



Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2024 – 10:35 AM EDT

PDB ID : 2BP2
Title : THE STRUCTURE OF BOVINE PANCREATIC PROPHOSPHOLIPASE
A2 AT 3.0 ANGSTROMS RESOLUTION
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Deposited on : 1981-06-05
Resolution : 3.00 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	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

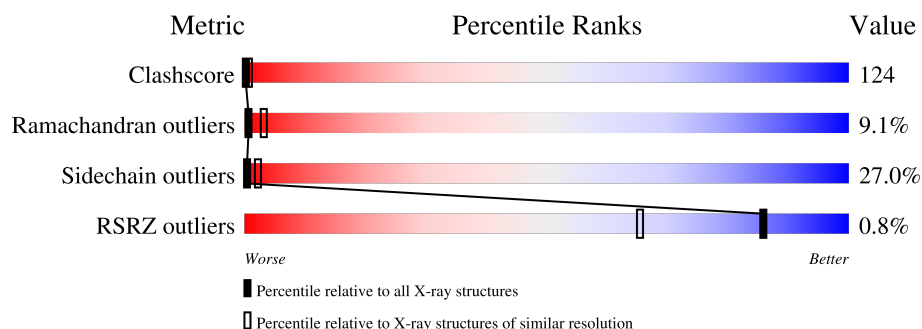
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)
RSRZ outliers	127900	1990 (3.00-3.00)

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	130	<div> <div></div> <div>22%</div> <div>34%</div> <div>31%</div> <div>8%</div> <div>5%</div> </div>

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 956 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.

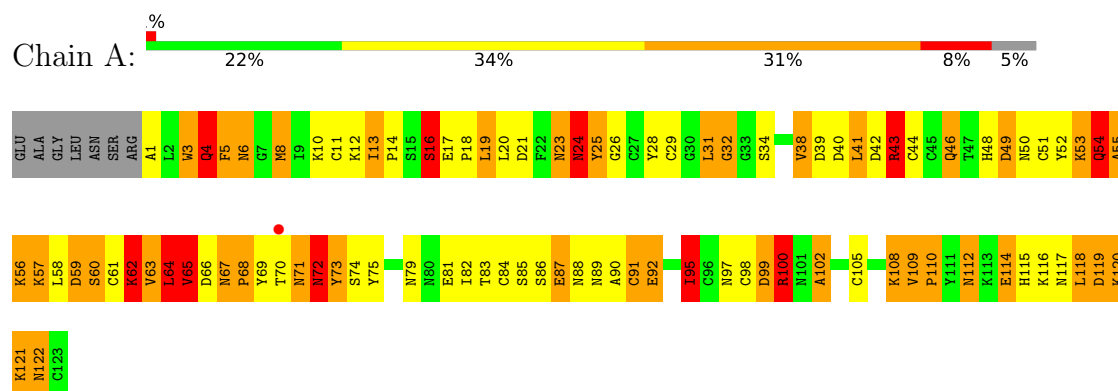
- Molecule 1 is a protein called PHOSPHOLIPASE A2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	123	956	586	164	191	15	0	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: PHOSPHOLIPASE A2



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	46.95Å 46.95Å 102.04Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	(Not available) – 3.00 7.02 – 2.98	Depositor EDS
% Data completeness (in resolution range)	(Not available) ((Not available)-3.00) 99.1 (7.02-2.98)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$	-	Xtriage
Refinement program	PROLSQ	Depositor
R, R_{free}	0.219 , (Not available) 0.211 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	31.0	Xtriage
Anisotropy	0.485	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.43 , 140.9	EDS
L-test for twinning ¹	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.059 for -h,-k,l	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	956	wwPDB-VP
Average B, all atoms (Å ²)	13.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.83% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Theoretical 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 [i](#)

5.1 Standard geometry [i](#)

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	1.44	5/975 (0.5%)	2.55	49/1317 (3.7%)

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	A	0	1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	3	TRP	NE1-CE2	-7.56	1.27	1.37
1	A	114	GLU	CD-OE2	-6.40	1.18	1.25
1	A	91	CYS	CB-SG	5.57	1.91	1.82
1	A	3	TRP	CD1-NE1	5.47	1.47	1.38
1	A	28	TYR	CD1-CE1	5.07	1.47	1.39

All (49) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	43	ARG	CD-NE-CZ	28.70	163.79	123.60
1	A	100	ARG	NE-CZ-NH2	18.49	129.54	120.30
1	A	87	GLU	OE1-CD-OE2	15.09	141.41	123.30
1	A	100	ARG	NE-CZ-NH1	-13.88	113.36	120.30
1	A	42	ASP	CB-CG-OD1	12.39	129.45	118.30
1	A	59	ASP	CB-CG-OD1	10.05	127.35	118.30
1	A	4	GLN	CA-CB-CG	9.63	134.59	113.40
1	A	71	ASN	CA-CB-CG	-8.94	93.73	113.40
1	A	24	ASN	CA-CB-CG	-8.74	94.18	113.40
1	A	87	GLU	CG-CD-OE1	-8.62	101.07	118.30
1	A	92	GLU	OE1-CD-OE2	8.45	133.43	123.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	49	ASP	CB-CG-OD1	8.39	125.85	118.30
1	A	39	ASP	CB-CG-OD2	-7.99	111.11	118.30
1	A	25	TYR	CB-CG-CD1	-7.72	116.37	121.00
1	A	64	LEU	N-CA-C	-7.32	91.23	111.00
1	A	114	GLU	OE1-CD-OE2	7.27	132.03	123.30
1	A	4	GLN	CA-C-O	-7.26	104.85	120.10
1	A	92	GLU	CG-CD-OE1	-7.25	103.81	118.30
1	A	97	ASN	CA-CB-CG	-7.23	97.50	113.40
1	A	4	GLN	N-CA-CB	7.15	123.47	110.60
1	A	122	ASN	CA-CB-CG	-7.11	97.77	113.40
1	A	46	GLN	CA-CB-CG	-6.89	98.23	113.40
1	A	39	ASP	CB-CG-OD1	6.74	124.37	118.30
1	A	54	GLN	CA-CB-CG	-6.43	99.25	113.40
1	A	71	ASN	N-CA-CB	6.38	122.08	110.60
1	A	67	ASN	CA-CB-CG	-6.26	99.62	113.40
1	A	13	ILE	CA-CB-CG2	-6.18	98.54	110.90
1	A	40	ASP	CB-CG-OD2	-6.16	112.75	118.30
1	A	43	ARG	NE-CZ-NH2	6.14	123.37	120.30
1	A	64	LEU	CA-C-N	-6.09	103.80	117.20
1	A	89	ASN	CA-CB-CG	-5.93	100.36	113.40
1	A	87	GLU	CA-CB-CG	-5.90	100.41	113.40
1	A	89	ASN	OD1-CG-ND2	5.67	134.94	121.90
1	A	3	TRP	NE1-CE2-CZ2	-5.55	124.29	130.40
1	A	99	ASP	CB-CG-OD2	-5.52	113.33	118.30
1	A	100	ARG	CD-NE-CZ	-5.52	115.87	123.60
1	A	5	PHE	N-CA-CB	5.51	120.52	110.60
1	A	99	ASP	CB-CG-OD1	5.44	123.19	118.30
1	A	38	VAL	CA-CB-CG2	5.43	119.05	110.90
1	A	79	ASN	CB-CG-OD1	5.33	132.26	121.60
1	A	72	ASN	CA-CB-CG	-5.32	101.69	113.40
1	A	16	SER	C-N-CA	5.29	134.92	121.70
1	A	110	PRO	CA-N-CD	-5.21	104.20	111.50
1	A	95	ILE	CA-CB-CG2	-5.14	100.61	110.90
1	A	42	ASP	CB-CG-OD2	-5.14	113.67	118.30
1	A	112	ASN	N-CA-CB	-5.13	101.37	110.60
1	A	65	VAL	N-CA-CB	-5.10	100.28	111.50
1	A	73	TYR	CB-CG-CD1	-5.09	117.94	121.00
1	A	102	ALA	CB-CA-C	-5.07	102.50	110.10

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	21	ASP	Mainchain

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	956	0	885	228	11
All	All	956	0	885	228	11

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 124.

All (228) 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:43:ARG:HH12	1:A:46:GLN:CG	1.04	1.55
1:A:43:ARG:NH1	1:A:46:GLN:CB	1.74	1.50
1:A:43:ARG:NH1	1:A:46:GLN:CG	1.73	1.45
1:A:67:ASN:CB	1:A:70:THR:OG1	1.78	1.28
1:A:41:LEU:HD12	1:A:41:LEU:O	1.35	1.26
1:A:43:ARG:NH1	1:A:46:GLN:HB3	1.36	1.23
1:A:62:LYS:O	1:A:65:VAL:CA	1.87	1.22
1:A:58:LEU:HD12	1:A:61:CYS:SG	1.82	1.20
1:A:60:SER:O	1:A:64:LEU:CD1	1.93	1.16
1:A:64:LEU:O	1:A:65:VAL:HG13	1.40	1.16
1:A:3:TRP:CH2	1:A:4:GLN:HG3	1.81	1.16
1:A:3:TRP:CZ2	1:A:4:GLN:HG3	1.82	1.14
1:A:64:LEU:C	1:A:65:VAL:HG13	1.62	1.13
1:A:5:PHE:CD1	1:A:99:ASP:HB3	1.86	1.10
1:A:64:LEU:C	1:A:65:VAL:CG1	2.21	1.09
1:A:67:ASN:HB3	1:A:70:THR:OG1	0.89	1.06
1:A:43:ARG:NH1	1:A:46:GLN:HG2	1.45	1.05
1:A:62:LYS:O	1:A:65:VAL:HA	1.50	1.04
1:A:43:ARG:HH11	1:A:46:GLN:CB	1.47	1.04
1:A:13:ILE:HG22	1:A:16:SER:HB2	1.39	1.03
1:A:56:LYS:HD3	1:A:68:PRO:HG3	1.38	1.03
1:A:13:ILE:O	1:A:16:SER:HB2	1.59	1.02

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:55:ALA:O	1:A:58:LEU:HG	1.57	1.02
1:A:121:LYS:HA	1:A:121:LYS:HE3	1.43	1.01
1:A:59:ASP:O	1:A:62:LYS:HB2	1.61	1.00
1:A:60:SER:O	1:A:64:LEU:HD11	1.61	0.97
1:A:55:ALA:O	1:A:58:LEU:CG	2.14	0.95
1:A:43:ARG:HH12	1:A:46:GLN:CB	1.54	0.94
1:A:62:LYS:O	1:A:65:VAL:N	2.01	0.93
1:A:1:ALA:N	1:A:71:ASN:O	2.02	0.93
1:A:8:MET:HE1	1:A:99:ASP:HB2	1.48	0.92
1:A:3:TRP:HE1	1:A:72:ASN:ND2	1.67	0.92
1:A:13:ILE:HG22	1:A:13:ILE:O	1.71	0.91
1:A:60:SER:C	1:A:64:LEU:HD11	1.89	0.91
1:A:41:LEU:O	1:A:44:CYS:HB2	1.70	0.91
1:A:1:ALA:H1	1:A:71:ASN:N	1.70	0.90
1:A:67:ASN:OD1	1:A:70:THR:N	2.05	0.90
1:A:121:LYS:HE3	1:A:121:LYS:CA	2.02	0.88
1:A:38:VAL:HG11	1:A:118:LEU:CD1	2.05	0.87
1:A:3:TRP:HE1	1:A:72:ASN:HD21	0.90	0.87
1:A:52:TYR:OH	1:A:99:ASP:OD2	1.92	0.86
1:A:119:ASP:OD1	1:A:121:LYS:HB2	1.74	0.86
1:A:5:PHE:HE1	1:A:99:ASP:O	1.58	0.86
1:A:56:LYS:HD3	1:A:68:PRO:CG	2.05	0.86
1:A:8:MET:CE	1:A:99:ASP:HB2	2.06	0.85
1:A:1:ALA:H1	1:A:71:ASN:H	1.24	0.85
1:A:41:LEU:HD12	1:A:41:LEU:C	1.96	0.85
1:A:43:ARG:HH12	1:A:46:GLN:HG2	0.69	0.85
1:A:66:ASP:CG	1:A:67:ASN:N	2.30	0.83
1:A:5:PHE:CE1	1:A:99:ASP:HB3	2.12	0.83
1:A:43:ARG:HH11	1:A:46:GLN:HB3	0.68	0.83
1:A:119:ASP:O	1:A:121:LYS:N	2.12	0.83
1:A:1:ALA:CB	1:A:71:ASN:O	2.27	0.83
1:A:120:LYS:N	1:A:120:LYS:HD3	1.93	0.83
1:A:59:ASP:HA	1:A:62:LYS:HD3	1.60	0.82
1:A:58:LEU:O	1:A:62:LYS:HD2	1.80	0.81
1:A:49:ASP:O	1:A:53:LYS:HG2	1.80	0.80
1:A:67:ASN:HB3	1:A:70:THR:HG1	1.47	0.78
1:A:3:TRP:CE2	1:A:4:GLN:HG3	2.19	0.77
1:A:4:GLN:N	1:A:4:GLN:OE1	2.16	0.77
1:A:5:PHE:CD1	1:A:99:ASP:CB	2.68	0.76
1:A:60:SER:O	1:A:64:LEU:HG	1.86	0.76
1:A:60:SER:O	1:A:64:LEU:CG	2.33	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:26:GLY:HA2	1:A:118:LEU:HD12	1.65	0.76
1:A:6:ASN:ND2	1:A:19:LEU:HG	2.02	0.74
1:A:60:SER:O	1:A:64:LEU:HD12	1.87	0.74
1:A:3:TRP:CZ2	1:A:4:GLN:CG	2.68	0.74
1:A:119:ASP:OD1	1:A:121:LYS:N	2.19	0.74
1:A:67:ASN:OD1	1:A:70:THR:HG23	1.89	0.73
1:A:67:ASN:HB3	1:A:70:THR:CB	2.16	0.73
1:A:13:ILE:O	1:A:16:SER:CB	2.35	0.73
1:A:58:LEU:CD1	1:A:61:CYS:SG	2.71	0.73
1:A:1:ALA:N	1:A:71:ASN:N	2.36	0.72
1:A:3:TRP:CZ3	1:A:4:GLN:HG3	2.24	0.72
1:A:1:ALA:HB3	1:A:71:ASN:O	1.87	0.72
1:A:10:LYS:HG2	1:A:18:PRO:HD3	1.71	0.72
1:A:43:ARG:CZ	1:A:46:GLN:HG2	2.20	0.72
1:A:60:SER:C	1:A:64:LEU:CD1	2.53	0.71
1:A:121:LYS:HA	1:A:121:LYS:CE	2.20	0.71
1:A:43:ARG:NH1	1:A:46:GLN:CD	2.42	0.71
1:A:64:LEU:C	1:A:66:ASP:N	2.40	0.70
1:A:13:ILE:CG2	1:A:16:SER:HB2	2.20	0.70
1:A:24:ASN:HB3	1:A:117:ASN:ND2	2.07	0.69
1:A:5:PHE:CE1	1:A:99:ASP:O	2.45	0.69
1:A:121:LYS:HE3	1:A:121:LYS:N	2.06	0.69
1:A:43:ARG:NH1	1:A:46:GLN:OE1	2.26	0.69
1:A:1:ALA:H3	1:A:71:ASN:C	1.96	0.68
1:A:43:ARG:CZ	1:A:46:GLN:OE1	2.41	0.68
1:A:119:ASP:OD1	1:A:119:ASP:O	2.12	0.67
1:A:6:ASN:HD22	1:A:19:LEU:CD2	2.08	0.67
1:A:118:LEU:HD22	1:A:122:ASN:HB2	1.78	0.65
1:A:23:ASN:ND2	1:A:24:ASN:H	1.94	0.65
1:A:5:PHE:HD1	1:A:99:ASP:HB3	1.56	0.65
1:A:55:ALA:O	1:A:58:LEU:CB	2.45	0.64
1:A:63:VAL:HG23	1:A:63:VAL:O	1.98	0.63
1:A:56:LYS:CD	1:A:68:PRO:HG3	2.22	0.63
1:A:10:LYS:HD2	1:A:18:PRO:CD	2.29	0.63
1:A:59:ASP:HA	1:A:62:LYS:CD	2.27	0.63
1:A:120:LYS:C	1:A:121:LYS:HE3	2.19	0.63
1:A:55:ALA:C	1:A:58:LEU:HG	2.19	0.63
1:A:61:CYS:O	1:A:66:ASP:N	2.33	0.62
1:A:43:ARG:NH1	1:A:46:GLN:CA	2.61	0.61
1:A:55:ALA:HA	1:A:58:LEU:HD11	1.83	0.60
1:A:51:CYS:O	1:A:54:GLN:HB3	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:105:CYS:O	1:A:108:LYS:HB2	2.01	0.60
1:A:114:GLU:OE2	1:A:114:GLU:N	2.33	0.60
1:A:24:ASN:HB3	1:A:116:LYS:HZ2	1.67	0.59
1:A:62:LYS:C	1:A:65:VAL:H	2.04	0.59
1:A:12:LYS:C	1:A:14:PRO:HD3	2.22	0.59
1:A:38:VAL:HG11	1:A:118:LEU:HD11	1.83	0.59
1:A:10:LYS:CG	1:A:18:PRO:HD3	2.31	0.59
1:A:122:ASN:N	1:A:122:ASN:ND2	2.45	0.59
1:A:59:ASP:C	1:A:62:LYS:HB2	2.23	0.58
1:A:12:LYS:O	1:A:14:PRO:HD3	2.03	0.58
1:A:64:LEU:O	1:A:65:VAL:CG1	2.30	0.58
1:A:8:MET:O	1:A:11:CYS:HB3	2.04	0.58
1:A:55:ALA:HA	1:A:58:LEU:CD1	2.34	0.57
1:A:24:ASN:HB2	1:A:116:LYS:NZ	2.19	0.57
1:A:6:ASN:HD22	1:A:19:LEU:HD23	1.69	0.57
1:A:13:ILE:N	1:A:14:PRO:HD3	2.15	0.57
1:A:10:LYS:HD2	1:A:18:PRO:HD2	1.87	0.56
1:A:38:VAL:HG21	1:A:115:HIS:HB3	1.86	0.56
1:A:4:GLN:OE1	1:A:4:GLN:CA	2.53	0.56
1:A:24:ASN:CB	1:A:116:LYS:HZ2	2.19	0.56
1:A:43:ARG:HH12	1:A:46:GLN:CA	2.18	0.55
1:A:73:TYR:HB3	1:A:92:GLU:HB2	1.88	0.55
1:A:71:ASN:OD1	1:A:92:GLU:HB2	2.07	0.55
1:A:3:TRP:CH2	1:A:4:GLN:CG	2.74	0.55
1:A:102:ALA:O	1:A:105:CYS:HB3	2.07	0.55
1:A:71:ASN:OD1	1:A:92:GLU:CB	2.56	0.54
1:A:64:LEU:N	1:A:65:VAL:HG12	2.22	0.54
1:A:10:LYS:CG	1:A:18:PRO:CD	2.86	0.54
1:A:73:TYR:OH	1:A:99:ASP:OD1	2.18	0.54
1:A:43:ARG:HH12	1:A:46:GLN:C	2.11	0.53
1:A:64:LEU:C	1:A:66:ASP:H	2.09	0.53
1:A:5:PHE:CE1	1:A:99:ASP:CB	2.87	0.53
1:A:24:ASN:O	1:A:116:LYS:HD2	2.07	0.53
1:A:38:VAL:HG11	1:A:118:LEU:HD13	1.85	0.53
1:A:119:ASP:OD1	1:A:119:ASP:C	2.45	0.53
1:A:10:LYS:HD2	1:A:18:PRO:HD3	1.89	0.53
1:A:24:ASN:CB	1:A:116:LYS:NZ	2.71	0.53
1:A:3:TRP:CE2	1:A:4:GLN:CG	2.90	0.53
1:A:67:ASN:OD1	1:A:69:TYR:C	2.47	0.52
1:A:8:MET:CE	1:A:99:ASP:CB	2.83	0.52
1:A:38:VAL:HG23	1:A:115:HIS:ND1	2.25	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:119:ASP:OD1	1:A:121:LYS:CB	2.53	0.51
1:A:23:ASN:ND2	1:A:24:ASN:N	2.57	0.51
1:A:3:TRP:CG	1:A:4:GLN:N	2.78	0.51
1:A:25:TYR:HB3	1:A:29:CYS:HB2	1.91	0.51
1:A:10:LYS:HD2	1:A:17:GLU:HA	1.92	0.51
1:A:19:LEU:HD23	1:A:19:LEU:N	2.26	0.51
1:A:19:LEU:O	1:A:23:ASN:HB2	2.10	0.51
1:A:118:LEU:HD22	1:A:122:ASN:CB	2.40	0.51
1:A:1:ALA:HB3	1:A:4:GLN:HE22	1.76	0.51
1:A:68:PRO:O	1:A:71:ASN:N	2.41	0.51
1:A:68:PRO:C	1:A:70:THR:H	2.15	0.50
1:A:66:ASP:OD2	1:A:67:ASN:N	2.44	0.50
1:A:4:GLN:NE2	1:A:72:ASN:HD22	2.09	0.50
1:A:10:LYS:CD	1:A:18:PRO:CD	2.89	0.50
1:A:38:VAL:HG21	1:A:115:HIS:CB	2.41	0.50
1:A:119:ASP:C	1:A:121:LYS:H	2.14	0.50
1:A:6:ASN:ND2	1:A:19:LEU:CG	2.74	0.49
1:A:6:ASN:HD22	1:A:19:LEU:CG	2.26	0.49
1:A:66:ASP:CG	1:A:71:ASN:HD22	2.16	0.49
1:A:67:ASN:O	1:A:71:ASN:N	2.45	0.49
1:A:64:LEU:C	1:A:65:VAL:HG12	2.21	0.49
1:A:95:ILE:O	1:A:95:ILE:CG2	2.60	0.49
1:A:55:ALA:O	1:A:58:LEU:HB2	2.11	0.49
1:A:60:SER:CA	1:A:64:LEU:HD11	2.42	0.48
1:A:6:ASN:O	1:A:10:LYS:HG3	2.14	0.48
1:A:17:GLU:O	1:A:20:LEU:HB2	2.13	0.48
1:A:1:ALA:CA	1:A:71:ASN:O	2.61	0.48
1:A:38:VAL:HG21	1:A:115:HIS:CG	2.48	0.48
1:A:84:CYS:SG	1:A:100:ARG:HG3	2.53	0.48
1:A:59:ASP:O	1:A:62:LYS:CB	2.49	0.47
1:A:43:ARG:CZ	1:A:46:GLN:CG	2.79	0.47
1:A:55:ALA:O	1:A:58:LEU:CD1	2.62	0.47
1:A:54:GLN:OE1	1:A:54:GLN:HA	1.96	0.47
1:A:119:ASP:C	1:A:121:LYS:N	2.66	0.47
1:A:64:LEU:O	1:A:66:ASP:N	2.48	0.46
1:A:8:MET:HE3	1:A:99:ASP:HB2	1.92	0.46
1:A:63:VAL:O	1:A:64:LEU:HD23	2.14	0.46
1:A:62:LYS:O	1:A:65:VAL:CB	2.61	0.46
1:A:10:LYS:CD	1:A:18:PRO:HD3	2.46	0.46
1:A:4:GLN:NE2	1:A:72:ASN:ND2	2.65	0.45
1:A:108:LYS:N	1:A:108:LYS:HD3	2.27	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:52:TYR:O	1:A:56:LYS:HG3	2.16	0.45
1:A:46:GLN:O	1:A:46:GLN:HG3	2.17	0.45
1:A:65:VAL:CG2	1:A:65:VAL:O	2.64	0.45
1:A:68:PRO:C	1:A:70:THR:N	2.70	0.45
1:A:6:ASN:HB2	1:A:19:LEU:HD21	1.99	0.45
1:A:43:ARG:NH2	1:A:46:GLN:HG2	2.31	0.45
1:A:50:ASN:O	1:A:54:GLN:HB2	2.17	0.45
1:A:105:CYS:O	1:A:109:VAL:HG22	2.17	0.45
1:A:66:ASP:CG	1:A:67:ASN:H	2.15	0.44
1:A:108:LYS:HA	1:A:108:LYS:HD2	1.30	0.44
1:A:55:ALA:CA	1:A:58:LEU:HG	2.48	0.44
1:A:66:ASP:OD2	1:A:67:ASN:C	2.56	0.44
1:A:23:ASN:HD22	1:A:23:ASN:HA	1.42	0.44
1:A:71:ASN:OD1	1:A:72:ASN:N	2.50	0.44
1:A:24:ASN:HB3	1:A:117:ASN:HD22	1.82	0.43
1:A:118:LEU:HD23	1:A:119:ASP:H	1.83	0.43
1:A:68:PRO:HA	1:A:95:ILE:HG13	2.00	0.43
1:A:88:ASN:HD22	1:A:88:ASN:HA	1.57	0.42
1:A:119:ASP:O	1:A:122:ASN:N	2.38	0.42
1:A:3:TRP:CD2	1:A:4:GLN:N	2.87	0.42
1:A:4:GLN:HB3	1:A:75:TYR:HE2	1.84	0.42
1:A:31:LEU:C	1:A:32:GLY:O	2.58	0.42
1:A:43:ARG:HA	1:A:46:GLN:HB3	2.01	0.42
1:A:52:TYR:HE1	1:A:98:CYS:HB3	1.83	0.42
1:A:73:TYR:HB3	1:A:92:GLU:O	2.19	0.42
1:A:48:HIS:HB2	1:A:102:ALA:HB2	2.01	0.42
1:A:54:GLN:OE1	1:A:54:GLN:CA	2.57	0.42
1:A:12:LYS:HA	1:A:12:LYS:HD2	1.66	0.42
1:A:46:GLN:CG	1:A:46:GLN:O	2.64	0.42
1:A:67:ASN:CG	1:A:70:THR:HG23	2.40	0.42
1:A:116:LYS:HZ2	1:A:117:ASN:ND2	2.17	0.42
1:A:10:LYS:CD	1:A:18:PRO:HD2	2.49	0.41
1:A:67:ASN:O	1:A:70:THR:C	2.58	0.41
1:A:65:VAL:O	1:A:65:VAL:HG22	2.20	0.41
1:A:66:ASP:OD2	1:A:68:PRO:N	2.54	0.41
1:A:85:SER:OG	1:A:87:GLU:HB2	2.20	0.41
1:A:43:ARG:NH1	1:A:46:GLN:C	2.72	0.41
1:A:82:ILE:HG22	1:A:83:THR:N	2.35	0.41
1:A:52:TYR:CE1	1:A:98:CYS:HB3	2.56	0.40
1:A:58:LEU:HB2	1:A:61:CYS:SG	2.61	0.40
1:A:25:TYR:C	1:A:25:TYR:CD1	2.95	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:57:LYS:O	1:A:62:LYS:HE2	2.21	0.40

All (11) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:63:VAL:CB	1:A:114:GLU:CG[2_655]	1.35	0.85
1:A:64:LEU:CD2	1:A:115:HIS:CE1[2_655]	1.42	0.78
1:A:63:VAL:CG2	1:A:115:HIS:CD2[2_655]	1.46	0.74
1:A:63:VAL:CG2	1:A:115:HIS:NE2[2_655]	1.57	0.63
1:A:63:VAL:CG1	1:A:114:GLU:CB[2_655]	1.72	0.48
1:A:63:VAL:CG1	1:A:114:GLU:CG[2_655]	1.74	0.46
1:A:63:VAL:CA	1:A:114:GLU:CG[2_655]	1.98	0.22
1:A:63:VAL:CG2	1:A:115:HIS:CG[2_655]	1.99	0.21
1:A:64:LEU:CD2	1:A:115:HIS:NE2[2_655]	2.06	0.14
1:A:63:VAL:CG2	1:A:115:HIS:CE1[2_655]	2.12	0.08
1:A:75:TYR:OH	1:A:87:GLU:OE2[5_556]	2.12	0.08

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	121/130 (93%)	100 (83%)	10 (8%)	11 (9%)	1 3

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	120	LYS
1	A	32	GLY
1	A	54	GLN
1	A	62	LYS
1	A	63	VAL

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Mol	Chain	Res	Type
1	A	91	CYS
1	A	112	ASN
1	A	90	ALA
1	A	55	ALA
1	A	65	VAL
1	A	68	PRO

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	111/116 (96%)	81 (73%)	30 (27%)	0 2

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	GLN
1	A	6	ASN
1	A	8	MET
1	A	16	SER
1	A	19	LEU
1	A	23	ASN
1	A	24	ASN
1	A	31	LEU
1	A	34	SER
1	A	41	LEU
1	A	43	ARG
1	A	53	LYS
1	A	56	LYS
1	A	57	LYS
1	A	60	SER
1	A	62	LYS
1	A	64	LEU
1	A	65	VAL
1	A	72	ASN
1	A	74	SER

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Mol	Chain	Res	Type
1	A	81	GLU
1	A	86	SER
1	A	95	ILE
1	A	100	ARG
1	A	108	LYS
1	A	109	VAL
1	A	110	PRO
1	A	118	LEU
1	A	119	ASP
1	A	121	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	6	ASN
1	A	23	ASN
1	A	72	ASN
1	A	88	ASN
1	A	117	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](#)

There are no ligands in this entry.

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](#)

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	123/130 (94%)	-0.69	1 (0%) 86 65	2, 10, 28, 41	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	70	THR	3.0

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 monosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.