



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 08:16 AM UTC

PDB ID : 150L / pdb_0000150L
Title : CONSERVATION OF SOLVENT-BINDING SITES IN 10 CRYSTAL FORMS OF T4 LYSOZYME
Authors : Faber, H.R.; Matthews, B.W.
Deposited on : 1994-01-25
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

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

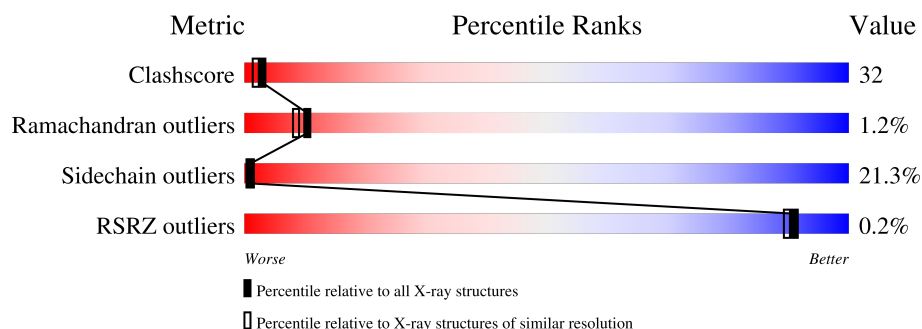
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (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 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	% 41% 41% 16% .
1	B	164	41% 38% 18% ..
1	C	164	57% 29% 10% ..
1	D	164	37% 44% 17% ..

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5327 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
			1303	821	235	241	6			
1	B	162	Total	C	N	O	S	0	0	0
			1284	809	233	236	6			
1	C	162	Total	C	N	O	S	0	0	0
			1288	811	234	237	6			
1	D	162	Total	C	N	O	S	0	0	0
			1288	811	234	237	6			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	6	ILE	MET	conflict	UNP P00720
B	6	ILE	MET	conflict	UNP P00720
C	6	ILE	MET	conflict	UNP P00720
D	6	ILE	MET	conflict	UNP P00720

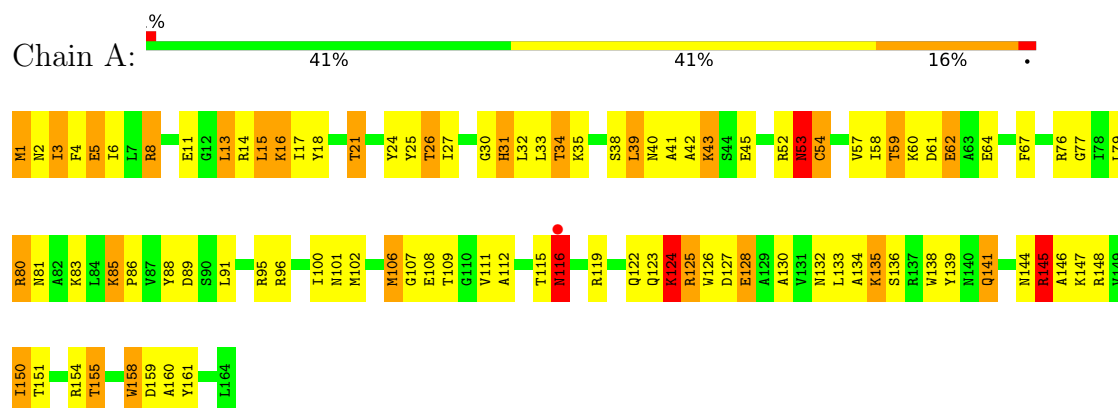
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	58	Total	O	0	0
			58	58		
2	B	21	Total	O	0	0
			21	21		
2	C	49	Total	O	0	0
			49	49		
2	D	36	Total	O	0	0
			36	36		

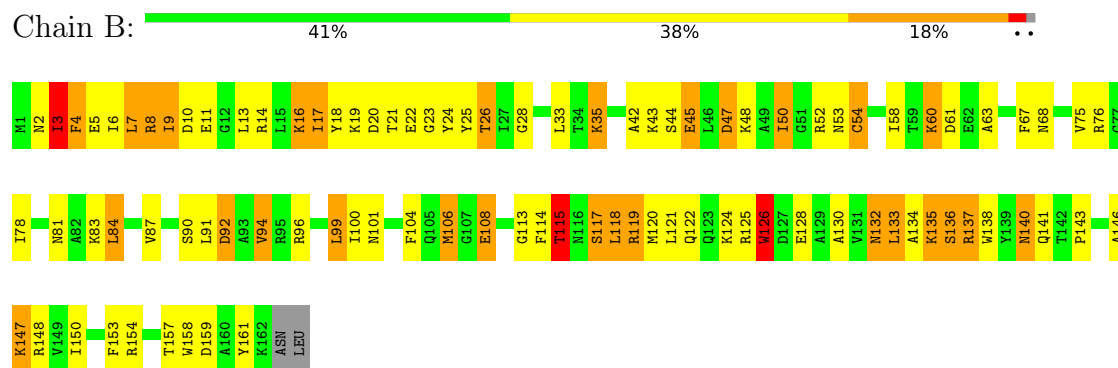
3 Residue-property plots

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.

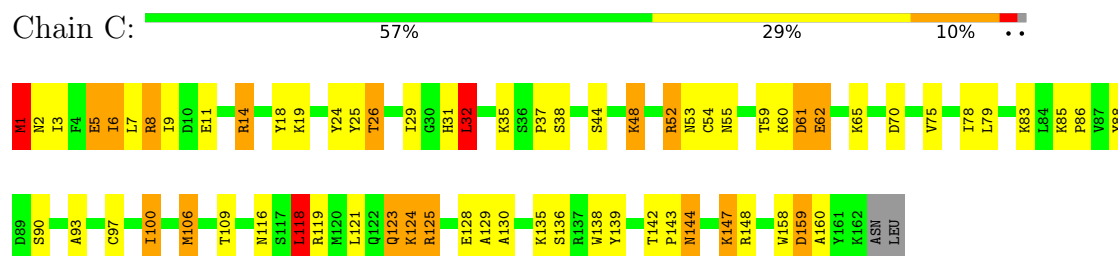
• Molecule 1: T4 LYSOZYME



• Molecule 1: T4 LYSOZYME



• Molecule 1: T4 LYSOZYME



• Molecule 1: T4 LYSOZYME

Chain D: 37% 44% 17% ..



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	72.20Å 73.80Å 150.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	6.00 – 2.20 6.00 – 2.22	Depositor EDS
% Data completeness (in resolution range)	(Not available) (6.00-2.20) 80.7 (6.00-2.22)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.26 (at 2.22Å)	Xtriage
Refinement program	TNT	Depositor
R, R_{free}	0.210 , (Not available) 0.209 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	24.0	Xtriage
Anisotropy	0.072	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.46 , 138.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.036 for k,h,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5327	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.65% 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

5.1 Standard geometry

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	1.43	3/1323 (0.2%)	1.80	23/1782 (1.3%)
1	B	1.36	1/1304 (0.1%)	1.77	16/1758 (0.9%)
1	C	1.54	5/1308 (0.4%)	1.86	29/1763 (1.6%)
1	D	1.44	3/1308 (0.2%)	1.82	19/1763 (1.1%)
All	All	1.44	12/5243 (0.2%)	1.81	87/7066 (1.2%)

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	100	ILE	CA-CB	-7.02	1.45	1.54
1	A	3	ILE	N-CA	-6.07	1.39	1.46
1	C	97	CYS	CA-CB	6.04	1.63	1.53
1	D	152	THR	CA-C	-6.00	1.44	1.52
1	C	62	GLU	N-CA	-5.80	1.39	1.46
1	D	89	ASP	C-N	-5.46	1.26	1.34
1	D	37	PRO	CA-C	-5.31	1.47	1.53
1	A	3	ILE	CA-CB	-5.13	1.48	1.54
1	C	129	ALA	N-CA	-5.11	1.40	1.46
1	B	108	GLU	CD-OE2	5.11	1.35	1.25
1	C	2	ASN	CA-C	5.10	1.60	1.53
1	C	130	ALA	CA-C	-5.09	1.46	1.52

All (87) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	3	ILE	N-CA-C	11.18	123.37	110.62
1	B	94	VAL	N-CA-CB	9.35	120.89	110.51
1	D	97	CYS	CB-CA-C	8.28	125.95	110.63
1	A	101	ASN	CA-CB-CG	8.25	120.85	112.60
1	A	53	ASN	CA-CB-CG	7.53	120.13	112.60
1	A	89	ASP	CA-CB-CG	7.50	120.10	112.60
1	D	44	SER	N-CA-CB	7.39	120.99	110.12
1	D	37	PRO	N-CA-CB	7.39	107.42	102.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	92	ASP	N-CA-C	-7.29	101.13	110.53
1	D	78	ILE	CB-CA-C	-7.07	102.46	112.22
1	C	6	ILE	CA-C-N	-6.99	109.68	120.31
1	C	6	ILE	C-N-CA	-6.99	109.68	120.31
1	C	97	CYS	CB-CA-C	6.92	123.96	110.67
1	D	105	GLN	N-CA-C	6.75	118.72	111.36
1	B	54	CYS	N-CA-C	6.73	118.69	111.36
1	C	70	ASP	CA-CB-CG	-6.58	106.02	112.60
1	A	126	TRP	N-CA-C	6.46	118.86	111.11
1	B	4	PHE	CA-C-N	6.37	129.10	120.44
1	B	4	PHE	C-N-CA	6.37	129.10	120.44
1	A	145	ARG	CD-NE-CZ	6.25	133.14	124.40
1	D	37	PRO	CB-CA-C	-6.20	104.77	112.89
1	D	51	GLY	CA-C-N	6.16	130.28	122.16
1	D	51	GLY	C-N-CA	6.16	130.28	122.16
1	D	91	LEU	N-CA-C	6.13	118.62	110.53
1	C	9	ILE	CB-CA-C	-6.11	103.89	112.14
1	C	1	MET	CA-C-N	-6.03	109.99	120.97
1	C	1	MET	C-N-CA	-6.03	109.99	120.97
1	A	34	THR	CA-C-N	-6.02	112.61	122.54
1	A	34	THR	C-N-CA	-6.02	112.61	122.54
1	A	53	ASN	CB-CA-C	6.01	120.40	110.78
1	C	123	GLN	CB-CG-CD	-6.00	102.41	112.60
1	D	42	ALA	N-CA-CB	5.96	118.66	110.01
1	C	3	ILE	N-CA-CB	5.96	117.92	110.47
1	B	21	THR	N-CA-C	-5.94	104.88	111.71
1	B	26	THR	N-CA-CB	-5.85	100.72	111.37
1	C	61	ASP	CA-C-N	-5.77	112.55	120.28
1	C	61	ASP	C-N-CA	-5.77	112.55	120.28
1	D	41	ALA	N-CA-C	-5.74	105.10	111.36
1	A	116	ASN	O-C-N	5.67	127.92	122.03
1	A	106	MET	N-CA-C	5.59	120.22	113.17
1	C	93	ALA	N-CA-C	5.58	117.45	111.36
1	D	155	THR	N-CA-C	5.57	118.20	111.40
1	C	37	PRO	N-CA-CB	5.56	106.43	102.65
1	C	2	ASN	CB-CA-C	5.55	121.05	109.68
1	D	69	GLN	N-CA-CB	5.53	119.58	110.39
1	C	142	THR	CA-C-N	5.53	124.72	118.97
1	C	142	THR	C-N-CA	5.53	124.72	118.97
1	A	159	ASP	N-CA-CB	5.53	118.18	109.94
1	B	126	TRP	N-CA-C	5.49	117.97	111.33
1	A	26	THR	N-CA-CB	-5.48	102.72	111.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	144	ASN	N-CA-C	5.46	116.91	111.07
1	D	89	ASP	CB-CA-C	-5.45	102.32	110.88
1	C	6	ILE	O-C-N	-5.44	116.58	121.91
1	C	136	SER	N-CA-C	5.44	117.90	110.55
1	A	145	ARG	NE-CZ-NH1	5.43	126.93	121.50
1	B	76	ARG	N-CA-CB	5.41	118.56	110.22
1	B	3	ILE	N-CA-C	5.39	115.53	110.30
1	C	31	HIS	CA-CB-CG	-5.36	108.44	113.80
1	B	83	LYS	CA-C-N	-5.35	111.68	121.52
1	B	83	LYS	C-N-CA	-5.35	111.68	121.52
1	D	161	TYR	N-CA-CB	5.34	119.80	110.98
1	D	5	GLU	CA-C-O	5.34	126.21	120.55
1	A	31	HIS	CA-CB-CG	-5.30	108.50	113.80
1	C	11	GLU	CB-CA-C	-5.26	101.08	110.70
1	A	8	ARG	N-CA-C	-5.22	105.67	111.36
1	C	32	LEU	N-CA-CB	5.22	117.67	109.69
1	C	65	LYS	CB-CA-C	-5.21	102.15	110.79
1	A	43	LYS	N-CA-CB	5.19	118.21	110.22
1	B	115	THR	CB-CA-C	5.18	119.09	110.81
1	A	40	ASN	N-CA-CB	5.17	117.45	109.91
1	B	47	ASP	CA-C-N	5.15	127.64	120.63
1	B	47	ASP	C-N-CA	5.15	127.64	120.63
1	D	76	ARG	CB-CA-C	5.14	119.32	110.79
1	D	110	GLY	N-CA-C	-5.13	106.28	112.49
1	A	158	TRP	N-CA-C	-5.12	106.88	112.72
1	D	115	THR	N-CA-C	5.10	121.67	110.80
1	B	94	VAL	N-CA-C	5.09	115.71	110.36
1	C	124	LYS	N-CA-CB	5.08	119.08	110.49
1	A	125	ARG	CB-CA-C	-5.08	104.56	112.07
1	C	159	ASP	CA-C-N	-5.07	112.49	120.60
1	C	159	ASP	C-N-CA	-5.07	112.49	120.60
1	C	118	LEU	N-CA-C	-5.06	105.94	111.82
1	A	30	GLY	CA-C-N	-5.05	114.40	121.72
1	A	30	GLY	C-N-CA	-5.05	114.40	121.72
1	C	8	ARG	N-CA-CB	5.04	117.45	109.94
1	A	128	GLU	N-CA-C	-5.01	105.71	111.07
1	A	159	ASP	CA-CB-CG	-5.00	107.60	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	1303	0	1325	91	0
1	B	1284	0	1302	110	0
1	C	1288	0	1308	47	1
1	D	1288	0	1308	92	1
2	A	58	0	0	6	0
2	B	21	0	0	2	0
2	C	49	0	0	0	0
2	D	36	0	0	2	0
All	All	5327	0	5243	334	1

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 (334) 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:1:MET:HG2	1:A:6:ILE:HD11	1.37	1.05
1:C:75:VAL:HG22	1:C:100:ILE:HD11	1.40	1.03
1:A:125:ARG:HB3	1:A:128:GLU:HG3	1.38	1.01
1:D:155:THR:HG22	1:D:157:THR:HG23	1.43	0.98
1:A:139:TYR:HE1	1:A:147:LYS:HG2	1.33	0.93
1:D:24:TYR:CE1	1:D:35:LYS:HG2	2.04	0.92
1:D:134:ALA:HA	1:D:139:TYR:CD2	2.05	0.92
1:D:17:ILE:HD13	1:D:33:LEU:HD23	1.53	0.91
1:A:116:ASN:H	1:A:116:ASN:ND2	1.51	0.90
1:D:60:LYS:H	1:D:60:LYS:HD2	1.37	0.88
1:B:2:ASN:HB3	1:B:5:GLU:CG	2.06	0.86
1:A:116:ASN:H	1:A:116:ASN:HD22	1.24	0.85
1:D:120:MET:CE	1:D:128:GLU:HB3	2.06	0.85
1:C:125:ARG:HG2	1:C:128:GLU:CD	2.02	0.85
1:B:148:ARG:HD2	1:B:161:TYR:CE2	2.12	0.85
1:D:88:TYR:CZ	1:D:96:ARG:HD3	2.12	0.84
1:B:2:ASN:HB3	1:B:5:GLU:HG3	1.62	0.82
1:D:16:LYS:HG3	1:D:57:VAL:HG22	1.60	0.81
1:D:52:ARG:HH21	1:D:54:CYS:HB3	1.46	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:16:LYS:HG3	1:D:57:VAL:CG2	2.11	0.80
1:A:139:TYR:CE1	1:A:147:LYS:HG2	2.15	0.80
1:B:50:ILE:CG2	1:B:52:ARG:HG2	2.12	0.80
1:D:120:MET:HE3	1:D:128:GLU:HB3	1.63	0.80
1:D:52:ARG:HH21	1:D:54:CYS:CB	1.94	0.79
1:A:116:ASN:ND2	1:A:116:ASN:N	2.29	0.79
1:A:1:MET:HA	1:A:5:GLU:OE1	1.84	0.78
1:A:1:MET:HG2	1:A:6:ILE:CD1	2.15	0.76
1:B:50:ILE:HG22	1:B:52:ARG:HG2	1.67	0.76
1:A:24:TYR:CZ	1:A:35:LYS:HD3	2.22	0.75
1:B:3:ILE:HG23	1:B:67:PHE:HE2	1.52	0.75
1:D:24:TYR:CD1	1:D:35:LYS:HG2	2.22	0.74
1:B:137:ARG:HA	1:B:140:ASN:HB2	1.69	0.74
1:B:20:ASP:HB3	1:B:24:TYR:H	1.51	0.74
1:D:60:LYS:H	1:D:60:LYS:CD	2.01	0.74
1:D:88:TYR:CE2	1:D:96:ARG:HD3	2.23	0.73
1:B:20:ASP:CB	1:B:24:TYR:H	2.01	0.73
1:B:124:LYS:HB3	1:B:126:TRP:CZ2	2.23	0.73
1:A:116:ASN:HD22	1:A:116:ASN:N	1.86	0.72
1:A:2:ASN:H	1:A:5:GLU:CD	1.98	0.71
1:D:116:ASN:O	1:D:120:MET:HG2	1.91	0.71
1:D:60:LYS:O	1:D:63:ALA:HB3	1.90	0.71
1:C:52:ARG:HG3	1:C:53:ASN:N	2.06	0.70
1:A:79:LEU:HD23	1:A:85:LYS:HD2	1.73	0.70
1:C:19:LYS:HE3	1:C:25:TYR:CZ	2.27	0.70
1:D:34:THR:CG2	1:D:42:ALA:HB2	2.22	0.69
1:B:87:VAL:HG12	1:B:91:LEU:HD11	1.72	0.69
1:B:20:ASP:HB3	1:B:23:GLY:H	1.57	0.69
1:B:45:GLU:HG2	2:B:170:HOH:O	1.92	0.68
1:D:34:THR:HG22	1:D:42:ALA:HB2	1.74	0.68
1:D:16:LYS:HG2	1:D:56:GLY:O	1.93	0.68
1:D:132:ASN:O	1:D:135:LYS:HG2	1.93	0.68
1:C:26:THR:HG23	1:C:32:LEU:HA	1.76	0.67
1:B:125:ARG:HD2	1:B:128:GLU:OE1	1.95	0.67
1:A:125:ARG:HE	1:A:128:GLU:CD	2.03	0.67
1:B:75:VAL:HG11	1:D:37:PRO:HG3	1.77	0.66
1:B:146:ALA:O	1:B:150:ILE:HG13	1.95	0.66
1:A:2:ASN:N	1:A:5:GLU:OE2	2.29	0.66
1:D:116:ASN:HB3	1:D:119:ARG:HH21	1.60	0.66
1:C:5:GLU:OE2	1:C:8:ARG:NH2	2.30	0.65
1:B:113:GLY:O	1:B:115:THR:HG23	1.97	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:117:SER:HB2	1:B:133:LEU:HD21	1.78	0.65
1:A:91:LEU:HD22	1:A:95:ARG:HB3	1.78	0.65
1:D:114:PHE:O	1:D:118:LEU:HD22	1.97	0.65
1:A:116:ASN:O	1:A:119:ARG:HB2	1.97	0.64
1:A:127:ASP:OD1	1:A:154:ARG:HG3	1.97	0.64
1:A:146:ALA:O	1:A:150:ILE:HD12	1.99	0.63
1:A:151:THR:O	1:A:155:THR:HG23	1.98	0.63
1:B:2:ASN:HD22	1:B:5:GLU:HG2	1.63	0.62
1:D:59:THR:N	1:D:62:GLU:OE1	2.29	0.62
1:B:50:ILE:HG21	1:B:54:CYS:SG	2.40	0.62
1:B:3:ILE:HG23	1:B:67:PHE:CE2	2.35	0.62
1:B:24:TYR:CD1	1:B:35:LYS:HB3	2.34	0.62
1:D:85:LYS:HG3	1:D:85:LYS:O	1.93	0.62
1:C:125:ARG:HG2	1:C:128:GLU:OE2	2.00	0.61
1:A:21:THR:HG22	1:C:1:MET:O	2.01	0.61
1:C:143:PRO:O	1:C:147:LYS:HG2	2.00	0.61
1:B:6:ILE:HD13	1:B:158:TRP:CD2	2.36	0.61
1:B:78:ILE:HD12	1:B:100:ILE:CD1	2.31	0.61
1:B:137:ARG:HA	1:B:140:ASN:CB	2.30	0.61
1:C:52:ARG:HG2	1:C:54:CYS:SG	2.41	0.60
1:D:54:CYS:HB3	1:D:57:VAL:O	2.00	0.60
1:B:117:SER:CB	1:B:133:LEU:HD21	2.32	0.60
1:B:124:LYS:HD2	1:B:126:TRP:CZ2	2.36	0.60
1:B:7:LEU:O	1:B:11:GLU:HB2	2.02	0.59
1:A:54:CYS:HB3	1:A:57:VAL:O	2.01	0.59
1:B:147:LYS:HB3	2:B:183:HOH:O	2.01	0.59
1:B:9:ILE:HG22	1:B:10:ASP:N	2.16	0.59
1:C:75:VAL:CG2	1:C:100:ILE:HD11	2.23	0.59
1:D:120:MET:HB2	1:D:129:ALA:HB2	1.84	0.59
1:A:83:LYS:HD3	1:A:112:ALA:HB1	1.85	0.59
1:D:105:GLN:HB2	1:D:145:ARG:NH2	2.18	0.59
1:B:137:ARG:HH11	1:B:137:ARG:CG	2.17	0.58
1:D:26:THR:HG22	1:D:27:ILE:N	2.17	0.57
1:D:155:THR:CG2	1:D:157:THR:HG23	2.26	0.57
1:A:13:LEU:HD12	1:A:14:ARG:N	2.19	0.57
1:A:13:LEU:C	1:A:14:ARG:HG2	2.29	0.57
1:B:52:ARG:NH1	1:B:54:CYS:HA	2.20	0.57
1:C:85:LYS:HB3	1:C:86:PRO:HD3	1.86	0.57
1:A:4:PHE:CE2	1:A:64:GLU:HG3	2.40	0.57
1:B:154:ARG:NH1	1:B:154:ARG:HG2	2.20	0.56
1:D:17:ILE:HD13	1:D:33:LEU:CD2	2.33	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1:MET:CE	1:A:161:TYR:HB3	2.36	0.56
1:C:1:MET:HE2	1:C:5:GLU:HB3	1.86	0.56
1:A:106:MET:HE1	1:A:138:TRP:CD1	2.41	0.56
1:B:6:ILE:HD13	1:B:158:TRP:CE2	2.41	0.56
1:B:91:LEU:HB2	1:B:96:ARG:HG2	1.88	0.56
1:B:154:ARG:HG2	1:B:154:ARG:HH11	1.71	0.56
1:A:52:ARG:HG2	1:A:54:CYS:SG	2.46	0.56
1:A:79:LEU:CD2	1:A:85:LYS:HD2	2.35	0.56
1:B:3:ILE:O	1:B:3:ILE:HD12	2.06	0.55
1:B:78:ILE:HD12	1:B:100:ILE:HD11	1.89	0.55
1:B:137:ARG:O	1:B:141:GLN:N	2.33	0.55
1:B:24:TYR:CE1	1:B:35:LYS:HE2	2.41	0.55
1:B:22:GLU:HB2	1:B:24:TYR:CE2	2.42	0.55
1:D:50:ILE:HG22	1:D:52:ARG:HB2	1.89	0.55
1:B:148:ARG:HD2	1:B:161:TYR:CD2	2.42	0.54
1:B:50:ILE:HG21	1:B:52:ARG:HG2	1.86	0.54
1:D:11:GLU:OE1	1:D:145:ARG:NH1	2.41	0.54
1:B:137:ARG:HH11	1:B:137:ARG:HG3	1.73	0.54
1:B:20:ASP:HB3	1:B:23:GLY:N	2.23	0.53
1:C:24:TYR:CZ	1:C:35:LYS:HD3	2.43	0.53
1:D:155:THR:HG21	1:D:157:THR:OG1	2.08	0.53
1:B:137:ARG:HG3	1:B:137:ARG:NH1	2.23	0.53
1:D:134:ALA:HA	1:D:139:TYR:CE2	2.43	0.53
1:D:144:ASN:O	1:D:147:LYS:HB2	2.08	0.53
1:A:123:GLN:HE22	1:A:125:ARG:NH1	2.07	0.53
1:C:125:ARG:HB3	1:C:128:GLU:HB2	1.91	0.53
1:B:47:ASP:OD1	1:B:54:CYS:N	2.37	0.52
1:A:115:THR:HG22	1:A:119:ARG:NH2	2.25	0.52
1:D:2:ASN:O	1:D:5:GLU:HB2	2.10	0.52
1:D:6:ILE:HD13	1:D:158:TRP:CD2	2.45	0.52
1:A:134:ALA:C	1:A:136:SER:H	2.18	0.52
1:B:96:ARG:O	1:B:100:ILE:HG12	2.09	0.52
1:C:139:TYR:OH	1:C:147:LYS:HD2	2.10	0.52
1:B:94:VAL:HG22	1:B:158:TRP:NE1	2.25	0.52
1:B:25:TYR:O	1:B:33:LEU:HB2	2.10	0.52
1:A:24:TYR:CE1	1:A:35:LYS:HD3	2.45	0.52
1:B:125:ARG:HB3	1:B:128:GLU:HB2	1.91	0.51
1:C:143:PRO:O	1:C:147:LYS:N	2.43	0.51
1:B:24:TYR:CE1	1:B:35:LYS:HB3	2.45	0.51
1:D:61:ASP:O	1:D:65:LYS:HD3	2.11	0.51
1:D:159:ASP:OD1	1:D:159:ASP:N	2.43	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:3:ILE:HB	1:D:20:ASP:O	2.10	0.51
1:B:3:ILE:HG21	1:D:19:LYS:HB2	1.92	0.51
1:A:148:ARG:HG2	1:A:160:ALA:HB1	1.93	0.51
1:C:123:GLN:OE1	1:C:125:ARG:NH1	2.42	0.51
1:A:116:ASN:ND2	2:A:194:HOH:O	2.42	0.51
1:B:124:LYS:HD2	1:B:126:TRP:CH2	2.45	0.51
1:A:85:LYS:HB3	1:A:86:PRO:HD3	1.92	0.51
1:A:130:ALA:HB3	2:A:198:HOH:O	2.09	0.51
1:A:139:TYR:OH	1:A:147:LYS:HE3	2.11	0.51
1:B:14:ARG:NH1	1:B:18:TYR:CD2	2.79	0.51
1:C:125:ARG:HG2	1:C:128:GLU:OE1	2.11	0.50
1:D:114:PHE:C	1:D:118:LEU:HD22	2.36	0.50
1:A:59:THR:OG1	1:A:62:GLU:OE2	2.29	0.50
1:D:106:MET:HG3	1:D:110:GLY:HA3	1.92	0.50
1:D:52:ARG:HH22	1:D:58:ILE:HA	1.76	0.50
1:A:139:TYR:CE1	1:A:147:LYS:HE2	2.46	0.50
1:A:160:ALA:HB3	2:A:216:HOH:O	2.11	0.50
1:A:1:MET:HE2	1:A:161:TYR:CB	2.42	0.50
1:B:133:LEU:O	1:B:136:SER:OG	2.30	0.50
1:A:6:ILE:HG23	1:A:161:TYR:CE2	2.47	0.50
1:A:58:ILE:HB	1:A:62:GLU:HG3	1.94	0.49
1:A:15:LEU:HB3	1:A:58:ILE:O	2.12	0.49
1:B:118:LEU:O	1:B:122:GLN:N	2.42	0.49
1:B:150:ILE:O	1:B:153:PHE:HB2	2.12	0.49
1:D:46:LEU:HD13	1:D:54:CYS:HB2	1.94	0.49
1:A:141:GLN:HG2	2:A:202:HOH:O	2.12	0.49
1:B:124:LYS:HA	1:B:126:TRP:CH2	2.47	0.49
1:B:18:TYR:CE1	1:B:26:THR:HG22	2.48	0.49
1:A:59:THR:H	1:A:62:GLU:HG3	1.78	0.49
1:B:20:ASP:HB2	1:B:24:TYR:H	1.76	0.48
1:A:116:ASN:HA	1:A:119:ARG:NH1	2.28	0.48
1:A:11:GLU:CD	1:A:145:ARG:HH22	2.20	0.48
1:A:123:GLN:NE2	1:A:125:ARG:HH11	2.10	0.48
1:B:137:ARG:CA	1:B:140:ASN:HB2	2.42	0.48
1:A:1:MET:HE2	1:A:161:TYR:HB3	1.94	0.48
1:A:26:THR:HG22	1:A:27:ILE:N	2.28	0.48
1:B:13:LEU:HD21	1:B:60:LYS:HG2	1.95	0.48
1:C:85:LYS:HB3	1:C:86:PRO:CD	2.43	0.48
1:D:149:VAL:HG21	2:D:186:HOH:O	2.14	0.48
1:A:26:THR:HG23	1:A:31:HIS:O	2.14	0.47
1:A:124:LYS:HE2	2:A:188:HOH:O	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:5:GLU:HG3	1:C:8:ARG:HH21	1.79	0.47
1:D:95:ARG:HD3	1:D:153:PHE:O	2.13	0.47
1:B:78:ILE:CD1	1:B:100:ILE:HD13	2.45	0.47
1:B:2:ASN:HB3	1:B:5:GLU:HG2	1.95	0.47
1:C:54:CYS:O	1:C:55:ASN:C	2.57	0.47
1:A:139:TYR:CE1	1:A:147:LYS:CE	2.98	0.47
1:C:106:MET:CE	1:C:138:TRP:CD1	2.98	0.47
1:A:144:ASN:HB2	2:A:203:HOH:O	2.14	0.47
1:D:34:THR:HG22	1:D:42:ALA:CB	2.41	0.47
1:D:52:ARG:NH2	1:D:54:CYS:HB3	2.23	0.47
1:B:157:THR:C	1:B:159:ASP:H	2.21	0.47
1:C:119:ARG:HD3	1:C:123:GLN:HE22	1.80	0.47
1:D:85:LYS:N	1:D:86:PRO:HD2	2.30	0.47
1:A:4:PHE:HD2	1:A:67:PHE:CD2	2.33	0.46
1:D:133:LEU:HB3	1:D:150:ILE:HD11	1.97	0.46
1:B:114:PHE:O	1:B:118:LEU:HD22	2.16	0.46
1:D:52:ARG:NH2	1:D:54:CYS:SG	2.88	0.46
1:D:79:LEU:HD12	1:D:79:LEU:HA	1.49	0.46
1:D:111:VAL:O	1:D:118:LEU:HD21	2.16	0.46
1:B:136:SER:O	1:B:140:ASN:HB2	2.16	0.46
1:D:24:TYR:CB	1:D:32:LEU:HD11	2.46	0.46
1:A:146:ALA:C	1:A:150:ILE:HD12	2.41	0.46
1:B:121:LEU:HD23	1:B:121:LEU:HA	1.68	0.46
1:A:45:GLU:OE1	1:A:45:GLU:HA	2.16	0.46
1:B:22:GLU:CB	1:B:24:TYR:CE2	2.98	0.46
1:C:59:THR:N	1:C:62:GLU:OE1	2.45	0.46
1:D:50:ILE:CG2	1:D:52:ARG:HB2	2.46	0.46
1:B:18:TYR:CZ	1:B:26:THR:CG2	2.99	0.46
1:C:24:TYR:HB3	1:C:32:LEU:CD1	2.46	0.46
1:D:99:LEU:HA	1:D:99:LEU:HD12	1.60	0.45
1:B:60:LYS:O	1:B:63:ALA:HB3	2.16	0.45
1:D:124:LYS:HG2	1:D:126:TRP:CZ2	2.52	0.45
1:A:52:ARG:CG	1:A:53:ASN:N	2.79	0.45
1:A:115:THR:CG2	1:A:119:ARG:NH2	2.79	0.45
1:D:26:THR:CG2	1:D:27:ILE:N	2.79	0.45
1:A:88:TYR:CD1	1:A:96:ARG:HG2	2.52	0.45
1:B:24:TYR:CE1	1:B:35:LYS:CE	2.99	0.45
1:C:1:MET:CE	1:C:5:GLU:HB3	2.46	0.45
1:D:85:LYS:N	1:D:86:PRO:CD	2.79	0.45
1:A:16:LYS:HB3	1:A:16:LYS:HE2	1.51	0.45
1:B:22:GLU:HB2	1:B:24:TYR:CD2	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:124:LYS:CD	1:B:126:TRP:CH2	2.99	0.45
1:A:1:MET:HE1	1:A:161:TYR:HB3	1.98	0.45
1:A:39:LEU:O	1:A:43:LYS:HG3	2.17	0.45
1:A:125:ARG:NH2	1:A:128:GLU:OE1	2.33	0.45
1:C:14:ARG:NH2	1:C:18:TYR:HB3	2.32	0.45
1:C:61:ASP:O	1:C:62:GLU:C	2.56	0.45
1:D:52:ARG:NE	1:D:54:CYS:SG	2.89	0.45
1:A:85:LYS:N	1:A:86:PRO:HD2	2.31	0.45
1:B:20:ASP:HB2	1:B:24:TYR:O	2.17	0.45
1:C:6:ILE:CD1	1:C:158:TRP:CD2	2.99	0.45
1:D:19:LYS:HE2	1:D:25:TYR:CZ	2.52	0.45
1:D:131:VAL:CG2	1:D:132:ASN:N	2.79	0.45
1:A:81:ASN:OD1	1:A:83:LYS:N	2.48	0.45
1:B:6:ILE:CD1	1:B:158:TRP:CD2	3.00	0.45
1:B:125:ARG:HD2	1:B:128:GLU:CD	2.42	0.45
1:D:145:ARG:HH11	1:D:145:ARG:HD3	1.61	0.45
1:B:135:LYS:HB2	1:B:135:LYS:HE2	1.58	0.44
1:B:137:ARG:NH2	1:B:140:ASN:OD1	2.50	0.44
1:D:59:THR:OG1	1:D:62:GLU:OE1	2.29	0.44
1:B:16:LYS:CB	1:B:16:LYS:NZ	2.80	0.44
1:D:84:LEU:C	1:D:86:PRO:HD2	2.43	0.44
1:C:29:ILE:O	1:C:29:ILE:HG22	2.17	0.44
1:C:48:LYS:HB2	1:C:48:LYS:HE2	1.35	0.44
1:D:46:LEU:HD23	1:D:46:LEU:HA	1.64	0.44
1:D:95:ARG:HG2	1:D:153:PHE:HA	1.99	0.44
1:B:81:ASN:HB3	1:B:84:LEU:HB2	1.98	0.44
1:C:6:ILE:HD11	1:C:158:TRP:CD2	2.52	0.44
1:D:27:ILE:CD1	1:D:58:ILE:HD13	2.48	0.44
1:A:13:LEU:O	1:A:14:ARG:HG2	2.17	0.44
1:A:62:GLU:H	1:A:62:GLU:HG2	1.53	0.44
1:A:132:ASN:O	1:A:133:LEU:C	2.60	0.43
1:B:106:MET:HE2	1:B:138:TRP:CD1	2.53	0.43
1:A:102:MET:HE3	1:A:138:TRP:CH2	2.52	0.43
1:D:133:LEU:HD23	1:D:133:LEU:HA	1.89	0.43
1:B:148:ARG:HD2	1:B:161:TYR:CZ	2.51	0.43
1:B:3:ILE:CG2	1:D:19:LYS:HB2	2.48	0.43
1:C:118:LEU:HD12	1:C:121:LEU:HD12	1.99	0.43
1:C:144:ASN:O	1:C:148:ARG:HG3	2.19	0.43
1:D:1:MET:HE2	1:D:161:TYR:CB	2.48	0.43
1:D:43:LYS:NZ	1:D:55:ASN:ND2	2.66	0.43
1:A:2:ASN:OD1	1:A:4:PHE:N	2.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:123:GLN:O	1:A:125:ARG:N	2.51	0.43
1:A:158:TRP:CE3	1:A:158:TRP:HA	2.53	0.43
1:B:35:LYS:HB3	1:B:35:LYS:HE2	1.62	0.43
1:C:79:LEU:HD21	1:C:88:TYR:CD2	2.53	0.43
1:D:6:ILE:HD13	1:D:158:TRP:CE3	2.53	0.43
1:D:11:GLU:OE1	1:D:11:GLU:HA	2.19	0.43
1:A:107:GLY:O	1:A:111:VAL:HG23	2.19	0.43
1:B:120:MET:HG2	1:B:125:ARG:HH11	1.84	0.43
1:B:18:TYR:CZ	1:B:26:THR:HG22	2.53	0.42
1:B:20:ASP:OD2	1:B:24:TYR:HD2	2.02	0.42
1:B:52:ARG:O	1:B:52:ARG:HG3	2.14	0.42
1:D:143:PRO:O	1:D:147:LYS:HG3	2.19	0.42
1:B:130:ALA:O	1:B:134:ALA:HB2	2.19	0.42
1:C:24:TYR:CD1	1:C:35:LYS:HA	2.54	0.42
1:A:88:TYR:CE1	1:A:96:ARG:HG2	2.55	0.42
1:C:6:ILE:HD13	1:C:6:ILE:HG21	1.75	0.42
1:D:81:ASN:HB3	1:D:84:LEU:HB2	2.00	0.42
1:A:11:GLU:CD	1:A:145:ARG:NH2	2.78	0.42
1:D:24:TYR:CZ	1:D:35:LYS:HG2	2.53	0.42
1:A:18:TYR:O	1:A:25:TYR:HA	2.20	0.42
1:B:141:GLN:C	1:B:143:PRO:HD3	2.45	0.42
1:C:118:LEU:HD12	1:C:118:LEU:HA	1.82	0.42
1:C:159:ASP:O	1:C:160:ALA:C	2.60	0.42
1:B:3:ILE:HG21	1:D:19:LYS:CB	2.49	0.42
1:B:20:ASP:N	1:B:24:TYR:O	2.52	0.42
1:B:99:LEU:HD23	1:B:99:LEU:HA	1.58	0.42
1:C:6:ILE:HD13	1:C:158:TRP:CE3	2.54	0.42
1:D:144:ASN:HB2	2:D:190:HOH:O	2.20	0.42
1:C:106:MET:HE2	1:C:138:TRP:CD1	2.55	0.42
1:D:52:ARG:HG2	1:D:53:ASN:H	1.83	0.42
1:A:17:ILE:HG22	1:A:25:TYR:HD1	1.85	0.42
1:C:52:ARG:NE	1:C:54:CYS:SG	2.83	0.42
1:A:1:MET:HE2	1:A:6:ILE:HD13	2.01	0.41
1:D:1:MET:HE2	1:D:161:TYR:HB2	2.02	0.41
1:D:121:LEU:HD12	1:D:121:LEU:HA	1.86	0.41
1:A:79:LEU:HD23	1:A:79:LEU:HA	1.87	0.41
1:A:106:MET:CE	1:A:138:TRP:CD1	3.02	0.41
1:B:120:MET:HA	1:B:125:ARG:NH1	2.35	0.41
1:A:41:ALA:O	1:A:42:ALA:C	2.61	0.41
1:B:124:LYS:CB	1:B:126:TRP:CZ2	3.00	0.41
1:D:33:LEU:HD12	1:D:33:LEU:HA	1.77	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:17:ILE:CG2	1:A:25:TYR:CD1	3.03	0.41
1:A:34:THR:OG1	1:A:35:LYS:N	2.53	0.41
1:A:77:GLY:O	1:A:80:ARG:HG2	2.20	0.41
1:B:14:ARG:O	1:B:28:GLY:N	2.53	0.41
1:D:16:LYS:HG3	1:D:57:VAL:HG23	1.96	0.41
1:D:134:ALA:CB	1:D:139:TYR:CE2	3.04	0.41
1:B:3:ILE:HD12	1:B:3:ILE:C	2.43	0.41
1:C:116:ASN:O	1:C:119:ARG:HB3	2.21	0.41
1:A:17:ILE:HG12	1:A:33:LEU:CD1	2.51	0.41
1:B:42:ALA:O	1:B:43:LYS:C	2.62	0.41
1:B:133:LEU:H	1:B:133:LEU:HG	1.57	0.41
1:C:119:ARG:CG	1:C:123:GLN:NE2	2.84	0.41
1:D:3:ILE:HD13	1:D:3:ILE:HG21	1.71	0.41
1:D:34:THR:HG22	1:D:42:ALA:HA	2.03	0.41
1:D:133:LEU:HB3	1:D:150:ILE:CD1	2.51	0.41
1:B:8:ARG:HH11	1:B:8:ARG:HG2	1.85	0.41
1:B:10:ASP:OD2	1:B:101:ASN:ND2	2.46	0.40
1:B:119:ARG:O	1:B:119:ARG:HG3	2.11	0.40
1:A:13:LEU:HD12	1:A:13:LEU:C	2.47	0.40
1:D:66:LEU:O	1:D:67:PHE:C	2.63	0.40
1:B:4:PHE:O	1:B:8:ARG:N	2.49	0.40
1:B:67:PHE:O	1:B:68:ASN:C	2.63	0.40
1:B:78:ILE:CD1	1:B:100:ILE:CD1	2.99	0.40
1:A:122:GLN:C	1:A:124:LYS:H	2.29	0.40
1:B:2:ASN:ND2	1:B:5:GLU:HG2	2.34	0.40
1:B:60:LYS:HB3	1:B:60:LYS:HE3	1.33	0.40
1:C:1:MET:HB2	1:C:6:ILE:HG13	2.04	0.40
1:C:106:MET:HE1	1:C:138:TRP:CD1	2.56	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:19:LYS:NZ	1:D:72:ASP:OD1[3_645]	2.06	0.14

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%)	152 (94%)	8 (5%)	2 (1%)	10	8
1	B	160/164 (98%)	140 (88%)	16 (10%)	4 (2%)	4	2
1	C	160/164 (98%)	148 (92%)	11 (7%)	1 (1%)	21	23
1	D	160/164 (98%)	146 (91%)	13 (8%)	1 (1%)	21	23
All	All	642/656 (98%)	586 (91%)	48 (8%)	8 (1%)	10	8

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	124	LYS
1	B	53	ASN
1	A	135	LYS
1	C	124	LYS
1	B	132	ASN
1	B	9	ILE
1	D	20	ASP
1	B	17	ILE

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	137/138 (99%)	108 (79%)	29 (21%)	1	1
1	B	134/138 (97%)	101 (75%)	33 (25%)	1	0

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	135/138 (98%)	114 (84%)	21 (16%)	2	2
1	D	135/138 (98%)	103 (76%)	32 (24%)	1	0
All	All	541/552 (98%)	426 (79%)	115 (21%)	1	1

All (115) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	3	ILE
1	A	5	GLU
1	A	8	ARG
1	A	13	LEU
1	A	15	LEU
1	A	16	LYS
1	A	21	THR
1	A	32	LEU
1	A	38	SER
1	A	39	LEU
1	A	53	ASN
1	A	54	CYS
1	A	59	THR
1	A	60	LYS
1	A	61	ASP
1	A	62	GLU
1	A	76	ARG
1	A	80	ARG
1	A	85	LYS
1	A	108	GLU
1	A	109	THR
1	A	116	ASN
1	A	124	LYS
1	A	135	LYS
1	A	141	GLN
1	A	145	ARG
1	A	150	ILE
1	A	155	THR
1	B	3	ILE
1	B	7	LEU
1	B	8	ARG
1	B	16	LYS
1	B	17	ILE

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Mol	Chain	Res	Type
1	B	19	LYS
1	B	35	LYS
1	B	44	SER
1	B	45	GLU
1	B	48	LYS
1	B	50	ILE
1	B	58	ILE
1	B	60	LYS
1	B	61	ASP
1	B	84	LEU
1	B	90	SER
1	B	92	ASP
1	B	99	LEU
1	B	104	PHE
1	B	106	MET
1	B	108	GLU
1	B	115	THR
1	B	117	SER
1	B	118	LEU
1	B	119	ARG
1	B	126	TRP
1	B	132	ASN
1	B	133	LEU
1	B	135	LYS
1	B	136	SER
1	B	137	ARG
1	B	140	ASN
1	B	147	LYS
1	C	1	MET
1	C	5	GLU
1	C	7	LEU
1	C	14	ARG
1	C	26	THR
1	C	32	LEU
1	C	38	SER
1	C	44	SER
1	C	48	LYS
1	C	52	ARG
1	C	60	LYS
1	C	78	ILE
1	C	83	LYS
1	C	90	SER

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Mol	Chain	Res	Type
1	C	100	ILE
1	C	106	MET
1	C	109	THR
1	C	118	LEU
1	C	125	ARG
1	C	135	LYS
1	C	147	LYS
1	D	3	ILE
1	D	5	GLU
1	D	9	ILE
1	D	11	GLU
1	D	14	ARG
1	D	16	LYS
1	D	17	ILE
1	D	29	ILE
1	D	32	LEU
1	D	33	LEU
1	D	43	LYS
1	D	46	LEU
1	D	50	ILE
1	D	55	ASN
1	D	58	ILE
1	D	65	LYS
1	D	69	GLN
1	D	78	ILE
1	D	79	LEU
1	D	80	ARG
1	D	83	LYS
1	D	100	ILE
1	D	103	VAL
1	D	104	PHE
1	D	106	MET
1	D	109	THR
1	D	115	THR
1	D	118	LEU
1	D	121	LEU
1	D	128	GLU
1	D	131	VAL
1	D	154	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	68	ASN
1	A	69	GLN
1	A	116	ASN
1	A	123	GLN
1	A	141	GLN
1	A	144	ASN
1	A	163	ASN
1	B	2	ASN
1	B	68	ASN
1	B	132	ASN
1	B	144	ASN
1	C	53	ASN
1	C	140	ASN
1	C	141	GLN
1	C	144	ASN
1	D	55	ASN
1	D	69	GLN
1	D	132	ASN
1	D	141	GLN
1	D	144	ASN

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 ⓘ

There are no ligands in this entry.

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 [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%)	-0.35	1 (0%) 85 83	12, 29, 52, 59	0
1	B	162/164 (98%)	-0.45	0 100 100	18, 37, 58, 71	0
1	C	162/164 (98%)	-0.55	0 100 100	15, 27, 49, 63	0
1	D	162/164 (98%)	-0.56	0 100 100	13, 35, 55, 65	0
All	All	650/656 (99%)	-0.48	1 (0%) 91 90	12, 33, 55, 71	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	116	ASN	2.1

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.