

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

#### Oct 18, 2023 – 05:06 AM EDT

PDB ID : 1SKG

Title : Structure-based rational drug design: Crystal structure of the complex formed

between Phospholipase A2 and a pentapeptide Val-Ala-Phe-Arg-Ser

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Deposited on : 2004-03-04

Resolution : 1.21 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 (i)) were used in the production of this report:

MolProbity : 4.02b-467

Mogul : 1.8.5 (274361), CSD as541be (2020)

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)

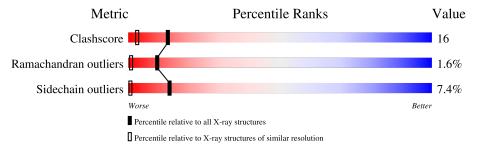
Validation Pipeline (wwPDB-VP) : 2.36

# 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 1.21 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	$(\# \mathrm{Entries})$	$(\#  ext{Entries},  ext{ resolution range}( ext{Å}))$
Clashscore	141614	1294 (1.24-1.20)
Ramachandran outliers	138981	1251 (1.24-1.20)
Sidechain outliers	138945	1250 (1.24-1.20)

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%

Note EDS was not executed.

Mol	Chain	Length	Q	uality of chain		
1	A	121	80	0%	17%	
2	В	5	40%	40%	20%	_



# 2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 1258 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.

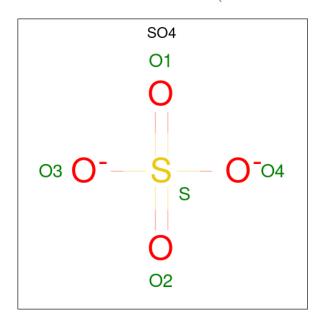
• Molecule 1 is a protein called Phospholipase A2.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	A	121	Total 944	C 591	N 159	O 178	S 16	0	0	0

• Molecule 2 is a protein called VAFRS.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	В	5	Total 40	C 26	N 8	O 6	0	0	0

• Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0

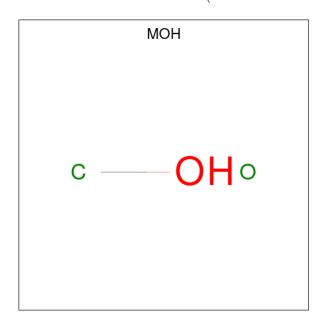
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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0

 $\bullet$  Molecule 4 is METHANOL (three-letter code: MOH) (formula:  $\mathrm{CH_{4}O}).$ 



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 2 1 1	0	0
4	A	1	Total C O 2 1 1	0	0

• Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	248	Total O 248 248	0	0
5	В	2	Total O 2 2	0	0

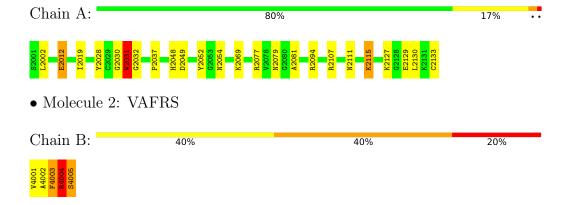


# 3 Residue-property plots (i)

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. 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





# 4 Data and refinement statistics (i)

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

Property	Value	Source	
Space group	P 43	Depositor	
Cell constants	52.35Å 52.35Å 47.83Å	Depositor	
a, b, c, $\alpha$ , $\beta$ , $\gamma$	90.00° 90.00° 90.00°	Depositor	
Resolution (Å)	20.00 - 1.21	Depositor	
% Data completeness	97.0 (20.00-1.21)	Depositor	
(in resolution range)	31.0 (20.00 1.21)		
$R_{merge}$	(Not available)	Depositor	
$R_{sym}$	0.08	Depositor	
Refinement program	SHELXL-97	Depositor	
$R, R_{free}$	0.141 , 0.183	Depositor	
Estimated twinning fraction	No twinning to report.	Xtriage	
Total number of atoms	1258	wwPDB-VP	
Average B, all atoms (Å <sup>2</sup> )	22.0	wwPDB-VP	



# 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: SO4, MOH

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	Bo	nd lengths	Bond angles		
		RMSZ	# Z  > 5	RMSZ	# Z  > 5	
1	A	0.77	1/964 (0.1%)	1.26	7/1292 (0.5%)	
2	В	0.74	0/40	1.68	1/52 (1.9%)	
All	All	0.77	1/1004 (0.1%)	1.28	8/1344 (0.6%)	

#### All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	$\mathbf{Z}$	Observed(A)	$\operatorname{Ideal}( ext{\AA})$
1	A	2133	CYS	C-OXT	8.70	1.39	1.23

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	$\mathbf{Z}$	$Observed(^o)$	$\mathbf{Ideal}(^o)$
1	A	2094	ARG	NE-CZ-NH1	-8.44	116.08	120.30
1	A	2052	TYR	CB-CG-CD2	-8.00	116.20	121.00
1	A	2012	GLU	CG-CD-OE2	7.99	134.28	118.30
1	A	2012	GLU	CG-CD-OE1	-6.17	105.97	118.30
1	A	2054	ASN	CB-CG-ND2	5.99	131.07	116.70
1	A	2052	TYR	CB-CG-CD1	5.80	124.48	121.00
1	A	2054	ASN	CA-CB-CG	5.39	125.25	113.40
2	В	4004	ARG	C-N-CA	5.21	134.74	121.70

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	asvmmetric	unit.	whereas S	Svmm-	Clashes	lists s	vmmetr	v-related	clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	944	0	904	26	0
2	В	40	0	40	25	0
3	A	20	0	0	0	0
4	A	4	0	0	0	0
5	A	248	0	0	3	0
5	В	2	0	0	1	0
All	All	1258	0	944	32	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 16.

All (32) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic	Clash
710011-1	1100111-2	$\operatorname{distance} \left( \operatorname{\AA} \right)$	overlap (Å)
1:A:2019:ILE:HD11	2:B:4001:VAL:HA	1.46	0.97
2:B:4004:ARG:HD2	2:B:4004:ARG:H	1.30	0.96
1:A:2031:TRP:HA	2:B:4004:ARG:HA	1.62	0.81
1:A:2002:LEU:HD11	2:B:4002:ALA:HB3	1.67	0.75
2:B:4004:ARG:HD3	5:B:9220:HOH:O	1.85	0.75
1:A:2032:GLY:O	2:B:4005:SER:HA	1.89	0.72
1:A:2002:LEU:HD21	2:B:4002:ALA:N	2.06	0.71
1:A:2049:ASP:OD1	2:B:4004:ARG:HG3	1.93	0.68
1:A:2002:LEU:HD21	2:B:4002:ALA:H	1.59	0.66
1:A:2002:LEU:HG	2:B:4001:VAL:HB	1.77	0.66
1:A:2032:GLY:H	2:B:4004:ARG:HB2	1.64	0.62
5:A:9214:HOH:O	2:B:4002:ALA:HB2	2.04	0.58
1:A:2002:LEU:CD2	2:B:4001:VAL:HB	2.34	0.58
1:A:2002:LEU:HD11	2:B:4002:ALA:CB	2.34	0.56
1:A:2002:LEU:CG	2:B:4001:VAL:HB	2.37	0.55
1:A:2028:TYR:HA	1:A:2032:GLY:HA2	1.87	0.55
1:A:2028:TYR:HA	1:A:2032:GLY:CA	2.37	0.54
1:A:2019:ILE:CD1	2:B:4001:VAL:HA	2.30	0.54
1:A:2002:LEU:CD1	2:B:4002:ALA:HB3	2.38	0.53
1:A:2069:LYS:HE2	2:B:4003:PHE:CD2	2.43	0.53
2:B:4002:ALA:O	2:B:4003:PHE:HB3	2.08	0.53
1:A:2069:LYS:HE2	2:B:4003:PHE:HD2	1.73	0.52
1:A:2030:GLY:HA2	2:B:4002:ALA:HB1	1.96	0.48
1:A:2130:LEU:HD23	5:A:9072:HOH:O	2.17	0.45
1:A:2012:GLU:HG3	1:A:2107:ARG:HB2	2.00	0.43
1:A:2048:HIS:HE1	2:B:4004:ARG:NH2	2.17	0.43

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Atom-1	Atom-2	Interatomic	Clash	
Atom-1	Atom-2	${ m distance} ({ m \AA})$	overlap (Å)	
2:B:4004:ARG:HD2	2:B:4004:ARG:N	2.10	0.43	
1:A:2077:ARG:HA	1:A:2081:ALA:O	2.20	0.42	
1:A:2037:PRO:HD3	1:A:2130:LEU:HD13	2.02	0.42	
1:A:2115:LYS:HB3	1:A:2115:LYS:HE2	1.71	0.42	
1:A:2032:GLY:N	2:B:4004:ARG:HB2	2.31	0.41	
5:A:9077:HOH:O	2:B:4001:VAL:HG13	2.20	0.41	

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	119/121 (98%)	115 (97%)	2 (2%)	2 (2%)	9 1
2	В	3/5~(60%)	0	3 (100%)	0	100 100
All	All	122/126~(97%)	115 (94%)	5 (4%)	2 (2%)	9 1

#### All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	2079	ASN
1	A	2031	TRP

#### 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	104/104 (100%)	99 (95%)	5 (5%)	25 2		
2	В	4/4 (100%)	1 (25%)	3 (75%)	0 0		
All	All	108/108 (100%)	100 (93%)	8 (7%)	13 1		

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2031	TRP
1	A	2111	ASN
1	A	2115	LYS
1	A	2127	LYS
1	A	2129	GLU
2	В	4003	PHE
2	В	4004	ARG
2	В	4005	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	2111	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)

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

#### 5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry (i)

6 ligands 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	Tuno	Chain	Chain Res		Bond lengths			Bond angles		
MIOI	Type	Chain	nes	Link	Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	SO4	A	3001	-	4,4,4	0.32	0	6,6,6	0.59	0
3	SO4	A	3004	-	4,4,4	0.23	0	6,6,6	0.35	0
3	SO4	A	3003	-	4,4,4	0.23	0	6,6,6	0.22	0
4	MOH	A	5679	-	1,1,1	0.30	0	-		
4	MOH	A	5678	-	1,1,1	0.28	0	-		
3	SO4	A	3002	-	4,4,4	0.30	0	6,6,6	0.31	0

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 (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)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

## 6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

#### 6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

#### 6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.

