



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 11:17 AM UTC

PDB ID : 1FL9 / pdb_00001fl9
Title : THE YJEE PROTEIN
Authors : Teplyakov, A.; Gilliland, G.L.; Structure 2 Function Project (S2F)
Deposited on : 2000-08-13
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) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

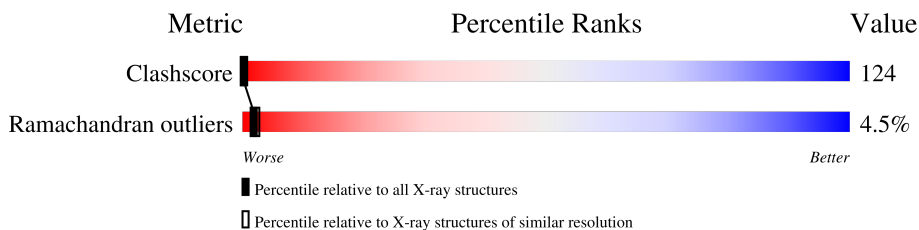
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 Å.

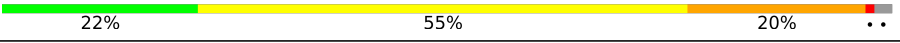
Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	161	
1	B	161	
1	C	161	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3799 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 HYPOTHETICAL PROTEIN HI0065.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	157	Total	C	N	O	S	0	0	0
			1255	806	201	241	7			
1	B	157	Total	C	N	O	S	0	0	0
			1255	806	201	241	7			
1	C	157	Total	C	N	O	S	0	0	0
			1255	806	201	241	7			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	SEE REMARK 999	UNP P44492
B	-1	SER	-	SEE REMARK 999	UNP P44492
C	0	HIS	-	SEE REMARK 999	UNP P44492

- Molecule 2 is MERCURY (II) ION (CCD ID: HG) (formula: Hg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Hg	0	0
			1	1		
2	B	1	Total	Hg	0	0
			1	1		
2	C	1	Total	Hg	0	0
			1	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	12	Total	O	0	0
			12	12		
3	B	7	Total	O	0	0
			7	7		

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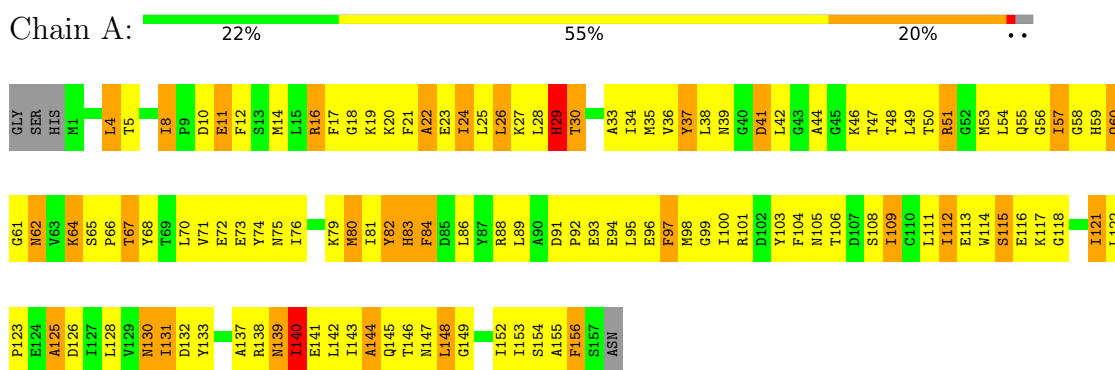
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	12	Total	O	0	0
			12	12		

3 Residue-property plots

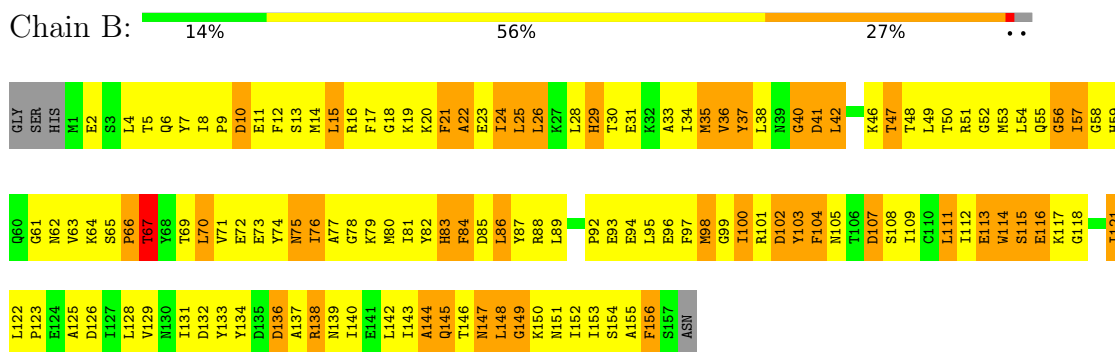
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

Note EDS was not executed.

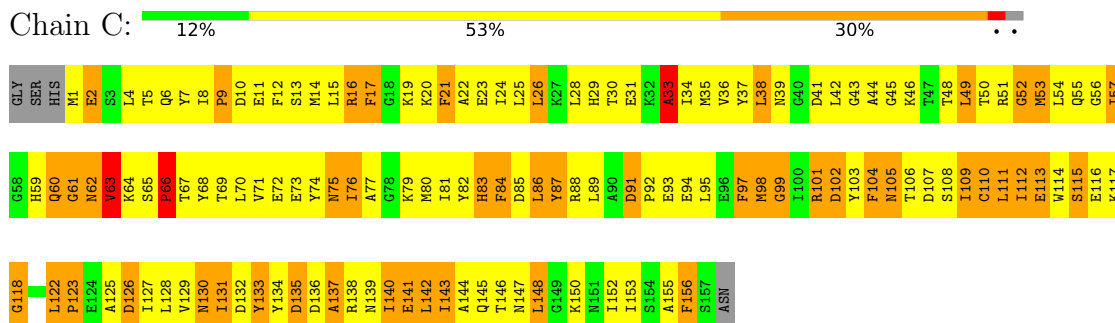
• Molecule 1: HYPOTHETICAL PROTEIN HI0065



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• Molecule 1: HYPOTHETICAL PROTEIN HI0065



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	76.40 Å 73.00 Å 96.80 Å 90.00° 110.20° 90.00°	Depositor
Resolution (Å)	10.00 – 2.50	Depositor
% Data completeness (in resolution range)	95.0 (10.00-2.50)	Depositor
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	REFMAC	Depositor
R, R_{free}	0.220 , 0.285	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3799	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: HG

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	0.93	0/1279	2.10	53/1726 (3.1%)
1	B	1.07	3/1279 (0.2%)	2.43	83/1726 (4.8%)
1	C	1.07	2/1279 (0.2%)	2.28	86/1726 (5.0%)
All	All	1.03	5/3837 (0.1%)	2.27	222/5178 (4.3%)

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	C	0	3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	113	GLU	CD-OE1	-7.01	1.12	1.25
1	B	66	PRO	N-CA	-5.36	1.40	1.47
1	B	100	ILE	N-CA	-5.33	1.40	1.46
1	B	25	LEU	N-CA	-5.07	1.40	1.46
1	C	113	GLU	CD-OE2	5.06	1.34	1.25

All (222) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	100	ILE	N-CA-C	17.80	128.70	110.72
1	B	102	ASP	CA-CB-CG	16.99	129.59	112.60
1	B	66	PRO	N-CA-C	15.84	133.66	113.86
1	A	156	PHE	CA-CB-CG	14.79	128.59	113.80
1	C	91	ASP	N-CA-C	14.00	120.82	108.22

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	21	PHE	CA-CB-CG	11.73	125.53	113.80
1	B	76	ILE	CB-CA-C	11.41	122.28	111.44
1	B	83	HIS	CA-CB-CG	11.06	124.86	113.80
1	A	16	ARG	CD-NE-CZ	11.04	139.85	124.40
1	C	136	ASP	N-CA-C	-10.41	100.89	114.31
1	A	16	ARG	N-CA-CB	10.30	124.95	109.91
1	B	33	ALA	N-CA-C	10.27	123.59	110.24
1	C	31	GLU	CA-C-O	9.71	130.08	119.40
1	B	65	SER	CA-C-N	9.64	129.30	119.56
1	B	65	SER	C-N-CA	9.64	129.30	119.56
1	C	141	GLU	N-CA-C	9.12	122.87	112.57
1	B	33	ALA	CA-C-N	9.07	134.79	121.18
1	B	33	ALA	C-N-CA	9.07	134.79	121.18
1	B	105	ASN	N-CA-C	8.99	129.95	110.80
1	B	75	ASN	CA-C-N	-8.96	114.03	123.08
1	B	75	ASN	C-N-CA	-8.96	114.03	123.08
1	B	99	GLY	CA-C-N	8.81	132.93	120.42
1	B	99	GLY	C-N-CA	8.81	132.93	120.42
1	C	101	ARG	N-CA-C	-8.77	101.30	111.03
1	B	24	ILE	CA-C-N	8.73	132.50	120.63
1	B	24	ILE	C-N-CA	8.73	132.50	120.63
1	C	17	PHE	CA-CB-CG	-8.54	105.26	113.80
1	B	156	PHE	CA-CB-CG	-8.50	105.30	113.80
1	B	103	TYR	N-CA-C	8.44	124.13	111.04
1	C	61	GLY	N-CA-C	-8.44	98.13	112.30
1	C	91	ASP	CA-C-N	8.10	127.78	119.19
1	C	91	ASP	C-N-CA	8.10	127.78	119.19
1	C	112	ILE	N-CA-C	8.08	121.30	108.85
1	B	25	LEU	N-CA-C	8.07	119.98	111.03
1	B	24	ILE	CB-CA-C	-8.06	101.49	112.04
1	C	140	ILE	N-CA-C	8.00	119.36	108.17
1	A	84	PHE	CA-CB-CG	-7.89	105.91	113.80
1	B	57	ILE	CB-CA-C	7.64	121.75	111.97
1	B	67	THR	N-CA-C	7.60	127.00	110.80
1	A	24	ILE	N-CA-C	-7.60	102.93	110.30
1	B	40	GLY	N-CA-C	7.57	122.19	110.91
1	C	136	ASP	CA-C-O	7.55	127.15	118.77
1	C	26	LEU	N-CA-C	-7.52	103.90	113.23
1	B	144	ALA	O-C-N	7.50	131.54	122.84
1	A	29	HIS	CA-CB-CG	7.43	121.23	113.80
1	C	123	PRO	N-CD-CG	7.43	114.34	103.20
1	A	109	ILE	CB-CA-C	-7.39	102.37	111.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	102	ASP	CA-CB-CG	7.39	119.99	112.60
1	C	111	LEU	CA-C-O	7.38	130.77	120.52
1	B	101	ARG	N-CA-CB	7.37	120.67	109.91
1	A	97	PHE	N-CA-C	7.36	120.75	111.69
1	A	125	ALA	N-CA-C	7.36	121.47	109.76
1	C	62	ASN	CA-C-N	-7.36	110.14	121.18
1	C	62	ASN	C-N-CA	-7.36	110.14	121.18
1	A	16	ARG	CG-CD-NE	7.33	128.13	112.00
1	A	140	ILE	CB-CA-C	-7.31	99.30	111.29
1	B	156	PHE	N-CA-C	7.30	121.09	111.75
1	C	85	ASP	CA-CB-CG	-7.20	105.40	112.60
1	C	63	VAL	N-CA-CB	7.13	118.74	110.82
1	C	97	PHE	CA-CB-CG	-7.12	106.68	113.80
1	A	121	ILE	N-CA-C	-7.11	103.25	111.00
1	C	60	GLN	OE1-CD-NE2	-7.10	115.50	122.60
1	A	39	ASN	N-CA-C	7.10	119.08	107.23
1	A	24	ILE	N-CA-CB	7.03	118.28	110.62
1	A	26	LEU	N-CA-C	6.90	121.15	112.87
1	C	135	ASP	O-C-N	6.90	128.55	121.79
1	B	121	ILE	CB-CA-C	-6.89	103.26	111.94
1	C	75	ASN	N-CA-CB	6.88	121.89	110.47
1	B	147	ASN	OD1-CG-ND2	6.83	129.43	122.60
1	B	114	TRP	N-CA-C	6.76	125.20	110.80
1	B	149	GLY	N-CA-C	-6.75	104.39	113.24
1	C	123	PRO	CA-N-CD	-6.73	102.58	112.00
1	C	140	ILE	CA-C-N	6.73	133.19	122.83
1	C	140	ILE	C-N-CA	6.73	133.19	122.83
1	B	78	GLY	N-CA-C	-6.69	106.22	114.92
1	A	62	ASN	CA-CB-CG	-6.64	105.96	112.60
1	C	33	ALA	CA-C-O	6.64	130.00	120.51
1	C	53	MET	CA-C-O	-6.64	114.03	121.00
1	B	12	PHE	CA-CB-CG	-6.63	107.17	113.80
1	C	66	PRO	N-CA-C	6.59	126.05	112.47
1	A	16	ARG	N-CA-C	-6.57	103.81	110.97
1	B	98	MET	N-CA-C	6.56	120.99	112.92
1	C	62	ASN	CA-C-O	-6.53	113.60	121.05
1	C	16	ARG	CD-NE-CZ	-6.53	115.26	124.40
1	C	66	PRO	CA-C-N	6.50	133.95	121.54
1	C	66	PRO	C-N-CA	6.50	133.95	121.54
1	B	26	LEU	N-CA-CB	6.46	120.30	110.28
1	B	10	ASP	CA-CB-CG	6.39	118.99	112.60
1	A	112	ILE	N-CA-C	6.35	118.62	108.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	105	ASN	N-CA-C	6.28	119.21	109.60
1	C	83	HIS	CA-CB-CG	-6.26	107.54	113.80
1	B	22	ALA	N-CA-C	6.25	120.11	112.23
1	A	96	GLU	CA-C-N	6.24	131.26	120.58
1	A	96	GLU	C-N-CA	6.24	131.26	120.58
1	B	131	ILE	N-CA-C	6.23	116.90	108.17
1	A	66	PRO	CA-C-N	6.21	130.35	121.71
1	A	66	PRO	C-N-CA	6.21	130.35	121.71
1	A	139	ASN	CA-C-O	-6.19	114.40	121.58
1	B	29	HIS	CA-CB-CG	-6.18	107.62	113.80
1	B	136	ASP	CB-CA-C	-6.17	101.02	111.26
1	A	51	ARG	CB-CA-C	-6.15	101.14	110.92
1	C	110	CYS	O-C-N	-6.13	116.23	123.22
1	C	137	ALA	N-CA-C	6.12	118.61	110.53
1	B	83	HIS	CA-C-O	6.10	126.99	120.46
1	B	42	LEU	N-CA-C	-6.10	101.15	110.30
1	A	130	ASN	OD1-CG-ND2	6.07	128.67	122.60
1	B	83	HIS	CB-CA-C	6.06	120.66	110.79
1	A	84	PHE	CB-CA-C	-6.03	100.85	110.14
1	C	84	PHE	O-C-N	6.02	130.31	123.27
1	A	97	PHE	CA-CB-CG	-6.01	107.78	113.80
1	B	47	THR	N-CA-CB	6.01	119.82	110.44
1	C	133	TYR	N-CA-C	6.01	118.54	109.23
1	C	65	SER	CA-C-N	5.99	127.33	119.84
1	C	65	SER	C-N-CA	5.99	127.33	119.84
1	A	109	ILE	N-CA-CB	5.98	119.39	112.15
1	A	84	PHE	N-CA-C	5.96	118.16	108.26
1	C	76	ILE	CB-CA-C	-5.93	101.67	110.33
1	C	61	GLY	CA-C-O	-5.93	115.22	122.75
1	C	112	ILE	N-CA-CB	-5.92	99.65	111.91
1	C	84	PHE	CA-C-N	-5.91	114.39	123.14
1	C	84	PHE	C-N-CA	-5.91	114.39	123.14
1	B	111	LEU	CA-C-N	-5.91	115.86	123.19
1	B	111	LEU	C-N-CA	-5.91	115.86	123.19
1	A	131	ILE	N-CA-C	5.90	116.11	107.37
1	A	104	PHE	CA-C-O	5.89	125.38	118.90
1	C	16	ARG	NE-CZ-NH1	-5.88	115.62	121.50
1	A	145	GLN	CB-CG-CD	-5.85	102.66	112.60
1	C	110	CYS	CA-C-N	5.82	131.33	122.95
1	C	110	CYS	C-N-CA	5.82	131.33	122.95
1	A	60	GLN	CB-CG-CD	5.80	122.46	112.60
1	C	122	LEU	CA-C-N	5.79	127.07	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	122	LEU	C-N-CA	5.79	127.07	119.84
1	B	35	MET	N-CA-C	5.78	119.14	109.95
1	C	99	GLY	CA-C-N	5.77	129.77	120.82
1	C	99	GLY	C-N-CA	5.77	129.77	120.82
1	C	86	LEU	N-CA-C	-5.74	106.18	112.72
1	B	86	LEU	CB-CA-C	5.74	118.80	111.40
1	C	155	ALA	N-CA-C	-5.71	105.20	111.82
1	A	57	ILE	CB-CA-C	-5.70	105.32	110.91
1	C	130	ASN	CA-CB-CG	5.70	118.30	112.60
1	C	61	GLY	CA-C-N	-5.69	113.29	121.42
1	C	61	GLY	C-N-CA	-5.69	113.29	121.42
1	B	102	ASP	CA-C-O	-5.69	112.36	118.79
1	B	37	TYR	CA-CB-CG	-5.68	103.67	113.90
1	B	62	ASN	CB-CA-C	5.67	119.09	109.80
1	C	126	ASP	N-CA-C	-5.66	104.12	111.02
1	C	68	TYR	CA-C-N	5.65	130.25	120.58
1	C	68	TYR	C-N-CA	5.65	130.25	120.58
1	B	70	LEU	CA-C-N	-5.65	115.14	122.99
1	B	70	LEU	C-N-CA	-5.65	115.14	122.99
1	C	84	PHE	CA-CB-CG	-5.65	108.15	113.80
1	B	104	PHE	O-C-N	-5.64	114.68	122.46
1	C	88	ARG	CD-NE-CZ	5.62	132.28	124.40
1	B	102	ASP	CB-CA-C	-5.59	100.80	109.80
1	B	84	PHE	CA-CB-CG	5.59	119.39	113.80
1	A	8	ILE	N-CA-CB	-5.57	105.62	111.36
1	A	82	TYR	N-CA-C	5.56	117.46	108.34
1	C	52	GLY	N-CA-C	-5.56	106.06	112.73
1	C	77	ALA	N-CA-C	5.56	122.64	110.80
1	B	65	SER	N-CA-C	5.52	116.92	109.24
1	B	136	ASP	N-CA-CB	5.50	121.30	110.63
1	A	144	ALA	N-CA-CB	-5.49	102.50	110.84
1	C	63	VAL	N-CA-C	-5.49	101.59	108.84
1	B	107	ASP	CA-CB-CG	5.48	118.08	112.60
1	C	49	LEU	N-CA-C	-5.47	105.22	111.07
1	B	105	ASN	O-C-N	-5.47	115.32	122.59
1	B	41	ASP	CA-C-N	-5.46	112.58	121.26
1	B	41	ASP	C-N-CA	-5.46	112.58	121.26
1	C	31	GLU	CB-CA-C	-5.42	104.01	112.31
1	B	136	ASP	CA-CB-CG	5.42	118.02	112.60
1	A	91	ASP	N-CA-C	5.39	116.73	109.24
1	A	57	ILE	N-CA-C	5.38	116.64	112.12
1	A	67	THR	CB-CA-C	-5.36	102.40	112.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	67	THR	N-CA-C	5.36	122.22	110.80
1	A	16	ARG	CA-CB-CG	5.36	124.81	114.10
1	B	100	ILE	CA-C-O	-5.35	115.18	120.85
1	A	51	ARG	CD-NE-CZ	5.32	131.85	124.40
1	A	36	VAL	N-CA-CB	5.32	119.25	112.34
1	C	88	ARG	CA-CB-CG	5.30	124.70	114.10
1	B	113	GLU	O-C-N	-5.30	116.50	123.02
1	A	80	MET	CA-C-O	-5.28	115.87	121.94
1	B	147	ASN	O-C-N	5.27	129.40	122.33
1	A	22	ALA	N-CA-C	5.27	117.11	111.36
1	B	15	LEU	CA-C-N	-5.27	113.22	120.28
1	B	15	LEU	C-N-CA	-5.27	113.22	120.28
1	A	28	LEU	CA-C-O	5.26	126.11	120.38
1	B	66	PRO	CA-C-O	-5.25	111.65	119.00
1	A	148	LEU	N-CA-C	-5.23	105.64	112.23
1	B	148	LEU	CA-C-O	5.23	125.72	119.60
1	C	105	ASN	CA-CB-CG	-5.23	107.37	112.60
1	A	83	HIS	CA-C-O	-5.22	115.47	121.47
1	C	21	PHE	N-CA-C	5.21	116.65	110.97
1	C	97	PHE	N-CA-C	5.20	117.74	111.40
1	B	42	LEU	CB-CA-C	5.19	117.27	110.06
1	C	135	ASP	CA-CB-CG	-5.19	107.41	112.60
1	A	37	TYR	N-CA-C	5.18	117.18	108.73
1	A	64	LYS	CB-CA-C	-5.17	101.83	109.90
1	B	145	GLN	CA-C-O	5.17	124.89	118.43
1	B	93	GLU	CA-C-N	5.12	128.10	120.31
1	B	93	GLU	C-N-CA	5.12	128.10	120.31
1	C	87	TYR	N-CA-C	5.12	116.67	111.14
1	B	134	TYR	CA-CB-CG	5.12	123.11	113.90
1	C	131	ILE	N-CA-C	5.11	119.97	109.34
1	C	2	GLU	CA-C-N	5.10	127.95	120.71
1	C	2	GLU	C-N-CA	5.10	127.95	120.71
1	B	138	ARG	CA-C-N	-5.09	115.82	123.00
1	B	138	ARG	C-N-CA	-5.09	115.82	123.00
1	C	134	TYR	CA-CB-CG	-5.09	104.74	113.90
1	B	26	LEU	CB-CA-C	-5.09	100.90	110.67
1	C	141	GLU	CB-CA-C	-5.08	103.18	111.06
1	A	30	THR	N-CA-C	5.07	117.41	108.75
1	B	56	GLY	CA-C-N	-5.07	114.18	120.56
1	B	56	GLY	C-N-CA	-5.07	114.18	120.56
1	A	41	ASP	N-CA-C	-5.06	102.79	110.28
1	C	104	PHE	CA-C-N	-5.06	113.75	121.74

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	104	PHE	C-N-CA	-5.06	113.75	121.74
1	C	148	LEU	N-CA-C	-5.05	106.97	113.23
1	B	36	VAL	N-CA-C	5.04	115.53	108.17
1	B	103	TYR	CA-CB-CG	-5.04	104.83	113.90
1	C	109	ILE	CA-C-O	-5.04	114.48	120.78
1	C	98	MET	CA-C-O	-5.02	115.53	120.70
1	C	76	ILE	CA-C-O	-5.00	115.06	120.36

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	101	ARG	Mainchain
1	C	143	ILE	Mainchain
1	C	63	VAL	Mainchain

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	1255	0	1244	254	0
1	B	1255	0	1244	383	0
1	C	1255	0	1244	344	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
3	A	12	0	0	1	0
3	B	7	0	0	1	0
3	C	12	0	0	3	0
All	All	3799	0	3732	928	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 124.

All (928) 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:79:LYS:HD3	1:A:109:ILE:CD1	1.66	1.24
1:B:89:LEU:HD11	1:B:95:LEU:CD2	1.67	1.23
1:B:21:PHE:HA	1:B:156:PHE:CZ	1.75	1.21
1:B:7:TYR:CE1	1:B:137:ALA:HB1	1.78	1.19
1:A:71:VAL:HG13	1:A:84:PHE:CE1	1.77	1.18
1:A:26:LEU:HA	1:A:79:LYS:CE	1.75	1.17
1:C:8:ILE:HD12	1:C:140:ILE:HG12	1.19	1.17
1:C:42:LEU:HD21	1:C:114:TRP:CZ2	1.80	1.16
1:C:57:ILE:CG2	1:C:76:ILE:HG21	1.74	1.15
1:B:121:ILE:HD12	1:B:121:ILE:N	1.61	1.15
1:A:16:ARG:NH2	1:C:135:ASP:HB2	1.59	1.15
1:B:121:ILE:H	1:B:121:ILE:CD1	1.54	1.14
1:B:20:LYS:O	1:B:24:ILE:HD13	1.45	1.14
1:C:8:ILE:CG2	1:C:138:ARG:HB2	1.78	1.13
1:C:35:MET:HE2	1:C:123:PRO:HD2	1.26	1.13
1:B:14:MET:HE2	1:B:49:LEU:HG	1.12	1.12
1:B:139:ASN:HD21	1:C:51:ARG:NH2	1.46	1.12
1:B:146:THR:HG22	1:B:147:ASN:N	1.61	1.11
1:B:37:TYR:CE1	1:B:125:ALA:HB2	1.87	1.10
1:B:7:TYR:HE1	1:B:137:ALA:HB1	0.99	1.09
1:C:35:MET:CE	1:C:104:PHE:HE1	1.64	1.09
1:A:16:ARG:HH21	1:C:135:ASP:HB2	1.11	1.09
1:C:94:GLU:HA	1:C:97:PHE:CD2	1.88	1.08
1:C:35:MET:HE1	1:C:104:PHE:HE1	1.15	1.08
1:C:57:ILE:HG22	1:C:76:ILE:HG21	1.13	1.08
1:C:8:ILE:HG22	1:C:138:ARG:HB2	1.29	1.07
1:B:69:THR:HG22	1:B:70:LEU:H	0.91	1.06
1:A:24:ILE:HD11	1:A:156:PHE:HD1	1.10	1.06
1:B:20:LYS:HD2	1:B:24:ILE:HD11	1.36	1.05
1:A:88:ARG:HG3	1:A:88:ARG:HH11	1.18	1.05
1:C:35:MET:CE	1:C:104:PHE:CE1	2.40	1.05
1:C:42:LEU:HD21	1:C:114:TRP:HZ2	0.91	1.05
1:B:64:LYS:HG3	1:B:72:GLU:HG2	1.39	1.04
1:B:95:LEU:HD13	1:B:122:LEU:CD1	1.87	1.04
1:B:82:TYR:HE2	1:B:103:TYR:HA	1.19	1.04
1:B:20:LYS:CD	1:B:24:ILE:HD11	1.88	1.04
1:A:73:GLU:HB3	1:A:82:TYR:CD1	1.93	1.03
1:A:24:ILE:HG21	1:A:155:ALA:CB	1.87	1.03
1:A:79:LYS:CD	1:A:109:ILE:HD11	1.89	1.03
1:A:146:THR:HG22	1:A:147:ASN:H	1.24	1.03
1:B:89:LEU:HD11	1:B:95:LEU:HD21	1.38	1.03
1:C:8:ILE:CD1	1:C:140:ILE:HG12	1.87	1.03

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:34:ILE:HG22	1:C:110:CYS:H	1.20	1.03
1:C:50:THR:HG22	1:C:54:LEU:HD13	1.38	1.02
1:B:95:LEU:HD13	1:B:122:LEU:HD12	1.40	1.02
1:B:6:GLN:H	1:B:140:ILE:HG22	1.19	1.02
1:C:131:ILE:HD13	1:C:131:ILE:N	1.72	1.02
1:A:35:MET:HE2	1:A:35:MET:N	1.75	1.01
1:B:146:THR:CG2	1:B:147:ASN:H	1.73	1.01
1:B:146:THR:HG22	1:B:147:ASN:H	1.10	1.01
1:C:35:MET:HE1	1:C:104:PHE:CE1	1.94	1.01
1:B:2:GLU:OE1	1:B:2:GLU:HA	1.59	1.01
1:A:17:PHE:CE2	1:A:156:PHE:HE2	1.78	1.00
1:A:24:ILE:HD11	1:A:156:PHE:CD1	1.95	1.00
1:B:82:TYR:CE2	1:B:103:TYR:HA	1.95	1.00
1:C:42:LEU:CD2	1:C:114:TRP:HZ2	1.74	1.00
1:B:118:GLY:HA3	1:B:122:LEU:HD12	1.43	1.00
1:A:26:LEU:CA	1:A:79:LYS:HE3	1.91	0.99
1:C:24:ILE:CD1	1:C:152:ILE:HG23	1.91	0.99
1:A:73:GLU:HB3	1:A:82:TYR:HD1	1.19	0.99
1:C:15:LEU:HD21	1:C:48:THR:HG22	1.42	0.99
1:B:89:LEU:O	1:B:117:LYS:HG2	1.62	0.98
1:A:89:LEU:HD11	1:A:94:GLU:HB2	1.45	0.98
1:C:15:LEU:HD21	1:C:48:THR:CG2	1.93	0.98
1:A:12:PHE:CE1	1:C:137:ALA:HB1	1.98	0.98
1:B:19:LYS:HE3	1:B:23:GLU:CD	1.88	0.98
1:A:17:PHE:CE2	1:A:156:PHE:CE2	2.52	0.97
1:B:133:TYR:HA	1:C:12:PHE:CZ	1.99	0.97
1:A:17:PHE:HE2	1:A:156:PHE:HE2	1.02	0.97
1:B:69:THR:CG2	1:B:70:LEU:H	1.77	0.97
1:C:29:HIS:ND1	1:C:30:THR:N	2.12	0.97
1:B:103:TYR:CD1	1:B:103:TYR:N	2.27	0.96
1:A:132:ASP:HB2	1:A:139:ASN:CB	1.95	0.96
1:B:125:ALA:O	1:B:145:GLN:NE2	1.98	0.96
1:C:99:GLY:HA2	1:C:103:TYR:HB2	1.46	0.96
1:A:33:ALA:HB1	1:A:108:SER:O	1.65	0.96
1:B:69:THR:HG22	1:B:70:LEU:N	1.73	0.96
1:B:148:LEU:HD23	1:B:149:GLY:H	1.30	0.96
1:C:41:ASP:HB3	1:C:133:TYR:HE1	1.31	0.95
1:B:7:TYR:HE1	1:B:137:ALA:CB	1.79	0.95
1:C:144:ALA:HB2	1:C:153:ILE:HG13	1.48	0.95
1:B:82:TYR:OH	1:B:102:ASP:HB2	1.66	0.95
1:B:95:LEU:N	1:B:95:LEU:HD23	1.80	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:21:PHE:CE1	1:C:25:LEU:HD11	2.01	0.95
1:B:139:ASN:HD21	1:C:51:ARG:HH22	1.15	0.94
1:A:71:VAL:HG22	1:A:84:PHE:CZ	2.02	0.94
1:B:139:ASN:ND2	1:C:51:ARG:NH2	2.14	0.94
1:A:24:ILE:HG21	1:A:155:ALA:HB1	1.47	0.94
1:A:26:LEU:HA	1:A:79:LYS:HE3	0.94	0.94
1:B:132:ASP:O	1:B:139:ASN:N	2.01	0.94
1:C:94:GLU:HA	1:C:97:PHE:HD2	1.28	0.94
1:B:66:PRO:HG3	1:B:72:GLU:HB2	1.49	0.93
1:A:67:THR:HG22	1:A:68:TYR:N	1.83	0.93
1:A:71:VAL:CG2	1:A:84:PHE:CZ	2.52	0.93
1:B:89:LEU:CD1	1:B:95:LEU:CD2	2.46	0.93
1:C:104:PHE:HZ	1:C:122:LEU:HD13	1.34	0.93
1:B:89:LEU:HD11	1:B:95:LEU:HD23	1.49	0.93
1:A:132:ASP:HB2	1:A:139:ASN:HB3	1.48	0.92
1:A:71:VAL:HG22	1:A:84:PHE:CE2	2.04	0.92
1:B:37:TYR:HD2	1:B:115:SER:HB3	1.33	0.92
1:B:7:TYR:CE1	1:B:137:ALA:CB	2.53	0.92
1:C:30:THR:HG21	1:C:107:ASP:H	1.33	0.92
1:B:148:LEU:HA	1:B:151:ASN:ND2	1.83	0.92
1:C:21:PHE:HE1	1:C:25:LEU:HD11	1.30	0.92
1:B:89:LEU:HD21	1:B:95:LEU:HD21	1.50	0.91
1:B:64:LYS:HG3	1:B:72:GLU:CG	2.00	0.91
1:A:71:VAL:CG1	1:A:84:PHE:CE1	2.53	0.91
1:A:79:LYS:HD3	1:A:109:ILE:HD11	0.92	0.91
1:C:37:TYR:CD1	1:C:125:ALA:HB2	2.05	0.91
1:B:37:TYR:CD2	1:B:115:SER:HB3	2.05	0.90
1:B:121:ILE:HD12	1:B:121:ILE:H	0.75	0.90
1:A:60:GLN:HA	1:C:6:GLN:NE2	1.87	0.90
1:C:57:ILE:HG22	1:C:76:ILE:CG2	2.00	0.90
1:C:50:THR:CG2	1:C:54:LEU:HD13	2.01	0.90
1:A:10:ASP:OD1	1:A:11:GLU:N	2.06	0.89
1:C:24:ILE:HD11	1:C:152:ILE:HG23	1.54	0.89
1:B:51:ARG:HA	1:B:63:VAL:HG11	1.55	0.89
1:C:144:ALA:HB2	1:C:153:ILE:CG1	2.03	0.89
1:B:26:LEU:HD22	1:B:77:ALA:HB3	1.55	0.88
1:B:103:TYR:HD1	1:B:103:TYR:H	1.20	0.88
1:B:95:LEU:CD1	1:B:122:LEU:HD12	2.02	0.88
1:B:6:GLN:HB2	1:B:140:ILE:CG2	2.04	0.88
1:B:67:THR:HB	1:B:85:ASP:OD2	1.72	0.88
1:B:7:TYR:O	1:B:9:PRO:HD3	1.74	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:14:MET:HE2	1:B:49:LEU:CG	2.01	0.88
1:B:136:ASP:H	1:C:16:ARG:HH12	1.22	0.88
1:C:15:LEU:CD2	1:C:48:THR:HG22	2.04	0.88
1:A:86:LEU:HD21	1:A:95:LEU:HD21	1.55	0.87
1:C:127:ILE:HG22	1:C:128:LEU:H	1.39	0.87
1:B:20:LYS:HD2	1:B:24:ILE:CD1	2.04	0.87
1:B:9:PRO:HG2	1:C:56:GLY:HA2	1.56	0.87
1:A:88:ARG:O	1:A:88:ARG:HD3	1.75	0.86
1:C:144:ALA:HB2	1:C:153:ILE:CD1	2.05	0.86
1:A:60:GLN:HA	1:C:6:GLN:HE22	1.40	0.86
1:A:86:LEU:CD2	1:A:89:LEU:HD22	2.05	0.86
1:C:70:LEU:O	1:C:71:VAL:CG2	2.23	0.86
1:B:6:GLN:HB2	1:B:140:ILE:HG21	1.58	0.85
1:A:35:MET:HB3	1:A:125:ALA:HA	1.58	0.85
1:A:71:VAL:HG13	1:A:84:PHE:HE1	1.40	0.85
1:A:25:LEU:HD23	1:A:34:ILE:HD13	1.57	0.85
1:C:125:ALA:HB1	1:C:127:ILE:O	1.77	0.85
1:C:15:LEU:HD23	1:C:15:LEU:N	1.92	0.85
1:C:41:ASP:HB3	1:C:133:TYR:CE1	2.11	0.85
1:A:24:ILE:CG2	1:A:155:ALA:HB1	2.07	0.84
1:B:103:TYR:N	1:B:103:TYR:HD1	1.73	0.84
1:A:25:LEU:CD2	1:A:34:ILE:HD13	2.06	0.84
1:B:21:PHE:HZ	1:B:36:VAL:HG11	1.40	0.83
1:B:6:GLN:N	1:B:140:ILE:HG22	1.92	0.83
1:C:6:GLN:HG3	1:C:17:PHE:CE1	2.13	0.83
1:A:118:GLY:HA3	1:A:122:LEU:CD1	2.08	0.83
1:C:79:LYS:HG2	1:C:109:ILE:HG13	1.61	0.83
1:B:54:LEU:HD12	1:B:74:TYR:CE2	2.14	0.83
1:B:139:ASN:ND2	1:C:51:ARG:HH22	1.76	0.83
1:B:133:TYR:O	1:B:133:TYR:CD1	2.32	0.82
1:C:1:MET:HB2	1:C:150:LYS:HE2	1.60	0.82
1:C:24:ILE:HD12	1:C:152:ILE:HG23	1.60	0.82
1:C:144:ALA:CB	1:C:153:ILE:CD1	2.57	0.82
1:B:18:GLY:O	1:B:53:MET:HG3	1.78	0.82
1:B:70:LEU:O	1:B:84:PHE:HA	1.80	0.82
1:B:8:ILE:HD12	1:B:14:MET:SD	2.20	0.82
1:B:2:GLU:HB2	1:B:144:ALA:CB	2.09	0.81
1:B:19:LYS:HB2	1:B:56:GLY:HA3	1.62	0.81
1:B:57:ILE:HG23	1:B:76:ILE:HG21	1.62	0.81
1:C:1:MET:CB	1:C:150:LYS:HE2	2.10	0.81
1:B:42:LEU:HG	1:B:114:TRP:HH2	1.45	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:21:PHE:HA	1:B:156:PHE:HZ	1.35	0.81
1:C:70:LEU:C	1:C:71:VAL:HG23	2.05	0.81
1:A:94:GLU:HA	1:A:97:PHE:CD2	2.15	0.81
1:B:133:TYR:HA	1:C:12:PHE:HZ	1.43	0.81
1:A:14:MET:HE1	1:A:140:ILE:HD11	1.61	0.81
1:B:97:PHE:N	1:B:97:PHE:HD1	1.79	0.81
1:B:95:LEU:HD23	1:B:95:LEU:H	1.43	0.81
1:B:19:LYS:HE3	1:B:23:GLU:OE2	1.80	0.80
1:B:89:LEU:CD1	1:B:95:LEU:HD21	2.10	0.80
1:C:59:HIS:CE1	1:C:64:LYS:HE3	2.17	0.80
1:B:92:PRO:HA	1:B:95:LEU:HG	1.64	0.80
1:B:14:MET:CE	1:B:49:LEU:HG	2.06	0.80
1:A:37:TYR:CE2	1:A:115:SER:HB2	2.17	0.80
1:B:24:ILE:CD1	1:B:24:ILE:H	1.95	0.80
1:A:33:ALA:CB	1:A:108:SER:O	2.30	0.79
1:C:82:TYR:HB3	1:C:84:PHE:HE1	1.47	0.79
1:B:53:MET:HE1	1:B:111:LEU:HD21	1.64	0.79
1:C:4:LEU:HD13	1:C:5:THR:H	1.46	0.79
1:C:94:GLU:CA	1:C:97:PHE:HD2	1.95	0.79
1:B:8:ILE:CD1	1:B:14:MET:SD	2.71	0.78
1:B:24:ILE:HG23	1:B:155:ALA:HB1	1.64	0.78
1:B:69:THR:HG22	1:B:71:VAL:H	1.48	0.78
1:B:9:PRO:CG	1:C:56:GLY:HA2	2.12	0.78
1:B:136:ASP:N	1:C:16:ARG:HH12	1.80	0.78
1:A:14:MET:HG2	1:A:48:THR:HG22	1.64	0.78
1:C:1:MET:HG2	1:C:145:GLN:HA	1.65	0.78
1:A:57:ILE:HG22	1:A:76:ILE:HG21	1.65	0.78
1:A:14:MET:HG2	1:A:48:THR:CG2	2.14	0.78
1:A:47:THR:CG2	1:A:48:THR:N	2.46	0.78
1:A:140:ILE:HG22	1:A:140:ILE:O	1.82	0.77
1:B:19:LYS:NZ	1:B:56:GLY:O	2.12	0.77
1:C:51:ARG:O	1:C:55:GLN:HB2	1.84	0.77
1:B:2:GLU:HB2	1:B:144:ALA:HB3	1.64	0.77
1:B:4:LEU:HD23	1:B:153:ILE:HG23	1.65	0.77
1:B:148:LEU:HD23	1:B:149:GLY:N	1.99	0.77
1:B:136:ASP:H	1:C:16:ARG:NH1	1.83	0.77
1:C:37:TYR:CE1	1:C:125:ALA:HB2	2.20	0.77
1:B:82:TYR:CE2	1:B:103:TYR:CG	2.73	0.77
1:B:89:LEU:HD12	1:B:94:GLU:HB2	1.67	0.77
1:C:53:MET:SD	1:C:111:LEU:HD11	2.25	0.77
1:B:4:LEU:HD21	1:C:62:ASN:HD21	1.50	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:35:MET:CE	1:C:123:PRO:HD2	2.10	0.77
1:A:57:ILE:CG2	1:A:76:ILE:HG21	2.16	0.76
1:C:109:ILE:O	1:C:109:ILE:HG22	1.85	0.76
1:B:54:LEU:HD12	1:B:74:TYR:CD2	2.20	0.76
1:B:25:LEU:HD13	1:B:109:ILE:HD13	1.65	0.76
1:A:79:LYS:CD	1:A:109:ILE:CD1	2.58	0.76
1:C:128:LEU:O	1:C:142:LEU:HA	1.84	0.76
1:B:20:LYS:HD3	1:B:24:ILE:HD11	1.67	0.76
1:B:34:ILE:HG22	1:B:34:ILE:O	1.85	0.76
1:C:104:PHE:CZ	1:C:122:LEU:HD13	2.20	0.76
1:A:19:LYS:HG3	1:A:56:GLY:HA3	1.68	0.76
1:C:30:THR:CG2	1:C:107:ASP:H	1.97	0.76
1:C:34:ILE:CG2	1:C:110:CYS:H	1.96	0.76
1:C:84:PHE:CE2	1:C:98:MET:HE3	2.20	0.76
1:A:19:LYS:HG3	1:A:56:GLY:CA	2.16	0.76
1:B:25:LEU:HD13	1:B:109:ILE:CD1	2.16	0.75
1:B:97:PHE:N	1:B:97:PHE:CD1	2.54	0.75
1:C:70:LEU:O	1:C:71:VAL:HG23	1.86	0.75
1:B:118:GLY:HA3	1:B:122:LEU:CD1	2.16	0.75
1:C:41:ASP:O	1:C:44:ALA:N	2.19	0.75
1:B:2:GLU:HG2	1:B:150:LYS:HG3	1.69	0.75
1:C:15:LEU:HA	1:C:52:GLY:HA3	1.69	0.75
1:C:144:ALA:CB	1:C:153:ILE:HG13	2.15	0.75
1:C:30:THR:HG21	1:C:107:ASP:N	2.00	0.75
1:A:8:ILE:HG22	1:A:8:ILE:O	1.87	0.74
1:B:82:TYR:CE2	1:B:103:TYR:CA	2.70	0.74
1:A:47:THR:HG22	1:A:48:THR:H	1.52	0.74
1:B:14:MET:HE3	1:B:14:MET:HA	1.69	0.74
1:B:26:LEU:HD23	1:B:79:LYS:HB2	1.69	0.74
1:A:118:GLY:HA3	1:A:122:LEU:HD12	1.69	0.74
1:B:95:LEU:HD13	1:B:122:LEU:HD11	1.67	0.74
1:C:8:ILE:HD12	1:C:140:ILE:CG1	2.09	0.74
1:B:84:PHE:HD1	1:B:84:PHE:H	1.35	0.74
1:B:143:ILE:HG22	1:B:144:ALA:O	1.87	0.74
1:B:2:GLU:O	1:B:143:ILE:HA	1.87	0.74
1:C:29:HIS:CG	1:C:30:THR:H	2.06	0.74
1:B:54:LEU:CD1	1:B:74:TYR:CD2	2.71	0.73
1:A:17:PHE:HE2	1:A:156:PHE:CE2	1.92	0.73
1:C:1:MET:CG	1:C:145:GLN:HA	2.18	0.73
1:A:37:TYR:CD1	1:A:125:ALA:HB2	2.24	0.73
1:B:21:PHE:CA	1:B:156:PHE:CZ	2.65	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:131:ILE:HD12	1:C:140:ILE:HD12	1.70	0.73
1:B:7:TYR:OH	1:B:137:ALA:HB2	1.88	0.73
1:B:19:LYS:HD2	1:B:56:GLY:CA	2.19	0.73
1:B:151:ASN:O	1:B:155:ALA:HB2	1.88	0.73
1:C:148:LEU:C	1:C:148:LEU:HD23	2.14	0.73
1:B:89:LEU:CD2	1:B:95:LEU:HD21	2.18	0.73
1:B:133:TYR:HB2	1:B:138:ARG:HG3	1.71	0.73
1:A:35:MET:CB	1:A:125:ALA:HA	2.18	0.73
1:C:35:MET:HE2	1:C:104:PHE:CE1	2.21	0.73
1:A:144:ALA:HB1	1:A:149:GLY:C	2.13	0.72
1:B:19:LYS:HB2	1:B:52:GLY:O	1.89	0.72
1:A:4:LEU:HD12	1:A:153:ILE:HG23	1.70	0.72
1:A:24:ILE:CD1	1:A:156:PHE:HD1	1.98	0.72
1:B:54:LEU:HD23	1:B:54:LEU:N	2.03	0.72
1:C:1:MET:HB2	1:C:144:ALA:O	1.87	0.72
1:B:95:LEU:HD12	1:B:121:ILE:HD13	1.70	0.72
1:A:25:LEU:HD21	1:A:34:ILE:HG21	1.72	0.72
1:A:101:ARG:HB2	1:A:103:TYR:CE1	2.25	0.72
1:B:151:ASN:O	1:B:155:ALA:CB	2.38	0.72
1:B:84:PHE:CD1	1:B:84:PHE:N	2.55	0.71
1:C:83:HIS:HD2	1:C:84:PHE:N	1.88	0.71
1:C:91:ASP:HB2	1:C:92:PRO:HD2	1.73	0.71
1:A:95:LEU:O	1:A:98:MET:HE3	1.91	0.71
1:C:129:VAL:HG12	1:C:131:ILE:HD11	1.72	0.71
1:C:102:ASP:O	1:C:103:TYR:CD1	2.44	0.70
1:C:49:LEU:O	1:C:53:MET:CB	2.39	0.70
1:A:35:MET:HE2	1:A:35:MET:CA	2.21	0.70
1:C:94:GLU:HA	1:C:97:PHE:CE2	2.26	0.70
1:A:83:HIS:NE2	1:A:113:GLU:OE2	2.25	0.70
1:C:8:ILE:HG21	1:C:138:ARG:HB2	1.73	0.70
1:C:84:PHE:HE2	1:C:98:MET:HE3	1.55	0.70
1:C:103:TYR:O	1:C:110:CYS:SG	2.50	0.70
1:C:70:LEU:O	1:C:71:VAL:HG22	1.91	0.70
1:C:8:ILE:HG22	1:C:8:ILE:O	1.90	0.70
1:B:24:ILE:HD13	1:B:24:ILE:H	1.56	0.69
1:C:83:HIS:HD2	1:C:84:PHE:H	1.39	0.69
1:C:144:ALA:HB2	1:C:153:ILE:HD11	1.74	0.69
1:A:12:PHE:CD1	1:C:137:ALA:HB1	2.28	0.69
1:B:37:TYR:CE1	1:B:125:ALA:CB	2.73	0.69
1:B:21:PHE:HD1	1:B:156:PHE:HE2	1.40	0.69
1:A:118:GLY:HA3	1:A:122:LEU:HD13	1.75	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:89:LEU:CD1	1:B:94:GLU:HB2	2.23	0.69
1:C:66:PRO:HG3	1:C:72:GLU:HB2	1.73	0.69
1:B:104:PHE:CE2	1:B:123:PRO:CD	2.76	0.69
1:C:75:ASN:C	1:C:76:ILE:HD12	2.18	0.69
1:C:84:PHE:HB3	1:C:86:LEU:HD13	1.73	0.69
1:C:46:LYS:HD2	1:C:113:GLU:HG2	1.74	0.69
1:C:62:ASN:N	1:C:62:ASN:HD22	1.90	0.68
1:B:82:TYR:CE2	1:B:103:TYR:CD2	2.82	0.68
1:B:95:LEU:HA	1:B:98:MET:HE2	1.74	0.68
1:C:74:TYR:O	1:C:80:MET:HA	1.93	0.68
1:B:14:MET:HB3	1:B:48:THR:HG22	1.76	0.68
1:A:19:LYS:NZ	1:A:56:GLY:O	2.26	0.68
1:A:37:TYR:CE2	1:A:115:SER:CB	2.77	0.68
1:C:84:PHE:HZ	1:C:103:TYR:CD2	2.09	0.68
1:B:7:TYR:HB3	1:C:55:GLN:HB3	1.76	0.68
1:B:8:ILE:HG22	1:B:8:ILE:O	1.93	0.68
1:B:21:PHE:CA	1:B:156:PHE:HZ	2.03	0.68
1:B:73:GLU:HG2	1:B:82:TYR:CE1	2.28	0.68
1:B:86:LEU:HD13	1:B:122:LEU:CD1	2.24	0.68
1:B:104:PHE:CE2	1:B:123:PRO:HD3	2.28	0.68
1:C:57:ILE:HG21	1:C:76:ILE:HG21	1.73	0.68
1:C:70:LEU:C	1:C:71:VAL:CG2	2.66	0.68
1:A:50:THR:HG21	1:A:83:HIS:CE1	2.28	0.68
1:A:74:TYR:CD1	1:A:81:ILE:HG22	2.28	0.68
1:C:11:GLU:HA	1:C:138:ARG:HH12	1.58	0.68
1:C:131:ILE:CD1	1:C:131:ILE:H	2.03	0.68
1:B:136:ASP:CA	1:C:16:ARG:HH12	2.06	0.68
1:C:8:ILE:CD1	1:C:140:ILE:CG1	2.69	0.68
1:A:4:LEU:HD23	1:A:5:THR:N	2.08	0.68
1:A:71:VAL:HG11	1:A:103:TYR:HE2	1.58	0.68
1:B:2:GLU:O	1:B:143:ILE:HG23	1.94	0.68
1:A:19:LYS:CG	1:A:56:GLY:HA3	2.23	0.67
1:A:61:GLY:H	1:C:6:GLN:HE21	1.40	0.67
1:C:49:LEU:HD22	1:C:53:MET:CE	2.24	0.67
1:A:24:ILE:HG21	1:A:155:ALA:HB3	1.73	0.67
1:A:16:ARG:NH2	1:C:135:ASP:CB	2.50	0.67
1:C:144:ALA:HB3	1:C:153:ILE:HD12	1.77	0.67
1:C:5:THR:HA	1:C:140:ILE:O	1.95	0.67
1:A:118:GLY:CA	1:A:122:LEU:HD12	2.26	0.66
1:B:26:LEU:O	1:B:79:LYS:NZ	2.28	0.66
1:B:86:LEU:HD13	1:B:122:LEU:HD11	1.76	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:89:LEU:HD21	1:B:95:LEU:CD2	2.25	0.66
1:C:127:ILE:HG22	1:C:128:LEU:N	2.09	0.66
1:B:35:MET:CE	1:B:112:ILE:HD11	2.25	0.66
1:A:71:VAL:HG21	1:A:84:PHE:CZ	2.29	0.66
1:A:88:ARG:HH11	1:A:88:ARG:CG	2.01	0.66
1:B:37:TYR:CD2	1:B:115:SER:CB	2.78	0.66
1:A:10:ASP:CG	1:A:11:GLU:H	2.02	0.66
1:C:35:MET:HE2	1:C:123:PRO:CD	2.15	0.66
1:C:128:LEU:HB2	1:C:143:ILE:O	1.96	0.66
1:A:146:THR:HG22	1:A:147:ASN:N	2.04	0.66
1:B:4:LEU:N	1:B:142:LEU:O	2.29	0.66
1:B:25:LEU:CD1	1:B:109:ILE:HD13	2.26	0.66
1:B:25:LEU:HD21	1:B:34:ILE:HG21	1.76	0.66
1:C:148:LEU:HD23	1:C:148:LEU:O	1.96	0.66
1:B:24:ILE:CD1	1:B:24:ILE:N	2.58	0.66
1:B:5:THR:HG23	1:B:5:THR:O	1.95	0.66
1:B:137:ALA:H	1:C:16:ARG:NH1	1.94	0.66
1:A:132:ASP:HB2	1:A:139:ASN:HB2	1.78	0.65
1:B:149:GLY:HA2	1:B:152:ILE:HD12	1.79	0.65
1:C:142:LEU:N	1:C:142:LEU:HD23	2.11	0.65
1:B:57:ILE:CG2	1:B:76:ILE:HG21	2.26	0.65
1:C:82:TYR:HB3	1:C:84:PHE:CE1	2.29	0.65
1:A:74:TYR:CD1	1:A:74:TYR:N	2.65	0.65
1:C:36:VAL:C	1:C:37:TYR:HD1	2.05	0.65
1:C:102:ASP:HA	1:C:105:ASN:ND2	2.12	0.65
1:A:89:LEU:HD11	1:A:94:GLU:CB	2.24	0.65
1:B:37:TYR:CE2	1:B:115:SER:CB	2.79	0.65
1:B:84:PHE:HD1	1:B:84:PHE:N	1.90	0.65
1:A:46:LYS:HB2	3:A:206:HOH:O	1.97	0.65
1:C:6:GLN:O	1:C:139:ASN:HA	1.97	0.65
1:C:53:MET:SD	1:C:111:LEU:CD1	2.84	0.65
1:C:89:LEU:HD23	1:C:117:LYS:HB3	1.79	0.65
1:C:106:THR:HG23	1:C:107:ASP:OD1	1.97	0.65
1:A:34:ILE:C	1:A:35:MET:HE2	2.22	0.64
1:B:11:GLU:O	1:B:15:LEU:HD12	1.96	0.64
1:B:6:GLN:HG3	1:B:17:PHE:CE1	2.33	0.64
1:C:4:LEU:CD1	1:C:5:THR:H	2.09	0.64
1:C:35:MET:HG2	1:C:123:PRO:CG	2.27	0.64
1:A:34:ILE:HD12	1:A:109:ILE:HG12	1.78	0.64
1:A:86:LEU:HD21	1:A:89:LEU:HD22	1.78	0.64
1:A:89:LEU:O	1:A:117:LYS:HG2	1.97	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:21:PHE:HE2	1:A:142:LEU:HD21	1.63	0.64
1:C:34:ILE:HA	1:C:126:ASP:OD2	1.98	0.64
1:C:146:THR:HG22	1:C:147:ASN:N	2.11	0.64
1:A:12:PHE:CE1	1:C:137:ALA:CB	2.77	0.64
1:A:57:ILE:CG2	1:A:76:ILE:CG2	2.76	0.64
1:A:115:SER:O	1:A:115:SER:OG	2.15	0.64
1:B:57:ILE:HG23	1:B:76:ILE:CG2	2.27	0.64
1:C:14:MET:HG2	1:C:138:ARG:NH1	2.13	0.64
1:C:144:ALA:HB3	1:C:153:ILE:CD1	2.28	0.64
1:A:35:MET:HB3	1:A:125:ALA:CA	2.28	0.63
1:C:49:LEU:O	1:C:53:MET:HB3	1.98	0.63
1:C:89:LEU:HD22	1:C:95:LEU:HD21	1.80	0.63
1:B:6:GLN:HG3	1:B:17:PHE:HE1	1.62	0.63
1:A:89:LEU:CD1	1:A:94:GLU:HB2	2.27	0.63
1:B:40:GLY:HA3	1:B:46:LYS:HD3	1.80	0.63
1:B:115:SER:OG	1:B:116:GLU:N	2.31	0.63
1:A:4:LEU:O	1:A:141:GLU:HA	1.98	0.63
1:C:35:MET:HG2	1:C:123:PRO:HG2	1.81	0.63
1:B:20:LYS:O	1:B:24:ILE:CD1	2.35	0.62
1:B:19:LYS:CE	1:B:23:GLU:OE2	2.47	0.62
1:B:126:ASP:HB3	1:B:146:THR:OG1	1.99	0.62
1:B:146:THR:HG21	1:B:148:LEU:HD22	1.79	0.62
1:A:8:ILE:O	1:A:8:ILE:CG2	2.48	0.62
1:A:8:ILE:CG2	1:A:138:ARG:HB2	2.28	0.62
1:B:148:LEU:O	1:B:152:ILE:HG13	1.98	0.62
1:C:21:PHE:CD1	1:C:21:PHE:C	2.77	0.62
1:B:37:TYR:CD1	1:B:125:ALA:HB2	2.34	0.62
1:B:139:ASN:HD21	1:C:51:ARG:HH21	1.43	0.62
1:C:41:ASP:CB	1:C:133:TYR:HE1	2.11	0.62
1:B:24:ILE:CG2	1:B:155:ALA:HB1	2.29	0.62
1:B:55:GLN:HG2	1:B:59:HIS:O	1.99	0.62
1:B:95:LEU:CD2	1:B:95:LEU:N	2.57	0.62
1:A:17:PHE:CE2	1:A:156:PHE:CZ	2.87	0.62
1:B:82:TYR:CZ	1:B:103:TYR:CD2	2.88	0.62
1:A:47:THR:CG2	1:A:48:THR:H	2.09	0.62
1:B:21:PHE:CZ	1:B:36:VAL:HG11	2.31	0.62
1:B:146:THR:HB	1:B:148:LEU:HD23	1.80	0.62
1:C:57:ILE:CG2	1:C:76:ILE:CG2	2.67	0.62
1:A:75:ASN:OD1	1:A:75:ASN:C	2.43	0.62
1:B:100:ILE:HD13	1:B:100:ILE:N	2.14	0.62
1:C:74:TYR:N	1:C:74:TYR:CD1	2.68	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:2:GLU:O	1:C:153:ILE:CD1	2.48	0.61
1:A:59:HIS:HE1	1:A:62:ASN:O	1.82	0.61
1:A:117:LYS:N	1:A:117:LYS:HD2	2.16	0.61
1:B:6:GLN:HB2	1:B:140:ILE:HG22	1.82	0.61
1:B:152:ILE:HA	1:B:155:ALA:HB3	1.83	0.61
1:C:60:GLN:H	1:C:60:GLN:NE2	1.98	0.61
1:B:37:TYR:CD1	1:B:125:ALA:CB	2.83	0.61
1:C:24:ILE:HG13	1:C:152:ILE:HD12	1.82	0.61
1:A:60:GLN:CD	1:C:20:LYS:HD3	2.26	0.61
1:B:2:GLU:HB2	1:B:144:ALA:HB2	1.83	0.61
1:B:8:ILE:N	1:B:138:ARG:O	2.30	0.61
1:C:83:HIS:CD2	1:C:84:PHE:N	2.66	0.61
1:C:126:ASP:O	1:C:145:GLN:N	2.30	0.61
1:B:25:LEU:HD21	1:B:34:ILE:HD13	1.83	0.61
1:A:47:THR:HG23	1:A:48:THR:N	2.14	0.61
1:A:121:ILE:HG22	1:A:122:LEU:N	2.16	0.61
1:C:144:ALA:CB	1:C:153:ILE:HD11	2.29	0.61
1:A:64:LYS:CB	1:A:72:GLU:OE1	2.49	0.61
1:B:151:ASN:O	1:B:155:ALA:N	2.30	0.60
1:A:88:ARG:HG3	1:A:88:ARG:NH1	1.97	0.60
1:B:29:HIS:CG	1:B:30:THR:N	2.67	0.60
1:C:95:LEU:HD11	1:C:118:GLY:HA2	1.82	0.60
1:A:57:ILE:HG22	1:A:76:ILE:CG2	2.30	0.60
1:A:61:GLY:H	1:C:6:GLN:NE2	1.98	0.60
1:B:26:LEU:HD22	1:B:77:ALA:CB	2.30	0.60
1:B:126:ASP:O	1:B:145:GLN:N	2.34	0.60
1:C:132:ASP:O	1:C:138:ARG:HA	2.01	0.60
1:A:24:ILE:CD1	1:A:156:PHE:CD1	2.78	0.60
1:B:19:LYS:HE3	1:B:23:GLU:OE1	2.01	0.60
1:B:36:VAL:HG12	1:B:129:VAL:HG23	1.83	0.60
1:B:69:THR:HG22	1:B:71:VAL:N	2.15	0.60
1:B:88:ARG:HG3	1:B:88:ARG:HH11	1.67	0.60
1:C:153:ILE:O	1:C:156:PHE:HB2	2.01	0.60
1:B:2:GLU:OE1	1:B:2:GLU:CA	2.41	0.60
1:B:82:TYR:HE2	1:B:103:TYR:CA	2.05	0.60
1:B:137:ALA:O	1:C:12:PHE:HE1	1.85	0.60
1:A:71:VAL:CG2	1:A:84:PHE:CE2	2.80	0.60
1:B:20:LYS:O	1:B:21:PHE:C	2.44	0.60
1:B:133:TYR:CB	1:B:138:ARG:HG3	2.31	0.60
1:C:84:PHE:HB2	1:C:112:ILE:HG12	1.84	0.60
1:A:25:LEU:CD2	1:A:34:ILE:HG21	2.32	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:53:MET:C	1:B:54:LEU:HD23	2.27	0.59
1:B:73:GLU:HG2	1:B:82:TYR:HE1	1.67	0.59
1:A:71:VAL:HG11	1:A:103:TYR:CE2	2.37	0.59
1:C:94:GLU:CA	1:C:97:PHE:CD2	2.71	0.59
1:C:129:VAL:CG1	1:C:131:ILE:HD11	2.32	0.59
1:A:5:THR:HB	1:A:141:GLU:HG3	1.82	0.59
1:A:24:ILE:HD13	1:A:155:ALA:CB	2.32	0.59
1:C:28:LEU:HD13	1:C:152:ILE:CD1	2.32	0.59
1:C:112:ILE:HD11	1:C:122:LEU:HD12	1.83	0.59
1:A:24:ILE:CG2	1:A:155:ALA:CB	2.68	0.59
1:A:34:ILE:HG12	1:A:148:LEU:HD21	1.84	0.59
1:B:82:TYR:CE1	1:B:103:TYR:CE2	2.91	0.59
1:C:8:ILE:HB	1:C:138:ARG:O	2.03	0.59
1:B:84:PHE:CZ	1:B:104:PHE:HE1	2.19	0.59
1:B:143:ILE:CG2	1:B:144:ALA:N	2.66	0.59
1:B:19:LYS:HD2	1:B:56:GLY:C	2.27	0.59
1:B:82:TYR:CD2	1:B:103:TYR:CD2	2.90	0.59
1:C:80:MET:SD	1:C:82:TYR:OH	2.56	0.58
1:A:88:ARG:O	1:A:88:ARG:CD	2.51	0.58
1:A:121:ILE:CG2	1:A:122:LEU:N	2.67	0.58
1:C:59:HIS:HE1	1:C:64:LYS:HE3	1.64	0.58
1:A:14:MET:CG	1:A:48:THR:HG22	2.33	0.58
1:B:35:MET:HE2	1:B:112:ILE:HD11	1.85	0.58
1:B:86:LEU:CD1	1:B:122:LEU:CD1	2.81	0.58
1:A:12:PHE:CD1	1:C:137:ALA:CB	2.86	0.58
1:C:35:MET:HB3	1:C:125:ALA:HA	1.85	0.58
1:C:37:TYR:OH	1:C:123:PRO:O	2.21	0.58
1:C:131:ILE:CD1	1:C:140:ILE:HD12	2.34	0.58
1:B:19:LYS:HA	1:B:53:MET:HA	1.86	0.58
1:B:25:LEU:HD22	1:B:25:LEU:O	2.04	0.58
1:B:41:ASP:O	1:B:42:LEU:C	2.46	0.58
1:A:24:ILE:HD13	1:A:155:ALA:HB3	1.86	0.58
1:C:73:GLU:HA	1:C:81:ILE:O	2.04	0.58
1:B:21:PHE:HD1	1:B:156:PHE:CE2	2.22	0.57
1:A:14:MET:O	1:A:18:GLY:N	2.37	0.57
1:C:73:GLU:C	1:C:74:TYR:CD1	2.82	0.57
1:B:86:LEU:CD1	1:B:122:LEU:HD11	2.34	0.57
1:C:62:ASN:N	1:C:62:ASN:ND2	2.52	0.57
1:A:4:LEU:C	1:A:4:LEU:CD2	2.77	0.57
1:A:4:LEU:HD23	1:A:4:LEU:C	2.30	0.57
1:A:139:ASN:O	1:A:140:ILE:HG13	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:143:ILE:HG22	1:B:144:ALA:N	2.18	0.57
1:C:1:MET:HG2	1:C:145:GLN:HG3	1.86	0.57
1:B:24:ILE:N	1:B:24:ILE:HD12	2.18	0.57
1:C:2:GLU:O	1:C:153:ILE:HD12	2.04	0.57
1:A:64:LYS:HB2	1:A:72:GLU:OE1	2.04	0.57
1:A:132:ASP:CB	1:A:139:ASN:HB3	2.30	0.57
1:B:10:ASP:OD1	1:B:11:GLU:N	2.35	0.57
1:C:64:LYS:H	1:C:64:LYS:HD2	1.70	0.57
1:A:29:HIS:HE1	1:A:106:THR:HG23	1.69	0.57
1:B:104:PHE:HE2	1:B:123:PRO:HD3	1.68	0.57
1:B:139:ASN:ND2	1:C:51:ARG:HH21	1.97	0.57
1:C:8:ILE:N	1:C:138:ARG:O	2.38	0.57
1:B:75:ASN:HD21	1:B:80:MET:HE2	1.69	0.56
1:B:37:TYR:CD1	1:B:37:TYR:N	2.74	0.56
1:B:115:SER:O	1:B:117:LYS:N	2.37	0.56
1:A:26:LEU:HG	1:A:79:LYS:HD2	1.87	0.56
1:C:131:ILE:N	1:C:131:ILE:CD1	2.40	0.56
1:A:76:ILE:N	1:A:79:LYS:O	2.33	0.56
1:A:152:ILE:O	1:A:155:ALA:HB3	2.05	0.56
1:B:133:TYR:CA	1:C:12:PHE:CZ	2.82	0.56
1:C:76:ILE:HD12	1:C:76:ILE:N	2.21	0.56
1:B:51:ARG:HG3	1:B:63:VAL:CG1	2.36	0.56
1:C:112:ILE:CD1	1:C:122:LEU:HD12	2.34	0.56
1:A:21:PHE:HE2	1:A:142:LEU:CD2	2.18	0.56
1:A:153:ILE:O	1:A:155:ALA:N	2.39	0.56
1:B:75:ASN:C	1:B:76:ILE:HG12	2.31	0.56
1:B:51:ARG:HG3	1:B:63:VAL:HG13	1.87	0.56
1:B:4:LEU:HB3	1:B:142:LEU:HB2	1.88	0.55
1:B:55:GLN:HA	1:B:59:HIS:H	1.70	0.55
1:C:22:ALA:HA	1:C:25:LEU:HD12	1.86	0.55
1:A:29:HIS:CG	1:A:30:THR:H	2.24	0.55
1:B:51:ARG:HA	1:B:63:VAL:CG1	2.34	0.55
1:C:82:TYR:CD2	1:C:103:TYR:O	2.58	0.55
1:C:93:GLU:O	1:C:97:PHE:CE2	2.59	0.55
1:C:94:GLU:C	1:C:97:PHE:HD2	2.13	0.55
1:B:25:LEU:O	1:B:25:LEU:CD2	2.54	0.55
1:B:146:THR:CG2	1:B:147:ASN:N	2.33	0.55
1:C:112:ILE:HD11	1:C:122:LEU:CD1	2.37	0.55
1:C:131:ILE:HG23	1:C:140:ILE:CD1	2.36	0.55
1:A:70:LEU:O	1:A:84:PHE:HA	2.07	0.55
1:C:1:MET:HB3	1:C:150:LYS:NZ	2.22	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:133:TYR:O	1:B:133:TYR:CG	2.57	0.55
1:B:148:LEU:HA	1:B:151:ASN:HD21	1.65	0.55
1:A:38:LEU:HB2	1:A:113:GLU:HG3	1.88	0.55
1:A:50:THR:HG21	1:A:83:HIS:NE2	2.22	0.55
1:A:133:TYR:O	1:A:133:TYR:CD1	2.60	0.55
1:B:21:PHE:CB	1:B:156:PHE:HZ	2.19	0.55
1:C:1:MET:HB3	1:C:150:LYS:HE2	1.86	0.55
1:C:131:ILE:HG23	1:C:140:ILE:HD11	1.86	0.55
1:B:8:ILE:HB	1:B:138:ARG:HB2	1.89	0.55
1:B:63:VAL:HG23	1:B:63:VAL:O	2.06	0.55
1:B:136:ASP:OD2	1:C:16:ARG:NH1	2.39	0.55
1:A:74:TYR:HD1	1:A:81:ILE:HG22	1.72	0.55
1:A:126:ASP:O	1:A:149:GLY:HA3	2.07	0.55
1:B:92:PRO:CA	1:B:95:LEU:HG	2.34	0.55
1:C:10:ASP:HB3	3:C:202:HOH:O	2.05	0.55
1:C:6:GLN:HG3	1:C:17:PHE:HE1	1.68	0.55
1:B:22:ALA:HB2	1:B:53:MET:HG2	1.88	0.54
1:B:148:LEU:O	1:B:152:ILE:CD1	2.56	0.54
1:B:7:TYR:N	1:C:55:GLN:OE1	2.26	0.54
1:C:21:PHE:CD1	1:C:25:LEU:HD11	2.43	0.54
1:C:59:HIS:ND1	1:C:63:VAL:CG2	2.71	0.54
1:B:5:THR:HA	1:B:140:ILE:O	2.07	0.54
1:A:46:LYS:O	1:A:50:THR:OG1	2.24	0.54
1:B:24:ILE:CG2	1:B:155:ALA:CB	2.86	0.54
1:A:60:GLN:CA	1:C:6:GLN:HE22	2.17	0.53
1:B:8:ILE:HG21	1:B:138:ARG:HD3	1.90	0.53
1:B:14:MET:CE	1:B:14:MET:HA	2.38	0.53
1:B:133:TYR:CA	1:C:12:PHE:HZ	2.16	0.53
1:A:62:ASN:ND2	1:C:4:LEU:HD22	2.23	0.53
1:B:136:ASP:CB	1:C:16:ARG:HH12	2.20	0.53
1:C:131:ILE:HD12	1:C:140:ILE:CD1	2.38	0.53
1:A:144:ALA:CB	1:A:149:GLY:C	2.80	0.53
1:C:146:THR:HG22	1:C:147:ASN:H	1.74	0.53
1:A:17:PHE:CD1	1:A:17:PHE:C	2.87	0.53
1:A:142:LEU:HB3	1:A:153:ILE:HD12	1.91	0.53
1:A:14:MET:HE1	1:A:140:ILE:CD1	2.34	0.53
1:B:96:GLU:HA	1:B:96:GLU:OE1	2.09	0.53
1:C:33:ALA:HB2	1:C:104:PHE:O	2.08	0.53
1:A:88:ARG:CG	1:A:88:ARG:NH1	2.65	0.53
1:B:9:PRO:HB3	1:C:19:LYS:CE	2.39	0.53
1:C:8:ILE:CG2	1:C:138:ARG:CB	2.70	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:14:MET:HE3	1:A:14:MET:HA	1.91	0.53
1:A:101:ARG:HB2	1:A:103:TYR:CD1	2.44	0.53
1:B:75:ASN:ND2	1:B:80:MET:HE2	2.23	0.53
1:C:24:ILE:CG1	1:C:152:ILE:HD12	2.39	0.53
1:C:50:THR:CG2	1:C:54:LEU:CD1	2.82	0.53
1:C:105:ASN:O	1:C:108:SER:HB2	2.09	0.53
1:A:49:LEU:HD12	1:A:131:ILE:HD11	1.91	0.53
1:B:29:HIS:ND1	1:B:30:THR:N	2.57	0.53
1:A:92:PRO:O	1:A:95:LEU:N	2.40	0.53
1:B:42:LEU:HG	1:B:114:TRP:CH2	2.36	0.53
1:B:128:LEU:CD1	1:B:145:GLN:OE1	2.57	0.53
1:C:1:MET:HB3	1:C:150:LYS:CE	2.38	0.53
1:B:86:LEU:HD22	1:B:89:LEU:HD22	1.91	0.52
1:C:87:TYR:O	1:C:87:TYR:CD1	2.62	0.52
1:B:67:THR:HA	1:B:85:ASP:HB3	1.91	0.52
1:C:49:LEU:HD13	1:C:129:VAL:HG11	1.91	0.52
1:B:38:LEU:CD2	1:B:129:VAL:HG11	2.39	0.52
1:B:54:LEU:N	1:B:54:LEU:CD2	2.71	0.52
1:C:2:GLU:HB2	1:C:150:LYS:HD2	1.90	0.52
1:A:64:LYS:HB3	1:A:72:GLU:OE1	2.09	0.52
1:B:6:GLN:NE2	1:C:60:GLN:HA	2.24	0.52
1:B:47:THR:HG22	1:B:63:VAL:HG21	1.92	0.52
1:C:144:ALA:CB	1:C:153:ILE:CG1	2.77	0.52
1:A:16:ARG:O	1:A:20:LYS:HG3	2.10	0.52
1:B:84:PHE:CE1	1:B:112:ILE:HG12	2.44	0.52
1:C:21:PHE:C	1:C:21:PHE:HD1	2.17	0.52
1:C:89:LEU:CD2	1:C:95:LEU:HD21	2.39	0.52
1:A:80:MET:O	1:A:81:ILE:HG13	2.10	0.52
1:C:24:ILE:HD12	1:C:152:ILE:CD1	2.40	0.52
1:A:47:THR:O	1:A:48:THR:C	2.53	0.51
1:A:71:VAL:CG1	1:A:82:TYR:HB3	2.41	0.51
1:B:2:GLU:CB	1:B:144:ALA:HB2	2.39	0.51
1:B:13:SER:C	1:B:15:LEU:N	2.66	0.51
1:C:64:LYS:HD2	1:C:74:TYR:CE2	2.44	0.51
1:A:101:ARG:CB	1:A:103:TYR:CE1	2.94	0.51
1:B:37:TYR:CE2	1:B:115:SER:HB2	2.44	0.51
1:C:35:MET:CG	1:C:123:PRO:HG2	2.40	0.51
1:C:37:TYR:HD1	1:C:37:TYR:N	2.08	0.51
1:C:39:ASN:O	1:C:130:ASN:HA	2.11	0.51
1:A:89:LEU:HD21	1:A:95:LEU:HG	1.93	0.51
1:B:95:LEU:CD1	1:B:118:GLY:CA	2.88	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:97:PHE:HD1	1:B:97:PHE:H	1.51	0.51
1:C:49:LEU:HD22	1:C:53:MET:HE3	1.92	0.51
1:A:73:GLU:HB3	1:A:82:TYR:CE1	2.43	0.51
1:B:50:THR:HG22	1:B:54:LEU:HG	1.92	0.51
1:A:34:ILE:HG23	1:A:148:LEU:CD2	2.40	0.51
1:B:52:GLY:O	1:B:56:GLY:HA3	2.11	0.51
1:B:2:GLU:O	1:B:143:ILE:CA	2.57	0.51
1:B:2:GLU:CB	1:B:144:ALA:CB	2.87	0.51
1:A:4:LEU:HD13	1:A:142:LEU:HD12	1.92	0.51
1:A:23:GLU:O	1:A:27:LYS:HG3	2.11	0.51
1:B:35:MET:HE1	1:B:112:ILE:HD11	1.91	0.51
1:B:140:ILE:HG22	1:B:140:ILE:O	2.10	0.51
1:C:115:SER:HB3	3:C:201:HOH:O	2.11	0.51
1:A:116:GLU:HB2	1:A:117:LYS:HD2	1.93	0.50
1:B:5:THR:O	1:B:5:THR:CG2	2.59	0.50
1:B:8:ILE:HD13	1:B:14:MET:HA	1.94	0.50
1:A:8:ILE:HG22	1:A:138:ARG:HB2	1.92	0.50
1:C:35:MET:HB3	1:C:37:TYR:HE1	1.76	0.50
1:B:9:PRO:HG3	1:C:56:GLY:HA2	1.91	0.50
1:B:82:TYR:CZ	1:B:103:TYR:CE2	2.99	0.50
1:C:37:TYR:CD1	1:C:37:TYR:N	2.80	0.50
1:B:7:TYR:OH	1:B:137:ALA:CB	2.59	0.50
1:B:20:LYS:HG3	1:B:21:PHE:N	2.25	0.50
1:C:49:LEU:CD1	1:C:129:VAL:HG11	2.41	0.50
1:B:14:MET:HB3	1:B:48:THR:CG2	2.41	0.50
1:B:126:ASP:HB3	1:B:146:THR:CB	2.42	0.50
1:B:150:LYS:O	1:B:154:SER:OG	2.30	0.50
1:B:66:PRO:O	1:B:69:THR:HB	2.12	0.50
1:C:94:GLU:O	1:C:97:PHE:HD2	1.94	0.50
1:A:54:LEU:CD2	1:A:81:ILE:HG21	2.41	0.50
1:B:84:PHE:CE2	1:B:104:PHE:HE1	2.29	0.50
1:B:148:LEU:CD2	1:B:149:GLY:N	2.73	0.50
1:B:84:PHE:CE2	1:B:104:PHE:CE1	3.00	0.50
1:C:11:GLU:O	1:C:14:MET:HB2	2.11	0.50
1:B:18:GLY:HA2	1:B:49:LEU:CD2	2.42	0.49
1:C:14:MET:CG	1:C:138:ARG:NH1	2.75	0.49
1:A:16:ARG:HH21	1:C:135:ASP:CB	2.03	0.49
1:B:37:TYR:N	1:B:37:TYR:HD1	2.10	0.49
1:B:92:PRO:CB	1:B:121:ILE:HD11	2.42	0.49
1:B:10:ASP:CG	1:B:11:GLU:H	2.19	0.49
1:C:127:ILE:CG2	1:C:128:LEU:H	2.18	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:24:ILE:O	1:A:25:LEU:C	2.54	0.49
1:C:1:MET:CB	1:C:150:LYS:CE	2.87	0.49
1:C:89:LEU:HD11	1:C:91:ASP:O	2.12	0.49
1:C:146:THR:HG22	1:C:148:LEU:H	1.78	0.49
1:A:21:PHE:CE2	1:A:142:LEU:HD21	2.45	0.49
1:B:37:TYR:HE2	1:B:115:SER:CB	2.23	0.49
1:C:64:LYS:HD2	1:C:74:TYR:HE2	1.77	0.49
1:C:146:THR:O	1:C:150:LYS:N	2.32	0.49
1:A:22:ALA:O	1:A:25:LEU:HB2	2.12	0.49
1:C:34:ILE:O	1:C:110:CYS:HB2	2.12	0.49
1:C:76:ILE:HD13	1:C:81:ILE:HD11	1.95	0.49
1:B:95:LEU:CD1	1:B:118:GLY:HA2	2.43	0.49
1:C:44:ALA:HA	1:C:133:TYR:CD1	2.47	0.49
1:A:44:ALA:HA	1:A:133:TYR:CD2	2.47	0.49
1:B:38:LEU:HD23	1:B:129:VAL:CG1	2.42	0.49
1:B:139:ASN:OD1	1:B:140:ILE:N	2.46	0.49
1:A:57:ILE:HG21	1:A:76:ILE:HG21	1.92	0.49
1:A:57:ILE:HG21	1:A:76:ILE:CG2	2.41	0.49
1:A:146:THR:CG2	1:A:147:ASN:H	2.06	0.49
1:B:2:GLU:HB3	1:B:153:ILE:HD12	1.95	0.49
1:B:104:PHE:CE2	1:B:123:PRO:HD2	2.48	0.49
1:C:17:PHE:C	1:C:17:PHE:CD2	2.90	0.49
1:C:36:VAL:CG2	1:C:127:ILE:HD13	2.43	0.49
1:A:141:GLU:HB3	1:A:143:ILE:HG13	1.94	0.48
1:B:19:LYS:HD2	1:B:56:GLY:O	2.13	0.48
1:B:25:LEU:HD21	1:B:34:ILE:CD1	2.44	0.48
1:C:26:LEU:HD11	1:C:57:ILE:HG23	1.95	0.48
1:C:62:ASN:HB2	3:C:210:HOH:O	2.12	0.48
1:B:19:LYS:CB	1:B:52:GLY:O	2.60	0.48
1:B:21:PHE:CD2	1:B:53:MET:SD	3.06	0.48
1:C:60:GLN:CD	1:C:60:GLN:C	2.81	0.48
1:C:146:THR:CG2	1:C:147:ASN:N	2.76	0.48
1:B:8:ILE:HB	1:B:138:ARG:O	2.14	0.48
1:B:92:PRO:HB3	1:B:121:ILE:HD11	1.94	0.48
1:C:22:ALA:O	1:C:25:LEU:HB2	2.14	0.48
1:C:38:LEU:HD11	1:C:111:LEU:HD21	1.93	0.48
1:B:146:THR:HG22	1:B:148:LEU:H	1.79	0.48
1:C:84:PHE:CD2	1:C:98:MET:HE3	2.49	0.48
1:C:54:LEU:HD11	1:C:111:LEU:HD13	1.96	0.48
1:C:69:THR:O	1:C:70:LEU:CB	2.62	0.48
1:A:64:LYS:HB2	1:A:64:LYS:HE3	1.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:131:ILE:O	1:A:131:ILE:HG22	2.14	0.48
1:B:104:PHE:HZ	1:B:122:LEU:HD23	1.79	0.48
1:C:59:HIS:ND1	1:C:63:VAL:HG22	2.28	0.48
1:B:7:TYR:CD2	1:C:52:GLY:HA2	2.49	0.48
1:C:102:ASP:HA	1:C:105:ASN:HD21	1.77	0.48
1:A:8:ILE:HG21	1:A:138:ARG:HB2	1.95	0.47
1:A:37:TYR:CE1	1:A:125:ALA:HB2	2.48	0.47
1:B:89:LEU:CD2	1:B:95:LEU:CD2	2.89	0.47
1:B:137:ALA:H	1:C:16:ARG:CZ	2.26	0.47
1:B:146:THR:HB	1:B:148:LEU:CD2	2.42	0.47
1:C:9:PRO:HD2	1:C:13:SER:OG	2.14	0.47
1:C:14:MET:HG2	1:C:138:ARG:HH11	1.76	0.47
1:C:60:GLN:C	1:C:61:GLY:O	2.53	0.47
1:A:35:MET:CA	1:A:35:MET:CE	2.83	0.47
1:A:139:ASN:O	1:A:140:ILE:CG1	2.61	0.47
1:B:21:PHE:HZ	1:B:36:VAL:CG1	2.19	0.47
1:C:49:LEU:O	1:C:53:MET:HB2	2.14	0.47
1:C:95:LEU:CD1	1:C:118:GLY:HA2	2.43	0.47
1:A:139:ASN:C	1:A:140:ILE:HG13	2.39	0.47
1:B:9:PRO:HB3	1:C:19:LYS:NZ	2.30	0.47
1:B:89:LEU:CG	1:B:95:LEU:HD21	2.43	0.47
1:B:133:TYR:CB	1:B:138:ARG:HA	2.44	0.47
1:C:82:TYR:HE2	1:C:108:SER:OG	1.97	0.47
1:C:23:GLU:HA	1:C:26:LEU:HD12	1.96	0.47
1:A:64:LYS:O	1:A:65:SER:C	2.57	0.47
1:B:29:HIS:CG	1:B:30:THR:H	2.33	0.47
1:B:82:TYR:CD2	1:B:103:TYR:CB	2.98	0.47
1:C:26:LEU:HD11	1:C:57:ILE:HD13	1.96	0.47
1:A:51:ARG:NH2	1:C:139:ASN:HD21	2.13	0.47
1:A:53:MET:HE2	1:A:53:MET:HB2	1.65	0.47
1:A:55:GLN:C	1:A:57:ILE:H	2.22	0.47
1:C:37:TYR:CD1	1:C:125:ALA:CB	2.89	0.47
1:B:21:PHE:HB2	1:B:156:PHE:HZ	1.80	0.47
1:B:54:LEU:O	1:B:59:HIS:HB2	2.14	0.47
1:A:126:ASP:OD1	1:A:126:ASP:N	2.47	0.46
1:B:25:LEU:CD2	1:B:25:LEU:C	2.88	0.46
1:A:84:PHE:HB2	1:A:112:ILE:HA	1.97	0.46
1:A:118:GLY:CA	1:A:122:LEU:CD1	2.87	0.46
1:A:130:ASN:O	1:A:140:ILE:HA	2.15	0.46
1:B:95:LEU:CD1	1:B:118:GLY:HA3	2.45	0.46
1:A:25:LEU:HD22	1:A:34:ILE:HD13	1.90	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:36:VAL:HG12	1:B:129:VAL:CG2	2.45	0.46
1:B:81:ILE:CG2	1:B:111:LEU:HD12	2.46	0.46
1:C:34:ILE:HG22	1:C:34:ILE:O	2.16	0.46
1:C:50:THR:HG23	1:C:54:LEU:HD13	1.94	0.46
1:A:92:PRO:O	1:A:93:GLU:C	2.57	0.46
1:A:118:GLY:O	1:A:122:LEU:HB3	2.15	0.46
1:B:10:ASP:O	1:B:13:SER:N	2.42	0.46
1:B:19:LYS:NZ	1:B:23:GLU:OE2	2.48	0.46
1:B:69:THR:C	1:B:70:LEU:HG	2.40	0.46
1:B:82:TYR:CE2	1:B:103:TYR:CB	2.98	0.46
1:B:74:TYR:HB2	1:B:81:ILE:HB	1.97	0.46
1:A:25:LEU:CD2	1:A:34:ILE:CD1	2.89	0.46
1:C:21:PHE:CE2	1:C:129:VAL:CG2	2.99	0.46
1:A:26:LEU:CA	1:A:79:LYS:CE	2.69	0.45
1:C:14:MET:O	1:C:15:LEU:C	2.57	0.45
1:A:33:ALA:O	1:A:35:MET:CE	2.65	0.45
1:B:7:TYR:CZ	1:B:137:ALA:CB	2.99	0.45
1:B:20:LYS:CD	1:B:24:ILE:CD1	2.74	0.45
1:C:82:TYR:CE2	1:C:103:TYR:HA	2.51	0.45
1:B:8:ILE:O	1:B:8:ILE:CG2	2.63	0.45
1:B:38:LEU:HD23	1:B:129:VAL:HG11	1.97	0.45
1:B:95:LEU:HD12	1:B:122:LEU:HD12	1.96	0.45
1:B:126:ASP:HB3	1:B:146:THR:HB	1.97	0.45
1:C:21:PHE:CD1	1:C:21:PHE:O	2.69	0.45
1:A:25:LEU:HD22	1:A:34:ILE:CD1	2.47	0.45
1:A:54:LEU:HD22	1:A:81:ILE:HG21	1.98	0.45
1:A:73:GLU:H	1:A:73:GLU:HG3	1.40	0.45
1:B:37:TYR:O	1:B:129:VAL:HB	2.16	0.45
1:C:28:LEU:HD13	1:C:152:ILE:HD11	1.98	0.45
1:C:35:MET:HG2	1:C:123:PRO:CD	2.46	0.45
1:C:81:ILE:H	1:C:81:ILE:HG12	1.39	0.45
1:C:82:TYR:HB2	1:C:110:CYS:HA	1.99	0.45
1:A:75:ASN:HB2	1:A:80:MET:HE2	1.99	0.45
1:A:98:MET:O	1:A:101:ARG:N	2.50	0.45
1:B:14:MET:CG	1:B:48:THR:HG22	2.47	0.45
1:B:115:SER:O	1:B:116:GLU:C	2.59	0.45
1:A:34:ILE:HG23	1:A:148:LEU:HD23	1.98	0.45
1:A:24:ILE:HG21	1:A:24:ILE:HD13	1.66	0.45
1:A:60:GLN:O	1:A:60:GLN:HG3	2.17	0.45
1:A:99:GLY:C	1:A:101:ARG:H	2.24	0.45
1:C:35:MET:SD	1:C:110:CYS:HB3	2.56	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:86:LEU:HD13	1:B:122:LEU:HD13	1.97	0.45
1:C:22:ALA:O	1:C:23:GLU:C	2.58	0.45
1:A:117:LYS:HD2	1:A:117:LYS:H	1.81	0.45
1:A:128:LEU:N	1:A:143:ILE:O	2.46	0.45
1:C:76:ILE:N	1:C:76:ILE:CD1	2.80	0.45
1:B:86:LEU:CD1	1:B:122:LEU:HD13	2.46	0.44
1:B:148:LEU:O	1:B:152:ILE:CG1	2.65	0.44
1:C:36:VAL:HG22	1:C:127:ILE:HD13	1.99	0.44
1:A:73:GLU:C	1:A:74:TYR:CD1	2.96	0.44
1:B:13:SER:C	1:B:15:LEU:H	2.26	0.44
1:B:19:LYS:CA	1:B:52:GLY:O	2.65	0.44
1:B:64:LYS:HD3	1:B:64:LYS:C	2.42	0.44
1:C:35:MET:SD	1:C:104:PHE:HE1	2.38	0.44
1:C:87:TYR:HB2	1:C:114:TRP:CD1	2.53	0.44
1:B:118:GLY:O	1:B:122:LEU:HB2	2.18	0.44
1:B:146:THR:HG21	1:B:148:LEU:CD2	2.44	0.44
1:C:11:GLU:HA	1:C:138:ARG:NH1	2.30	0.44
1:A:122:LEU:HA	1:A:123:PRO:HD3	1.77	0.44
1:A:22:ALA:HA	1:A:25:LEU:HB2	1.99	0.44
1:A:59:HIS:CE1	1:A:62:ASN:O	2.67	0.44
1:C:23:GLU:O	1:C:26:LEU:HB2	2.17	0.44
1:C:98:MET:HG2	1:C:99:GLY:N	2.33	0.44
1:A:86:LEU:CD2	1:A:89:LEU:CD2	2.89	0.44
1:B:8:ILE:HD13	1:B:14:MET:CA	2.48	0.44
1:B:18:GLY:O	1:B:21:PHE:HB3	2.18	0.44
1:B:30:THR:O	1:B:107:ASP:HA	2.18	0.44
1:B:140:ILE:CG2	1:B:140:ILE:O	2.63	0.44
1:C:142:LEU:HD23	1:C:142:LEU:H	1.81	0.44
1:A:30:THR:HG21	1:A:34:ILE:HG13	2.00	0.43
1:A:97:PHE:N	1:A:97:PHE:CD1	2.76	0.43
1:C:84:PHE:HB2	1:C:112:ILE:CG1	2.47	0.43
1:C:87:TYR:CD1	1:C:87:TYR:C	2.90	0.43
1:A:79:LYS:HD3	1:A:109:ILE:HD12	1.81	0.43
1:A:144:ALA:HB1	1:A:149:GLY:HA3	2.00	0.43
1:B:87:TYR:O	1:B:88:ARG:C	2.61	0.43
1:B:128:LEU:HD11	1:B:145:GLN:OE1	2.18	0.43
1:C:37:TYR:CG	1:C:125:ALA:HB2	2.49	0.43
1:C:59:HIS:ND1	1:C:63:VAL:HG23	2.32	0.43
1:A:71:VAL:CG1	1:A:84:PHE:CZ	2.98	0.43
1:A:71:VAL:HA	1:A:83:HIS:O	2.18	0.43
1:A:100:ILE:O	1:A:100:ILE:HG22	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:144:ALA:HB1	1:A:149:GLY:CA	2.46	0.43
1:B:25:LEU:CD2	1:B:34:ILE:HD13	2.48	0.43
1:B:63:VAL:O	1:B:63:VAL:CG2	2.67	0.43
1:C:86:LEU:O	1:C:117:LYS:HG3	2.18	0.43
1:C:130:ASN:HB2	1:C:141:GLU:CB	2.48	0.43
1:A:60:GLN:O	1:A:60:GLN:CG	2.66	0.43
1:A:61:GLY:N	1:C:6:GLN:HE21	2.13	0.43
1:C:13:SER:O	1:C:14:MET:C	2.62	0.43
1:C:22:ALA:O	1:C:25:LEU:N	2.52	0.43
1:A:4:LEU:O	1:A:142:LEU:N	2.38	0.43
1:B:8:ILE:CG2	1:B:138:ARG:HD3	2.48	0.43
1:B:146:THR:CG2	1:B:148:LEU:CD2	2.96	0.43
1:C:35:MET:HG3	1:C:37:TYR:CE1	2.53	0.43
1:C:75:ASN:O	1:C:76:ILE:HD12	2.18	0.43
1:A:74:TYR:HB2	1:A:81:ILE:CG2	2.49	0.43
1:C:37:TYR:CE1	1:C:125:ALA:CB	2.98	0.43
1:B:26:LEU:HD11	1:B:57:ILE:HG13	1.99	0.43
1:B:51:ARG:C	1:B:53:MET:N	2.75	0.43
1:C:50:THR:HG23	1:C:54:LEU:CD1	2.49	0.43
1:C:84:PHE:CZ	1:C:103:TYR:CD2	2.99	0.43
1:A:41:ASP:O	1:A:42:LEU:C	2.61	0.43
1:B:82:TYR:CD2	1:B:103:TYR:HB3	2.53	0.43
1:C:116:GLU:H	1:C:116:GLU:HG2	1.67	0.43
1:C:146:THR:CG2	1:C:147:ASN:H	2.31	0.43
1:A:44:ALA:O	1:A:131:ILE:HG21	2.18	0.43
1:B:88:ARG:HG3	1:B:88:ARG:NH1	2.29	0.43
1:C:8:ILE:HD11	1:C:140:ILE:HG12	1.90	0.43
1:A:53:MET:O	1:A:57:ILE:HG13	2.18	0.43
1:A:83:HIS:CE1	1:A:113:GLU:OE2	2.72	0.43
1:B:66:PRO:HG2	1:B:83:HIS:O	2.19	0.43
1:C:41:ASP:O	1:C:43:GLY:N	2.52	0.43
1:C:150:LYS:HB2	1:C:150:LYS:HE3	1.77	0.43
1:B:132:ASP:HB2	3:B:206:HOH:O	2.18	0.42
1:B:133:TYR:CB	1:B:138:ARG:CG	2.96	0.42
1:C:7:TYR:HA	1:C:139:ASN:HD22	1.84	0.42
1:C:37:TYR:O	1:C:39:ASN:N	2.51	0.42
1:C:45:GLY:O	1:C:46:LYS:C	2.62	0.42
1:C:131:ILE:HD12	1:C:131:ILE:HG23	1.73	0.42
1:B:6:GLN:O	1:B:139:ASN:OD1	2.37	0.42
1:B:47:THR:HG23	1:B:50:THR:HB	2.00	0.42
1:B:108:SER:O	1:B:109:ILE:HG12	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:34:ILE:CG2	1:C:110:CYS:N	2.75	0.42
1:B:18:GLY:HA2	1:B:49:LEU:HD23	2.01	0.42
1:B:82:TYR:CE1	1:B:103:TYR:CD2	3.07	0.42
1:C:2:GLU:CB	1:C:150:LYS:HD2	2.49	0.42
1:C:60:GLN:NE2	1:C:60:GLN:N	2.66	0.42
1:B:7:TYR:CE1	1:C:15:LEU:HB3	2.54	0.42
1:B:84:PHE:HE2	1:B:104:PHE:CE1	2.38	0.42
1:B:133:TYR:HA	1:C:12:PHE:CE1	2.50	0.42
1:C:102:ASP:CB	1:C:105:ASN:HD22	2.32	0.42
1:B:10:ASP:OD1	1:B:11:GLU:HB3	2.20	0.42
1:B:18:GLY:CA	1:B:49:LEU:HD23	2.50	0.42
1:B:26:LEU:HD23	1:B:79:LYS:CB	2.46	0.42
1:B:37:TYR:HE2	1:B:115:SER:HB2	1.82	0.42
1:C:28:LEU:HD13	1:C:152:ILE:HD13	2.01	0.42
1:C:80:MET:CE	1:C:82:TYR:OH	2.67	0.42
1:A:144:ALA:HB2	1:A:149:GLY:O	2.20	0.42
1:C:35:MET:HG3	1:C:37:TYR:HE1	1.84	0.42
1:A:98:MET:O	1:A:101:ARG:HB2	2.19	0.42
1:B:2:GLU:O	1:B:144:ALA:N	2.52	0.42
1:C:122:LEU:N	1:C:122:LEU:HD23	2.35	0.42
1:A:50:THR:O	1:A:53:MET:HB3	2.19	0.42
1:A:144:ALA:CB	1:A:149:GLY:O	2.68	0.42
1:B:136:ASP:H	1:C:16:ARG:CZ	2.33	0.42
1:A:44:ALA:HB1	1:A:133:TYR:CG	2.55	0.41
1:A:98:MET:O	1:A:101:ARG:CB	2.68	0.41
1:B:55:GLN:NE2	1:B:61:GLY:O	2.51	0.41
1:B:137:ALA:H	1:C:16:ARG:HH12	1.68	0.41
1:C:17:PHE:CE1	1:C:156:PHE:HE2	2.38	0.41
1:C:80:MET:H	1:C:80:MET:HG3	1.67	0.41
1:A:8:ILE:O	1:A:137:ALA:HB1	2.20	0.41
1:A:14:MET:HE2	1:A:49:LEU:HG	2.02	0.41
1:B:133:TYR:O	1:B:133:TYR:CE1	2.71	0.41
1:A:23:GLU:H	1:A:23:GLU:HG3	1.65	0.41
1:C:15:LEU:HD22	1:C:52:GLY:CA	2.50	0.41
1:A:67:THR:HG22	1:A:68:TYR:CD1	2.56	0.41
1:B:24:ILE:O	1:B:28:LEU:HG	2.19	0.41
1:C:8:ILE:HG21	1:C:138:ARG:CB	2.42	0.41
1:B:92:PRO:HA	1:B:95:LEU:CG	2.41	0.41
1:A:141:GLU:HB3	1:A:143:ILE:CG1	2.51	0.41
1:B:53:MET:O	1:B:57:ILE:HB	2.21	0.41
1:C:91:ASP:HB2	1:C:92:PRO:CD	2.48	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:7:TYR:CZ	1:B:137:ALA:HB2	2.55	0.41
1:B:82:TYR:CD2	1:B:103:TYR:CG	3.07	0.41
1:C:5:THR:HG22	1:C:6:GLN:N	2.36	0.41
1:B:14:MET:CB	1:B:48:THR:HG22	2.45	0.40
1:B:85:ASP:HA	1:B:113:GLU:HB3	2.02	0.40
1:B:87:TYR:C	1:B:89:LEU:N	2.77	0.40
1:B:142:LEU:HD23	1:B:142:LEU:HA	1.91	0.40
1:C:5:THR:CG2	1:C:6:GLN:N	2.85	0.40
1:C:22:ALA:HA	1:C:25:LEU:CD1	2.51	0.40
1:A:4:LEU:CD1	1:A:153:ILE:HG23	2.45	0.40
1:A:76:ILE:HG13	1:A:81:ILE:HD12	2.02	0.40
1:A:83:HIS:O	1:A:84:PHE:CD1	2.74	0.40
1:A:111:LEU:HD23	1:A:111:LEU:H	1.86	0.40
1:A:114:TRP:O	1:A:116:GLU:N	2.55	0.40
1:A:152:ILE:O	1:A:153:ILE:C	2.64	0.40
1:B:87:TYR:CD1	1:B:114:TRP:CZ3	3.09	0.40
1:C:29:HIS:CG	1:C:30:THR:N	2.72	0.40
1:C:128:LEU:HD13	1:C:145:GLN:OE1	2.21	0.40
1:C:148:LEU:C	1:C:148:LEU:CD2	2.86	0.40
1:A:60:GLN:CA	1:C:6:GLN:NE2	2.69	0.40
1:B:4:LEU:CD2	1:C:62:ASN:HD21	2.29	0.40
1:B:136:ASP:CB	1:C:16:ARG:NH1	2.82	0.40
1:B:15:LEU:O	1:B:16:ARG:C	2.61	0.40
1:B:37:TYR:HE1	1:B:125:ALA:HB2	1.67	0.40
1:B:70:LEU:N	1:B:70:LEU:HD12	2.37	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	155/161 (96%)	118 (76%)	30 (19%)	7 (4%)	2 2

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	155/161 (96%)	127 (82%)	23 (15%)	5 (3%)	3	4
1	C	155/161 (96%)	116 (75%)	30 (19%)	9 (6%)	1	1
All	All	465/483 (96%)	361 (78%)	83 (18%)	21 (4%)	2	2

All (21) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	116	GLU
1	C	115	SER
1	A	115	SER
1	A	140	ILE
1	A	154	SER
1	B	31	GLU
1	C	38	LEU
1	C	118	GLY
1	A	4	LEU
1	B	58	GLY
1	C	33	ALA
1	C	57	ILE
1	C	142	LEU
1	B	67	THR
1	B	115	SER
1	C	9	PRO
1	C	66	PRO
1	A	11	GLU
1	A	29	HIS
1	C	156	PHE
1	A	58	GLY

5.3.2 Protein sidechains ⓘ

There are no protein residues with a non-rotameric sidechain to report in this entry.

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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 3 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.