



# Full wwPDB X-ray Structure Validation Report ⓘ

Apr 30, 2024 – 02:08 PM JST

PDB ID : 8JA4  
Title : Structure of the alginate epimerase/lyase  
Authors : Fujiwara, T.  
Deposited on : 2023-05-05  
Resolution : 2.04 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)

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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>.

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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 : 2.36.2  
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) : 2.36.2

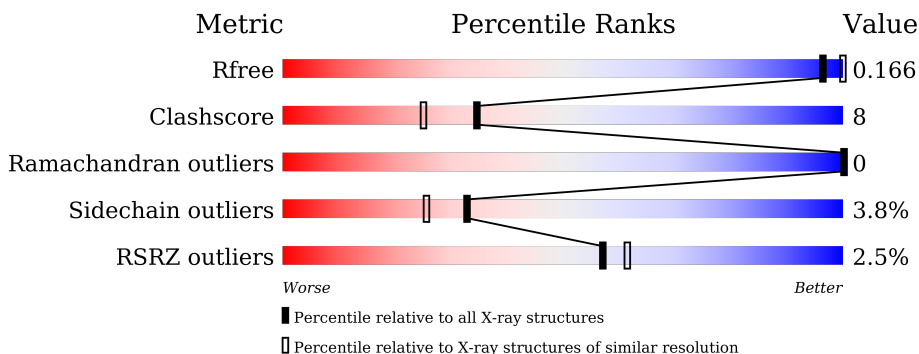
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.04 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 of 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	493	

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	PEG	A	509	-	-	X	-
4	PGE	A	510	-	X	-	X

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 3840 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 mannuronan 5-epimerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	480	3630	2209	633	782	6	0	8	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	486	LEU	-	expression tag	UNP A0A0C4WKK2
A	487	GLU	-	expression tag	UNP A0A0C4WKK2
A	488	HIS	-	expression tag	UNP A0A0C4WKK2
A	489	HIS	-	expression tag	UNP A0A0C4WKK2
A	490	HIS	-	expression tag	UNP A0A0C4WKK2
A	491	HIS	-	expression tag	UNP A0A0C4WKK2
A	492	HIS	-	expression tag	UNP A0A0C4WKK2
A	493	HIS	-	expression tag	UNP A0A0C4WKK2

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca) (labeled as "Ligand of Interest" by depositor).

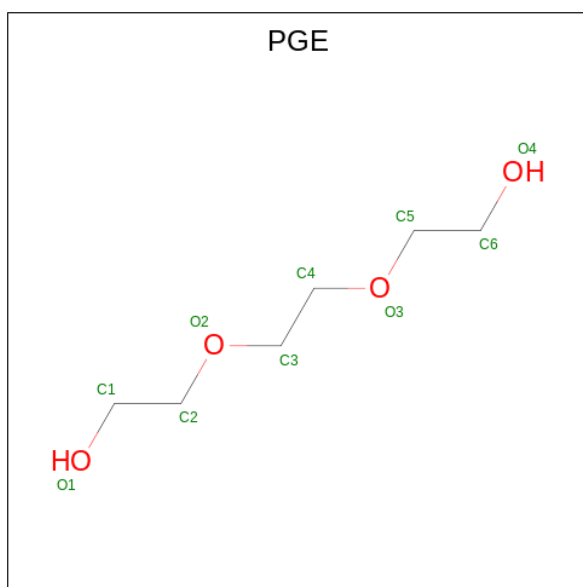
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	6	Total	Ca	0	0
			6	6		

- Molecule 3 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C<sub>4</sub>H<sub>10</sub>O<sub>3</sub>).



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

- Molecule 4 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula:  $C_6H_{14}O_4$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			10	6	4		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
4	A	1	10	6	4	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	A	163	163	163	0	0

### 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: mannuronan 5-epimerase



## 4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	122.52Å 122.52Å 119.80Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.51 – 2.04 48.51 – 2.04	Depositor EDS
% Data completeness (in resolution range)	100.0 (48.51-2.04) 100.0 (48.51-2.04)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.74 (at 2.05Å)	Xtrriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, $R_{free}$	0.167 , 0.172 0.164 , 0.166	Depositor DCC
$R_{free}$ test set	2113 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	48.3	Xtrriage
Anisotropy	0.065	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 47.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.019 for h,-h-k,-l	Xtrriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	3840	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	60.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.36% 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, PGE, PEG

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	0.64	0/3702	0.82	2/5005 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	387	ASP	N-CA-CB	-7.67	96.80	110.60
1	A	387	ASP	N-CA-C	6.34	128.11	111.00

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	3630	0	3380	59	0
2	A	6	0	0	0	0
3	A	21	0	30	9	0
4	A	20	0	22	0	0
5	A	163	0	0	1	0
All	All	3840	0	3432	59	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 8.



All (59) 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:383:ASN:OD1	1:A:385:GLU:HG3	1.53	1.09
1:A:102:ILE:HG13	1:A:141:VAL:HG21	1.36	1.05
1:A:195:ARG:HD3	3:A:507:PEG:H31	1.58	0.85
1:A:66:ASP:OD2	1:A:103:ARG:HD2	1.84	0.78
1:A:382:ARG:HG2	1:A:388:THR:HA	1.67	0.75
1:A:89:ILE:HD12	1:A:107:LEU:HD21	1.72	0.71
1:A:454:ASN:C	1:A:455:LEU:HD12	2.10	0.71
1:A:382:ARG:NH1	1:A:388:THR:HB	2.06	0.70
1:A:382:ARG:CG	1:A:388:THR:HA	2.21	0.70
1:A:308:ASP:H	3:A:508:PEG:H42	1.57	0.69
1:A:263[A]:GLN:HG3	1:A:286:LEU:HB2	1.76	0.68
1:A:455:LEU:HD12	1:A:455:LEU:N	2.12	0.64
1:A:330:VAL:HB	1:A:348:ALA:HB3	1.81	0.61
1:A:108:ASP:OD1	5:A:601:HOH:O	2.16	0.61
1:A:102:ILE:HG13	1:A:141:VAL:CG2	2.23	0.61
1:A:199:ASN:HD21	3:A:509:PEG:H12	1.68	0.59
1:A:428:GLU:N	1:A:428:GLU:OE2	2.36	0.59
1:A:263[A]:GLN:NE2	1:A:287:ASP:OD1	2.39	0.55
1:A:89:ILE:CD1	1:A:107:LEU:HD21	2.35	0.55
1:A:383:ASN:OD1	1:A:385:GLU:CG	2.43	0.54
1:A:308:ASP:H	3:A:508:PEG:C4	2.20	0.54
1:A:195:ARG:HG3	3:A:507:PEG:H12	1.90	0.53
1:A:199:ASN:ND2	3:A:509:PEG:H12	2.22	0.53
1:A:161:THR:HG21	1:A:163:ARG:NH1	2.23	0.53
1:A:467:ASP:OD1	1:A:470:GLN:HG3	2.09	0.53
1:A:140:ARG:HA	1:A:164:ASP:O	2.09	0.53
1:A:105:LEU:HD12	1:A:105:LEU:C	2.30	0.52
1:A:89:ILE:HD12	1:A:107:LEU:CD2	2.38	0.51
1:A:185:PHE:HB2	1:A:209:LEU:HD23	1.93	0.49
1:A:201:VAL:HG11	3:A:509:PEG:H11	1.93	0.49
1:A:308:ASP:N	3:A:508:PEG:H42	2.25	0.49
1:A:215:TYR:HA	1:A:245:HIS:O	2.13	0.49
1:A:382:ARG:CZ	1:A:388:THR:HB	2.42	0.48
1:A:397:PHE:CE2	1:A:404:ILE:HG13	2.49	0.48
1:A:299:LYS:HD2	1:A:478:ASP:HB3	1.96	0.47
1:A:105:LEU:HD12	1:A:105:LEU:O	2.14	0.47
1:A:125:TYR:CE1	1:A:126:ILE:HD12	2.50	0.46
1:A:149:TYR:CE1	1:A:172:LYS:HE2	2.51	0.45
1:A:455:LEU:N	1:A:455:LEU:CD1	2.77	0.45
1:A:148:ARG:HG3	1:A:172:LYS:HD3	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:76:LYS:HG3	1:A:108:ASP:HB3	1.99	0.45
1:A:154:HIS:HA	1:A:178:ASP:HB3	1.99	0.44
1:A:45:ARG:HB3	1:A:78:VAL:HG12	2.00	0.44
1:A:199:ASN:HD21	3:A:509:PEG:C1	2.30	0.44
1:A:58:MET:CE	1:A:92:LYS:HB2	2.48	0.43
1:A:74:VAL:HG22	1:A:106:THR:HB	2.00	0.43
1:A:382:ARG:HH11	1:A:388:THR:HB	1.82	0.43
1:A:72:GLU:OE2	1:A:72:GLU:HA	2.19	0.43
1:A:213:VAL:HA	1:A:243:ALA:O	2.19	0.43
1:A:378:GLN:HE22	1:A:445:ASN:ND2	2.17	0.42
1:A:196:HIS:HD1	1:A:220:ASN:H	1.66	0.42
1:A:201:VAL:HG22	1:A:202:THR:N	2.34	0.42
1:A:268:TYR:HA	1:A:291:HIS:O	2.20	0.42
1:A:389:SER:O	1:A:390:ARG:HB2	2.20	0.42
1:A:429:ASP:OD2	1:A:429:ASP:N	2.46	0.42
1:A:146:VAL:HG22	1:A:170:ASN:OD1	2.20	0.42
1:A:235:TYR:HA	1:A:258:SER:O	2.20	0.42
1:A:77:LEU:HD23	1:A:114:THR:HG21	2.00	0.42
1:A:161:THR:HG22	1:A:163:ARG:HG2	2.04	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	A	486/493 (99%)	461 (95%)	25 (5%)	0	<a href="#">100</a> <a href="#">100</a>

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	A	376/380 (99%)	362 (96%)	14 (4%)	34 27

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	PHE
1	A	6	LYS
1	A	38	ARG
1	A	50	ASP
1	A	60	LYS
1	A	66	ASP
1	A	82	ASP
1	A	130	ASP
1	A	154	HIS
1	A	163	ARG
1	A	172	LYS
1	A	179	PHE
1	A	382	ARG
1	A	389	SER

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

Mol	Chain	Res	Type
1	A	83	GLN
1	A	199	ASN
1	A	225	GLN
1	A	316	ASN
1	A	378	GLN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 11 ligands modelled in this entry, 6 are monoatomic - leaving 5 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
3	PEG	A	508	-	6,6,6	0.10	0	5,5,5	0.10	0
4	PGE	A	511	-	9,9,9	0.45	0	8,8,8	0.75	0
3	PEG	A	507	-	6,6,6	0.12	0	5,5,5	0.07	0
4	PGE	A	510	-	9,9,9	3.27	5 (55%)	8,8,8	4.64	5 (62%)
3	PEG	A	509	-	6,6,6	0.11	0	5,5,5	0.10	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	PEG	A	508	-	-	1/4/4/4	-
4	PGE	A	511	-	-	3/7/7/7	-
3	PEG	A	507	-	-	0/4/4/4	-
4	PGE	A	510	-	-	6/7/7/7	-
3	PEG	A	509	-	-	1/4/4/4	-

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	510	PGE	C4-C3	-4.76	1.24	1.49
4	A	510	PGE	O2-C3	-4.68	1.21	1.42
4	A	510	PGE	O1-C1	-4.53	1.18	1.42
4	A	510	PGE	O2-C2	-4.40	1.22	1.42
4	A	510	PGE	O4-C6	-2.85	1.27	1.42

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	510	PGE	O2-C3-C4	9.59	153.61	110.39
4	A	510	PGE	O3-C5-C6	5.07	132.34	110.07
4	A	510	PGE	C5-O3-C4	4.56	133.03	113.29
4	A	510	PGE	O3-C4-C3	4.29	129.74	110.39
4	A	510	PGE	C3-O2-C2	-2.73	101.47	113.29

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	511	PGE	C1-C2-O2-C3
4	A	510	PGE	C3-C4-O3-C5
4	A	510	PGE	O3-C5-C6-O4
4	A	511	PGE	O3-C5-C6-O4
4	A	510	PGE	C4-C3-O2-C2
4	A	510	PGE	C6-C5-O3-C4
3	A	509	PEG	C4-C3-O2-C2
3	A	508	PEG	C1-C2-O2-C3
4	A	510	PGE	O1-C1-C2-O2
4	A	510	PGE	O2-C3-C4-O3
4	A	511	PGE	O2-C3-C4-O3

There are no ring outliers.

3 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	508	PEG	3	0
3	A	507	PEG	2	0
3	A	509	PEG	4	0

## 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	480/493 (97%)	0.18	12 (2%) 57 61	35, 55, 91, 122	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1	MET	6.2
1	A	386	GLY	6.1
1	A	388	THR	4.1
1	A	10	ALA	3.8
1	A	387	ASP	3.7
1	A	16	SER	2.8
1	A	15	LYS	2.6
1	A	428	GLU	2.4
1	A	13	ASP	2.3
1	A	28	ALA	2.2
1	A	21	ALA	2.2
1	A	71	GLY	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](#)

There are no monosaccharides in this entry.

### 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.

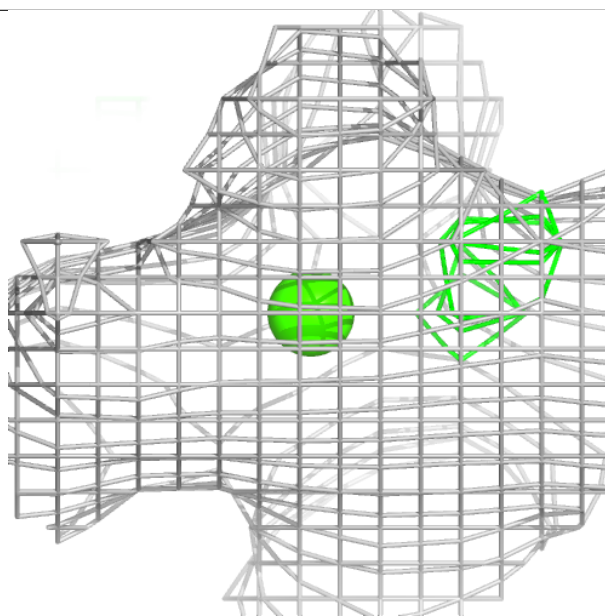
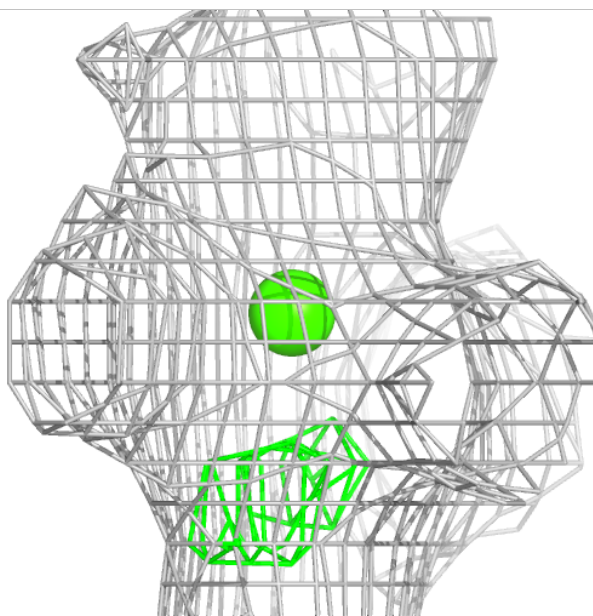
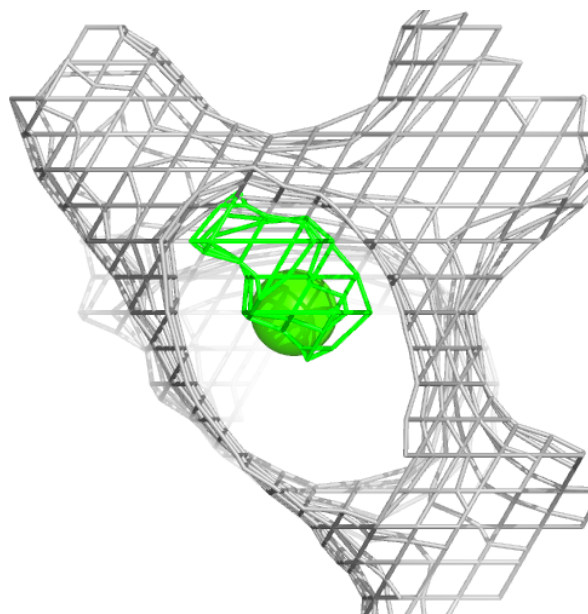
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
4	PGE	A	510	10/10	0.65	0.43	42,51,61,63	10
4	PGE	A	511	10/10	0.68	0.15	84,89,95,98	0
3	PEG	A	507	7/7	0.73	0.33	47,54,60,61	7
3	PEG	A	509	7/7	0.79	0.31	60,72,80,81	0
2	CA	A	506	1/1	0.85	0.09	91,91,91,91	1
3	PEG	A	508	7/7	0.89	0.26	41,51,61,72	7
2	CA	A	501	1/1	0.94	0.09	60,60,60,60	1
2	CA	A	503	1/1	0.99	0.07	41,41,41,41	0
2	CA	A	505	1/1	0.99	0.05	55,55,55,55	0
2	CA	A	504	1/1	1.00	0.07	49,49,49,49	0
2	CA	A	502	1/1	1.00	0.10	42,42,42,42	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.



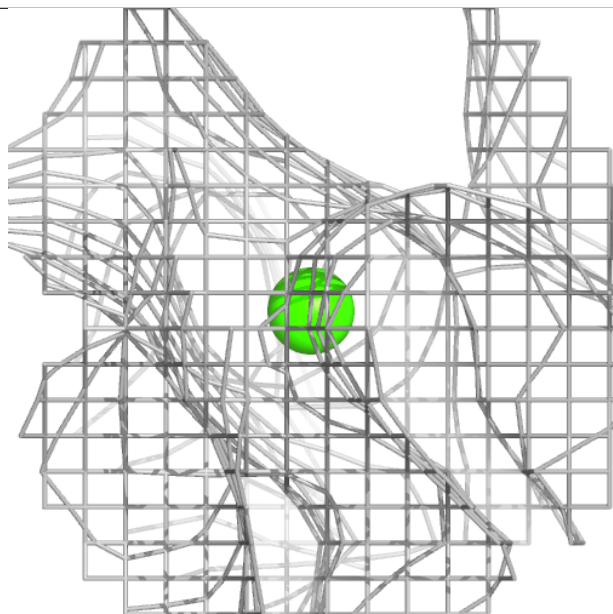
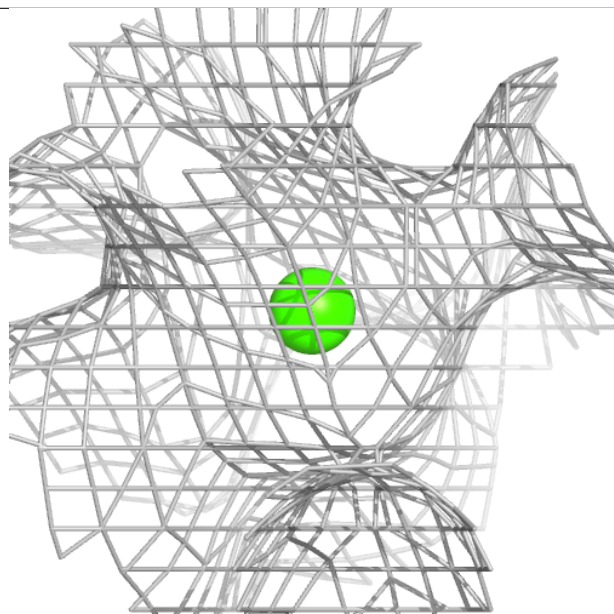
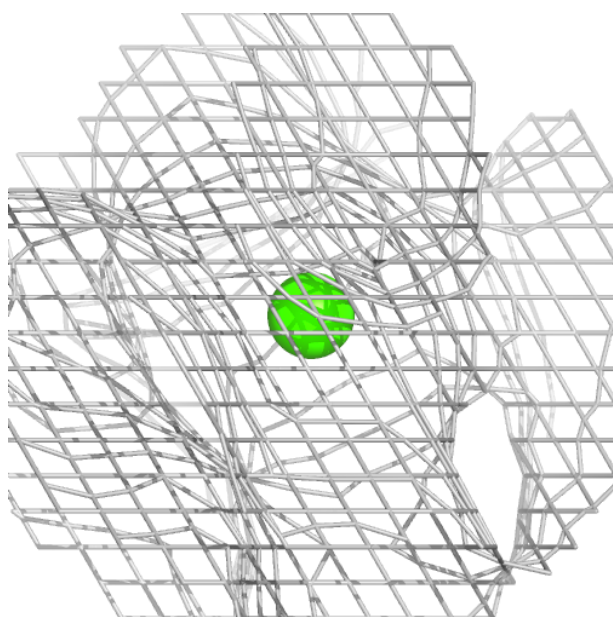
**Electron density around CA A 506:**

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



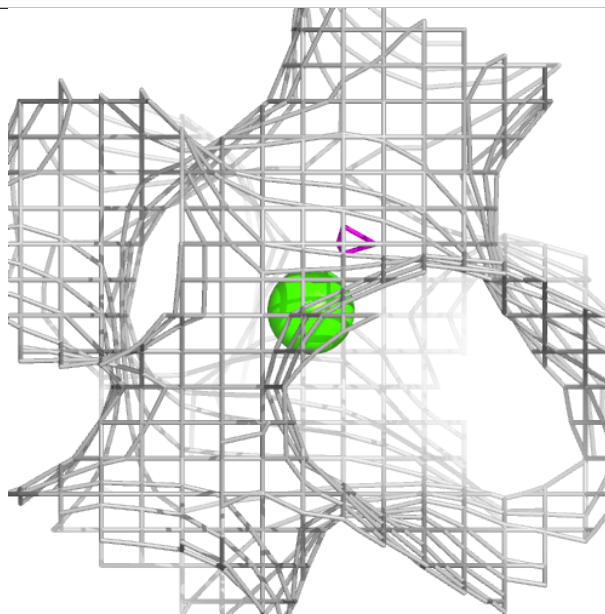
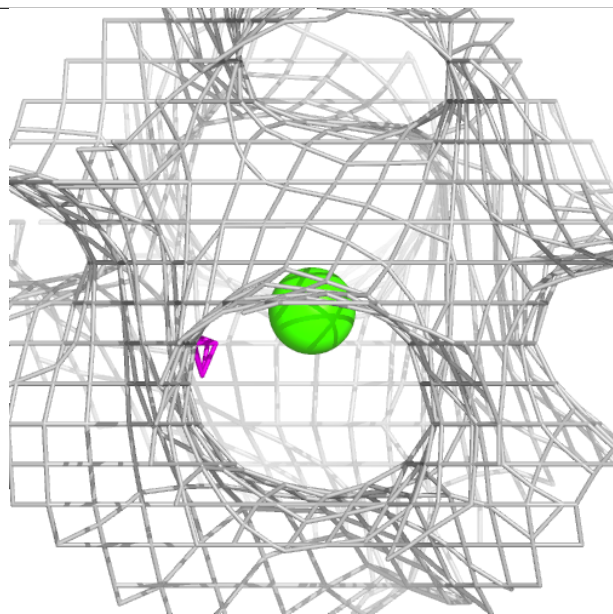
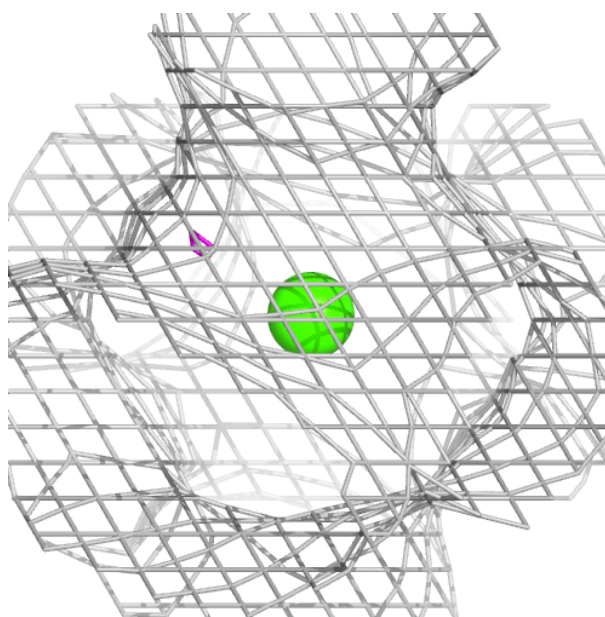
**Electron density around CA A 501:**

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



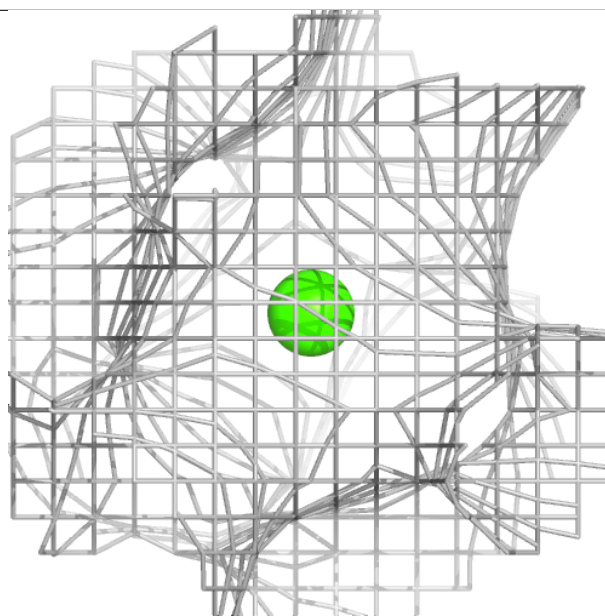
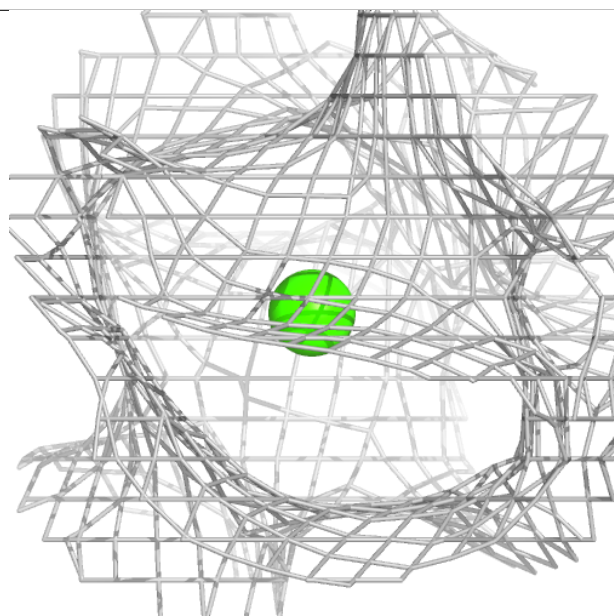
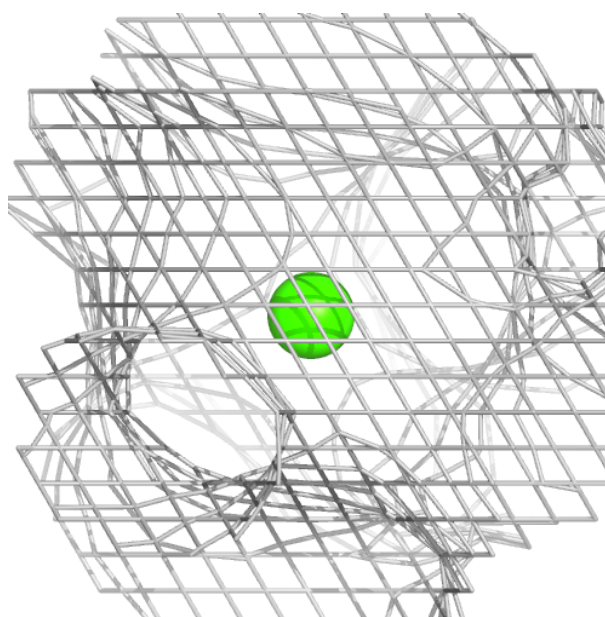
**Electron density around CA A 503:**

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



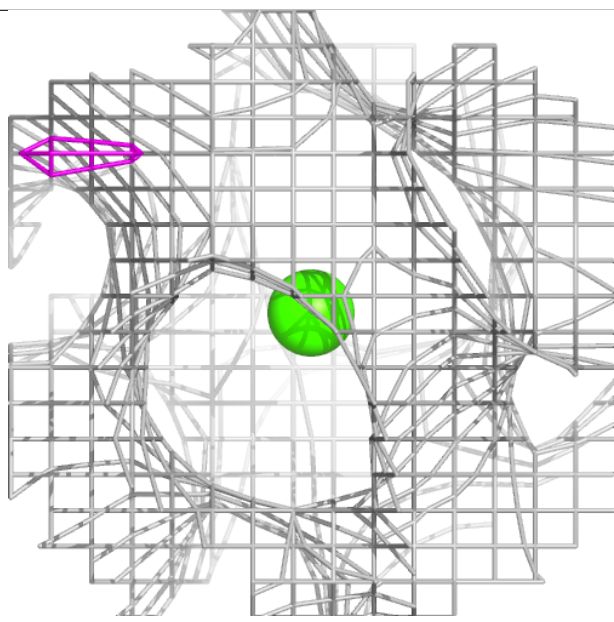
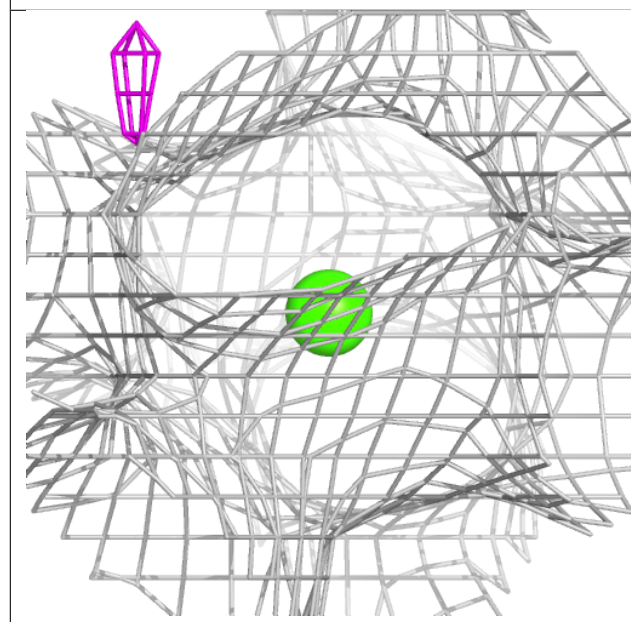
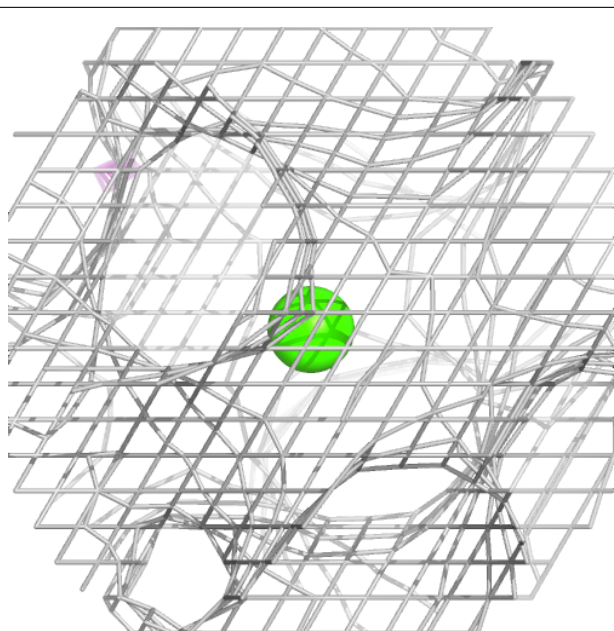
**Electron density around CA A 505:**

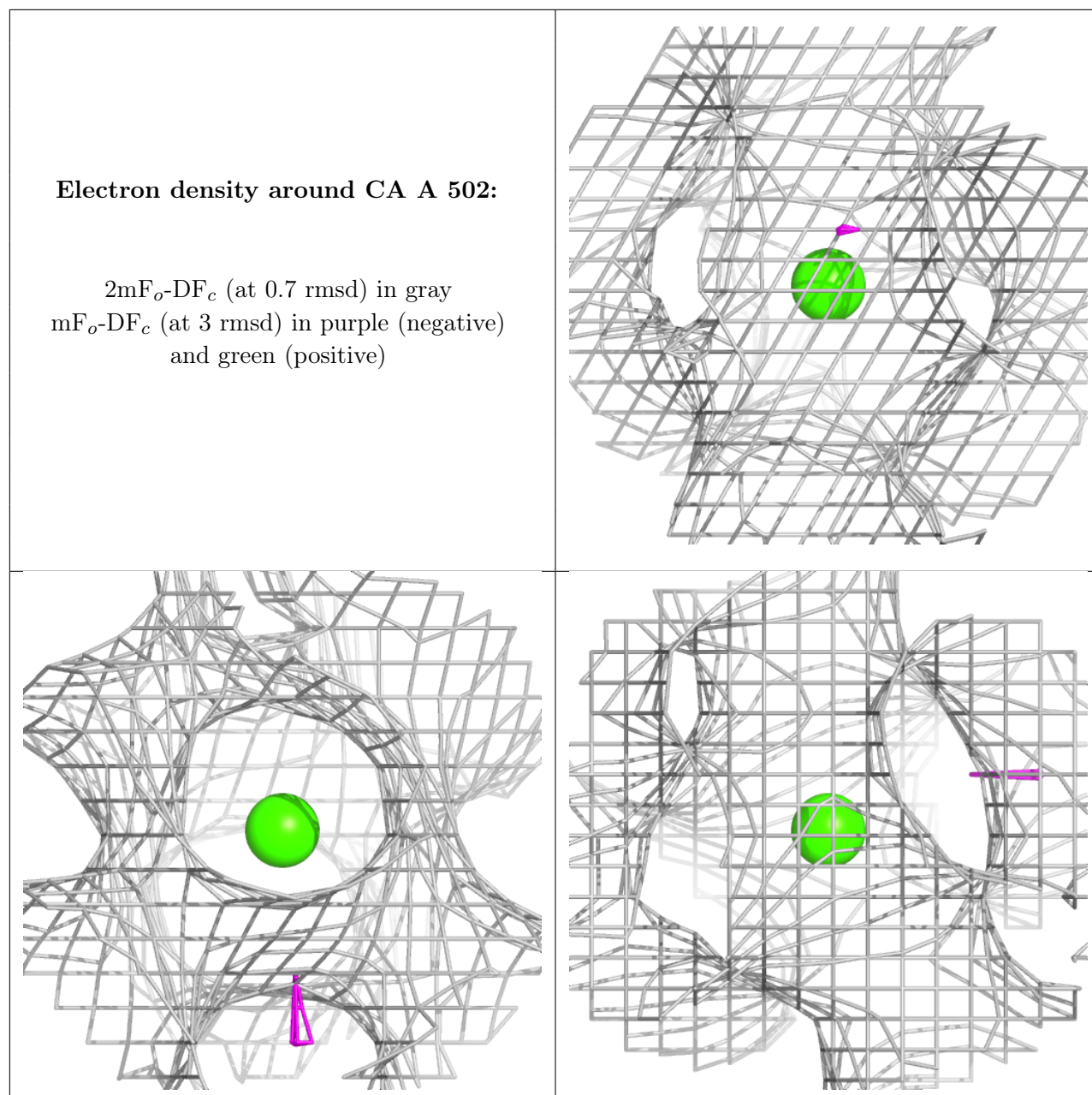
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



**Electron density around CA A 504:**

$2mF_o-DF_c$  (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.