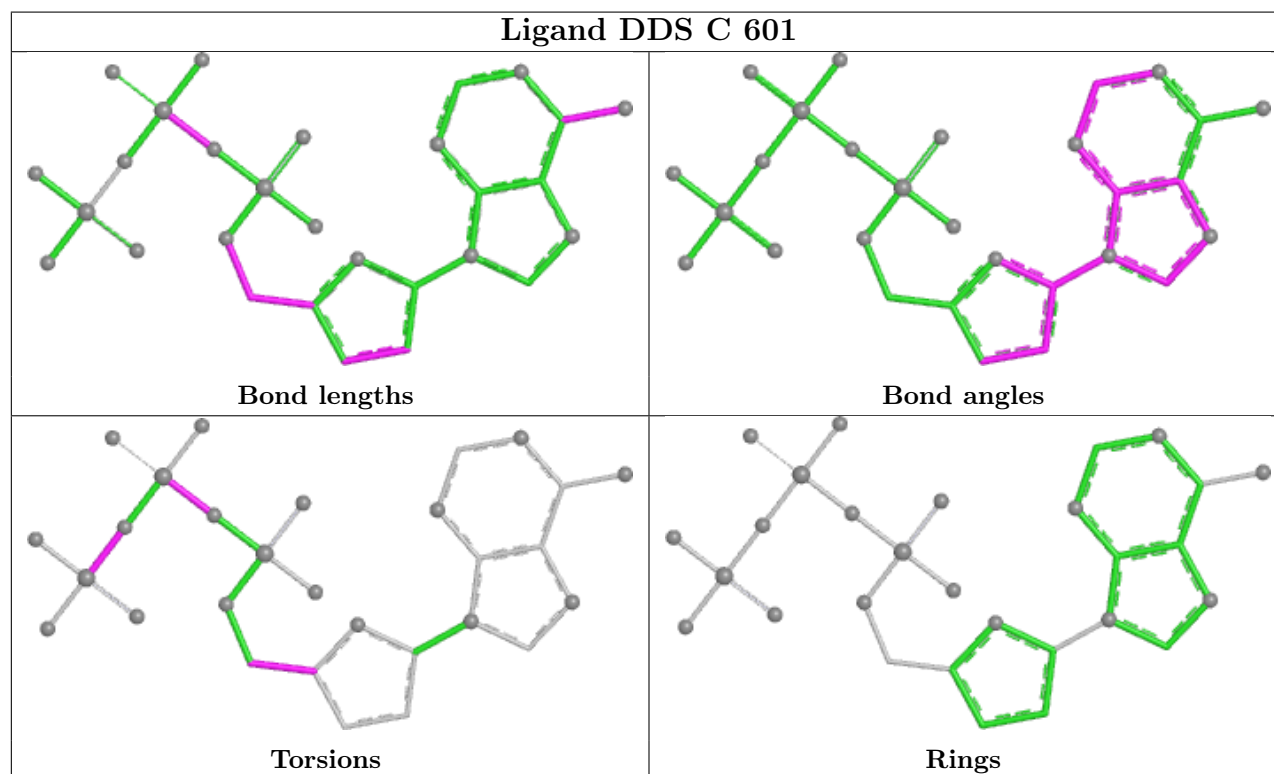
Mol	Chain	Res	Type	Atoms
7	A	603	DDS	C5'-O5'-PA-O1A
7	A	603	DDS	C3'-C4'-C5'-O5'
7	A	603	DDS	O4'-C4'-C5'-O5'
7	C	601	DDS	C3'-C4'-C5'-O5'
7	C	601	DDS	O4'-C4'-C5'-O5'
7	C	601	DDS	PB-O3B-PG-O1G
7	A	603	DDS	PA-O3A-PB-O2B
7	C	601	DDS	PA-O3A-PB-O2B
8	B	503	EDO	O1-C1-C2-O2
8	D	502	EDO	O1-C1-C2-O2
8	D	503	EDO	O1-C1-C2-O2
7	A	603	DDS	PA-O3A-PB-O1B
7	C	601	DDS	PA-O3A-PB-O1B

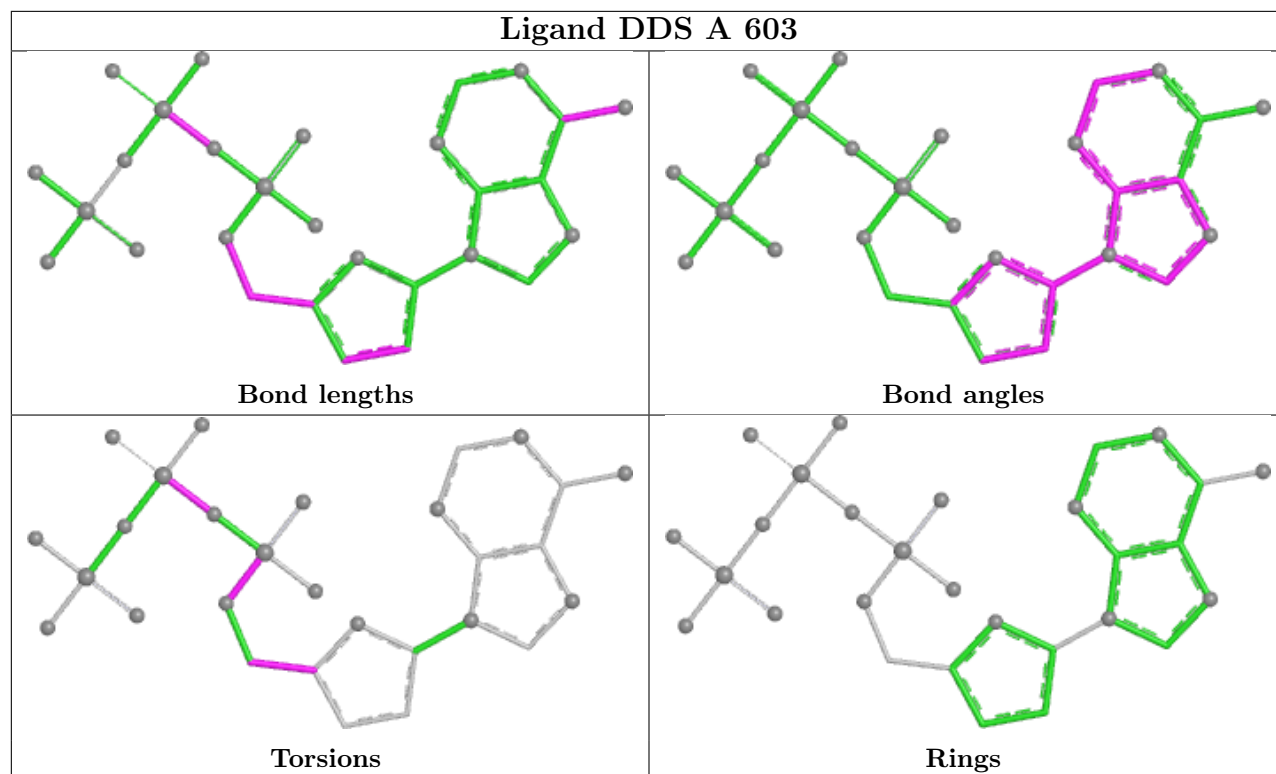
There are no ring outliers.

5 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	A	604	EDO	1	0
8	B	504	EDO	1	0
7	C	601	DDS	2	0
8	D	503	EDO	1	0
7	A	603	DDS	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.





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	556/557 (99%)	-0.07	14 (2%) 58 55	31, 85, 182, 297	0
1	C	556/557 (99%)	0.06	15 (2%) 56 53	35, 92, 196, 259	0
2	B	411/428 (96%)	-0.30	13 (3%) 50 46	34, 69, 132, 213	0
2	D	412/428 (96%)	0.02	18 (4%) 39 35	40, 83, 162, 226	0
3	E	24/27 (88%)	0.13	0 100 100	80, 121, 199, 256	0
3	T	24/27 (88%)	-0.08	0 100 100	75, 115, 214, 240	0
4	F	19/21 (90%)	-0.13	0 100 100	64, 103, 209, 212	0
4	P	19/21 (90%)	-0.24	1 (5%) 32 28	69, 105, 182, 220	0
All	All	2021/2066 (97%)	-0.06	61 (3%) 52 49	31, 83, 180, 297	0

All (61) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	74	LEU	4.7
2	B	231	GLY	3.8
1	C	61	PHE	3.6
1	A	73	LYS	3.6
2	D	217	PRO	3.4
2	D	420	PRO	3.4
2	B	90	VAL	3.3
1	C	280	SER	3.2
2	D	359	GLY	3.2
2	B	423	VAL	3.0
2	D	232	TYR	3.0
2	D	423	VAL	2.9
1	C	5	ILE	2.8
2	D	418	ASN	2.8
2	D	422	LEU	2.8
1	A	4	PRO	2.8

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Mol	Chain	Res	Type	RSRZ
2	B	422	LEU	2.6
1	C	277	ARG	2.6
2	D	214	LEU	2.6
1	A	133	PRO	2.5
2	D	421	PRO	2.5
2	D	360	ALA	2.5
1	A	131	THR	2.5
2	D	241	VAL	2.5
1	A	109	LEU	2.4
2	B	5	ILE	2.4
1	C	24	TRP	2.4
2	D	426	TRP	2.4
1	C	-1	MET	2.4
2	B	232	TYR	2.4
2	B	87	PHE	2.4
1	A	74	LEU	2.4
1	A	132	ILE	2.4
1	C	264	LEU	2.4
1	A	135	ILE	2.4
1	A	38	CYS	2.4
2	B	230	MET	2.3
1	C	142	ILE	2.3
1	A	247	PRO	2.3
2	D	88	TRP	2.3
2	D	95	PRO	2.3
2	D	87	PHE	2.3
1	C	274	ILE	2.2
1	C	299	ALA	2.2
4	P	803	DC	2.2
2	D	94	ILE	2.2
2	B	88	TRP	2.2
2	B	213	GLY	2.2
2	D	216	THR	2.2
2	B	86	ASP	2.2
1	A	358	ARG	2.1
1	A	37	ILE	2.1
1	C	34	LEU	2.1
2	B	228	LEU	2.1
1	C	293	ILE	2.1
1	A	60	VAL	2.1
1	C	73	LYS	2.0
2	B	266	TRP	2.0

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Mol	Chain	Res	Type	RSRZ
1	C	295	LEU	2.0
1	A	24	TRP	2.0
2	D	427	TYR	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	0.87	0.12	111,128,154,182	0
4	MRG	P	817	26/27	0.91	0.10	95,122,147,149	0

6.3 Carbohydrates [i](#)

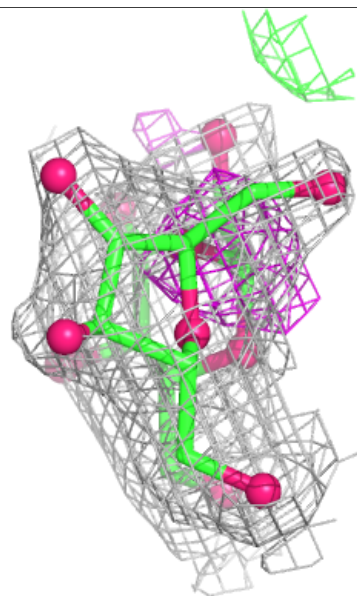
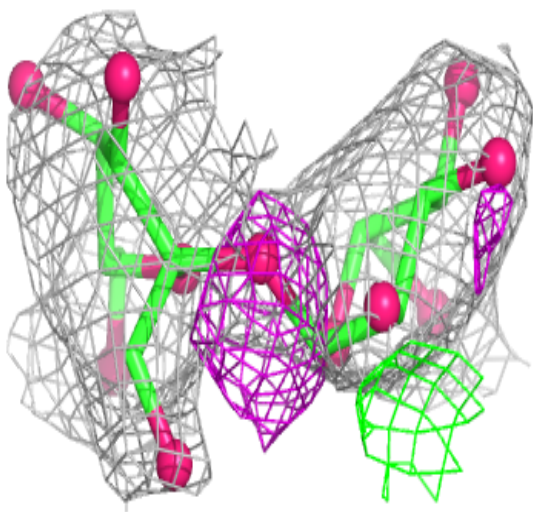
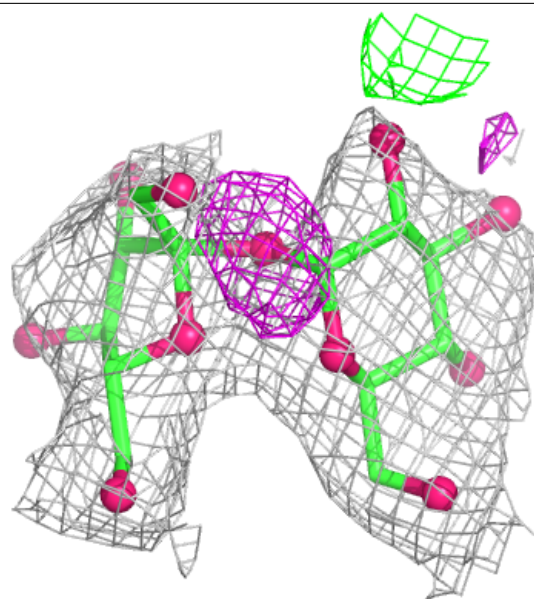
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

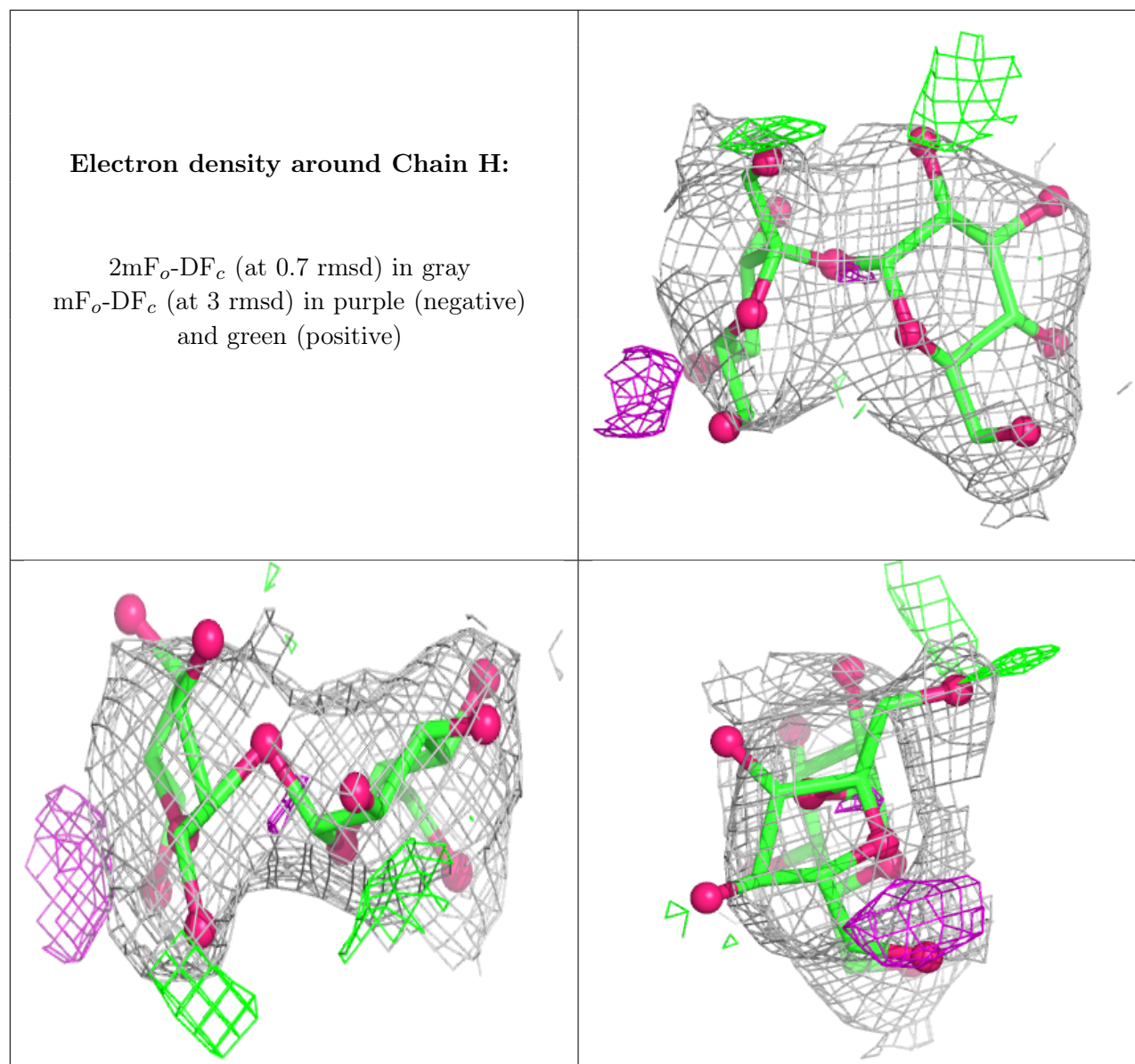
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	GLC	G	1	11/12	-	-	79,105,117,129	0
5	FRU	G	2	12/12	-	-	84,128,135,141	0
5	GLC	H	1	11/12	-	-	80,109,129,134	0
5	FRU	H	2	12/12	-	-	125,146,150,155	0

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 ⓘ

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.

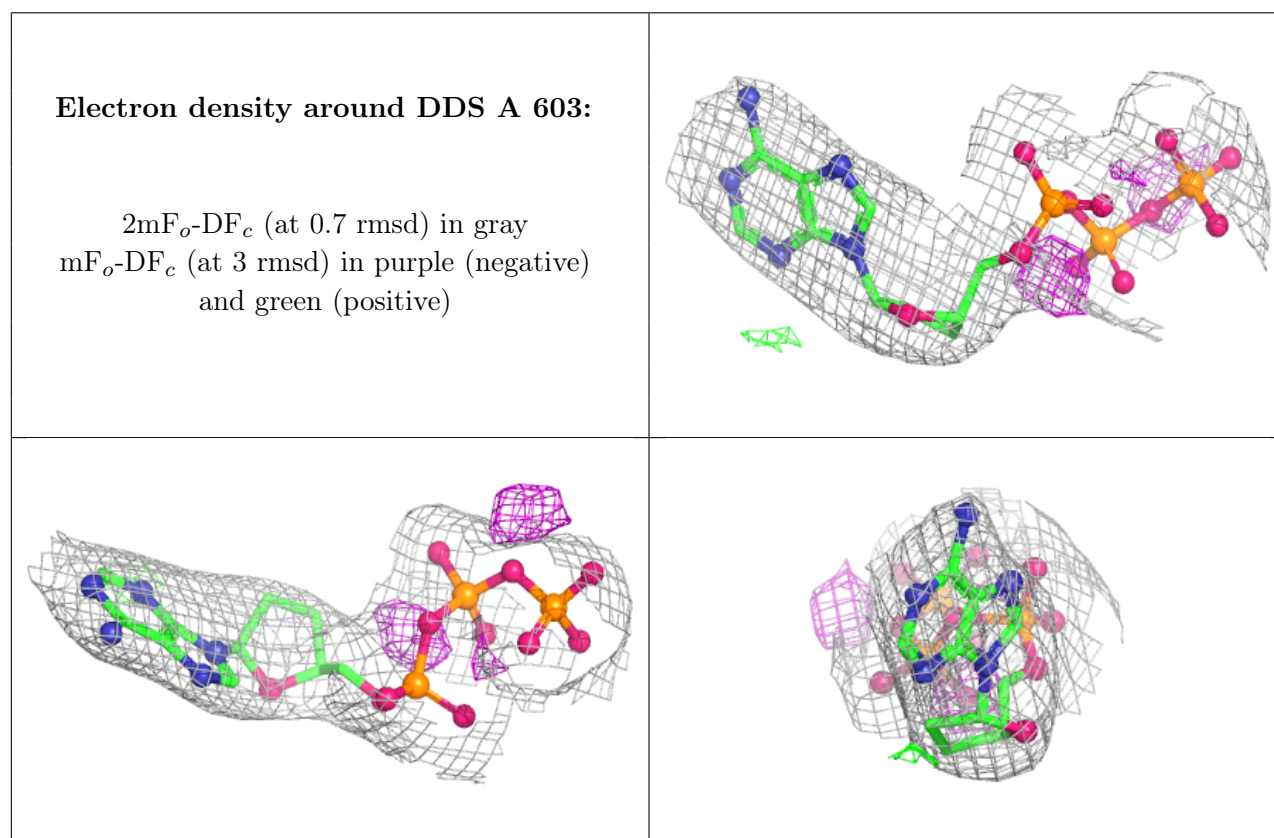
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
8	EDO	A	604	4/4	0.77	0.19	79,84,94,96	0
8	EDO	T	801	4/4	0.77	0.23	82,92,100,101	0
8	EDO	D	502	4/4	0.78	0.28	91,93,93,96	0
6	MG	C	603	1/1	0.80	0.22	107,107,107,107	0
8	EDO	B	504	4/4	0.82	0.22	73,80,89,93	0

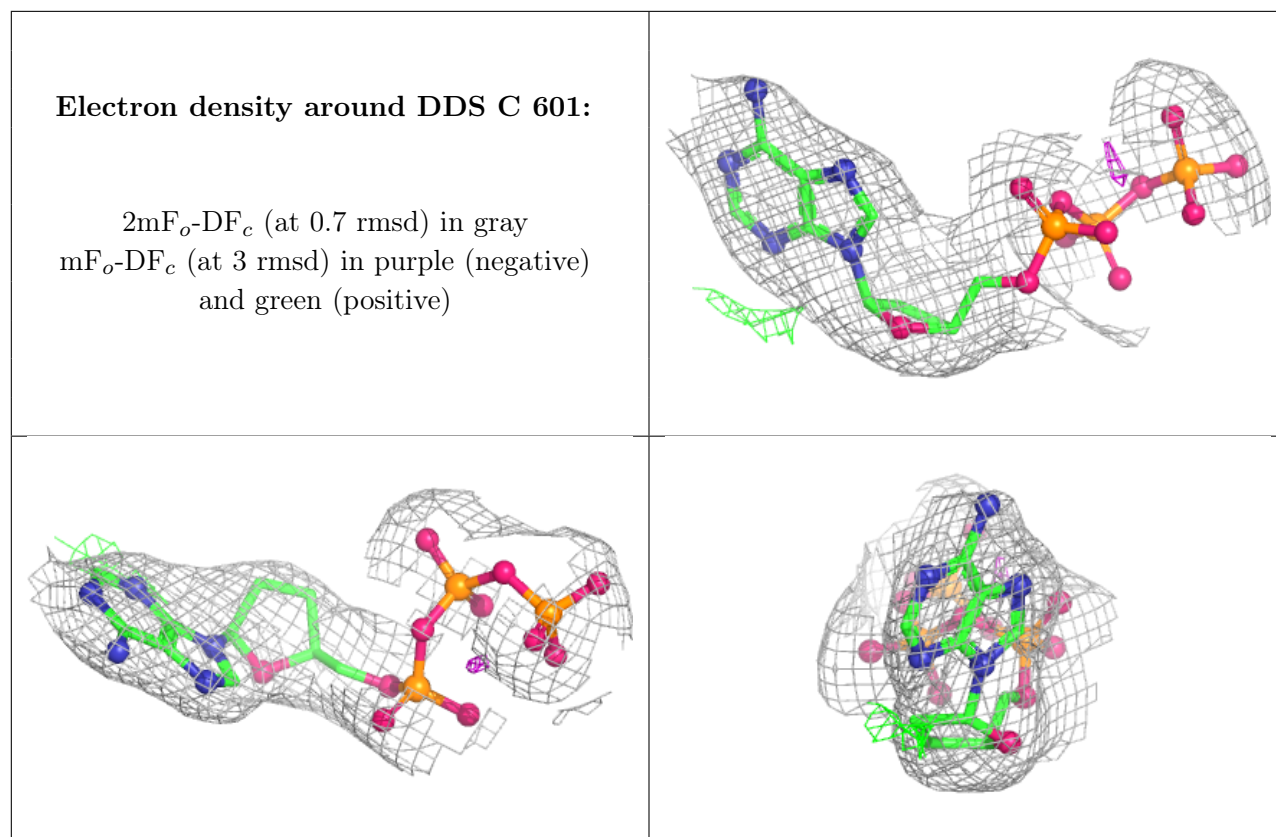
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
8	EDO	B	503	4/4	0.87	0.22	64,67,73,81	0
6	MG	A	602	1/1	0.87	0.14	117,117,117,117	0
9	SO4	B	502	5/5	0.89	0.20	150,152,154,161	0
8	EDO	D	503	4/4	0.90	0.19	66,70,77,82	0
7	DDS	A	603	29/29	0.90	0.09	66,92,155,158	0
7	DDS	C	601	29/29	0.93	0.07	66,81,131,138	0
6	MG	C	602	1/1	0.94	0.06	109,109,109,109	0
6	MG	A	601	1/1	0.97	0.05	159,159,159,159	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.