



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:16 PM UTC

PDB ID : 1L97 / pdb_00001l97
Title : STRUCTURE OF A HINGE-BENDING BACTERIOPHAGE T4
LYSOZYME MUTANT, ILE3-> PRO
Authors : Dixon, M.; Shewchuk, L.; Matthews, B.W.
Deposited on : 1992-02-11
Resolution : 2.00 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	:	4-5-2 with Phenix2.0
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

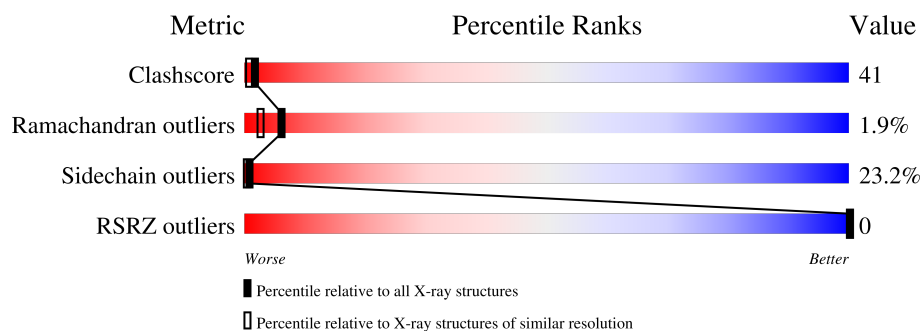
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	164	
1	B	164	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2817 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 T4 LYSOZYME.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	164	Total	C	N	O	S	0	0	0
			1308	822	238	241	7			
1	B	164	Total	C	N	O	S	0	0	0
			1308	822	238	241	7			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	3	PRO	ILE	conflict	UNP P00720
B	3	PRO	ILE	conflict	UNP P00720

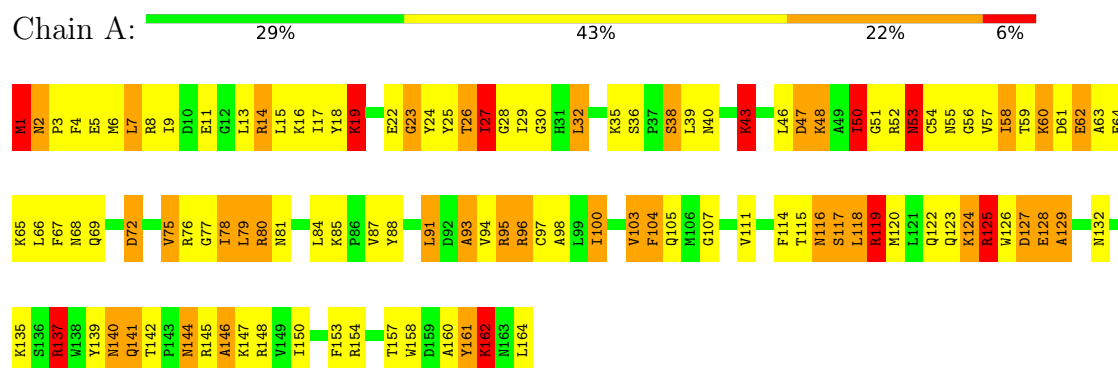
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	103	Total	O	0	0
			103	103		
2	B	98	Total	O	0	0
			98	98		

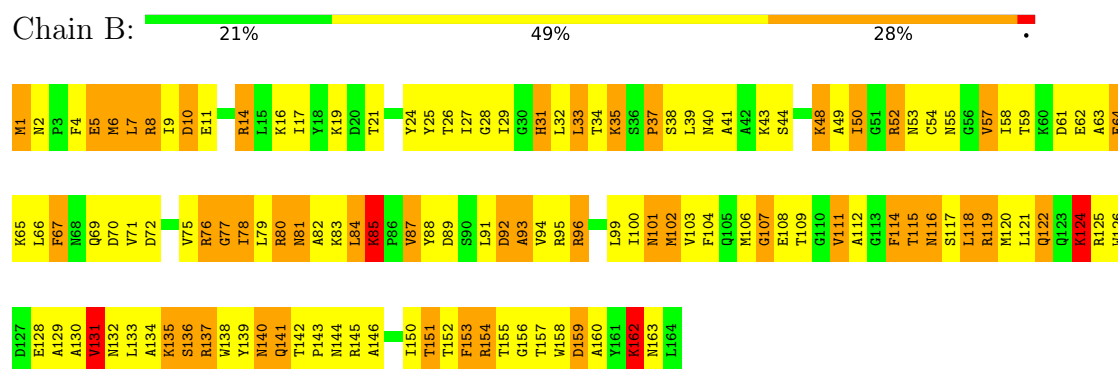
3 Residue-property plots

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: T4 LYSOZYME



• Molecule 1: T4 LYSOZYME



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	86.50Å 96.60Å 39.20Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	(Not available) – 2.00 64.44 – 2.00	Depositor EDS
% Data completeness (in resolution range)	(Not available) ((Not available)-2.00) 70.8 (64.44-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.33 (at 1.99Å)	Xtriage
Refinement program	TNT	Depositor
R, R_{free}	0.167 , (Not available) 0.170 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	17.7	Xtriage
Anisotropy	0.349	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 134.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	2817	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.18% 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 [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	1.66	17/1329 (1.3%)	2.20	54/1789 (3.0%)
1	B	1.55	6/1329 (0.5%)	2.14	59/1789 (3.3%)
All	All	1.61	23/2658 (0.9%)	2.17	113/3578 (3.2%)

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	116	ASN	CA-C	-7.97	1.42	1.52
1	A	103	VAL	C-N	-7.75	1.23	1.33
1	A	8	ARG	NE-CZ	7.47	1.41	1.33
1	A	26	THR	CA-C	-6.20	1.44	1.52
1	A	7	LEU	CA-C	-5.96	1.44	1.52
1	A	32	LEU	CA-C	-5.92	1.45	1.52
1	A	28	GLY	N-CA	5.81	1.53	1.45
1	A	129	ALA	N-CA	-5.71	1.39	1.46
1	A	63	ALA	C-N	-5.55	1.26	1.33
1	A	61	ASP	CA-C	-5.54	1.45	1.52
1	A	146	ALA	N-CA	5.48	1.53	1.46
1	B	37	PRO	CA-C	-5.46	1.44	1.52
1	A	26	THR	C-N	-5.46	1.28	1.33
1	A	27	ILE	C-N	5.42	1.41	1.33
1	B	92	ASP	CG-OD2	5.41	1.35	1.25
1	A	11	GLU	CD-OE1	5.34	1.35	1.25
1	B	54	CYS	C-O	-5.27	1.17	1.24
1	A	154	ARG	C-O	-5.21	1.18	1.24
1	B	159	ASP	CG-OD1	5.19	1.35	1.25
1	B	50	ILE	CA-CB	-5.18	1.47	1.54
1	A	64	GLU	CD-OE2	5.05	1.34	1.25
1	A	25	TYR	CA-C	-5.03	1.46	1.52
1	B	64	GLU	CA-C	-5.01	1.46	1.52

All (113) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	50	ILE	CB-CA-C	10.44	126.36	111.65
1	A	2	ASN	N-CA-C	-9.86	99.26	108.13
1	A	53	ASN	CA-CB-CG	9.74	122.34	112.60
1	B	124	LYS	CB-CA-C	8.80	123.60	111.63
1	A	2	ASN	CA-C-N	8.62	128.35	119.56
1	A	2	ASN	C-N-CA	8.62	128.35	119.56
1	A	137	ARG	NE-CZ-NH2	-7.90	112.09	119.20
1	B	67	PHE	CA-CB-CG	-7.67	106.13	113.80
1	A	140	ASN	CA-CB-CG	-7.48	105.12	112.60
1	B	31	HIS	CA-CB-CG	-7.36	106.44	113.80
1	A	68	ASN	CA-CB-CG	7.30	119.90	112.60
1	B	114	PHE	CA-CB-CG	-7.24	106.56	113.80
1	A	19	LYS	CA-C-N	-7.17	109.50	122.12
1	A	19	LYS	C-N-CA	-7.17	109.50	122.12
1	B	101	ASN	CA-C-N	6.97	129.92	120.44
1	B	101	ASN	C-N-CA	6.97	129.92	120.44
1	A	38	SER	N-CA-C	6.95	120.32	109.96
1	A	47	ASP	CA-CB-CG	-6.95	105.65	112.60
1	B	141	GLN	CA-C-N	6.94	129.50	122.00
1	B	141	GLN	C-N-CA	6.94	129.50	122.00
1	A	96	ARG	CA-C-N	6.88	129.50	120.28
1	A	96	ARG	C-N-CA	6.88	129.50	120.28
1	A	91	LEU	N-CA-C	6.70	120.41	109.76
1	A	14	ARG	NE-CZ-NH2	-6.52	113.33	119.20
1	A	142	THR	CA-C-N	6.48	125.71	118.97
1	A	142	THR	C-N-CA	6.48	125.71	118.97
1	A	162	LYS	N-CA-C	6.41	118.27	111.28
1	A	23	GLY	CA-C-N	6.40	129.96	120.87
1	A	23	GLY	C-N-CA	6.40	129.96	120.87
1	B	10	ASP	N-CA-C	6.38	118.11	111.03
1	B	115	THR	CA-C-N	6.35	129.61	120.79
1	B	115	THR	C-N-CA	6.35	129.61	120.79
1	B	91	LEU	N-CA-CB	6.34	119.40	109.83
1	A	144	ASN	CA-CB-CG	-6.30	106.30	112.60
1	B	162	LYS	N-CA-C	6.24	121.07	112.90
1	B	57	VAL	CA-C-N	-6.09	114.82	123.17
1	B	57	VAL	C-N-CA	-6.09	114.82	123.17
1	B	71	VAL	N-CA-C	6.09	116.25	110.53
1	B	49	ALA	CA-C-O	6.05	126.93	120.70
1	A	67	PHE	CA-CB-CG	-5.96	107.84	113.80
1	B	142	THR	CB-CA-C	-5.95	103.26	110.81
1	A	1	MET	CB-CA-C	5.94	121.39	110.10
1	B	52	ARG	CA-C-N	-5.94	113.98	122.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	52	ARG	C-N-CA	-5.94	113.98	122.77
1	B	33	LEU	N-CA-CB	5.91	118.45	110.59
1	A	39	LEU	N-CA-C	-5.89	104.77	111.07
1	B	131	VAL	N-CA-CB	-5.87	103.14	110.47
1	B	135	LYS	N-CA-CB	5.76	118.84	110.26
1	A	64	GLU	CA-C-O	5.72	126.48	120.42
1	B	37	PRO	N-CA-CB	5.71	109.25	103.25
1	B	159	ASP	CA-CB-CG	5.71	118.31	112.60
1	A	137	ARG	CA-C-O	5.71	126.39	119.31
1	A	72	ASP	CA-CB-CG	-5.63	106.97	112.60
1	B	94	VAL	N-CA-C	5.62	116.27	110.36
1	A	19	LYS	O-C-N	-5.61	116.67	123.29
1	A	64	GLU	N-CA-CB	-5.59	101.88	110.16
1	A	161	TYR	CB-CA-C	-5.59	100.42	109.13
1	B	144	ASN	CA-C-N	5.59	128.22	120.29
1	B	144	ASN	C-N-CA	5.59	128.22	120.29
1	A	125	ARG	CA-C-N	5.57	132.19	121.54
1	A	125	ARG	C-N-CA	5.57	132.19	121.54
1	A	93	ALA	N-CA-C	5.56	122.65	110.80
1	A	127	ASP	N-CA-CB	5.56	119.67	110.33
1	A	96	ARG	N-CA-C	-5.55	105.23	111.28
1	B	116	ASN	CB-CA-C	-5.53	102.13	110.92
1	B	140	ASN	CA-CB-CG	5.48	118.08	112.60
1	A	48	LYS	O-C-N	5.47	128.00	122.09
1	B	118	LEU	N-CA-C	-5.47	105.31	111.28
1	B	55	ASN	CA-CB-CG	-5.46	107.14	112.60
1	B	107	GLY	CA-C-N	5.46	131.96	121.54
1	B	107	GLY	C-N-CA	5.46	131.96	121.54
1	B	144	ASN	CA-CB-CG	5.46	118.06	112.60
1	A	43	LYS	O-C-N	5.42	127.72	122.09
1	A	119	ARG	CA-C-N	5.41	127.47	120.44
1	A	119	ARG	C-N-CA	5.41	127.47	120.44
1	B	155	THR	CA-CB-OG1	-5.40	101.50	109.60
1	B	81	ASN	CA-CB-CG	-5.39	107.20	112.60
1	B	91	LEU	CA-C-N	5.39	129.54	121.72
1	B	91	LEU	C-N-CA	5.39	129.54	121.72
1	A	114	PHE	N-CA-CB	-5.37	104.17	110.35
1	B	40	ASN	O-C-N	5.36	129.62	122.43
1	B	77	GLY	CA-C-N	5.35	127.39	120.70
1	B	77	GLY	C-N-CA	5.35	127.39	120.70
1	A	96	ARG	O-C-N	5.32	127.75	122.12
1	B	85	LYS	N-CA-CB	5.31	120.11	110.08

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	24	TYR	N-CA-C	5.29	118.62	110.42
1	B	122	GLN	CB-CG-CD	-5.28	103.63	112.60
1	A	62	GLU	N-CA-C	-5.27	105.70	111.82
1	B	124	LYS	CA-C-O	5.25	127.49	120.98
1	A	160	ALA	N-CA-C	5.16	119.21	113.01
1	B	61	ASP	CA-C-O	5.16	125.89	120.42
1	A	54	CYS	CB-CA-C	-5.16	102.09	110.85
1	A	95	ARG	CD-NE-CZ	-5.14	117.20	124.40
1	A	30	GLY	CA-C-N	-5.14	114.12	122.34
1	A	30	GLY	C-N-CA	-5.14	114.12	122.34
1	B	163	ASN	CA-CB-CG	5.14	117.74	112.60
1	A	118	LEU	CA-C-N	5.13	127.40	120.38
1	A	118	LEU	C-N-CA	5.13	127.40	120.38
1	B	153	PHE	CA-C-N	5.12	127.42	120.65
1	B	153	PHE	C-N-CA	5.12	127.42	120.65
1	B	78	ILE	N-CA-C	-5.10	105.35	110.30
1	B	131	VAL	N-CA-C	-5.09	104.82	110.62
1	B	37	PRO	O-C-N	5.09	129.51	122.64
1	B	53	ASN	CB-CA-C	5.06	120.14	111.68
1	B	96	ARG	CA-C-N	5.06	127.06	120.28
1	B	96	ARG	C-N-CA	5.06	127.06	120.28
1	B	116	ASN	N-CA-C	5.03	117.16	111.02
1	A	26	THR	CA-C-N	-5.02	116.49	122.11
1	A	26	THR	C-N-CA	-5.02	116.49	122.11
1	A	75	VAL	N-CA-C	-5.02	105.60	110.42
1	A	68	ASN	OD1-CG-ND2	-5.02	117.58	122.60
1	B	116	ASN	CA-C-N	5.01	126.99	120.28
1	B	116	ASN	C-N-CA	5.01	126.99	120.28

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	1308	0	1330	90	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1308	0	1330	128	0
2	A	103	0	0	13	0
2	B	98	0	0	12	0
All	All	2817	0	2660	216	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 41.

All (216) 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:B:114:PHE:HB2	1:B:118:LEU:HD11	1.38	1.04
1:B:84:LEU:HD21	1:B:112:ALA:HB2	1.40	0.99
1:A:1:MET:HE3	1:A:5:GLU:HB3	1.48	0.96
1:B:99:LEU:O	1:B:103:VAL:HG22	1.70	0.92
1:A:125:ARG:HG2	1:A:128:GLU:OE1	1.70	0.91
1:B:124:LYS:HB2	1:B:126:TRP:CZ2	2.07	0.90
1:A:17:ILE:HD12	1:A:56:GLY:HA2	1.55	0.88
1:B:77:GLY:HA2	1:B:80:ARG:HD3	1.55	0.88
1:B:118:LEU:HB2	2:B:166:HOH:O	1.77	0.84
1:B:64:GLU:O	1:B:67:PHE:HB3	1.78	0.83
1:A:120:MET:HE1	1:A:132:ASN:CB	2.08	0.82
1:A:94:VAL:O	1:A:97:CYS:HB2	1.81	0.81
1:A:157:THR:HG22	2:A:257:HOH:O	1.80	0.81
1:B:119:ARG:HH11	1:B:119:ARG:HG3	1.46	0.80
1:A:93:ALA:HA	1:A:96:ARG:HG3	1.62	0.80
1:A:124:LYS:HD2	2:A:200:HOH:O	1.82	0.80
1:A:53:ASN:HD22	1:A:53:ASN:H	1.30	0.80
1:B:124:LYS:HB2	1:B:126:TRP:CH2	2.17	0.79
1:B:93:ALA:O	1:B:96:ARG:HB2	1.82	0.79
1:B:2:ASN:OD1	1:B:5:GLU:HB2	1.81	0.79
1:B:84:LEU:CD2	1:B:112:ALA:HB2	2.13	0.79
1:B:145:ARG:HD2	2:B:261:HOH:O	1.84	0.76
1:A:105:GLN:O	1:A:105:GLN:HG2	1.85	0.76
1:B:116:ASN:O	1:B:119:ARG:HB3	1.86	0.75
1:B:2:ASN:ND2	1:B:4:PHE:HB3	2.01	0.75
1:B:76:ARG:O	1:B:80:ARG:HD2	1.87	0.75
1:A:107:GLY:O	1:A:111:VAL:HG23	1.88	0.74
1:B:92:ASP:OD2	1:B:95:ARG:HD2	1.88	0.74
1:B:75:VAL:O	1:B:79:LEU:HG	1.89	0.72
1:A:119:ARG:O	1:A:123:GLN:HG2	1.90	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:139:TYR:OH	1:A:147:LYS:HE3	1.89	0.71
1:A:79:LEU:HA	1:A:85:LYS:HG3	1.74	0.69
1:B:151:THR:HA	1:B:154:ARG:HB2	1.75	0.69
1:A:76:ARG:O	1:A:80:ARG:HD3	1.93	0.69
1:A:65:LYS:NZ	2:A:254:HOH:O	2.27	0.68
1:A:16:LYS:HG3	1:A:57:VAL:HG22	1.76	0.67
1:A:137:ARG:O	1:A:141:GLN:HB2	1.94	0.67
1:A:120:MET:HE1	1:A:132:ASN:HB2	1.75	0.67
1:A:140:ASN:ND2	2:A:266:HOH:O	2.27	0.66
1:A:120:MET:HE3	1:A:129:ALA:HA	1.77	0.66
1:B:5:GLU:HA	1:B:5:GLU:OE1	1.96	0.66
1:B:114:PHE:CB	1:B:118:LEU:HD11	2.21	0.66
1:B:10:ASP:OD2	1:B:101:ASN:ND2	2.28	0.65
1:A:46:LEU:O	1:A:50:ILE:HG23	1.97	0.64
1:B:106:MET:CG	1:B:111:VAL:HG23	2.28	0.64
1:B:120:MET:HE1	1:B:132:ASN:ND2	2.13	0.64
1:B:81:ASN:O	1:B:85:LYS:HB3	1.98	0.63
1:A:14:ARG:HG3	1:A:18:TYR:CE1	2.34	0.63
1:B:89:ASP:HB3	2:B:186:HOH:O	1.97	0.63
1:B:1:MET:HB3	1:B:158:TRP:CD2	2.34	0.63
1:B:106:MET:SD	1:B:111:VAL:HG23	2.39	0.63
1:B:75:VAL:HG22	1:B:100:ILE:HD11	1.79	0.63
1:A:1:MET:CE	1:A:5:GLU:HB3	2.27	0.62
1:A:85:LYS:NZ	2:A:211:HOH:O	2.31	0.62
1:B:72:ASP:HA	2:B:225:HOH:O	1.99	0.62
1:B:117:SER:HA	1:B:120:MET:HE3	1.82	0.61
1:A:27:ILE:HD13	1:A:58:ILE:HG12	1.83	0.61
1:B:119:ARG:HH11	1:B:119:ARG:CG	2.12	0.61
1:A:53:ASN:H	1:A:53:ASN:ND2	1.96	0.61
1:B:146:ALA:O	1:B:150:ILE:HD12	2.00	0.61
1:B:43:LYS:NZ	2:B:256:HOH:O	2.28	0.60
1:B:1:MET:HG3	1:B:2:ASN:N	2.17	0.60
1:B:38:SER:OG	1:B:41:ALA:HB2	2.02	0.59
1:B:100:ILE:HA	1:B:103:VAL:CG2	2.33	0.59
1:A:116:ASN:O	1:A:120:MET:HG3	2.03	0.58
1:B:26:THR:HG22	1:B:27:ILE:N	2.19	0.58
1:B:95:ARG:HG2	1:B:153:PHE:HA	1.85	0.58
1:B:1:MET:H3	1:B:158:TRP:CD1	2.22	0.58
1:B:79:LEU:O	1:B:85:LYS:HD2	2.04	0.58
1:B:106:MET:HE2	1:B:138:TRP:CD1	2.39	0.58
1:A:81:ASN:HB3	1:A:84:LEU:HB2	1.86	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:57:VAL:HG12	1:B:58:ILE:N	2.19	0.57
1:B:1:MET:HG3	1:B:2:ASN:H	1.68	0.56
1:B:151:THR:CA	1:B:154:ARG:HB2	2.34	0.56
1:A:1:MET:HE2	1:A:9:ILE:CD1	2.36	0.56
1:B:11:GLU:CD	1:B:145:ARG:HH12	2.14	0.55
1:B:158:TRP:HB2	2:B:245:HOH:O	2.05	0.55
1:B:134:ALA:HA	1:B:139:TYR:CD2	2.42	0.55
1:A:38:SER:OG	1:A:40:ASN:HB3	2.07	0.54
1:A:120:MET:CE	1:A:129:ALA:HA	2.36	0.54
1:A:24:TYR:CE1	1:A:35:LYS:HB3	2.43	0.54
1:B:135:LYS:O	1:B:136:SER:O	2.25	0.54
1:B:7:LEU:HD12	1:B:11:GLU:HB2	1.90	0.54
1:A:19:LYS:HG2	1:A:23:GLY:O	2.06	0.54
1:B:130:ALA:HA	1:B:133:LEU:HD12	1.89	0.54
1:A:81:ASN:O	1:A:85:LYS:HB2	2.07	0.54
1:B:66:LEU:O	1:B:69:GLN:HB2	2.08	0.54
1:A:1:MET:HE2	1:A:9:ILE:HD12	1.90	0.54
1:A:117:SER:HA	1:A:120:MET:HE2	1.88	0.53
1:A:158:TRP:O	1:A:162:LYS:HG2	2.09	0.53
1:A:87:VAL:HG11	1:A:118:LEU:HD22	1.91	0.53
1:A:43:LYS:O	1:A:46:LEU:HB3	2.09	0.52
1:A:75:VAL:HG11	1:B:37:PRO:HG3	1.90	0.52
1:B:131:VAL:HG11	2:B:242:HOH:O	2.10	0.52
1:A:17:ILE:CD1	1:A:56:GLY:HA2	2.35	0.52
1:B:133:LEU:O	1:B:136:SER:OG	2.27	0.52
1:B:14:ARG:HG3	1:B:14:ARG:NH1	2.24	0.51
1:B:75:VAL:CG2	1:B:100:ILE:HD11	2.40	0.51
1:B:44:SER:O	1:B:48:LYS:HG3	2.11	0.51
1:B:128:GLU:O	1:B:131:VAL:HG12	2.10	0.51
1:A:26:THR:HG22	1:A:27:ILE:N	2.26	0.50
1:B:92:ASP:CG	1:B:95:ARG:HH11	2.19	0.50
1:B:120:MET:HE1	1:B:132:ASN:CG	2.36	0.50
1:B:114:PHE:HB2	1:B:118:LEU:CD1	2.27	0.49
1:B:157:THR:O	2:B:258:HOH:O	2.20	0.49
1:B:29:ILE:O	1:B:29:ILE:HG22	2.11	0.49
1:B:38:SER:O	1:B:41:ALA:HB3	2.12	0.49
1:B:137:ARG:O	1:B:141:GLN:N	2.43	0.49
1:A:100:ILE:O	1:A:104:PHE:HB2	2.12	0.49
1:B:87:VAL:HA	1:B:122:GLN:HG2	1.93	0.49
1:B:1:MET:CG	1:B:2:ASN:N	2.75	0.49
1:B:151:THR:HA	1:B:154:ARG:CB	2.42	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:103:VAL:O	1:B:107:GLY:N	2.46	0.49
1:A:120:MET:HE1	1:A:132:ASN:HB3	1.90	0.49
1:B:100:ILE:HA	1:B:103:VAL:HG23	1.95	0.48
1:A:116:ASN:HA	1:A:119:ARG:HD3	1.94	0.48
1:A:125:ARG:NH1	2:A:207:HOH:O	2.47	0.48
1:B:137:ARG:C	1:B:141:GLN:HG3	2.39	0.48
1:B:27:ILE:CG1	1:B:28:GLY:H	2.27	0.48
1:B:80:ARG:HG2	2:B:195:HOH:O	2.14	0.48
1:B:151:THR:C	1:B:154:ARG:HB2	2.38	0.48
1:A:119:ARG:NE	2:A:242:HOH:O	2.29	0.48
1:A:162:LYS:CB	1:A:162:LYS:NZ	2.77	0.47
1:B:78:ILE:HA	1:B:84:LEU:HD12	1.95	0.47
1:B:27:ILE:HG12	1:B:28:GLY:H	1.79	0.47
1:A:27:ILE:CD1	1:A:58:ILE:HG12	2.43	0.47
1:B:78:ILE:HG21	1:B:88:TYR:CD1	2.49	0.47
1:B:150:ILE:O	1:B:152:THR:N	2.47	0.47
1:A:91:LEU:O	1:A:96:ARG:NE	2.45	0.47
1:A:148:ARG:HB3	1:A:161:TYR:CZ	2.50	0.47
1:B:120:MET:CE	1:B:132:ASN:ND2	2.76	0.47
1:A:47:ASP:O	1:A:51:GLY:N	2.37	0.47
1:B:103:VAL:O	1:B:107:GLY:HA2	2.15	0.47
1:B:119:ARG:HG3	1:B:119:ARG:NH1	2.18	0.47
1:B:111:VAL:O	1:B:114:PHE:CD1	2.68	0.47
1:B:59:THR:OG1	1:B:62:GLU:HG3	2.15	0.46
1:A:98:ALA:HB3	1:A:153:PHE:CE1	2.49	0.46
1:A:87:VAL:HG11	1:A:118:LEU:CD2	2.45	0.46
1:B:25:TYR:HB2	1:B:34:THR:HG22	1.97	0.46
1:B:125:ARG:HD2	1:B:128:GLU:OE2	2.16	0.46
1:B:159:ASP:HA	1:B:162:LYS:HG3	1.97	0.46
1:A:14:ARG:CB	1:A:18:TYR:CE1	2.99	0.46
1:A:4:PHE:CE1	1:A:29:ILE:HD11	2.50	0.46
1:B:26:THR:CG2	1:B:27:ILE:N	2.79	0.46
1:A:6:MET:HE1	1:A:98:ALA:HA	1.98	0.45
1:A:146:ALA:O	1:A:150:ILE:HD12	2.16	0.45
1:B:106:MET:HG3	1:B:111:VAL:HG23	1.99	0.45
1:A:115:THR:O	1:A:119:ARG:CD	2.64	0.45
1:B:4:PHE:CE1	1:B:29:ILE:CD1	3.00	0.45
1:A:79:LEU:O	1:A:85:LYS:HD2	2.16	0.45
1:A:14:ARG:HB2	1:A:18:TYR:HE1	1.82	0.45
1:A:14:ARG:CG	1:A:18:TYR:CE1	2.99	0.45
1:A:123:GLN:C	1:A:124:LYS:HG3	2.40	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:120:MET:H	1:B:120:MET:HG3	1.58	0.44
1:A:132:ASN:O	1:A:135:LYS:HB2	2.17	0.44
1:B:14:ARG:HH11	1:B:14:ARG:CG	2.30	0.44
1:A:66:LEU:HD23	1:A:66:LEU:HA	1.85	0.44
1:A:55:ASN:CA	2:A:265:HOH:O	2.66	0.44
1:B:7:LEU:O	1:B:11:GLU:N	2.45	0.44
1:A:60:LYS:HE3	1:A:60:LYS:HB3	1.82	0.44
1:A:26:THR:CG2	1:A:27:ILE:N	2.80	0.44
1:A:75:VAL:HG12	1:A:79:LEU:HD11	2.00	0.44
1:A:144:ASN:HB2	2:A:229:HOH:O	2.17	0.44
1:B:7:LEU:HD13	1:B:7:LEU:HA	1.76	0.44
1:A:161:TYR:O	1:A:164:LEU:HB2	2.17	0.43
1:A:72:ASP:OD1	1:B:19:LYS:NZ	2.47	0.43
1:B:75:VAL:HG12	1:B:79:LEU:HD11	2.00	0.43
1:B:152:THR:O	1:B:156:GLY:N	2.41	0.43
1:B:11:GLU:CG	2:B:177:HOH:O	2.66	0.43
1:B:118:LEU:O	1:B:122:GLN:HG3	2.19	0.43
1:A:65:LYS:NZ	1:A:69:GLN:OE1	2.30	0.43
1:B:139:TYR:CD1	1:B:139:TYR:C	2.97	0.43
1:B:66:LEU:O	1:B:69:GLN:N	2.44	0.43
1:A:15:LEU:O	1:A:57:VAL:HG13	2.19	0.43
1:A:137:ARG:HD2	1:A:141:GLN:HE22	1.83	0.43
1:B:6:MET:HB2	1:B:158:TRP:CZ3	2.53	0.43
1:B:103:VAL:O	1:B:107:GLY:CA	2.67	0.43
1:B:139:TYR:CD1	1:B:139:TYR:O	2.72	0.43
1:A:115:THR:O	1:A:119:ARG:HD3	2.19	0.43
1:B:102:MET:HB3	1:B:111:VAL:HG21	2.01	0.43
1:A:52:ARG:HH21	1:A:62:GLU:CD	2.26	0.42
1:B:83:LYS:HB2	1:B:83:LYS:HE2	1.90	0.42
1:B:121:LEU:HD23	1:B:121:LEU:HA	1.96	0.42
1:B:129:ALA:O	1:B:133:LEU:HG	2.18	0.42
1:A:95:ARG:HH11	1:A:95:ARG:HD2	1.50	0.42
1:B:150:ILE:O	1:B:154:ARG:HB2	2.19	0.42
1:A:56:GLY:N	2:A:265:HOH:O	2.51	0.42
1:B:81:ASN:OD1	1:B:82:ALA:N	2.52	0.42
1:A:93:ALA:O	1:A:96:ARG:HB2	2.18	0.42
1:B:31:HIS:HD1	1:B:70:ASP:CG	2.28	0.42
1:B:38:SER:OG	1:B:41:ALA:CB	2.68	0.42
1:A:55:ASN:HA	2:A:265:HOH:O	2.20	0.42
1:B:38:SER:CB	1:B:41:ALA:HB2	2.49	0.42
1:B:108:GLU:HB3	1:B:109:THR:H	1.63	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:88:TYR:CE1	1:A:96:ARG:HD3	2.55	0.41
1:B:14:ARG:HG3	1:B:14:ARG:HH11	1.84	0.41
1:B:117:SER:HB3	1:B:133:LEU:HD21	2.02	0.41
1:A:2:ASN:HA	1:A:3:PRO:HD3	1.69	0.41
1:B:83:LYS:N	1:B:83:LYS:HD3	2.36	0.41
1:A:16:LYS:HB3	1:A:16:LYS:HE2	1.87	0.41
1:A:52:ARG:CZ	2:A:232:HOH:O	2.68	0.41
1:B:78:ILE:HG21	1:B:78:ILE:HD13	1.80	0.41
1:B:138:TRP:HA	1:B:141:GLN:NE2	2.36	0.41
1:B:8:ARG:NH1	2:B:228:HOH:O	2.53	0.41
1:B:150:ILE:O	1:B:151:THR:C	2.64	0.41
1:B:78:ILE:O	1:B:85:LYS:HB3	2.21	0.41
1:B:32:LEU:HD21	1:B:35:LYS:HD2	2.02	0.41
1:A:84:LEU:HD23	1:A:84:LEU:HA	1.91	0.41
1:A:55:ASN:N	2:A:265:HOH:O	2.54	0.41
1:A:77:GLY:O	1:A:78:ILE:C	2.63	0.41
1:B:2:ASN:HD21	1:B:4:PHE:HB3	1.84	0.41
1:B:63:ALA:O	1:B:67:PHE:HB2	2.21	0.41
1:A:84:LEU:O	1:A:85:LYS:C	2.64	0.40
1:B:129:ALA:O	1:B:132:ASN:CB	2.69	0.40
1:B:138:TRP:HA	1:B:141:GLN:CD	2.46	0.40
1:B:160:ALA:CB	2:B:258:HOH:O	2.70	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	162/164 (99%)	140 (86%)	21 (13%)	1 (1%)	21	17
1	B	162/164 (99%)	136 (84%)	21 (13%)	5 (3%)	3	1
All	All	324/328 (99%)	276 (85%)	42 (13%)	6 (2%)	6	3

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	93	ALA
1	B	115	THR
1	B	136	SER
1	B	151	THR
1	A	126	TRP
1	B	143	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	138/138 (100%)	106 (77%)	32 (23%)	1	0
1	B	138/138 (100%)	106 (77%)	32 (23%)	1	0
All	All	276/276 (100%)	212 (77%)	64 (23%)	1	0

All (64) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	7	LEU
1	A	13	LEU
1	A	19	LYS
1	A	22	GLU
1	A	27	ILE
1	A	32	LEU
1	A	36	SER
1	A	43	LYS
1	A	48	LYS
1	A	50	ILE
1	A	53	ASN
1	A	58	ILE
1	A	59	THR
1	A	60	LYS
1	A	78	ILE
1	A	79	LEU

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Mol	Chain	Res	Type
1	A	80	ARG
1	A	100	ILE
1	A	103	VAL
1	A	104	PHE
1	A	117	SER
1	A	119	ARG
1	A	122	GLN
1	A	124	LYS
1	A	125	ARG
1	A	127	ASP
1	A	128	GLU
1	A	137	ARG
1	A	141	GLN
1	A	145	ARG
1	A	162	LYS
1	B	1	MET
1	B	5	GLU
1	B	6	MET
1	B	7	LEU
1	B	8	ARG
1	B	9	ILE
1	B	14	ARG
1	B	16	LYS
1	B	17	ILE
1	B	21	THR
1	B	33	LEU
1	B	35	LYS
1	B	39	LEU
1	B	48	LYS
1	B	50	ILE
1	B	52	ARG
1	B	65	LYS
1	B	76	ARG
1	B	80	ARG
1	B	84	LEU
1	B	85	LYS
1	B	87	VAL
1	B	102	MET
1	B	104	PHE
1	B	111	VAL
1	B	119	ARG
1	B	124	LYS

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Mol	Chain	Res	Type
1	B	131	VAL
1	B	137	ARG
1	B	140	ASN
1	B	154	ARG
1	B	162	LYS

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

Mol	Chain	Res	Type
1	A	40	ASN
1	A	53	ASN
1	A	116	ASN
1	A	141	GLN
1	B	40	ASN
1	B	68	ASN
1	B	140	ASN
1	B	141	GLN
1	B	144	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 oligosaccharides 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 ⓘ

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	164/164 (100%)	-0.79	0 100 100	5, 16, 26, 43	0
1	B	164/164 (100%)	-0.50	0 100 100	8, 21, 34, 60	0
All	All	328/328 (100%)	-0.65	0 100 100	5, 19, 32, 60	0

There are no RSRZ outliers to report.

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.