



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 07:04 PM UTC

PDB ID : 7LVH / pdb_00007lvh
Title : CRYSTAL STRUCTURE OF AP2 ASSOCIATED KINASE 1 ISOFORM 1
COMPLEXED WITH LIGAND N-[3-METHOXY-4-(1,3-OXAZOL-5-YL)PH
ENYL]-3-(PROPAN-2-YL)PIPERIDINE-2-CARBOXAMIDE
Authors : Muckelbauer, J.K.
Deposited on : 2021-02-25
Resolution : 2.65 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

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

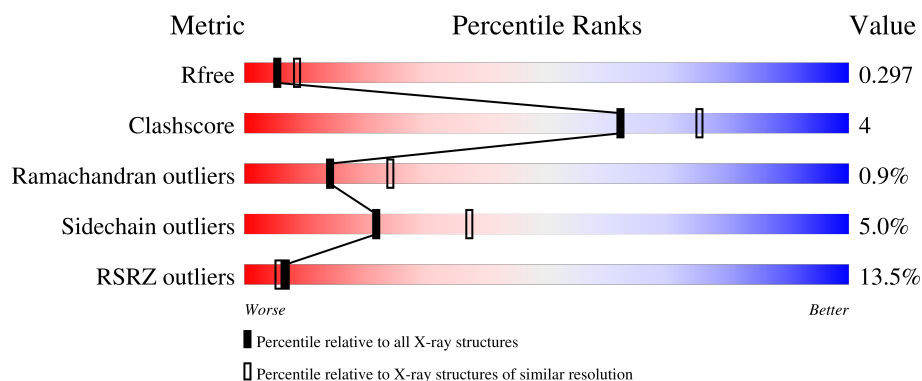
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.65 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}	180053	1110 (2.66-2.66)
Clashscore	190562	1141 (2.66-2.66)
Ramachandran outliers	187476	1126 (2.66-2.66)
Sidechain outliers	187428	1126 (2.66-2.66)
RSRZ outliers	180081	1110 (2.66-2.66)

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 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	12% 79% 8% 11%
1	B	318	12% 76% 10% 12%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4305 atoms, of which 49 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 AP2-associated protein kinase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	284	Total	C	N	O	S	0	0	0
			2097	1340	343	395	19			
1	B	281	Total	C	N	O	S	0	0	0
			2084	1332	344	389	19			

There are 26 discrepancies between the modelled and reference sequences:

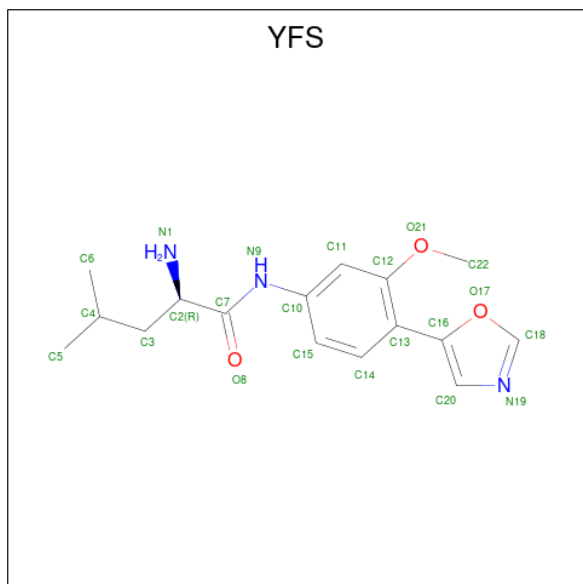
Chain	Residue	Modelled	Actual	Comment	Reference
A	13	MET	-	initiating methionine	UNP Q3UHJ0
A	14	HIS	-	expression tag	UNP Q3UHJ0
A	15	HIS	-	expression tag	UNP Q3UHJ0
A	16	HIS	-	expression tag	UNP Q3UHJ0
A	17	HIS	-	expression tag	UNP Q3UHJ0
A	18	HIS	-	expression tag	UNP Q3UHJ0
A	19	HIS	-	expression tag	UNP Q3UHJ0
A	20	LEU	-	expression tag	UNP Q3UHJ0
A	21	VAL	-	expression tag	UNP Q3UHJ0
A	22	PRO	-	expression tag	UNP Q3UHJ0
A	23	ARG	-	expression tag	UNP Q3UHJ0
A	24	GLY	-	expression tag	UNP Q3UHJ0
A	25	SER	-	expression tag	UNP Q3UHJ0
B	13	MET	-	initiating methionine	UNP Q3UHJ0
B	14	HIS	-	expression tag	UNP Q3UHJ0
B	15	HIS	-	expression tag	UNP Q3UHJ0
B	16	HIS	-	expression tag	UNP Q3UHJ0
B	17	HIS	-	expression tag	UNP Q3UHJ0
B	18	HIS	-	expression tag	UNP Q3UHJ0
B	19	HIS	-	expression tag	UNP Q3UHJ0
B	20	LEU	-	expression tag	UNP Q3UHJ0
B	21	VAL	-	expression tag	UNP Q3UHJ0
B	22	PRO	-	expression tag	UNP Q3UHJ0
B	23	ARG	-	expression tag	UNP Q3UHJ0
B	24	GLY	-	expression tag	UNP Q3UHJ0

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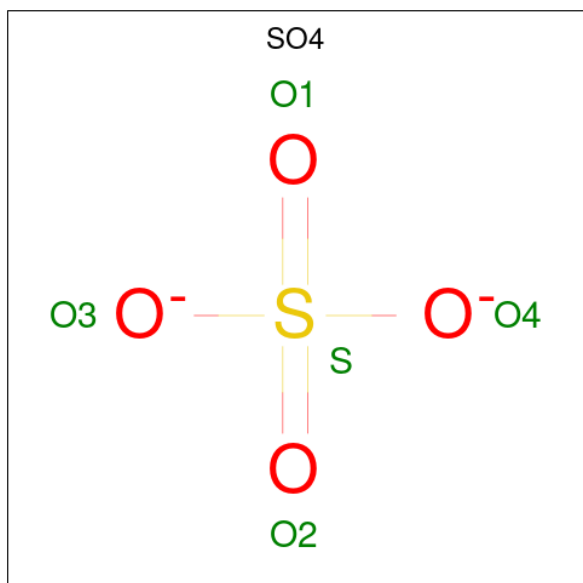
Chain	Residue	Modelled	Actual	Comment	Reference
B	25	SER	-	expression tag	UNP Q3UHQ0

- Molecule 2 is N-[3-methoxy-4-(1,3-oxazol-5-yl)phenyl]-D-leucinamide (CCD ID: YFS) (formula: $C_{16}H_{21}N_3O_3$).



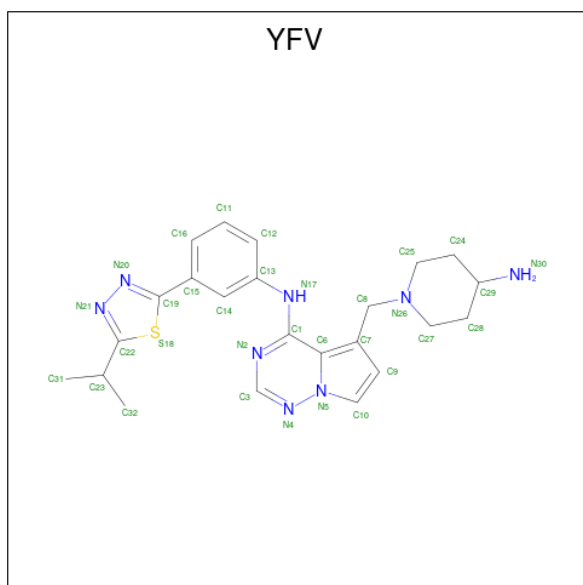
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	H	N	O		
2	A	1	43	16	21	3	3	0	0

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O_4S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0

- Molecule 4 is 5-[(4-aminopiperidin-1-yl)methyl]-N-{3-[5-(propan-2-yl)-1,3,4-thiadiazol-2-yl]phenyl}pyrrolo[2,1-f][1,2,4]triazin-4-amine (CCD ID: YFV) (formula: C₂₃H₂₈N₈S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total C H N S 60 23 28 8 1	0	0

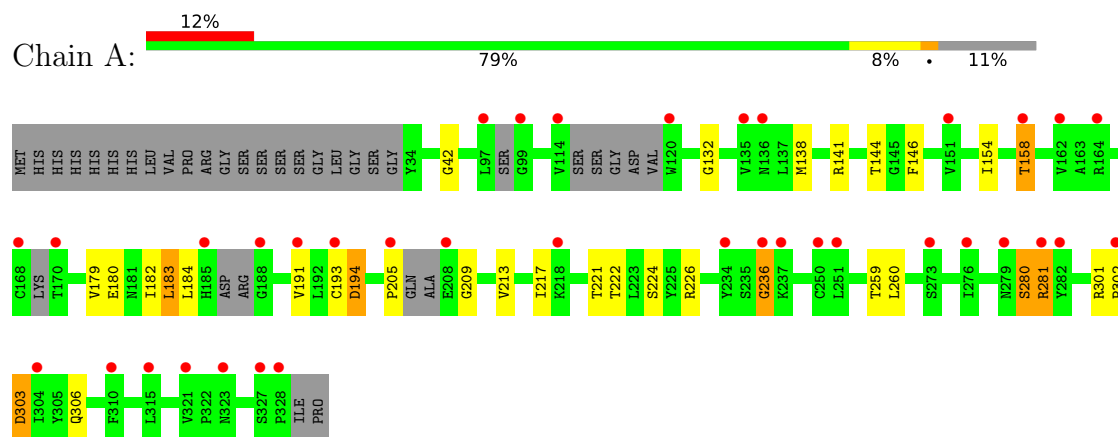
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	5	Total O 5 5	0	0
5	B	6	Total O 6 6	0	0

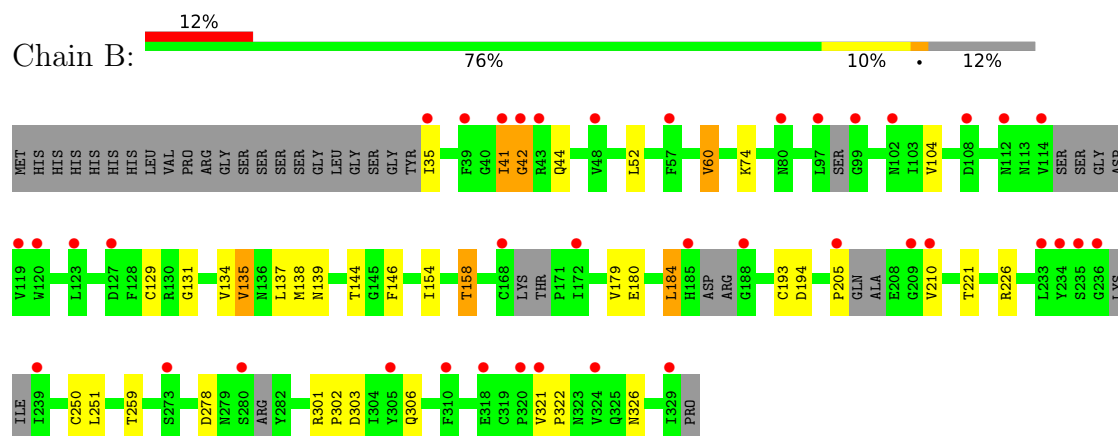
3 Residue-property plots [i](#)

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: AP2-associated protein kinase 1



- Molecule 1: AP2-associated protein kinase 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	91.51Å 91.51Å 172.15Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.67 – 2.65 30.67 – 2.65	Depositor EDS
% Data completeness (in resolution range)	100.0 (30.67-2.65) 99.9 (30.67-2.65)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.11 (at 2.65Å)	Xtriage
Refinement program	BUSTER 2.11.5	Depositor
R, R_{free}	0.265 , 0.299 0.271 , 0.297	Depositor DCC
R_{free} test set	1267 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å ²)	61.9	Xtriage
Anisotropy	0.080	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 40.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.018 for -h,-k,l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	4305	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.89% 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: YFS, SO4, YFV

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	0.89	0/2136	1.39	10/2908 (0.3%)
1	B	0.87	0/2120	1.39	10/2878 (0.3%)
All	All	0.88	0/4256	1.39	20/5786 (0.3%)

There are no bond length outliers.

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	280	SER	CA-C-N	7.87	135.87	121.70
1	A	280	SER	C-N-CA	7.87	135.87	121.70
1	A	236	GLY	CA-C-N	7.74	135.63	121.70
1	A	236	GLY	C-N-CA	7.74	135.63	121.70
1	A	42	GLY	CA-C-N	6.20	132.86	121.70
1	A	42	GLY	C-N-CA	6.20	132.86	121.70
1	A	146	PHE	CA-CB-CG	5.96	119.76	113.80
1	B	42	GLY	CA-C-N	5.92	132.36	121.70
1	B	42	GLY	C-N-CA	5.92	132.36	121.70
1	B	131	GLY	N-CA-C	-5.84	105.06	112.54
1	B	146	PHE	CA-CB-CG	5.83	119.63	113.80
1	B	184	LEU	CA-C-N	5.80	132.14	121.70
1	B	184	LEU	C-N-CA	5.80	132.14	121.70
1	B	129	CYS	CA-C-N	5.79	128.04	120.28
1	B	129	CYS	C-N-CA	5.79	128.04	120.28
1	B	250	CYS	CA-C-N	5.22	127.54	120.44
1	B	250	CYS	C-N-CA	5.22	127.54	120.44
1	A	184	LEU	CA-C-N	5.12	130.91	121.70
1	A	184	LEU	C-N-CA	5.12	130.91	121.70
1	A	303	ASP	CA-CB-CG	5.07	117.67	112.60

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	2097	0	1910	15	0
1	B	2084	0	1927	20	0
2	A	22	21	0	0	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
4	B	32	28	0	0	0
5	A	5	0	0	0	0
5	B	6	0	0	0	0
All	All	4256	49	3837	33	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 (33) 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:B:135:VAL:HA	1:B:138:MET:HE3	1.79	0.64
1:A:205:PRO:HB3	1:A:213:VAL:HG11	1.79	0.64
1:B:154:ILE:O	1:B:158:THR:HG23	2.02	0.60
1:B:221:THR:HG21	1:B:226:ARG:HG3	1.84	0.60
1:A:154:ILE:O	1:A:158:THR:HG23	2.02	0.59
1:A:221:THR:HG22	1:A:226:ARG:HE	1.69	0.58
1:A:141:ARG:NH2	1:B:322:PRO:HG3	2.23	0.53
1:B:42:GLY:HA3	1:B:44:GLN:H	1.74	0.53
1:A:141:ARG:HH22	1:B:322:PRO:HG3	1.74	0.51
1:B:52:LEU:HB2	1:B:60:VAL:HG23	1.93	0.51
1:A:183:LEU:HD11	1:A:193:CYS:HB3	1.92	0.50
1:B:41:ILE:O	1:B:44:GLN:HB2	2.12	0.49
1:A:224:SER:CB	1:A:260:LEU:HD13	2.43	0.49
1:B:137:LEU:HD11	1:B:184:LEU:HD22	1.94	0.49
1:A:224:SER:HB3	1:A:260:LEU:HD13	1.95	0.49
1:B:104:VAL:HG22	1:B:193:CYS:HA	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:221:THR:CG2	1:A:226:ARG:HE	2.26	0.48
1:A:301:ARG:HG3	1:A:302:PRO:HD2	1.96	0.48
1:B:221:THR:CG2	1:B:226:ARG:HG3	2.45	0.47
1:B:134:VAL:HG12	1:B:138:MET:HE2	1.96	0.46
1:B:301:ARG:HG3	1:B:302:PRO:HD2	1.97	0.46
1:A:303:ASP:H	1:A:306:GLN:HE21	1.64	0.45
1:A:280:SER:HA	1:A:281:ARG:CB	2.47	0.44
1:B:42:GLY:CA	1:B:44:GLN:H	2.32	0.42
1:B:60:VAL:HG12	1:B:74:LYS:HG3	2.01	0.42
1:A:217:ILE:O	1:A:221:THR:HB	2.19	0.42
1:B:179:VAL:CG1	1:B:251:LEU:HD12	2.50	0.41
1:B:179:VAL:HG12	1:B:251:LEU:HD12	2.02	0.41
1:A:209:GLY:O	1:A:213:VAL:HG12	2.20	0.41
1:B:205:PRO:HB2	1:B:210:VAL:HG22	2.03	0.41
1:B:303:ASP:H	1:B:306:GLN:HE21	1.68	0.41
1:B:135:VAL:HA	1:B:138:MET:CE	2.50	0.40
1:A:138:MET:HE1	1:A:179:VAL:HG11	2.03	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	272/318 (86%)	255 (94%)	13 (5%)	4 (2%)	8	13
1	B	265/318 (83%)	254 (96%)	10 (4%)	1 (0%)	30	45
All	All	537/636 (84%)	509 (95%)	23 (4%)	5 (1%)	14	24

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	194	ASP

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Mol	Chain	Res	Type
1	A	132	GLY
1	A	236	GLY
1	B	194	ASP
1	A	281	ARG

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	208/283 (74%)	199 (96%)	9 (4%)	26	44
1	B	211/283 (75%)	199 (94%)	12 (6%)	18	32
All	All	419/566 (74%)	398 (95%)	21 (5%)	22	37

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	144	THR
1	A	158	THR
1	A	180	GLU
1	A	182	ILE
1	A	183	LEU
1	A	191	VAL
1	A	194	ASP
1	A	222	THR
1	A	259	THR
1	B	35	ILE
1	B	41	ILE
1	B	60	VAL
1	B	135	VAL
1	B	139	ASN
1	B	144	THR
1	B	158	THR
1	B	180	GLU
1	B	259	THR
1	B	278	ASP

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Mol	Chain	Res	Type
1	B	321	VAL
1	B	326	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	166	HIS
1	A	174	HIS
1	A	203	GLN
1	A	232	ASN
1	A	306	GLN
1	B	166	HIS
1	B	306	GLN
1	B	325	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](#)

4 ligands are 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	YFS	A	401	-	22,23,23	0.21	0	24,31,31	0.35	0
3	SO4	B	402	-	4,4,4	0.31	0	6,6,6	0.18	0
3	SO4	A	402	-	4,4,4	0.23	0	6,6,6	0.34	0
4	YFV	B	401	-	34,36,36	0.47	0	39,51,51	0.72	1 (2%)

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	YFS	A	401	-	-	0/18/18/18	0/2/2/2
4	YFV	B	401	-	-	4/16/26/26	0/5/5/5

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	401	YFV	C13-N17-C1	2.68	134.57	128.41

There are no chirality outliers.

All (4) torsion outliers are listed below:

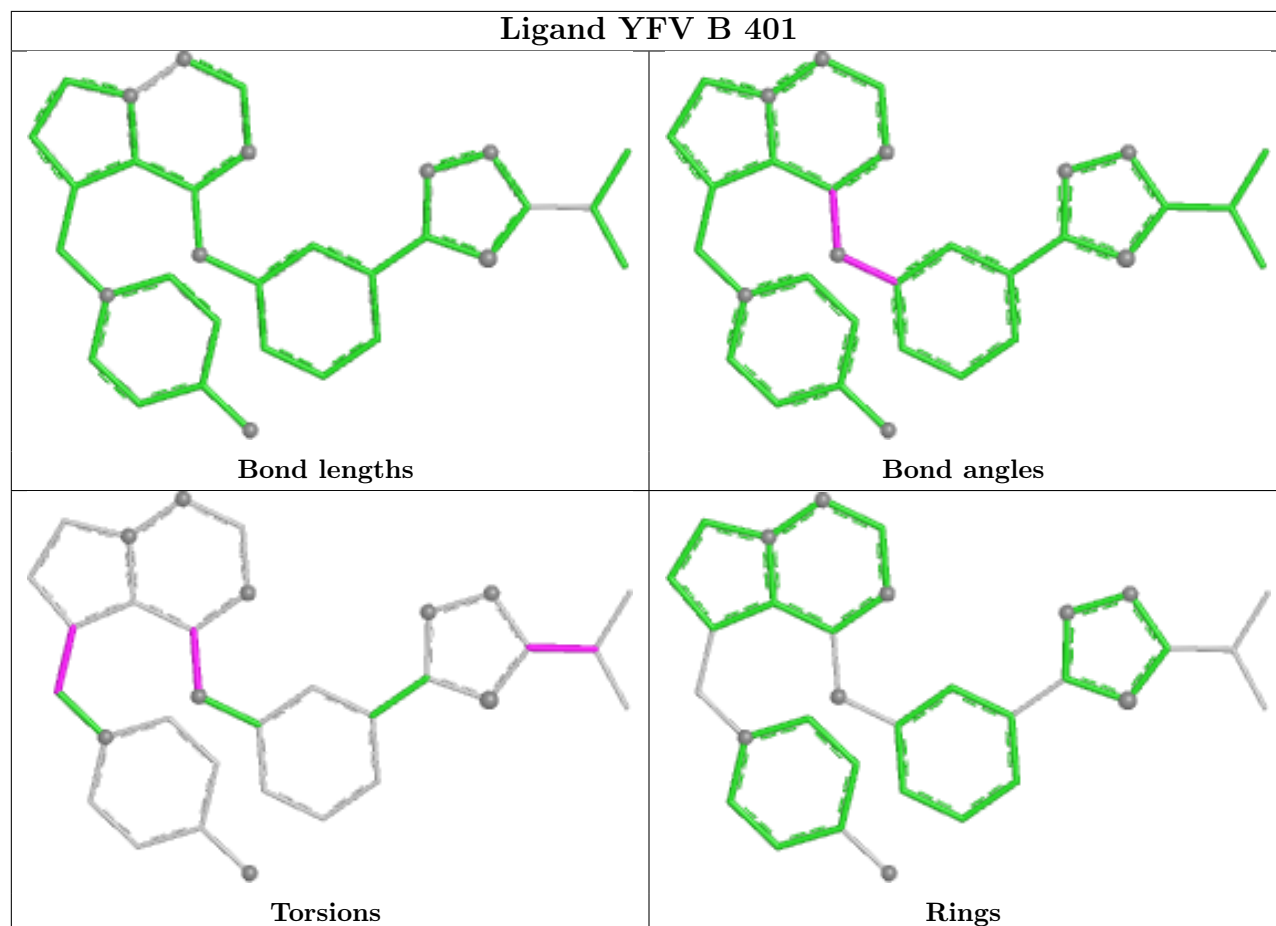
Mol	Chain	Res	Type	Atoms
4	B	401	YFV	N2-C1-N17-C13
4	B	401	YFV	S18-C22-C23-C32
4	B	401	YFV	C9-C7-C8-N26
4	B	401	YFV	C6-C1-N17-C13

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, 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	284/318 (89%)	1.10	37 (13%) 7 6	38, 60, 81, 93	0
1	B	281/318 (88%)	0.96	39 (13%) 6 5	37, 59, 83, 110	0
All	All	565/636 (88%)	1.03	76 (13%) 7 5	37, 60, 83, 110	0

All (76) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	114	VAL	5.6
1	A	188	GLY	5.4
1	A	185	HIS	4.8
1	A	276	ILE	4.6
1	A	114	VAL	4.5
1	B	43	ARG	4.2
1	A	208	GLU	4.1
1	B	280	SER	4.1
1	A	97	LEU	4.1
1	A	250	CYS	4.0
1	A	205	PRO	4.0
1	B	41	ILE	4.0
1	B	97	LEU	3.7
1	B	329	ILE	3.7
1	A	193	CYS	3.6
1	B	185	HIS	3.6
1	A	315	LEU	3.6
1	A	168	CYS	3.6
1	B	321	VAL	3.5
1	B	318	GLU	3.4
1	A	282	TYR	3.4
1	A	304	ILE	3.3
1	B	236	GLY	3.3
1	B	168	CYS	3.2

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Mol	Chain	Res	Type	RSRZ
1	B	120	TRP	3.2
1	B	119	VAL	3.2
1	B	42	GLY	3.1
1	B	234	TYR	3.1
1	B	35	ILE	3.0
1	A	99	GLY	3.0
1	B	112	ASN	2.9
1	A	170	THR	2.8
1	B	57	PHE	2.8
1	A	279	ASN	2.7
1	A	327	SER	2.7
1	B	210	VAL	2.6
1	B	99	GLY	2.5
1	A	191	VAL	2.5
1	A	251	LEU	2.5
1	A	328	PRO	2.5
1	A	151	VAL	2.5
1	A	162	VAL	2.5
1	A	281	ARG	2.5
1	B	102	ASN	2.4
1	B	108	ASP	2.4
1	A	310	PHE	2.4
1	B	324	VAL	2.4
1	A	136	ASN	2.4
1	B	127	ASP	2.4
1	A	135	VAL	2.4
1	B	239	ILE	2.3
1	B	188	GLY	2.3
1	A	321	VAL	2.3
1	A	158	THR	2.3
1	B	205	PRO	2.3
1	B	209	GLY	2.3
1	B	39	PHE	2.3
1	A	237	LYS	2.2
1	A	273	SER	2.2
1	B	273	SER	2.2
1	A	120	TRP	2.2
1	A	302	PRO	2.2
1	A	236	GLY	2.2
1	B	123	LEU	2.2
1	A	164	ARG	2.1
1	A	323	ASN	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	80	ASN	2.1
1	B	233	LEU	2.1
1	B	235	SER	2.1
1	A	234	TYR	2.1
1	A	218	LYS	2.1
1	B	305	TYR	2.1
1	B	320	PRO	2.1
1	B	172	ILE	2.1
1	B	310	PHE	2.0
1	B	48	VAL	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](#)

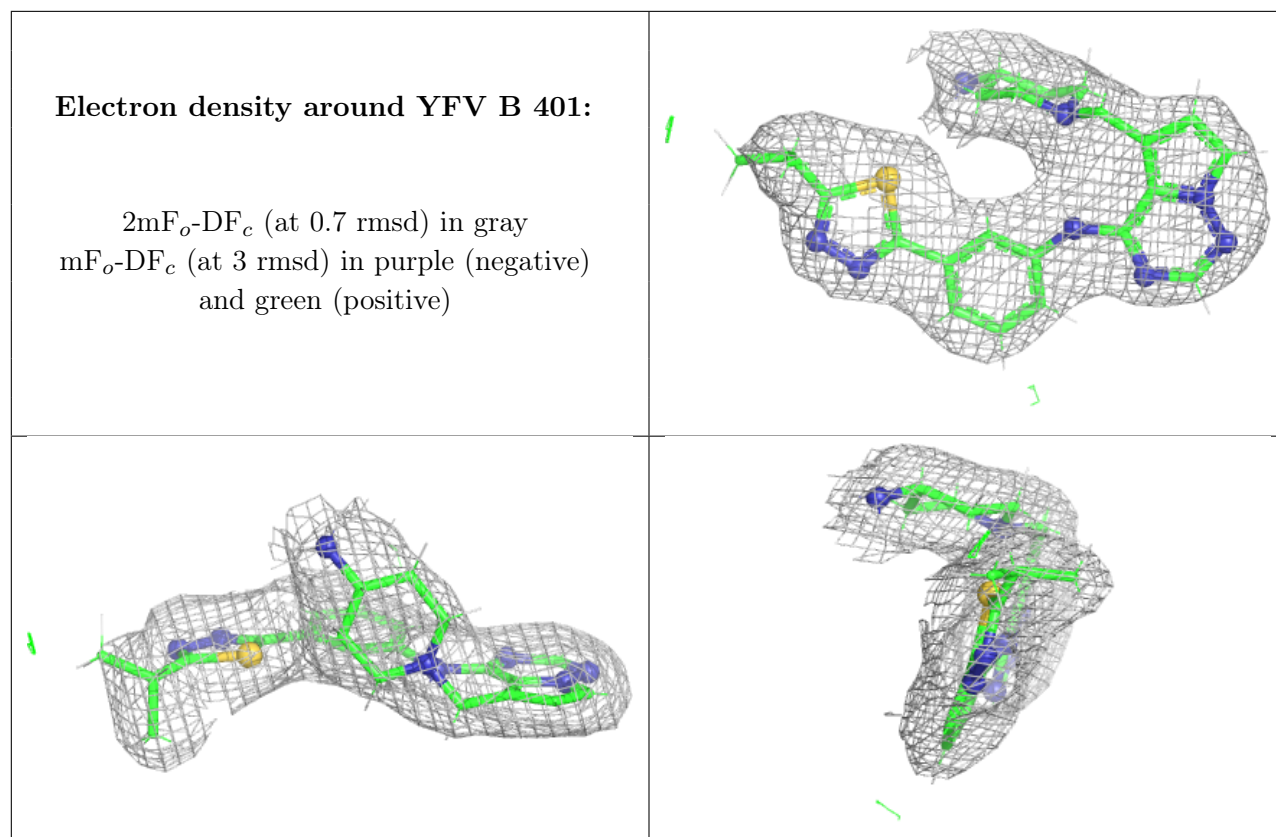
There are no oligosaccharides in this entry.

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(Å ²)	Q<0.9
2	YFS	A	401	22/22	0.88	0.11	39,48,55,58	0
4	YFV	B	401	32/32	0.91	0.10	50,55,71,72	0
3	SO4	A	402	5/5	0.94	0.10	57,60,62,63	0
3	SO4	B	402	5/5	0.96	0.09	77,77,78,79	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.