Full wwPDB X-ray Structure Validation Report (i)

Jan 2, 2024 – 11:39 pm GMT

PDB ID	:	5F3G
Title	:	Crystal structure of human KDM4A in complex with compound 53a
Authors	:	Le Bihan, Y.-V.; Westwood, I.M.; van Montfort, R.L.M.
Deposited on	:	2015-12-02
Resolution	:	2.50 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 (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (i)) were used in the production of this report:

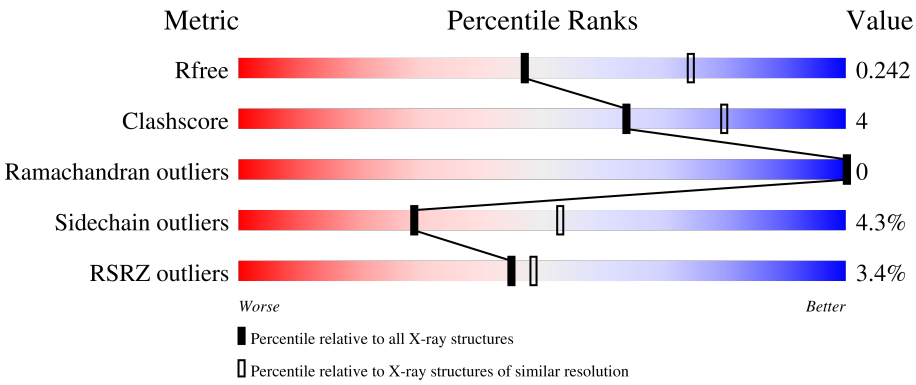
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	360	2% 81% 13% 6%
1	B	360	3% 84% 9% 7%
1	C	360	4% 81% 12% • 6%
1	D	360	4% 82% 12% • 6%

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 11155 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 Lysine-specific demethylase 4A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	339	2717	1765	448	488	16	0	1	0
1	B	335	2628	1707	435	472	14	0	2	0
1	C	337	2687	1753	436	483	15	0	2	0
1	D	339	2629	1717	422	476	14	0	0	0

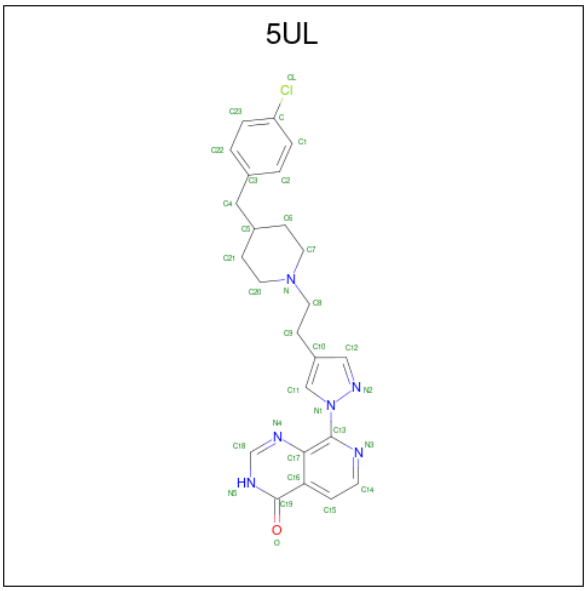
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	expression tag	UNP O75164
B	0	SER	-	expression tag	UNP O75164
C	0	SER	-	expression tag	UNP O75164
D	0	SER	-	expression tag	UNP O75164

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

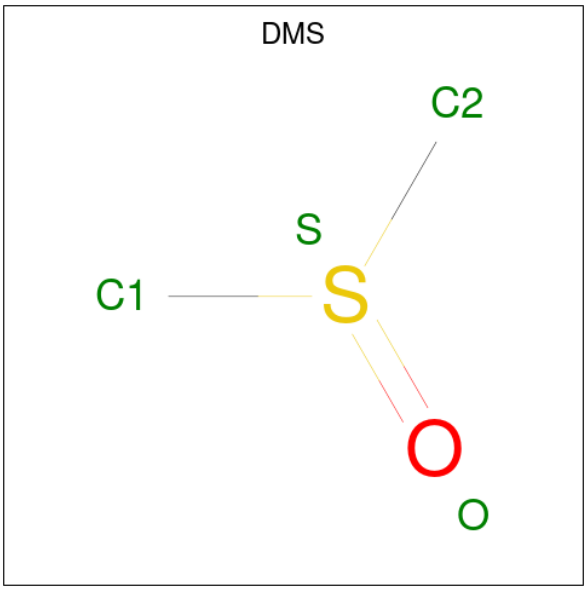
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Zn		
2	A	2	2	2	0	0
2	B	2	2	2	0	0
2	C	2	2	2	0	0
2	D	2	2	2	0	0

- Molecule 3 is 8-[4-[2-[4-[(4-chlorophenyl)methyl]piperidin-1-yl]ethyl]pyrazol-1-yl]-3 {H}-pyrido[3,4-d]pyrimidin-4-one (three-letter code: 5UL) (formula: $C_{24}H_{25}ClN_6O$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C Cl N O 32 24 1 6 1	0	0
3	B	1	Total C N O 21 14 6 1	0	0
3	C	1	Total C N O 25 18 6 1	0	0
3	D	1	Total C N O 16 10 5 1	0	0

- Molecule 4 is DIMETHYL SULFOXIDE (three-letter code: DMS) (formula: C_2H_6OS).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	O	S		
4	A	1	4	2	1	1	0	0
			Total	C	O	S		
4	A	1	4	2	1	1	0	0
			Total	C	O	S		
4	C	1	4	2	1	1	0	0
			Total	C	O	S		
4	C	1	4	2	1	1	0	0
			Total	C	O	S		
4	C	1	4	2	1	1	0	0
			Total	C	O	S		
4	D	1	4	2	1	1	0	0
			Total	C	O	S		
4	D	1	4	2	1	1	0	0

- Molecule 5 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Cl		
5	D	1	1	1	0	0

- Molecule 6 is water.

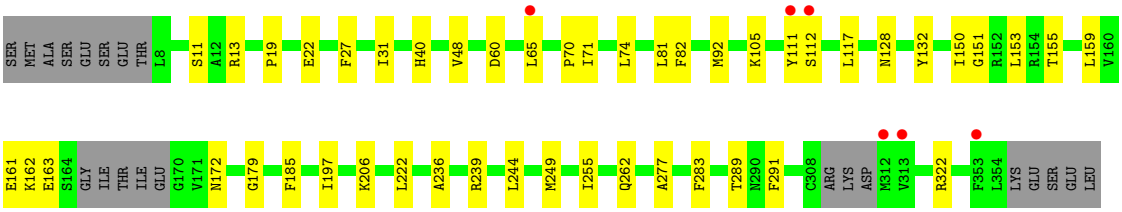
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
6	A	90	90	90	0	0
			Total	O		
6	B	107	107	107	0	0
			Total	O		
6	C	84	84	84	0	0
			Total	O		
6	D	82	82	82	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

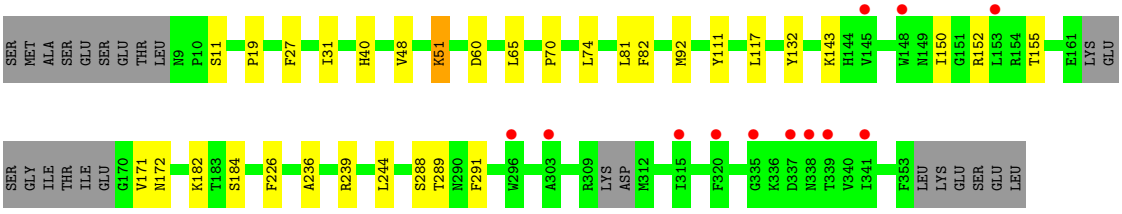
• Molecule 1: Lysine-specific demethylase 4A

Chain A:



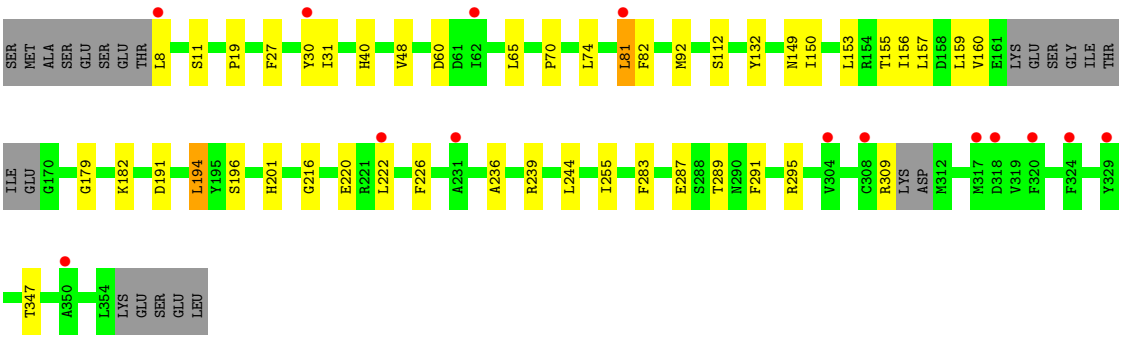
• Molecule 1: Lysine-specific demethylase 4A

Chain B:



• Molecule 1: Lysine-specific demethylase 4A

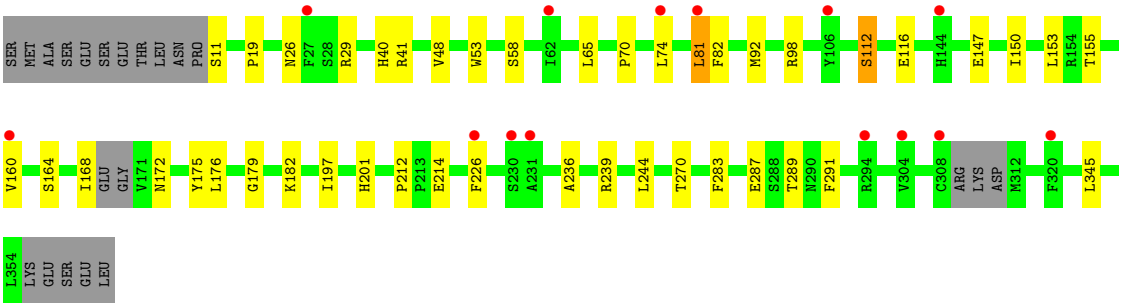
Chain C:



• Molecule 1: Lysine-specific demethylase 4A

Chain D:

4%
82%
12%
6%


SER MET ALA SER GLU SER GLU THR LEU ASN PRO S11 P19 N26 F27 S28 R29 H40 R41 V48 W53 S58 I62 L65 P70 L74 L81 F82 M92 R98 Y106 S112 E116 H144 E147 I150 L153 R154 T155
V160 S164 I168 GLU GLY V171 N172 Y175 L176 G179 K182 I197 H201 P212 P213 E214 F226 S230 A231 A236 R239 L244 T270 F283 E287 S288 T289 N290 F291 R294 V304 C308 ARG LYS ASP M312 F320 L345
L354 LYS GLU SER GLU LEU

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	57.27Å 100.57Å 142.59Å 90.00° 99.43° 90.00°	Depositor
Resolution (Å)	42.50 – 2.50 42.50 – 2.50	Depositor EDS
% Data completeness (in resolution range)	97.7 (42.50-2.50) 88.2 (42.50-2.50)	Depositor EDS
R_{merge}	0.20	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.54 (at 2.51Å)	Xtriage
Refinement program	BUSTER 2.10.2	Depositor
R, R_{free}	0.203 , 0.246 0.207 , 0.242	Depositor DCC
R_{free} test set	2576 reflections (4.77%)	wwPDB-VP
Wilson B-factor (Å^2)	42.1	Xtriage
Anisotropy	0.400	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 54.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.39$, $< L^2 > = 0.22$	Xtriage
Estimated twinning fraction	0.136 for h,-k,-h-l	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	11155	wwPDB-VP
Average B, all atoms (Å^2)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 18.46% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: 5UL, CL, DMS, ZN

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	0.53	0/2802	0.68	0/3809
1	B	0.53	0/2713	0.67	0/3702
1	C	0.55	0/2772	0.68	0/3772
1	D	0.53	0/2713	0.69	0/3702
All	All	0.54	0/11000	0.68	0/14985

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	2717	0	2532	22	0
1	B	2628	0	2354	14	0
1	C	2687	0	2481	20	0
1	D	2629	0	2372	21	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
3	A	32	0	25	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	21	0	12	3	0
3	C	25	0	19	2	0
3	D	16	0	6	2	0
4	A	8	0	12	0	0
4	C	12	0	18	0	0
4	D	8	0	12	0	0
5	D	1	0	0	0	0
6	A	90	0	0	3	0
6	B	107	0	0	1	0
6	C	84	0	0	1	0
6	D	82	0	0	3	0
All	All	11155	0	9843	86	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 4.

All (86) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:51:LYS:H	1:B:51:LYS:HD2	1.17	1.06
1:A:161:GLU:OE2	1:A:172:ASN:HB2	1.63	0.97
3:D:403:5UL:N4	3:D:403:5UL:H9	2.07	0.70
3:B:403:5UL:H10	6:B:537:HOH:O	1.91	0.69
1:D:92:MET:HE2	6:D:506:HOH:O	1.92	0.69
1:A:161:GLU:OE2	1:A:172:ASN:CB	2.40	0.69
1:C:70:PRO:HG3	1:C:92:MET:HG2	1.75	0.67
3:A:403:5UL:N4	3:A:403:5UL:H9	2.13	0.64
3:B:403:5UL:H9	3:B:403:5UL:N4	2.13	0.62
3:C:403:5UL:N4	3:C:403:5UL:H9	2.13	0.62
1:B:51:LYS:H	1:B:51:LYS:CD	1.99	0.62
1:C:150:ILE:HG23	1:C:289:THR:HG22	1.82	0.62
1:B:51:LYS:HD2	1:B:51:LYS:N	2.02	0.60
1:D:214:GLU:HG2	6:D:502:HOH:O	2.02	0.59
1:C:19:PRO:HB3	1:C:30:TYR:CZ	2.37	0.59
1:A:111:TYR:CD1	1:A:117:LEU:HD13	2.39	0.57
1:A:82:PHE:HB2	1:A:244:LEU:HB2	1.85	0.57
1:B:82:PHE:HB2	1:B:244:LEU:HB2	1.86	0.57
1:C:82:PHE:HB2	1:C:244:LEU:HB2	1.86	0.57
1:D:82:PHE:HB2	1:D:244:LEU:HB2	1.86	0.57
1:B:111:TYR:CD1	1:B:117:LEU:HD13	2.40	0.56
3:D:403:5UL:N4	3:D:403:5UL:C11	2.68	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:236:ALA:HB1	1:B:239:ARG:HG3	1.89	0.55
1:B:27:PHE:O	1:B:31:ILE:HG12	2.07	0.55
1:C:160:VAL:O	1:C:160:VAL:HG12	2.06	0.55
3:A:403:5UL:N4	3:A:403:5UL:C11	2.70	0.55
1:C:149:ASN:O	1:C:153:LEU:HD23	2.07	0.54
1:A:236:ALA:HB1	1:A:239:ARG:HG3	1.89	0.54
3:B:403:5UL:N4	3:B:403:5UL:C11	2.72	0.53
3:C:403:5UL:N4	3:C:403:5UL:C11	2.71	0.53
1:D:70:PRO:HG3	1:D:92:MET:HG2	1.91	0.53
1:B:150:ILE:HG23	1:B:289:THR:HG22	1.92	0.52
1:A:22:GLU:HB3	6:A:523:HOH:O	2.10	0.52
1:C:19:PRO:HD2	1:C:48:VAL:O	2.10	0.52
1:B:70:PRO:HG3	1:B:92:MET:HG2	1.93	0.51
1:D:236:ALA:HB1	1:D:239:ARG:HG3	1.94	0.50
1:A:185:PHE:O	1:A:277:ALA:HB1	2.12	0.49
1:C:236:ALA:HB1	1:C:239:ARG:HG3	1.94	0.49
1:D:150:ILE:HG23	1:D:289:THR:HG22	1.94	0.49
1:A:206:LYS:HD2	1:A:262:GLN:OE1	2.13	0.48
1:D:26:ASN:HB3	1:D:29:ARG:HB3	1.94	0.48
1:D:112:SER:OG	1:D:116:GLU:OE1	2.26	0.48
1:C:309:ARG:HD2	6:C:572:HOH:O	2.13	0.48
1:C:155:THR:HG21	1:C:291:PHE:HB2	1.94	0.48
1:B:155:THR:HG21	1:B:291:PHE:HB2	1.95	0.48
1:D:155:THR:HG21	1:D:291:PHE:HB2	1.95	0.47
1:D:19:PRO:HD2	1:D:48:VAL:O	2.15	0.47
1:A:70:PRO:HG3	1:A:92:MET:HG2	1.96	0.47
1:A:19:PRO:HD2	1:A:48:VAL:O	2.15	0.46
1:D:81:LEU:HD11	1:D:226:PHE:CG	2.52	0.45
1:B:19:PRO:HD2	1:B:48:VAL:O	2.17	0.45
1:D:176:LEU:HD23	1:D:287:GLU:HG3	1.97	0.45
1:A:40:HIS:H	1:A:40:HIS:CD2	2.33	0.45
1:D:179:GLY:O	1:D:283:PHE:HA	2.17	0.45
1:C:216:GLY:O	1:C:220:GLU:HG3	2.17	0.44
1:A:27:PHE:O	1:A:31:ILE:HG12	2.17	0.44
1:A:153:LEU:HD11	1:A:197:ILE:HG21	2.00	0.44
1:B:143:LYS:O	1:B:152:ARG:NH1	2.44	0.44
1:A:27:PHE:CD2	1:A:153:LEU:HD22	2.52	0.44
1:D:153:LEU:HD11	1:D:197:ILE:HG21	2.00	0.44
1:A:155:THR:HG21	1:A:291:PHE:HB2	2.00	0.43
1:A:159:LEU:HD23	1:A:162:LYS:HD3	2.01	0.43
1:D:40:HIS:CD2	1:D:40:HIS:H	2.34	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:81:LEU:HD11	1:C:226:PHE:CG	2.53	0.43
1:C:27:PHE:O	1:C:31:ILE:HG12	2.18	0.43
1:B:81:LEU:HD11	1:B:226:PHE:CG	2.54	0.42
1:B:40:HIS:CD2	1:B:40:HIS:H	2.36	0.42
1:A:249:MET:HE2	6:A:577:HOH:O	2.18	0.42
1:D:160:VAL:O	1:D:164:SER:HB2	2.19	0.42
1:A:179:GLY:O	1:A:283:PHE:HA	2.20	0.42
1:C:222:LEU:HD22	1:C:255:ILE:HG13	2.02	0.42
1:A:150:ILE:HG23	1:A:289:THR:HG22	2.01	0.42
1:C:179:GLY:O	1:C:283:PHE:HA	2.20	0.42
1:C:156:ILE:O	1:C:159:LEU:HG	2.19	0.41
1:D:41:ARG:O	1:D:212:PRO:HB3	2.20	0.41
1:C:40:HIS:CD2	1:C:40:HIS:H	2.37	0.41
1:A:105:LYS:O	1:A:128:ASN:ND2	2.46	0.41
1:A:222:LEU:HD22	1:A:255:ILE:HG13	2.03	0.41
1:C:201:HIS:HE2	1:C:287:GLU:HB2	1.85	0.41
1:D:201:HIS:HE2	1:D:287:GLU:HB2	1.86	0.41
1:D:345:LEU:HD22	6:D:572:HOH:O	2.19	0.41
1:A:151:GLY:HA3	6:A:525:HOH:O	2.20	0.41
1:D:53:TRP:NE1	1:D:147:GLU:HB2	2.36	0.41
1:C:295:ARG:HB2	1:C:347:THR:HA	2.03	0.40
1:C:157:LEU:HD13	1:C:194:LEU:HD21	2.03	0.40
1:D:168:ILE:H	1:D:172:ASN:HD21	1.69	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	334/360 (93%)	327 (98%)	7 (2%)	0	100	100
1	B	331/360 (92%)	325 (98%)	6 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	333/360 (92%)	322 (97%)	11 (3%)	0	100	100
1	D	333/360 (92%)	322 (97%)	11 (3%)	0	100	100
All	All	1331/1440 (92%)	1296 (97%)	35 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	272/316 (86%)	261 (96%)	11 (4%)	31	56
1	B	246/316 (78%)	235 (96%)	11 (4%)	27	51
1	C	263/316 (83%)	251 (95%)	12 (5%)	27	50
1	D	250/316 (79%)	240 (96%)	10 (4%)	31	56
All	All	1031/1264 (82%)	987 (96%)	44 (4%)	29	53

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

Mol	Chain	Res	Type
1	A	11	SER
1	A	13	ARG
1	A	60	ASP
1	A	65	LEU
1	A	71	ILE
1	A	74	LEU
1	A	81	LEU
1	A	112	SER
1	A	132	TYR
1	A	163	GLU
1	A	322	ARG
1	B	11	SER
1	B	51	LYS
1	B	60	ASP

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Mol	Chain	Res	Type
1	B	65	LEU
1	B	74	LEU
1	B	132	TYR
1	B	171	VAL
1	B	172	ASN
1	B	182	LYS
1	B	184	SER
1	B	288	SER
1	C	8	LEU
1	C	11	SER
1	C	60	ASP
1	C	65	LEU
1	C	74	LEU
1	C	81	LEU
1	C	112	SER
1	C	132	TYR
1	C	182	LYS
1	C	191	ASP
1	C	194	LEU
1	C	196	SER
1	D	11	SER
1	D	58	SER
1	D	65	LEU
1	D	74	LEU
1	D	81	LEU
1	D	98	ARG
1	D	112	SER
1	D	175	TYR
1	D	182	LYS
1	D	270	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	40	HIS
1	B	40	HIS
1	B	128	ASN
1	C	40	HIS
1	C	86	ASN
1	D	40	HIS
1	D	172	ASN
1	D	215	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 20 ligands modelled in this entry, 9 are monoatomic - leaving 11 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	$\#\|Z\| > 2$
4	DMS	A	404	-	3,3,3	0.26	0	3,3,3	0.35	0
4	DMS	C	404	-	3,3,3	0.33	0	3,3,3	0.11	0
3	5UL	D	403	2	17,18,36	0.70	0	18,25,50	0.94	1 (5%)
4	DMS	D	405	-	3,3,3	0.15	0	3,3,3	0.54	0
4	DMS	C	405	-	3,3,3	0.28	0	3,3,3	0.19	0
3	5UL	B	403	2	22,23,36	0.65	0	22,31,50	0.84	1 (4%)
4	DMS	A	405	-	3,3,3	0.19	0	3,3,3	0.24	0
4	DMS	C	406	-	3,3,3	0.28	0	3,3,3	0.41	0
3	5UL	C	403	2	27,28,36	0.56	0	31,39,50	0.74	1 (3%)
4	DMS	D	404	-	3,3,3	0.26	0	3,3,3	0.23	0
3	5UL	A	403	2	35,36,36	0.52	0	42,50,50	0.64	1 (2%)

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
3	5UL	B	403	2	-	1/5/9/23	0/3/3/5
3	5UL	D	403	2	-	0/0/4/23	0/3/3/5
3	5UL	C	403	2	-	1/5/19/23	0/4/4/5
3	5UL	A	403	2	-	2/9/23/23	0/5/5/5

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	403	5UL	C17-C13-N3	3.23	125.89	118.67
3	C	403	5UL	C17-C13-N3	3.07	125.53	118.67
3	D	403	5UL	C17-C13-N3	3.03	125.45	118.67
3	B	403	5UL	C17-C13-N3	2.83	125.00	118.67

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	403	5UL	C3-C4-C5-C6
3	B	403	5UL	C9-C8-N-C7
3	C	403	5UL	C9-C8-N-C7
3	A	403	5UL	C3-C4-C5-C21

There are no ring outliers.

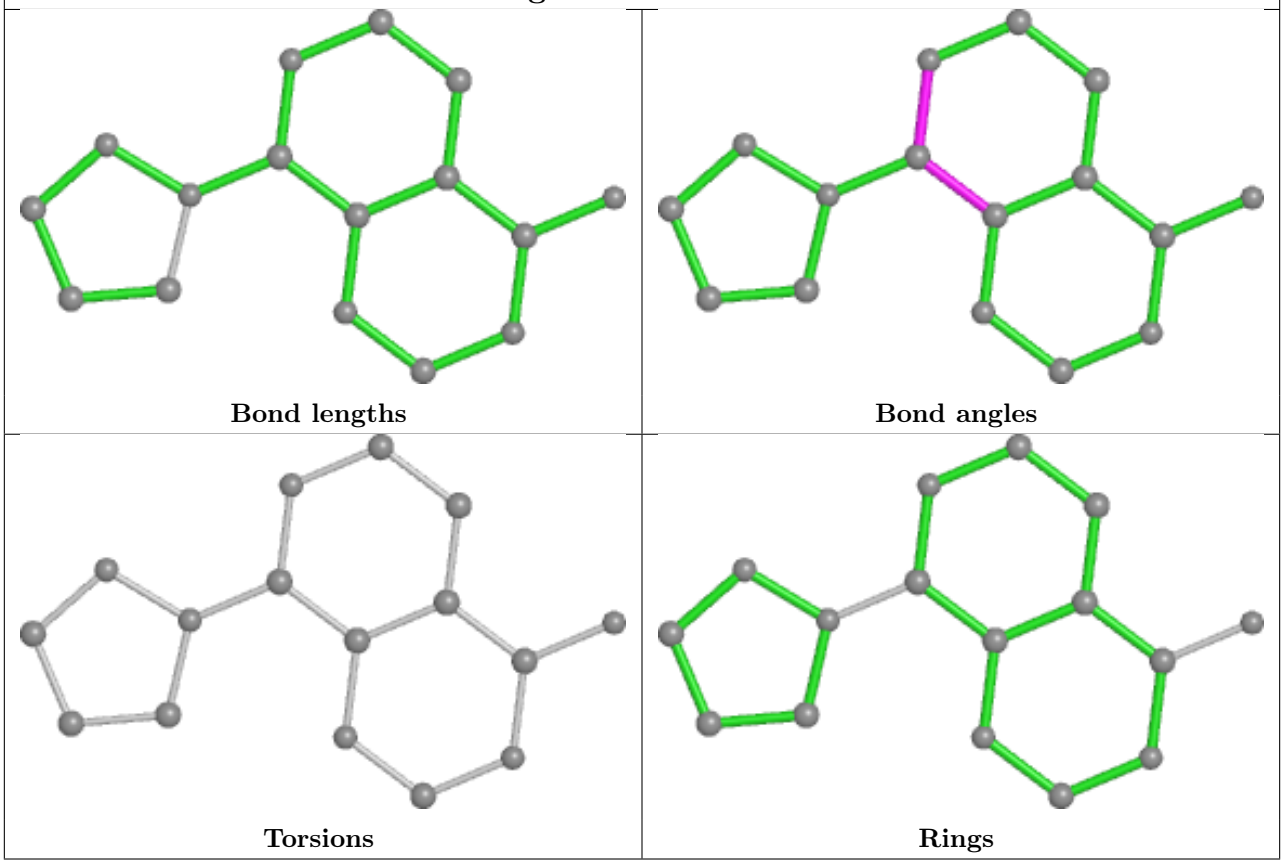
4 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	403	5UL	2	0
3	B	403	5UL	3	0
3	C	403	5UL	2	0
3	A	403	5UL	2	0

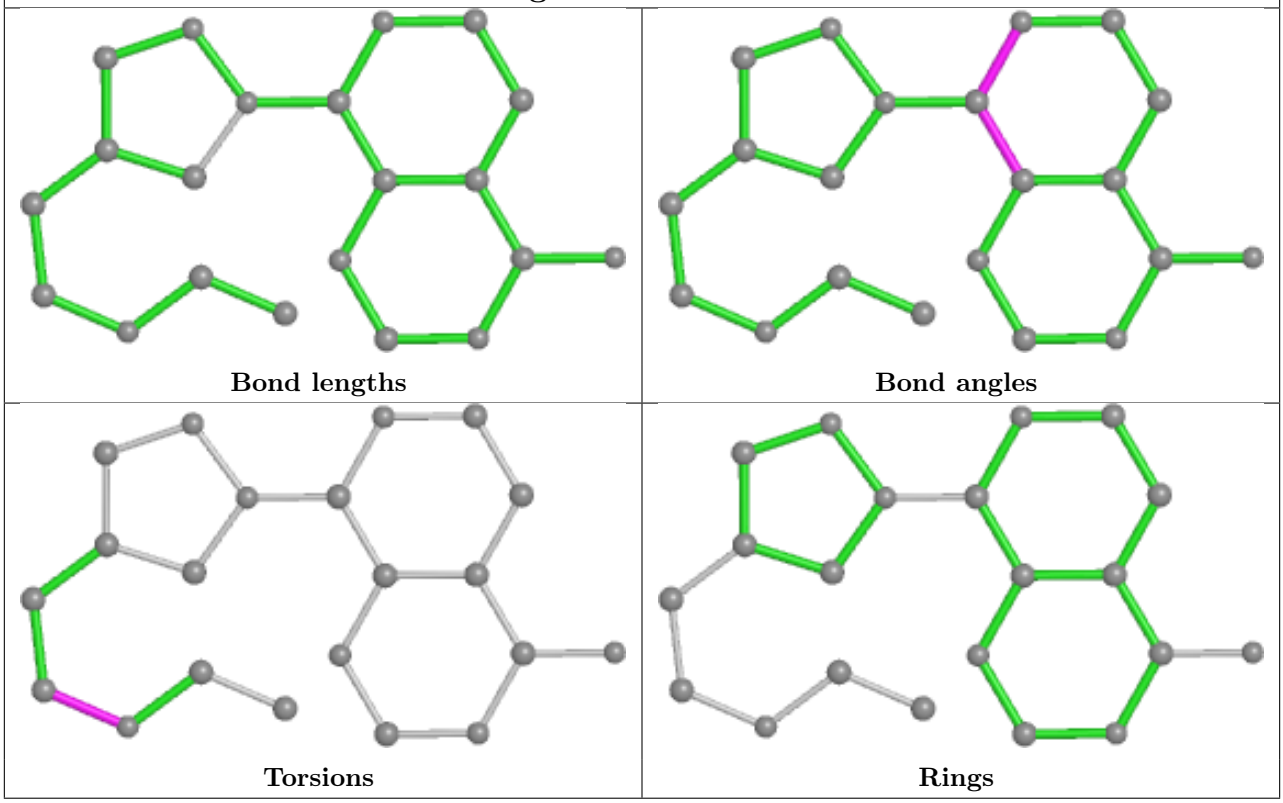
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 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.

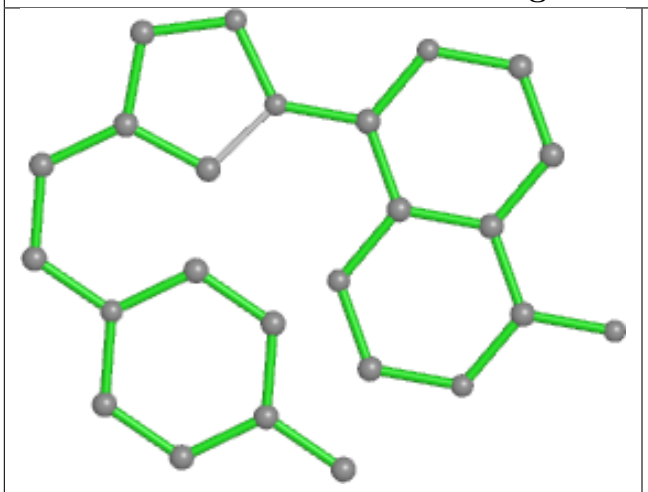
Ligand 5UL D 403



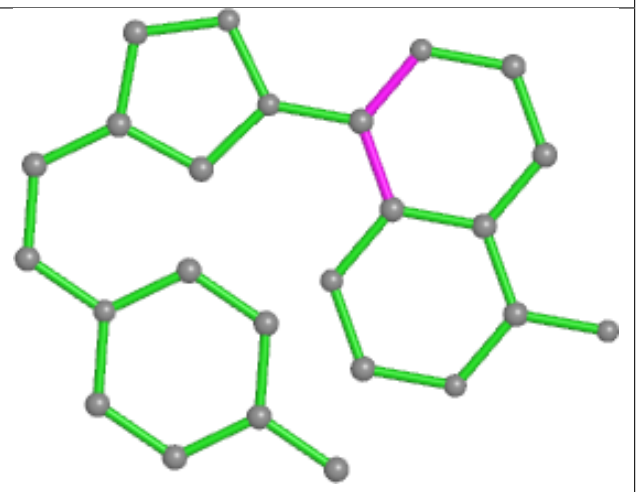
Ligand 5UL B 403



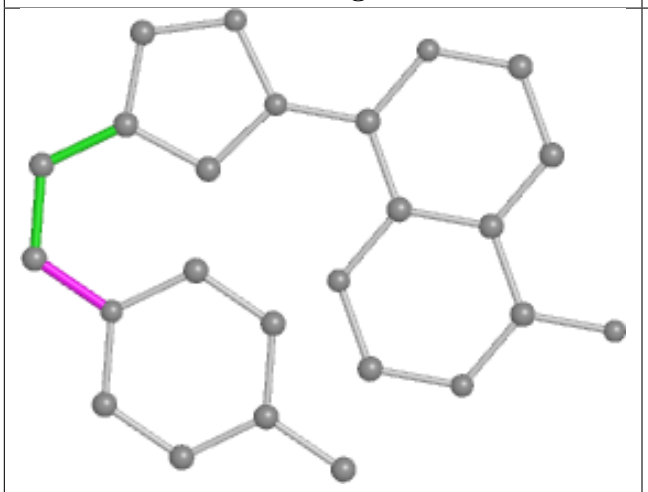
Ligand 5UL C 403



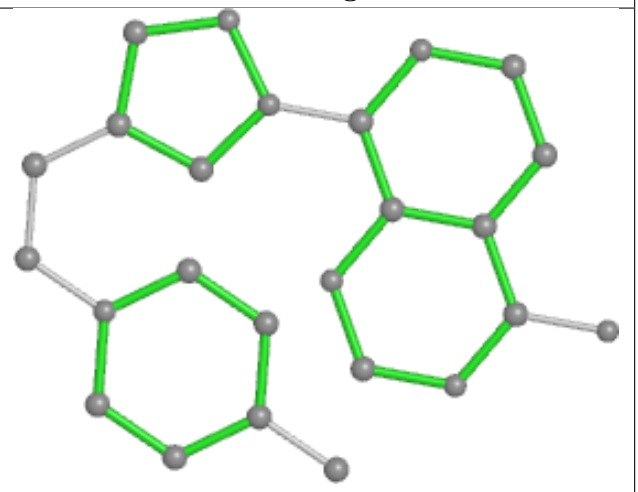
Bond lengths



Bond angles

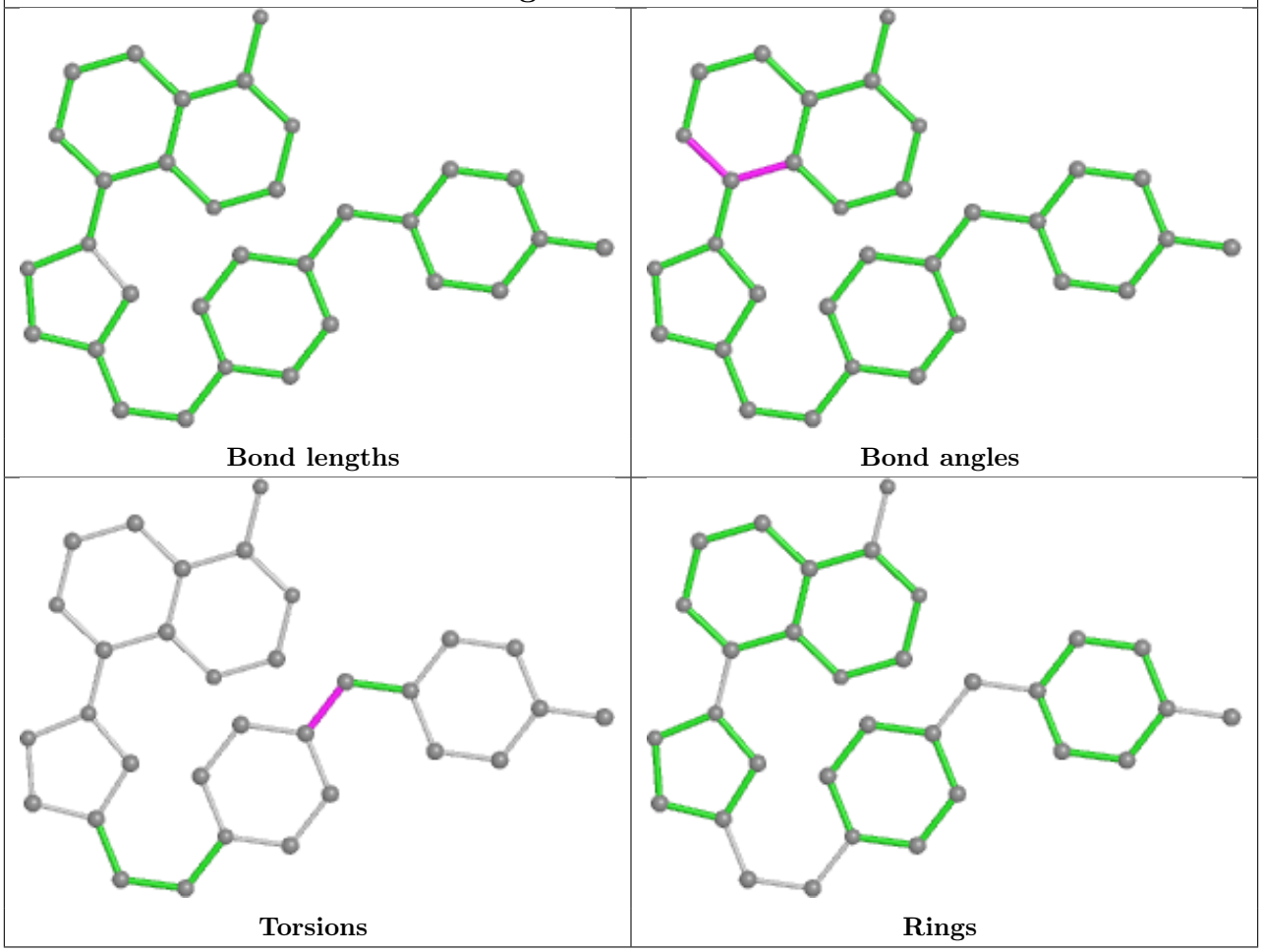


Torsions



Rings

Ligand 5UL A 403



5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	A	339/360 (94%)	0.06	6 (1%)	68	71	26, 46, 74, 97	0
1	B	335/360 (93%)	0.23	12 (3%)	42	46	24, 55, 103, 117	0
1	C	337/360 (93%)	0.13	14 (4%)	36	39	19, 48, 85, 105	0
1	D	339/360 (94%)	0.28	14 (4%)	37	40	20, 56, 92, 117	0
All	All	1350/1440 (93%)	0.17	46 (3%)	45	48	19, 51, 91, 117	0

All (46) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	303	ALA	5.4
1	B	339	THR	4.0
1	D	106	TYR	3.7
1	D	294	ARG	3.5
1	B	145	VAL	3.5
1	A	112	SER	3.5
1	C	317	MET	3.3
1	D	144	HIS	3.3
1	B	337	ASP	3.1
1	B	320	PHE	3.0
1	A	312	MET	3.0
1	C	222	LEU	3.0
1	A	313	VAL	2.7
1	C	8	LEU	2.7
1	C	308	CYS	2.7
1	D	231	ALA	2.7
1	B	153	LEU	2.7
1	D	230	SER	2.6
1	D	308	CYS	2.5
1	D	226	PHE	2.5
1	C	329	TYR	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	338	ASN	2.5
1	D	304	VAL	2.4
1	C	81	LEU	2.4
1	C	318	ASP	2.4
1	D	62	ILE	2.4
1	C	30	TYR	2.4
1	D	160	VAL	2.3
1	D	74	LEU	2.3
1	B	341	ILE	2.3
1	B	315	ILE	2.2
1	C	62	ILE	2.2
1	A	353	PHE	2.2
1	D	27	PHE	2.2
1	B	296	TRP	2.2
1	D	320	PHE	2.2
1	D	81	LEU	2.1
1	B	148	TRP	2.1
1	B	335	GLY	2.1
1	C	320	PHE	2.1
1	C	304	VAL	2.1
1	A	65	LEU	2.1
1	C	231	ALA	2.1
1	C	324	PHE	2.1
1	A	111	TYR	2.0
1	C	350	ALA	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 monosaccharides in this entry.

6.4 Ligands (i)

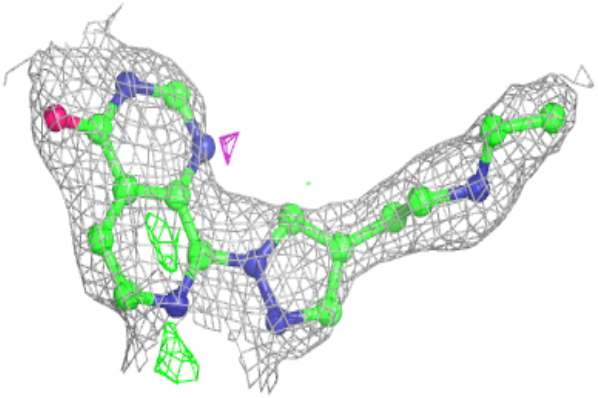
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

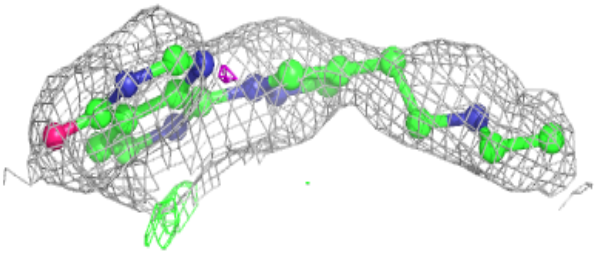
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
4	DMS	D	404	4/4	0.81	0.22	104,105,105,105	0
5	CL	D	406	1/1	0.82	0.14	74,74,74,74	0
4	DMS	C	406	4/4	0.85	0.25	74,74,75,75	4
4	DMS	C	404	4/4	0.86	0.35	44,49,50,50	4
3	5UL	B	403	21/32	0.91	0.17	41,50,52,53	0
4	DMS	A	404	4/4	0.91	0.38	54,55,55,56	4
3	5UL	C	403	25/32	0.93	0.21	27,34,60,60	25
3	5UL	A	403	32/32	0.93	0.21	32,38,61,66	32
2	ZN	B	402	1/1	0.93	0.07	105,105,105,105	0
4	DMS	C	405	4/4	0.94	0.16	69,71,71,72	0
3	5UL	D	403	16/32	0.94	0.20	57,59,66,66	0
4	DMS	D	405	4/4	0.96	0.18	43,44,45,47	4
4	DMS	A	405	4/4	0.96	0.14	58,59,59,60	4
2	ZN	A	402	1/1	0.97	0.10	13,13,13,13	1
2	ZN	D	402	1/1	0.97	0.10	54,54,54,54	1
2	ZN	C	402	1/1	0.99	0.08	55,55,55,55	1
2	ZN	D	401	1/1	0.99	0.13	44,44,44,44	0
2	ZN	B	401	1/1	0.99	0.12	50,50,50,50	0
2	ZN	A	401	1/1	0.99	0.10	42,42,42,42	0
2	ZN	C	401	1/1	0.99	0.16	43,43,43,43	0

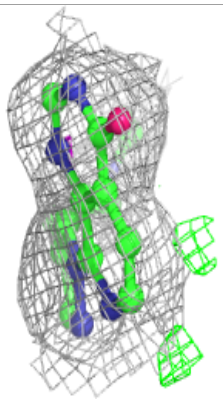
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around 5UL B 403:

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

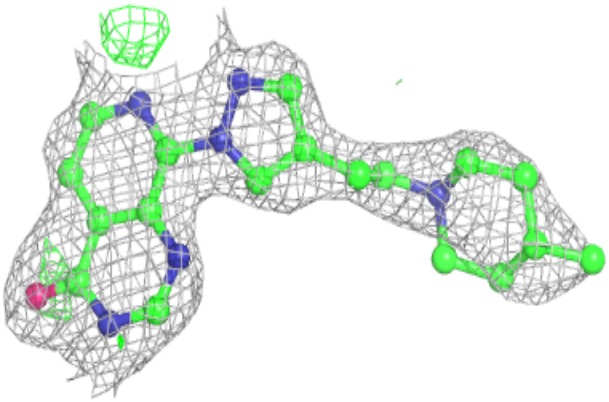


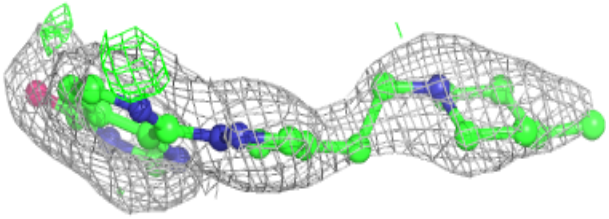


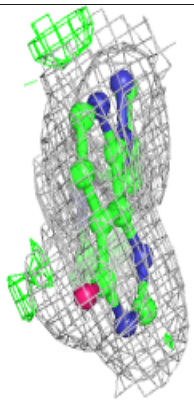


Electron density around 5UL C 403:

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

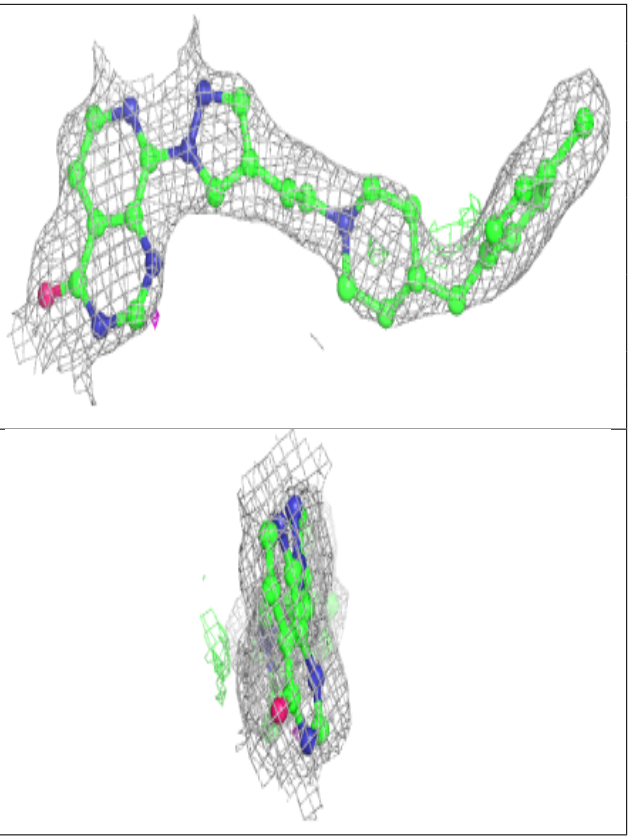






Electron density around 5UL A 403:

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



Electron density around 5UL D 403:

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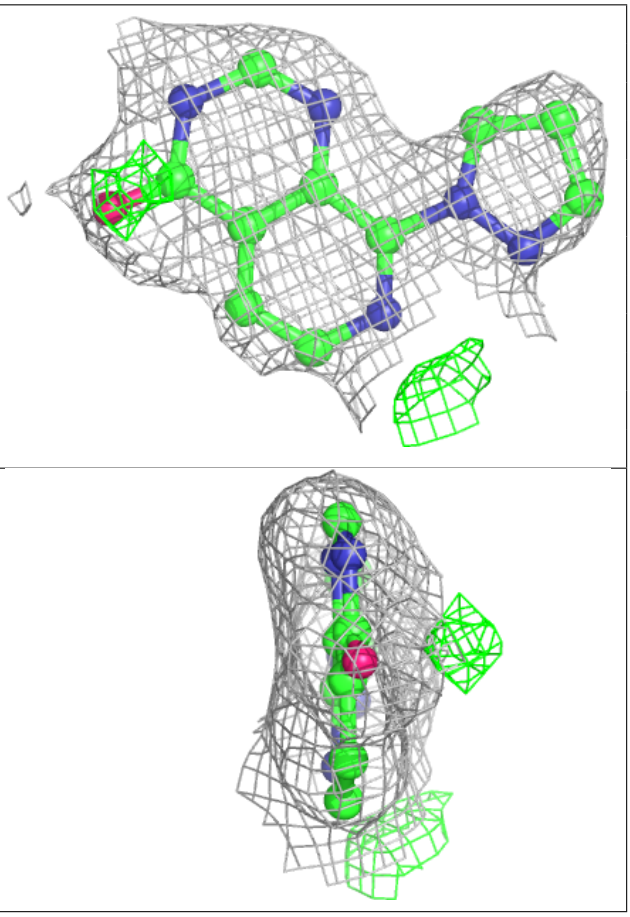
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers (i)

There are no such residues in this entry.