



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:01 AM UTC

PDB ID : 1R4Q / pdb_00001r4q
Title : Shiga toxin
Authors : Fraser, M.E.; Fujinaga, M.; Cherney, M.M.; Melton-Celsa, A.R.; Twiddy, E.M.; O'Brien, A.D.; James, M.N.G.
Deposited on : 2003-10-07
Resolution : 2.50 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

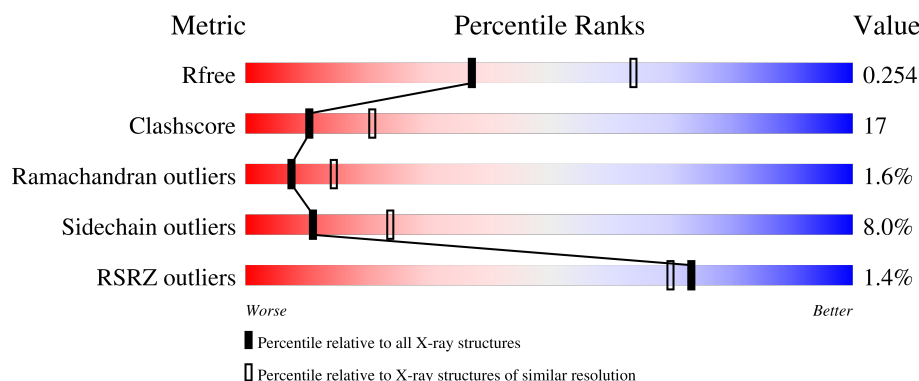
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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))
R_{free}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	293	2% 62% 26% 6% 6%
1	L	293	2% 61% 28% 6% 5%
2	B	69	61% 33% 6%
2	C	69	% 55% 41% .
2	D	69	67% 30% .

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Mol	Chain	Length	Quality of chain
2	E	69	75% 20%
2	F	69	4% 62% 28% 7%
2	G	69	74% 23%
2	H	69	75% 19%
2	I	69	% 70% 26%
2	J	69	% 57% 38% 6%
2	K	69	3% 51% 42%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9725 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 SHT cytotoxin A subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	276	Total	C	N	O	S	0	0	0
			2136	1334	386	407	9			
1	L	277	Total	C	N	O	S	0	0	0
			2143	1338	387	409	9			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	168	ALA	PRO	see remark 999	UNP Q7BQ99
L	168	ALA	PRO	see remark 999	UNP Q7BQ99

- Molecule 2 is a protein called Shigella toxin chain B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	69	Total	C	N	O	S	0	0	0
			540	339	90	108	3			
2	C	69	Total	C	N	O	S	0	0	0
			540	339	90	108	3			
2	D	69	Total	C	N	O	S	0	0	0
			540	339	90	108	3			
2	E	69	Total	C	N	O	S	0	0	0
			540	339	90	108	3			
2	F	69	Total	C	N	O	S	0	0	0
			540	339	90	108	3			
2	G	69	Total	C	N	O	S	0	0	0
			540	339	90	108	3			
2	H	69	Total	C	N	O	S	0	0	0
			540	339	90	108	3			
2	I	69	Total	C	N	O	S	0	0	0
			540	339	90	108	3			
2	J	69	Total	C	N	O	S	0	0	0
			540	339	90	108	3			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	K	69	Total	C	N	O	S	0	0	0
			540	339	90	108	3			

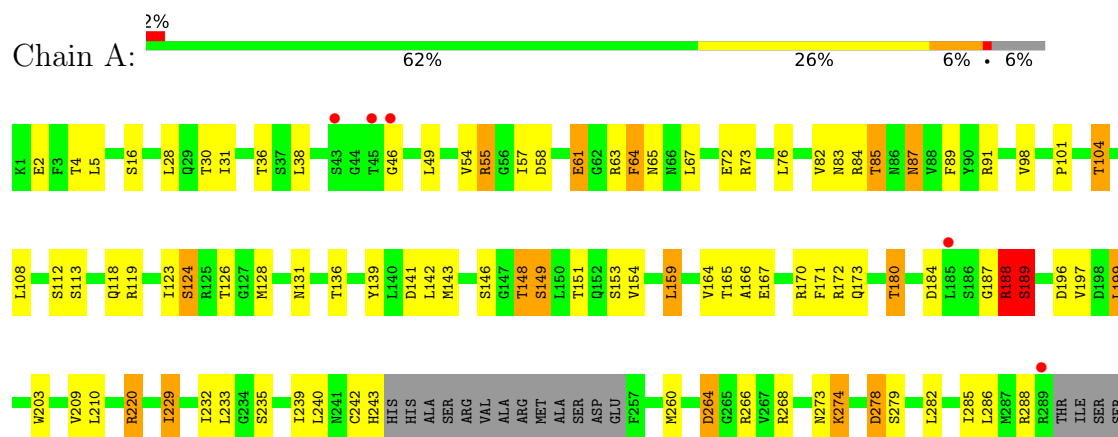
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	9	Total	O	0	0
			9	9		
3	L	6	Total	O	0	0
			6	6		
3	B	1	Total	O	0	0
			1	1		
3	C	1	Total	O	0	0
			1	1		
3	D	1	Total	O	0	0
			1	1		
3	E	4	Total	O	0	0
			4	4		
3	F	2	Total	O	0	0
			2	2		
3	G	10	Total	O	0	0
			10	10		
3	H	5	Total	O	0	0
			5	5		
3	I	1	Total	O	0	0
			1	1		
3	J	2	Total	O	0	0
			2	2		
3	K	4	Total	O	0	0
			4	4		

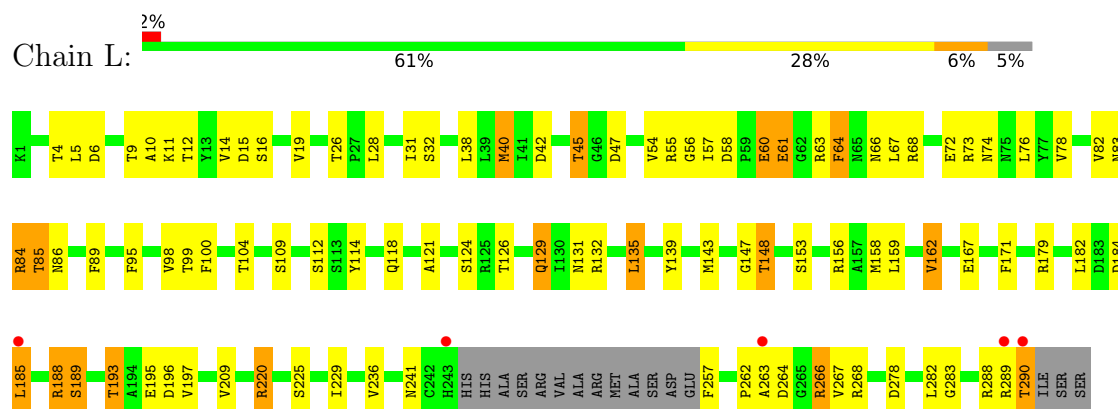
3 Residue-property plots [i](#)

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.

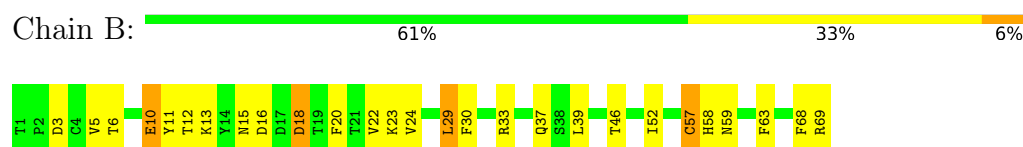
• Molecule 1: SHT cytotoxin A subunit



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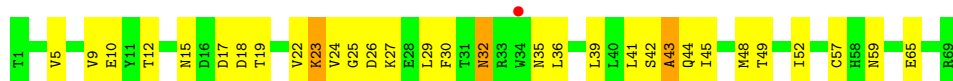


• Molecule 2: Shigella toxin chain B



• Molecule 2: Shigella toxin chain B

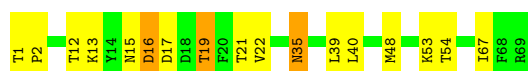




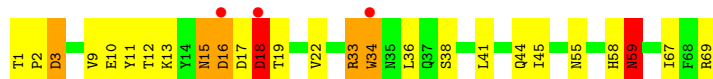
- Molecule 2: Shigella toxin chain B



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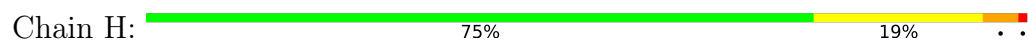
- Molecule 2: Shigella toxin chain B



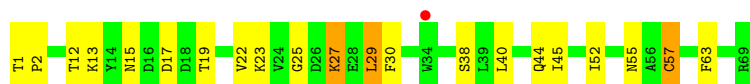
- Molecule 2: Shigella toxin chain B



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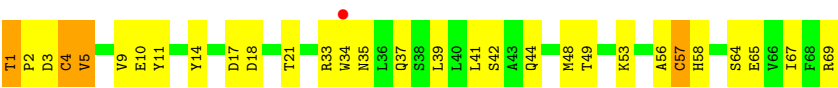


- Molecule 2: Shigella toxin chain B

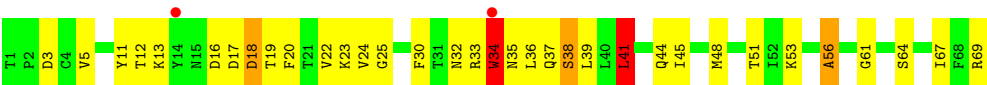


- Molecule 2: Shigella toxin chain B





● Molecule 2: Shigella toxin chain B



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	133.00Å 147.18Å 82.85Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	98.60 – 2.50 98.60 – 2.50	Depositor EDS
% Data completeness (in resolution range)	(Not available) (98.60-2.50) 94.0 (98.60-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.56 (at 2.51Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.199 , 0.257 0.198 , 0.254	Depositor DCC
R_{free} test set	2747 reflections (5.13%)	wwPDB-VP
Wilson B-factor (Å ²)	38.6	Xtriage
Anisotropy	0.214	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 35.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9725	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 3.39% 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	0.94	2/2171 (0.1%)	1.27	19/2941 (0.6%)
1	L	0.87	2/2178 (0.1%)	1.24	20/2951 (0.7%)
2	B	0.78	0/549	1.10	2/742 (0.3%)
2	C	0.82	0/549	1.15	3/742 (0.4%)
2	D	0.96	0/549	1.17	1/742 (0.1%)
2	E	1.10	0/549	1.21	2/742 (0.3%)
2	F	1.06	2/549 (0.4%)	1.28	4/742 (0.5%)
2	G	1.03	0/549	1.18	0/742
2	H	0.96	0/549	1.31	6/742 (0.8%)
2	I	0.84	1/549 (0.2%)	1.02	0/742
2	J	0.87	0/549	1.20	5/742 (0.7%)
2	K	1.02	1/549 (0.2%)	1.25	4/742 (0.5%)
All	All	0.93	8/9839 (0.1%)	1.22	66/13312 (0.5%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	229	ILE	CA-CB	6.33	1.62	1.54
2	F	16	ASP	CA-C	5.79	1.58	1.52
2	K	34	TRP	CB-CG	5.75	1.68	1.50
1	L	45	THR	CA-CB	5.72	1.61	1.52
1	A	197	VAL	CA-CB	5.55	1.60	1.54
1	L	40	MET	SD-CE	5.36	1.93	1.79
2	F	59	ASN	CG-OD1	5.18	1.33	1.23
2	I	45	ILE	CA-CB	5.05	1.61	1.54

All (66) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	104	THR	N-CA-C	-8.90	96.73	110.17
1	L	73	ARG	N-CA-C	8.64	120.31	111.07
2	J	1	THR	CA-C-N	7.88	127.81	119.85
2	J	1	THR	C-N-CA	7.88	127.81	119.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	73	ARG	N-CA-C	7.69	123.12	112.04
2	F	15	ASN	N-CA-C	7.63	121.69	109.79
1	A	278	ASP	N-CA-C	-7.58	97.23	109.96
1	L	257	PHE	CA-C-N	7.28	127.69	119.83
1	L	257	PHE	C-N-CA	7.28	127.69	119.83
1	L	148	THR	N-CA-C	7.26	122.55	112.45
2	F	34	TRP	N-CA-C	7.18	120.02	111.33
1	L	209	VAL	N-CA-C	7.15	119.16	111.58
2	J	4	CYS	N-CA-C	7.05	121.50	112.34
1	A	112	SER	N-CA-C	7.00	121.05	112.23
2	F	16	ASP	N-CA-C	-6.84	103.83	112.93
1	L	100	PHE	CA-C-N	6.75	127.15	119.93
1	L	100	PHE	C-N-CA	6.75	127.15	119.93
1	L	159	LEU	N-CA-C	-6.71	104.05	111.36
1	A	209	VAL	N-CA-C	6.66	117.45	110.72
2	H	1	THR	CA-C-N	6.53	126.49	119.76
2	H	1	THR	C-N-CA	6.53	126.49	119.76
2	B	3	ASP	N-CA-C	-6.39	102.05	110.43
1	A	279	SER	N-CA-C	6.37	119.21	111.82
1	A	87	ASN	N-CA-C	6.36	120.62	112.86
1	L	135	LEU	N-CA-C	-6.34	105.41	113.01
2	K	34	TRP	CA-CB-CG	6.27	125.50	113.60
2	E	16	ASP	N-CA-C	6.24	118.89	111.71
1	L	104	THR	N-CA-C	-6.21	99.78	109.72
1	A	164	VAL	CA-C-N	-6.20	115.06	122.44
1	A	164	VAL	C-N-CA	-6.20	115.06	122.44
1	L	278	ASP	N-CA-C	-6.19	100.47	110.32
1	L	112	SER	N-CA-C	6.10	120.45	112.89
1	L	182	LEU	N-CA-C	-6.03	105.57	113.17
1	L	147	GLY	N-CA-C	5.98	120.54	112.17
1	A	159	LEU	N-CA-C	-5.88	104.87	111.28
1	L	74	ASN	N-CA-C	5.88	117.68	111.28
1	A	180	THR	N-CA-C	5.83	119.49	112.38
1	A	210	LEU	CA-C-N	-5.78	113.73	119.56
1	A	210	LEU	C-N-CA	-5.78	113.73	119.56
2	C	43	ALA	N-CA-C	-5.75	105.01	111.28
1	L	264	ASP	N-CA-C	5.71	118.87	110.30
1	L	171	PHE	N-CA-C	5.66	118.00	107.99
2	F	3	ASP	N-CA-C	-5.61	101.98	110.28
1	L	197	VAL	N-CA-C	-5.58	105.06	110.42
1	A	232	ILE	CB-CA-C	-5.46	104.87	112.02
2	K	39	LEU	N-CA-C	-5.44	105.25	111.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L	267	VAL	N-CA-C	5.43	116.39	108.58
1	A	149	SER	N-CA-C	5.41	118.28	110.28
2	C	57	CYS	N-CA-C	5.35	118.31	108.58
1	A	136	THR	N-CA-C	-5.34	105.36	111.07
1	A	146	SER	N-CA-C	-5.31	101.22	109.24
2	K	51	THR	N-CA-C	-5.29	99.79	108.41
2	B	18	ASP	N-CA-C	5.29	119.31	112.86
2	E	19	THR	N-CA-C	-5.27	101.12	109.76
2	K	41	LEU	CA-CB-CG	-5.23	98.00	116.30
2	H	33	ARG	N-CA-C	-5.20	100.60	108.67
2	H	45	ILE	CB-CA-C	-5.19	105.13	112.14
2	H	43	ALA	N-CA-C	-5.18	105.72	111.36
1	A	264	ASP	N-CA-C	5.16	121.78	110.80
2	D	37	GLN	N-CA-C	5.14	116.58	110.97
1	A	210	LEU	N-CA-C	5.14	121.17	109.81
2	H	45	ILE	N-CA-C	5.10	115.82	110.62
2	C	15	ASN	N-CA-C	5.09	118.05	110.52
1	L	56	GLY	N-CA-C	5.08	122.89	115.08
2	J	9	VAL	CA-C-N	5.05	127.31	120.65
2	J	9	VAL	C-N-CA	5.05	127.31	120.65

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	2136	0	2129	72	0
1	L	2143	0	2136	78	0
2	B	540	0	527	21	0
2	C	540	0	527	23	0
2	D	540	0	527	17	0
2	E	540	0	527	15	0
2	F	540	0	527	25	0
2	G	540	0	527	15	0
2	H	540	0	527	15	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	I	540	0	527	18	0
2	J	540	0	527	24	0
2	K	540	0	527	46	0
3	A	9	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
3	E	4	0	0	0	0
3	F	2	0	0	0	0
3	G	10	0	0	0	0
3	H	5	0	0	0	0
3	I	1	0	0	0	0
3	J	2	0	0	0	0
3	K	4	0	0	1	0
3	L	6	0	0	0	0
All	All	9725	0	9535	324	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 17.

All (324) 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)
2:J:39:LEU:HD22	2:K:41:LEU:HD22	1.41	1.01
2:F:16:ASP:C	2:F:18:ASP:H	1.64	1.01
2:D:69:ARG:HG3	2:D:69:ARG:HH11	1.28	0.98
1:L:10:ALA:HB1	1:L:179:ARG:HG2	1.47	0.97
2:K:12:THR:HG22	2:K:22:VAL:HG12	1.45	0.94
1:L:121:ALA:HB1	1:L:156:ARG:HG3	1.55	0.88
1:A:188:ARG:O	1:A:189:SER:HB3	1.73	0.88
2:J:65:GLU:HG3	2:K:16:ASP:HB2	1.59	0.85
1:L:60:GLU:HG2	1:L:61:GLU:HG2	1.61	0.83
2:I:1:THR:HG23	2:I:2:PRO:HD2	1.62	0.82
1:L:40:MET:HE2	1:L:241:ASN:HA	1.64	0.79
2:D:69:ARG:HG3	2:D:69:ARG:NH1	1.94	0.79
2:F:12:THR:HG22	2:F:22:VAL:HG12	1.64	0.77
2:E:67:ILE:HG12	2:F:13:LYS:HG2	1.66	0.76
1:L:185:LEU:HD12	1:L:185:LEU:H	1.49	0.75
2:K:17:ASP:O	2:K:18:ASP:HB2	1.86	0.75
1:L:60:GLU:H	1:L:60:GLU:CD	1.95	0.74
1:A:188:ARG:HB3	1:A:188:ARG:HH11	1.52	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:1:THR:O	2:G:53:LYS:HE3	1.86	0.73
1:A:61:GLU:OE1	1:A:63:ARG:NH2	2.22	0.73
2:E:22:VAL:HG11	2:E:40:LEU:HD13	1.70	0.72
1:A:170:ARG:HD3	1:A:203:TRP:CE2	2.25	0.72
1:A:188:ARG:HH11	1:A:188:ARG:CB	2.03	0.71
2:K:5:VAL:HG21	2:K:24:VAL:O	1.89	0.71
1:A:54:VAL:HG12	1:A:57:ILE:HD11	1.73	0.71
2:B:22:VAL:HG12	2:B:24:VAL:HG23	1.73	0.71
1:L:15:ASP:O	1:L:19:VAL:HG23	1.89	0.71
1:L:114:TYR:HB3	1:L:118:GLN:HE21	1.56	0.70
1:L:220:ARG:HB3	1:L:220:ARG:NH1	2.06	0.70
2:K:33:ARG:HH12	2:K:64:SER:HB3	1.54	0.70
1:L:10:ALA:HB1	1:L:179:ARG:CG	2.21	0.69
2:K:17:ASP:OD1	2:K:19:THR:HG22	1.93	0.69
1:A:84:ARG:HH11	1:A:84:ARG:HG3	1.58	0.69
2:E:13:LYS:HE3	2:E:15:ASN:OD1	1.93	0.69
2:C:5:VAL:HG21	2:C:24:VAL:O	1.93	0.68
1:L:188:ARG:HH11	1:L:188:ARG:HA	1.59	0.68
1:L:68:ARG:HD2	1:L:84:ARG:HD3	1.75	0.68
2:B:16:ASP:O	2:F:33:ARG:NH2	2.26	0.68
2:I:13:LYS:HE3	2:I:15:ASN:OD1	1.96	0.66
1:A:118:GLN:HE21	1:A:124:SER:HA	1.59	0.66
1:A:184:ASP:OD2	1:A:188:ARG:HG3	1.96	0.66
2:F:18:ASP:HB3	2:K:32:ASN:CB	2.27	0.65
2:F:34:TRP:CZ2	2:K:34:TRP:HB3	2.31	0.65
2:I:27:LYS:HB3	2:I:29:LEU:HD21	1.78	0.65
1:L:131:ASN:HB3	1:L:189:SER:CB	2.27	0.65
1:A:171:PHE:CD1	1:A:199:LEU:HD13	2.32	0.65
2:J:39:LEU:HD22	2:K:41:LEU:CD2	2.20	0.65
2:D:1:THR:HG21	2:D:54:THR:HA	1.80	0.64
2:B:13:LYS:HD3	2:B:15:ASN:OD1	1.98	0.64
1:A:54:VAL:CG1	1:A:57:ILE:HD11	2.27	0.64
2:I:23:LYS:NZ	2:I:25:GLY:O	2.30	0.64
1:L:95:PHE:HB3	1:L:98:VAL:CG2	2.26	0.64
2:G:32:ASN:C	2:G:32:ASN:HD22	2.05	0.63
2:F:16:ASP:C	2:F:18:ASP:N	2.43	0.63
2:K:53:LYS:HD2	2:K:67:ILE:HD11	1.80	0.63
1:A:184:ASP:CG	1:A:188:ARG:HG3	2.25	0.62
2:F:18:ASP:HB3	2:K:32:ASN:HB2	1.81	0.62
2:H:5:VAL:HG22	2:H:24:VAL:HG12	1.81	0.62
1:A:84:ARG:HG3	1:A:84:ARG:NH1	2.12	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:1:THR:CG2	2:E:2:PRO:HD2	2.29	0.62
2:K:33:ARG:O	2:K:37:GLN:HG3	1.99	0.62
1:L:54:VAL:CG1	1:L:57:ILE:HD11	2.29	0.61
1:L:185:LEU:HD12	1:L:185:LEU:N	2.16	0.61
2:F:34:TRP:CH2	2:K:34:TRP:HB3	2.35	0.61
1:A:57:ILE:N	1:A:57:ILE:HD12	2.16	0.61
1:A:188:ARG:O	1:A:189:SER:CB	2.48	0.61
1:A:58:ASP:HB2	1:A:64:PHE:CE2	2.36	0.61
2:B:12:THR:HB	2:B:20:PHE:CZ	2.36	0.61
2:C:43:ALA:HA	2:C:48:MET:HE2	1.82	0.61
1:L:188:ARG:HA	1:L:188:ARG:NH1	2.16	0.61
2:I:17:ASP:OD1	2:I:19:THR:HG23	1.99	0.61
1:A:188:ARG:HA	1:A:188:ARG:NH1	2.16	0.60
2:J:56:ALA:O	2:J:58:HIS:N	2.31	0.60
2:K:33:ARG:NH1	2:K:64:SER:HB3	2.15	0.60
1:A:31:ILE:HG23	1:A:229:ILE:HD11	1.82	0.60
1:L:6:ASP:HB3	1:L:12:THR:HB	1.83	0.60
2:K:11:TYR:CE2	2:K:23:LYS:HD2	2.36	0.60
2:E:1:THR:HG21	2:E:54:THR:HA	1.84	0.60
1:L:95:PHE:HB3	1:L:98:VAL:HG21	1.84	0.60
1:L:67:LEU:HD12	1:L:68:ARG:H	1.67	0.60
1:L:5:LEU:HD12	1:L:16:SER:CB	2.32	0.60
1:L:193:THR:HG23	1:L:196:ASP:OD1	2.02	0.59
1:A:242:CYS:O	1:A:243:HIS:C	2.46	0.59
1:L:78:VAL:HG21	1:L:162:VAL:HG22	1.85	0.59
2:E:1:THR:HG23	2:E:2:PRO:HD2	1.85	0.59
2:G:45:ILE:CG1	2:K:48:MET:HE3	2.33	0.59
1:L:5:LEU:HD12	1:L:16:SER:HB2	1.85	0.58
1:L:220:ARG:HB3	1:L:220:ARG:HH11	1.69	0.58
1:L:158:MET:O	1:L:162:VAL:HB	2.03	0.58
1:A:54:VAL:HG21	1:A:67:LEU:HD23	1.85	0.58
2:G:45:ILE:HG12	2:K:48:MET:HE3	1.84	0.58
2:F:15:ASN:HB2	2:F:19:THR:O	2.04	0.57
2:F:16:ASP:O	2:F:18:ASP:N	2.36	0.57
1:L:167:GLU:HA	1:L:167:GLU:OE1	2.03	0.57
1:A:141:ASP:OD2	1:A:153:SER:HB3	2.04	0.57
1:L:4:THR:HG23	1:L:55:ARG:HG3	1.87	0.57
1:L:10:ALA:CB	1:L:179:ARG:HG2	2.30	0.57
2:F:41:LEU:O	2:F:45:ILE:HG13	2.04	0.57
1:L:63:ARG:HH11	1:L:63:ARG:HB3	1.69	0.57
2:J:4:CYS:O	2:J:5:VAL:HB	2.04	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:288:ARG:HG2	2:K:38:SER:HB3	1.86	0.56
2:D:69:ARG:NH1	2:D:69:ARG:CG	2.67	0.56
1:A:49:LEU:HD11	1:A:98:VAL:HG13	1.86	0.56
1:L:64:PHE:N	1:L:64:PHE:CD1	2.74	0.56
2:J:53:LYS:HE2	2:J:67:ILE:HD11	1.87	0.56
2:H:33:ARG:HG2	2:H:33:ARG:HH11	1.71	0.56
2:H:30:PHE:CD1	2:H:30:PHE:C	2.84	0.56
2:G:32:ASN:C	2:G:32:ASN:ND2	2.61	0.56
1:A:83:ASN:OD1	1:A:85:THR:HB	2.06	0.56
1:A:167:GLU:OE1	1:A:167:GLU:HA	2.06	0.56
2:J:65:GLU:HG2	2:K:16:ASP:N	2.21	0.55
1:A:82:VAL:HG22	1:A:89:PHE:HB2	1.88	0.55
2:K:12:THR:HG22	2:K:22:VAL:CG1	2.28	0.55
1:A:5:LEU:CD2	1:A:165:THR:HG22	2.37	0.54
1:A:220:ARG:NH1	1:A:220:ARG:HB3	2.22	0.54
2:K:3:ASP:OD2	3:K:72:HOH:O	2.18	0.54
1:L:68:ARG:HB3	1:L:82:VAL:HB	1.88	0.54
2:B:13:LYS:HG3	2:F:67:ILE:HG12	1.89	0.54
2:B:58:HIS:O	2:B:59:ASN:C	2.49	0.54
2:I:29:LEU:HD12	2:I:52:ILE:HG21	1.89	0.54
2:C:27:LYS:HA	2:C:59:ASN:HD21	1.72	0.54
2:G:22:VAL:HG13	2:G:24:VAL:HG23	1.89	0.54
2:C:9:VAL:CG1	2:C:44:GLN:HG3	2.38	0.54
2:K:20:PHE:O	2:K:30:PHE:HB2	2.08	0.54
2:J:11:TYR:HA	2:J:44:GLN:NE2	2.22	0.53
2:F:9:VAL:HB	2:F:44:GLN:HG3	1.90	0.53
1:L:72:GLU:O	1:L:76:LEU:HA	2.08	0.53
1:L:38:LEU:CD1	1:L:236:VAL:HG12	2.38	0.53
2:B:10:GLU:OE1	2:B:23:LYS:NZ	2.38	0.53
2:E:22:VAL:CG1	2:E:40:LEU:HD13	2.38	0.53
2:I:27:LYS:HB3	2:I:29:LEU:CD2	2.38	0.53
1:L:14:VAL:HG21	1:L:179:ARG:HD3	1.89	0.53
1:L:266:ARG:NH1	1:L:266:ARG:HG3	2.24	0.53
1:L:185:LEU:H	1:L:185:LEU:CD1	2.20	0.53
1:L:266:ARG:HG3	1:L:266:ARG:HH11	1.73	0.53
1:A:220:ARG:HB3	1:A:220:ARG:HH11	1.73	0.53
2:K:69:ARG:HG2	2:K:69:ARG:HH11	1.73	0.53
2:I:55:ASN:C	2:I:57:CYS:H	2.18	0.52
2:J:67:ILE:HG12	2:K:13:LYS:HB3	1.92	0.52
1:A:288:ARG:HD3	2:F:38:SER:HB3	1.92	0.52
1:L:54:VAL:HG13	1:L:57:ILE:HD11	1.90	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:17:ASP:OD1	2:E:19:THR:HG23	2.09	0.52
1:A:118:GLN:NE2	1:A:124:SER:HA	2.25	0.52
1:A:128:MET:HE1	1:A:159:LEU:HB3	1.92	0.52
1:A:142:LEU:CD2	1:A:154:VAL:HG13	2.40	0.52
2:F:18:ASP:HB3	2:K:32:ASN:HB3	1.91	0.52
1:A:142:LEU:HD23	1:A:154:VAL:HG13	1.91	0.51
1:L:57:ILE:HD12	1:L:57:ILE:N	2.26	0.51
2:J:65:GLU:CG	2:K:16:ASP:HB2	2.36	0.51
1:L:57:ILE:O	1:L:57:ILE:HG22	2.10	0.51
2:G:32:ASN:ND2	2:G:33:ARG:HD3	2.26	0.51
1:L:66:ASN:HB2	1:L:143:MET:HE2	1.93	0.51
2:H:32:ASN:N	2:H:32:ASN:OD1	2.43	0.51
1:A:151:THR:OG1	1:A:154:VAL:HG23	2.11	0.50
2:E:39:LEU:HD11	2:F:41:LEU:HD13	1.93	0.50
2:K:5:VAL:CG2	2:K:25:GLY:HA3	2.41	0.50
1:L:290:THR:HG22	2:K:34:TRP:CD1	2.47	0.50
2:C:27:LYS:HA	2:C:59:ASN:ND2	2.25	0.50
2:B:68:PHE:HB2	2:C:12:THR:OG1	2.11	0.50
2:G:20:PHE:HB2	2:G:37:GLN:HG2	1.94	0.50
2:I:27:LYS:CB	2:I:29:LEU:HD21	2.41	0.50
1:A:123:ILE:HD11	1:A:128:MET:HE2	1.92	0.49
2:C:36:LEU:HD23	2:C:39:LEU:HD12	1.93	0.49
1:L:83:ASN:OD1	1:L:85:THR:HB	2.13	0.49
2:J:53:LYS:HB2	2:J:65:GLU:HB2	1.93	0.49
2:G:32:ASN:HD21	2:G:33:ARG:HD3	1.77	0.49
2:I:30:PHE:O	2:I:63:PHE:CD2	2.65	0.49
1:L:195:GLU:CD	1:L:195:GLU:H	2.20	0.49
2:I:1:THR:CG2	2:I:2:PRO:HD2	2.39	0.49
2:J:34:TRP:O	2:J:37:GLN:HG3	2.12	0.49
2:D:12:THR:HG22	2:D:22:VAL:HG23	1.95	0.49
2:J:53:LYS:HE2	2:J:67:ILE:CD1	2.43	0.49
1:A:268:ARG:NH1	2:C:45:ILE:O	2.46	0.48
2:D:69:ARG:HH11	2:D:69:ARG:CG	2.09	0.48
1:A:240:LEU:HD11	1:A:260:MET:HB2	1.95	0.48
2:C:19:THR:HG22	2:C:32:ASN:HB3	1.94	0.48
1:L:95:PHE:HB3	1:L:98:VAL:HG23	1.96	0.48
2:B:22:VAL:CG1	2:B:24:VAL:HG23	2.42	0.48
1:A:5:LEU:HD22	1:A:165:THR:HG22	1.95	0.48
1:L:11:LYS:O	1:L:15:ASP:OD2	2.32	0.48
1:L:288:ARG:CG	2:K:38:SER:HB3	2.43	0.48
2:C:35:ASN:OD1	2:C:35:ASN:N	2.44	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:39:LEU:CD1	2:F:41:LEU:HD13	2.44	0.48
1:A:166:ALA:O	1:A:170:ARG:HG3	2.14	0.48
2:B:30:PHE:CD1	2:B:30:PHE:C	2.92	0.47
1:A:58:ASP:HB2	1:A:64:PHE:HE2	1.76	0.47
2:C:43:ALA:HA	2:C:48:MET:CE	2.45	0.47
2:B:33:ARG:NH1	2:J:17:ASP:HB2	2.29	0.47
2:G:48:MET:HE2	2:H:45:ILE:HG13	1.95	0.47
2:J:33:ARG:HH12	2:K:18:ASP:CG	2.22	0.47
2:C:36:LEU:HD23	2:C:36:LEU:HA	1.78	0.47
2:G:18:ASP:HB3	2:G:34:TRP:CH2	2.49	0.47
1:A:108:LEU:HA	1:A:148:THR:O	2.15	0.47
2:D:1:THR:HG22	2:D:2:PRO:HD2	1.95	0.47
2:I:44:GLN:O	2:I:44:GLN:HG2	2.14	0.47
2:B:46:THR:HG21	2:C:45:ILE:HD13	1.96	0.47
1:L:85:THR:HG22	1:L:86:ASN:N	2.29	0.47
1:A:84:ARG:HH11	1:A:84:ARG:CG	2.27	0.46
1:L:262:PRO:O	1:L:263:ALA:C	2.58	0.46
2:H:20:PHE:CG	2:H:37:GLN:HG2	2.50	0.46
2:I:17:ASP:CG	2:I:19:THR:HG23	2.40	0.46
1:A:28:LEU:HD12	1:A:38:LEU:HD23	1.97	0.46
1:L:28:LEU:HD23	1:L:28:LEU:HA	1.70	0.46
1:L:31:ILE:HD12	1:L:38:LEU:CD2	2.46	0.46
2:G:48:MET:HE2	2:H:45:ILE:CG1	2.45	0.46
1:L:290:THR:CG2	2:K:34:TRP:CD1	2.99	0.46
2:J:33:ARG:NH1	2:K:18:ASP:OD1	2.45	0.46
1:A:123:ILE:HD11	1:A:128:MET:CE	2.46	0.46
1:A:170:ARG:HD3	1:A:203:TRP:CZ2	2.50	0.45
2:H:5:VAL:CG2	2:H:24:VAL:HG12	2.45	0.45
2:J:1:THR:HA	2:J:2:PRO:HD3	1.78	0.45
1:A:58:ASP:HB2	1:A:64:PHE:CD2	2.51	0.45
1:A:57:ILE:HD12	1:A:57:ILE:H	1.79	0.45
2:G:32:ASN:HD22	2:G:33:ARG:N	2.14	0.45
2:B:11:TYR:CD1	2:B:11:TYR:C	2.94	0.45
2:B:20:PHE:CD1	2:B:37:GLN:HG2	2.52	0.45
2:E:12:THR:HA	2:E:21:THR:O	2.16	0.45
1:A:172:ARG:HD2	1:A:233:LEU:O	2.17	0.45
1:L:47:ASP:OD2	1:L:72:GLU:OE2	2.34	0.45
2:D:1:THR:CG2	2:D:2:PRO:HD2	2.46	0.45
2:C:29:LEU:HD13	2:C:52:ILE:HG21	1.99	0.45
2:C:23:LYS:HE2	2:C:25:GLY:O	2.16	0.45
2:J:48:MET:HE1	2:K:44:GLN:NE2	2.32	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:184:ASP:OD2	1:L:188:ARG:HG3	2.17	0.44
2:B:57:CYS:O	2:B:57:CYS:SG	2.76	0.44
2:C:17:ASP:O	2:C:18:ASP:HB2	2.17	0.44
2:D:46:THR:OG1	2:D:48:MET:HG3	2.17	0.44
2:I:1:THR:HG23	2:I:2:PRO:CD	2.40	0.44
1:A:91:ARG:NH1	1:A:98:VAL:O	2.48	0.44
1:L:288:ARG:HG2	2:K:38:SER:CB	2.47	0.44
2:K:17:ASP:O	2:K:18:ASP:CB	2.62	0.44
2:G:66:VAL:HG12	2:G:67:ILE:N	2.33	0.44
1:L:61:GLU:HB2	1:L:63:ARG:HG3	2.00	0.44
2:C:30:PHE:CD1	2:C:30:PHE:C	2.96	0.44
1:L:282:LEU:O	1:L:283:GLY:C	2.60	0.44
2:C:12:THR:HG22	2:C:22:VAL:HG12	2.00	0.44
1:A:273:ASN:O	1:A:274:LYS:HB2	2.18	0.43
1:L:4:THR:CG2	1:L:55:ARG:HG3	2.48	0.43
1:L:11:LYS:HG3	1:L:15:ASP:OD2	2.18	0.43
2:E:48:MET:HE3	2:E:48:MET:HB2	1.70	0.43
1:A:54:VAL:HG12	1:A:57:ILE:CD1	2.46	0.43
1:L:290:THR:CG2	2:K:34:TRP:HD1	2.31	0.43
2:B:63:PHE:CD1	2:B:63:PHE:C	2.95	0.43
2:F:1:THR:CG2	2:F:2:PRO:HD2	2.49	0.43
1:A:128:MET:HE1	1:A:159:LEU:HD13	1.99	0.43
2:J:4:CYS:O	2:J:4:CYS:SG	2.75	0.43
2:I:22:VAL:HG11	2:I:40:LEU:HD13	2.00	0.43
2:D:26:ASP:OD1	2:D:27:LYS:HD3	2.18	0.43
2:H:11:TYR:CD1	2:H:11:TYR:C	2.95	0.43
2:K:69:ARG:HG2	2:K:69:ARG:NH1	2.33	0.43
1:A:188:ARG:HH11	1:A:188:ARG:CA	2.31	0.43
1:L:32:SER:O	1:L:229:ILE:HD13	2.19	0.43
1:L:289:ARG:O	1:L:289:ARG:HG3	2.18	0.43
1:A:286:LEU:HD23	1:A:286:LEU:HA	1.86	0.43
2:B:29:LEU:HD13	2:B:52:ILE:HG21	2.01	0.43
1:L:57:ILE:O	1:L:57:ILE:CG2	2.67	0.42
2:G:37:GLN:OE1	2:K:35:ASN:ND2	2.38	0.42
1:A:266:ARG:HE	1:A:266:ARG:HB3	1.48	0.42
2:F:10:GLU:O	2:F:11:TYR:HB3	2.19	0.42
1:A:188:ARG:CB	1:A:188:ARG:NH1	2.79	0.42
1:L:139:TYR:CE1	1:L:143:MET:HG3	2.54	0.42
2:H:67:ILE:HG12	2:I:13:LYS:HG3	2.01	0.42
2:I:55:ASN:C	2:I:57:CYS:N	2.77	0.42
2:K:17:ASP:CG	2:K:19:THR:HG22	2.44	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:188:ARG:NH1	1:A:188:ARG:CA	2.82	0.42
1:L:60:GLU:CD	1:L:60:GLU:N	2.68	0.42
1:L:63:ARG:HB3	1:L:63:ARG:NH1	2.35	0.42
2:D:13:LYS:HB3	2:D:21:THR:HB	2.02	0.42
1:A:229:ILE:HD12	1:A:229:ILE:HA	1.86	0.42
1:L:89:PHE:C	1:L:89:PHE:CD1	2.97	0.42
1:A:30:THR:HB	1:A:31:ILE:HD12	2.00	0.42
2:F:2:PRO:HD2	2:F:55:ASN:ND2	2.35	0.42
2:J:42:SER:HB3	2:K:45:ILE:CD1	2.50	0.42
2:K:56:ALA:O	2:K:61:GLY:HA3	2.18	0.42
1:A:173:GLN:NE2	1:A:196:ASP:OD1	2.44	0.42
1:L:45:THR:HG23	1:L:45:THR:O	2.19	0.42
1:L:64:PHE:N	1:L:64:PHE:HD1	2.16	0.42
1:L:131:ASN:HB3	1:L:189:SER:HB3	2.00	0.42
2:E:48:MET:HE1	2:F:45:ILE:HG12	2.01	0.42
1:L:220:ARG:HD2	1:L:225:SER:HB2	2.00	0.42
1:A:139:TYR:CE1	1:A:143:MET:HG3	2.55	0.42
1:L:31:ILE:HD12	1:L:38:LEU:HD22	2.02	0.42
1:A:5:LEU:HD12	1:A:16:SER:HB3	2.02	0.42
2:C:32:ASN:C	2:C:32:ASN:HD22	2.28	0.42
2:D:58:HIS:O	2:D:59:ASN:C	2.63	0.41
2:E:35:ASN:HD21	2:K:34:TRP:HZ3	1.65	0.41
2:F:69:ARG:HG2	2:F:69:ARG:HH11	1.85	0.41
2:H:12:THR:HG22	2:H:22:VAL:HG23	2.03	0.41
1:A:38:LEU:HD21	1:A:239:ILE:HD13	2.01	0.41
1:A:187:GLY:O	1:A:188:ARG:O	2.39	0.41
1:A:285:ILE:HD12	2:F:45:ILE:HG22	2.02	0.41
2:B:18:ASP:OD2	2:J:33:ARG:CZ	2.68	0.41
2:E:53:LYS:O	2:E:54:THR:HB	2.20	0.41
2:C:65:GLU:HA	2:D:14:TYR:O	2.20	0.41
2:D:20:PHE:CD1	2:D:37:GLN:HG2	2.55	0.41
2:D:41:LEU:HA	2:D:41:LEU:HD12	1.79	0.41
2:K:12:THR:CG2	2:K:22:VAL:HG12	2.33	0.41
1:L:67:LEU:HD12	1:L:68:ARG:N	2.32	0.41
2:D:11:TYR:CD1	2:D:11:TYR:C	2.99	0.41
2:H:67:ILE:HG12	2:I:13:LYS:CG	2.51	0.41
1:A:55:ARG:HH22	1:A:84:ARG:HH12	1.68	0.41
1:A:63:ARG:O	1:A:65:ASN:N	2.50	0.41
2:J:14:TYR:CE1	2:J:18:ASP:HA	2.56	0.41
2:K:20:PHE:O	2:K:30:PHE:HA	2.21	0.41
1:L:290:THR:HG21	2:K:34:TRP:HD1	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:1:THR:HA	2:H:2:PRO:HD3	1.75	0.41
2:H:66:VAL:HG12	2:H:67:ILE:N	2.35	0.41
2:B:39:LEU:CD2	2:C:41:LEU:HD22	2.51	0.41
1:A:148:THR:HG22	1:A:149:SER:N	2.36	0.41
1:L:129:GLN:NE2	1:L:189:SER:OG	2.54	0.41
2:F:58:HIS:O	2:F:59:ASN:C	2.64	0.41
2:H:11:TYR:OH	2:H:28:GLU:OE1	2.37	0.41
1:A:57:ILE:H	1:A:57:ILE:CD1	2.34	0.41
1:A:131:ASN:HD22	1:A:189:SER:HB2	1.86	0.41
2:C:26:ASP:OD2	2:C:27:LYS:HE3	2.21	0.40
2:D:16:ASP:C	2:D:18:ASP:H	2.29	0.40
1:L:9:THR:C	1:L:11:LYS:N	2.76	0.40
2:B:13:LYS:CD	2:B:15:ASN:OD1	2.69	0.40
2:C:41:LEU:O	2:C:45:ILE:HG13	2.21	0.40
1:A:72:GLU:O	1:A:76:LEU:HA	2.22	0.40
2:J:4:CYS:O	2:J:57:CYS:SG	2.80	0.40
2:B:13:LYS:HE2	2:B:15:ASN:OD1	2.21	0.40
1:A:57:ILE:N	1:A:57:ILE:CD1	2.82	0.40
2:J:14:TYR:CZ	2:J:18:ASP:HA	2.56	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	272/293 (93%)	254 (93%)	12 (4%)	6 (2%)	5	9
1	L	273/293 (93%)	245 (90%)	23 (8%)	5 (2%)	6	12
2	B	67/69 (97%)	62 (92%)	4 (6%)	1 (2%)	8	16
2	C	67/69 (97%)	62 (92%)	5 (8%)	0	100	100
2	D	67/69 (97%)	60 (90%)	7 (10%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	E	67/69 (97%)	65 (97%)	2 (3%)	0	100	100
2	F	67/69 (97%)	61 (91%)	4 (6%)	2 (3%)	3	5
2	G	67/69 (97%)	64 (96%)	3 (4%)	0	100	100
2	H	67/69 (97%)	66 (98%)	1 (2%)	0	100	100
2	I	67/69 (97%)	63 (94%)	3 (4%)	1 (2%)	8	16
2	J	67/69 (97%)	62 (92%)	3 (4%)	2 (3%)	3	5
2	K	67/69 (97%)	60 (90%)	5 (8%)	2 (3%)	3	5
All	All	1215/1276 (95%)	1124 (92%)	72 (6%)	19 (2%)	7	14

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	64	PHE
1	A	188	ARG
1	A	189	SER
2	F	18	ASP
2	J	57	CYS
2	K	18	ASP
1	A	46	GLY
1	L	132	ARG
1	L	189	SER
2	B	57	CYS
2	F	17	ASP
1	A	264	ASP
1	L	42	ASP
1	L	84	ARG
2	K	56	ALA
1	L	109	SER
2	I	57	CYS
2	J	5	VAL
1	A	113	SER

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	238/252 (94%)	216 (91%)	22 (9%)	8	19
1	L	239/252 (95%)	218 (91%)	21 (9%)	9	20
2	B	61/61 (100%)	56 (92%)	5 (8%)	10	23
2	C	61/61 (100%)	56 (92%)	5 (8%)	10	23
2	D	61/61 (100%)	58 (95%)	3 (5%)	22	45
2	E	61/61 (100%)	59 (97%)	2 (3%)	33	61
2	F	61/61 (100%)	56 (92%)	5 (8%)	10	23
2	G	61/61 (100%)	55 (90%)	6 (10%)	7	16
2	H	61/61 (100%)	59 (97%)	2 (3%)	33	61
2	I	61/61 (100%)	57 (93%)	4 (7%)	15	32
2	J	61/61 (100%)	53 (87%)	8 (13%)	4	8
2	K	61/61 (100%)	57 (93%)	4 (7%)	15	32
All	All	1087/1114 (98%)	1000 (92%)	87 (8%)	11	24

All (87) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	GLU
1	A	4	THR
1	A	36	THR
1	A	55	ARG
1	A	61	GLU
1	A	85	THR
1	A	87	ASN
1	A	101	PRO
1	A	104	THR
1	A	119	ARG
1	A	124	SER
1	A	126	THR
1	A	148	THR
1	A	180	THR
1	A	188	ARG
1	A	189	SER
1	A	199	LEU
1	A	220	ARG
1	A	235	SER
1	A	274	LYS
1	A	278	ASP
1	A	282	LEU

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Mol	Chain	Res	Type
1	L	26	THR
1	L	58	ASP
1	L	60	GLU
1	L	61	GLU
1	L	64	PHE
1	L	85	THR
1	L	99	THR
1	L	124	SER
1	L	126	THR
1	L	129	GLN
1	L	135	LEU
1	L	148	THR
1	L	153	SER
1	L	162	VAL
1	L	185	LEU
1	L	188	ARG
1	L	193	THR
1	L	220	ARG
1	L	266	ARG
1	L	268	ARG
1	L	290	THR
2	B	5	VAL
2	B	6	THR
2	B	10	GLU
2	B	29	LEU
2	B	69	ARG
2	C	10	GLU
2	C	23	LYS
2	C	32	ASN
2	C	42	SER
2	C	49	THR
2	D	29	LEU
2	D	53	LYS
2	D	69	ARG
2	E	16	ASP
2	E	35	ASN
2	F	3	ASP
2	F	18	ASP
2	F	33	ARG
2	F	36	LEU
2	F	59	ASN
2	G	21	THR

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Mol	Chain	Res	Type
2	G	22	VAL
2	G	32	ASN
2	G	50	VAL
2	G	54	THR
2	G	69	ARG
2	H	1	THR
2	H	22	VAL
2	I	12	THR
2	I	27	LYS
2	I	29	LEU
2	I	38	SER
2	J	3	ASP
2	J	10	GLU
2	J	21	THR
2	J	35	ASN
2	J	41	LEU
2	J	49	THR
2	J	64	SER
2	J	69	ARG
2	K	34	TRP
2	K	36	LEU
2	K	38	SER
2	K	41	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	118	GLN
1	A	214	HIS
1	A	273	ASN
1	L	66	ASN
1	L	97	HIS
1	L	118	GLN
1	L	129	GLN
1	L	145	HIS
1	L	202	ASN
1	L	273	ASN
2	B	37	GLN
2	B	55	ASN
2	C	37	GLN
2	C	59	ASN
2	D	15	ASN

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Mol	Chain	Res	Type
2	D	55	ASN
2	E	35	ASN
2	E	37	GLN
2	F	35	ASN
2	F	55	ASN
2	G	32	ASN
2	G	58	HIS
2	H	44	GLN
2	H	55	ASN
2	I	37	GLN
2	I	55	ASN
2	J	58	HIS

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 [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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	276/293 (94%)	-0.47	5 (1%) 67 64	18, 37, 71, 86	0
1	L	277/293 (94%)	-0.25	5 (1%) 67 64	30, 49, 82, 93	0
2	B	69/69 (100%)	-0.53	0 100 100	28, 45, 56, 66	0
2	C	69/69 (100%)	-0.65	1 (1%) 73 70	26, 36, 49, 63	0
2	D	69/69 (100%)	-0.61	0 100 100	20, 37, 52, 60	0
2	E	69/69 (100%)	-0.88	0 100 100	15, 24, 40, 56	0
2	F	69/69 (100%)	-0.34	3 (4%) 40 35	26, 38, 67, 81	0
2	G	69/69 (100%)	-1.04	0 100 100	17, 22, 35, 43	0
2	H	69/69 (100%)	-0.77	0 100 100	19, 31, 43, 52	0
2	I	69/69 (100%)	-0.32	1 (1%) 73 70	33, 50, 64, 72	0
2	J	69/69 (100%)	-0.29	1 (1%) 73 70	33, 47, 62, 66	0
2	K	69/69 (100%)	-0.31	2 (2%) 53 49	23, 42, 75, 87	0
All	All	1243/1276 (97%)	-0.48	18 (1%) 73 70	15, 40, 68, 93	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	L	290	THR	4.8
2	K	34	TRP	4.3
2	K	14	TYR	3.2
2	F	18	ASP	3.2
2	F	16	ASP	3.1
1	A	185	LEU	3.0
1	A	289	ARG	2.7
1	A	45	THR	2.7
1	A	46	GLY	2.6
2	F	34	TRP	2.6
2	C	34	TRP	2.5

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Mol	Chain	Res	Type	RSRZ
2	J	34	TRP	2.3
1	A	43	SER	2.2
1	L	185	LEU	2.1
1	L	289	ARG	2.1
2	I	34	TRP	2.1
1	L	243	HIS	2.1
1	L	263	ALA	2.0

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.