



# Full wwPDB X-ray Structure Validation Report

Oct 16, 2021 – 11:24 PM EDT

PDB ID : 1SQM  
Title : STRUCTURE OF [R563A] LEUKOTRIENE A4 HYDROLASE  
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Deposited on : 2004-03-19  
Resolution : 2.30 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](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  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtrriage (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) : 2.23.2

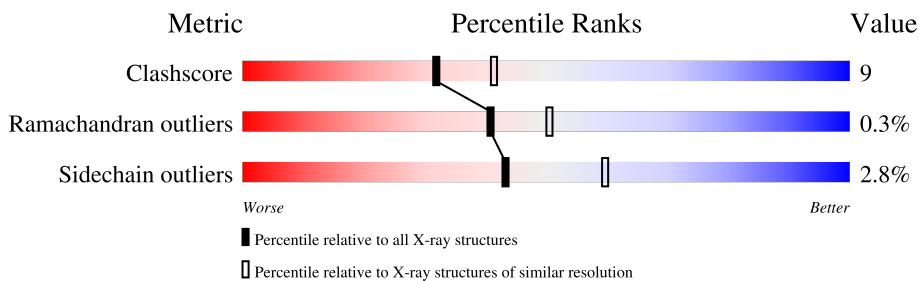
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.30 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	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)

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  $\geq 3$ , 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  $\leq 5\%$ .

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	610	

## 2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 5480 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 LEUKOTRIENE A-4 HYDROLASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	610	4871	3127	808	915	21	0	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	563	ALA	ARG	engineered mutation	UNP P09960

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		

- Molecule 3 is YTTERBIUM (III) ION (three-letter code: YB) (formula: Yb).

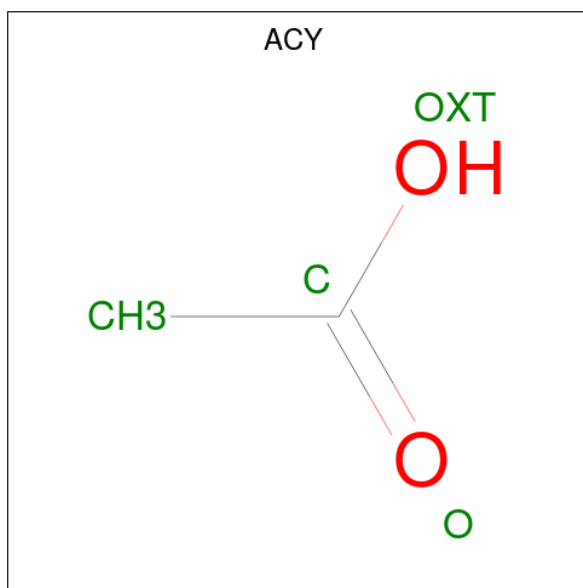
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Yb	0	0
			1	1		

- Molecule 4 is IMIDAZOLE (three-letter code: IMD) (formula: C<sub>3</sub>H<sub>5</sub>N<sub>2</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	N	0	0
			5	3	2		

- Molecule 5 is ACETIC ACID (three-letter code: ACY) (formula: C<sub>2</sub>H<sub>4</sub>O<sub>2</sub>).



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

- Molecule 6 is water.

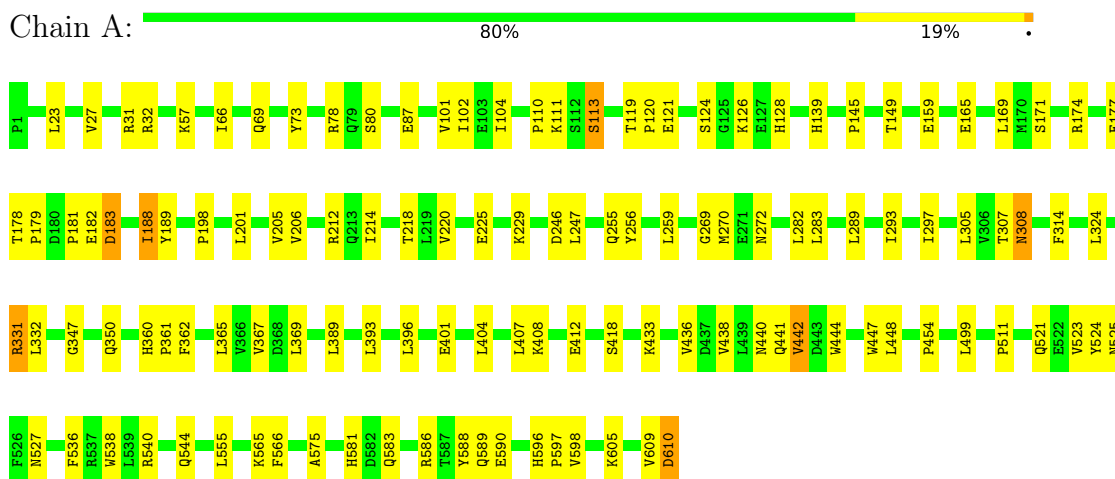
<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>		<b>ZeroOcc</b>	<b>AltConf</b>
6	A	598	Total 598	O 598	0	0

### 3 Residue-property plots

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: LEUKOTRIENE A-4 HYDROLASE



## 4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	78.00Å 86.85Å 98.95Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	16.48 – 2.30	Depositor
% Data completeness (in resolution range)	(Not available) (16.48-2.30)	Depositor
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	CNS	Depositor
R, $R_{free}$	0.190 , 0.230	Depositor
Estimated twinning fraction	No twinning to report.	Xtrriage
Total number of atoms	5480	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	26.0	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: ZN, YB, IMD, ACY

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	0.56	0/4996	0.74	0/6791

There are no bond length outliers.

There are no bond angle outliers.

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	4871	0	4829	86	0
2	A	1	0	0	0	0
3	A	1	0	0	0	0
4	A	5	0	5	0	0
5	A	4	0	3	1	0
6	A	598	0	0	10	0
All	All	5480	0	4837	86	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.

All (86) 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:360:HIS:HD2	1:A:362:PHE:H	1.08	0.92
1:A:527:ASN:ND2	1:A:538:TRP:HE1	1.71	0.87
1:A:360:HIS:CD2	1:A:362:PHE:H	1.97	0.80
1:A:527:ASN:HD22	1:A:538:TRP:HE1	1.28	0.76
1:A:159:GLU:HB3	1:A:188:ILE:HD11	1.68	0.74
1:A:521:GLN:HE22	1:A:527:ASN:H	1.37	0.72
1:A:174:ARG:HD2	6:A:3051:HOH:O	1.90	0.71
1:A:324:LEU:HG	1:A:407:LEU:HD21	1.74	0.68
1:A:111:LYS:HG2	1:A:111:LYS:O	1.97	0.64
1:A:360:HIS:HD2	1:A:362:PHE:N	1.90	0.64
1:A:347:GLY:HA2	1:A:350:GLN:HE21	1.65	0.62
1:A:396:LEU:HG	1:A:438:VAL:CG1	2.31	0.60
1:A:269:GLY:O	5:A:3001:ACY:H1	2.01	0.60
1:A:78:ARG:NH2	6:A:3210:HOH:O	2.36	0.59
1:A:270:MET:HA	1:A:270:MET:HE2	1.85	0.58
1:A:293:ILE:HG22	1:A:297:ILE:HD12	1.87	0.56
1:A:588:TYR:CE2	1:A:609:VAL:HG21	2.40	0.56
1:A:555:LEU:HD11	1:A:583:GLN:HG2	1.88	0.55
1:A:178:THR:O	1:A:188:ILE:HG22	2.07	0.55
1:A:182:GLU:O	1:A:183:ASP:HB2	2.07	0.54
1:A:401:GLU:HG3	6:A:3204:HOH:O	2.07	0.54
1:A:27:VAL:HG21	1:A:206:VAL:HG21	1.90	0.53
1:A:282:LEU:HA	1:A:289:LEU:HD12	1.91	0.52
1:A:586:ARG:O	1:A:590:GLU:HG3	2.10	0.52
1:A:32:ARG:HB2	1:A:110:PRO:HB3	1.90	0.52
1:A:31:ARG:HG2	1:A:31:ARG:HH21	1.75	0.51
1:A:565:LYS:HE3	1:A:566:PHE:CZ	2.46	0.51
1:A:433:LYS:O	1:A:436:VAL:HG13	2.11	0.51
1:A:605:LYS:NZ	1:A:605:LYS:HB3	2.26	0.51
1:A:159:GLU:CB	1:A:188:ILE:HD11	2.40	0.51
1:A:225:GLU:H	1:A:225:GLU:CD	2.15	0.51
1:A:229:LYS:HG3	6:A:3087:HOH:O	2.11	0.50
1:A:596:HIS:HE1	6:A:3002:HOH:O	1.94	0.50
1:A:179:PRO:O	1:A:181:PRO:HD3	2.11	0.50
1:A:57:LYS:NZ	1:A:78:ARG:HH22	2.10	0.50
1:A:120:PRO:O	1:A:126:LYS:HA	2.12	0.50
1:A:523:VAL:HG23	1:A:524:TYR:CD1	2.46	0.49
1:A:307:THR:O	1:A:418:SER:HA	2.12	0.49
1:A:104:ILE:HG21	1:A:145:PRO:HG2	1.95	0.49
1:A:293:ILE:HG22	1:A:297:ILE:CD1	2.42	0.49
1:A:609:VAL:O	1:A:610:ASP:HB2	2.13	0.49
1:A:111:LYS:O	1:A:111:LYS:CG	2.62	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:169:LEU:HB2	1:A:205:VAL:CG1	2.44	0.47
1:A:229:LYS:HE2	1:A:283:LEU:HB3	1.96	0.47
1:A:586:ARG:NH1	1:A:590:GLU:OE2	2.47	0.47
1:A:246:ASP:OD2	1:A:331:ARG:NH2	2.47	0.47
1:A:367:VAL:HG23	1:A:369:LEU:HG	1.97	0.46
1:A:128:HIS:CE1	1:A:165:GLU:HG3	2.50	0.46
1:A:442:VAL:HG22	1:A:444:TRP:CE2	2.50	0.46
1:A:447:TRP:CZ3	1:A:454:PRO:HB3	2.51	0.46
1:A:499:LEU:HD22	6:A:3500:HOH:O	2.16	0.46
1:A:171:SER:HB3	1:A:272:ASN:HB3	1.96	0.45
1:A:139:HIS:HA	6:A:3477:HOH:O	2.16	0.45
1:A:218:THR:HG23	1:A:256:TYR:O	2.17	0.45
1:A:57:LYS:HE3	1:A:113:SER:OG	2.16	0.45
1:A:174:ARG:CD	6:A:3051:HOH:O	2.58	0.45
1:A:596:HIS:CD2	1:A:598:VAL:H	2.35	0.45
1:A:360:HIS:CD2	1:A:361:PRO:HD2	2.52	0.45
1:A:124:SER:HB3	1:A:225:GLU:CG	2.47	0.44
1:A:305:LEU:HD23	1:A:305:LEU:HA	1.87	0.44
1:A:586:ARG:NH1	6:A:3556:HOH:O	2.50	0.44
1:A:255:GLN:HG3	1:A:256:TYR:N	2.32	0.44
1:A:119:THR:OG1	1:A:121:GLU:HG2	2.18	0.44
1:A:214:ILE:HD13	1:A:220:VAL:CG2	2.48	0.44
1:A:408:LYS:O	1:A:412:GLU:HG3	2.18	0.43
1:A:121:GLU:H	1:A:121:GLU:CD	2.21	0.43
1:A:511:PRO:HB3	1:A:544:GLN:O	2.18	0.43
1:A:360:HIS:HA	1:A:361:PRO:HD3	1.88	0.43
1:A:389:LEU:O	1:A:393:LEU:HG	2.19	0.43
1:A:596:HIS:CG	1:A:597:PRO:HD2	2.53	0.43
1:A:447:TRP:CH2	1:A:454:PRO:HB3	2.54	0.43
1:A:589:GLN:OE1	1:A:589:GLN:HA	2.18	0.42
1:A:365:LEU:HD23	1:A:448:LEU:HD23	2.02	0.42
1:A:121:GLU:HA	1:A:126:LYS:HD2	2.01	0.42
1:A:308:ASN:OD1	1:A:314:PHE:HA	2.18	0.42
1:A:536:PHE:CZ	1:A:540:ARG:HD2	2.55	0.42
1:A:188:ILE:HD12	1:A:189:TYR:N	2.35	0.41
1:A:544:GLN:NE2	6:A:3136:HOH:O	2.52	0.41
1:A:73:TYR:HA	1:A:87:GLU:O	2.20	0.41
1:A:177:GLU:O	1:A:178:THR:CG2	2.69	0.41
1:A:198:PRO:HD2	1:A:201:LEU:HD12	2.03	0.41
1:A:404:LEU:HD23	1:A:404:LEU:HA	1.87	0.40
1:A:575:ALA:O	1:A:581:HIS:HD2	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:440:ASN:HD22	1:A:440:ASN:HA	1.73	0.40
1:A:66:ILE:CD1	1:A:102:ILE:HG12	2.52	0.40
1:A:205:VAL:HG11	1:A:259: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	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	608/610 (100%)	584 (96%)	22 (4%)	2 (0%)	41 50

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	80	SER
1	A	183	ASP

### 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	541/541 (100%)	526 (97%)	15 (3%)	43 60

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	23	LEU
1	A	69	GLN
1	A	101	VAL
1	A	113	SER
1	A	149	THR
1	A	188	ILE
1	A	212	ARG
1	A	247	LEU
1	A	308	ASN
1	A	331	ARG
1	A	332	LEU
1	A	441	GLN
1	A	442	VAL
1	A	525	ASN
1	A	610	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	69	GLN
1	A	213	GLN
1	A	226	GLN
1	A	341	ASN
1	A	350	GLN
1	A	360	HIS
1	A	440	ASN
1	A	466	ASN
1	A	521	GLN
1	A	525	ASN
1	A	527	ASN
1	A	544	GLN
1	A	581	HIS
1	A	596	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](#)

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](#)

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 for Mogul analysis.

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	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
5	ACY	A	3001	2	1,3,3	2.57	1 (100%)	0,3,3	-	-
4	IMD	A	2002	-	3,5,5	0.60	0	4,5,5	0.49	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
4	IMD	A	2002	-	-	-	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	3001	ACY	CH3-C	2.57	1.52	1.48

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	3001	ACY	1	0

## 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.