



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 7, 2020 – 04:36 PM BST

PDB ID : 6CXE
Title : Structure of alpha-GSA[26,6P] bound by CD1d and in complex with the Va14Vb8.2 TCR
Authors : Wang, J.; Zajonc, D.
Deposited on : 2018-04-02
Resolution : 2.05 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.13.1
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) : 2.13.1

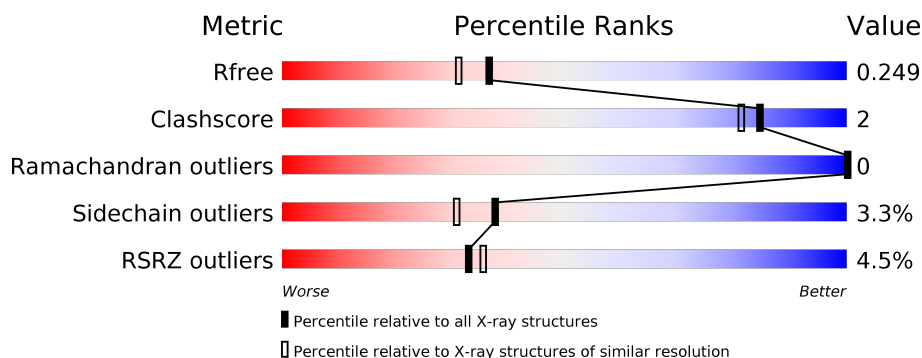
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.05 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	1692 (2.04-2.04)
Clashscore	141614	1773 (2.04-2.04)
Ramachandran outliers	138981	1752 (2.04-2.04)
Sidechain outliers	138945	1752 (2.04-2.04)
RSRZ outliers	127900	1672 (2.04-2.04)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	C	209	<div> <div>5%</div> <div> <div></div> <div>93%</div> <div></div> </div> <div></div> </div>
2	D	241	<div> <div></div> <div> <div></div> <div>93%</div> <div></div> </div> <div>5%</div> </div>
3	A	285	<div> <div>8%</div> <div> <div></div> <div>83%</div> <div></div> </div> <div>11% 6%</div> </div>
4	B	99	<div> <div>2%</div> <div> <div></div> <div>90%</div> <div></div> </div> <div>6%</div> </div>
5	E	3	<div> <div></div> <div>100%</div> <div></div> </div>

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 6787 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 Chimeric T cell antigen receptor alpha chain Va14,Va24,Ja18.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	202	Total	C	N	O	S	0	0	0
			1536	950	262	316	8			

- Molecule 2 is a protein called Chimeric T cell antigen receptor beta chain Vb8.2, vb11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	239	Total	C	N	O	S	0	0	0
			1872	1175	333	358	6			

- Molecule 3 is a protein called Antigen-presenting glycoprotein CD1d1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	269	Total	C	N	O	S	0	0	0
			2165	1381	372	399	13			

There are 6 discrepancies between the modelled and reference sequences:

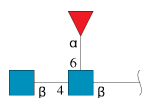
Chain	Residue	Modelled	Actual	Comment	Reference
A	280	HIS	-	expression tag	UNP A0A0R4J090
A	281	HIS	-	expression tag	UNP A0A0R4J090
A	282	HIS	-	expression tag	UNP A0A0R4J090
A	283	HIS	-	expression tag	UNP A0A0R4J090
A	284	HIS	-	expression tag	UNP A0A0R4J090
A	285	HIS	-	expression tag	UNP A0A0R4J090

- Molecule 4 is a protein called Beta-2-microglobulin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	B	97	Total	C	N	O	S	0	0	0
			801	512	136	147	6			

- Molecule 5 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[al

pha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose.

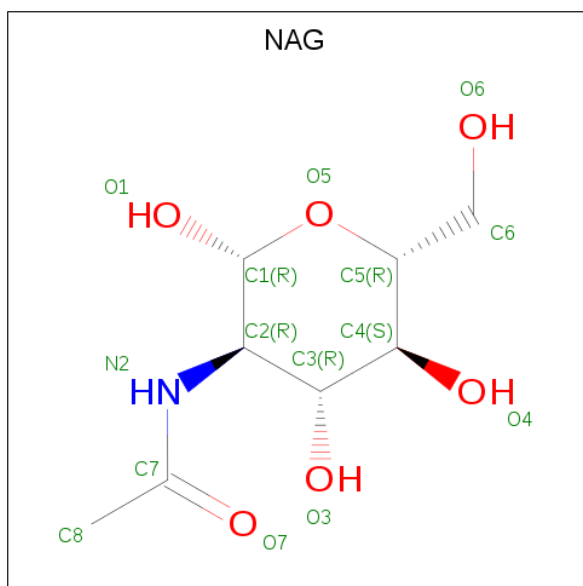


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	E	3	Total	C	N	O	0	0	0
			38	22	2	14			

- Molecule 6 is SODIUM ION (three-letter code: NA) (formula: Na).

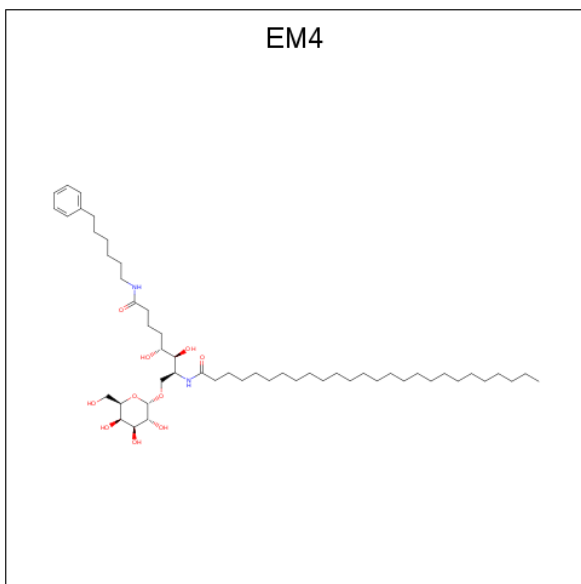
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	Na	0	0
			1	1		
6	D	2	Total	Na	0	0
			2	2		
6	C	1	Total	Na	0	0
			1	1		

- Molecule 7 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C₈H₁₅NO₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	1	Total	C	N	O	0	0
			14	8	1	5		
7	A	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 8 is N-[(2S,3S,4R)-3,4-dihydroxy-8-oxo-8-[(6-phenylhexyl)amino]-1-{[(2S,3R,4S,5R,6R)-3,4,5-trihydroxy-6-(hydroxymethyl)tetrahydro-2H-pyran-2-yl]oxy}octan-2-yl]hexacosamide (three-letter code: EM4) (formula: C₅₂H₉₄N₂O₁₀).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	A	1	Total	C	N	O	0	0
			64	52	2	10		

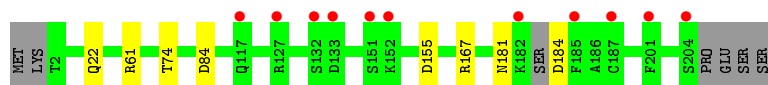
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	C	66	Total	O	0	0
			66	66		
9	D	111	Total	O	0	0
			111	111		
9	A	86	Total	O	0	0
			86	86		
9	B	16	Total	O	0	0
			16	16		

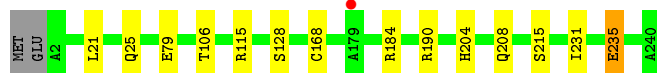
3 Residue-property plots

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.

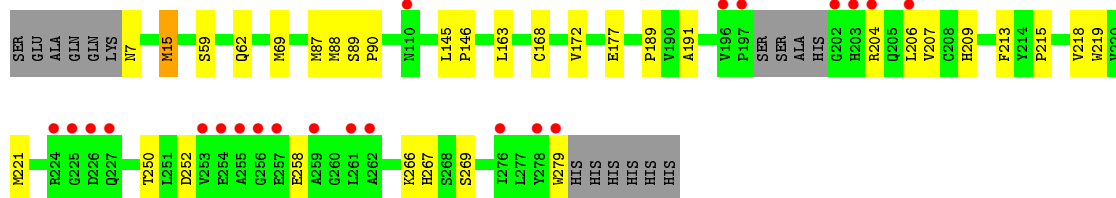
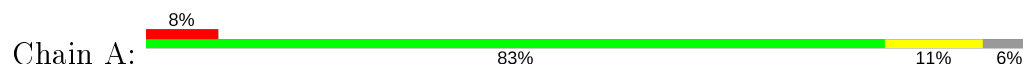
- Molecule 1: Chimeric T cell antigen receptor alpha chain Va14,Va24,Ja18



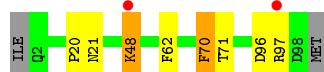
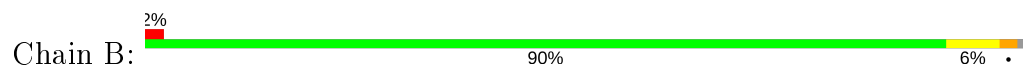
- Molecule 2: Chimeric T cell antigen receptor beta chain Vb8.2, vb11



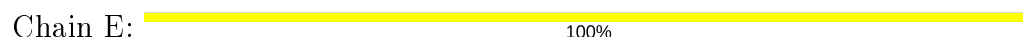
- Molecule 3: Antigen-presenting glycoprotein CD1d1



- Molecule 4: Beta-2-microglobulin



- Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose



MAG1
MAG2
FUC3

4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	78.44Å 189.07Å 150.91Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.05 44.41 – 2.05	Depositor EDS
% Data completeness (in resolution range)	95.9 (50.00-2.05) 96.0 (44.41-2.05)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.65 (at 2.05Å)	Xtriage
Refinement program	REFMAC 5.8.0155	Depositor
R, R_{free}	0.209 , 0.248 0.211 , 0.249	Depositor DCC
R_{free} test set	1388 reflections (2.05%)	wwPDB-VP
Wilson B-factor (Å ²)	41.5	Xtriage
Anisotropy	0.143	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 40.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6787	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.21% 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: FUC, NA, NAG, EM4

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	C	0.48	0/1563	0.68	0/2128
2	D	0.50	0/1923	0.63	0/2622
3	A	0.50	0/2228	0.63	0/3028
4	B	0.49	0/827	0.63	0/1123
All	All	0.49	0/6541	0.64	0/8901

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	C	1536	0	1440	2	0
2	D	1872	0	1774	3	0
3	A	2165	0	2071	21	0
4	B	801	0	774	4	0
5	E	38	0	34	0	0
6	A	1	0	0	0	0
6	C	1	0	0	0	0
6	D	2	0	0	0	0
7	A	28	0	26	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	A	64	0	0	0	0
9	A	86	0	0	0	0
9	B	16	0	0	0	0
9	C	66	0	0	0	0
9	D	111	0	0	0	0
All	All	6787	0	6119	29	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 2.

All (29) 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)
3:A:59:SER:H	3:A:62:GLN:HE21	1.30	0.79
3:A:59:SER:H	3:A:62:GLN:NE2	1.94	0.66
3:A:15:MET:HG2	4:B:62:PHE:HE2	1.60	0.66
3:A:88:MET:HE3	3:A:146:PRO:HD3	1.79	0.65
2:D:204:HIS:NE2	2:D:235:GLU:HG2	2.12	0.64
3:A:215:PRO:O	3:A:267:HIS:HE1	1.83	0.60
3:A:267:HIS:CD2	3:A:269:SER:H	2.21	0.59
3:A:219:TRP:HB3	3:A:266:LYS:HB2	1.86	0.57
4:B:48:LYS:H	4:B:48:LYS:HD3	1.70	0.57
4:B:20:PRO:HA	4:B:71:THR:HG22	1.89	0.54
3:A:87:MET:SD	3:A:88:MET:HE1	2.51	0.50
3:A:267:HIS:HD2	3:A:269:SER:OG	1.94	0.50
3:A:258:GLU:HB3	3:A:279:TRP:CD1	2.48	0.48
3:A:59:SER:N	3:A:62:GLN:HE21	2.04	0.48
4:B:21:ASN:HB3	4:B:70:PHE:CE1	2.49	0.47
3:A:219:TRP:CZ2	3:A:221:MET:HG3	2.49	0.47
1:C:22:GLN:OE1	1:C:74:THR:HG22	2.15	0.46
3:A:89:SER:HA	3:A:90:PRO:HA	1.74	0.45
2:D:208:GLN:HG3	2:D:231:ILE:HG23	1.98	0.45
3:A:189:PRO:HB3	3:A:213:PHE:HB3	1.99	0.43
3:A:218:VAL:HG23	3:A:267:HIS:HB2	2.00	0.43
1:C:61:ARG:NH2	1:C:84:ASP:OD2	2.40	0.42
3:A:168:CYS:O	3:A:172:VAL:HG23	2.19	0.42
3:A:69:MET:CE	3:A:163:LEU:HD21	2.49	0.42
3:A:88:MET:CE	3:A:146:PRO:HD3	2.48	0.41
3:A:191:ALA:HA	3:A:209:HIS:O	2.21	0.41
3:A:207:VAL:HG22	3:A:250:THR:HG22	2.02	0.41
3:A:145:LEU:HB3	3:A:146:PRO:HD3	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:21:LEU:HD22	2:D:106:THR:HG21	2.02	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	C	198/209 (95%)	191 (96%)	7 (4%)	0	100	100
2	D	237/241 (98%)	235 (99%)	2 (1%)	0	100	100
3	A	265/285 (93%)	261 (98%)	4 (2%)	0	100	100
4	B	95/99 (96%)	92 (97%)	3 (3%)	0	100	100
All	All	795/834 (95%)	779 (98%)	16 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	C	174/188 (93%)	170 (98%)	4 (2%)	50	44
2	D	202/208 (97%)	193 (96%)	9 (4%)	27	20
3	A	235/249 (94%)	229 (97%)	6 (3%)	46	39
4	B	91/93 (98%)	87 (96%)	4 (4%)	28	21

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	702/738 (95%)	679 (97%)	23 (3%)	38	31

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	155	ASP
1	C	167	ARG
1	C	181	ASN
1	C	184	ASP
2	D	25	GLN
2	D	79	GLU
2	D	115	ARG
2	D	128	SER
2	D	168	CYS
2	D	184	ARG
2	D	190	ARG
2	D	215	SER
2	D	235	GLU
3	A	7	ASN
3	A	15	MET
3	A	177	GLU
3	A	204	ARG
3	A	206	LEU
3	A	252	ASP
4	B	48	LYS
4	B	70	PHE
4	B	96	ASP
4	B	97	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
2	D	230	GLN
3	A	62	GLN
3	A	267	HIS

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 ⓘ

3 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	NAG	E	1	3,5	14,14,15	0.55	0	17,19,21	1.17	3 (17%)
5	NAG	E	2	5	14,14,15	0.53	0	17,19,21	1.31	1 (5%)
5	FUC	E	3	5	10,10,11	0.91	0	14,14,16	1.07	1 (7%)

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	NAG	E	1	3,5	-	1/6/23/26	0/1/1/1
5	NAG	E	2	5	-	1/6/23/26	0/1/1/1
5	FUC	E	3	5	-	-	0/1/1/1

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	E	2	NAG	C1-O5-C5	4.53	118.33	112.19
5	E	1	NAG	O5-C5-C6	2.33	110.86	107.20
5	E	1	NAG	C1-O5-C5	2.30	115.31	112.19
5	E	1	NAG	C1-C2-N2	2.15	114.16	110.49
5	E	3	FUC	O5-C5-C6	2.13	111.90	107.33

There are no chirality outliers.

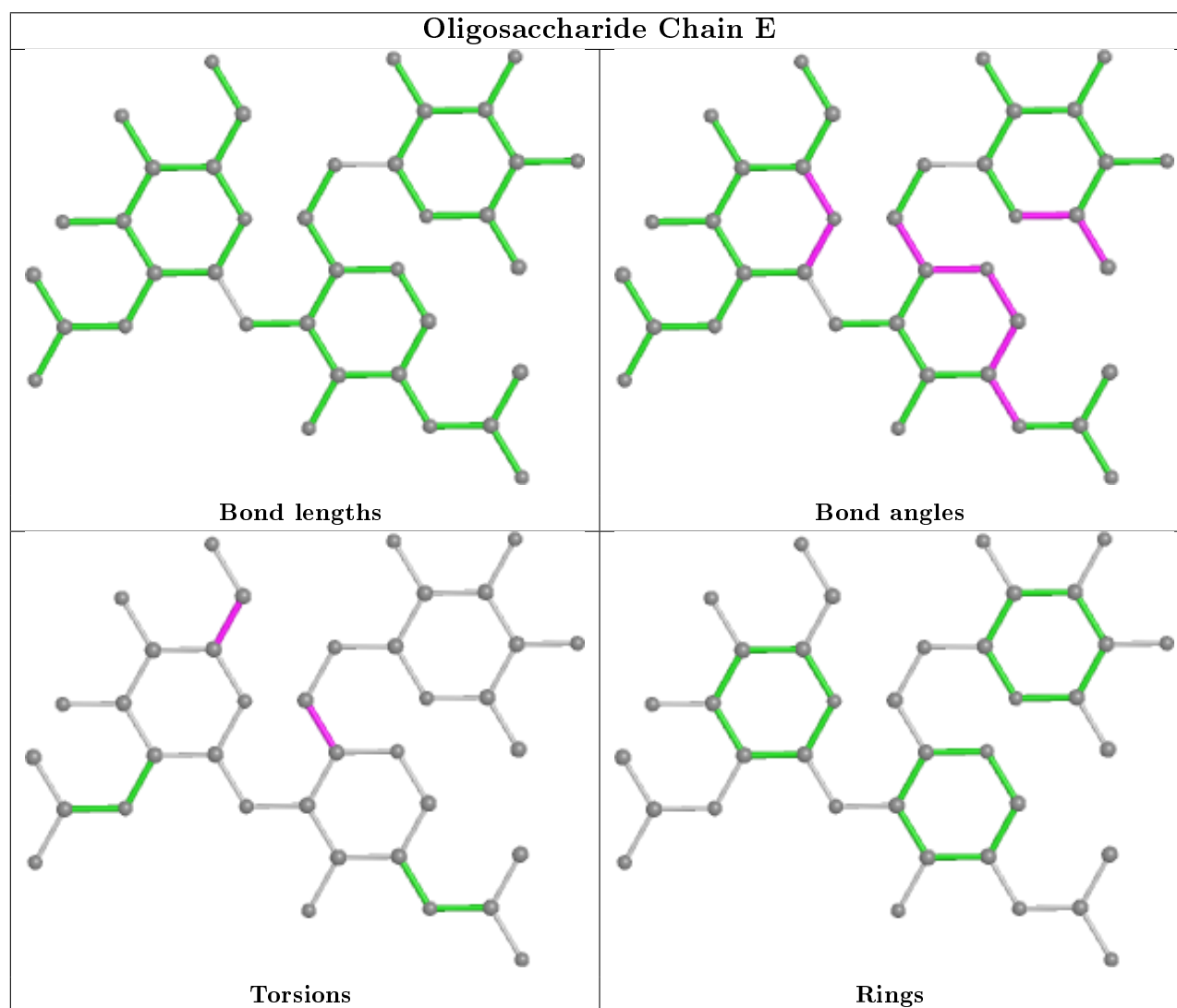
All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	E	2	NAG	O5-C5-C6-O6
5	E	1	NAG	O5-C5-C6-O6

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 7 ligands modelled in this entry, 4 are monoatomic - leaving 3 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$
7	NAG	A	301	3	14,14,15	0.60	0	17,19,21	1.16	2 (11%)
8	EM4	A	306	-	65,65,65	0.73	1 (1%)	73,77,77	0.82	1 (1%)
7	NAG	A	302	3	14,14,15	0.75	0	17,19,21	1.20	1 (5%)

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
7	NAG	A	301	3	-	0/6/23/26	0/1/1/1
8	EM4	A	306	-	-	18/60/80/80	0/2/2/2
7	NAG	A	302	3	-	0/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	A	306	EM4	O9-C31	2.19	1.47	1.41

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	A	302	NAG	C1-O5-C5	4.01	117.63	112.19
7	A	301	NAG	C1-O5-C5	3.14	116.44	112.19
8	A	306	EM4	C31-O9-C30	2.52	118.64	113.69
7	A	301	NAG	O5-C5-C6	2.37	110.93	107.20

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
8	A	306	EM4	C17-C18-C19-N1
8	A	306	EM4	C5-C6-C7-C40
8	A	306	EM4	C46-C47-C48-C49

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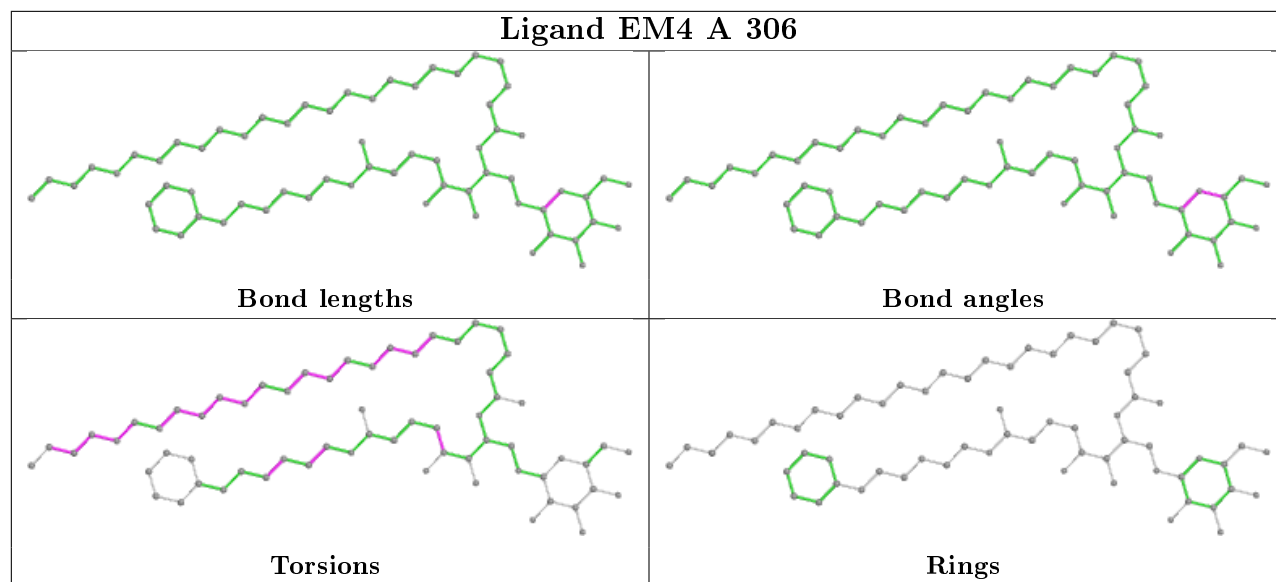
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Mol	Chain	Res	Type	Atoms
8	A	306	EM4	C41-C42-C43-C35
8	A	306	EM4	C48-C49-C50-C51
8	A	306	EM4	C47-C48-C49-C50
8	A	306	EM4	C43-C35-C36-C37
8	A	306	EM4	C7-C40-C41-C42
8	A	306	EM4	C36-C37-C38-C39
8	A	306	EM4	C38-C39-C44-C45
8	A	306	EM4	C49-C50-C51-C52
8	A	306	EM4	C36-C35-C43-C42
8	A	306	EM4	C44-C45-C46-C47
8	A	306	EM4	C23-C24-C25-O2
8	A	306	EM4	C41-C40-C7-C6
8	A	306	EM4	C15-C16-C17-C18
8	A	306	EM4	C37-C38-C39-C44
8	A	306	EM4	C39-C44-C45-C46

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 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	C	202/209 (96%)	0.15	11 (5%) 25 28	30, 46, 83, 104	0
2	D	239/241 (99%)	-0.22	1 (0%) 92 93	31, 42, 61, 87	0
3	A	269/285 (94%)	0.37	22 (8%) 11 12	32, 44, 95, 123	0
4	B	97/99 (97%)	0.11	2 (2%) 63 67	37, 53, 81, 94	0
All	All	807/834 (96%)	0.11	36 (4%) 33 35	30, 44, 84, 123	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	152	LYS	7.9
3	A	203	HIS	5.8
3	A	254	GLU	4.8
3	A	204	ARG	4.1
3	A	276	ILE	3.8
3	A	110	ASN	3.6
3	A	253	VAL	3.5
3	A	256	GLY	3.5
3	A	255	ALA	3.5
3	A	279	TRP	3.4
1	C	132	SER	3.3
3	A	202	GLY	3.2
3	A	226	ASP	3.1
1	C	117	GLN	3.1
2	D	179	ALA	3.1
3	A	278	TYR	3.0
3	A	262	ALA	2.9
3	A	259	ALA	2.9
3	A	225	GLY	2.9
1	C	182	LYS	2.8
3	A	196	VAL	2.8

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Mol	Chain	Res	Type	RSRZ
3	A	261	LEU	2.8
1	C	187	CYS	2.6
4	B	97	ARG	2.6
3	A	227	GLN	2.4
1	C	151	SER	2.4
3	A	224	ARG	2.4
4	B	48	LYS	2.3
3	A	197	PRO	2.3
3	A	206	LEU	2.2
1	C	201	PHE	2.2
1	C	133	ASP	2.2
1	C	204	SER	2.2
1	C	185	PHE	2.1
1	C	127	ARG	2.1
3	A	257	GLU	2.1

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

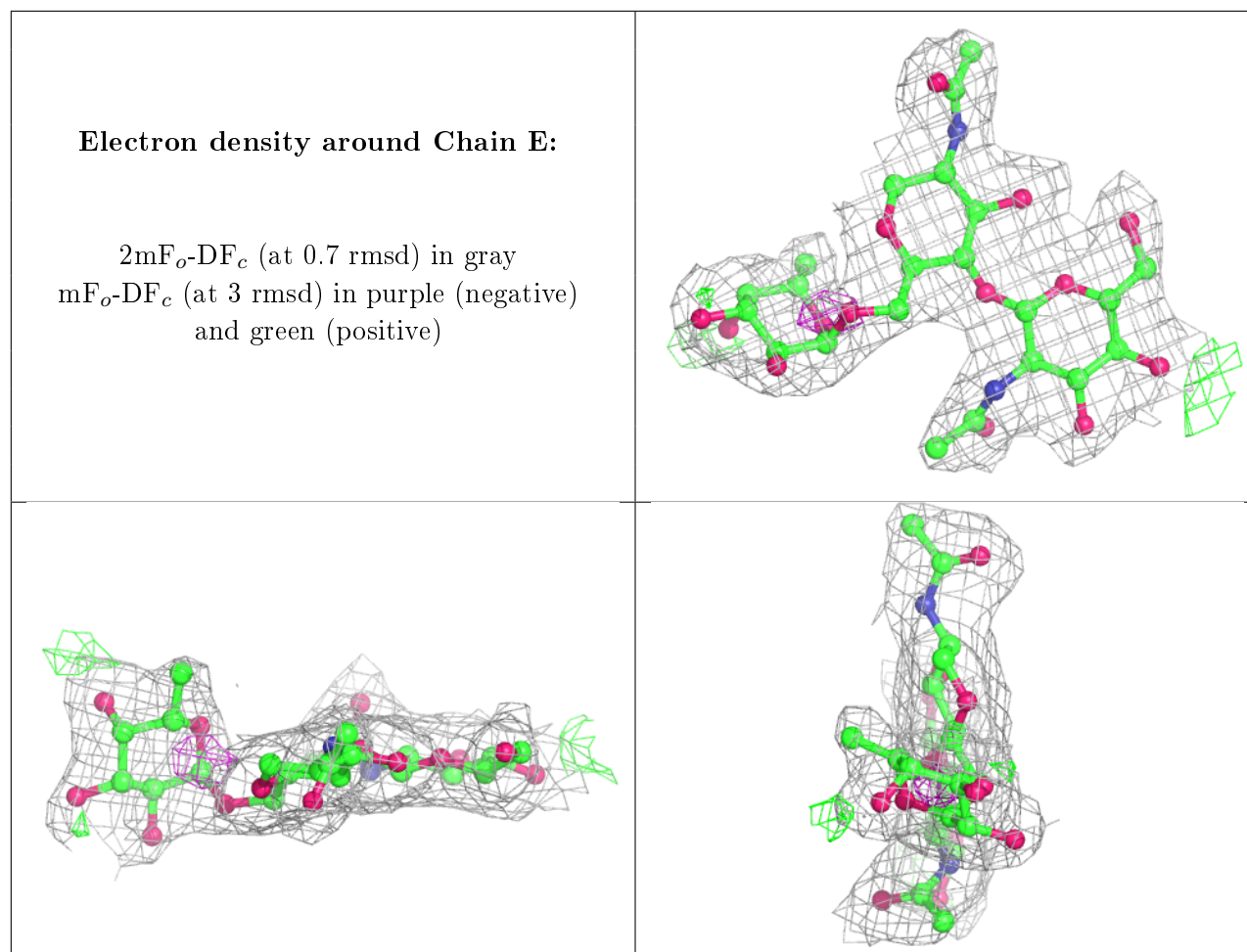
There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	FUC	E	3	10/11	0.73	0.18	68,71,74,75	0
5	NAG	E	2	14/15	0.93	0.13	59,66,70,73	0
5	NAG	E	1	14/15	0.95	0.09	42,48,60,67	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.



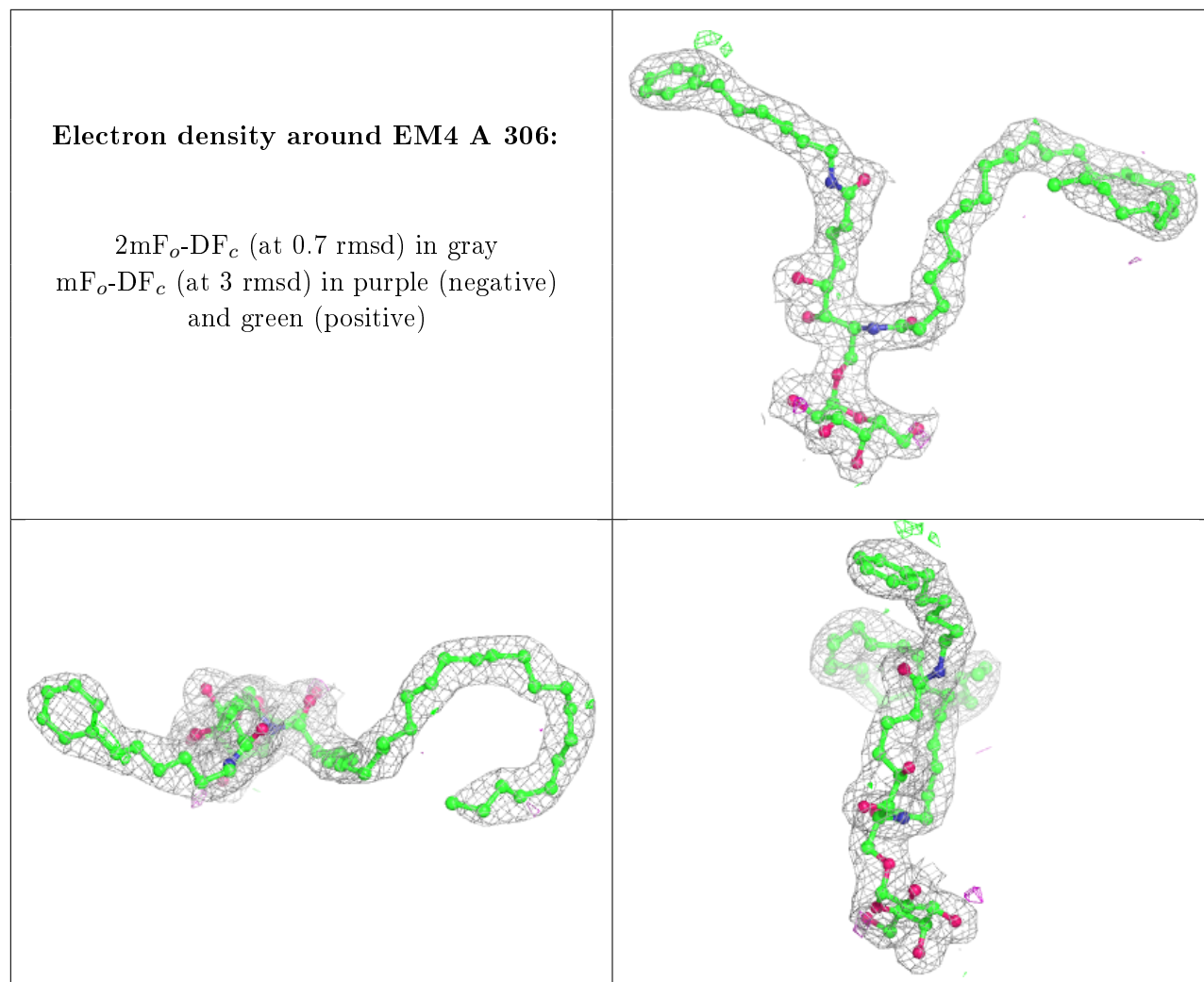
6.4 Ligands ⓘ

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
6	NA	D	302	1/1	0.89	0.16	64,64,64,64	0
7	NAG	A	301	14/15	0.90	0.16	61,67,71,75	0
6	NA	C	301	1/1	0.91	0.08	53,53,53,53	0
7	NAG	A	302	14/15	0.92	0.10	44,46,50,50	0
8	EM4	A	306	64/64	0.96	0.17	31,43,53,57	0
6	NA	D	301	1/1	0.98	0.09	43,43,43,43	0
6	NA	A	307	1/1	0.99	0.09	46,46,46,46	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 ⓘ

There are no such residues in this entry.