



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 11:10 PM UTC

PDB ID : 1PZ7 / pdb_00001pz7
Title : Modulation of agrin function by alternative splicing and Ca²⁺ binding
Authors : Stetefeld, J.; Alexandrescu, A.T.; Maciejewski, M.W.; Jenny, M.; Rathgeb-Szabo, K.; Schulthess, T.; Landwehr, R.; Frank, S.; Ruegg, M.A.; Kammerer, R.A.
Deposited on : 2003-07-10
Resolution : 1.42 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

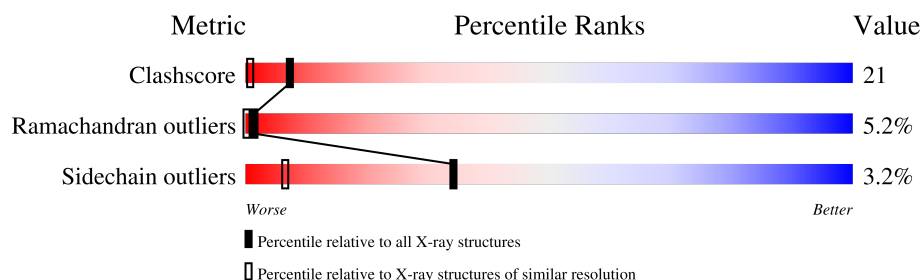
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.42 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	4154 (1.44-1.40)
Ramachandran outliers	187476	4083 (1.44-1.40)
Sidechain outliers	187428	4082 (1.44-1.40)

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	204	 66% 20% 5% • 8%
1	B	204	 73% 14% • • 8%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3375 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 Agrin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	188	Total	C	N	O	S	0	0	0
			1445	910	252	277	6			
1	B	187	Total	C	N	O	S	0	0	0
			1443	908	251	278	6			

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

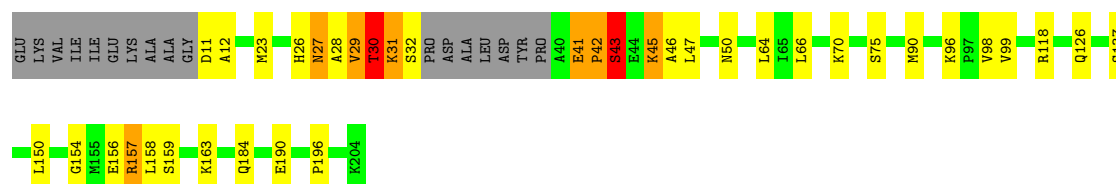
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Ca	0	0
			1	1		
2	B	1	Total	Ca	0	0
			1	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	274	Total	O	0	0
			274	274		
3	B	211	Total	O	0	0
			211	211		

Note EDS was not executed.

Chain A: 66% 20% 5% 8%



4 Data and refinement statistics

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

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	101.33Å 57.54Å 74.62Å 90.00° 127.10° 90.00°	Depositor
Resolution (Å)	20.00 – 1.42	Depositor
% Data completeness (in resolution range)	100.0 (20.00-1.42)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.04	Depositor
Refinement program	REFMAC 5.0	Depositor
R, R_{free}	0.220 , 0.240	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3375	wwPDB-VP
Average B, all atoms (Å ²)	14.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

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

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	0.43	0/1473	0.83	2/1997 (0.1%)
1	B	0.38	0/1471	0.81	2/1993 (0.1%)
All	All	0.41	0/2944	0.82	4/3990 (0.1%)

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	B	0	1

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	31	LYS	N-CA-C	-8.52	96.03	109.23
1	A	30	THR	N-CA-C	7.12	120.24	109.63
1	B	75	SER	N-CA-C	5.41	118.29	110.28
1	A	75	SER	N-CA-C	5.20	117.83	110.50

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	30	THR	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	1445	0	1427	66	0
1	B	1443	0	1424	52	1
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	274	0	0	8	1
3	B	211	0	0	9	0
All	All	3375	0	2851	118	1

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 21.

All (118) 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:29:VAL:HG21	1:A:184:GLN:HG2	1.22	1.16
1:A:30:THR:HG23	1:A:31:LYS:H	1.18	1.06
1:A:8:ALA:HB2	1:A:13:GLU:HG3	1.44	0.95
1:A:41:GLU:HB3	1:A:42:PRO:HD3	1.48	0.93
1:A:30:THR:HG23	1:A:31:LYS:N	1.79	0.90
1:A:160:VAL:HG11	3:A:815:HOH:O	1.75	0.86
1:A:29:VAL:HG11	1:A:184:GLN:HE21	1.41	0.86
1:A:29:VAL:HG21	1:A:184:GLN:CG	2.09	0.81
1:B:32:SER:O	1:B:45:LYS:HG2	1.81	0.80
1:A:30:THR:CG2	1:A:31:LYS:N	2.45	0.78
1:A:29:VAL:CG2	1:A:184:GLN:HG2	2.08	0.78
1:A:41:GLU:CB	1:A:42:PRO:HD3	2.14	0.77
1:A:30:THR:O	1:A:32:SER:N	2.17	0.77
1:B:50:ASN:HD22	1:B:118:ARG:HE	1.35	0.75
1:B:32:SER:O	1:B:47:LEU:HD21	1.86	0.74
1:A:160:VAL:C	3:A:1056:HOH:O	2.29	0.74
1:A:31:LYS:HE2	1:A:31:LYS:HA	1.71	0.71
1:B:45:LYS:CE	1:B:46:ALA:H	2.03	0.70
1:B:163:LYS:NZ	3:B:819:HOH:O	2.20	0.69
1:B:45:LYS:HE3	1:B:46:ALA:H	1.58	0.69
1:A:7:LYS:HB3	1:A:201:CYS:O	1.92	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:29:VAL:HG11	1:A:184:GLN:NE2	2.08	0.68
1:B:26:HIS:O	1:B:28:ALA:N	2.27	0.67
1:A:24:GLU:HB2	1:A:151:TRP:CZ3	2.30	0.67
1:B:29:VAL:O	1:B:30:THR:CB	2.43	0.66
1:A:50:ASN:HD22	1:A:118:ARG:HE	1.43	0.66
1:A:32:SER:CB	1:A:182:ASP:O	2.44	0.65
1:A:26:HIS:O	1:A:28:ALA:N	2.31	0.64
1:A:32:SER:HB3	1:A:182:ASP:O	1.98	0.63
1:B:41:GLU:HB3	1:B:42:PRO:HD2	1.81	0.62
1:A:30:THR:C	1:A:32:SER:H	2.08	0.62
1:A:7:LYS:HD3	1:A:9:ALA:HB3	1.81	0.61
1:B:45:LYS:CE	1:B:46:ALA:N	2.63	0.61
1:A:121:ARG:NH2	3:A:888:HOH:O	2.28	0.61
1:B:28:ALA:O	1:B:29:VAL:HG23	2.01	0.60
1:A:40:ALA:O	1:A:41:GLU:HB3	2.01	0.60
1:B:27:ASN:HB3	1:B:32:SER:HB3	1.84	0.60
1:A:32:SER:HB2	1:A:182:ASP:HB3	1.84	0.59
1:B:41:GLU:CB	1:B:42:PRO:HD2	2.32	0.59
1:A:8:ALA:HB3	1:A:200:HIS:ND1	2.18	0.58
1:A:31:LYS:HA	1:A:31:LYS:CE	2.32	0.58
1:B:90:MET:HG2	1:B:99:VAL:HG22	1.84	0.58
1:A:27:ASN:O	1:A:27:ASN:CG	2.46	0.57
1:B:31:LYS:O	1:B:32:SER:CB	2.52	0.57
1:A:22:TYR:HD2	1:A:151:TRP:CE3	2.23	0.57
1:B:29:VAL:O	1:B:30:THR:OG1	2.21	0.57
1:B:70:LYS:O	1:B:157:ARG:NH1	2.39	0.56
1:A:29:VAL:HG11	1:A:184:GLN:HB2	1.88	0.55
1:B:29:VAL:HG21	1:B:184:GLN:HG2	1.88	0.55
1:B:31:LYS:O	1:B:32:SER:HB3	2.06	0.55
1:A:31:LYS:HE2	1:A:31:LYS:CA	2.31	0.55
1:B:23:MET:HE3	1:B:196:PRO:HG3	1.89	0.55
1:B:45:LYS:HE2	1:B:46:ALA:H	1.72	0.54
1:B:45:LYS:CD	3:B:911:HOH:O	2.56	0.54
1:A:88:GLN:HE21	1:A:101:ARG:HH21	1.55	0.54
1:A:23:MET:HG3	1:A:196:PRO:HG3	1.90	0.54
1:A:32:SER:HB2	1:A:182:ASP:O	2.07	0.54
1:A:24:GLU:HB2	1:A:151:TRP:CH2	2.44	0.53
1:B:41:GLU:CG	1:B:42:PRO:HD2	2.39	0.53
1:B:64:LEU:HD23	1:B:154:GLY:HA2	1.92	0.52
1:A:50:ASN:ND2	1:A:118:ARG:HH21	2.07	0.52
1:B:45:LYS:HE2	1:B:46:ALA:N	2.24	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:22:TYR:CE2	1:A:151:TRP:CZ3	2.99	0.51
1:B:50:ASN:ND2	1:B:118:ARG:HH11	2.09	0.51
1:B:45:LYS:HE3	1:B:46:ALA:N	2.23	0.51
1:B:156:GLU:O	1:B:158:LEU:N	2.44	0.51
1:A:70:LYS:NZ	3:A:980:HOH:O	2.44	0.50
1:B:41:GLU:CB	1:B:42:PRO:CD	2.90	0.50
1:B:11:ASP:N	3:B:732:HOH:O	2.45	0.50
1:A:30:THR:C	1:A:32:SER:N	2.67	0.49
1:A:182:ASP:O	1:A:183:ARG:HB2	2.12	0.49
1:A:8:ALA:HB3	1:A:200:HIS:CE1	2.47	0.49
1:A:41:GLU:CB	1:A:42:PRO:CD	2.88	0.49
1:B:50:ASN:ND2	1:B:118:ARG:HE	2.08	0.49
1:B:26:HIS:O	1:B:27:ASN:C	2.56	0.48
1:B:158:LEU:HD13	3:B:861:HOH:O	2.13	0.48
1:A:66:LEU:HG	1:A:150:LEU:HD11	1.96	0.48
1:A:8:ALA:CB	1:A:13:GLU:HG3	2.31	0.47
1:B:23:MET:HE3	1:B:196:PRO:CG	2.44	0.47
1:B:45:LYS:HD2	3:B:911:HOH:O	2.13	0.47
1:B:66:LEU:HG	1:B:150:LEU:HD11	1.97	0.46
1:A:41:GLU:HB3	1:A:42:PRO:CD	2.32	0.46
1:B:42:PRO:HA	3:B:887:HOH:O	2.15	0.46
1:B:45:LYS:HE3	1:B:45:LYS:HA	1.98	0.46
1:A:45:LYS:NZ	1:A:45:LYS:HB3	2.31	0.46
1:A:198:ILE:C	1:A:199:LEU:HD12	2.41	0.45
1:B:41:GLU:HG2	1:B:42:PRO:CD	2.46	0.45
1:A:27:ASN:ND2	1:A:148:GLY:HA3	2.32	0.45
1:A:7:LYS:HA	1:A:14:ALA:HB3	1.97	0.45
1:A:22:TYR:CD2	1:A:151:TRP:CE3	3.03	0.45
1:A:158:LEU:O	1:A:161:ALA:HB3	2.17	0.45
1:A:45:LYS:O	1:A:145:ASP:HA	2.17	0.44
1:A:187:HIS:O	1:A:191:ASP:HB2	2.17	0.44
1:B:45:LYS:HE3	1:B:45:LYS:CA	2.47	0.44
1:B:158:LEU:HD21	3:B:799:HOH:O	2.17	0.44
1:A:10:GLY:O	1:A:11:ASP:HB2	2.18	0.43
1:A:120:GLN:NE2	3:A:1042:HOH:O	2.51	0.43
1:B:98:VAL:HB	1:B:137:SER:HA	1.99	0.43
1:A:22:TYR:CD2	1:A:151:TRP:CZ3	3.07	0.43
1:B:27:ASN:CB	1:B:32:SER:HB3	2.47	0.43
1:A:8:ALA:HB2	1:A:13:GLU:CG	2.33	0.43
1:A:8:ALA:HA	1:A:13:GLU:HA	2.00	0.42
1:A:44:GLU:HA	3:A:1053:HOH:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:183:ARG:N	3:A:909:HOH:O	2.49	0.42
1:B:90:MET:CG	1:B:99:VAL:HG22	2.47	0.42
1:A:197:THR:HG22	1:A:199:LEU:CD1	2.49	0.42
1:B:126:GLN:NE2	3:B:864:HOH:O	2.52	0.41
1:A:40:ALA:O	1:A:41:GLU:CB	2.66	0.41
1:B:32:SER:HA	3:B:914:HOH:O	2.20	0.41
1:B:190:GLU:H	1:B:190:GLU:CD	2.28	0.41
1:B:42:PRO:O	1:B:43:SER:CB	2.68	0.41
1:B:28:ALA:O	1:B:29:VAL:CG2	2.68	0.41
1:B:41:GLU:HG2	1:B:42:PRO:HD2	2.03	0.41
1:A:187:HIS:HB2	1:A:191:ASP:CG	2.46	0.41
1:A:191:ASP:OD2	3:A:1059:HOH:O	2.22	0.41
1:B:28:ALA:O	1:B:29:VAL:CB	2.69	0.41
1:A:62:GLN:NE2	1:A:84:ASP:H	2.20	0.40
1:B:29:VAL:O	1:B:30:THR:HB	2.20	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:31:LYS:NZ	3:A:1069:HOH:O[2_555]	1.91	0.29

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	184/204 (90%)	169 (92%)	6 (3%)	9 (5%)	1	0
1	B	183/204 (90%)	170 (93%)	3 (2%)	10 (6%)	1	0
All	All	367/408 (90%)	339 (92%)	9 (2%)	19 (5%)	1	0

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	27	ASN
1	A	31	LYS
1	A	41	GLU
1	A	42	PRO
1	B	27	ASN
1	B	29	VAL
1	B	30	THR
1	A	8	ALA
1	B	12	ALA
1	A	11	ASP
1	B	43	SER
1	B	159	SER
1	A	161	ALA
1	B	157	ARG
1	B	42	PRO
1	B	41	GLU
1	A	10	GLY
1	B	96	LYS
1	A	96	LYS

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	154/168 (92%)	146 (95%)	8 (5%)	21	2
1	B	155/168 (92%)	153 (99%)	2 (1%)	61	29
All	All	309/336 (92%)	299 (97%)	10 (3%)	34	6

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	LYS
1	A	27	ASN
1	A	30	THR
1	A	31	LYS
1	A	32	SER

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Mol	Chain	Res	Type
1	A	45	LYS
1	A	66	LEU
1	A	186	LEU
1	B	43	SER
1	B	45	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	26	HIS
1	A	27	ASN
1	A	48	GLN
1	A	50	ASN
1	A	62	GLN
1	A	88	GLN
1	A	120	GLN
1	A	126	GLN
1	A	143	GLN
1	A	184	GLN
1	B	27	ASN
1	B	50	ASN
1	B	62	GLN
1	B	120	GLN
1	B	126	GLN
1	B	143	GLN
1	B	162	HIS
1	B	195	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

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.