



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 6, 2020 – 12:10 PM JST

PDB ID : 6LGG  
Title : Bombyx mori GH13 sucrose hydrolase mutant E322Q complexed with sucrose  
Authors : Miyazaki, T.  
Deposited on : 2019-12-05  
Resolution : 1.84 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 : 2.14.6  
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) : 2.14.6

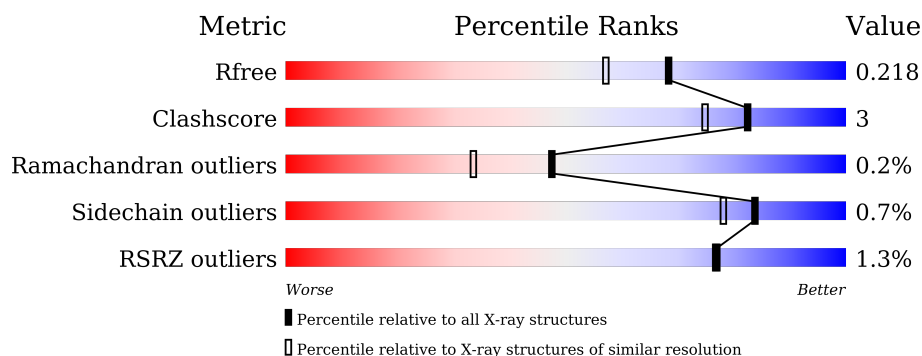
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.84 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	4003 (1.86-1.82)
Clashscore	141614	4233 (1.86-1.82)
Ramachandran outliers	138981	4185 (1.86-1.82)
Sidechain outliers	138945	4186 (1.86-1.82)
RSRZ outliers	127900	3957 (1.86-1.82)

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	598	<div> <div>0.1%</div> <div>89% 5% 6%</div> </div>
1	B	598	<div> <div>2%</div> <div>87% 8% 5%</div> </div>
2	E	2	<div> <div>100%</div> </div>
2	F	2	<div> <div>100%</div> </div>

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 10069 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 Sucrose hydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	564	Total	C	N	O	S	0	0	0
			4607	2953	764	876	14			
1	B	569	Total	C	N	O	S	0	5	0
			4670	2995	770	890	15			

There are 44 discrepancies between the modelled and reference sequences:

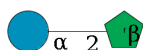
Chain	Residue	Modelled	Actual	Comment	Reference
A	9	MET	-	expression tag	UNP A0A077JI83
A	10	GLY	-	expression tag	UNP A0A077JI83
A	11	SER	-	expression tag	UNP A0A077JI83
A	12	SER	-	expression tag	UNP A0A077JI83
A	13	HIS	-	expression tag	UNP A0A077JI83
A	14	HIS	-	expression tag	UNP A0A077JI83
A	15	HIS	-	expression tag	UNP A0A077JI83
A	16	HIS	-	expression tag	UNP A0A077JI83
A	17	HIS	-	expression tag	UNP A0A077JI83
A	18	HIS	-	expression tag	UNP A0A077JI83
A	19	SER	-	expression tag	UNP A0A077JI83
A	20	SER	-	expression tag	UNP A0A077JI83
A	21	GLY	-	expression tag	UNP A0A077JI83
A	22	LEU	-	expression tag	UNP A0A077JI83
A	23	VAL	-	expression tag	UNP A0A077JI83
A	24	PRO	-	expression tag	UNP A0A077JI83
A	25	ARG	-	expression tag	UNP A0A077JI83
A	26	GLY	-	expression tag	UNP A0A077JI83
A	27	SER	-	expression tag	UNP A0A077JI83
A	28	HIS	-	expression tag	UNP A0A077JI83
A	29	MET	-	expression tag	UNP A0A077JI83
A	322	GLN	GLU	engineered mutation	UNP A0A077JI83
B	9	MET	-	expression tag	UNP A0A077JI83
B	10	GLY	-	expression tag	UNP A0A077JI83
B	11	SER	-	expression tag	UNP A0A077JI83

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Chain	Residue	Modelled	Actual	Comment	Reference
B	12	SER	-	expression tag	UNP A0A077JI83
B	13	HIS	-	expression tag	UNP A0A077JI83
B	14	HIS	-	expression tag	UNP A0A077JI83
B	15	HIS	-	expression tag	UNP A0A077JI83
B	16	HIS	-	expression tag	UNP A0A077JI83
B	17	HIS	-	expression tag	UNP A0A077JI83
B	18	HIS	-	expression tag	UNP A0A077JI83
B	19	SER	-	expression tag	UNP A0A077JI83
B	20	SER	-	expression tag	UNP A0A077JI83
B	21	GLY	-	expression tag	UNP A0A077JI83
B	22	LEU	-	expression tag	UNP A0A077JI83
B	23	VAL	-	expression tag	UNP A0A077JI83
B	24	PRO	-	expression tag	UNP A0A077JI83
B	25	ARG	-	expression tag	UNP A0A077JI83
B	26	GLY	-	expression tag	UNP A0A077JI83
B	27	SER	-	expression tag	UNP A0A077JI83
B	28	HIS	-	expression tag	UNP A0A077JI83
B	29	MET	-	expression tag	UNP A0A077JI83
B	322	GLN	GLU	engineered mutation	UNP A0A077JI83

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



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	E	2	Total	C	O	0	0	0
			23	12	11			
2	F	2	Total	C	O	0	0	0
			23	12	11			

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg) (labeled as "Ligand of Interest" by author).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	2	Total	Mg	0	0
			2	2		
3	A	2	Total	Mg	0	0
			2	2		

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca) (labeled as "Ligand of

Interest" by author).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total 1	Ca 1	0	0
4	A	1	Total 1	Ca 1	0	0

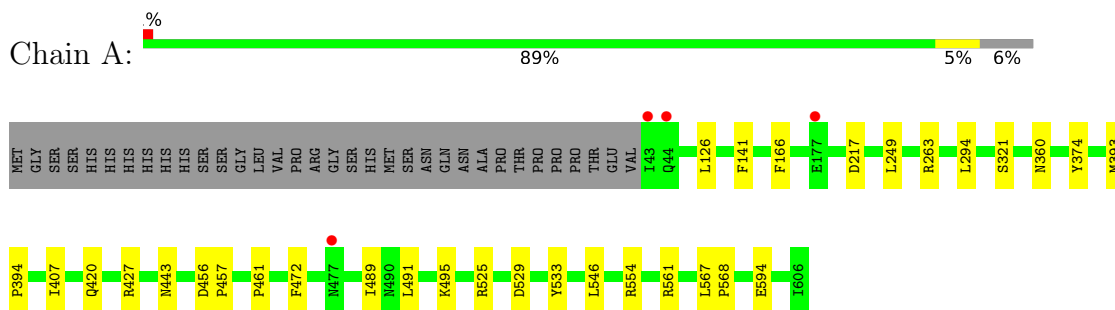
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	314	Total 314	O 314	0	0
5	B	426	Total 426	O 426	0	0

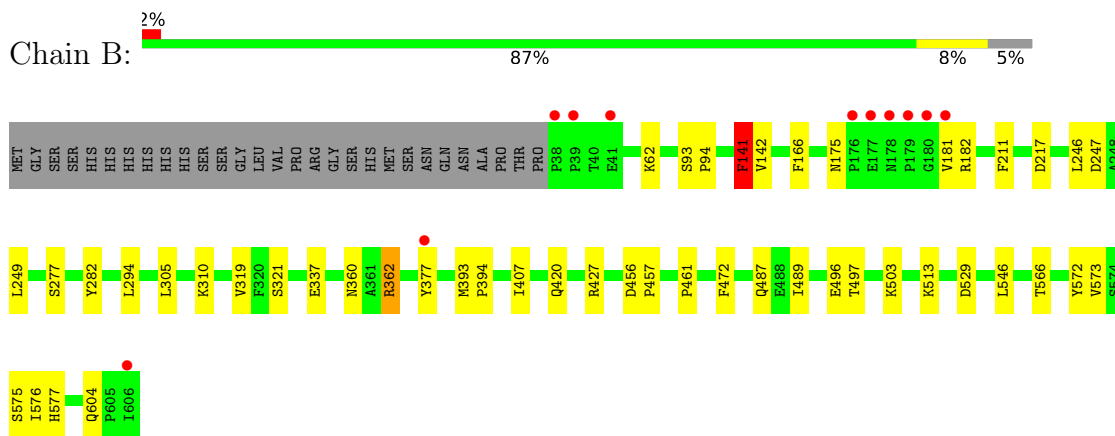
### 3 Residue-property plots

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.

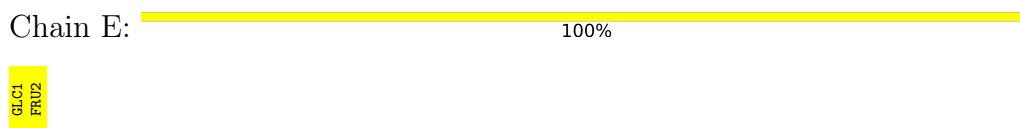
- Molecule 1: Sucrose hydrolase



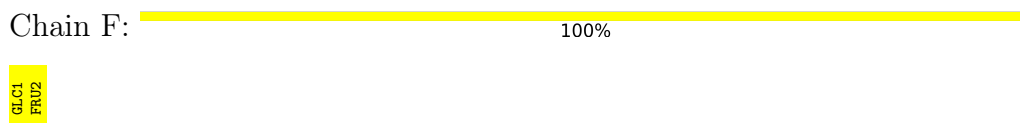
- Molecule 1: Sucrose hydrolase



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



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



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	64.70Å 127.90Å 154.30Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.28 – 1.84 49.23 – 1.84	Depositor EDS
% Data completeness (in resolution range)	99.8 (49.28-1.84) 99.9 (49.23-1.84)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.13 (at 1.84Å)	Xtriage
Refinement program	REFMAC 5.8.0257	Depositor
R, $R_{free}$	0.181 , 0.211 0.189 , 0.218	Depositor DCC
$R_{free}$ test set	5621 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	23.5	Xtriage
Anisotropy	0.095	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 36.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	10069	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.15% 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

### 5.1 Standard geometry

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

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	0.71	0/4748	0.81	0/6463
1	B	0.74	0/4825	0.85	5/6570 (0.1%)
All	All	0.73	0/9573	0.83	5/13033 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	362	ARG	NE-CZ-NH2	-7.07	116.76	120.30
1	B	141	PHE	CB-CA-C	7.02	124.44	110.40
1	B	362	ARG	NE-CZ-NH1	6.18	123.39	120.30
1	B	211	PHE	CB-CA-C	5.16	120.73	110.40
1	B	337	GLU	CB-CA-C	5.12	120.63	110.40

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4607	0	4329	18	0
1	B	4670	0	4398	30	0
2	E	23	0	21	0	0
2	F	23	0	21	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	2	0	0	0	0
3	B	2	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	314	0	0	3	0
5	B	426	0	0	7	0
All	All	10069	0	8769	46	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 3.

All (46) 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:393:MET:HB3	1:A:394:PRO:HD3	1.78	0.65
1:B:141:PHE:CD2	1:B:246:LEU:CD2	2.85	0.60
1:B:141:PHE:CD2	1:B:246:LEU:HD21	2.39	0.58
1:B:393:MET:HB3	1:B:394:PRO:HD3	1.86	0.58
1:B:142:VAL:HA	1:B:247:ASP:HB3	1.84	0.57
1:B:141:PHE:HD2	1:B:246:LEU:HD21	1.70	0.57
1:B:249:LEU:HD12	1:B:321:SER:HB2	1.88	0.56
1:A:427:ARG:CZ	1:A:489:ILE:HD11	2.37	0.54
1:A:461:PRO:HB3	1:A:472:PHE:CG	2.43	0.54
1:A:554:ARG:HD2	1:A:594:GLU:OE2	2.08	0.54
1:B:305:LEU:HD22	1:B:319[B]:VAL:HG23	1.90	0.53
1:B:497:THR:O	1:B:503:LYS:HE2	2.08	0.52
1:A:141:PHE:HA	5:A:1032:HOH:O	2.09	0.51
1:A:407:ILE:HG21	1:A:546:LEU:HD13	1.93	0.51
1:B:572:TYR:CD2	1:B:573:VAL:HG23	2.46	0.51
1:B:513:LYS:HG3	5:B:828:HOH:O	2.11	0.51
1:B:575:SER:HB3	5:B:1099:HOH:O	2.12	0.50
1:A:561:ARG:NH2	1:B:377:TYR:CD1	2.80	0.49
1:A:294:LEU:C	1:A:294:LEU:HD12	2.33	0.49
1:B:487:GLN:CG	5:B:835:HOH:O	2.61	0.48
1:B:360:ASN:HB2	1:B:529:ASP:O	2.13	0.48
1:B:141:PHE:CD2	1:B:246:LEU:HD22	2.50	0.47
1:B:62:LYS:HE2	5:B:945:HOH:O	2.13	0.46
1:B:277:SER:HA	1:B:282:TYR:CD2	2.50	0.46
1:A:491:LEU:HG	1:A:495:LYS:HD2	1.98	0.45
1:A:525:ARG:HB3	1:A:533:TYR:CZ	2.52	0.45
1:A:360:ASN:HB2	1:A:529:ASP:O	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:407:ILE:HG21	1:B:546:LEU:HD13	1.99	0.45
1:B:427:ARG:CZ	1:B:489:ILE:HD11	2.46	0.45
1:B:566:THR:OG1	1:B:604:GLN:HG2	2.16	0.44
1:B:456:ASP:N	1:B:457:PRO:CD	2.81	0.44
1:A:456:ASP:N	1:A:457:PRO:CD	2.80	0.44
1:A:249:LEU:HD12	1:A:321:SER:HB2	1.99	0.43
1:A:561:ARG:HD2	5:A:890:HOH:O	2.18	0.43
1:B:496:GLU:HG3	5:B:961:HOH:O	2.18	0.43
1:A:567:LEU:HA	1:A:568:PRO:C	2.39	0.42
1:B:461:PRO:HB3	1:B:472:PHE:CG	2.53	0.42
1:B:294:LEU:C	1:B:294:LEU:HD12	2.39	0.42
1:B:503:LYS:NZ	5:B:820:HOH:O	2.52	0.42
1:B:182:ARG:HD3	5:B:858:HOH:O	2.19	0.42
1:A:443:ASN:ND2	5:A:815:HOH:O	2.48	0.42
1:A:374:TYR:HA	1:B:362:ARG:HD3	2.03	0.41
1:B:175:ASN:ND2	1:B:181:VAL:HG13	2.35	0.41
1:A:126:LEU:C	1:A:126:LEU:HD23	2.41	0.40
1:B:576:ILE:HG13	1:B:577:HIS:CD2	2.57	0.40
1:B:93:SER:HB3	1:B:94:PRO:CD	2.51	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	562/598 (94%)	549 (98%)	12 (2%)	1 (0%)	47 33
1	B	572/598 (96%)	557 (97%)	14 (2%)	1 (0%)	47 33
All	All	1134/1196 (95%)	1106 (98%)	26 (2%)	2 (0%)	47 33

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	420	GLN
1	B	420	GLN

### 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	494/524 (94%)	491 (99%)	3 (1%)	86	82
1	B	504/524 (96%)	500 (99%)	4 (1%)	81	75
All	All	998/1048 (95%)	991 (99%)	7 (1%)	84	78

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	166	PHE
1	A	217	ASP
1	A	263	ARG
1	B	141	PHE
1	B	166	PHE
1	B	217	ASP
1	B	310	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	191	GLN
1	A	391	ASN
1	A	577	HIS
1	B	391	ASN
1	B	577	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 ⓘ

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$
2	GLC	E	1	2	11,11,12	0.63	0	15,15,17	1.76	3 (20%)
2	FRU	E	2	2	11,12,12	1.12	1 (9%)	10,18,18	1.18	1 (10%)
2	GLC	F	1	2	11,11,12	0.76	0	15,15,17	2.12	3 (20%)
2	FRU	F	2	2	11,12,12	1.60	3 (27%)	10,18,18	1.48	1 (10%)

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	GLC	E	1	2	-	0/2/19/22	0/1/1/1
2	FRU	E	2	2	-	0/5/24/24	0/1/1/1
2	GLC	F	1	2	-	0/2/19/22	0/1/1/1
2	FRU	F	2	2	-	0/5/24/24	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	2	FRU	O2-C2	3.13	1.46	1.40
2	F	2	FRU	C1-C2	2.49	1.56	1.52
2	E	2	FRU	O2-C2	2.48	1.45	1.40
2	F	2	FRU	O5-C2	2.07	1.46	1.43

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	1	GLC	C1-O5-C5	6.76	121.35	112.19
2	E	1	GLC	C1-O5-C5	4.62	118.45	112.19
2	F	2	FRU	O1-C1-C2	3.05	118.36	111.86
2	E	1	GLC	C2-C3-C4	-2.80	106.05	110.89
2	F	1	GLC	C2-C3-C4	-2.70	106.23	110.89
2	F	1	GLC	C1-C2-C3	-2.30	106.83	109.67
2	E	1	GLC	O5-C5-C4	2.28	116.38	110.83
2	E	2	FRU	O3-C3-C4	-2.10	106.08	113.32

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 6 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 [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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	564/598 (94%)	-0.27	4 (0%) 87 87	17, 29, 47, 75	0
1	B	569/598 (95%)	-0.36	11 (1%) 66 65	15, 23, 41, 82	0
All	All	1133/1196 (94%)	-0.32	15 (1%) 77 77	15, 26, 44, 82	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	38	PRO	5.3
1	A	43	ILE	5.2
1	B	181	VAL	5.2
1	B	180	GLY	4.7
1	B	176	PRO	3.5
1	B	606	ILE	3.2
1	B	179	PRO	3.1
1	B	377	TYR	2.9
1	B	177	GLU	2.8
1	B	178	ASN	2.7
1	B	41	GLU	2.6
1	A	477	ASN	2.5
1	B	39	PRO	2.4
1	A	44	GLN	2.3
1	A	177	GLU	2.2

### 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	FRU	E	2	12/12	0.94	0.09	23,24,26,26	0
2	FRU	F	2	12/12	0.96	0.09	21,23,26,27	0
2	GLC	F	1	11/12	0.97	0.11	16,18,21,23	0
2	GLC	E	1	11/12	0.98	0.09	21,22,23,24	0

### 6.4 Ligands [i](#)

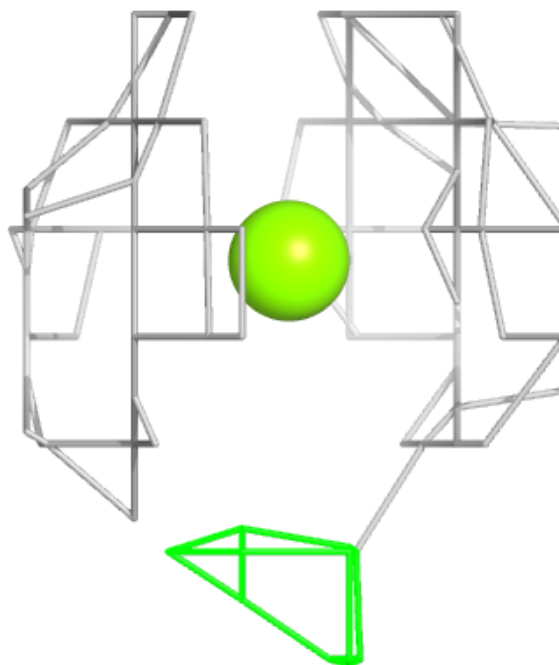
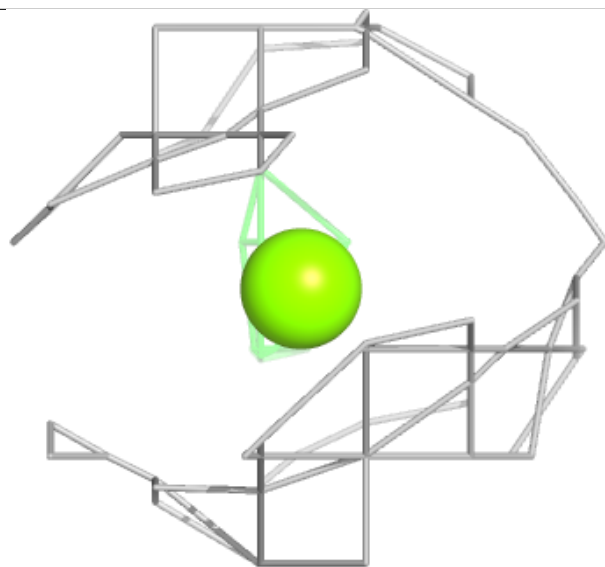
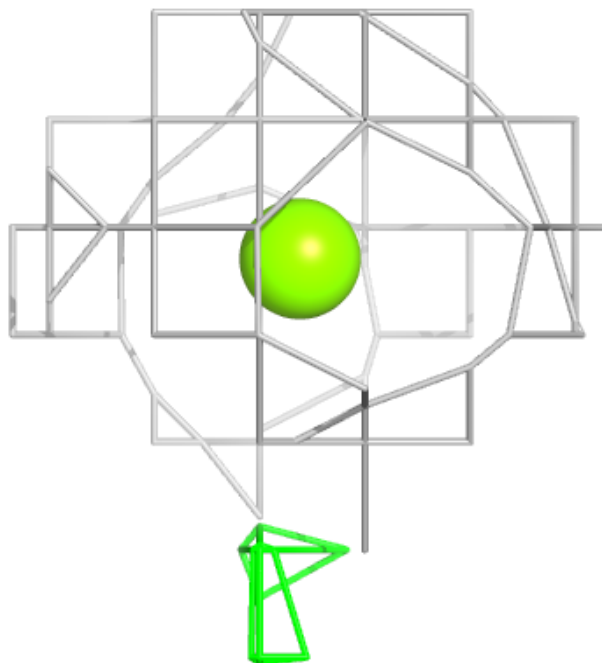
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<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	MG	A	701	1/1	0.95	0.07	32,32,32,32	0
3	MG	A	703	1/1	0.96	0.13	35,35,35,35	0
3	MG	B	703	1/1	0.99	0.10	32,32,32,32	0
3	MG	B	701	1/1	0.99	0.04	19,19,19,19	0
4	CA	B	702	1/1	1.00	0.05	18,18,18,18	0
4	CA	A	702	1/1	1.00	0.05	25,25,25,25	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 MG A 701:**

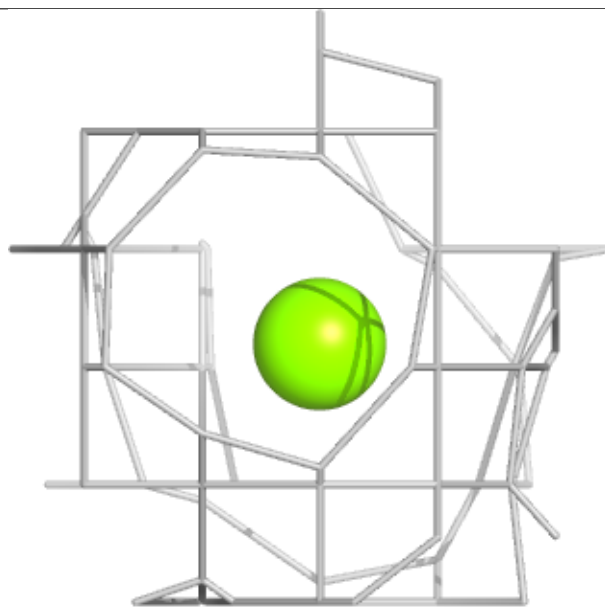
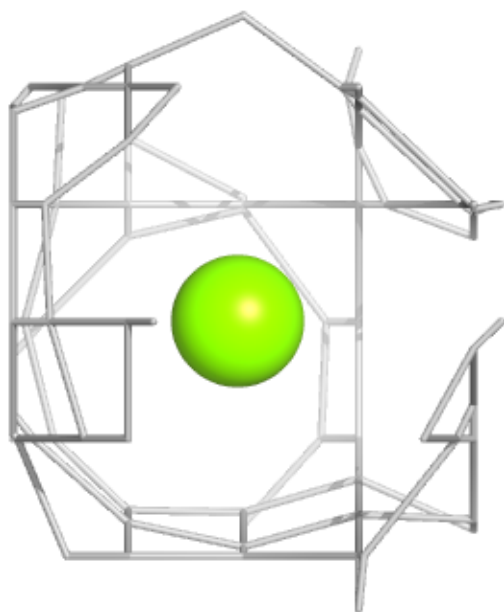
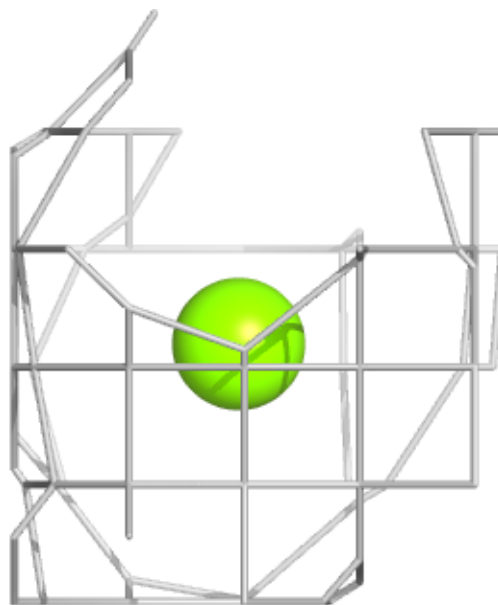
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





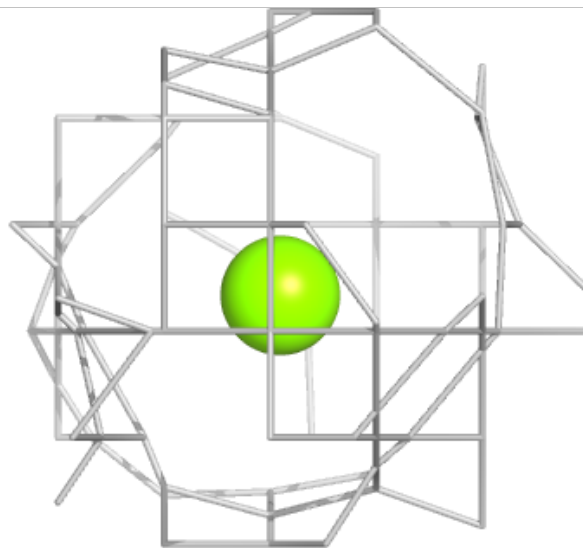
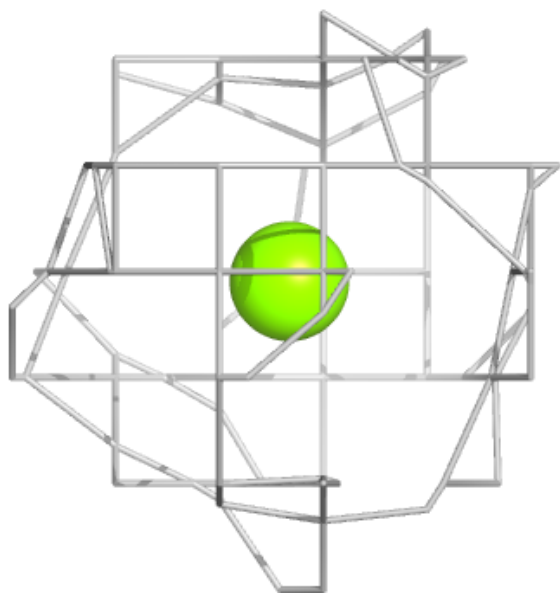
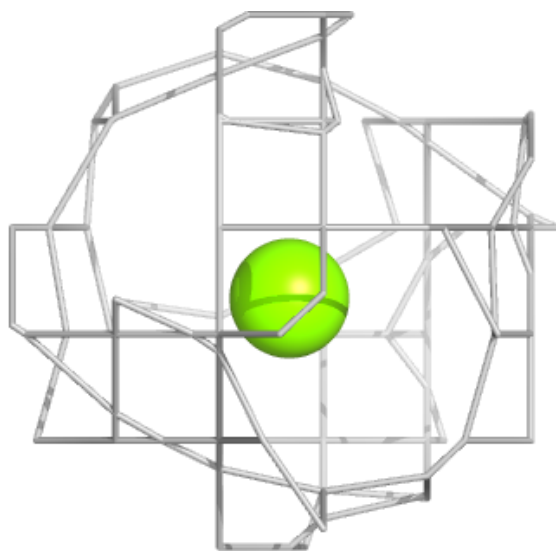
**Electron density around MG A 703:**

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



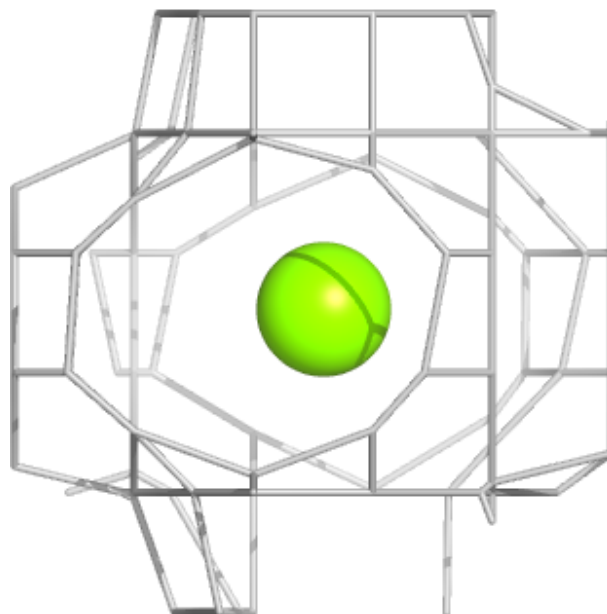
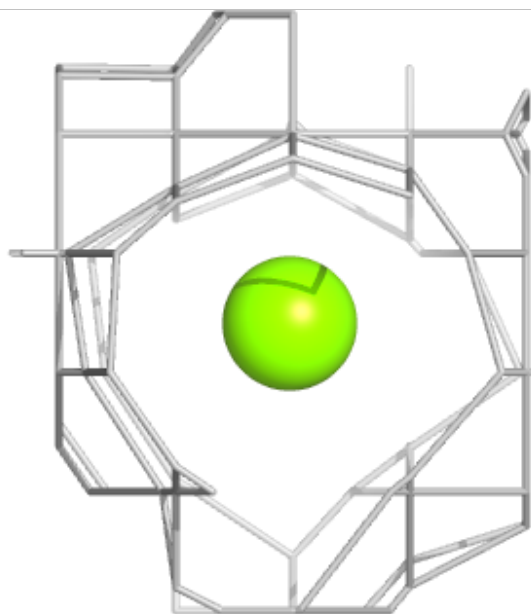
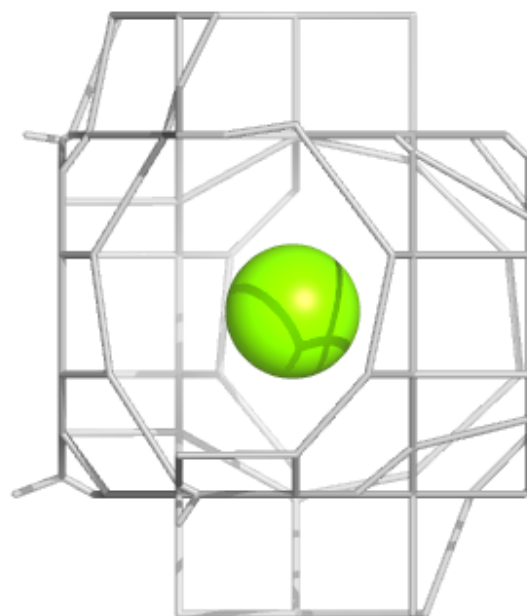
**Electron density around MG B 703:**

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



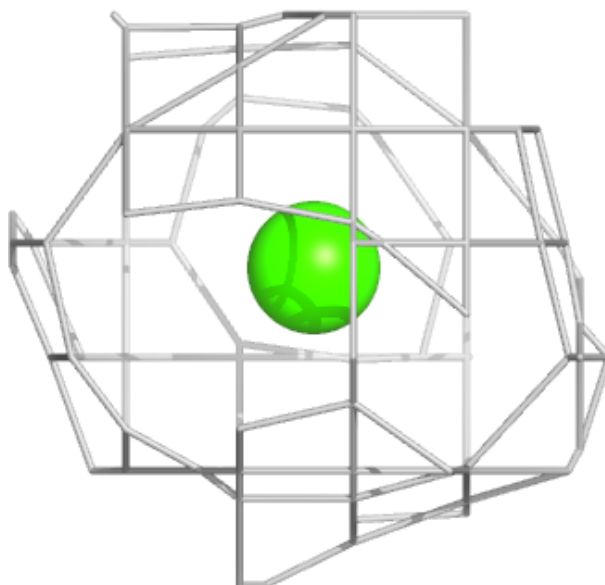
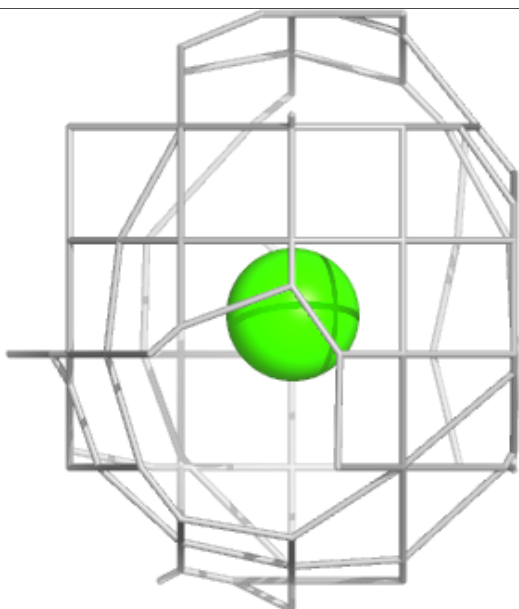
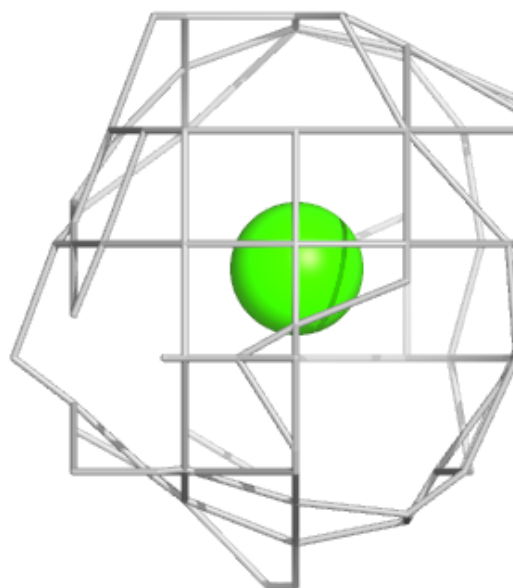
**Electron density around MG B 701:**

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



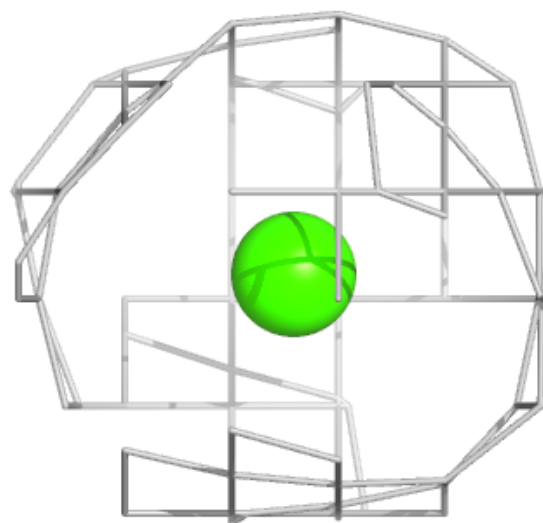
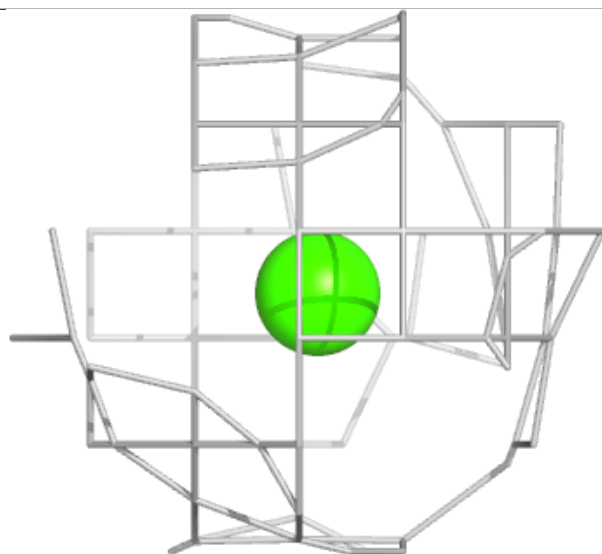
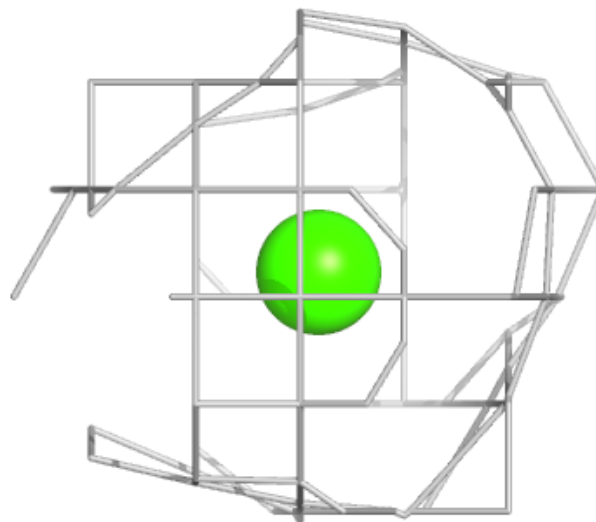
**Electron density around CA B 702:**

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



**Electron density around CA A 702:**

$2mF_o-DF_c$  (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.