



# wwPDB X-ray Structure Validation Summary Report ⓘ

May 29, 2020 – 07:14 am BST

PDB ID : 5H7D  
Title : Crystal structure of the YgjG-protein A-Zpa963-calmodulin complex  
Authors : Youn, S.J.; Kwon, N.Y.; Lee, J.H.; Kim, J.H.; Lee, H.; Lee, J.O.  
Deposited on : 2016-11-17  
Resolution : 2.57 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	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)	:	2.11

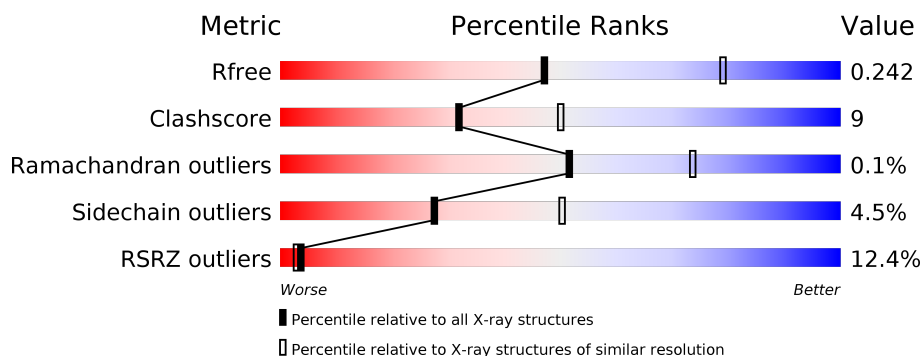
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.57 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))
$R_{free}$	130704	3676 (2.60-2.56)
Clashscore	141614	4049 (2.60-2.56)
Ramachandran outliers	138981	3979 (2.60-2.56)
Sidechain outliers	138945	3979 (2.60-2.56)
RSRZ outliers	127900	3614 (2.60-2.56)

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 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  $\leq 5\%$ . 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	499	<div> <div>5%</div> <div>81%17%..</div> </div>
1	B	499	<div> <div>8%</div> <div>83%15%..</div> </div>
1	C	499	<div> <div>6%</div> <div>80%18%..</div> </div>
1	D	499	<div> <div>6%</div> <div>82%15%..</div> </div>
1	I	499	<div> <div>5%</div> <div>82%16%..</div> </div>
1	J	499	<div> <div>7%</div> <div>81%17%..</div> </div>

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Mol	Chain	Length	Quality of chain
1	M	499	<div><div></div><div>6%</div><div>84%</div><div>14%</div><div>••</div></div>
1	N	499	<div><div></div><div>8%</div><div>83%</div><div>15%</div><div>••</div></div>
2	E	120	<div><div></div><div>25%</div><div>64%</div><div>24%</div><div>6%</div><div>6%</div></div>
2	F	120	<div><div></div><div>49%</div><div>62%</div><div>26%</div><div>5%</div><div>•</div><div>7%</div></div>
2	G	120	<div><div></div><div>18%</div><div>63%</div><div>25%</div><div>6%</div><div>6%</div></div>
2	H	120	<div><div></div><div>50%</div><div>63%</div><div>26%</div><div>•</div><div>7%</div></div>
2	K	120	<div><div></div><div>19%</div><div>62%</div><div>28%</div><div>5%</div><div>6%</div></div>
2	L	120	<div><div></div><div>48%</div><div>68%</div><div>22%</div><div>•</div><div>7%</div></div>
2	O	120	<div><div></div><div>22%</div><div>63%</div><div>25%</div><div>6%</div><div>6%</div></div>
2	P	120	<div><div></div><div>60%</div><div>63%</div><div>22%</div><div>8%</div><div>7%</div></div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 37132 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 Putrescine aminotransferase,Immunoglobulin G-binding protein A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	495	Total	C	N	O	S	0	0	0
			3775	2396	651	706	22			
1	B	495	Total	C	N	O	S	0	0	0
			3775	2396	651	706	22			
1	C	495	Total	C	N	O	S	0	0	0
			3775	2396	651	706	22			
1	D	495	Total	C	N	O	S	0	0	0
			3775	2396	651	706	22			
1	I	495	Total	C	N	O	S	0	0	0
			3775	2396	651	706	22			
1	J	495	Total	C	N	O	S	0	0	0
			3775	2396	651	706	22			
1	M	495	Total	C	N	O	S	0	0	0
			3775	2396	651	706	22			
1	N	495	Total	C	N	O	S	0	0	0
			3775	2396	651	706	22			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	3	GLY	-	expression tag	UNP P42588
A	4	SER	-	expression tag	UNP P42588
A	5	HIS	-	expression tag	UNP P42588
A	6	MET	-	expression tag	UNP P42588
A	456	VAL	ASN	engineered mutation	UNP P38507
A	474	ALA	GLY	engineered mutation	UNP P38507
B	3	GLY	-	expression tag	UNP P42588
B	4	SER	-	expression tag	UNP P42588
B	5	HIS	-	expression tag	UNP P42588
B	6	MET	-	expression tag	UNP P42588
B	456	VAL	ASN	engineered mutation	UNP P38507
B	474	ALA	GLY	engineered mutation	UNP P38507

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Chain	Residue	Modelled	Actual	Comment	Reference
C	3	GLY	-	expression tag	UNP P42588
C	4	SER	-	expression tag	UNP P42588
C	5	HIS	-	expression tag	UNP P42588
C	6	MET	-	expression tag	UNP P42588
C	456	VAL	ASN	engineered mutation	UNP P38507
C	474	ALA	GLY	engineered mutation	UNP P38507
D	3	GLY	-	expression tag	UNP P42588
D	4	SER	-	expression tag	UNP P42588
D	5	HIS	-	expression tag	UNP P42588
D	6	MET	-	expression tag	UNP P42588
D	456	VAL	ASN	engineered mutation	UNP P38507
D	474	ALA	GLY	engineered mutation	UNP P38507
I	3	GLY	-	expression tag	UNP P42588
I	4	SER	-	expression tag	UNP P42588
I	5	HIS	-	expression tag	UNP P42588
I	6	MET	-	expression tag	UNP P42588
I	456	VAL	ASN	engineered mutation	UNP P38507
I	474	ALA	GLY	engineered mutation	UNP P38507
J	3	GLY	-	expression tag	UNP P42588
J	4	SER	-	expression tag	UNP P42588
J	5	HIS	-	expression tag	UNP P42588
J	6	MET	-	expression tag	UNP P42588
J	456	VAL	ASN	engineered mutation	UNP P38507
J	474	ALA	GLY	engineered mutation	UNP P38507
M	3	GLY	-	expression tag	UNP P42588
M	4	SER	-	expression tag	UNP P42588
M	5	HIS	-	expression tag	UNP P42588
M	6	MET	-	expression tag	UNP P42588
M	456	VAL	ASN	engineered mutation	UNP P38507
M	474	ALA	GLY	engineered mutation	UNP P38507
N	3	GLY	-	expression tag	UNP P42588
N	4	SER	-	expression tag	UNP P42588
N	5	HIS	-	expression tag	UNP P42588
N	6	MET	-	expression tag	UNP P42588
N	456	VAL	ASN	engineered mutation	UNP P38507
N	474	ALA	GLY	engineered mutation	UNP P38507

- Molecule 2 is a protein called Zpa963, Calmodulin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	113	Total	C	N	O	S	0	0	0
			872	544	143	181	4			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	F	112	Total 857	C 536	N 139	O 178	S 4	0	0	0
2	G	113	Total 872	C 544	N 143	O 181	S 4	0	0	0
2	H	112	Total 857	C 536	N 139	O 178	S 4	0	0	0
2	K	113	Total 872	C 544	N 143	O 181	S 4	0	0	0
2	L	112	Total 857	C 536	N 139	O 178	S 4	0	0	0
2	O	113	Total 872	C 544	N 143	O 181	S 4	0	0	0
2	P	112	Total 857	C 536	N 139	O 178	S 4	0	0	0

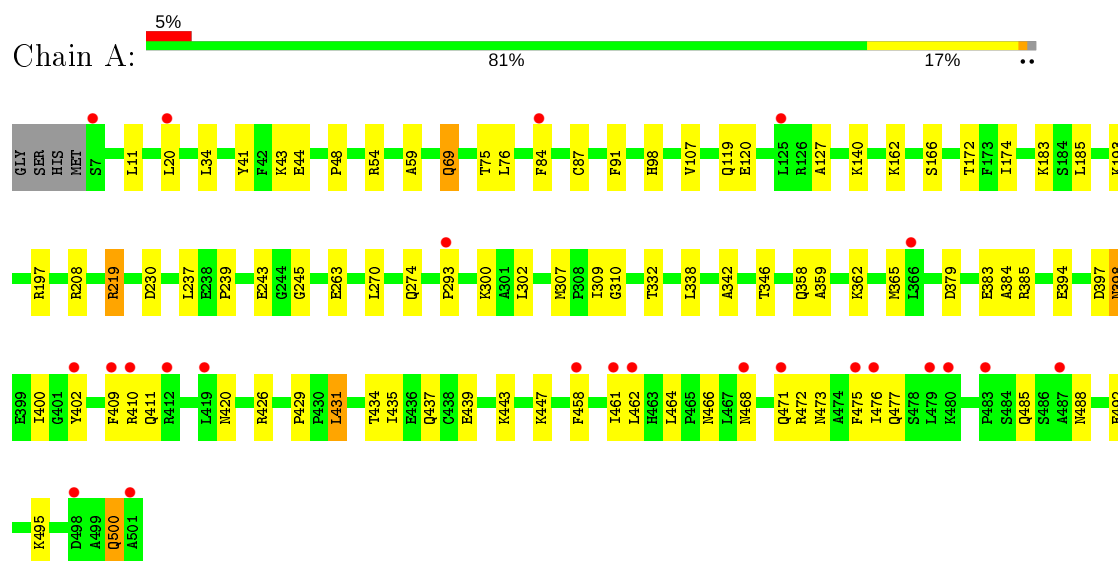
- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	P	2	Total 2	Ca 2	0	0
3	G	2	Total 2	Ca 2	0	0
3	K	2	Total 2	Ca 2	0	0
3	E	2	Total 2	Ca 2	0	0
3	H	2	Total 2	Ca 2	0	0
3	O	2	Total 2	Ca 2	0	0
3	L	2	Total 2	Ca 2	0	0
3	F	2	Total 2	Ca 2	0	0

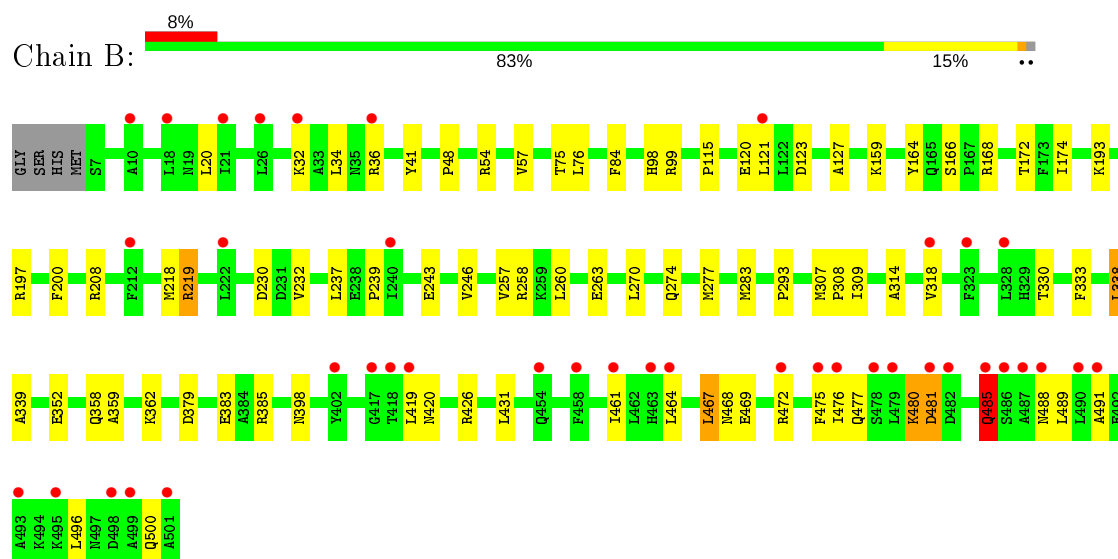
### 3 Residue-property plots [i](#)

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.

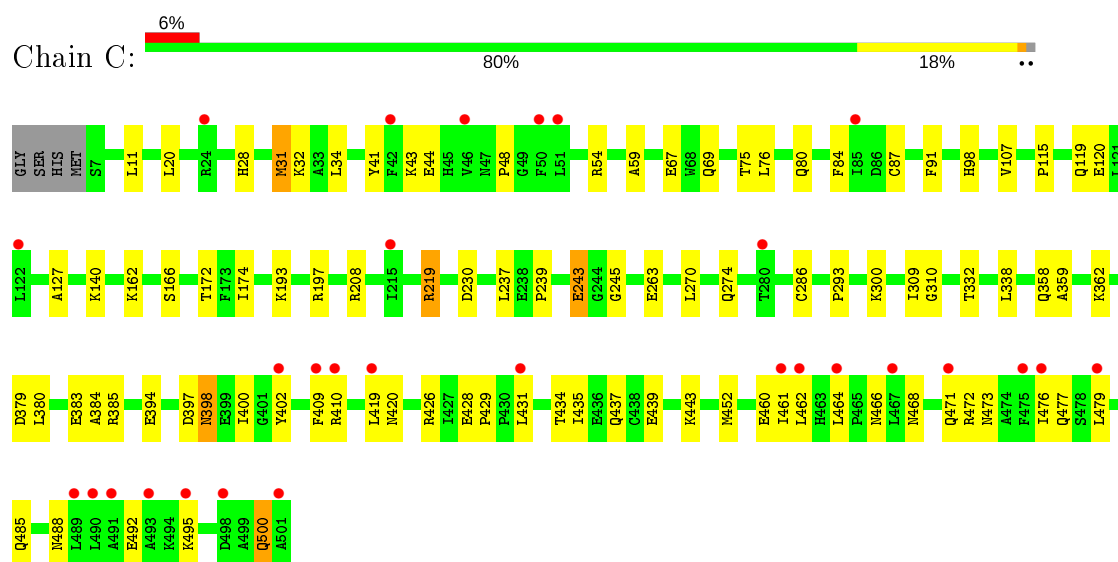
- Molecule 1: Putrescine aminotransferase,Immunoglobulin G-binding protein A



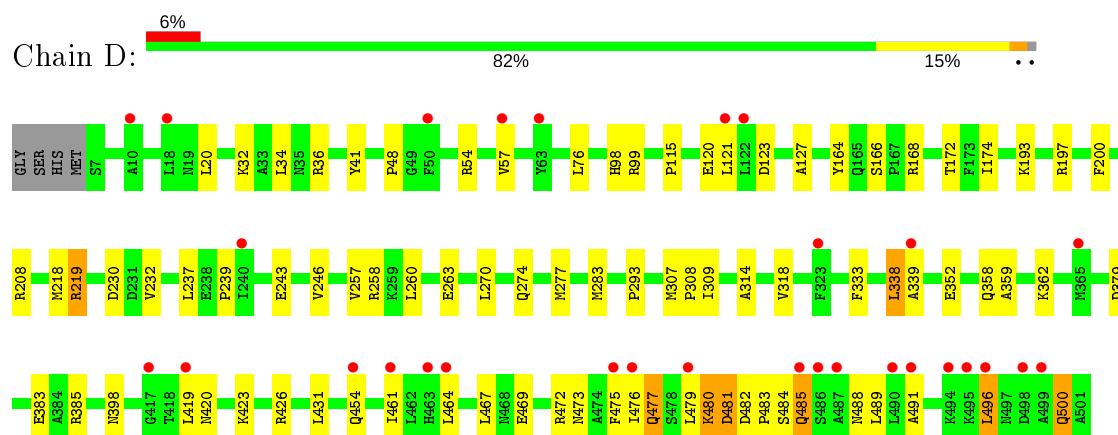
- Molecule 1: Putrescine aminotransferase,Immunoglobulin G-binding protein A



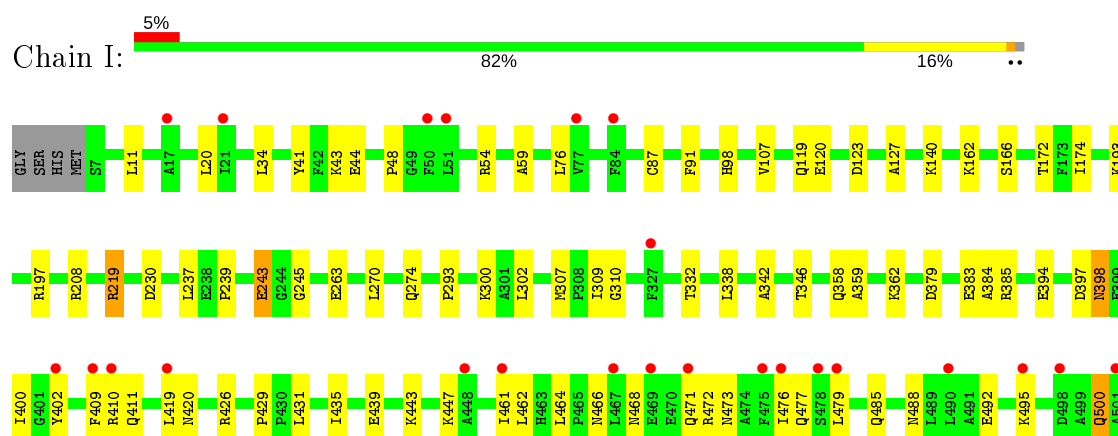
- Molecule 1: Putrescine aminotransferase,Immunoglobulin G-binding protein A



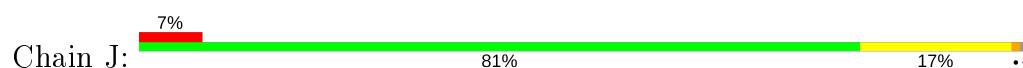
- Molecule 1: Putrescine aminotransferase,Immunoglobulin G-binding protein A



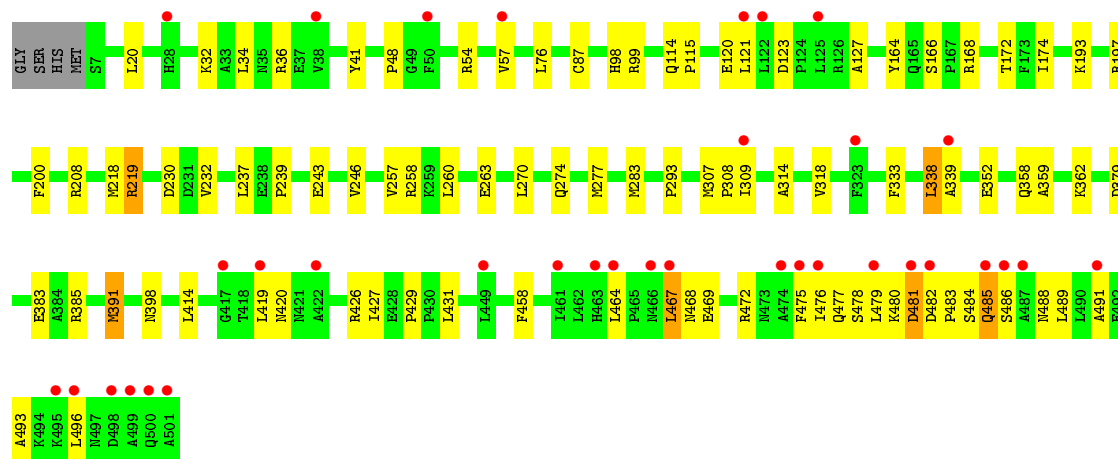
- Molecule 1: Putrescine aminotransferase,Immunoglobulin G-binding protein A



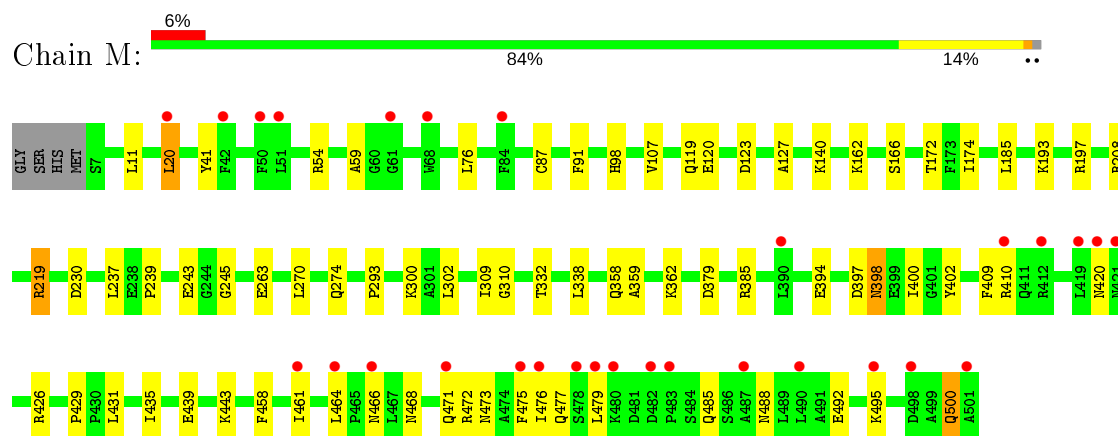
- Molecule 1: Putrescine aminotransferase,Immunoglobulin G-binding protein A



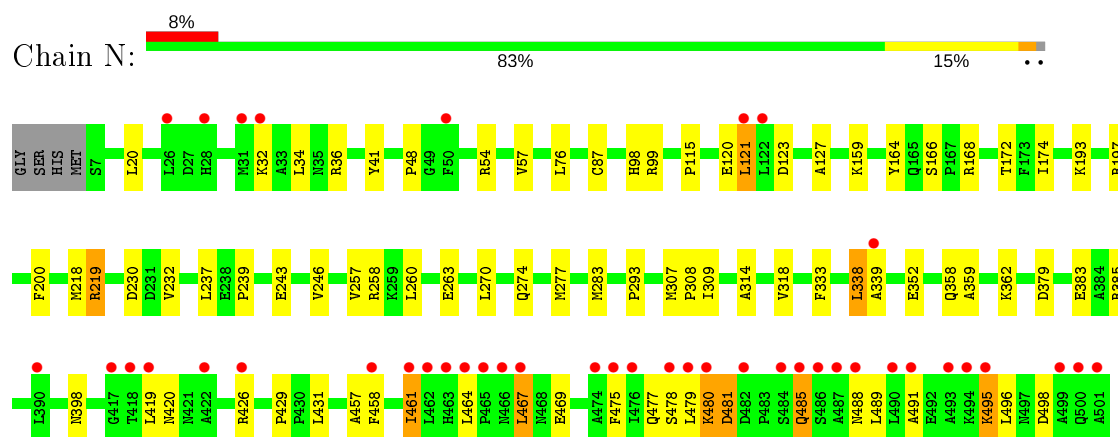




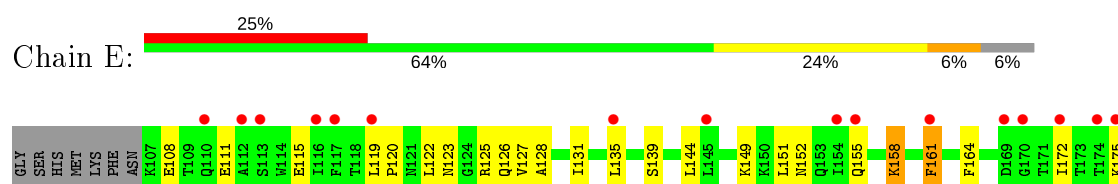
- Molecule 1: Putrescine aminotransferase,Immunoglobulin G-binding protein A

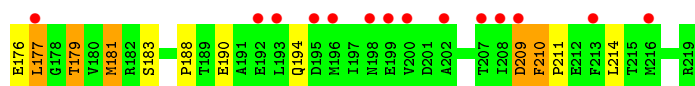


- Molecule 1: Putrescine aminotransferase,Immunoglobulin G-binding protein A

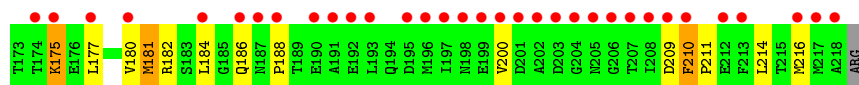
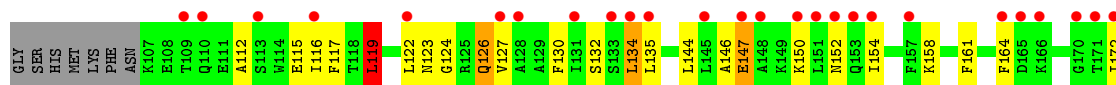


- Molecule 2: Zpa963,Calmodulin

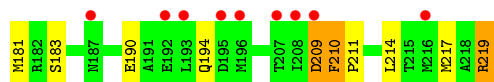
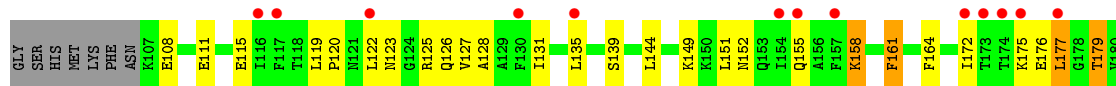




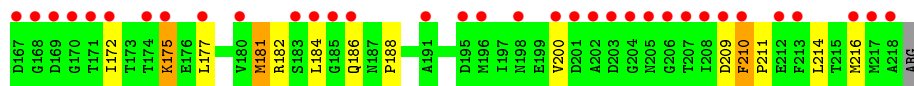
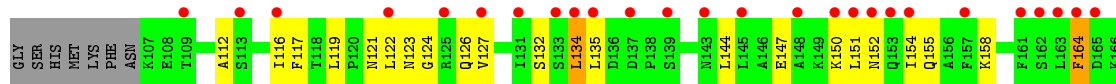
• Molecule 2: Zpa963,Calmodulin



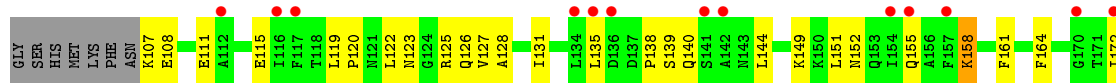
• Molecule 2: Zpa963,Calmodulin



• Molecule 2: Zpa963,Calmodulin

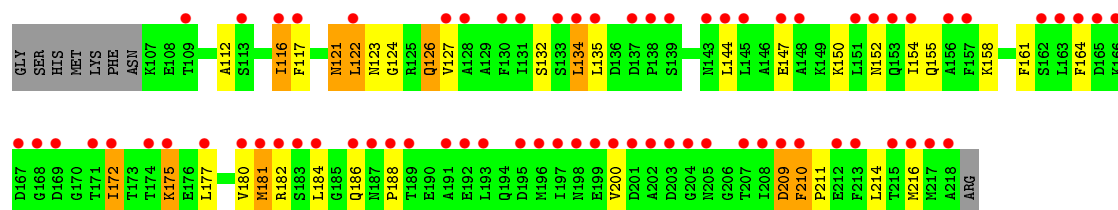


• Molecule 2: Zpa963,Calmodulin



• Molecule 2: Zpa963,Calmodulin





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	79.49Å 155.38Å 155.49Å 84.84° 89.68° 89.56°	Depositor
Resolution (Å)	50.00 – 2.57 47.70 – 2.57	Depositor EDS
% Data completeness (in resolution range)	89.5 (50.00-2.57) 89.4 (47.70-2.57)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.64 (at 2.58Å)	Xtriage
Refinement program	REFMAC 5.8.0155	Depositor
R, $R_{free}$	0.204 , 0.241 0.204 , 0.242	Depositor DCC
$R_{free}$ test set	21075 reflections (10.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	62.1	Xtriage
Anisotropy	0.014	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 38.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.397 for h,-k,-l 0.387 for -h,l,k 0.439 for -h,-l,-k	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	37132	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	88.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.72% 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: 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	0/3845	0.52	0/5203
1	B	0.47	0/3845	0.57	0/5203
1	C	0.45	1/3845 (0.0%)	0.54	1/5203 (0.0%)
1	D	0.47	0/3845	0.56	0/5203
1	I	0.44	0/3845	0.54	0/5203
1	J	0.46	0/3845	0.57	0/5203
1	M	0.44	0/3845	0.53	1/5203 (0.0%)
1	N	0.47	0/3845	0.56	0/5203
2	E	0.58	0/884	0.69	0/1194
2	F	0.66	0/869	0.86	3/1175 (0.3%)
2	G	0.58	0/884	0.68	0/1194
2	H	0.65	0/869	0.82	1/1175 (0.1%)
2	K	0.57	0/884	0.69	0/1194
2	L	0.63	0/869	0.84	2/1175 (0.2%)
2	O	0.57	0/884	0.70	0/1194
2	P	0.63	0/869	0.84	1/1175 (0.1%)
All	All	0.49	1/37772 (0.0%)	0.60	9/51100 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	460	GLU	CD-OE2	5.33	1.31	1.25

The worst 5 of 9 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	31	MET	CG-SD-CE	-7.73	87.83	100.20
2	F	119	LEU	CB-CG-CD2	5.82	120.89	111.00
2	P	184	LEU	CA-CB-CG	5.53	128.01	115.30
2	L	163	LEU	CB-CG-CD1	5.52	120.38	111.00
2	L	184	LEU	CA-CB-CG	5.51	127.97	115.30

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	3775	0	3801	61	0
1	B	3775	0	3801	60	0
1	C	3775	0	3801	65	0
1	D	3775	0	3801	63	0
1	I	3775	0	3801	58	0
1	J	3775	0	3801	61	0
1	M	3775	0	3801	54	0
1	N	3775	0	3801	56	0
2	E	872	0	846	32	0
2	F	857	0	829	35	0
2	G	872	0	846	32	0
2	H	857	0	829	31	0
2	K	872	0	846	32	0
2	L	857	0	829	23	0
2	O	872	0	846	34	0
2	P	857	0	829	34	0
3	E	2	0	0	0	0
3	F	2	0	0	0	0
3	G	2	0	0	0	0
3	H	2	0	0	0	0
3	K	2	0	0	0	0
3	L	2	0	0	0	0
3	O	2	0	0	0	0
3	P	2	0	0	0	0
All	All	37132	0	37108	660	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 9.

The worst 5 of 660 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:461:ILE:HD11	1:A:475:PHE:HB3	1.28	1.14
1:M:461:ILE:HD11	1:M:475:PHE:HB3	1.30	1.12
1:D:485:GLN:HB2	1:D:489:LEU:HG	1.30	1.11
1:B:461:ILE:HD11	1:B:475:PHE:HB3	1.26	1.08
1:N:485:GLN:HB2	1:N:489:LEU:HG	1.35	1.07

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	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	493/499 (99%)	472 (96%)	21 (4%)	0	100	100
1	B	493/499 (99%)	470 (95%)	22 (4%)	1 (0%)	47	69
1	C	493/499 (99%)	473 (96%)	19 (4%)	1 (0%)	47	69
1	D	493/499 (99%)	470 (95%)	23 (5%)	0	100	100
1	I	493/499 (99%)	472 (96%)	21 (4%)	0	100	100
1	J	493/499 (99%)	469 (95%)	24 (5%)	0	100	100
1	M	493/499 (99%)	472 (96%)	21 (4%)	0	100	100
1	N	493/499 (99%)	470 (95%)	23 (5%)	0	100	100
2	E	111/120 (92%)	105 (95%)	5 (4%)	1 (1%)	17	34
2	F	110/120 (92%)	103 (94%)	7 (6%)	0	100	100
2	G	111/120 (92%)	104 (94%)	6 (5%)	1 (1%)	17	34
2	H	110/120 (92%)	102 (93%)	8 (7%)	0	100	100
2	K	111/120 (92%)	105 (95%)	5 (4%)	1 (1%)	17	34
2	L	110/120 (92%)	103 (94%)	7 (6%)	0	100	100
2	O	111/120 (92%)	104 (94%)	6 (5%)	1 (1%)	17	34
2	P	110/120 (92%)	103 (94%)	7 (6%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	4828/4952 (98%)	4597 (95%)	225 (5%)	6 (0%)	51 73

5 of 6 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	485	GLN
2	E	139	SER
2	G	139	SER
2	K	139	SER
2	O	139	SER

### 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	396/399 (99%)	385 (97%)	11 (3%)	43 67
1	B	396/399 (99%)	382 (96%)	14 (4%)	36 59
1	C	396/399 (99%)	386 (98%)	10 (2%)	47 70
1	D	396/399 (99%)	380 (96%)	16 (4%)	31 55
1	I	396/399 (99%)	387 (98%)	9 (2%)	50 73
1	J	396/399 (99%)	382 (96%)	14 (4%)	36 59
1	M	396/399 (99%)	388 (98%)	8 (2%)	55 76
1	N	396/399 (99%)	380 (96%)	16 (4%)	31 55
2	E	95/101 (94%)	84 (88%)	11 (12%)	5 9
2	F	93/101 (92%)	85 (91%)	8 (9%)	10 19
2	G	95/101 (94%)	82 (86%)	13 (14%)	3 6
2	H	93/101 (92%)	86 (92%)	7 (8%)	13 26
2	K	95/101 (94%)	84 (88%)	11 (12%)	5 9
2	L	93/101 (92%)	86 (92%)	7 (8%)	13 26
2	O	95/101 (94%)	84 (88%)	11 (12%)	5 9
2	P	93/101 (92%)	82 (88%)	11 (12%)	5 9

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	3920/4000 (98%)	3743 (96%)	177 (4%)	27 50

5 of 177 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	G	210	PHE
1	J	48	PRO
2	O	183	SER
2	G	219	ARG
2	H	210	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 34 such sidechains are listed below:

Mol	Chain	Res	Type
2	G	110	GLN
2	H	126	GLN
2	O	155	GLN
2	G	152	ASN
2	E	152	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 16 ligands modelled in this entry, 16 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 ⓘ

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(Å²)	Q<0.9
1	A	495/499 (99%)	0.61	24 (4%)	30	26	45, 69, 137, 163	0
1	B	495/499 (99%)	0.77	40 (8%)	12	9	43, 62, 126, 159	0
1	C	495/499 (99%)	0.61	29 (5%)	22	19	42, 65, 86, 98	0
1	D	495/499 (99%)	0.75	30 (6%)	21	17	44, 62, 129, 162	0
1	I	495/499 (99%)	0.51	24 (4%)	30	26	44, 68, 130, 155	0
1	J	495/499 (99%)	0.62	35 (7%)	16	13	44, 61, 128, 162	0
1	M	495/499 (99%)	0.57	29 (5%)	22	19	46, 69, 137, 164	0
1	N	495/499 (99%)	0.74	42 (8%)	10	8	44, 63, 128, 160	0
2	E	113/120 (94%)	1.39	30 (26%)	0	0	103, 125, 147, 155	0
2	F	112/120 (93%)	3.49	59 (52%)	0	0	163, 203, 250, 261	0
2	G	113/120 (94%)	1.04	22 (19%)	1	0	97, 119, 151, 160	0
2	H	112/120 (93%)	2.63	60 (53%)	0	0	147, 183, 233, 246	0
2	K	113/120 (94%)	1.18	23 (20%)	1	0	97, 117, 142, 153	0
2	L	112/120 (93%)	2.96	57 (50%)	0	0	166, 201, 241, 251	0
2	O	113/120 (94%)	1.21	26 (23%)	0	0	103, 127, 156, 162	0
2	P	112/120 (93%)	3.03	72 (64%)	0	0	170, 192, 227, 237	0
All	All	4860/4952 (98%)	0.92	602 (12%)	4	3	42, 70, 188, 261	0

The worst 5 of 602 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	L	208	ILE	21.6
2	F	200	VAL	20.5
2	F	208	ILE	16.9
2	P	208	ILE	16.2
2	F	205	ASN	14.1

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

There are no non-standard protein/DNA/RNA residues in this entry.

## 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 6.4 Ligands [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( $\text{\AA}^2$ )	Q<0.9
3	CA	O	502	1/1	0.56	0.12	88,88,88,88	0
3	CA	E	502	1/1	0.77	0.13	86,86,86,86	0
3	CA	P	502	1/1	0.83	0.41	84,84,84,84	0
3	CA	O	501	1/1	0.83	0.10	79,79,79,79	0
3	CA	H	502	1/1	0.83	0.54	83,83,83,83	0
3	CA	G	502	1/1	0.83	0.18	84,84,84,84	0
3	CA	G	501	1/1	0.83	0.10	82,82,82,82	0
3	CA	P	501	1/1	0.85	0.24	89,89,89,89	0
3	CA	K	501	1/1	0.85	0.09	80,80,80,80	0
3	CA	L	501	1/1	0.87	0.22	88,88,88,88	0
3	CA	K	502	1/1	0.89	0.20	86,86,86,86	0
3	CA	H	501	1/1	0.90	0.22	90,90,90,90	0
3	CA	F	501	1/1	0.91	0.19	91,91,91,91	0
3	CA	E	501	1/1	0.91	0.12	82,82,82,82	0
3	CA	L	502	1/1	0.94	0.52	82,82,82,82	0
3	CA	F	502	1/1	0.94	0.53	87,87,87,87	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.