



Full wwPDB NMR Structure Validation Report ⓘ

Mar 8, 2026 – 12:36 PM UTC

PDB ID : 2022 / pdb_00002o22
Title : Solution structure of the anti-apoptotic protein Bcl-2 in complex with an acyl-sulfonamide-based ligand
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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)
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
wwPDB-RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
wwPDB-ShiftChecker	:	v1.2
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

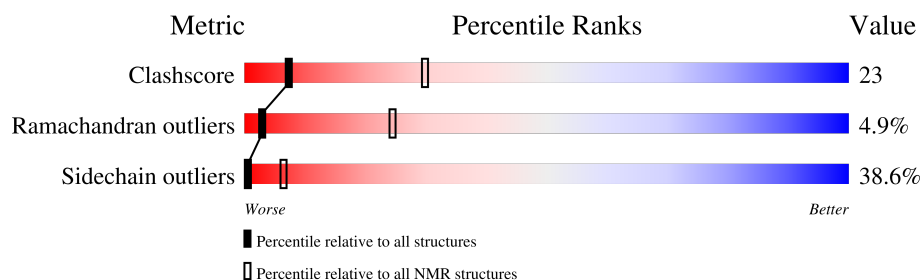
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment was not calculated.

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)	NMR archive (#Entries)
Clashscore	229148	14424
Ramachandran outliers	224038	12848
Sidechain outliers	223484	12823

The table below summarises the geometric issues observed across the polymeric chains and their fit to the experimental data. The red, orange, yellow and green segments indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A cyan segment indicates the fraction of residues that are not part of the well-defined cores, and 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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	164	42% 44% 13% .

2 Ensemble composition and analysis ⓘ

This entry contains 1 models. Identification of well-defined residues and clustering analysis are not possible.

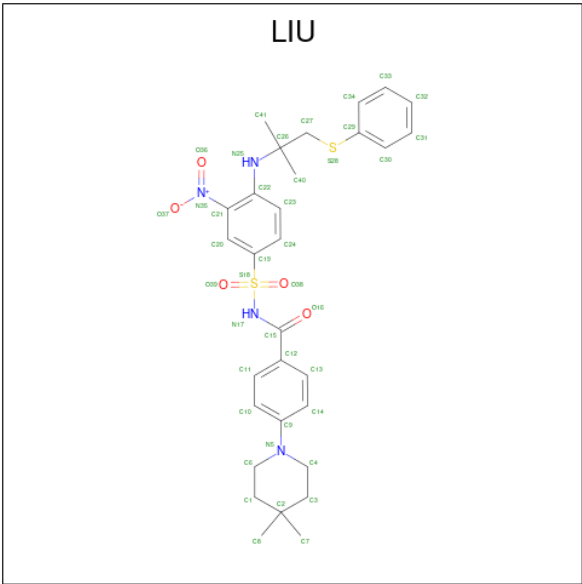
3 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 2688 atoms, of which 1297 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Apoptosis regulator Bcl-2.

Mol	Chain	Residues	Atoms						Trace
1	A	164	Total	C	H	N	O	S	0
			2611	847	1261	242	254	7	

- Molecule 2 is N-[(4-{[1,1-dimethyl-2-(phenylthio)ethyl]amino}-3-nitrophenyl)sulfonyl]-4-(4,4-dimethylpiperidin-1-yl)benzamide (CCD ID: LIU) (formula: C₃₀H₃₆N₄O₅S₂).

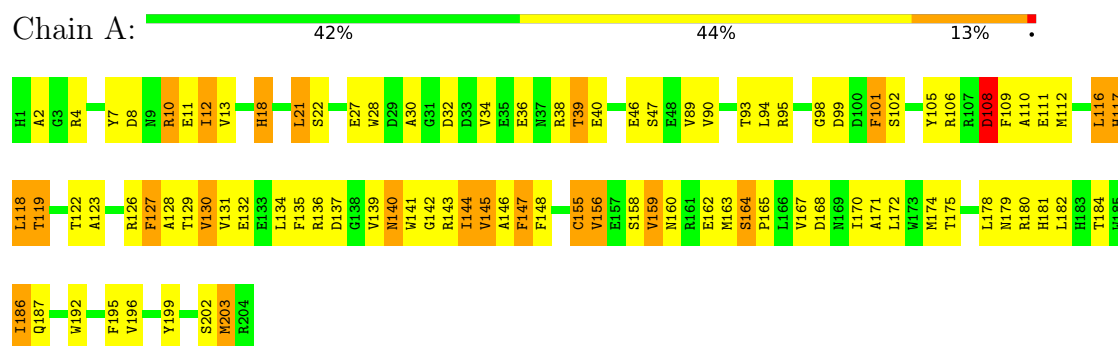


Mol	Chain	Residues	Atoms					
2	A	1	Total	C	H	N	O	S
			77	30	36	4	5	2

4 Residue-property plots

These plots are provided for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic is the same as shown in the summary in section 1 of this report. The second graphic shows the sequence where residues are colour-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. Stretches of 2 or more consecutive residues without any outliers are shown as green connectors. Residues which are classified as ill-defined in the NMR ensemble, are shown in cyan with an underline colour-coded according to the previous scheme. Residues which were present in the experimental sample, but not modelled in the final structure are shown in grey.

• Molecule 1: Apoptosis regulator Bcl-2



5 Refinement protocol and experimental data overview ⓘ

The models were refined using the following method: *simulated annealing*.

Of the ? calculated structures, 1 were deposited, based on the following criterion: ?.

The following table shows the software used for structure solution, optimisation and refinement.

Software name	Classification	Version
CNS	refinement	

No chemical shift data was provided.

6 Model quality ⓘ

6.1 Standard geometry ⓘ

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

There are no covalent bond-length or bond-angle outliers.

There are no bond-length outliers.

There are no bond-angle outliers.

There are no chirality outliers.

There are no planarity outliers.

6.2 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 each 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 averaged over the ensemble.

Mol	Chain	Non-H	H(model)	H(added)	Clashes
1	A	1350	1261	1261	61
2	A	41	36	36	9
All	All	1391	1297	1297	63

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 23.

All clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)
1:A:141:TRP:O	1:A:145:VAL:HG12	0.80	1.76
1:A:145:VAL:HG11	2:A:1000:LIU:H33	0.73	1.61
1:A:28:TRP:CE2	1:A:30:ALA:HB2	0.70	2.21
1:A:171:ALA:O	1:A:175:THR:HG23	0.67	1.88
1:A:159:VAL:HG12	1:A:164:SER:HA	0.66	1.68
1:A:156:VAL:O	1:A:159:VAL:HG22	0.65	1.91
1:A:119:THR:O	1:A:123:ALA:HB2	0.64	1.93
1:A:10:ARG:HA	1:A:175:THR:HG21	0.62	1.71
1:A:186:ILE:HD12	1:A:192:TRP:CD1	0.62	2.30
1:A:182:LEU:O	1:A:186:ILE:HG23	0.61	1.95
1:A:118:LEU:O	1:A:119:THR:HG23	0.59	1.97
1:A:164:SER:O	1:A:167:VAL:HG13	0.55	2.02

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Atom-1	Atom-2	Clash(Å)	Distance(Å)
1:A:199:TYR:CD1	2:A:1000:LIU:H402	0.54	2.38
1:A:145:VAL:HG11	2:A:1000:LIU:C33	0.53	2.33
1:A:135:PHE:CD2	1:A:181:HIS:CD2	0.53	2.96
1:A:142:GLY:O	1:A:145:VAL:HG13	0.52	2.05
1:A:116:LEU:HD23	1:A:117:HIS:N	0.52	2.19
1:A:21:LEU:HD22	1:A:28:TRP:HB2	0.52	1.82
1:A:126:ARG:O	1:A:130:VAL:HG13	0.52	2.04
1:A:135:PHE:CG	1:A:135:PHE:O	0.52	2.63
1:A:39:THR:HG22	1:A:40:GLU:HG3	0.51	1.81
1:A:98:GLY:HA2	1:A:145:VAL:HG23	0.50	1.84
1:A:105:TYR:CE2	2:A:1000:LIU:H10	0.50	2.42
1:A:186:ILE:HD12	1:A:192:TRP:HD1	0.50	1.65
1:A:164:SER:O	1:A:167:VAL:HG22	0.49	2.07
1:A:178:LEU:C	1:A:178:LEU:HD23	0.49	2.32
1:A:18:HIS:CD2	1:A:28:TRP:CZ3	0.49	3.00
1:A:135:PHE:CZ	1:A:181:HIS:CE1	0.49	3.00
1:A:28:TRP:CD2	1:A:30:ALA:HB2	0.49	2.43
1:A:123:ALA:HA	1:A:126:ARG:HG2	0.48	1.84
1:A:135:PHE:CD1	1:A:181:HIS:NE2	0.47	2.83
1:A:127:PHE:CD2	1:A:128:ALA:N	0.46	2.83
1:A:164:SER:N	1:A:165:PRO:CD	0.46	2.79
1:A:101:PHE:CD2	1:A:102:SER:N	0.46	2.83
1:A:140:ASN:O	1:A:144:ILE:HD12	0.46	2.10
2:A:1000:LIU:H412	2:A:1000:LIU:H23	0.46	1.87
1:A:108:ASP:O	1:A:112:MET:HE3	0.46	2.11
1:A:109:PHE:CD1	1:A:110:ALA:N	0.46	2.84
1:A:34:VAL:O	1:A:34:VAL:HG23	0.45	2.10
1:A:186:ILE:CG1	1:A:187:GLN:N	0.45	2.79
1:A:93:THR:HG22	1:A:195:PHE:CZ	0.45	2.47
1:A:148:PHE:CZ	1:A:178:LEU:CD1	0.45	3.00
1:A:135:PHE:CE1	1:A:181:HIS:CE1	0.44	3.05
1:A:7:TYR:CB	1:A:192:TRP:CE3	0.43	3.02
1:A:181:HIS:CD2	1:A:182:LEU:HD12	0.43	2.48
1:A:146:ALA:HB1	2:A:1000:LIU:C3	0.43	2.43
1:A:147:PHE:HE1	1:A:178:LEU:HD12	0.43	1.72
1:A:108:ASP:OD2	1:A:109:PHE:CD2	0.43	2.72
1:A:147:PHE:CD1	1:A:148:PHE:N	0.43	2.87
1:A:155:CYS:O	1:A:167:VAL:HG12	0.42	2.14
2:A:1000:LIU:H11	2:A:1000:LIU:H31	0.42	1.92
1:A:127:PHE:CG	1:A:128:ALA:N	0.42	2.85
1:A:46:GLU:O	1:A:89:VAL:HG22	0.42	2.15
1:A:135:PHE:CG	1:A:181:HIS:NE2	0.41	2.88

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Atom-1	Atom-2	Clash(Å)	Distance(Å)
1:A:142:GLY:HA3	2:A:1000:LIU:N17	0.41	2.30
1:A:28:TRP:CZ2	1:A:30:ALA:HB2	0.41	2.50
1:A:28:TRP:CD1	1:A:28:TRP:C	0.41	2.98
1:A:178:LEU:O	1:A:182:LEU:HB2	0.41	2.16
1:A:12:ILE:CG2	1:A:13:VAL:N	0.41	2.83
1:A:135:PHE:CD2	1:A:135:PHE:O	0.40	2.74
1:A:7:TYR:HB2	1:A:192:TRP:CE3	0.40	2.51
1:A:130:VAL:CG2	1:A:131:VAL:N	0.40	2.84
1:A:142:GLY:HA3	2:A:1000:LIU:HN17	0.40	1.77

6.3 Torsion angles [i](#)

6.3.1 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 PDB entries followed by that with respect to all NMR entries. 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	162/164 (99%)	135 (83%)	19 (12%)	8 (5%)	3	24
All	All	162/164 (99%)	135 (83%)	19 (12%)	8 (5%)	3	24

All 8 Ramachandran outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type
1	A	2	ALA
1	A	27	GLU
1	A	36	GLU
1	A	39	THR
1	A	108	ASP
1	A	162	GLU
1	A	202	SER
1	A	203	MET

6.3.2 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 PDB entries followed by that with respect to all NMR entries. 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	140/140 (100%)	86 (61%)	54 (39%)	0 6
All	All	140/140 (100%)	86 (61%)	54 (39%)	0 6

All 54 residues with a non-rotameric sidechain are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type
1	A	4	ARG
1	A	8	ASP
1	A	10	ARG
1	A	11	GLU
1	A	12	ILE
1	A	18	HIS
1	A	21	LEU
1	A	22	SER
1	A	32	ASP
1	A	38	ARG
1	A	47	SER
1	A	90	VAL
1	A	94	LEU
1	A	95	ARG
1	A	99	ASP
1	A	101	PHE
1	A	106	ARG
1	A	108	ASP
1	A	111	GLU
1	A	116	LEU
1	A	117	HIS
1	A	118	LEU
1	A	119	THR
1	A	122	THR
1	A	127	PHE
1	A	129	THR
1	A	130	VAL
1	A	132	GLU
1	A	134	LEU
1	A	136	ARG
1	A	137	ASP
1	A	139	VAL
1	A	140	ASN
1	A	143	ARG

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Mol	Chain	Res	Type
1	A	144	ILE
1	A	145	VAL
1	A	147	PHE
1	A	155	CYS
1	A	156	VAL
1	A	158	SER
1	A	159	VAL
1	A	160	ASN
1	A	163	MET
1	A	164	SER
1	A	168	ASP
1	A	170	ILE
1	A	172	LEU
1	A	174	MET
1	A	179	ASN
1	A	180	ARG
1	A	184	THR
1	A	186	ILE
1	A	196	VAL
1	A	203	MET

6.3.3 RNA [i](#)

There are no RNA molecules in this entry.

6.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.6 Ligand geometry [i](#)

1 ligand is modelled in this entry.

In the following table, the Counts columns list the number of bonds for which Mogul statistics could be retrieved, the number of bonds that are observed in the model and the number of bonds 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 is the number of standard

deviations the observed value is removed from the expected value. A bond length with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the average root-mean-square of all Z scores of the bond lengths.

Mol	Type	Chain	Res	Link	Bond lengths		
					Counts	RMSZ	#Z>2
2	LIU	A	1000	-	42,44,44	3.42	9 (21%)

In the following table, the Counts columns list the number of angles for which Mogul statistics could be retrieved, the number of angles that are observed in the model and the number of 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 angle is the number of standard deviations the observed value is removed from the expected value. A bond angle with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the average root-mean-square of all Z scores of the bond angles.

Mol	Type	Chain	Res	Link	Bond angles		
					Counts	RMSZ	#Z>2
2	LIU	A	1000	-	57,65,65	1.90	13 (22%)

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	LIU	A	1000	-	-	0,32,46,46	0,4,4,4

All bond outliers are listed below. They are sorted according to the Z-score.

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1000	LIU	O38-S18	14.07	1.59	1.43
2	A	1000	LIU	O39-S18	13.57	1.59	1.43
2	A	1000	LIU	C15-N17	4.39	1.33	1.39
2	A	1000	LIU	C22-N25	3.92	1.49	1.39
2	A	1000	LIU	C9-N5	3.91	1.49	1.38
2	A	1000	LIU	C21-N35	3.73	1.52	1.45
2	A	1000	LIU	S18-N17	2.74	1.70	1.64
2	A	1000	LIU	C6-N5	2.30	1.50	1.46
2	A	1000	LIU	C4-N5	2.22	1.50	1.46

All angle outliers are listed below. They are sorted according to the Z-score.

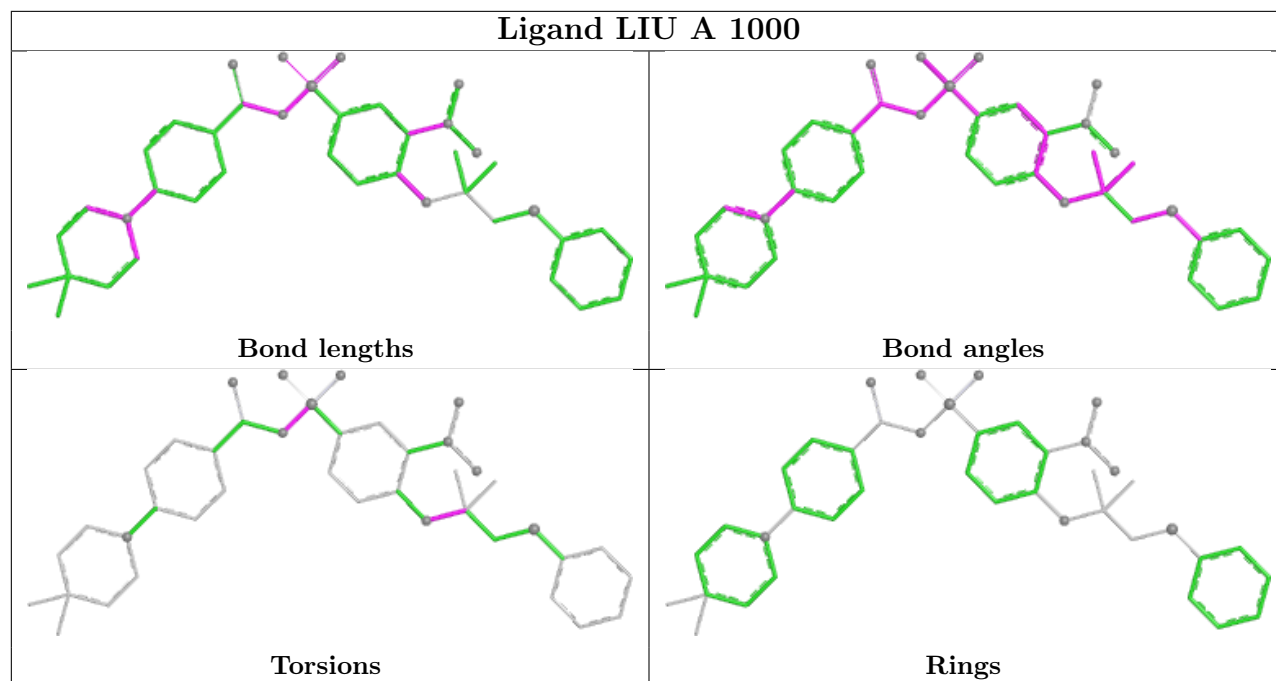
Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1000	LIU	O39-S18-O38	6.68	111.40	119.52
2	A	1000	LIU	C12-C15-N17	5.09	122.14	116.08
2	A	1000	LIU	O16-C15-N17	4.38	115.78	121.16
2	A	1000	LIU	C21-C22-N25	3.85	119.51	123.39
2	A	1000	LIU	O38-S18-C19	3.47	112.36	107.98
2	A	1000	LIU	C19-S18-N17	3.25	111.25	106.06
2	A	1000	LIU	C27-S28-C29	3.23	109.67	103.04
2	A	1000	LIU	C20-C21-C22	2.94	118.87	121.55
2	A	1000	LIU	C41-C26-C40	2.68	105.48	109.55
2	A	1000	LIU	O39-S18-C19	2.67	111.35	107.98
2	A	1000	LIU	C41-C26-N25	2.38	113.26	107.60
2	A	1000	LIU	C4-N5-C9	2.31	124.41	118.11
2	A	1000	LIU	C26-N25-C22	2.30	137.78	125.30

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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.



6.7 Other polymers [i](#)

There are no such molecules in this entry.

6.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

7 Chemical shift validation

No chemical shift data were provided