



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 20, 2017 – 10:45 PM EST

PDB ID : 176L  
Title : PROTEIN FLEXIBILITY AND ADAPTABILITY SEEN IN 25 CRYSTAL FORMS OF T4 LYSOZYME  
Authors : Zhang, X.-J.; Weaver, L.; Dubose, R.; Matthews, B.W.  
Deposited on : unknown  
Resolution : 2.20 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](mailto: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.

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

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20030345

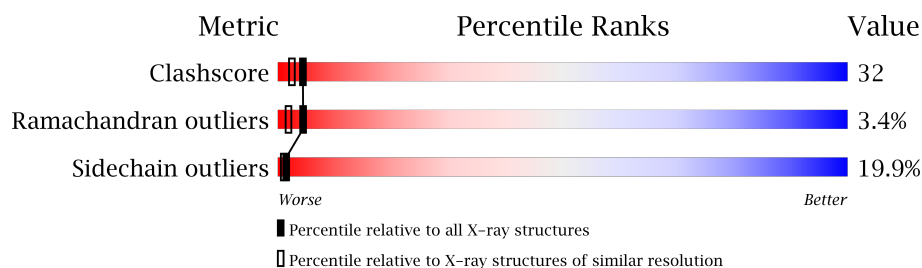
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.20 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	112137	4730 (2.20-2.20)
Ramachandran outliers	110173	4656 (2.20-2.20)
Sidechain outliers	110143	4657 (2.20-2.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	164	 37% 42% 18% .
1	B	164	 45% 41% 13% .

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2640 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
			1307	819	239	244	5			
1	B	164	Total	C	N	O	S	0	0	0
			1307	819	239	244	5			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	32	THR	LEU	CONFLICT	UNP P00720
A	34	LYS	THR	CONFLICT	UNP P00720
A	35	VAL	LYS	CONFLICT	UNP P00720
A	36	ASP	SER	CONFLICT	UNP P00720
A	37	GLY	PRO	CONFLICT	UNP P00720
A	38	ASN	SER	CONFLICT	UNP P00720
A	39	SER	LEU	CONFLICT	UNP P00720
A	54	THR	CYS	CONFLICT	UNP P00720
A	97	ALA	CYS	CONFLICT	UNP P00720
B	32	THR	LEU	CONFLICT	UNP P00720
B	34	LYS	THR	CONFLICT	UNP P00720
B	35	VAL	LYS	CONFLICT	UNP P00720
B	36	ASP	SER	CONFLICT	UNP P00720
B	37	GLY	PRO	CONFLICT	UNP P00720
B	38	ASN	SER	CONFLICT	UNP P00720
B	39	SER	LEU	CONFLICT	UNP P00720
B	54	THR	CYS	CONFLICT	UNP P00720
B	97	ALA	CYS	CONFLICT	UNP P00720

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Cl	0	0
			1	1		

- Molecule 3 is water.

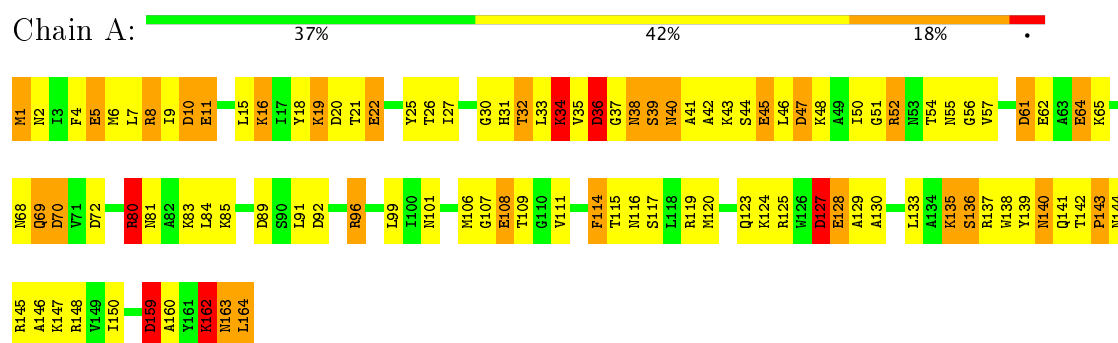
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	7	Total 7	O 7	0	0
3	B	18	Total 18	O 18	0	0

### 3 Residue-property plots

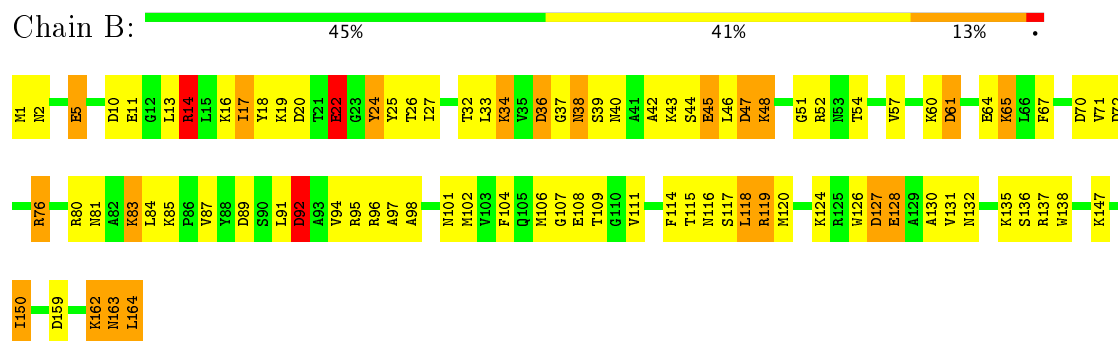
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are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

Note EDS was not executed.

#### • Molecule 1: T4 LYSOZYME



#### • Molecule 1: T4 LYSOZYME



## 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, $\alpha$ , $\beta$ , $\gamma$	49.60Å 127.10Å 29.10Å 90.00° 98.40° 90.00°	Depositor
Resolution (Å)	14.00 – 2.20	Depositor
% Data completeness (in resolution range)	(Not available) (14.00-2.20)	Depositor
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	TNT	Depositor
R, $R_{free}$	0.183 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2640	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	31.0	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: CL

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	1.00	7/1326 (0.5%)	1.47	19/1784 (1.1%)
1	B	1.01	5/1326 (0.4%)	1.45	22/1784 (1.2%)
All	All	1.01	12/2652 (0.5%)	1.46	41/3568 (1.1%)

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	11	GLU	CD-OE2	6.97	1.33	1.25
1	A	45	GLU	CD-OE1	6.70	1.33	1.25
1	A	62	GLU	CD-OE2	6.07	1.32	1.25
1	A	22	GLU	CD-OE2	5.93	1.32	1.25
1	B	108	GLU	CD-OE2	5.88	1.32	1.25
1	B	22	GLU	CD-OE2	5.79	1.32	1.25
1	B	11	GLU	CD-OE2	5.71	1.31	1.25
1	B	45	GLU	CD-OE2	5.65	1.31	1.25
1	A	108	GLU	CD-OE2	5.64	1.31	1.25
1	B	128	GLU	CD-OE2	5.24	1.31	1.25
1	A	128	GLU	CD-OE2	5.14	1.31	1.25
1	A	5	GLU	CD-OE2	5.07	1.31	1.25

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	61	ASP	CB-CG-OD2	-10.56	108.80	118.30
1	B	89	ASP	CB-CG-OD2	-8.73	110.44	118.30
1	B	92	ASP	CB-CG-OD2	-8.64	110.52	118.30
1	A	142	THR	C-N-CD	-7.87	103.28	120.60
1	B	119	ARG	NE-CZ-NH1	7.72	124.16	120.30
1	A	61	ASP	CB-CG-OD1	7.55	125.09	118.30
1	B	89	ASP	CB-CG-OD1	7.41	124.97	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	72	ASP	CB-CG-OD2	-7.29	111.74	118.30
1	B	72	ASP	CB-CG-OD2	-6.83	112.15	118.30
1	A	159	ASP	CB-CG-OD2	-6.74	112.23	118.30
1	A	10	ASP	CB-CG-OD1	6.49	124.14	118.30
1	B	127	ASP	CB-CG-OD2	-6.30	112.63	118.30
1	A	92	ASP	CB-CG-OD2	-6.05	112.85	118.30
1	A	70	ASP	CB-CG-OD1	6.02	123.72	118.30
1	B	80	ARG	NE-CZ-NH1	6.00	123.30	120.30
1	B	47	ASP	CB-CG-OD2	-5.98	112.92	118.30
1	B	159	ASP	CB-CG-OD1	5.93	123.63	118.30
1	B	61	ASP	CB-CG-OD2	-5.80	113.08	118.30
1	B	127	ASP	CB-CG-OD1	5.73	123.46	118.30
1	B	159	ASP	CB-CG-OD2	-5.73	113.14	118.30
1	A	47	ASP	CB-CG-OD2	-5.66	113.20	118.30
1	B	20	ASP	CB-CG-OD1	5.58	123.33	118.30
1	B	119	ARG	NE-CZ-NH2	-5.58	117.51	120.30
1	B	20	ASP	CB-CG-OD2	-5.55	113.30	118.30
1	A	89	ASP	CB-CG-OD1	5.54	123.29	118.30
1	B	137	ARG	NE-CZ-NH1	5.53	123.06	120.30
1	A	47	ASP	CB-CG-OD1	5.49	123.24	118.30
1	A	70	ASP	CB-CG-OD2	-5.47	113.37	118.30
1	B	10	ASP	CB-CG-OD2	-5.46	113.39	118.30
1	A	89	ASP	CB-CG-OD2	-5.37	113.46	118.30
1	B	96	ARG	NE-CZ-NH1	5.33	122.97	120.30
1	A	10	ASP	CB-CG-OD2	-5.33	113.50	118.30
1	A	80	ARG	NE-CZ-NH1	5.32	122.96	120.30
1	A	127	ASP	CB-CG-OD2	-5.27	113.56	118.30
1	B	61	ASP	CB-CG-OD1	5.25	123.02	118.30
1	B	47	ASP	CB-CG-OD1	5.23	123.01	118.30
1	B	14	ARG	NE-CZ-NH2	-5.20	117.70	120.30
1	A	20	ASP	CB-CG-OD1	5.12	122.91	118.30
1	A	164	LEU	N-CA-CB	5.10	120.61	110.40
1	B	92	ASP	CB-CG-OD1	5.05	122.84	118.30
1	A	36	ASP	CB-CG-OD2	5.01	122.81	118.30

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	1307	0	1324	95	0
1	B	1307	0	1324	72	0
2	A	1	0	0	0	0
3	A	7	0	0	0	0
3	B	18	0	0	1	0
All	All	2640	0	2648	167	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 32.

All (167) 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:144:ASN:HA	1:A:147:LYS:HE3	1.39	1.02
1:B:19:LYS:HD3	1:B:25:TYR:CE1	2.00	0.96
1:A:162:LYS:HE3	1:A:163:ASN:HB2	1.44	0.96
1:A:96:ARG:HB2	1:A:96:ARG:HH11	1.27	0.95
1:A:34:LYS:H	1:A:34:LYS:HD3	1.34	0.92
1:B:124:LYS:HD3	1:B:126:TRP:CZ2	2.03	0.92
1:A:35:VAL:CG2	1:A:45:GLU:HG2	2.06	0.85
1:B:34:LYS:HD3	1:B:34:LYS:H	1.40	0.84
1:A:162:LYS:HE3	1:A:163:ASN:CB	2.07	0.84
1:B:13:LEU:HD12	1:B:14:ARG:N	1.95	0.81
1:B:124:LYS:HD3	1:B:126:TRP:HZ2	1.44	0.81
1:A:25:TYR:O	1:A:32:THR:HG22	1.82	0.80
1:B:34:LYS:CD	1:B:34:LYS:H	1.96	0.79
1:B:13:LEU:HD12	1:B:14:ARG:H	1.51	0.76
1:A:143:PRO:O	1:A:147:LYS:HB2	1.86	0.75
1:A:120:MET:CE	1:A:128:GLU:HB3	2.16	0.75
1:A:61:ASP:O	1:A:65:LYS:HG2	1.87	0.74
1:A:116:ASN:O	1:A:120:MET:HG3	1.88	0.74
1:A:5:GLU:O	1:A:9:ILE:HD12	1.89	0.72
1:A:34:LYS:CD	1:A:34:LYS:H	1.99	0.72
1:A:96:ARG:CB	1:A:96:ARG:HH11	2.01	0.71
1:B:163:ASN:HD22	1:B:164:LEU:N	1.89	0.71
1:B:61:ASP:O	1:B:65:LYS:HE3	1.91	0.70
1:A:35:VAL:HG21	1:A:45:GLU:HG2	1.71	0.70
1:A:107:GLY:O	1:A:111:VAL:HG23	1.93	0.69
1:A:144:ASN:HA	1:A:147:LYS:CE	2.20	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:76:ARG:HB3	1:B:76:ARG:HH11	1.59	0.67
1:B:163:ASN:O	1:B:164:LEU:HB2	1.94	0.67
1:A:16:LYS:HE2	1:A:57:VAL:CG2	2.25	0.66
1:B:76:ARG:CB	1:B:76:ARG:HH11	2.09	0.66
1:B:2:ASN:N	1:B:5:GLU:OE1	2.28	0.66
1:B:45:GLU:O	1:B:48:LYS:HB3	1.97	0.65
1:A:38:ASN:OD1	1:A:40:ASN:HB2	1.97	0.64
1:A:81:ASN:OD1	1:A:83:LYS:N	2.30	0.64
1:B:38:ASN:OD1	1:B:39:SER:N	2.30	0.64
1:B:61:ASP:O	1:B:65:LYS:HG2	1.98	0.63
1:A:162:LYS:CG	1:A:163:ASN:H	2.11	0.63
1:A:120:MET:HE2	1:A:128:GLU:HB3	1.80	0.63
1:A:19:LYS:HD3	1:A:25:TYR:CE1	2.34	0.63
1:B:81:ASN:O	1:B:85:LYS:HB3	2.00	0.62
1:A:65:LYS:O	1:A:69:GLN:HG2	1.98	0.62
1:B:1:MET:HG3	1:B:5:GLU:OE1	2.00	0.62
1:B:45:GLU:HA	1:B:45:GLU:OE1	1.98	0.62
1:B:116:ASN:CG	1:B:119:ARG:HH22	2.03	0.62
1:A:18:TYR:O	1:A:26:THR:N	2.28	0.61
1:A:38:ASN:O	1:A:41:ALA:HB3	2.01	0.61
1:B:162:LYS:O	1:B:163:ASN:ND2	2.34	0.61
1:A:80:ARG:HB2	1:A:80:ARG:CZ	2.29	0.61
1:A:120:MET:HE1	1:A:128:GLU:HB3	1.83	0.60
1:A:2:ASN:N	1:A:5:GLU:OE1	2.31	0.60
1:B:38:ASN:OD1	1:B:40:ASN:N	2.29	0.60
1:B:76:ARG:NH1	1:B:76:ARG:HB3	2.17	0.60
1:B:39:SER:O	1:B:42:ALA:HB3	2.02	0.60
1:A:16:LYS:HA	1:A:56:GLY:O	2.02	0.59
1:B:19:LYS:HD3	1:B:25:TYR:CD1	2.36	0.59
1:A:81:ASN:OD1	1:A:84:LEU:N	2.36	0.59
1:B:52:ARG:O	1:B:54:THR:HG23	2.02	0.59
1:A:47:ASP:OD1	1:A:54:THR:HG23	2.03	0.58
1:A:1:MET:HA	1:A:5:GLU:OE1	2.04	0.58
1:A:36:ASP:CG	1:A:37:GLY:H	2.03	0.57
1:B:81:ASN:OD1	1:B:83:LYS:N	2.29	0.57
1:B:81:ASN:HB3	1:B:84:LEU:HB2	1.87	0.57
1:B:94:VAL:O	1:B:97:ALA:HB3	2.05	0.56
1:B:36:ASP:OD1	1:B:37:GLY:N	2.28	0.56
1:B:36:ASP:CG	1:B:37:GLY:H	2.09	0.56
1:B:120:MET:HE1	1:B:128:GLU:HB3	1.87	0.55
1:A:11:GLU:OE1	1:A:30:GLY:HA3	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:16:LYS:HG3	1:B:57:VAL:CG2	2.37	0.55
1:A:137:ARG:HB3	1:A:141:GLN:OE1	2.05	0.55
1:A:31:HIS:ND1	1:A:70:ASP:OD2	2.29	0.54
1:B:131:VAL:HG12	1:B:132:ASN:N	2.23	0.54
1:A:25:TYR:CD2	1:A:42:ALA:HB2	2.43	0.54
1:B:25:TYR:HE2	1:B:37:GLY:O	1.90	0.54
1:A:36:ASP:OD1	1:A:37:GLY:N	2.27	0.54
1:B:127:ASP:O	1:B:130:ALA:HB3	2.08	0.54
1:B:44:SER:O	1:B:48:LYS:HB2	2.08	0.53
1:A:11:GLU:OE2	1:A:145:ARG:NH1	2.36	0.53
1:A:31:HIS:HD1	1:A:70:ASP:CG	2.12	0.53
1:A:163:ASN:O	1:A:164:LEU:HB2	2.09	0.52
1:A:143:PRO:O	1:A:147:LYS:HE2	2.10	0.52
1:A:52:ARG:HG3	1:A:52:ARG:HH11	1.74	0.52
1:A:25:TYR:HE2	1:A:37:GLY:O	1.92	0.51
1:B:116:ASN:ND2	1:B:119:ARG:HH22	2.08	0.51
1:B:16:LYS:HG3	1:B:57:VAL:HG22	1.93	0.51
1:A:120:MET:HE2	1:A:128:GLU:CB	2.41	0.51
1:A:52:ARG:O	1:A:54:THR:HG23	2.11	0.51
1:B:98:ALA:O	1:B:101:ASN:HB3	2.11	0.51
1:A:47:ASP:O	1:A:51:GLY:N	2.43	0.50
1:B:67:PHE:O	1:B:71:VAL:HG23	2.10	0.50
1:B:107:GLY:O	1:B:111:VAL:HG23	2.12	0.50
1:B:163:ASN:ND2	1:B:164:LEU:O	2.45	0.50
1:B:138:TRP:CZ3	1:B:150:ILE:HD11	2.47	0.50
1:A:16:LYS:HE2	1:A:57:VAL:HG23	1.92	0.49
1:A:114:PHE:N	1:A:114:PHE:CD1	2.80	0.49
1:A:96:ARG:NH1	1:A:96:ARG:HB2	2.10	0.49
1:A:91:LEU:HB2	1:A:96:ARG:HG2	1.94	0.49
1:B:114:PHE:CD1	1:B:114:PHE:N	2.81	0.49
1:B:120:MET:CE	1:B:128:GLU:HB3	2.42	0.49
1:A:146:ALA:O	1:A:150:ILE:N	2.36	0.48
1:B:26:THR:HG22	1:B:27:ILE:N	2.27	0.48
1:A:137:ARG:O	1:A:141:GLN:N	2.47	0.48
1:A:34:LYS:N	1:A:34:LYS:CD	2.73	0.48
1:A:114:PHE:N	1:A:114:PHE:HD1	2.11	0.48
1:B:18:TYR:O	1:B:26:THR:N	2.29	0.48
1:A:140:ASN:HD22	1:A:140:ASN:HA	1.44	0.47
1:A:162:LYS:HG3	1:A:163:ASN:H	1.79	0.47
1:A:15:LEU:O	1:A:57:VAL:HA	2.15	0.47
1:B:22:GLU:HG2	1:B:24:TYR:CE2	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:159:ASP:O	1:A:162:LYS:HG2	2.14	0.47
1:A:123:GLN:OE1	1:A:125:ARG:NH2	2.48	0.47
1:A:8:ARG:NH2	1:A:64:GLU:OE2	2.36	0.47
1:A:38:ASN:OD1	1:A:40:ASN:N	2.26	0.47
1:A:52:ARG:HH11	1:A:52:ARG:CG	2.27	0.47
1:B:138:TRP:HZ3	1:B:150:ILE:HD11	1.81	0.46
1:A:39:SER:C	1:A:43:LYS:HD2	2.37	0.45
1:A:114:PHE:HD2	1:A:133:LEU:HD23	1.80	0.45
1:A:138:TRP:HZ3	1:A:150:ILE:HD11	1.81	0.45
1:A:46:LEU:HD23	1:A:56:GLY:HA2	1.98	0.45
1:B:17:ILE:HG22	1:B:17:ILE:O	2.09	0.45
1:A:7:LEU:CD2	1:A:145:ARG:NH2	2.80	0.45
1:A:2:ASN:OD1	1:A:4:PHE:HB2	2.17	0.45
1:B:51:GLY:O	1:B:52:ARG:HB3	2.17	0.45
1:A:120:MET:HB2	1:A:129:ALA:HB2	1.99	0.44
1:A:65:LYS:H	1:A:65:LYS:HG2	1.47	0.44
1:B:70:ASP:HB3	1:B:104:PHE:CE2	2.52	0.44
1:A:2:ASN:OD1	1:A:4:PHE:N	2.51	0.44
1:A:50:ILE:HD13	1:A:50:ILE:HA	1.78	0.43
1:A:99:LEU:HA	1:A:99:LEU:HD12	1.78	0.43
1:B:22:GLU:HB3	1:B:24:TYR:CE2	2.52	0.43
1:A:10:ASP:C	1:A:11:GLU:HG2	2.38	0.43
1:B:92:ASP:OD1	1:B:95:ARG:HD2	2.18	0.43
1:A:123:GLN:HB2	1:A:125:ARG:HG3	1.99	0.43
1:A:6:MET:CE	1:A:101:ASN:HD22	2.31	0.43
1:A:162:LYS:CE	1:A:163:ASN:H	2.31	0.43
1:B:163:ASN:ND2	1:B:164:LEU:N	2.62	0.43
1:A:119:ARG:O	1:A:123:GLN:HG3	2.18	0.43
1:A:33:LEU:O	1:A:35:VAL:N	2.51	0.43
1:B:24:TYR:CD1	1:B:24:TYR:N	2.82	0.42
1:A:162:LYS:HE3	1:A:163:ASN:H	1.84	0.42
1:B:43:LYS:O	1:B:46:LEU:HB3	2.19	0.42
1:A:120:MET:H	1:A:120:MET:HG3	1.62	0.42
1:A:19:LYS:NZ	1:A:25:TYR:CE1	2.85	0.42
1:B:33:LEU:HD22	1:B:34:LYS:HZ3	1.85	0.42
1:B:76:ARG:NH1	1:B:76:ARG:CB	2.78	0.42
1:B:34:LYS:N	1:B:34:LYS:HE2	2.35	0.41
1:A:148:ARG:HG3	1:A:160:ALA:HB1	2.02	0.41
1:B:84:LEU:O	1:B:87:VAL:HB	2.20	0.41
1:A:114:PHE:CD2	1:A:133:LEU:HD23	2.55	0.41
1:A:26:THR:HG22	1:A:27:ILE:N	2.36	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:35:VAL:HG23	1:A:45:GLU:HG2	1.94	0.41
1:B:120:MET:HE1	1:B:128:GLU:CB	2.50	0.41
1:A:135:LYS:O	1:A:136:SER:O	2.39	0.41
1:B:120:MET:CE	1:B:128:GLU:CB	2.98	0.41
1:B:47:ASP:OD1	1:B:54:THR:HG23	2.19	0.41
1:A:39:SER:O	1:A:43:LYS:HD2	2.20	0.41
1:B:60:LYS:HD2	3:B:181:HOH:O	2.20	0.41
1:A:33:LEU:HD23	1:A:33:LEU:HA	1.94	0.41
1:B:22:GLU:HG2	1:B:24:TYR:CZ	2.56	0.41
1:B:26:THR:CG2	1:B:27:ILE:N	2.84	0.41
1:B:114:PHE:O	1:B:118:LEU:HG	2.20	0.41
1:A:162:LYS:CG	1:A:163:ASN:N	2.80	0.40
1:B:1:MET:HA	1:B:5:GLU:OE1	2.21	0.40
1:A:127:ASP:O	1:A:130:ALA:HB3	2.21	0.40
1:B:87:VAL:O	1:B:91:LEU:HG	2.20	0.40
1:A:38:ASN:OD1	1:A:39:SER:N	2.54	0.40
1:B:102:MET:O	1:B:106:MET:HG2	2.21	0.40
1:A:139:TYR:CD1	1:A:139:TYR:C	2.94	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	162/164 (99%)	138 (85%)	16 (10%)	8 (5%)	2	1
1	B	162/164 (99%)	147 (91%)	12 (7%)	3 (2%)	9	6
All	All	324/328 (99%)	285 (88%)	28 (9%)	11 (3%)	4	2

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	36	ASP
1	A	136	SER
1	A	162	LYS
1	B	38	ASN
1	A	34	LYS
1	A	38	ASN
1	A	135	LYS
1	A	143	PRO
1	B	36	ASP
1	B	136	SER
1	A	115	THR

### 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	136/136 (100%)	105 (77%)	31 (23%)	1	1
1	B	136/136 (100%)	113 (83%)	23 (17%)	2	2
All	All	272/272 (100%)	218 (80%)	54 (20%)	1	1

All (54) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	8	ARG
1	A	16	LYS
1	A	19	LYS
1	A	21	THR
1	A	22	GLU
1	A	32	THR
1	A	34	LYS
1	A	39	SER
1	A	40	ASN
1	A	44	SER
1	A	48	LYS
1	A	52	ARG

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Mol	Chain	Res	Type
1	A	55	ASN
1	A	64	GLU
1	A	68	ASN
1	A	69	GLN
1	A	80	ARG
1	A	85	LYS
1	A	96	ARG
1	A	106	MET
1	A	108	GLU
1	A	109	THR
1	A	114	PHE
1	A	117	SER
1	A	124	LYS
1	A	127	ASP
1	A	140	ASN
1	A	159	ASP
1	A	162	LYS
1	A	163	ASN
1	B	5	GLU
1	B	14	ARG
1	B	17	ILE
1	B	22	GLU
1	B	24	TYR
1	B	32	THR
1	B	34	LYS
1	B	48	LYS
1	B	64	GLU
1	B	65	LYS
1	B	76	ARG
1	B	83	LYS
1	B	92	ASP
1	B	109	THR
1	B	115	THR
1	B	117	SER
1	B	118	LEU
1	B	135	LYS
1	B	147	LYS
1	B	150	ILE
1	B	162	LYS
1	B	163	ASN
1	B	164	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	53	ASN
1	A	140	ASN
1	B	53	ASN
1	B	116	ASN
1	B	163	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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 1 ligands modelled in this entry, 1 is 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 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.