

# Full wwPDB X-ray Structure Validation Report (i)

Jan 22, 2023 – 12:18 am GMT

PDB ID : 8B58

Title : Crystal Structure of Cyclophilin TgCyp23 from Toxoplasma gondii in complex

with Cyclosporin A

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Deposited on : 2022-09-22

Resolution : 1.10 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 (i) 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 (1)) were used in the production of this report:

 $Mol Probity \quad : \quad 4.02b\text{--}467$ 

Mogul : 1.8.4, CSD as541be (2020)

Xtriage (Phenix) : 1.13 EDS : 2.31.3

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

 $Refmac \quad : \quad 5.8.0158$ 

CCP4 : 7.0.044 (Gargrove)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

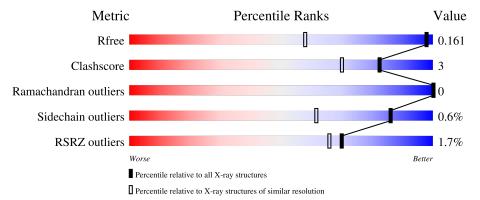
Validation Pipeline (wwPDB-VP) : 2.31.3

## 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 1.10 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	Similar resolution
Metric	$(\# \mathrm{Entries})$	$(\#  ext{Entries},  ext{ resolution range}( ext{Å}))$
$R_{free}$	130704	1619 (1.14-1.06)
Clashscore	141614	1671 (1.14-1.06)
Ramachandran outliers	138981	1615 (1.14-1.06)
Sidechain outliers	138945	1613 (1.14-1.06)
RSRZ outliers	127900	1588 (1.14-1.06)

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 <=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	211	2%	91%	• 6%
1	В	211	.%	92%	• 5%
2	С	11	27%	55%	18%
2	D	11	9%	91%	



## 2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 3824 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 Peptidyl-prolyl cis-trans isomerase.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
1	Δ	199	Total	С	N	О	S	0	0	0
1	A 199	199	1593	1012	282	291	8		9	
1	D	200	Total	С	N	О	S	0	Q	0
1	Ъ	В 200	1593	1010	281	294	8		8	

• Molecule 2 is a protein called Cyclosporin A.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	С	11	Total 85			0	0	0
2	D	11	Total 85		N 11	0	0	0

• Molecule 3 is water.

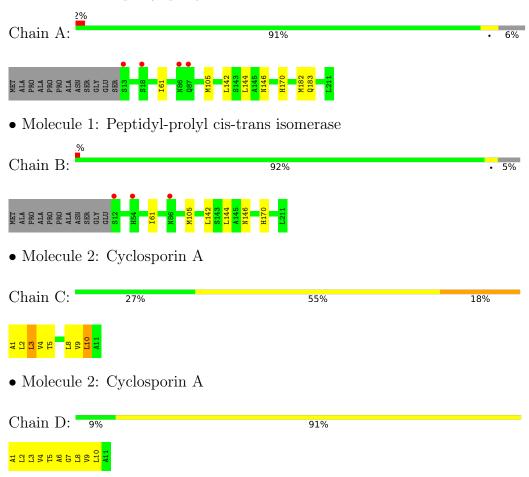
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	221	Total O 221 221	0	0
3	В	230	Total O 230 230	0	0
3	С	7	Total O 7 7	0	0
3	D	10	Total O 10 10	0	0



## 3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: Peptidyl-prolyl cis-trans isomerase





## 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants	38.40Å 119.42Å 46.35Å	Donositor
a, b, c, $\alpha$ , $\beta$ , $\gamma$	90.00° 103.62° 90.00°	Depositor
Resolution (Å)	45.09 - 1.10	Depositor
rtesolution (A)	45.05 - 1.10	EDS
% Data completeness	99.8 (45.09-1.10)	Depositor
(in resolution range)	99.8 (45.05-1.10)	EDS
$R_{merge}$	0.04	Depositor
$R_{sym}$	(Not available)	Depositor
$< I/\sigma(I) > 1$	1.80 (at 1.10Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
P. P.	0.133 , $0.153$	Depositor
$R, R_{free}$	0.142 , 0.161	DCC
$R_{free}$ test set	8309  reflections  (5.09%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	12.3	Xtriage
Anisotropy	0.141	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$ , $B_{sol}(Å^2)$	0.34, 42.4	EDS
L-test for twinning <sup>2</sup>	$ < L >=0.50, < L^2>=0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.98	EDS
Total number of atoms	3824	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 6.22% of the height of the origin peak. No significant pseudotranslation is detected.

<sup>&</sup>lt;sup>2</sup>Theoretical values of <|L|>,  $<L^2>$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



<sup>&</sup>lt;sup>1</sup>Intensities estimated from amplitudes.

## 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: ABA, BMT, SAR, DAL, MVA, MLE

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).

Mal	Chain	Bond	lengths	Bond angles		
Mol	Chain	RMSZ	# Z  > 5	RMSZ	# Z  > 5	
1	A	0.55	0/1648	0.76	1/2220 (0.0%)	
1	В	0.53	0/1643	0.75	0/2214	
2	С	0.87	0/10	1.36	0/11	
2	D	0.58	0/10	0.72	0/11	
All	All	0.54	0/3311	0.76	1/4456 (0.0%)	

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
1	A	182	MET	CG-SD-CE	5.50	109.00	100.20

There are no chirality outliers.

There are no planarity outliers.

#### 5.2 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 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	1593	0	1587	5	0
1	В	1593	0	1574	4	0
2	С	85	0	110	4	0
2	D	85	0	110	6	0
3	A	221	0	0	2	0
3	В	230	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	С	7	0	0	0	0
3	D	10	0	0	0	0
All	All	3824	0	3381	17	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 3.

All (17) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic	Clash	
7100111-1	7100111-2	$\operatorname{distance}\left( \operatorname{\AA}\right)$	overlap(A)	
1:B:170:HIS:HE1	3:B:478:HOH:O	1.79	0.66	
1:A:170:HIS:HE1	3:A:469:HOH:O	1.82	0.61	
1:A:144:LEU:O	1:A:170:HIS:HD2	1.83	0.61	
2:D:1:DAL:C	2:D:3:MLE:HN1	2.33	0.59	
1:B:144:LEU:O	1:B:170:HIS:HD2	1.88	0.55	
2:D:9:VAL:HA	2:D:10:MLE:HN1	1.72	0.47	
2:D:6:ABA:C	2:D:8:MLE:HN1	2.45	0.46	
2:C:9:VAL:HA	2:C:10:MLE:HN1	1.69	0.45	
1:A:146:ASN:O	2:C:5:BMT:HA	2.16	0.45	
2:C:1:DAL:HA	2:C:2:MLE:HN1	1.71	0.45	
1:A:183:GLN:NE2	3:A:303:HOH:O	2.49	0.44	
2:C:1:DAL:C	2:C:3:MLE:HN1	2.47	0.44	
2:D:1:DAL:HA	2:D:2:MLE:HN1	1.70	0.43	
1:B:146:ASN:O	2:D:5:BMT:HA	2.19	0.43	
1:A:61:ILE:HG21	1:A:142[B]:LEU:HD22	2.03	0.41	
2:D:2:MLE:N	2:D:3:MLE:HN1	2.35	0.41	
1:B:61:ILE:HG21	1:B:142:LEU:HD22	2.03	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	in Analysed Favoured Allowed		Outliers	Perce	entiles	
1	A	$206/211\ (98\%)$	203 (98%)	3 (2%)	0	100	100
1	В	$206/211\ (98\%)$	202 (98%)	4 (2%)	0	100	100
2	C	1/11~(9%)	1 (100%)	0	0	100	100
2	D	1/11~(9%)	1 (100%)	0	0	100	100
All	All	414/444~(93%)	407 (98%)	7 (2%)	0	100	100

There are no Ramachandran outliers to report.

#### 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	Percen	tiles
1	A	176/175 (101%)	175 (99%)	1 (1%)	86	61
1	В	176/175~(101%)	175 (99%)	1 (1%)	86	61
2	$\mathbf{C}$	1/1 (100%)	1 (100%)	0	100	100
2	D	1/1 (100%)	1 (100%)	0	100	100
All	All	354/352 (101%)	352 (99%)	2 (1%)	86	61

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

Mol	Chain	Res	Type
1	A	105	MET
1	В	105	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	170	HIS
1	A	183	GLN
1	В	103	ASN
1	В	170	HIS



#### 5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

18 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).

N T 1	TD.	aı ·	ъ	т. 1	Вс	ond leng	ths	В	ond ang	gles
Mol	Type	Chain	Res	Link	Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	MLE	D	2	2	7,8,9	0.46	0	6,9,11	0.73	0
2	MVA	D	4	2	6,7,8	0.54	0	7,8,10	1.38	2 (28%)
2	ABA	С	6	2	4,5,6	0.43	0	1,5,7	0.72	0
2	BMT	D	5	2	11,12,13	0.39	0	12,14,16	0.80	0
2	ABA	D	6	2	4,5,6	0.44	0	1,5,7	0.42	0
2	MLE	D	3	2	7,8,9	0.56	0	6,9,11	0.91	0
2	DAL	D	1	2	3,4,5	0.74	0	2,4,6	1.19	0
2	MLE	D	10	2	7,8,9	0.47	0	6,9,11	0.75	0
2	MLE	С	8	2	7,8,9	1.14	1 (14%)	6,9,11	1.18	1 (16%)
2	BMT	С	5	2	11,12,13	0.52	0	12,14,16	0.71	0
2	MLE	С	10	2	7,8,9	0.47	0	6,9,11	0.98	1 (16%)
2	MVA	С	4	2	6,7,8	0.40	0	7,8,10	1.27	1 (14%)
2	MLE	D	8	2	7,8,9	0.61	0	6,9,11	0.75	0
2	MLE	С	3	2	7,8,9	0.43	0	6,9,11	1.02	1 (16%)
2	DAL	С	1	2	3,4,5	0.80	0	2,4,6	1.14	0
2	SAR	D	7	2	4,4,5	0.93	0	1,3,5	2.16	1 (100%)
2	MLE	С	2	2	7,8,9	0.40	0	6,9,11	0.84	0
2	SAR	С	7	2	4,4,5	0.75	0	1,3,5	1.58	0

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	MLE	D	2	2	-	0/5/8/10	-
2	MVA	D	4	2	-	1/6/8/10	-
2	ABA	С	6	2	-	0/3/4/6	-
2	BMT	D	5	2	-	1/13/16/18	-
2	ABA	D	6	2	-	0/3/4/6	-
2	MLE	D	3	2	-	0/5/8/10	-
2	DAL	D	1	2	-	0/0/2/4	-
2	MLE	D	10	2	-	0/5/8/10	-
2	MLE	С	8	2	-	0/5/8/10	-
2	BMT	С	5	2	-	1/13/16/18	-
2	MLE	С	10	2	-	0/5/8/10	-
2	MVA	С	4	2	-	3/6/8/10	-
2	MLE	D	8	2	-	1/5/8/10	-
2	MLE	С	3	2	-	0/5/8/10	-
2	DAL	С	1	2	-	0/0/2/4	-
2	SAR	D	7	2	-	1/1/2/3	-
2	MLE	С	2	2	-	0/5/8/10	-
2	SAR	С	7	2	-	1/1/2/3	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(\text{\AA})$	$Ideal(\AA)$
2	С	8	MLE	O-C	2.30	1.29	1.19

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	$\mathbf{Z}$	$Observed(^o)$	$Ideal(^{o})$
2	С	8	MLE	CN-N-CA	2.18	120.41	113.64
2	D	7	SAR	O-C-CA	-2.16	119.17	125.42
2	С	3	MLE	O-C-CA	-2.13	119.19	124.78
2	С	4	MVA	CB-CA-N	2.13	113.95	111.17
2	D	4	MVA	CB-CA-N	2.08	113.88	111.17
2	D	4	MVA	CG1-CB-CA	-2.06	108.05	111.21
2	С	10	MLE	O-C-CA	-2.03	119.46	124.78

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	8	MLE	O-C-CA-CB
2	С	5	BMT	CB-CA-N-CN
2	D	5	BMT	CB-CA-N-CN

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Mol	Chain	Res	Type	Atoms
2	С	4	MVA	CB-CA-N-CN
2	D	4	MVA	CB-CA-N-CN
2	С	4	MVA	O-C-CA-CB
2	С	7	SAR	C-CA-N-CN
2	D	7	SAR	C-CA-N-CN
2	С	4	MVA	N-CA-CB-CG2

There are no ring outliers.

12 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	2	MLE	2	0
2	D	5	BMT	1	0
2	D	6	ABA	1	0
2	D	3	MLE	2	0
2	D	1	DAL	2	0
2	D	10	MLE	1	0
2	С	5	BMT	1	0
2	С	10	MLE	1	0
2	D	8	MLE	1	0
2	С	3	MLE	1	0
2	C	1	DAL	2	0
2	С	2	MLE	1	0

### 5.5 Carbohydrates (i)

There are no monosaccharides 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,  $95^{th}$  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>	$\# \mathrm{RSRZ}{>}2$	$OWAB(\AA^2)$	Q < 0.9
1	A	199/211 (94%)	-0.02	4 (2%) 65 60	11, 16, 30, 43	0
1	В	200/211 (94%)	-0.07	3 (1%) 73 69	10, 16, 31, 47	0
2	С	2/11 (18%)	0.52	0 100 100	20, 20, 20, 24	0
2	D	2/11 (18%)	-0.24	0 100 100	16, 16, 16, 19	0
All	All	403/444 (90%)	-0.04	7 (1%) 70 66	10, 16, 30, 47	0

#### All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	В	86	ASN	4.7
1	В	12	SER	3.4
1	A	87	GLN	3.0
1	В	54[A]	HIS	2.9
1	A	13	SER	2.8
1	A	18	SER	2.7
1	A	86	ASN	2.7

### 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,  $95^{th}$  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	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
2	MLE	С	8	9/10	0.94	0.10	16,20,22,25	0
2	MLE	D	8	9/10	0.95	0.08	14,16,24,24	0
2	ABA	D	6	6/7	0.96	0.07	12,13,14,17	0
2	MLE	D	10	9/10	0.96	0.07	14,17,20,20	0
2	MLE	D	2	9/10	0.97	0.08	11,11,13,13	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q < 0.9
2	SAR	D	7	5/6	0.97	0.05	15,15,18,20	0
2	MLE	С	3	9/10	0.97	0.08	12,13,17,17	0
2	MVA	D	4	8/9	0.97	0.09	10,10,11,11	0
2	MLE	С	10	9/10	0.97	0.08	14,17,21,22	0
2	ABA	С	6	6/7	0.97	0.08	11,12,14,17	0
2	DAL	С	1	5/6	0.98	0.06	13,15,16,17	0
2	DAL	D	1	5/6	0.98	0.07	12,13,14,15	0
2	SAR	С	7	5/6	0.98	0.05	15,16,21,27	0
2	MLE	D	3	9/10	0.98	0.07	10,11,14,14	0
2	MVA	С	4	8/9	0.98	0.07	11,12,12,12	0
2	MLE	С	2	9/10	0.98	0.07	13,14,16,20	0
2	BMT	С	5	13/14	0.98	0.08	11,12,15,17	0
2	BMT	D	5	13/14	0.98	0.06	11,12,14,16	0

### 6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

## 6.4 Ligands (i)

There are no ligands in this entry.

### 6.5 Other polymers (i)

There are no such residues in this entry.

