



Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2020 – 12:56 am BST

PDB ID : 2RD4
Title : Design of specific inhibitors of Phospholipase A2: Crystal structure of the complex of phospholipase A2 with pentapeptide Leu-Val-Phe-Phe-Ala at 2.9 Å resolution
Authors : Mirza, Z.; Kaur, A.; Singh, N.; Sinha, M.; Sharma, S.; Srinivasan, A.; Kaur, P.; Singh, T.P.
Deposited on : 2007-09-21
Resolution : 2.97 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 ⓘ symbol.

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

MolProbity : 4.02b-467
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

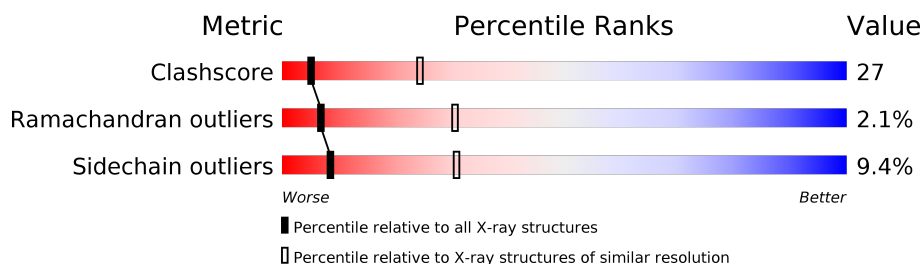
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.97 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	141614	3103 (3.00-2.96)
Ramachandran outliers	138981	2993 (3.00-2.96)
Sidechain outliers	138945	2996 (3.00-2.96)

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 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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	119	
2	B	119	
3	C	5	

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 1940 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 Phospholipase A2 isoform 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	119	Total	C	N	O	S	0	0	0
			916	553	166	182	15			

- Molecule 2 is a protein called Phospholipase A2 isoform 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	119	Total	C	N	O	S	0	0	0
			922	557	168	182	15			

- Molecule 3 is a protein called pentapeptide inhibitor.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	5	Total	C	N	O	0	0	0
			42	32	5	5			

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Ca	0	0
			1	1		
4	A	1	Total	Ca	0	0
			1	1		

- Molecule 5 is water.

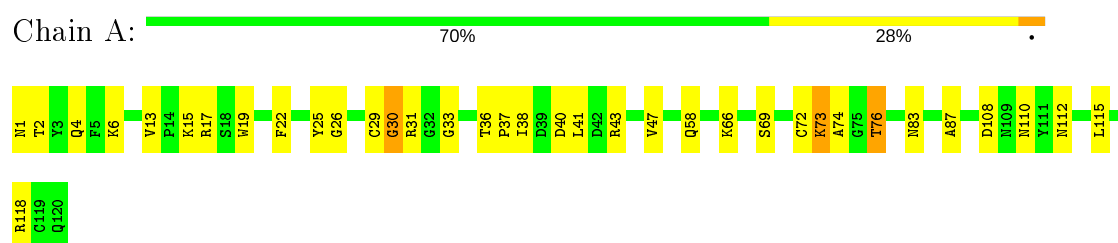
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	30	Total	O	0	0
			30	30		
5	B	28	Total	O	0	0
			28	28		

3 Residue-property plots

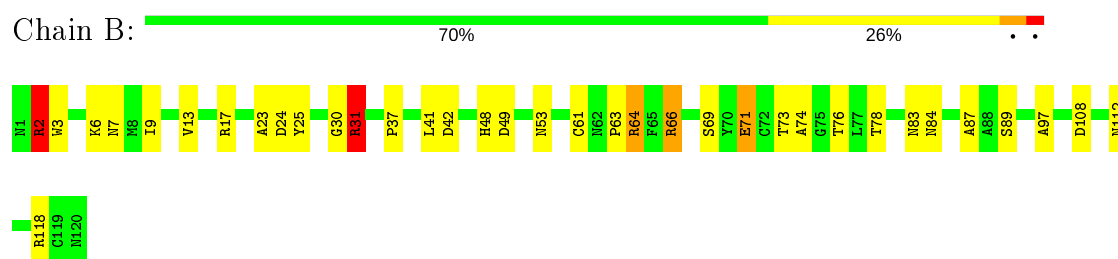
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded according 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. 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: Phospholipase A2 isoform 1



- Molecule 2: Phospholipase A2 isoform 2



- Molecule 3: pentapeptide inhibitor



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α , β , γ	65.54Å 65.54Å 58.55Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.97	Depositor
% Data completeness (in resolution range)	98.7 (20.00-2.97)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.18	Depositor
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.251 , 0.274	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	1940	wwPDB-VP
Average B, all atoms (Å ²)	30.0	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:
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	0.36	0/935	0.63	0/1262
2	B	0.37	0/943	0.63	0/1279
3	C	2.11	2/43 (4.7%)	5.65	13/57 (22.8%)
All	All	0.48	2/1921 (0.1%)	1.05	13/2598 (0.5%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
2	B	0	1
3	C	0	4
All	All	0	5

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	5	PHE	C-N	9.37	1.55	1.34
3	C	3	VAL	CA-CB	5.77	1.66	1.54

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	2	LEU	CB-CA-C	18.72	145.77	110.20
3	C	3	VAL	CA-C-N	15.45	151.19	117.20
3	C	3	VAL	CA-C-O	-14.83	88.95	120.10
3	C	2	LEU	CA-CB-CG	13.94	147.37	115.30
3	C	3	VAL	O-C-N	-12.91	102.05	122.70
3	C	5	PHE	C-N-CA	10.52	147.99	121.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	4	PHE	N-CA-CB	10.23	129.02	110.60
3	C	2	LEU	N-CA-C	-8.57	87.86	111.00
3	C	2	LEU	O-C-N	-8.45	109.19	122.70
3	C	6	ALA	N-CA-CB	6.22	118.81	110.10
3	C	3	VAL	N-CA-CB	5.56	123.74	111.50
3	C	5	PHE	O-C-N	5.38	131.31	122.70
3	C	5	PHE	CA-C-N	-5.17	105.83	117.20

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	2	ARG	Sidechain
3	C	2	LEU	Mainchain
3	C	3	VAL	Mainchain,Peptide
3	C	4	PHE	Peptide

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	916	0	827	34	0
2	B	922	0	818	65	0
3	C	42	0	42	27	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	30	0	0	4	0
5	B	28	0	0	2	0
All	All	1940	0	1687	96	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 27.

All (96) 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)
2:B:2:ARG:HB2	2:B:2:ARG:NH1	1.27	1.46
1:A:33:GLY:HA3	2:B:31:ARG:HH12	1.07	1.14
2:B:2:ARG:HG2	3:C:2:LEU:HD22	1.21	1.13
2:B:2:ARG:CB	2:B:2:ARG:NH1	2.16	1.07
2:B:31:ARG:HH11	3:C:4:PHE:HB3	1.24	0.97
2:B:2:ARG:CG	3:C:2:LEU:HD22	1.94	0.97
2:B:9:ILE:HD12	3:C:5:PHE:HZ	1.32	0.94
1:A:33:GLY:HA3	2:B:31:ARG:NH1	1.83	0.92
2:B:2:ARG:HB2	2:B:2:ARG:CZ	2.04	0.87
2:B:118:ARG:HG2	2:B:118:ARG:HH11	1.37	0.87
2:B:2:ARG:NH1	2:B:64:ARG:HG3	1.89	0.86
1:A:73:LYS:HA	1:A:73:LYS:HE2	1.59	0.85
2:B:2:ARG:HB2	2:B:2:ARG:HH11	1.04	0.84
2:B:31:ARG:HD2	3:C:3:VAL:HG12	1.57	0.84
1:A:2:THR:HG22	5:A:429:HOH:O	1.78	0.81
2:B:2:ARG:CZ	2:B:2:ARG:CB	2.57	0.81
2:B:3:TRP:CZ3	2:B:7:ASN:ND2	2.48	0.80
3:C:4:PHE:CD1	3:C:5:PHE:HB3	2.17	0.80
2:B:48:HIS:ND1	3:C:6:ALA:HB1	2.00	0.77
1:A:1:ASN:OD1	1:A:4:GLN:HG3	1.86	0.75
2:B:2:ARG:HG2	3:C:2:LEU:CD2	2.09	0.75
2:B:9:ILE:HD12	3:C:5:PHE:CZ	2.20	0.75
2:B:2:ARG:CZ	2:B:2:ARG:HA	2.19	0.73
1:A:38:ILE:HD12	1:A:110:ASN:CG	2.10	0.72
2:B:31:ARG:NH1	3:C:4:PHE:HB3	2.02	0.72
2:B:6:LYS:HZ1	3:C:4:PHE:HZ	1.38	0.71
2:B:49:ASP:HB3	5:B:410:HOH:O	1.90	0.70
2:B:37:PRO:HA	2:B:42:ASP:OD2	1.92	0.69
1:A:38:ILE:CD1	1:A:110:ASN:HB3	2.23	0.69
2:B:48:HIS:ND1	3:C:6:ALA:CB	2.55	0.69
1:A:38:ILE:HD12	1:A:110:ASN:HB3	1.76	0.68
2:B:31:ARG:CD	3:C:4:PHE:HA	2.23	0.67
1:A:73:LYS:O	1:A:76:THR:HG23	1.94	0.67
2:B:9:ILE:CD1	3:C:5:PHE:HZ	2.07	0.66
2:B:31:ARG:HD3	3:C:4:PHE:HA	1.77	0.66
1:A:26:GLY:HA2	1:A:38:ILE:HD13	1.77	0.65
2:B:2:ARG:HH12	2:B:64:ARG:HG3	1.60	0.64
2:B:2:ARG:CZ	2:B:2:ARG:CA	2.76	0.63
1:A:38:ILE:HD12	1:A:110:ASN:CB	2.28	0.63
2:B:66:ARG:HH11	2:B:66:ARG:HG2	1.63	0.62
2:B:2:ARG:HB3	3:C:2:LEU:HD13	1.81	0.62
1:A:69:SER:N	1:A:83:ASN:OD1	2.30	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:26:GLY:HA2	1:A:38:ILE:CD1	2.29	0.62
1:A:112:ASN:OD1	2:B:112:ASN:ND2	2.32	0.61
2:B:9:ILE:CD1	3:C:5:PHE:CZ	2.84	0.60
2:B:2:ARG:HD3	2:B:6:LYS:HE3	1.83	0.59
2:B:66:ARG:NH1	2:B:66:ARG:HG2	2.17	0.58
2:B:23:ALA:O	2:B:30:GLY:HA3	2.03	0.58
1:A:33:GLY:CA	2:B:31:ARG:HH12	1.98	0.58
2:B:118:ARG:HG2	2:B:118:ARG:NH1	2.12	0.58
2:B:13:VAL:HG12	2:B:17:ARG:HB2	1.86	0.58
2:B:6:LYS:NZ	3:C:4:PHE:HZ	2.02	0.57
3:C:4:PHE:HD1	3:C:5:PHE:CB	2.20	0.55
3:C:4:PHE:CD1	3:C:5:PHE:CB	2.89	0.55
1:A:66:LYS:HE3	1:A:87:ALA:HB2	1.90	0.54
1:A:118:ARG:HD2	5:A:417:HOH:O	2.08	0.53
2:B:49:ASP:OD1	3:C:6:ALA:HB3	2.10	0.51
2:B:69:SER:N	2:B:83:ASN:OD1	2.43	0.51
2:B:25:TYR:CZ	2:B:41:LEU:HD23	2.45	0.51
1:A:25:TYR:CZ	1:A:41:LEU:HD23	2.46	0.50
2:B:2:ARG:CA	2:B:2:ARG:NH1	2.74	0.50
1:A:31:ARG:NH1	2:B:24:ASP:OD1	2.45	0.50
1:A:43:ARG:O	1:A:47:VAL:HG23	2.11	0.49
2:B:2:ARG:HD3	3:C:4:PHE:CZ	2.47	0.49
1:A:73:LYS:O	1:A:74:ALA:HB3	2.13	0.49
2:B:2:ARG:HA	2:B:2:ARG:NH2	2.26	0.49
3:C:3:VAL:O	3:C:4:PHE:CB	2.59	0.49
1:A:74:ALA:HB3	1:A:76:THR:HG23	1.95	0.48
1:A:17:ARG:NH1	1:A:17:ARG:HG2	2.28	0.47
2:B:61:CYS:O	2:B:63:PRO:HD3	2.13	0.47
2:B:83:ASN:HB3	2:B:87:ALA:HB3	1.95	0.47
1:A:13:VAL:HG12	1:A:13:VAL:O	2.16	0.46
2:B:6:LYS:HG2	3:C:5:PHE:HE1	1.81	0.46
2:B:71:GLU:OE1	2:B:78:THR:HB	2.15	0.46
1:A:33:GLY:CA	2:B:31:ARG:NH1	2.67	0.46
1:A:83:ASN:HB3	1:A:87:ALA:CB	2.47	0.45
2:B:2:ARG:CB	2:B:2:ARG:HH11	1.97	0.45
2:B:31:ARG:HD2	3:C:4:PHE:N	2.31	0.45
2:B:31:ARG:HD3	3:C:4:PHE:CB	2.46	0.45
1:A:22:PHE:O	1:A:29:CYS:HB3	2.17	0.45
1:A:6:LYS:HE3	5:A:412:HOH:O	2.16	0.45
2:B:48:HIS:CD2	2:B:97:ALA:HB2	2.53	0.44
2:B:76:THR:HG23	5:B:418:HOH:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:31:ARG:HD2	3:C:4:PHE:HA	1.99	0.43
1:A:38:ILE:HD11	1:A:110:ASN:HB3	1.99	0.43
2:B:73:THR:O	2:B:74:ALA:C	2.57	0.42
1:A:30:GLY:O	2:B:31:ARG:NE	2.52	0.42
2:B:24:ASP:OD2	2:B:112:ASN:OD1	2.38	0.42
2:B:2:ARG:NH1	2:B:64:ARG:CG	2.74	0.42
1:A:36:THR:HA	1:A:37:PRO:HD3	1.95	0.42
1:A:115:LEU:HG	5:A:405:HOH:O	2.19	0.41
2:B:66:ARG:HH11	2:B:66:ARG:CG	2.32	0.41
1:A:25:TYR:O	1:A:29:CYS:HB2	2.20	0.41
2:B:2:ARG:HH12	2:B:64:ARG:CG	2.31	0.41
2:B:30:GLY:O	2:B:31:ARG:O	2.39	0.40
1:A:17:ARG:HH11	1:A:17:ARG:HG2	1.85	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	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	117/119 (98%)	106 (91%)	10 (8%)	1 (1%)	17	53
2	B	117/119 (98%)	104 (89%)	12 (10%)	1 (1%)	17	53
3	C	3/5 (60%)	0	0	3 (100%)	0	0
All	All	237/243 (98%)	210 (89%)	22 (9%)	5 (2%)	7	30

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	31	ARG
3	C	3	VAL
3	C	4	PHE
3	C	5	PHE

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Mol	Chain	Res	Type
1	A	30	GLY

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	99/99 (100%)	91 (92%)	8 (8%)	11	37
2	B	99/99 (100%)	90 (91%)	9 (9%)	9	32
3	C	4/4 (100%)	2 (50%)	2 (50%)	0	0
All	All	202/202 (100%)	183 (91%)	19 (9%)	8	31

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	LYS
1	A	19	TRP
1	A	40	ASP
1	A	58	GLN
1	A	72	CYS
1	A	73	LYS
1	A	76	THR
1	A	108	ASP
2	B	2	ARG
2	B	31	ARG
2	B	53	ASN
2	B	64	ARG
2	B	66	ARG
2	B	71	GLU
2	B	84	ASN
2	B	89	SER
2	B	108	ASP
3	C	4	PHE
3	C	5	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	10	GLN
1	A	58	GLN
1	A	65	GLN
2	B	53	ASN
2	B	112	ASN
2	B	116	GLN
2	B	120	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 2 ligands modelled in this entry, 2 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

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.