



Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 09:15 am BST

PDB ID : 3UDL
Title : 3-heterocyclyl quinolone bound to HCV NS5B
Authors : Somoza, J.R.
Deposited on : 2011-10-28
Resolution : 2.17 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 : 2.11
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

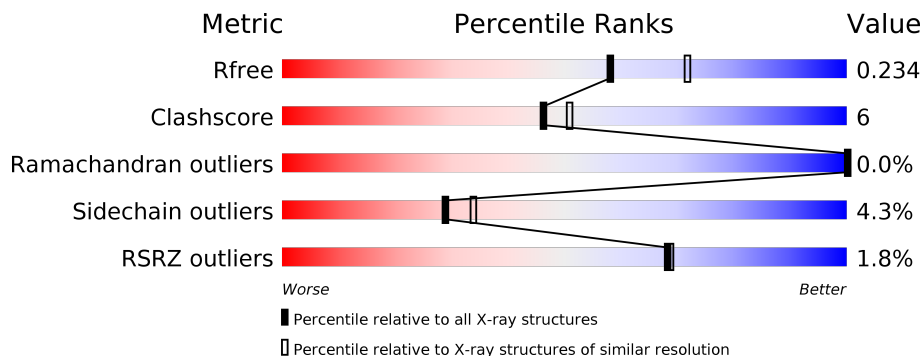
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.17 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	6864 (2.20-2.16)
Clashscore	141614	7689 (2.20-2.16)
Ramachandran outliers	138981	7564 (2.20-2.16)
Sidechain outliers	138945	7564 (2.20-2.16)
RSRZ outliers	127900	6738 (2.20-2.16)

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 ≥ 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	576	
1	B	576	
1	C	576	
1	D	576	

2 Entry composition i

There are 3 unique types of molecules in this entry. The entry contains 18412 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 HCV NS5B polymerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	555	4319	2725	761	802	31	0	0	0
1	B	555	4319	2725	761	802	31	0	0	0
1	C	555	4319	2725	761	802	31	0	0	0
1	D	555	4319	2725	761	802	31	0	0	0

There are 24 discrepancies between the modelled and reference sequences:

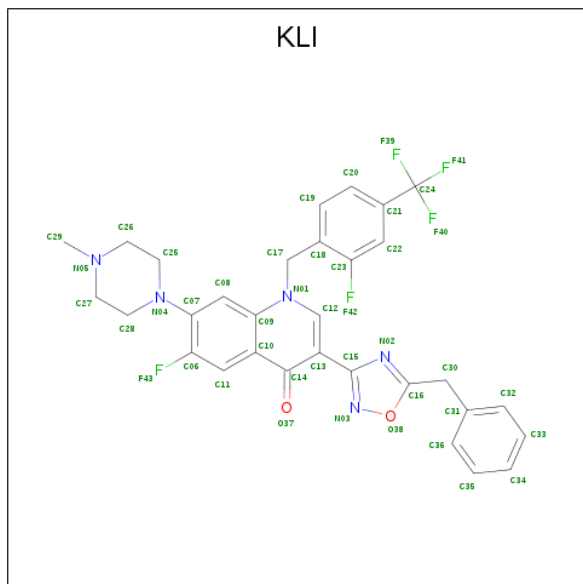
Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	HIS	-	EXPRESSION TAG	UNP Q99AU2
A	-4	HIS	-	EXPRESSION TAG	UNP Q99AU2
A	-3	HIS	-	EXPRESSION TAG	UNP Q99AU2
A	-2	HIS	-	EXPRESSION TAG	UNP Q99AU2
A	-1	HIS	-	EXPRESSION TAG	UNP Q99AU2
A	0	HIS	-	EXPRESSION TAG	UNP Q99AU2
B	-5	HIS	-	EXPRESSION TAG	UNP Q99AU2
B	-4	HIS	-	EXPRESSION TAG	UNP Q99AU2
B	-3	HIS	-	EXPRESSION TAG	UNP Q99AU2
B	-2	HIS	-	EXPRESSION TAG	UNP Q99AU2
B	-1	HIS	-	EXPRESSION TAG	UNP Q99AU2
B	0	HIS	-	EXPRESSION TAG	UNP Q99AU2
C	-5	HIS	-	EXPRESSION TAG	UNP Q99AU2
C	-4	HIS	-	EXPRESSION TAG	UNP Q99AU2
C	-3	HIS	-	EXPRESSION TAG	UNP Q99AU2
C	-2	HIS	-	EXPRESSION TAG	UNP Q99AU2
C	-1	HIS	-	EXPRESSION TAG	UNP Q99AU2
C	0	HIS	-	EXPRESSION TAG	UNP Q99AU2
D	-5	HIS	-	EXPRESSION TAG	UNP Q99AU2
D	-4	HIS	-	EXPRESSION TAG	UNP Q99AU2
D	-3	HIS	-	EXPRESSION TAG	UNP Q99AU2

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-2	HIS	-	EXPRESSION TAG	UNP Q99AU2
D	-1	HIS	-	EXPRESSION TAG	UNP Q99AU2
D	0	HIS	-	EXPRESSION TAG	UNP Q99AU2

- Molecule 2 is 3-(5-benzyl-1,2,4-oxadiazol-3-yl)-6-fluoro-1-[2-fluoro-4-(trifluoromethyl)benzyl]-7-(4-methylpiperazin-1-yl)quinolin-4(1H)-one (three-letter code: KLI) (formula: C₃₁H₂₆F₅N₅O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
			Total	C	F	N			O
2	A	1	Total	C	F	N	O	0	0
			43	31	5	5	2		
2	B	1	Total	C	F	N	O	0	0
			43	31	5	5	2		
2	C	1	Total	C	F	N	O	0	0
			43	31	5	5	2		
2	D	1	Total	C	F	N	O	0	0
			43	31	5	5	2		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	274	Total	O	0	0
			274	274		
3	B	248	Total	O	0	0
			248	248		
3	C	231	Total	O	0	0
			231	231		

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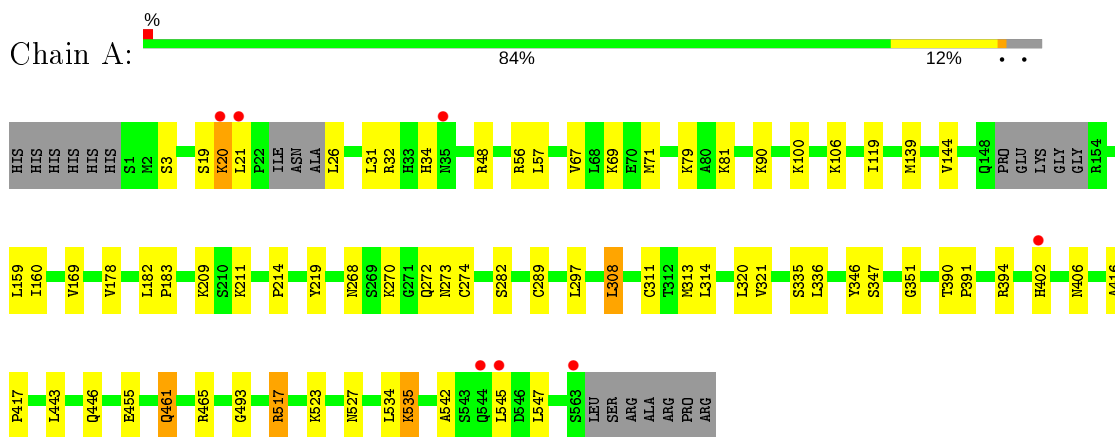
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	211	Total 211	O 211	0	0

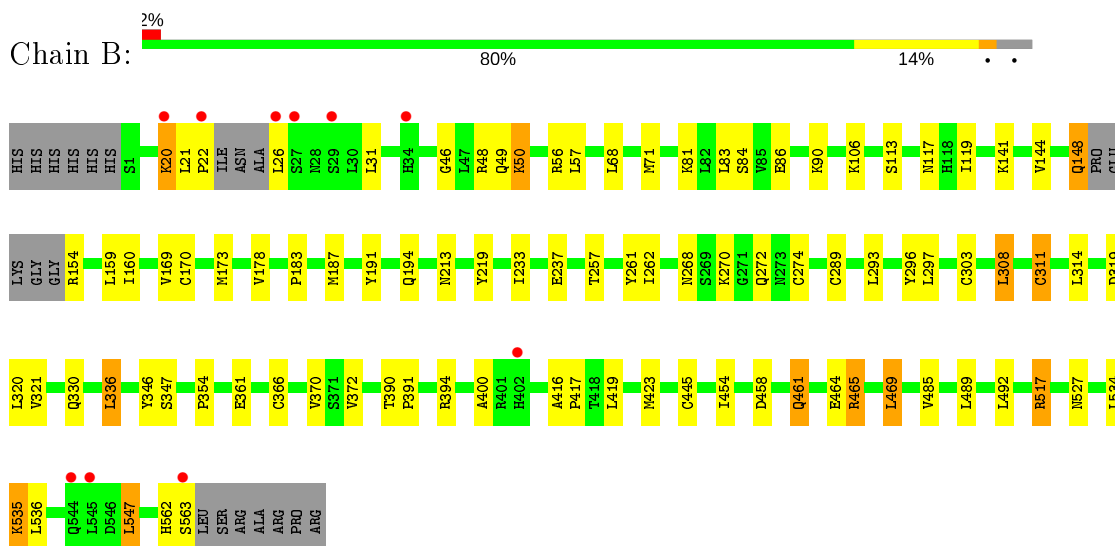
3 Residue-property plots i

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.

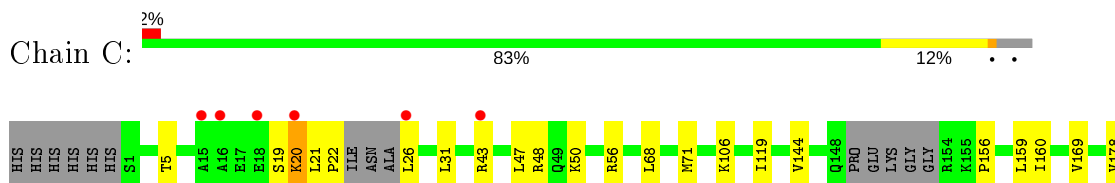
- Molecule 1: HCV NS5B polymerase

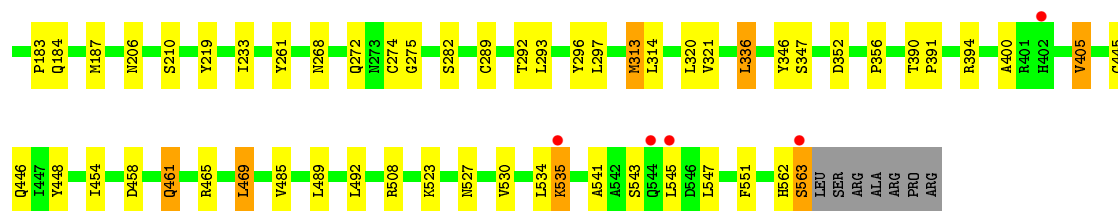


- Molecule 1: HCV NS5B polymerase

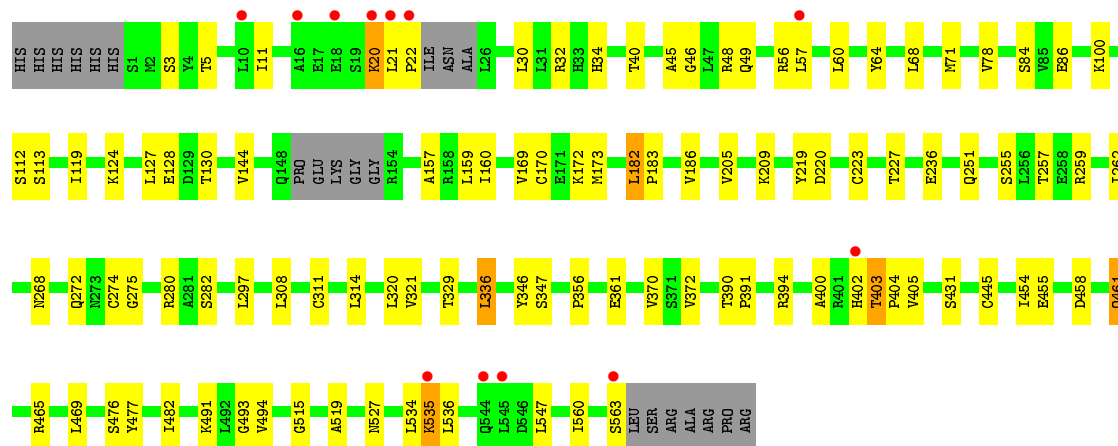
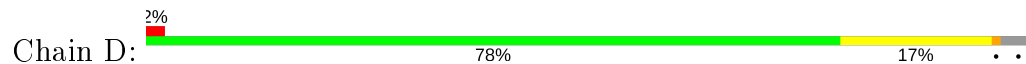


- Molecule 1: HCV NS5B polymerase





● Molecule 1: HCV NS5B polymerase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	101.75Å 101.86Å 250.87Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.91 – 2.17 49.91 – 2.17	Depositor EDS
% Data completeness (in resolution range)	92.9 (49.91-2.17) 92.4 (49.91-2.17)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.38 (at 2.18Å)	Xtrriage
Refinement program	PHENIX 1.7_650	Depositor
R, R_{free}	0.191 , 0.243 0.184 , 0.234	Depositor DCC
R_{free} test set	13245 reflections (9.98%)	wwPDB-VP
Wilson B-factor (Å ²)	30.4	Xtrriage
Anisotropy	0.367	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 34.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.087 for k,h,-l	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	18412	wwPDB-VP
Average B, all atoms (Å ²)	36.0	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 20.64 % 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 8.3176e-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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.22	0/4412	0.41	0/5985
1	B	0.23	0/4412	0.41	0/5985
1	C	0.23	0/4412	0.40	0/5985
1	D	0.22	0/4412	0.41	0/5985
All	All	0.23	0/17648	0.41	0/23940

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	4319	0	4334	39	0
1	B	4319	0	4336	63	0
1	C	4319	0	4334	57	0
1	D	4319	0	4334	60	0
2	A	43	0	26	0	0
2	B	43	0	26	2	0
2	C	43	0	26	0	0
2	D	43	0	26	1	0
3	A	274	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	248	0	0	2	0
3	C	231	0	0	1	0
3	D	211	0	0	1	0
All	All	18412	0	17442	220	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 6.

All (220) 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:D:71:MET:HE1	1:D:297:LEU:HD13	1.49	0.94
1:D:268:ASN:HD21	1:D:272:GLN:HE21	1.18	0.91
1:B:187:MET:HE3	1:B:293:LEU:HA	1.60	0.83
1:B:187:MET:HE3	1:B:293:LEU:HD23	1.60	0.83
1:C:527:ASN:HD21	1:C:534:LEU:H	1.23	0.83
1:D:46:GLY:HA2	1:D:49:GLN:HE21	1.42	0.83
1:D:527:ASN:HD21	1:D:534:LEU:H	1.26	0.83
1:B:187:MET:CE	1:B:293:LEU:HA	2.09	0.83
1:C:187:MET:CE	1:C:293:LEU:HA	2.11	0.79
1:A:527:ASN:HD21	1:A:534:LEU:H	1.31	0.78
1:A:79:LYS:HE3	1:A:81:LYS:HE2	1.65	0.76
1:B:21:LEU:HD12	1:B:22:PRO:HD2	1.67	0.75
1:B:535:LYS:H	1:B:535:LYS:HD2	1.51	0.75
1:A:144:VAL:HB	1:A:394:ARG:HG2	1.70	0.73
1:B:527:ASN:HD21	1:B:534:LEU:H	1.37	0.72
1:B:148:GLN:HA	1:B:148:GLN:HE21	1.53	0.72
1:B:20:LYS:N	1:B:20:LYS:HD2	2.05	0.72
1:B:20:LYS:H	1:B:20:LYS:HD2	1.55	0.71
1:C:458:ASP:HA	1:C:461:GLN:HE21	1.57	0.69
1:D:78:VAL:HG21	1:D:182:LEU:HD23	1.74	0.69
1:D:68:LEU:HA	1:D:71:MET:HE2	1.74	0.69
1:D:130:THR:HG22	1:D:251:GLN:HE22	1.57	0.68
1:B:187:MET:HE1	1:B:296:TYR:HB3	1.75	0.67
1:C:268:ASN:HD21	1:C:272:GLN:HE21	1.40	0.67
1:B:71:MET:HE2	1:B:297:LEU:HD13	1.77	0.67
1:B:119:ILE:HD13	1:B:169:VAL:HG11	1.77	0.66
1:C:22:PRO:HG2	1:C:400:ALA:HB1	1.77	0.66
1:C:523:LYS:HG3	1:C:534:LEU:HD23	1.75	0.66
1:A:119:ILE:HD13	1:A:169:VAL:HG11	1.77	0.65
1:C:68:LEU:HA	1:C:71:MET:HE2	1.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:160:ILE:HD12	1:C:282:SER:OG	1.95	0.65
1:C:31:LEU:HD11	1:C:492:LEU:HD22	1.79	0.64
1:C:535:LYS:HD2	1:C:535:LYS:H	1.64	0.63
1:C:19:SER:HB3	1:C:43:ARG:HH21	1.65	0.62
1:C:187:MET:CE	1:C:293:LEU:HD23	2.30	0.62
1:A:406:ASN:ND2	1:A:443:LEU:HB3	2.16	0.61
1:B:465:ARG:HH11	1:B:547:LEU:HD22	1.66	0.60
1:D:144:VAL:HB	1:D:394:ARG:HG2	1.81	0.60
1:B:423:MET:HB2	2:B:571:KLI:H34	1.84	0.59
1:C:535:LYS:CD	1:C:535:LYS:H	2.16	0.59
1:C:144:VAL:HB	1:C:394:ARG:HG2	1.85	0.58
1:A:268:ASN:HB3	1:A:274:CYS:SG	2.44	0.58
1:A:314:LEU:HB3	1:A:321:VAL:HG13	1.84	0.57
1:C:178:VAL:HG23	3:C:704:HOH:O	2.05	0.57
1:C:187:MET:HE3	1:C:293:LEU:HA	1.85	0.57
1:C:21:LEU:HD12	1:C:22:PRO:HD2	1.86	0.57
1:C:336:LEU:HD12	1:C:356:PRO:HD3	1.86	0.57
1:D:535:LYS:O	1:D:536:LEU:HB2	2.04	0.56
1:C:48:ARG:HG2	1:C:159:LEU:HG	1.88	0.56
1:D:390:THR:HB	1:D:391:PRO:HD3	1.86	0.56
1:C:187:MET:HE3	1:C:293:LEU:HD23	1.88	0.56
1:D:160:ILE:HD12	1:D:282:SER:OG	2.05	0.56
1:B:86:GLU:O	1:B:90:LYS:HE3	2.07	0.55
1:C:219:TYR:HB3	1:C:320:LEU:HD23	1.89	0.55
1:B:464:GLU:HG3	1:B:469:LEU:HD13	1.88	0.55
1:D:458:ASP:O	1:D:461:GLN:HG2	2.06	0.55
1:A:406:ASN:HD22	1:A:443:LEU:HB3	1.71	0.55
1:C:390:THR:HB	1:C:391:PRO:HD3	1.88	0.55
1:B:187:MET:CE	1:B:293:LEU:HD23	2.34	0.54
1:B:314:LEU:HB3	1:B:321:VAL:CG1	2.37	0.54
1:C:268:ASN:HB3	1:C:274:CYS:SG	2.47	0.54
1:A:535:LYS:HD2	1:A:535:LYS:H	1.73	0.54
1:B:314:LEU:HB3	1:B:321:VAL:HG13	1.90	0.54
1:D:11:ILE:HD12	1:D:45:ALA:HB1	1.90	0.53
1:D:130:THR:HG22	1:D:251:GLN:NE2	2.23	0.53
1:D:48:ARG:HG2	1:D:159:LEU:HG	1.89	0.53
1:D:3:SER:HB3	1:D:56:ARG:HD2	1.89	0.53
1:C:268:ASN:HD21	1:C:272:GLN:NE2	2.04	0.53
1:D:455:GLU:HB3	3:D:834:HOH:O	2.09	0.53
1:D:68:LEU:HA	1:D:71:MET:CE	2.37	0.53
1:C:119:ILE:HD13	1:C:169:VAL:HG11	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:117:ASN:HB2	3:B:699:HOH:O	2.09	0.53
1:B:308:LEU:HB3	1:B:311:CYS:SG	2.49	0.52
1:A:314:LEU:HB3	1:A:321:VAL:CG1	2.38	0.52
1:D:314:LEU:HB3	1:D:321:VAL:CG1	2.39	0.52
1:A:390:THR:HB	1:A:391:PRO:HD3	1.91	0.52
1:B:46:GLY:O	1:B:50:LYS:HD2	2.09	0.52
1:D:182:LEU:HD22	1:D:186:VAL:HG23	1.91	0.52
1:B:268:ASN:HB3	1:B:274:CYS:SG	2.50	0.52
1:C:187:MET:HE1	1:C:296:TYR:HB3	1.90	0.52
1:C:20:LYS:O	1:C:20:LYS:HD2	2.09	0.52
1:D:268:ASN:HB3	1:D:274:CYS:SG	2.50	0.52
1:B:144:VAL:HB	1:B:394:ARG:HG2	1.92	0.51
1:A:270:LYS:HB2	1:A:272:GLN:HE21	1.74	0.51
1:B:390:THR:HB	1:B:391:PRO:HD3	1.92	0.51
1:B:31:LEU:HD11	1:B:492:LEU:HD22	1.92	0.51
1:A:160:ILE:HD12	1:A:282:SER:OG	2.11	0.51
1:A:67:VAL:O	1:A:71:MET:HG3	2.11	0.51
1:B:303:CYS:HG	1:B:311:CYS:HG	1.55	0.51
1:B:535:LYS:HG2	1:B:536:LEU:N	2.26	0.51
1:B:46:GLY:HA2	1:B:49:GLN:HE21	1.75	0.50
1:C:314:LEU:HB3	1:C:321:VAL:CG1	2.40	0.50
1:D:130:THR:CG2	1:D:251:GLN:HE22	2.24	0.50
1:D:477:TYR:HB2	1:D:482:ILE:HD11	1.94	0.50
1:D:71:MET:HE1	1:D:297:LEU:CD1	2.33	0.50
1:C:445:CYS:SG	1:C:454:ILE:HD12	2.52	0.50
1:A:219:TYR:HB3	1:A:320:LEU:HD23	1.94	0.50
1:C:535:LYS:N	1:C:535:LYS:HD2	2.26	0.50
1:D:336:LEU:HD12	1:D:356:PRO:HD3	1.94	0.50
1:A:21:LEU:HD23	1:A:34:HIS:HB2	1.94	0.49
1:A:71:MET:HE2	1:A:297:LEU:HD13	1.94	0.49
1:B:48:ARG:HG2	1:B:159:LEU:HG	1.94	0.49
1:B:170:CYS:HA	1:B:173:MET:CE	2.42	0.49
1:B:183:PRO:HG3	1:B:289:CYS:SG	2.53	0.49
1:D:84:SER:OG	1:D:86:GLU:HG2	2.12	0.49
1:C:314:LEU:HB3	1:C:321:VAL:HG13	1.93	0.49
1:D:205:VAL:HG12	1:D:209:LYS:HE2	1.95	0.49
1:D:403:THR:HG22	1:D:404:PRO:HD2	1.94	0.49
1:D:60:LEU:HD13	1:D:64:TYR:CE2	2.48	0.48
1:D:172:LYS:HE3	1:D:560:ILE:HD13	1.95	0.48
1:B:68:LEU:HA	1:B:71:MET:HE2	1.96	0.48
1:C:68:LEU:HA	1:C:71:MET:CE	2.42	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:515:GLY:HA2	1:D:519:ALA:HB2	1.96	0.48
1:D:22:PRO:HG2	1:D:400:ALA:HB1	1.95	0.48
1:B:178:VAL:HG23	3:B:730:HOH:O	2.12	0.47
1:C:187:MET:HE2	1:C:293:LEU:HD23	1.95	0.47
1:D:445:CYS:SG	1:D:454:ILE:HD12	2.54	0.47
1:B:257:THR:O	1:B:262:ILE:HG23	2.15	0.47
1:B:535:LYS:CD	1:B:535:LYS:H	2.25	0.47
1:C:71:MET:HE1	1:C:297:LEU:HD22	1.96	0.47
1:A:178:VAL:HG23	3:A:734:HOH:O	2.14	0.47
1:A:346:TYR:O	1:A:347:SER:HB3	2.15	0.47
1:A:535:LYS:CD	1:A:535:LYS:H	2.28	0.47
1:B:84:SER:OG	1:B:86:GLU:HG2	2.15	0.47
1:D:361:GLU:HG3	1:D:370:VAL:O	2.15	0.47
1:D:32:ARG:HG2	1:D:493:GLY:O	2.15	0.46
1:A:523:LYS:HG3	1:A:534:LEU:HD23	1.97	0.46
1:D:236:GLU:OE2	1:D:280:ARG:NH2	2.35	0.46
1:B:219:TYR:HB3	1:B:320:LEU:HD23	1.97	0.46
1:D:127:LEU:HA	1:D:251:GLN:HG3	1.98	0.46
1:B:187:MET:HE1	1:B:296:TYR:CB	2.44	0.46
1:C:508:ARG:CZ	1:C:530:VAL:HG11	2.46	0.46
1:C:233:ILE:HD13	1:C:261:TYR:O	2.15	0.46
1:A:517:ARG:H	1:A:517:ARG:HG2	1.40	0.45
1:B:535:LYS:O	1:B:536:LEU:HB2	2.15	0.45
1:C:20:LYS:C	1:C:20:LYS:HD2	2.37	0.45
1:C:183:PRO:HG3	1:C:289:CYS:SG	2.56	0.45
1:C:313:MET:HE2	1:C:320:LEU:HD11	1.98	0.45
1:D:257:THR:O	1:D:262:ILE:HG23	2.15	0.45
1:A:183:PRO:HG3	1:A:289:CYS:SG	2.57	0.45
1:B:170:CYS:HA	1:B:173:MET:HE3	1.98	0.45
1:B:233:ILE:HD13	1:B:261:TYR:O	2.16	0.45
1:C:47:LEU:HB2	1:C:156:PRO:HB3	1.99	0.45
1:A:71:MET:HE3	1:A:71:MET:HB2	1.72	0.45
1:B:191:TYR:O	1:B:194:GLN:HG2	2.17	0.45
1:D:71:MET:CE	1:D:297:LEU:HD13	2.35	0.45
1:C:405:VAL:HG21	1:C:446:GLN:HG2	1.99	0.45
1:A:461:GLN:H	1:A:461:GLN:HE21	1.65	0.45
1:C:187:MET:HE1	1:C:296:TYR:CB	2.47	0.45
1:C:71:MET:HE2	1:C:297:LEU:HD13	1.98	0.45
1:D:314:LEU:HB3	1:D:321:VAL:HG13	1.99	0.45
1:A:3:SER:HB3	1:A:56:ARG:HD2	1.98	0.44
1:C:71:MET:CE	1:C:297:LEU:HD13	2.46	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:517:ARG:HG2	1:B:517:ARG:H	1.54	0.44
1:C:485:VAL:O	1:C:489:LEU:HG	2.17	0.44
1:D:20:LYS:HD2	1:D:20:LYS:O	2.18	0.44
1:D:535:LYS:H	1:D:535:LYS:HG2	1.51	0.44
1:D:46:GLY:HA2	1:D:49:GLN:NE2	2.21	0.44
1:C:458:ASP:CA	1:C:461:GLN:HE21	2.29	0.44
1:C:448:TYR:CE2	1:C:551:PHE:HD1	2.35	0.44
1:D:346:TYR:O	1:D:347:SER:HB3	2.17	0.44
1:A:182:LEU:HB3	1:A:183:PRO:HD3	1.99	0.44
1:A:461:GLN:HB2	1:A:542:ALA:HA	2.00	0.44
1:A:19:SER:H	1:A:20:LYS:HD2	1.81	0.43
1:A:211:LYS:HB2	1:A:214:PRO:HB3	2.00	0.43
1:B:336:LEU:HD22	1:B:354:PRO:HG2	2.00	0.43
1:C:458:ASP:HA	1:C:461:GLN:NE2	2.28	0.43
1:D:255:SER:OG	1:D:259:ARG:HD3	2.17	0.43
1:B:445:CYS:SG	1:B:454:ILE:HD12	2.59	0.43
1:B:22:PRO:HG2	1:B:400:ALA:HB1	2.01	0.43
1:B:416:ALA:N	1:B:417:PRO:CD	2.81	0.43
1:B:562:HIS:O	1:B:563:SER:HB3	2.18	0.43
2:B:571:KLI:H28	2:B:571:KLI:F43	2.07	0.43
1:D:119:ILE:HD13	1:D:169:VAL:HG11	2.00	0.43
1:B:535:LYS:HG2	1:B:536:LEU:H	1.83	0.43
1:D:476:SER:HA	2:D:571:KLI:O37	2.18	0.43
1:C:465:ARG:HG3	1:C:545:LEU:HD12	2.01	0.43
1:D:30:LEU:O	1:D:494:VAL:HG22	2.19	0.43
1:B:237:GLU:HG3	1:B:257:THR:OG1	2.18	0.43
1:C:469:LEU:HD12	1:C:469:LEU:HA	1.88	0.42
1:D:21:LEU:HD23	1:D:34:HIS:CB	2.50	0.42
1:A:31:LEU:HD23	1:A:31:LEU:C	2.39	0.42
1:C:562:HIS:O	1:C:563:SER:HB2	2.20	0.42
1:B:141:LYS:HE3	1:B:141:LYS:HB3	1.87	0.42
1:B:469:LEU:HA	1:B:469:LEU:HD12	1.85	0.42
1:D:219:TYR:HB3	1:D:320:LEU:HD23	2.00	0.42
1:A:48:ARG:HG2	1:A:159:LEU:HG	2.02	0.42
1:D:40:THR:HB	1:D:157:ALA:HB2	2.00	0.42
1:A:20:LYS:N	1:A:20:LYS:HD2	2.35	0.41
1:D:100:LYS:HE2	1:D:100:LYS:HB3	1.87	0.41
1:A:455:GLU:HB3	3:A:724:HOH:O	2.20	0.41
1:B:361:GLU:HG3	1:B:370:VAL:O	2.20	0.41
1:C:5:THR:O	1:C:275:GLY:HA3	2.19	0.41
1:D:21:LEU:HD23	1:D:34:HIS:HB2	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:83:LEU:HB2	1:B:173:MET:HA	2.02	0.41
1:A:270:LYS:HD2	1:A:272:GLN:NE2	2.35	0.41
1:B:71:MET:HB2	1:B:71:MET:HE3	1.81	0.41
1:C:71:MET:HE1	1:C:297:LEU:HB2	2.02	0.41
1:B:465:ARG:NH1	1:B:547:LEU:HD22	2.34	0.41
1:C:346:TYR:O	1:C:347:SER:HB3	2.20	0.41
1:A:139:MET:HB2	1:A:160:ILE:CG2	2.51	0.41
1:D:124:LYS:HE3	1:D:128:GLU:OE1	2.21	0.41
1:A:32:ARG:HG2	1:A:493:GLY:O	2.21	0.41
1:A:416:ALA:N	1:A:417:PRO:CD	2.83	0.41
1:B:346:TYR:O	1:B:347:SER:HB3	2.20	0.41
1:B:458:ASP:O	1:B:461:GLN:HG2	2.21	0.41
1:C:206:ASN:O	1:C:210:SER:HB3	2.21	0.41
1:D:170:CYS:HA	1:D:173:MET:HE3	2.03	0.41
1:D:5:THR:O	1:D:275:GLY:HA3	2.20	0.41
1:C:541:ALA:C	1:C:543:SER:H	2.23	0.41
1:B:268:ASN:HD21	1:B:272:GLN:NE2	2.18	0.41
1:A:308:LEU:CD1	1:A:335:SER:HB3	2.51	0.41
1:B:187:MET:CE	1:B:296:TYR:HB3	2.47	0.40
1:D:491:LYS:HE2	1:D:491:LYS:HB3	1.92	0.40
1:C:187:MET:HE1	1:C:292:THR:O	2.21	0.40
1:D:182:LEU:N	1:D:183:PRO:CD	2.84	0.40
1:D:223:CYS:O	1:D:227:THR:HG23	2.20	0.40
1:D:469:LEU:HD12	1:D:469:LEU:HA	1.93	0.40
1:B:319:ASP:CG	1:B:366:CYS:H	2.25	0.40
1:B:485:VAL:O	1:B:489:LEU:HG	2.20	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	549/576 (95%)	533 (97%)	15 (3%)	1 (0%)	47	52
1	B	549/576 (95%)	538 (98%)	11 (2%)	0	100	100
1	C	549/576 (95%)	534 (97%)	15 (3%)	0	100	100
1	D	549/576 (95%)	531 (97%)	18 (3%)	0	100	100
All	All	2196/2304 (95%)	2136 (97%)	59 (3%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	351	GLY

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	474/491 (96%)	453 (96%)	21 (4%)	28	33
1	B	474/491 (96%)	449 (95%)	25 (5%)	22	25
1	C	474/491 (96%)	459 (97%)	15 (3%)	39	47
1	D	474/491 (96%)	454 (96%)	20 (4%)	30	35
All	All	1896/1964 (96%)	1815 (96%)	81 (4%)	29	34

All (81) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	LYS
1	A	26	LEU
1	A	57	LEU
1	A	69	LYS
1	A	90	LYS
1	A	100	LYS
1	A	106	LYS
1	A	209	LYS
1	A	273	ASN
1	A	308	LEU

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Mol	Chain	Res	Type
1	A	311	CYS
1	A	313	MET
1	A	336	LEU
1	A	402	HIS
1	A	446	GLN
1	A	461	GLN
1	A	465	ARG
1	A	517	ARG
1	A	535	LYS
1	A	545	LEU
1	A	547	LEU
1	B	20	LYS
1	B	26	LEU
1	B	50	LYS
1	B	56	ARG
1	B	57	LEU
1	B	81	LYS
1	B	106	LYS
1	B	113	SER
1	B	148	GLN
1	B	154	ARG
1	B	160	ILE
1	B	213	ASN
1	B	270	LYS
1	B	308	LEU
1	B	311	CYS
1	B	330	GLN
1	B	336	LEU
1	B	372	VAL
1	B	419	LEU
1	B	461	GLN
1	B	465	ARG
1	B	469	LEU
1	B	517	ARG
1	B	535	LYS
1	B	547	LEU
1	C	20	LYS
1	C	26	LEU
1	C	50	LYS
1	C	56	ARG
1	C	106	LYS
1	C	184	GLN

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Mol	Chain	Res	Type
1	C	313	MET
1	C	336	LEU
1	C	352	ASP
1	C	405	VAL
1	C	461	GLN
1	C	469	LEU
1	C	535	LYS
1	C	547	LEU
1	C	563	SER
1	D	20	LYS
1	D	57	LEU
1	D	112	SER
1	D	113	SER
1	D	182	LEU
1	D	220	ASP
1	D	308	LEU
1	D	311	CYS
1	D	329	THR
1	D	336	LEU
1	D	372	VAL
1	D	402	HIS
1	D	403	THR
1	D	405	VAL
1	D	431	SER
1	D	461	GLN
1	D	465	ARG
1	D	535	LYS
1	D	547	LEU
1	D	563	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (40) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	49	GLN
1	A	272	GLN
1	A	273	ASN
1	A	406	ASN
1	A	436	GLN
1	A	461	GLN
1	A	475	HIS
1	A	483	ASN
1	A	514	GLN

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Mol	Chain	Res	Type
1	A	527	ASN
1	B	49	GLN
1	B	148	GLN
1	B	213	ASN
1	B	272	GLN
1	B	406	ASN
1	B	436	GLN
1	B	475	HIS
1	B	514	GLN
1	B	527	ASN
1	C	142	ASN
1	C	184	GLN
1	C	213	ASN
1	C	251	GLN
1	C	272	GLN
1	C	406	ASN
1	C	461	GLN
1	C	475	HIS
1	C	483	ASN
1	C	514	GLN
1	C	527	ASN
1	D	49	GLN
1	D	213	ASN
1	D	251	GLN
1	D	272	GLN
1	D	406	ASN
1	D	436	GLN
1	D	461	GLN
1	D	475	HIS
1	D	514	GLN
1	D	527	ASN

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](#)

4 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	KLI	D	571	-	42,48,48	1.78	5 (11%)	53,71,71	1.47	8 (15%)
2	KLI	B	571	-	42,48,48	1.81	5 (11%)	53,71,71	1.46	7 (13%)
2	KLI	C	571	-	42,48,48	1.74	5 (11%)	53,71,71	1.44	9 (16%)
2	KLI	A	571	-	42,48,48	1.77	5 (11%)	53,71,71	1.50	6 (11%)

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	KLI	D	571	-	-	4/21/32/32	0/6/6/6
2	KLI	B	571	-	-	1/21/32/32	0/6/6/6
2	KLI	C	571	-	-	4/21/32/32	0/6/6/6
2	KLI	A	571	-	-	2/21/32/32	0/6/6/6

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	571	KLI	C12-N01	8.56	1.39	1.33
2	A	571	KLI	C12-N01	8.36	1.39	1.33
2	D	571	KLI	C12-N01	8.33	1.39	1.33
2	C	571	KLI	C12-N01	8.06	1.39	1.33
2	B	571	KLI	C11-C06	3.68	1.41	1.35
2	D	571	KLI	C11-C06	3.63	1.40	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	571	KLI	C11-C06	3.60	1.40	1.35
2	A	571	KLI	C11-C06	3.54	1.40	1.35
2	D	571	KLI	C08-C07	2.39	1.42	1.38
2	C	571	KLI	C08-C07	2.36	1.42	1.38
2	B	571	KLI	C08-C07	2.32	1.42	1.38
2	A	571	KLI	C08-C07	2.29	1.42	1.38
2	C	571	KLI	C18-C23	2.28	1.42	1.38
2	D	571	KLI	C18-C23	2.27	1.42	1.38
2	B	571	KLI	C18-C23	2.23	1.42	1.38
2	D	571	KLI	C12-C13	-2.19	1.35	1.39
2	A	571	KLI	C18-C23	2.18	1.41	1.38
2	C	571	KLI	C12-C13	-2.16	1.35	1.39
2	A	571	KLI	C12-C13	-2.09	1.35	1.39
2	B	571	KLI	C12-C13	-2.09	1.35	1.39

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	571	KLI	C12-C13-C14	-5.78	117.06	119.97
2	A	571	KLI	C12-C13-C14	-5.44	117.24	119.97
2	B	571	KLI	C12-C13-C14	-5.37	117.27	119.97
2	C	571	KLI	C12-C13-C14	-5.23	117.34	119.97
2	A	571	KLI	C18-C17-N01	-3.51	107.34	113.23
2	B	571	KLI	C13-C15-N02	-3.31	119.46	123.67
2	B	571	KLI	F43-C06-C07	3.24	121.36	118.42
2	C	571	KLI	C13-C15-N02	-3.20	119.59	123.67
2	A	571	KLI	C14-C10-C09	3.12	121.31	118.59
2	C	571	KLI	C14-C10-C09	3.01	121.21	118.59
2	D	571	KLI	C14-C10-C09	3.01	121.21	118.59
2	D	571	KLI	C28-C27-N05	-2.94	107.48	110.80
2	D	571	KLI	C13-C15-N02	-2.88	120.00	123.67
2	D	571	KLI	C18-C17-N01	-2.86	108.44	113.23
2	A	571	KLI	C13-C15-N02	-2.78	120.13	123.67
2	C	571	KLI	F43-C06-C07	2.77	120.94	118.42
2	A	571	KLI	C22-C23-C18	-2.73	120.39	123.98
2	B	571	KLI	C14-C10-C09	2.73	120.96	118.59
2	D	571	KLI	C22-C23-C18	-2.60	120.57	123.98
2	B	571	KLI	C22-C23-C18	-2.57	120.60	123.98
2	A	571	KLI	F43-C06-C07	2.55	120.74	118.42
2	B	571	KLI	C25-N04-C07	2.53	122.28	116.27
2	C	571	KLI	C28-C27-N05	-2.53	107.95	110.80
2	C	571	KLI	C22-C23-C18	-2.36	120.88	123.98

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	571	KLI	C18-C17-N01	-2.36	109.28	113.23
2	D	571	KLI	F43-C06-C07	2.31	120.52	118.42
2	C	571	KLI	F39-C24-C21	-2.23	108.04	112.93
2	B	571	KLI	F39-C24-C21	-2.12	108.28	112.93
2	D	571	KLI	F39-C24-C21	-2.01	108.51	112.93
2	C	571	KLI	C25-N04-C07	2.01	121.03	116.27

There are no chirality outliers.

All (11) torsion outliers are listed below:

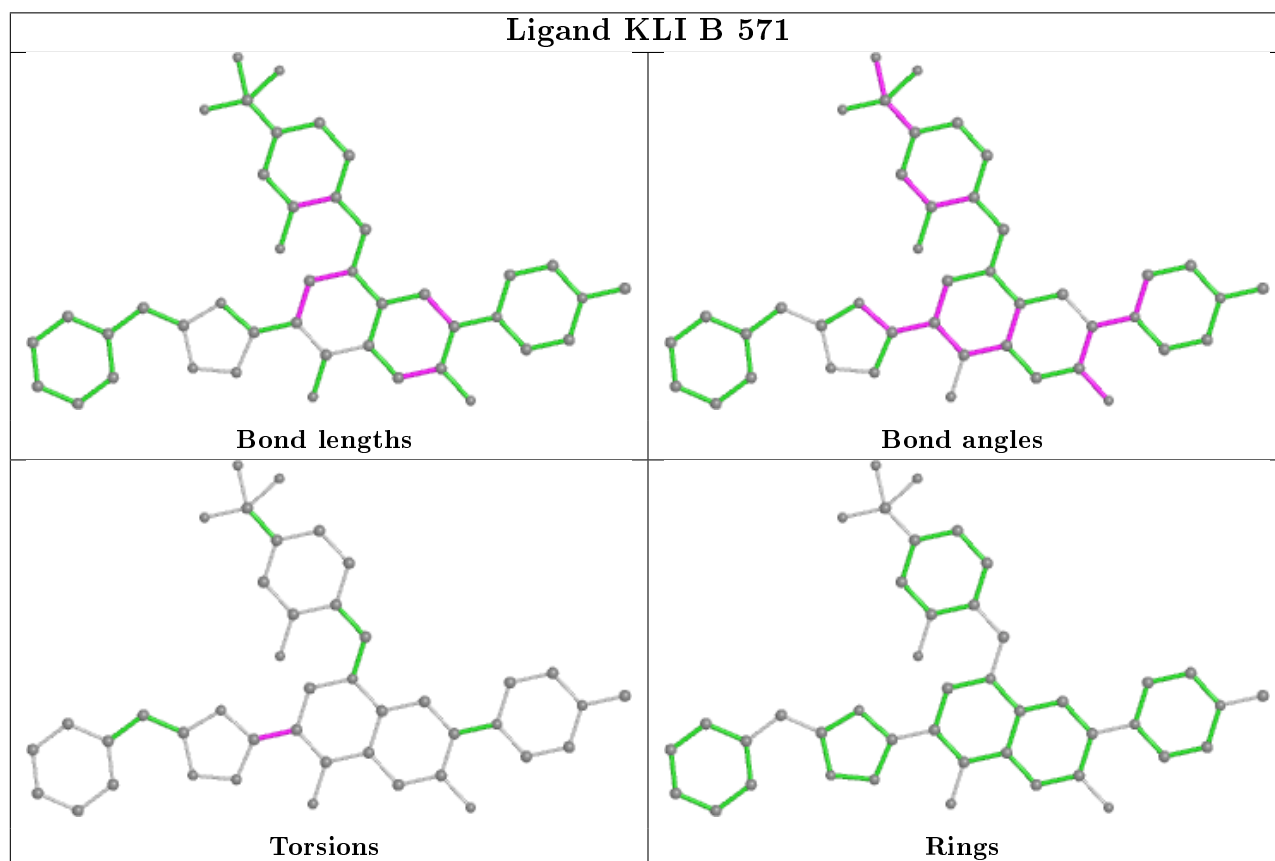
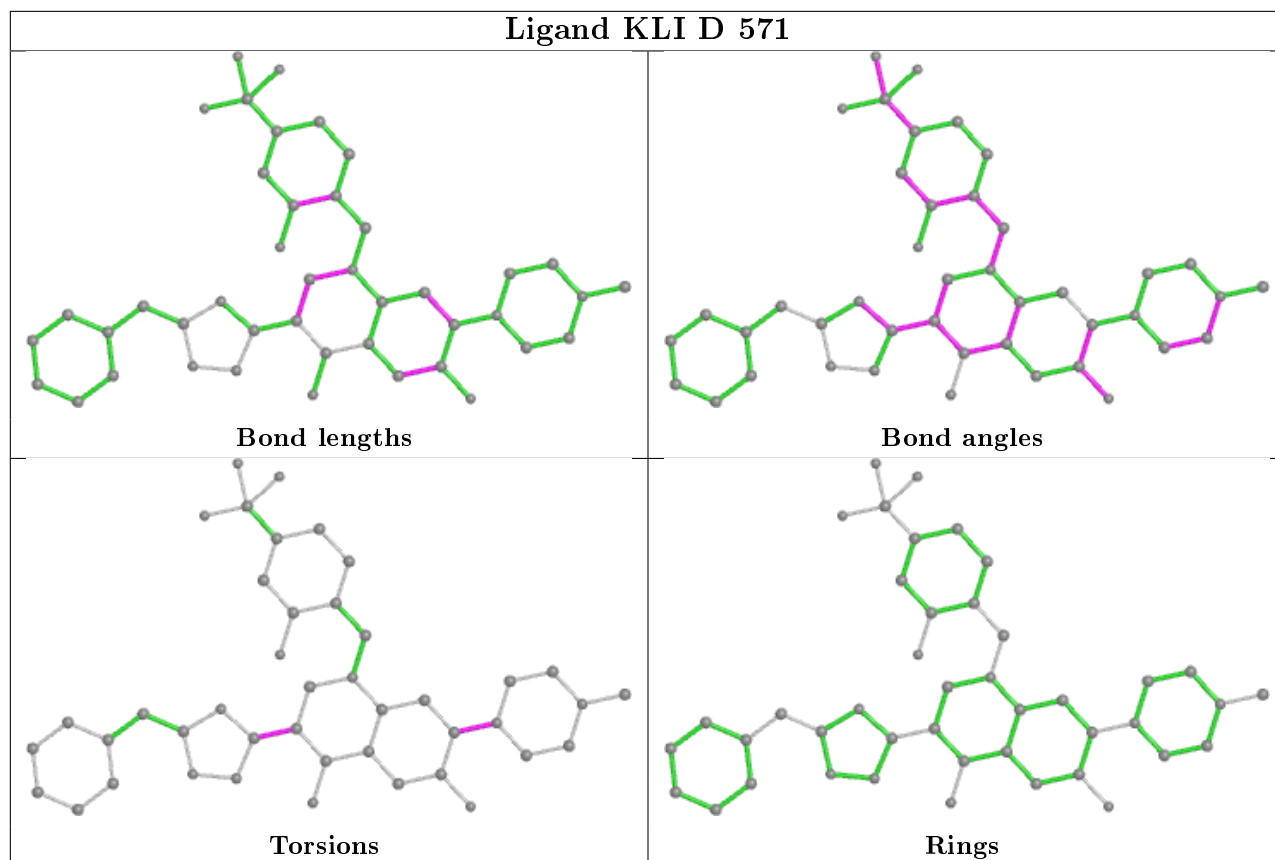
Mol	Chain	Res	Type	Atoms
2	C	571	KLI	C06-C07-N04-C28
2	D	571	KLI	C06-C07-N04-C28
2	C	571	KLI	C06-C07-N04-C25
2	D	571	KLI	C06-C07-N04-C25
2	B	571	KLI	C14-C13-C15-N02
2	C	571	KLI	C14-C13-C15-N02
2	C	571	KLI	C08-C07-N04-C28
2	A	571	KLI	C06-C07-N04-C28
2	D	571	KLI	C08-C07-N04-C28
2	D	571	KLI	C14-C13-C15-N02
2	A	571	KLI	C14-C13-C15-N02

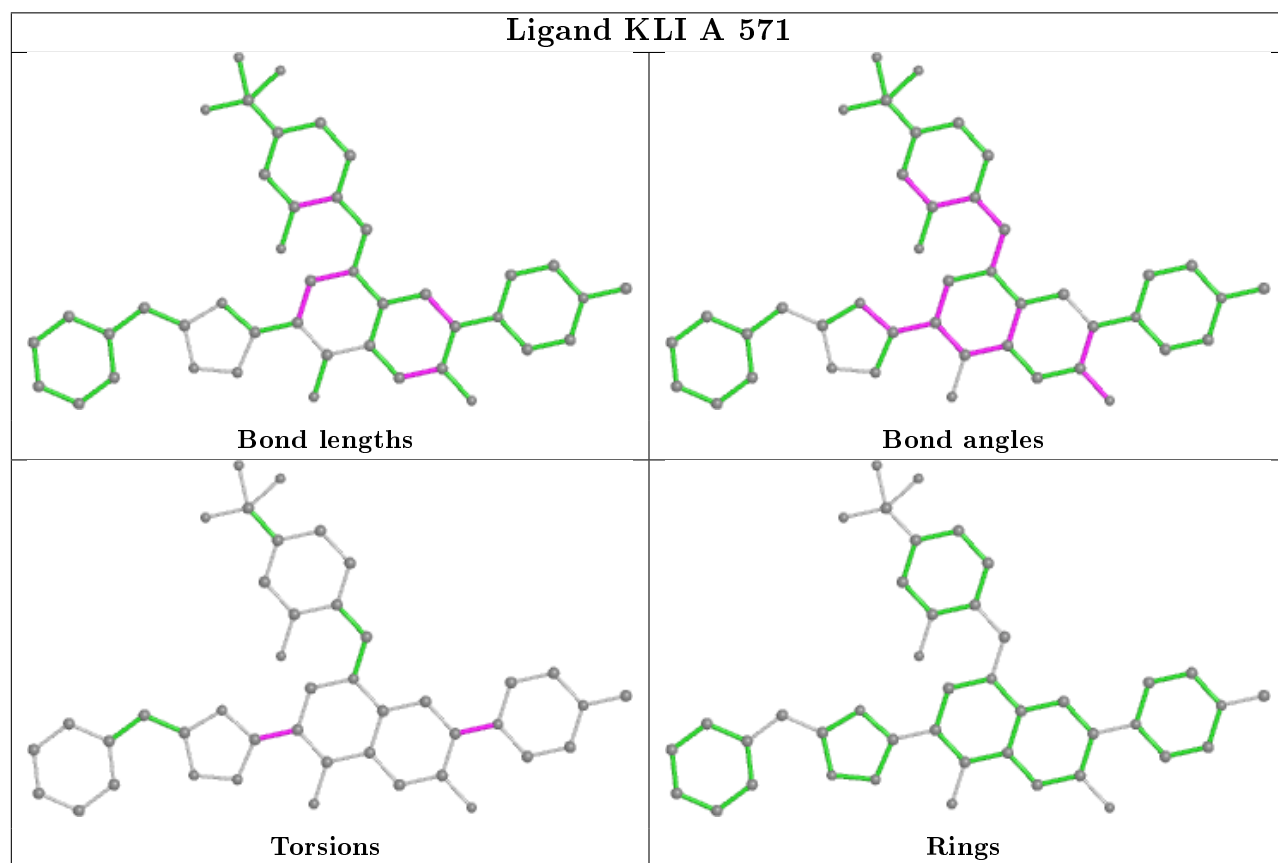
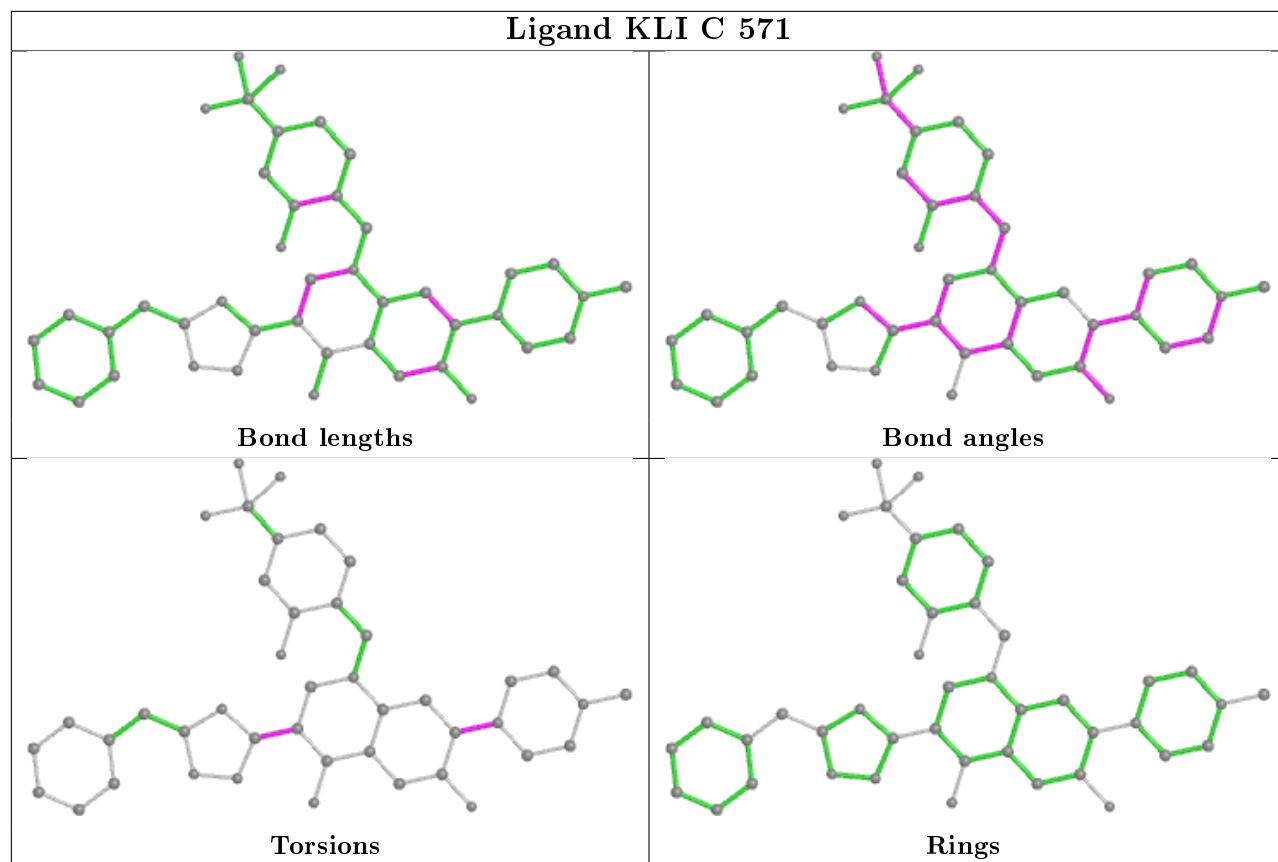
There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	571	KLI	1	0
2	B	571	KLI	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	555/576 (96%)	-0.20	7 (1%) 77 77	21, 33, 57, 82	0
1	B	555/576 (96%)	-0.17	10 (1%) 68 69	21, 33, 59, 81	0
1	C	555/576 (96%)	-0.15	11 (1%) 65 66	21, 34, 59, 85	0
1	D	555/576 (96%)	-0.13	12 (2%) 62 62	20, 36, 61, 85	0
All	All	2220/2304 (96%)	-0.16	40 (1%) 68 69	20, 34, 59, 85	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	563	SER	6.5
1	C	563	SER	5.3
1	C	16	ALA	4.5
1	C	26	LEU	4.2
1	C	544	GLN	3.9
1	C	535	LYS	3.9
1	B	26	LEU	3.8
1	D	545	LEU	3.7
1	B	544	GLN	3.7
1	A	544	GLN	3.6
1	D	544	GLN	3.6
1	B	545	LEU	3.5
1	B	563	SER	3.4
1	A	402	HIS	3.3
1	D	20	LYS	3.3
1	A	545	LEU	3.3
1	C	545	LEU	3.2
1	C	402	HIS	2.9
1	D	21	LEU	2.9
1	B	20	LYS	2.8
1	B	402	HIS	2.7

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Mol	Chain	Res	Type	RSRZ
1	C	15	ALA	2.6
1	D	563	SER	2.6
1	B	34	HIS	2.6
1	D	57	LEU	2.6
1	B	29	SER	2.6
1	D	10	LEU	2.5
1	C	20	LYS	2.5
1	B	27	SER	2.5
1	D	18	GLU	2.4
1	D	535	LYS	2.3
1	A	21	LEU	2.3
1	A	35	ASN	2.2
1	B	22	PRO	2.1
1	C	18	GLU	2.1
1	D	16	ALA	2.1
1	D	22	PRO	2.1
1	C	43	ARG	2.0
1	A	20	LYS	2.0
1	D	402	HIS	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](#)

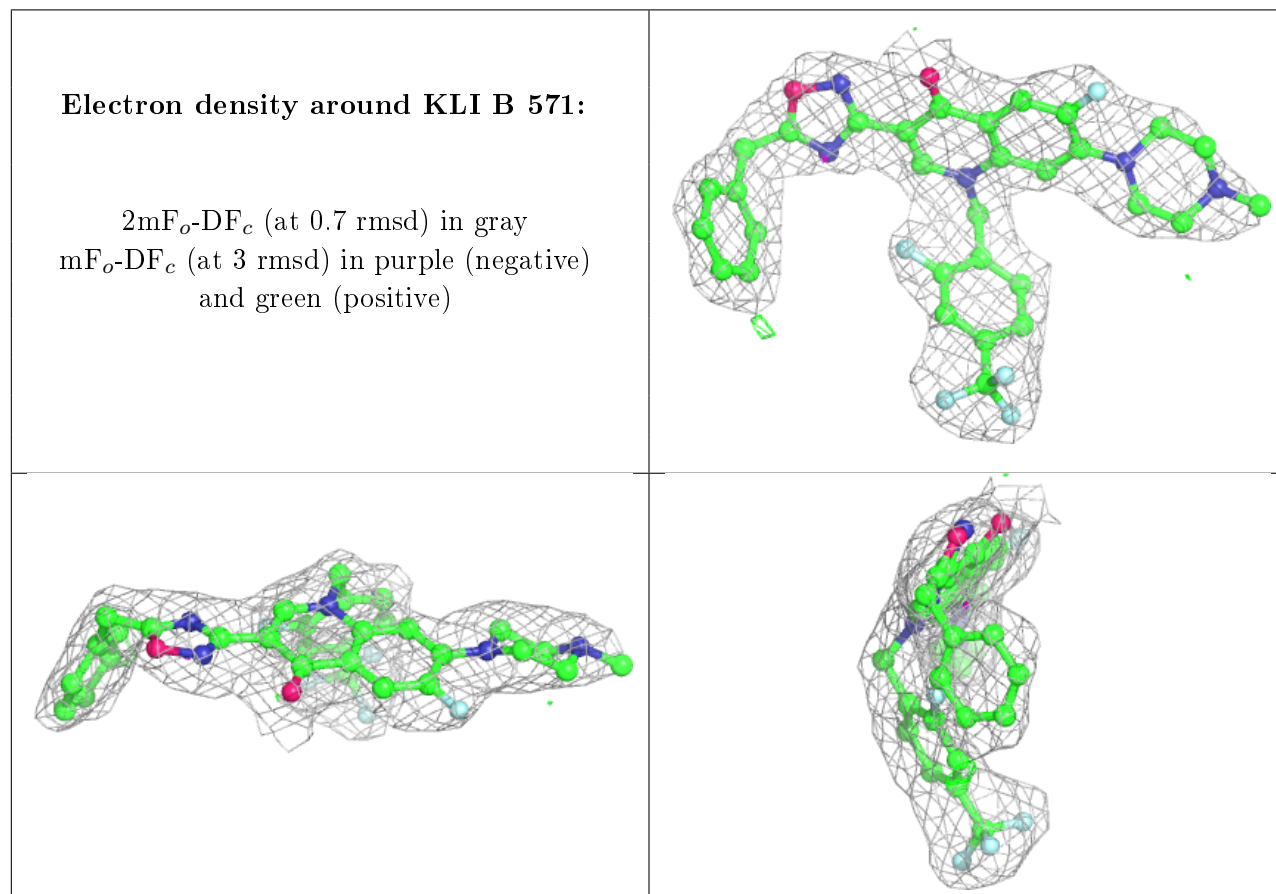
There are no carbohydrates in this entry.

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, 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.

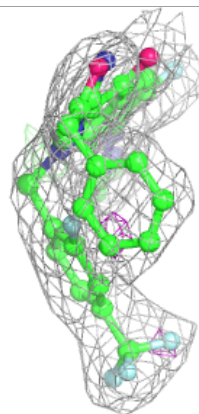
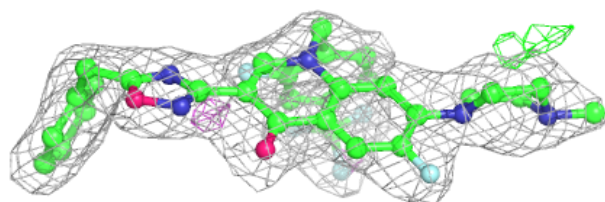
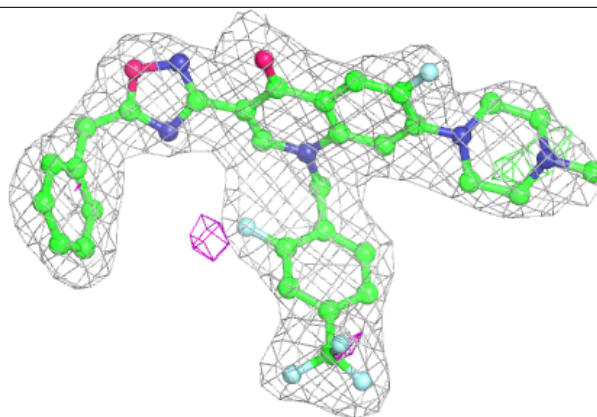
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	KLI	B	571	43/43	0.92	0.12	30,46,55,66	0
2	KLI	A	571	43/43	0.92	0.12	27,41,49,55	0
2	KLI	C	571	43/43	0.93	0.12	30,42,53,61	0
2	KLI	D	571	43/43	0.94	0.11	27,41,52,61	0

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.

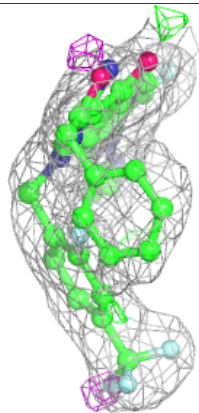
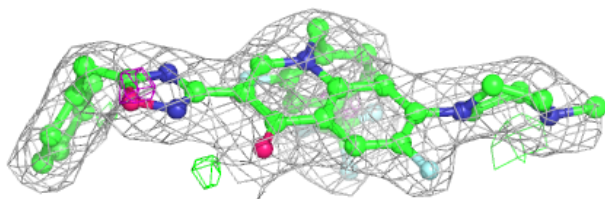
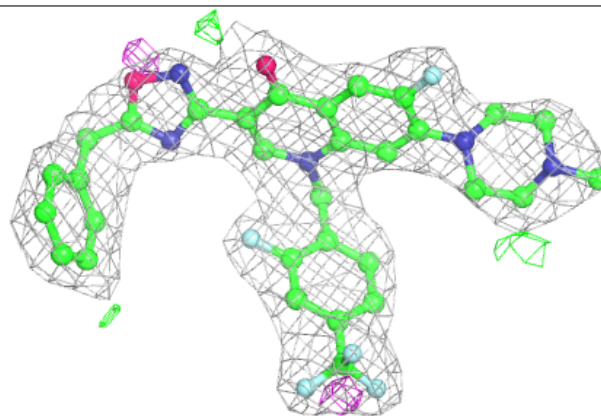


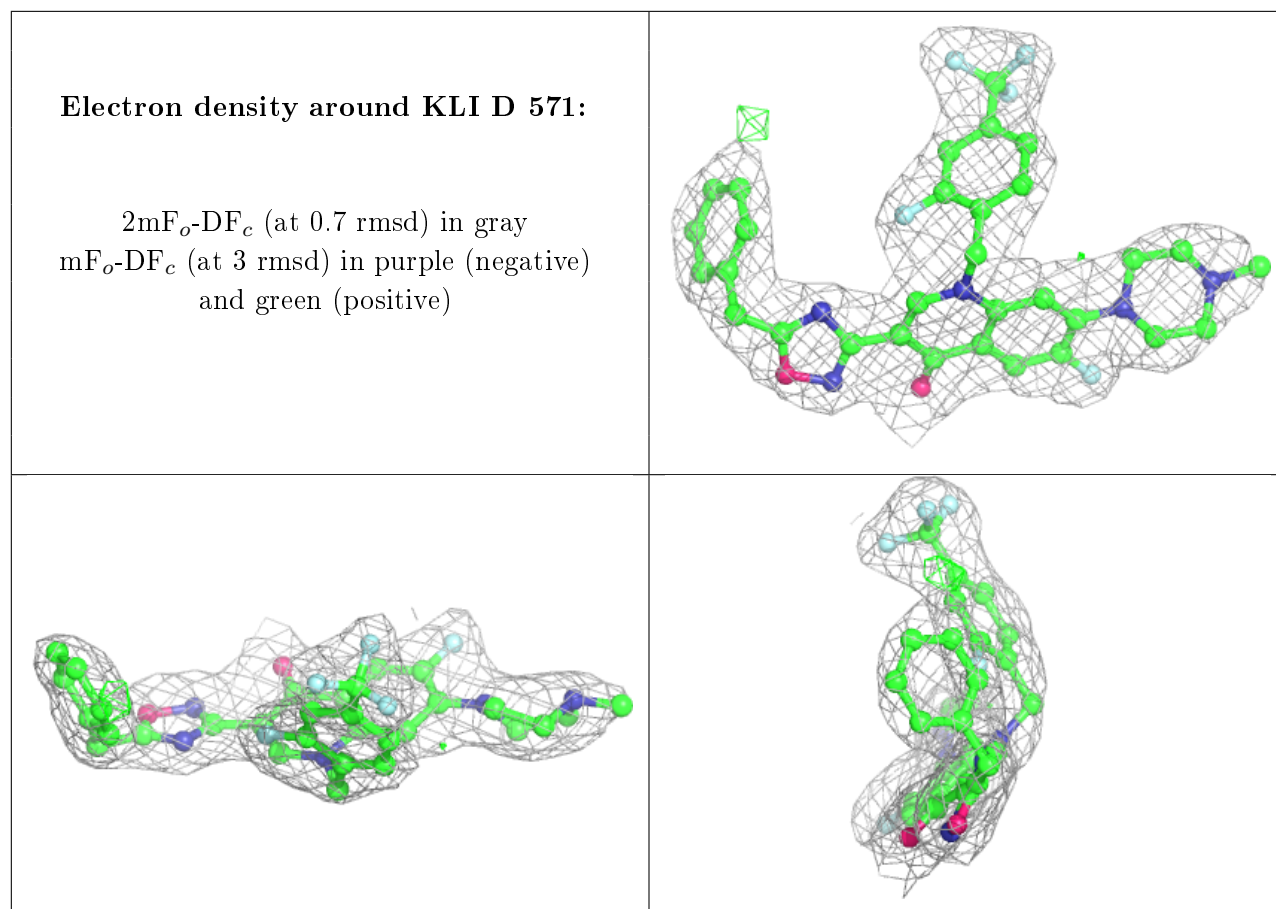
Electron density around KLI A 571:

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

**Electron density around KLI C 571:**

$2mF_o-DF_c$ (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.