



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:57 PM GMT

PDB ID : 1I22
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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

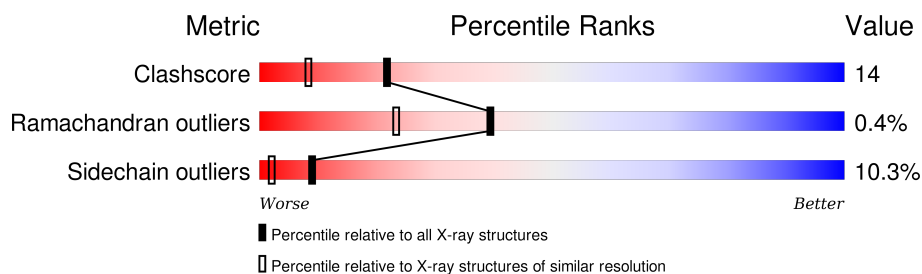
1 Overall quality at a glance

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	102246	5383 (1.80-1.80)
Ramachandran outliers	100387	5320 (1.80-1.80)
Sidechain outliers	100360	5319 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	130	
1	B	130	
1	C	130	
1	D	130	

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	UNP P61626
A	86	ASP	GLN	ENGINEERED	UNP P61626
A	92	ASP	ALA	ENGINEERED	UNP P61626
B	83	LYS	ALA	ENGINEERED	UNP P61626
B	86	ASP	GLN	ENGINEERED	UNP P61626
B	92	ASP	ALA	ENGINEERED	UNP P61626
C	83	LYS	ALA	ENGINEERED	UNP P61626
C	86	ASP	GLN	ENGINEERED	UNP P61626
C	92	ASP	ALA	ENGINEERED	UNP P61626
D	83	LYS	ALA	ENGINEERED	UNP P61626
D	86	ASP	GLN	ENGINEERED	UNP P61626
D	92	ASP	ALA	ENGINEERED	UNP P61626

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

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

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

- Molecule 3 is water.

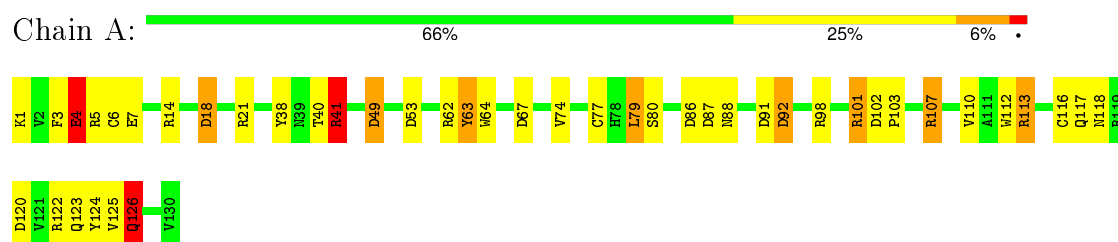
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	87	Total	O	0	0
			87	87		
3	B	83	Total	O	0	0
			83	83		
3	C	61	Total	O	0	0
			61	61		
3	D	86	Total	O	0	0
			86	86		

3 Residue-property plots [i](#)

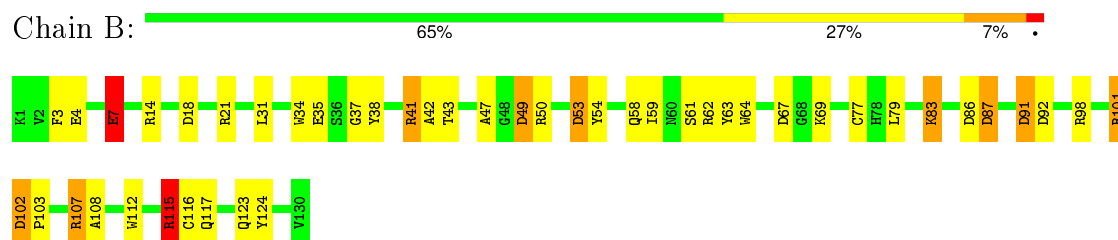
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according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

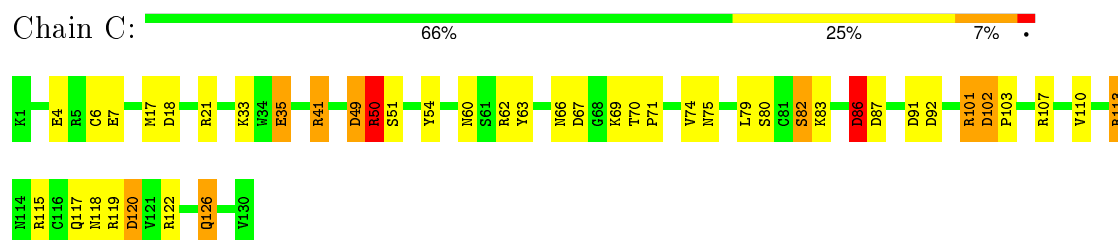
• Molecule 1: LYSOZYME C



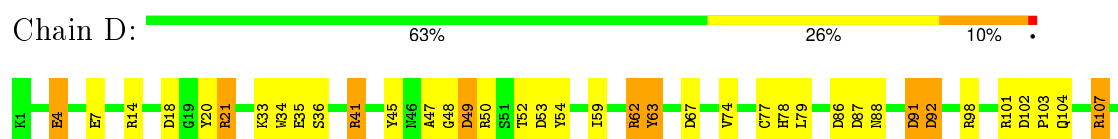
• Molecule 1: LYSOZYME C

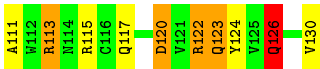


• Molecule 1: LYSOZYME C



• Molecule 1: LYSOZYME C





4 Data and refinement statistics

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

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	62.00 Å 113.00 Å 41.20 Å 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	1.07	2/1055 (0.2%)	1.78	30/1423 (2.1%)
1	B	1.04	3/1055 (0.3%)	1.73	28/1423 (2.0%)
1	C	1.04	3/1055 (0.3%)	1.77	31/1423 (2.2%)
1	D	1.08	2/1055 (0.2%)	1.86	36/1423 (2.5%)
All	All	1.06	10/4220 (0.2%)	1.78	125/5692 (2.2%)

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	4	GLU	CD-OE2	9.87	1.36	1.25
1	B	4	GLU	CD-OE2	8.64	1.35	1.25
1	A	4	GLU	CD-OE2	8.27	1.34	1.25
1	C	4	GLU	CD-OE2	7.83	1.34	1.25
1	B	7	GLU	CD-OE2	6.21	1.32	1.25
1	D	7	GLU	CD-OE2	6.07	1.32	1.25
1	A	7	GLU	CD-OE2	5.93	1.32	1.25
1	C	7	GLU	CD-OE2	5.90	1.32	1.25
1	B	35	GLU	CD-OE2	5.05	1.31	1.25
1	C	35	GLU	CD-OE1	-5.02	1.20	1.25

All (125) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	67	ASP	CB-CG-OD2	-12.09	107.42	118.30
1	A	67	ASP	CB-CG-OD1	11.84	128.96	118.30
1	D	14	ARG	NE-CZ-NH1	11.42	126.01	120.30
1	D	67	ASP	CB-CG-OD1	11.37	128.53	118.30
1	D	92	ASP	CB-CG-OD2	-10.57	108.79	118.30
1	D	87	ASP	CB-CG-OD2	-10.29	109.03	118.30
1	D	18	ASP	CB-CG-OD1	9.79	127.11	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	87	ASP	CB-CG-OD2	-9.70	109.57	118.30
1	C	50	ARG	NE-CZ-NH1	9.56	125.08	120.30
1	B	67	ASP	CB-CG-OD2	-9.53	109.72	118.30
1	D	14	ARG	NE-CZ-NH2	-9.45	115.58	120.30
1	A	53	ASP	CB-CG-OD2	-9.18	110.04	118.30
1	B	87	ASP	CB-CG-OD2	-8.98	110.21	118.30
1	D	91	ASP	CB-CG-OD2	-8.90	110.29	118.30
1	D	86	ASP	CB-CG-OD2	-8.85	110.34	118.30
1	B	107	ARG	NE-CZ-NH1	8.83	124.71	120.30
1	C	120	ASP	CB-CG-OD1	8.63	126.07	118.30
1	B	98	ARG	NE-CZ-NH2	-8.49	116.06	120.30
1	B	107	ARG	NE-CZ-NH2	-8.38	116.11	120.30
1	C	18	ASP	CB-CG-OD1	8.38	125.84	118.30
1	D	120	ASP	CB-CG-OD1	8.33	125.80	118.30
1	C	87	ASP	CB-CG-OD1	8.25	125.73	118.30
1	A	124	TYR	CB-CG-CD1	-8.12	116.13	121.00
1	A	87	ASP	CB-CG-OD2	-7.99	111.11	118.30
1	D	86	ASP	CB-CG-OD1	7.83	125.34	118.30
1	C	91	ASP	CB-CG-OD2	-7.80	111.28	118.30
1	A	63	TYR	CB-CA-C	-7.78	94.84	110.40
1	C	54	TYR	CB-CG-CD1	-7.77	116.34	121.00
1	C	113	ARG	NE-CZ-NH2	-7.76	116.42	120.30
1	D	59	ILE	CA-CB-CG1	-7.71	96.36	111.00
1	B	98	ARG	NE-CZ-NH1	7.66	124.13	120.30
1	A	98	ARG	NE-CZ-NH2	7.64	124.12	120.30
1	D	63	TYR	CB-CA-C	-7.63	95.14	110.40
1	D	18	ASP	CB-CG-OD2	-7.58	111.48	118.30
1	C	92	ASP	CB-CG-OD1	7.38	124.94	118.30
1	A	49	ASP	CB-CG-OD1	7.25	124.83	118.30
1	C	67	ASP	CB-CG-OD1	7.20	124.78	118.30
1	C	86	ASP	CB-CG-OD2	-7.18	111.84	118.30
1	A	120	ASP	CB-CG-OD2	-7.16	111.86	118.30
1	C	18	ASP	CB-CG-OD2	-7.16	111.86	118.30
1	A	120	ASP	CB-CG-OD1	7.13	124.72	118.30
1	B	53	ASP	CB-CG-OD1	7.11	124.70	118.30
1	B	87	ASP	CB-CG-OD1	7.02	124.62	118.30
1	C	102	ASP	CB-CG-OD1	7.02	124.61	118.30
1	D	36	SER	N-CA-CB	-6.96	100.06	110.50
1	D	92	ASP	CB-CG-OD1	6.92	124.53	118.30
1	B	18	ASP	CB-CG-OD2	-6.88	112.11	118.30
1	C	101	ARG	NE-CZ-NH1	6.88	123.74	120.30
1	B	115	ARG	CB-CA-C	-6.75	96.90	110.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	67	ASP	CB-CG-OD2	-6.73	112.24	118.30
1	A	5	ARG	NE-CZ-NH1	6.65	123.63	120.30
1	D	87	ASP	CB-CG-OD1	6.65	124.29	118.30
1	A	5	ARG	NE-CZ-NH2	-6.62	116.99	120.30
1	B	53	ASP	CB-CG-OD2	-6.50	112.44	118.30
1	D	53	ASP	CB-CG-OD1	6.42	124.08	118.30
1	D	122	ARG	NE-CZ-NH1	6.41	123.50	120.30
1	D	104	GLN	CB-CA-C	-6.40	97.59	110.40
1	C	54	TYR	CZ-CE2-CD2	-6.39	114.05	119.80
1	B	42	ALA	CB-CA-C	-6.33	100.60	110.10
1	D	113	ARG	NE-CZ-NH2	-6.27	117.17	120.30
1	A	41	ARG	CD-NE-CZ	6.26	132.36	123.60
1	D	113	ARG	CB-CA-C	-6.26	97.89	110.40
1	C	91	ASP	CB-CG-OD1	6.25	123.92	118.30
1	A	53	ASP	CB-CG-OD1	6.21	123.89	118.30
1	B	18	ASP	CB-CG-OD1	6.21	123.88	118.30
1	B	91	ASP	CB-CG-OD2	-6.19	112.72	118.30
1	A	67	ASP	CB-CG-OD2	-6.19	112.73	118.30
1	C	120	ASP	CB-CG-OD2	-6.17	112.74	118.30
1	A	86	ASP	CB-CG-OD1	6.16	123.85	118.30
1	A	87	ASP	CB-CG-OD1	6.13	123.82	118.30
1	C	82	SER	CB-CA-C	-6.10	98.51	110.10
1	B	67	ASP	CB-CG-OD1	6.10	123.79	118.30
1	D	107	ARG	NE-CZ-NH1	-6.09	117.25	120.30
1	A	18	ASP	CB-CG-OD1	6.07	123.76	118.30
1	D	107	ARG	CD-NE-CZ	-6.05	115.12	123.60
1	A	79	LEU	N-CA-CB	-6.05	98.31	110.40
1	D	91	ASP	CB-CG-OD1	5.96	123.67	118.30
1	A	113	ARG	CB-CA-C	-5.90	98.59	110.40
1	A	122	ARG	NE-CZ-NH1	5.83	123.21	120.30
1	D	63	TYR	CB-CG-CD1	-5.75	117.55	121.00
1	B	92	ASP	CB-CG-OD1	5.75	123.48	118.30
1	D	120	ASP	CB-CG-OD2	-5.74	113.14	118.30
1	C	54	TYR	CB-CG-CD2	5.68	124.41	121.00
1	B	101	ARG	NE-CZ-NH1	5.65	123.12	120.30
1	C	49	ASP	CB-CG-OD1	5.62	123.36	118.30
1	A	3	PHE	CB-CA-C	-5.60	99.20	110.40
1	D	98	ARG	NE-CZ-NH1	5.58	123.09	120.30
1	C	86	ASP	CB-CG-OD1	5.57	123.31	118.30
1	A	126	GLN	CA-CB-CG	-5.55	101.19	113.40
1	A	63	TYR	CB-CG-CD1	-5.54	117.68	121.00
1	D	98	ARG	CD-NE-CZ	5.53	131.34	123.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	14	ARG	NE-CZ-NH1	5.52	123.06	120.30
1	C	115	ARG	CB-CA-C	-5.51	99.37	110.40
1	A	92	ASP	CB-CG-OD2	-5.50	113.35	118.30
1	C	119	ARG	CB-CA-C	-5.50	99.41	110.40
1	A	122	ARG	NE-CZ-NH2	-5.47	117.56	120.30
1	B	53	ASP	CA-CB-CG	-5.47	101.36	113.40
1	B	107	ARG	CD-NE-CZ	-5.46	115.96	123.60
1	A	38	TYR	N-CA-CB	5.45	120.41	110.60
1	B	21	ARG	CD-NE-CZ	5.43	131.21	123.60
1	D	126	GLN	N-CA-CB	5.43	120.38	110.60
1	B	49	ASP	CB-CG-OD2	-5.42	113.42	118.30
1	D	49	ASP	CB-CG-OD2	-5.42	113.42	118.30
1	A	101	ARG	CA-CB-CG	-5.42	101.48	113.40
1	A	18	ASP	CB-CG-OD2	-5.40	113.44	118.30
1	D	78	HIS	CA-CB-CG	-5.40	104.42	113.60
1	C	79	LEU	CB-CA-C	-5.39	99.95	110.20
1	B	86	ASP	CB-CG-OD2	-5.38	113.46	118.30
1	B	63	TYR	N-CA-C	5.36	125.48	111.00
1	C	17	MET	CG-SD-CE	-5.34	91.65	100.20
1	B	34	TRP	CA-CB-CG	-5.33	103.57	113.70
1	B	108	ALA	CB-CA-C	-5.28	102.18	110.10
1	C	92	ASP	CB-CG-OD2	-5.28	113.55	118.30
1	D	14	ARG	CD-NE-CZ	5.26	130.96	123.60
1	C	113	ARG	N-CA-CB	-5.20	101.24	110.60
1	C	35	GLU	OE1-CD-OE2	-5.13	117.14	123.30
1	D	77	CYS	N-CA-CB	5.11	119.81	110.60
1	C	6	CYS	CA-CB-SG	-5.09	104.84	114.00
1	C	63	TYR	CB-CG-CD1	-5.07	117.96	121.00
1	A	86	ASP	CB-CG-OD2	-5.07	113.74	118.30
1	A	6	CYS	CA-CB-SG	-5.06	104.89	114.00
1	D	113	ARG	CB-CG-CD	-5.01	98.57	111.60
1	B	14	ARG	CD-NE-CZ	5.01	130.61	123.60
1	B	50	ARG	NE-CZ-NH1	5.00	122.80	120.30
1	D	4	GLU	N-CA-CB	5.00	119.60	110.60

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	27	0
1	B	1035	0	995	31	0
1	C	1035	0	995	26	0
1	D	1035	0	995	40	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	114	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 14.

All (114) 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:OE1	1:D:126:GLN:HA	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:HG2	1:A:41:ARG:HH21	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:102:ASP:CB	1:B:103:PRO:HD2	2.30	0.62
1:B:62:ARG:HB2	1:B:62:ARG:NH1	2.14	0.62
1:A:126:GLN:OE1	1:A:126:GLN:HA	1.98	0.61
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:D:101:ARG:HD2	3:D:332:HOH:O	2.02	0.60
1:A:4:GLU:HB2	3:A:418:HOH:O	2.02	0.60
1:D:41:ARG:HH21	1:D:41:ARG:HG3	1.66	0.59
1:C:102:ASP:HB3	1:C:103:PRO:HD2	1.85	0.59
1:A:107:ARG:HH21	1:A:107:ARG:CG	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:C:126:GLN:OE1	1:C:126:GLN:HA	2.06	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
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:D:49:ASP:O	1:D:50:ARG:HB2	2.11	0.51
1:B:3:PHE:HD1	1:B:7:GLU:HG2	1.76	0.50
1:B:91:ASP:N	1:B:91:ASP:OD2	2.45	0.50
1:C:83:LYS:O	1:C:86:ASP:OD2	2.29	0.50
1:A:107:ARG:HH21	1:A:107:ARG:HG3	1.76	0.50
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:D:107:ARG:NE	3:D:391:HOH:O	2.31	0.49
1:B:112:TRP:CD1	1:B:116:CYS:HB2	2.47	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:D:62:ARG:O	1:D:62:ARG:HG3	2.10	0.48
1:D:20:TYR:CE2	1:D:21:ARG:HG2	2.48	0.48
1:C:51:SER:HA	1:C:70:THR:HG23	1.95	0.48
1:A:21:ARG:HH12	1:D:21:ARG:NH1	2.11	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:117:GLN:O	1:A:118:ASN:HB2	2.13	0.47
1:A:1:LYS:N	1:A:40:THR:OG1	2.41	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: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:107:ARG:O	1:C:113:ARG:NH2	2.49	0.45
1:C:51:SER:CB	1:C:60:ASN:HD21	2.29	0.45
1:A:49:ASP:CB	1:A:62:ARG:HH12	2.25	0.45
1:A:107:ARG:NH2	1:A:107:ARG:CG	2.79	0.45
1:D:50:ARG:HG2	1:D:50:ARG:HH11	1.80	0.45
1:B:41:ARG:NH2	1:B:41:ARG:CG	2.79	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
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:HH21	1:B:41:ARG:CG	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: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:D:120:ASP:OD2	1:D:122:ARG:NH1	2.42	0.43
1:D:49:ASP:HB2	1:D:62:ARG:NH1	2.26	0.43
1:B:102:ASP:HB3	1:B:103:PRO:HD2	1.99	0.43
1:B:115:ARG:HH12	1:D:45:TYR:CB	2.32	0.43
1:A:21:ARG:NH2	1:D:21:ARG:HH11	2.15	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:C:62:ARG:O	1:C:62:ARG:HG3	2.18	0.43
1:B:49:ASP:CG	1:B:62:ARG:HH12	2.22	0.43
1:B:123:GLN:HG3	1:B:124:TYR:N	2.33	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:C:35:GLU:OE1	1:C:35:GLU:HA	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:HD2	1:A:113:ARG:HH11	1.73	0.41
1:B:115:ARG:NH1	1:D:45:TYR:CB	2.80	0.41
1:B:83:LYS:CA	1:B:83:LYS:HE2	2.41	0.41
1:D:21:ARG:HA	1:D:21:ARG:HD2	1.67	0.41
1:C:66:ASN:HB2	1:C:75:ASN:ND2	2.36	0.40
1:B:102:ASP:CB	1:B:103:PRO:CD	2.98	0.40
1:B:102:ASP:CG	1:B:103:PRO:HD2	2.42	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%)	24	8
1	C	128/130 (98%)	123 (96%)	5 (4%)	0	100	100
1	D	128/130 (98%)	124 (97%)	3 (2%)	1 (1%)	24	8
All	All	512/520 (98%)	491 (96%)	19 (4%)	2 (0%)	39	23

All (2) Ramachandran outliers are listed below:

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

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	107/107 (100%)	95 (89%)	12 (11%)	7	1
1	B	107/107 (100%)	96 (90%)	11 (10%)	9	2
1	C	107/107 (100%)	98 (92%)	9 (8%)	14	4
1	D	107/107 (100%)	95 (89%)	12 (11%)	7	1
All	All	428/428 (100%)	384 (90%)	44 (10%)	9	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	92	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	115	ARG
1	C	33	LYS

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Mol	Chain	Res	Type
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	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	88	ASN
1	D	91	ASP
1	D	92	ASP
1	D	123	GLN
1	D	126	GLN
1	D	130	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	104	GLN
1	A	118	ASN
1	B	104	GLN
1	B	118	ASN
1	C	88	ASN
1	C	118	ASN
1	D	88	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 [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

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 [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 ⓘ

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.