



wwPDB X-ray Structure Validation Summary Report ⓘ

Sep 30, 2026 – 04:06 PM EDT

PDB ID : 9YHP / pdb_00009yhp
Title : FEM1B with small molecule activator
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Deposited on : 2025-09-30
Resolution : 3.50 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 Phenix2.2
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.2
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 : 9.0.015 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.51

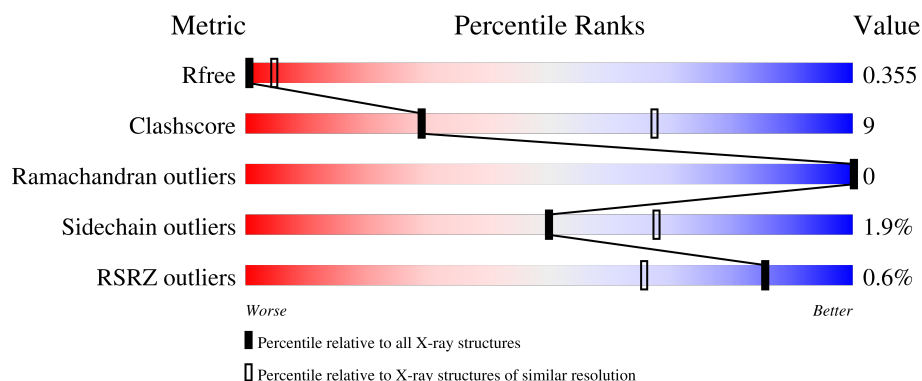
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.50 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	1085 (3.54-3.46)
Clashscore	190562	1140 (3.54-3.46)
Ramachandran outliers	187476	1113 (3.54-3.46)
Sidechain outliers	187428	1114 (3.54-3.46)
RSRZ outliers	180081	1084 (3.54-3.46)

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	379	 81% 17% ..
1	B	379	 78% 21% ..
2	E	31	 48% 52%
3	J	66	 62% 38%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	A1CW4	B	401	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6578 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 Protein fem-1 homolog B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	377	Total	C	N	O	S	0	0	0
			2949	1855	535	544	15			
1	B	376	Total	C	N	O	S	0	0	0
			2944	1852	534	543	15			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	ALA	-	expression tag	UNP Q9Z2G0
A	0	ALA	-	expression tag	UNP Q9Z2G0
B	-1	ALA	-	expression tag	UNP Q9Z2G0
B	0	ALA	-	expression tag	UNP Q9Z2G0

- Molecule 2 is a protein called Folliculin-interacting protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	15	Total	C	N	O	S	0	0	0
			115	69	23	20	3			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	1	GLY	-	expression tag	UNP Q68FD7

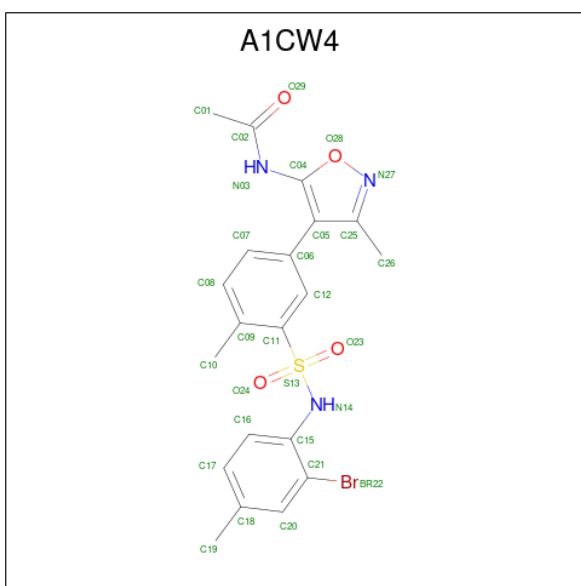
- Molecule 3 is a protein called Mitochondrial import receptor subunit TOM20 homolog.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	J	66	Total	C	N	O	S	0	0	0
			517	334	82	99	2			

- Molecule 4 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total Zn 1 1	0	0
4	B	2	Total Zn 2 2	0	0
4	E	1	Total Zn 1 1	0	0

- Molecule 5 is N-[(4M)-4-{3-[(2-bromo-4-methylphenyl)sulfamoyl]-4-methylphenyl}-3-methyl-1,2-oxazol-5-yl]acetamide (CCD ID: A1CW4) (formula: C₂₀H₂₀BrN₃O₄S) (labeled as "Ligand of Interest" by depositor).

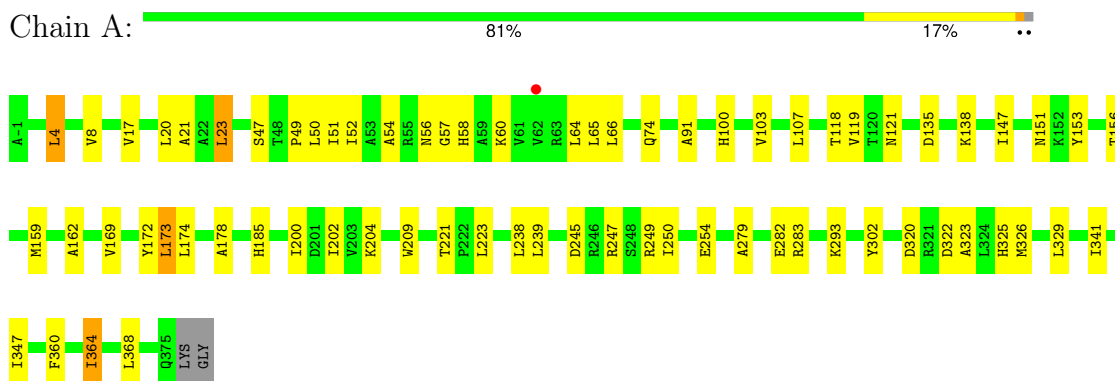


Mol	Chain	Residues	Atoms							ZeroOcc	AltConf
5	B	1	Total	Br	C	H	N	O	S	0	0
			49	1	20	20	3	4	1		

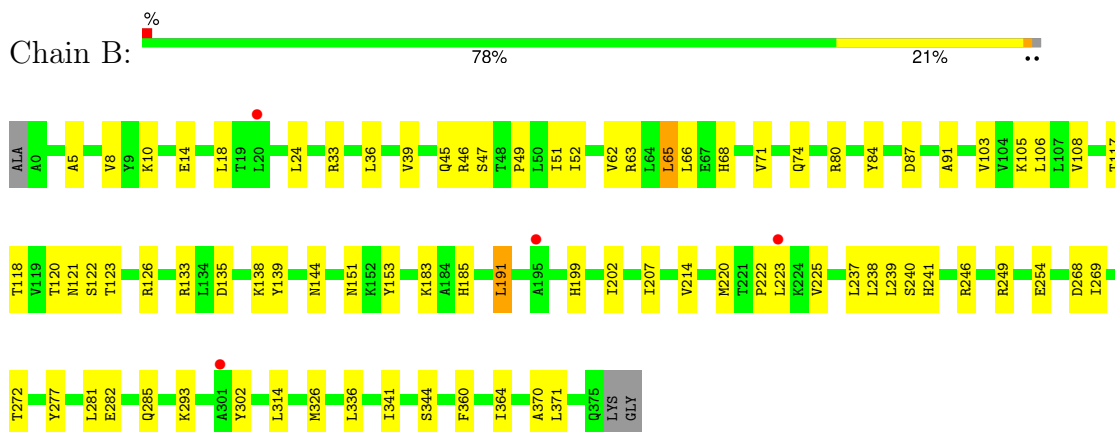
3 Residue-property plots [i](#)

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.

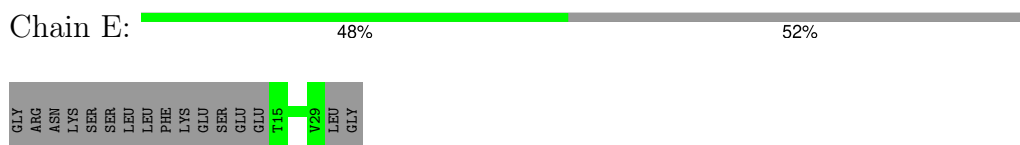
• Molecule 1: Protein fem-1 homolog B



• Molecule 1: Protein fem-1 homolog B



• Molecule 2: Folliculin-interacting protein 1



• Molecule 3: Mitochondrial import receptor subunit TOM20 homolog



D62	F69	E72	E73	L76	G77	E78	E79	L80	L81	A82	Q83	G89	H92	L93	T94	N95	A96	P103	Q104	Q105	L106	L107	L110	T113	L114	P115	P116	Q120	P121	L122	L123	T124	K125	L126	P127
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4 Data and refinement statistics

Property	Value	Source
Space group	P 41 3 2	Depositor
Cell constants a, b, c, α , β , γ	185.01Å 185.01Å 185.01Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.45 – 3.50 49.45 – 3.50	Depositor EDS
% Data completeness (in resolution range)	99.9 (49.45-3.50) 99.9 (49.45-3.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.34 (at 3.48Å)	Xtriage
Refinement program	PHENIX 1.20.1-4487	Depositor
R, R_{free}	0.300 , 0.333 0.322 , 0.355	Depositor DCC
R_{free} test set	690 reflections (4.85%)	wwPDB-VP
Wilson B-factor (Å ²)	117.2	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 105.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.42$, $\langle L^2 \rangle = 0.24$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	6578	wwPDB-VP
Average B, all atoms (Å ²)	132.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.02% 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: ZN, A1CW4

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	0.11	0/3003	0.32	0/4070
1	B	0.13	0/2998	0.34	0/4063
2	E	0.11	0/118	0.28	0/160
3	J	0.14	0/526	0.34	0/715
All	All	0.12	0/6645	0.33	0/9008

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	2949	0	2939	42	0
1	B	2944	0	2934	62	0
2	E	115	0	102	0	0
3	J	517	0	523	15	0
4	A	1	0	0	0	0
4	B	2	0	0	0	0
4	E	1	0	0	0	0
5	B	29	20	0	11	0
All	All	6558	20	6498	113	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 9.

The worst 5 of 113 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:126:ARG:HH21	5:B:401:A1CW4:C07	1.68	1.03
1:B:151:ASN:ND2	5:B:401:A1CW4:C26	2.21	1.01
1:B:151:ASN:HD22	5:B:401:A1CW4:C26	1.76	0.95
1:A:254:GLU:OE2	1:A:293:LYS:NZ	2.09	0.86
1:B:126:ARG:NH2	5:B:401:A1CW4:C07	2.45	0.78

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	375/379 (99%)	364 (97%)	11 (3%)	0	100	100
1	B	374/379 (99%)	361 (96%)	13 (4%)	0	100	100
2	E	13/31 (42%)	13 (100%)	0	0	100	100
3	J	64/66 (97%)	61 (95%)	3 (5%)	0	100	100
All	All	826/855 (97%)	799 (97%)	27 (3%)	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	305/306 (100%)	298 (98%)	7 (2%)	44	64
1	B	305/306 (100%)	300 (98%)	5 (2%)	55	70
2	E	14/29 (48%)	14 (100%)	0	100	100
3	J	57/57 (100%)	56 (98%)	1 (2%)	51	69
All	All	681/698 (98%)	668 (98%)	13 (2%)	50	68

5 of 13 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	65	LEU
1	B	66	LEU
3	J	72	GLU
1	B	144	ASN
1	B	191	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	110	HIS
1	A	166	HIS
1	B	121	ASN
1	B	155	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](#)

Of 5 ligands modelled in this entry, 4 are monoatomic - leaving 1 for Mogul analysis.

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	$\# Z > 2$
5	A1CW4	B	401	-	30,31,31	1.96	6 (20%)	40,46,46	2.57	18 (45%)

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
5	A1CW4	B	401	-	-	5/19/19/19	0/3/3/3

The worst 5 of 6 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	401	A1CW4	C06-C05	-5.26	1.39	1.49
5	B	401	A1CW4	S13-N14	5.11	1.71	1.63
5	B	401	A1CW4	O28-N27	-3.46	1.36	1.42
5	B	401	A1CW4	C11-S13	-3.40	1.72	1.77
5	B	401	A1CW4	C02-N03	2.93	1.43	1.37

The worst 5 of 18 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	401	A1CW4	C16-C15-C21	-5.50	111.15	117.58
5	B	401	A1CW4	C08-C09-C11	5.21	121.76	116.27
5	B	401	A1CW4	C12-C11-C09	-5.06	114.40	121.63
5	B	401	A1CW4	BR22-C21-C20	-4.05	110.89	118.28
5	B	401	A1CW4	BR22-C21-C15	3.79	123.07	119.82

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	401	A1CW4	C15-N14-S13-C11

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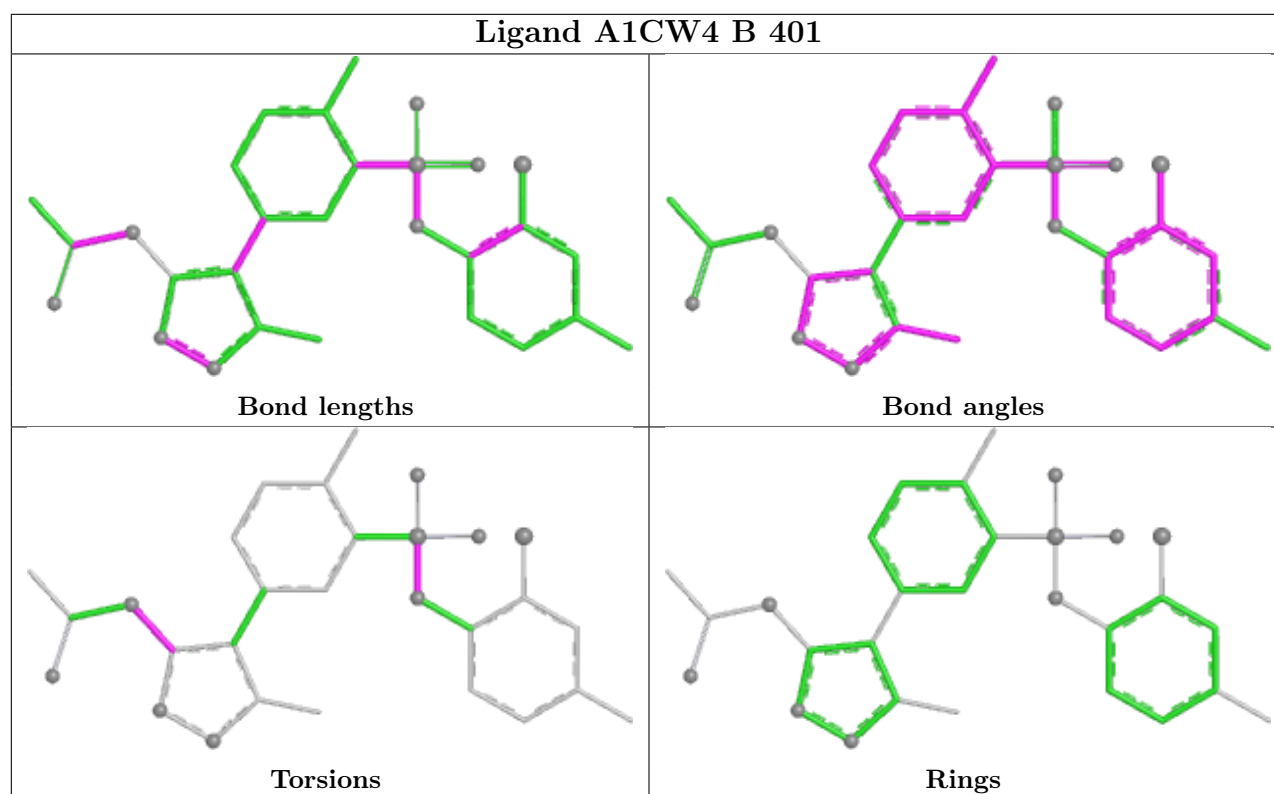
Mol	Chain	Res	Type	Atoms
5	B	401	A1CW4	C15-N14-S13-O23
5	B	401	A1CW4	C15-N14-S13-O24
5	B	401	A1CW4	O28-C04-N03-C02
5	B	401	A1CW4	C05-C04-N03-C02

There are no ring outliers.

1 monomer is involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	401	A1CW4	11	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 [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	377/379 (99%)	-0.11	1 (0%) 90 76	80, 133, 179, 205	0
1	B	376/379 (99%)	-0.10	4 (1%) 78 54	84, 118, 156, 178	0
2	E	15/31 (48%)	-0.21	0 100 100	128, 145, 157, 160	0
3	J	66/66 (100%)	-0.11	0 100 100	172, 201, 216, 222	0
All	All	834/855 (97%)	-0.11	5 (0%) 85 65	80, 129, 199, 222	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	223	LEU	3.7
1	B	20	LEU	3.0
1	A	62	VAL	2.8
1	B	301	ALA	2.3
1	B	195	ALA	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,

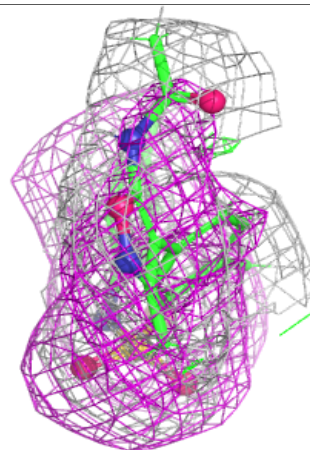
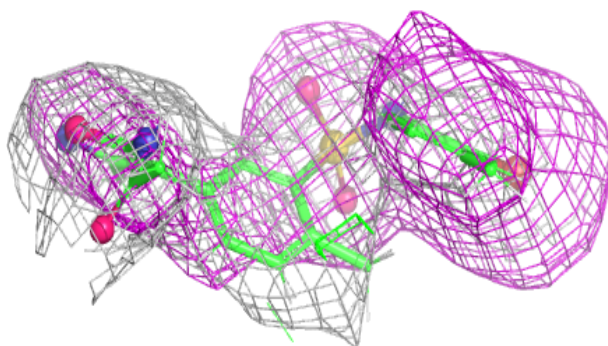
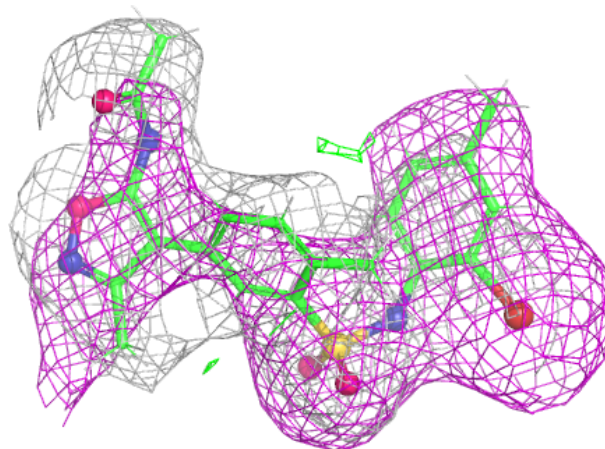
median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	ZN	B	402	1/1	0.72	0.06	261,261,261,261	0
5	A1CW4	B	401	29/29	0.76	0.15	20,20,20,20	0
4	ZN	B	403	1/1	0.80	0.06	241,241,241,241	0
4	ZN	E	601	1/1	0.98	0.04	156,156,156,156	0
4	ZN	A	401	1/1	0.98	0.04	121,121,121,121	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 A1CW4 B 401:

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 ⓘ

There are no such residues in this entry.