



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:56 AM UTC

PDB ID : 7AYM / pdb_00007aym
Title : Structure of DDR2 Kinase domain in complex with IBZ3
Authors : Nawrotek, A.; Talagas, A.; Vuillard, L.M.; Miallau, L.
Deposited on : 2020-11-12
Resolution : 2.12 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

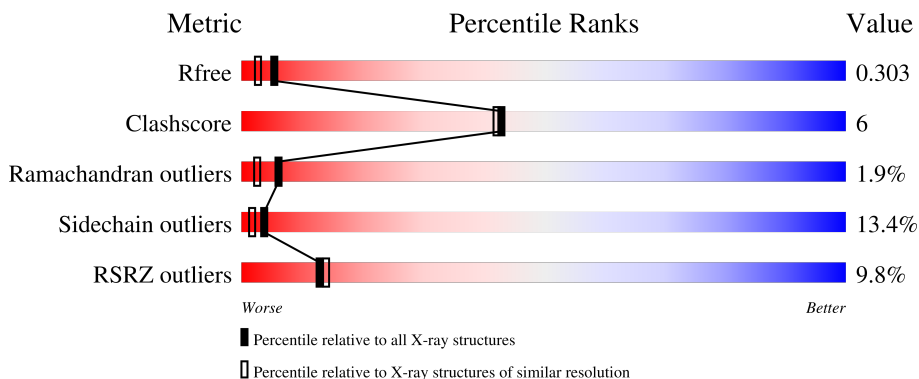
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.12 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	8290 (2.14-2.10)
Clashscore	190562	8817 (2.14-2.10)
Ramachandran outliers	187476	8738 (2.14-2.10)
Sidechain outliers	187428	8739 (2.14-2.10)
RSRZ outliers	180081	8294 (2.14-2.10)

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	337	8% 58% 20% • 18%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2293 atoms, of which 20 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 Discoidin domain-containing receptor 2, Epithelial discoidin domain-containing receptor 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	275	Total	C	N	O	S	0	1	0
			2234	1424	385	407	18			

There are 30 discrepancies between the modelled and reference sequences:

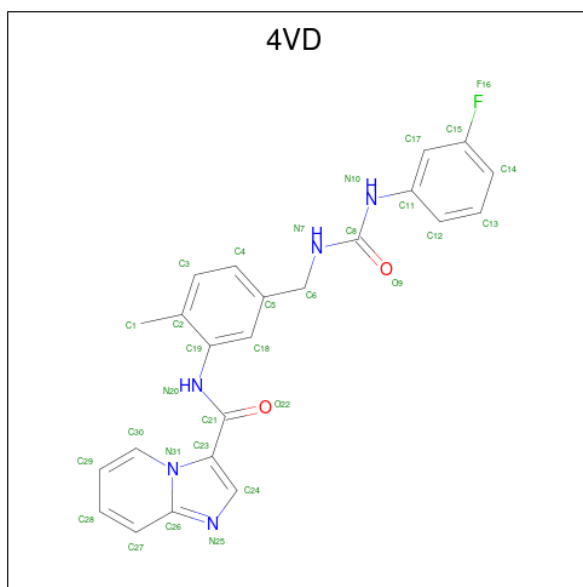
Chain	Residue	Modelled	Actual	Comment	Reference
A	530	MET	-	initiating methionine	UNP Q16832
A	531	GLY	-	expression tag	UNP Q16832
A	532	SER	-	expression tag	UNP Q16832
A	533	SER	-	expression tag	UNP Q16832
A	534	HIS	-	expression tag	UNP Q16832
A	535	HIS	-	expression tag	UNP Q16832
A	536	HIS	-	expression tag	UNP Q16832
A	537	HIS	-	expression tag	UNP Q16832
A	538	HIS	-	expression tag	UNP Q16832
A	539	HIS	-	expression tag	UNP Q16832
A	540	ASP	-	expression tag	UNP Q16832
A	541	TYR	-	expression tag	UNP Q16832
A	542	ASP	-	expression tag	UNP Q16832
A	543	ILE	-	expression tag	UNP Q16832
A	544	PRO	-	expression tag	UNP Q16832
A	545	THR	-	expression tag	UNP Q16832
A	546	THR	-	expression tag	UNP Q16832
A	547	GLU	-	expression tag	UNP Q16832
A	548	ASN	-	expression tag	UNP Q16832
A	549	LEU	-	expression tag	UNP Q16832
A	550	TYR	-	expression tag	UNP Q16832
A	551	PHE	-	expression tag	UNP Q16832
A	552	GLN	-	expression tag	UNP Q16832
A	553	GLY	-	expression tag	UNP Q16832
A	554	PRO	-	expression tag	UNP Q16832
A	555	ARG	-	expression tag	UNP Q16832

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
A	556	VAL	-	expression tag	UNP Q16832
A	557	ASP	-	expression tag	UNP Q16832
A	835	GLU	LYS	engineered mutation	UNP Q16832
A	856	ARG	LEU	engineered mutation	UNP Q16832

- Molecule 2 is N-[5-({[(3-fluorophenyl)carbamoyl]amino}methyl)-2-methylphenyl]imidazo[1,2-a]pyridine-3-carboxamide (CCD ID: 4VD) (formula: C₂₃H₂₀FN₅O₂) (labeled as "Ligand of Interest" by depositor).



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	41.75Å 85.36Å 89.54Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	61.78 – 2.12 61.78 – 2.12	Depositor EDS
% Data completeness (in resolution range)	57.0 (61.78-2.12) 48.1 (61.78-2.12)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.32 (at 2.10Å)	Xtriage
Refinement program	BUSTER 2.11.7	Depositor
R, R_{free}	0.236 , 0.289 0.249 , 0.303	Depositor DCC
R_{free} test set	515 reflections (2.64%)	wwPDB-VP
Wilson B-factor (Å ²)	21.2	Xtriage
Anisotropy	0.258	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 43.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.037 for -h,l,k	Xtriage
F_o, F_c correlation	0.85	EDS
Total number of atoms	2293	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.24% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 4VD

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	0.91	0/2285	1.45	19/3083 (0.6%)

There are no bond length outliers.

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	661	ASP	CA-CB-CG	8.04	120.64	112.60
1	A	844	ASP	CA-C-N	6.09	128.72	120.38
1	A	844	ASP	C-N-CA	6.09	128.72	120.38
1	A	715	PHE	CA-CB-CG	5.89	119.69	113.80
1	A	778	ASP	CA-C-N	5.68	127.72	120.56
1	A	778	ASP	C-N-CA	5.68	127.72	120.56
1	A	707	MET	CA-C-N	5.65	127.85	120.28
1	A	707	MET	C-N-CA	5.65	127.85	120.28
1	A	831	ASP	CA-C-N	5.52	127.68	120.28
1	A	831	ASP	C-N-CA	5.52	127.68	120.28
1	A	730	ASN	CA-C-N	5.40	130.22	122.40
1	A	730	ASN	C-N-CA	5.40	130.22	122.40
1	A	787	TRP	CA-C-N	5.38	127.80	120.54
1	A	787	TRP	C-N-CA	5.38	127.80	120.54
1	A	816	ASP	CA-CB-CG	5.20	117.80	112.60
1	A	777	SER	CA-C-N	5.09	127.53	120.29
1	A	777	SER	C-N-CA	5.09	127.53	120.29
1	A	798	TYR	CA-C-N	5.00	126.94	120.44
1	A	798	TYR	C-N-CA	5.00	126.94	120.44

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	2234	0	2216	28	0
2	A	31	20	20	1	0
3	A	8	0	0	1	0
All	All	2273	20	2236	28	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 6.

All (28) 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:701:THR:HG22	1:A:858:LEU:HB3	1.49	0.92
1:A:798:TYR:HB3	1:A:801:LEU:HD12	1.70	0.73
1:A:834:TYR:HD1	1:A:837:MET:HE2	1.58	0.69
1:A:697:LYS:O	1:A:701:THR:HG23	1.94	0.67
1:A:575:GLY:HA2	1:A:611:ARG:HG2	1.88	0.55
1:A:844:ASP:HB2	1:A:847:ASN:HD22	1.71	0.55
1:A:834:TYR:CD1	1:A:837:MET:HE2	2.39	0.55
1:A:740[A]:MET:HG3	1:A:757:VAL:HG21	1.93	0.51
1:A:669:HIS:CD2	1:A:689:PRO:HB2	2.48	0.49
1:A:739:GLY:HA2	3:A:1005:HOH:O	2.12	0.49
1:A:740[A]:MET:HE2	1:A:749:TYR:HB2	1.95	0.49
1:A:800:GLN:HE21	1:A:801:LEU:HG	1.78	0.48
1:A:802:SER:H	1:A:805:GLN:HB2	1.80	0.46
1:A:758:LEU:HD22	1:A:763:MET:HE1	1.96	0.46
1:A:632:LEU:HB3	1:A:637:ILE:HG21	1.97	0.45
1:A:637:ILE:HD11	1:A:706:GLY:O	2.18	0.44
1:A:661:ASP:OD1	1:A:663:ASN:HB3	2.17	0.44
1:A:644:CYS:HB2	1:A:651:CYS:HB2	2.00	0.44
1:A:656:TYR:CE2	1:A:658:GLU:HA	2.53	0.43
1:A:831:ASP:O	1:A:835:GLU:HG2	2.19	0.43
1:A:751:ARG:HH21	1:A:753:GLN:HA	1.84	0.43
1:A:836:LEU:HD13	1:A:857:LEU:HD12	1.99	0.43
1:A:740[B]:MET:HG3	1:A:749:TYR:CG	2.53	0.43
1:A:736:ALA:HA	2:A:901:4VD:H44	2.00	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:559:PRO:HB2	1:A:562:LEU:HD13	2.02	0.42
1:A:842:ARG:HH21	1:A:847:ASN:HD22	1.68	0.41
1:A:559:PRO:O	1:A:562:LEU:HB2	2.21	0.41
1:A:812:GLU:OE1	1:A:819:ARG:HB2	2.20	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	268/337 (80%)	250 (93%)	13 (5%)	5 (2%)	6 2

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	599	ALA
1	A	614	ALA
1	A	741	SER
1	A	769	LEU
1	A	575	GLY

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	247/295 (84%)	213 (86%)	34 (14%)	3 2

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	568	LYS
1	A	581	GLU
1	A	609	MET
1	A	613	ASP
1	A	615	ASN
1	A	624	LYS
1	A	631	ARG
1	A	632	LEU
1	A	633	LYS
1	A	643	VAL
1	A	645	ILE
1	A	646	THR
1	A	670	GLN
1	A	690	THR
1	A	691	VAL
1	A	705	SER
1	A	708	LYS
1	A	713	LEU
1	A	727	VAL
1	A	740[A]	MET
1	A	740[B]	MET
1	A	751	ARG
1	A	757	VAL
1	A	767	SER
1	A	768	ILE
1	A	769	LEU
1	A	779	VAL
1	A	794	GLN
1	A	799	SER
1	A	820	GLN
1	A	821	THR
1	A	837	MET
1	A	842	ARG
1	A	852	GLN

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	847	ASN

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](#)

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	4VD	A	901	-	34,34,34	0.20	0	46,47,47	0.52	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	4VD	A	901	-	-	0/17/17/17	0/4/4/4

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

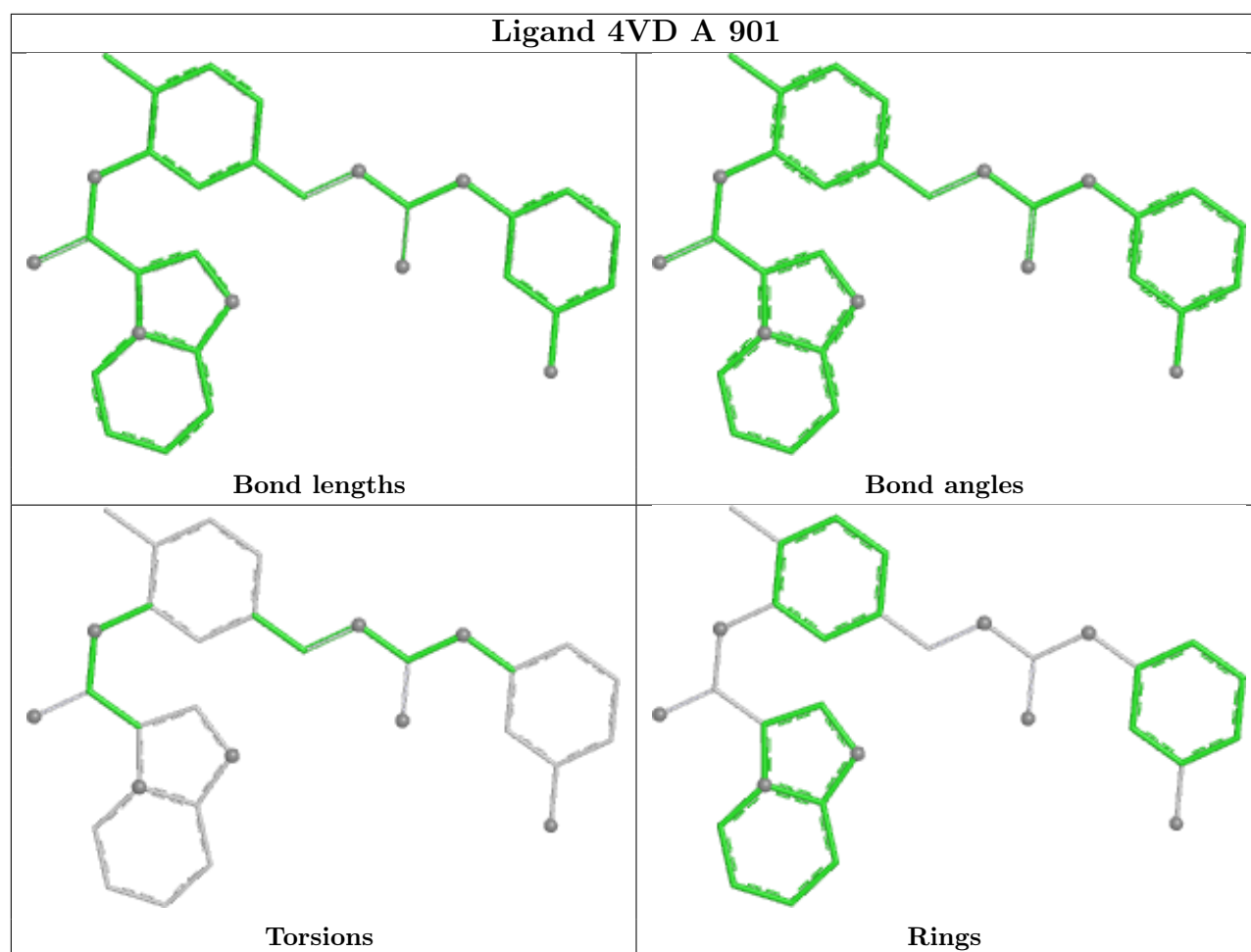
There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	901	4VD	1	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, 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	275/337 (81%)	0.71	27 (9%) 13 14	7, 23, 47, 67	4 (1%)

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	574	PHE	4.6
1	A	572	GLY	3.6
1	A	807	ILE	3.6
1	A	740[A]	MET	3.4
1	A	599	ALA	3.0
1	A	768	ILE	2.8
1	A	638	ILE	2.8
1	A	644	CYS	2.8
1	A	614	ALA	2.7
1	A	741	SER	2.7
1	A	670	GLN	2.7
1	A	817	GLN	2.7
1	A	646	THR	2.6
1	A	752	ILE	2.6
1	A	600	ASN	2.6
1	A	611	ARG	2.5
1	A	755	ARG	2.5
1	A	617	ASN	2.5
1	A	601	GLN	2.3
1	A	743	ASN	2.2
1	A	557	ASP	2.2
1	A	598	SER	2.2
1	A	612	ALA	2.2
1	A	627	LYS	2.2
1	A	575	GLY	2.1
1	A	742	ARG	2.1
1	A	818	GLY	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](#)

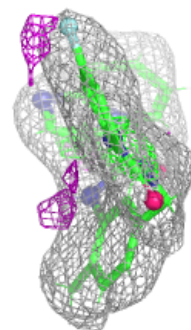
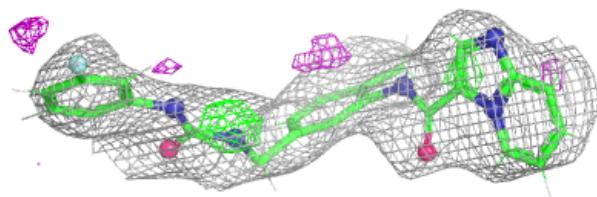
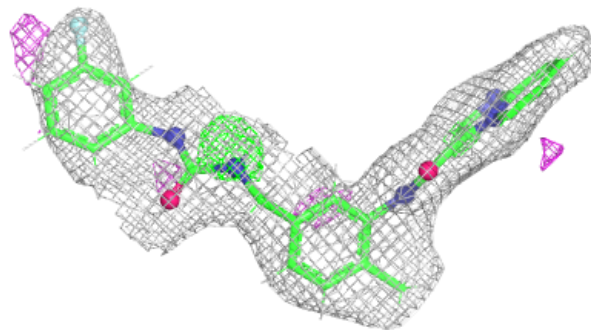
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	4VD	A	901	31/31	0.86	0.22	47,54,54,54	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 4VD A 901:

$2mF_o-DF_c$ (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.