



Full wwPDB NMR Structure Validation Report ⓘ

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PDB ID : 1004
Title : P395S mutant of the p85 regulatory subunit of the N-terminal src homology 2 domain of PI3-Kinase complexed to a peptide derived from PDGFr
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This is a Full wwPDB NMR Structure Validation Report for a publicly released PDB entry.

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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>.

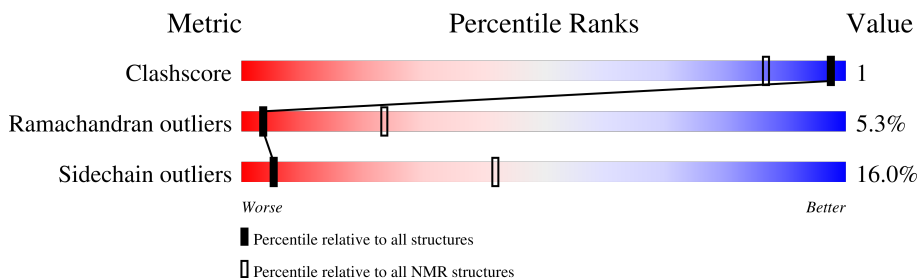
1 Overall quality at a glance

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	158937	12864
Ramachandran outliers	154571	11451
Sidechain outliers	154315	11428

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	111	
2	B	8	

2 Ensemble composition and analysis

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

3 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 1910 atoms, of which 938 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Phosphatidylinositol 3-kinase regulatory alpha subunit.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	111	1779	569	874	158	175	3	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	60	SER	LEU	conflict	UNP P23727
A	75	SER	PRO	engineered mutation	UNP P23727

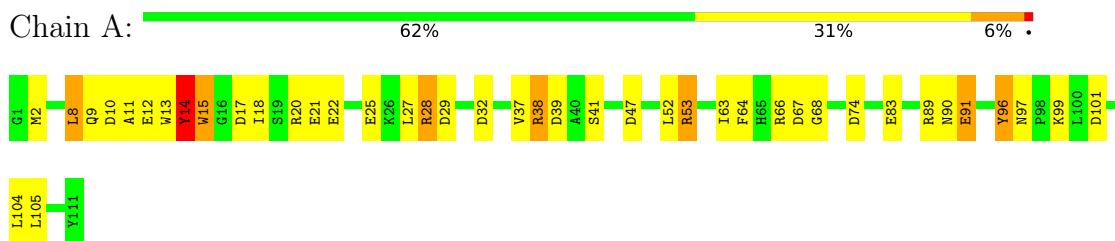
- Molecule 2 is a protein called 8-mer peptide from PDGFr.

Mol	Chain	Residues	Atoms						Trace	
			Total	C	H	N	O	P		S
2	B	8	131	42	64	8	15	1	1	0

4 Residue-property plots [i](#)

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: Phosphatidylinositol 3-kinase regulatory alpha subunit



- Molecule 2: 8-mer peptide from PDGFr



5 Refinement protocol and experimental data overview

The models were refined using the following method: *distance geometry, 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
DYANA	structure solution	1.5
Discover	refinement	97.0
TALOS	structure solution	1999.019.15.47

No chemical shift data was provided.

6 Model quality i

6.1 Standard geometry i

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

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 (average) 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	1.52	11/924 (1.2%)	1.74	36/1245 (2.9%)
2	B	1.37	1/50 (2.0%)	1.61	0/66 (0.0%)
All	All	1.51	12/974 (1.2%)	1.73	36/1311 (2.7%)

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	Planarity
1	A	0	2
All	All	0	2

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

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	12	GLU	CD-OE2	10.36	1.37	1.25
1	A	83	GLU	CD-OE2	10.30	1.36	1.25
1	A	21	GLU	CD-OE2	10.28	1.36	1.25
1	A	22	GLU	CD-OE2	10.09	1.36	1.25
1	A	25	GLU	CD-OE2	10.05	1.36	1.25
1	A	91	GLU	CD-OE2	10.04	1.36	1.25
1	A	32	ASP	CG-OD2	5.25	1.37	1.25
1	A	74	ASP	CG-OD2	5.25	1.37	1.25
1	A	39	ASP	CG-OD2	5.13	1.37	1.25
2	B	184	ASP	CG-OD2	5.10	1.37	1.25
1	A	29	ASP	CG-OD2	5.06	1.36	1.25
1	A	10	ASP	CG-OD2	5.02	1.36	1.25

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	28	ARG	NE-CZ-NH1	8.77	124.68	120.30
1	A	89	ARG	NE-CZ-NH1	8.71	124.66	120.30
1	A	66	ARG	NE-CZ-NH1	8.50	124.55	120.30
1	A	53	ARG	NE-CZ-NH1	8.34	124.47	120.30
1	A	38	ARG	NE-CZ-NH1	7.83	124.22	120.30
1	A	20	ARG	NE-CZ-NH1	7.66	124.13	120.30
1	A	89	ARG	NE-CZ-NH2	-7.43	116.59	120.30
1	A	38	ARG	NE-CZ-NH2	-7.40	116.60	120.30
1	A	28	ARG	NE-CZ-NH2	-7.01	116.79	120.30
1	A	39	ASP	CB-CG-OD2	-6.98	112.02	118.30
1	A	47	ASP	CB-CG-OD2	-6.97	112.02	118.30
1	A	20	ARG	NE-CZ-NH2	-6.69	116.95	120.30
1	A	17	ASP	CB-CG-OD2	-6.54	112.42	118.30
1	A	67	ASP	CB-CG-OD2	-6.53	112.43	118.30
1	A	101	ASP	CB-CG-OD2	-6.37	112.56	118.30
1	A	29	ASP	CB-CG-OD2	-6.28	112.65	118.30
1	A	74	ASP	CB-CG-OD2	-6.26	112.67	118.30
1	A	17	ASP	CB-CG-OD1	6.24	123.91	118.30
1	A	67	ASP	CB-CG-OD1	6.20	123.88	118.30
1	A	15	TRP	CD1-NE1-CE2	-6.15	103.47	109.00
1	A	32	ASP	CB-CG-OD2	-6.10	112.81	118.30
1	A	96	TYR	CB-CG-CD2	-6.09	117.35	121.00
1	A	39	ASP	CB-CG-OD1	6.02	123.72	118.30
1	A	47	ASP	CB-CG-OD1	6.02	123.71	118.30
1	A	101	ASP	CB-CG-OD1	5.99	123.69	118.30
1	A	66	ARG	NE-CZ-NH2	-5.82	117.39	120.30
1	A	11	ALA	CB-CA-C	5.61	118.52	110.10
1	A	13	TRP	CD1-NE1-CE2	-5.61	103.95	109.00
1	A	14	TYR	CA-CB-CG	5.51	123.87	113.40
1	A	14	TYR	CB-CG-CD2	-5.50	117.70	121.00
1	A	10	ASP	CB-CG-OD2	-5.48	113.37	118.30
1	A	74	ASP	CB-CG-OD1	5.29	123.06	118.30
1	A	10	ASP	CB-CG-OD1	5.27	123.04	118.30
1	A	29	ASP	CB-CG-OD1	5.21	122.99	118.30
1	A	32	ASP	CB-CG-OD1	5.20	122.98	118.30
1	A	53	ARG	NE-CZ-NH2	-5.07	117.76	120.30

There are no chirality outliers.

All planar outliers are listed below.

Mol	Chain	Res	Type	Group
1	A	53	ARG	Sidechain
1	A	96	TYR	Sidechain

6.2 Too-close contacts [i](#)

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	905	874	876	1
All	All	972	938	936	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 1.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)
1:A:14:TYR:HA	1:A:37:VAL:HG21	0.55	1.79

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	109/111 (98%)	70 (64%)	33 (30%)	6 (6%)	3	23
2	B	5/8 (62%)	3 (60%)	2 (40%)	0 (0%)	100	100
All	All	114/119 (96%)	73 (64%)	35 (31%)	6 (5%)	3	23

All 6 Ramachandran outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type
1	A	8	LEU
1	A	27	LEU
1	A	63	ILE
1	A	64	PHE
1	A	68	GLY
1	A	97	ASN

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	99/99 (100%)	84 (85%)	15 (15%)	6	44
2	B	7/7 (100%)	5 (71%)	2 (29%)	2	18
All	All	106/106 (100%)	89 (84%)	17 (16%)	5	42

All 17 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	2	MET
1	A	8	LEU
1	A	9	GLN
1	A	14	TYR
1	A	15	TRP
1	A	18	ILE
1	A	28	ARG
1	A	38	ARG
1	A	41	SER
1	A	52	LEU
1	A	90	ASN
1	A	91	GLU
1	A	99	LYS
1	A	104	LEU
1	A	105	LEU
2	B	184	ASP
2	B	189	LEU

6.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue 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	PTR	B	185	2	15,16,17	1.44	2 (13%)

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	PTR	B	185	2	19,22,24	1.07	1 (5%)

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	PTR	B	185	2	-	0,10,11,13	0,1,1,1

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	B	185	PTR	P-O1P	3.22	1.60	1.50
2	B	185	PTR	P-OH	3.20	1.64	1.59

All angle outliers are listed below.

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	185	PTR	OH-CZ-CE1	2.18	125.70	119.23

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.6 Ligand geometry [i](#)

There are no ligands in this entry.

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