



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 11:00 PM UTC

PDB ID : 7QXD / pdb_00007qxd
Title : Recognition of Staphylococcus aureus wall teichoic acid analogue SA475 (compound 2) by Fab4497
Authors : Soriano-Maldonado, P.; van Raaij, M.J.
Deposited on : 2022-01-26
Resolution : 1.65 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

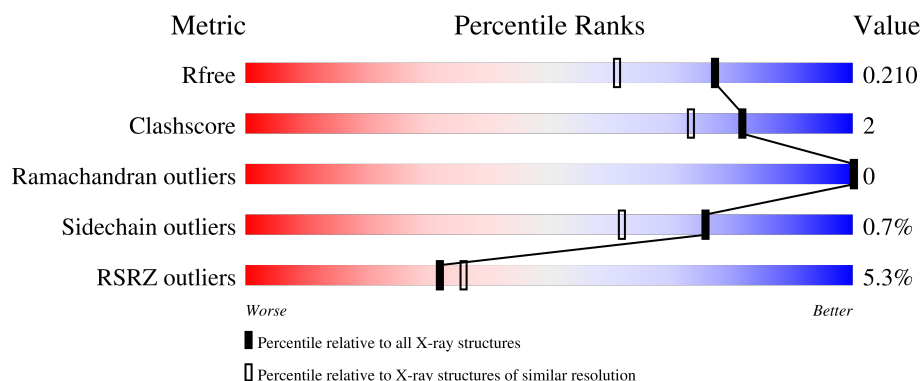
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.65 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	1141 (1.64-1.64)
Clashscore	190562	1171 (1.64-1.64)
Ramachandran outliers	187476	1151 (1.64-1.64)
Sidechain outliers	187428	1150 (1.64-1.64)
RSRZ outliers	180081	1141 (1.64-1.64)

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	HHH	236	4% 84% 7% 9%
1	KKK	236	7% 84% 7% 9%
2	LLL	220	5% 93% 6%
2	MMM	220	4% 94% 5%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 7623 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.

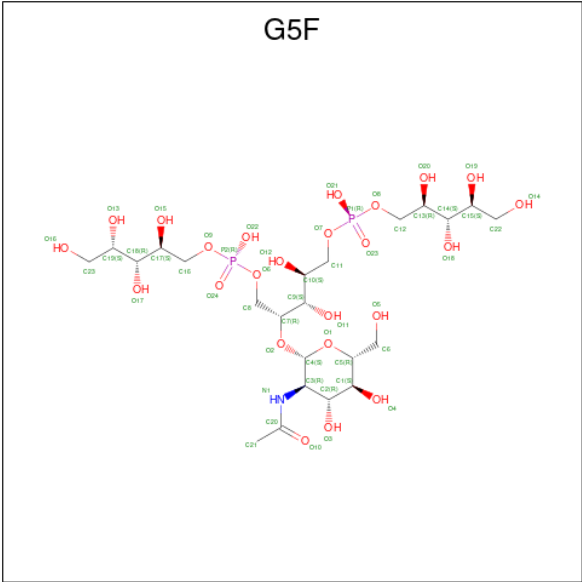
- Molecule 1 is a protein called Antibody Fab 4497 heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	HHH	214	Total	C	N	O	S	0	3	0
			1608	1014	272	316	6			
1	KKK	215	Total	C	N	O	S	0	2	0
			1614	1017	274	317	6			

- Molecule 2 is a protein called Antibody Fab 4497 light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	LLL	219	Total	C	N	O	S	0	1	0
			1721	1080	296	341	4			
2	MMM	219	Total	C	N	O	S	0	1	0
			1726	1083	299	340	4			

- Molecule 3 is [(2 {S},3 {S},4 {R})-4-[(2 {S},3 {R},4 {R},5 {S},6 {R})-3-acetamido-6-(hydroxymethyl)-4,5-bis(oxidanyl)oxan-2-yl]oxy-5-[[[(2 {S},3 {R})-2,3-bis(oxidanyl)butoxy]-oxidanyl-phosphoryl]oxy-2,3-bis(oxidanyl)pentyl] [(2 {R},3 {S},4 {S})-2,3,4,5-tetrakis(oxidanyl)pentyl] hydrogen phosphate (CCD ID: G5F) (formula: C₂₃H₄₇NO₂₄P₂) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	HHH	1	Total 47	C 22	N 1	O 22	P 2	0	0
3	MMM	1	Total 49	C 23	N 1	O 23	P 2	0	0

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



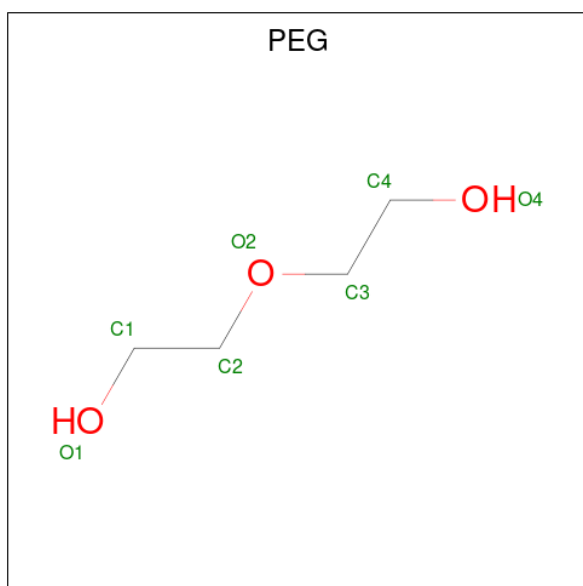
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	LLL	1	Total	C	O	0	0
			6	3	3		
4	KKK	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	LLL	1	Total	O	S	0	0
			5	4	1		
5	MMM	1	Total	O	S	0	0
			5	4	1		

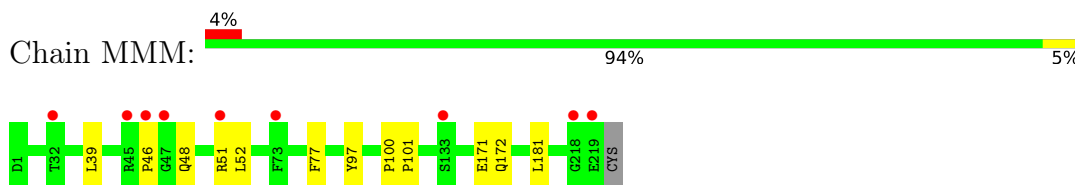
- Molecule 6 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: C₄H₁₀O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	KKK	1	Total	C	O	0	0
			7	4	3		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	HHH	195	Total 197	O 197	0	2
7	LLL	176	Total 178	O 178	0	2
7	KKK	188	Total 192	O 192	0	4
7	MMM	259	Total 262	O 262	0	3



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	64.39Å 113.37Å 154.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	90.02 – 1.65 90.02 – 1.65	Depositor EDS
% Data completeness (in resolution range)	93.0 (90.02-1.65) 93.0 (90.02-1.65)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.64 (at 1.65Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.175 , 0.202 0.186 , 0.210	Depositor DCC
R_{free} test set	6292 reflections (4.60%)	wwPDB-VP
Wilson B-factor (Å ²)	23.3	Xtriage
Anisotropy	0.230	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 34.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	7623	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.19% 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: GOL, G5F, PEG, SO4

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	HHH	0.99	0/1654	1.12	0/2251
1	KKK	0.98	1/1657 (0.1%)	1.12	2/2256 (0.1%)
2	LLL	0.93	0/1763	1.10	1/2393 (0.0%)
2	MMM	0.96	0/1768	1.11	2/2399 (0.1%)
All	All	0.96	1/6842 (0.0%)	1.11	5/9299 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	KKK	100	ASP	N-CA	8.29	1.51	1.46

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	KKK	130	SER	CA-C-N	5.26	127.64	120.54
1	KKK	130	SER	C-N-CA	5.26	127.64	120.54
2	LLL	172	GLN	CB-CA-C	5.14	117.91	109.90
2	MMM	172	GLN	CB-CA-C	5.13	117.54	109.84
2	MMM	48	GLN	CB-CG-CD	-5.03	104.04	112.60

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	HHH	1608	0	1567	8	0
1	KKK	1614	0	1578	8	1
2	LLL	1721	0	1680	8	1
2	MMM	1726	0	1688	9	0
3	HHH	47	0	0	0	0
3	MMM	49	0	0	0	0
4	KKK	6	0	8	2	0
4	LLL	6	0	8	0	0
5	LLL	5	0	0	0	0
5	MMM	5	0	0	0	0
6	KKK	7	0	10	0	0
7	HHH	197	0	0	0	0
7	KKK	192	0	0	2	0
7	LLL	178	0	0	0	0
7	MMM	262	0	0	1	0
All	All	7623	0	6539	33	1

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 2.

All (33) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:KKK:104:ASP:OD1	2:MMM:51[A]:ARG:NH2	2.08	0.84
2:LLL:155:LYS:NZ	2:LLL:201:GLU:OE1	2.24	0.70
1:HHH:20:LEU:HG	1:HHH:83:MET:HE2	1.75	0.67
4:KKK:302:GOL:H11	7:KKK:441:HOH:O	1.96	0.66
2:MMM:51[A]:ARG:HG3	7:MMM:519:HOH:O	1.99	0.60
4:KKK:302:GOL:C1	7:KKK:441:HOH:O	2.51	0.58
1:HHH:181:LEU:C	1:HHH:181:LEU:HD12	2.34	0.53
1:KKK:68:PHE:O	1:KKK:69:ILE:HD13	2.09	0.52
2:MMM:51[A]:ARG:HH11	2:MMM:52:LEU:H	1.55	0.52
2:MMM:51[A]:ARG:NH1	2:MMM:52:LEU:H	2.07	0.52
2:LLL:100:PRO:HA	2:LLL:101:PRO:C	2.35	0.52
1:KKK:181:LEU:C	1:KKK:181:LEU:HD12	2.37	0.49
1:HHH:202:ASN:ND2	1:HHH:209:LYS:HE2	2.29	0.48
2:LLL:39:LEU:HD22	2:LLL:77:PHE:CG	2.48	0.48
2:MMM:39:LEU:HD22	2:MMM:77:PHE:CG	2.49	0.47
2:MMM:39:LEU:HD22	2:MMM:77:PHE:CB	2.44	0.47
2:MMM:100:PRO:HA	2:MMM:101:PRO:C	2.38	0.47
1:KKK:22:CYS:HB3	1:KKK:79:LEU:HB3	1.97	0.46
1:HHH:22:CYS:HB3	1:HHH:79:LEU:HB3	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:LLL:205:GLN:O	1:KKK:1:GLU:HB2	2.16	0.45
1:HHH:51:THR:OG1	1:HHH:55:GLY:HA2	2.17	0.44
2:LLL:84:LEU:HD11	2:LLL:110:LEU:HD21	1.99	0.44
1:HHH:98:ARG:CZ	1:HHH:100[A]:ASP:OD1	2.65	0.44
1:HHH:83:MET:HB3	1:HHH:86:LEU:HD21	1.99	0.43
2:LLL:181:LEU:C	2:LLL:181:LEU:HD23	2.43	0.43
1:KKK:83:MET:HE2	1:KKK:86:LEU:HD21	1.99	0.43
2:LLL:89:VAL:HG21	2:LLL:172:GLN:HB3	2.01	0.42
1:KKK:48:ILE:HD13	1:KKK:48:ILE:HA	1.86	0.42
2:MMM:181:LEU:C	2:MMM:181:LEU:HD23	2.45	0.42
2:LLL:39:LEU:HD22	2:LLL:77:PHE:CB	2.51	0.41
1:HHH:204:LYS:N	1:HHH:205:PRO:CD	2.84	0.41
2:MMM:46:PRO:HG2	2:MMM:171:GLU:CD	2.46	0.41
1:KKK:204:LYS:N	1:KKK:205:PRO:CD	2.84	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:LLL:33:SER:O	1:KKK:31:SER:OG[3_544]	2.10	0.10

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	HHH	213/236 (90%)	207 (97%)	6 (3%)	0	100	100
1	KKK	213/236 (90%)	207 (97%)	6 (3%)	0	100	100
2	LLL	218/220 (99%)	215 (99%)	3 (1%)	0	100	100
2	MMM	218/220 (99%)	215 (99%)	3 (1%)	0	100	100
All	All	862/912 (94%)	844 (98%)	18 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	HHH	179/194 (92%)	177 (99%)	2 (1%)	65	44
1	KKK	181/194 (93%)	180 (99%)	1 (1%)	78	66
2	LLL	196/196 (100%)	195 (100%)	1 (0%)	81	71
2	MMM	196/196 (100%)	195 (100%)	1 (0%)	81	71
All	All	752/780 (96%)	747 (99%)	5 (1%)	76	62

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	HHH	50	PHE
1	HHH	132	LYS
2	LLL	97	TYR
1	KKK	50	PHE
2	MMM	97	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

7 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	G5F	HHH	301	-	47,47,50	0.44	0	64,67,71	0.79	3 (4%)
3	G5F	MMM	301	-	49,49,50	0.48	0	67,70,71	0.81	2 (2%)
6	PEG	KKK	301	-	6,6,6	0.36	0	5,5,5	0.30	0
4	GOL	KKK	302	-	5,5,5	0.06	0	5,5,5	0.29	0
4	GOL	LLL	301	-	5,5,5	0.23	0	5,5,5	0.69	0
5	SO4	LLL	302	-	4,4,4	0.18	0	6,6,6	0.12	0
5	SO4	MMM	302	-	4,4,4	0.27	0	6,6,6	0.26	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
3	G5F	HHH	301	-	-	7/56/76/82	0/1/1/1
3	G5F	MMM	301	-	-	7/60/80/82	0/1/1/1
6	PEG	KKK	301	-	-	1/4/4/4	-
4	GOL	LLL	301	-	-	4/4/4/4	-
4	GOL	KKK	302	-	-	2/4/4/4	-

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	MMM	301	G5F	O17-C18-C17	3.04	115.85	108.93
3	HHH	301	G5F	C4-C3-N1	-2.47	106.77	110.92
3	HHH	301	G5F	O21-P1-O23	2.27	123.01	112.44
3	MMM	301	G5F	O22-P2-O24	2.08	122.14	112.44
3	HHH	301	G5F	O9-C16-C17	2.05	114.83	109.36

There are no chirality outliers.

All (21) torsion outliers are listed below:

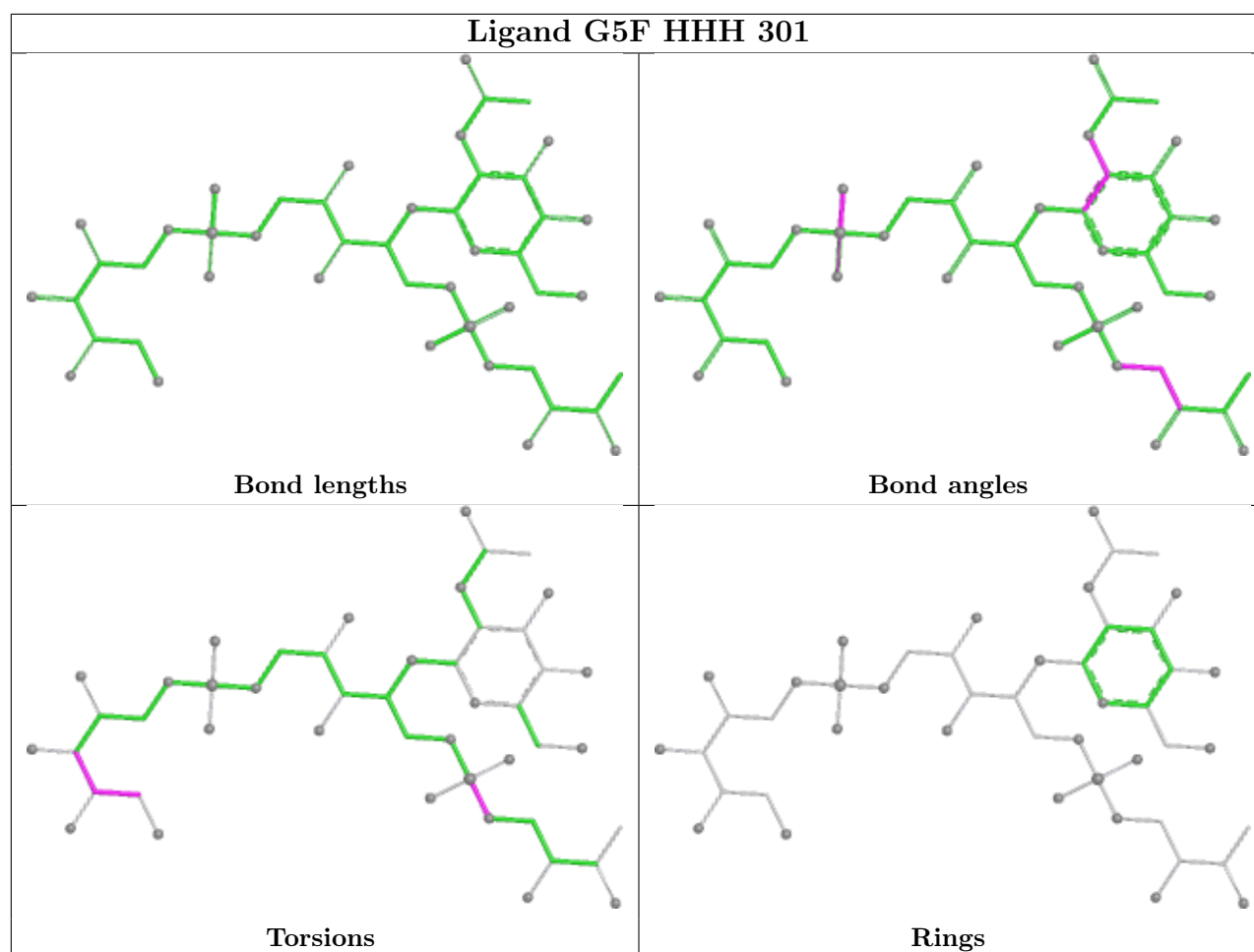
Mol	Chain	Res	Type	Atoms
3	HHH	301	G5F	O19-C15-C22-O14
3	MMM	301	G5F	C16-C17-C18-O17
3	MMM	301	G5F	O15-C17-C18-C19
3	MMM	301	G5F	O15-C17-C18-O17
3	MMM	301	G5F	C17-C18-C19-O13
4	LLL	301	GOL	O1-C1-C2-C3
4	LLL	301	GOL	C1-C2-C3-O3
3	HHH	301	G5F	C14-C15-C22-O14
3	MMM	301	G5F	C16-C17-C18-C19
4	KKK	302	GOL	C1-C2-C3-O3
3	MMM	301	G5F	C17-C18-C19-C23
3	MMM	301	G5F	O17-C18-C19-O13
4	LLL	301	GOL	O2-C2-C3-O3
3	HHH	301	G5F	C13-C14-C15-O19
6	KKK	301	PEG	C1-C2-O2-C3
4	KKK	302	GOL	O2-C2-C3-O3
3	HHH	301	G5F	C16-O9-P2-O22
3	HHH	301	G5F	C16-O9-P2-O6
4	LLL	301	GOL	O1-C1-C2-O2
3	HHH	301	G5F	O18-C14-C15-O19
3	HHH	301	G5F	O18-C14-C15-C22

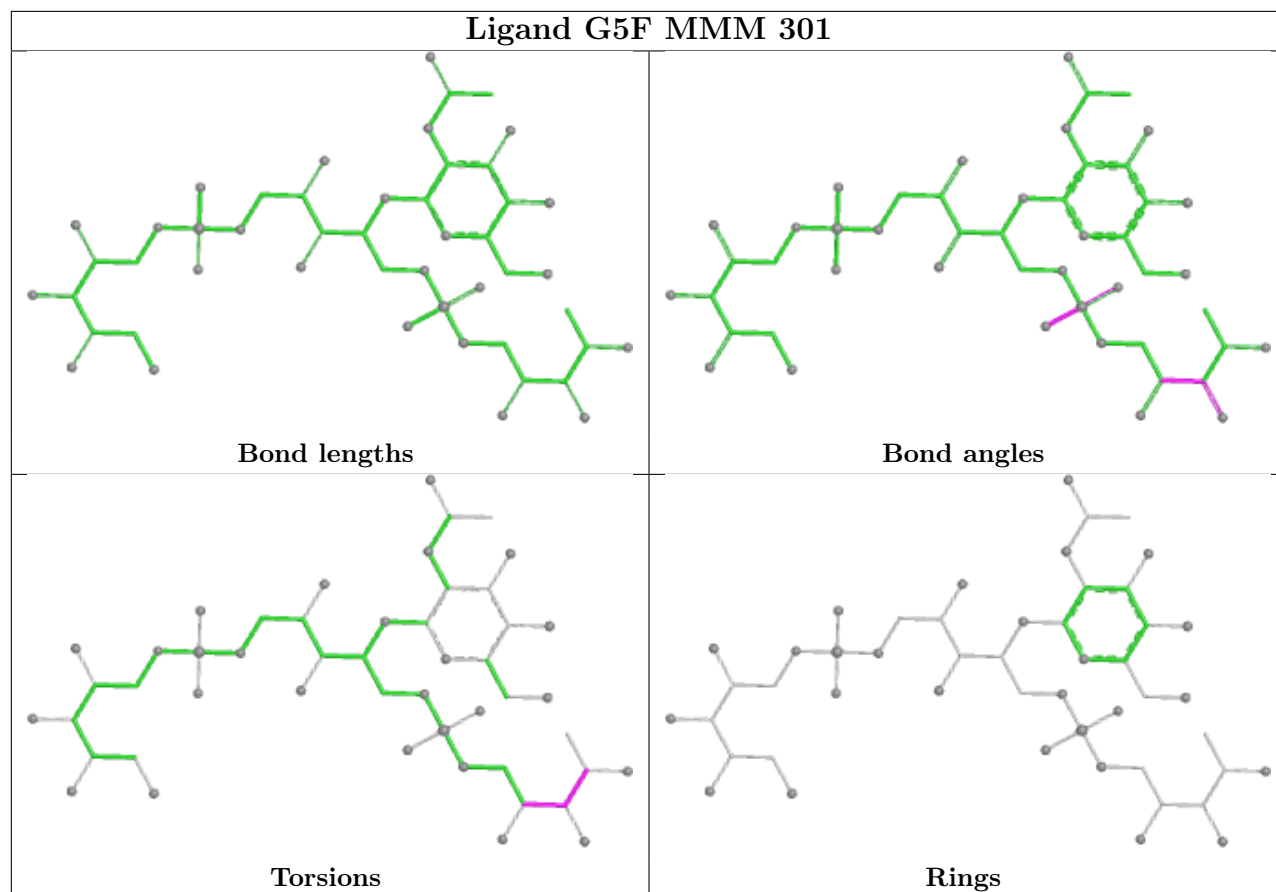
There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	KKK	302	GOL	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	HHH	214/236 (90%)	0.23	9 (4%)	40	45	15, 24, 41, 84	3 (1%)
1	KKK	215/236 (91%)	0.33	17 (7%)	18	21	11, 24, 50, 88	2 (0%)
2	LLL	219/220 (99%)	0.22	11 (5%)	34	39	17, 26, 47, 77	1 (0%)
2	MMM	219/220 (99%)	-0.01	9 (4%)	41	46	13, 21, 39, 62	1 (0%)
All	All	867/912 (95%)	0.19	46 (5%)	32	36	11, 24, 46, 88	7 (0%)

All (46) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	KKK	136	GLY	6.9
2	MMM	46	PRO	5.7
1	KKK	132	LYS	5.0
1	HHH	136	GLY	4.8
1	KKK	26	GLY	4.6
1	KKK	218	SER	4.1
1	KKK	137	GLY	3.9
1	HHH	132	LYS	3.9
1	KKK	131	SER	3.5
1	HHH	193	GLY	3.2
1	KKK	29	PHE	3.2
2	LLL	160	LEU	3.2
1	KKK	28	SER	3.1
1	KKK	1	GLU	3.0
2	MMM	47	GLY	2.9
1	HHH	194	THR	2.8
1	KKK	25	SER	2.8
1	KKK	2	VAL	2.8
2	LLL	73	PHE	2.7
2	MMM	73	PHE	2.7
2	MMM	133	SER	2.6

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Mol	Chain	Res	Type	RSRZ
2	MMM	219	GLU	2.6
2	LLL	218	GLY	2.5
2	LLL	219	GLU	2.5
1	KKK	130	SER	2.5
1	KKK	194	THR	2.4
2	MMM	51[A]	ARG	2.4
1	KKK	27	PHE	2.4
2	LLL	133	SER	2.4
2	LLL	208	SER	2.3
1	KKK	76	LYS	2.3
1	KKK	30	ASN	2.3
1	HHH	137	GLY	2.3
1	HHH	1	GLU	2.3
1	HHH	131	SER	2.3
2	LLL	207	LEU	2.2
2	LLL	196	LYS	2.2
2	MMM	218	GLY	2.2
2	MMM	32	THR	2.2
1	HHH	217	LYS	2.1
2	LLL	158	ASN	2.1
1	KKK	69	ILE	2.1
2	MMM	45	ARG	2.1
1	HHH	76	LYS	2.1
2	LLL	159	ALA	2.0
2	LLL	132	LYS	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

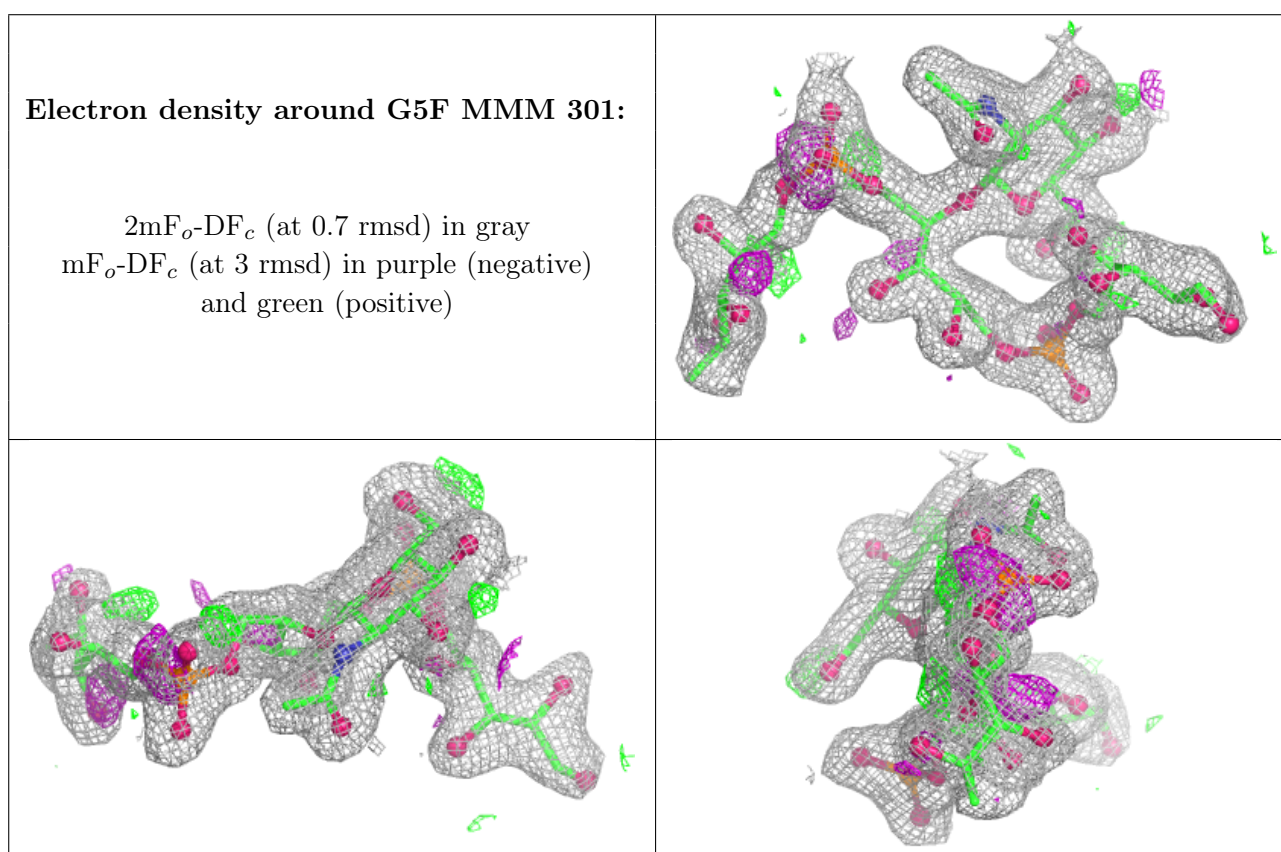
There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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

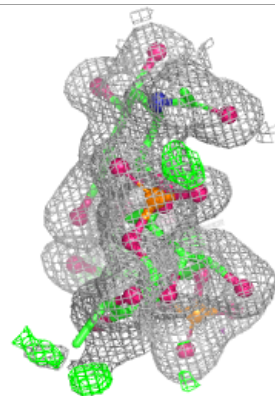
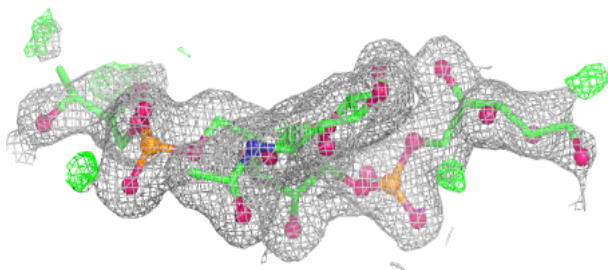
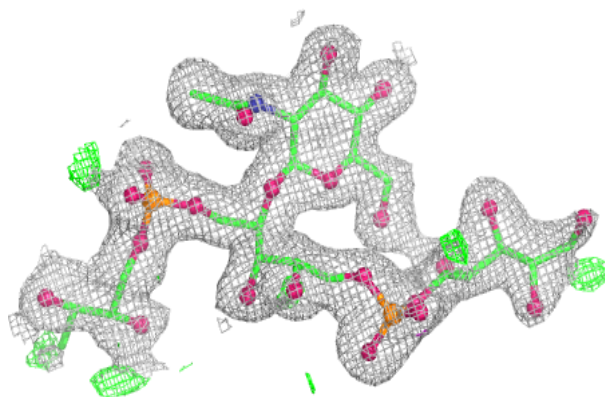
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
6	PEG	KKK	301	7/7	0.74	0.25	55,63,77,77	0
5	SO4	MMM	302	5/5	0.79	0.15	47,59,67,71	0
4	GOL	KKK	302	6/6	0.87	0.14	44,48,50,55	0
4	GOL	LLL	301	6/6	0.91	0.12	27,35,48,52	0
5	SO4	LLL	302	5/5	0.91	0.10	39,44,51,52	0
3	G5F	MMM	301	49/50	0.95	0.09	22,29,47,50	0
3	G5F	HHH	301	47/50	0.96	0.08	20,29,55,58	0

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



Electron density around G5F HHH 301:

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



6.5 Other polymers [i](#)

There are no such residues in this entry.