



# Full wwPDB X-ray Structure Validation Report ⓘ

May 23, 2020 – 03:41 pm BST

PDB ID : 5H0B  
Title : Crystal structure of HCK complexed with a pyrrolo-pyrimidine inhibitor (S)-2-(((1r,4S)-4-(4-amino-5-(4-phenoxyphenyl)-7H-pyrrolo[2,3-d]pyrimidin-7-yl)cyclohexyl)amino)-4-methylpentanoic acid  
Authors : Tomabechi, Y.; Kukimoto-Niino, M.; Shirouzu, M.  
Deposited on : 2016-10-04  
Resolution : 1.65 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 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 : 2.11  
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) : 2.11

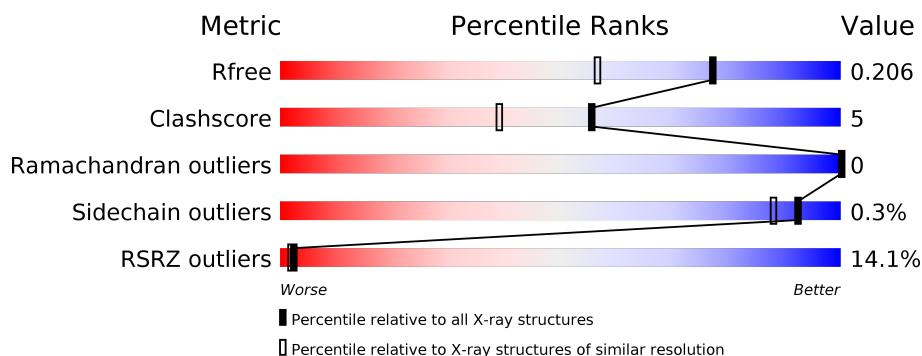
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.65 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	1827 (1.66-1.66)
Clashscore	141614	1931 (1.66-1.66)
Ramachandran outliers	138981	1891 (1.66-1.66)
Sidechain outliers	138945	1891 (1.66-1.66)
RSRZ outliers	127900	1791 (1.66-1.66)

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	454	<div> <div>14%</div> <div>88%</div> <div>11%</div> <div>.</div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4033 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.

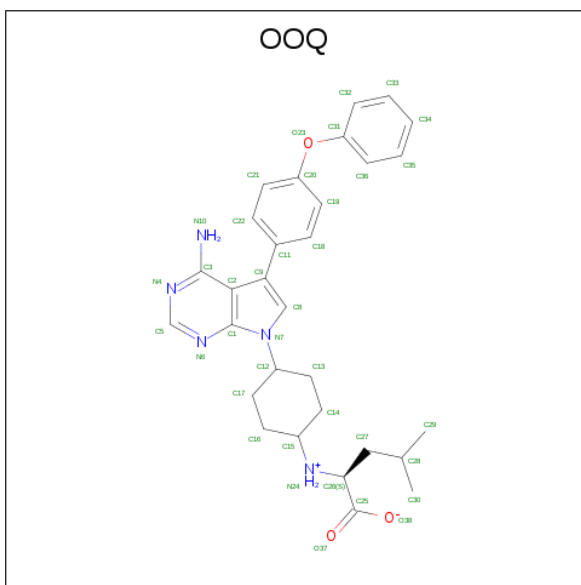
- Molecule 1 is a protein called Tyrosine-protein kinase HCK.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	P	S			
1	A	447	3613	2306	609	677	1	20	0	0	0

There are 11 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	78	GLY	-	expression tag	UNP P08631
A	79	ALA	-	expression tag	UNP P08631
A	80	MET	-	expression tag	UNP P08631
A	81	GLY	-	expression tag	UNP P08631
A	82	SER	-	expression tag	UNP P08631
A	83	GLY	-	expression tag	UNP P08631
A	84	ILE	-	expression tag	UNP P08631
A	85	ARG	-	linker	UNP P08631
A	528	GLU	GLN	engineered mutation	UNP P08631
A	529	GLU	GLN	engineered mutation	UNP P08631
A	530	ILE	GLN	engineered mutation	UNP P08631

- Molecule 2 is (2 {S})-2-[[4-[4-azanyl-5-(4-phenoxyphenyl)pyrrolo[2,3-d]pyrimidin-7-yl]cyclohexyl]azaniumyl]-4-methyl-pentanoate (three-letter code: OOQ) (formula: C<sub>30</sub>H<sub>35</sub>N<sub>5</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			38	30	5	3		

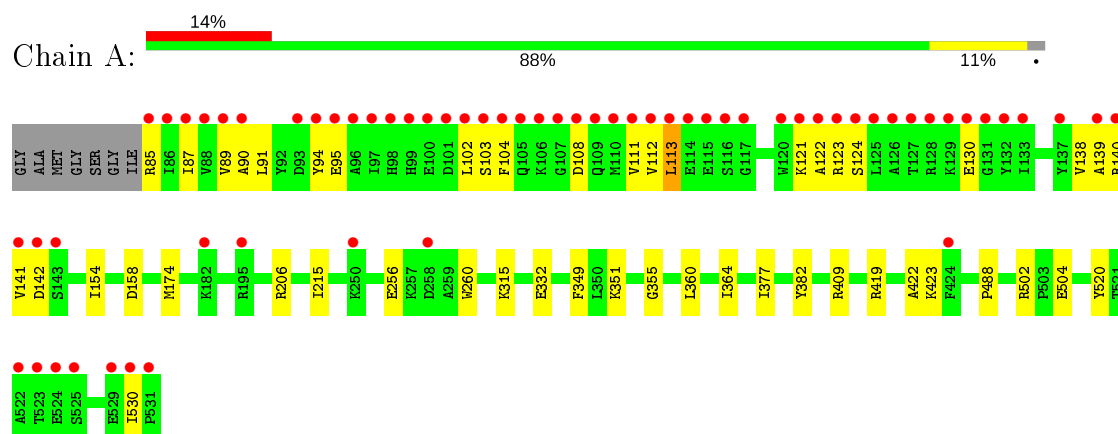
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	382	Total O 382 382	0	0

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

These plots are drawn for all protein, RNA and DNA 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: Tyrosine-protein kinase HCK



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	43.18 Å 85.58 Å 130.28 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.73 – 1.65 38.73 – 1.65	Depositor EDS
% Data completeness (in resolution range)	97.6 (38.73-1.65) 97.6 (38.73-1.65)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.26 (at 1.65 Å)	Xtriage
Refinement program	PHENIX 1.9_1690	Depositor
R, $R_{free}$	0.180 , 0.205 0.184 , 0.206	Depositor DCC
$R_{free}$ test set	2000 reflections (3.48%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	24.2	Xtriage
Anisotropy	0.261	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 52.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	4033	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.73% 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 ⓘ

### 5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: OOQ, 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	0.35	0/3682	0.50	0/4972

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 ⓘ

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	3613	0	3574	37	0
2	A	38	0	0	0	0
3	A	382	0	0	14	1
All	All	4033	0	3574	37	1

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 (37) 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:419:ARG:NH2	3:A:702:HOH:O	2.05	0.88

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:409:ARG:NH2	3:A:701:HOH:O	1.98	0.83
1:A:332:GLU:OE1	3:A:703:HOH:O	2.05	0.73
1:A:206:ARG:NH2	3:A:707:HOH:O	2.23	0.71
1:A:85:ARG:N	3:A:708:HOH:O	2.23	0.71
1:A:351:LYS:NZ	3:A:711:HOH:O	2.25	0.70
1:A:140:ARG:NH2	3:A:709:HOH:O	2.24	0.70
1:A:158:ASP:OD1	3:A:704:HOH:O	2.13	0.65
1:A:94:TYR:H	1:A:104:PHE:HB2	1.62	0.65
1:A:256:GLU:OE1	3:A:705:HOH:O	2.15	0.64
1:A:140:ARG:CZ	3:A:709:HOH:O	2.49	0.59
1:A:113:LEU:N	1:A:121:LYS:O	2.35	0.58
1:A:123:ARG:HG3	1:A:130:GLU:HG2	1.87	0.57
1:A:140:ARG:NH1	3:A:709:HOH:O	2.37	0.56
1:A:89:VAL:HG23	1:A:108:ASP:O	2.06	0.56
1:A:94:TYR:O	1:A:104:PHE:N	2.35	0.55
1:A:94:TYR:HB3	1:A:104:PHE:HD2	1.72	0.55
1:A:112:VAL:HA	1:A:122:ALA:HA	1.88	0.54
1:A:154:ILE:HD11	1:A:174:MET:SD	2.47	0.54
1:A:91:LEU:HD11	1:A:139:ALA:HB2	1.90	0.54
1:A:502:ARG:NH2	1:A:504:GLU:OE1	2.42	0.53
1:A:364:ILE:HD12	1:A:520:TYR:OH	2.10	0.51
1:A:90:ALA:HA	1:A:138:VAL:HG12	1.91	0.51
1:A:87:ILE:HG23	1:A:141:VAL:HG11	1.91	0.51
1:A:215:ILE:HG22	1:A:530:ILE:HG21	1.92	0.50
1:A:141:VAL:HG13	1:A:142:ASP:OD1	2.12	0.50
1:A:102:LEU:HG	1:A:103:SER:H	1.77	0.49
1:A:422:ALA:HA	3:A:716:HOH:O	2.13	0.49
1:A:85:ARG:HH11	1:A:111:VAL:HG21	1.80	0.46
1:A:349:PHE:O	1:A:355:GLY:HA3	2.16	0.46
1:A:377:ILE:HG23	1:A:382:TYR:HB3	1.99	0.44
1:A:360:LEU:HD21	1:A:488:PRO:HD3	2.02	0.42
1:A:423:LYS:HD3	3:A:778:HOH:O	2.19	0.41
1:A:95:GLU:OE2	3:A:706:HOH:O	2.22	0.41
1:A:142:ASP:N	1:A:142:ASP:OD1	2.54	0.41
1:A:124:SER:H	1:A:130:GLU:HG3	1.86	0.41
1:A:260:TRP:CZ2	1:A:315:LYS:HD3	2.57	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.



Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:858:HOH:O	3:A:919:HOH:O[3_644]	2.19	0.01

## 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	444/454 (98%)	428 (96%)	16 (4%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	390/393 (99%)	389 (100%)	1 (0%)	92	88

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	113	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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	PTR	A	527	1	15,16,17	1.28	1 (6%)	19,22,24	0.67	0

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	527	1	-	0/10/11/13	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	527	PTR	OH-CZ	-4.37	1.30	1.40

There are no bond angle outliers.

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 carbohydrates 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	OOQ	A	601	-	37,42,42	0.96	1 (2%)	41,59,59	1.16	5 (12%)

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	OOQ	A	601	-	-	0/16/46/46	0/5/5/5

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601	OOQ	C9-C2	4.10	1.48	1.41

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	OOQ	C9-C2-C1	-3.35	104.86	107.54
2	A	601	OOQ	C8-N7-C12	2.95	128.10	125.48
2	A	601	OOQ	N6-C5-N4	-2.70	124.46	128.68
2	A	601	OOQ	C2-C3-N4	-2.21	116.94	122.53
2	A	601	OOQ	C31-O23-C20	2.15	123.84	118.80

There are no chirality outliers.

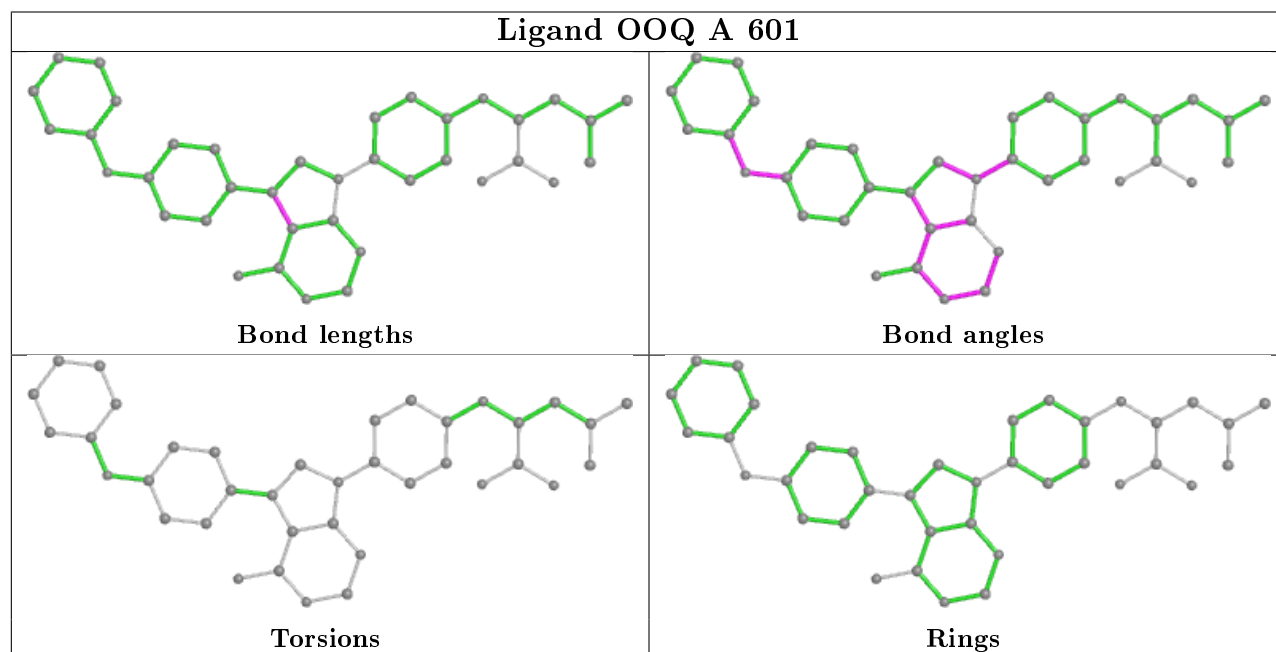
There are no torsion outliers.

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, 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	446/454 (98%)	0.89	63 (14%) 2 2	16, 29, 97, 121	0

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	127	THR	17.9
1	A	131	GLY	13.3
1	A	112	VAL	12.6
1	A	86	ILE	12.4
1	A	126	ALA	11.9
1	A	99	HIS	11.0
1	A	116	SER	10.1
1	A	97	ILE	10.0
1	A	93	ASP	9.4
1	A	105	GLN	9.3
1	A	125	LEU	9.1
1	A	129	LYS	9.0
1	A	88	VAL	8.3
1	A	102	LEU	8.2
1	A	132	TYR	7.5
1	A	142	ASP	7.2
1	A	522	ALA	7.2
1	A	106	LYS	7.1
1	A	523	THR	6.8
1	A	111	VAL	6.7
1	A	141	VAL	6.6
1	A	87	ILE	6.6
1	A	122	ALA	6.6
1	A	104	PHE	6.5
1	A	114	GLU	6.2
1	A	120	TRP	6.1
1	A	113	LEU	6.1

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	A	123	ARG	6.1
1	A	128	ARG	6.0
1	A	110	MET	5.8
1	A	108	ASP	5.7
1	A	117	GLY	5.7
1	A	85	ARG	5.5
1	A	133	ILE	5.4
1	A	531	PRO	5.4
1	A	103	SER	5.1
1	A	130	GLU	5.0
1	A	143	SER	4.7
1	A	109	GLN	4.6
1	A	94	TYR	4.5
1	A	121	LYS	4.3
1	A	98	HIS	4.3
1	A	90	ALA	4.0
1	A	530	ILE	3.8
1	A	529	GLU	3.6
1	A	101	ASP	3.5
1	A	89	VAL	3.3
1	A	195	ARG	3.2
1	A	140	ARG	3.2
1	A	96	ALA	3.1
1	A	100	GLU	3.1
1	A	525	SER	3.1
1	A	524	GLU	3.1
1	A	124	SER	3.0
1	A	95	GLU	3.0
1	A	115	GLU	2.8
1	A	107	GLY	2.7
1	A	182	LYS	2.6
1	A	258	ASP	2.5
1	A	139	ALA	2.5
1	A	250	LYS	2.5
1	A	424	PHE	2.3
1	A	137	TYR	2.1

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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( $\text{\AA}^2$ )	Q<0.9
1	PTR	A	527	16/17	0.97	0.11	28,36,40,45	0

### 6.3 Carbohydrates

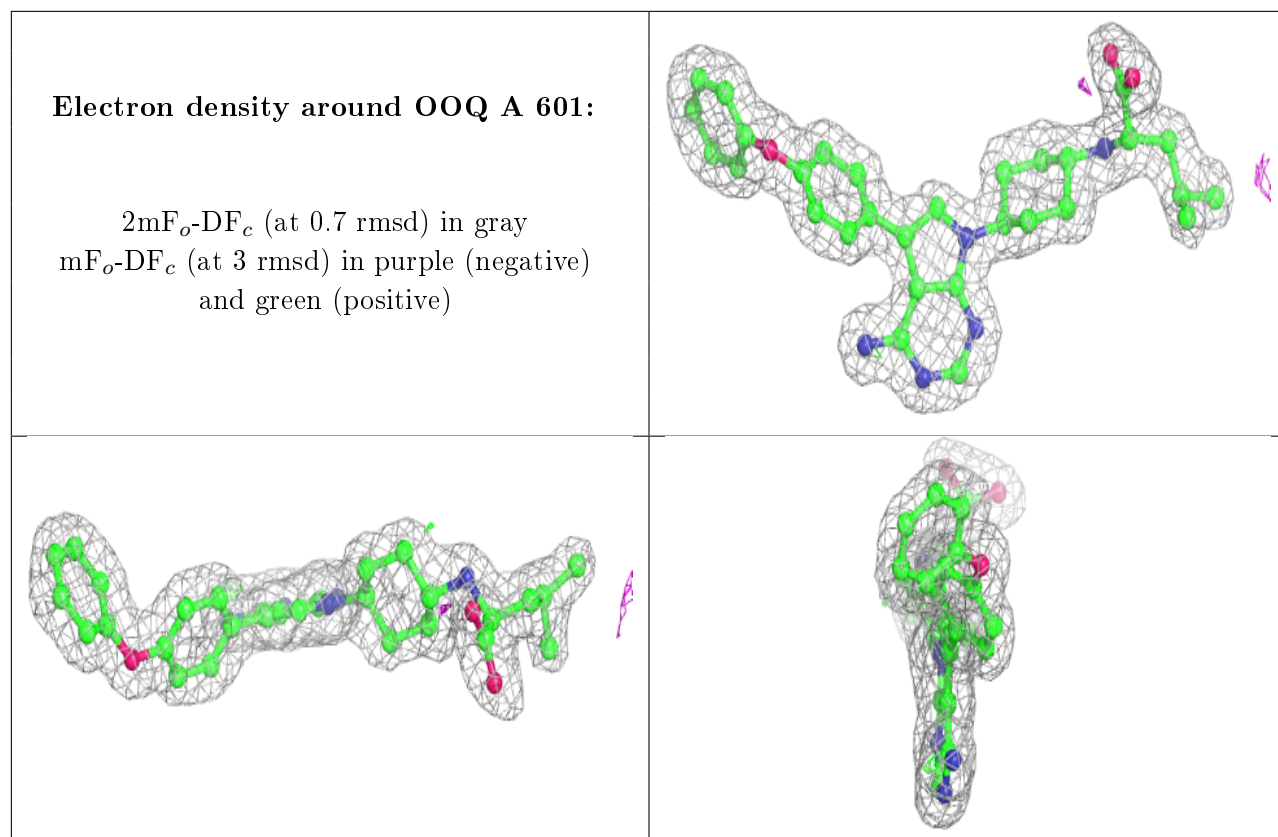
There are no carbohydrates in this entry.

### 6.4 Ligands

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( $\text{\AA}^2$ )	Q<0.9
2	OOQ	A	601	38/38	0.95	0.08	17,21,31,34	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.