



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 01:49 PM GMT

PDB ID : 1PIP
Title : CRYSTAL STRUCTURE OF PAPAIN-SUCCINYL-GLN-VAL-VAL-ALA-ALA-P-NITROANILIDE COMPLEX AT 1.7 ANGSTROMS RESOLUTION: NONCOVALENT BINDING MODE OF A COMMON SEQUENCE OF ENDOGENOUS THIOL PROTEASE INHIBITORS
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Deposited on : 1992-10-03
Resolution : 1.70 Å (reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

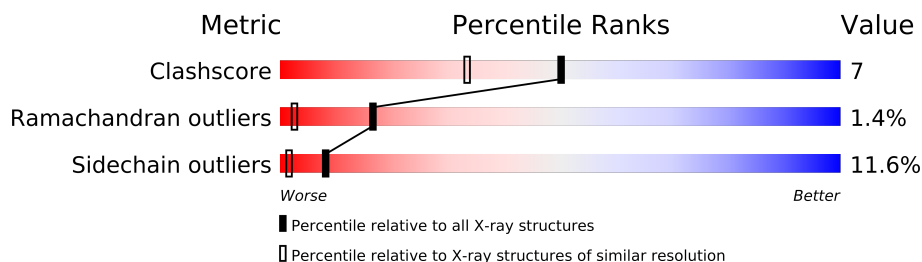
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.15 2013
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	21963
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP)	:	stable22683

1 Overall quality at a glance

The reported resolution of this entry is 1.70 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	79885	2929 (1.70-1.70)
Ramachandran outliers	78287	2878 (1.70-1.70)
Sidechain outliers	78261	2878 (1.70-1.70)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	212	
2	B	7	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 1705 atoms, of which 0 are hydrogen and 0 are deuterium.

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 PAPAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	212	Total	C	N	O	S	0	0	0
			1655	1050	293	305	7			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	47	GLN	GLU	CONFLICT	UNP P00784
A	118	GLN	GLU	CONFLICT	UNP P00784
A	135	GLN	GLU	CONFLICT	UNP P00784

- Molecule 2 is a protein called SUCCINYL-GLN-VAL-VAL-ALA-ALA-P-NITROANILIDE.

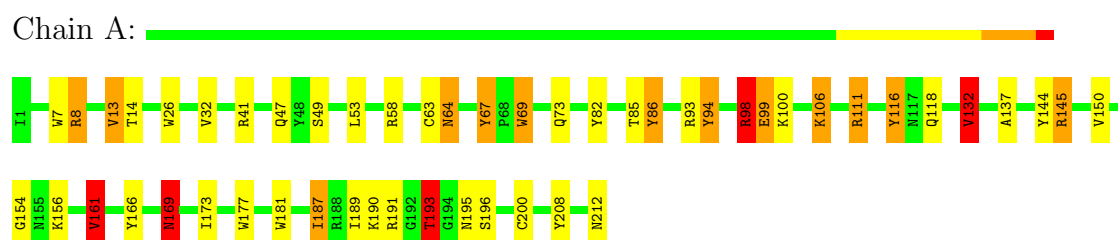
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	7	Total	C	N	O	0	0	0
			50	31	8	11			

3 Residue-property plots

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: PAPAIN



• Molecule 2: SUCCINYL-GLN-VAL-VAL-ALA-ALA-P-NITROANILIDE



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	43.09Å 102.30Å 49.69Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 1.70	Depositor
% Data completeness (in resolution range)	(Not available) (10.00-1.70)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.196 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	1705	wwPDB-VP
Average B, all atoms (Å ²)	15.0	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: SIN, NIT

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	0.90	0/1699	1.63	37/2306 (1.6%)
2	B	0.83	0/32	3.10	4/43 (9.3%)
All	All	0.90	0/1731	1.67	41/2349 (1.7%)

There are no bond length outliers.

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	98	ARG	NE-CZ-NH1	8.96	124.78	120.30
1	A	69	TRP	CD1-CG-CD2	8.33	112.96	106.30
2	B	401	VAL	CA-CB-CG2	-8.20	98.60	110.90
1	A	7	TRP	CD1-CG-CD2	7.57	112.35	106.30
1	A	111	ARG	NE-CZ-NH2	-7.45	116.57	120.30
1	A	116	TYR	CB-CG-CD2	-7.09	116.75	121.00
1	A	181	TRP	CD1-CG-CD2	6.89	111.81	106.30
1	A	69	TRP	CE2-CD2-CG	-6.89	101.79	107.30
1	A	7	TRP	CE2-CD2-CG	-6.84	101.83	107.30
1	A	177	TRP	CA-C-N	6.80	129.81	116.20
1	A	98	ARG	NE-CZ-NH2	-6.80	116.90	120.30
1	A	145	ARG	NE-CZ-NH1	6.79	123.69	120.30
1	A	177	TRP	O-C-N	-6.70	111.82	123.20
1	A	94	TYR	CA-CB-CG	6.66	126.05	113.40
1	A	64	ASN	CA-CB-CG	6.66	128.05	113.40
1	A	177	TRP	CD1-CG-CD2	6.60	111.58	106.30
2	B	403	ALA	N-CA-C	6.42	128.33	111.00
1	A	8	ARG	NE-CZ-NH2	-6.40	117.10	120.30
1	A	161	VAL	CG1-CB-CG2	6.29	120.97	110.90
1	A	177	TRP	CE2-CD2-CG	-6.20	102.34	107.30
1	A	132	VAL	N-CA-CB	-6.08	98.13	111.50
1	A	26	TRP	CD1-CG-CD2	6.04	111.13	106.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	63	CYS	O-C-N	-6.01	113.08	122.70
1	A	181	TRP	CE2-CD2-CG	-5.99	102.51	107.30
1	A	26	TRP	CE2-CD2-CG	-5.86	102.61	107.30
1	A	208	TYR	CB-CG-CD2	-5.83	117.50	121.00
1	A	161	VAL	N-CA-CB	-5.76	98.83	111.50
2	B	401	VAL	CA-CB-CG1	5.67	119.40	110.90
1	A	7	TRP	CG-CD1-NE1	-5.66	104.44	110.10
1	A	67	TYR	CB-CG-CD1	-5.64	117.61	121.00
1	A	193	THR	N-CA-CB	-5.60	99.65	110.30
1	A	166	TYR	CB-CG-CD2	-5.56	117.66	121.00
1	A	181	TRP	CG-CD1-NE1	-5.56	104.54	110.10
1	A	93	ARG	NE-CZ-NH1	5.54	123.07	120.30
1	A	69	TRP	CG-CD1-NE1	-5.29	104.81	110.10
1	A	106	LYS	CA-CB-CG	5.20	124.84	113.40
1	A	169	ASN	CB-CG-ND2	5.13	129.01	116.70
1	A	145	ARG	NE-CZ-NH2	-5.10	117.75	120.30
1	A	94	TYR	CB-CG-CD1	5.08	124.05	121.00
2	B	400	GLN	O-C-N	5.04	130.77	122.70
1	A	49	SER	N-CA-CB	-5.01	102.99	110.50

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1655	0	1596	24	1
2	B	50	0	43	7	1
All	All	1705	0	1639	25	1

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 7.

All (25) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:137:ALA:HA	2:B:219:NIT:ON1	1.49	1.09
1:A:156:LYS:HE3	2:B:213:SIN:H22	1.44	1.00
1:A:144:TYR:CE2	1:A:187:ILE:HG12	2.33	0.64
1:A:173:ILE:HD12	1:A:189:ILE:HG13	1.85	0.58
1:A:156:LYS:CE	2:B:213:SIN:H22	2.25	0.56
1:A:132:VAL:HG12	1:A:161:VAL:HG12	1.88	0.56
1:A:82:TYR:HB2	1:A:85:THR:HG23	1.90	0.52
2:B:400:GLN:O	2:B:401:VAL:HB	2.09	0.52
1:A:150:VAL:HG12	1:A:190:LYS:HG3	1.92	0.51
1:A:137:ALA:HB2	2:B:401:VAL:HG21	1.93	0.51
1:A:118:GLN:NE2	1:A:191:ARG:HH11	2.09	0.50
1:A:137:ALA:CB	2:B:401:VAL:HG21	2.42	0.50
1:A:98:ARG:HB2	1:A:98:ARG:HH11	1.79	0.47
1:A:137:ALA:CA	2:B:219:NIT:ON1	2.42	0.47
1:A:161:VAL:HG13	1:A:173:ILE:HG23	1.97	0.46
1:A:116:TYR:HB3	1:A:196:SER:OG	2.15	0.46
1:A:169:ASN:HD22	1:A:169:ASN:H	1.63	0.45
1:A:154:GLY:O	1:A:200:CYS:HA	2.18	0.43
1:A:144:TYR:CZ	1:A:187:ILE:HG12	2.54	0.41
1:A:14:THR:HG21	1:A:47:GLN:HG2	2.02	0.41
1:A:67:TYR:HB3	1:A:69:TRP:NE1	2.36	0.41
1:A:99:GLU:HG2	1:A:100:LYS:HG3	2.02	0.41
1:A:193:THR:HG22	1:A:195:ASN:H	1.86	0.41
1:A:86:TYR:CD2	1:A:100:LYS:HD2	2.56	0.40
1:A:8:ARG:HA	1:A:13:VAL:HG13	2.03	0.40

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	Distance(Å)	Clash(Å)
1:A:73:GLN:CG	2:B:213:SIN:C1[2.765]	2.06	0.14

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

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	210/212 (99%)	203 (97%)	7 (3%)	0	100	100
2	B	3/7 (43%)	0	0	3 (100%)	0	0
All	All	213/219 (97%)	203 (95%)	7 (3%)	3 (1%)	16	2

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	401	VAL
2	B	403	ALA
2	B	402	VAL

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	170/170 (100%)	151 (89%)	19 (11%)	9	2
2	B	3/3 (100%)	2 (67%)	1 (33%)	0	0
All	All	173/173 (100%)	153 (88%)	20 (12%)	8	1

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	VAL
1	A	32	VAL
1	A	41	ARG
1	A	53	LEU
1	A	58	ARG
1	A	64	ASN
1	A	86	TYR
1	A	94	TYR
1	A	98	ARG
1	A	99	GLU
1	A	106	LYS
1	A	111	ARG
1	A	132	VAL
1	A	145	ARG

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Mol	Chain	Res	Type
1	A	161	VAL
1	A	169	ASN
1	A	187	ILE
1	A	193	THR
1	A	212	ASN
2	B	401	VAL

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	64	ASN
1	A	169	ASN
1	A	212	ASN

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

There are no ligands in this entry.

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.