



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 02:15 AM UTC

PDB ID : 8TRY / pdb_00008try
Title : Crystal Structure of Mtb Pks13 Thioesterase domain in complex with inhibitor X20348
Authors : Krieger, I.V.; Sacchettini, J.C.
Deposited on : 2023-08-10
Resolution : 2.35 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 types of validation reports are described at

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

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtrriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

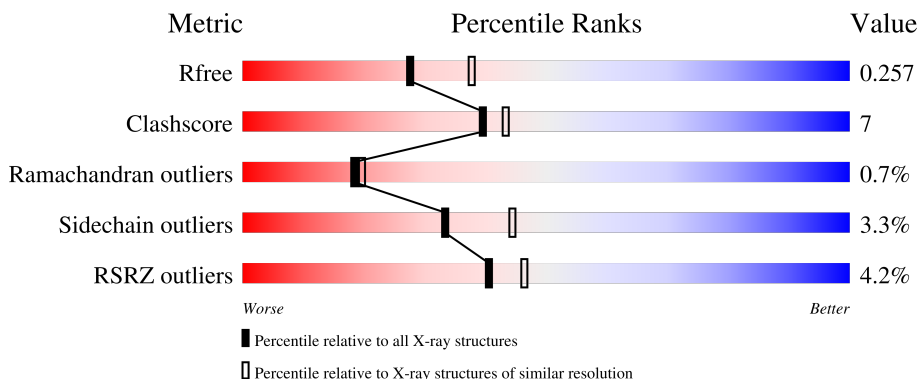
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION


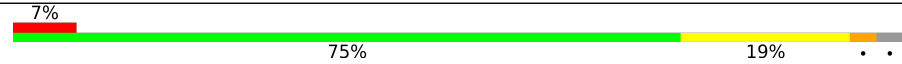
The reported resolution of this entry is 2.35 Å.

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}	180053	1596 (2.36-2.36)
Clashscore	190562	1663 (2.36-2.36)
Ramachandran outliers	187476	1646 (2.36-2.36)
Sidechain outliers	187428	1646 (2.36-2.36)
RSRZ outliers	180081	1598 (2.36-2.36)

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 $\leq 5\%$. 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	286	
1	B	286	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4582 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.

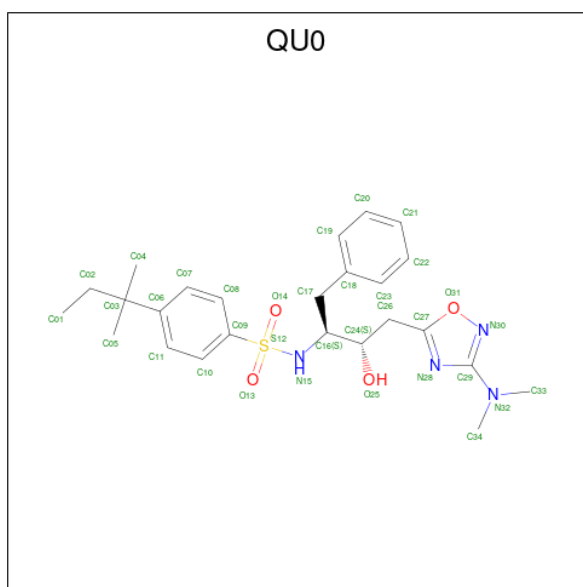
- Molecule 1 is a protein called Polyketide synthase Pks13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	277	Total	C	N	O	S	0	0	0
			2177	1381	379	411	6			
1	B	277	Total	C	N	O	S	0	0	0
			2177	1381	379	411	6			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1448	SER	-	expression tag	UNP I6X8D2
A	1449	ASN	-	expression tag	UNP I6X8D2
A	1450	ALA	-	expression tag	UNP I6X8D2
B	1448	SER	-	expression tag	UNP I6X8D2
B	1449	ASN	-	expression tag	UNP I6X8D2
B	1450	ALA	-	expression tag	UNP I6X8D2

- Molecule 2 is N-{(2S,3S)-4-[3-(dimethylamino)-1,2,4-oxadiazol-5-yl]-3-hydroxy-1-phenylbutan-2-yl}-4-(2-methylbutan-2-yl)benzene-1-sulfonamide (CCD ID: QU0) (formula: C₂₅H₃₄N₄O₄S) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
2	A	1	Total	C	N	O	S	0	0
			34	25	4	4	1		
2	A	1	Total	C	N	O	S	0	0
			34	25	4	4	1		
2	B	1	Total	C	N	O	S	0	0
			34	25	4	4	1		
2	B	1	Total	C	N	O	S	0	0
			34	25	4	4	1		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	51	Total	O	0	0
			51	51		
4	B	31	Total	O	0	0
			31	31		

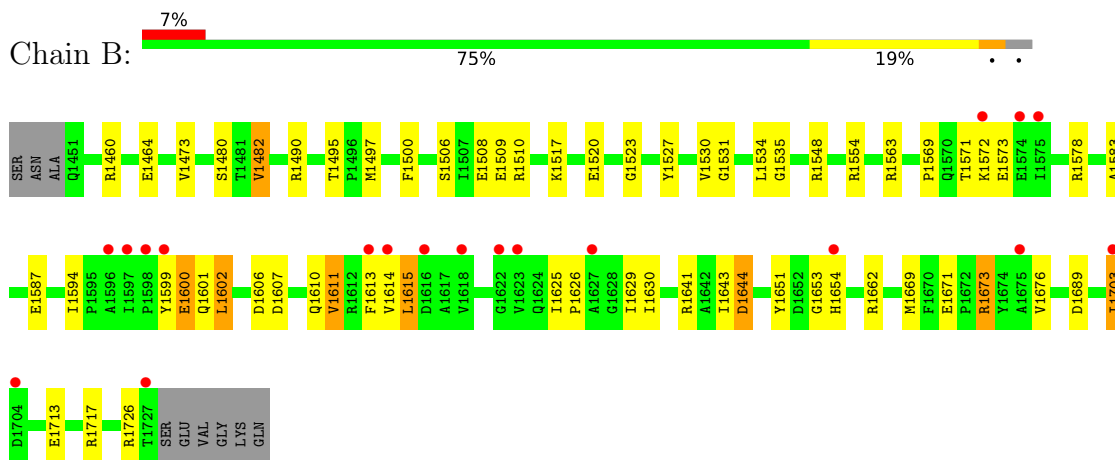
3 Residue-property plots [i](#)

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: Polyketide synthase Pks13



- Molecule 1: Polyketide synthase Pks13



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	54.86Å 96.62Å 130.97Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.36 – 2.35 48.36 – 2.35	Depositor EDS
% Data completeness (in resolution range)	84.6 (48.36-2.35) 84.6 (48.36-2.35)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.55 (at 2.34Å)	Xtrriage
Refinement program	PDB-REDO	Depositor
R, R_{free}	0.196 , 0.256 0.209 , 0.257	Depositor DCC
R_{free} test set	1217 reflections (4.09%)	wwPDB-VP
Wilson B-factor (Å ²)	37.3	Xtrriage
Anisotropy	0.567	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 34.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4582	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.89% 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: QU0, SO4

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	0.89	2/2227 (0.1%)	1.08	6/3024 (0.2%)
1	B	0.86	2/2227 (0.1%)	1.07	7/3024 (0.2%)
All	All	0.87	4/4454 (0.1%)	1.08	13/6048 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	8
1	B	0	11
All	All	0	19

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1497	MET	SD-CE	7.53	1.98	1.79
1	B	1669	MET	SD-CE	7.40	1.98	1.79
1	B	1497	MET	SD-CE	6.47	1.95	1.79
1	A	1644	ASP	CG-OD1	5.52	1.35	1.25

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1608	GLU	CB-CG-CD	7.57	125.46	112.60
1	B	1644	ASP	CA-CB-CG	7.27	119.87	112.60
1	A	1606	ASP	CA-CB-CG	6.33	118.93	112.60
1	B	1500	PHE	CA-CB-CG	-6.04	107.76	113.80
1	B	1606	ASP	CA-CB-CG	6.00	118.60	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1689	ASP	CA-CB-CG	5.55	118.15	112.60
1	B	1689	ASP	CA-CB-CG	5.43	118.03	112.60
1	B	1713	GLU	CB-CG-CD	5.34	121.68	112.60
1	A	1625	ILE	CA-C-O	5.24	122.60	119.19
1	B	1703	ILE	N-CA-CB	-5.24	104.10	112.07
1	A	1607	ASP	N-CA-C	5.16	116.91	111.28
1	A	1508	GLU	CB-CG-CD	5.12	121.31	112.60
1	B	1594	ILE	CB-CA-C	5.12	115.96	110.53

There are no chirality outliers.

All (19) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1457	ARG	Sidechain
1	A	1510	ARG	Sidechain
1	A	1548	ARG	Sidechain
1	A	1554	ARG	Sidechain
1	A	1563	ARG	Sidechain
1	A	1634	ARG	Sidechain
1	A	1641	ARG	Sidechain
1	A	1717	ARG	Sidechain
1	B	1490	ARG	Sidechain
1	B	1510	ARG	Sidechain
1	B	1548	ARG	Sidechain
1	B	1554	ARG	Sidechain
1	B	1563	ARG	Sidechain
1	B	1578	ARG	Sidechain
1	B	1641	ARG	Sidechain
1	B	1662	ARG	Sidechain
1	B	1673	ARG	Sidechain
1	B	1717	ARG	Sidechain
1	B	1726	ARG	Sidechain

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	2177	0	2127	35	0
1	B	2177	0	2127	25	0
2	A	68	0	0	0	0
2	B	68	0	0	0	0
3	B	10	0	0	0	0
4	A	51	0	0	7	0
4	B	31	0	0	3	0
All	All	4582	0	4254	59	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 7.

All (59) 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:1600:GLU:O	1:A:1601:GLN:HG2	1.92	0.70
1:B:1600:GLU:O	1:B:1601:GLN:HG2	1.96	0.66
1:A:1641:ARG:O	1:A:1645:THR:HG23	1.96	0.66
1:A:1480:SER:OG	1:A:1482:VAL:HG13	1.97	0.64
1:B:1599:TYR:O	1:B:1600:GLU:HB2	2.01	0.60
1:A:1608:GLU:HG3	4:A:1934:HOH:O	2.01	0.59
1:A:1512:GLN:NE2	4:A:1903:HOH:O	2.37	0.57
1:A:1611:VAL:HG22	4:A:1934:HOH:O	2.04	0.57
1:A:1678:GLN:HG3	4:A:1920:HOH:O	2.05	0.56
1:A:1451:GLN:C	1:A:1452:ILE:HD12	2.32	0.55
1:A:1498:TYR:CG	1:A:1521:MET:HE1	2.42	0.55
1:B:1625:ILE:HG22	1:B:1630:ILE:HG13	1.89	0.55
1:A:1569:PRO:O	1:A:1570:GLN:HG2	2.08	0.54
1:B:1626:PRO:HG2	1:B:1629:ILE:HD12	1.89	0.53
1:A:1473:VAL:HA	1:A:1530:VAL:O	2.09	0.53
1:A:1583:ALA:O	1:A:1587:GLU:HG3	2.09	0.52
1:B:1615:LEU:HD21	1:B:1630:ILE:HG23	1.91	0.52
1:A:1618:VAL:HG11	1:A:1625:ILE:HD11	1.92	0.52
1:A:1548:ARG:HA	1:B:1482:VAL:HG22	1.91	0.52
1:B:1473:VAL:HA	1:B:1530:VAL:O	2.10	0.51
1:B:1583:ALA:O	1:B:1587:GLU:HG3	2.10	0.51
1:A:1569:PRO:O	1:A:1570:GLN:CB	2.58	0.51
1:A:1534:LEU:HB2	1:A:1643:ILE:HD11	1.93	0.50
1:B:1506:SER:OG	1:B:1509:GLU:HG3	2.12	0.49
1:B:1480:SER:OG	1:B:1482:VAL:HG13	2.14	0.48
1:B:1534:LEU:HB2	1:B:1643:ILE:HD11	1.96	0.48
1:A:1460:ARG:NH1	1:A:1495:THR:O	2.46	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1615:LEU:CD2	4:A:1950:HOH:O	2.62	0.47
1:B:1517:LYS:HD2	4:B:1919:HOH:O	2.15	0.47
1:A:1615:LEU:HD22	4:A:1950:HOH:O	2.15	0.46
1:A:1626:PRO:HG2	1:A:1629:ILE:HD12	1.96	0.46
1:A:1451:GLN:HG2	1:A:1452:ILE:HD12	1.98	0.46
1:B:1610:GLN:O	1:B:1614:VAL:HG23	2.16	0.45
1:B:1599:TYR:O	1:B:1600:GLU:CB	2.64	0.45
1:A:1659:MET:HE3	1:A:1679:PRO:O	2.16	0.45
1:B:1607:ASP:O	1:B:1611:VAL:HG13	2.17	0.45
1:A:1522:GLN:O	1:A:1522:GLN:HG3	2.17	0.45
1:B:1654:HIS:HE1	4:B:1927:HOH:O	1.99	0.45
1:A:1607:ASP:O	1:A:1611:VAL:HG13	2.17	0.44
1:A:1498:TYR:CD1	1:A:1521:MET:HE1	2.53	0.44
1:B:1569:PRO:HB2	1:B:1571:THR:HG23	1.98	0.44
1:A:1523:GLY:O	1:A:1527:TYR:OH	2.35	0.43
1:A:1625:ILE:HG22	1:A:1630:ILE:HG13	2.00	0.43
1:B:1531:GLY:HA3	1:B:1535:GLY:C	2.44	0.43
1:A:1563:ARG:NE	4:A:1906:HOH:O	2.51	0.43
1:A:1569:PRO:O	1:A:1570:GLN:CG	2.66	0.43
1:B:1602:LEU:HD12	1:B:1613:PHE:CD1	2.53	0.43
1:B:1654:HIS:CE1	4:B:1927:HOH:O	2.72	0.43
1:B:1460:ARG:NH1	1:B:1495:THR:O	2.48	0.43
1:B:1625:ILE:CG2	1:B:1630:ILE:HG13	2.48	0.43
1:A:1703:ILE:HA	1:A:1708:ILE:HD13	2.00	0.42
1:B:1671:GLU:OE2	1:B:1673:ARG:HD3	2.20	0.42
1:B:1651:TYR:CE2	1:B:1653:GLY:HA3	2.55	0.42
1:A:1569:PRO:O	1:A:1570:GLN:HB2	2.20	0.42
1:B:1523:GLY:O	1:B:1527:TYR:OH	2.36	0.41
1:A:1469:VAL:HA	1:A:1470:PRO:HD3	1.95	0.41
1:A:1452:ILE:HD12	1:A:1452:ILE:N	2.36	0.41
1:A:1531:GLY:HA3	1:A:1535:GLY:C	2.47	0.40
1:A:1651:TYR:CE2	1:A:1653:GLY:HA3	2.56	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	275/286 (96%)	263 (96%)	10 (4%)	2 (1%)	18	20
1	B	275/286 (96%)	268 (98%)	5 (2%)	2 (1%)	18	20
All	All	550/572 (96%)	531 (96%)	15 (3%)	4 (1%)	18	20

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1569	PRO
1	A	1570	GLN
1	B	1602	LEU
1	B	1600	GLU

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	224/231 (97%)	220 (98%)	4 (2%)	51	66
1	B	224/231 (97%)	213 (95%)	11 (5%)	22	28
All	All	448/462 (97%)	433 (97%)	15 (3%)	33	44

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1482	VAL
1	A	1568	ILE
1	A	1611	VAL
1	A	1662	ARG
1	B	1464	GLU
1	B	1482	VAL
1	B	1508	GLU
1	B	1520	GLU

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Mol	Chain	Res	Type
1	B	1572	LYS
1	B	1573	GLU
1	B	1611	VAL
1	B	1615	LEU
1	B	1644	ASP
1	B	1676	VAL
1	B	1703	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	1451	GLN
1	B	1601	GLN
1	B	1721	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

6 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	QU0	A	1801	-	36,36,36	2.58	10 (27%)	46,52,52	3.42	13 (28%)
2	QU0	B	1802	-	36,36,36	2.32	9 (25%)	46,52,52	2.98	14 (30%)
3	SO4	B	1804	-	4,4,4	0.29	0	6,6,6	0.12	0
3	SO4	B	1803	-	4,4,4	0.27	0	6,6,6	0.19	0
2	QU0	A	1802	-	36,36,36	2.70	8 (22%)	46,52,52	2.95	11 (23%)
2	QU0	B	1801	-	36,36,36	2.23	11 (30%)	46,52,52	3.32	14 (30%)

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	QU0	A	1801	-	-	5/36/36/36	0/3/3/3
2	QU0	B	1802	-	-	3/36/36/36	0/3/3/3
2	QU0	B	1801	-	-	0/36/36/36	0/3/3/3
2	QU0	A	1802	-	-	0/36/36/36	0/3/3/3

All (38) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1801	QU0	S12-N15	8.23	1.75	1.61
2	B	1801	QU0	S12-N15	7.74	1.74	1.61
2	A	1802	QU0	O14-S12	7.64	1.52	1.43
2	B	1802	QU0	S12-N15	7.09	1.73	1.61
2	A	1802	QU0	S12-N15	6.57	1.72	1.61
2	A	1801	QU0	C29-N30	6.16	1.35	1.30
2	A	1802	QU0	C09-S12	5.76	1.85	1.76
2	A	1802	QU0	O13-S12	5.69	1.50	1.43
2	A	1802	QU0	C29-N30	4.75	1.34	1.30
2	A	1801	QU0	C27-N28	4.72	1.34	1.29
2	A	1801	QU0	C09-S12	4.72	1.83	1.76
2	B	1802	QU0	C27-N28	4.68	1.34	1.29
2	B	1801	QU0	C29-N30	4.68	1.34	1.30
2	B	1801	QU0	O14-S12	4.63	1.48	1.43
2	A	1802	QU0	O31-N30	-4.49	1.34	1.42
2	B	1802	QU0	O13-S12	4.31	1.48	1.43
2	A	1801	QU0	C29-N32	4.30	1.45	1.35
2	A	1802	QU0	C29-N32	4.26	1.45	1.35
2	B	1802	QU0	C29-N30	4.16	1.33	1.30
2	B	1802	QU0	O31-N30	-4.16	1.34	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1802	QU0	O14-S12	4.12	1.48	1.43
2	B	1802	QU0	C29-N32	4.11	1.45	1.35
2	B	1802	QU0	C09-S12	4.08	1.82	1.76
2	B	1801	QU0	C29-N32	4.06	1.45	1.35
2	A	1801	QU0	O14-S12	4.01	1.48	1.43
2	A	1801	QU0	O13-S12	4.01	1.48	1.43
2	A	1802	QU0	C27-N28	3.36	1.32	1.29
2	B	1801	QU0	C27-N28	3.00	1.32	1.29
2	B	1801	QU0	O31-N30	-2.96	1.37	1.42
2	B	1801	QU0	C26-C24	-2.81	1.49	1.53
2	A	1801	QU0	C24-C16	-2.79	1.49	1.53
2	B	1801	QU0	C09-S12	2.75	1.80	1.76
2	B	1801	QU0	O13-S12	2.70	1.46	1.43
2	B	1802	QU0	O31-C27	-2.61	1.31	1.34
2	A	1801	QU0	C10-C09	2.26	1.42	1.38
2	B	1801	QU0	O31-C27	-2.16	1.32	1.34
2	A	1801	QU0	O31-N30	-2.14	1.38	1.42
2	B	1801	QU0	C24-C16	-2.13	1.50	1.53

All (52) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1801	QU0	O31-N30-C29	14.47	110.22	101.09
2	A	1801	QU0	O31-N30-C29	13.43	109.56	101.09
2	A	1801	QU0	O14-S12-O13	-10.38	106.91	119.52
2	B	1802	QU0	O14-S12-O13	-10.02	107.35	119.52
2	B	1802	QU0	O31-N30-C29	9.45	107.05	101.09
2	A	1802	QU0	O14-S12-O13	-9.44	108.06	119.52
2	A	1802	QU0	O31-N30-C29	8.71	106.59	101.09
2	B	1801	QU0	O14-S12-O13	-8.65	109.02	119.52
2	A	1801	QU0	N28-C29-N30	-8.00	106.14	118.63
2	B	1801	QU0	N28-C29-N30	-7.73	106.56	118.63
2	A	1802	QU0	O14-S12-N15	6.40	118.50	106.88
2	B	1802	QU0	N28-C29-N30	-6.39	108.64	118.63
2	A	1801	QU0	O14-S12-C09	6.25	115.87	107.98
2	A	1802	QU0	N28-C29-N30	-6.10	109.09	118.63
2	A	1802	QU0	O31-C27-C26	5.72	124.71	117.11
2	B	1801	QU0	O31-C27-C26	5.57	124.50	117.11
2	B	1802	QU0	O31-C27-N28	-5.21	109.40	113.62
2	A	1802	QU0	C27-O31-N30	5.17	108.92	106.10
2	A	1802	QU0	O31-C27-N28	-5.13	109.47	113.62
2	B	1802	QU0	C27-O31-N30	5.12	108.89	106.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1801	QU0	O13-S12-N15	5.04	116.03	106.88
2	A	1801	QU0	O31-C27-C26	4.78	123.46	117.11
2	B	1802	QU0	O31-C27-C26	4.46	123.03	117.11
2	B	1801	QU0	C18-C17-C16	-4.39	105.96	113.40
2	B	1801	QU0	C27-O31-N30	-4.21	103.81	106.10
2	A	1801	QU0	O31-C27-N28	-4.21	110.21	113.62
2	B	1802	QU0	O13-S12-C09	4.17	113.23	107.98
2	B	1802	QU0	C10-C09-S12	-3.67	115.72	119.76
2	A	1801	QU0	C18-C17-C16	-3.44	107.56	113.40
2	B	1801	QU0	C10-C09-S12	-3.42	115.99	119.76
2	B	1801	QU0	C08-C09-S12	3.37	123.47	119.76
2	A	1801	QU0	C07-C06-C03	-3.32	115.53	121.52
2	B	1802	QU0	C18-C17-C16	-3.21	107.96	113.40
2	A	1802	QU0	C16-N15-S12	-3.11	115.59	121.96
2	B	1802	QU0	C11-C06-C03	-3.11	115.90	121.52
2	B	1801	QU0	O14-S12-N15	3.08	112.47	106.88
2	A	1801	QU0	C11-C06-C07	3.04	122.56	118.03
2	A	1802	QU0	C11-C06-C03	-2.77	116.52	121.52
2	B	1802	QU0	N32-C29-N28	2.71	126.87	119.34
2	B	1801	QU0	C09-S12-N15	2.68	111.47	107.79
2	A	1802	QU0	C11-C06-C07	2.68	122.02	118.03
2	B	1801	QU0	C11-C06-C03	-2.55	116.91	121.52
2	A	1801	QU0	N32-C29-N28	2.55	126.41	119.34
2	A	1801	QU0	C02-C03-C06	2.47	117.06	110.02
2	B	1802	QU0	C08-C09-S12	2.29	122.28	119.76
2	B	1801	QU0	N32-C29-N28	2.25	125.58	119.34
2	A	1801	QU0	C17-C16-C24	-2.20	108.08	111.66
2	B	1801	QU0	C23-C18-C19	2.15	121.42	118.23
2	A	1802	QU0	C17-C16-N15	2.12	113.26	110.26
2	B	1802	QU0	O13-S12-N15	2.07	110.64	106.88
2	B	1801	QU0	O25-C24-C16	2.03	113.89	109.88
2	B	1802	QU0	C34-N32-C33	2.00	122.28	115.87

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1801	QU0	C01-C02-C03-C04
2	A	1801	QU0	C01-C02-C03-C05
2	A	1801	QU0	C01-C02-C03-C06
2	B	1802	QU0	C02-C03-C06-C07
2	A	1801	QU0	C16-N15-S12-O14

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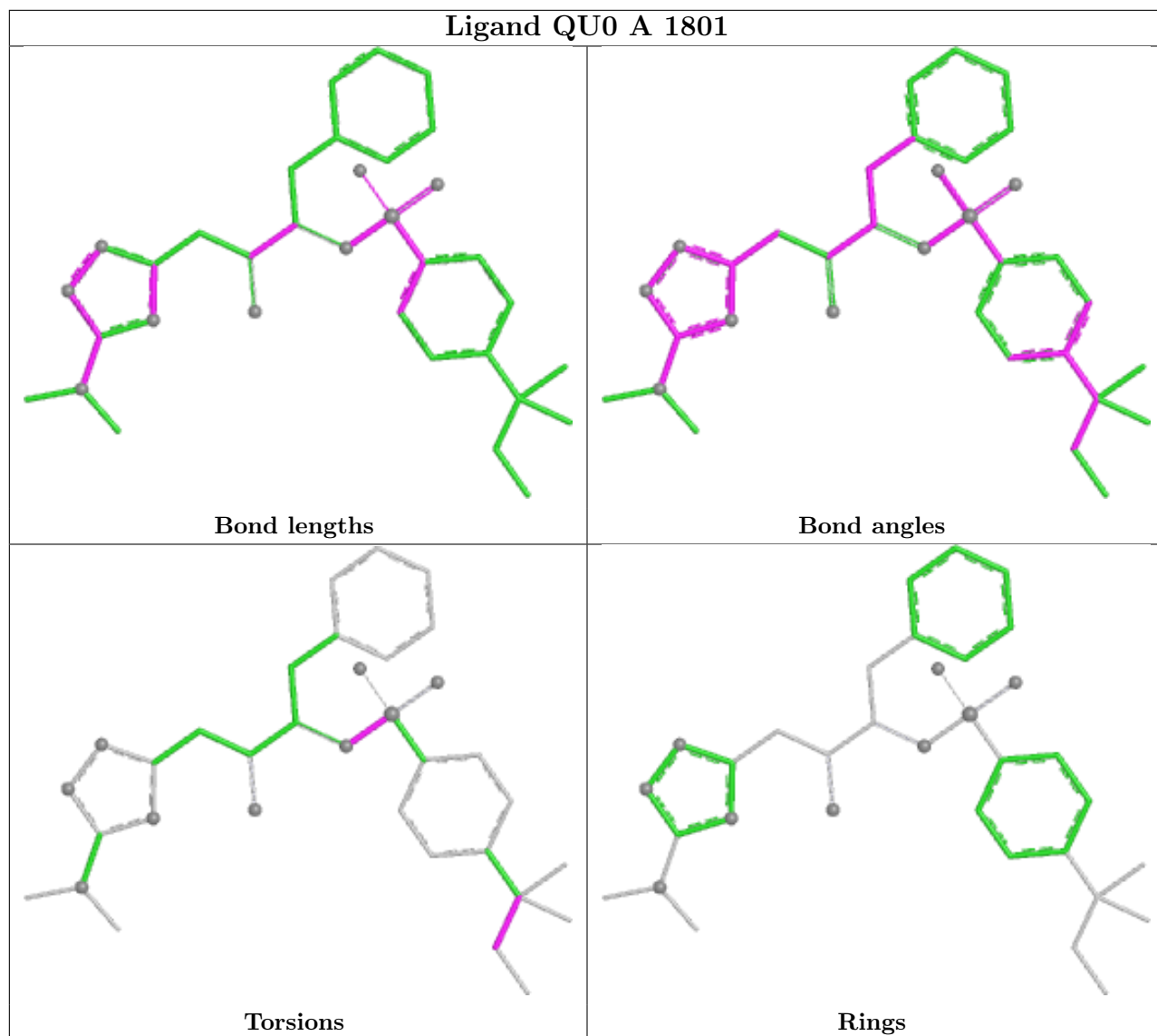
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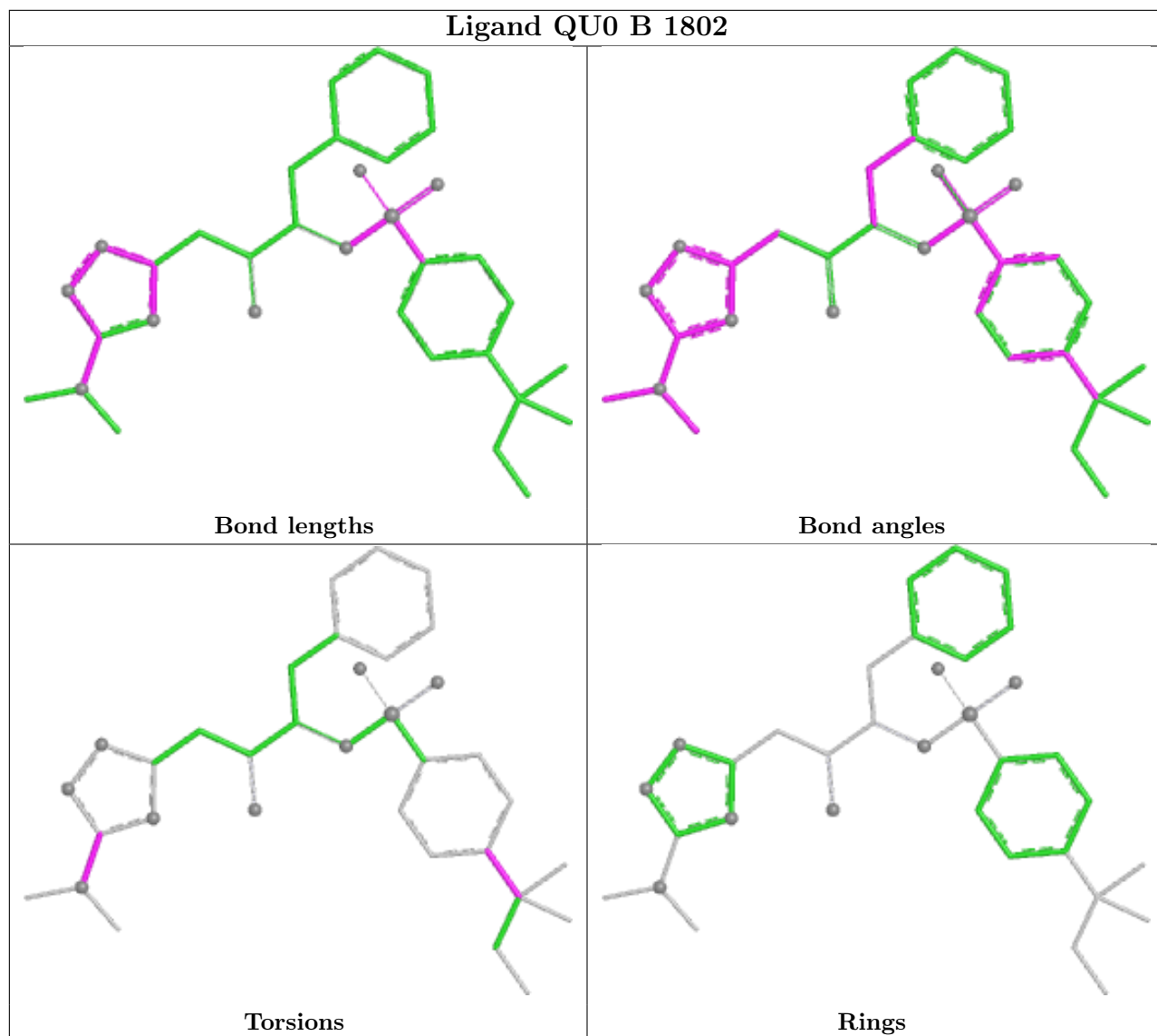
Mol	Chain	Res	Type	Atoms
2	B	1802	QU0	C02-C03-C06-C11
2	A	1801	QU0	C16-N15-S12-O13
2	B	1802	QU0	N30-C29-N32-C34

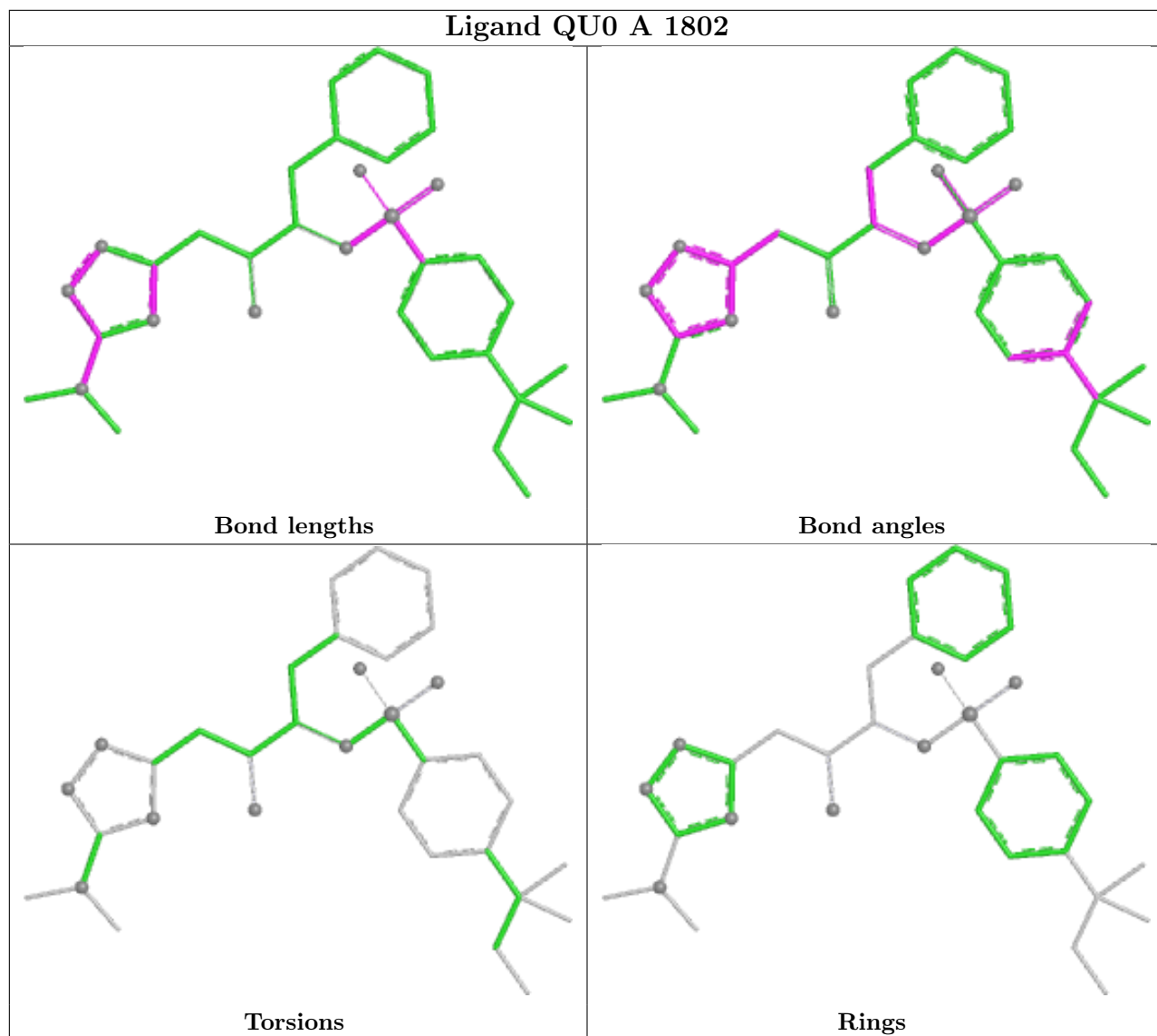
There are no ring outliers.

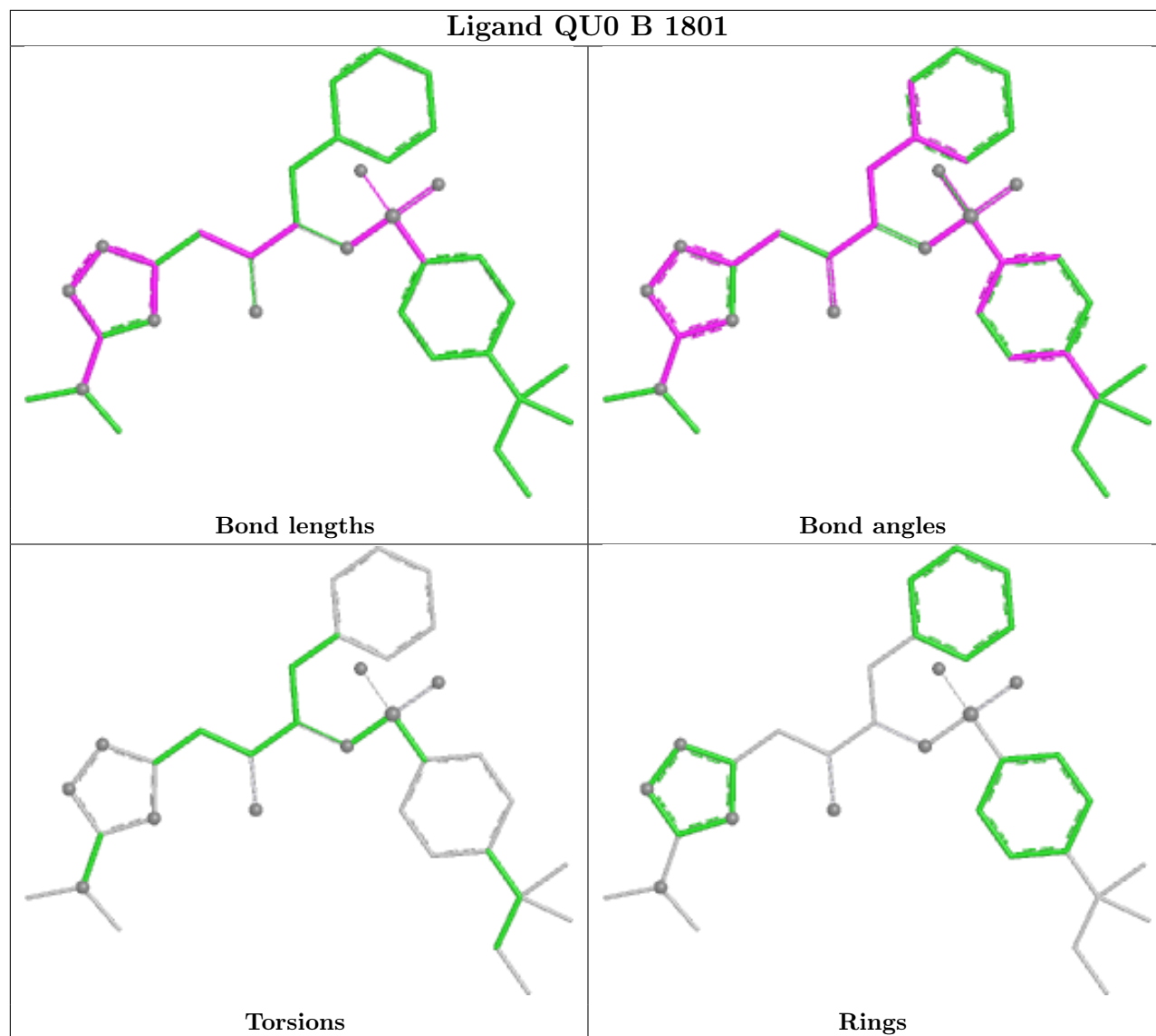
No monomer is involved in short contacts.

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	277/286 (96%)	0.12	4 (1%) 73 78	26, 41, 70, 107	0
1	B	277/286 (96%)	0.43	19 (6%) 23 26	25, 50, 88, 125	0
All	All	554/572 (96%)	0.28	23 (4%) 40 46	25, 45, 83, 125	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1453	ASP	4.0
1	B	1574	GLU	3.8
1	B	1572	LYS	3.4
1	B	1599	TYR	3.4
1	B	1598	PRO	3.1
1	A	1727	THR	3.0
1	B	1622	GLY	3.0
1	B	1675	ALA	3.0
1	B	1575	ILE	2.8
1	A	1644	ASP	2.8
1	B	1704	ASP	2.7
1	B	1616	ASP	2.5
1	B	1618	VAL	2.5
1	B	1597	ILE	2.5
1	B	1614	VAL	2.3
1	B	1623	VAL	2.2
1	A	1620	GLN	2.2
1	B	1596	ALA	2.2
1	B	1727	THR	2.2
1	B	1613	PHE	2.1
1	B	1627	ALA	2.1
1	B	1703	ILE	2.0
1	B	1654	HIS	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

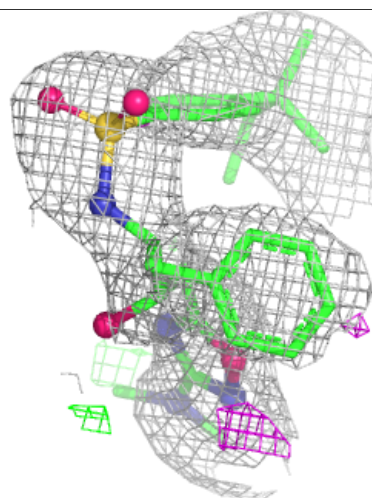
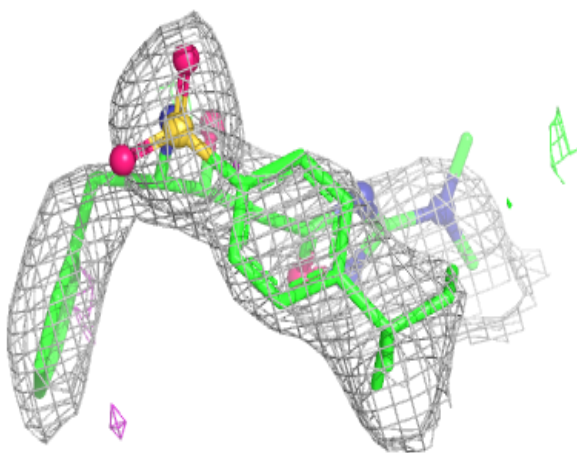
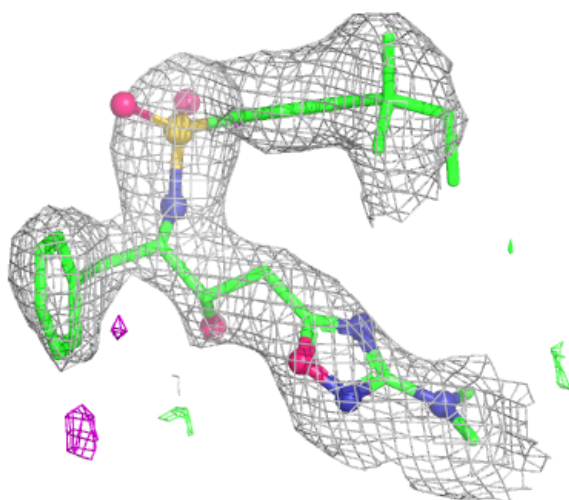
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

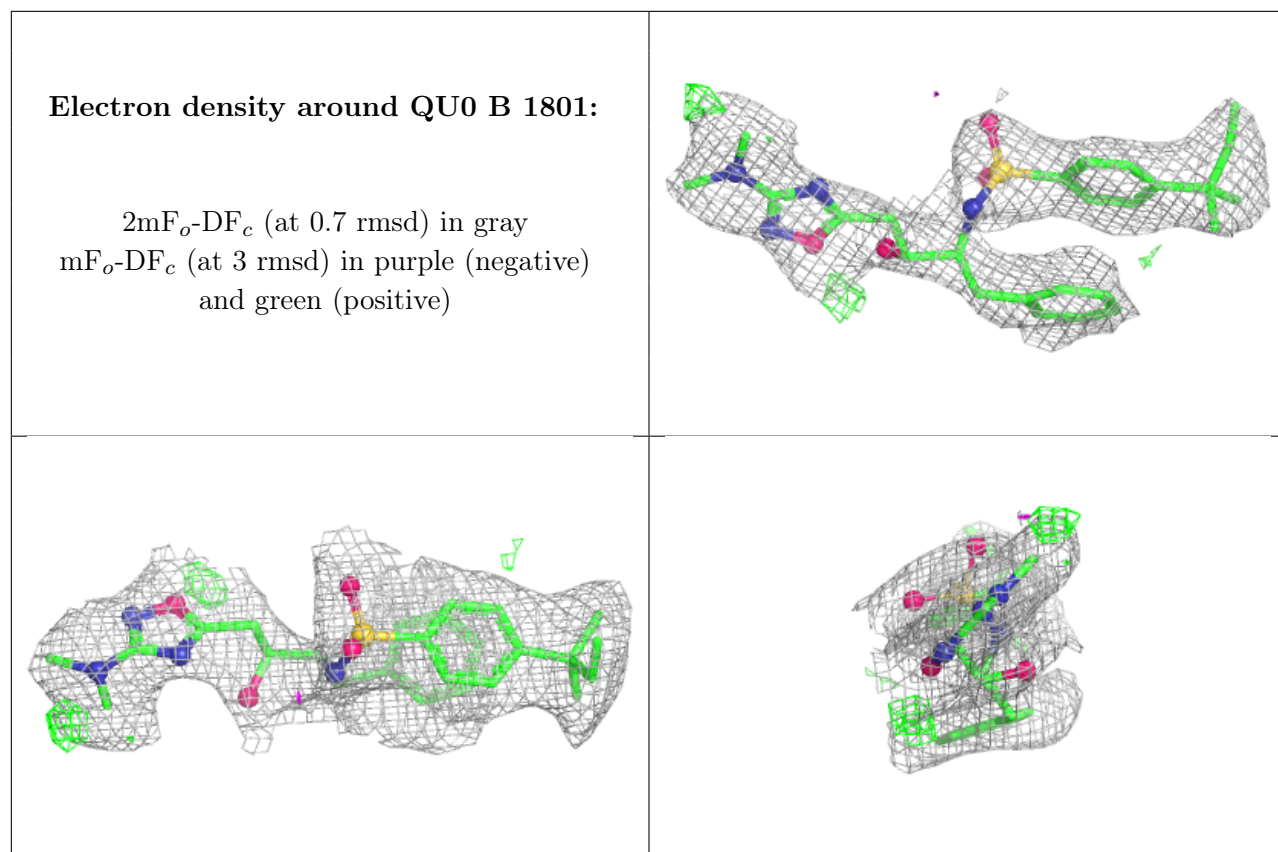
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	SO4	B	1804	5/5	0.78	0.12	80,89,99,104	0
3	SO4	B	1803	5/5	0.79	0.10	69,83,96,114	0
2	QU0	B	1802	34/34	0.93	0.14	40,58,82,86	0
2	QU0	B	1801	34/34	0.94	0.10	36,44,54,57	0
2	QU0	A	1802	34/34	0.95	0.10	31,44,69,76	0
2	QU0	A	1801	34/34	0.96	0.08	26,34,47,49	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 QU0 B 1802:

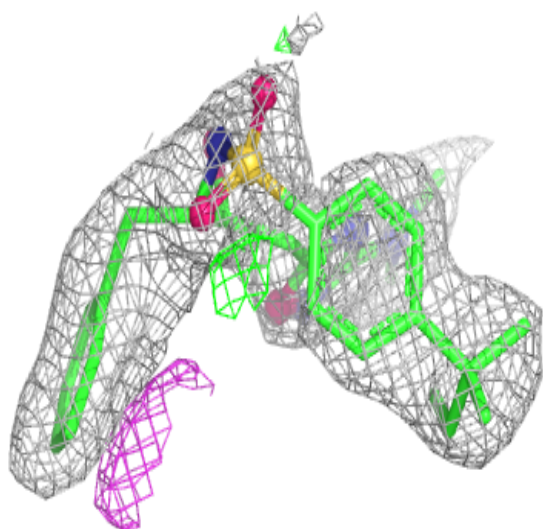
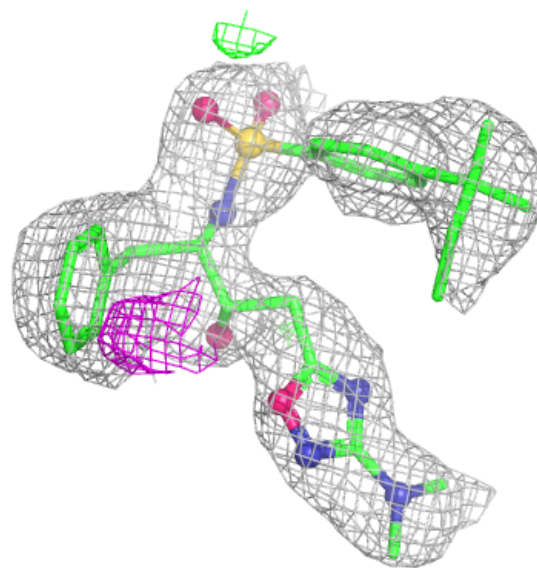
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

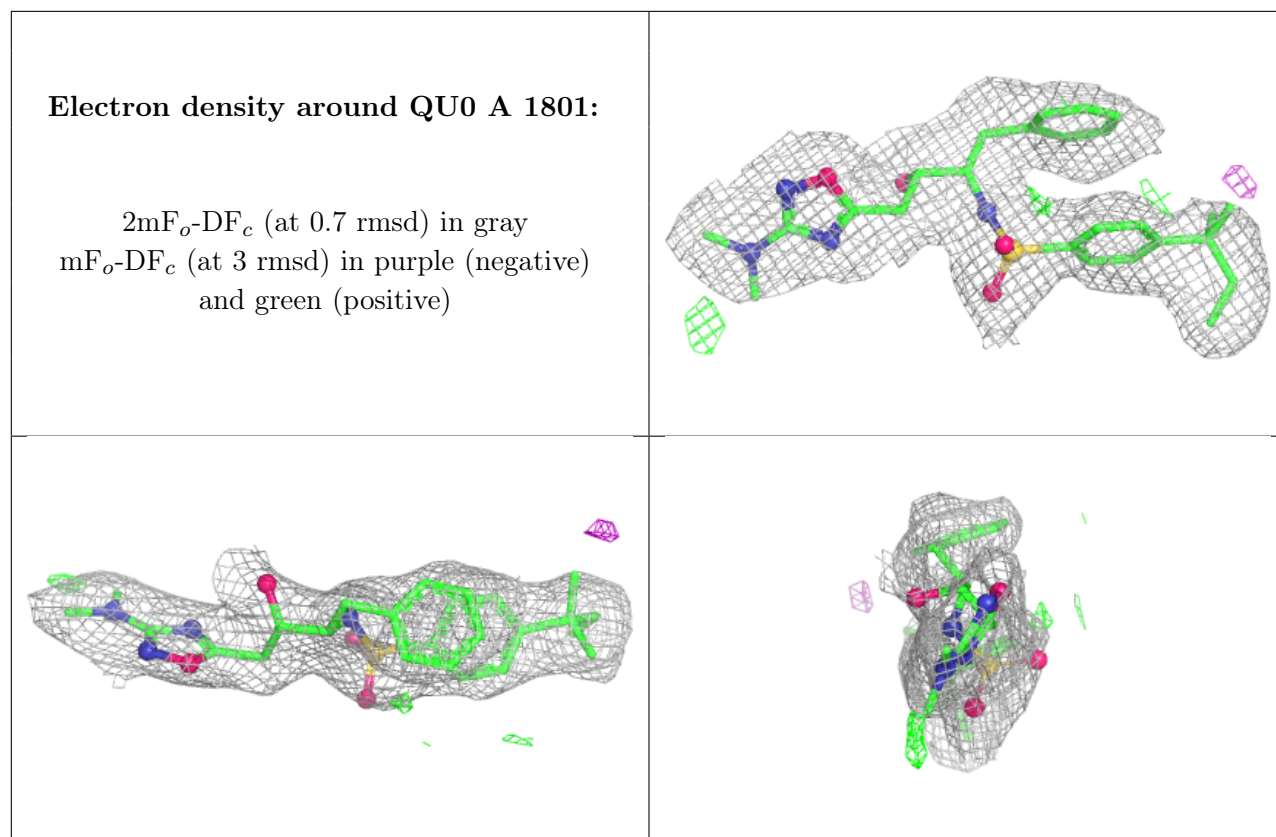




Electron density around QU0 A 1802:

$2mF_o-DF_c$ (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.