



# Full wwPDB X-ray Structure Validation Report ⓘ

May 25, 2020 – 04:31 am BST

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 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 : 2.11  
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) : 2.11

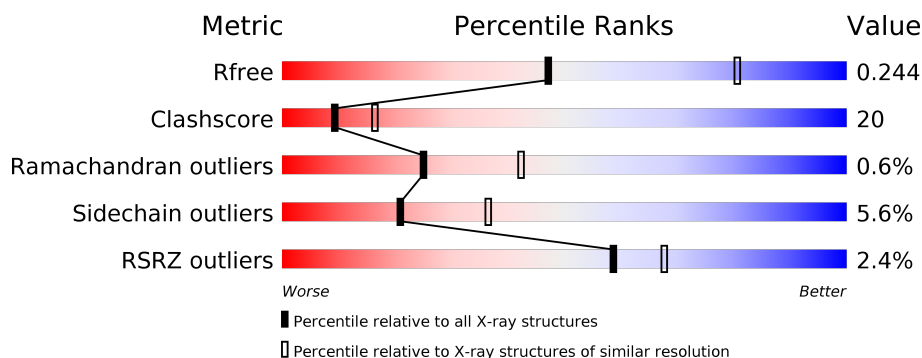
# 1 Overall quality at a glance

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 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	549	<div> <div>3%</div> <div> <div></div> <div>57%</div> <div>33%</div> <div>• 6%</div> </div> </div>
1	B	549	<div> <div>%</div> <div> <div></div> <div>61%</div> <div>30%</div> <div>• 7%</div> </div> </div>

## 2 Entry composition

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-monooxygenase.

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

There are 60 discrepancies between the modelled and reference sequences:

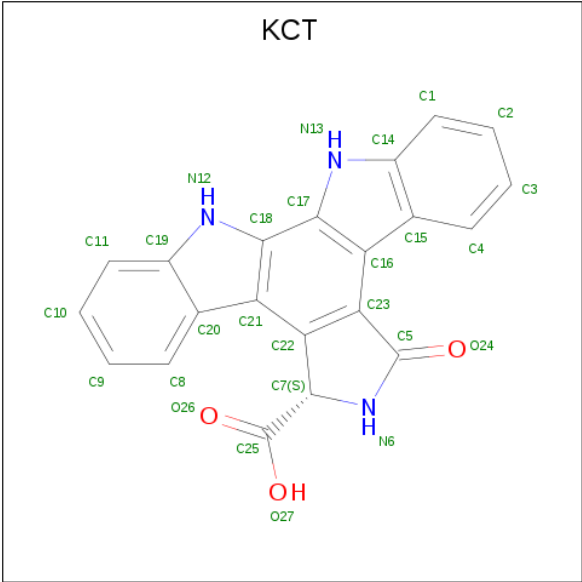
Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	EXPRESSION TAG	UNP Q8KI25
A	-18	GLY	-	EXPRESSION TAG	UNP Q8KI25
A	-17	SER	-	EXPRESSION TAG	UNP Q8KI25
A	-16	SER	-	EXPRESSION TAG	UNP Q8KI25
A	-15	HIS	-	EXPRESSION TAG	UNP Q8KI25
A	-14	HIS	-	EXPRESSION TAG	UNP Q8KI25
A	-13	HIS	-	EXPRESSION TAG	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: C<sub>21</sub>H<sub>13</sub>N<sub>3</sub>O<sub>3</sub>).



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

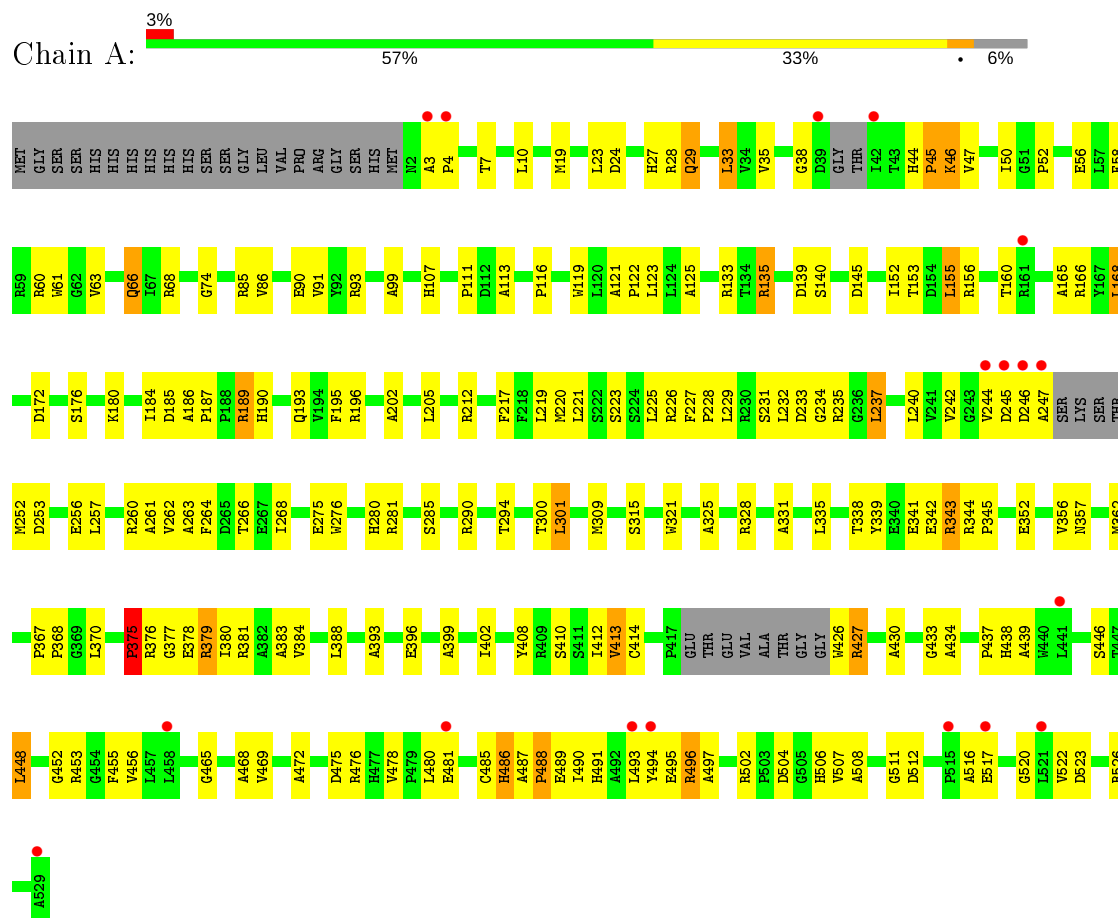
- Molecule 3 is water.

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

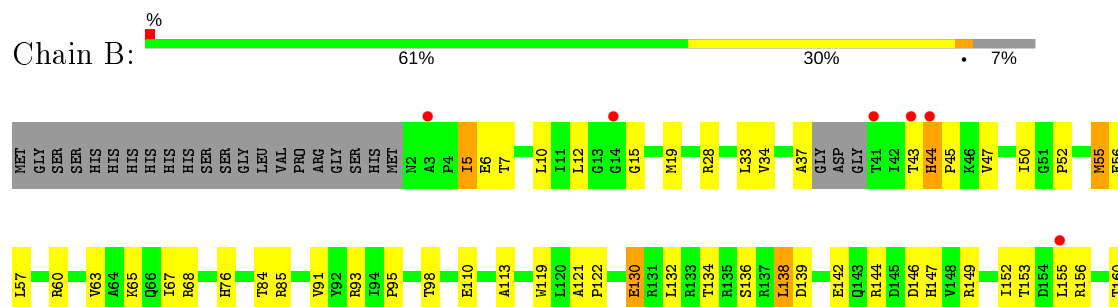
### 3 Residue-property plots [i](#)

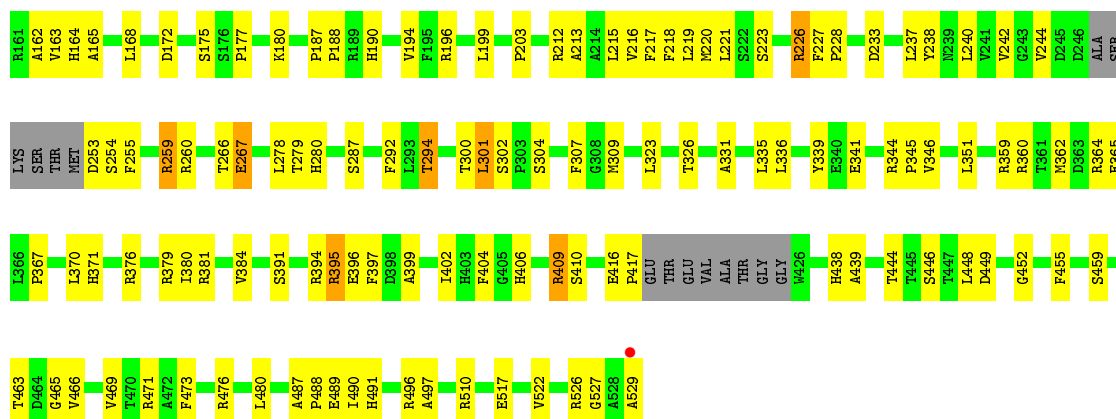
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	63.18 Å 77.70 Å 123.14 Å 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
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.38 (at 2.77 Å)	Xtriage
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 <sup>2</sup> )	62.9	Xtriage
Anisotropy	0.172	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 41.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	7928	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	61.0	wwPDB-VP

Xtriage'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.*

<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: 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.

All (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
1	A	496	ARG	N-CA-C	-5.31	96.66	111.00
1	A	140	SER	N-CA-CB	-5.29	102.57	110.50
1	A	139	ASP	CB-CA-C	5.15	120.69	110.40
1	A	486	HIS	N-CA-C	5.12	124.81	111.00
1	A	413	VAL	N-CA-C	-5.11	97.21	111.00

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	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
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.

All (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
1:B:44:HIS:HB2	1:B:45:PRO:HD2	1.40	0.99
1:A:46:LYS:O	1:A:47:VAL:CG2	2.13	0.96
1:B:44:HIS:CB	1:B:45:PRO:HD2	1.93	0.94
1:B:43:THR:O	1:B:44:HIS:HB2	1.67	0.94
1:A:489:GLU:HG2	1:A:490:ILE:HD12	1.50	0.93
1:A:315:SER:HA	1:A:343:ARG:HD2	1.50	0.92
1:A:46:LYS:O	1:A:47:VAL:HG22	1.68	0.91
1:A:496:ARG:HD2	1:A:511:GLY:HA2	1.53	0.91
1:A:86:VAL:HG21	1:A:229:LEU:HD13	1.54	0.89
1:B:194:VAL:O	1:B:244:VAL:HG22	1.72	0.86
1:A:219:LEU:O	1:A:226:ARG:HA	1.76	0.84
1:A:46:LYS:C	1:A:47:VAL:CG2	2.45	0.83
1:B:522:VAL:O	1:B:526:ARG:HG2	1.79	0.82
1:A:344:ARG:HB3	1:A:345:PRO:HD3	1.62	0.81
1:B:130:GLU:H	1:B:130:GLU:CD	1.85	0.79
1:A:66:GLN:H	1:A:66:GLN:NE2	1.80	0.79
1:A:10:LEU:HD23	1:A:168:LEU:HD12	1.66	0.78
1:A:135:ARG:HD2	1:A:156:ARG:HE	1.49	0.77
1:A:52:PRO:O	1:A:56:GLU:HG3	1.85	0.77
1:B:409:ARG:HA	1:B:409:ARG:HE	1.51	0.76
1:A:121:ALA:HB3	1:A:122:PRO:HD3	1.69	0.74
1:B:365:GLU:O	1:B:367:PRO:HD3	1.85	0.74
1:B:175:SER:O	1:B:177:PRO:HD3	1.87	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:189:ARG:HH11	1:A:189:ARG:HG2	1.53	0.73
1:A:437:PRO:HG2	1:A:494:TYR:CE1	2.25	0.72
1:B:138:LEU:HD12	1:B:152:ILE:HD11	1.70	0.72
1:B:5:ILE:HG22	1:B:162:ALA:O	1.90	0.72
1:B:52:PRO:O	1:B:56:GLU:HG3	1.89	0.72
1:B:194:VAL:HB	1:B:244:VAL:CG2	2.20	0.72
1:A:66:GLN:N	1:A:66:GLN:HE21	1.88	0.71
1:A:135:ARG:HG3	1:A:135:ARG:HH11	1.56	0.70
1:A:66:GLN:H	1:A:66:GLN:HE21	1.37	0.70
1:B:91:VAL:HG21	1:B:220:MET:CE	2.21	0.70
1:B:144:ARG:NH2	1:B:149:ARG:HD2	2.06	0.70
1:A:24:ASP:HB2	1:A:61:TRP:CZ2	2.27	0.70
1:A:156:ARG:HG2	1:A:156:ARG:HH11	1.58	0.69
1:A:275:GLU:HG2	1:A:276:TRP:N	2.06	0.69
1:B:60:ARG:HD2	1:B:452:GLY:HA3	1.74	0.69
1:A:46:LYS:O	1:A:47:VAL:HG23	1.90	0.68
1:A:217:PHE:O	1:A:228:PRO:HA	1.94	0.68
1:B:130:GLU:N	1:B:130:GLU:CD	2.47	0.67
1:B:476:ARG:HH11	1:B:476:ARG:HG3	1.60	0.67
1:A:244:VAL:HG12	1:A:245:ASP:O	1.94	0.67
1:B:399:ALA:O	1:B:402:ILE:HG22	1.96	0.66
1:B:43:THR:O	1:B:44:HIS:CB	2.40	0.66
1:A:91:VAL:HG21	1:A:220:MET:CE	2.25	0.66
1:A:502:ARG:HG3	1:A:506:HIS:O	1.96	0.66
1:B:28:ARG:HH11	1:B:28:ARG:HG2	1.60	0.66
1:A:399:ALA:HB1	1:A:402:ILE:CD1	2.09	0.66
1:A:487:ALA:HB1	1:A:489:GLU:OE1	1.96	0.65
1:B:394:ARG:HB3	1:B:394:ARG:NH1	2.12	0.65
1:B:5:ILE:C	1:B:5:ILE:HD13	2.17	0.65
1:B:227:PHE:HB3	1:B:228:PRO:HD2	1.79	0.64
1:A:60:ARG:HD3	1:A:452:GLY:HA3	1.80	0.64
1:B:212:ARG:NH1	1:B:212:ARG:HB3	2.11	0.64
1:A:522:VAL:O	1:A:526:ARG:HG2	1.97	0.64
1:A:91:VAL:HG21	1:A:220:MET:HE2	1.80	0.64
1:A:375:PRO:HA	1:A:378:GLU:HB2	1.78	0.64
1:B:5:ILE:HG23	1:B:163:VAL:HA	1.79	0.64
1:A:227:PHE:HB3	1:A:228:PRO:HD2	1.80	0.63
1:B:344:ARG:HB3	1:B:345:PRO:HD3	1.78	0.63
1:A:47:VAL:HA	2:A:600:KCT:O27	1.98	0.63
1:B:489:GLU:N	1:B:489:GLU:OE1	2.31	0.62
1:A:221:LEU:O	1:A:221:LEU:HD13	1.99	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:262:VAL:HG12	1:A:264:PHE:H	1.64	0.61
1:B:199:LEU:HD11	1:B:237:LEU:HD22	1.81	0.61
1:A:10:LEU:HD23	1:A:168:LEU:CD1	2.30	0.61
1:A:184:ILE:HD13	1:A:285:SER:HB2	1.84	0.60
1:A:344:ARG:NH2	3:A:736:HOH:O	2.34	0.60
1:A:135:ARG:O	1:A:155:LEU:HB2	2.02	0.60
1:A:388:LEU:CD2	1:A:393:ALA:HB2	2.32	0.60
1:A:33:LEU:HD21	1:A:133:ARG:HG2	1.84	0.59
1:A:402:ILE:N	1:A:402:ILE:HD12	2.18	0.59
1:A:315:SER:CA	1:A:343:ARG:HD2	2.27	0.59
1:B:364:ARG:HG2	1:B:364:ARG:HH11	1.67	0.59
1:A:153:THR:OG1	1:A:160:THR:HG22	2.03	0.58
1:B:370:LEU:HD22	1:B:381:ARG:HG2	1.85	0.58
1:A:189:ARG:NH1	3:A:735:HOH:O	2.37	0.58
1:B:5:ILE:HD13	1:B:5:ILE:O	2.03	0.58
1:B:12:LEU:HD22	1:B:172:ASP:HB3	1.86	0.58
1:A:294:THR:HG22	1:A:339:TYR:OH	2.04	0.57
1:B:144:ARG:HH22	1:B:149:ARG:HD2	1.68	0.57
1:A:35:VAL:HG11	1:A:152:ILE:HG12	1.87	0.57
1:B:476:ARG:HG3	1:B:476:ARG:NH1	2.20	0.57
1:B:394:ARG:CB	1:B:394:ARG:NH1	2.67	0.57
1:A:488:PRO:N	1:A:489:GLU:OE1	2.37	0.57
1:B:446:SER:O	1:B:449:ASP:HB2	2.05	0.57
1:B:394:ARG:CB	1:B:394:ARG:HH11	2.18	0.57
1:A:517:GLU:HG2	1:A:520:GLY:H	1.69	0.57
1:B:255:PHE:CE1	1:B:259:ARG:HD2	2.39	0.56
1:B:190:HIS:HB2	1:B:280:HIS:CG	2.40	0.56
1:A:220:MET:HE3	1:A:370:LEU:HD21	1.88	0.56
1:A:376:ARG:O	1:A:380:ILE:HG13	2.05	0.56
1:A:107:HIS:HB2	3:A:726:HOH:O	2.04	0.56
1:B:360:ARG:NE	1:B:391:SER:O	2.38	0.56
1:A:489:GLU:OE1	1:A:489:GLU:N	2.39	0.56
1:B:294:THR:HG23	1:B:339:TYR:OH	2.05	0.56
1:B:84:THR:O	1:B:221:LEU:HG	2.05	0.56
1:A:68:ARG:NH2	1:A:111:PRO:O	2.33	0.56
1:A:185:ASP:OD2	1:A:187:PRO:HD3	2.05	0.56
1:A:455:PHE:O	1:A:480:LEU:HD12	2.06	0.56
1:B:341:GLU:OE2	1:B:410:SER:HB2	2.06	0.56
1:B:300:THR:C	1:B:301:LEU:HD23	2.26	0.55
1:B:6:GLU:OE2	1:B:164:HIS:ND1	2.39	0.55
1:A:28:ARG:CZ	1:A:321:TRP:HB3	2.36	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:7:THR:O	1:A:165:ALA:HA	2.07	0.55
1:A:223:SER:O	1:A:226:ARG:NH1	2.40	0.55
1:B:28:ARG:NH1	1:B:28:ARG:HG2	2.19	0.55
1:A:342:GLU:HB3	1:A:408:TYR:CD1	2.41	0.55
1:B:255:PHE:HE1	1:B:259:ARG:HD2	1.72	0.55
1:B:47:VAL:HG21	1:B:309:MET:SD	2.47	0.54
1:A:172:ASP:OD1	1:A:176:SER:OG	2.21	0.54
1:A:341:GLU:OE2	1:A:410:SER:HB2	2.06	0.54
1:B:194:VAL:HB	1:B:244:VAL:HG21	1.89	0.54
1:B:190:HIS:HB2	1:B:280:HIS:CD2	2.43	0.54
1:B:7:THR:O	1:B:165:ALA:HA	2.08	0.54
1:B:55:MET:HB3	1:B:110:GLU:HB3	1.90	0.53
1:B:144:ARG:CZ	1:B:149:ARG:HD2	2.38	0.53
1:B:226:ARG:NH2	1:B:362:MET:O	2.41	0.53
1:A:412:ILE:O	1:A:412:ILE:HG13	2.07	0.53
1:A:19:MET:HE3	1:A:125:ALA:HA	1.91	0.53
1:B:527:GLY:C	1:B:529:ALA:H	2.11	0.53
1:A:50:ILE:HG12	1:A:309:MET:CE	2.39	0.53
1:B:217:PHE:O	1:B:228:PRO:HA	2.09	0.53
1:A:352:GLU:O	1:A:356:VAL:HG23	2.09	0.52
1:A:427:ARG:HG2	1:A:427:ARG:HH21	1.75	0.52
1:A:3:ALA:HB1	1:A:4:PRO:HD2	1.91	0.52
1:A:375:PRO:C	1:A:377:GLY:N	2.62	0.52
1:B:219:LEU:O	1:B:226:ARG:HA	2.10	0.52
1:B:279:THR:OG1	1:B:300:THR:HG21	2.09	0.52
1:A:244:VAL:C	1:A:245:ASP:O	2.46	0.52
1:B:346:VAL:HG13	1:B:406:HIS:CD2	2.45	0.52
1:B:394:ARG:HB2	1:B:394:ARG:HH11	1.73	0.51
1:A:502:ARG:HB2	1:A:504:ASP:OD1	2.09	0.51
1:A:190:HIS:HB2	1:A:280:HIS:CG	2.45	0.51
1:B:93:ARG:O	1:B:95:PRO:HD3	2.11	0.51
1:B:91:VAL:HG21	1:B:220:MET:HE2	1.89	0.51
1:B:142:GLU:OE2	1:B:144:ARG:NH1	2.44	0.50
1:A:453:ARG:HB2	3:A:720:HOH:O	2.12	0.50
1:B:215:LEU:HG	1:B:216:VAL:HG23	1.92	0.50
1:B:359:ARG:HH11	1:B:359:ARG:HG3	1.76	0.50
1:A:196:ARG:HB2	1:A:242:VAL:CG1	2.42	0.50
1:A:328:ARG:HH11	1:A:328:ARG:HG3	1.76	0.50
1:A:456:VAL:HG22	1:A:481:GLU:HB2	1.92	0.50
1:B:395:ARG:CG	1:B:395:ARG:HH21	2.24	0.50
1:B:438:HIS:CG	1:B:439:ALA:N	2.80	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:219:LEU:O	1:A:226:ARG:CA	2.55	0.50
1:A:378:GLU:HA	1:A:381:ARG:NH1	2.26	0.50
1:A:502:ARG:HG2	1:A:508:ALA:HB2	1.94	0.49
1:A:375:PRO:O	1:A:377:GLY:N	2.45	0.49
1:A:202:ALA:HB1	1:A:205:LEU:HB2	1.95	0.49
1:B:473:PHE:CD1	1:B:480:LEU:HD22	2.47	0.49
1:A:257:LEU:O	1:A:260:ARG:HB2	2.13	0.49
1:B:370:LEU:CD2	1:B:381:ARG:HG2	2.42	0.49
1:A:497:ALA:HB3	1:A:512:ASP:HA	1.95	0.49
1:A:221:LEU:HD12	1:A:225:LEU:HD12	1.95	0.49
1:B:438:HIS:HA	1:B:448:LEU:HD13	1.95	0.49
1:A:27:HIS:HB3	3:A:705:HOH:O	2.11	0.49
1:A:465:GLY:O	1:A:468:ALA:HB3	2.13	0.48
1:A:375:PRO:C	1:A:377:GLY:H	2.15	0.48
1:B:331:ALA:HB1	1:B:335:LEU:HD22	1.93	0.48
1:A:189:ARG:HG2	1:A:189:ARG:NH1	2.22	0.48
1:A:38:GLY:HA3	1:A:135:ARG:HH12	1.79	0.48
1:B:68:ARG:NH1	1:B:113:ALA:HB2	2.28	0.48
1:B:119:TRP:O	1:B:122:PRO:HD2	2.14	0.48
1:A:50:ILE:HG12	1:A:309:MET:HE3	1.96	0.48
1:A:44:HIS:O	1:A:45:PRO:O	2.32	0.48
1:B:394:ARG:CZ	1:B:394:ARG:HB3	2.43	0.48
1:B:444:THR:HG22	1:B:444:THR:O	2.14	0.48
1:B:47:VAL:HG13	1:B:47:VAL:O	2.12	0.48
1:A:74:GLY:HA2	1:A:99:ALA:HB3	1.95	0.47
1:B:409:ARG:NE	1:B:409:ARG:HA	2.26	0.47
1:A:335:LEU:O	1:A:338:THR:OG1	2.24	0.47
1:A:472:ALA:O	1:A:475:ASP:HB2	2.13	0.47
1:B:37:ALA:HA	1:B:136:SER:O	2.13	0.47
1:B:144:ARG:HG3	1:B:147:HIS:NE2	2.29	0.47
1:A:156:ARG:NH1	1:A:156:ARG:HG2	2.27	0.47
1:A:266:THR:O	1:A:268:ILE:HG13	2.14	0.47
1:A:246:ASP:O	1:A:247:ALA:HB3	2.13	0.47
1:A:66:GLN:N	1:A:66:GLN:NE2	2.51	0.47
1:A:402:ILE:HD12	1:A:402:ILE:H	1.80	0.47
1:A:370:LEU:HD12	1:A:381:ARG:HG2	1.97	0.47
1:A:437:PRO:HB2	1:A:493:LEU:HD21	1.96	0.47
1:A:523:ASP:OD1	1:A:526:ARG:NH1	2.48	0.47
1:A:38:GLY:HA3	1:A:135:ARG:NH1	2.30	0.47
1:A:135:ARG:HH11	1:A:135:ARG:CG	2.25	0.46
1:B:138:LEU:HA	1:B:152:ILE:HD13	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:15:GLY:O	1:B:19:MET:HG2	2.16	0.46
1:B:459:SER:HA	1:B:497:ALA:O	2.14	0.46
1:A:193:GLN:HE22	1:A:362:MET:CE	2.28	0.46
1:A:339:TYR:O	1:A:343:ARG:HB2	2.15	0.46
1:A:344:ARG:HB3	1:A:345:PRO:CD	2.41	0.46
1:A:485:CYS:SG	1:A:486:HIS:N	2.89	0.46
1:A:119:TRP:O	1:A:122:PRO:HD2	2.16	0.46
1:A:68:ARG:NH1	1:A:113:ALA:HB2	2.31	0.46
1:B:260:ARG:HG3	1:B:260:ARG:HH11	1.80	0.46
1:B:346:VAL:HG13	1:B:406:HIS:NE2	2.31	0.46
1:A:488:PRO:HA	1:A:491:HIS:HB3	1.97	0.46
1:B:307:PHE:CE1	1:B:402:ILE:HG21	2.51	0.46
1:A:388:LEU:HD23	1:A:393:ALA:HB2	1.98	0.46
1:B:376:ARG:O	1:B:380:ILE:HG13	2.16	0.46
1:A:35:VAL:HG23	1:A:35:VAL:O	2.16	0.45
1:B:203:PRO:HD2	1:B:267:GLU:HG2	1.98	0.45
1:A:380:ILE:O	1:A:383:ALA:HB3	2.15	0.45
1:B:218:PHE:HE1	1:B:397:PHE:HZ	1.63	0.45
1:A:186:ALA:O	1:A:281:ARG:CZ	2.65	0.45
1:B:10:LEU:O	1:B:168:LEU:HD12	2.16	0.45
1:A:253:ASP:O	1:A:257:LEU:HG	2.17	0.45
1:B:85:ARG:HG2	1:B:221:LEU:HD21	1.99	0.45
1:B:489:GLU:HG2	1:B:490:ILE:N	2.32	0.45
1:A:116:PRO:HD2	1:A:119:TRP:CE3	2.51	0.45
1:B:279:THR:O	1:B:300:THR:HG22	2.17	0.45
1:B:91:VAL:HG21	1:B:220:MET:HE3	1.97	0.45
1:B:156:ARG:HB3	1:B:156:ARG:NH1	2.32	0.45
1:A:357:ASN:CG	1:A:396:GLU:HG3	2.37	0.45
1:B:121:ALA:HB3	1:B:122:PRO:HD3	1.98	0.45
1:A:325:ALA:HB1	1:A:331:ALA:HB2	1.98	0.45
1:A:135:ARG:NH1	1:A:135:ARG:HG3	2.27	0.44
1:A:233:ASP:O	1:A:235:ARG:HG3	2.17	0.44
1:A:375:PRO:O	1:A:376:ARG:C	2.55	0.44
1:A:379:ARG:HD3	1:A:379:ARG:HA	1.76	0.44
1:B:371:HIS:CD2	1:B:371:HIS:N	2.85	0.44
1:B:63:VAL:O	1:B:67:ILE:HG13	2.18	0.44
1:A:28:ARG:NH2	1:A:321:TRP:HB3	2.32	0.44
1:B:359:ARG:HG3	1:B:359:ARG:NH1	2.33	0.44
1:A:300:THR:C	1:A:301:LEU:HD23	2.38	0.44
1:B:119:TRP:CH2	1:B:233:ASP:HB2	2.53	0.44
1:B:196:ARG:HG3	1:B:244:VAL:CG1	2.48	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:412:ILE:O	1:A:413:VAL:HG22	2.17	0.44
1:B:138:LEU:HD12	1:B:152:ILE:CD1	2.45	0.44
1:A:172:ASP:OD1	1:A:172:ASP:N	2.45	0.44
1:B:199:LEU:HD12	1:B:238:TYR:O	2.17	0.44
1:A:195:PHE:HB2	1:A:276:TRP:CE2	2.53	0.43
1:A:413:VAL:HG11	1:A:433:GLY:CA	2.48	0.43
1:A:58:PHE:HB3	1:A:63:VAL:HG23	1.99	0.43
1:B:5:ILE:CG2	1:B:162:ALA:O	2.62	0.43
1:B:278:LEU:HD21	1:B:302:SER:HB2	2.00	0.43
1:A:438:HIS:ND1	1:A:446:SER:OG	2.48	0.43
1:A:85:ARG:CZ	1:A:263:ALA:HA	2.47	0.43
1:B:144:ARG:NH1	1:B:149:ARG:HD2	2.33	0.43
1:B:212:ARG:HH11	1:B:212:ARG:HB3	1.81	0.43
1:A:495:GLU:HB3	1:A:496:ARG:HH12	1.84	0.43
1:A:29:GLN:HB2	1:A:328:ARG:HH22	1.84	0.43
1:B:93:ARG:NH2	1:B:213:ALA:O	2.51	0.43
1:A:152:ILE:N	1:A:152:ILE:HD12	2.33	0.43
1:B:156:ARG:HH11	1:B:156:ARG:HB3	1.84	0.43
1:B:416:GLU:HB3	1:B:417:PRO:HD2	2.00	0.43
1:B:146:ASP:N	1:B:146:ASP:OD1	2.50	0.43
1:A:487:ALA:HB1	1:A:489:GLU:CD	2.38	0.43
1:B:446:SER:O	1:B:449:ASP:N	2.45	0.43
1:B:180:LYS:HA	1:B:180:LYS:HD2	1.72	0.42
1:B:50:ILE:HD12	1:B:67:ILE:HG21	2.01	0.42
1:A:378:GLU:OE1	1:A:381:ARG:NH1	2.48	0.42
1:B:76:HIS:O	1:B:98:THR:HG22	2.19	0.42
1:B:212:ARG:CB	1:B:212:ARG:HH11	2.32	0.42
1:A:237:LEU:HD23	1:A:237:LEU:HA	1.80	0.42
1:A:252:MET:N	1:A:256:GLU:OE1	2.53	0.42
1:A:90:GLU:CD	1:A:93:ARG:HE	2.22	0.42
1:B:346:VAL:HG22	1:B:406:HIS:CD2	2.54	0.42
1:B:463:THR:O	1:B:466:VAL:HG23	2.20	0.42
1:A:232:LEU:HD12	1:A:237:LEU:HD13	2.02	0.42
1:A:23:LEU:O	1:A:27:HIS:HD2	2.03	0.42
1:B:226:ARG:NH2	1:B:362:MET:HA	2.34	0.42
1:B:381:ARG:HA	1:B:384:VAL:HG22	2.02	0.42
1:A:301:LEU:HD23	1:A:301:LEU:N	2.35	0.41
1:A:378:GLU:HA	1:A:381:ARG:HH12	1.85	0.41
1:A:219:LEU:HD11	1:A:261:ALA:HB1	2.01	0.41
1:A:219:LEU:HA	1:A:219:LEU:HD23	1.79	0.41
1:A:28:ARG:C	1:A:29:GLN:HG2	2.40	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:412:ILE:C	1:A:413:VAL:CG2	2.88	0.41
1:B:496:ARG:HG3	1:B:510:ARG:HD2	2.03	0.41
1:B:34:VAL:O	1:B:132:LEU:HD12	2.20	0.41
1:B:402:ILE:HA	1:B:402:ILE:HD12	1.92	0.41
1:A:196:ARG:HG3	1:A:244:VAL:CG2	2.50	0.41
1:B:187:PRO:HA	1:B:188:PRO:HD3	1.77	0.41
1:B:455:PHE:O	1:B:480:LEU:HA	2.19	0.41
1:B:487:ALA:HA	1:B:488:PRO:HD2	1.85	0.41
1:B:266:THR:OG1	1:B:267:GLU:N	2.54	0.41
1:B:351:LEU:C	1:B:351:LEU:HD23	2.41	0.41
1:A:135:ARG:NH1	1:A:135:ARG:CG	2.81	0.41
1:A:384:VAL:O	1:A:388:LEU:HB2	2.20	0.41
1:A:448:LEU:HD12	1:A:448:LEU:HA	1.86	0.41
1:B:326:THR:HA	1:B:331:ALA:HB3	2.03	0.41
1:B:395:ARG:CG	1:B:395:ARG:NH2	2.82	0.41
1:B:465:GLY:O	1:B:469:VAL:HG23	2.19	0.41
1:A:46:LYS:HA	1:A:46:LYS:HD2	1.83	0.41
1:A:430:ALA:O	1:A:507:VAL:HG21	2.20	0.41
1:A:231:SER:OG	1:A:234:GLY:HA2	2.21	0.41
1:B:196:ARG:HB2	1:B:242:VAL:HG22	2.03	0.41
1:B:57:LEU:HA	1:B:57:LEU:HD23	1.80	0.41
1:B:153:THR:HA	1:B:160:THR:HG22	2.03	0.41
1:A:328:ARG:HH11	1:A:328:ARG:CG	2.32	0.40
1:A:367:PRO:HD2	1:A:370:LEU:HD22	2.03	0.40
1:A:434:ALA:O	1:A:506:HIS:HA	2.21	0.40
1:B:323:LEU:HD23	1:B:336:LEU:HD21	2.03	0.40
1:A:166:ARG:O	1:A:290:ARG:NH1	2.51	0.40
1:B:52:PRO:HB2	1:B:404:PHE:HZ	1.87	0.40
1:A:476:ARG:HB2	1:A:478:VAL:HG23	2.03	0.40
1:A:439:ALA:HB1	1:A:490:ILE:HG13	2.04	0.40
1:B:223:SER:O	1:B:226:ARG:HD3	2.22	0.40
1:B:287:SER:HB3	1:B:292:PHE:CD2	2.56	0.40
1:A:24:ASP:HB2	1:A:61:TRP:HZ2	1.80	0.40
1:B:396:GLU:O	2:B:600:KCT:H10	2.21	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	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

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	44	HIS
1	A	155	LEU
1	A	45	PRO
1	B	304	SER
1	A	368	PRO
1	A	488	PRO

### 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	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

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	GLN
1	A	33	LEU
1	A	66	GLN
1	A	123	LEU
1	A	135	ARG
1	A	145	ASP
1	A	168	LEU
1	A	180	LYS
1	A	189	ARG
1	A	212	ARG
1	A	237	LEU
1	A	240	LEU
1	A	301	LEU
1	A	343	ARG
1	A	375	PRO
1	A	379	ARG
1	A	414	CYS
1	A	426	TRP
1	A	427	ARG
1	A	448	LEU
1	A	469	VAL
1	B	5	ILE
1	B	33	LEU
1	B	55	MET
1	B	65	LYS
1	B	130	GLU
1	B	134	THR
1	B	138	LEU
1	B	139	ASP
1	B	155	LEU
1	B	226	ARG
1	B	240	LEU
1	B	253	ASP
1	B	254	SER
1	B	259	ARG
1	B	267	GLU
1	B	294	THR
1	B	301	LEU
1	B	379	ARG
1	B	395	ARG
1	B	409	ARG
1	B	471	ARG
1	B	491	HIS

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Mol	Chain	Res	Type
1	B	517	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	27	HIS
1	A	66	GLN
1	A	193	GLN
1	B	193	GLN
1	B	280	HIS
1	B	371	HIS
1	B	406	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 carbohydrates 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	A	600	-	22,32,32	1.94	8 (36%)	27,50,50	2.41	8 (29%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	KCT	B	600	-	22,32,32	2.01	7 (31%)	27,50,50	2.64	9 (33%)

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

All (15) 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	-3.07	1.47	1.52
2	B	600	KCT	C3-C4	2.79	1.43	1.36
2	A	600	KCT	O24-C5	2.71	1.28	1.23
2	A	600	KCT	C3-C4	2.55	1.42	1.36
2	A	600	KCT	C2-C1	2.54	1.42	1.36
2	B	600	KCT	C9-C8	2.52	1.42	1.36
2	B	600	KCT	C10-C11	2.50	1.42	1.36
2	A	600	KCT	C10-C11	2.48	1.42	1.36
2	B	600	KCT	C2-C1	2.41	1.42	1.36
2	A	600	KCT	C7-N6	2.27	1.48	1.46
2	A	600	KCT	C9-C8	2.22	1.41	1.36
2	B	600	KCT	C22-C7	-2.18	1.48	1.52

All (17) 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	C22-C21-C18	-3.59	119.25	122.71
2	B	600	KCT	C22-C21-C18	-3.59	119.25	122.71
2	A	600	KCT	C4-C15-C14	3.40	122.68	118.17
2	B	600	KCT	C4-C15-C14	3.11	122.29	118.17
2	A	600	KCT	C22-C23-C5	-2.73	105.68	108.21
2	B	600	KCT	C22-C23-C5	-2.66	105.74	108.21
2	B	600	KCT	O24-C5-C23	-2.59	126.25	129.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	600	KCT	C8-C20-C19	2.52	121.51	118.17
2	A	600	KCT	O24-C5-C23	-2.51	126.34	129.32
2	B	600	KCT	C21-C20-C19	2.45	108.74	106.09
2	A	600	KCT	C21-C20-C19	2.41	108.69	106.09
2	A	600	KCT	O24-C5-N6	-2.31	123.98	125.98
2	B	600	KCT	C8-C20-C19	2.27	121.18	118.17
2	B	600	KCT	O24-C5-N6	-2.16	124.10	125.98
2	B	600	KCT	C22-C23-C16	2.16	123.39	120.47

There are no chirality outliers.

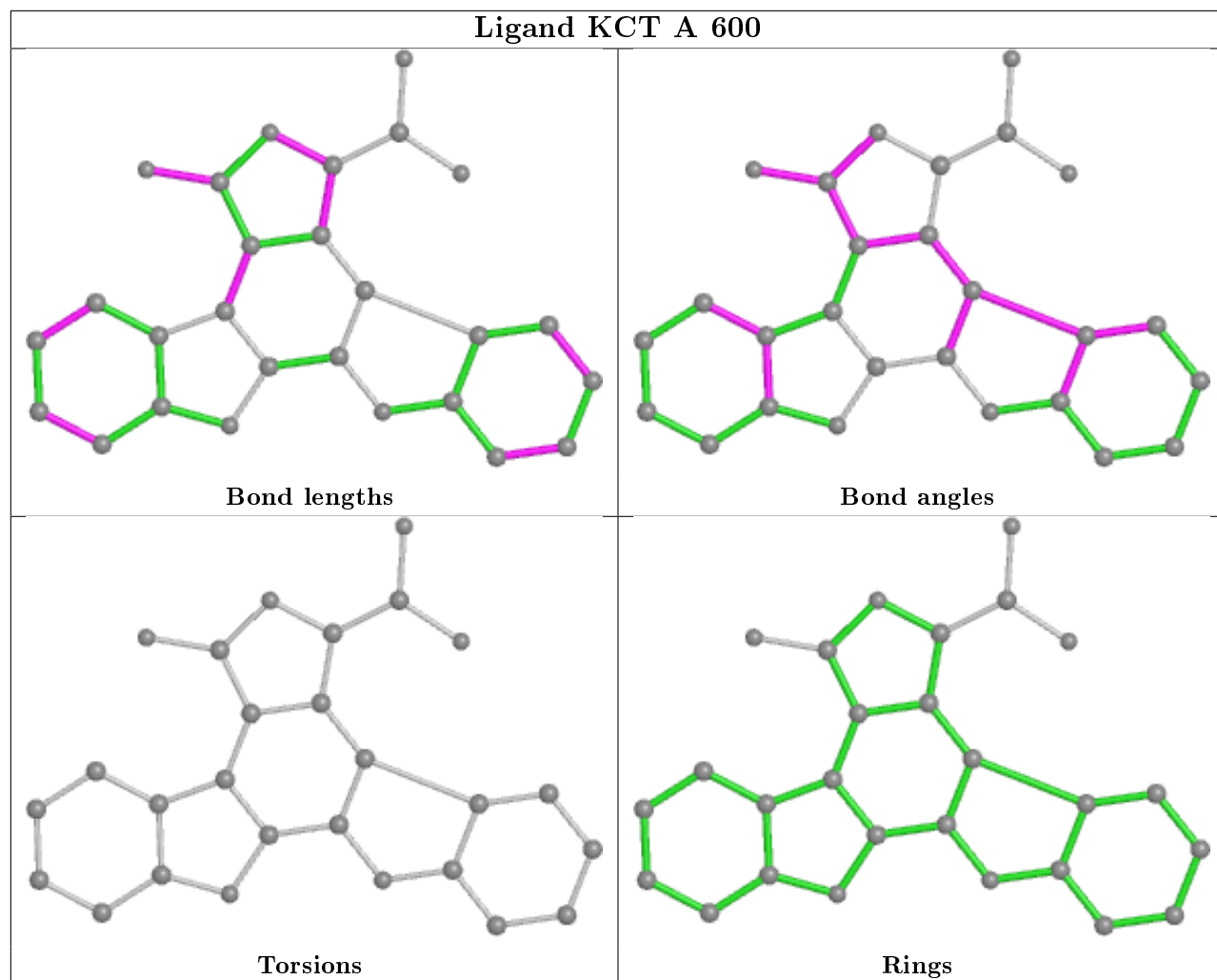
There are no torsion outliers.

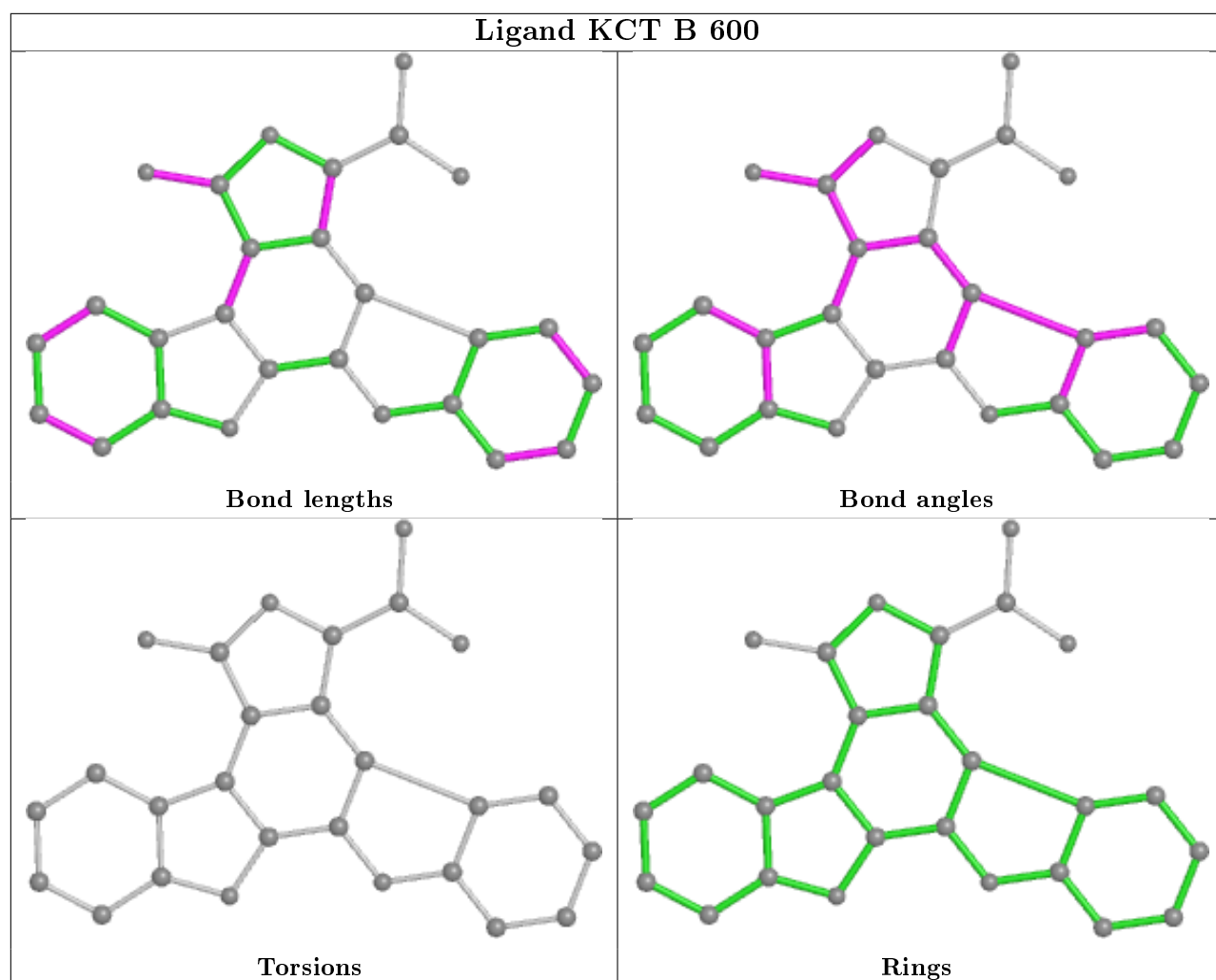
There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	600	KCT	1	0
2	B	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 ⓘ

### 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, 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	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

All (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
1	A	245	ASP	3.3
1	A	517	GLU	3.1
1	A	4	PRO	2.9
1	B	529	ALA	2.9
1	A	42	ILE	2.9
1	B	43	THR	2.7
1	A	161	ARG	2.7
1	B	3	ALA	2.6
1	A	244	VAL	2.6
1	B	155	LEU	2.5
1	A	458	LEU	2.5
1	B	14	GLY	2.4
1	A	521	LEU	2.4
1	B	41	THR	2.3
1	A	247	ALA	2.3
1	A	515	PRO	2.2
1	A	481	GLU	2.1
1	A	441	LEU	2.1
1	A	39	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	493	LEU	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 carbohydrates in this entry.

## 6.4 Ligands [i](#)

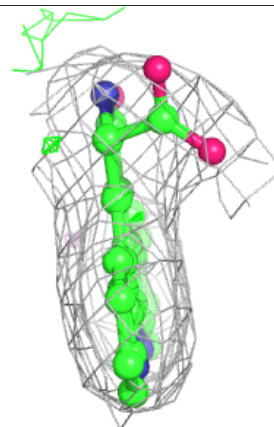
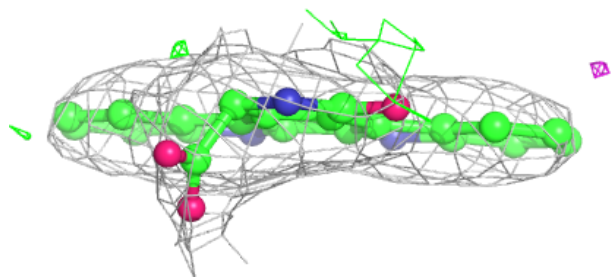
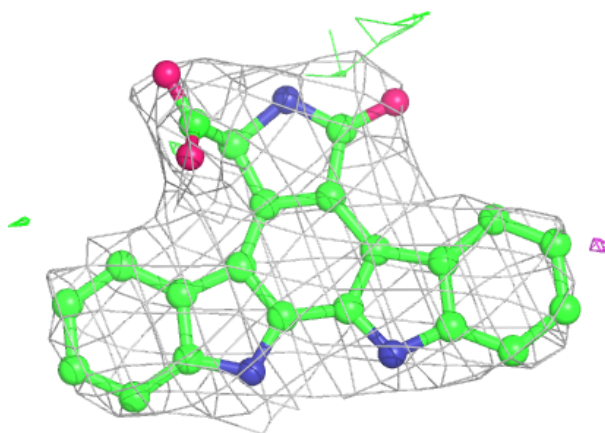
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

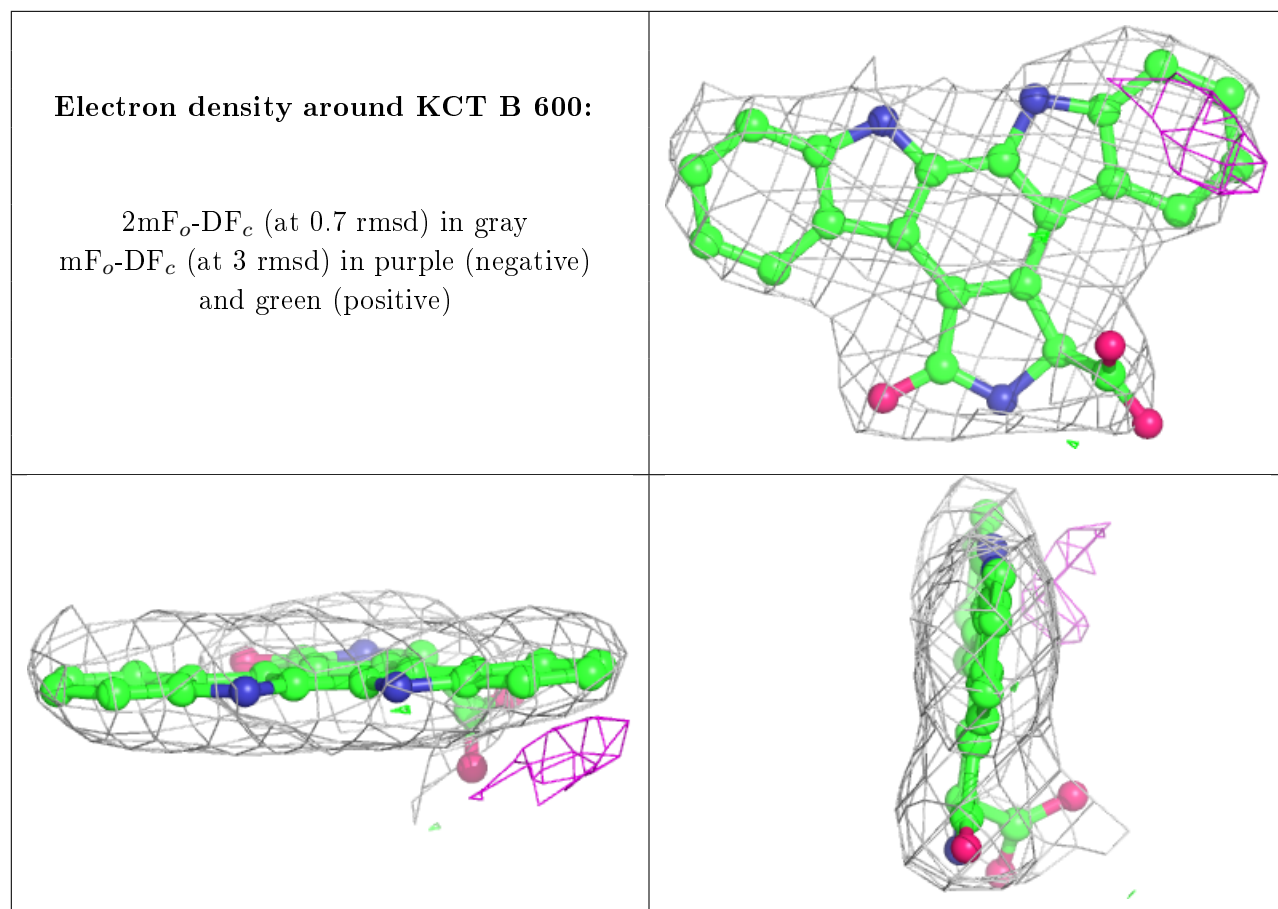
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	KCT	A	600	27/27	0.93	0.20	69,71,80,84	0
2	KCT	B	600	27/27	0.93	0.26	59,64,75,80	0

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.

**Electron density around KCT A 600:**

$2mF_o-DF_c$  (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.