



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 21, 2020 – 02:19 PM BST

PDB ID : 5KN2  
Title : Native bovine skeletal calsequestrin, high-Ca<sup>2+</sup> form  
Authors : Lewis, K.M.; Byrd, S.S.; Kang, C.  
Deposited on : 2016-06-27  
Resolution : 2.60 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](mailto: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.

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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  
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

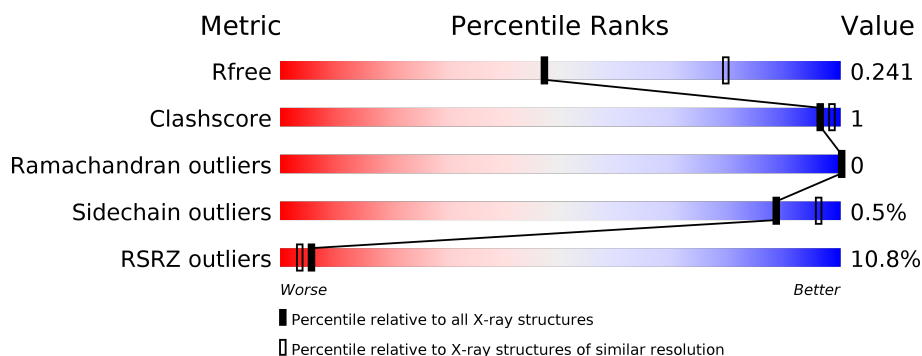
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.60 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	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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  $\geq 3$ , 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	361	<div> <div>2%</div> <div>95%</div> <div>5%</div> <div>• •</div> </div>
1	B	361	<div> <div>18%</div> <div>93%</div> <div>5%</div> <div>•</div> </div>
1	C	361	<div> <div>12%</div> <div>94%</div> <div>5%</div> <div>• •</div> </div>
2	D	2	<div> <div>50%</div> <div>50%</div> </div>
2	E	2	<div> <div>50%</div> <div>50%</div> </div>
2	F	2	<div> <div>50%</div> <div>50%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	CA	A	406	-	-	-	X
3	CA	A	407	-	-	-	X
3	CA	A	408	-	-	-	X
3	CA	A	409	-	-	-	X
3	CA	A	410	-	-	-	X
3	CA	C	410	-	-	-	X

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 16462 atoms, of which 7878 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 Calsequestrin.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	352	Total	C	H	N	O	S	0	0	0
			5448	1809	2618	420	595	6			
1	B	351	Total	C	H	N	O	S	0	0	0
			5395	1794	2585	420	590	6			
1	C	352	Total	C	H	N	O	S	0	0	0
			5412	1799	2596	420	591	6			

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



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	2	Total	C	H	N	O	0	0	0
			54	16	26	2	10			
2	E	2	Total	C	H	N	O	0	0	0
			55	16	27	2	10			
2	F	2	Total	C	H	N	O	0	0	0
			54	16	26	2	10			

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	14	Total	Ca	0	0
			14	14		
3	A	13	Total	Ca	0	0
			13	13		
3	C	12	Total	Ca	0	0
			12	12		

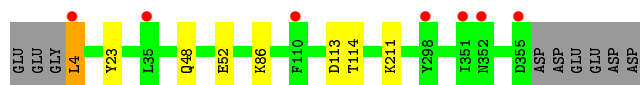
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	5	Total	O	0	0
			5	5		

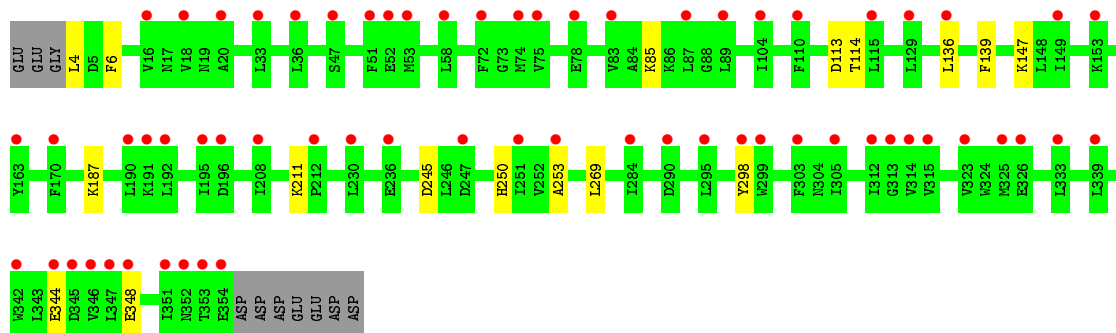
### 3 Residue-property plots [i](#)

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.

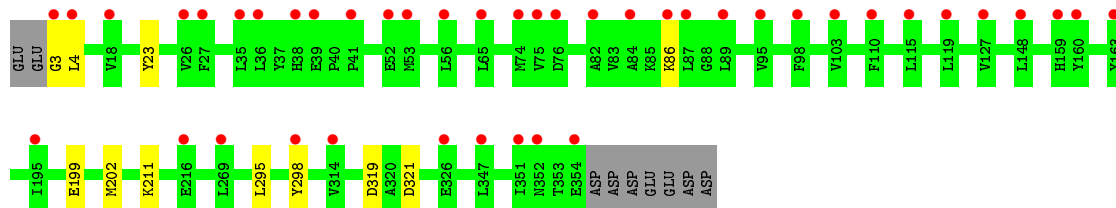
- Molecule 1: Calsequestrin



- Molecule 1: Calsequestrin



- Molecule 1: Calsequestrin



- Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose





- Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain E:  50% 50%



- Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F:  50% 50%



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	130.36Å 169.19Å 155.48Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.95 – 2.60 49.95 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.7 (49.95-2.60) 92.8 (49.95-2.60)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.23 (at 2.61Å)	Xtriage
Refinement program	PHENIX 1.10 _2155	Depositor
R, $R_{free}$	0.205 , 0.241 0.206 , 0.241	Depositor DCC
$R_{free}$ test set	2000 reflections (3.79%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	75.2	Xtriage
Anisotropy	0.212	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 63.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	16462	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	117.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.98% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: CA, NAG

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	0.34	0/2895	0.52	0/3941
1	B	0.32	0/2875	0.52	0/3913
1	C	0.32	0/2881	0.52	0/3922
All	All	0.33	0/8651	0.52	0/11776

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	A	2830	2618	2618	4	0
1	B	2810	2585	2586	10	0
1	C	2816	2596	2595	6	0
2	D	28	26	25	0	0
2	E	28	27	25	0	0
2	F	28	26	25	0	0
3	A	13	0	0	0	0
3	B	14	0	0	0	0
3	C	12	0	0	0	0
4	A	5	0	0	1	0
All	All	8584	7878	7874	17	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 1.

All (17) 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:A:23:TYR:OH	1:A:86:LYS:NZ	2.27	0.65
4:A:501:HOH:O	1:B:85:LYS:NZ	2.36	0.58
1:B:253:ALA:CB	1:B:269:LEU:HD11	2.43	0.48
1:B:139:PHE:O	1:B:147:LYS:NZ	2.41	0.48
1:A:4:LEU:HD13	1:A:4:LEU:O	2.15	0.47
1:B:344:GLU:O	1:B:348:GLU:HG3	2.15	0.46
1:B:6:PHE:HA	1:C:295:LEU:HD21	1.98	0.45
1:B:4:LEU:HA	1:C:298:TYR:OH	2.17	0.44
1:B:245:ASP:HA	1:B:250:HIS:CE1	2.52	0.44
1:C:319:ASP:OD2	1:C:321:ASP:HB2	2.18	0.43
1:A:48:GLN:O	1:A:52:GLU:HG3	2.19	0.43
1:B:136:LEU:HD21	1:B:187:LYS:HE3	2.01	0.41
1:A:113:ASP:OD1	1:A:114:THR:N	2.51	0.41
1:B:298:TYR:OH	1:C:3:GLY:HA3	2.20	0.41
1:C:23:TYR:OH	1:C:86:LYS:NZ	2.35	0.41
1:C:199:GLU:HB3	1:C:202:MET:HG3	2.01	0.41
1:B:113:ASP:OD1	1:B:114:THR:N	2.53	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	350/361 (97%)	343 (98%)	7 (2%)	0	100	100
1	B	349/361 (97%)	342 (98%)	7 (2%)	0	100	100
1	C	350/361 (97%)	343 (98%)	7 (2%)	0	100	100
All	All	1049/1083 (97%)	1028 (98%)	21 (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	A	308/323 (95%)	306 (99%)	2 (1%)	86	95
1	B	303/323 (94%)	302 (100%)	1 (0%)	92	98
1	C	304/323 (94%)	302 (99%)	2 (1%)	84	94
All	All	915/969 (94%)	910 (100%)	5 (0%)	88	96

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

Mol	Chain	Res	Type
1	A	4	LEU
1	A	211	LYS
1	B	211	LYS
1	C	4	LEU
1	C	211	LYS

Some 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 ⓘ

6 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	$\# Z  > 2$
2	NAG	D	1	1,2	14,14,15	0.75	1 (7%)	17,19,21	0.73	1 (5%)
2	NAG	D	2	2	14,14,15	0.27	0	17,19,21	0.62	0
2	NAG	E	1	1,2	14,14,15	0.69	1 (7%)	17,19,21	0.68	1 (5%)
2	NAG	E	2	2	14,14,15	0.51	0	17,19,21	0.49	0
2	NAG	F	1	1,2	14,14,15	0.58	0	17,19,21	0.72	1 (5%)
2	NAG	F	2	2	14,14,15	0.41	0	17,19,21	0.47	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
2	NAG	D	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	D	2	2	-	1/6/23/26	0/1/1/1
2	NAG	E	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	E	2	2	-	2/6/23/26	0/1/1/1
2	NAG	F	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	F	2	2	-	0/6/23/26	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	1	NAG	O5-C1	2.44	1.47	1.43
2	E	1	NAG	O5-C1	2.13	1.47	1.43

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	1	NAG	C1-O5-C5	2.48	115.55	112.19
2	E	1	NAG	C1-O5-C5	2.44	115.49	112.19
2	D	1	NAG	C1-O5-C5	2.28	115.28	112.19

There are no chirality outliers.

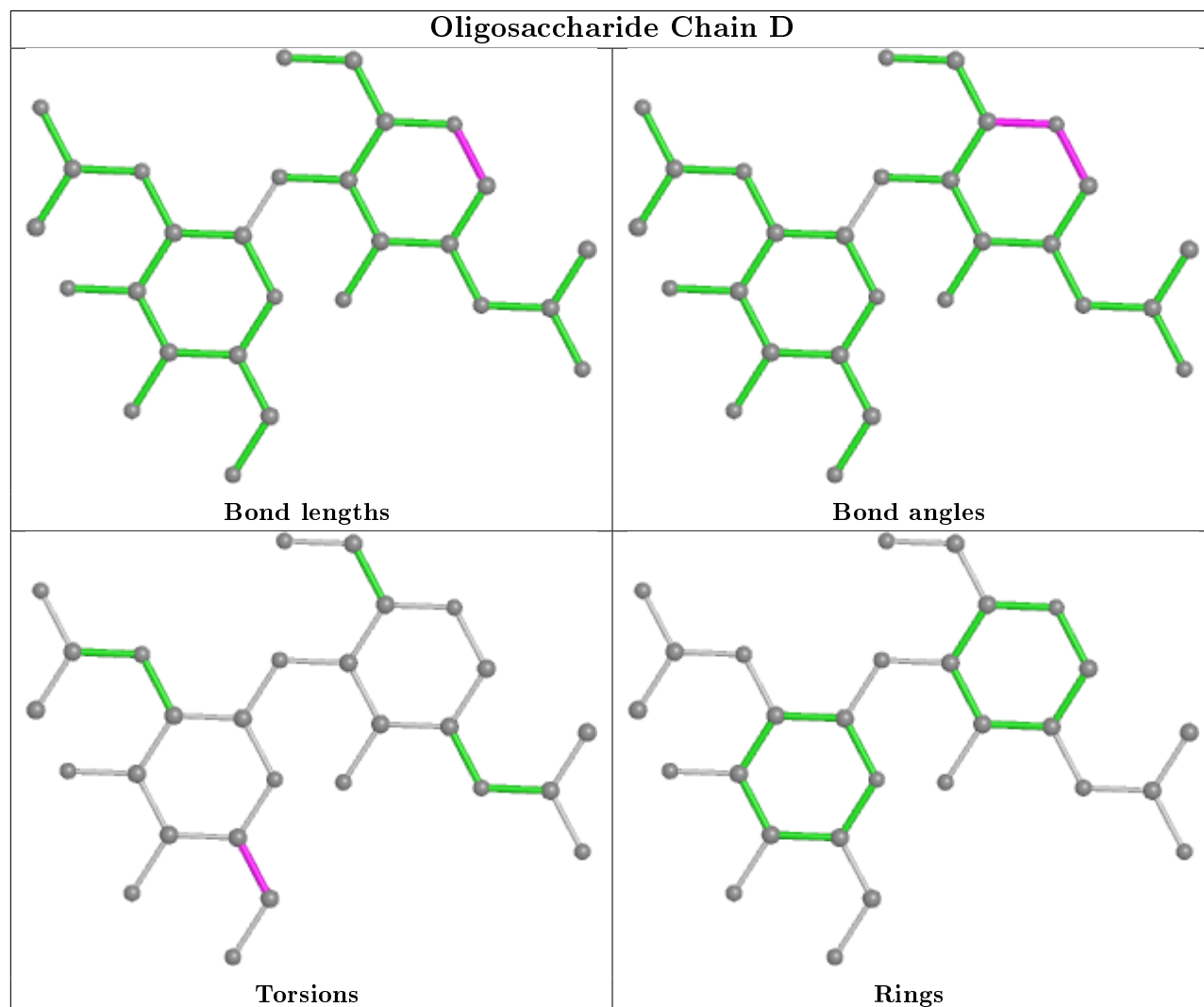
All (7) torsion outliers are listed below:

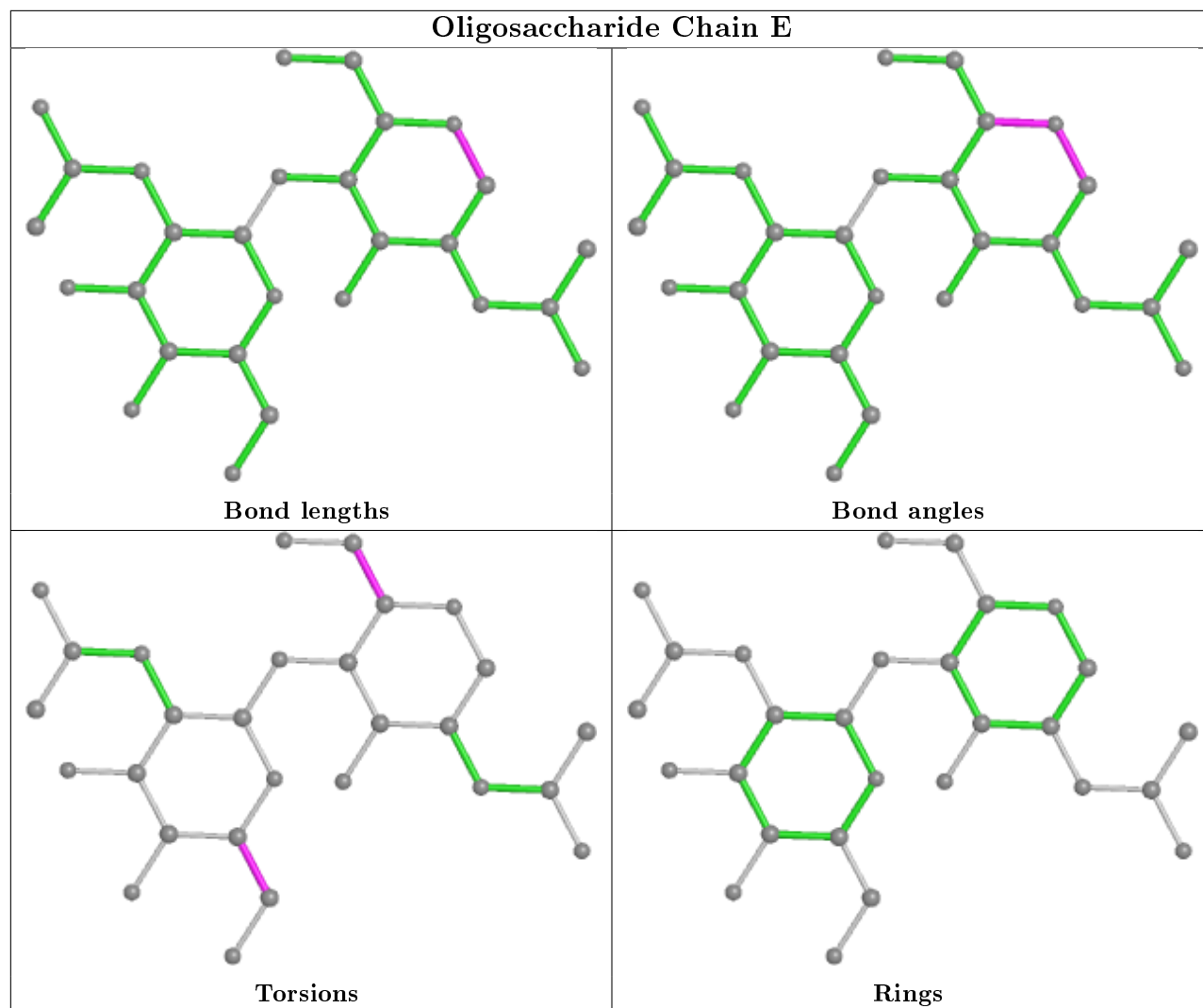
Mol	Chain	Res	Type	Atoms
2	F	1	NAG	O5-C5-C6-O6
2	F	1	NAG	C4-C5-C6-O6
2	E	1	NAG	C4-C5-C6-O6
2	E	2	NAG	C4-C5-C6-O6
2	E	1	NAG	O5-C5-C6-O6
2	E	2	NAG	O5-C5-C6-O6
2	D	2	NAG	C4-C5-C6-O6

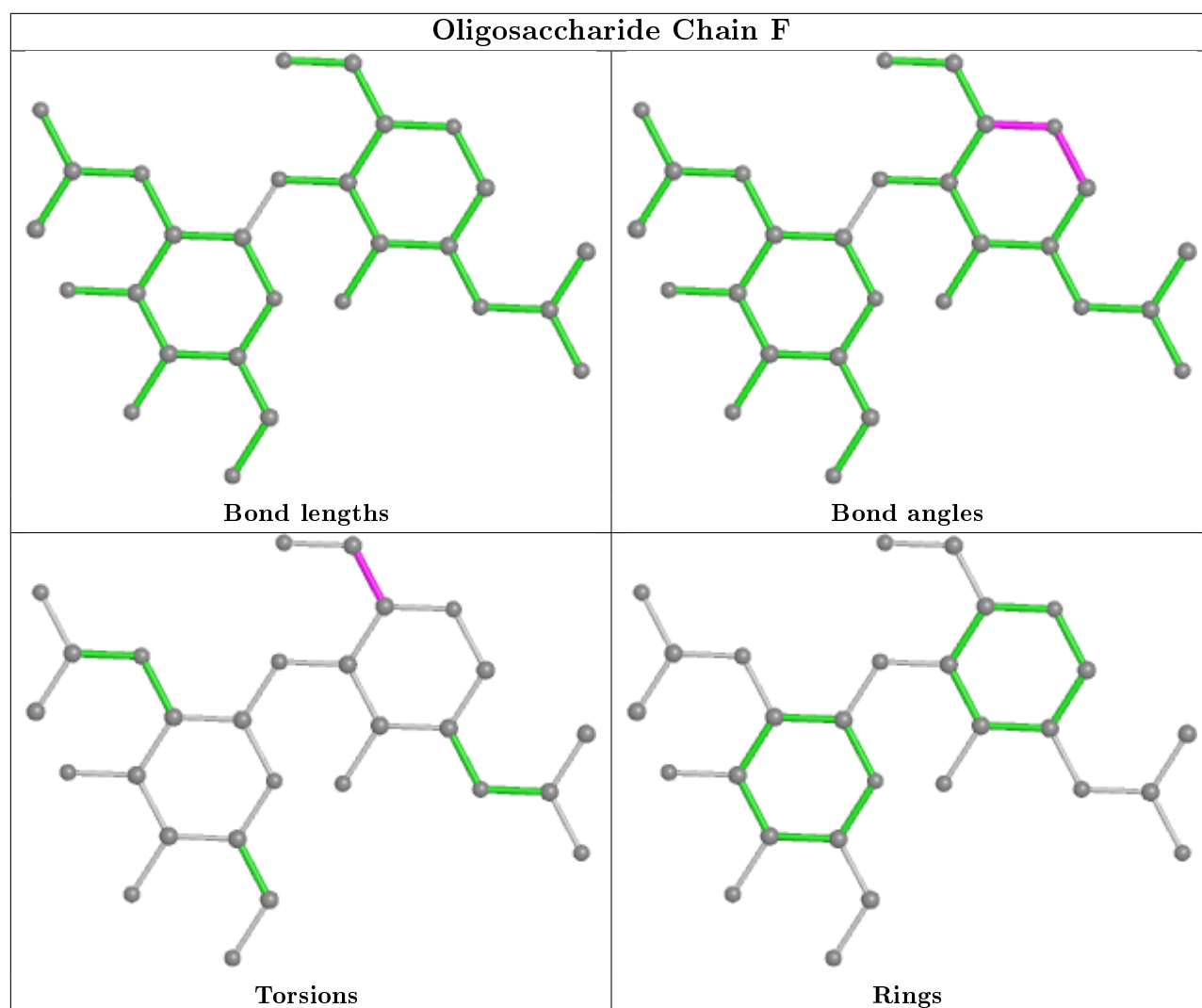
There are no ring outliers.

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 39 ligands modelled in this entry, 39 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	352/361 (97%)	0.28	7 (1%) 65 60	56, 82, 124, 170	0
1	B	351/361 (97%)	0.90	64 (18%) 1 0	74, 122, 165, 230	0
1	C	352/361 (97%)	0.73	43 (12%) 4 2	69, 112, 154, 170	0
All	All	1055/1083 (97%)	0.63	114 (10%) 5 3	56, 107, 159, 230	0

All (114) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	348	GLU	6.5
1	B	149	ILE	6.2
1	B	347	LEU	5.6
1	B	314	VAL	5.6
1	B	351	ILE	5.3
1	B	18	VAL	5.2
1	B	346	VAL	5.1
1	B	354	GLU	5.0
1	C	18	VAL	5.0
1	C	351	ILE	4.9
1	B	75	VAL	4.8
1	C	89	LEU	4.7
1	B	312	ILE	4.5
1	B	20	ALA	4.5
1	B	74	MET	4.4
1	B	352	ASN	4.4
1	B	323	VAL	4.3
1	C	103	VAL	4.2
1	A	351	ILE	4.0
1	C	36	LEU	4.0
1	B	353	THR	3.9
1	B	325	MET	3.8
1	B	298	TYR	3.8

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Mol	Chain	Res	Type	RSRZ
1	C	39	GLU	3.8
1	A	352	ASN	3.8
1	B	251	ILE	3.8
1	C	163	TYR	3.7
1	B	339	LEU	3.7
1	C	84	ALA	3.6
1	B	190	LEU	3.6
1	C	41	PRO	3.5
1	B	83	VAL	3.4
1	B	51	PHE	3.2
1	B	303	PHE	3.2
1	C	75	VAL	3.2
1	A	4	LEU	3.2
1	B	72	PHE	3.1
1	C	38	HIS	3.1
1	B	284	ILE	3.1
1	C	4	LEU	3.1
1	C	65	LEU	3.1
1	C	98	PHE	3.1
1	B	208	ILE	3.1
1	C	35	LEU	3.0
1	C	95	VAL	3.0
1	C	26	VAL	3.0
1	B	315	VAL	3.0
1	B	196	ASP	2.9
1	B	192	LEU	2.9
1	B	153	LYS	2.9
1	B	253	ALA	2.9
1	C	110	PHE	2.8
1	B	313	GLY	2.8
1	B	78	GLU	2.8
1	B	136	LEU	2.8
1	B	290	ASP	2.8
1	C	86	LYS	2.7
1	B	163	TYR	2.7
1	B	87	LEU	2.7
1	B	342	TRP	2.7
1	B	16	VAL	2.6
1	B	170	PHE	2.6
1	C	53	MET	2.6
1	C	74	MET	2.6
1	B	305	ILE	2.6

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Mol	Chain	Res	Type	RSRZ
1	C	269	LEU	2.5
1	C	52	GLU	2.5
1	C	82	ALA	2.5
1	B	247	ASP	2.5
1	B	33	LEU	2.5
1	B	129	LEU	2.5
1	C	27	PHE	2.5
1	B	344	GLU	2.4
1	B	333	LEU	2.4
1	B	299	TRP	2.4
1	C	354	GLU	2.4
1	C	326	GLU	2.4
1	C	352	ASN	2.4
1	B	36	LEU	2.4
1	A	298	TYR	2.4
1	B	115	LEU	2.4
1	C	87	LEU	2.3
1	C	195	ILE	2.3
1	B	191	LYS	2.3
1	B	195	ILE	2.3
1	B	345	ASP	2.3
1	C	115	LEU	2.3
1	B	230	LEU	2.3
1	C	347	LEU	2.3
1	C	298	TYR	2.3
1	B	104	ILE	2.3
1	B	110	PHE	2.2
1	A	35	LEU	2.2
1	B	58	LEU	2.2
1	B	89	LEU	2.2
1	C	159	HIS	2.2
1	C	148	LEU	2.2
1	C	216	GLU	2.2
1	C	314	VAL	2.2
1	C	76	ASP	2.2
1	B	47	SER	2.2
1	B	326	GLU	2.2
1	A	110	PHE	2.2
1	C	56	LEU	2.1
1	B	236	GLU	2.1
1	B	212	PRO	2.1
1	C	3	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	160	TYR	2.1
1	B	295	LEU	2.0
1	B	52	GLU	2.0
1	C	119	LEU	2.0
1	A	355	ASP	2.0
1	B	53	MET	2.0
1	C	127	VAL	2.0

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

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

## 6.3 Carbohydrates [i](#)

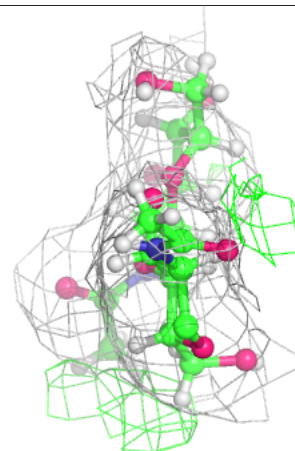
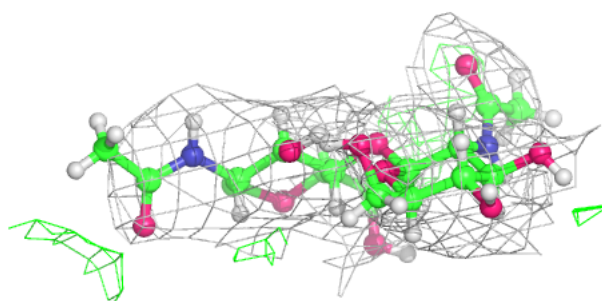
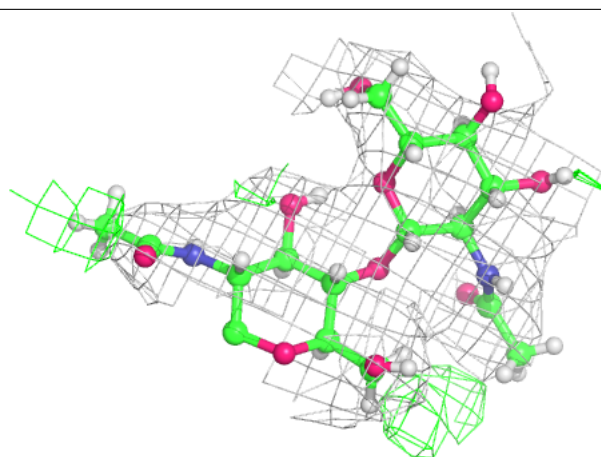
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	NAG	F	2	14/15	0.80	0.26	113,153,186,191	0
2	NAG	E	1	14/15	0.82	0.22	137,167,201,219	0
2	NAG	D	2	14/15	0.82	0.23	105,136,166,194	0
2	NAG	E	2	14/15	0.84	0.34	157,189,220,227	0
2	NAG	F	1	14/15	0.87	0.20	105,138,172,183	0
2	NAG	D	1	14/15	0.91	0.13	95,117,141,169	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.

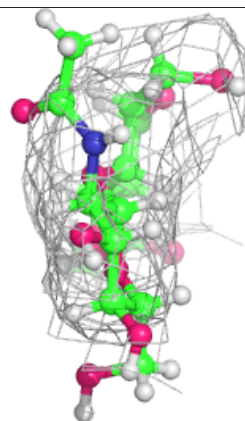
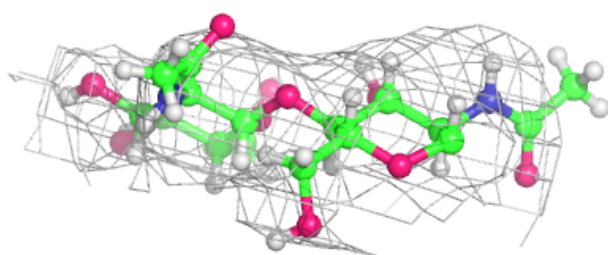
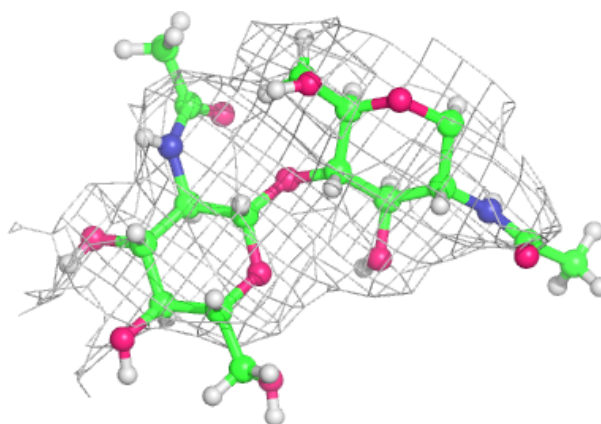
**Electron density around Chain D:**

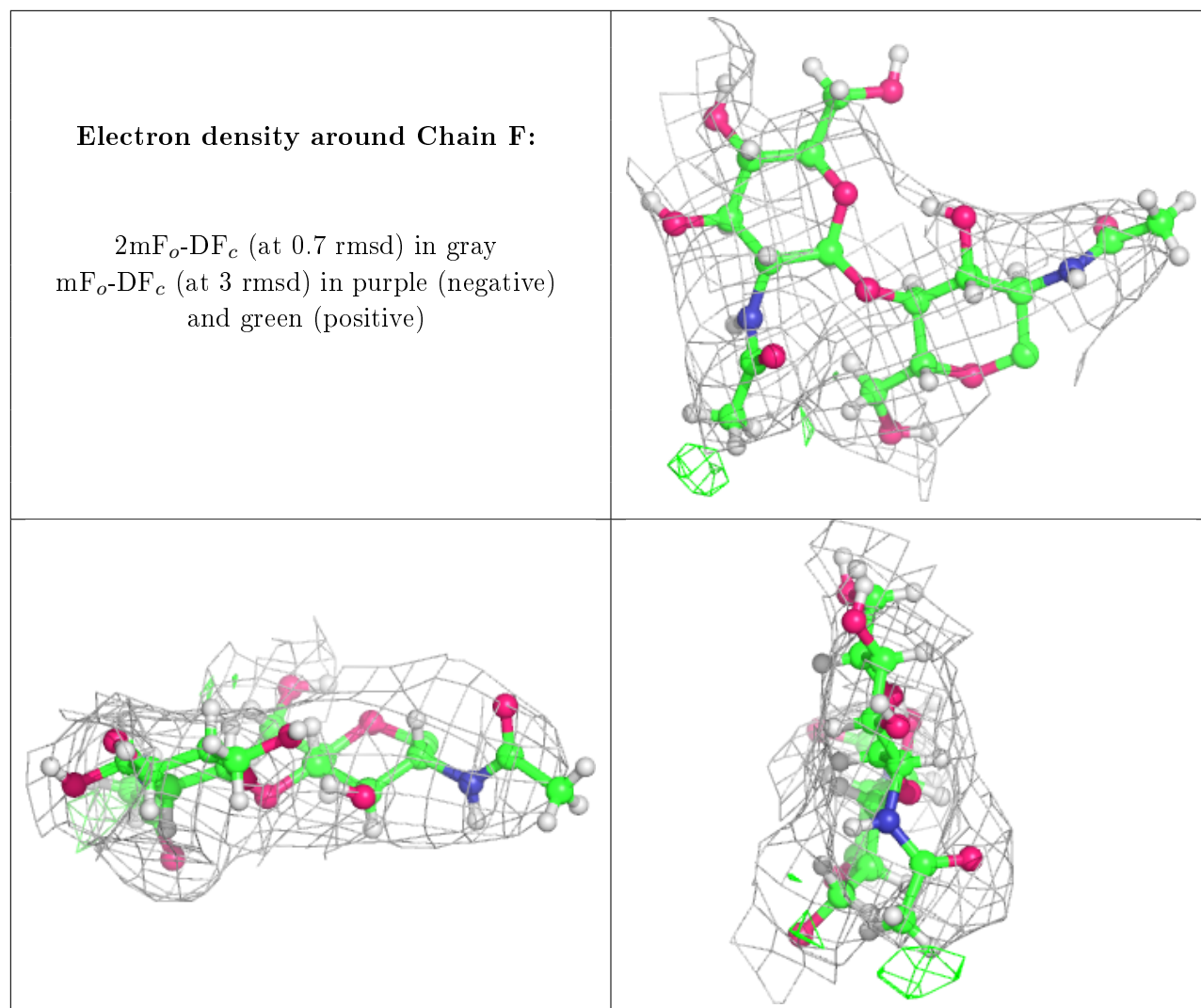
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



**Electron density around Chain E:**

$2mF_o-DF_c$  (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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
3	CA	C	410	1/1	0.31	0.61	157,157,157,157	0
3	CA	A	408	1/1	0.54	0.59	128,128,128,128	0
3	CA	A	410	1/1	0.60	0.52	147,147,147,147	0
3	CA	B	416	1/1	0.62	0.24	150,150,150,150	0
3	CA	A	407	1/1	0.66	1.21	141,141,141,141	0
3	CA	A	409	1/1	0.69	0.65	201,201,201,201	0
3	CA	C	414	1/1	0.74	0.22	177,177,177,177	0
3	CA	B	413	1/1	0.74	0.15	140,140,140,140	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	CA	A	414	1/1	0.76	0.16	127,127,127,127	0
3	CA	A	406	1/1	0.77	0.63	133,133,133,133	0
3	CA	A	415	1/1	0.81	0.17	150,150,150,150	0
3	CA	B	412	1/1	0.81	0.46	120,120,120,120	0
3	CA	B	403	1/1	0.82	1.33	189,189,189,189	0
3	CA	B	409	1/1	0.84	0.17	105,105,105,105	0
3	CA	C	404	1/1	0.87	0.17	104,104,104,104	0
3	CA	B	411	1/1	0.88	0.70	138,138,138,138	0
3	CA	B	405	1/1	0.90	0.27	111,111,111,111	0
3	CA	B	407	1/1	0.91	0.10	134,134,134,134	0
3	CA	C	406	1/1	0.91	1.17	149,149,149,149	0
3	CA	C	403	1/1	0.92	0.14	147,147,147,147	0
3	CA	A	412	1/1	0.92	0.33	89,89,89,89	0
3	CA	C	412	1/1	0.92	0.27	81,81,81,81	0
3	CA	B	404	1/1	0.93	0.58	140,140,140,140	0
3	CA	C	408	1/1	0.93	0.24	119,119,119,119	0
3	CA	B	410	1/1	0.94	0.32	115,115,115,115	0
3	CA	C	413	1/1	0.94	0.12	132,132,132,132	0
3	CA	C	409	1/1	0.94	0.08	147,147,147,147	0
3	CA	A	411	1/1	0.94	0.17	109,109,109,109	0
3	CA	B	406	1/1	0.94	0.24	121,121,121,121	0
3	CA	B	408	1/1	0.95	0.14	129,129,129,129	0
3	CA	A	413	1/1	0.95	0.30	89,89,89,89	0
3	CA	A	404	1/1	0.96	0.27	89,89,89,89	0
3	CA	B	415	1/1	0.96	0.07	149,149,149,149	0
3	CA	C	411	1/1	0.97	0.27	92,92,92,92	0
3	CA	C	407	1/1	0.97	0.50	138,138,138,138	0
3	CA	B	414	1/1	0.97	0.06	140,140,140,140	0
3	CA	A	403	1/1	0.97	0.33	165,165,165,165	0
3	CA	C	405	1/1	0.98	0.26	101,101,101,101	0
3	CA	A	405	1/1	0.99	0.44	107,107,107,107	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.