



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:04 PM UTC

PDB ID : 6Q74 / pdb_00006q74
Title : PI3K delta in complex with 1benzylN[5(3,6dihydro2Hpyran4yl)2methoxyipyridin3yl]2methyl1Himidazole4sulfonamide
Authors : Convery, M.A.; Rowland, P.; Down, K.; Barton, N.
Deposited on : 2018-12-12
Resolution : 2.48 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)
Xtriage (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

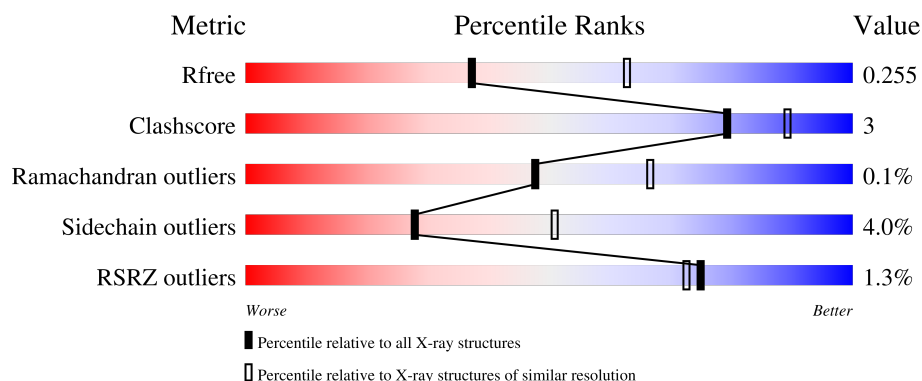
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

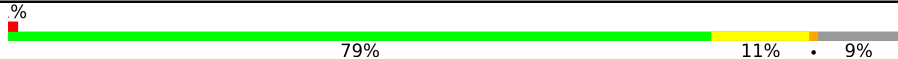
The reported resolution of this entry is 2.48 Å.

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	7589 (2.50-2.46)
Clashscore	190562	8295 (2.50-2.46)
Ramachandran outliers	187476	8164 (2.50-2.46)
Sidechain outliers	187428	8166 (2.50-2.46)
RSRZ outliers	180081	7593 (2.50-2.46)

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	940	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 7120 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.

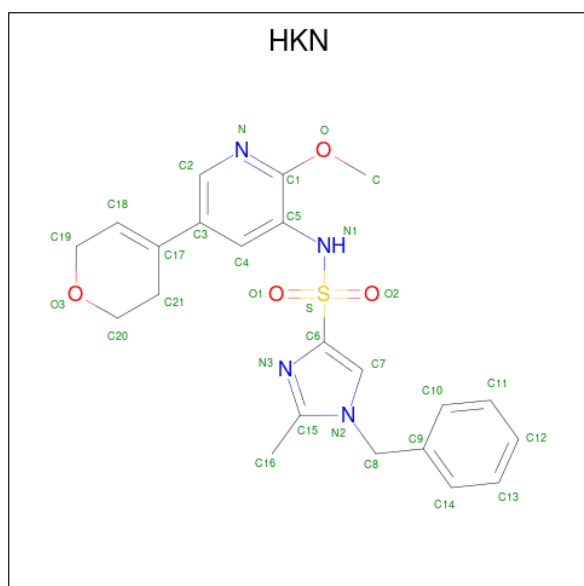
- Molecule 1 is a protein called Phosphatidylinositol 4,5-bisphosphate 3-kinase catalytic sub-unit delta isoform.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	854	Total	C	N	O	S	0	0	0
			6869	4392	1172	1251	54			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	105	GLY	-	expression tag	UNP O35904
A	508	GLN	-	insertion	UNP O35904

- Molecule 2 is {N}-[5-(3,6-dihydro-2 {H}-pyran-4-yl)-2-methoxy-pyridin-3-yl]-2-methyl-1-(phenylmethyl)imidazole-4-sulfonamide (CCD ID: HKN) (formula: C₂₂H₂₄N₄O₄S) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			31	22	4	4	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	220	Total 220	O 220	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	141.28Å 63.96Å 116.22Å 90.00° 102.16° 90.00°	Depositor
Resolution (Å)	56.81 – 2.48 56.81 – 2.48	Depositor EDS
% Data completeness (in resolution range)	85.3 (56.81-2.48) 85.2 (56.81-2.48)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.99 (at 2.48Å)	Xtriage
Refinement program	BUSTER 2.11.7	Depositor
R, R_{free}	0.174 , 0.249 0.181 , 0.255	Depositor DCC
R_{free} test set	1474 reflections (4.05%)	wwPDB-VP
Wilson B-factor (Å ²)	65.9	Xtriage
Anisotropy	0.039	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 79.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7120	wwPDB-VP
Average B, all atoms (Å ²)	82.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.33% 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: HKN

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	0.88	1/7019 (0.0%)	1.40	25/9475 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	757	LYS	CA-CB	5.67	1.56	1.52

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	915	PHE	CA-CB-CG	9.40	123.20	113.80
1	A	284	ALA	CA-C-N	6.56	129.07	120.28
1	A	284	ALA	C-N-CA	6.56	129.07	120.28
1	A	427	ASP	CA-CB-CG	6.50	119.10	112.60
1	A	205	PHE	CA-CB-CG	5.93	119.73	113.80
1	A	932	PHE	CA-CB-CG	5.88	119.68	113.80
1	A	488	LEU	CA-C-N	5.61	128.06	120.38
1	A	488	LEU	C-N-CA	5.61	128.06	120.38
1	A	951	GLU	N-CA-C	-5.59	104.88	110.97
1	A	656	HIS	CA-C-N	5.43	126.29	122.60
1	A	656	HIS	C-N-CA	5.43	126.29	122.60
1	A	634	LEU	CA-C-N	5.29	125.85	119.98
1	A	634	LEU	C-N-CA	5.29	125.85	119.98
1	A	135	GLU	CB-CG-CD	5.27	121.56	112.60
1	A	456	ALA	CA-C-N	5.27	125.11	120.10
1	A	456	ALA	C-N-CA	5.27	125.11	120.10
1	A	729	SER	CA-C-N	5.16	130.14	122.82
1	A	729	SER	C-N-CA	5.16	130.14	122.82
1	A	865	GLY	CA-C-N	5.10	127.83	120.38
1	A	865	GLY	C-N-CA	5.10	127.83	120.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	243	VAL	N-CA-C	-5.09	102.17	108.89
1	A	1022	ALA	CA-C-N	5.04	127.30	120.44
1	A	1022	ALA	C-N-CA	5.04	127.30	120.44
1	A	109	VAL	CA-C-N	5.00	126.98	120.28
1	A	109	VAL	C-N-CA	5.00	126.98	120.28

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	6869	0	6835	37	0
2	A	31	0	0	0	0
3	A	220	0	0	0	0
All	All	7120	0	6835	37	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 3.

All (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:216:LEU:HD12	1:A:256:LEU:HD11	1.78	0.64
1:A:758:PRO:HB3	1:A:779:LYS:HG3	1.80	0.63
1:A:932:PHE:HB3	1:A:1022:ALA:HB2	1.78	0.63
1:A:603:LEU:HD22	1:A:607:GLU:HB3	1.83	0.58
1:A:333:VAL:HG11	1:A:341:LEU:HD11	1.86	0.57
1:A:515:LEU:HD11	1:A:538:VAL:HG22	1.88	0.55
1:A:583:LEU:HD11	1:A:600:LEU:HD11	1.89	0.55
1:A:713:GLU:O	1:A:717:MET:HG3	2.07	0.54
1:A:971:GLY:HA3	1:A:1004:LEU:HD21	1.89	0.53
1:A:621:TYR:CZ	1:A:983:ALA:HB2	2.44	0.53
1:A:328:ILE:HD11	1:A:474:TYR:HB2	1.90	0.52
1:A:390:LEU:HB2	1:A:425:LEU:HD21	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:534:MET:O	1:A:538:VAL:HG23	2.11	0.51
1:A:609:PHE:HE1	1:A:646:PHE:CD2	2.29	0.51
1:A:637:ALA:HB1	1:A:644:GLY:HA2	1.94	0.50
1:A:512:ARG:HG3	1:A:542:PHE:HZ	1.77	0.50
1:A:340:LYS:HD3	1:A:362:GLU:HB3	1.94	0.48
1:A:341:LEU:HG	1:A:365:VAL:HG22	1.95	0.48
1:A:698:PHE:HZ	1:A:714:MET:HB3	1.79	0.48
1:A:116:GLN:HG2	1:A:683:MET:HE1	1.95	0.48
1:A:788:MET:O	1:A:792:GLN:HG3	2.15	0.47
1:A:859:LEU:HD21	1:A:901:ILE:HD11	1.97	0.47
1:A:133:ASP:HB3	1:A:136:VAL:HB	1.97	0.47
1:A:432:LEU:HD23	1:A:478:VAL:HG21	1.98	0.45
1:A:695:LEU:HD23	1:A:759:LEU:HD13	1.99	0.45
1:A:943:GLN:HG2	1:A:949:ASN:HB3	2.00	0.44
1:A:784:LEU:HD12	1:A:823:GLY:HA3	2.00	0.43
1:A:789:LEU:HD13	1:A:981:MET:HE2	2.00	0.43
1:A:417:PRO:HB3	1:A:452:LEU:HD23	2.01	0.43
1:A:268:SER:OG	1:A:270:LEU:HD12	2.20	0.42
1:A:971:GLY:HA2	1:A:1002:LEU:HD13	2.01	0.42
1:A:703:SER:HB3	1:A:711:THR:HG21	2.02	0.42
1:A:936:TYR:HA	1:A:939:VAL:HB	2.02	0.41
1:A:630:THR:HG22	1:A:634:LEU:HD12	2.03	0.41
1:A:154:ARG:HD2	1:A:165:TYR:CE1	2.56	0.40
1:A:341:LEU:HD13	1:A:369:PRO:HG3	2.03	0.40
1:A:961:GLU:O	1:A:965:THR:HG23	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	838/940 (89%)	813 (97%)	24 (3%)	1 (0%)	48	66

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	328	ILE

5.3.2 Protein sidechains ⓘ

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	752/827 (91%)	722 (96%)	30 (4%)	28	51

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	154	ARG
1	A	166	SER
1	A	174	SER
1	A	192	VAL
1	A	195	LYS
1	A	210	LYS
1	A	283	LEU
1	A	342	VAL
1	A	370	VAL
1	A	394	LEU
1	A	436	GLU
1	A	445	VAL
1	A	471	LEU
1	A	483	VAL
1	A	570	SER
1	A	684	LYS
1	A	687	GLU
1	A	690	SER
1	A	696	ASN
1	A	699	VAL
1	A	779	LYS
1	A	787	ASP
1	A	855	LEU
1	A	897	ASP

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Mol	Chain	Res	Type
1	A	915	PHE
1	A	935	THR
1	A	947	THR
1	A	951	GLU
1	A	998	LEU
1	A	1017	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	187	ASN
1	A	247	HIS
1	A	351	ASN
1	A	508	GLN
1	A	539	GLN
1	A	610	GLN
1	A	696	ASN
1	A	732	GLN
1	A	836	ASN
1	A	918	ASN
1	A	943	GLN
1	A	944	GLN
1	A	948	ASN
1	A	949	ASN
1	A	970	HIS
1	A	976	HIS
1	A	996	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

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	HKN	A	4001	-	33,34,34	0.57	0	40,48,48	0.60	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
2	HKN	A	4001	-	-	2/19/29/29	0/4/4/4

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

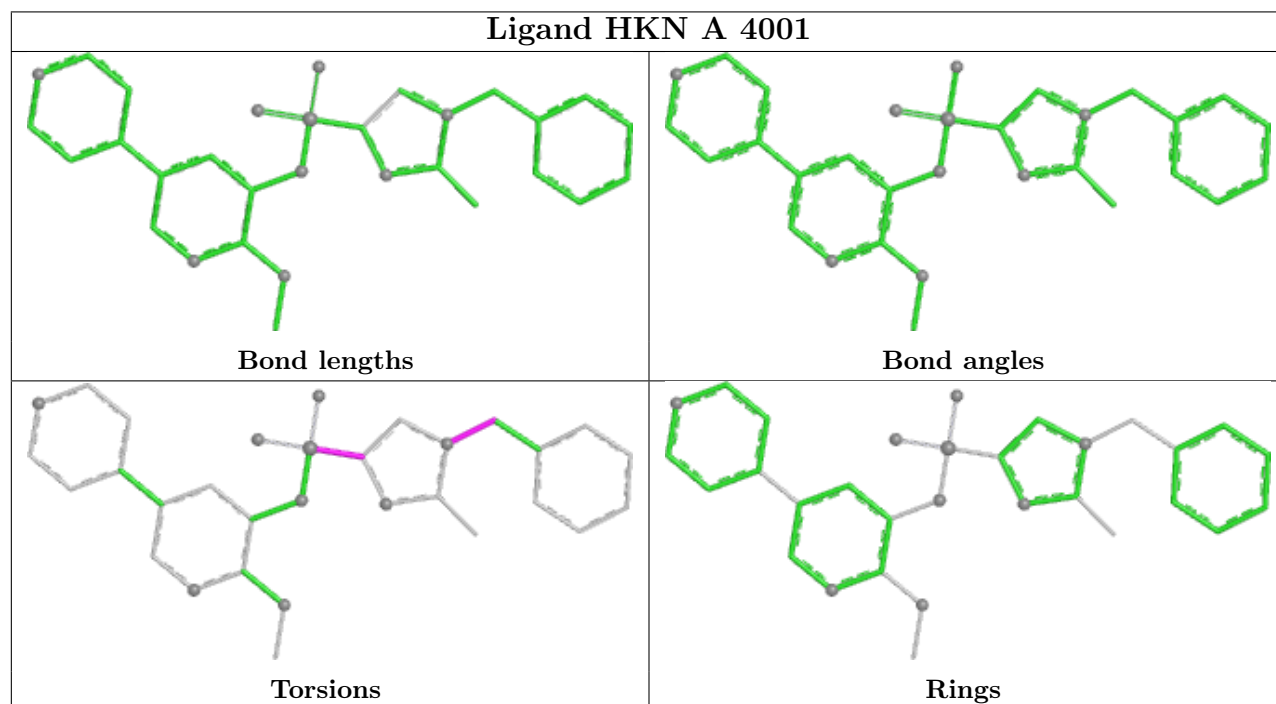
Mol	Chain	Res	Type	Atoms
2	A	4001	HKN	N3-C6-S-O1
2	A	4001	HKN	C9-C8-N2-C15

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 [i](#)

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

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	854/940 (90%)	-0.00	11 (1%) 75 72	43, 78, 128, 175	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	934	LEU	3.6
1	A	933	ILE	3.1
1	A	521	GLY	3.0
1	A	931	PRO	2.8
1	A	932	PHE	2.8
1	A	445	VAL	2.6
1	A	398	VAL	2.6
1	A	177	GLY	2.5
1	A	452	LEU	2.2
1	A	107	ASP	2.2
1	A	938	PHE	2.1

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.

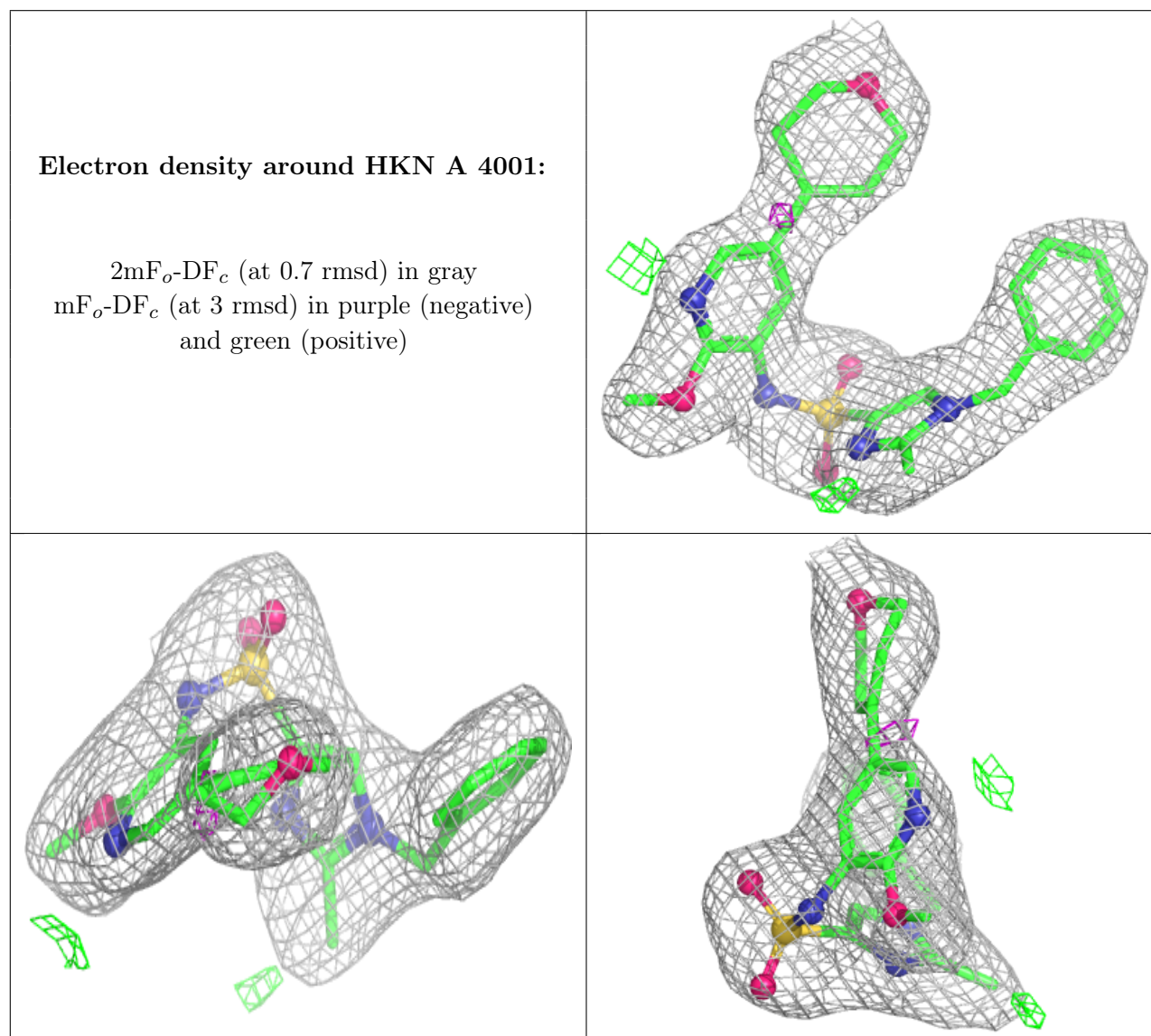
6.4 Ligands [i](#)

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(Å ²)	Q<0.9
2	HKN	A	4001	31/31	0.96	0.08	47,65,72,74	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 ⓘ

There are no such residues in this entry.