



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 15, 2020 – 10:08 am BST

PDB ID : 5FTV
Title : Pseudomonas aeruginosa RmlA in complex with allosteric inhibitor
Authors : Alphey, M.S.; Tran, F.; Westwood, N.J.; Naismith, J.H.
Deposited on : 2016-01-18
Resolution : 2.21 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

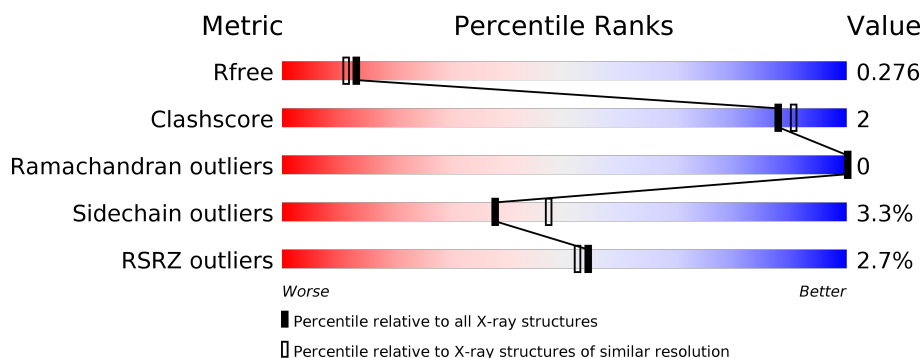
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.21 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	5912 (2.24-2.20)
Clashscore	141614	6646 (2.24-2.20)
Ramachandran outliers	138981	6543 (2.24-2.20)
Sidechain outliers	138945	6544 (2.24-2.20)
RSRZ outliers	127900	5797 (2.24-2.20)

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	303	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 1%, green 98%);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> % 90% 7% • </div> </div>
1	B	303	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 1%, green 98%);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> % 90% 6% • </div> </div>
1	C	303	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 6%, orange 1%, yellow 1%, green 92%);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> 6% 91% 7% • </div> </div>
1	D	303	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 2%, orange 1%, yellow 1%, green 96%);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> 2% 87% 9% • </div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 9879 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 GLUCOSE-1-PHOSPHATE THYMIDYLYLTRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	296	Total	C	N	O	S	0	1	0
			2331	1490	398	439	4			
1	B	292	Total	C	N	O	S	0	1	0
			2291	1466	386	435	4			
1	C	297	Total	C	N	O	S	0	0	0
			2335	1492	400	439	4			
1	D	292	Total	C	N	O	S	0	0	0
			2285	1462	385	434	4			

There are 40 discrepancies between the modelled and reference sequences:

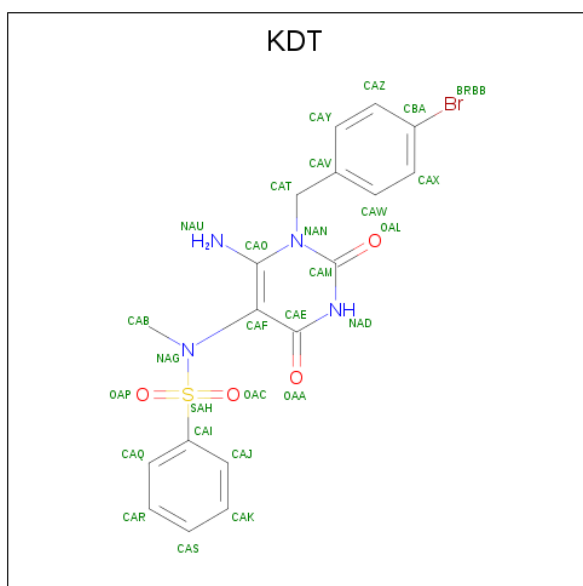
Chain	Residue	Modelled	Actual	Comment	Reference
A	-9	HIS	-	expression tag	UNP Q9HU22
A	-8	HIS	-	expression tag	UNP Q9HU22
A	-7	HIS	-	expression tag	UNP Q9HU22
A	-6	HIS	-	expression tag	UNP Q9HU22
A	-5	HIS	-	expression tag	UNP Q9HU22
A	-4	HIS	-	expression tag	UNP Q9HU22
A	-3	GLY	-	expression tag	UNP Q9HU22
A	-2	SER	-	expression tag	UNP Q9HU22
A	-1	MET	-	expression tag	UNP Q9HU22
A	0	ALA	-	expression tag	UNP Q9HU22
B	-9	HIS	-	expression tag	UNP Q9HU22
B	-8	HIS	-	expression tag	UNP Q9HU22
B	-7	HIS	-	expression tag	UNP Q9HU22
B	-6	HIS	-	expression tag	UNP Q9HU22
B	-5	HIS	-	expression tag	UNP Q9HU22
B	-4	HIS	-	expression tag	UNP Q9HU22
B	-3	GLY	-	expression tag	UNP Q9HU22
B	-2	SER	-	expression tag	UNP Q9HU22
B	-1	MET	-	expression tag	UNP Q9HU22
B	0	ALA	-	expression tag	UNP Q9HU22

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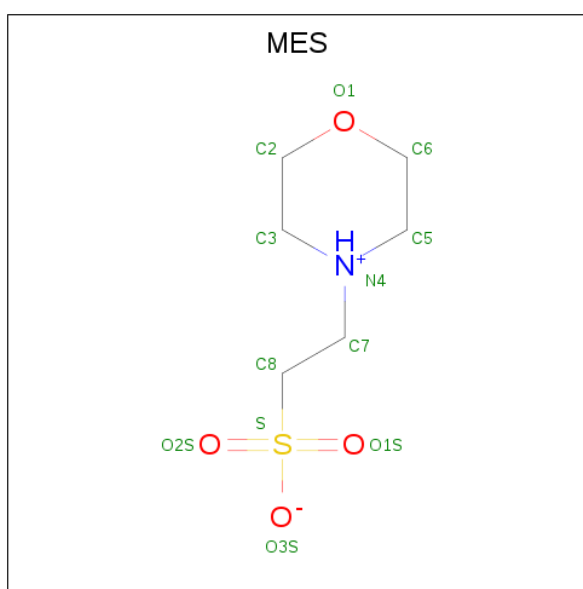
Chain	Residue	Modelled	Actual	Comment	Reference
C	-9	HIS	-	expression tag	UNP Q9HU22
C	-8	HIS	-	expression tag	UNP Q9HU22
C	-7	HIS	-	expression tag	UNP Q9HU22
C	-6	HIS	-	expression tag	UNP Q9HU22
C	-5	HIS	-	expression tag	UNP Q9HU22
C	-4	HIS	-	expression tag	UNP Q9HU22
C	-3	GLY	-	expression tag	UNP Q9HU22
C	-2	SER	-	expression tag	UNP Q9HU22
C	-1	MET	-	expression tag	UNP Q9HU22
C	0	ALA	-	expression tag	UNP Q9HU22
D	-9	HIS	-	expression tag	UNP Q9HU22
D	-8	HIS	-	expression tag	UNP Q9HU22
D	-7	HIS	-	expression tag	UNP Q9HU22
D	-6	HIS	-	expression tag	UNP Q9HU22
D	-5	HIS	-	expression tag	UNP Q9HU22
D	-4	HIS	-	expression tag	UNP Q9HU22
D	-3	GLY	-	expression tag	UNP Q9HU22
D	-2	SER	-	expression tag	UNP Q9HU22
D	-1	MET	-	expression tag	UNP Q9HU22
D	0	ALA	-	expression tag	UNP Q9HU22

- Molecule 2 is N-(6-AMINO-1-(4-BROMOBENZYL)-2,4-DIOXO-1,2,3,4-TETRAHYDROPYRIMIDIN-5-YL)-N-METHYLBENZENESULFONAMIDE (three-letter code: KDT) (formula: C₁₈H₁₇BrN₄O₄S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total 28	Br 1	C 18	N 4	O 4	S 1	0	0
2	B	1	Total 28	Br 1	C 18	N 4	O 4	S 1	0	0
2	C	1	Total 28	Br 1	C 18	N 4	O 4	S 1	0	0
2	D	1	Total 28	Br 1	C 18	N 4	O 4	S 1	0	0

- Molecule 3 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula: C₆H₁₃NO₄S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
3	B	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
3	C	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
3	D	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Cl	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total 1	Cl 1	0	0
4	D	1	Total 1	Cl 1	0	0
4	C	1	Total 1	Cl 1	0	0

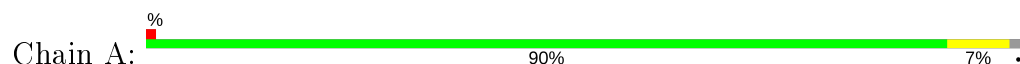
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	128	Total 128	O 128	0	0
5	B	112	Total 112	O 112	0	0
5	C	123	Total 123	O 123	0	0
5	D	110	Total 110	O 110	0	0

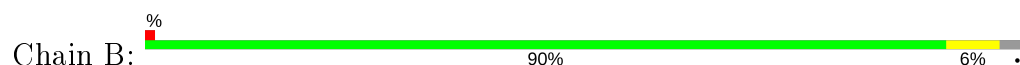
3 Residue-property plots [i](#)

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.

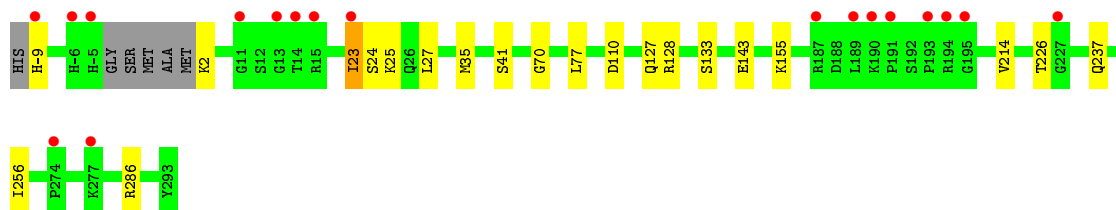
- Molecule 1: GLUCOSE-1-PHOSPHATE THYMIDYLTRANSFERASE



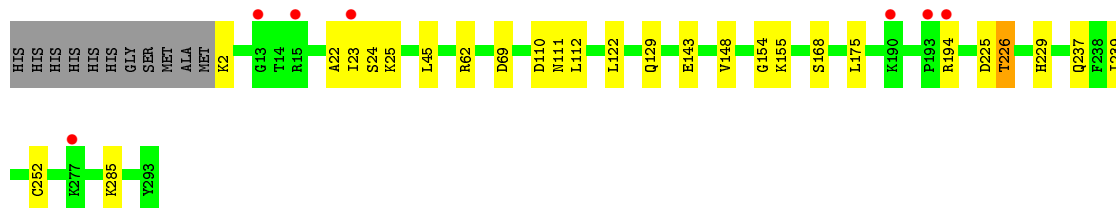
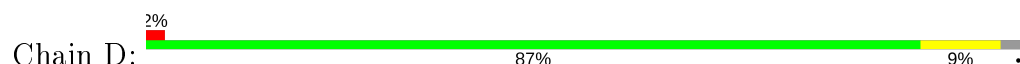
- Molecule 1: GLUCOSE-1-PHOSPHATE THYMIDYLTRANSFERASE



- Molecule 1: GLUCOSE-1-PHOSPHATE THYMIDYLTRANSFERASE



- Molecule 1: GLUCOSE-1-PHOSPHATE THYMIDYLTRANSFERASE



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	64.26 Å 154.35 Å 134.81 Å 90.00° 92.40° 90.00°	Depositor
Resolution (Å)	134.69 – 2.21 27.82 – 2.21	Depositor EDS
% Data completeness (in resolution range)	98.9 (134.69-2.21) 99.0 (27.82-2.21)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.38 (at 2.22 Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.227 , 0.275 0.233 , 0.276	Depositor DCC
R_{free} test set	3245 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	14.8	Xtriage
Anisotropy	0.260	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 41.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.078 for h,-k,-l	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	9879	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.29% 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: KDT, MES, CL

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	0.63	0/2387	0.76	2/3237 (0.1%)
1	B	0.62	0/2344	0.81	5/3180 (0.2%)
1	C	0.61	0/2389	0.80	2/3240 (0.1%)
1	D	0.58	0/2335	0.79	1/3168 (0.0%)
All	All	0.61	0/9455	0.79	10/12825 (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	B	128	ARG	NE-CZ-NH1	6.67	123.64	120.30
1	D	69	ASP	CB-CG-OD1	5.57	123.31	118.30
1	A	204	ARG	NE-CZ-NH1	5.51	123.06	120.30
1	A	23	ILE	CB-CA-C	-5.42	100.77	111.60
1	B	3	ARG	NE-CZ-NH1	5.36	122.98	120.30
1	C	286	ARG	NE-CZ-NH1	5.26	122.93	120.30
1	B	151	ASP	CB-CG-OD1	5.11	122.90	118.30
1	B	225	ASP	CB-CG-OD1	5.11	122.90	118.30
1	C	23	ILE	CB-CA-C	-5.08	101.45	111.60
1	B	232	LEU	CA-CB-CG	5.02	126.84	115.30

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	2331	0	2310	12	0
1	B	2291	0	2279	9	0
1	C	2335	0	2309	12	0
1	D	2285	0	2272	12	0
2	A	28	0	0	0	0
2	B	28	0	0	0	0
2	C	28	0	0	0	0
2	D	28	0	0	1	0
3	A	12	0	13	1	0
3	B	12	0	13	0	0
3	C	12	0	13	1	0
3	D	12	0	13	1	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
5	A	128	0	0	0	0
5	B	112	0	0	1	0
5	C	123	0	0	2	0
5	D	110	0	0	0	0
All	All	9879	0	9222	38	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 2.

All (38) 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:154:GLY:O	3:A:1295:MES:H32	1.93	0.69
1:B:101:ASN:HB2	5:B:2010:HOH:O	1.93	0.67
1:A:226:THR:O	1:A:226:THR:HG22	2.01	0.60
1:C:237:GLN:NE2	1:D:237:GLN:OE1	2.22	0.59
1:C:226:THR:HG22	1:C:226:THR:O	2.04	0.58
5:C:2015:HOH:O	1:D:229:HIS:ND1	2.33	0.56
1:B:25:LYS:HE3	1:B:226:THR:HG22	1.89	0.54
1:B:226:THR:O	1:B:226:THR:HG22	2.09	0.52
1:C:25:LYS:NZ	1:C:110:ASP:OD2	2.33	0.52
1:A:23:ILE:HG23	1:B:23:ILE:HD12	1.92	0.51
1:A:233:LEU:HD11	1:B:237[A]:GLN:HG3	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:25:LYS:NZ	1:D:110:ASP:OD2	2.45	0.48
1:C:27:LEU:HD22	1:D:22:ALA:HB1	1.95	0.48
1:B:54:ILE:HA	1:B:80:ALA:O	2.14	0.48
1:C:214:VAL:O	3:C:1295:MES:H71	2.14	0.47
1:C:35:MET:CE	1:C:226:THR:HG21	2.45	0.47
1:A:24:SER:HB3	1:A:59:ASP:OD2	2.16	0.46
1:C:237:GLN:HE21	1:D:237:GLN:CD	2.11	0.45
1:A:23:ILE:HG23	1:B:23:ILE:CD1	2.46	0.45
1:D:122:LEU:CD2	1:D:175:LEU:HD21	2.47	0.44
1:C:128:ARG:NH1	1:C:133:SER:OG	2.50	0.44
1:A:234:GLU:HA	1:A:237[A]:GLN:HE21	1.82	0.44
1:B:88:LEU:HD13	1:B:108:LEU:HD21	1.99	0.44
1:D:148:VAL:O	1:D:168:SER:OG	2.29	0.42
1:C:41:SER:HB3	1:C:256:ILE:CD1	2.49	0.42
1:D:154:GLY:O	3:D:1295:MES:H32	2.20	0.42
1:D:45:LEU:HD22	2:D:1294:KDT:CAJ	2.49	0.42
1:C:237:GLN:NE2	1:D:237:GLN:HG2	2.35	0.42
1:B:35:MET:O	1:B:39:PRO:HD2	2.20	0.41
1:C:127:GLN:NE2	5:C:2053:HOH:O	2.47	0.41
1:A:225:ASP:HB3	1:A:231:SER:HB2	2.02	0.41
1:A:128:ARG:HB2	1:A:128:ARG:HE	1.64	0.41
1:D:226:THR:O	1:D:226:THR:CG2	2.69	0.41
1:A:56:THR:HB	1:A:58:GLN:OE1	2.21	0.41
1:C:70:GLY:HA3	1:C:77:LEU:HG	2.02	0.41
1:A:55:SER:O	1:A:81:VAL:HA	2.21	0.40
1:A:60:THR:N	1:A:61:PRO:CD	2.83	0.40
1:D:112:LEU:HD21	1:D:239:ILE:HD11	2.02	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	293/303 (97%)	287 (98%)	6 (2%)	0	100	100
1	B	291/303 (96%)	282 (97%)	9 (3%)	0	100	100
1	C	293/303 (97%)	284 (97%)	9 (3%)	0	100	100
1	D	290/303 (96%)	285 (98%)	5 (2%)	0	100	100
All	All	1167/1212 (96%)	1138 (98%)	29 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	244/248 (98%)	236 (97%)	8 (3%)	38	47
1	B	240/248 (97%)	235 (98%)	5 (2%)	53	65
1	C	244/248 (98%)	238 (98%)	6 (2%)	47	58
1	D	239/248 (96%)	226 (95%)	13 (5%)	22	25
All	All	967/992 (98%)	935 (97%)	32 (3%)	38	47

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	-9	HIS
1	A	23	ILE
1	A	24	SER
1	A	111	ASN
1	A	152	GLN
1	A	155	LYS
1	A	194	ARG
1	A	232	LEU
1	B	2	LYS
1	B	62	ARG
1	B	111	ASN
1	B	155	LYS
1	B	232	LEU

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Mol	Chain	Res	Type
1	C	-9	HIS
1	C	2	LYS
1	C	23	ILE
1	C	24	SER
1	C	143	GLU
1	C	155	LYS
1	D	2	LYS
1	D	23	ILE
1	D	24	SER
1	D	62	ARG
1	D	111	ASN
1	D	129	GLN
1	D	143	GLU
1	D	155	LYS
1	D	194	ARG
1	D	225	ASP
1	D	226	THR
1	D	252	CYS
1	D	285	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	-9	HIS
1	B	152	GLN
1	C	-6	HIS
1	D	64	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 carbohydrates in this entry.

5.6 Ligand geometry

Of 12 ligands modelled in this entry, 4 are monoatomic - leaving 8 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	KDT	A	1294	-	28,30,30	2.73	5 (17%)	32,44,44	2.17	8 (25%)
2	KDT	C	1294	-	28,30,30	2.84	5 (17%)	32,44,44	2.41	9 (28%)
2	KDT	B	1294	-	28,30,30	2.85	6 (21%)	32,44,44	2.38	9 (28%)
3	MES	C	1295	-	12,12,12	1.89	1 (8%)	14,16,16	1.85	3 (21%)
3	MES	B	1295	-	12,12,12	2.08	1 (8%)	14,16,16	1.86	5 (35%)
2	KDT	D	1294	-	28,30,30	3.02	7 (25%)	32,44,44	2.45	12 (37%)
3	MES	D	1295	-	12,12,12	1.97	1 (8%)	14,16,16	1.81	2 (14%)
3	MES	A	1295	-	12,12,12	1.97	2 (16%)	14,16,16	2.26	3 (21%)

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	KDT	A	1294	-	-	4/20/20/20	0/3/3/3
2	KDT	C	1294	-	-	3/20/20/20	0/3/3/3
2	KDT	B	1294	-	-	3/20/20/20	0/3/3/3
3	MES	C	1295	-	-	0/6/14/14	0/1/1/1
3	MES	B	1295	-	-	0/6/14/14	0/1/1/1
2	KDT	D	1294	-	-	4/20/20/20	0/3/3/3
3	MES	D	1295	-	-	3/6/14/14	0/1/1/1
3	MES	A	1295	-	-	5/6/14/14	0/1/1/1

All (28) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1294	KDT	OAP-SAH	9.03	1.53	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	1294	KDT	OAP-SAH	8.81	1.53	1.43
2	C	1294	KDT	OAC-SAH	8.51	1.53	1.43
2	D	1294	KDT	OAC-SAH	7.94	1.52	1.43
2	A	1294	KDT	OAC-SAH	7.84	1.52	1.43
2	A	1294	KDT	OAP-SAH	7.37	1.51	1.43
2	B	1294	KDT	CAF-NAG	-7.32	1.35	1.45
2	D	1294	KDT	CAF-NAG	-7.31	1.36	1.45
2	D	1294	KDT	OAP-SAH	6.96	1.51	1.43
2	A	1294	KDT	CAF-NAG	-6.92	1.36	1.45
2	B	1294	KDT	OAC-SAH	6.73	1.51	1.43
3	B	1295	MES	C8-S	-6.57	1.68	1.77
3	D	1295	MES	C8-S	-6.39	1.68	1.77
2	C	1294	KDT	CAF-NAG	-6.29	1.37	1.45
3	C	1295	MES	C8-S	-6.00	1.69	1.77
3	A	1295	MES	C8-S	-5.89	1.69	1.77
2	D	1294	KDT	SAH-NAG	-5.46	1.57	1.64
2	D	1294	KDT	CAE-NAD	5.05	1.41	1.33
2	B	1294	KDT	CAI-SAH	-4.71	1.69	1.76
2	A	1294	KDT	CAE-NAD	4.31	1.40	1.33
2	C	1294	KDT	CAT-CAV	-3.78	1.42	1.51
2	D	1294	KDT	CAT-CAV	-3.51	1.42	1.51
2	B	1294	KDT	CAT-CAV	-3.16	1.43	1.51
2	A	1294	KDT	CAT-CAV	-2.83	1.44	1.51
2	D	1294	KDT	BRBB-CBA	2.57	1.95	1.90
2	C	1294	KDT	CAE-NAD	2.38	1.37	1.33
2	B	1294	KDT	CAE-NAD	2.21	1.36	1.33
3	A	1295	MES	C7-N4	2.05	1.52	1.47

All (51) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1294	KDT	CAE-NAD-CAM	7.32	121.32	115.14
2	C	1294	KDT	OAC-SAH-OAP	-6.49	109.00	119.52
2	D	1294	KDT	CAE-CAF-CAO	6.34	120.54	114.57
3	A	1295	MES	O3S-S-C8	6.11	115.65	105.77
2	A	1294	KDT	CAE-NAD-CAM	6.03	120.23	115.14
2	B	1294	KDT	CAE-NAD-CAM	5.83	120.06	115.14
2	B	1294	KDT	OAC-SAH-OAP	-5.60	110.45	119.52
2	A	1294	KDT	CAE-CAF-CAO	5.35	119.61	114.57
2	B	1294	KDT	OAC-SAH-NAG	5.05	111.89	106.77
2	D	1294	KDT	CAE-NAD-CAM	4.97	119.33	115.14
3	D	1295	MES	O3S-S-C8	4.83	113.58	105.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	1294	KDT	OAC-SAH-OAP	-4.65	111.99	119.52
2	B	1294	KDT	CAE-CAF-CAO	4.62	118.92	114.57
2	B	1294	KDT	OAP-SAH-NAG	4.54	111.38	106.77
3	C	1295	MES	O3S-S-C8	4.50	113.05	105.77
2	D	1294	KDT	OAP-SAH-CAI	-4.45	102.42	108.05
2	C	1294	KDT	CAE-CAF-CAO	4.30	118.63	114.57
2	C	1294	KDT	OAC-SAH-NAG	4.08	110.91	106.77
2	D	1294	KDT	OAP-SAH-NAG	4.05	110.88	106.77
2	A	1294	KDT	OAC-SAH-OAP	-4.02	113.02	119.52
3	B	1295	MES	O1S-S-C8	3.83	111.53	106.92
2	A	1294	KDT	OAP-SAH-NAG	3.82	110.65	106.77
3	A	1295	MES	C2-C3-N4	3.82	115.90	110.10
2	A	1294	KDT	BRBB-CBA-CAX	3.43	124.28	119.30
2	C	1294	KDT	OAC-SAH-CAI	3.17	112.06	108.05
2	B	1294	KDT	OAP-SAH-CAI	-3.17	104.04	108.05
2	D	1294	KDT	OAC-SAH-CAI	3.12	112.00	108.05
2	C	1294	KDT	OAP-SAH-NAG	3.07	109.89	106.77
3	C	1295	MES	O2S-S-C8	-3.04	103.25	106.92
3	A	1295	MES	O1S-S-C8	-3.02	103.28	106.92
3	C	1295	MES	O1S-S-C8	3.01	110.54	106.92
2	D	1294	KDT	CAI-SAH-NAG	2.74	110.54	106.63
2	C	1294	KDT	BRBB-CBA-CAX	2.73	123.27	119.30
2	A	1294	KDT	CAB-NAG-SAH	2.73	121.44	116.83
2	D	1294	KDT	CAB-NAG-SAH	2.73	121.44	116.83
3	B	1295	MES	O3S-S-C8	2.67	110.09	105.77
2	D	1294	KDT	CAV-CAT-NAN	2.65	116.74	112.63
3	B	1295	MES	C2-C3-N4	2.64	114.11	110.10
3	D	1295	MES	C2-C3-N4	2.60	114.05	110.10
2	D	1294	KDT	CAF-CAE-NAD	-2.38	118.25	123.14
2	A	1294	KDT	OAC-SAH-NAG	2.36	109.16	106.77
2	D	1294	KDT	BRBB-CBA-CAX	2.34	122.69	119.30
2	A	1294	KDT	CAV-CAT-NAN	2.33	116.24	112.63
2	B	1294	KDT	NAU-CAO-NAN	2.22	123.12	120.12
2	B	1294	KDT	CAT-CAV-CAY	2.20	124.58	120.40
3	B	1295	MES	C6-O1-C2	2.17	117.13	109.89
2	D	1294	KDT	OAC-SAH-NAG	2.13	108.93	106.77
3	B	1295	MES	O1-C2-C3	2.11	116.44	111.80
2	C	1294	KDT	CAB-NAG-SAH	2.10	120.38	116.83
2	B	1294	KDT	OAC-SAH-CAI	2.10	110.70	108.05
2	C	1294	KDT	CAF-CAE-NAD	-2.08	118.87	123.14

There are no chirality outliers.

All (22) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1294	KDT	CAB-NAG-SAH-OAP
2	A	1294	KDT	CAB-NAG-SAH-OAC
2	C	1294	KDT	CAB-NAG-SAH-OAP
2	C	1294	KDT	CAB-NAG-SAH-OAC
2	D	1294	KDT	CAB-NAG-SAH-OAP
2	D	1294	KDT	CAB-NAG-SAH-OAC
3	A	1295	MES	C8-C7-N4-C5
3	D	1295	MES	C7-C8-S-O3S
3	A	1295	MES	C7-C8-S-O3S
2	B	1294	KDT	CAB-NAG-SAH-OAP
3	A	1295	MES	C8-C7-N4-C3
3	D	1295	MES	C7-C8-S-O1S
3	D	1295	MES	C7-C8-S-O2S
3	A	1295	MES	C7-C8-S-O1S
3	A	1295	MES	C7-C8-S-O2S
2	D	1294	KDT	NAN-CAT-CAV-CAW
2	D	1294	KDT	NAN-CAT-CAV-CAY
2	B	1294	KDT	NAN-CAT-CAV-CAW
2	B	1294	KDT	NAN-CAT-CAV-CAY
2	A	1294	KDT	NAN-CAT-CAV-CAW
2	A	1294	KDT	NAN-CAT-CAV-CAY
2	C	1294	KDT	NAN-CAT-CAV-CAW

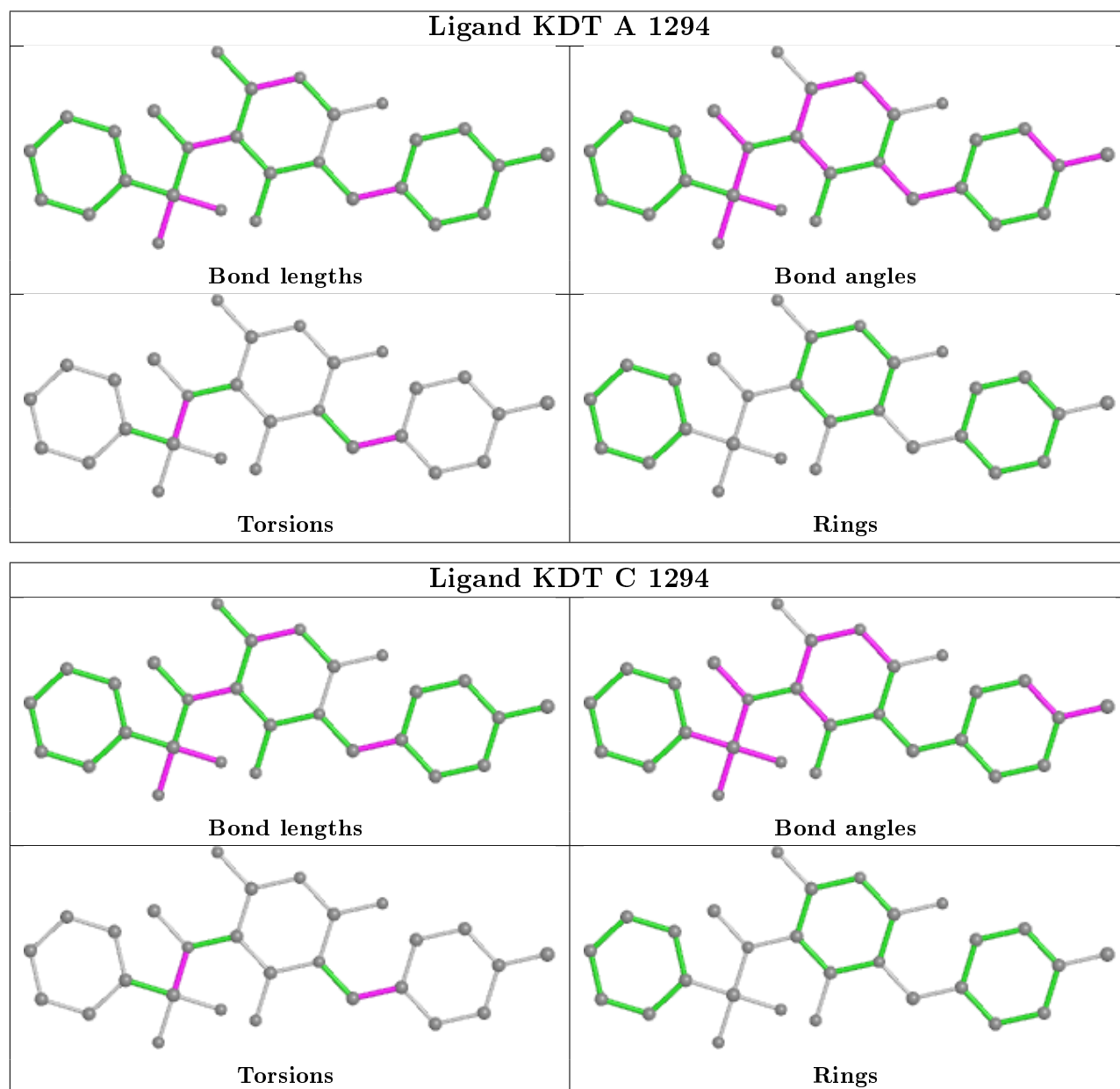
There are no ring outliers.

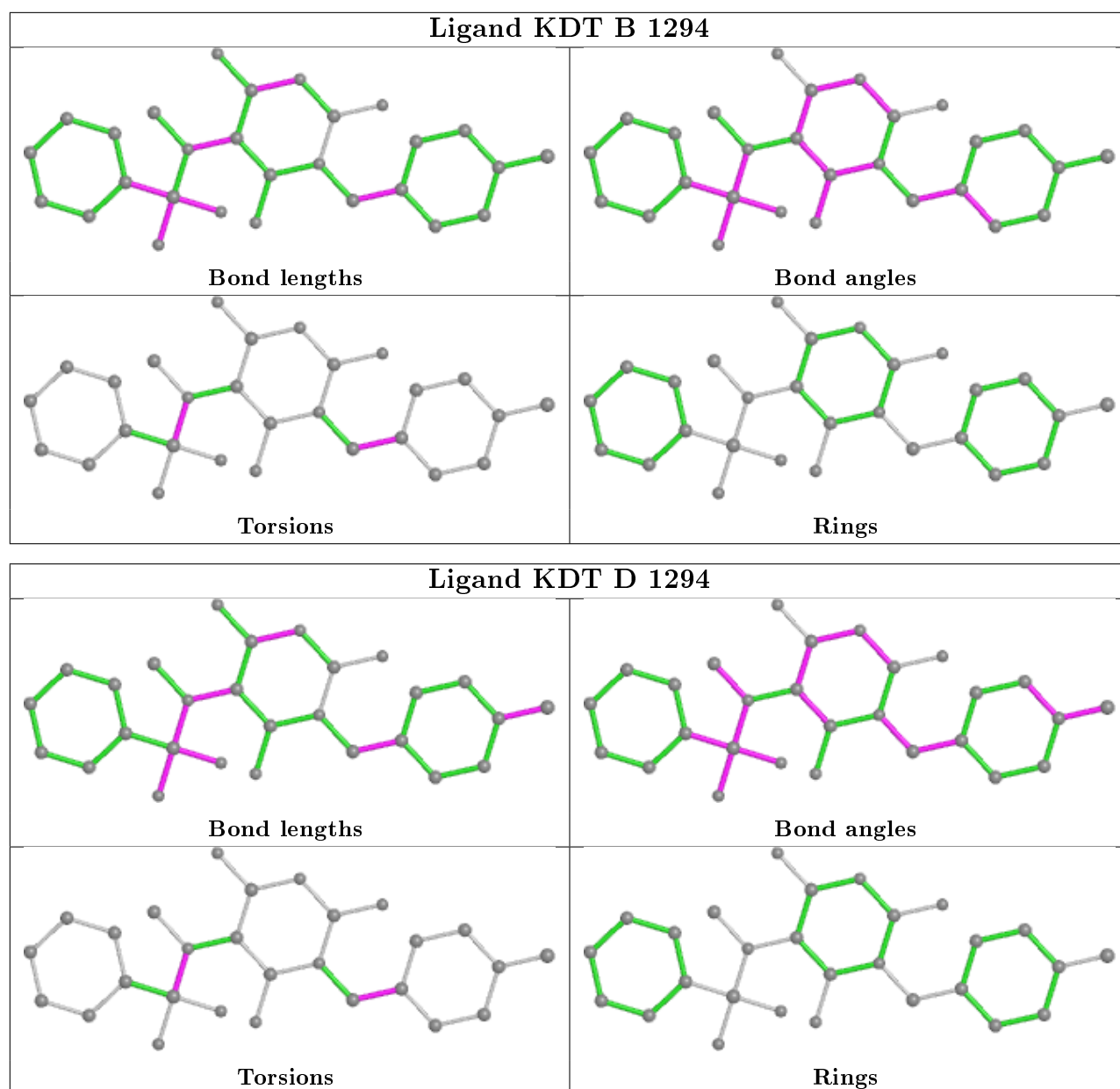
4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	1295	MES	1	0
2	D	1294	KDT	1	0
3	D	1295	MES	1	0
3	A	1295	MES	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, 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	296/303 (97%)	-0.31	4 (1%) 75 73	10, 21, 41, 78	0
1	B	292/303 (96%)	-0.17	3 (1%) 82 81	14, 24, 42, 63	0
1	C	297/303 (98%)	-0.06	18 (6%) 21 19	10, 22, 62, 94	0
1	D	292/303 (96%)	-0.13	7 (2%) 59 57	14, 26, 49, 85	0
All	All	1177/1212 (97%)	-0.17	32 (2%) 54 52	10, 23, 49, 94	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	23	ILE	4.5
1	C	189	LEU	4.5
1	C	15	ARG	3.5
1	B	193	PRO	3.4
1	D	190	LYS	3.4
1	D	194	ARG	3.3
1	C	13	GLY	3.1
1	C	14	THR	3.1
1	C	195	GLY	3.1
1	D	193	PRO	3.1
1	C	187	ARG	2.9
1	C	191	PRO	2.9
1	B	275	LEU	2.7
1	C	190	LYS	2.7
1	C	193	PRO	2.7
1	A	277	LYS	2.7
1	C	11	GLY	2.7
1	D	13	GLY	2.7
1	C	194	ARG	2.6
1	C	-9	HIS	2.6
1	D	277	LYS	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	11	GLY	2.6
1	A	193	PRO	2.5
1	C	-6	HIS	2.5
1	C	277	LYS	2.3
1	B	195	GLY	2.3
1	C	-5	HIS	2.3
1	A	194	ARG	2.2
1	C	227	GLY	2.2
1	D	23	ILE	2.2
1	D	15	ARG	2.1
1	C	274	PRO	2.1

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](#)

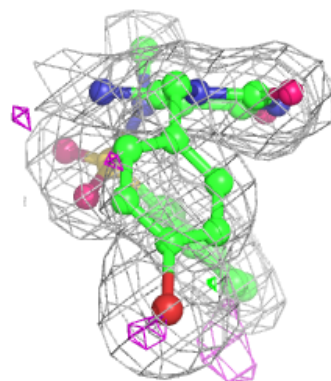
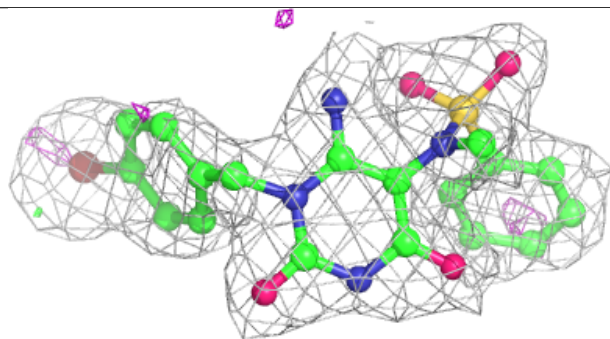
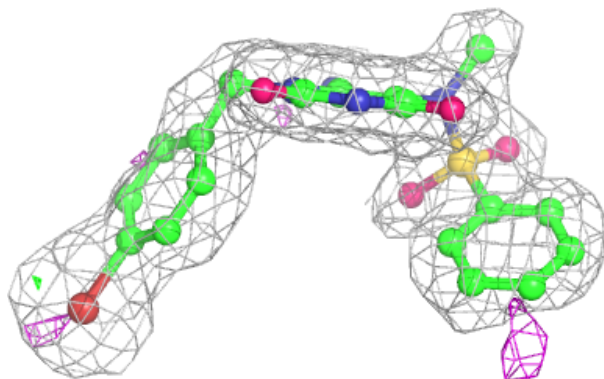
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	MES	B	1295	12/12	0.89	0.18	42,49,51,53	0
3	MES	D	1295	12/12	0.91	0.16	42,45,50,52	0
4	CL	D	1296	1/1	0.93	0.08	38,38,38,38	0
3	MES	A	1295	12/12	0.94	0.14	36,44,45,45	0
4	CL	C	1296	1/1	0.96	0.12	37,37,37,37	0
3	MES	C	1295	12/12	0.96	0.13	30,36,37,39	0
4	CL	B	1296	1/1	0.97	0.05	34,34,34,34	0
2	KDT	B	1294	28/28	0.98	0.10	15,18,28,29	0
2	KDT	A	1294	28/28	0.98	0.09	13,14,21,22	0
2	KDT	C	1294	28/28	0.98	0.10	12,13,19,21	0
2	KDT	D	1294	28/28	0.98	0.09	18,20,28,30	0
4	CL	A	1296	1/1	0.99	0.09	28,28,28,28	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.

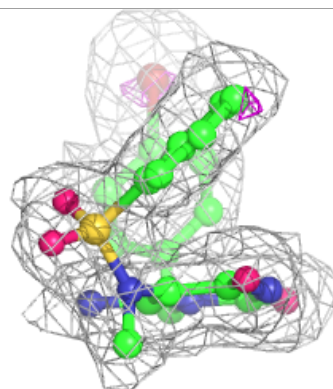
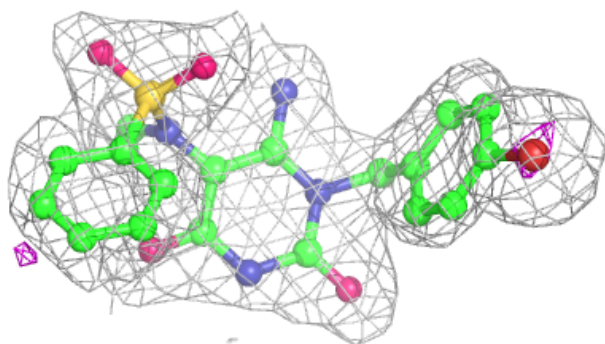
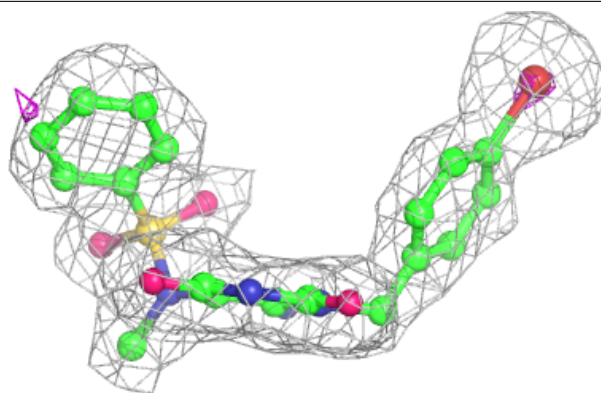
Electron density around KDT B 1294:

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

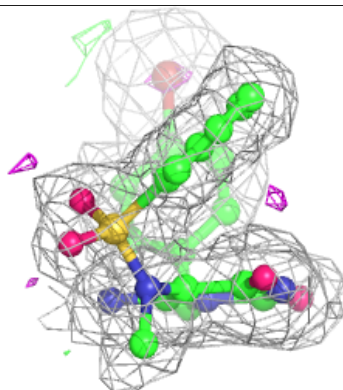
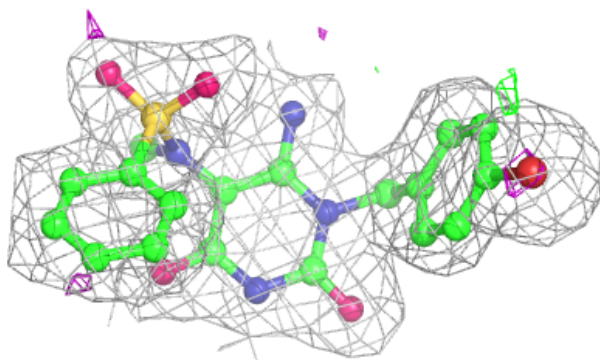
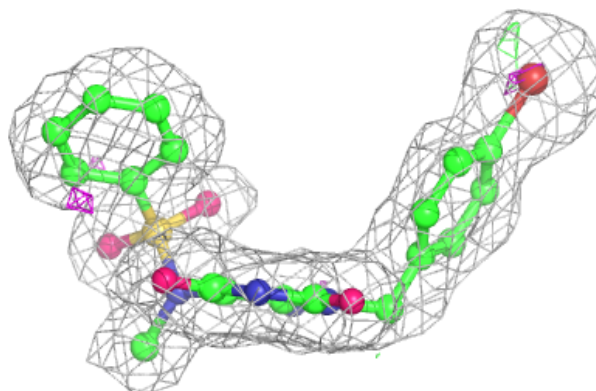


Electron density around KDT A 1294:

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

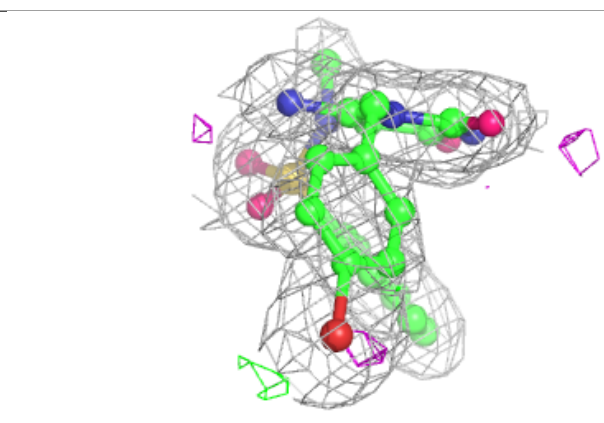
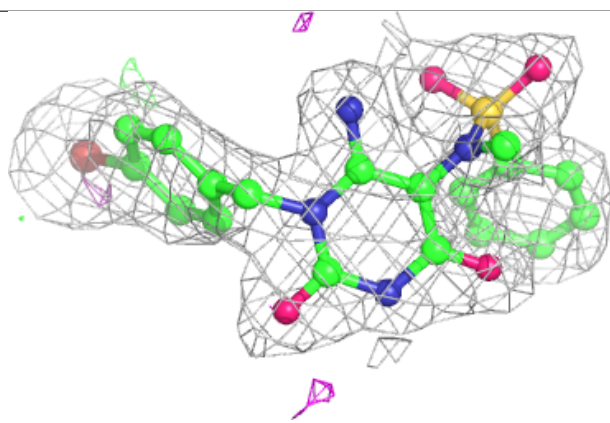
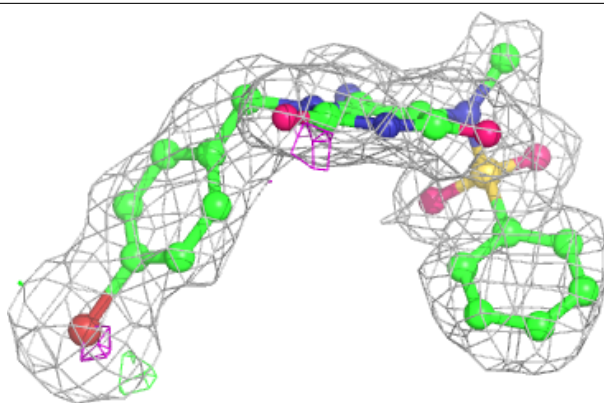
**Electron density around KDT C 1294:**

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



Electron density around KDT D 1294:

$2mF_o-DF_c$ (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.