



# Full wwPDB X-ray Structure Validation Report ⓘ

May 21, 2020 – 02:56 am BST

PDB ID : 4E73  
Title : Crystal structure of JNK1beta-JIP in complex with an azaquinolone inhibitor  
Authors : Lukacs, C.M.; Janson, C.A.  
Deposited on : 2012-03-16  
Resolution : 2.27 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.

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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

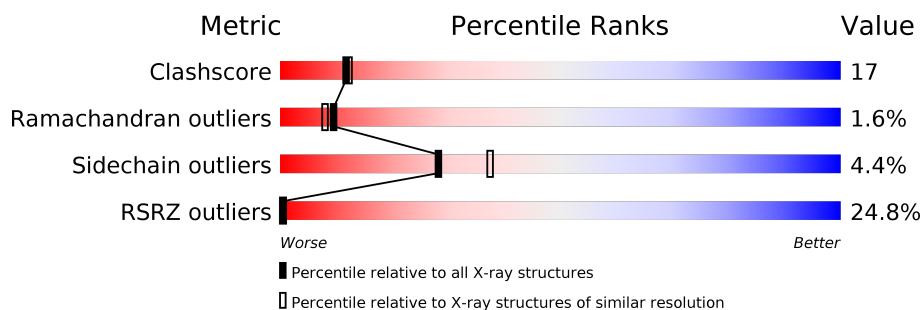
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.27 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))
Clashscore	141614	7711 (2.30-2.26)
Ramachandran outliers	138981	7597 (2.30-2.26)
Sidechain outliers	138945	7598 (2.30-2.26)
RSRZ outliers	127900	6849 (2.30-2.26)

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	369	<div> <div>21%</div> <div>59%</div> <div>23%</div> <div>• •</div> <div>14%</div> </div>
2	B	11	<div> <div>18%</div> <div>55%</div> <div>27%</div> <div>18%</div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 2784 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 Mitogen-activated protein kinase 8.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	318	2581	1659	435	467	20	0	1	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	364	HIS	-	EXPRESSION TAG	UNP P45983
A	365	HIS	-	EXPRESSION TAG	UNP P45983
A	366	HIS	-	EXPRESSION TAG	UNP P45983
A	367	HIS	-	EXPRESSION TAG	UNP P45983
A	368	HIS	-	EXPRESSION TAG	UNP P45983
A	369	HIS	-	EXPRESSION TAG	UNP P45983

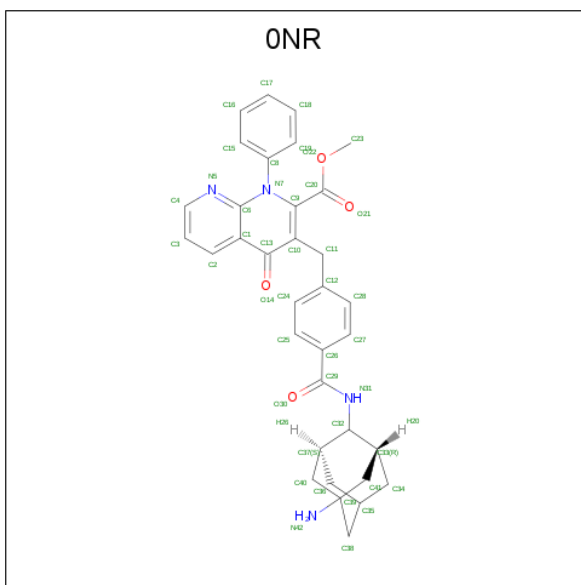
- Molecule 2 is a protein called C-Jun-amino-terminal kinase-interacting protein 1.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	B	9	72	46	14	12	0	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	153	LYS	ARG	CONFLICT	UNP Q9UQF2

- Molecule 3 is methyl 3-(4-{[(1R,2S,3S,5S,7s)-5-aminotricyclo[3.3.1.1<sup>3,7</sup>]dec-2-yl]carbamoyl}benzyl)-4-oxo-1-phenyl-1,4-dihydro-1,8-naphthyridine-2-carboxylate (three-letter code: 0NR) (formula: C<sub>34</sub>H<sub>34</sub>N<sub>4</sub>O<sub>4</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			42	34	4	4		

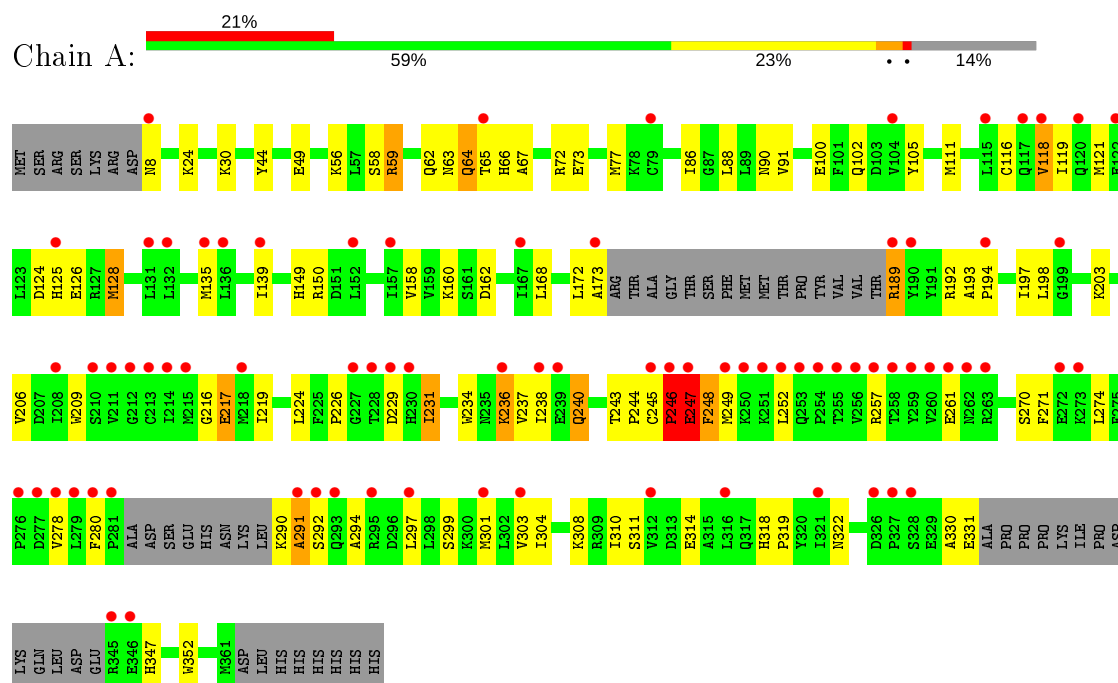
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	88	Total O 88 88	0	0
4	B	1	Total O 1 1	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: Mitogen-activated protein kinase 8



- Molecule 2: C-Jun-amino-terminal kinase-interacting protein 1



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	62.90Å 79.84Å 85.52Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	27.70 – 2.27 27.69 – 2.27	Depositor EDS
% Data completeness (in resolution range)	98.9 (27.70-2.27) 99.0 (27.69-2.27)	Depositor EDS
$R_{merge}$	0.04	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.72 (at 2.26Å)	Xtriage
Refinement program	CNX 2005	Depositor
R, $R_{free}$	0.247 , 0.282 0.248 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	48.8	Xtriage
Anisotropy	0.122	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 61.2	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	0.93	EDS
Total number of atoms	2784	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.73% 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: 0NR

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	0.32	0/2635	0.48	0/3555
2	B	0.33	0/73	0.63	0/98
All	All	0.32	0/2708	0.49	0/3653

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

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	246	PRO	Peptide

### 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	2581	0	2601	94	0
2	B	72	0	82	3	0
3	A	42	0	34	0	0
4	A	88	0	0	4	0
4	B	1	0	0	0	0
All	All	2784	0	2717	94	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 17.

All (94) 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:59:ARG:NH1	1:A:59:ARG:HB3	1.83	0.94
1:A:245:CYS:HB2	1:A:246:PRO:HD2	1.53	0.90
1:A:245:CYS:SG	1:A:247:GLU:HB3	2.14	0.88
1:A:231:ILE:H	1:A:231:ILE:HD13	1.42	0.84
1:A:63:ASN:HD22	1:A:65:THR:H	1.24	0.83
1:A:119:ILE:HD11	1:A:217:GLU:HB3	1.61	0.81
1:A:189:ARG:HD3	1:A:197:ILE:HG22	1.64	0.80
1:A:59:ARG:HB3	1:A:59:ARG:HH11	1.50	0.77
1:A:240:GLN:NE2	1:A:274:LEU:HD22	2.08	0.69
1:A:59:ARG:HB2	1:A:62:GLN:HB2	1.73	0.68
1:A:226:PRO:HD2	1:A:236:LYS:HD3	1.77	0.67
1:A:249:MET:O	1:A:257:ARG:HD2	1.95	0.67
1:A:216:GLY:HA3	1:A:224:LEU:HD11	1.77	0.66
1:A:63:ASN:ND2	1:A:65:THR:H	1.94	0.65
1:A:290:LYS:O	1:A:292:SER:N	2.31	0.64
1:A:271:PHE:CZ	1:A:299:SER:HA	2.31	0.64
1:A:249:MET:HA	1:A:252:LEU:HD12	1.81	0.62
1:A:118:VAL:O	1:A:121:MET:HG2	1.99	0.62
1:A:197:ILE:HG13	1:A:198:LEU:CD1	2.28	0.62
1:A:189:ARG:HG2	1:A:192:ARG:HD2	1.80	0.62
1:A:247:GLU:O	1:A:248:PHE:HB2	2.00	0.61
1:A:197:ILE:HG13	1:A:198:LEU:HD12	1.82	0.61
1:A:160:LYS:HB3	1:A:162:ASP:OD1	2.01	0.60
1:A:322:ASN:HB2	4:A:581:HOH:O	2.03	0.59
1:A:56:LYS:HE3	1:A:105:TYR:OH	2.03	0.58
1:A:58:SER:C	1:A:59:ARG:HG2	2.22	0.58
1:A:135[A]:MET:O	1:A:139:ILE:HG13	2.03	0.58
1:A:319:PRO:HA	1:A:322:ASN:HD21	1.69	0.57
1:A:194:PRO:O	1:A:198:LEU:HD13	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:245:CYS:HG	1:A:247:GLU:HB3	1.68	0.57
1:A:73:GLU:O	1:A:77:MET:HG3	2.04	0.57
1:A:67:ALA:HB1	1:A:352:TRP:HB3	1.87	0.57
1:A:118:VAL:HG21	2:B:160:LEU:HD21	1.85	0.56
1:A:319:PRO:HA	1:A:322:ASN:ND2	2.21	0.56
1:A:59:ARG:CZ	1:A:59:ARG:HB3	2.36	0.56
1:A:172:LEU:O	1:A:173:ALA:CB	2.54	0.55
1:A:63:ASN:HD22	1:A:65:THR:N	2.01	0.55
1:A:249:MET:HA	1:A:252:LEU:CD1	2.37	0.55
1:A:257:ARG:O	1:A:261:GLU:HB2	2.08	0.53
1:A:135[B]:MET:O	1:A:139:ILE:HG13	2.09	0.53
1:A:63:ASN:HD21	1:A:65:THR:HB	1.73	0.53
1:A:116:CYS:O	1:A:119:ILE:HG22	2.08	0.53
1:A:198:LEU:HD11	1:A:234:TRP:CE3	2.45	0.52
1:A:126:GLU:HG2	2:B:157:PRO:HG3	1.90	0.52
1:A:297:LEU:HD11	1:A:301:MET:HE2	1.92	0.51
1:A:111:MET:HG3	1:A:158:VAL:HG23	1.93	0.51
1:A:100:GLU:O	1:A:102:GLN:HG2	2.10	0.50
1:A:244:PRO:O	1:A:249:MET:HE2	2.11	0.50
1:A:59:ARG:HG3	1:A:66:HIS:CE1	2.46	0.49
1:A:270:SER:O	1:A:274:LEU:HG	2.12	0.49
1:A:203:LYS:O	1:A:206:VAL:HG12	2.12	0.49
1:A:172:LEU:O	1:A:173:ALA:HB3	2.13	0.48
1:A:234:TRP:O	1:A:238:ILE:HG12	2.12	0.48
1:A:310:ILE:HG12	1:A:311:SER:N	2.28	0.48
1:A:247:GLU:O	1:A:248:PHE:CB	2.62	0.48
1:A:229:ASP:HB3	1:A:231:ILE:HD11	1.96	0.47
1:A:231:ILE:H	1:A:231:ILE:CD1	2.11	0.47
1:A:330:ALA:O	1:A:331:GLU:HG3	2.14	0.47
1:A:318:HIS:CG	1:A:319:PRO:HD2	2.50	0.47
1:A:246:PRO:O	1:A:248:PHE:N	2.49	0.46
1:A:124:ASP:O	1:A:128:MET:HB2	2.15	0.46
1:A:72:ARG:HD3	4:A:583:HOH:O	2.16	0.46
1:A:72:ARG:NE	4:A:588:HOH:O	2.26	0.46
1:A:189:ARG:CG	1:A:192:ARG:HD2	2.45	0.45
1:A:119:ILE:CD1	1:A:217:GLU:HB3	2.38	0.45
1:A:245:CYS:HB2	1:A:246:PRO:CD	2.33	0.45
1:A:149:HIS:O	1:A:150:ARG:HB2	2.16	0.45
1:A:72:ARG:CZ	1:A:173:ALA:HB3	2.46	0.45
1:A:237:VAL:HG12	1:A:304:ILE:HD11	1.98	0.44
1:A:231:ILE:HD13	1:A:231:ILE:N	2.23	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:229:ASP:HB3	1:A:231:ILE:CD1	2.48	0.44
1:A:125:HIS:CD2	1:A:290:LYS:HB2	2.53	0.44
1:A:63:ASN:ND2	1:A:65:THR:N	2.64	0.43
1:A:189:ARG:NH1	1:A:192:ARG:HH11	2.15	0.43
1:A:193:ALA:HA	1:A:209:TRP:CD1	2.54	0.43
1:A:310:ILE:HD11	1:A:314:GLU:HB3	2.01	0.43
1:A:56:LYS:HE3	1:A:105:TYR:HH	1.83	0.43
1:A:240:GLN:HB3	1:A:240:GLN:HE21	1.52	0.43
1:A:126:GLU:HG2	2:B:157:PRO:CG	2.49	0.42
1:A:44:TYR:OH	1:A:49:GLU:HA	2.19	0.42
1:A:8:ASN:O	1:A:24:LYS:HE3	2.19	0.42
1:A:219:ILE:HD13	1:A:280:PHE:CZ	2.55	0.42
1:A:280:PHE:CD1	1:A:291:ALA:HA	2.54	0.42
1:A:86:ILE:HG13	1:A:168:LEU:HD12	2.03	0.41
1:A:88:LEU:HD11	1:A:91:VAL:CG2	2.50	0.41
1:A:280:PHE:HE1	1:A:294:ALA:HB3	1.86	0.41
1:A:189:ARG:CD	1:A:197:ILE:HG22	2.43	0.41
1:A:280:PHE:CE1	1:A:294:ALA:HB3	2.56	0.41
1:A:64:GLN:NE2	1:A:347:HIS:O	2.53	0.41
1:A:303:VAL:HG11	1:A:308:LYS:HB2	2.03	0.41
1:A:245:CYS:CB	1:A:246:PRO:HD2	2.33	0.40
1:A:203:LYS:NZ	4:A:586:HOH:O	2.53	0.40
1:A:243:THR:HG22	1:A:244:PRO:O	2.22	0.40
1:A:88:LEU:HD11	1:A:91:VAL:HG22	2.03	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	311/369 (84%)	293 (94%)	14 (4%)	4 (1%)	12 11

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	7/11 (64%)	6 (86%)	0	1 (14%)	0	0
All	All	318/380 (84%)	299 (94%)	14 (4%)	5 (2%)	9	8

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	248	PHE
1	A	291	ALA
2	B	155	LYS
1	A	246	PRO
1	A	247	GLU

### 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	285/331 (86%)	272 (95%)	13 (5%)	27	35
2	B	9/11 (82%)	9 (100%)	0	100	100
All	All	294/342 (86%)	281 (96%)	13 (4%)	28	37

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	LYS
1	A	59	ARG
1	A	64	GLN
1	A	90	ASN
1	A	118	VAL
1	A	128	MET
1	A	189	ARG
1	A	217	GLU
1	A	231	ILE
1	A	236	LYS
1	A	240	GLN

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Mol	Chain	Res	Type
1	A	247	GLU
1	A	278	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	9	ASN
1	A	27	GLN
1	A	62	GLN
1	A	63	ASN
1	A	90	ASN
1	A	240	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 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	$\# Z  > 2$
3	0NR	A	401	-	45,48,48	2.00	20 (44%)	53,72,72	1.16	6 (11%)

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
3	0NR	A	401	-	-	2/22/53/53	0/8/7/7

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	401	0NR	C13-C10	3.82	1.52	1.41
3	A	401	0NR	C4-N5	3.17	1.38	1.32
3	A	401	0NR	C41-C39	3.11	1.58	1.53
3	A	401	0NR	C19-C8	3.09	1.44	1.38
3	A	401	0NR	C15-C8	3.04	1.44	1.38
3	A	401	0NR	C6-N5	2.93	1.39	1.35
3	A	401	0NR	C38-C39	2.87	1.57	1.53
3	A	401	0NR	C3-C2	2.87	1.43	1.36
3	A	401	0NR	C40-C39	2.74	1.57	1.53
3	A	401	0NR	C27-C26	2.65	1.43	1.39
3	A	401	0NR	C28-C27	2.48	1.43	1.38
3	A	401	0NR	C29-N31	2.35	1.39	1.34
3	A	401	0NR	C28-C12	2.23	1.43	1.38
3	A	401	0NR	C41-C33	2.22	1.57	1.53
3	A	401	0NR	C11-C10	2.16	1.56	1.52
3	A	401	0NR	C25-C26	2.12	1.42	1.39
3	A	401	0NR	C40-C37	2.09	1.57	1.53
3	A	401	0NR	C34-C33	2.08	1.58	1.53
3	A	401	0NR	C39-N42	-2.07	1.42	1.49
3	A	401	0NR	C24-C12	2.06	1.43	1.38

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	401	0NR	C1-C6-N5	-3.22	119.65	124.62
3	A	401	0NR	C4-N5-C6	3.05	120.55	116.77
3	A	401	0NR	C19-C8-N7	2.76	121.80	119.24
3	A	401	0NR	C10-C11-C12	-2.35	108.75	114.10
3	A	401	0NR	O22-C20-O21	2.26	127.87	123.45
3	A	401	0NR	C11-C10-C13	-2.26	115.68	119.29

There are no chirality outliers.

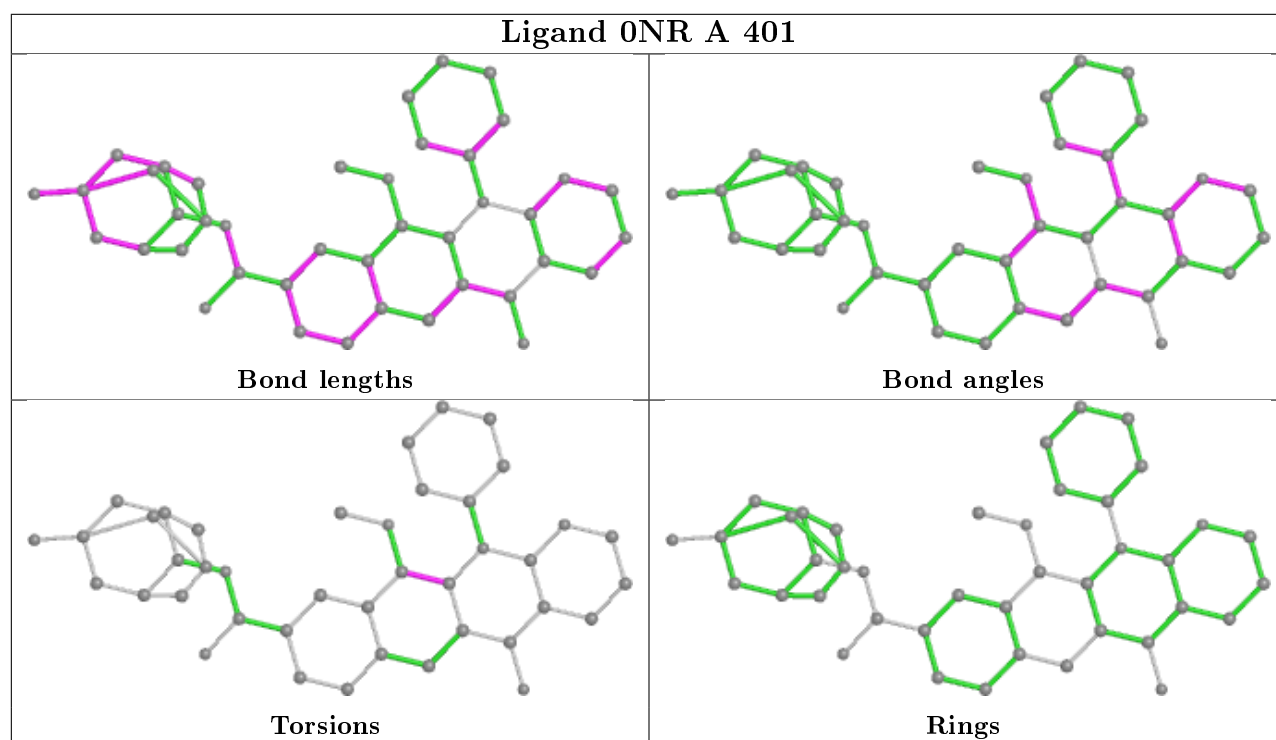
All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	401	0NR	O22-C20-C9-N7
3	A	401	0NR	O22-C20-C9-C10

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 ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues

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	318/369 (86%)	1.15	79 (24%) <b>0</b> <b>0</b>	29, 56, 99, 113	0
2	B	9/11 (81%)	1.02	2 (22%) <b>0</b> <b>1</b>	66, 67, 80, 82	0
All	All	327/380 (86%)	1.14	81 (24%) <b>0</b> <b>0</b>	29, 56, 99, 113	0

All (81) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	253	GLN	7.8
1	A	291	ALA	7.7
1	A	258	THR	6.5
1	A	278	VAL	6.4
1	A	256	VAL	6.4
1	A	229	ASP	6.1
1	A	247	GLU	5.9
1	A	211	VAL	5.9
1	A	254	PRO	5.4
1	A	292	SER	5.1
1	A	250	LYS	5.1
1	A	345	ARG	5.0
1	A	135[A]	MET	4.6
1	A	228	THR	4.6
1	A	280	PHE	4.6
1	A	279	LEU	4.5
1	A	273	LYS	4.4
1	A	328	SER	4.4
1	A	246	PRO	4.3
1	A	152	LEU	4.3
1	A	326	ASP	4.2
1	A	281	PRO	4.1
1	A	214	ILE	4.1
1	A	189	ARG	3.9

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Mol	Chain	Res	Type	RSRZ
1	A	8	ASN	3.9
1	A	276	PRO	3.8
1	A	131	LEU	3.7
1	A	136	LEU	3.5
1	A	249	MET	3.5
1	A	262	ASN	3.5
1	A	227	GLY	3.5
1	A	139	ILE	3.4
1	A	132	LEU	3.3
1	A	115	LEU	3.2
1	A	65	THR	3.2
1	A	157	ILE	3.2
1	A	190	TYR	3.1
1	A	120	GLN	3.1
1	A	210	SER	3.1
1	A	245	CYS	3.0
1	A	208	ILE	2.9
1	A	79	CYS	2.8
1	A	239	GLU	2.7
1	A	117	GLN	2.7
1	A	301	MET	2.7
1	A	303	VAL	2.7
1	A	230	HIS	2.6
1	A	260	VAL	2.6
1	A	295	ARG	2.6
1	A	125	HIS	2.6
1	A	213	CYS	2.6
1	A	327	PRO	2.6
1	A	118	VAL	2.5
1	A	257	ARG	2.5
1	A	122	GLU	2.5
1	A	261	GLU	2.5
2	B	158	THR	2.5
1	A	173	ALA	2.5
1	A	199	GLY	2.5
1	A	215	MET	2.4
1	A	194	PRO	2.4
1	A	236	LYS	2.4
1	A	255	THR	2.3
1	A	293	GLN	2.3
1	A	252	LEU	2.3
1	A	212	GLY	2.3

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Mol	Chain	Res	Type	RSRZ
2	B	155	LYS	2.3
1	A	167	ILE	2.3
1	A	104	VAL	2.3
1	A	272	GLU	2.2
1	A	218	MET	2.2
1	A	259	TYR	2.2
1	A	321	ILE	2.2
1	A	263	ARG	2.1
1	A	312	VAL	2.1
1	A	316	LEU	2.1
1	A	346	GLU	2.1
1	A	277	ASP	2.1
1	A	297	LEU	2.1
1	A	238	ILE	2.0
1	A	251	LYS	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](#)

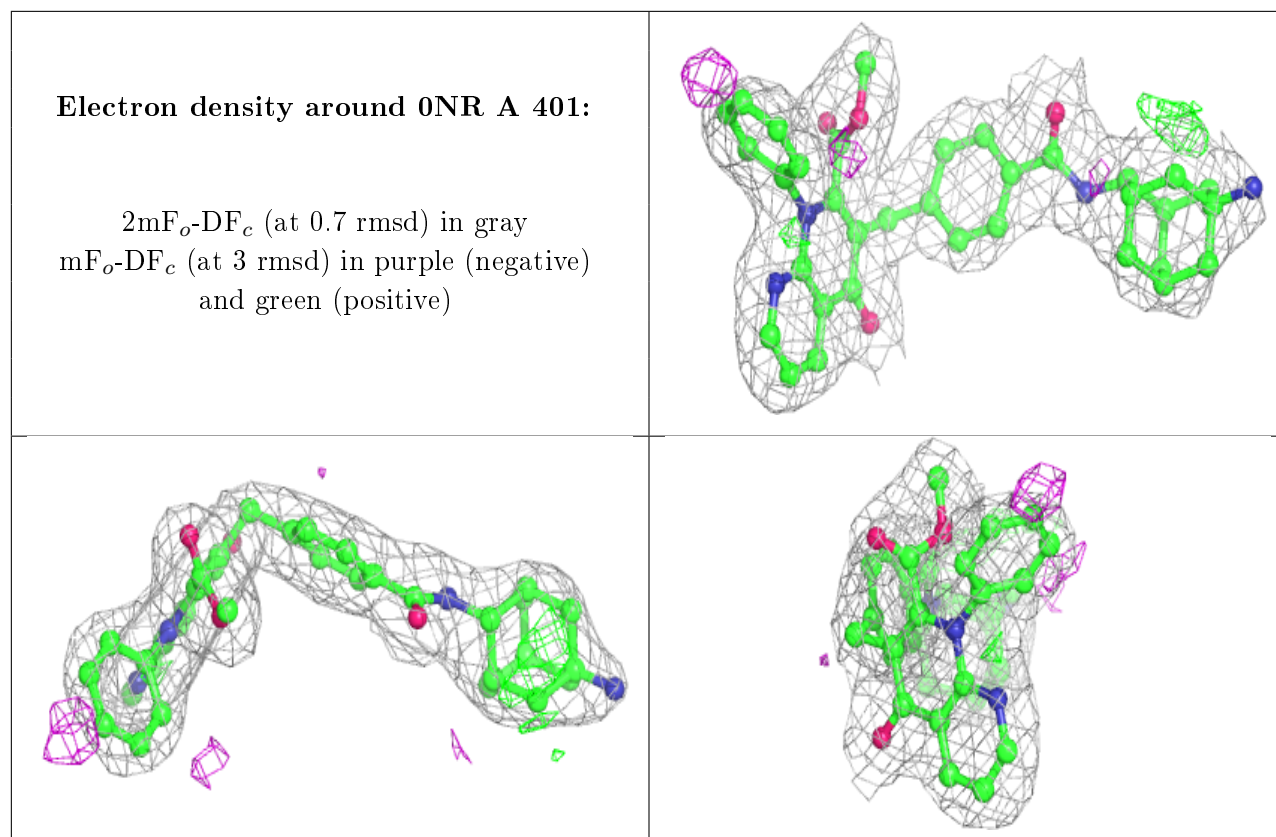
There are no carbohydrates in this entry.

## 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, 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
3	0NR	A	401	42/42	0.93	0.15	24,39,63,63	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.