



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 16, 2022 – 01:06 pm BST

PDB ID : 7Z4R
Title : Plasmodium falciparum pyruvate kinase mutant - C343A
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Deposited on : 2022-03-04
Resolution : 2.00 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at <http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4.02b-467
Xtrriage (Phenix) : 1.13
EDS : 2.29
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0267
CCP4 : 7.1.010 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.29

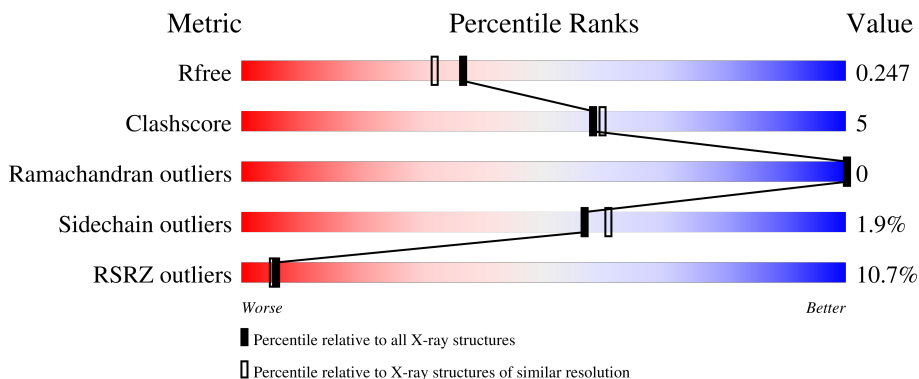
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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

Mol	Chain	Length	Quality of chain
1	A	519	

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 3943 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 Pyruvate kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	494	3767	2372	652	713	30	0	2	0

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	343	ALA	CYS	engineered mutation	UNP C6KTA4
A	512	LEU	-	expression tag	UNP C6KTA4
A	513	GLU	-	expression tag	UNP C6KTA4
A	514	HIS	-	expression tag	UNP C6KTA4
A	515	HIS	-	expression tag	UNP C6KTA4
A	516	HIS	-	expression tag	UNP C6KTA4
A	517	HIS	-	expression tag	UNP C6KTA4
A	518	HIS	-	expression tag	UNP C6KTA4
A	519	HIS	-	expression tag	UNP C6KTA4

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mg	0	0
			1	1		

- Molecule 3 is POTASSIUM ION (three-letter code: K) (formula: K) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	K	0	0
			1	1		

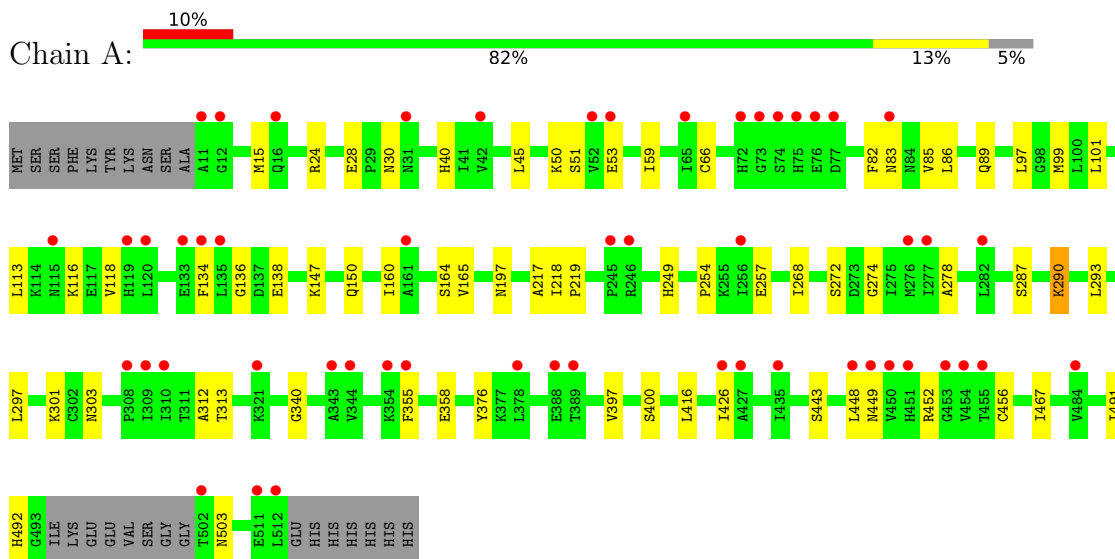
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	174	Total 174	O 174	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: Pyruvate kinase



4 Data and refinement statistics i

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	89.42Å 109.13Å 134.10Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	32.30 – 2.00 33.52 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.2 (32.30-2.00) 99.2 (33.52-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.07 (at 2.00Å)	Xtrriage
Refinement program	PHENIX 1.19.2_4158, PHENIX 1.19.2_4158	Depositor
R, R_{free}	0.215 , 0.248 0.212 , 0.247	Depositor DCC
R_{free} test set	3547 reflections (8.00%)	wwPDB-VP
Wilson B-factor (Å ²)	49.9	Xtrriage
Anisotropy	0.147	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	(Not available) , (Not available)	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3943	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.18% 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: MG, K

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	0.27	0/3818	0.48	0/5154

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	3767	0	3930	37	0
2	A	1	0	0	0	0
3	A	1	0	0	0	0
4	A	174	0	0	1	0
All	All	3943	0	3930	37	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 5.

All (37) 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:89:GLN:HE22	1:A:97:LEU:H	1.44	0.65
1:A:268:ILE:O	1:A:272:SER:OG	2.17	0.60
1:A:113:LEU:HD13	1:A:118:VAL:HG13	1.84	0.59
1:A:355:PHE:HB3	1:A:358:GLU:HB2	1.85	0.58
1:A:45:LEU:HD11	1:A:66:CYS:HB3	1.89	0.54
1:A:40:HIS:ND1	1:A:449:ASN:O	2.38	0.53
1:A:83:ASN:HA	1:A:86:LEU:HD12	1.90	0.53
1:A:467:ILE:HD12	1:A:467:ILE:H	1.73	0.53
1:A:138:GLU:H	1:A:138:GLU:CD	2.16	0.49
1:A:312:ALA:HB1	1:A:313:THR:HA	1.94	0.49
1:A:376:TYR:OH	1:A:452:ARG:NH1	2.46	0.48
1:A:449:ASN:ND2	4:A:713:HOH:O	2.44	0.48
1:A:24:ARG:NH2	1:A:28:GLU:OE2	2.48	0.47
1:A:249:HIS:HB3	1:A:443:SER:OG	2.15	0.47
1:A:297:LEU:HD11	1:A:301:LYS:HE2	1.96	0.47
1:A:82:PHE:O	1:A:86:LEU:HG	2.15	0.46
1:A:82:PHE:CE2	1:A:86:LEU:HD21	2.51	0.46
1:A:147:LYS:HG2	1:A:150:GLN:HB2	1.98	0.46
1:A:257:GLU:HG2	1:A:278:ALA:HB3	1.98	0.45
1:A:287:SER:HB2	1:A:290:LYS:HD3	1.98	0.45
1:A:85:VAL:HG11	1:A:99:MET:HE1	1.98	0.45
1:A:254:PRO:HD2	1:A:274:GLY:O	2.17	0.45
1:A:416:LEU:HB2	1:A:492:HIS:HB3	1.99	0.44
1:A:59:ILE:HD11	1:A:97:LEU:HD11	2.00	0.43
1:A:101:LEU:HD13	1:A:217:ALA:HB2	1.99	0.43
1:A:53:GLU:H	1:A:53:GLU:CD	2.21	0.43
1:A:303:ASN:HD21	1:A:340:GLY:HA3	1.83	0.43
1:A:160:ILE:HB	1:A:165:VAL:HG13	2.01	0.43
1:A:426:ILE:HD11	1:A:491:ILE:HG21	1.99	0.43
1:A:218:ILE:HB	1:A:219:PRO:HD3	2.01	0.42
1:A:290:LYS:HG3	1:A:293:LEU:HD12	2.01	0.42
1:A:51:SER:HB2	1:A:53:GLU:OE2	2.19	0.42
1:A:312:ALA:HB1	1:A:313:THR:HG22	2.02	0.42
1:A:448:LEU:HB2	1:A:456:CYS:SG	2.59	0.42
1:A:82:PHE:HE2	1:A:86:LEU:HD21	1.85	0.41
1:A:134:PHE:CE2	1:A:136:GLY:HA2	2.56	0.41
1:A:397:VAL:HG21	1:A:503:ASN:HB3	2.03	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	492/519 (95%)	482 (98%)	10 (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	A	425/445 (96%)	417 (98%)	8 (2%)	57 61

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	MET
1	A	30	ASN
1	A	50	LYS
1	A	116	LYS
1	A	164	SER
1	A	197	ASN
1	A	290	LYS
1	A	400	SER

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

Mol	Chain	Res	Type
1	A	30	ASN
1	A	89	GLN
1	A	119	HIS

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 2 ligands modelled in this entry, 2 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

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	A	494/519 (95%)	0.67	53 (10%) 6 5	32, 51, 76, 99	0

All (53) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	11	ALA	9.2
1	A	389	THR	6.7
1	A	12	GLY	6.5
1	A	512	LEU	4.7
1	A	119	HIS	4.3
1	A	74	SER	3.9
1	A	343	ALA	3.7
1	A	309	ILE	3.6
1	A	246	ARG	3.6
1	A	134	PHE	3.5
1	A	83	ASN	3.3
1	A	450	VAL	3.1
1	A	77	ASP	2.9
1	A	42	VAL	2.9
1	A	76	GLU	2.9
1	A	427	ALA	2.8
1	A	282	LEU	2.8
1	A	321	LYS	2.8
1	A	65	ILE	2.8
1	A	16	GLN	2.8
1	A	135	LEU	2.7
1	A	448	LEU	2.7
1	A	453	GLY	2.7
1	A	354	LYS	2.7
1	A	72	HIS	2.6
1	A	120	LEU	2.6
1	A	310	ILE	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	511	GLU	2.6
1	A	454	VAL	2.5
1	A	502	THR	2.5
1	A	435	ILE	2.5
1	A	484	VAL	2.5
1	A	355	PHE	2.5
1	A	75	HIS	2.4
1	A	455	THR	2.4
1	A	256	ILE	2.4
1	A	73	GLY	2.3
1	A	161	ALA	2.3
1	A	53	GLU	2.3
1	A	133	GLU	2.3
1	A	52	VAL	2.3
1	A	276	MET	2.2
1	A	115	ASN	2.2
1	A	426	ILE	2.2
1	A	344	VAL	2.2
1	A	449	ASN	2.2
1	A	245	PRO	2.2
1	A	388	GLU	2.1
1	A	308	PRO	2.1
1	A	277	ILE	2.1
1	A	451	HIS	2.1
1	A	378	LEU	2.1
1	A	31	ASN	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 [\(i\)](#)

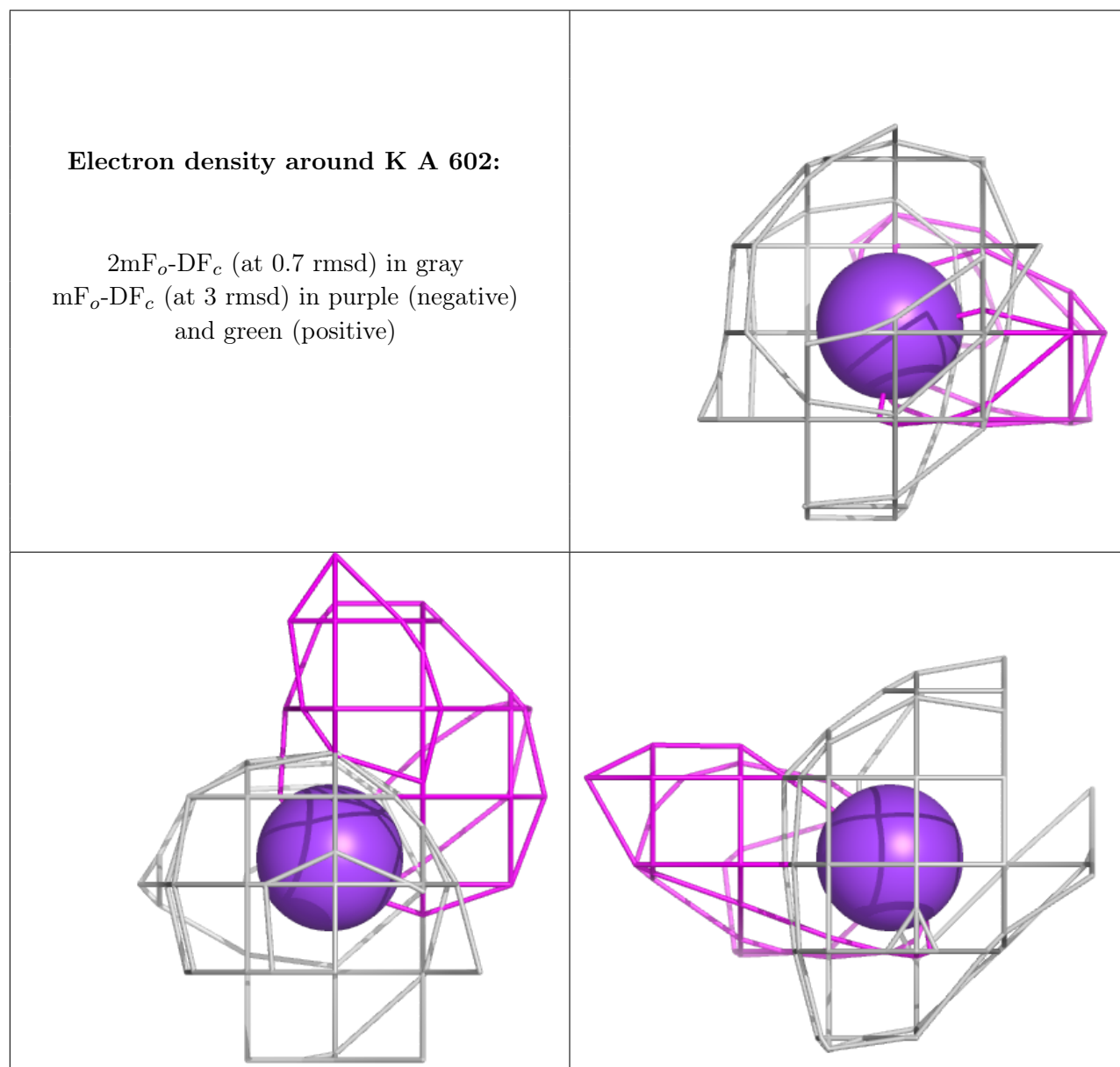
There are no monosaccharides in this entry.

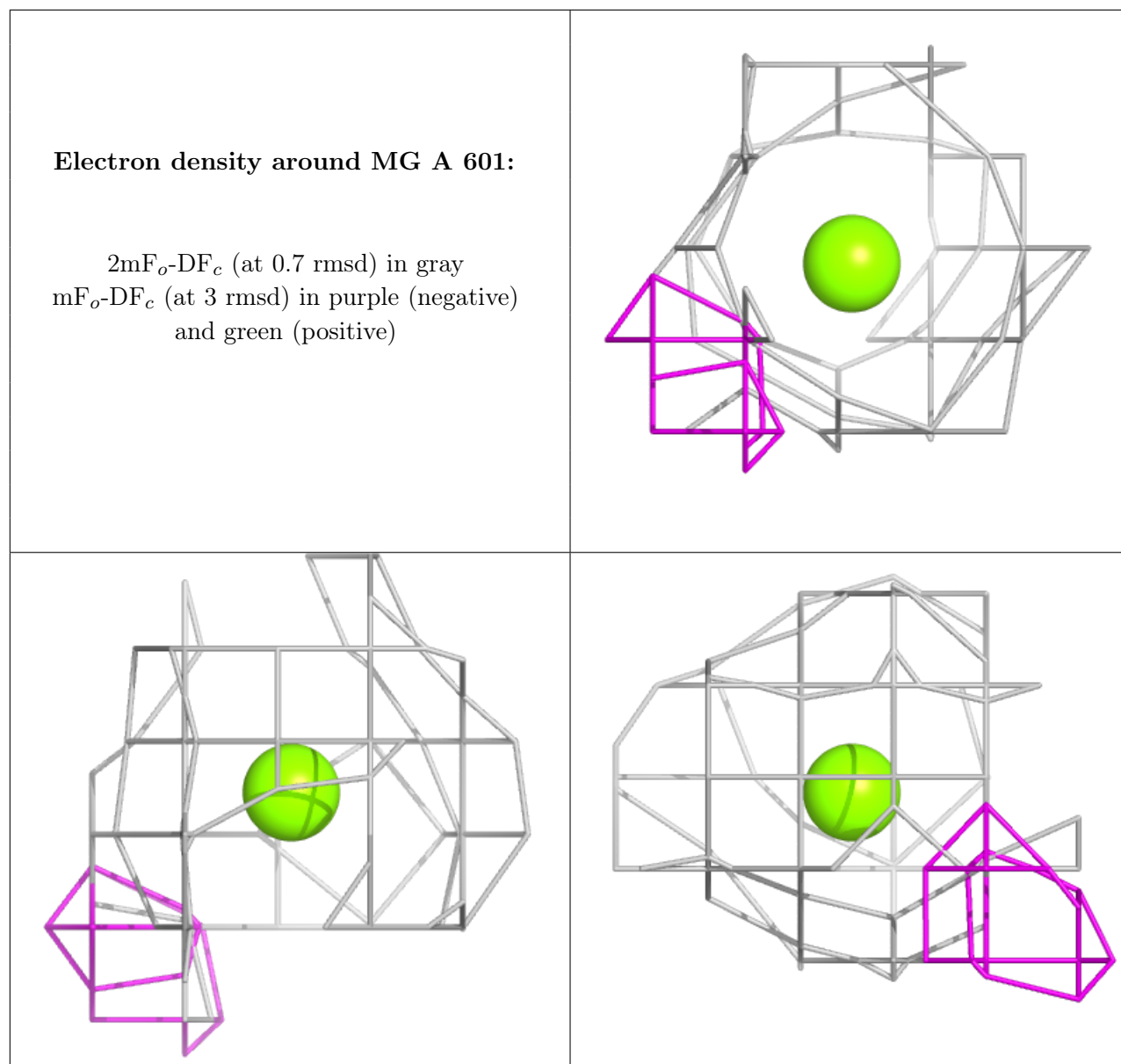
6.4 Ligands [\(i\)](#)

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, 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
3	K	A	602	1/1	0.83	0.15	80,80,80,80	0
2	MG	A	601	1/1	0.87	0.12	69,69,69,69	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 [i](#)

There are no such residues in this entry.