Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 11:29 PM GMT

PDB ID	:	1XJD
Title	:	Crystal Structure of PKC-theta complexed with Staurosporine at 2A resolution
Authors	:	Xu, Z.B.
Deposited on	:	2004-09-23
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
http://wwpdb.org/validation/2016/XrayValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

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

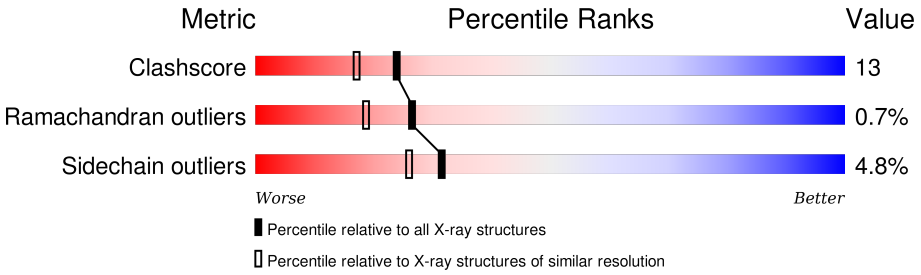
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

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	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (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 on the lower bar indicate the fraction of residues that contain outliers for >=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 <=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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	345	63% 15% • 18%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2502 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.

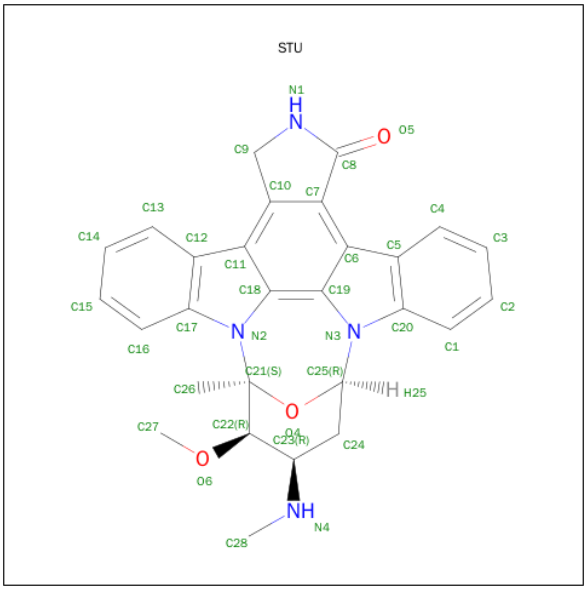
- Molecule 1 is a protein called Protein kinase C, theta type.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	P	S			
1	A	282	2352	1525	388	421	2	16	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	538	TPO	THR	MODIFIED RESIDUE	UNP Q04759
A	695	SEP	SER	MODIFIED RESIDUE	UNP Q04759

- Molecule 2 is STAUROSPORINE (three-letter code: STU) (formula: $C_{28}H_{26}N_4O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
2	A	1	35	28	4	3	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	115	Total	O	0	0
			115	115		

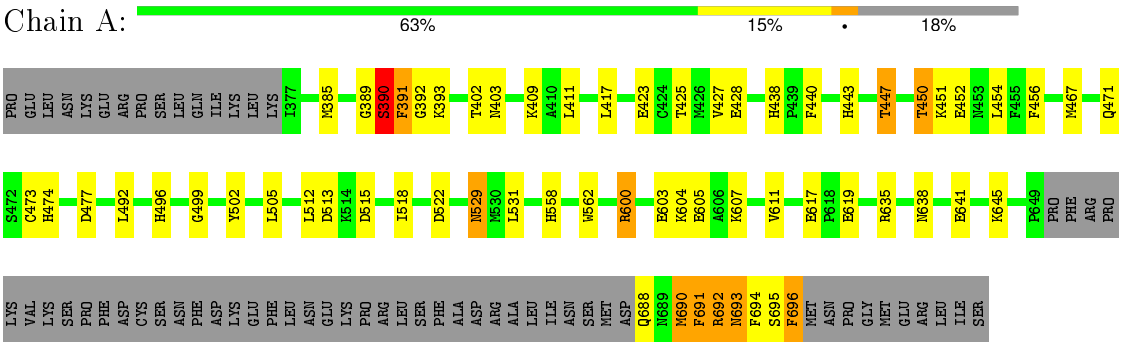
3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

Note EDS was not executed.

- Molecule 1: Protein kinase C, theta type

Chain A:



4 Data and refinement statistics (i)

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

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	139.57Å 42.40Å 67.68Å 90.00° 116.24° 90.00°	Depositor
Resolution (Å)	19.93 – 2.00	Depositor
% Data completeness (in resolution range)	83.2 (19.93-2.00)	Depositor
R_{merge}	0.06	Depositor
R_{sym}	0.04	Depositor
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.201 , 0.216	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2502	wwPDB-VP
Average B, all atoms (Å^2)	29.0	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: TPO, STU, SEP

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	0.41	0/2390	0.64	0/3210

There are no bond length outliers.

There are no bond angle outliers.

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	2352	0	2304	62	1
2	A	35	0	26	1	0
3	A	115	0	0	4	1
All	All	2502	0	2330	62	2

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 (62) 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:385:MET:CE	1:A:393:LYS:HD2	2.04	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:417:LEU:HD21	1:A:690:MET:HG3	1.59	0.84
1:A:385:MET:HE3	1:A:393:LYS:HD2	1.59	0.84
1:A:603:GLU:HG3	3:A:103:HOH:O	1.82	0.78
1:A:447:THR:HG22	1:A:696:PHE:N	1.98	0.77
1:A:392:GLY:HA3	1:A:411:LEU:HD23	1.67	0.75
1:A:450:THR:CG2	1:A:451:LYS:N	2.52	0.72
1:A:447:THR:N	1:A:696:PHE:O	2.22	0.72
1:A:392:GLY:CA	1:A:411:LEU:HD23	2.19	0.72
1:A:450:THR:HG22	1:A:451:LYS:N	2.07	0.70
1:A:696:PHE:C	1:A:696:PHE:CD1	2.69	0.65
1:A:450:THR:CG2	1:A:451:LYS:H	2.08	0.65
1:A:417:LEU:CD2	1:A:690:MET:HG3	2.25	0.65
1:A:391:PHE:HZ	1:A:425:THR:HA	1.62	0.64
1:A:391:PHE:CZ	1:A:425:THR:HA	2.33	0.64
1:A:450:THR:HG22	1:A:452:GLU:H	1.65	0.61
1:A:391:PHE:C	1:A:411:LEU:HD21	2.24	0.58
1:A:438:HIS:O	1:A:443:HIS:HE1	1.88	0.57
1:A:450:THR:HG23	1:A:451:LYS:H	1.70	0.56
1:A:391:PHE:O	1:A:409:LYS:HE3	2.06	0.55
1:A:499:GLY:HA3	1:A:531:LEU:HD21	1.89	0.54
1:A:454:LEU:HD22	1:A:694:PHE:CD2	2.43	0.54
1:A:447:THR:HG23	1:A:695:SEP:HA	1.89	0.54
1:A:411:LEU:HD12	1:A:456:PHE:HE1	1.72	0.53
1:A:423:GLU:O	1:A:427:VAL:HG23	2.08	0.53
1:A:389:GLY:O	1:A:390:SER:C	2.46	0.53
1:A:417:LEU:HG	1:A:691:PHE:HZ	1.74	0.53
1:A:513:ASP:HB3	1:A:515:ASP:OD1	2.09	0.52
1:A:402:THR:O	1:A:403:ASN:HB2	2.10	0.51
1:A:447:THR:HG22	1:A:696:PHE:HB3	1.92	0.51
1:A:605:GLU:HG2	3:A:103:HOH:O	2.09	0.51
1:A:391:PHE:O	1:A:411:LEU:HD21	2.11	0.51
1:A:392:GLY:HA3	1:A:411:LEU:CD2	2.40	0.50
1:A:391:PHE:O	1:A:409:LYS:CE	2.59	0.50
1:A:385:MET:HE3	1:A:393:LYS:CD	2.38	0.50
1:A:467:MET:O	1:A:471:GLN:HG3	2.12	0.49
1:A:696:PHE:O	1:A:696:PHE:HD1	1.96	0.48
1:A:605:GLU:HG3	1:A:635:ARG:NH2	2.29	0.48
1:A:447:THR:CG2	1:A:694:PHE:O	2.63	0.47
1:A:447:THR:HG21	1:A:694:PHE:O	2.14	0.47
1:A:438:HIS:CD2	1:A:440:PHE:H	2.32	0.47
1:A:496:HIS:CE1	1:A:558:HIS:HB2	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:492:LEU:HD11	1:A:505:LEU:HD22	1.98	0.46
1:A:391:PHE:C	1:A:411:LEU:CD2	2.84	0.45
1:A:390:SER:HB2	1:A:522:ASP:OD1	2.16	0.45
1:A:392:GLY:N	1:A:411:LEU:HD23	2.32	0.45
1:A:617:GLU:OE1	1:A:619:GLU:HG3	2.17	0.45
1:A:607:LYS:O	1:A:611:VAL:HG23	2.17	0.44
1:A:450:THR:HG22	1:A:452:GLU:N	2.31	0.44
1:A:512:LEU:HD23	1:A:518:ILE:HG22	1.99	0.44
1:A:558:HIS:HD2	1:A:562:TRP:HE1	1.66	0.44
1:A:409:LYS:NZ	1:A:428:GLU:OE1	2.51	0.43
1:A:389:GLY:HA3	2:A:200:STU:H16	2.02	0.41
1:A:502:TYR:HE2	1:A:505:LEU:HA	1.85	0.41
1:A:499:GLY:CA	1:A:531:LEU:HD21	2.51	0.41
1:A:645:LYS:HA	3:A:34:HOH:O	2.20	0.41
1:A:454:LEU:HD22	1:A:694:PHE:CE2	2.55	0.41
1:A:600:ARG:HD3	3:A:42:HOH:O	2.20	0.41
1:A:638:ASN:CG	1:A:641:GLU:HG3	2.40	0.41
1:A:473:CYS:O	1:A:474:HIS:HB2	2.20	0.41
1:A:529:ASN:HD22	1:A:529:ASN:HA	1.57	0.41
1:A:692:ARG:O	1:A:693:ASN:HB2	2.21	0.40

All (2) 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)
3:A:114:HOH:O	3:A:114:HOH:O[2_656]	1.77	0.43
1:A:477:ASP:OD1	1:A:477:ASP:OD1[2_656]	2.18	0.02

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	276/345 (80%)	262 (95%)	12 (4%)	2 (1%)	26 19

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	690	MET
1	A	390	SER

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	252/312 (81%)	240 (95%)	12 (5%)	31 26

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	390	SER
1	A	391	PHE
1	A	447	THR
1	A	450	THR
1	A	529	ASN
1	A	600	ARG
1	A	604	LYS
1	A	688	GLN
1	A	691	PHE
1	A	692	ARG
1	A	693	ASN
1	A	696	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	404	GLN
1	A	438	HIS

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Mol	Chain	Res	Type
1	A	443	HIS
1	A	529	ASN
1	A	554	GLN
1	A	558	HIS
1	A	582	GLN
1	A	595	ASN
1	A	630	GLN
1	A	688	GLN
1	A	693	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)

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	$\#\|Z\| > 2$
1	TPO	A	538	1	8,10,11	2.07	1 (12%)	7,14,16	1.52	0
1	SEP	A	695	1	8,9,10	2.20	1 (12%)	8,12,14	3.28	3 (37%)

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
1	TPO	A	538	1	-	0/8/11/13	0/0/0/0
1	SEP	A	695	1	-	0/6/8/10	0/0/0/0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	695	SEP	P-OG	-5.93	1.40	1.60
1	A	538	TPO	P-OG1	-5.47	1.43	1.60

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	695	SEP	O2P-P-OG	-5.94	89.47	106.56
1	A	695	SEP	O3P-P-O1P	-2.00	104.14	110.58
1	A	695	SEP	O3P-P-OG	5.76	123.14	106.56

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
1	A	695	SEP	1	0

5.5 Carbohydrates (i)

There are no carbohydrates in this entry.

5.6 Ligand geometry (i)

1 ligand is 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	STU	A	200	-	27,42,42	4.49	22 (81%)	23,68,68	1.31	3 (13%)

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	STU	A	200	-	-	0/4/42/42	0/0/8/8

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	200	STU	C28-N4	2.34	1.53	1.46
2	A	200	STU	C26-C21	2.68	1.55	1.51
2	A	200	STU	C15-C16	3.00	1.43	1.36
2	A	200	STU	C13-C12	3.32	1.47	1.41
2	A	200	STU	C4-C5	3.35	1.47	1.41
2	A	200	STU	C3-C4	3.58	1.44	1.36
2	A	200	STU	C16-C17	3.70	1.48	1.41
2	A	200	STU	C22-C23	3.95	1.56	1.52
2	A	200	STU	C14-C13	4.03	1.45	1.36
2	A	200	STU	C2-C1	4.35	1.46	1.36
2	A	200	STU	C1-C20	4.40	1.50	1.41
2	A	200	STU	O6-C22	4.49	1.51	1.42
2	A	200	STU	C3-C2	4.54	1.49	1.38
2	A	200	STU	C10-C11	4.79	1.50	1.42
2	A	200	STU	C14-C15	5.03	1.50	1.38
2	A	200	STU	C24-C23	5.42	1.62	1.53
2	A	200	STU	C6-C19	5.97	1.50	1.42
2	A	200	STU	C5-C20	6.02	1.51	1.41
2	A	200	STU	C12-C17	6.07	1.51	1.41
2	A	200	STU	C24-C25	6.54	1.62	1.51
2	A	200	STU	C11-C18	7.02	1.52	1.42
2	A	200	STU	C7-C6	8.96	1.57	1.43

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	200	STU	C24-C23-N4	-2.51	106.48	112.16
2	A	200	STU	C1-C20-N3	-2.27	129.45	132.18
2	A	200	STU	C11-C12-C17	2.95	109.60	106.37

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	A	200	STU	1	0

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 will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates

EDS was not executed - this section will therefore be empty.

6.4 Ligands

EDS was not executed - this section will therefore be empty.

6.5 Other polymers

EDS was not executed - this section will therefore be empty.