



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 04:18 AM UTC

PDB ID : 6QHR / pdb_00006qhr
Title : Structure of the mitogen activated kinase kinase 7 in complex with pyrazolopyrimidine 1m
Authors : Wolle, P.; Mueller, M.P.; Rauh, D.
Deposited on : 2019-01-17
Resolution : 2.52 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

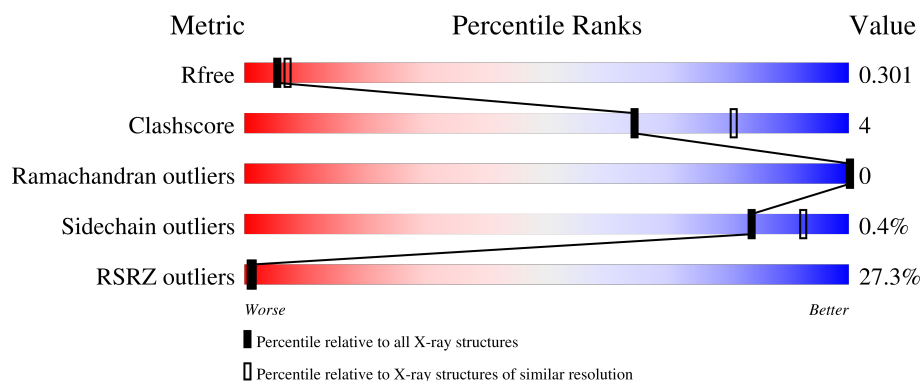
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.52 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	7383 (2.54-2.50)
Clashscore	190562	8079 (2.54-2.50)
Ramachandran outliers	187476	7944 (2.54-2.50)
Sidechain outliers	187428	7946 (2.54-2.50)
RSRZ outliers	180081	7387 (2.54-2.50)

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	318	25% 82% 8% 10%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2299 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.

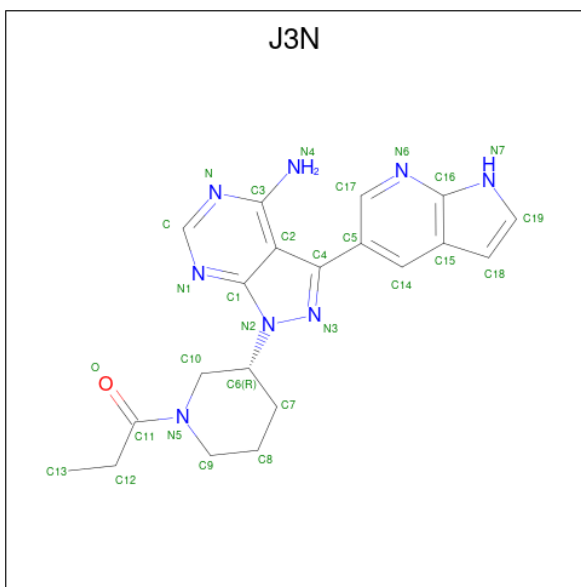
- Molecule 1 is a protein called Dual specificity mitogen-activated protein kinase kinase 7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	286	2178	1387	375	397	19	0	0	0

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	107	MET	-	initiating methionine	UNP O14733
A	108	GLY	-	expression tag	UNP O14733
A	109	HIS	-	expression tag	UNP O14733
A	110	HIS	-	expression tag	UNP O14733
A	111	HIS	-	expression tag	UNP O14733
A	112	HIS	-	expression tag	UNP O14733
A	113	HIS	-	expression tag	UNP O14733
A	114	HIS	-	expression tag	UNP O14733
A	115	SER	-	expression tag	UNP O14733
A	116	ALA	-	expression tag	UNP O14733
A	117	LYS	-	expression tag	UNP O14733
A	118	GLN	-	expression tag	UNP O14733
A	119	THR	-	expression tag	UNP O14733
A	120	GLY	-	expression tag	UNP O14733
A	121	TYR	-	expression tag	UNP O14733
A	122	LEU	-	expression tag	UNP O14733
A	123	THR	-	expression tag	UNP O14733
A	124	ILE	-	expression tag	UNP O14733
A	125	GLY	-	expression tag	UNP O14733
A	126	GLY	-	expression tag	UNP O14733

- Molecule 2 is 1-[(3 {R})-3-[4-azanyl-3-(1 {H}-pyrrolo[2,3-b]pyridin-5-yl)pyrazolo[3,4-d]pyrimidin-1-yl]piperidin-1-yl]propan-1-one (CCD ID: J3N) (formula: C₂₀H₂₂N₈O).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			29	20	8	1		

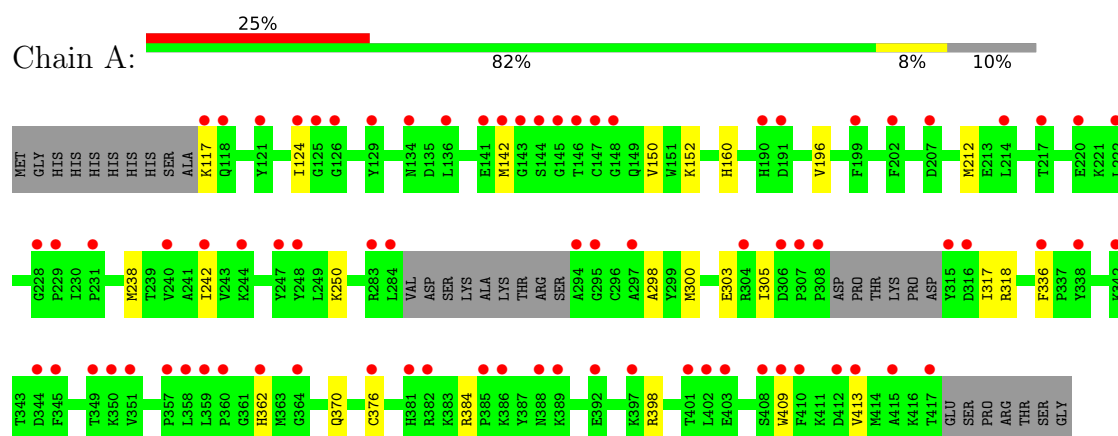
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	92	Total	O	0	0
			92	92		

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: Dual specificity mitogen-activated protein kinase kinase 7



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	62.10Å 69.00Å 84.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.16 – 2.52 46.16 – 2.52	Depositor EDS
% Data completeness (in resolution range)	98.5 (46.16-2.52) 98.9 (46.16-2.52)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.40	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.71 (at 2.39Å)	Xtriage
Refinement program	PHENIX (1.14_3260: ???)	Depositor
R, R_{free}	0.237 , 0.294 0.243 , 0.301	Depositor DCC
R_{free} test set	1026 reflections (6.94%)	wwPDB-VP
Wilson B-factor (Å ²)	38.4	Xtriage
Anisotropy	0.308	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 48.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.82	EDS
Total number of atoms	2299	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.67% 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: J3N

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	0.15	0/2221	0.34	0/3002

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	2178	0	2104	17	0
2	A	29	0	0	1	0
3	A	92	0	0	3	0
All	All	2299	0	2104	17	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 4.

All (17) 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:212:MET:HE2	2:A:501:J3N:N4	2.14	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:370:GLN:NE2	3:A:603:HOH:O	2.30	0.60
1:A:298:ALA:HB1	1:A:336:PHE:CZ	2.39	0.57
1:A:117:LYS:N	3:A:610:HOH:O	2.40	0.54
1:A:317:ILE:HG23	1:A:318:ARG:HG2	1.90	0.54
1:A:124:ILE:HG21	1:A:160:HIS:ND1	2.24	0.53
1:A:398:ARG:NH2	3:A:602:HOH:O	2.24	0.49
1:A:303:GLU:OE2	1:A:384:ARG:NH2	2.46	0.49
1:A:142:MET:HE2	1:A:152:LYS:HB2	1.95	0.48
1:A:142:MET:HE3	1:A:150:VAL:HG12	1.97	0.47
1:A:196:VAL:HB	1:A:212:MET:HE3	1.97	0.47
1:A:376:CYS:O	1:A:384:ARG:HD2	2.19	0.42
1:A:238:MET:O	1:A:242:ILE:HG12	2.19	0.42
1:A:300:MET:HE2	1:A:305:ILE:HG12	2.01	0.42
1:A:409:TRP:O	1:A:413:VAL:HG23	2.19	0.41
1:A:250:LYS:HE3	1:A:317:ILE:HD11	2.03	0.41
1:A:303:GLU:CD	1:A:384:ARG:HH22	2.28	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	280/318 (88%)	269 (96%)	11 (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	223/277 (80%)	222 (100%)	1 (0%)	84	92

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

Mol	Chain	Res	Type
1	A	362	HIS

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	149	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](#)

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	J3N	A	501	1	33,33,33	1.82	6 (18%)	48,48,48	3.53	21 (43%)

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	J3N	A	501	1	-	10/14/30/30	0/5/5/5

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	J3N	C2-C3	6.06	1.49	1.42
2	A	501	J3N	C2-C1	5.39	1.47	1.40
2	A	501	J3N	C15-C16	3.72	1.47	1.42
2	A	501	J3N	C4-N3	2.54	1.36	1.33
2	A	501	J3N	C2-C4	2.37	1.50	1.44
2	A	501	J3N	C1-N2	-2.36	1.33	1.36

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	J3N	C15-C16-N6	-11.13	120.06	126.66
2	A	501	J3N	C15-C16-N7	-7.33	104.06	108.24
2	A	501	J3N	C2-C1-N1	-7.32	119.20	126.97
2	A	501	J3N	N1-C1-N2	6.66	134.73	126.05
2	A	501	J3N	C16-C15-C18	6.52	110.16	106.48
2	A	501	J3N	C2-C4-N3	-5.87	106.16	110.19
2	A	501	J3N	N7-C16-N6	5.67	135.88	126.69
2	A	501	J3N	C16-N7-C19	4.96	110.97	108.11
2	A	501	J3N	C4-N3-N2	4.81	109.50	106.97
2	A	501	J3N	C6-C10-N5	-4.34	103.66	109.46
2	A	501	J3N	C7-C6-N2	-4.29	107.82	111.88
2	A	501	J3N	C8-C9-N5	-4.02	102.51	110.67
2	A	501	J3N	C-N1-C1	3.74	120.96	111.83
2	A	501	J3N	C5-C4-N3	3.62	123.62	118.91
2	A	501	J3N	C15-C18-C19	-3.53	102.75	107.05
2	A	501	J3N	N1-C-N	-3.35	123.50	128.58
2	A	501	J3N	C13-C12-C11	-2.60	108.14	112.69
2	A	501	J3N	C14-C5-C17	2.56	120.56	117.92
2	A	501	J3N	C1-N2-N3	2.42	112.45	110.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	J3N	C15-C14-C5	-2.31	117.69	121.32
2	A	501	J3N	C6-N2-N3	-2.19	114.75	119.89

There are no chirality outliers.

All (10) torsion outliers are listed below:

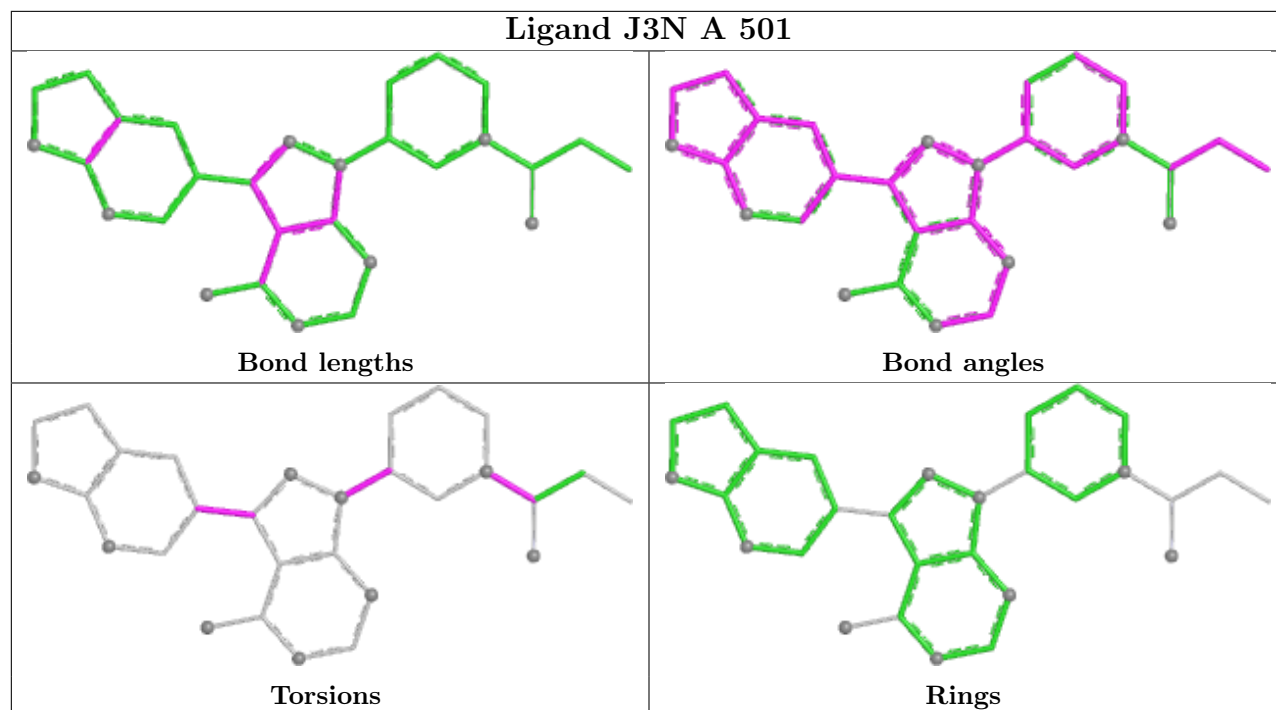
Mol	Chain	Res	Type	Atoms
2	A	501	J3N	O-C11-N5-C10
2	A	501	J3N	C12-C11-N5-C10
2	A	501	J3N	N3-C4-C5-C17
2	A	501	J3N	O-C11-N5-C9
2	A	501	J3N	C12-C11-N5-C9
2	A	501	J3N	C10-C6-N2-C1
2	A	501	J3N	C2-C4-C5-C17
2	A	501	J3N	C7-C6-N2-C1
2	A	501	J3N	C2-C4-C5-C14
2	A	501	J3N	N3-C4-C5-C14

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	J3N	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	286/318 (89%)	1.56	78 (27%) 1 1	20, 35, 53, 72	0

All (78) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	125	GLY	6.2
1	A	284	LEU	6.2
1	A	144	SER	6.1
1	A	315	TYR	5.0
1	A	388	ASN	4.8
1	A	316	ASP	4.5
1	A	295	GLY	4.0
1	A	307	PRO	4.0
1	A	358	LEU	3.9
1	A	143	GLY	3.8
1	A	146	THR	3.7
1	A	126	GLY	3.7
1	A	349	THR	3.6
1	A	417	THR	3.5
1	A	217	THR	3.5
1	A	386	LYS	3.4
1	A	117	LYS	3.3
1	A	403	GLU	3.3
1	A	360	PRO	3.3
1	A	294	ALA	3.2
1	A	283	ARG	3.2
1	A	308	PRO	3.2
1	A	359	LEU	3.1
1	A	412	ASP	2.9
1	A	148	GLY	2.9
1	A	297	ALA	2.8
1	A	145	GLY	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	121	TYR	2.8
1	A	244	LYS	2.8
1	A	392	GLU	2.8
1	A	191	ASP	2.6
1	A	336	PHE	2.6
1	A	248	TYR	2.6
1	A	350	LYS	2.6
1	A	397	LYS	2.6
1	A	362	HIS	2.6
1	A	136	LEU	2.6
1	A	214	LEU	2.5
1	A	344	ASP	2.5
1	A	124	ILE	2.5
1	A	410	PHE	2.5
1	A	147	CYS	2.5
1	A	415	ALA	2.5
1	A	247	TYR	2.4
1	A	222	LEU	2.4
1	A	357	PRO	2.4
1	A	364	GLY	2.4
1	A	304	ARG	2.4
1	A	385	PRO	2.3
1	A	401	THR	2.3
1	A	409	TRP	2.4
1	A	190	HIS	2.3
1	A	338	TYR	2.3
1	A	199	PHE	2.3
1	A	351	VAL	2.3
1	A	376	CYS	2.3
1	A	242	ILE	2.2
1	A	202	PHE	2.2
1	A	402	LEU	2.2
1	A	207	ASP	2.2
1	A	129	TYR	2.2
1	A	345	PHE	2.2
1	A	382	ARG	2.2
1	A	240	VAL	2.1
1	A	142	MET	2.1
1	A	231	PRO	2.1
1	A	342	LYS	2.1
1	A	228	GLY	2.1
1	A	306	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	134	ASN	2.1
1	A	229	PRO	2.1
1	A	413	VAL	2.0
1	A	141	GLU	2.0
1	A	408	SER	2.0
1	A	118	GLN	2.0
1	A	220	GLU	2.0
1	A	389	LYS	2.0
1	A	381	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](#)

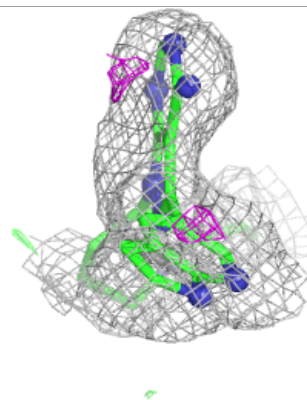
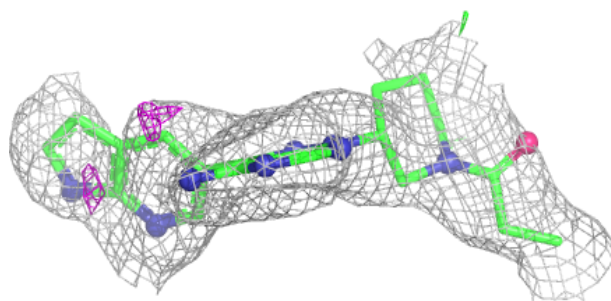
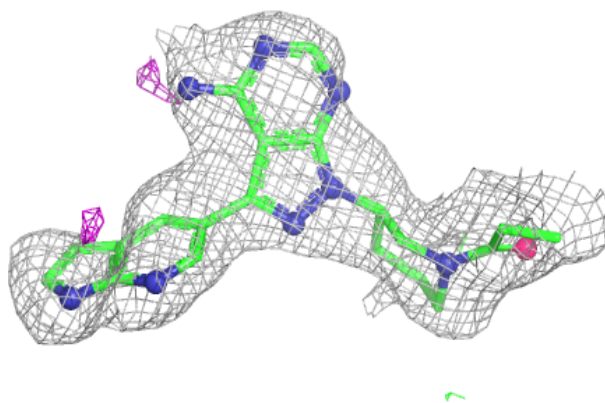
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	J3N	A	501	29/29	0.80	0.18	33,39,51,59	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 J3N A 501:

$2mF_o - DF_c$ (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.