



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:57 AM UTC

PDB ID : 1I22 / pdb_00001i22
Title : MUTANT HUMAN LYSOZYME (A83K/Q86D/A92D)
Authors : Kuroki, R.
Deposited on : 2001-02-05
Resolution : 1.80 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
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

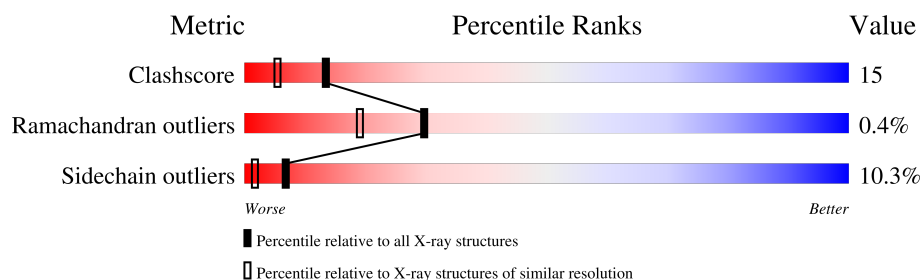
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	130	 67% 24% 7% •
1	B	130	 64% 27% 8% •
1	C	130	 66% 24% 10%
1	D	130	 62% 28% 10%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4461 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 LYSOZYME C.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	130	Total	C	N	O	S	0	0	0
			1035	636	200	189	10			
1	B	130	Total	C	N	O	S	0	0	0
			1035	636	200	189	10			
1	C	130	Total	C	N	O	S	0	0	0
			1035	636	200	189	10			
1	D	130	Total	C	N	O	S	0	0	0
			1035	636	200	189	10			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	83	LYS	ALA	engineered mutation	UNP P61626
A	86	ASP	GLN	engineered mutation	UNP P61626
A	92	ASP	ALA	engineered mutation	UNP P61626
B	83	LYS	ALA	engineered mutation	UNP P61626
B	86	ASP	GLN	engineered mutation	UNP P61626
B	92	ASP	ALA	engineered mutation	UNP P61626
C	83	LYS	ALA	engineered mutation	UNP P61626
C	86	ASP	GLN	engineered mutation	UNP P61626
C	92	ASP	ALA	engineered mutation	UNP P61626
D	83	LYS	ALA	engineered mutation	UNP P61626
D	86	ASP	GLN	engineered mutation	UNP P61626
D	92	ASP	ALA	engineered mutation	UNP P61626

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Ca	0	0
			1	1		
2	B	1	Total	Ca	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	C	1	Total 1	Ca 1	0	0
2	D	1	Total 1	Ca 1	0	0

- Molecule 3 is water.

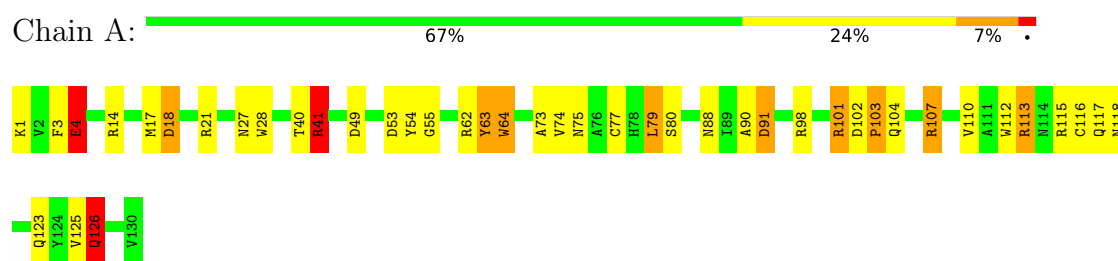
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	87	Total 87	O 87	0	0
3	B	83	Total 83	O 83	0	0
3	C	61	Total 61	O 61	0	0
3	D	86	Total 86	O 86	0	0

3 Residue-property plots [i](#)

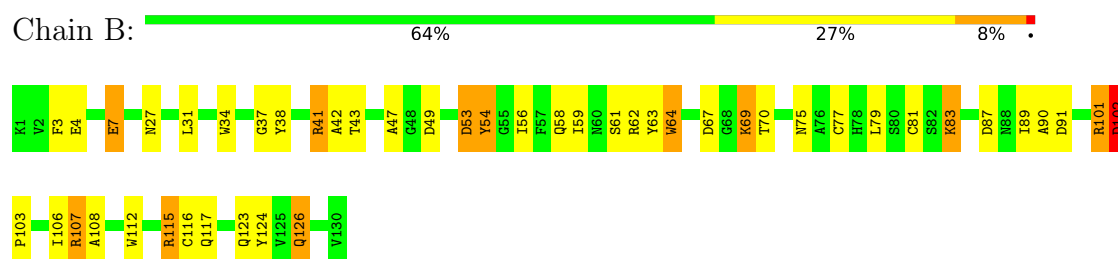
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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.

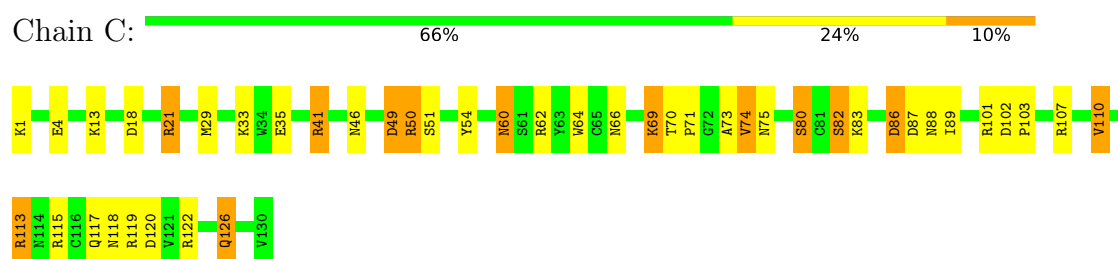
• Molecule 1: LYSOZYME C



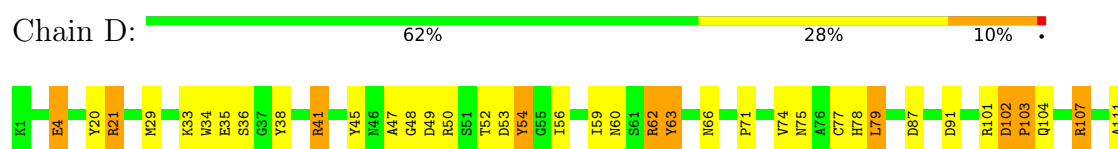
• Molecule 1: LYSOZYME C



• Molecule 1: LYSOZYME C



• Molecule 1: LYSOZYME C



W112	R113	N114	R115	C116	Q117	N118	R119	D120	V121	R122	Q123	Y124	V125	Q126	G127	C128	G129	V130
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4 Data and refinement statistics

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

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	62.00Å 113.00Å 41.20Å 90.00° 74.60° 90.00°	Depositor
Resolution (Å)	20.00 – 1.80	Depositor
% Data completeness (in resolution range)	77.0 (20.00-1.80)	Depositor
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	TNT	Depositor
R, R_{free}	0.172 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4461	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CA

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	1.38	0/1055	1.96	25/1423 (1.8%)
1	B	1.33	3/1055 (0.3%)	1.99	28/1423 (2.0%)
1	C	1.34	5/1055 (0.5%)	1.95	21/1423 (1.5%)
1	D	1.34	1/1055 (0.1%)	2.02	34/1423 (2.4%)
All	All	1.35	9/4220 (0.2%)	1.98	108/5692 (1.9%)

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	69	LYS	CA-C	-6.53	1.45	1.53
1	B	53	ASP	CA-C	-5.98	1.45	1.52
1	D	4	GLU	CD-OE2	5.87	1.36	1.25
1	C	74	VAL	CA-CB	-5.43	1.48	1.54
1	C	69	LYS	N-CA	-5.41	1.41	1.46
1	B	115	ARG	CA-C	-5.35	1.46	1.52
1	C	29	MET	CA-CB	5.25	1.61	1.53
1	C	49	ASP	CA-C	-5.19	1.47	1.52
1	B	4	GLU	CD-OE2	5.16	1.35	1.25

All (108) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	53	ASP	CA-CB-CG	-11.24	101.36	112.60
1	B	63	TYR	N-CA-C	9.43	125.48	113.88
1	D	78	HIS	CA-CB-CG	-9.38	104.42	113.80
1	C	88	ASN	CA-CB-CG	-8.87	103.73	112.60
1	A	63	TYR	CB-CA-C	-8.76	94.84	109.03
1	D	59	ILE	CA-CB-CG1	-8.26	96.36	110.40
1	D	63	TYR	CB-CA-C	-8.26	95.14	109.02
1	B	115	ARG	CB-CA-C	-7.99	96.90	110.64

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	59	ILE	N-CA-CB	-7.87	101.95	110.53
1	A	53	ASP	CA-CB-CG	-7.77	104.83	112.60
1	B	27	ASN	CA-CB-CG	-7.60	105.00	112.60
1	D	36	SER	N-CA-CB	-7.56	100.06	110.95
1	A	79	LEU	N-CA-CB	-7.08	98.31	111.13
1	D	126	GLN	N-CA-CB	7.03	120.38	109.48
1	A	64	TRP	N-CA-C	6.99	121.97	112.68
1	D	75	ASN	N-CA-C	6.96	120.99	111.39
1	D	113	ARG	CB-CA-C	-6.89	97.89	110.63
1	A	27	ASN	CA-CB-CG	-6.83	105.77	112.60
1	B	89	ILE	N-CA-C	6.78	118.67	112.29
1	C	64	TRP	N-CA-C	6.63	121.50	112.68
1	A	54	TYR	N-CA-C	6.63	120.20	109.40
1	D	107	ARG	CD-NE-CZ	-6.62	115.12	124.40
1	A	90	ALA	CA-C-N	-6.61	109.44	120.68
1	A	90	ALA	C-N-CA	-6.61	109.44	120.68
1	D	118	ASN	CA-CB-CG	6.59	119.19	112.60
1	C	82	SER	CB-CA-C	-6.55	98.51	110.63
1	A	126	GLN	CA-CB-CG	-6.45	101.19	114.10
1	A	103	PRO	N-CA-C	6.43	122.33	113.65
1	B	4	GLU	N-CA-C	-6.41	100.13	110.32
1	A	101	ARG	CA-CB-CG	-6.31	101.48	114.10
1	A	113	ARG	CB-CA-C	-6.29	98.59	110.67
1	D	128	CYS	N-CA-C	-6.24	105.71	113.38
1	D	104	GLN	CB-CA-C	-6.21	97.59	109.95
1	C	119	ARG	CB-CA-C	-6.15	99.41	110.36
1	D	102	ASP	CA-C-N	6.15	126.33	119.32
1	D	102	ASP	C-N-CA	6.15	126.33	119.32
1	A	73	ALA	CA-C-O	-6.13	114.82	120.95
1	D	4	GLU	N-CA-CB	6.13	119.60	110.29
1	D	4	GLU	CB-CG-CD	6.12	123.00	112.60
1	C	115	ARG	CB-CA-C	-6.10	99.37	109.99
1	D	104	GLN	CB-CG-CD	-6.07	102.28	112.60
1	B	90	ALA	CA-C-N	-6.04	110.57	121.14
1	B	90	ALA	C-N-CA	-6.04	110.57	121.14
1	C	113	ARG	N-CA-CB	-6.04	101.24	110.12
1	C	60	ASN	N-CA-C	6.03	119.01	109.96
1	B	107	ARG	CD-NE-CZ	-6.03	115.96	124.40
1	D	87	ASP	N-CA-C	6.03	120.64	113.16
1	C	87	ASP	CA-C-N	-6.00	114.78	122.77
1	C	87	ASP	C-N-CA	-6.00	114.78	122.77
1	D	71	PRO	CB-CA-C	-6.00	102.06	110.63

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	4	GLU	N-CA-C	-5.97	101.64	110.48
1	C	18	ASP	CA-CB-CG	-5.96	106.64	112.60
1	B	102	ASP	CA-C-N	5.94	126.09	119.32
1	B	102	ASP	C-N-CA	5.94	126.09	119.32
1	C	89	ILE	N-CA-C	5.94	117.20	111.91
1	B	64	TRP	N-CA-C	5.90	120.53	112.68
1	B	115	ARG	CA-C-N	-5.87	110.33	121.54
1	B	115	ARG	C-N-CA	-5.87	110.33	121.54
1	D	54	TYR	N-CA-C	5.82	118.89	109.40
1	C	82	SER	CA-CB-OG	-5.76	99.58	111.10
1	C	73	ALA	CA-C-O	-5.76	114.71	121.16
1	A	18	ASP	CA-CB-CG	-5.70	106.90	112.60
1	A	41	ARG	CD-NE-CZ	5.69	132.36	124.40
1	A	55	GLY	O-C-N	5.69	128.31	122.91
1	D	79	LEU	N-CA-CB	-5.65	101.01	111.13
1	B	126	GLN	N-CA-CB	5.65	118.36	109.83
1	C	80	SER	CB-CA-C	5.59	118.97	109.80
1	D	77	CYS	N-CA-CB	5.59	119.81	110.41
1	A	91	ASP	N-CA-CB	5.55	118.85	110.30
1	D	113	ARG	CB-CG-CD	-5.53	98.57	111.30
1	B	75	ASN	N-CA-C	5.52	119.25	111.52
1	A	104	GLN	CB-CG-CD	-5.51	103.23	112.60
1	B	56	ILE	CB-CA-C	5.50	120.08	112.05
1	B	42	ALA	CB-CA-C	-5.49	100.60	109.72
1	B	90	ALA	O-C-N	-5.48	116.31	122.12
1	D	66	ASN	CA-CB-CG	-5.48	107.12	112.60
1	D	53	ASP	CA-CB-CG	-5.47	107.13	112.60
1	A	98	ARG	NE-CZ-NH2	5.47	124.12	119.20
1	D	60	ASN	N-CA-C	5.42	118.09	109.96
1	B	101	ARG	N-CA-C	-5.40	106.53	113.23
1	C	113	ARG	CA-C-O	5.32	126.19	120.55
1	A	3	PHE	CB-CA-C	-5.31	99.20	109.76
1	B	70	THR	CB-CA-C	-5.28	104.07	110.15
1	B	34	TRP	CA-CB-CG	-5.28	103.57	113.60
1	D	112	TRP	CA-C-O	5.25	126.17	120.55
1	C	46	ASN	CA-CB-CG	-5.25	107.35	112.60
1	C	54	TYR	N-CA-C	5.24	118.00	109.24
1	B	54	TYR	N-CA-C	5.24	117.93	109.40
1	D	103	PRO	N-CA-C	5.21	121.52	113.75
1	B	63	TYR	O-C-N	5.21	129.34	122.36
1	C	87	ASP	CB-CA-C	5.20	119.69	109.72
1	D	124	TYR	CA-C-N	-5.18	114.97	122.69

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	124	TYR	C-N-CA	-5.18	114.97	122.69
1	D	56	ILE	N-CA-C	5.18	116.53	110.62
1	B	81	CYS	CA-C-O	5.17	125.94	120.10
1	A	75	ASN	N-CA-C	5.16	121.78	110.80
1	D	4	GLU	N-CA-C	-5.12	102.90	110.48
1	A	3	PHE	CA-C-N	-5.12	113.32	122.07
1	A	3	PHE	C-N-CA	-5.12	113.32	122.07
1	A	115	ARG	CB-CA-C	-5.10	100.82	109.38
1	C	1	LYS	CB-CA-C	-5.09	100.42	110.10
1	B	42	ALA	N-CA-C	5.08	117.60	110.23
1	B	101	ARG	N-CA-CB	5.06	118.33	110.44
1	C	74	VAL	N-CA-CB	5.05	119.32	112.35
1	B	69	LYS	CB-CA-C	-5.03	102.22	110.72
1	D	38	TYR	CA-C-N	-5.01	116.10	122.77
1	D	38	TYR	C-N-CA	-5.01	116.10	122.77
1	A	4	GLU	CB-CG-CD	5.01	121.11	112.60

There are no chirality outliers.

There are no planarity outliers.

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	1035	0	995	28	0
1	B	1035	0	995	34	0
1	C	1035	0	995	26	0
1	D	1035	0	995	41	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	87	0	0	4	0
3	B	83	0	0	0	0
3	C	61	0	0	0	0
3	D	86	0	0	4	0
All	All	4461	0	3980	119	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 15.

All (119) 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:C:49:ASP:HB2	1:C:62:ARG:HH12	1.26	0.96
1:B:115:ARG:NH2	1:D:45:TYR:HB3	1.85	0.91
1:C:50:ARG:HG3	1:C:50:ARG:HH11	1.35	0.88
1:B:83:LYS:HA	1:B:83:LYS:HE2	1.60	0.83
1:B:115:ARG:HH22	1:D:45:TYR:HB3	1.45	0.82
1:C:50:ARG:HG3	1:C:50:ARG:NH1	1.95	0.77
1:D:126:GLN:HA	1:D:126:GLN:OE1	1.84	0.77
1:B:115:ARG:CZ	1:D:45:TYR:HB3	2.15	0.77
1:D:49:ASP:HB2	1:D:62:ARG:HH12	1.50	0.76
1:C:49:ASP:HB2	1:C:62:ARG:NH1	2.01	0.76
1:C:21:ARG:HE	1:C:101:ARG:HG2	1.50	0.76
1:A:49:ASP:HB2	1:A:62:ARG:HH12	1.50	0.75
1:A:41:ARG:HH21	1:A:41:ARG:HG2	1.50	0.75
1:A:49:ASP:HB2	1:A:62:ARG:NH1	2.03	0.73
1:D:63:TYR:HA	1:D:74:VAL:CG2	2.23	0.68
1:C:41:ARG:HH21	1:C:41:ARG:HG2	1.59	0.68
1:C:86:ASP:OD2	1:C:86:ASP:N	2.27	0.67
1:C:41:ARG:HG2	1:C:41:ARG:NH2	2.09	0.66
1:D:49:ASP:CB	1:D:62:ARG:HH12	2.08	0.65
1:B:107:ARG:NH2	1:B:117:GLN:OE1	2.29	0.65
1:B:115:ARG:NH1	1:D:45:TYR:HB3	2.12	0.64
1:D:123:GLN:NE2	3:D:448:HOH:O	2.26	0.64
1:B:62:ARG:NH1	1:B:62:ARG:HB2	2.14	0.62
1:B:102:ASP:CB	1:B:103:PRO:HD2	2.30	0.62
1:A:125:VAL:HG23	1:A:126:GLN:HE22	1.65	0.60
1:A:21:ARG:HH22	1:D:21:ARG:HH11	1.48	0.60
1:A:4:GLU:HB2	3:A:418:HOH:O	2.02	0.60
1:D:101:ARG:HD2	3:D:332:HOH:O	2.02	0.60
1:D:41:ARG:HG3	1:D:41:ARG:HH21	1.66	0.59
1:C:102:ASP:HB3	1:C:103:PRO:HD2	1.85	0.59
1:A:107:ARG:CG	1:A:107:ARG:HH21	2.17	0.58
1:A:112:TRP:CD1	1:A:116:CYS:HB2	2.40	0.57
1:C:49:ASP:CB	1:C:62:ARG:HH12	2.11	0.56
1:D:107:ARG:NH2	1:D:117:GLN:OE1	2.40	0.54
1:A:21:ARG:HE	1:A:101:ARG:HG2	1.72	0.53
1:A:123:GLN:HG2	3:A:469:HOH:O	2.07	0.53
1:D:41:ARG:HH21	1:D:41:ARG:CG	2.22	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:123:GLN:HG2	1:D:124:TYR:N	2.24	0.52
1:D:63:TYR:HA	1:D:74:VAL:HG23	1.91	0.51
1:D:34:TRP:CZ3	1:D:115:ARG:NH1	2.79	0.51
1:D:49:ASP:OD2	1:D:62:ARG:NH1	2.44	0.51
1:B:3:PHE:HD1	1:B:7:GLU:HG2	1.76	0.50
1:C:83:LYS:O	1:C:86:ASP:OD2	2.29	0.50
1:A:126:GLN:OE1	1:A:126:GLN:HA	1.98	0.50
1:A:107:ARG:HH21	1:A:107:ARG:HG3	1.76	0.50
1:C:126:GLN:HA	1:C:126:GLN:OE1	2.06	0.49
1:D:45:TYR:CE1	1:D:47:ALA:HA	2.47	0.49
1:A:77:CYS:HB3	3:A:502:HOH:O	2.12	0.49
1:B:112:TRP:CD1	1:B:116:CYS:HB2	2.47	0.49
1:D:49:ASP:O	1:D:50:ARG:HB2	2.11	0.49
1:D:107:ARG:NE	3:D:391:HOH:O	2.31	0.49
1:C:110:VAL:O	1:C:113:ARG:HB2	2.13	0.48
1:C:62:ARG:NH2	1:C:71:PRO:HD2	2.28	0.48
1:B:115:ARG:HH12	1:D:45:TYR:HB3	1.78	0.48
1:A:18:ASP:HB3	3:A:515:HOH:O	2.14	0.48
1:C:51:SER:HA	1:C:70:THR:HG23	1.95	0.48
1:D:20:TYR:CE2	1:D:21:ARG:HG2	2.48	0.48
1:A:21:ARG:HH12	1:D:21:ARG:NH1	2.11	0.47
1:A:1:LYS:N	1:A:40:THR:OG1	2.41	0.47
1:A:117:GLN:O	1:A:118:ASN:HB2	2.13	0.47
1:B:49:ASP:CG	1:B:62:ARG:HH12	2.22	0.47
1:A:102:ASP:HB3	1:A:103:PRO:CD	2.44	0.47
1:B:49:ASP:OD2	1:B:62:ARG:NH1	2.48	0.47
1:D:122:ARG:O	1:D:126:GLN:NE2	2.48	0.46
1:D:41:ARG:CG	1:D:41:ARG:NH2	2.79	0.46
1:B:54:TYR:OH	1:B:67:ASP:OD2	2.28	0.46
1:A:49:ASP:HB2	1:A:62:ARG:CZ	2.46	0.46
1:D:52:THR:HB	1:D:54:TYR:CE2	2.51	0.46
1:C:51:SER:CB	1:C:60:ASN:HD21	2.29	0.45
1:C:107:ARG:O	1:C:113:ARG:NH2	2.49	0.45
1:B:102:ASP:CG	1:B:103:PRO:HD2	2.42	0.45
1:A:49:ASP:CB	1:A:62:ARG:HH12	2.25	0.45
1:A:107:ARG:CG	1:A:107:ARG:NH2	2.79	0.45
1:B:41:ARG:CG	1:B:41:ARG:NH2	2.79	0.45
1:D:50:ARG:HH11	1:D:50:ARG:HG2	1.80	0.45
1:D:50:ARG:HG2	1:D:50:ARG:NH1	2.32	0.45
1:C:51:SER:HA	1:C:70:THR:CG2	2.47	0.45
1:C:41:ARG:HH21	1:C:41:ARG:CG	2.23	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:59:ILE:CG2	1:B:64:TRP:HB2	2.47	0.44
1:A:125:VAL:HG23	1:A:126:GLN:NE2	2.29	0.44
1:B:41:ARG:CG	1:B:41:ARG:HH21	2.31	0.44
1:C:120:ASP:OD2	1:C:122:ARG:NH1	2.51	0.44
1:C:51:SER:HB2	1:C:60:ASN:HD21	1.82	0.44
1:C:117:GLN:O	1:C:118:ASN:HB2	2.18	0.44
1:B:91:ASP:N	1:B:91:ASP:OD2	2.45	0.44
1:B:53:ASP:HB3	1:B:58:GLN:HB3	2.00	0.44
1:D:35:GLU:OE2	1:D:111:ALA:HB3	2.18	0.44
1:A:17:MET:HE3	1:A:28:TRP:CD2	2.53	0.43
1:D:120:ASP:OD2	1:D:122:ARG:NH1	2.42	0.43
1:B:102:ASP:HB3	1:B:103:PRO:HD2	1.99	0.43
1:D:49:ASP:HB2	1:D:62:ARG:NH1	2.26	0.43
1:A:21:ARG:NH2	1:D:21:ARG:HH11	2.15	0.43
1:B:115:ARG:HH12	1:D:45:TYR:CB	2.32	0.43
1:A:107:ARG:NH2	1:A:117:GLN:OE1	2.51	0.43
1:B:64:TRP:O	1:B:77:CYS:HB2	2.19	0.43
1:C:102:ASP:HB3	1:C:103:PRO:CD	2.48	0.43
1:B:54:TYR:HE2	1:B:61:SER:HB3	1.84	0.43
1:B:123:GLN:HG3	1:B:124:TYR:N	2.33	0.43
1:C:62:ARG:O	1:C:62:ARG:HG3	2.18	0.43
1:B:3:PHE:CD1	1:B:7:GLU:HG2	2.54	0.43
1:A:63:TYR:HB2	1:A:64:TRP:CD1	2.53	0.42
1:A:102:ASP:HB3	1:A:103:PRO:HD2	2.00	0.42
1:B:41:ARG:HH21	1:B:41:ARG:HG3	1.84	0.42
1:B:62:ARG:HB2	1:B:62:ARG:HH11	1.82	0.42
1:D:102:ASP:HB3	1:D:103:PRO:CD	2.49	0.42
1:D:107:ARG:NH1	3:D:391:HOH:O	2.40	0.42
1:B:37:GLY:O	1:B:38:TYR:HB2	2.20	0.41
1:A:113:ARG:HH11	1:A:113:ARG:HD2	1.73	0.41
1:B:115:ARG:NH1	1:D:45:TYR:CB	2.80	0.41
1:C:35:GLU:OE1	1:C:35:GLU:HA	2.18	0.41
1:B:49:ASP:CG	1:B:62:ARG:NH1	2.79	0.41
1:B:83:LYS:HE2	1:B:83:LYS:CA	2.41	0.41
1:B:106:ILE:C	1:B:108:ALA:H	2.28	0.41
1:D:21:ARG:HD2	1:D:21:ARG:HA	1.67	0.41
1:C:66:ASN:HB2	1:C:75:ASN:ND2	2.36	0.40
1:D:29:MET:HE3	1:D:29:MET:HB3	1.93	0.40
1:B:102:ASP:CB	1:B:103:PRO:CD	2.98	0.40
1:D:62:ARG:O	1:D:62:ARG:HG3	2.10	0.40
1:D:113:ARG:HH11	1:D:113:ARG:HD2	1.65	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	128/130 (98%)	124 (97%)	4 (3%)	0	100	100
1	B	128/130 (98%)	120 (94%)	7 (6%)	1 (1%)	16	6
1	C	128/130 (98%)	123 (96%)	5 (4%)	0	100	100
1	D	128/130 (98%)	124 (97%)	3 (2%)	1 (1%)	16	6
All	All	512/520 (98%)	491 (96%)	19 (4%)	2 (0%)	30	19

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	48	GLY
1	B	47	ALA

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	107/107 (100%)	96 (90%)	11 (10%)	7	2
1	B	107/107 (100%)	96 (90%)	11 (10%)	7	2
1	C	107/107 (100%)	95 (89%)	12 (11%)	6	1
1	D	107/107 (100%)	97 (91%)	10 (9%)	8	2
All	All	428/428 (100%)	384 (90%)	44 (10%)	7	2

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	GLU
1	A	14	ARG
1	A	41	ARG
1	A	74	VAL
1	A	79	LEU
1	A	80	SER
1	A	88	ASN
1	A	91	ASP
1	A	107	ARG
1	A	110	VAL
1	A	126	GLN
1	B	7	GLU
1	B	31	LEU
1	B	41	ARG
1	B	43	THR
1	B	69	LYS
1	B	79	LEU
1	B	83	LYS
1	B	87	ASP
1	B	101	ARG
1	B	102	ASP
1	B	126	GLN
1	C	13	LYS
1	C	21	ARG
1	C	33	LYS
1	C	41	ARG
1	C	50	ARG
1	C	69	LYS
1	C	74	VAL
1	C	80	SER
1	C	82	SER
1	C	86	ASP
1	C	110	VAL
1	C	126	GLN
1	D	4	GLU
1	D	21	ARG
1	D	33	LYS
1	D	41	ARG
1	D	62	ARG
1	D	79	LEU
1	D	91	ASP
1	D	123	GLN

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Mol	Chain	Res	Type
1	D	126	GLN
1	D	130	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	58	GLN
1	A	78	HIS
1	A	104	GLN
1	A	118	ASN
1	B	88	ASN
1	B	104	GLN
1	B	118	ASN
1	C	88	ASN
1	C	118	ASN
1	D	44	ASN
1	D	118	ASN
1	D	126	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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

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

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 is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section is therefore empty.

6.4 Ligands ⓘ

EDS was not executed - this section is therefore empty.

6.5 Other polymers ⓘ

EDS was not executed - this section is therefore empty.