



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 15, 2022 – 02:04 PM EDT

PDB ID : 7UYR
Title : Crystal structure of TYK2 kinase domain in complex with compound 12
Authors : Toms, A.V.; Leit, S.; Greenwood, J.R.; Mondal, S.; Carriero, S.; Dahlgren, M.; Harriman, G.C.; Kennedy-Smith, J.J.; Kapeller, R.; Lawson, J.P.; Romero, D.L.; Shelley, M.; Wester, R.T.; Westlin, W.; Mc Elwee, J.J.; Miao, W.; Edmondson, S.D.; Masee, C.E.
Deposited on : 2022-05-07
Resolution : 2.15 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.29
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) : 2.29

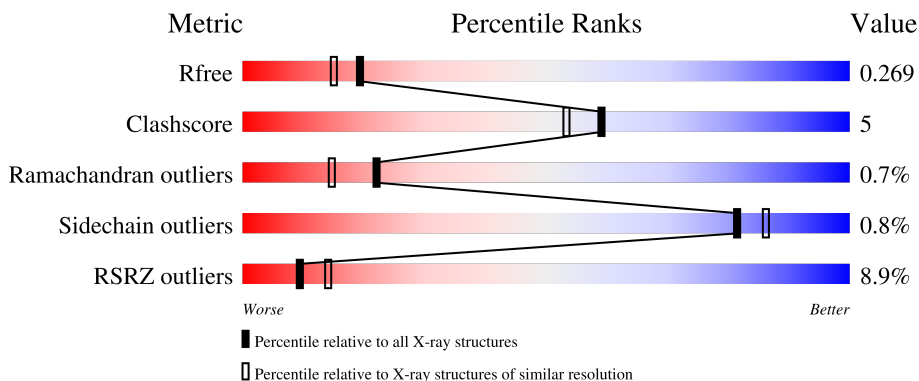
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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	1479 (2.16-2.16)
Clashscore	141614	1585 (2.16-2.16)
Ramachandran outliers	138981	1560 (2.16-2.16)
Sidechain outliers	138945	1559 (2.16-2.16)
RSRZ outliers	127900	1456 (2.16-2.16)

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	289	

2 Entry composition [i](#)

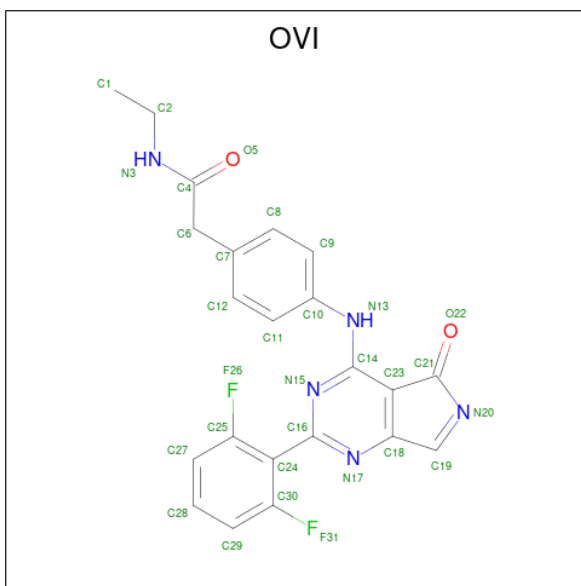
There are 3 unique types of molecules in this entry. The entry contains 2464 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Non-receptor tyrosine-protein kinase TYK2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
			Total	C	N	O	P				S
1	A	282	2337	1499	399	422	2	15	75	5	0

- Molecule 2 is 2-(4-{[2-(2,6-difluorophenyl)-5-oxo-5H-pyrrolo[3,4-d]pyrimidin-4-yl]amino}phenyl)-N-ethylacetamide (three-letter code: OVI) (formula: C₂₂H₁₇F₂N₅O₂) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	F	N	O		
2	A	1	44	32	2	7	3	0	1

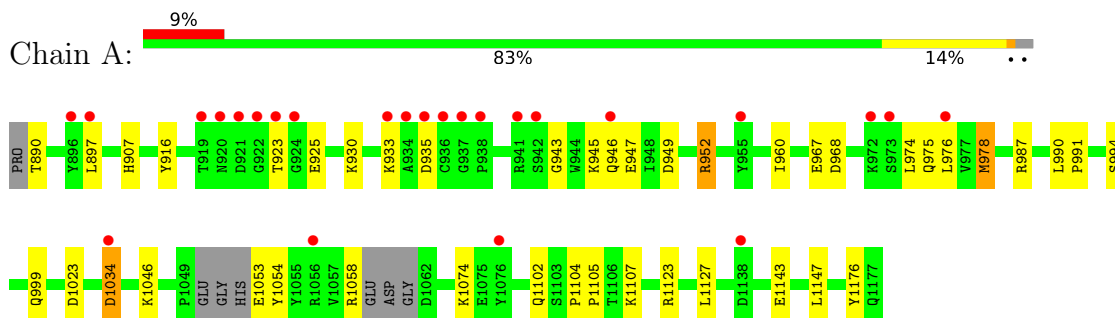
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	83	Total	O	0	0
			83	83		

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: Non-receptor tyrosine-protein kinase TYK2



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	36.19Å 74.24Å 106.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	34.30 – 2.15 34.28 – 2.15	Depositor EDS
% Data completeness (in resolution range)	95.5 (34.30-2.15) 95.5 (34.28-2.15)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.43 (at 2.16Å)	Xtrriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.200 , 0.259 0.209 , 0.269	Depositor DCC
R_{free} test set	1426 reflections (9.14%)	wwPDB-VP
Wilson B-factor (Å ²)	34.9	Xtrriage
Anisotropy	0.179	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 41.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	2464	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.59% 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: OVI, PTR

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	1.12	15/2361 (0.6%)	0.99	14/3190 (0.4%)

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	1

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	925	GLU	CB-CG	-19.09	1.15	1.52
1	A	933	LYS	CB-CG	-16.19	1.08	1.52
1	A	967	GLU	CB-CG	13.71	1.78	1.52
1	A	987	ARG	CD-NE	-12.22	1.25	1.46
1	A	1058	ARG	CB-CG	-12.07	1.20	1.52
1	A	1123	ARG	NE-CZ	11.58	1.48	1.33
1	A	890	THR	CB-CG2	-11.08	1.15	1.52
1	A	935	ASP	CB-CG	9.20	1.71	1.51
1	A	945	LYS	CE-NZ	-7.32	1.30	1.49
1	A	1074	LYS	CE-NZ	-6.86	1.31	1.49
1	A	1053	GLU	CG-CD	-6.71	1.41	1.51
1	A	923	THR	CB-OG1	6.41	1.56	1.43
1	A	923	THR	CB-CG2	-5.97	1.32	1.52
1	A	978[A]	MET	C-O	5.74	1.34	1.23
1	A	978[B]	MET	C-O	5.74	1.34	1.23

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	952	ARG	NE-CZ-NH1	-13.51	113.55	120.30
1	A	952	ARG	NE-CZ-NH2	11.28	125.94	120.30
1	A	890	THR	OG1-CB-CG2	10.08	133.19	110.00
1	A	933	LYS	CA-CB-CG	9.27	133.79	113.40
1	A	935	ASP	CB-CG-OD1	6.45	124.11	118.30
1	A	925	GLU	CA-CB-CG	6.35	127.38	113.40
1	A	952	ARG	CD-NE-CZ	6.28	132.39	123.60
1	A	923	THR	CA-CB-OG1	-6.10	96.18	109.00
1	A	949	ASP	CB-CG-OD1	6.09	123.78	118.30
1	A	1123	ARG	NE-CZ-NH2	5.92	123.26	120.30
1	A	987	ARG	CG-CD-NE	5.91	124.22	111.80
1	A	890	THR	CA-CB-OG1	-5.62	97.20	109.00
1	A	949	ASP	CB-CG-OD2	-5.25	113.58	118.30
1	A	935	ASP	CB-CG-OD2	-5.08	113.72	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	952	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	2337	0	2289	21	0
2	A	44	0	0	0	0
3	A	83	0	0	3	0
All	All	2464	0	2289	21	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 5.

All (21) 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:1102[B]:GLN:NE2	3:A:1301:HOH:O	1.94	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:976[A]:LEU:HG	1:A:978[A]:MET:SD	2.14	0.86
1:A:897:LEU:HD12	1:A:916:TYR:CD1	2.26	0.71
1:A:897:LEU:HD12	1:A:916:TYR:HD1	1.59	0.67
1:A:947:GLU:OE2	3:A:1302:HOH:O	2.10	0.67
1:A:1046:LYS:NZ	3:A:1303:HOH:O	2.37	0.56
1:A:999:GLN:NE2	1:A:1034:ASP:O	2.40	0.54
1:A:943:GLY:O	1:A:946:GLN:HB3	2.10	0.52
1:A:976[A]:LEU:CG	1:A:978[A]:MET:SD	2.95	0.51
1:A:1104:PRO:HB2	1:A:1105:PRO:HD3	1.92	0.50
1:A:930[A]:LYS:HD2	1:A:976[A]:LEU:HD23	1.93	0.49
1:A:930[A]:LYS:CD	1:A:976[A]:LEU:HD23	2.44	0.47
1:A:1102[A]:GLN:NE2	1:A:1107:LYS:HE2	2.29	0.47
1:A:960:ILE:HG21	1:A:978[B]:MET:CE	2.45	0.47
1:A:968:ASP:HB2	1:A:975:GLN:HG3	1.99	0.45
1:A:1147:LEU:HD23	1:A:1147:LEU:C	2.36	0.44
1:A:907:HIS:NE2	1:A:1023:ASP:OD2	2.48	0.43
1:A:930[A]:LYS:HE3	1:A:978[A]:MET:HG3	1.99	0.43
1:A:1143:GLU:OE2	1:A:1176:TYR:OH	2.22	0.42
1:A:990:LEU:N	1:A:991:PRO:CD	2.82	0.41
1:A:960:ILE:HG21	1:A:978[B]:MET:HE2	2.01	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	279/289 (96%)	270 (97%)	7 (2%)	2 (1%)	22 15

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	994	SER

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Mol	Chain	Res	Type
1	A	1034	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	247/252 (98%)	245 (99%)	2 (1%)	81 86

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	974	LEU
1	A	1127	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1150	ASN
1	A	1173	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](#)

2 non-standard protein/DNA/RNA residues 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	$\# Z > 2$
1	PTR	A	1055	1	15,16,17	0.63	0	19,22,24	0.57	0
1	PTR	A	1054	1	15,16,17	0.55	0	19,22,24	0.72	1 (5%)

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	PTR	A	1055	1	-	0/10/11/13	0/1/1/1
1	PTR	A	1054	1	-	0/10/11/13	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	1054	PTR	CB-CA-C	-2.25	107.25	111.47

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 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	OVI	A	1201[A]	-	33,34,34	2.50	9 (27%)	40,48,48	1.99	10 (25%)
2	OVI	A	1201[B]	-	33,34,34	2.50	9 (27%)	40,48,48	2.36	16 (40%)

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	OVI	A	1201[A]	-	-	1/15/24/24	0/4/4/4
2	OVI	A	1201[B]	-	-	2/15/24/24	0/4/4/4

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1201[A]	OVI	C6-C7	-8.40	1.37	1.51
2	A	1201[B]	OVI	C6-C7	-8.13	1.38	1.51
2	A	1201[A]	OVI	C24-C16	-6.23	1.39	1.49
2	A	1201[B]	OVI	C24-C16	-6.23	1.39	1.49
2	A	1201[A]	OVI	C18-C19	-6.18	1.34	1.42
2	A	1201[B]	OVI	C18-C19	-6.18	1.34	1.42
2	A	1201[A]	OVI	C23-C18	-4.04	1.36	1.41
2	A	1201[B]	OVI	C23-C18	-4.04	1.36	1.41
2	A	1201[A]	OVI	C14-N15	3.25	1.40	1.34
2	A	1201[B]	OVI	C14-N15	3.25	1.40	1.34
2	A	1201[A]	OVI	C23-C14	-2.66	1.36	1.42
2	A	1201[B]	OVI	C23-C14	-2.66	1.36	1.42
2	A	1201[A]	OVI	C21-N20	-2.56	1.33	1.37
2	A	1201[B]	OVI	C21-N20	-2.56	1.33	1.37
2	A	1201[B]	OVI	C10-N13	-2.56	1.35	1.40
2	A	1201[A]	OVI	C16-N15	2.47	1.39	1.34
2	A	1201[B]	OVI	C16-N15	2.47	1.39	1.34
2	A	1201[A]	OVI	C10-N13	-2.45	1.35	1.40

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1201[B]	OVI	N13-C14-N15	5.60	127.45	119.12

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1201[A]	OVI	N13-C14-N15	5.01	126.58	119.12
2	A	1201[A]	OVI	C25-C24-C30	4.81	119.28	114.56
2	A	1201[B]	OVI	C25-C24-C30	4.81	119.28	114.56
2	A	1201[B]	OVI	C6-C7-C8	-4.07	115.07	120.89
2	A	1201[B]	OVI	C6-C7-C12	3.98	126.59	120.89
2	A	1201[B]	OVI	C6-C4-N3	-3.97	110.83	116.19
2	A	1201[A]	OVI	C27-C25-C24	-3.87	119.22	123.94
2	A	1201[B]	OVI	C27-C25-C24	-3.87	119.22	123.94
2	A	1201[A]	OVI	C18-N17-C16	3.58	119.70	116.30
2	A	1201[B]	OVI	C18-N17-C16	3.58	119.70	116.30
2	A	1201[A]	OVI	N17-C16-N15	-3.26	119.96	125.23
2	A	1201[B]	OVI	N17-C16-N15	-3.26	119.96	125.23
2	A	1201[A]	OVI	C24-C16-N17	3.11	120.78	116.70
2	A	1201[B]	OVI	C24-C16-N17	3.11	120.78	116.70
2	A	1201[A]	OVI	O22-C21-C23	3.11	127.03	122.65
2	A	1201[B]	OVI	O22-C21-C23	3.11	127.03	122.65
2	A	1201[B]	OVI	C11-C10-N13	3.03	130.79	120.64
2	A	1201[B]	OVI	C9-C10-N13	-2.60	111.93	120.64
2	A	1201[B]	OVI	C1-C2-N3	-2.53	102.80	111.52
2	A	1201[A]	OVI	C14-N15-C16	2.39	120.40	116.10
2	A	1201[B]	OVI	C14-N15-C16	2.39	120.40	116.10
2	A	1201[A]	OVI	C29-C30-C24	-2.31	121.11	123.94
2	A	1201[B]	OVI	C29-C30-C24	-2.31	121.11	123.94
2	A	1201[A]	OVI	C23-C14-N15	-2.11	117.12	122.22
2	A	1201[B]	OVI	C23-C14-N15	-2.11	117.12	122.22

There are no chirality outliers.

All (3) torsion outliers are listed below:

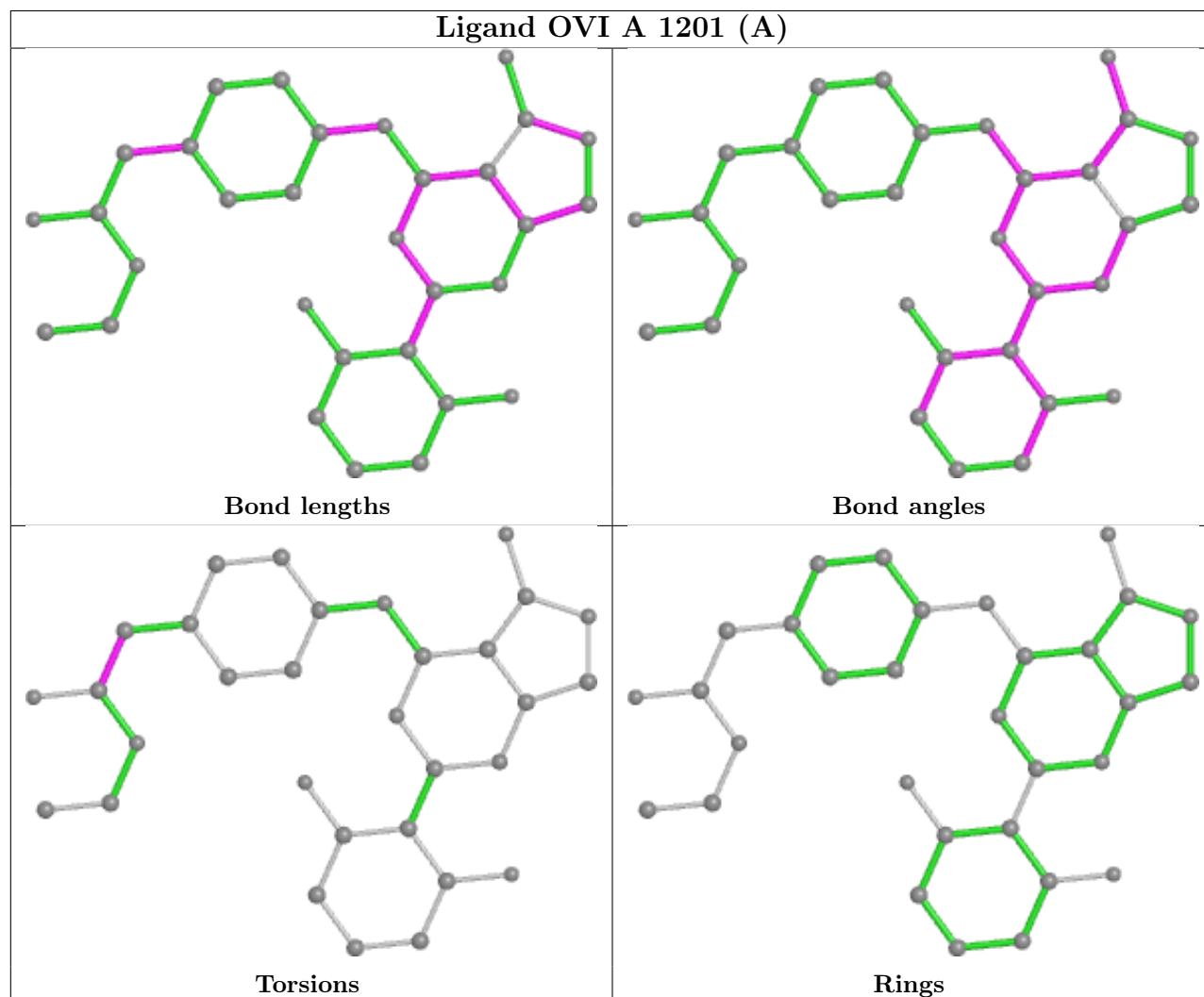
Mol	Chain	Res	Type	Atoms
2	A	1201[B]	OVI	C4-C6-C7-C12
2	A	1201[B]	OVI	C4-C6-C7-C8
2	A	1201[A]	OVI	O5-C4-C6-C7

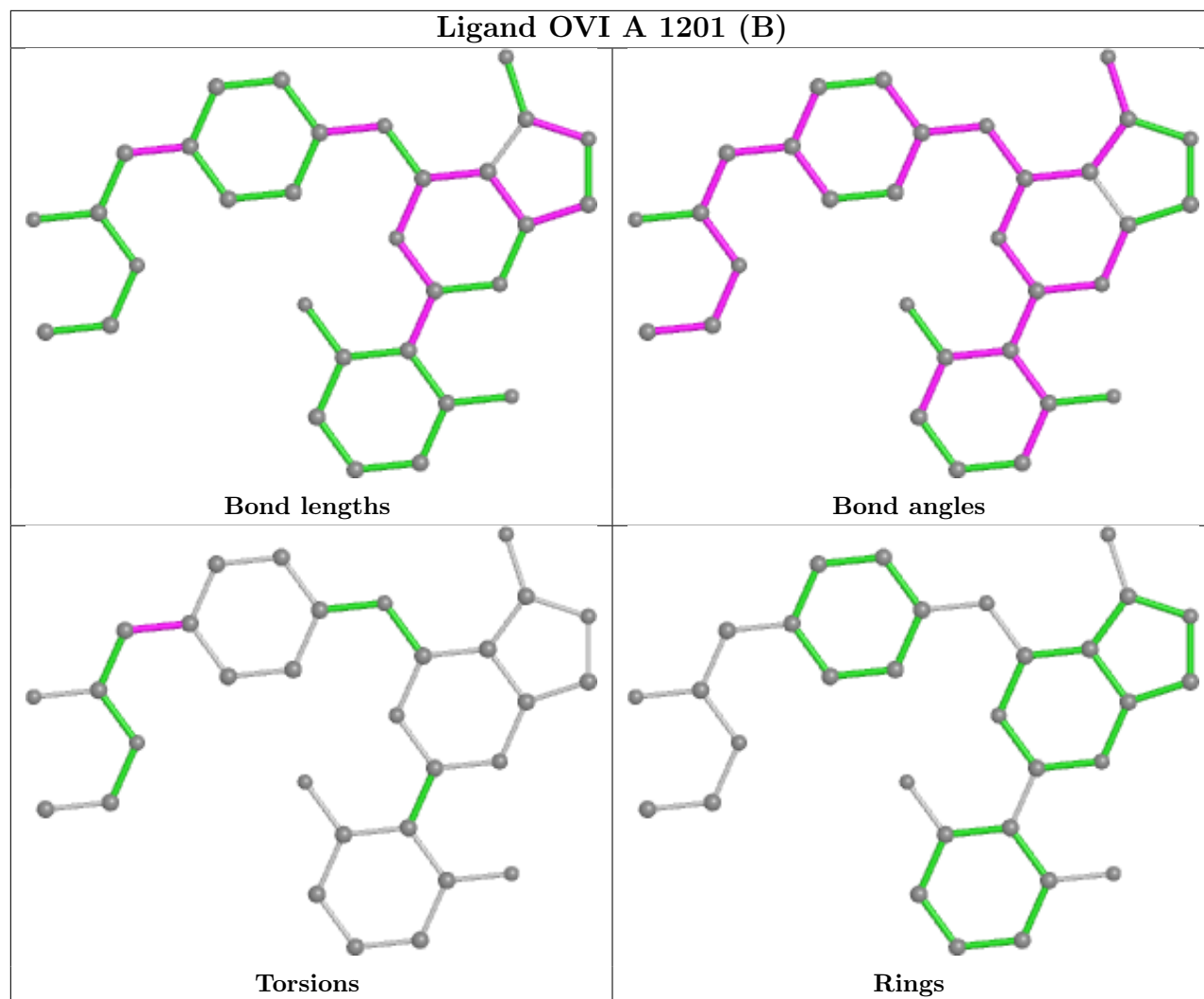
There are no ring outliers.

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	280/289 (96%)	0.55	25 (8%) 9 14	22, 39, 69, 84	24 (8%)

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	934	ALA	7.8
1	A	920	ASN	3.8
1	A	1056	ARG	3.8
1	A	923	THR	3.3
1	A	938	PRO	3.2
1	A	1034	ASP	3.1
1	A	1076	TYR	3.1
1	A	921	ASP	3.0
1	A	941	ARG	3.0
1	A	933	LYS	2.9
1	A	919	THR	2.8
1	A	937	GLY	2.6
1	A	955	TYR	2.6
1	A	922	GLY	2.5
1	A	896	TYR	2.5
1	A	942	SER	2.5
1	A	935	ASP	2.3
1	A	897	LEU	2.3
1	A	924	GLY	2.3
1	A	973	SER	2.2
1	A	946	GLN	2.2
1	A	1138	ASP	2.2
1	A	936	CYS	2.2
1	A	976[A]	LEU	2.1
1	A	972	LYS	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, 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
1	PTR	A	1055	16/17	0.74	0.23	52,72,88,89	0
1	PTR	A	1054	16/17	0.80	0.23	53,60,83,83	0

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

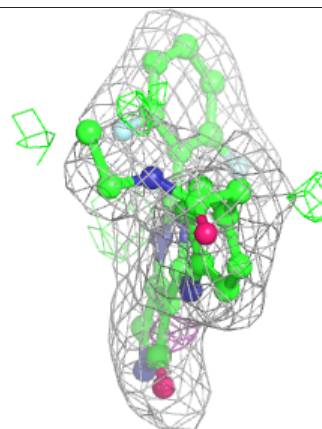
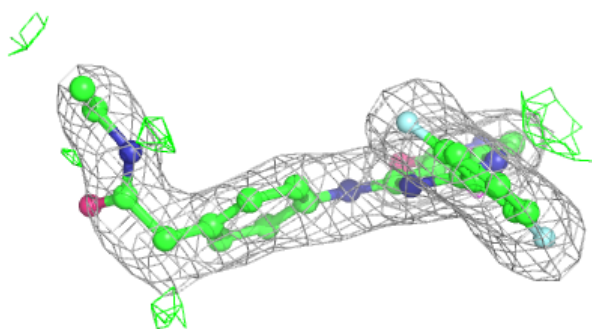
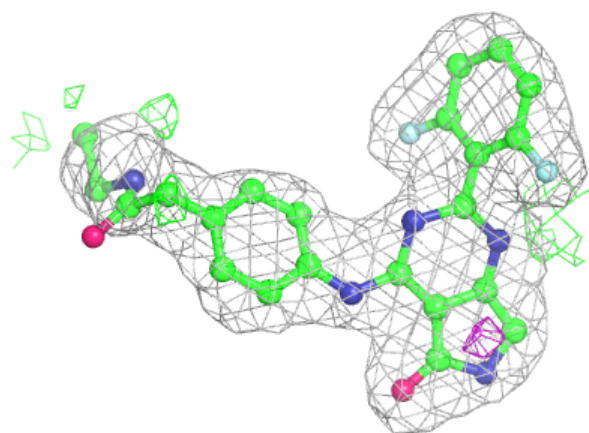
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

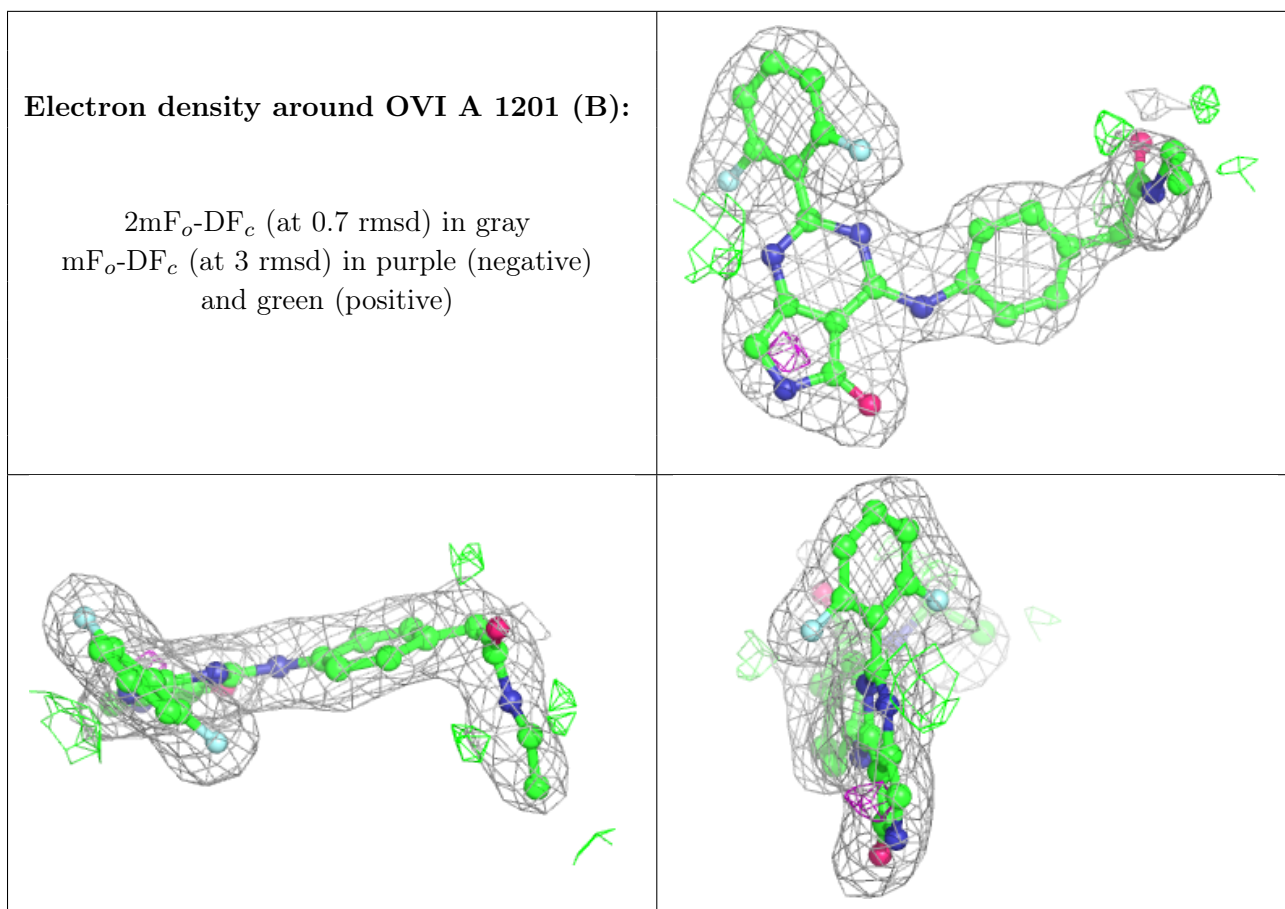
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	OVI	A	1201[A]	31/31	0.95	0.13	29,31,43,44	13
2	OVI	A	1201[B]	31/31	0.95	0.13	29,31,43,44	13

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 OVI A 1201 (A):

$2mF_o-DF_c$ (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.