



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 03:44 PM UTC

PDB ID : 1G73 / pdb_00001g73
Title : CRYSTAL STRUCTURE OF SMAC BOUND TO XIAP-BIR3 DOMAIN
Authors : Wu, G.; Chai, J.; Suber, T.L.; Wu, J.-W.; Shi, Y.
Deposited on : 2000-11-08
Resolution : 2.00 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
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

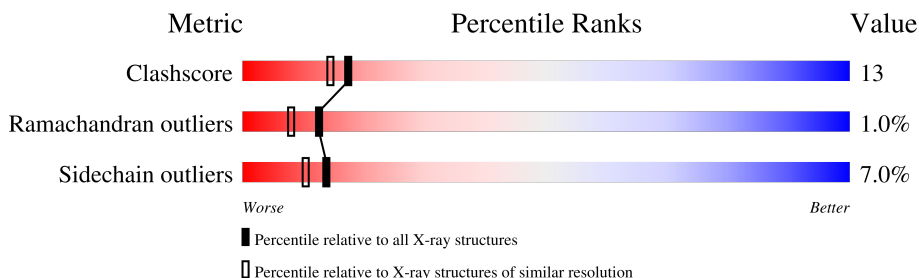
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	162	 67% 25% 5% . .
1	B	162	 71% 23% . .
2	C	121	 50% 21% . 26%
2	D	121	 60% 18% 7% 16%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4010 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 SECOND MITOCHONDRIA-DERIVED ACTIVATOR OF CASPASES.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	157	Total	C	N	O	S	0	0	0
			1223	761	206	251	5			
1	B	157	Total	C	N	O	S	0	0	0
			1223	761	206	251	5			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	33	ASP	PHE	engineered mutation	UNP Q9NR28
B	33	ASP	PHE	engineered mutation	UNP Q9NR28

- Molecule 2 is a protein called INHIBITORS OF APOPTOSIS-LIKE PROTEIN ILP.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	89	Total	C	N	O	S	0	0	0
			729	471	123	131	4			
2	D	102	Total	C	N	O	S	0	0	0
			833	533	141	154	5			

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Zn	0	0
			1	1		
3	D	1	Total	Zn	0	0
			1	1		

N341	L342	H343	L344	T345	H346	T356	E357	LYS
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4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	48.10Å 52.90Å 67.10Å 100.00° 104.10° 94.00°	Depositor
Resolution (Å)	20.00 – 2.00	Depositor
% Data completeness (in resolution range)	94.8 (20.00-2.00)	Depositor
R_{merge}	0.04	Depositor
R_{sym}	0.04	Depositor
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.227 , 0.268	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4010	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	0.64	0/1239	0.98	3/1680 (0.2%)
1	B	0.63	0/1239	1.03	2/1680 (0.1%)
2	C	0.64	0/754	0.93	2/1021 (0.2%)
2	D	0.62	0/859	1.12	7/1164 (0.6%)
All	All	0.63	0/4091	1.02	14/5545 (0.3%)

There are no bond length outliers.

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	344	LEU	N-CA-C	-10.11	94.69	110.36
2	D	311	LYS	N-CA-C	-7.28	99.14	110.14
2	D	320	HIS	N-CA-C	-6.43	104.19	111.07
2	D	279	VAL	N-CA-C	-6.10	100.70	108.93
2	D	346	HIS	N-CA-C	6.02	119.45	111.75
1	B	78	TRP	N-CA-C	5.99	117.89	111.36
1	A	54	THR	N-CA-C	-5.76	104.92	111.14
1	A	11	HIS	N-CA-C	-5.58	103.75	111.28
2	C	329	TYR	N-CA-C	-5.52	104.90	111.69
2	D	356	THR	N-CA-C	-5.47	106.66	113.38
2	D	323	TRP	N-CA-C	5.31	118.92	112.23
2	C	336	GLN	N-CA-C	5.15	116.98	111.36
1	A	77	VAL	N-CA-C	-5.12	105.72	110.53
1	B	148	LYS	N-CA-C	-5.06	105.65	111.07

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	1223	0	1221	38	0
1	B	1223	0	1221	24	0
2	C	729	0	682	22	0
2	D	833	0	782	22	0
3	B	1	0	0	0	0
3	D	1	0	0	0	0
All	All	4010	0	3906	99	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 13.

All (99) 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:6:GLN:O	1:A:7:LYS:HB2	1.70	0.88
1:A:114:GLU:O	1:A:118:GLN:HG2	1.79	0.82
1:A:138:GLN:O	1:A:142:VAL:HG13	1.86	0.76
1:A:4:ILE:HG12	2:D:308:THR:OG1	1.86	0.75
1:A:123:GLN:HG2	1:A:124:ALA:N	2.01	0.75
1:A:73:GLU:O	1:A:77:VAL:HG23	1.90	0.71
2:D:266:GLU:HA	2:D:269:ILE:HD11	1.71	0.70
2:C:321:ALA:HA	2:C:330:LEU:HD21	1.76	0.67
2:D:311:LYS:HB2	2:D:314:GLU:HG3	1.76	0.67
2:C:308:THR:HG22	2:C:309:ASP:H	1.61	0.66
2:D:318:GLU:HG2	2:D:342:ILE:HD12	1.79	0.64
2:C:297:LYS:HG2	2:C:308:THR:HG23	1.82	0.62
1:A:2:VAL:HG13	1:A:3:PRO:HD2	1.83	0.61
2:C:262:MET:HE1	2:C:271:THR:HG21	1.82	0.60
2:D:344:LEU:O	2:D:345:THR:HB	2.01	0.59
2:C:294:GLU:HG3	2:C:297:LYS:HD2	1.86	0.58
2:D:258:ARG:HD3	2:D:304:GLY:HA3	1.86	0.58
1:B:4:ILE:HG12	2:C:308:THR:OG1	2.04	0.57
1:B:44:ALA:HB2	1:B:98:LEU:HD12	1.87	0.57
2:C:262:MET:HE1	2:C:271:THR:CG2	2.35	0.57
1:B:95:TYR:C	1:B:95:TYR:CD1	2.82	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:95:TYR:C	1:A:95:TYR:CD1	2.83	0.56
1:A:23:VAL:O	1:A:27:THR:HG23	2.06	0.56
1:B:42:ILE:HG23	1:B:142:VAL:HG11	1.88	0.55
2:C:272:PHE:HB3	2:C:281:LYS:HD2	1.88	0.55
1:A:123:GLN:CG	1:A:124:ALA:N	2.68	0.54
2:D:342:ILE:O	2:D:346:HIS:HB3	2.07	0.54
2:D:265:TYR:O	2:D:269:ILE:HG13	2.07	0.54
1:B:48:TYR:O	1:B:52:VAL:HG23	2.07	0.54
2:D:356:THR:O	2:D:357:GLU:HB2	2.09	0.53
1:B:118:GLN:HA	1:B:118:GLN:OE1	2.09	0.53
2:D:308:THR:HG22	2:D:309:ASP:H	1.74	0.52
1:B:89:THR:O	1:B:93:GLN:HG3	2.09	0.52
2:D:308:THR:HG22	2:D:309:ASP:N	2.25	0.51
1:A:7:LYS:HA	1:A:7:LYS:HE2	1.93	0.50
1:A:99:GLU:OE2	2:C:260:PRO:HD2	2.12	0.50
1:A:103:MET:HE2	2:C:260:PRO:HG3	1.93	0.50
1:A:55:LEU:HD11	1:A:81:ILE:HG23	1.94	0.50
1:B:123:GLN:HG2	1:B:124:ALA:N	2.27	0.49
1:B:69:MET:HG3	1:B:74:GLU:HB2	1.94	0.49
1:A:21:ARG:NE	1:A:21:ARG:HA	2.27	0.49
2:D:321:ALA:HA	2:D:330:LEU:HD21	1.94	0.48
1:A:38:THR:HG21	1:A:135:VAL:HG22	1.96	0.48
2:D:258:ARG:HH11	2:D:304:GLY:HA3	1.78	0.48
1:A:85:ARG:HG2	1:A:85:ARG:HH11	1.79	0.47
1:A:19:MET:CE	1:B:26:VAL:HG11	2.43	0.47
1:A:153:LEU:O	1:A:157:GLN:HG2	2.14	0.47
1:B:58:LEU:HD21	1:B:80:VAL:HG12	1.96	0.47
2:D:280:ASN:OD1	2:D:282:GLU:HB2	2.15	0.47
2:D:344:LEU:O	2:D:345:THR:CB	2.63	0.47
1:B:20:ARG:CG	1:B:20:ARG:HH11	2.28	0.47
1:A:6:GLN:O	1:A:7:LYS:CB	2.55	0.47
1:A:103:MET:HE2	2:C:260:PRO:CG	2.45	0.47
1:A:95:TYR:HD1	1:A:95:TYR:O	1.98	0.46
1:B:123:GLN:H	1:B:123:GLN:CD	2.24	0.46
1:A:6:GLN:H	1:A:6:GLN:NE2	2.13	0.46
2:C:321:ALA:O	2:C:343:HIS:HE1	1.99	0.46
1:B:23:VAL:O	1:B:27:THR:HG23	2.16	0.46
2:D:284:LEU:HD23	2:D:316:PRO:HB3	1.98	0.46
2:C:292:LEU:HD12	2:C:297:LYS:HB3	1.97	0.46
1:A:76:GLU:O	1:A:80:VAL:HG23	2.16	0.45
2:C:264:ASP:O	2:C:268:ARG:HG3	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:137:LEU:O	1:A:141:GLU:HG3	2.16	0.45
1:B:95:TYR:C	1:B:95:TYR:HD1	2.23	0.45
2:C:294:GLU:HG3	2:C:294:GLU:O	2.17	0.45
1:A:75:ASP:O	1:A:79:GLN:HG2	2.17	0.44
1:B:110:GLU:HG2	1:B:132:ILE:HD13	1.99	0.44
2:C:328:LYS:HE3	2:C:331:LEU:HD21	2.00	0.44
1:A:65:LEU:HB3	1:A:69:MET:HG3	1.98	0.44
2:D:341:ASN:O	2:D:344:LEU:O	2.36	0.44
2:D:342:ILE:C	2:D:344:LEU:O	2.60	0.44
2:D:287:ALA:O	2:D:320:HIS:CE1	2.71	0.44
2:D:342:ILE:O	2:D:344:LEU:O	2.36	0.44
1:A:72:GLU:HG2	1:A:73:GLU:N	2.32	0.43
1:B:23:VAL:HG11	1:B:116:ALA:HA	2.00	0.43
1:A:99:GLU:O	1:A:103:MET:HG2	2.19	0.43
1:A:41:LEU:O	1:A:45:ILE:HG13	2.17	0.43
2:C:327:CYS:SG	2:C:329:TYR:HB3	2.59	0.43
1:A:102:TRP:O	1:A:106:VAL:HG23	2.18	0.43
1:A:2:VAL:HG13	1:A:3:PRO:CD	2.48	0.42
2:C:262:MET:CE	2:C:271:THR:HG21	2.46	0.42
1:B:122:ASP:O	1:B:125:SER:HB3	2.20	0.42
1:B:21:ARG:HD3	1:B:21:ARG:HA	1.74	0.42
2:D:345:THR:HG22	2:D:346:HIS:N	2.33	0.42
1:B:59:TYR:HE1	1:B:157:GLN:HE22	1.66	0.42
2:C:334:LYS:HA	2:C:334:LYS:HD2	1.95	0.42
1:B:2:VAL:HG13	1:B:3:PRO:HD2	2.02	0.41
2:C:321:ALA:HA	2:C:330:LEU:CD2	2.48	0.41
1:B:65:LEU:O	1:B:69:MET:HE2	2.20	0.41
1:B:66:LEU:HA	1:B:69:MET:HE2	2.02	0.41
1:A:6:GLN:HB2	1:A:7:LYS:H	1.47	0.41
1:A:97:LYS:HE2	1:A:97:LYS:HB3	1.77	0.41
2:C:338:TYR:O	2:C:342:ILE:HG12	2.21	0.41
2:C:307:LEU:HD11	2:C:323:TRP:CD1	2.56	0.41
1:A:4:ILE:HG12	2:D:308:THR:HG1	1.82	0.40
1:B:58:LEU:HD12	1:B:58:LEU:HA	1.93	0.40
1:A:2:VAL:HG12	1:A:3:PRO:O	2.21	0.40
1:A:19:MET:O	1:A:23:VAL:HG23	2.22	0.40
1:A:144:GLN:O	1:A:148:LYS:HG2	2.21	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	155/162 (96%)	148 (96%)	4 (3%)	3 (2%)	6	3
1	B	155/162 (96%)	148 (96%)	6 (4%)	1 (1%)	21	17
2	C	87/121 (72%)	84 (97%)	2 (2%)	1 (1%)	11	7
2	D	100/121 (83%)	98 (98%)	2 (2%)	0	100	100
All	All	497/566 (88%)	478 (96%)	14 (3%)	5 (1%)	12	8

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	7	LYS
1	A	71	SER
1	B	6	GLN
1	A	8	SER
2	C	312	PRO

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	133/138 (96%)	123 (92%)	10 (8%)	12	9
1	B	133/138 (96%)	123 (92%)	10 (8%)	12	9
2	C	74/105 (70%)	70 (95%)	4 (5%)	20	17
2	D	87/105 (83%)	81 (93%)	6 (7%)	14	11
All	All	427/486 (88%)	397 (93%)	30 (7%)	14	10

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	GLN
1	A	7	LYS
1	A	21	ARG
1	A	58	LEU
1	A	87	GLU
1	A	95	TYR
1	A	135	VAL
1	A	137	LEU
1	A	142	VAL
1	A	157	GLN
1	B	2	VAL
1	B	53	TYR
1	B	58	LEU
1	B	87	GLU
1	B	94	GLU
1	B	95	TYR
1	B	135	VAL
1	B	137	LEU
1	B	142	VAL
1	B	145	LEU
2	C	279	VAL
2	C	307	LEU
2	C	308	THR
2	C	341	ASN
2	D	269	ILE
2	D	274	THR
2	D	279	VAL
2	D	307	LEU
2	D	328	LYS
2	D	345	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	6	GLN
1	A	93	GLN
1	A	143	HIS
1	B	36	GLN
1	B	79	GLN
1	B	93	GLN
1	B	131	HIS

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Mol	Chain	Res	Type
1	B	133	GLN
1	B	143	HIS
2	C	259	ASN
2	C	343	HIS
2	D	259	ASN
2	D	346	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.