



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 24, 2022 – 08:13 PM JST

PDB ID : 7E06
Title : Trans-3/4-proline-hydroxylase H11 with AKG
Authors : Gong, W.G.; Yang, L.Y.
Deposited on : 2021-01-27
Resolution : 1.97 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 : 2.26
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) : 2.26

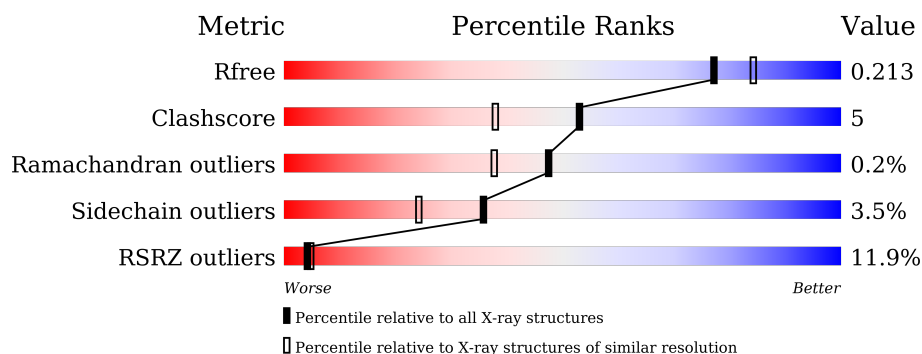
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.97 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	11647 (2.00-1.96)
Clashscore	141614	1014 (1.98-1.98)
Ramachandran outliers	138981	1006 (1.98-1.98)
Sidechain outliers	138945	1006 (1.98-1.98)
RSRZ outliers	127900	11410 (2.00-1.96)

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	268	<div> <div>9%</div> <div>85%</div> <div>7% • 6%</div> </div>
1	B	268	<div> <div>14%</div> <div>83%</div> <div>13% • •</div> </div>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4572 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.

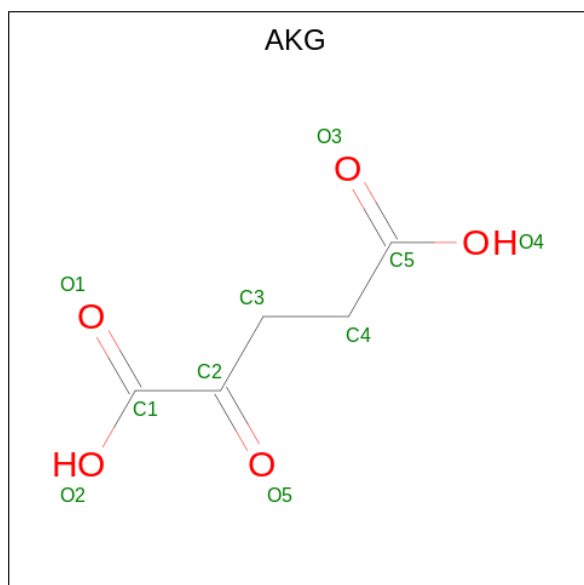
- Molecule 1 is a protein called Phytanoyl-CoA dioxygenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	251	Total	C	N	O	S	0	1	0
			2000	1270	357	367	6			
1	B	261	Total	C	N	O	S	0	1	0
			2079	1321	368	384	6			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLN	-	expression tag	UNP S5TUM1
A	0	ILE	-	expression tag	UNP S5TUM1
B	-1	GLN	-	expression tag	UNP S5TUM1
B	0	ILE	-	expression tag	UNP S5TUM1

- Molecule 2 is 2-OXOGLUTARIC ACID (three-letter code: AKG) (formula: $C_5H_6O_5$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			10	5	5		
2	B	1	Total	C	O	0	0
			10	5	5		

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Fe	0	0
			1	1		
3	B	1	Total	Fe	0	0
			1	1		

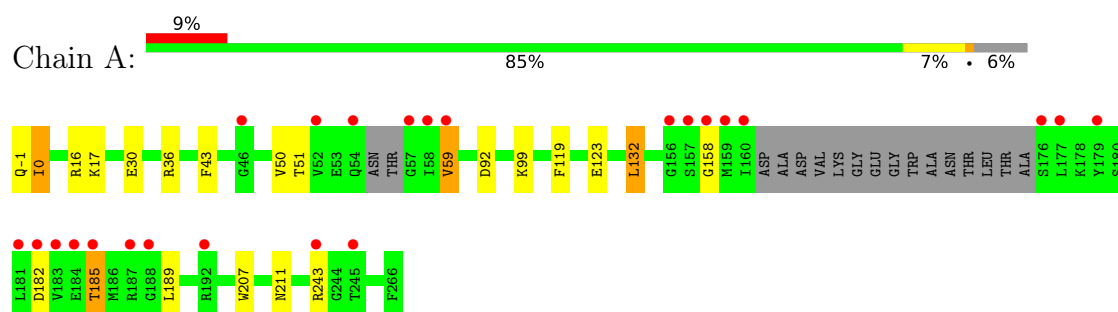
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	256	Total	O	0	0
			256	256		
4	B	215	Total	O	0	0
			215	215		

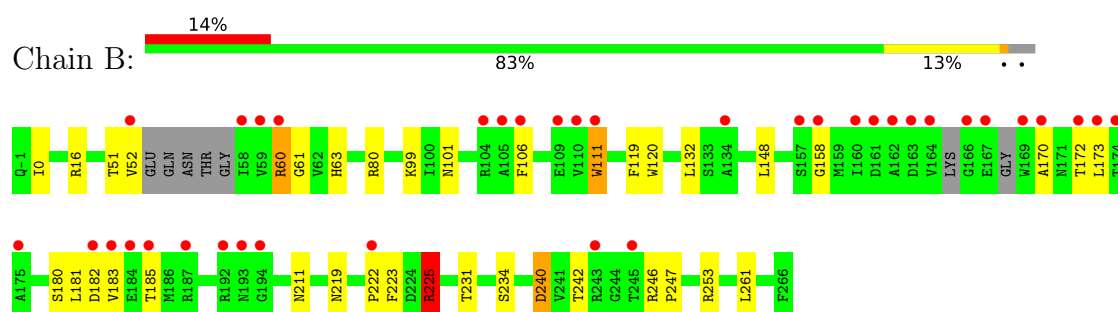
3 Residue-property plots [i](#)

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: Phytanoyl-CoA dioxygenase



• Molecule 1: Phytanoyl-CoA dioxygenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	106.76Å 106.76Å 144.01Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	85.76 – 1.97 39.79 – 1.97	Depositor EDS
% Data completeness (in resolution range)	99.2 (85.76-1.97) 99.2 (39.79-1.97)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	10.49 (at 1.98Å)	Xtrriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.169 , 0.204 0.181 , 0.213	Depositor DCC
R_{free} test set	2882 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	28.3	Xtrriage
Anisotropy	0.022	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 50.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4572	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 20.75 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 8.1213e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: FE, AKG

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	1.17	2/2052 (0.1%)	1.03	4/2784 (0.1%)
1	B	1.19	1/2132 (0.0%)	1.23	10/2896 (0.3%)
All	All	1.18	3/4184 (0.1%)	1.14	14/5680 (0.2%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	207	TRP	CE3-CZ3	5.30	1.47	1.38
1	A	30	GLU	CD-OE2	5.22	1.31	1.25
1	B	111	TRP	CB-CG	-5.13	1.41	1.50

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	80	ARG	NE-CZ-NH1	-19.43	110.58	120.30
1	B	80	ARG	NE-CZ-NH2	16.96	128.78	120.30
1	B	225	ARG	NE-CZ-NH1	10.56	125.58	120.30
1	B	80	ARG	CD-NE-CZ	9.46	136.85	123.60
1	B	225	ARG	NE-CZ-NH2	-8.70	115.95	120.30
1	B	80	ARG	CG-CD-NE	-8.69	93.55	111.80
1	B	80	ARG	CB-CG-CD	7.58	131.31	111.60
1	A	16	ARG	NE-CZ-NH1	7.12	123.86	120.30
1	A	36	ARG	NE-CZ-NH1	6.25	123.42	120.30
1	B	16	ARG	NE-CZ-NH1	6.04	123.32	120.30
1	A	92	ASP	CB-CG-OD2	-5.79	113.09	118.30
1	B	240	ASP	CB-CG-OD2	-5.54	113.31	118.30
1	A	132	LEU	CB-CG-CD2	-5.25	102.08	111.00
1	B	261	LEU	CB-CA-C	-5.16	100.40	110.20

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	2000	0	1927	12	0
1	B	2079	0	2004	31	0
2	A	10	0	4	0	0
2	B	10	0	4	2	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	256	0	0	6	0
4	B	215	0	0	4	0
All	All	4572	0	3939	43	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 (43) 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:B:225:ARG:O	1:B:225:ARG:HD2	1.83	0.79
1:A:0:ILE:HG22	4:A:803:HOH:O	1.86	0.75
1:A:0:ILE:CG2	4:A:803:HOH:O	2.38	0.71
1:B:111:TRP:CZ3	1:B:172:THR:O	2.47	0.67
1:A:50:VAL:HG11	1:A:59:VAL:HG12	1.81	0.62
1:B:240:ASP:OD1	1:B:242:THR:HB	1.99	0.62
1:B:111:TRP:CH2	1:B:172:THR:O	2.52	0.61
1:B:63:HIS:ND1	4:B:601:HOH:O	2.30	0.59
1:B:101:ASN:OD1	1:B:173:LEU:CD2	2.50	0.59
1:A:119:PHE:CE1	1:A:123:GLU:HG3	2.38	0.58
4:A:628:HOH:O	1:B:0:ILE:HG23	2.04	0.57
1:A:132:LEU:C	1:A:132:LEU:HD12	2.25	0.56
1:B:61:GLY:HA3	1:B:173:LEU:HD13	1.87	0.56
1:B:158:GLY:HA2	1:B:211:ASN:HD21	1.74	0.53
1:A:158:GLY:C	1:A:211:ASN:HD21	2.11	0.53
1:B:158:GLY:C	1:B:211:ASN:HD21	2.13	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:101:ASN:OD1	1:B:173:LEU:HD21	2.10	0.51
1:B:158:GLY:CA	1:B:211:ASN:HD21	2.24	0.51
1:B:253:ARG:NH1	4:B:604:HOH:O	2.44	0.50
1:B:101:ASN:OD1	1:B:173:LEU:HD22	2.10	0.50
1:B:158:GLY:HA2	1:B:211:ASN:ND2	2.27	0.49
1:B:106:PHE:CE2	1:B:222:PRO:HA	2.47	0.48
1:B:99:LYS:HE3	2:B:501:AKG:O1	2.14	0.48
4:A:628:HOH:O	1:B:0:ILE:CG2	2.61	0.47
1:B:172:THR:HG21	4:B:725:HOH:O	2.17	0.45
1:B:119:PHE:CD2	1:B:170:ALA:HB1	2.52	0.45
1:B:111:TRP:HH2	1:B:173:LEU:HD23	1.82	0.45
1:A:43:PHE:HB3	1:A:59:VAL:HG21	2.00	0.44
1:A:59:VAL:O	1:A:59:VAL:HG23	2.18	0.43
1:B:219:ASN:ND2	1:B:223:PHE:O	2.27	0.43
1:B:51:THR:HG23	4:B:601:HOH:O	2.19	0.43
1:B:132:LEU:C	1:B:132:LEU:HD12	2.39	0.43
1:B:111:TRP:CH2	1:B:173:LEU:HD23	2.54	0.42
1:A:50:VAL:HG12	1:A:51:THR:N	2.34	0.42
1:A:243:ARG:NH2	4:A:603:HOH:O	2.45	0.42
1:B:99:LYS:CE	2:B:501:AKG:O1	2.68	0.42
1:A:182:ASP:OD1	1:A:185:THR:N	2.40	0.41
1:B:111:TRP:CH2	1:B:172:THR:C	2.93	0.41
1:A:99:LYS:NZ	4:A:612:HOH:O	2.54	0.41
1:B:120:TRP:CH2	1:B:231:THR:HG21	2.55	0.41
1:B:111:TRP:CH2	1:B:173:LEU:HA	2.56	0.40
1:B:60:ARG:HD3	1:B:173:LEU:HB3	2.03	0.40
1:B:246:ARG:HB3	1:B:247:PRO:HD2	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	246/268 (92%)	242 (98%)	3 (1%)	1 (0%)	34	22
1	B	254/268 (95%)	245 (96%)	9 (4%)	0	100	100
All	All	500/536 (93%)	487 (97%)	12 (2%)	1 (0%)	47	38

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	59	VAL

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	210/225 (93%)	205 (98%)	5 (2%)	49	41
1	B	219/225 (97%)	209 (95%)	10 (5%)	27	14
All	All	429/450 (95%)	414 (96%)	15 (4%)	36	24

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	-1	GLN
1	A	0	ILE
1	A	17	LYS
1	A	185	THR
1	A	189	LEU
1	B	52	VAL
1	B	60	ARG
1	B	148	LEU
1	B	180	SER
1	B	181	LEU
1	B	182	ASP
1	B	183	VAL
1	B	185	THR
1	B	225	ARG
1	B	234	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	-1	GLN
1	A	41	GLN
1	A	63	HIS
1	A	211	ASN
1	B	41	GLN
1	B	63	HIS
1	B	97	GLN
1	B	121	HIS
1	B	193	ASN
1	B	211	ASN

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 ⓘ

There are no monosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 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	$\# Z > 2$
2	AKG	B	501	3	3,9,9	0.93	0	4,11,11	1.14	0
2	AKG	A	501	3	3,9,9	0.90	0	4,11,11	2.19	2 (50%)

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	AKG	B	501	3	-	0/3/9/9	-
2	AKG	A	501	3	-	0/3/9/9	-

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	AKG	O5-C2-C3	3.45	126.27	120.38
2	A	501	AKG	C4-C3-C2	2.31	118.12	113.14

There are no chirality outliers.

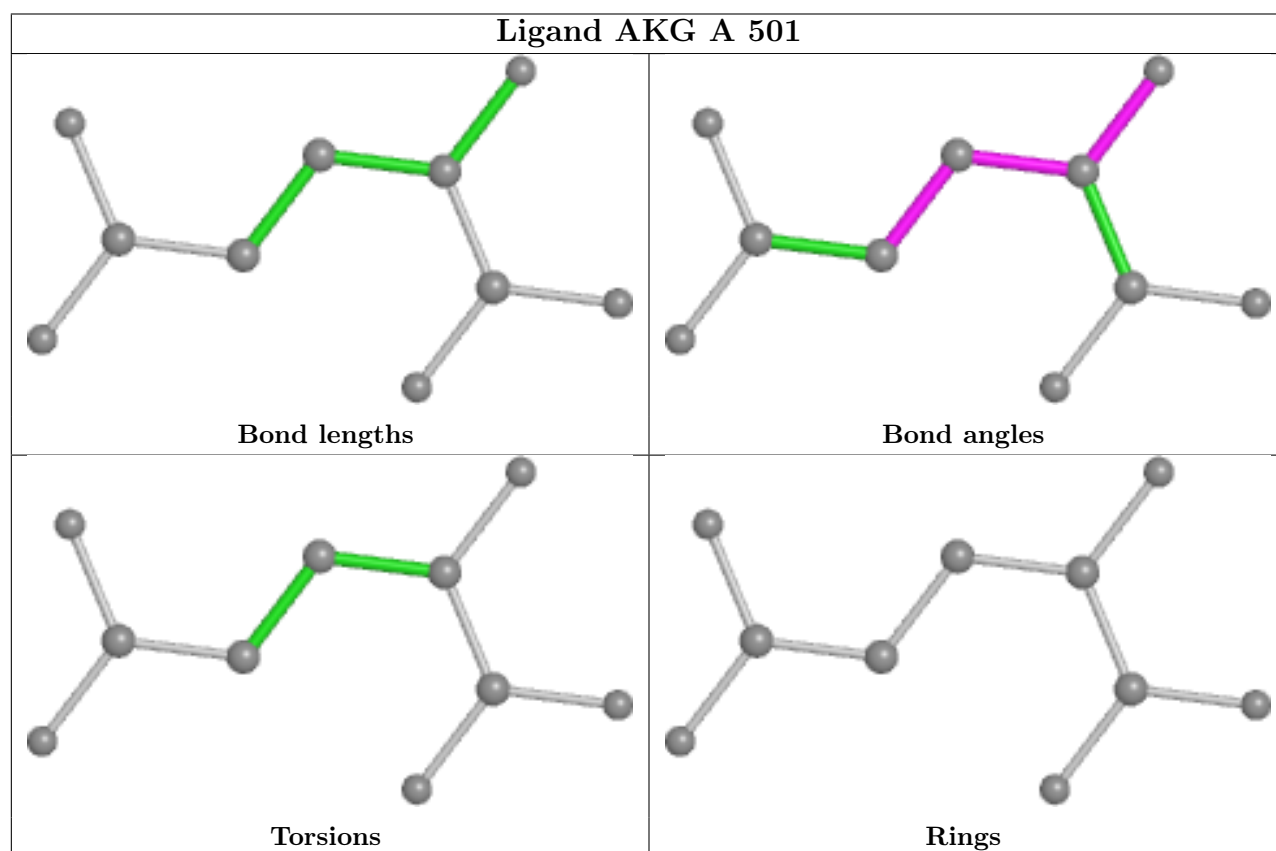
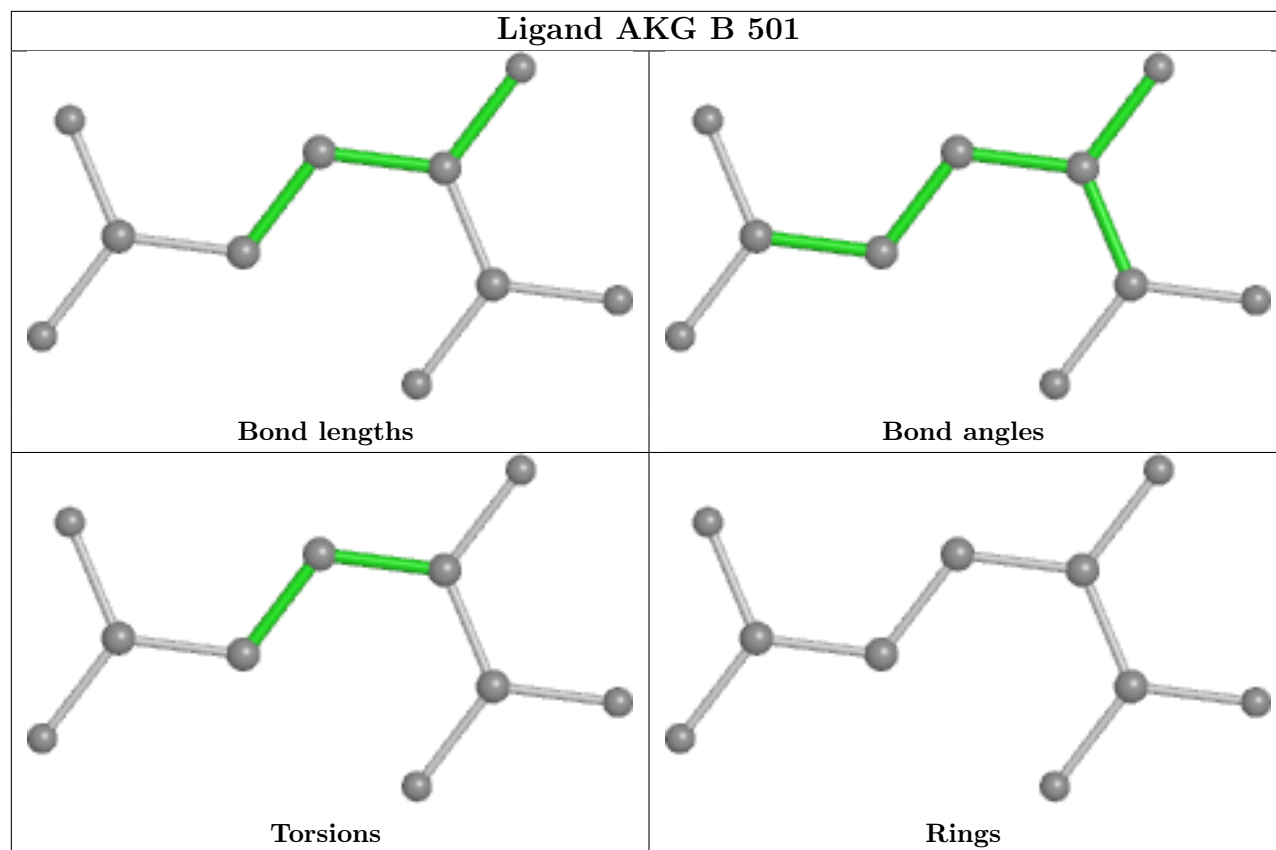
There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	501	AKG	2	0

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	A	251/268 (93%)	0.29	24 (9%) 8 9	19, 30, 77, 104	0
1	B	261/268 (97%)	0.50	37 (14%) 2 2	19, 33, 86, 114	0
All	All	512/536 (95%)	0.40	61 (11%) 4 5	19, 31, 83, 114	0

All (61) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	58	ILE	6.8
1	B	169	TRP	5.8
1	A	160	ILE	5.5
1	A	176	SER	5.4
1	B	160	ILE	5.4
1	B	183	VAL	5.4
1	A	57	GLY	5.3
1	B	105	ALA	5.1
1	B	106	PHE	5.1
1	B	164	VAL	4.9
1	B	167	GLU	4.9
1	B	161	ASP	4.8
1	B	166	GLY	4.6
1	B	173	LEU	4.5
1	A	158	GLY	4.4
1	B	163	ASP	4.3
1	A	159	MET	4.2
1	B	182	ASP	4.1
1	B	174	THR	4.1
1	A	184	GLU	4.0
1	A	58	ILE	3.9
1	B	172	THR	3.9
1	B	170	ALA	3.7
1	B	59	VAL	3.6

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Mol	Chain	Res	Type	RSRZ
1	A	185	THR	3.6
1	A	177	LEU	3.5
1	A	59	VAL	3.4
1	A	245	THR	3.3
1	B	52	VAL	3.3
1	A	182	ASP	3.3
1	A	157	SER	3.2
1	B	184	GLU	3.2
1	A	183	VAL	3.1
1	A	156	GLY	3.0
1	B	110	VAL	3.0
1	B	192	ARG	2.9
1	A	52	VAL	2.8
1	A	54	GLN	2.8
1	B	162	ALA	2.8
1	A	188	GLY	2.6
1	A	179	TYR	2.6
1	B	157	SER	2.6
1	A	46	GLY	2.6
1	A	181	LEU	2.6
1	B	222	PRO	2.6
1	B	111	TRP	2.4
1	B	104	ARG	2.4
1	B	109	GLU	2.4
1	A	192	ARG	2.4
1	B	243	ARG	2.4
1	B	187	ARG	2.3
1	A	243	ARG	2.3
1	B	134	ALA	2.3
1	B	193	ASN	2.2
1	B	158	GLY	2.2
1	B	185	THR	2.2
1	B	245	THR	2.2
1	B	175	ALA	2.2
1	A	187	ARG	2.1
1	B	60	ARG	2.1
1	B	194	GLY	2.0

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 [i](#)

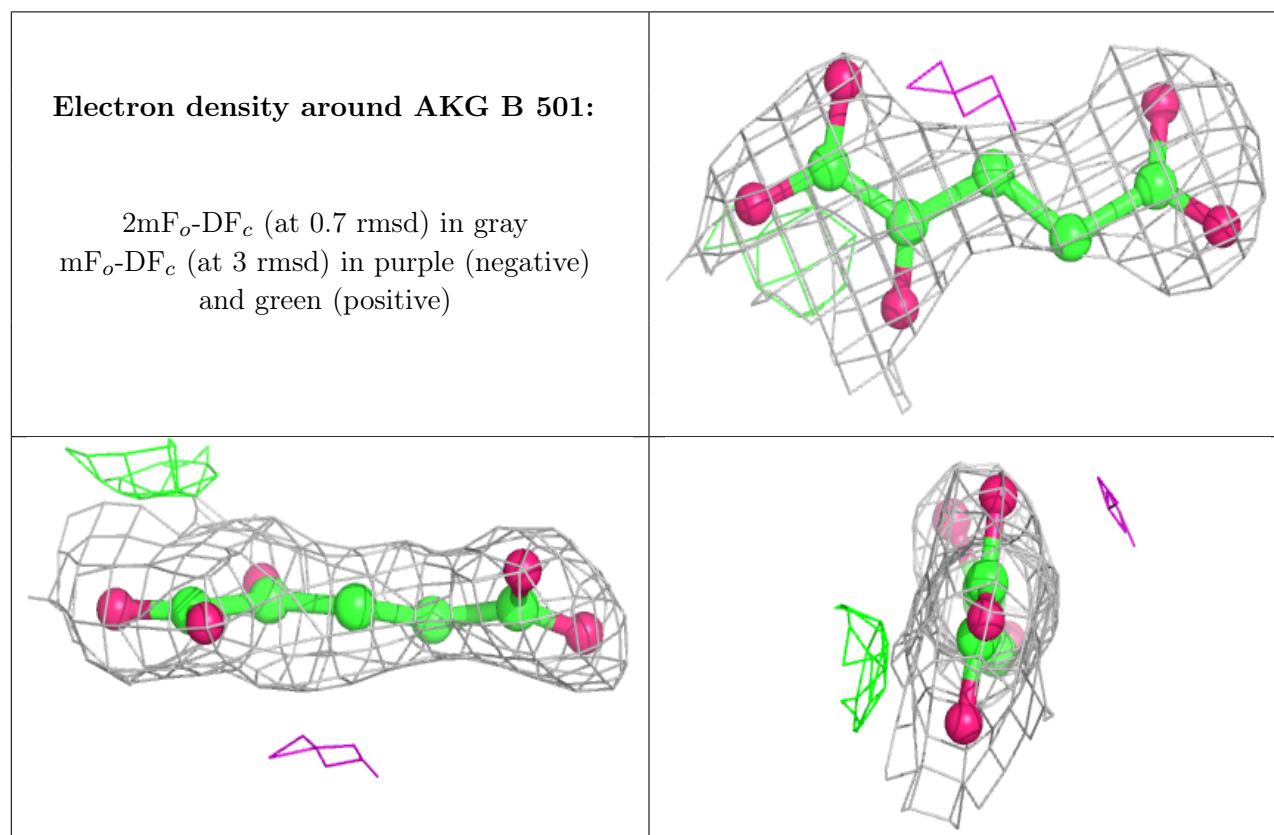
There are no monosaccharides in this entry.

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.

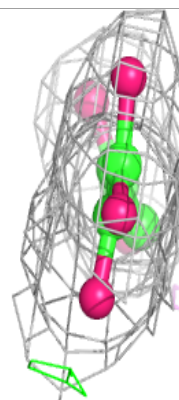
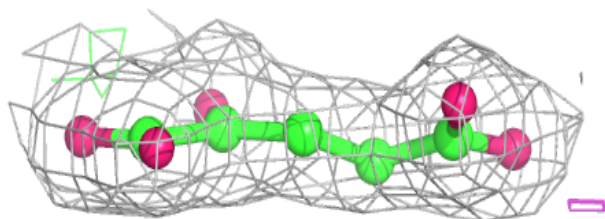
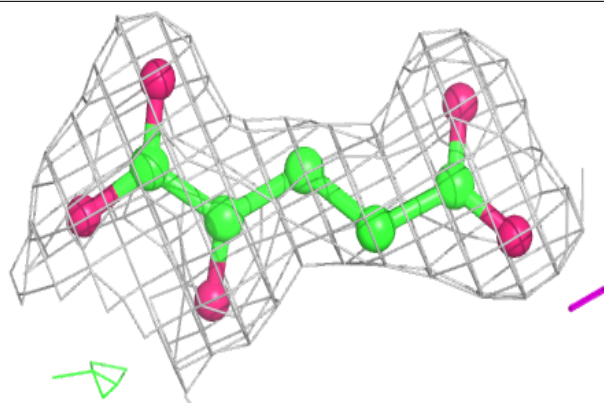
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	AKG	B	501	10/10	0.94	0.13	43,61,66,66	0
2	AKG	A	501	10/10	0.98	0.09	34,40,44,47	0
3	FE	A	502	1/1	0.99	0.14	18,18,18,18	1
3	FE	B	502	1/1	0.99	0.04	34,34,34,34	1

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.



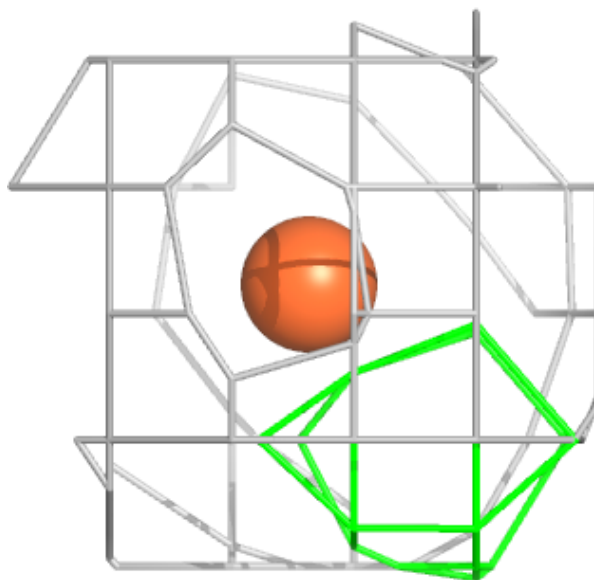
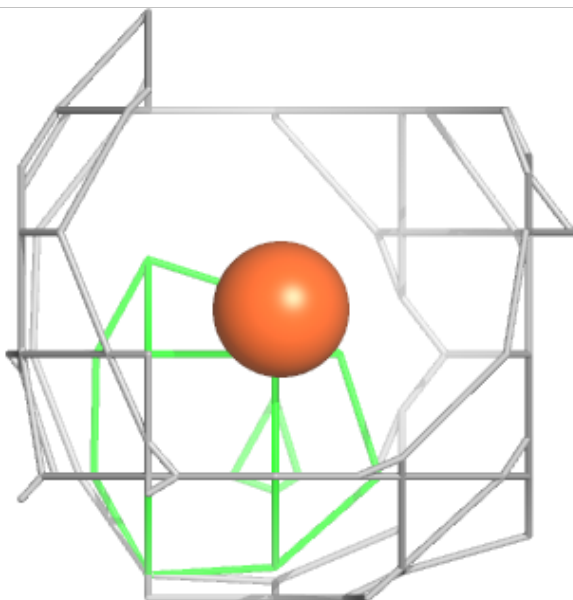
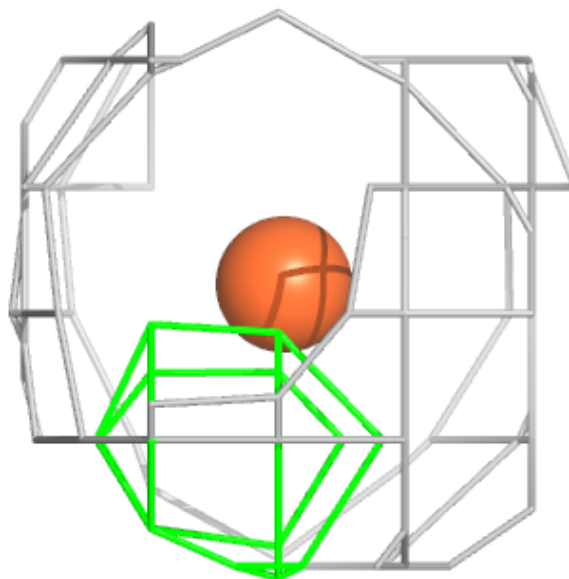
Electron density around AKG A 501:

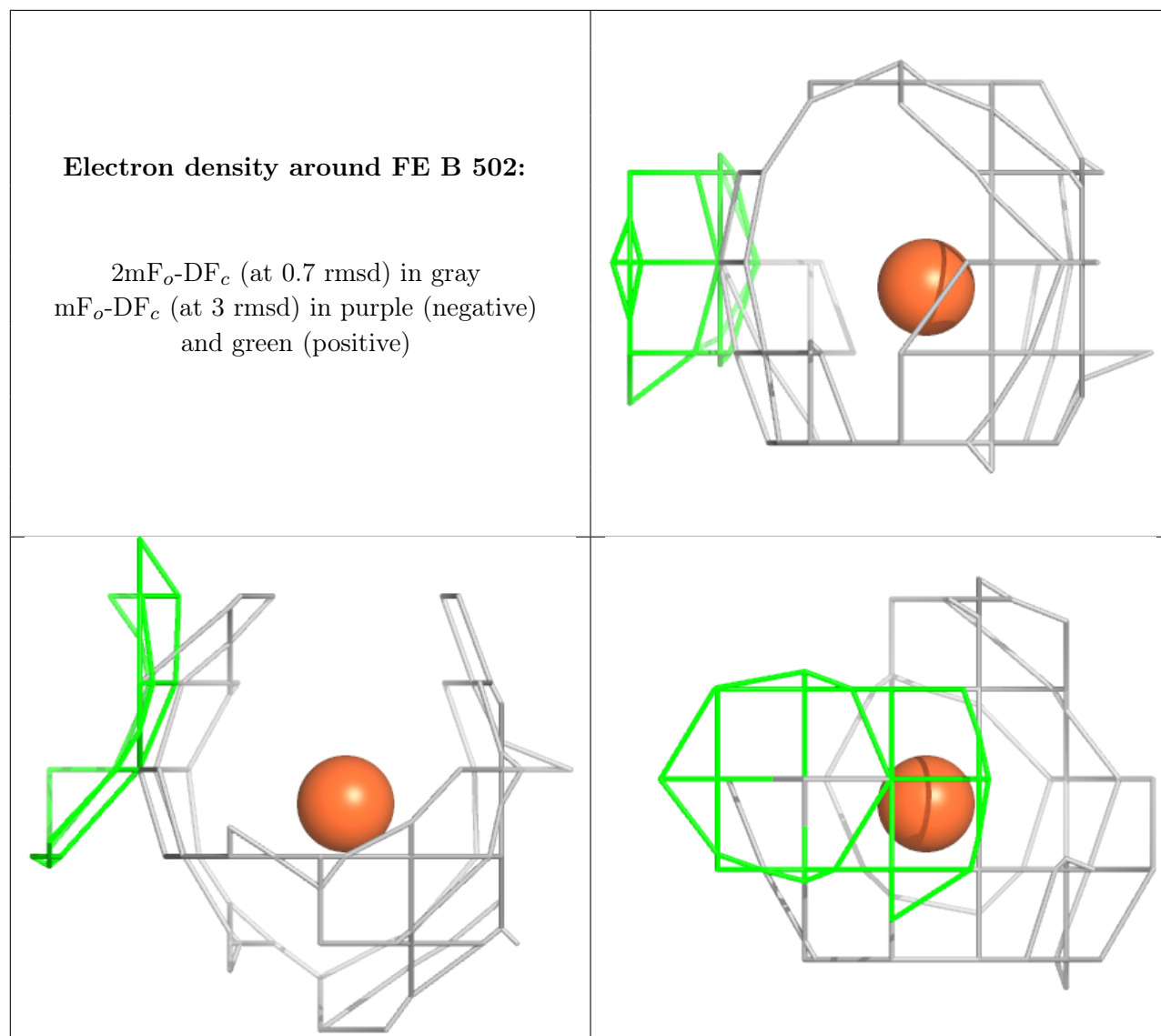
$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around FE A 502:

$2mF_o-DF_c$ (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.