



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 18, 2024 – 11:12 AM EST

PDB ID : 4EIQ
Title : Chromopyrrolic acid-soaked RebC-10x with bound 7-carboxy-K252c
Authors : Goldman, P.J.; Ryan, K.S.; Howard-Jones, A.R.; Hamill, M.J.; Elliott, S.J.;
Walsh, C.T.; Drennan, C.L.
Deposited on : 2012-04-05
Resolution : 2.76 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.36
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

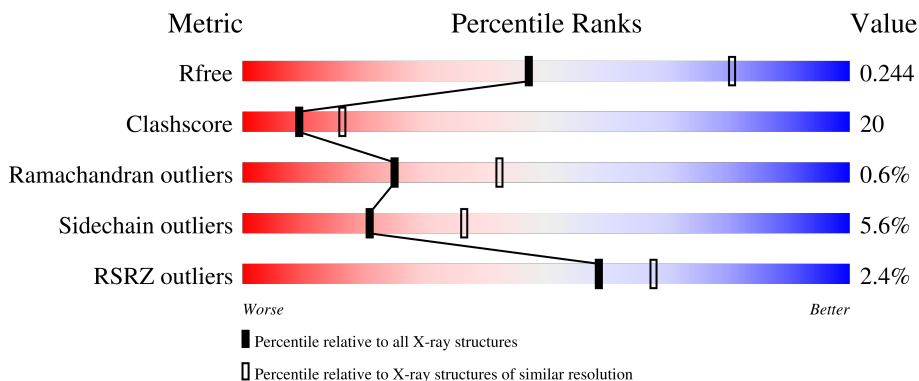
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.76 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	1235 (2.78-2.74)
Clashscore	141614	1277 (2.78-2.74)
Ramachandran outliers	138981	1257 (2.78-2.74)
Sidechain outliers	138945	1257 (2.78-2.74)
RSRZ outliers	127900	1207 (2.78-2.74)

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	549	 3% 57% 33% • 6%
1	B	549	 % 61% 30% • 7%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 7928 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 Putative FAD-monoxygenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	514	3925	2463	742	710	10	0	0	0
1	B	511	3887	2444	735	699	9	0	0	0

There are 60 discrepancies between the modelled and reference sequences:

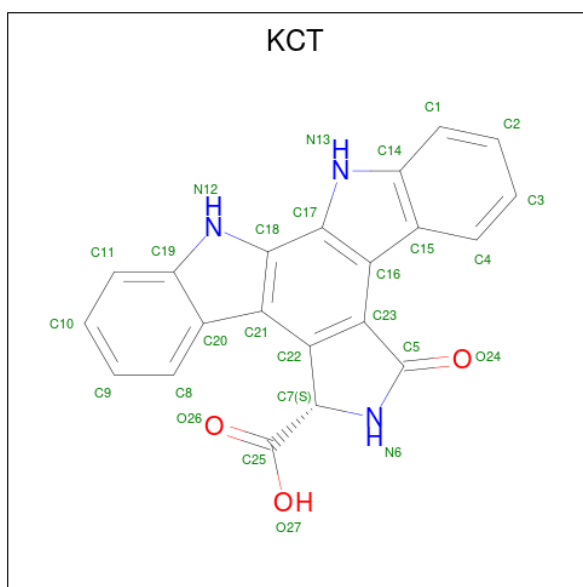
Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	expression tag	UNP Q8KI25
A	-18	GLY	-	expression tag	UNP Q8KI25
A	-17	SER	-	expression tag	UNP Q8KI25
A	-16	SER	-	expression tag	UNP Q8KI25
A	-15	HIS	-	expression tag	UNP Q8KI25
A	-14	HIS	-	expression tag	UNP Q8KI25
A	-13	HIS	-	expression tag	UNP Q8KI25
A	-12	HIS	-	expression tag	UNP Q8KI25
A	-11	HIS	-	expression tag	UNP Q8KI25
A	-10	HIS	-	expression tag	UNP Q8KI25
A	-9	SER	-	expression tag	UNP Q8KI25
A	-8	SER	-	expression tag	UNP Q8KI25
A	-7	GLY	-	expression tag	UNP Q8KI25
A	-6	LEU	-	expression tag	UNP Q8KI25
A	-5	VAL	-	expression tag	UNP Q8KI25
A	-4	PRO	-	expression tag	UNP Q8KI25
A	-3	ARG	-	expression tag	UNP Q8KI25
A	-2	GLY	-	expression tag	UNP Q8KI25
A	-1	SER	-	expression tag	UNP Q8KI25
A	0	HIS	-	expression tag	UNP Q8KI25
A	36	ASP	GLU	engineered mutation	UNP Q8KI25
A	37	ALA	GLN	engineered mutation	UNP Q8KI25
A	38	GLY	THR	engineered mutation	UNP Q8KI25
A	46	LYS	ARG	engineered mutation	UNP Q8KI25
A	48	SER	GLY	engineered mutation	UNP Q8KI25

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Chain	Residue	Modelled	Actual	Comment	Reference
A	117	ALA	GLN	engineered mutation	UNP Q8KI25
A	216	VAL	PHE	engineered mutation	UNP Q8KI25
A	231	SER	ALA	engineered mutation	UNP Q8KI25
A	239	ASN	ARG	engineered mutation	UNP Q8KI25
A	241	VAL	THR	engineered mutation	UNP Q8KI25
B	-19	MET	-	expression tag	UNP Q8KI25
B	-18	GLY	-	expression tag	UNP Q8KI25
B	-17	SER	-	expression tag	UNP Q8KI25
B	-16	SER	-	expression tag	UNP Q8KI25
B	-15	HIS	-	expression tag	UNP Q8KI25
B	-14	HIS	-	expression tag	UNP Q8KI25
B	-13	HIS	-	expression tag	UNP Q8KI25
B	-12	HIS	-	expression tag	UNP Q8KI25
B	-11	HIS	-	expression tag	UNP Q8KI25
B	-10	HIS	-	expression tag	UNP Q8KI25
B	-9	SER	-	expression tag	UNP Q8KI25
B	-8	SER	-	expression tag	UNP Q8KI25
B	-7	GLY	-	expression tag	UNP Q8KI25
B	-6	LEU	-	expression tag	UNP Q8KI25
B	-5	VAL	-	expression tag	UNP Q8KI25
B	-4	PRO	-	expression tag	UNP Q8KI25
B	-3	ARG	-	expression tag	UNP Q8KI25
B	-2	GLY	-	expression tag	UNP Q8KI25
B	-1	SER	-	expression tag	UNP Q8KI25
B	0	HIS	-	expression tag	UNP Q8KI25
B	36	ASP	GLU	engineered mutation	UNP Q8KI25
B	37	ALA	GLN	engineered mutation	UNP Q8KI25
B	38	GLY	THR	engineered mutation	UNP Q8KI25
B	46	LYS	ARG	engineered mutation	UNP Q8KI25
B	48	SER	GLY	engineered mutation	UNP Q8KI25
B	117	ALA	GLN	engineered mutation	UNP Q8KI25
B	216	VAL	PHE	engineered mutation	UNP Q8KI25
B	231	SER	ALA	engineered mutation	UNP Q8KI25
B	239	ASN	ARG	engineered mutation	UNP Q8KI25
B	241	VAL	THR	engineered mutation	UNP Q8KI25

- Molecule 2 is (5S)-7-oxo-6,7,12,13-tetrahydro-5H-indolo[2,3-a]pyrrolo[3,4-c]carbazole-5-carboxylic acid (three-letter code: KCT) (formula: C₂₁H₁₃N₃O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	
			Total	C	N			O
2	A	1	27	21	3	3	0	0
2	B	1	27	21	3	3	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	36	36	36	0	0
3	B	26	26	26	0	0

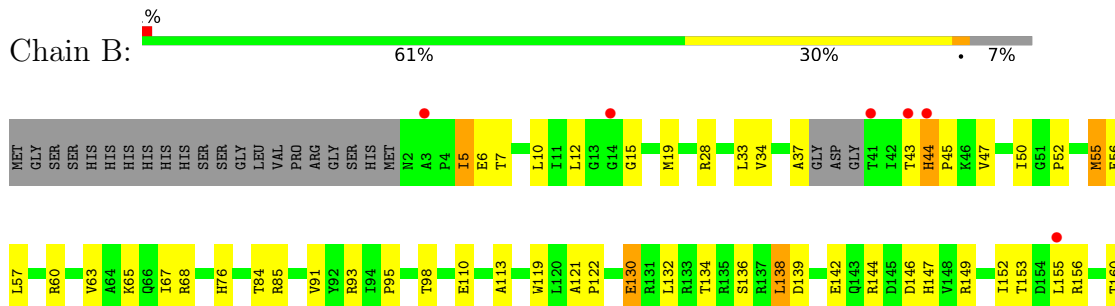
3 Residue-property plots [i](#)

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: Putative FAD-monooxygenase



• Molecule 1: Putative FAD-monooxygenase



R161	IYS	L366	T463
A162	SER	P367	D464
V163	THR	L370	G465
H164	MET	H371	V466
A165	D253	R376	V469
L168	F255	R379	T470
D172	R259	I380	R471
S175	R260	R381	A472
S176	T266	V384	F473
P177	E267	S391	R476
K180	L278	R394	L480
P187	T279	R395	A487
P188	H280	E396	P488
R189	S287	F397	E489
H190	F292	D398	I490
V194	L293	A399	H491
F195	T294	I402	R496
R196	T300	H403	A497
L199	L301	F404	B510
P203	S302	G405	E517
R212	S304	H406	V522
A213	F307	R409	R526
E214	G308	S410	G527
L215	M309	E416	A528
V216	L323	P417	A529
F217	L326	GLU	
F218	T326	THR	
L219	A331	GLU	
M220	L335	VAL	
L221	L336	ALA	
S222	L336	THR	
S223	Y339	GLY	
R226	E340	GLY	
F227	E341	W426	
P228	R344	H438	
D233	P345	A439	
L237	V346	T444	
Y238	L351	T445	
M239	R359	S446	
L240	G243	T447	
V241	T361	L448	
V242	R360	D449	
G243	M362	G452	
V244	D363	F455	
D245	R364	S459	
D246	E365		
ALA			
SER			

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	63.18Å 77.70Å 123.14Å 90.00° 98.78° 90.00°	Depositor
Resolution (Å)	47.30 – 2.76 47.34 – 2.76	Depositor EDS
% Data completeness (in resolution range)	94.4 (47.30-2.76) 94.6 (47.34-2.76)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.38 (at 2.77Å)	Xtrriage
Refinement program	CNS	Depositor
R, R_{free}	0.206 , 0.255 0.199 , 0.244	Depositor DCC
R_{free} test set	1458 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	62.9	Xtrriage
Anisotropy	0.172	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 41.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7928	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.18% 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: KCT

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	0.40	0/4020	0.71	10/5469 (0.2%)
1	B	0.42	0/3983	0.66	0/5423
All	All	0.41	0/8003	0.69	10/10892 (0.1%)

There are no bond length outliers.

The worst 5 of 10 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	486	HIS	CB-CA-C	-5.87	98.67	110.40
1	A	516	ALA	CB-CA-C	-5.63	101.66	110.10
1	A	46	LYS	CB-CA-C	-5.45	99.49	110.40
1	A	375	PRO	N-CA-C	5.43	126.21	112.10
1	A	516	ALA	N-CA-C	5.38	125.51	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	3925	0	3857	166	0
1	B	3887	0	3811	142	0
2	A	27	0	12	1	0
2	B	27	0	12	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	36	0	0	5	0
3	B	26	0	0	0	0
All	All	7928	0	7692	308	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 20.

The worst 5 of 308 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:399:ALA:HB1	1:A:402:ILE:HD13	1.14	1.13
1:B:44:HIS:HB3	1:B:45:PRO:CD	1.79	1.11
1:B:44:HIS:CB	1:B:45:PRO:CD	2.28	1.11
1:B:44:HIS:HB3	1:B:45:PRO:HD3	1.36	1.06
1:A:46:LYS:C	1:A:47:VAL:HG22	1.75	1.04

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	506/549 (92%)	454 (90%)	48 (10%)	4 (1%)	19	34
1	B	503/549 (92%)	471 (94%)	30 (6%)	2 (0%)	34	53
All	All	1009/1098 (92%)	925 (92%)	78 (8%)	6 (1%)	25	42

5 of 6 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	44	HIS
1	A	155	LEU

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Mol	Chain	Res	Type
1	A	45	PRO
1	B	304	SER
1	A	368	PRO

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	396/432 (92%)	375 (95%)	21 (5%)	22	38
1	B	390/432 (90%)	367 (94%)	23 (6%)	19	34
All	All	786/864 (91%)	742 (94%)	44 (6%)	21	36

5 of 44 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	139	ASP
1	B	267	GLU
1	B	155	LEU
1	B	253	ASP
1	B	301	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 7 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	193	GLN
1	B	280	HIS
1	B	406	HIS
1	B	371	HIS
1	A	193	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](#)

2 ligands 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	KCT	B	600	-	25,32,32	1.96	7 (28%)	31,50,50	2.54	10 (32%)
2	KCT	A	600	-	25,32,32	1.93	9 (36%)	31,50,50	2.43	9 (29%)

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	KCT	B	600	-	-	3/4/16/16	0/6/6/6
2	KCT	A	600	-	-	3/4/16/16	0/6/6/6

The worst 5 of 16 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	600	KCT	C23-C16	5.90	1.52	1.43
2	A	600	KCT	C23-C16	5.20	1.51	1.43
2	B	600	KCT	O24-C5	3.29	1.30	1.23
2	A	600	KCT	C22-C7	-3.11	1.47	1.52
2	B	600	KCT	C3-C4	2.79	1.43	1.36

The worst 5 of 19 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	600	KCT	C23-C22-C21	-10.91	116.70	122.04
2	A	600	KCT	C23-C22-C21	-9.13	117.57	122.04
2	A	600	KCT	C25-C7-N6	-4.89	105.46	112.10
2	A	600	KCT	C22-C21-C18	-3.59	119.25	122.71
2	B	600	KCT	C22-C21-C18	-3.59	119.25	122.71

There are no chirality outliers.

5 of 6 torsion outliers are listed below:

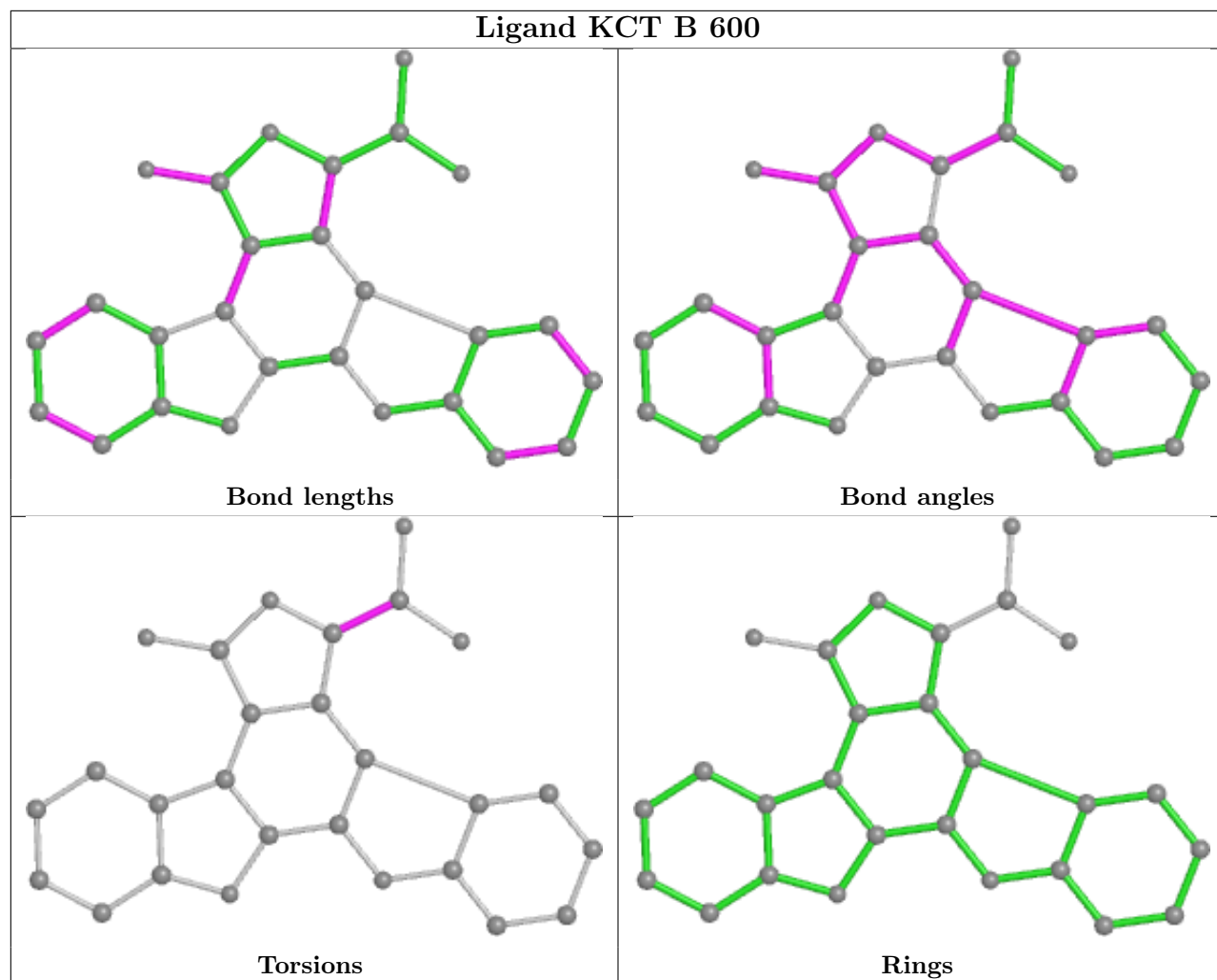
Mol	Chain	Res	Type	Atoms
2	A	600	KCT	O27-C25-C7-N6
2	A	600	KCT	O26-C25-C7-N6
2	A	600	KCT	O26-C25-C7-C22
2	B	600	KCT	O27-C25-C7-N6
2	B	600	KCT	O26-C25-C7-N6

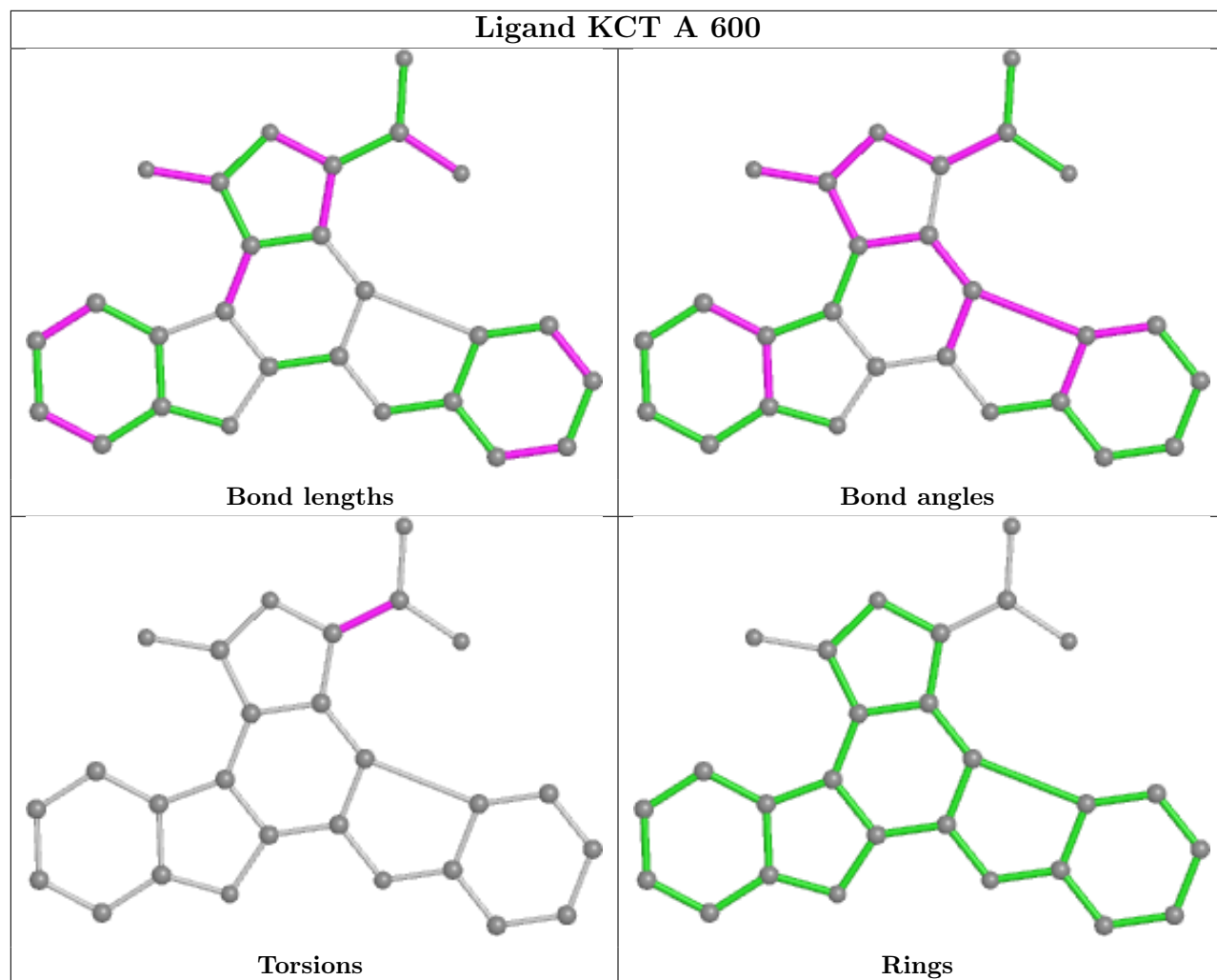
There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	600	KCT	1	0
2	A	600	KCT	1	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 [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, 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	514/549 (93%)	0.30	18 (3%) 44 52	45, 66, 91, 117	0
1	B	511/549 (93%)	0.10	7 (1%) 75 82	34, 52, 80, 108	0
All	All	1025/1098 (93%)	0.20	25 (2%) 59 68	34, 59, 88, 117	0

The worst 5 of 25 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	44	HIS	4.6
1	A	3	ALA	4.1
1	A	246	ASP	4.1
1	A	529	ALA	4.0
1	A	494	TYR	3.9

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](#)

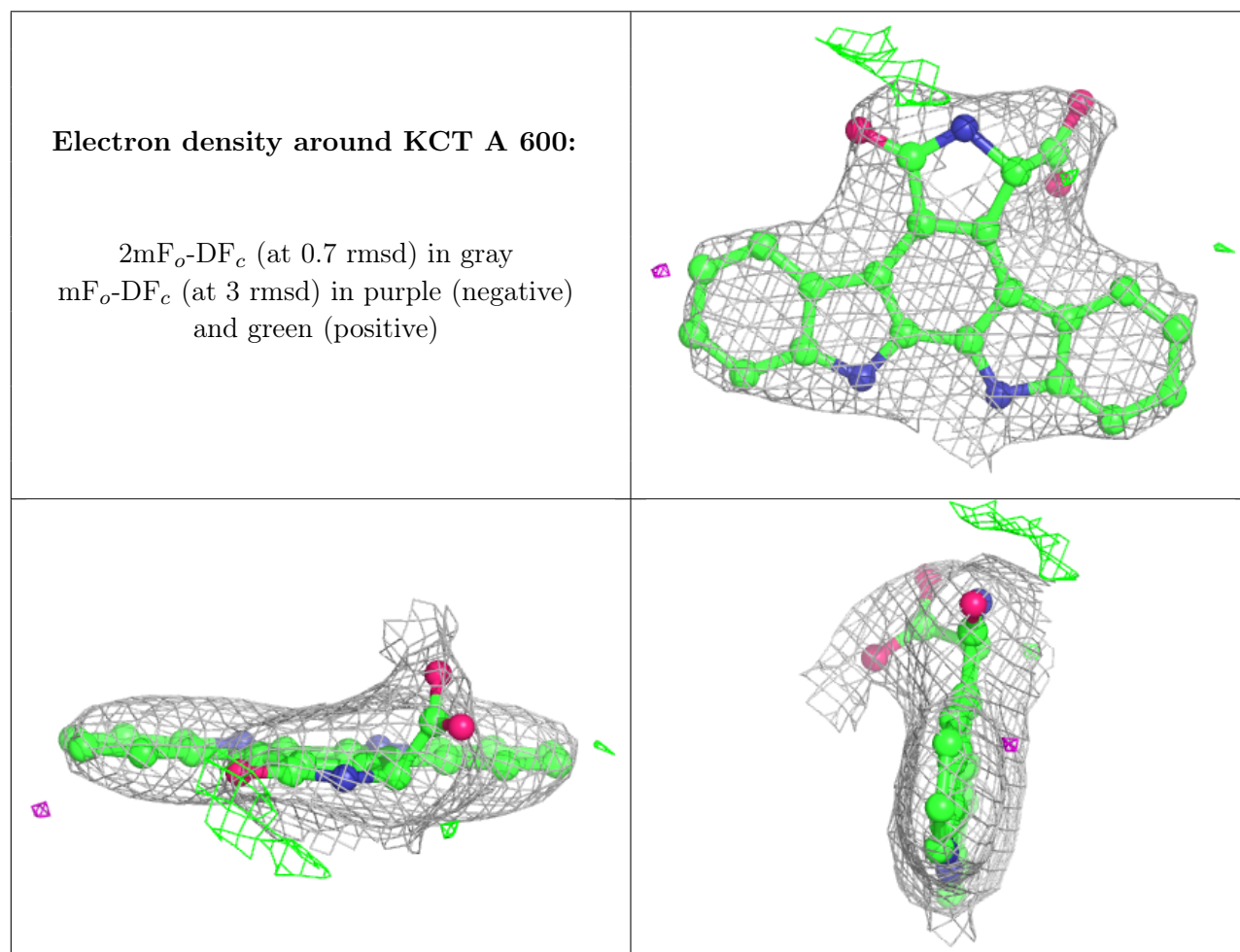
There are no monosaccharides in this entry.

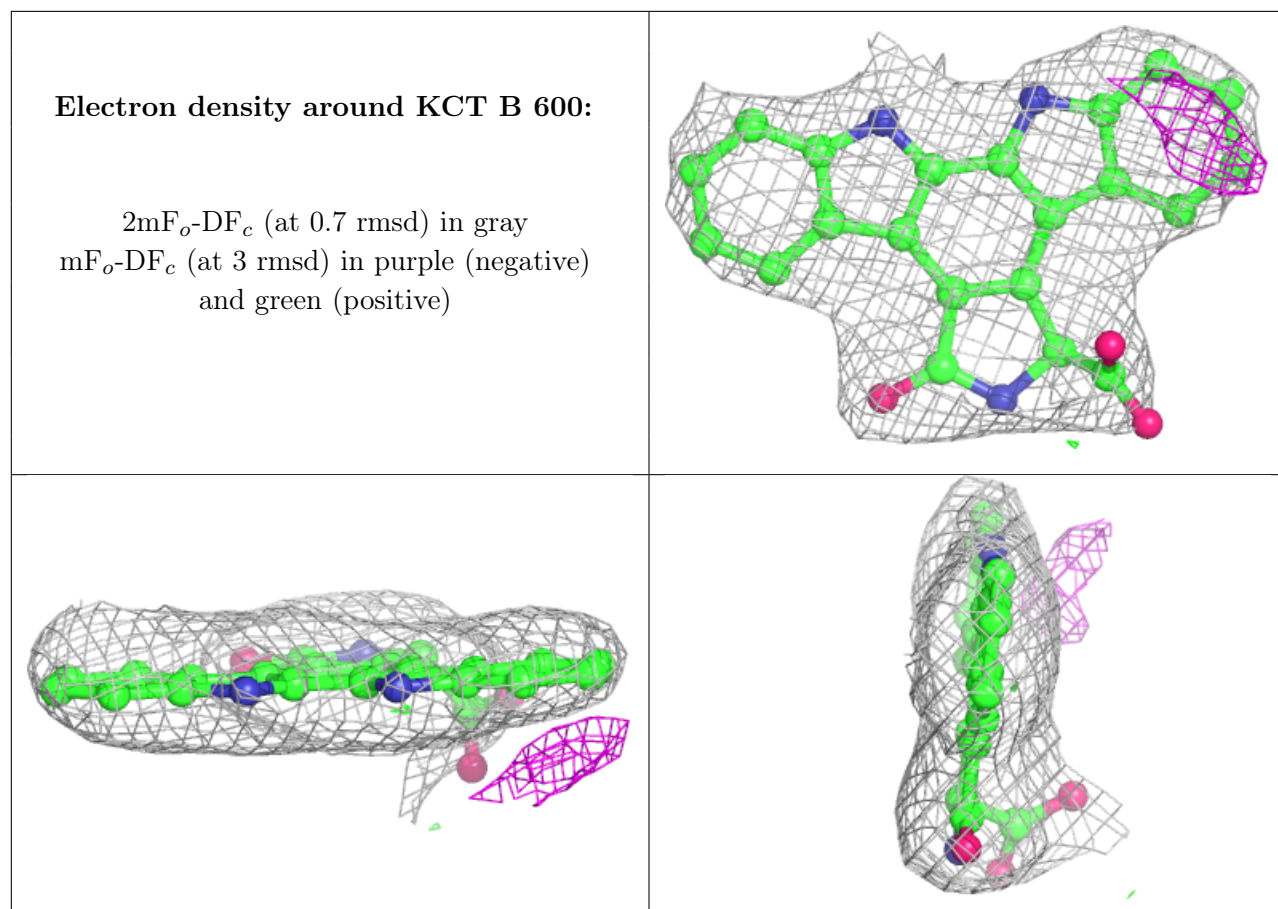
6.4 Ligands [i](#)

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
2	KCT	A	600	27/27	0.93	0.20	69,71,80,84	0
2	KCT	B	600	27/27	0.93	0.26	59,64,75,80	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.