WORLDWIDE wwPDB PROTEIN DATA BANK

Full wwPDB X-ray Structure Validation Report (i)

Jan 31, 2016 – 09:49 PM GMT

PDB ID	:	1QTH
Title	:	THE INTRODUCTION OF STRAIN AND ITS EFFECTS ON THE STRUC-TURE AND STABILITY OF T4 LYSOZYME
Authors	:	Liu, R.; Baase, W.A.; Matthews, B.W.
Deposited on	:	1999-06-28
Resolution	:	1.90 Å(reported)

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

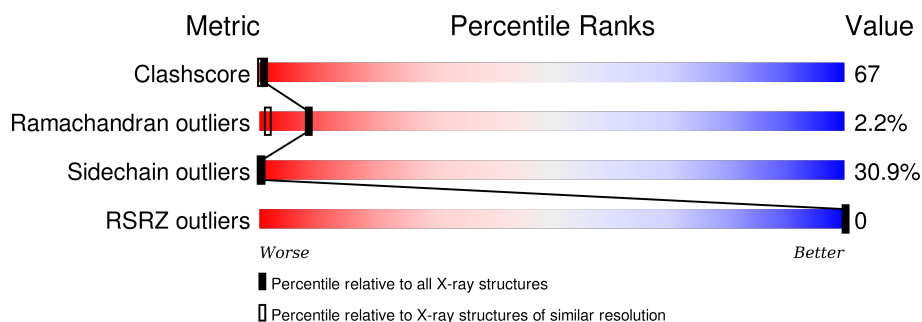|		
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
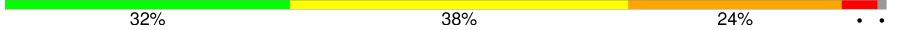*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.90 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	Percentile Ranks	Value
Clashscore		67
Ramachandran outliers		2.2%
Sidechain outliers		30.9%
RSRZ outliers		0

Worse ——— *Better*

■ Percentile relative to all X-ray structures
▯ Percentile relative to X-ray structures of similar resolution

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	102246	5398 (1.90-1.90)
Ramachandran outliers	100387	5338 (1.90-1.90)
Sidechain outliers	100360	5339 (1.90-1.90)
RSRZ outliers	91569	4766 (1.90-1.90)

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 <=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	21% 45% 26% 8% •
1	B	164	32% 38% 24% • •

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 2703 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.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	162	Total 1295	C 816	N 235	O 238	S 6	0	0	0
1	B	162	Total 1295	C 816	N 235	O 238	S 6	0	0	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	54	THR	CYS	ENGINEERED	UNP P00720
A	97	ALA	CYS	ENGINEERED	UNP P00720
A	98	MET	ALA	ENGINEERED	UNP P00720
B	54	THR	CYS	ENGINEERED	UNP P00720
B	97	ALA	CYS	ENGINEERED	UNP P00720
B	98	MET	ALA	ENGINEERED	UNP P00720
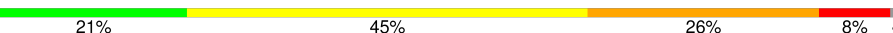
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	47	Total 47	O 47	0	0
2	B	66	Total 66	O 66	0	0

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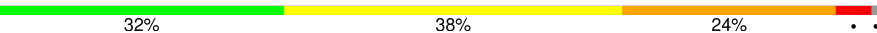
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according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

- Molecule 1: LYSOZYME

Chain A:
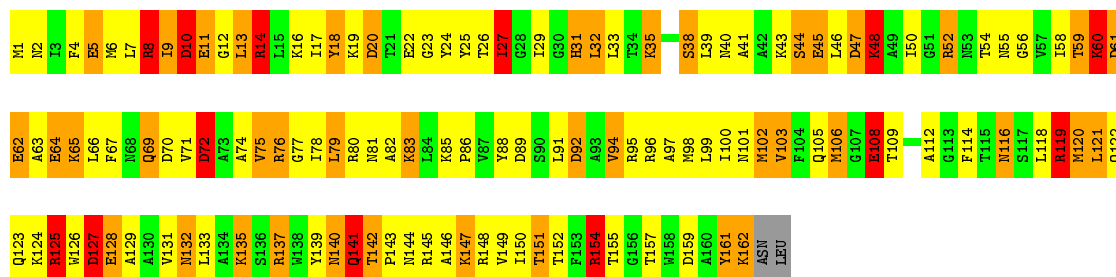
- Molecule 1: LYSOZYME

Chain B:

4   Data and refinement statistics (i)

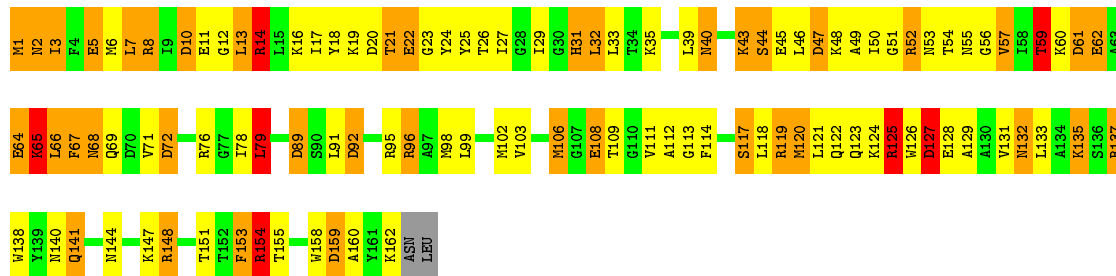| Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, α, β, γ	53.57Å　　53.57Å　　101.87Å 90.00°　　90.00°　　120.00°	Depositor
Resolution (Å)	20.00　–　1.90 46.39　–　1.76	Depositor EDS
% Data completeness (in resolution range)	85.3 (20.00-1.90) 73.7 (46.39-1.76)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.41 (at 1.76Å)	Xtriage
Refinement program	TNT	Depositor
R, R_{free}	0.189　,　(Not available) 0.203　,　(Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å^2)	25.5	Xtriage
Anisotropy	0.230	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.22 , 115.9	EDS
Estimated twinning fraction	0.106 for -h,-k,l 0.373 for h,-h-k,-l 0.107 for -k,-h,-l	Xtriage
L-test for twinning[2]	$< |L| > = 0.42, < L^2 > = 0.25$	Xtriage
Outliers	0 of 23736 reflections	Xtriage
F_o,F_c correlation	0.97	EDS
Total number of atoms	2703	wwPDB-VP
Average B, all atoms (Å^2)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.95% of the height of the origin peak. No significant pseudotranslation is detected.*

---

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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	1.09	8/1315 (0.6%)	1.68	27/1770 (1.5%)
1	B	1.12	8/1315 (0.6%)	1.76	33/1770 (1.9%)
All	All	1.11	16/2630 (0.6%)	1.72	60/3540 (1.7%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	1	0

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	45	GLU	CD-OE2	6.78	1.33	1.25
1	B	45	GLU	CD-OE2	6.25	1.32	1.25
1	A	5	GLU	CD-OE2	6.24	1.32	1.25
1	A	108	GLU	CD-OE2	6.16	1.32	1.25
1	B	128	GLU	CD-OE2	6.16	1.32	1.25
1	A	62	GLU	CD-OE2	6.10	1.32	1.25
1	A	11	GLU	CD-OE2	6.00	1.32	1.25
1	B	108	GLU	CD-OE2	5.99	1.32	1.25
1	A	128	GLU	CD-OE2	5.97	1.32	1.25
1	A	64	GLU	CD-OE2	5.93	1.32	1.25
1	B	22	GLU	CD-OE2	5.89	1.32	1.25
1	B	5	GLU	CD-OE2	5.72	1.31	1.25
1	B	11	GLU	CD-OE2	5.54	1.31	1.25
1	B	64	GLU	CD-OE2	5.22	1.31	1.25
1	B	62	GLU	CD-OE2	5.21	1.31	1.25
1	A	22	GLU	CD-OE2	5.17	1.31	1.25

All (60) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	154	ARG	NE-CZ-NH2	-10.77	114.92	120.30
1	B	154	ARG	NE-CZ-NH1	10.48	125.54	120.30
1	A	89	ASP	CB-CG-OD2	-9.22	110.00	118.30
1	B	92	ASP	CB-CG-OD1	8.91	126.32	118.30
1	B	89	ASP	CB-CG-OD1	8.80	126.22	118.30
1	B	89	ASP	CB-CG-OD2	-8.68	110.49	118.30
1	A	159	ASP	CB-CG-OD1	8.66	126.10	118.30
1	B	92	ASP	CB-CG-OD2	-8.66	110.50	118.30
1	A	119	ARG	NE-CZ-NH1	8.28	124.44	120.30
1	A	125	ARG	NE-CZ-NH2	-7.99	116.31	120.30
1	B	76	ARG	NE-CZ-NH1	7.75	124.18	120.30
1	A	127	ASP	CB-CG-OD2	-7.73	111.34	118.30
1	B	59	THR	N-CA-CB	-7.62	95.81	110.30
1	B	159	ASP	CB-CG-OD2	-7.59	111.47	118.30
1	B	127	ASP	CB-CG-OD2	-7.53	111.52	118.30
1	B	148	ARG	NE-CZ-NH2	-7.12	116.74	120.30
1	A	125	ARG	NE-CZ-NH1	7.08	123.84	120.30
1	B	14	ARG	NE-CZ-NH1	7.07	123.83	120.30
1	B	72	ASP	CB-CG-OD2	-6.71	112.26	118.30
1	A	47	ASP	CB-CG-OD2	-6.57	112.39	118.30
1	B	47	ASP	CB-CG-OD2	-6.48	112.47	118.30
1	B	72	ASP	CB-CG-OD1	6.48	124.13	118.30
1	A	92	ASP	CB-CG-OD2	-6.39	112.55	118.30
1	B	25	TYR	N-CA-CB	6.29	121.93	110.60
1	B	7	LEU	CA-CB-CG	-6.21	101.01	115.30
1	B	49	ALA	CB-CA-C	-6.13	100.91	110.10
1	B	65	LYS	N-CA-C	-6.10	94.53	111.00
1	A	127	ASP	CB-CG-OD1	6.09	123.78	118.30
1	B	159	ASP	CB-CG-OD1	6.08	123.77	118.30
1	A	154	ARG	NE-CZ-NH1	6.06	123.33	120.30
1	B	127	ASP	CB-CG-OD1	6.04	123.73	118.30
1	A	14	ARG	NE-CZ-NH1	6.01	123.31	120.30
1	B	148	ARG	NE-CZ-NH1	5.93	123.27	120.30
1	B	76	ARG	NE-CZ-NH2	-5.90	117.35	120.30
1	A	72	ASP	CB-CG-OD2	-5.83	113.05	118.30
1	A	10	ASP	CB-CG-OD2	-5.81	113.07	118.30
1	A	8	ARG	NE-CZ-NH2	-5.76	117.42	120.30
1	A	151	THR	CA-CB-CG2	-5.75	104.34	112.40
1	B	79	LEU	CB-CA-C	5.75	121.12	110.20
1	A	8	ARG	NE-CZ-NH1	5.73	123.16	120.30
1	A	89	ASP	CB-CG-OD1	5.72	123.45	118.30
1	B	125	ARG	NE-CZ-NH1	5.72	123.16	120.30
1	A	20	ASP	CB-CG-OD1	5.62	123.36	118.30

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	67	PHE	CB-CA-C	-5.61	99.18	110.40
1	B	79	LEU	CA-CB-CG	-5.61	102.40	115.30
1	A	70	ASP	CB-CG-OD2	-5.50	113.35	118.30
1	B	61	ASP	CB-CG-OD2	-5.47	113.38	118.30
1	A	31	HIS	CA-CB-CG	-5.36	104.50	113.60
1	B	31	HIS	CA-CB-CG	-5.35	104.50	113.60
1	A	48	LYS	N-CA-CB	5.33	120.19	110.60
1	B	72	ASP	CB-CA-C	5.33	121.06	110.40
1	A	18	TYR	CB-CG-CD2	-5.32	117.81	121.00
1	A	94	VAL	CA-CB-CG1	5.32	118.87	110.90
1	A	161	TYR	N-CA-CB	5.31	120.15	110.60
1	A	126	TRP	N-CA-C	5.25	125.17	111.00
1	A	47	ASP	CB-CG-OD1	5.20	122.98	118.30
1	B	53	ASN	N-CA-CB	5.12	119.82	110.60
1	A	79	LEU	CA-CB-CG	-5.11	103.55	115.30
1	B	61	ASP	CB-CG-OD1	5.07	122.86	118.30
1	B	10	ASP	CB-CG-OD1	5.04	122.84	118.30

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	141	GLN	CA

There are no planarity outliers.

5.2 Too-close contacts (i)

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	1295	0	1323	197	0
1	B	1295	0	1323	154	0
2	A	47	0	0	9	0
2	B	66	0	0	8	0
All	All	2703	0	2646	351	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 67.

All (351) 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:102:MET:HG2	1:A:106:MET:HE3	1.23	1.16
1:A:39:LEU:HD21	1:A:43:LYS:HE3	1.28	1.15
1:B:16:LYS:HG2	1:B:57:VAL:HG23	1.32	1.09
1:A:77:GLY:N	1:A:80:ARG:HH21	1.50	1.09
1:B:137:ARG:HH21	1:B:141:GLN:HB2	1.13	1.05
1:A:98:MET:HE2	1:A:149:VAL:HG13	1.31	1.05
1:A:154:ARG:HH11	1:A:154:ARG:HG2	1.19	1.05
1:B:59:THR:HG22	1:B:62:GLU:HB2	1.37	1.02
1:A:98:MET:HE1	1:A:149:VAL:HG22	1.40	1.02
1:A:46:LEU:HD23	1:A:56:GLY:HA2	1.43	0.98
1:B:14:ARG:HB2	1:B:18:TYR:CE2	1.98	0.98
1:B:154:ARG:HG2	1:B:154:ARG:HH11	1.28	0.95
1:A:137:ARG:NH2	1:A:141:GLN:HB2	1.86	0.90
1:A:88:TYR:HD1	1:A:99:LEU:HD23	1.38	0.86
1:B:119:ARG:HH11	1:B:120:MET:N	1.74	0.85
1:A:7:LEU:HD12	1:A:67:PHE:HE1	1.41	0.85
1:A:4:PHE:CE1	1:A:29:ILE:HD11	2.13	0.83
1:B:46:LEU:HD23	1:B:56:GLY:HA2	1.60	0.82
1:B:106:MET:HE1	1:B:138:TRP:CD1	2.15	0.82
1:B:89:ASP:HA	1:B:96:ARG:HH21	1.44	0.82
1:A:132:ASN:HD22	1:A:132:ASN:N	1.74	0.81
1:B:16:LYS:CG	1:B:57:VAL:HG23	2.09	0.81
1:A:29:ILE:HD13	2:A:180:HOH:O	1.81	0.80
1:A:146:ALA:O	1:A:150:ILE:HD12	1.82	0.80
1:B:24:TYR:CE2	1:B:35:LYS:HB3	2.16	0.80
1:A:6:MET:CE	1:A:101:ASN:HD22	1.94	0.79
1:B:119:ARG:NH1	1:B:120:MET:N	2.28	0.79
1:B:1:MET:HE3	1:B:5:GLU:HB2	1.62	0.79
1:B:137:ARG:NH2	1:B:141:GLN:HB2	1.94	0.78
1:A:29:ILE:HG21	2:A:180:HOH:O	1.82	0.78
1:A:7:LEU:HD12	1:A:67:PHE:CE1	2.19	0.78
1:B:151:THR:O	1:B:155:THR:HG23	1.83	0.77
1:A:77:GLY:N	1:A:80:ARG:NH2	2.29	0.77
1:A:102:MET:HG2	1:A:106:MET:CE	2.08	0.77
1:A:20:ASP:HA	2:A:375:HOH:O	1.84	0.77
1:A:76:ARG:HB3	1:A:80:ARG:CZ	2.14	0.77
1:B:59:THR:HG22	1:B:62:GLU:CB	2.15	0.77
1:B:1:MET:HE1	1:B:6:MET:HB2	1.67	0.76
1:B:6:MET:HB2	1:B:158:TRP:HZ3	1.50	0.76
1:A:137:ARG:NE	1:A:141:GLN:HG3	2.01	0.75

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:7:LEU:HB2	2:A:180:HOH:O	1.85	0.75
1:A:32:LEU:HD22	1:A:33:LEU:N	2.01	0.75
1:B:24:TYR:CD2	1:B:35:LYS:HA	2.22	0.74
1:B:16:LYS:HG2	1:B:57:VAL:CG2	2.16	0.74
1:B:119:ARG:NH1	1:B:119:ARG:HB3	2.03	0.74
1:A:98:MET:HE3	1:A:98:MET:HA	1.69	0.74
1:A:98:MET:HE2	1:A:149:VAL:CG1	2.16	0.74
1:B:19:LYS:HG3	1:B:23:GLY:O	1.88	0.73
1:B:39:LEU:HD11	1:B:43:LYS:HE3	1.71	0.73
1:B:31:HIS:CG	1:B:66:LEU:HD11	2.24	0.73
1:B:89:ASP:HA	1:B:96:ARG:NH2	2.03	0.73
1:B:65:LYS:O	1:B:65:LYS:HE3	1.89	0.72
1:B:46:LEU:CD2	1:B:56:GLY:HA2	2.18	0.72
1:A:137:ARG:HH21	1:A:141:GLN:HB2	1.54	0.72
1:A:46:LEU:CD2	1:A:56:GLY:HA2	2.17	0.72
1:B:3:ILE:HD12	1:B:6:MET:CE	2.20	0.72
1:B:89:ASP:CA	1:B:96:ARG:HH21	2.03	0.72
1:A:154:ARG:HG2	1:A:154:ARG:NH1	1.96	0.72
1:B:158:TRP:O	1:B:160:ALA:N	2.23	0.72
1:A:59:THR:O	1:A:60:LYS:C	2.27	0.72
1:B:154:ARG:HG2	1:B:154:ARG:NH1	1.99	0.71
1:A:96:ARG:O	1:A:100:ILE:HD12	1.89	0.71
1:B:2:ASN:N	1:B:2:ASN:OD1	2.23	0.71
1:A:74:ALA:O	1:A:78:ILE:HD12	1.90	0.71
1:A:108:GLU:HG2	1:A:109:THR:N	2.03	0.71
1:A:50:ILE:HG22	1:A:52:ARG:N	2.06	0.70
1:B:65:LYS:O	1:B:66:LEU:HB2	1.91	0.69
1:B:62:GLU:HA	1:B:65:LYS:CE	2.23	0.69
1:B:95:ARG:NH2	1:B:154:ARG:O	2.26	0.69
1:A:17:ILE:N	1:A:27:ILE:HD12	2.07	0.69
1:B:59:THR:CG2	1:B:62:GLU:HB2	2.21	0.69
1:B:62:GLU:HA	1:B:65:LYS:HE2	1.74	0.69
1:A:161:TYR:C	1:A:162:LYS:HE2	2.13	0.68
1:A:16:LYS:NZ	1:A:17:ILE:HD12	2.07	0.68
1:A:151:THR:O	1:A:155:THR:HG23	1.93	0.68
1:A:29:ILE:HD12	1:A:67:PHE:CG	2.28	0.68
1:A:50:ILE:CG2	1:A:52:ARG:HG3	2.24	0.68
1:B:66:LEU:N	1:B:69:GLN:OE1	2.27	0.68
1:A:148:ARG:HG3	1:A:148:ARG:HH11	1.59	0.67
1:A:7:LEU:CD1	1:A:67:PHE:HE1	2.07	0.67
1:A:98:MET:CE	1:A:149:VAL:HG13	2.19	0.67

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:6:MET:HE2	1:A:101:ASN:HD22	1.60	0.67
1:A:1:MET:HE3	2:A:226:HOH:O	1.94	0.67
1:B:66:LEU:O	1:B:69:GLN:HB2	1.95	0.66
1:A:59:THR:O	1:A:61:ASP:N	2.28	0.66
1:B:119:ARG:NH1	1:B:120:MET:HE3	2.10	0.66
1:A:88:TYR:CD1	1:A:99:LEU:HD23	2.26	0.66
1:A:144:ASN:O	1:A:148:ARG:NH1	2.29	0.66
1:A:85:LYS:N	1:A:86:PRO:HD2	2.11	0.66
1:A:46:LEU:O	1:A:50:ILE:HD13	1.95	0.66
1:B:62:GLU:O	1:B:65:LYS:NZ	2.27	0.66
1:B:127:ASP:OD1	1:B:154:ARG:HD2	1.96	0.65
1:B:14:ARG:HG2	1:B:14:ARG:HH11	1.59	0.65
1:A:38:SER:OG	1:A:41:ALA:N	2.29	0.65
1:B:92:ASP:O	1:B:96:ARG:HG3	1.96	0.65
1:A:62:GLU:O	1:A:65:LYS:NZ	2.27	0.65
1:A:132:ASN:N	1:A:132:ASN:ND2	2.44	0.65
1:B:119:ARG:HD2	1:B:123:GLN:NE2	2.12	0.65
1:A:148:ARG:NH1	1:A:148:ARG:HG3	2.10	0.65
1:A:162:LYS:HD3	2:A:192:HOH:O	1.97	0.64
1:A:98:MET:CE	1:A:149:VAL:HG22	2.24	0.64
1:A:128:GLU:O	1:A:132:ASN:ND2	2.30	0.64
1:A:137:ARG:CZ	1:A:141:GLN:HG3	2.27	0.64
1:B:55:ASN:ND2	2:B:354:HOH:O	2.29	0.64
1:B:6:MET:HB2	1:B:158:TRP:CZ3	2.30	0.64
1:B:67:PHE:CE1	1:B:71:VAL:HG21	2.33	0.64
1:B:32:LEU:HD22	1:B:33:LEU:N	2.12	0.64
1:B:40:ASN:O	1:B:44:SER:HB2	1.97	0.63
1:A:4:PHE:HE1	1:A:29:ILE:HD11	1.63	0.63
1:B:95:ARG:HG2	1:B:153:PHE:HD2	1.63	0.63
1:A:76:ARG:HA	1:A:79:LEU:HD12	1.80	0.63
1:A:7:LEU:O	1:A:11:GLU:HB2	1.99	0.63
1:B:39:LEU:O	1:B:43:LYS:HG2	1.99	0.62
1:B:129:ALA:O	1:B:133:LEU:HG	1.99	0.62
1:A:14:ARG:HB2	1:A:18:TYR:CE2	2.34	0.62
1:A:99:LEU:O	1:A:103:VAL:HG13	1.98	0.62
1:B:99:LEU:O	1:B:103:VAL:HG23	1.99	0.62
1:A:98:MET:HE1	1:A:149:VAL:CG2	2.25	0.62
1:A:98:MET:O	1:A:101:ASN:HB3	2.00	0.62
1:A:5:GLU:O	1:A:9:ILE:HG13	2.00	0.62
1:B:119:ARG:HH11	1:B:119:ARG:C	2.04	0.61
1:A:125:ARG:HG2	1:A:128:GLU:OE1	2.01	0.61

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:77:GLY:CA	1:A:80:ARG:HH21	2.13	0.60
1:B:51:GLY:HA2	2:B:351:HOH:O	2.00	0.60
1:A:116:ASN:O	1:A:120:MET:HE3	2.01	0.60
1:B:3:ILE:HA	1:B:6:MET:HE2	1.83	0.60
1:A:59:THR:H	1:A:62:GLU:CG	2.15	0.60
1:B:14:ARG:HB2	1:B:18:TYR:CD2	2.37	0.60
1:B:7:LEU:N	1:B:7:LEU:HD23	2.12	0.60
1:A:45:GLU:HA	1:A:48:LYS:HE3	1.84	0.59
1:A:77:GLY:H	1:A:80:ARG:HH21	1.45	0.59
1:A:20:ASP:O	1:A:23:GLY:N	2.32	0.59
1:A:59:THR:O	1:A:62:GLU:N	2.21	0.59
1:B:137:ARG:NE	1:B:137:ARG:O	2.36	0.59
1:B:67:PHE:CZ	1:B:71:VAL:HG21	2.38	0.59
1:A:132:ASN:HD22	1:A:132:ASN:H	1.48	0.58
1:B:24:TYR:HA	2:B:382:HOH:O	2.01	0.58
1:A:61:ASP:OD1	1:A:61:ASP:N	2.28	0.58
1:A:6:MET:HE1	1:A:101:ASN:HD22	1.66	0.58
1:B:127:ASP:O	1:B:131:VAL:HG23	2.03	0.58
1:A:65:LYS:HZ2	1:A:66:LEU:HG	1.69	0.58
1:A:8:ARG:HD3	1:A:8:ARG:C	2.23	0.58
1:A:102:MET:O	1:A:106:MET:HG3	2.04	0.58
1:B:22:GLU:N	1:B:22:GLU:OE2	2.37	0.58
1:A:39:LEU:HD21	1:A:43:LYS:CE	2.20	0.57
1:A:46:LEU:HD23	1:A:56:GLY:CA	2.27	0.57
1:A:10:ASP:N	1:A:10:ASP:OD1	2.36	0.57
1:A:24:TYR:CE2	1:A:35:LYS:HG3	2.39	0.57
1:B:65:LYS:HA	1:B:68:ASN:HB2	1.86	0.57
1:A:13:LEU:O	1:A:14:ARG:HG2	2.05	0.57
1:B:21:THR:HG23	2:B:206:HOH:O	2.04	0.57
1:A:16:LYS:HD2	1:A:17:ILE:H	1.70	0.57
1:A:16:LYS:HZ3	1:A:17:ILE:HD12	1.68	0.57
1:A:142:THR:HB	1:A:145:ARG:HB3	1.86	0.57
1:A:132:ASN:H	1:A:132:ASN:ND2	2.03	0.57
1:B:120:MET:HB3	1:B:125:ARG:HB3	1.84	0.57
1:B:98:MET:O	1:B:102:MET:HG3	2.03	0.57
1:A:92:ASP:OD2	1:A:95:ARG:HD2	2.05	0.57
1:B:27:ILE:O	1:B:31:HIS:HB3	2.04	0.56
1:B:10:ASP:OD1	1:B:148:ARG:NH2	2.29	0.56
1:A:76:ARG:O	1:A:79:LEU:N	2.38	0.56
1:A:98:MET:CE	1:A:98:MET:HA	2.33	0.56
1:B:119:ARG:HH12	1:B:120:MET:CG	2.19	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:6:MET:CB	1:B:158:TRP:HZ3	2.17	0.56
1:A:23:GLY:N	2:A:353:HOH:O	2.40	0.55
1:A:19:LYS:HD3	1:A:23:GLY:O	2.07	0.55
1:B:120:MET:CB	1:B:125:ARG:HB3	2.37	0.55
1:B:1:MET:HE1	1:B:6:MET:CB	2.36	0.54
1:A:6:MET:O	1:A:10:ASP:OD1	2.25	0.54
1:B:3:ILE:HD12	1:B:6:MET:HE3	1.87	0.54
1:B:3:ILE:O	1:B:7:LEU:HG	2.08	0.54
1:A:59:THR:H	1:A:62:GLU:HG3	1.72	0.54
1:A:63:ALA:O	1:A:66:LEU:HB2	2.08	0.54
1:A:24:TYR:CZ	1:A:35:LYS:HG3	2.43	0.54
1:A:17:ILE:CA	1:A:27:ILE:HD12	2.37	0.54
1:B:3:ILE:HA	1:B:6:MET:CE	2.37	0.54
1:A:114:PHE:O	1:A:118:LEU:HG	2.08	0.53
1:B:62:GLU:HA	1:B:65:LYS:NZ	2.24	0.53
1:B:119:ARG:HH11	1:B:119:ARG:HB3	1.74	0.53
1:B:119:ARG:HD2	1:B:123:GLN:CD	2.29	0.53
1:B:12:GLY:O	1:B:29:ILE:HA	2.09	0.52
1:A:32:LEU:HD22	1:A:33:LEU:H	1.74	0.52
1:A:75:VAL:O	1:A:79:LEU:HG	2.09	0.52
1:A:26:THR:HG22	1:A:27:ILE:N	2.25	0.52
1:A:148:ARG:CG	1:A:148:ARG:HH11	2.23	0.52
1:A:25:TYR:C	1:A:33:LEU:HD12	2.30	0.52
1:A:131:VAL:HG12	1:A:132:ASN:HD22	1.75	0.51
1:B:59:THR:HG22	1:B:62:GLU:N	2.25	0.51
1:A:98:MET:CE	1:A:149:VAL:HA	2.40	0.51
1:A:91:LEU:HD22	1:A:95:ARG:HB2	1.92	0.51
1:A:161:TYR:O	1:A:162:LYS:HE2	2.09	0.51
1:A:145:ARG:O	1:A:149:VAL:HG23	2.10	0.51
1:A:19:LYS:HB3	1:A:23:GLY:HA2	1.91	0.51
1:A:50:ILE:HD12	1:A:50:ILE:N	2.25	0.51
1:A:17:ILE:CD1	1:A:43:LYS:HG2	2.41	0.51
1:A:26:THR:HA	1:A:31:HIS:O	2.11	0.51
1:B:14:ARG:CG	1:B:14:ARG:HH11	2.24	0.51
1:B:31:HIS:ND1	1:B:66:LEU:HD11	2.25	0.51
1:A:91:LEU:HB2	1:A:96:ARG:HG2	1.93	0.51
1:A:62:GLU:O	1:A:63:ALA:C	2.48	0.50
1:A:105:GLN:HB2	1:A:145:ARG:NH2	2.26	0.50
1:B:3:ILE:HD12	1:B:6:MET:HE1	1.92	0.50
1:B:7:LEU:O	1:B:8:ARG:C	2.47	0.50
1:A:135:LYS:HE2	1:A:135:LYS:HA	1.93	0.50

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:39:LEU:CD2	1:A:43:LYS:HE3	2.19	0.50
1:B:137:ARG:CZ	1:B:137:ARG:HB3	2.41	0.50
1:A:67:PHE:CZ	1:A:71:VAL:HG21	2.46	0.50
1:A:88:TYR:CE1	1:A:96:ARG:HB3	2.46	0.50
1:B:102:MET:O	1:B:106:MET:HG2	2.11	0.50
1:B:20:ASP:OD1	1:B:22:GLU:HG2	2.12	0.50
1:A:52:ARG:O	1:A:54:THR:HG23	2.11	0.49
1:A:91:LEU:HD22	1:A:95:ARG:CB	2.42	0.49
1:B:95:ARG:HG2	1:B:153:PHE:CD2	2.43	0.49
1:A:24:TYR:HB3	1:A:32:LEU:HD21	1.94	0.49
1:A:66:LEU:O	1:A:69:GLN:N	2.46	0.49
1:B:119:ARG:NH1	1:B:120:MET:CE	2.76	0.49
1:A:16:LYS:HZ2	1:A:17:ILE:HD12	1.75	0.49
1:B:67:PHE:O	1:B:71:VAL:HG23	2.13	0.49
1:B:137:ARG:HH21	1:B:141:GLN:CB	2.04	0.49
1:B:17:ILE:HA	1:B:27:ILE:HD12	1.95	0.48
1:A:17:ILE:HA	1:A:27:ILE:HD12	1.94	0.48
1:B:13:LEU:O	1:B:14:ARG:HG2	2.12	0.48
1:A:99:LEU:O	1:A:100:ILE:C	2.48	0.48
1:A:17:ILE:HD12	1:A:56:GLY:HA3	1.95	0.48
1:A:31:HIS:CE1	1:A:66:LEU:HD22	2.49	0.48
1:A:29:ILE:HD12	1:A:67:PHE:CD1	2.48	0.48
1:B:52:ARG:HG2	2:B:371:HOH:O	2.13	0.48
1:A:27:ILE:HD11	1:A:56:GLY:O	2.13	0.48
1:B:17:ILE:HD12	1:B:43:LYS:HD3	1.95	0.48
1:A:121:LEU:CD2	1:A:129:ALA:CB	2.92	0.48
1:B:54:THR:HG22	2:B:371:HOH:O	2.13	0.48
1:B:67:PHE:CE1	1:B:71:VAL:CG2	2.96	0.48
1:A:139:TYR:CE1	1:A:147:LYS:HG3	2.49	0.48
1:A:44:SER:O	1:A:48:LYS:HG3	2.14	0.47
1:B:39:LEU:HD11	1:B:43:LYS:CE	2.43	0.47
1:B:59:THR:HG22	1:B:62:GLU:H	1.80	0.47
1:A:85:LYS:N	1:A:86:PRO:CD	2.76	0.47
1:A:121:LEU:CD2	1:A:129:ALA:HB1	2.44	0.47
1:A:46:LEU:HD12	1:A:50:ILE:HD13	1.97	0.47
1:A:76:ARG:CB	1:A:80:ARG:NH2	2.78	0.47
1:B:119:ARG:NH2	1:B:120:MET:CE	2.78	0.47
1:B:31:HIS:ND1	1:B:66:LEU:CD1	2.77	0.47
1:B:26:THR:CG2	1:B:27:ILE:N	2.78	0.46
1:A:137:ARG:CZ	1:A:141:GLN:HB2	2.45	0.46
1:A:16:LYS:HD2	1:A:17:ILE:N	2.31	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:106:MET:HG2	1:B:111:VAL:CG2	2.46	0.46
1:B:54:THR:CG2	2:B:371:HOH:O	2.63	0.46
1:A:50:ILE:HG22	1:A:52:ARG:H	1.80	0.46
1:A:72:ASP:OD2	1:A:76:ARG:NH1	2.49	0.46
1:B:26:THR:HG22	1:B:27:ILE:N	2.31	0.46
1:B:6:MET:C	1:B:7:LEU:HD23	2.36	0.46
1:A:50:ILE:HG21	1:A:54:THR:CG2	2.46	0.46
1:A:16:LYS:HE2	1:A:55:ASN:O	2.15	0.46
1:A:60:LYS:O	1:A:63:ALA:HB3	2.15	0.46
1:B:91:LEU:HD22	1:B:95:ARG:HB3	1.98	0.46
1:B:120:MET:HB3	1:B:125:ARG:CB	2.46	0.46
1:B:132:ASN:O	1:B:135:LYS:HG3	2.16	0.46
1:A:58:ILE:HA	1:A:62:GLU:OE1	2.17	0.45
1:B:62:GLU:CA	1:B:65:LYS:NZ	2.79	0.45
1:A:24:TYR:CD2	1:A:35:LYS:HA	2.51	0.45
1:A:143:PRO:O	1:A:147:LYS:HB2	2.16	0.45
1:B:121:LEU:HD23	1:B:121:LEU:HA	1.57	0.45
1:A:120:MET:N	1:A:120:MET:CE	2.79	0.45
1:B:106:MET:HG2	1:B:111:VAL:HG23	1.98	0.45
1:A:8:ARG:O	1:A:12:GLY:HA2	2.16	0.45
1:A:152:THR:HG23	1:A:157:THR:O	2.16	0.45
1:A:16:LYS:HD3	1:A:56:GLY:C	2.37	0.45
1:A:120:MET:O	1:A:121:LEU:C	2.52	0.45
1:B:154:ARG:NH1	1:B:154:ARG:CG	2.76	0.45
1:B:91:LEU:HD13	1:B:95:ARG:HB3	1.98	0.45
1:B:19:LYS:HG3	1:B:23:GLY:C	2.37	0.45
1:B:8:ARG:O	1:B:12:GLY:HA2	2.16	0.45
1:B:111:VAL:O	1:B:113:GLY:N	2.50	0.45
1:B:151:THR:O	1:B:151:THR:HG22	2.17	0.45
1:A:121:LEU:HD23	1:A:129:ALA:CB	2.47	0.45
1:A:41:ALA:O	1:A:45:GLU:HG2	2.17	0.45
1:A:32:LEU:HB3	2:A:389:HOH:O	2.17	0.45
1:B:114:PHE:O	1:B:118:LEU:HD12	2.17	0.45
1:A:39:LEU:C	1:A:39:LEU:HD23	2.37	0.44
1:A:133:LEU:HA	1:A:133:LEU:HD23	1.57	0.44
1:B:132:ASN:HA	1:B:132:ASN:HD22	1.46	0.44
1:A:16:LYS:HZ3	1:A:56:GLY:HA3	1.82	0.44
1:A:76:ARG:HB3	1:A:80:ARG:NH2	2.33	0.44
1:B:1:MET:HG2	1:B:5:GLU:CB	2.48	0.44
1:B:78:ILE:HG22	1:B:79:LEU:HD23	1.99	0.44
1:B:120:MET:HG2	1:B:120:MET:H	1.37	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:103:VAL:HA	1:B:111:VAL:HG21	2.00	0.44
1:A:121:LEU:HD23	1:A:129:ALA:HB2	1.99	0.44
1:A:5:GLU:HB3	2:A:226:HOH:O	2.17	0.44
1:A:26:THR:HG23	1:A:31:HIS:O	2.18	0.44
1:A:65:LYS:HB3	1:A:65:LYS:HE3	1.68	0.44
1:B:54:THR:O	1:B:55:ASN:HB3	2.17	0.44
1:A:81:ASN:OD1	1:A:82:ALA:N	2.51	0.44
1:A:162:LYS:HB2	1:A:162:LYS:HE3	1.35	0.43
1:A:127:ASP:HB3	1:A:154:ARG:CZ	2.49	0.43
1:B:119:ARG:NH2	1:B:120:MET:HE1	2.33	0.43
1:A:88:TYR:O	1:A:96:ARG:NE	2.51	0.43
1:B:119:ARG:CZ	1:B:119:ARG:HB3	2.47	0.43
1:A:112:ALA:HA	1:A:118:LEU:HD11	2.01	0.43
1:A:60:LYS:O	1:A:63:ALA:N	2.52	0.43
1:A:100:ILE:HG22	1:A:100:ILE:O	2.19	0.43
1:A:50:ILE:CD1	1:A:50:ILE:N	2.81	0.43
1:A:16:LYS:CE	1:A:56:GLY:HA3	2.48	0.43
1:A:75:VAL:CG1	1:A:76:ARG:N	2.82	0.43
1:B:1:MET:HA	1:B:5:GLU:HG3	2.00	0.43
1:B:117:SER:OG	1:B:132:ASN:HB3	2.18	0.43
1:B:17:ILE:N	1:B:27:ILE:HD12	2.34	0.42
1:A:121:LEU:HD21	1:A:129:ALA:HB1	2.01	0.42
1:A:17:ILE:HD11	1:A:56:GLY:HA2	2.01	0.42
1:B:119:ARG:CB	1:B:119:ARG:HH11	2.33	0.42
1:B:137:ARG:NE	1:B:137:ARG:CA	2.80	0.42
1:B:32:LEU:CD2	1:B:33:LEU:N	2.82	0.42
1:B:19:LYS:HA	1:B:24:TYR:O	2.19	0.42
1:B:91:LEU:HD22	1:B:95:ARG:CB	2.49	0.42
1:A:88:TYR:HE1	1:A:100:ILE:HD11	1.84	0.42
1:A:65:LYS:NZ	1:A:66:LEU:HG	2.33	0.42
1:A:81:ASN:OD1	1:A:83:LYS:N	2.46	0.42
1:A:64:GLU:HG2	1:A:64:GLU:H	1.71	0.42
1:A:127:ASP:O	1:A:131:VAL:HG23	2.20	0.42
1:B:14:ARG:NH1	1:B:14:ARG:CG	2.83	0.42
1:B:119:ARG:CZ	1:B:120:MET:HE3	2.49	0.42
1:A:6:MET:HE1	1:A:101:ASN:HB2	2.02	0.41
1:B:31:HIS:CG	1:B:66:LEU:CD1	3.01	0.41
1:A:140:ASN:C	1:A:141:GLN:HG2	2.40	0.41
1:B:65:LYS:O	1:B:66:LEU:CB	2.65	0.41
1:B:46:LEU:O	1:B:47:ASP:C	2.58	0.41
1:B:119:ARG:NH1	1:B:120:MET:H	2.13	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:119:ARG:O	1:A:122:GLN:HB3	2.20	0.41
1:B:64:GLU:HB2	2:B:243:HOH:O	2.20	0.41
1:A:98:MET:HE3	1:A:98:MET:CA	2.43	0.41
1:B:120:MET:HB3	1:B:120:MET:HE2	1.71	0.41
1:A:67:PHE:CE1	1:A:71:VAL:CG2	3.03	0.41
1:A:94:VAL:O	1:A:97:ALA:HB3	2.20	0.41
1:A:45:GLU:O	1:A:46:LEU:C	2.59	0.41
1:A:129:ALA:O	1:A:132:ASN:HB2	2.20	0.41
1:A:50:ILE:HG21	1:A:52:ARG:HG3	1.99	0.41
1:B:137:ARG:CZ	1:B:137:ARG:CB	2.99	0.41
1:A:105:GLN:O	1:A:105:GLN:HG2	2.20	0.41
1:A:145:ARG:HD2	1:A:145:ARG:HH11	1.64	0.41
1:A:6:MET:CE	1:A:98:MET:SD	3.09	0.41
1:B:13:LEU:HB2	1:B:29:ILE:HG12	2.03	0.41
1:B:120:MET:HE2	1:B:125:ARG:HG2	2.03	0.41
1:B:32:LEU:C	1:B:32:LEU:HD22	2.40	0.41
1:B:61:ASP:HA	1:B:64:GLU:HB3	2.02	0.41
1:A:59:THR:N	1:A:62:GLU:HG3	2.35	0.41
1:B:137:ARG:C	1:B:137:ARG:NE	2.74	0.41
1:B:120:MET:CB	1:B:125:ARG:CB	2.99	0.40
1:B:48:LYS:C	1:B:50:ILE:H	2.23	0.40
1:A:123:GLN:HB2	1:A:125:ARG:HD3	2.03	0.40
1:B:96:ARG:HH11	1:B:96:ARG:HD2	1.65	0.40
1:A:17:ILE:HG13	1:A:27:ILE:HD12	2.02	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	160/164 (98%)	140 (88%)	17 (11%)	3 (2%)	10	2

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	B	160/164 (98%)	141 (88%)	15 (9%)	4 (2%)	7 1
All	All	320/328 (98%)	281 (88%)	32 (10%)	7 (2%)	8 1

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	60	LYS
1	B	66	LEU
1	B	159	ASP
1	B	112	ALA
1	A	141	GLN
1	B	126	TRP
1	A	27	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	136/138 (99%)	92 (68%)	44 (32%)	0 0
1	B	136/138 (99%)	96 (71%)	40 (29%)	0 0
All	All	272/276 (99%)	188 (69%)	84 (31%)	0 0

All (84) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	ASN
1	A	8	ARG
1	A	9	ILE
1	A	10	ASP
1	A	13	LEU
1	A	14	ARG
1	A	27	ILE
1	A	32	LEU
1	A	35	LYS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	38	SER
1	A	40	ASN
1	A	44	SER
1	A	47	ASP
1	A	48	LYS
1	A	52	ARG
1	A	59	THR
1	A	60	LYS
1	A	61	ASP
1	A	65	LYS
1	A	69	GLN
1	A	72	ASP
1	A	75	VAL
1	A	76	ARG
1	A	83	LYS
1	A	102	MET
1	A	103	VAL
1	A	106	MET
1	A	108	GLU
1	A	116	ASN
1	A	119	ARG
1	A	120	MET
1	A	121	LEU
1	A	124	LYS
1	A	125	ARG
1	A	127	ASP
1	A	132	ASN
1	A	135	LYS
1	A	137	ARG
1	A	140	ASN
1	A	141	GLN
1	A	142	THR
1	A	147	LYS
1	A	154	ARG
1	A	162	LYS
1	B	1	MET
1	B	2	ASN
1	B	3	ILE
1	B	8	ARG
1	B	13	LEU
1	B	14	ARG
1	B	21	THR

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	32	LEU
1	B	40	ASN
1	B	43	LYS
1	B	44	SER
1	B	52	ARG
1	B	57	VAL
1	B	59	THR
1	B	60	LYS
1	B	65	LYS
1	B	68	ASN
1	B	72	ASP
1	B	79	LEU
1	B	96	ARG
1	B	106	MET
1	B	108	GLU
1	B	109	THR
1	B	117	SER
1	B	119	ARG
1	B	120	MET
1	B	122	GLN
1	B	124	LYS
1	B	125	ARG
1	B	127	ASP
1	B	132	ASN
1	B	135	LYS
1	B	137	ARG
1	B	140	ASN
1	B	141	GLN
1	B	144	ASN
1	B	147	LYS
1	B	153	PHE
1	B	154	ARG
1	B	162	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	40	ASN
1	A	101	ASN
1	A	122	GLN
1	A	132	ASN
1	B	40	ASN

Continued from previous page...

Mol	Chain	Res	Type
1	B	123	GLN
1	B	132	ASN
1	B	140	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)

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 (i)

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	162/164 (98%)	-0.73	0	100	100	22, 40, 65, 76	0
1	B	162/164 (98%)	-0.75	0	100	100	22, 38, 60, 78	0
All	All	324/328 (98%)	-0.74	0	100	100	22, 39, 63, 78	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 carbohydrates 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.