



# Full wwPDB X-ray Structure Validation Report ⓘ

Sep 17, 2020 – 07:21 PM BST

PDB ID : 6YKY  
Title : Biochemical, Cellular and Structural Characterization of Novel ERK3 Inhibitors  
Authors : Graedler, U.  
Deposited on : 2020-04-06  
Resolution : 2.52 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.14.5  
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) : 2.14.5

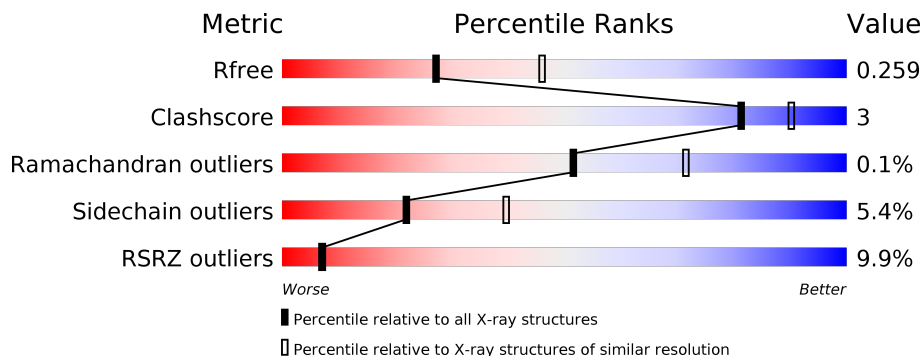
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.52 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	5743 (2.54-2.50)
Clashscore	141614	6463 (2.54-2.50)
Ramachandran outliers	138981	6335 (2.54-2.50)
Sidechain outliers	138945	6337 (2.54-2.50)
RSRZ outliers	127900	5630 (2.54-2.50)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	319	 8% 79% 9% • 12%
1	B	319	 10% 81% 9% • 9%
1	C	319	 7% 79% 9% • 11%
1	D	319	 12% 84% 8% • 6%

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9474 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.

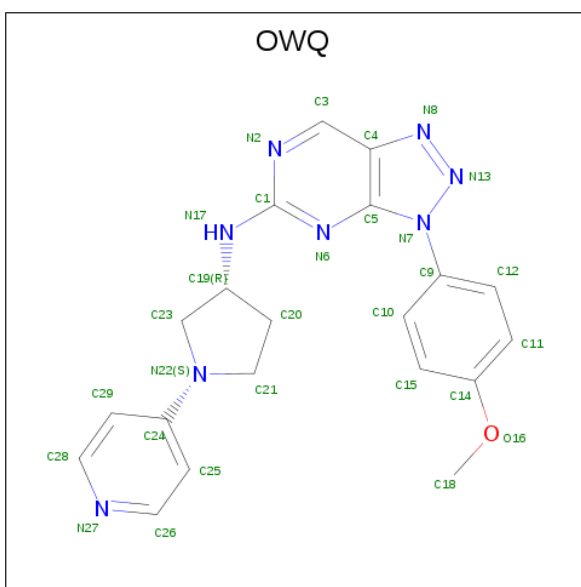
- Molecule 1 is a protein called Mitogen-activated protein kinase 6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	281	2231	1430	377	411	13	0	0	0
1	B	289	2297	1472	386	426	13	0	2	0
1	C	283	2244	1437	380	414	13	0	0	0
1	D	301	2397	1533	404	448	12	0	2	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	290	VAL	LEU	engineered mutation	UNP Q16659
B	290	VAL	LEU	engineered mutation	UNP Q16659
C	290	VAL	LEU	engineered mutation	UNP Q16659
D	290	VAL	LEU	engineered mutation	UNP Q16659

- Molecule 2 is 3-(4-methoxyphenyl)- {N}-[(3 {R})-1-pyridin-4-ylpyrrolidin-3-yl]-[1,2,3]triazolo[4,5-d]pyrimidin-5-amine (three-letter code: OWQ) (formula: C<sub>20</sub>H<sub>20</sub>N<sub>8</sub>O) (labeled as "Ligand of Interest" by author).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
2	A	1	29	20	8	1	0	0
2	B	1	29	20	8	1	0	0
2	C	1	29	20	8	1	0	0
2	D	1	29	20	8	1	0	0

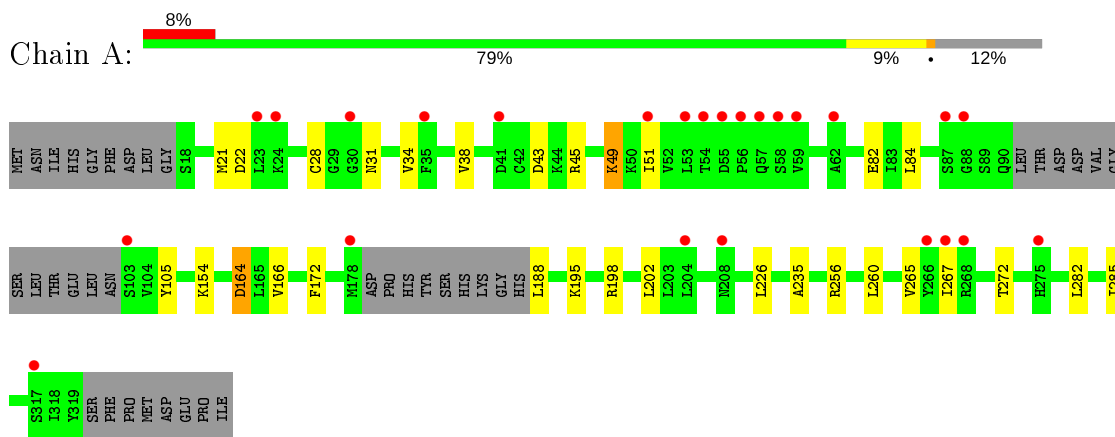
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	45	45	45	0	0
3	B	40	40	40	0	0
3	C	56	56	56	0	0
3	D	48	48	48	0	0

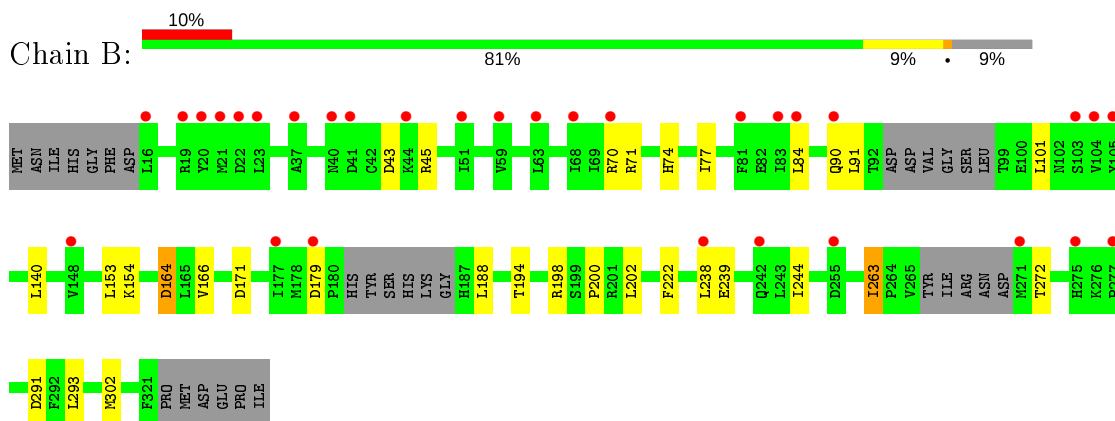
### 3 Residue-property plots [i](#)

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.

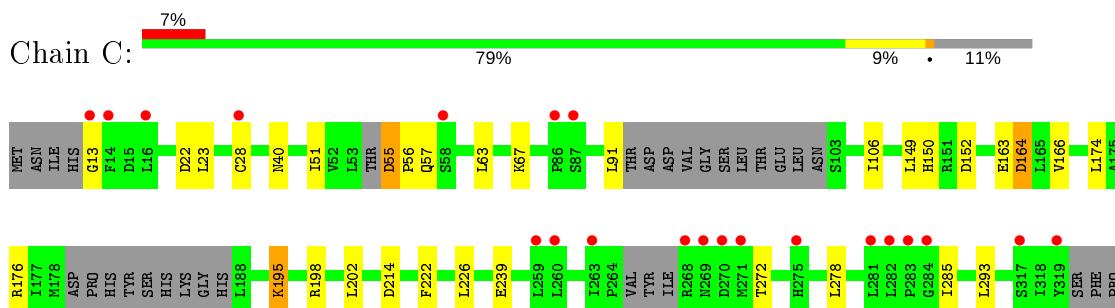
- Molecule 1: Mitogen-activated protein kinase 6



- Molecule 1: Mitogen-activated protein kinase 6




- Molecule 1: Mitogen-activated protein kinase 6



MET  
ASP  
GLU  
PRO  
ILE

• Molecule 1: Mitogen-activated protein kinase 6

Chain D:  12% 84% 8% 6%

MET ASP ILE HIS GLY F14 G17 S18 M21 D22 L23 L26 N31 G32 I33 V34 D39 N40 D41 C42 D43 K44 R45 K49 L53 T54 D55 V59 R70 R71 H74 I77 D94 V95 G96 S97 L98 T99 E100 L149 D164 I165 V166 I177

H181 K185 G186 H187 K195 R198 L202 L213 F222 E239 E252 E253 E254 D255 L259 L260 I263 P264 VAL TVR ILE ARG ASN ASP MET T272 E273 P274 E275 K276 P277 L278 T279 Q280 L281 L282 E288 A289 V290 D291 F292 L293 T298 F299 E305 F321

PRO  
MET  
ASP  
GLU  
PRO  
ILE

## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	65.98Å 100.88Å 195.52Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.08 – 2.52 48.08 – 2.52	Depositor EDS
% Data completeness (in resolution range)	82.2 (48.08-2.52) 82.2 (48.08-2.52)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	0.09	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.20 (at 2.51Å)	Xtrriage
Refinement program	BUSTER 2.11.7	Depositor
R, $R_{free}$	0.198 , 0.251 0.204 , 0.259	Depositor DCC
$R_{free}$ test set	1857 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	60.5	Xtrriage
Anisotropy	0.048	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 64.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	9474	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	74.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.18% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup> Intensities estimated from amplitudes.

<sup>2</sup> Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

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

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	0.51	0/2274	0.68	0/3080
1	B	0.51	0/2347	0.67	0/3178
1	C	0.50	0/2285	0.70	0/3089
1	D	0.49	0/2450	0.68	0/3320
All	All	0.50	0/9356	0.68	0/12667

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	2231	0	2253	15	0
1	B	2297	0	2324	11	0
1	C	2244	0	2262	12	0
1	D	2397	0	2402	12	0
2	A	29	0	0	0	0
2	B	29	0	0	0	0
2	C	29	0	0	1	0
2	D	29	0	0	0	0
3	A	45	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	40	0	0	1	0
3	C	56	0	0	0	0
3	D	48	0	0	0	0
All	All	9474	0	9241	48	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 3.

All (48) 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:53:LEU:HD13	1:D:59:VAL:HG22	1.68	0.75
1:D:149:LEU:HD23	1:D:177:ILE:HG12	1.68	0.75
1:A:282:LEU:HB3	1:A:285:ILE:HG12	1.81	0.62
1:B:302:MET:HB3	1:D:71:ARG:CZ	2.31	0.60
1:A:21:MET:HG3	1:A:38:VAL:HG13	1.83	0.59
1:B:140:LEU:HD21	1:B:153:LEU:HD11	1.85	0.58
1:A:154:LYS:HD2	1:C:163:GLU:HG2	1.86	0.56
1:A:31:ASN:HB2	1:A:49:LYS:HE3	1.87	0.56
1:C:150:HIS:O	1:C:214:ASP:OD1	2.24	0.55
1:A:226:LEU:HD13	1:A:285:ILE:HD13	1.89	0.54
1:C:51:ILE:HD11	1:C:106:ILE:HD11	1.89	0.54
1:B:164:ASP:HB2	1:B:166:VAL:HG23	1.90	0.53
1:C:55:ASP:HB3	1:C:56:PRO:HD3	1.90	0.53
1:D:164:ASP:HB2	1:D:166:VAL:HG23	1.91	0.52
1:C:164:ASP:HB2	1:C:166:VAL:HG23	1.91	0.51
1:D:222:PHE:HD2	1:D:293:LEU:HD13	1.75	0.51
1:A:31:ASN:HB3	1:A:51:ILE:HG12	1.93	0.50
1:C:152:ASP:HB2	1:C:174:LEU:HD12	1.93	0.50
1:B:222:PHE:HD2	1:B:293:LEU:HD13	1.76	0.49
1:A:164:ASP:HB2	1:A:166:VAL:HG23	1.95	0.49
1:B:291:ASP:HB3	3:B:509:HOH:O	2.13	0.47
1:A:235:ALA:HB3	2:C:401:OWQ:C26	2.45	0.46
1:B:238:LEU:CD1	1:B:263:ILE:HG21	2.46	0.46
1:C:13:GLY:HA3	1:C:22:ASP:HA	1.98	0.46
1:D:198:ARG:HG2	1:D:202:LEU:HD23	1.97	0.45
1:A:198:ARG:HG2	1:A:202:LEU:HD23	1.98	0.45
1:B:74:HIS:HB3	1:B:77:ILE:HG12	1.97	0.45
1:A:256:ARG:O	1:A:260:LEU:HD13	2.16	0.45
1:C:222:PHE:HD2	1:C:293:LEU:HD13	1.81	0.45
1:A:49:LYS:HE2	1:A:172:PHE:CD2	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:198:ARG:HG2	1:C:202:LEU:HD23	1.98	0.45
1:A:195:LYS:HA	1:A:198:ARG:HD2	1.99	0.44
1:B:200:PRO:HG3	1:B:244:ILE:HG21	2.00	0.44
1:A:34:VAL:HG22	1:A:49:LYS:HD3	2.00	0.44
1:D:181:HIS:O	1:D:185:LYS:HE3	2.18	0.44
1:D:34:VAL:HG22	1:D:49:LYS:HD2	2.00	0.43
1:B:154:LYS:HD3	1:B:194:THR:HG21	2.00	0.43
1:D:55:ASP:O	1:D:59:VAL:HG23	2.18	0.43
1:A:226:LEU:HB3	1:A:285:ILE:HG21	2.01	0.42
1:D:195:LYS:HA	1:D:198:ARG:HD2	2.01	0.42
1:C:195:LYS:HA	1:C:198:ARG:HD2	2.02	0.41
1:B:90:GLN:HE21	1:B:91:LEU:H	1.69	0.41
1:D:44:LYS:HD3	1:D:45:ARG:H	1.86	0.41
1:B:198:ARG:HG2	1:B:202:LEU:HD23	2.01	0.41
1:D:74:HIS:HB3	1:D:77:ILE:HG12	2.01	0.41
1:C:51:ILE:HD11	1:C:106:ILE:CD1	2.51	0.40
1:A:49:LYS:O	1:A:105:TYR:HA	2.21	0.40
1:C:226:LEU:HD13	1:C:285:ILE:HG21	2.03	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	275/319 (86%)	267 (97%)	8 (3%)	0	100	100
1	B	283/319 (89%)	275 (97%)	6 (2%)	2 (1%)	22	37
1	C	273/319 (86%)	266 (97%)	7 (3%)	0	100	100
1	D	299/319 (94%)	288 (96%)	11 (4%)	0	100	100
All	All	1130/1276 (89%)	1096 (97%)	32 (3%)	2 (0%)	51	67

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	171[A]	ASP
1	B	171[B]	ASP

### 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	246/283 (87%)	234 (95%)	12 (5%)	25	45
1	B	256/283 (90%)	244 (95%)	12 (5%)	26	46
1	C	247/283 (87%)	232 (94%)	15 (6%)	18	34
1	D	266/283 (94%)	249 (94%)	17 (6%)	17	31
All	All	1015/1132 (90%)	959 (94%)	56 (6%)	22	39

All (56) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	ASP
1	A	28	CYS
1	A	43	ASP
1	A	45	ARG
1	A	49	LYS
1	A	82	GLU
1	A	84	LEU
1	A	164	ASP
1	A	188	LEU
1	A	265	VAL
1	A	267	ILE
1	A	272	THR
1	B	43	ASP
1	B	45	ARG
1	B	70	ARG
1	B	71	ARG
1	B	84	LEU
1	B	101	LEU
1	B	164	ASP
1	B	179	ASP

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	B	188	LEU
1	B	239	GLU
1	B	263	ILE
1	B	272	THR
1	C	23	LEU
1	C	28	CYS
1	C	40	ASN
1	C	55	ASP
1	C	57	GLN
1	C	63	LEU
1	C	67	LYS
1	C	91	LEU
1	C	149	LEU
1	C	164	ASP
1	C	176	ARG
1	C	195	LYS
1	C	239	GLU
1	C	272	THR
1	C	278	LEU
1	D	21	MET
1	D	23	LEU
1	D	31	ASN
1	D	43	ASP
1	D	53	LEU
1	D	55	ASP
1	D	70	ARG
1	D	164	ASP
1	D	185	LYS
1	D	187[A]	HIS
1	D	187[B]	HIS
1	D	195	LYS
1	D	213	ILE
1	D	239	GLU
1	D	253	GLU
1	D	272	THR
1	D	275	HIS

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

<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	A	31	ASN
1	A	295	GLN

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Mol	Chain	Res	Type
1	B	40	ASN
1	B	90	GLN
1	B	295	GLN
1	C	295	GLN
1	D	295	GLN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

### 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

### 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

### 5.6 Ligand geometry [i](#)

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	OWQ	D	401	-	30,33,33	1.35	5 (16%)	36,46,46	2.48	12 (33%)
2	OWQ	A	401	-	30,33,33	1.16	3 (10%)	36,46,46	2.72	15 (41%)
2	OWQ	C	401	-	30,33,33	1.26	4 (13%)	36,46,46	2.20	11 (30%)
2	OWQ	B	401	-	30,33,33	1.47	5 (16%)	36,46,46	2.62	12 (33%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns.

'-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	OWQ	D	401	-	-	10/14/23/23	0/5/5/5
2	OWQ	A	401	-	-	8/14/23/23	0/5/5/5
2	OWQ	C	401	-	-	9/14/23/23	0/5/5/5
2	OWQ	B	401	-	-	8/14/23/23	0/5/5/5

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	OWQ	C1-N17	3.96	1.39	1.34
2	D	401	OWQ	C1-N17	3.21	1.38	1.34
2	D	401	OWQ	N8-N13	-3.15	1.28	1.34
2	C	401	OWQ	N8-N13	-3.08	1.28	1.34
2	B	401	OWQ	N8-N13	-3.08	1.28	1.34
2	B	401	OWQ	N13-N7	2.93	1.41	1.37
2	C	401	OWQ	N13-N7	2.89	1.41	1.37
2	C	401	OWQ	C1-N17	2.82	1.38	1.34
2	D	401	OWQ	N13-N7	2.80	1.41	1.37
2	A	401	OWQ	N8-N13	-2.78	1.29	1.34
2	A	401	OWQ	N13-N7	2.78	1.41	1.37
2	D	401	OWQ	C23-N22	2.67	1.50	1.46
2	B	401	OWQ	C23-N22	2.40	1.49	1.46
2	D	401	OWQ	C21-N22	2.40	1.50	1.47
2	B	401	OWQ	C21-N22	2.28	1.50	1.47
2	C	401	OWQ	C5-N6	-2.14	1.32	1.35
2	A	401	OWQ	C5-N6	-2.09	1.32	1.35

All (50) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	OWQ	C23-N22-C24	-8.22	110.94	122.72
2	B	401	OWQ	C21-N22-C24	-7.61	114.52	123.57
2	A	401	OWQ	C21-N22-C24	-6.48	115.85	123.57
2	D	401	OWQ	C21-N22-C24	-6.46	115.88	123.57
2	C	401	OWQ	C23-N22-C24	-6.22	113.81	122.72
2	B	401	OWQ	N2-C1-N6	-6.11	120.61	126.52
2	C	401	OWQ	C1-N6-C5	5.11	121.08	115.28
2	D	401	OWQ	C1-N6-C5	5.00	120.96	115.28
2	B	401	OWQ	C1-N6-C5	4.98	120.93	115.28
2	D	401	OWQ	N2-C1-N6	-4.93	121.76	126.52
2	A	401	OWQ	C1-N17-C19	-4.53	116.66	124.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	OWQ	C3-N2-C1	4.53	122.60	115.88
2	D	401	OWQ	C1-N17-C19	-4.30	117.05	124.31
2	A	401	OWQ	C9-N7-N13	4.27	126.80	119.95
2	B	401	OWQ	C23-N22-C24	-4.22	116.68	122.72
2	D	401	OWQ	C3-N2-C1	4.05	121.89	115.88
2	C	401	OWQ	N2-C1-N6	-4.04	122.61	126.52
2	A	401	OWQ	N8-N13-N7	4.04	110.86	106.37
2	B	401	OWQ	N8-N13-N7	3.95	110.77	106.37
2	D	401	OWQ	C23-N22-C24	-3.92	117.10	122.72
2	D	401	OWQ	N8-N13-N7	3.88	110.68	106.37
2	A	401	OWQ	C1-N6-C5	3.87	119.67	115.28
2	A	401	OWQ	C28-N27-C26	3.58	125.28	116.85
2	B	401	OWQ	C9-N7-N13	3.47	125.51	119.95
2	C	401	OWQ	N8-N13-N7	3.41	110.17	106.37
2	C	401	OWQ	C21-N22-C24	-3.40	119.52	123.57
2	B	401	OWQ	C20-C21-N22	3.40	107.34	103.35
2	B	401	OWQ	C21-N22-C23	-3.23	103.19	111.55
2	A	401	OWQ	N2-C1-N6	-3.10	123.52	126.52
2	D	401	OWQ	C9-N7-N13	3.09	124.89	119.95
2	C	401	OWQ	C1-N17-C19	-3.04	119.19	124.31
2	D	401	OWQ	C28-N27-C26	3.02	123.94	116.85
2	D	401	OWQ	C21-C20-C19	-2.99	99.74	104.08
2	C	401	OWQ	C3-N2-C1	2.92	120.21	115.88
2	A	401	OWQ	C3-N2-C1	2.91	120.20	115.88
2	C	401	OWQ	C9-N7-N13	2.87	124.54	119.95
2	A	401	OWQ	C25-C26-N27	-2.70	118.92	123.62
2	A	401	OWQ	C29-C28-N27	-2.68	118.96	123.62
2	C	401	OWQ	C28-N27-C26	2.67	123.14	116.85
2	D	401	OWQ	C25-C26-N27	-2.39	119.46	123.62
2	B	401	OWQ	N17-C1-N6	2.36	120.72	117.18
2	C	401	OWQ	C29-C28-N27	-2.34	119.55	123.62
2	A	401	OWQ	C29-C24-N22	-2.19	118.36	121.38
2	A	401	OWQ	C3-C4-N8	2.13	133.96	130.51
2	D	401	OWQ	C29-C28-N27	-2.09	119.98	123.62
2	B	401	OWQ	C18-O16-C14	2.09	122.04	117.51
2	A	401	OWQ	C9-N7-C5	-2.07	127.75	129.98
2	A	401	OWQ	C21-N22-C23	-2.03	106.31	111.55
2	C	401	OWQ	C21-N22-C23	-2.02	106.33	111.55
2	B	401	OWQ	C3-C4-N8	2.01	133.78	130.51

There are no chirality outliers.

All (35) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	401	OWQ	N2-C1-N17-C19
2	D	401	OWQ	C10-C9-N7-C5
2	D	401	OWQ	C12-C9-N7-C5
2	D	401	OWQ	C25-C24-N22-C23
2	D	401	OWQ	C29-C24-N22-C23
2	A	401	OWQ	C10-C9-N7-C5
2	A	401	OWQ	C12-C9-N7-C5
2	A	401	OWQ	C25-C24-N22-C23
2	A	401	OWQ	C29-C24-N22-C23
2	C	401	OWQ	C10-C9-N7-C5
2	C	401	OWQ	C10-C9-N7-N13
2	C	401	OWQ	C12-C9-N7-C5
2	C	401	OWQ	C12-C9-N7-N13
2	B	401	OWQ	C10-C9-N7-C5
2	B	401	OWQ	C12-C9-N7-C5
2	B	401	OWQ	C25-C24-N22-C21
2	B	401	OWQ	C29-C24-N22-C21
2	D	401	OWQ	C11-C14-O16-C18
2	D	401	OWQ	C15-C14-O16-C18
2	A	401	OWQ	C11-C14-O16-C18
2	C	401	OWQ	C25-C24-N22-C21
2	A	401	OWQ	C15-C14-O16-C18
2	C	401	OWQ	C29-C24-N22-C21
2	C	401	OWQ	C15-C14-O16-C18
2	C	401	OWQ	C11-C14-O16-C18
2	D	401	OWQ	C10-C9-N7-N13
2	D	401	OWQ	C12-C9-N7-N13
2	A	401	OWQ	C12-C9-N7-N13
2	B	401	OWQ	C12-C9-N7-N13
2	B	401	OWQ	C15-C14-O16-C18
2	B	401	OWQ	C11-C14-O16-C18
2	A	401	OWQ	C10-C9-N7-N13
2	B	401	OWQ	C10-C9-N7-N13
2	D	401	OWQ	N6-C1-N17-C19
2	C	401	OWQ	C29-C24-N22-C23

There are no ring outliers.

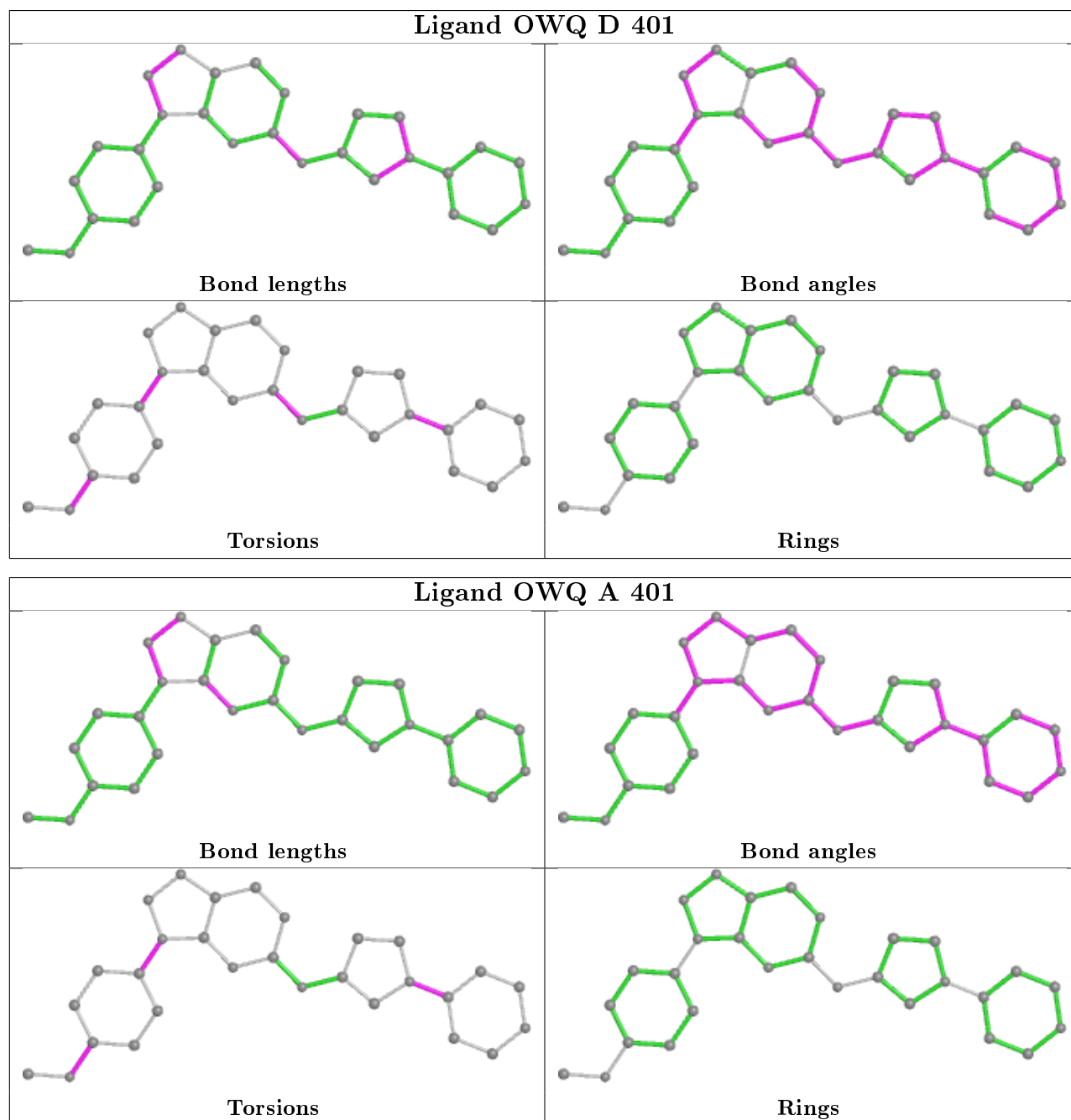
1 monomer is involved in 1 short contact:

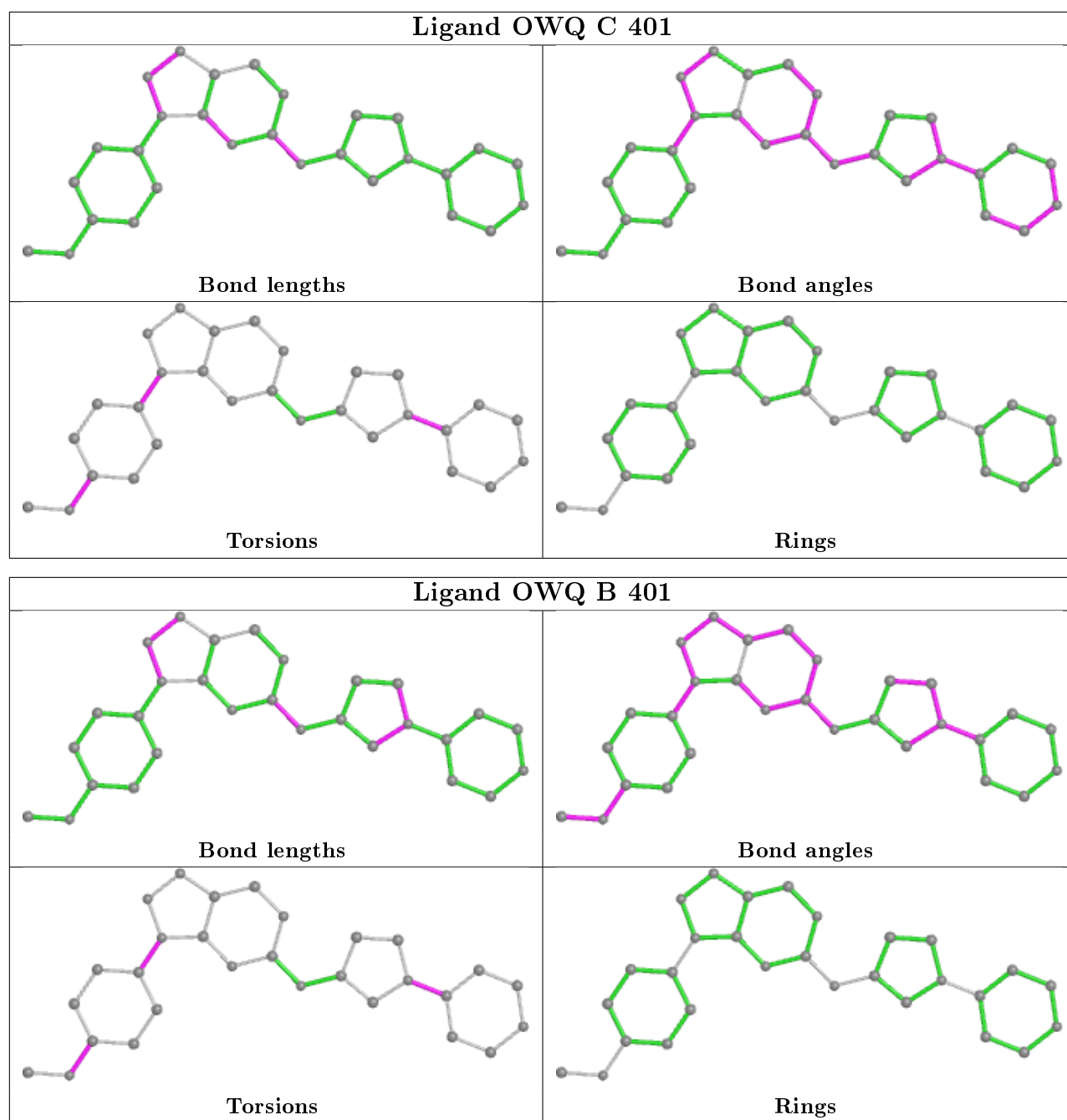
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	401	OWQ	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths,



bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight  $> 250$  and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





## 5.7 Other polymers [\(i\)](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [\(i\)](#)

There are no chain breaks in this entry.

## 6 Fit of model and data [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	281/319 (88%)	0.63	25 (8%) <b>9</b> <b>10</b>	33, 67, 123, 150	0
1	B	289/319 (90%)	0.60	31 (10%) <b>6</b> <b>6</b>	44, 73, 119, 143	0
1	C	283/319 (88%)	0.44	21 (7%) <b>14</b> <b>15</b>	34, 68, 113, 136	0
1	D	301/319 (94%)	0.67	37 (12%) <b>4</b> <b>4</b>	43, 74, 131, 151	0
All	All	1154/1276 (90%)	0.59	114 (9%) <b>7</b> <b>7</b>	33, 71, 121, 151	0

All (114) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	277	PRO	7.4
1	D	263	ILE	7.4
1	D	95	VAL	7.2
1	A	56	PRO	6.8
1	B	21	MET	6.4
1	C	284	GLY	6.0
1	B	23	LEU	5.9
1	D	94	ASP	5.6
1	A	58	SER	5.6
1	C	13	GLY	5.3
1	A	87	SER	5.2
1	B	105	TYR	4.9
1	D	275	HIS	4.7
1	D	96	GLY	4.7
1	A	59	VAL	4.6
1	A	204	LEU	4.5
1	A	267	ILE	4.5
1	D	278	LEU	4.4
1	B	83	ILE	4.3
1	A	55	ASP	4.3
1	B	84	LEU	4.3

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	A	103	SER	4.2
1	B	44	LYS	4.2
1	A	268	ARG	4.1
1	B	40	ASN	4.1
1	D	272	THR	4.1
1	C	263	ILE	4.1
1	A	266	TYR	4.0
1	B	90	GLN	4.0
1	B	19	ARG	3.9
1	A	57	GLN	3.9
1	C	259	LEU	3.8
1	C	282	LEU	3.8
1	B	68	ILE	3.7
1	A	178	MET	3.7
1	C	16	LEU	3.7
1	A	54	THR	3.6
1	A	53	LEU	3.6
1	B	20	TYR	3.6
1	D	18	SER	3.6
1	D	97	SER	3.5
1	D	259	LEU	3.5
1	B	63	LEU	3.5
1	C	268	ARG	3.4
1	C	271	MET	3.4
1	D	276	LYS	3.4
1	B	37	ALA	3.4
1	D	260	LEU	3.3
1	A	317	SER	3.2
1	C	269	ASN	3.2
1	D	282	LEU	3.2
1	D	280	GLN	3.2
1	B	104	VAL	3.1
1	D	274	PRO	3.0
1	D	99	THR	3.0
1	C	283	PRO	3.0
1	A	30	GLY	2.9
1	D	21	MET	2.9
1	C	14	PHE	2.9
1	D	273	GLU	2.9
1	D	33	LEU	2.9
1	B	271	MET	2.9
1	D	264	PRO	2.9

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	B	242	GLN	2.9
1	B	70	ARG	2.8
1	B	103	SER	2.8
1	B	59	VAL	2.8
1	C	270	ASP	2.8
1	D	291	ASP	2.8
1	B	275[A]	HIS	2.8
1	A	23	LEU	2.8
1	D	289	ALA	2.7
1	C	28	CYS	2.7
1	D	253	GLU	2.7
1	C	87	SER	2.6
1	B	81	PHE	2.6
1	D	298	THR	2.6
1	A	41	ASP	2.6
1	B	22	ASP	2.6
1	D	14	PHE	2.6
1	B	41	ASP	2.6
1	D	305	LEU	2.5
1	A	88	GLY	2.5
1	C	281	LEU	2.5
1	D	299	PHE	2.5
1	D	100	GLU	2.5
1	C	58	SER	2.5
1	D	39	ASP	2.5
1	D	252	HIS	2.5
1	D	255	ASP	2.5
1	A	62	ALA	2.4
1	A	275	HIS	2.4
1	C	317	SER	2.4
1	C	86	PRO	2.4
1	C	319	TYR	2.3
1	C	275	HIS	2.3
1	D	290	VAL	2.3
1	A	24	LYS	2.2
1	A	208	ASN	2.2
1	D	41	ASP	2.2
1	D	17	GLY	2.2
1	B	16	LEU	2.2
1	D	288	GLU	2.2
1	B	148	VAL	2.2
1	B	238	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	260	LEU	2.1
1	B	179	ASP	2.1
1	B	277	PRO	2.1
1	B	255	ASP	2.0
1	B	51	ILE	2.0
1	B	177	ILE	2.0
1	D	26	LEU	2.0
1	A	35	PHE	2.0
1	A	51	ILE	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 [\(i\)](#)

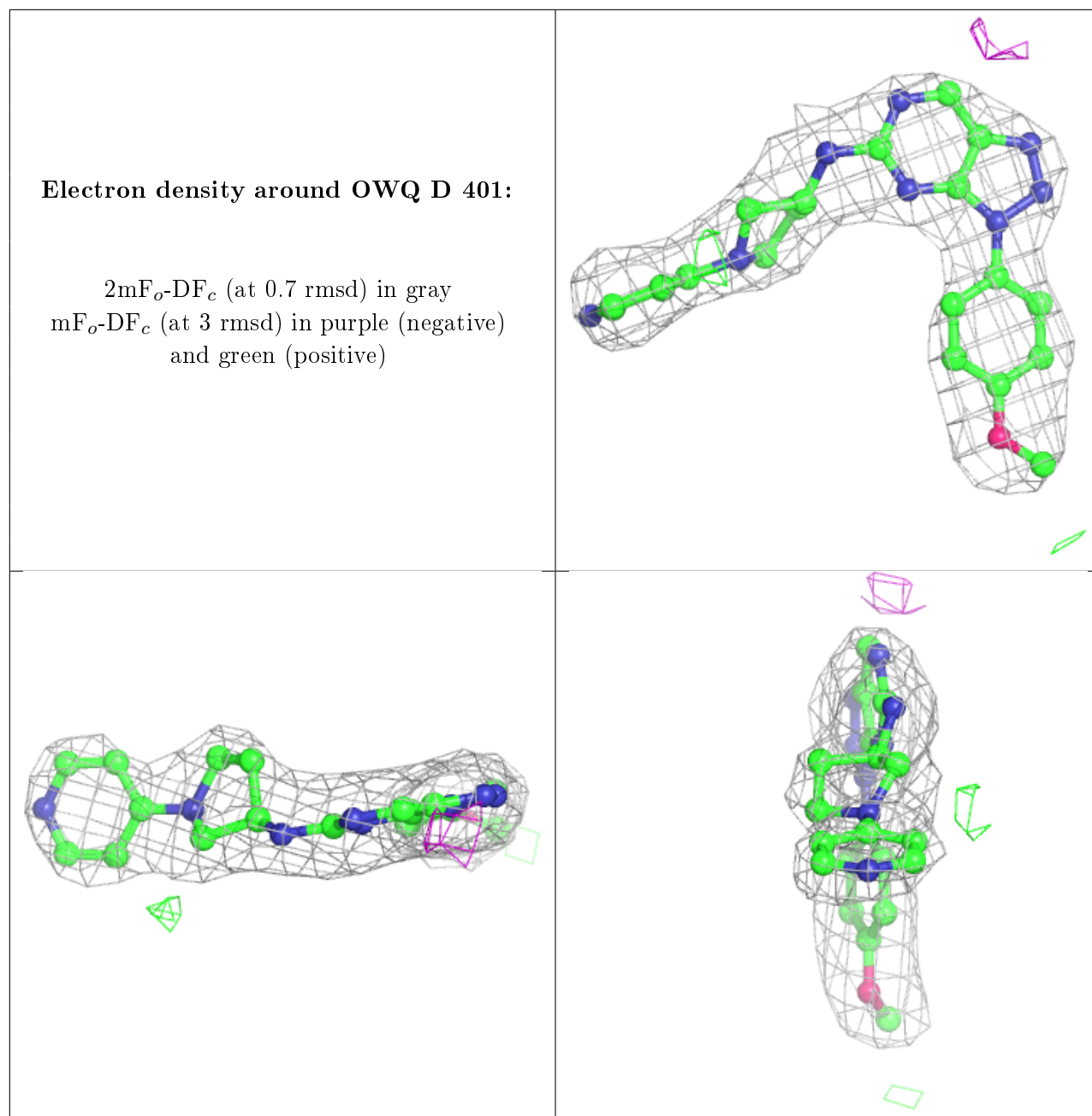
There are no monosaccharides in this entry.

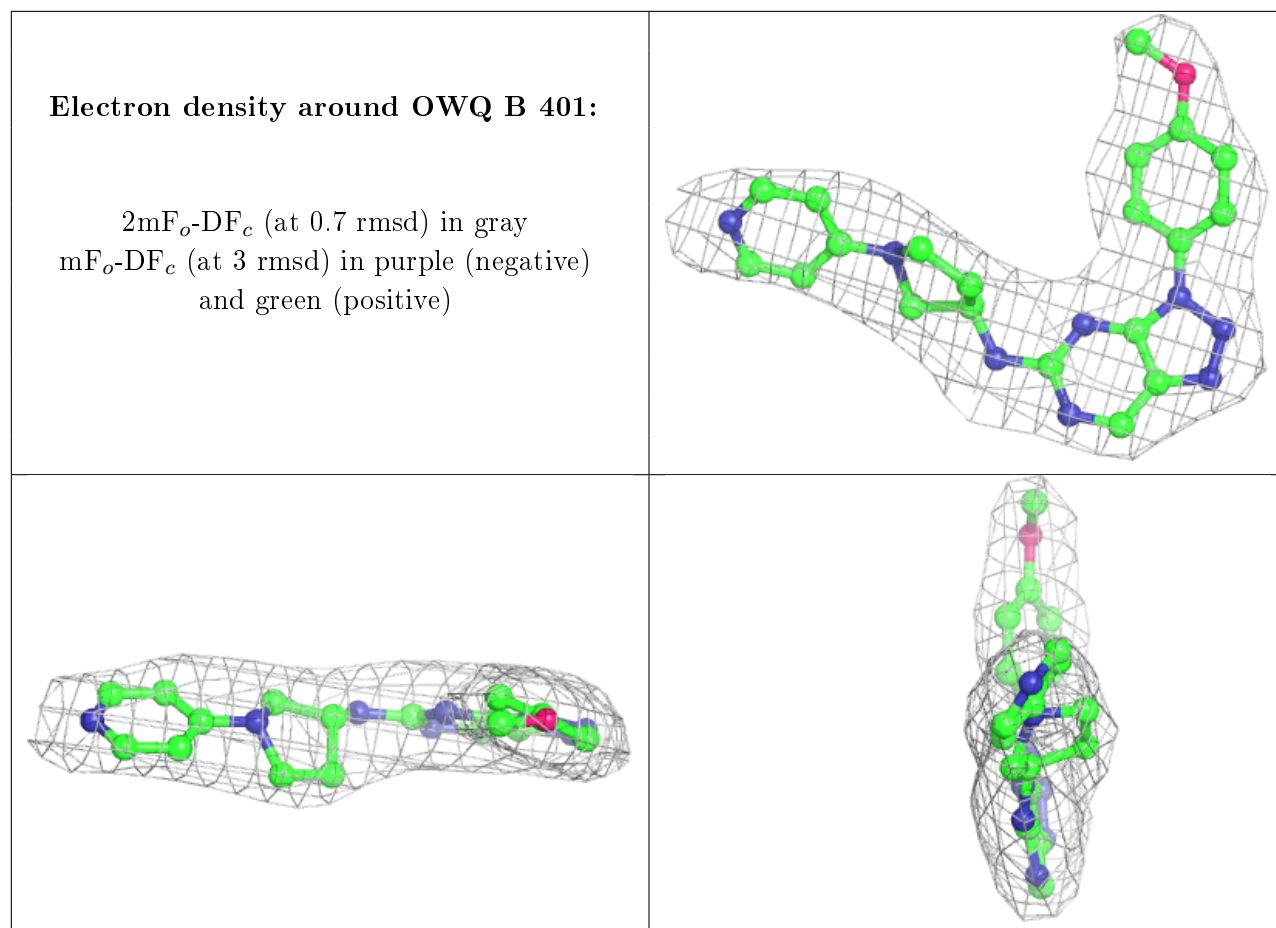
## 6.4 Ligands [\(i\)](#)

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, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

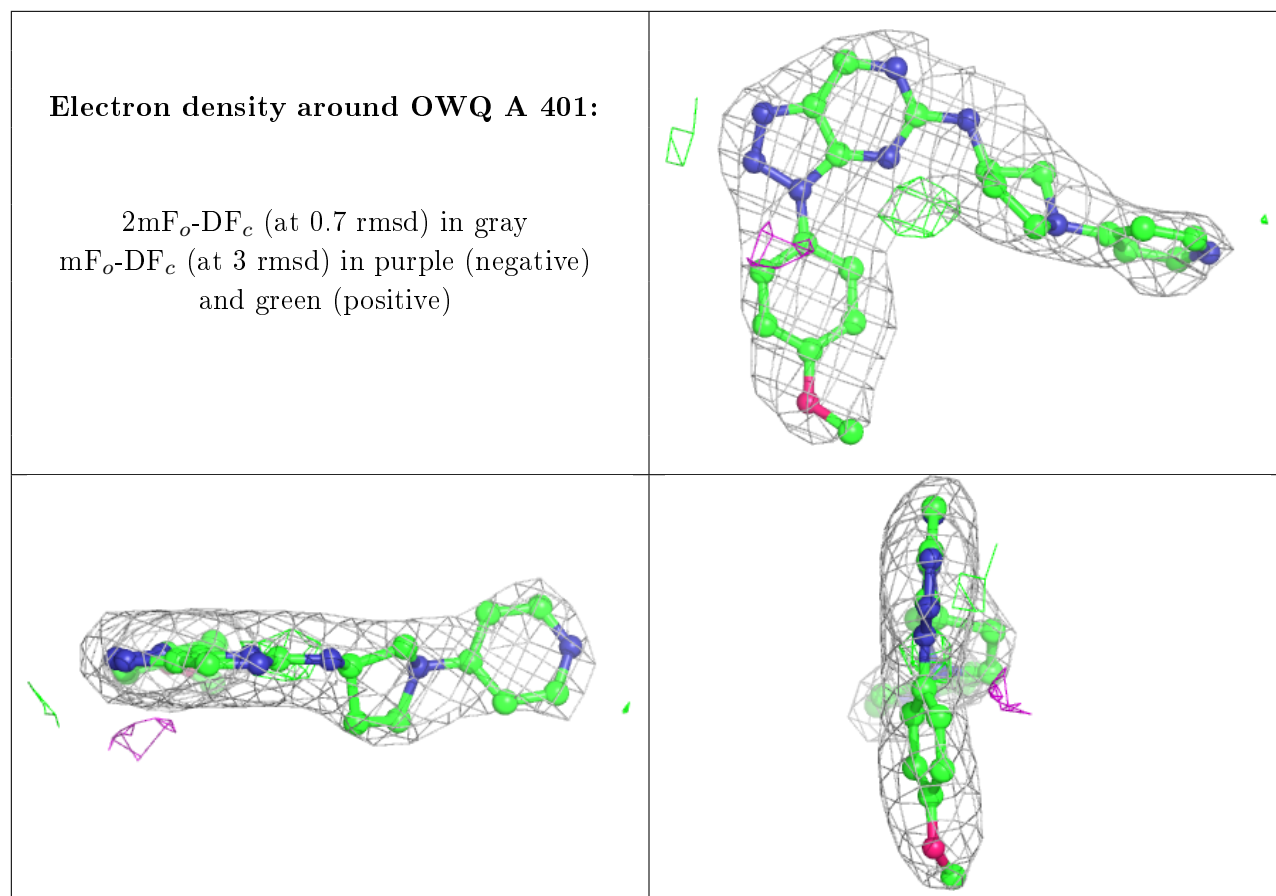
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	OWQ	D	401	29/29	0.94	0.16	59,65,81,82	0
2	OWQ	B	401	29/29	0.94	0.15	49,55,77,77	0
2	OWQ	A	401	29/29	0.95	0.17	47,59,81,83	0
2	OWQ	C	401	29/29	0.96	0.18	38,50,61,62	0

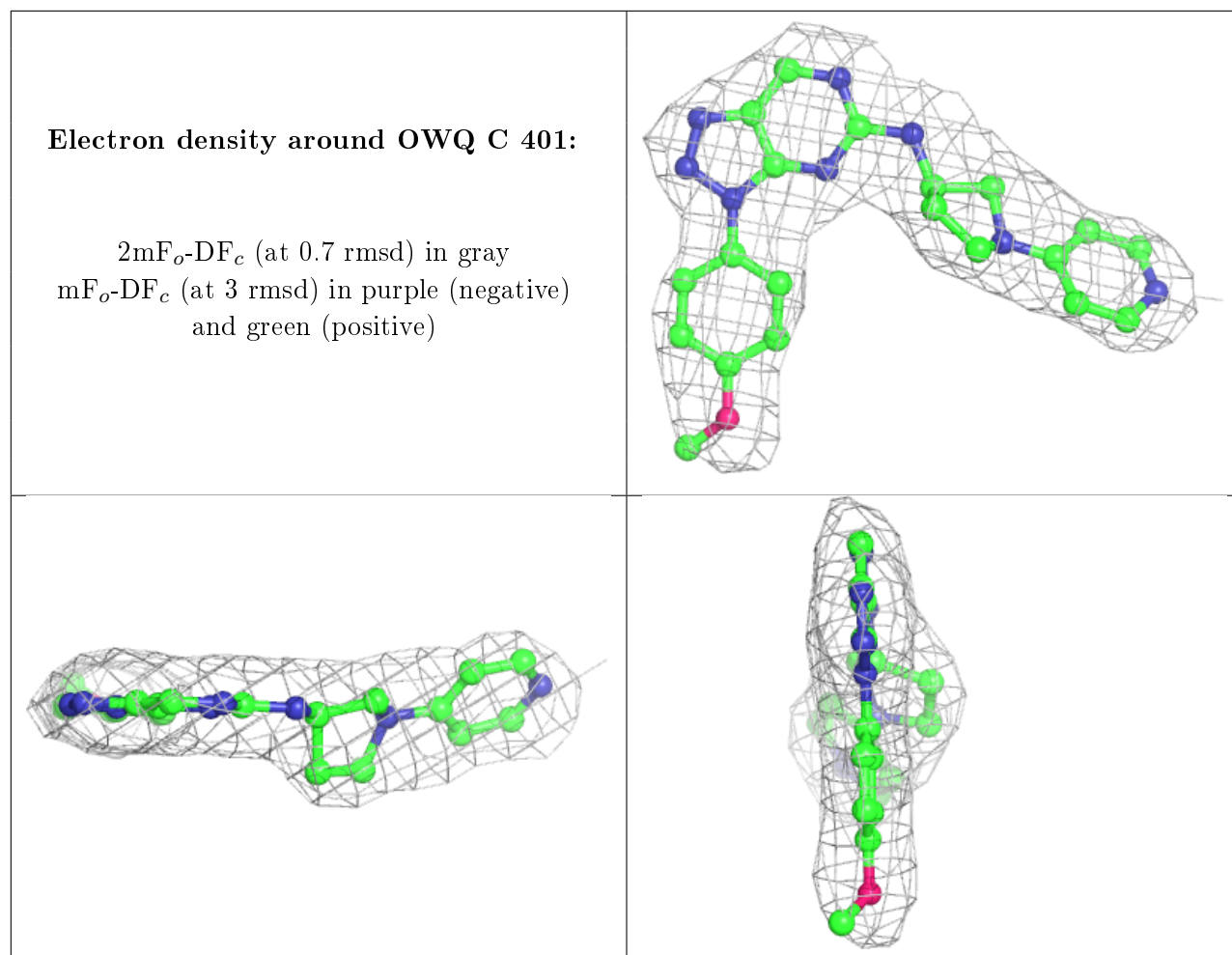
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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











## 6.5 Other polymers [i](#)

There are no such residues in this entry.