



# Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 01:35 PM GMT

PDB ID : 1PP2  
Title : THE REFINED CRYSTAL STRUCTURE OF DIMERIC PHOSPHOLIPASE  
A2 AT 2.5 ANGSTROMS. ACCESS TO A SHIELDED CATALYTIC CEN-  
TER  
Authors : Brunie, S.; Sigler, P.B.  
Deposited on : 1986-03-10  
Resolution : 2.50 Å(reported)

This is a full wwPDB 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 <http://wwpdb.org/ValidationPDFNotes.html>

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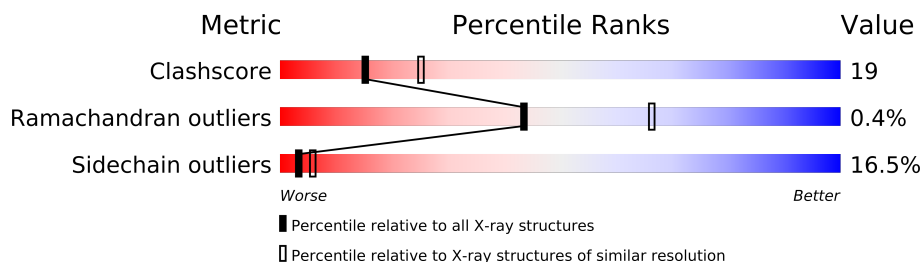
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 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 2.50 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	3562 (2.50-2.50)
Ramachandran outliers	78287	3480 (2.50-2.50)
Sidechain outliers	78261	3482 (2.50-2.50)

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  $\geq 3$ , 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	L	122	
1	R	122	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2029 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 CALCIUM-FREE PHOSPHOLIPASE A2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	R	122	Total	C	N	O	S	0	0	0
			946	591	154	186	15			
1	L	122	Total	C	N	O	S	0	0	0
			946	591	154	186	15			

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	L	74	Total	O	0	0
			74	74		
2	R	63	Total	O	0	0
			63	63		

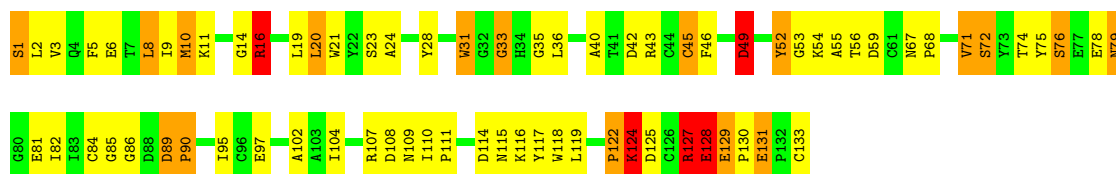
### 3 Residue-property plots

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.

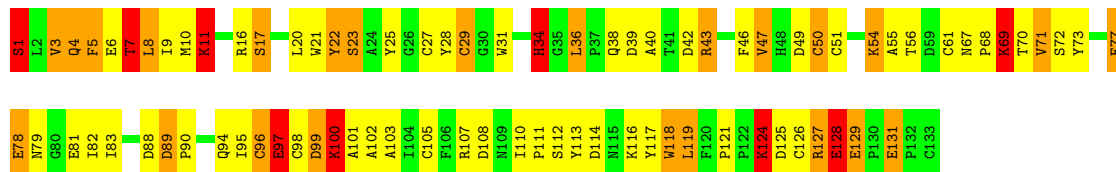
#### • Molecule 1: CALCIUM-FREE PHOSPHOLIPASE A2

Chain R: 



#### • Molecule 1: CALCIUM-FREE PHOSPHOLIPASE A2

Chain L: 



## 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, $\alpha$ , $\beta$ , $\gamma$	53.40Å 100.20Å 48.60Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	(Not available) – 2.50	Depositor
% Data completeness (in resolution range)	(Not available) ((Not available)-2.50)	Depositor
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	PROLSQ	Depositor
R, $R_{free}$	0.178 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2029	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	11.0	wwPDB-VP

## 5 Model quality

### 5.1 Standard geometry

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	L	1.55	3/973 (0.3%)	3.35	120/1320 (9.1%)
1	R	1.63	8/973 (0.8%)	3.29	107/1320 (8.1%)
All	All	1.59	11/1946 (0.6%)	3.32	227/2640 (8.6%)

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	L	0	1

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	R	1	SER	CB-OG	-7.45	1.32	1.42
1	L	6	GLU	CD-OE1	-6.71	1.18	1.25
1	L	1	SER	CB-OG	-6.38	1.33	1.42
1	R	6	GLU	CD-OE1	-6.26	1.18	1.25
1	R	43	ARG	CZ-NH2	6.20	1.41	1.33
1	R	86	GLY	N-CA	-5.74	1.37	1.46
1	R	28	TYR	CG-CD2	5.66	1.46	1.39
1	L	118	TRP	CG-CD1	5.60	1.44	1.36
1	R	108	ASP	C-N	-5.47	1.21	1.34
1	R	43	ARG	CD-NE	5.39	1.55	1.46
1	R	85	GLY	N-CA	5.31	1.54	1.46

All (227) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L	43	ARG	NE-CZ-NH1	35.69	138.15	120.30
1	L	16	ARG	NE-CZ-NH2	28.90	134.75	120.30
1	R	127	ARG	NE-CZ-NH1	24.91	132.75	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	R	49	ASP	CB-CG-OD1	-23.22	97.40	118.30
1	R	108	ASP	CB-CG-OD1	23.11	139.10	118.30
1	R	16	ARG	NE-CZ-NH1	21.54	131.07	120.30
1	L	88	ASP	CB-CG-OD1	20.70	136.93	118.30
1	R	128	GLU	OE1-CD-OE2	-19.14	100.33	123.30
1	L	89	ASP	CB-CG-OD1	15.87	132.58	118.30
1	L	127	ARG	NE-CZ-NH2	-15.35	112.63	120.30
1	R	108	ASP	CB-CG-OD2	-15.15	104.66	118.30
1	L	114	ASP	CB-CG-OD2	15.11	131.89	118.30
1	R	42	ASP	CB-CG-OD2	14.31	131.18	118.30
1	L	108	ASP	CB-CG-OD2	-13.92	105.77	118.30
1	R	128	GLU	CG-CD-OE1	13.92	146.13	118.30
1	R	16	ARG	NE-CZ-NH2	-13.69	113.46	120.30
1	R	127	ARG	NE-CZ-NH2	-13.42	113.59	120.30
1	L	129	GLU	CA-CB-CG	13.21	142.47	113.40
1	L	88	ASP	CB-CG-OD2	-12.78	106.80	118.30
1	R	16	ARG	CD-NE-CZ	12.77	141.48	123.60
1	L	16	ARG	NH1-CZ-NH2	-12.65	105.48	119.40
1	L	43	ARG	NH1-CZ-NH2	-12.39	105.77	119.40
1	R	97	GLU	OE1-CD-OE2	-12.23	108.62	123.30
1	L	97	GLU	CA-CB-CG	12.02	139.84	113.40
1	R	117	TYR	CB-CG-CD1	11.67	128.00	121.00
1	R	6	GLU	OE1-CD-OE2	-11.54	109.45	123.30
1	L	16	ARG	CG-CD-NE	11.42	135.78	111.80
1	R	55	ALA	CB-CA-C	-11.20	93.30	110.10
1	R	6	GLU	CG-CD-OE1	11.12	140.54	118.30
1	L	49	ASP	CB-CG-OD1	-10.64	108.73	118.30
1	R	125	ASP	CA-CB-CG	10.46	136.41	113.40
1	L	99	ASP	CB-CG-OD2	-10.46	108.89	118.30
1	L	43	ARG	NE-CZ-NH2	-10.40	115.10	120.30
1	L	108	ASP	CB-CG-OD1	10.39	127.65	118.30
1	R	125	ASP	CB-CG-OD1	10.36	127.62	118.30
1	R	28	TYR	CB-CG-CD2	-10.18	114.89	121.00
1	L	99	ASP	CB-CG-OD1	10.02	127.32	118.30
1	R	124	LYS	CB-CG-CD	9.89	137.31	111.60
1	L	98	CYS	CA-CB-SG	9.68	131.42	114.00
1	R	52	TYR	CB-CG-CD2	-9.63	115.22	121.00
1	L	124	LYS	CD-CE-NZ	-9.58	89.67	111.70
1	R	54	LYS	CA-CB-CG	9.55	134.41	113.40
1	L	43	ARG	CD-NE-CZ	9.47	136.86	123.60
1	L	117	TYR	CB-CG-CD1	9.32	126.59	121.00
1	R	49	ASP	OD1-CG-OD2	9.26	140.89	123.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L	61	CYS	O-C-N	9.19	137.41	122.70
1	R	28	TYR	CB-CG-CD1	9.07	126.44	121.00
1	L	25	TYR	CB-CG-CD1	8.92	126.35	121.00
1	R	124	LYS	CA-CB-CG	8.91	132.99	113.40
1	R	85	GLY	C-N-CA	8.88	140.95	122.30
1	R	20	LEU	CA-CB-CG	8.70	135.31	115.30
1	L	94	GLN	CB-CG-CD	8.62	134.02	111.60
1	R	129	GLU	CG-CD-OE2	-8.62	101.06	118.30
1	L	47	VAL	CA-CB-CG1	8.55	123.72	110.90
1	L	1	SER	N-CA-CB	-8.41	97.88	110.50
1	R	55	ALA	N-CA-CB	8.39	121.85	110.10
1	R	72	SER	N-CA-CB	-8.39	97.91	110.50
1	R	114	ASP	CB-CG-OD2	8.29	125.76	118.30
1	L	56	THR	CA-CB-CG2	8.15	123.81	112.40
1	L	11	LYS	CB-CG-CD	8.06	132.55	111.60
1	L	17	SER	CA-CB-OG	8.04	132.90	111.20
1	L	128	GLU	OE1-CD-OE2	8.02	132.92	123.30
1	R	128	GLU	O-C-N	-8.02	109.87	122.70
1	R	119	LEU	CB-CA-C	7.96	125.32	110.20
1	L	3	VAL	CA-CB-CG2	7.90	122.75	110.90
1	L	69	LYS	CA-CB-CG	7.84	130.65	113.40
1	L	49	ASP	OD1-CG-OD2	7.81	138.15	123.30
1	L	114	ASP	CB-CG-OD1	-7.78	111.30	118.30
1	L	22	TYR	CB-CG-CD2	-7.75	116.35	121.00
1	R	107	ARG	NE-CZ-NH1	-7.74	116.43	120.30
1	R	117	TYR	CG-CD1-CE1	7.71	127.47	121.30
1	L	8	LEU	CA-CB-CG	7.69	132.98	115.30
1	L	119	LEU	O-C-N	-7.62	110.50	122.70
1	R	54	LYS	CB-CG-CD	7.56	131.25	111.60
1	R	119	LEU	CA-CB-CG	7.55	132.66	115.30
1	L	29	CYS	CA-C-N	7.53	131.25	116.20
1	R	49	ASP	N-CA-CB	-7.50	97.11	110.60
1	R	108	ASP	C-N-CA	7.49	140.41	121.70
1	L	113	TYR	CG-CD1-CE1	7.45	127.26	121.30
1	R	89	ASP	CB-CG-OD1	7.44	125.00	118.30
1	L	78	GLU	C-N-CA	7.35	140.08	121.70
1	R	10	MET	N-CA-CB	-7.33	97.40	110.60
1	L	4	GLN	OE1-CD-NE2	-7.30	105.11	121.90
1	L	47	VAL	N-CA-CB	7.20	127.33	111.50
1	R	11	LYS	CD-CE-NZ	7.14	128.11	111.70
1	L	3	VAL	CB-CA-C	-7.14	97.84	111.40
1	L	6	GLU	OE1-CD-OE2	-7.08	114.80	123.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	R	125	ASP	OD1-CG-OD2	-7.08	109.84	123.30
1	R	46	PHE	O-C-N	7.04	133.97	122.70
1	L	43	ARG	O-C-N	-7.03	111.45	122.70
1	L	16	ARG	CD-NE-CZ	-6.96	113.86	123.60
1	L	16	ARG	O-C-N	6.94	133.80	122.70
1	R	79	ASN	N-CA-CB	-6.91	98.16	110.60
1	R	129	GLU	CG-CD-OE1	6.90	132.10	118.30
1	L	127	ARG	CB-CG-CD	-6.89	93.67	111.60
1	R	131	GLU	CA-CB-CG	6.88	128.54	113.40
1	R	8	LEU	O-C-N	-6.86	111.72	122.70
1	L	1	SER	CB-CA-C	-6.83	97.13	110.10
1	R	49	ASP	CB-CA-C	6.80	123.99	110.40
1	R	43	ARG	NE-CZ-NH1	-6.79	116.91	120.30
1	R	81	GLU	OE1-CD-OE2	6.71	131.35	123.30
1	L	116	LYS	CA-CB-CG	6.70	128.15	113.40
1	L	4	GLN	O-C-N	-6.64	112.08	122.70
1	L	125	ASP	CB-CG-OD2	6.63	124.27	118.30
1	R	108	ASP	CA-C-N	6.63	131.79	117.20
1	R	71	VAL	CA-CB-CG2	6.63	120.85	110.90
1	R	20	LEU	CB-CA-C	6.62	122.77	110.20
1	R	42	ASP	OD1-CG-OD2	-6.61	110.75	123.30
1	L	89	ASP	OD1-CG-OD2	-6.59	110.78	123.30
1	L	129	GLU	CB-CG-CD	-6.59	96.41	114.20
1	R	56	THR	CA-CB-CG2	6.58	121.61	112.40
1	R	116	LYS	CG-CD-CE	6.53	131.49	111.90
1	R	24	ALA	N-CA-CB	-6.52	100.98	110.10
1	L	101	ALA	CB-CA-C	-6.50	100.35	110.10
1	L	16	ARG	N-CA-CB	6.50	122.30	110.60
1	L	103	ALA	CB-CA-C	6.46	119.78	110.10
1	R	102	ALA	N-CA-CB	6.42	119.09	110.10
1	L	7	THR	CA-C-O	-6.42	106.62	120.10
1	R	128	GLU	CB-CA-C	6.38	123.17	110.40
1	R	117	TYR	CD1-CE1-CZ	-6.38	114.06	119.80
1	R	108	ASP	CA-C-O	-6.36	106.75	120.10
1	L	3	VAL	CG1-CB-CG2	-6.29	100.84	110.90
1	L	77	GLU	CA-C-O	-6.28	106.92	120.10
1	R	119	LEU	C-N-CA	6.27	137.37	121.70
1	R	45	CYS	O-C-N	-6.24	112.72	122.70
1	L	94	GLN	CG-CD-NE2	6.22	131.63	116.70
1	L	119	LEU	CA-CB-CG	6.22	129.60	115.30
1	L	107	ARG	NE-CZ-NH1	-6.21	117.20	120.30
1	R	109	ASN	O-C-N	6.18	132.59	122.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	R	54	LYS	CA-C-O	-6.18	107.12	120.10
1	L	55	ALA	N-CA-CB	6.13	118.68	110.10
1	L	113	TYR	CA-CB-CG	6.11	125.00	113.40
1	R	81	GLU	C-N-CA	6.10	136.95	121.70
1	L	116	LYS	CB-CA-C	6.05	122.50	110.40
1	L	97	GLU	O-C-N	-6.01	113.09	122.70
1	L	107	ARG	O-C-N	5.99	132.29	122.70
1	R	52	TYR	O-C-N	-5.99	113.02	123.20
1	L	34	HIS	CB-CA-C	-5.96	98.47	110.40
1	L	29	CYS	O-C-N	-5.94	113.10	123.20
1	R	115	ASN	O-C-N	-5.92	113.22	122.70
1	R	59	ASP	CB-CG-OD1	5.92	123.63	118.30
1	R	54	LYS	N-CA-CB	5.91	121.24	110.60
1	R	131	GLU	CG-CD-OE2	-5.90	106.50	118.30
1	L	23	SER	O-C-N	5.87	132.10	122.70
1	L	94	GLN	O-C-N	-5.87	113.30	122.70
1	L	113	TYR	CB-CG-CD1	5.85	124.51	121.00
1	R	10	MET	CB-CG-SD	5.85	129.94	112.40
1	R	3	VAL	CA-CB-CG2	5.84	119.67	110.90
1	R	42	ASP	CA-CB-CG	5.82	126.20	113.40
1	R	90	PRO	O-C-N	-5.81	113.40	122.70
1	R	116	LYS	CA-CB-CG	5.80	126.15	113.40
1	R	68	PRO	C-N-CA	5.78	136.15	121.70
1	R	122	PRO	O-C-N	5.77	131.94	122.70
1	L	3	VAL	N-CA-CB	5.76	124.17	111.50
1	L	5	PHE	CZ-CE2-CD2	-5.76	113.19	120.10
1	L	49	ASP	CB-CG-OD2	-5.75	113.12	118.30
1	R	128	GLU	N-CA-CB	-5.71	100.31	110.60
1	L	38	GLN	CG-CD-OE1	5.71	133.02	121.60
1	L	96	CYS	CA-CB-SG	5.70	124.25	114.00
1	L	119	LEU	CA-C-O	5.66	131.98	120.10
1	R	43	ARG	NE-CZ-NH2	5.65	123.12	120.30
1	R	31	TRP	CA-CB-CG	5.64	124.41	113.70
1	R	119	LEU	N-CA-CB	-5.64	99.12	110.40
1	R	40	ALA	CB-CA-C	5.64	118.56	110.10
1	L	51	CYS	CA-CB-SG	5.63	124.14	114.00
1	R	71	VAL	C-N-CA	5.60	135.70	121.70
1	R	117	TYR	CB-CA-C	5.60	121.60	110.40
1	R	67	ASN	CA-CB-CG	5.60	125.71	113.40
1	L	73	TYR	CD1-CE1-CZ	5.59	124.83	119.80
1	R	33	GLY	CA-C-O	5.58	130.65	120.60
1	L	10	MET	N-CA-CB	-5.56	100.59	110.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	R	133	CYS	CA-C-O	-5.56	108.43	120.10
1	R	59	ASP	CA-C-N	5.52	129.34	117.20
1	R	131	GLU	CG-CD-OE1	5.51	129.33	118.30
1	L	112	SER	N-CA-CB	5.51	118.77	110.50
1	L	4	GLN	C-N-CA	5.50	135.45	121.70
1	L	4	GLN	CG-CD-OE1	5.49	132.58	121.60
1	L	29	CYS	C-N-CA	5.48	133.81	122.30
1	L	5	PHE	CG-CD2-CE2	5.46	126.81	120.80
1	R	20	LEU	CA-C-N	5.46	129.21	117.20
1	R	16	ARG	O-C-N	5.45	131.41	122.70
1	L	102	ALA	O-C-N	5.45	131.41	122.70
1	R	20	LEU	CA-C-O	-5.43	108.69	120.10
1	L	128	GLU	CA-C-N	-5.43	105.26	117.20
1	R	97	GLU	O-C-N	-5.42	114.03	122.70
1	R	84	CYS	O-C-N	5.41	132.40	123.20
1	L	42	ASP	CB-CG-OD2	-5.41	113.43	118.30
1	L	10	MET	O-C-N	-5.41	114.05	122.70
1	R	85	GLY	N-CA-C	-5.39	99.63	113.10
1	L	99	ASP	C-N-CA	5.37	135.13	121.70
1	R	117	TYR	CG-CD2-CE2	5.36	125.59	121.30
1	L	108	ASP	CB-CA-C	-5.34	99.72	110.40
1	L	77	GLU	CA-C-N	5.34	128.95	117.20
1	L	119	LEU	CB-CA-C	5.32	120.31	110.20
1	R	1	SER	N-CA-C	5.32	125.35	111.00
1	L	113	TYR	CD1-CE1-CZ	-5.31	115.02	119.80
1	L	102	ALA	N-CA-CB	5.29	117.51	110.10
1	R	127	ARG	NH1-CZ-NH2	-5.28	113.59	119.40
1	R	75	TYR	CB-CG-CD1	-5.27	117.84	121.00
1	L	114	ASP	CB-CA-C	5.26	120.93	110.40
1	R	54	LYS	CA-C-N	5.26	128.78	117.20
1	R	115	ASN	C-N-CA	5.26	134.85	121.70
1	R	3	VAL	CG1-CB-CG2	-5.24	102.51	110.90
1	L	22	TYR	N-CA-CB	-5.24	101.16	110.60
1	L	7	THR	O-C-N	5.23	131.06	122.70
1	L	34	HIS	O-C-N	5.23	132.09	123.20
1	L	128	GLU	CA-C-O	5.21	131.04	120.10
1	R	117	TYR	CB-CG-CD2	-5.20	117.88	121.00
1	L	107	ARG	CA-CB-CG	5.20	124.84	113.40
1	L	27	CYS	CB-CA-C	-5.19	100.02	110.40
1	L	39	ASP	CB-CG-OD2	-5.18	113.64	118.30
1	L	105	CYS	N-CA-CB	5.18	119.92	110.60
1	L	125	ASP	OD1-CG-OD2	-5.17	113.47	123.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L	10	MET	CB-CA-C	5.17	120.73	110.40
1	R	124	LYS	C-N-CA	5.16	134.60	121.70
1	L	81	GLU	O-C-N	5.16	130.96	122.70
1	L	78	GLU	N-CA-CB	5.15	119.88	110.60
1	L	117	TYR	CB-CG-CD2	-5.15	117.91	121.00
1	L	99	ASP	CB-CA-C	5.15	120.69	110.40
1	L	100	LYS	CB-CA-C	-5.14	100.11	110.40
1	L	77	GLU	N-CA-C	-5.12	97.17	111.00
1	L	39	ASP	CB-CG-OD1	5.12	122.91	118.30
1	L	28	TYR	N-CA-CB	-5.11	101.40	110.60
1	L	50	CYS	CA-C-O	-5.06	109.47	120.10
1	L	10	MET	CA-CB-CG	-5.05	104.71	113.30
1	L	55	ALA	CA-C-O	-5.05	109.50	120.10
1	R	1	SER	CA-C-O	5.01	130.62	120.10

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	L	127	ARG	Sidechain

## 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	L	946	0	853	50	0
1	R	946	0	852	26	0
2	L	74	0	0	2	0
2	R	63	0	0	1	0
All	All	2029	0	1705	68	0

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 19.

All (68) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:R:118:TRP:HE1	1:L:124:LYS:NZ	1.49	1.09
1:L:36:LEU:HD12	1:L:128:GLU:HG3	1.42	1.01
1:L:11:LYS:HE3	1:L:82:ILE:HD11	1.46	0.94
1:R:53:GLY:HA2	1:L:70:THR:HG21	1.58	0.85
1:L:68:PRO:HA	1:L:95:ILE:HD13	1.61	0.83
1:R:36:LEU:N	1:R:128:GLU:OE2	2.15	0.80
1:L:124:LYS:H	1:L:124:LYS:CD	1.96	0.78
1:L:11:LYS:HE3	1:L:82:ILE:CD1	2.15	0.76
1:R:78:GLU:O	1:R:79:ASN:HB2	1.84	0.74
1:L:68:PRO:HG3	1:L:95:ILE:HD11	1.72	0.72
1:R:89:ASP:HB2	1:R:90:PRO:HD2	1.72	0.71
1:L:36:LEU:HD12	1:L:128:GLU:CG	2.17	0.71
1:L:124:LYS:H	1:L:124:LYS:HD3	1.56	0.70
1:L:11:LYS:CE	1:L:82:ILE:HD11	2.22	0.69
1:L:40:ALA:HA	1:L:43:ARG:HD2	1.74	0.69
1:L:121:PRO:HB3	1:L:124:LYS:HE2	1.76	0.67
1:L:5:PHE:CE2	1:L:9:ILE:HD11	2.30	0.67
1:R:49:ASP:OD1	2:R:573:HOH:O	2.14	0.65
1:L:34:HIS:HB2	1:L:126:CYS:O	1.97	0.65
1:R:110:ILE:N	1:R:111:PRO:CD	2.61	0.63
1:L:97:GLU:OE1	1:L:100:LYS:NZ	2.33	0.62
1:L:110:ILE:N	1:L:111:PRO:HD2	2.16	0.60
1:L:17:SER:O	1:L:21:TRP:HB2	2.02	0.60
1:R:23:SER:HA	1:L:31:TRP:CE3	2.38	0.59
1:L:5:PHE:CZ	1:L:9:ILE:HD11	2.37	0.59
1:L:3:VAL:O	1:L:3:VAL:HG12	2.02	0.58
1:L:4:GLN:OE1	1:L:72:SER:HA	2.04	0.57
1:L:124:LYS:CD	1:L:124:LYS:N	2.69	0.55
1:R:36:LEU:HB2	1:R:128:GLU:OE2	2.06	0.55
1:R:118:TRP:NE1	1:L:124:LYS:NZ	2.24	0.55
1:L:11:LYS:NZ	2:L:567:HOH:O	2.40	0.55
1:R:129:GLU:HG3	1:R:130:PRO:HD2	1.88	0.54
1:L:110:ILE:N	1:L:111:PRO:CD	2.72	0.52
1:R:45:CYS:O	1:R:49:ASP:HB2	2.10	0.51
1:R:124:LYS:O	1:R:127:ARG:HD2	2.11	0.51
1:R:31:TRP:CE3	1:L:23:SER:HA	2.46	0.51
1:L:3:VAL:O	1:L:7:THR:HG23	2.10	0.50
1:R:35:GLY:HA2	1:R:128:GLU:HB2	1.94	0.49
1:L:78:GLU:HG2	1:L:83:ILE:HD13	1.95	0.49
1:R:16:ARG:HB3	1:R:21:TRP:CD1	2.47	0.49
1:R:52:TYR:O	1:L:67:ASN:ND2	2.46	0.49
1:R:76:SER:O	1:R:82:ILE:HA	2.13	0.48
1:R:1:SER:N	1:R:71:VAL:O	2.23	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:L:121:PRO:CB	1:L:124:LYS:HD3	2.44	0.48
1:L:50:CYS:O	1:L:54:LYS:HD2	2.14	0.48
1:L:46:PHE:CZ	1:L:131:GLU:HG3	2.48	0.47
1:L:3:VAL:CG1	1:L:3:VAL:O	2.63	0.47
1:L:68:PRO:HG3	1:L:95:ILE:CD1	2.43	0.47
1:L:110:ILE:HG22	1:L:111:PRO:N	2.29	0.46
1:L:1:SER:N	1:L:71:VAL:O	2.38	0.46
1:L:68:PRO:CA	1:L:95:ILE:HD13	2.40	0.46
1:R:33:GLY:O	1:R:130:PRO:HG3	2.16	0.45
1:R:5:PHE:CE1	1:R:9:ILE:HD11	2.52	0.45
1:L:5:PHE:O	1:L:9:ILE:HG13	2.18	0.44
1:L:3:VAL:O	1:L:7:THR:CG2	2.65	0.44
1:L:1:SER:HB3	1:L:4:GLN:H	1.83	0.44
1:L:89:ASP:HA	1:L:90:PRO:HD2	1.81	0.44
1:R:14:GLY:O	1:R:16:ARG:HG3	2.17	0.44
1:R:104:ILE:HG21	1:R:104:ILE:HD13	1.83	0.43
1:L:69:LYS:HE3	2:L:535:HOH:O	2.20	0.42
1:L:96:CYS:O	1:L:99:ASP:HB2	2.20	0.42
1:R:127:ARG:O	1:R:128:GLU:C	2.59	0.42
1:L:70:THR:O	1:L:71:VAL:C	2.58	0.41
1:R:122:PRO:HG2	1:L:118:TRP:CZ2	2.55	0.41
1:R:118:TRP:CE2	1:L:124:LYS:NZ	2.88	0.40
1:L:110:ILE:HG22	1:L:111:PRO:HD3	2.03	0.40
1:L:22:TYR:O	1:L:29:CYS:HB3	2.22	0.40
1:L:110:ILE:HB	1:L:111:PRO:HD3	2.03	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	L	120/122 (98%)	111 (92%)	8 (7%)	1 (1%)	27	46
1	R	120/122 (98%)	113 (94%)	7 (6%)	0	100	100
All	All	240/244 (98%)	224 (93%)	15 (6%)	1 (0%)	43	66

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	L	79	ASN

### 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	L	103/103 (100%)	84 (82%)	19 (18%)	2	4
1	R	103/103 (100%)	88 (85%)	15 (15%)	5	8
All	All	206/206 (100%)	172 (84%)	34 (16%)	3	6

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	R	2	LEU
1	R	8	LEU
1	R	10	MET
1	R	16	ARG
1	R	19	LEU
1	R	20	LEU
1	R	49	ASP
1	R	72	SER
1	R	74	THR
1	R	76	SER
1	R	95	ILE
1	R	124	LYS
1	R	127	ARG
1	R	128	GLU
1	R	131	GLU
1	L	1	SER
1	L	7	THR
1	L	8	LEU
1	L	11	LYS
1	L	20	LEU
1	L	34	HIS
1	L	36	LEU
1	L	47	VAL

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Mol	Chain	Res	Type
1	L	54	LYS
1	L	69	LYS
1	L	71	VAL
1	L	77	GLU
1	L	97	GLU
1	L	100	LYS
1	L	119	LEU
1	L	124	LYS
1	L	128	GLU
1	L	129	GLU
1	L	131	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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.