



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 6, 2023 – 09:21 AM EST

PDB ID : 6WVF  
Title : E.coli DsbB C104S with ubiquinone  
Authors : Liu, S.; Li, W.  
Deposited on : 2020-05-05  
Resolution : 2.90 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.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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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 : 2.36  
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

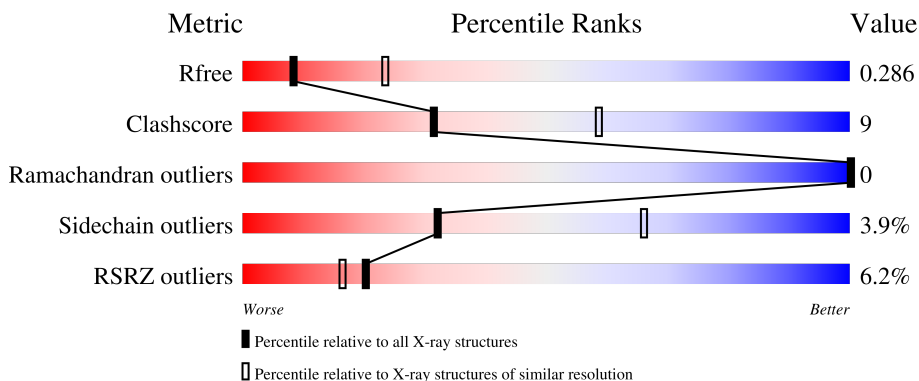
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.90 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	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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  $\geq 3$ , 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	399	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	UQ1	A	501	-	-	X	-

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3107 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.

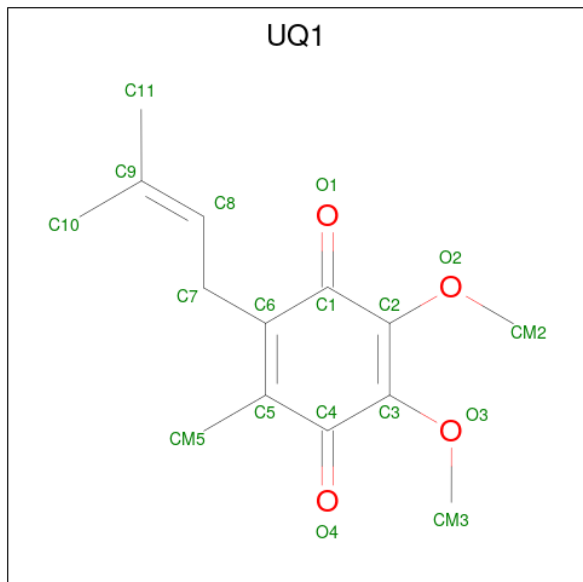
- Molecule 1 is a protein called Green fluorescent protein,Disulfide bond formation protein B,Green fluorescent protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	387	3088	2006	514	554	14	0	0	0

There are 25 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	30	ARG	SER	conflict	UNP P42212
A	39	ASN	TYR	conflict	UNP P42212
A	64	LEU	PHE	conflict	UNP P42212
A	66	CRO	SER	chromophore	UNP P42212
A	66	CRO	TYR	chromophore	UNP P42212
A	66	CRO	GLY	chromophore	UNP P42212
A	80	ARG	GLN	conflict	UNP P42212
A	99	SER	PHE	conflict	UNP P42212
A	105	THR	ASN	conflict	UNP P42212
A	147	ALA	CYS	engineered mutation	UNP P0A6M2
A	188	VAL	CYS	engineered mutation	UNP P0A6M2
A	243	SER	CYS	engineered mutation	UNP P0A6M2
A	314	THR	MET	conflict	UNP P42212
A	324	ALA	VAL	conflict	UNP P42212
A	332	VAL	ILE	conflict	UNP P42212
A	367	VAL	ALA	conflict	UNP P42212
A	393	HIS	-	expression tag	UNP P42212
A	394	HIS	-	expression tag	UNP P42212
A	395	HIS	-	expression tag	UNP P42212
A	396	HIS	-	expression tag	UNP P42212
A	397	HIS	-	expression tag	UNP P42212
A	398	HIS	-	expression tag	UNP P42212
A	399	HIS	-	expression tag	UNP P42212
A	400	HIS	-	expression tag	UNP P42212
A	401	HIS	-	expression tag	UNP P42212

- Molecule 2 is UBIQUINONE-1 (three-letter code: UQ1) (formula:  $C_{14}H_{18}O_4$ ) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	C O		
2	A	1	18	14 4	0	0

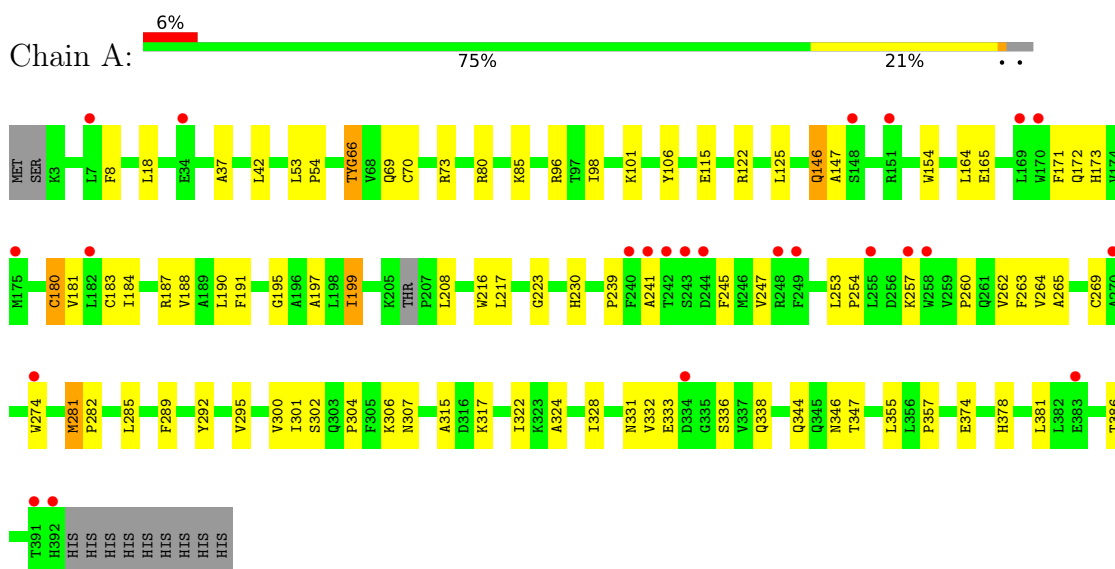
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	1	1	1	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: Green fluorescent protein, Disulfide bond formation protein B, Green fluorescent protein



## 4 Data and refinement statistics i

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	53.38Å 53.38Å 280.74Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	46.23 – 2.90 46.23 – 2.68	Depositor EDS
% Data completeness (in resolution range)	97.8 (46.23-2.90) 96.9 (46.23-2.68)	Depositor EDS
$R_{merge}$	0.18	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.05 (at 2.69Å)	Xtrriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, $R_{free}$	0.259 , 0.286 0.257 , 0.286	Depositor DCC
$R_{free}$ test set	629 reflections (4.65%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	75.7	Xtrriage
Anisotropy	0.403	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 67.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.047 for -h,-k,l	Xtrriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	3107	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	113.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.02% 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: CRO, UQ1

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	0.24	0/3145	0.42	0/4270

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	3088	0	3069	55	0
2	A	18	0	18	11	0
3	A	1	0	0	0	0
All	All	3107	0	3087	56	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 9.

All (56) 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:274:TRP:HB2	1:A:281:MET:HG2	1.61	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:183:CYS:SG	2:A:501:UQ1:O4	2.40	0.80
1:A:183:CYS:HG	2:A:501:UQ1:C4	1.99	0.75
1:A:183:CYS:SG	2:A:501:UQ1:C4	2.76	0.74
1:A:181:VAL:HG22	1:A:265:ALA:HB1	1.71	0.72
1:A:180:CYS:SG	1:A:181:VAL:N	2.64	0.70
1:A:183:CYS:SG	2:A:501:UQ1:HM52	2.34	0.68
1:A:98:ILE:HB	1:A:106:TYR:HB2	1.76	0.68
1:A:115:GLU:OE2	1:A:122:ARG:NH1	2.28	0.66
1:A:254:PRO:HB2	1:A:257:LYS:HB3	1.77	0.65
1:A:70:CYS:O	1:A:85:LYS:NZ	2.28	0.65
2:A:501:UQ1:O1	2:A:501:UQ1:H8	1.98	0.63
1:A:324:ALA:HB3	1:A:344:GLN:HB3	1.84	0.58
1:A:172:GLN:HE22	1:A:264:VAL:HG21	1.69	0.57
1:A:187:ARG:NH1	1:A:262:VAL:O	2.38	0.56
1:A:171:PHE:HD2	2:A:501:UQ1:HM21	1.73	0.53
1:A:154:TRP:HE3	1:A:197:ALA:HA	1.74	0.53
1:A:217:LEU:HD11	1:A:300:VAL:HG21	1.92	0.52
1:A:73:ARG:HB3	1:A:386:THR:HG22	1.92	0.52
1:A:322:ILE:HG12	1:A:346:ASN:HB2	1.92	0.51
1:A:146:GLN:HB2	1:A:307:ASN:HB3	1.92	0.51
1:A:172:GLN:NE2	1:A:264:VAL:HG21	2.26	0.50
1:A:230:HIS:CG	1:A:282:PRO:HB3	2.47	0.49
1:A:66:CRO:HG11	1:A:381:LEU:HD21	1.95	0.48
1:A:315:ALA:HB2	1:A:357:PRO:O	2.13	0.48
1:A:332:VAL:HB	1:A:336:SER:HB3	1.96	0.48
1:A:66:CRO:HB2	1:A:69:GLN:HE21	1.78	0.47
1:A:172:GLN:HE21	1:A:173:HIS:CE1	2.33	0.47
1:A:216:TRP:HH2	1:A:295:VAL:HG12	1.79	0.47
1:A:66:CRO:O2	1:A:96:ARG:NH2	2.46	0.46
1:A:188:VAL:HG22	1:A:263:PHE:HE2	1.80	0.46
1:A:147:ALA:O	1:A:331:ASN:ND2	2.49	0.46
1:A:66:CRO:C2	1:A:96:ARG:HH22	2.28	0.45
1:A:260:PRO:O	1:A:264:VAL:HG22	2.17	0.45
1:A:302:SER:O	1:A:306:LYS:N	2.50	0.45
1:A:223:GLY:HA3	1:A:289:PHE:CD1	2.52	0.45
1:A:18:LEU:HD11	1:A:125:LEU:HB2	1.99	0.44
1:A:195:GLY:O	1:A:199:ILE:HG22	2.18	0.44
1:A:183:CYS:SG	2:A:501:UQ1:CM5	3.06	0.44
1:A:80:ARG:HA	1:A:355:LEU:HD13	2.00	0.44
1:A:180:CYS:SG	1:A:269:CYS:N	2.90	0.44
1:A:216:TRP:CH2	1:A:295:VAL:HG12	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:8:PHE:HB3	1:A:37:ALA:HB3	1.99	0.44
1:A:164:LEU:HB3	2:A:501:UQ1:H102	2.01	0.43
1:A:165:GLU:HG2	1:A:190:LEU:HD23	2.01	0.43
1:A:184:ILE:HD12	1:A:263:PHE:HA	1.99	0.43
1:A:253:LEU:HD12	1:A:253:LEU:HA	1.86	0.42
1:A:239:PRO:C	1:A:241:ALA:H	2.21	0.42
1:A:285:LEU:HD22	2:A:501:UQ1:H103	2.01	0.42
1:A:374:GLU:HB3	1:A:378:HIS:CE1	2.54	0.42
1:A:183:CYS:SG	2:A:501:UQ1:C5	3.07	0.42
1:A:181:VAL:HG21	1:A:245:PHE:C	2.40	0.42
1:A:53:LEU:HA	1:A:54:PRO:HD3	1.95	0.42
1:A:301:ILE:O	1:A:304:PRO:HD2	2.20	0.41
1:A:101:LYS:O	1:A:338:GLN:NE2	2.54	0.41
1:A:171:PHE:CD2	2:A:501:UQ1:HM21	2.54	0.41

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	380/399 (95%)	363 (96%)	17 (4%)	0	<b>100</b> <b>100</b>

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	330/343 (96%)	317 (96%)	13 (4%)	32 66

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	42	LEU
1	A	146	GLN
1	A	180	CYS
1	A	191	PHE
1	A	199	ILE
1	A	208	LEU
1	A	247	VAL
1	A	281	MET
1	A	292	TYR
1	A	317	LYS
1	A	328	ILE
1	A	333	GLU
1	A	347	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	172	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](#)

1 non-standard protein/DNA/RNA residue is 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
1	CRO	A	66	1	23,23,24	2.72	7 (30%)	30,32,34	2.80	8 (26%)

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
1	CRO	A	66	1	-	7/12/31/32	0/2/2/2

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	66	CRO	C1-N2	7.15	1.42	1.32
1	A	66	CRO	CA2-C2	6.22	1.54	1.48
1	A	66	CRO	C1-N3	5.08	1.45	1.37
1	A	66	CRO	C2-N3	3.87	1.48	1.39
1	A	66	CRO	CG2-CB2	3.59	1.53	1.46
1	A	66	CRO	CB2-CA2	-2.81	1.32	1.35
1	A	66	CRO	CA2-N2	2.26	1.43	1.38

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	66	CRO	O2-C2-CA2	-8.22	126.34	130.96
1	A	66	CRO	CA2-C2-N3	8.04	107.17	103.37
1	A	66	CRO	C2-N3-C1	-5.05	105.41	107.97
1	A	66	CRO	CA2-N2-C1	4.92	109.40	105.77
1	A	66	CRO	CG2-CB2-CA2	-4.16	124.85	129.94
1	A	66	CRO	C2-CA2-N2	-3.86	106.23	108.93
1	A	66	CRO	CB2-CA2-C2	2.59	125.36	122.28
1	A	66	CRO	O3-C3-CA3	-2.12	119.99	126.39

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	66	CRO	N1-CA1-CB1-CG1
1	A	66	CRO	C1-CA1-CB1-CG1
1	A	66	CRO	N2-CA2-CB2-CG2
1	A	66	CRO	C3-CA3-N3-C2

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Mol	Chain	Res	Type	Atoms
1	A	66	CRO	C2-CA2-CB2-CG2
1	A	66	CRO	C1-CA1-CB1-OG1
1	A	66	CRO	N1-CA1-CB1-OG1

There are no ring outliers.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	66	CRO	4	0

## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

1 ligand is 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	UQ1	A	501	-	18,18,18	2.94	6 (33%)	22,25,25	1.61	5 (22%)

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	UQ1	A	501	-	-	4/9/33/33	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	UQ1	C6-C5	9.83	1.53	1.35
2	A	501	UQ1	C3-C2	4.11	1.53	1.36
2	A	501	UQ1	C6-C1	2.62	1.54	1.46
2	A	501	UQ1	O4-C4	-2.14	1.18	1.23
2	A	501	UQ1	O2-CM2	-2.12	1.40	1.45
2	A	501	UQ1	O3-CM3	-2.06	1.40	1.45

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	UQ1	C7-C6-C1	4.41	123.78	118.48
2	A	501	UQ1	CM5-C5-C6	-3.05	119.43	124.40
2	A	501	UQ1	C11-C9-C10	2.61	120.38	114.60
2	A	501	UQ1	C6-C5-C4	2.27	120.98	119.18
2	A	501	UQ1	C7-C8-C9	-2.02	120.94	127.26

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	UQ1	C1-C6-C7-C8
2	A	501	UQ1	C5-C6-C7-C8
2	A	501	UQ1	C4-C3-O3-CM3
2	A	501	UQ1	C2-C3-O3-CM3

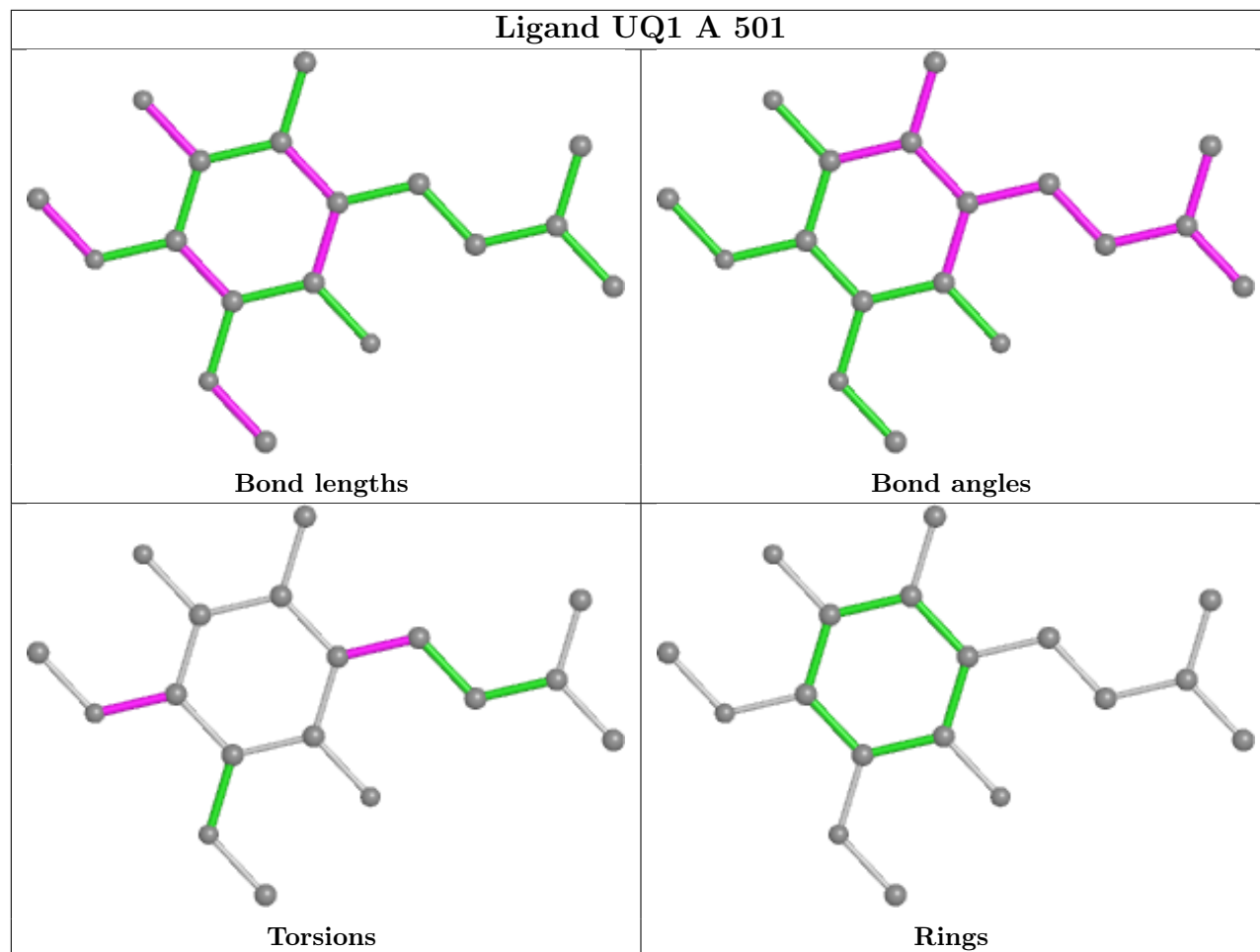
There are no ring outliers.

1 monomer is involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	UQ1	11	0

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

equivalents in the CSD to analyse the geometry.



## 5.7 Other polymers [i](#)

There are no such residues in this entry.

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

There are no chain breaks in this entry.

## 6 Fit of model and data

### 6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	386/399 (96%)	0.29	24 (6%) <span style="border: 1px solid red; padding: 2px;">20</span> <span style="border: 1px solid red; padding: 2px;">16</span>	68, 107, 167, 227	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	274	TRP	9.7
1	A	241	ALA	7.6
1	A	243	SER	5.3
1	A	391	THR	5.2
1	A	248	ARG	5.1
1	A	244	ASP	3.4
1	A	257	LYS	3.4
1	A	169	LEU	3.3
1	A	182	LEU	3.2
1	A	392	HIS	3.2
1	A	151	ARG	3.2
1	A	175	MET	3.0
1	A	34	GLU	2.9
1	A	170	TRP	2.8
1	A	255	LEU	2.6
1	A	7	LEU	2.5
1	A	383	GLU	2.4
1	A	148	SER	2.3
1	A	242	THR	2.3
1	A	334	ASP	2.1
1	A	240	PHE	2.1
1	A	249	PHE	2.1
1	A	270	ALA	2.1
1	A	258	TRP	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains [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
1	CRO	A	66	22/23	0.92	0.23	60,91,113,116	0

## 6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

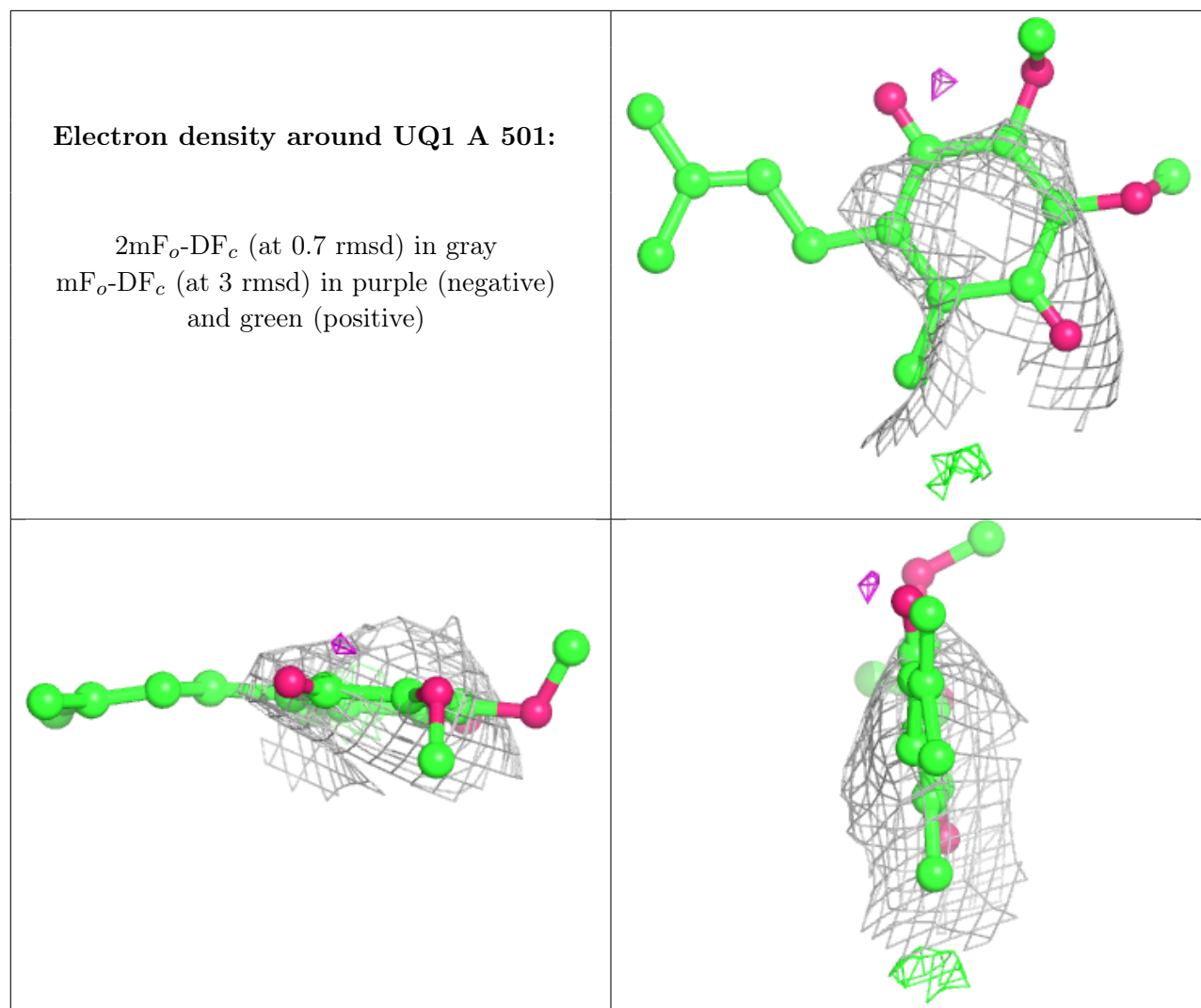
## 6.4 Ligands [i](#)

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	UQ1	A	501	18/18	0.80	0.79	127,152,162,164	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.