



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 03:13 AM UTC

PDB ID : 3MLM / pdb_00003mlm
Title : Crystal structure of Bn IV in complex with myristic acid: A Lys49 myotoxic phospholipase A2 from Bothrops neuwiedi venom
Authors : Delatorre, P.; Rocha, B.A.M.; Cavada, B.S.; Toyama, M.H.; Toyama, D.; Gadelha, C.A.A.
Deposited on : 2010-04-17
Resolution : 2.21 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 ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

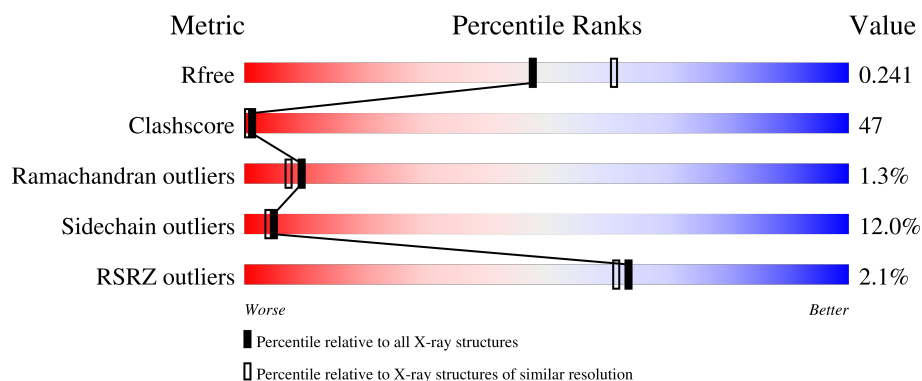
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.21 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 (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	7682 (2.24-2.20)
Clashscore	190562	8402 (2.24-2.20)
Ramachandran outliers	187476	8303 (2.24-2.20)
Sidechain outliers	187428	8304 (2.24-2.20)
RSRZ outliers	180081	7683 (2.24-2.20)

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 $\leq 5\%$. 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	121	
1	B	121	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	SO4	A	122	-	-	X	-
2	SO4	B	122	-	-	X	-
2	SO4	B	123	-	-	X	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 2034 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.

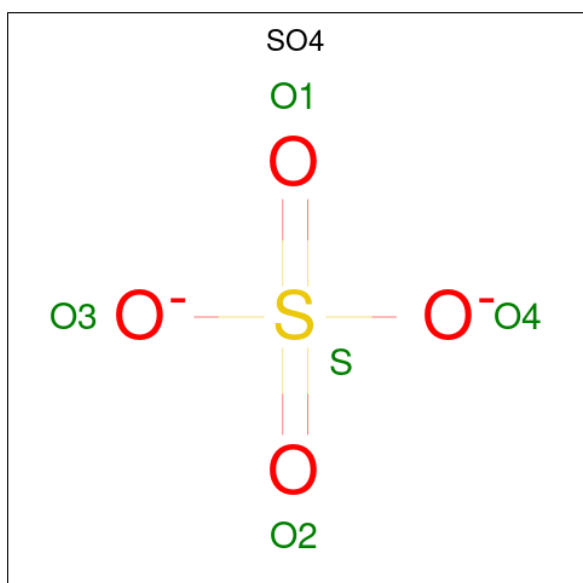
- Molecule 1 is a protein called BN-IV Lys-49 Phospholipase A2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	121	Total	C	N	O	S	0	0	0
			948	594	166	173	15			
1	B	121	Total	C	N	O	S	0	0	0
			948	594	166	173	15			

There are 6 discrepancies between the modelled and reference sequences:

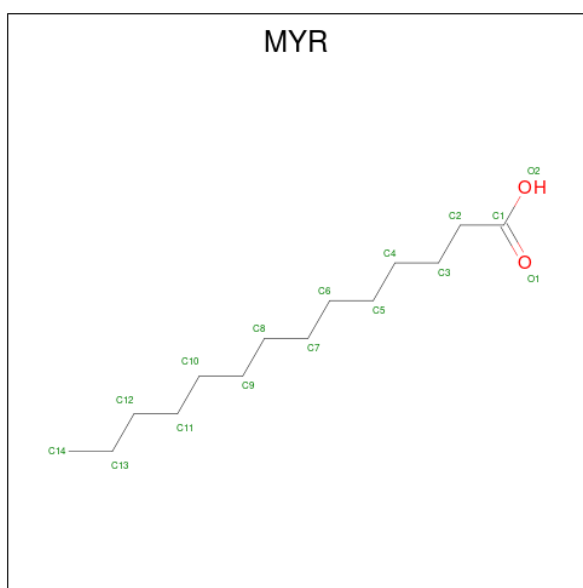
Chain	Residue	Modelled	Actual	Comment	Reference
A	2	LEU	-	variant	UNP Q9IAT9
A	35	GLY	GLN	variant	UNP Q9IAT9
A	54	ILE	LEU	variant	UNP Q9IAT9
B	2	LEU	-	variant	UNP Q9IAT9
B	35	GLY	GLN	variant	UNP Q9IAT9
B	54	ILE	LEU	variant	UNP Q9IAT9

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

- Molecule 3 is MYRISTIC ACID (CCD ID: MYR) (formula: $C_{14}H_{28}O_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 16 14 2	0	0
3	B	1	Total C O 16 14 2	0	0

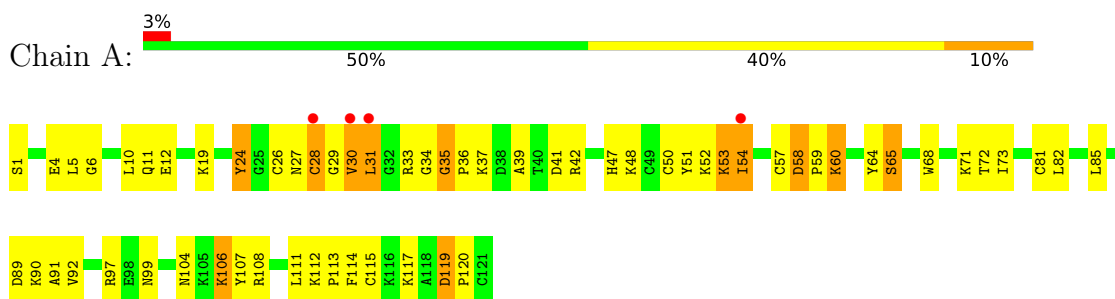
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	44	Total O 44 44	0	0
4	B	42	Total O 42 42	0	0

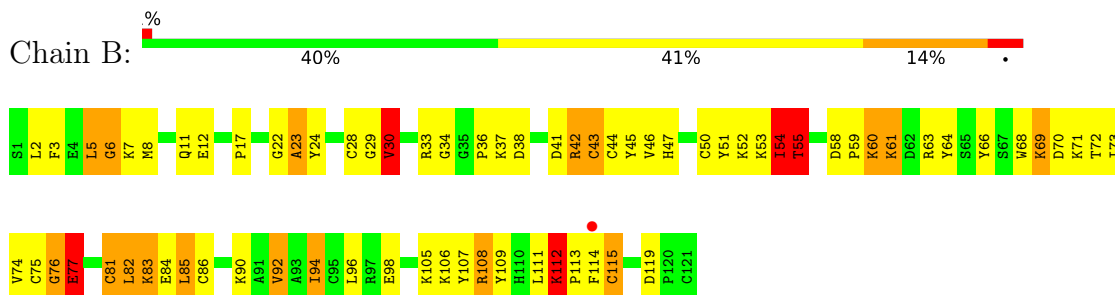
3 Residue-property plots

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 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: BN-IV Lys-49 Phospholipase A2



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4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	38.80Å 70.47Å 44.01Å 90.00° 109.32° 90.00°	Depositor
Resolution (Å)	32.49 – 2.21 32.49 – 2.21	Depositor EDS
% Data completeness (in resolution range)	96.4 (32.49-2.21) 96.8 (32.49-2.21)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.97 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.5.0066	Depositor
R, R_{free}	0.205 , 0.253 0.201 , 0.241	Depositor DCC
R_{free} test set	522 reflections (4.61%)	wwPDB-VP
Wilson B-factor (Å ²)	26.7	Xtriage
Anisotropy	0.215	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 45.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	2034	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.22% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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: MYR, SO4

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	1.49	2/968 (0.2%)	1.49	10/1295 (0.8%)
1	B	1.65	11/968 (1.1%)	1.58	15/1295 (1.2%)
All	All	1.57	13/1936 (0.7%)	1.54	25/2590 (1.0%)

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
1	B	0	2

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	77	GLU	CA-C	-8.02	1.42	1.52
1	B	41	ASP	C-O	-7.80	1.14	1.24
1	B	85	LEU	C-O	-6.56	1.16	1.24
1	B	94	ILE	CA-CB	-6.31	1.47	1.54
1	A	65	SER	CA-C	-6.21	1.45	1.52
1	B	53	LYS	C-O	-6.01	1.16	1.24
1	B	112	LYS	CA-C	5.41	1.59	1.52
1	B	54	ILE	C-O	-5.38	1.16	1.23
1	B	81	CYS	C-O	-5.20	1.18	1.24
1	B	23	ALA	CA-CB	-5.17	1.47	1.53
1	B	98	GLU	C-O	-5.14	1.17	1.24
1	B	22	GLY	C-O	-5.09	1.17	1.24
1	A	28	CYS	CB-SG	-5.03	1.64	1.81

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	28	CYS	N-CA-C	-11.35	90.01	108.52
1	B	30	VAL	N-CA-C	8.29	120.08	112.29
1	B	106	LYS	N-CA-C	-8.04	102.74	112.54
1	A	24	TYR	N-CA-C	7.57	120.83	109.25
1	A	119	ASP	CA-C-N	-7.53	111.86	119.92
1	A	119	ASP	C-N-CA	-7.53	111.86	119.92
1	B	22	GLY	N-CA-C	-6.53	106.26	114.16
1	A	64	TYR	CA-C-N	-6.52	112.58	122.81
1	A	64	TYR	C-N-CA	-6.52	112.58	122.81
1	B	50	CYS	N-CA-C	-6.19	104.54	111.28
1	A	10	LEU	N-CA-C	-6.05	104.60	111.14
1	B	3	PHE	N-CA-C	-5.70	105.15	111.36
1	B	5	LEU	N-CA-C	-5.49	104.98	110.97
1	B	55	THR	CB-CA-C	-5.48	103.02	111.74
1	B	76	GLY	CA-C-N	-5.48	111.07	121.54
1	B	76	GLY	C-N-CA	-5.48	111.07	121.54
1	B	43	CYS	N-CA-C	-5.30	105.50	111.28
1	B	112	LYS	CA-C-O	5.24	123.15	118.33
1	B	115	CYS	CA-CB-SG	-5.23	102.37	114.40
1	B	54	ILE	CA-C-N	5.12	131.19	123.04
1	B	54	ILE	C-N-CA	5.12	131.19	123.04
1	A	106	LYS	CA-C-N	-5.04	114.59	122.65
1	A	106	LYS	C-N-CA	-5.04	114.59	122.65
1	B	54	ILE	CB-CA-C	5.01	118.39	111.38
1	A	91	ALA	N-CA-C	-5.00	105.92	112.23

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	108	ARG	Sidechain
1	B	54	ILE	Peptide

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	948	0	936	79	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	948	0	938	105	0
2	A	10	0	0	4	0
2	B	10	0	0	7	0
3	A	16	0	27	2	0
3	B	16	0	27	8	0
4	A	44	0	0	2	0
4	B	42	0	0	8	0
All	All	2034	0	1928	182	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 47.

All (182) 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:29:GLY:O	1:A:30:VAL:CG2	1.70	1.39
1:A:31:LEU:HD22	1:A:31:LEU:O	1.35	1.24
1:A:31:LEU:HD21	1:A:112:LYS:HD3	1.30	1.12
1:B:105:LYS:HG2	1:B:108:ARG:NH1	1.61	1.12
1:B:55:THR:HA	4:B:158:HOH:O	1.53	1.08
1:B:105:LYS:HG2	1:B:108:ARG:HH11	1.16	1.08
1:B:69:LYS:CE	1:B:74:VAL:HG21	1.84	1.07
1:A:31:LEU:HD11	1:A:112:LYS:HE3	1.38	1.06
1:B:42:ARG:HG3	1:B:42:ARG:HH11	1.26	1.01
1:A:31:LEU:HD11	1:A:112:LYS:CE	1.90	1.01
1:A:31:LEU:HD11	1:A:112:LYS:CD	1.90	1.00
1:A:29:GLY:C	1:A:30:VAL:HG23	1.88	0.98
1:A:29:GLY:O	1:A:30:VAL:HG23	0.80	0.98
1:A:106:LYS:HE3	1:A:107:TYR:CE2	1.99	0.97
1:A:31:LEU:HD21	1:A:112:LYS:CD	1.97	0.95
1:A:31:LEU:O	1:A:31:LEU:CD2	2.15	0.94
1:B:69:LYS:HZ1	1:B:74:VAL:HG11	1.35	0.91
1:A:31:LEU:HD13	1:A:31:LEU:N	1.85	0.90
1:A:27:ASN:OD1	1:A:33:ARG:HB3	1.71	0.90
1:B:69:LYS:HE3	1:B:74:VAL:HG21	1.49	0.90
1:A:31:LEU:HD13	1:A:31:LEU:H	1.34	0.90
1:A:31:LEU:CD2	1:A:112:LYS:HD3	2.00	0.90
1:B:42:ARG:HH11	1:B:42:ARG:CG	1.87	0.88
1:B:6:GLY:HA3	3:B:125:MYR:H121	1.57	0.87
1:B:69:LYS:CD	1:B:74:VAL:HG21	2.07	0.84
1:B:42:ARG:HG3	1:B:42:ARG:NH1	1.87	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:42:ARG:HD2	4:B:137:HOH:O	1.78	0.82
1:B:69:LYS:NZ	1:B:74:VAL:HG11	1.94	0.82
1:B:61:LYS:N	1:B:61:LYS:HD2	1.96	0.80
1:A:31:LEU:N	1:A:31:LEU:CD1	2.44	0.80
1:B:69:LYS:HD2	1:B:74:VAL:HG21	1.64	0.79
1:B:42:ARG:CD	4:B:137:HOH:O	2.29	0.79
1:B:2:LEU:HD11	1:B:60:LYS:HD2	1.65	0.79
1:B:60:LYS:HB3	1:B:61:LYS:HD3	1.65	0.78
1:B:111:LEU:C	1:B:113:PRO:HD2	2.09	0.78
1:B:108:ARG:NH2	2:B:123:SO4:O1	2.19	0.74
1:B:60:LYS:C	1:B:61:LYS:HD2	2.13	0.73
1:A:30:VAL:O	1:A:30:VAL:HG12	1.88	0.72
1:A:59:PRO:HB3	1:A:85:LEU:CD1	2.21	0.71
1:B:105:LYS:HE2	1:B:108:ARG:HH12	1.55	0.71
1:A:51:TYR:O	1:A:54:ILE:HB	1.91	0.71
1:A:24:TYR:O	1:A:28:CYS:O	2.09	0.70
1:A:54:ILE:HG23	1:A:81:CYS:SG	2.31	0.70
1:B:61:LYS:N	1:B:61:LYS:CD	2.55	0.70
1:A:104:ASN:OD1	1:A:106:LYS:HB3	1.93	0.69
1:B:112:LYS:N	1:B:113:PRO:CD	2.54	0.69
1:A:31:LEU:CD1	1:A:112:LYS:HE3	2.19	0.69
1:A:33:ARG:NE	2:A:122:SO4:O1	2.21	0.69
1:B:60:LYS:CG	1:B:61:LYS:HZ2	2.06	0.69
1:A:59:PRO:HB3	1:A:85:LEU:HD13	1.75	0.68
1:B:54:ILE:HD11	1:B:84:GLU:HB3	1.76	0.68
1:A:112:LYS:O	1:A:115:CYS:HB2	1.96	0.66
1:B:6:GLY:CA	3:B:125:MYR:H121	2.24	0.66
1:A:31:LEU:CD1	1:A:112:LYS:CD	2.71	0.66
1:A:71:LYS:NZ	1:B:12:GLU:OE2	2.24	0.66
1:B:76:GLY:C	1:B:77:GLU:HG2	2.19	0.65
1:A:111:LEU:O	1:A:114:PHE:HB2	1.96	0.65
1:B:60:LYS:HB3	1:B:61:LYS:CD	2.27	0.65
1:B:54:ILE:CD1	1:B:84:GLU:HB3	2.28	0.64
1:B:69:LYS:HE3	1:B:74:VAL:CG2	2.26	0.63
1:A:27:ASN:OD1	1:A:33:ARG:CB	2.46	0.62
1:A:52:LYS:NZ	2:A:122:SO4:O4	2.19	0.62
1:A:24:TYR:O	1:A:28:CYS:HB2	2.00	0.61
1:B:33:ARG:NE	2:B:122:SO4:O3	2.25	0.61
1:A:6:GLY:HA3	3:A:124:MYR:H122	1.81	0.61
1:A:24:TYR:CD1	1:A:24:TYR:C	2.79	0.61
1:A:112:LYS:N	1:A:113:PRO:CD	2.64	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:5:LEU:HD11	1:B:92:VAL:HG11	1.83	0.60
1:B:64:TYR:CD1	1:B:86:CYS:HB2	2.36	0.60
1:A:36:PRO:CB	1:A:42:ARG:HG2	2.31	0.60
1:B:29:GLY:H	3:B:125:MYR:H22	1.67	0.59
1:B:77:GLU:OE2	1:B:82:LEU:HB3	2.01	0.59
1:A:30:VAL:O	1:A:30:VAL:CG1	2.51	0.58
1:B:111:LEU:C	1:B:113:PRO:CD	2.76	0.58
1:B:92:VAL:O	1:B:96:LEU:HG	2.03	0.58
1:B:64:TYR:CE1	1:B:86:CYS:HB2	2.39	0.58
1:B:68:TRP:CE2	1:B:71:LYS:HA	2.39	0.58
1:B:60:LYS:HG3	1:B:61:LYS:HZ2	1.68	0.57
1:B:60:LYS:CG	1:B:61:LYS:NZ	2.66	0.57
1:A:5:LEU:HD21	3:A:124:MYR:H21	1.86	0.57
1:A:57:CYS:N	1:A:81:CYS:SG	2.78	0.57
1:A:119:ASP:O	1:A:120:PRO:C	2.47	0.57
1:A:31:LEU:HD11	1:A:112:LYS:HD3	1.80	0.56
1:B:17:PRO:HB2	3:B:125:MYR:H92	1.87	0.56
1:B:47:HIS:CD2	1:B:92:VAL:HB	2.42	0.55
1:B:90:LYS:HE3	1:B:94:ILE:HD11	1.88	0.55
1:A:29:GLY:C	1:A:30:VAL:CG2	2.60	0.55
1:A:31:LEU:CD1	1:A:112:LYS:HD3	2.35	0.55
1:B:69:LYS:HD2	1:B:74:VAL:CG2	2.34	0.55
1:B:113:PRO:C	1:B:115:CYS:H	2.13	0.55
1:A:34:GLY:O	1:A:35:GLY:C	2.49	0.54
1:B:44:CYS:HB3	3:B:125:MYR:O2	2.06	0.54
1:B:105:LYS:CE	1:B:108:ARG:HH12	2.20	0.54
1:A:108:ARG:NH2	2:A:123:SO4:O4	2.30	0.54
1:B:105:LYS:HE3	2:B:123:SO4:O3	2.07	0.54
1:B:64:TYR:HB3	1:B:82:LEU:HG	1.90	0.54
1:A:106:LYS:HE3	1:A:107:TYR:CZ	2.43	0.54
1:A:112:LYS:N	1:A:113:PRO:HD3	2.23	0.53
1:A:24:TYR:C	1:A:28:CYS:O	2.51	0.53
1:B:24:TYR:HB3	1:B:28:CYS:HB2	1.90	0.53
1:A:114:PHE:HD1	4:A:139:HOH:O	1.90	0.53
1:B:52:LYS:NZ	2:B:122:SO4:O1	2.42	0.53
1:B:112:LYS:N	1:B:113:PRO:HD3	2.23	0.53
1:A:47:HIS:CD2	1:A:92:VAL:HB	2.44	0.52
1:A:47:HIS:HB2	1:A:92:VAL:CG2	2.39	0.52
1:B:24:TYR:OH	1:B:38:ASP:OD2	2.19	0.52
1:B:47:HIS:HB2	1:B:92:VAL:CG2	2.41	0.51
1:B:2:LEU:HD21	1:B:60:LYS:HD2	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:48:LYS:O	1:A:52:LYS:HG3	2.13	0.49
1:B:60:LYS:HB3	1:B:61:LYS:NZ	2.27	0.49
1:A:73:ILE:O	1:A:90:LYS:NZ	2.37	0.49
1:A:113:PRO:C	1:A:115:CYS:H	2.20	0.49
1:A:19:LYS:HZ2	1:A:108:ARG:NH1	2.11	0.49
1:B:24:TYR:CE1	1:B:107:TYR:HB2	2.48	0.49
1:A:31:LEU:HD11	1:A:112:LYS:HD2	1.85	0.48
1:A:59:PRO:HB3	1:A:85:LEU:HD11	1.95	0.48
1:B:29:GLY:H	3:B:125:MYR:C2	2.25	0.48
1:B:111:LEU:CB	1:B:113:PRO:HD2	2.44	0.48
1:A:36:PRO:HB3	1:A:42:ARG:HG2	1.95	0.48
1:A:52:LYS:NZ	2:A:122:SO4:S	2.83	0.48
1:B:23:ALA:C	4:B:147:HOH:O	2.56	0.48
1:B:105:LYS:CE	2:B:123:SO4:O3	2.61	0.48
1:B:54:ILE:HG23	1:B:81:CYS:SG	2.53	0.48
1:A:1:SER:OG	1:A:4:GLU:OE1	2.27	0.47
1:B:60:LYS:HB3	1:B:61:LYS:HZ3	1.80	0.47
1:A:11:GLN:HG2	4:B:164:HOH:O	2.15	0.47
1:A:58:ASP:OD2	1:A:60:LYS:N	2.45	0.47
1:A:31:LEU:CD2	1:A:112:LYS:CD	2.76	0.47
1:B:105:LYS:HE2	1:B:108:ARG:NH1	2.25	0.46
1:B:8:MET:HE2	1:B:73:ILE:HG21	1.97	0.46
1:B:42:ARG:O	1:B:43:CYS:C	2.52	0.46
1:B:60:LYS:HG2	1:B:61:LYS:NZ	2.30	0.46
1:B:52:LYS:NZ	2:B:122:SO4:S	2.86	0.46
1:B:2:LEU:CD1	1:B:60:LYS:HD2	2.42	0.45
1:B:7:LYS:O	1:B:11:GLN:HG3	2.16	0.45
1:B:42:ARG:NE	4:B:137:HOH:O	2.45	0.45
1:B:58:ASP:C	1:B:58:ASP:OD2	2.59	0.45
1:A:58:ASP:OD2	1:A:58:ASP:C	2.59	0.45
1:B:42:ARG:O	1:B:46:VAL:HG23	2.16	0.44
1:B:108:ARG:NH1	2:B:123:SO4:O1	2.50	0.44
1:A:68:TRP:CE2	1:A:71:LYS:HA	2.52	0.44
1:A:114:PHE:CD1	4:A:139:HOH:O	2.57	0.44
1:B:82:LEU:O	1:B:83:LYS:C	2.60	0.44
1:A:26:CYS:H	1:A:41:ASP:CG	2.25	0.44
1:B:69:LYS:O	1:B:72:THR:HB	2.17	0.44
1:B:83:LYS:HA	1:B:83:LYS:HD3	1.81	0.44
1:B:113:PRO:C	1:B:115:CYS:N	2.76	0.44
1:A:12:GLU:O	1:A:97:ARG:HG3	2.17	0.44
1:B:76:GLY:O	1:B:77:GLU:CB	2.59	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:69:LYS:NZ	1:B:74:VAL:HG21	2.31	0.44
1:B:51:TYR:O	1:B:54:ILE:HB	2.18	0.44
1:A:47:HIS:HB2	1:A:92:VAL:HG23	2.00	0.44
1:B:76:GLY:O	1:B:77:GLU:HB3	2.18	0.44
1:B:42:ARG:HB2	4:B:137:HOH:O	2.17	0.43
1:A:71:LYS:HD2	1:B:68:TRP:HH2	1.84	0.43
1:B:51:TYR:CE2	1:B:59:PRO:HB2	2.53	0.43
1:B:114:PHE:HB3	4:B:155:HOH:O	2.17	0.43
1:A:27:ASN:OD1	1:A:33:ARG:CA	2.67	0.43
1:A:58:ASP:HA	1:A:59:PRO:HD3	1.68	0.43
1:B:69:LYS:O	1:B:69:LYS:HG2	2.09	0.43
1:B:29:GLY:N	3:B:125:MYR:H22	2.33	0.43
1:B:34:GLY:O	1:B:36:PRO:HD3	2.18	0.43
1:B:108:ARG:O	1:B:109:TYR:HB2	2.18	0.43
1:B:74:VAL:C	1:B:75:CYS:O	2.60	0.42
1:B:112:LYS:O	1:B:115:CYS:HB3	2.18	0.42
1:B:47:HIS:HB2	1:B:92:VAL:HG23	2.01	0.42
1:B:45:TYR:OH	1:B:119:ASP:O	2.37	0.42
1:A:39:ALA:HB1	1:A:99:ASN:CG	2.45	0.42
1:A:36:PRO:CG	1:A:42:ARG:HG2	2.49	0.42
1:A:35:GLY:O	1:A:36:PRO:C	2.63	0.41
1:B:8:MET:HE2	1:B:73:ILE:CG2	2.50	0.41
1:B:75:CYS:HA	1:B:86:CYS:SG	2.60	0.41
1:B:59:PRO:HB3	1:B:85:LEU:HD13	2.02	0.41
1:B:30:VAL:HG22	1:B:30:VAL:O	2.21	0.41
1:B:105:LYS:CG	1:B:108:ARG:HH11	2.07	0.41
1:B:47:HIS:ND1	3:B:125:MYR:O1	2.48	0.41
1:A:50:CYS:O	1:A:53:LYS:HG2	2.21	0.41
1:A:89:ASP:O	1:A:92:VAL:HG12	2.20	0.41
1:A:19:LYS:NZ	1:A:108:ARG:NH1	2.69	0.41
1:B:54:ILE:HD13	1:B:84:GLU:HB3	2.03	0.41
1:A:5:LEU:HD11	1:A:92:VAL:HG11	2.02	0.40
1:B:11:GLN:HE22	1:B:66:TYR:HE2	1.69	0.40

There are no symmetry-related clashes.

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	119/121 (98%)	108 (91%)	9 (8%)	2 (2%)	7	5
1	B	119/121 (98%)	107 (90%)	11 (9%)	1 (1%)	16	15
All	All	238/242 (98%)	215 (90%)	20 (8%)	3 (1%)	9	7

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	30	VAL
1	A	35	GLY
1	B	6	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	104/104 (100%)	94 (90%)	10 (10%)	8	7
1	B	104/104 (100%)	89 (86%)	15 (14%)	3	2
All	All	208/208 (100%)	183 (88%)	25 (12%)	5	4

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	31	LEU
1	A	37	LYS
1	A	53	LYS

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Mol	Chain	Res	Type
1	A	54	ILE
1	A	58	ASP
1	A	60	LYS
1	A	65	SER
1	A	72	THR
1	A	82	LEU
1	A	117	LYS
1	B	30	VAL
1	B	37	LYS
1	B	42	ARG
1	B	54	ILE
1	B	55	THR
1	B	60	LYS
1	B	61	LYS
1	B	63	ARG
1	B	69	LYS
1	B	70	ASP
1	B	77	GLU
1	B	82	LEU
1	B	83	LYS
1	B	92	VAL
1	B	112	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	11	GLN
1	A	110	HIS
1	B	11	GLN

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 [i](#)

There are no oligosaccharides 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	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SO4	A	123	-	4,4,4	0.31	0	6,6,6	0.38	0
2	SO4	B	122	-	4,4,4	0.47	0	6,6,6	0.66	0
3	MYR	B	125	-	15,15,15	1.01	1 (6%)	15,15,15	1.57	2 (13%)
3	MYR	A	124	-	15,15,15	0.81	1 (6%)	15,15,15	1.01	0
2	SO4	B	123	-	4,4,4	0.25	0	6,6,6	0.10	0
2	SO4	A	122	-	4,4,4	0.51	0	6,6,6	0.92	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
3	MYR	A	124	-	-	10/13/13/13	-
3	MYR	B	125	-	-	10/13/13/13	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	124	MYR	O1-C1	2.43	1.30	1.22
3	B	125	MYR	C2-C1	2.09	1.55	1.50

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	125	MYR	O2-C1-O1	-4.67	111.33	123.33
3	B	125	MYR	O2-C1-C2	2.86	123.05	114.00

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	125	MYR	C1-C2-C3-C4
3	A	124	MYR	C9-C10-C11-C12
3	B	125	MYR	C9-C10-C11-C12
3	B	125	MYR	C2-C3-C4-C5
3	A	124	MYR	C11-C10-C9-C8
3	A	124	MYR	C5-C6-C7-C8
3	B	125	MYR	C11-C10-C9-C8
3	A	124	MYR	C2-C3-C4-C5
3	B	125	MYR	C3-C4-C5-C6
3	A	124	MYR	C10-C11-C12-C13
3	A	124	MYR	C3-C4-C5-C6
3	B	125	MYR	C5-C6-C7-C8
3	B	125	MYR	C6-C7-C8-C9
3	B	125	MYR	C11-C12-C13-C14
3	A	124	MYR	C11-C12-C13-C14
3	A	124	MYR	C1-C2-C3-C4
3	B	125	MYR	O2-C1-C2-C3
3	B	125	MYR	O1-C1-C2-C3
3	A	124	MYR	C7-C8-C9-C10
3	A	124	MYR	O2-C1-C2-C3

There are no ring outliers.

6 monomers are involved in 21 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	123	SO4	1	0
2	B	122	SO4	3	0
3	B	125	MYR	8	0
3	A	124	MYR	2	0
2	B	123	SO4	4	0
2	A	122	SO4	3	0

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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, 95th 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(Å ²)	Q<0.9
1	A	121/121 (100%)	0.13	4 (3%) 49 46	8, 22, 36, 41	0
1	B	121/121 (100%)	-0.05	1 (0%) 82 81	10, 20, 39, 44	0
All	All	242/242 (100%)	0.04	5 (2%) 63 61	8, 21, 38, 44	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	30	VAL	5.6
1	A	28	CYS	3.3
1	A	31	LEU	3.2
1	B	114	PHE	2.5
1	A	54	ILE	2.2

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 oligosaccharides 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, 95th 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	MYR	B	125	16/16	0.75	0.15	21,34,40,40	0
3	MYR	A	124	16/16	0.81	0.14	28,35,40,42	0
2	SO4	B	123	5/5	0.86	0.11	56,60,61,61	0
2	SO4	B	122	5/5	0.87	0.14	39,42,43,45	0
2	SO4	A	123	5/5	0.90	0.10	50,54,56,59	0
2	SO4	A	122	5/5	0.96	0.09	32,35,38,40	0

6.5 Other polymers [i](#)

There are no such residues in this entry.