



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2018 – 06:10 pm GMT

PDB ID : 1A1C  
Title : C-SRC (SH2 DOMAIN) COMPLEXED WITH ACE-PHOSPHOTYR-GLU-(  
N-ME(-(CH<sub>2</sub>)<sub>3</sub>-CYCLOPENTYL))  
Authors : Shewchuk, L.; Jordan, S.  
Deposited on : 1997-12-10  
Resolution : 2.40 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](mailto: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.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.7.3 (157068), CSD as539be (2018)  
Xtriage (Phenix) : 1.13  
EDS : trunk30967  
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)  
Refmac : 5.8.0158  
CCP4 : 7.0 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk30967

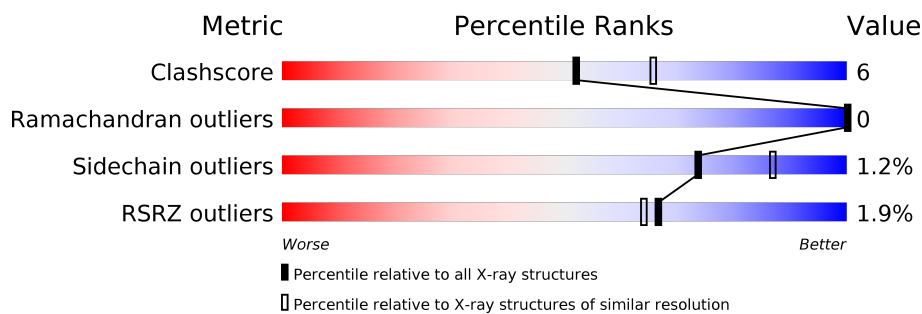
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.40 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	122126	3956 (2.40-2.40)
Ramachandran outliers	120053	3897 (2.40-2.40)
Sidechain outliers	120020	3898 (2.40-2.40)
RSRZ outliers	108989	3386 (2.40-2.40)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	107	<div> <div>3%</div> <div>73%</div> <div>21%</div> <div>5%</div> <div>.</div> </div>
1	B	107	<div> <div>%</div> <div>77%</div> <div>19%</div> <div>.</div> <div>.</div> </div>
2	C	4	<div> <div>50%</div> <div>50%</div> </div>
2	D	4	<div> <div>75%</div> <div>25%</div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2242 atoms, of which 479 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 C-SRC TYROSINE KINASE.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	105	Total	C	H	N	O	S	0	0	0
			1041	527	203	149	159	3			
1	B	104	Total	C	H	N	O	S	0	0	0
			995	510	188	142	152	3			

- Molecule 2 is a protein called ACE-PHOSPHOTYR-GLU-(N-ME(-(CH<sub>2</sub>)<sub>3</sub>-CYCLOPENTYL)).

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	C	4	Total	C	H	N	O	P	0	0	0
			40	25	2	3	9	1			
2	D	4	Total	C	H	N	O	P	0	0	0
			40	25	2	3	9	1			

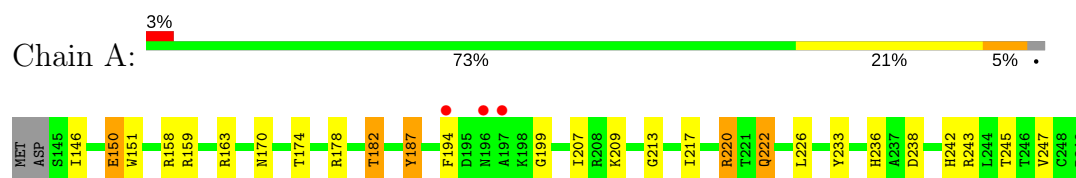
- Molecule 3 is water.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	21	Total	H	O	0	0
			63	42	21		
3	B	16	Total	H	O	0	0
			48	32	16		
3	C	3	Total	H	O	0	0
			9	6	3		
3	D	2	Total	H	O	0	0
			6	4	2		

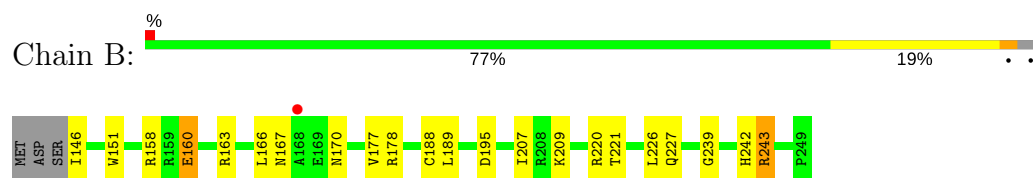
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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: C-SRC TYROSINE KINASE



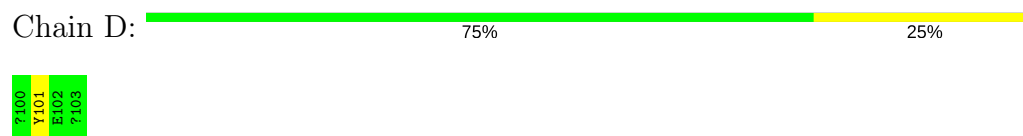
- Molecule 1: C-SRC TYROSINE KINASE



- Molecule 2: ACE-PHOSPHOTYR-GLU-(N-ME(-(CH<sub>2</sub>)<sub>3</sub>-CYCLOPENTYL))



- Molecule 2: ACE-PHOSPHOTYR-GLU-(N-ME(-(CH<sub>2</sub>)<sub>3</sub>-CYCLOPENTYL))



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	51.70Å 66.90Å 75.60Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	6.00 – 2.40 50.10 – 2.40	Depositor EDS
% Data completeness (in resolution range)	92.0 (6.00-2.40) 91.8 (50.10-2.40)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.23 (at 2.39Å)	Xtriage
Refinement program	X-PLOR 3.1	Depositor
R, $R_{free}$	0.186 , (Not available) 0.173 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	31.2	Xtriage
Anisotropy	0.179	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 71.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	2242	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.65% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: DIX, PTR, ACE

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	1.04	0/855	1.82	17/1153 (1.5%)
1	B	1.00	0/824	1.68	16/1116 (1.4%)
2	C	1.01	0/8	0.91	0/9
2	D	1.36	0/8	0.97	0/9
All	All	1.02	0/1695	1.75	33/2287 (1.4%)

There are no bond length outliers.

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	158	ARG	NE-CZ-NH2	-13.52	113.54	120.30
1	A	159	ARG	NE-CZ-NH2	-10.58	115.01	120.30
1	A	158	ARG	NE-CZ-NH1	10.12	125.36	120.30
1	A	220	ARG	NE-CZ-NH2	-9.56	115.52	120.30
1	A	220	ARG	NE-CZ-NH1	8.96	124.78	120.30
1	B	163	ARG	NE-CZ-NH2	-8.77	115.92	120.30
1	B	151	TRP	CD1-CG-CD2	8.24	112.89	106.30
1	A	178	ARG	NE-CZ-NH2	-7.97	116.31	120.30
1	B	151	TRP	CE2-CD2-CG	-7.86	101.01	107.30
1	B	243	ARG	NE-CZ-NH2	-7.83	116.39	120.30
1	A	151	TRP	CE2-CD2-CG	-7.63	101.19	107.30
1	B	160	GLU	CA-CB-CG	-7.43	97.06	113.40
1	A	151	TRP	CD1-CG-CD2	7.00	111.90	106.30
1	A	151	TRP	CB-CG-CD1	-6.96	117.95	127.00
1	A	182	THR	CA-CB-CG2	6.66	121.73	112.40
1	A	163	ARG	NE-CZ-NH2	-6.64	116.98	120.30
1	B	221	THR	CA-CB-CG2	6.46	121.45	112.40
1	A	151	TRP	CG-CD2-CE3	6.46	139.71	133.90
1	B	177	VAL	CG1-CB-CG2	-6.26	100.88	110.90
1	B	151	TRP	CB-CG-CD1	-6.21	118.93	127.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	209	LYS	CB-CG-CD	-5.87	96.35	111.60
1	A	163	ARG	NE-CZ-NH1	5.86	123.23	120.30
1	A	150	GLU	CA-CB-CG	5.84	126.25	113.40
1	B	178	ARG	NE-CZ-NH2	-5.84	117.38	120.30
1	B	158	ARG	NE-CZ-NH2	-5.77	117.42	120.30
1	B	166	LEU	CA-C-N	-5.74	104.57	117.20
1	B	220	ARG	NE-CZ-NH2	-5.67	117.46	120.30
1	A	217	ILE	O-C-N	-5.57	113.78	122.70
1	A	217	ILE	CA-C-N	5.50	129.31	117.20
1	B	151	TRP	CG-CD1-NE1	-5.49	104.61	110.10
1	B	151	TRP	CG-CD2-CE3	5.30	138.67	133.90
1	B	239	GLY	O-C-N	-5.20	114.39	122.70
1	A	187	TYR	CB-CG-CD2	-5.19	117.89	121.00

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	838	203	809	14	0
1	B	807	188	759	6	1
2	C	38	2	34	1	0
2	D	38	2	34	1	0
3	A	21	42	0	0	1
3	B	16	32	0	0	0
3	C	3	6	0	0	0
3	D	2	4	0	0	0
All	All	1763	479	1636	20	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 6.

All (20) 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:242:HIS:HD2	1:B:243:ARG:O	1.75	0.69
1:A:222:GLN:HE21	1:A:222:GLN:N	1.92	0.67
1:A:242:HIS:HD2	1:A:243:ARG:O	1.82	0.63
1:B:167:ASN:HB3	1:B:170:ASN:ND2	2.15	0.60
1:A:222:GLN:NE2	1:A:222:GLN:N	2.50	0.59
1:A:209:LYS:HE3	1:A:213:GLY:O	2.02	0.59
1:A:146:ILE:CD1	1:A:226:LEU:HD12	2.39	0.53
1:A:170:ASN:HB3	1:A:174:THR:OG1	2.12	0.49
1:A:182:THR:HG21	1:B:160:GLU:HG2	1.96	0.48
1:B:189:LEU:HB2	1:B:207:ILE:HD11	1.95	0.47
1:A:243:ARG:NH2	1:A:245:THR:HG22	2.29	0.47
1:A:242:HIS:CD2	1:A:243:ARG:O	2.67	0.46
1:A:150:GLU:HG3	1:A:247:VAL:HG11	1.98	0.44
1:A:233:TYR:HA	1:A:236:HIS:O	2.17	0.44
2:C:102:GLU:HA	2:C:103:DIX:HM1	1.80	0.44
1:B:188:CYS:SG	2:D:101:PTR:HE2	2.59	0.43
1:A:220:ARG:HD3	1:A:238:ASP:OD2	2.19	0.43
1:A:194:PHE:HA	1:A:199:GLY:O	2.20	0.42
1:B:146:ILE:CD1	1:B:226:LEU:HD12	2.50	0.42
1:A:187:TYR:HB2	1:A:207:ILE:HB	2.02	0.42

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:227:GLN:HE21	3:A:539:HOH:H2[4_456]	1.30	0.30

## 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	103/107 (96%)	99 (96%)	4 (4%)	0	<b>100</b> <b>100</b>

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	102/107 (95%)	98 (96%)	4 (4%)	0	100	100
2	C	1/4 (25%)	1 (100%)	0	0	100	100
2	D	1/4 (25%)	1 (100%)	0	0	100	100
All	All	207/222 (93%)	199 (96%)	8 (4%)	0	100	100

There are no Ramachandran outliers to report.

### 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	89/94 (95%)	88 (99%)	1 (1%)	76	88
1	B	82/94 (87%)	81 (99%)	1 (1%)	74	87
2	C	1/1 (100%)	1 (100%)	0	100	100
2	D	1/1 (100%)	1 (100%)	0	100	100
All	All	173/190 (91%)	171 (99%)	2 (1%)	74	87

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	222	GLN
1	B	195	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	222	GLN
1	A	242	HIS
1	B	242	HIS

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues are modelled in this entry.

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	$\# Z  > 2$
2	PTR	C	101	2	16,16,17	2.67	3 (18%)	21,22,24	1.00	0
2	PTR	D	101	2	16,16,17	2.24	3 (18%)	21,22,24	1.48	5 (23%)

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	C	101	2	-	0/9/11/13	0/1/1/1
2	PTR	D	101	2	-	0/9/11/13	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	101	PTR	OH-CZ	-4.56	1.30	1.40
2	C	101	PTR	OH-CZ	-2.91	1.34	1.40
2	C	101	PTR	CB-CA	-2.38	1.48	1.53
2	D	101	PTR	CB-CA	-2.14	1.49	1.53
2	D	101	PTR	P-OH	6.93	1.72	1.59
2	C	101	PTR	P-OH	9.51	1.76	1.59

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	101	PTR	O2P-P-OH	-3.13	94.92	105.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	101	PTR	O3P-P-OH	-2.62	96.65	105.50
2	D	101	PTR	CG-CB-CA	-2.25	109.76	114.23
2	D	101	PTR	O2P-P-O1P	2.49	120.32	110.60
2	D	101	PTR	O3P-P-O2P	3.07	119.73	107.59

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	101	PTR	1	0

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

There are no ligands in this entry.

## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	105/107 (98%)	-0.40	3 (2%) 51 49	10, 24, 58, 72	0
1	B	104/107 (97%)	-0.45	1 (0%) 82 80	8, 23, 51, 67	0
2	C	1/4 (25%)	-1.00	0 100 100	23, 23, 23, 23	0
2	D	1/4 (25%)	-0.93	0 100 100	29, 29, 29, 29	0
All	All	211/222 (95%)	-0.43	4 (1%) 66 64	8, 24, 58, 72	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	194	PHE	3.2
1	B	168	ALA	2.5
1	A	197	ALA	2.4
1	A	196	ASN	2.1

### 6.2 Non-standard residues in protein, DNA, RNA chains [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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
2	PTR	D	101	16/17	0.96	0.14	0,28,36,39	0
2	PTR	C	101	16/17	0.98	0.10	0,13,18,21	0

### 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 6.4 Ligands

There are no ligands in this entry.

## 6.5 Other polymers

There are no such residues in this entry.