



# Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 04:10 PM GMT

PDB ID : 2ZBE  
Title : Calcium pump crystal structure with bound BeF<sub>3</sub> in the absence of calcium and TG  
Authors : Toyoshima, C.; Ogawa, H.; Norimatsu, Y.  
Deposited on : 2007-10-20  
Resolution : 3.80 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

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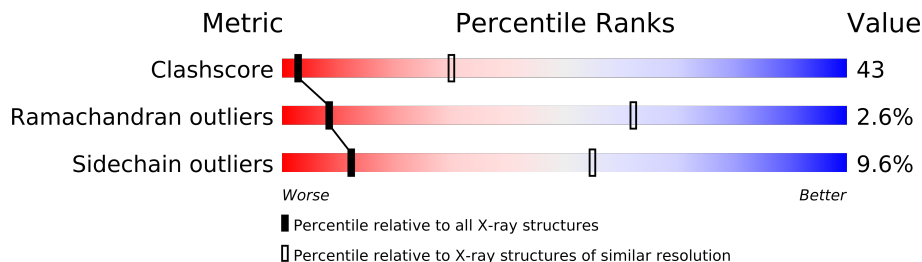
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 21963  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 3.80 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	79885	1100 (4.10-3.50)
Ramachandran outliers	78287	1050 (4.10-3.50)
Sidechain outliers	78261	1042 (4.10-3.50)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	995	
1	B	995	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 15358 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Sarcoplasmic/endoplasmicreticulum calcium ATPase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	995	Total	C	N	O	S	1	0	0
			7674	4878	1287	1452	57			
1	B	995	Total	C	N	O	S	1	0	0
			7674	4878	1287	1452	57			

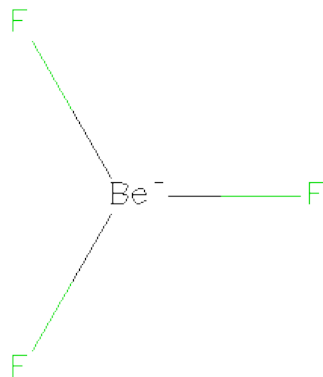
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	994	GLY	-	SEE REMARK 999	UNP P04191
B	994	GLY	-	SEE REMARK 999	UNP P04191

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 3 is BERYLLIUM TRIFLUORIDE ION (three-letter code: BEF) (formula: BeF<sub>3</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	Be	F	0	0
			4	1	3		
3	B	1	Total	Be	F	0	0
			4	1	3		

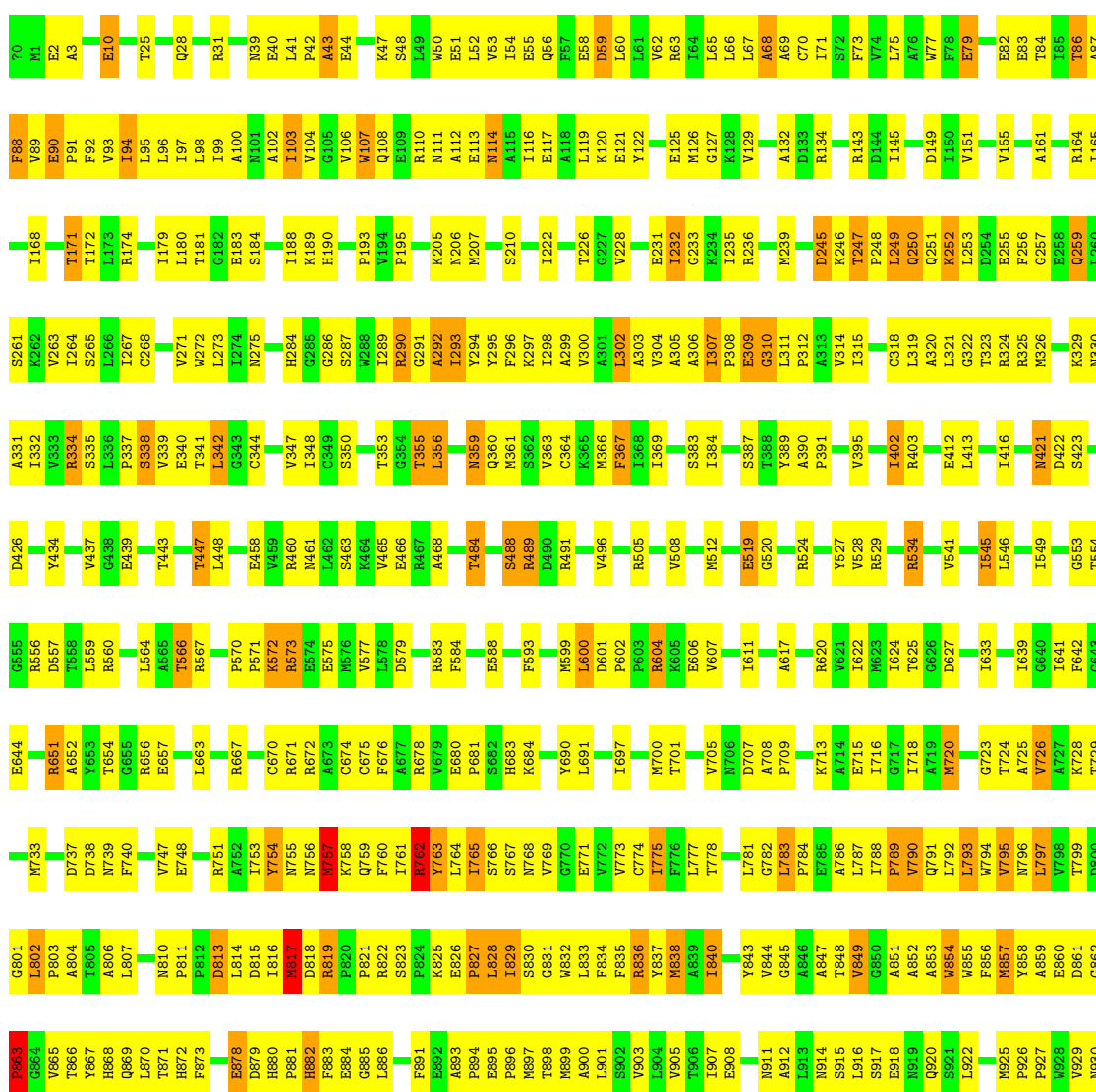
### 3 Residue-property plots

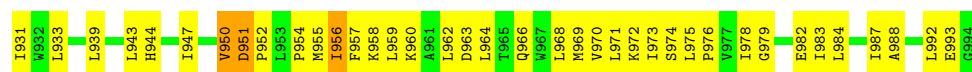
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

Note EDS was not executed.

- Molecule 1: Sarcoplasmic/endoplasmicreticulum calcium ATPase 1

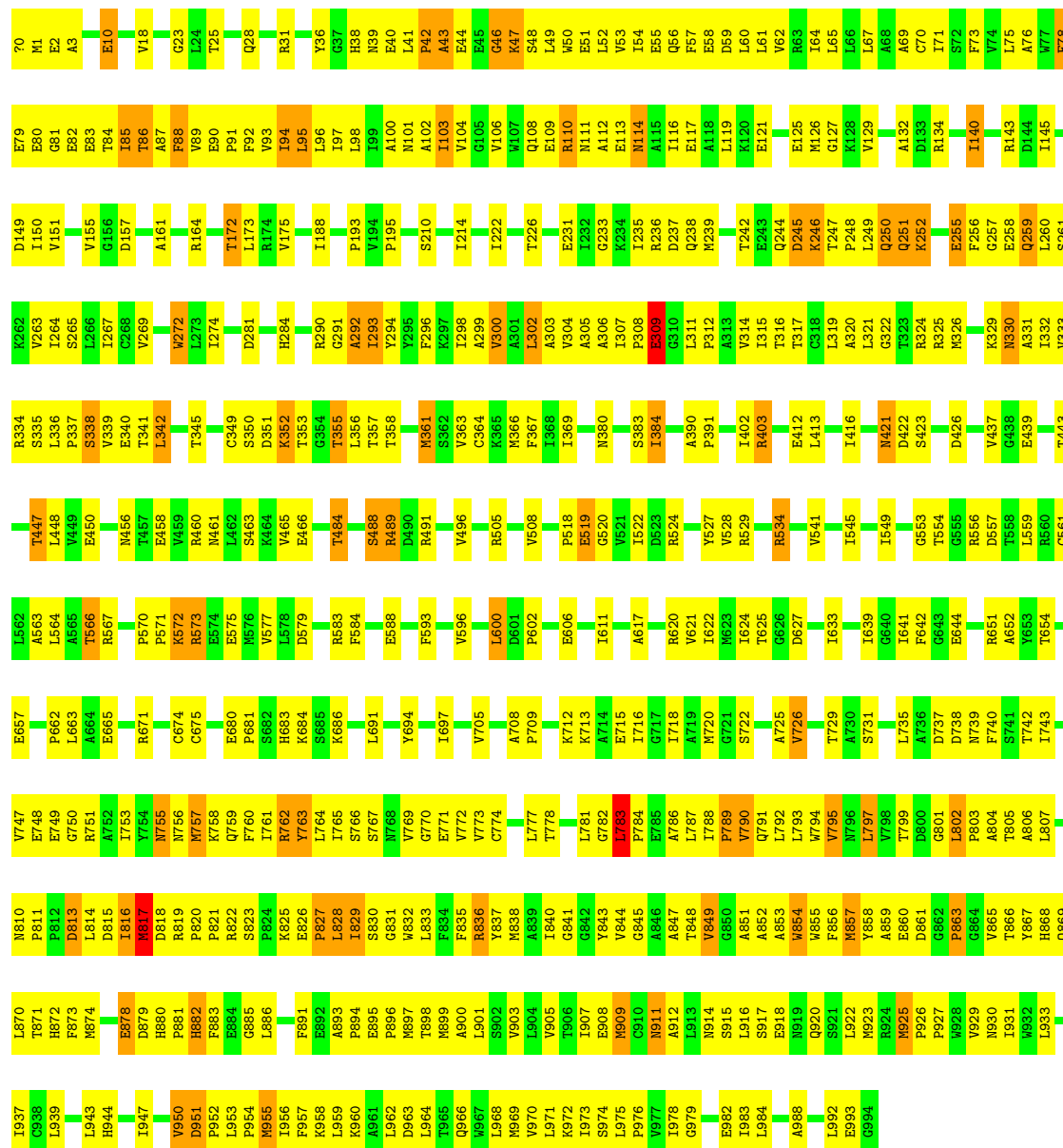
Chain A:





• Molecule 1: Sarcoplasmic/endoplasmicreticulum calcium ATPase 1

Chain B:



## 4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	133.87Å 91.59Å 248.46Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	12.00 – 3.80	Depositor
% Data completeness (in resolution range)	90.3 (12.00-3.80)	Depositor
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	CNS 1.2	Depositor
R, $R_{free}$	0.293 , 0.327	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	15358	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	163.0	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: MG, BEF, ACE

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	0.74	2/7813 (0.0%)	0.89	25/10594 (0.2%)
1	B	0.75	1/7813 (0.0%)	0.86	15/10594 (0.1%)
All	All	0.74	3/15626 (0.0%)	0.88	40/21188 (0.2%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	B	0	2
All	All	0	4

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	107	TRP	CB-CG	5.95	1.60	1.50
1	A	863	PRO	C-N	-5.75	1.22	1.33
1	B	561	CYS	CB-SG	-5.35	1.73	1.81

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	671	ARG	NE-CZ-NH1	9.46	125.03	120.30
1	A	822	ARG	NE-CZ-NH2	7.72	124.16	120.30
1	A	651	ARG	NE-CZ-NH2	7.24	123.92	120.30
1	A	762	ARG	NE-CZ-NH2	7.24	123.92	120.30
1	A	290	ARG	NE-CZ-NH2	7.05	123.82	120.30
1	A	560	ARG	NE-CZ-NH2	6.99	123.80	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	836	ARG	NE-CZ-NH2	6.97	123.79	120.30
1	A	751	ARG	NE-CZ-NH2	6.95	123.78	120.30
1	A	678	ARG	NE-CZ-NH2	6.74	123.67	120.30
1	A	656	ARG	NE-CZ-NH2	6.67	123.63	120.30
1	A	819	ARG	NE-CZ-NH2	6.66	123.63	120.30
1	A	667	ARG	NE-CZ-NH2	6.65	123.62	120.30
1	A	604	ARG	NE-CZ-NH2	6.47	123.53	120.30
1	B	822	ARG	NE-CZ-NH2	6.43	123.52	120.30
1	A	672	ARG	NE-CZ-NH2	6.42	123.51	120.30
1	B	366	MET	CG-SD-CE	6.20	110.12	100.20
1	B	838	MET	CG-SD-CE	6.16	110.05	100.20
1	A	366	MET	CG-SD-CE	6.15	110.04	100.20
1	A	757	MET	CG-SD-CE	6.04	109.87	100.20
1	A	838	MET	CG-SD-CE	6.03	109.85	100.20
1	A	817	MET	CG-SD-CE	5.99	109.78	100.20
1	A	361	MET	CG-SD-CE	5.96	109.75	100.20
1	B	857	MET	CG-SD-CE	5.92	109.68	100.20
1	B	309	GLU	O-C-N	-5.92	113.14	123.20
1	B	126	MET	CG-SD-CE	5.91	109.65	100.20
1	A	599	MET	CG-SD-CE	5.79	109.46	100.20
1	B	361	MET	CG-SD-CE	5.75	109.39	100.20
1	A	720	MET	CG-SD-CE	5.73	109.37	100.20
1	A	857	MET	CG-SD-CE	5.65	109.25	100.20
1	B	836	ARG	NE-CZ-NH2	5.65	123.12	120.30
1	B	955	MET	CG-SD-CE	5.60	109.17	100.20
1	A	733	MET	CG-SD-CE	5.60	109.16	100.20
1	A	126	MET	CG-SD-CE	5.47	108.95	100.20
1	A	700	MET	CG-SD-CE	5.46	108.94	100.20
1	B	909	MET	CG-SD-CE	5.46	108.93	100.20
1	B	923	MET	CG-SD-CE	5.40	108.83	100.20
1	B	874	MET	CG-SD-CE	5.32	108.70	100.20
1	B	817	MET	CG-SD-CE	5.31	108.69	100.20
1	B	274	ILE	O-C-N	-5.27	114.26	122.70
1	B	925	MET	CG-SD-CE	5.03	108.24	100.20

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	310	GLY	Mainchain
1	A	68	ALA	Mainchain
1	B	284	HIS	Mainchain

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Mol	Chain	Res	Type	Group
1	B	380	ASN	Mainchain

## 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	7674	0	7765	645	0
1	B	7674	0	7765	689	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	4	0	0	1	0
3	B	4	0	0	0	0
All	All	15358	0	15530	1334	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 43.

All (1334) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:757:MET:CE	1:A:761:ILE:HD11	1.60	1.30
1:B:762:ARG:HE	1:B:833:LEU:HD21	1.02	1.18
1:B:802:LEU:HD13	1:B:939:LEU:CD2	1.74	1.17
1:A:247:THR:OG1	1:A:250:GLN:HB3	1.45	1.17
1:A:802:LEU:HD13	1:A:939:LEU:CD2	1.76	1.15
1:B:802:LEU:HD13	1:B:939:LEU:HD23	1.23	1.15
1:B:735:LEU:HD11	1:B:743:ILE:HD11	1.23	1.12
1:A:762:ARG:HE	1:A:833:LEU:HD21	1.00	1.12
1:A:758:LYS:HA	1:A:761:ILE:HD12	1.32	1.11
1:A:802:LEU:HD13	1:A:939:LEU:HD23	1.25	1.09
1:A:898:THR:HA	1:A:959:LEU:HD22	1.33	1.09
1:A:248:PRO:HG2	1:A:341:THR:CG2	1.83	1.09
1:A:248:PRO:HG2	1:A:341:THR:HG21	1.27	1.08
1:A:757:MET:HE2	1:A:761:ILE:HD11	1.13	1.08
1:A:758:LYS:HB2	1:A:828:LEU:HD11	1.30	1.08
1:A:803:PRO:O	1:A:806:ALA:HB3	1.52	1.08

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:823:SER:HB3	1:A:826:GLU:HB2	1.35	1.07
1:B:549:ILE:HD11	1:B:596:VAL:HG21	1.37	1.07
1:B:803:PRO:O	1:B:806:ALA:HB3	1.55	1.06
1:B:823:SER:HB3	1:B:826:GLU:HB2	1.37	1.06
1:B:758:LYS:HB2	1:B:828:LEU:HD11	1.33	1.06
1:A:762:ARG:NE	1:A:833:LEU:HD21	1.69	1.05
1:B:898:THR:HA	1:B:959:LEU:HD22	1.33	1.05
1:B:855:TRP:HA	1:B:859:ALA:CB	1.87	1.05
1:A:855:TRP:HA	1:A:859:ALA:HB2	1.06	1.04
1:B:762:ARG:NE	1:B:833:LEU:HD21	1.70	1.04
1:B:735:LEU:CD1	1:B:743:ILE:HD11	1.88	1.04
1:B:267:ILE:HG12	1:B:302:LEU:HD22	1.40	1.04
1:A:341:THR:HG22	1:A:716:ILE:HD11	1.38	1.02
1:A:855:TRP:HA	1:A:859:ALA:CB	1.88	1.02
1:B:855:TRP:HA	1:B:859:ALA:HB2	1.05	1.02
1:A:899:MET:HE1	1:A:970:VAL:HG23	1.42	1.01
1:B:899:MET:HE1	1:B:970:VAL:HG23	1.41	1.00
1:B:342:LEU:HG	1:B:716:ILE:HG21	1.44	0.99
1:A:760:PHE:HB3	1:A:807:LEU:HD12	1.44	0.99
1:B:783:LEU:HD12	1:B:783:LEU:H	1.27	0.98
1:B:311:LEU:HD11	1:B:761:ILE:CD1	1.92	0.98
1:A:849:VAL:O	1:A:853:ALA:HB3	1.63	0.97
1:B:786:ALA:O	1:B:787:LEU:HG	1.65	0.96
1:A:120:LYS:HE2	1:A:728:LYS:NZ	1.80	0.96
1:B:963:ASP:HB3	1:B:966:GLN:HG3	1.48	0.96
1:A:963:ASP:HB3	1:A:966:GLN:HG3	1.48	0.95
1:B:617:ALA:O	1:B:821:PRO:HD3	1.66	0.95
1:B:311:LEU:HB3	1:B:312:PRO:HD3	1.46	0.95
1:B:49:LEU:HD13	1:B:110:ARG:HE	1.30	0.95
1:B:573:ARG:HD2	1:B:573:ARG:H	1.32	0.95
1:B:760:PHE:HB3	1:B:807:LEU:HD12	1.48	0.95
1:B:261:SER:HA	1:B:264:ILE:HD12	1.49	0.95
1:A:866:THR:H	1:A:869:GLN:HE21	1.12	0.94
1:A:622:ILE:HD12	1:A:674:CYS:HA	1.46	0.94
1:A:786:ALA:O	1:A:787:LEU:HG	1.67	0.94
1:B:866:THR:H	1:B:869:GLN:HE21	1.14	0.93
1:A:248:PRO:CG	1:A:341:THR:HG21	1.97	0.93
1:B:909:MET:SD	1:B:937:ILE:HD12	2.09	0.93
1:A:783:LEU:HD12	1:A:783:LEU:H	1.32	0.92
1:B:311:LEU:HD11	1:B:761:ILE:HD12	1.51	0.92
1:A:762:ARG:HE	1:A:833:LEU:CD2	1.82	0.91
1:A:788:ILE:HB	1:A:789:PRO:HD2	1.52	0.91

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:757:MET:HE1	1:A:761:ILE:HD11	1.52	0.91
1:A:573:ARG:H	1:A:573:ARG:HD2	1.36	0.91
1:B:762:ARG:HE	1:B:833:LEU:CD2	1.83	0.90
1:A:59:ASP:OD2	1:A:62:VAL:HG23	1.72	0.90
1:A:849:VAL:O	1:A:853:ALA:CB	2.20	0.90
1:B:50:TRP:CD1	1:B:54:ILE:HD11	2.07	0.90
1:B:49:LEU:CD1	1:B:110:ARG:HE	1.84	0.89
1:B:844:VAL:HG22	1:B:907:ILE:HG21	1.52	0.89
1:A:762:ARG:HB3	1:A:837:TYR:CZ	2.07	0.89
1:B:762:ARG:HB3	1:B:837:TYR:CZ	2.07	0.89
1:B:80:GLU:HG2	1:B:82:GLU:HG2	1.52	0.89
1:A:395:VAL:HG12	1:A:402:ILE:HD11	1.55	0.89
1:B:622:ILE:HD12	1:B:674:CYS:HA	1.52	0.89
1:B:756:ASN:O	1:B:759:GLN:HB2	1.72	0.88
1:A:757:MET:HG2	1:A:760:PHE:CZ	2.08	0.88
1:A:777:LEU:O	1:A:781:LEU:HB3	1.73	0.88
1:A:899:MET:SD	1:A:962:LEU:HD13	2.13	0.88
1:B:383:SER:C	1:B:384:ILE:HD12	1.94	0.87
1:B:69:ALA:O	1:B:73:PHE:HB2	1.74	0.87
1:B:245:ASP:O	1:B:246:LYS:HG2	1.74	0.87
1:B:57:PHE:CZ	1:B:102:ALA:HB2	2.10	0.86
1:B:899:MET:SD	1:B:962:LEU:HD13	2.16	0.86
1:A:248:PRO:CG	1:A:341:THR:CG2	2.54	0.86
1:B:855:TRP:CA	1:B:859:ALA:HB2	2.00	0.86
1:A:899:MET:CE	1:A:970:VAL:HG23	2.06	0.86
1:A:765:ILE:HD13	1:A:765:ILE:O	1.76	0.86
1:A:895:GLU:HG3	1:A:960:LYS:HD3	1.56	0.86
1:A:120:LYS:HE2	1:A:728:LYS:HZ2	1.40	0.86
1:A:556:ARG:HG2	1:A:644:GLU:HG3	1.56	0.86
1:A:844:VAL:HG22	1:A:907:ILE:HG21	1.56	0.86
1:B:898:THR:CA	1:B:959:LEU:HD22	2.06	0.85
1:B:518:PRO:O	1:B:522:ILE:HD13	1.75	0.85
1:A:898:THR:CA	1:A:959:LEU:HD22	2.06	0.85
1:B:757:MET:HG2	1:B:760:PHE:CZ	2.12	0.84
1:A:39:ASN:OD1	1:A:226:THR:HB	1.75	0.84
1:B:895:GLU:HG3	1:B:960:LYS:HD3	1.57	0.84
1:B:787:LEU:HD11	1:B:900:ALA:HB1	1.59	0.84
1:A:567:ARG:HH11	1:A:570:PRO:HA	1.42	0.84
1:A:843:TYR:OH	1:A:976:PRO:HG2	1.78	0.84
1:A:865:VAL:HG13	1:A:869:GLN:HG3	1.59	0.84
1:B:342:LEU:HD11	1:B:718:ILE:HD11	1.60	0.84
1:B:816:ILE:HA	1:B:819:ARG:HH21	1.39	0.84

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:899:MET:CE	1:B:970:VAL:HG23	2.07	0.84
1:A:950:VAL:HG12	1:A:952:PRO:HD2	1.58	0.84
1:A:833:LEU:HA	1:A:836:ARG:HB2	1.60	0.84
1:B:857:MET:HB2	1:B:858:TYR:CD1	2.13	0.83
1:A:52:LEU:HD12	1:A:106:VAL:HG13	1.60	0.83
1:B:567:ARG:HH11	1:B:570:PRO:HA	1.43	0.83
1:B:421:ASN:ND2	1:B:423:SER:H	1.77	0.83
1:B:740:PHE:O	1:B:743:ILE:HD13	1.79	0.83
1:B:843:TYR:OH	1:B:976:PRO:HG2	1.78	0.83
1:A:857:MET:HB2	1:A:858:TYR:CD1	2.14	0.83
1:A:95:LEU:O	1:A:99:ILE:HD12	1.77	0.83
1:A:421:ASN:ND2	1:A:423:SER:H	1.77	0.83
1:A:245:ASP:O	1:A:246:LYS:HG3	1.77	0.83
1:B:157:ASP:HB2	1:B:214:ILE:HD13	1.61	0.83
1:B:833:LEU:HA	1:B:836:ARG:HB2	1.61	0.82
1:A:541:VAL:O	1:A:545:ILE:HD13	1.79	0.82
1:B:777:LEU:O	1:B:781:LEU:HB3	1.79	0.82
1:A:802:LEU:CD1	1:A:939:LEU:HG	2.10	0.82
1:B:788:ILE:HB	1:B:789:PRO:HD2	1.58	0.82
1:A:519:GLU:H	1:A:519:GLU:CD	1.83	0.82
1:B:802:LEU:CD1	1:B:939:LEU:HG	2.09	0.82
1:A:757:MET:HE2	1:A:761:ILE:CD1	2.04	0.81
1:A:878:GLU:HG3	1:A:879:ASP:H	1.45	0.81
1:B:248:PRO:HG2	1:B:341:THR:CG2	2.11	0.81
1:A:773:VAL:O	1:A:777:LEU:HG	1.80	0.81
1:B:773:VAL:O	1:B:777:LEU:HG	1.81	0.81
1:A:757:MET:CE	1:A:761:ILE:CD1	2.54	0.81
1:B:849:VAL:O	1:B:853:ALA:HB3	1.80	0.81
1:A:758:LYS:CB	1:A:828:LEU:HD11	2.10	0.81
1:B:245:ASP:O	1:B:246:LYS:CG	2.29	0.81
1:A:865:VAL:HG13	1:A:869:GLN:CG	2.11	0.80
1:B:363:VAL:HG11	1:B:448:LEU:HD22	1.64	0.79
1:B:802:LEU:HD13	1:B:939:LEU:CG	2.12	0.79
1:A:899:MET:HE1	1:A:966:GLN:O	1.81	0.79
1:B:878:GLU:HG3	1:B:879:ASP:H	1.45	0.79
1:A:802:LEU:HD13	1:A:939:LEU:CG	2.13	0.79
1:A:50:TRP:CD1	1:A:54:ILE:HD11	2.17	0.79
1:B:311:LEU:HD11	1:B:761:ILE:HD13	1.64	0.79
1:A:350:SER:OG	1:A:356:LEU:HD13	1.82	0.79
1:B:899:MET:HA	1:B:962:LEU:HD11	1.64	0.79
1:A:832:TRP:O	1:A:836:ARG:HG3	1.83	0.79
1:B:898:THR:OG1	1:B:959:LEU:HB3	1.83	0.79

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:898:THR:OG1	1:A:959:LEU:HB3	1.83	0.79
1:A:855:TRP:CA	1:A:859:ALA:HB2	2.01	0.78
1:B:787:LEU:HD21	1:B:901:LEU:N	1.97	0.78
1:B:899:MET:HE1	1:B:966:GLN:O	1.83	0.78
1:B:484:THR:HB	1:B:496:VAL:HG12	1.64	0.78
1:A:899:MET:HA	1:A:962:LEU:HD11	1.64	0.78
1:A:484:THR:HB	1:A:496:VAL:HG12	1.64	0.78
1:B:248:PRO:HG2	1:B:341:THR:HG23	1.64	0.78
1:A:771:GLU:O	1:A:775:ILE:HD13	1.83	0.78
1:A:852:ALA:HB1	1:A:896:PRO:O	1.84	0.78
1:A:488:SER:HB3	1:A:491:ARG:HH21	1.49	0.78
1:A:787:LEU:HD11	1:A:900:ALA:HB1	1.66	0.77
1:B:926:PRO:O	1:B:929:VAL:HG23	1.85	0.77
1:B:39:ASN:OD1	1:B:226:THR:HB	1.83	0.77
1:B:897:MET:HE1	1:B:958:LYS:HB3	1.67	0.77
1:B:519:GLU:CD	1:B:519:GLU:H	1.84	0.77
1:B:140:ILE:HD13	1:B:140:ILE:H	1.50	0.77
1:B:23:GLY:HA2	1:B:150:ILE:HD13	1.65	0.77
1:B:786:ALA:O	1:B:897:MET:HA	1.84	0.77
1:A:851:ALA:HB2	1:A:973:ILE:HG13	1.67	0.77
1:B:832:TRP:O	1:B:836:ARG:HG3	1.84	0.77
1:B:852:ALA:HB1	1:B:896:PRO:O	1.86	0.76
1:B:855:TRP:O	1:B:859:ALA:HB3	1.85	0.76
1:A:249:LEU:HD11	1:A:253:LEU:HD21	1.64	0.76
1:A:421:ASN:HD22	1:A:423:SER:H	1.33	0.76
1:A:112:ALA:O	1:A:116:ILE:HG12	1.85	0.76
1:A:807:LEU:O	1:A:810:ASN:HB2	1.84	0.76
1:A:99:ILE:O	1:A:103:ILE:HG22	1.86	0.75
1:A:95:LEU:HA	1:A:98:LEU:HD12	1.66	0.75
1:B:245:ASP:C	1:B:246:LYS:HG3	2.05	0.75
1:A:787:LEU:HD21	1:A:901:LEU:N	2.01	0.75
1:A:903:VAL:HG22	1:A:970:VAL:HG13	1.68	0.75
1:A:866:THR:H	1:A:869:GLN:NE2	1.85	0.75
1:B:851:ALA:HB2	1:B:973:ILE:HG13	1.68	0.75
1:B:807:LEU:O	1:B:810:ASN:HB2	1.86	0.74
1:B:54:ILE:O	1:B:58:GLU:HG3	1.87	0.74
1:B:488:SER:HB3	1:B:491:ARG:HH21	1.51	0.74
1:A:326:MET:HB3	1:A:331:ALA:HB3	1.67	0.74
1:A:708:ALA:HB3	1:A:709:PRO:HD3	1.68	0.74
1:B:75:LEU:HD11	1:B:300:VAL:HB	1.68	0.74
1:A:914:ASN:ND2	1:A:922:LEU:HD11	2.03	0.74
1:A:65:LEU:HD13	1:A:307:ILE:HG21	1.70	0.74

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:950:VAL:HG12	1:B:952:PRO:HD2	1.70	0.74
1:B:308:PRO:HG3	1:B:765:ILE:HG22	1.70	0.74
1:A:786:ALA:O	1:A:897:MET:HA	1.88	0.74
1:A:897:MET:HE1	1:A:958:LYS:HB3	1.70	0.74
1:B:573:ARG:H	1:B:573:ARG:CD	2.01	0.74
1:B:421:ASN:HD22	1:B:423:SER:H	1.33	0.73
1:B:40:GLU:OE1	1:B:143:ARG:HB2	1.88	0.73
1:A:171:THR:HG22	1:A:172:THR:HG23	1.71	0.73
1:A:852:ALA:HA	1:A:899:MET:CG	2.19	0.73
1:B:758:LYS:CB	1:B:828:LEU:HD11	2.15	0.73
1:B:788:ILE:HG13	1:B:791:GLN:HB2	1.70	0.73
1:B:749:GLU:O	1:B:753:ILE:HD13	1.89	0.73
1:A:102:ALA:O	1:A:106:VAL:HG23	1.89	0.73
1:B:914:ASN:ND2	1:B:922:LEU:HD11	2.02	0.73
1:B:245:ASP:C	1:B:246:LYS:CG	2.57	0.72
1:B:611:ILE:HD12	1:B:621:VAL:HG11	1.70	0.72
1:A:866:THR:N	1:A:869:GLN:HE21	1.85	0.72
1:A:100:ALA:O	1:A:104:VAL:HG23	1.89	0.72
1:B:899:MET:CE	1:B:966:GLN:O	2.36	0.72
1:B:903:VAL:HG22	1:B:970:VAL:HG13	1.71	0.72
1:A:51:GLU:HA	1:A:54:ILE:HD12	1.70	0.72
1:A:790:VAL:HG12	1:A:790:VAL:O	1.89	0.72
1:A:364:CYS:O	1:A:384:ILE:N	2.21	0.72
1:A:899:MET:CE	1:A:966:GLN:O	2.38	0.72
1:B:866:THR:H	1:B:869:GLN:NE2	1.86	0.72
1:B:852:ALA:HA	1:B:899:MET:CG	2.19	0.71
1:B:76:ALA:HB1	1:B:88:PHE:HA	1.72	0.71
1:B:311:LEU:CB	1:B:312:PRO:HD3	2.21	0.71
1:B:802:LEU:N	1:B:803:PRO:HD2	2.06	0.71
1:A:855:TRP:O	1:A:859:ALA:HB3	1.89	0.71
1:B:57:PHE:CZ	1:B:102:ALA:CB	2.73	0.71
1:A:926:PRO:O	1:A:929:VAL:HG23	1.89	0.71
1:B:326:MET:HB3	1:B:331:ALA:HB3	1.71	0.71
1:B:51:GLU:HA	1:B:54:ILE:HD12	1.70	0.71
1:B:611:ILE:HD12	1:B:621:VAL:HG21	1.72	0.71
1:B:349:CYS:HB3	1:B:624:ILE:CD1	2.20	0.70
1:A:88:PHE:O	1:A:91:PRO:HD2	1.92	0.70
1:A:112:ALA:HB1	1:A:334:ARG:HG2	1.74	0.70
1:B:790:VAL:HG12	1:B:790:VAL:O	1.92	0.70
1:A:573:ARG:CD	1:A:573:ARG:H	2.05	0.70
1:B:127:GLY:O	1:B:140:ILE:HD13	1.90	0.70
1:B:844:VAL:CG2	1:B:907:ILE:HG21	2.21	0.70

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:40:GLU:OE1	1:A:143:ARG:HB2	1.91	0.70
1:B:829:ILE:HD12	1:B:829:ILE:H	1.56	0.70
1:A:817:MET:O	1:A:819:ARG:N	2.24	0.70
1:B:866:THR:N	1:B:869:GLN:HE21	1.88	0.70
1:B:311:LEU:HB3	1:B:312:PRO:CD	2.20	0.69
1:B:349:CYS:HB3	1:B:624:ILE:HD13	1.74	0.69
1:A:181:THR:HB	1:A:183:GLU:HG3	1.74	0.69
1:B:355:THR:HG23	1:B:720:MET:HG2	1.74	0.69
1:A:42:PRO:O	1:A:43:ALA:C	2.31	0.69
1:A:802:LEU:CD1	1:A:939:LEU:CG	2.70	0.69
1:A:792:LEU:HA	1:A:795:VAL:CG2	2.23	0.69
1:B:129:VAL:HG22	1:B:140:ILE:HD12	1.72	0.69
1:B:342:LEU:CG	1:B:716:ILE:HG21	2.21	0.69
1:B:817:MET:O	1:B:819:ARG:N	2.24	0.69
1:B:852:ALA:HA	1:B:899:MET:HB3	1.75	0.69
1:A:784:PRO:CD	1:A:870:LEU:HD11	2.22	0.69
1:B:802:LEU:CD1	1:B:939:LEU:CG	2.69	0.69
1:A:788:ILE:HB	1:A:789:PRO:CD	2.22	0.69
1:B:895:GLU:CG	1:B:960:LYS:HD3	2.22	0.69
1:A:75:LEU:HD12	1:A:300:VAL:HB	1.74	0.69
1:A:828:LEU:HG	1:A:829:ILE:HG13	1.75	0.69
1:B:762:ARG:NH2	1:B:918:GLU:HG3	2.08	0.69
1:B:157:ASP:O	1:B:214:ILE:HD12	1.93	0.69
1:B:61:LEU:CD1	1:B:311:LEU:HD23	2.23	0.69
1:B:757:MET:O	1:B:761:ILE:HG12	1.92	0.69
1:B:853:ALA:O	1:B:856:PHE:HB2	1.93	0.69
1:B:412:GLU:OE1	1:B:529:ARG:HD2	1.93	0.69
1:A:844:VAL:CG2	1:A:907:ILE:HG21	2.22	0.68
1:B:299:ALA:HA	1:B:302:LEU:HD11	1.73	0.68
1:B:56:GLN:HG3	1:B:106:VAL:HG23	1.75	0.68
1:B:802:LEU:CD1	1:B:939:LEU:CD2	2.65	0.68
1:A:944:HIS:O	1:A:947:ILE:HG22	1.93	0.68
1:B:308:PRO:CG	1:B:765:ILE:HG22	2.23	0.68
1:A:895:GLU:CG	1:A:960:LYS:HD3	2.23	0.68
1:B:792:LEU:HA	1:B:795:VAL:CG2	2.23	0.68
1:A:931:ILE:HD12	1:A:931:ILE:H	1.59	0.68
1:A:489:ARG:CD	1:A:489:ARG:H	2.07	0.68
1:B:129:VAL:HG11	1:B:145:ILE:HD12	1.76	0.68
1:B:849:VAL:O	1:B:853:ALA:CB	2.41	0.68
1:B:49:LEU:CD1	1:B:110:ARG:HG3	2.23	0.68
1:A:111:ASN:O	1:A:114:ASN:HB2	1.92	0.68
1:B:67:LEU:O	1:B:71:ILE:HG12	1.94	0.68

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:311:LEU:HG	1:A:315:ILE:HD13	1.75	0.68
1:B:61:LEU:HD11	1:B:311:LEU:HD23	1.74	0.68
1:A:963:ASP:H	1:A:966:GLN:HE21	1.42	0.68
1:B:767:SER:O	1:B:771:GLU:HG3	1.93	0.68
1:B:898:THR:OG1	1:B:959:LEU:CB	2.42	0.68
1:A:899:MET:SD	1:A:962:LEU:CD1	2.82	0.68
1:A:412:GLU:OE1	1:A:529:ARG:HD2	1.93	0.68
1:A:802:LEU:N	1:A:803:PRO:HD2	2.09	0.67
1:B:489:ARG:CD	1:B:489:ARG:H	2.07	0.67
1:A:600:LEU:HD13	1:A:600:LEU:O	1.94	0.67
1:B:931:ILE:H	1:B:931:ILE:HD12	1.59	0.67
1:A:769:VAL:O	1:A:773:VAL:HG23	1.95	0.67
1:A:817:MET:C	1:A:819:ARG:H	1.97	0.67
1:A:69:ALA:O	1:A:73:PHE:HB2	1.94	0.67
1:A:762:ARG:NH2	1:A:918:GLU:HG3	2.10	0.67
1:B:817:MET:C	1:B:819:ARG:H	1.97	0.67
1:A:767:SER:O	1:A:771:GLU:HG3	1.93	0.67
1:B:784:PRO:CD	1:B:870:LEU:HD11	2.24	0.67
1:A:261:SER:O	1:A:264:ILE:HG12	1.93	0.67
1:A:829:ILE:H	1:A:829:ILE:HD12	1.59	0.67
1:B:786:ALA:HB2	1:B:849:VAL:HG13	1.77	0.67
1:A:120:LYS:NZ	1:A:723:GLY:O	2.24	0.67
1:A:788:ILE:CD1	1:A:791:GLN:HB2	2.24	0.66
1:A:786:ALA:HB2	1:A:849:VAL:HG13	1.77	0.66
1:A:898:THR:OG1	1:A:959:LEU:CB	2.43	0.66
1:A:65:LEU:HG	1:A:94:ILE:HD11	1.77	0.66
1:B:172:THR:HG23	1:B:489:ARG:HD2	1.77	0.66
1:B:252:LYS:HE3	1:B:252:LYS:HA	1.77	0.66
1:A:950:VAL:O	1:A:954:PRO:HD2	1.94	0.66
1:B:651:ARG:HG3	1:B:651:ARG:HH11	1.59	0.66
1:A:308:PRO:HB2	1:A:764:LEU:HD23	1.76	0.66
1:A:774:CYS:SG	1:A:786:ALA:HB3	2.35	0.66
1:A:334:ARG:HB3	1:A:334:ARG:HH11	1.60	0.66
1:A:622:ILE:CD1	1:A:674:CYS:HA	2.23	0.66
1:A:247:THR:HG23	1:A:250:GLN:HG2	1.78	0.66
1:B:863:PRO:HB2	1:B:865:VAL:HG23	1.76	0.66
1:A:567:ARG:NH1	1:A:571:PRO:HD3	2.10	0.66
1:B:86:THR:HA	1:B:89:VAL:HG23	1.77	0.66
1:A:639:ILE:HD11	1:A:641:ILE:HD12	1.77	0.66
1:B:944:HIS:O	1:B:947:ILE:HG22	1.96	0.66
1:B:893:ALA:HB1	1:B:895:GLU:OE1	1.95	0.66
1:B:897:MET:CE	1:B:958:LYS:HB3	2.25	0.66

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:963:ASP:H	1:B:966:GLN:HE21	1.44	0.65
1:A:852:ALA:HA	1:A:899:MET:HB3	1.78	0.65
1:A:54:ILE:O	1:A:58:GLU:HG3	1.96	0.65
1:A:680:GLU:HB3	1:A:681:PRO:HD2	1.77	0.65
1:A:651:ARG:HH11	1:A:651:ARG:HG3	1.61	0.65
1:B:314:VAL:HG11	1:B:804:ALA:HB1	1.76	0.65
1:A:762:ARG:HH22	1:A:918:GLU:HA	1.61	0.65
1:B:567:ARG:NH1	1:B:571:PRO:HD3	2.12	0.65
1:B:247:THR:HG21	1:B:337:PRO:HB2	1.79	0.65
1:B:611:ILE:CD1	1:B:621:VAL:HG11	2.25	0.65
1:A:784:PRO:HD2	1:A:870:LEU:HD11	1.77	0.65
1:B:899:MET:SD	1:B:962:LEU:CD1	2.84	0.65
1:A:25:THR:HA	1:A:132:ALA:HB3	1.77	0.65
1:A:803:PRO:O	1:A:806:ALA:CB	2.40	0.65
1:A:232:ILE:HD13	1:A:232:ILE:O	1.97	0.65
1:A:259:GLN:O	1:A:263:VAL:HG23	1.96	0.65
1:A:760:PHE:CB	1:A:807:LEU:HD12	2.25	0.64
1:B:832:TRP:CZ3	1:B:835:PHE:CD2	2.85	0.64
1:A:97:ILE:O	1:A:100:ALA:HB3	1.98	0.64
1:A:113:GLU:O	1:A:117:GLU:HG2	1.97	0.64
1:B:252:LYS:HA	1:B:252:LYS:CE	2.28	0.64
1:A:300:VAL:HA	1:A:303:ALA:HB3	1.80	0.64
1:B:55:GLU:HA	1:B:58:GLU:OE1	1.97	0.64
1:A:600:LEU:O	1:A:602:PRO:HD3	1.97	0.64
1:A:857:MET:HB2	1:A:858:TYR:CE1	2.32	0.64
1:A:795:VAL:HG22	1:A:901:LEU:HD11	1.80	0.64
1:B:303:ALA:O	1:B:306:ALA:HB3	1.98	0.64
1:A:853:ALA:O	1:A:856:PHE:HB2	1.97	0.64
1:A:342:LEU:HG	1:A:716:ILE:HG21	1.78	0.64
1:A:286:GLY:CA	1:A:290:ARG:HB3	2.27	0.64
1:A:300:VAL:HA	1:A:303:ALA:CB	2.28	0.64
1:B:857:MET:HB2	1:B:858:TYR:CE1	2.32	0.64
1:B:322:GLY:CA	1:B:325:ARG:HH21	2.11	0.64
1:A:893:ALA:HB1	1:A:895:GLU:OE1	1.96	0.64
1:B:214:ILE:H	1:B:214:ILE:HD12	1.62	0.64
1:A:600:LEU:HD13	1:A:600:LEU:C	2.19	0.64
1:B:715:GLU:O	1:B:716:ILE:HD13	1.97	0.63
1:B:778:THR:HB	1:B:783:LEU:O	1.98	0.63
1:B:567:ARG:NH1	1:B:570:PRO:HA	2.13	0.63
1:B:329:LYS:O	1:B:330:ASN:HB2	1.97	0.63
1:A:311:LEU:HG	1:A:315:ILE:CD1	2.28	0.63
1:B:298:ILE:HD12	1:B:299:ALA:N	2.14	0.63

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:322:GLY:N	1:A:325:ARG:HH21	1.97	0.63
1:A:783:LEU:HA	1:A:870:LEU:HG	1.81	0.63
1:A:832:TRP:CZ3	1:A:835:PHE:CD2	2.86	0.63
1:B:803:PRO:O	1:B:806:ALA:CB	2.42	0.63
1:B:781:LEU:HD23	1:B:782:GLY:N	2.13	0.63
1:B:298:ILE:O	1:B:302:LEU:HG	1.98	0.63
1:B:49:LEU:HD13	1:B:110:ARG:NE	2.10	0.63
1:A:784:PRO:CD	1:A:870:LEU:CD1	2.76	0.63
1:B:49:LEU:HD11	1:B:110:ARG:HG3	1.80	0.63
1:B:384:ILE:N	1:B:384:ILE:HD12	2.13	0.63
1:A:577:VAL:HG11	1:A:583:ARG:NH1	2.14	0.63
1:B:335:SER:HB3	1:B:338:SER:OG	1.98	0.63
1:B:774:CYS:SG	1:B:786:ALA:HB3	2.38	0.63
1:B:917:SER:OG	1:B:920:GLN:HB2	1.98	0.63
1:B:788:ILE:HB	1:B:789:PRO:CD	2.29	0.63
1:B:49:LEU:CD1	1:B:110:ARG:NE	2.60	0.63
1:A:55:GLU:HA	1:A:58:GLU:OE1	1.99	0.63
1:B:129:VAL:HG22	1:B:140:ILE:CD1	2.28	0.63
1:B:157:ASP:O	1:B:214:ILE:CD1	2.46	0.62
1:B:795:VAL:HG22	1:B:901:LEU:HD11	1.80	0.62
1:A:395:VAL:O	1:A:402:ILE:HD13	1.99	0.62
1:A:369:ILE:H	1:A:369:ILE:HD12	1.64	0.62
1:B:75:LEU:CD1	1:B:300:VAL:HB	2.29	0.62
1:A:620:ARG:NH2	1:A:622:ILE:HD11	2.13	0.62
1:B:622:ILE:CD1	1:B:674:CYS:HA	2.28	0.62
1:B:42:PRO:O	1:B:43:ALA:C	2.38	0.62
1:B:577:VAL:HG11	1:B:583:ARG:NH1	2.14	0.62
1:B:150:ILE:HD12	1:B:150:ILE:N	2.14	0.62
1:B:242:THR:O	1:B:712:LYS:HE2	1.99	0.62
1:A:188:ILE:HD12	1:A:188:ILE:N	2.15	0.62
1:A:2:GLU:O	1:A:3:ALA:HB3	2.00	0.62
1:A:866:THR:HG23	1:A:869:GLN:NE2	2.15	0.62
1:A:527:TYR:HB3	1:A:534:ARG:HG3	1.82	0.62
1:B:828:LEU:HG	1:B:829:ILE:HG13	1.80	0.62
1:B:969:MET:O	1:B:973:ILE:HG12	1.99	0.62
1:B:784:PRO:HD3	1:B:870:LEU:CD1	2.30	0.62
1:A:606:GLU:H	1:A:606:GLU:CD	2.04	0.62
1:A:121:GLU:OE2	1:A:121:GLU:O	2.16	0.61
1:B:319:LEU:HD21	1:B:337:PRO:HA	1.81	0.61
1:B:783:LEU:H	1:B:783:LEU:CD1	2.07	0.61
1:B:762:ARG:HH22	1:B:918:GLU:HA	1.63	0.61
1:A:248:PRO:CG	1:A:341:THR:HG23	2.30	0.61

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:952:PRO:O	1:B:956:ILE:HG12	2.00	0.61
1:A:917:SER:OG	1:A:920:GLN:HB2	1.99	0.61
1:A:795:VAL:O	1:A:799:THR:HB	2.01	0.61
1:A:897:MET:CE	1:A:958:LYS:HB3	2.30	0.61
1:B:769:VAL:O	1:B:773:VAL:HG23	2.00	0.61
1:B:781:LEU:C	1:B:781:LEU:HD23	2.21	0.61
1:A:120:LYS:HE2	1:A:728:LYS:HZ1	1.64	0.61
1:B:83:GLU:O	1:B:87:ALA:HB2	2.00	0.61
1:A:792:LEU:HA	1:A:795:VAL:HG23	1.82	0.61
1:A:969:MET:O	1:A:973:ILE:HG12	2.00	0.61
1:B:907:ILE:HD11	1:B:974:SER:HA	1.81	0.61
1:B:527:TYR:HB3	1:B:534:ARG:HG3	1.83	0.61
1:A:795:VAL:CG2	1:A:901:LEU:HD11	2.30	0.61
1:B:518:PRO:O	1:B:522:ILE:CD1	2.46	0.61
1:B:59:ASP:HB3	1:B:62:VAL:HG23	1.83	0.61
1:B:639:ILE:HD11	1:B:641:ILE:HD12	1.82	0.61
1:A:907:ILE:HD11	1:A:974:SER:HA	1.82	0.61
1:A:67:LEU:O	1:A:71:ILE:HG13	2.01	0.61
1:A:572:LYS:N	1:A:572:LYS:HD2	2.16	0.61
1:B:311:LEU:CD1	1:B:761:ILE:CD1	2.74	0.61
1:A:275:ASN:HA	1:A:295:TYR:OH	2.00	0.61
1:A:802:LEU:CD1	1:A:939:LEU:CD2	2.67	0.61
1:A:775:ILE:N	1:A:775:ILE:CD1	2.63	0.61
1:A:412:GLU:OE2	1:A:566:THR:HG21	2.01	0.61
1:A:121:GLU:O	1:A:121:GLU:CD	2.39	0.61
1:B:2:GLU:O	1:B:3:ALA:HB3	2.01	0.61
1:A:802:LEU:N	1:A:803:PRO:CD	2.64	0.61
1:A:759:GLN:HA	1:A:762:ARG:HD2	1.83	0.60
1:B:792:LEU:HA	1:B:795:VAL:HG23	1.83	0.60
1:A:528:VAL:HG21	1:A:541:VAL:HG11	1.82	0.60
1:A:520:GLY:O	1:A:524:ARG:HG3	2.00	0.60
1:A:315:ILE:O	1:A:319:LEU:HB2	2.01	0.60
1:B:813:ASP:OD2	1:B:813:ASP:N	2.29	0.60
1:B:867:TYR:O	1:B:871:THR:HG23	2.01	0.60
1:A:755:ASN:O	1:A:759:GLN:HG2	2.01	0.60
1:A:784:PRO:HD3	1:A:870:LEU:CD1	2.32	0.60
1:B:351:ASP:HB3	1:B:355:THR:OG1	2.01	0.60
1:B:572:LYS:HD2	1:B:572:LYS:N	2.16	0.60
1:B:749:GLU:HG3	1:B:753:ILE:HD13	1.81	0.60
1:B:968:LEU:O	1:B:972:LYS:HG2	2.01	0.60
1:A:310:GLY:O	1:A:314:VAL:HG23	2.02	0.60
1:A:778:THR:HB	1:A:783:LEU:O	2.01	0.60

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:795:VAL:CG2	1:B:901:LEU:HD11	2.30	0.60
1:A:617:ALA:O	1:A:821:PRO:HD3	2.01	0.60
1:B:802:LEU:N	1:B:803:PRO:CD	2.63	0.60
1:A:781:LEU:HD23	1:A:782:GLY:N	2.17	0.60
1:A:914:ASN:C	1:A:916:LEU:H	2.05	0.60
1:B:95:LEU:HD23	1:B:96:LEU:N	2.16	0.60
1:A:786:ALA:C	1:A:787:LEU:HG	2.21	0.60
1:B:786:ALA:C	1:B:787:LEU:HG	2.21	0.60
1:B:784:PRO:HD2	1:B:870:LEU:HD11	1.81	0.60
1:B:421:ASN:C	1:B:421:ASN:HD22	2.05	0.60
1:B:319:LEU:HD23	1:B:336:LEU:HB3	1.84	0.60
1:A:567:ARG:NH1	1:A:570:PRO:HA	2.13	0.60
1:A:326:MET:C	1:A:331:ALA:HB3	2.22	0.60
1:A:119:LEU:O	1:A:122:TYR:HB3	2.02	0.60
1:B:760:PHE:CB	1:B:807:LEU:HD12	2.29	0.59
1:A:489:ARG:NE	1:A:489:ARG:H	1.99	0.59
1:B:577:VAL:HG11	1:B:583:ARG:HH11	1.67	0.59
1:B:852:ALA:HA	1:B:899:MET:CB	2.32	0.59
1:B:865:VAL:HG13	1:B:869:GLN:HG3	1.83	0.59
1:A:94:ILE:O	1:A:98:LEU:HG	2.02	0.59
1:A:577:VAL:HG11	1:A:583:ARG:HH11	1.67	0.59
1:A:271:VAL:HG12	1:A:275:ASN:HD21	1.67	0.59
1:B:489:ARG:NE	1:B:489:ARG:H	2.00	0.59
1:A:303:ALA:O	1:A:306:ALA:HB3	2.03	0.59
1:B:855:TRP:CA	1:B:859:ALA:CB	2.73	0.59
1:B:57:PHE:CE1	1:B:102:ALA:HB2	2.36	0.59
1:B:826:GLU:HG2	1:B:827:PRO:HD2	1.85	0.59
1:B:759:GLN:HA	1:B:762:ARG:HD2	1.84	0.59
1:B:784:PRO:CD	1:B:870:LEU:CD1	2.80	0.59
1:A:95:LEU:HA	1:A:98:LEU:CD1	2.32	0.59
1:B:708:ALA:HB3	1:B:709:PRO:HD3	1.85	0.59
1:A:762:ARG:HB3	1:A:837:TYR:CE2	2.38	0.59
1:A:247:THR:HG23	1:A:250:GLN:CG	2.31	0.59
1:A:869:GLN:OE1	1:A:883:PHE:HB3	2.02	0.59
1:B:214:ILE:N	1:B:214:ILE:HD12	2.17	0.59
1:A:188:ILE:H	1:A:188:ILE:HD12	1.67	0.59
1:B:317:THR:O	1:B:321:LEU:HG	2.03	0.59
1:B:369:ILE:HD12	1:B:369:ILE:H	1.66	0.59
1:A:826:GLU:HG2	1:A:827:PRO:HD2	1.85	0.59
1:A:369:ILE:HD13	1:A:529:ARG:O	2.03	0.59
1:B:18:VAL:HG21	1:B:150:ILE:HD11	1.85	0.59
1:A:334:ARG:CB	1:A:334:ARG:HH11	2.16	0.59

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:369:ILE:HG12	1:A:528:VAL:CG1	2.33	0.58
1:A:334:ARG:NH1	1:A:334:ARG:HB3	2.18	0.58
1:B:950:VAL:O	1:B:954:PRO:HD2	2.02	0.58
1:B:322:GLY:HA2	1:B:325:ARG:HH21	1.68	0.58
1:A:311:LEU:O	1:A:315:ILE:HD13	2.03	0.58
1:A:413:LEU:HG	1:A:564:LEU:HD12	1.85	0.58
1:B:421:ASN:HD22	1:B:422:ASP:N	2.01	0.58
1:A:65:LEU:O	1:A:68:ALA:HB3	2.04	0.58
1:B:89:VAL:O	1:B:93:VAL:HG23	2.04	0.58
1:A:120:LYS:CE	1:A:728:LYS:NZ	2.63	0.58
1:B:869:GLN:OE1	1:B:883:PHE:HB3	2.03	0.58
1:B:53:VAL:HG22	1:B:106:VAL:HG21	1.85	0.58
1:A:41:LEU:HD11	1:A:233:GLY:HA2	1.85	0.58
1:A:247:THR:HG1	1:A:250:GLN:HB3	1.63	0.58
1:A:968:LEU:O	1:A:972:LYS:HG2	2.03	0.58
1:A:90:GLU:HG3	1:A:91:PRO:N	2.18	0.58
1:A:863:PRO:HB2	1:A:865:VAL:HG23	1.84	0.58
1:B:556:ARG:HG2	1:B:644:GLU:HG3	1.85	0.58
1:B:121:GLU:O	1:B:121:GLU:HG2	2.02	0.58
1:A:814:LEU:H	1:A:814:LEU:HD23	1.68	0.58
1:A:355:THR:HG23	1:A:720:MET:HG2	1.85	0.58
1:A:963:ASP:N	1:A:966:GLN:HE21	2.01	0.58
1:A:964:LEU:O	1:A:968:LEU:HG	2.03	0.58
1:B:865:VAL:HG13	1:B:869:GLN:CG	2.33	0.58
1:B:364:CYS:O	1:B:384:ILE:HD13	2.03	0.58
1:A:421:ASN:HD22	1:A:422:ASP:N	2.02	0.58
1:A:725:ALA:O	1:A:729:THR:HG23	2.04	0.58
1:B:113:GLU:O	1:B:117:GLU:HG2	2.03	0.58
1:A:801:GLY:C	1:A:803:PRO:HD2	2.24	0.58
1:B:964:LEU:O	1:B:968:LEU:HG	2.03	0.58
1:A:308:PRO:HG3	1:A:765:ILE:HG12	1.84	0.58
1:B:340:GLU:HG2	1:B:750:GLY:O	2.04	0.58
1:A:783:LEU:N	1:A:783:LEU:HD12	2.12	0.58
1:A:184:SER:HB3	1:A:627:ASP:OD2	2.04	0.58
1:B:833:LEU:HD12	1:B:836:ARG:HD3	1.86	0.58
1:B:783:LEU:HA	1:B:870:LEU:HG	1.86	0.58
1:B:663:LEU:CD1	1:B:663:LEU:H	2.17	0.58
1:B:783:LEU:N	1:B:783:LEU:HD12	2.09	0.57
1:A:39:ASN:O	1:A:40:GLU:HB3	2.02	0.57
1:B:342:LEU:HG	1:B:716:ILE:CG2	2.29	0.57
1:A:338:SER:HA	1:A:341:THR:HB	1.86	0.57
1:A:952:PRO:O	1:A:956:ILE:HG12	2.04	0.57

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:788:ILE:HD12	1:A:791:GLN:HB2	1.86	0.57
1:A:852:ALA:HA	1:A:899:MET:HG2	1.86	0.57
1:B:898:THR:OG1	1:B:959:LEU:CA	2.52	0.57
1:A:847:ALA:HB1	1:A:903:VAL:HG11	1.85	0.57
1:A:715:GLU:HB3	1:A:716:ILE:HD12	1.86	0.57
1:B:364:CYS:C	1:B:384:ILE:HD13	2.25	0.57
1:B:914:ASN:C	1:B:916:LEU:H	2.06	0.57
1:A:39:ASN:O	1:A:143:ARG:HA	2.05	0.57
1:A:760:PHE:O	1:A:764:LEU:HB2	2.05	0.57
1:B:795:VAL:O	1:B:799:THR:HB	2.04	0.57
1:B:91:PRO:O	1:B:94:ILE:HG22	2.04	0.57
1:A:833:LEU:HD12	1:A:836:ARG:HD3	1.87	0.57
1:A:341:THR:CG2	1:A:716:ILE:HD11	2.24	0.57
1:B:413:LEU:HG	1:B:564:LEU:HD12	1.86	0.57
1:A:898:THR:HA	1:A:959:LEU:CD2	2.22	0.57
1:B:528:VAL:HG21	1:B:541:VAL:HG11	1.85	0.57
1:A:867:TYR:O	1:A:871:THR:HG23	2.04	0.56
1:B:898:THR:HA	1:B:959:LEU:CD2	2.23	0.56
1:B:384:ILE:N	1:B:384:ILE:CD1	2.67	0.56
1:B:412:GLU:OE2	1:B:566:THR:HG21	2.04	0.56
1:A:315:ILE:CD1	1:A:760:PHE:HE1	2.18	0.56
1:A:782:GLY:HA3	1:A:871:THR:HB	1.88	0.56
1:B:520:GLY:O	1:B:524:ARG:HG3	2.04	0.56
1:A:756:ASN:O	1:A:759:GLN:HB2	2.05	0.56
1:B:762:ARG:HB3	1:B:837:TYR:CE2	2.40	0.56
1:B:749:GLU:HG3	1:B:753:ILE:CD1	2.34	0.56
1:A:802:LEU:HD11	1:A:939:LEU:HG	1.87	0.56
1:A:865:VAL:HG12	1:A:866:THR:O	2.05	0.56
1:A:311:LEU:CD2	1:A:761:ILE:HG23	2.36	0.56
1:B:311:LEU:O	1:B:315:ILE:HG12	2.06	0.56
1:B:620:ARG:NH2	1:B:622:ILE:HD11	2.19	0.56
1:B:606:GLU:CD	1:B:606:GLU:H	2.07	0.56
1:B:847:ALA:HB1	1:B:903:VAL:HG11	1.86	0.56
1:A:402:ILE:HD13	1:A:402:ILE:H	1.70	0.56
1:A:423:SER:HB3	1:A:437:VAL:O	2.05	0.56
1:B:866:THR:HG23	1:B:869:GLN:NE2	2.20	0.56
1:B:423:SER:HB3	1:B:437:VAL:O	2.04	0.56
1:B:369:ILE:HD13	1:B:529:ARG:O	2.05	0.56
1:B:633:ILE:HD11	1:B:652:ALA:CB	2.36	0.56
1:B:41:LEU:HD11	1:B:233:GLY:HA2	1.87	0.56
1:A:852:ALA:HA	1:A:899:MET:CB	2.34	0.56
1:A:364:CYS:HA	1:A:384:ILE:HB	1.87	0.56

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:651:ARG:HG3	1:B:651:ARG:NH1	2.21	0.56
1:A:360:GLN:HB3	1:A:389:TYR:CE2	2.40	0.56
1:B:463:SER:OG	1:B:465:VAL:HG22	2.05	0.56
1:B:600:LEU:HD13	1:B:600:LEU:C	2.25	0.56
1:B:814:LEU:H	1:B:814:LEU:HD23	1.70	0.56
1:A:840:ILE:CD1	1:A:840:ILE:N	2.69	0.56
1:B:825:LYS:O	1:B:826:GLU:C	2.42	0.56
1:A:963:ASP:OD2	1:A:964:LEU:N	2.38	0.56
1:A:355:THR:HG21	1:A:701:THR:HG22	1.88	0.56
1:A:825:LYS:O	1:A:826:GLU:C	2.43	0.56
1:A:781:LEU:C	1:A:781:LEU:HD23	2.26	0.56
1:A:421:ASN:C	1:A:421:ASN:HD22	2.07	0.56
1:B:157:ASP:CB	1:B:214:ILE:HD13	2.35	0.56
1:B:239:MET:CE	1:B:708:ALA:HB1	2.36	0.56
1:B:342:LEU:HD11	1:B:718:ILE:CD1	2.35	0.56
1:B:832:TRP:CZ3	1:B:835:PHE:HD2	2.23	0.56
1:B:792:LEU:HA	1:B:795:VAL:HG21	1.88	0.56
1:B:963:ASP:N	1:B:966:GLN:HE21	2.03	0.56
1:B:267:ILE:HD13	1:B:302:LEU:HD13	1.88	0.56
1:A:318:CYS:O	1:A:321:LEU:HB2	2.06	0.56
1:A:771:GLU:HG2	1:A:792:LEU:HD13	1.88	0.55
1:B:816:ILE:HA	1:B:819:ARG:NH2	2.16	0.55
1:A:309:GLU:OE2	1:A:796:ASN:HB2	2.06	0.55
1:A:759:GLN:HE22	1:A:762:ARG:HH11	1.54	0.55
1:B:848:THR:O	1:B:852:ALA:HB3	2.06	0.55
1:B:121:GLU:CG	1:B:121:GLU:O	2.54	0.55
1:A:355:THR:CG2	1:A:701:THR:HG22	2.36	0.55
1:A:832:TRP:CZ3	1:A:835:PHE:HD2	2.24	0.55
1:A:795:VAL:HA	1:A:799:THR:OG1	2.07	0.55
1:B:311:LEU:CD1	1:B:761:ILE:HD12	2.30	0.55
1:B:852:ALA:HA	1:B:899:MET:HG2	1.89	0.55
1:B:705:VAL:HG22	1:B:726:VAL:HG21	1.89	0.55
1:B:315:ILE:HD11	1:B:757:MET:CE	2.37	0.55
1:A:879:ASP:OD1	1:A:882:HIS:HB2	2.07	0.55
1:A:249:LEU:O	1:A:252:LYS:HB2	2.06	0.55
1:A:600:LEU:CD1	1:A:600:LEU:C	2.75	0.55
1:B:557:ASP:HB3	1:B:559:LEU:HG	1.89	0.55
1:A:757:MET:HG2	1:A:760:PHE:HZ	1.65	0.55
1:A:792:LEU:HA	1:A:795:VAL:HG21	1.87	0.55
1:A:899:MET:SD	1:A:962:LEU:HD22	2.46	0.55
1:B:771:GLU:HG2	1:B:792:LEU:HD13	1.89	0.55
1:A:329:LYS:O	1:A:330:ASN:HB2	2.07	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:662:PRO:HG2	1:B:665:GLU:HG2	1.87	0.55
1:B:762:ARG:HH22	1:B:918:GLU:HG3	1.72	0.55
1:A:894:PRO:HB2	1:A:959:LEU:C	2.26	0.55
1:A:866:THR:HG23	1:A:869:GLN:HE21	1.72	0.55
1:B:39:ASN:O	1:B:143:ARG:HA	2.07	0.55
1:A:463:SER:OG	1:A:465:VAL:HG22	2.06	0.55
1:A:848:THR:O	1:A:852:ALA:HB3	2.07	0.54
1:A:41:LEU:HB3	1:A:44:GLU:OE2	2.07	0.54
1:A:460:ARG:CZ	1:A:461:ASN:HD21	2.21	0.54
1:A:757:MET:O	1:A:761:ILE:HG13	2.07	0.54
1:B:248:PRO:HA	1:B:251:GLN:HB2	1.88	0.54
1:B:333:VAL:HG11	1:B:339:VAL:HG22	1.89	0.54
1:B:899:MET:SD	1:B:962:LEU:HD22	2.47	0.54
1:B:248:PRO:O	1:B:249:LEU:C	2.43	0.54
1:A:120:LYS:HE3	1:A:728:LYS:HE3	1.89	0.54
1:B:879:ASP:OD1	1:B:882:HIS:HB2	2.07	0.54
1:B:239:MET:HE2	1:B:708:ALA:HB1	1.89	0.54
1:A:355:THR:HG23	1:A:720:MET:CG	2.36	0.54
1:B:340:GLU:HA	1:B:750:GLY:HA2	1.88	0.54
1:B:759:GLN:HE22	1:B:762:ARG:HH11	1.55	0.54
1:A:792:LEU:CA	1:A:795:VAL:HG23	2.38	0.54
1:A:898:THR:OG1	1:A:959:LEU:CA	2.55	0.54
1:B:722:SER:OG	1:B:738:ASP:OD2	2.20	0.54
1:A:315:ILE:HD11	1:A:760:PHE:HE1	1.71	0.54
1:A:315:ILE:CD1	1:A:760:PHE:CE1	2.91	0.54
1:B:338:SER:HA	1:B:341:THR:HB	1.90	0.54
1:B:815:ASP:O	1:B:819:ARG:NE	2.41	0.54
1:A:782:GLY:O	1:A:783:LEU:O	2.26	0.54
1:B:787:LEU:CD1	1:B:900:ALA:HB1	2.34	0.54
1:A:914:ASN:HD22	1:A:922:LEU:HD11	1.73	0.54
1:B:369:ILE:HG12	1:B:528:VAL:CG1	2.38	0.54
1:B:912:ALA:HB1	1:B:933:LEU:HD11	1.90	0.54
1:B:755:ASN:O	1:B:759:GLN:HG2	2.07	0.54
1:A:897:MET:HE1	1:A:958:LYS:HE3	1.90	0.54
1:B:857:MET:CE	1:B:867:TYR:HA	2.38	0.54
1:A:165:ILE:HG21	1:A:168:ILE:HD11	1.89	0.54
1:A:789:PRO:C	1:A:791:GLN:H	2.11	0.54
1:B:112:ALA:O	1:B:116:ILE:HD12	2.07	0.54
1:B:140:ILE:N	1:B:140:ILE:HD13	2.20	0.54
1:A:180:LEU:O	1:A:180:LEU:HD12	2.08	0.54
1:A:762:ARG:NH2	1:A:833:LEU:CD1	2.71	0.53
1:B:319:LEU:HD23	1:B:336:LEU:C	2.28	0.53

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:116:ILE:HD11	1:B:334:ARG:HG2	1.90	0.53
1:A:840:ILE:HD13	1:A:840:ILE:H	1.74	0.53
1:A:546:LEU:HA	1:A:549:ILE:HD12	1.90	0.53
1:B:979:GLY:O	1:B:983:ILE:HG12	2.08	0.53
1:B:761:ILE:O	1:B:765:ILE:HG12	2.08	0.53
1:B:795:VAL:HA	1:B:799:THR:OG1	2.08	0.53
1:A:326:MET:HB3	1:A:331:ALA:CB	2.37	0.53
1:B:914:ASN:HD22	1:B:922:LEU:HD11	1.70	0.53
1:B:792:LEU:CA	1:B:795:VAL:HG23	2.39	0.53
1:B:460:ARG:CZ	1:B:461:ASN:HD21	2.21	0.53
1:A:189:LYS:HE3	1:A:207:MET:O	2.09	0.53
1:A:979:GLY:O	1:A:983:ILE:HG12	2.08	0.53
1:A:395:VAL:CG1	1:A:402:ILE:HD11	2.35	0.53
1:A:249:LEU:HB2	1:A:340:GLU:OE1	2.08	0.53
1:A:86:THR:O	1:A:89:VAL:HB	2.08	0.53
1:A:129:VAL:HG12	1:A:151:VAL:HG22	1.90	0.53
1:A:161:ALA:HA	1:A:210:SER:HB2	1.91	0.53
1:B:847:ALA:HA	1:B:973:ILE:HG21	1.90	0.53
1:B:62:VAL:CG1	1:B:98:LEU:HD22	2.39	0.53
1:A:847:ALA:HA	1:A:973:ILE:CG2	2.38	0.53
1:A:66:LEU:HD13	1:A:98:LEU:HD13	1.91	0.53
1:B:129:VAL:HG12	1:B:151:VAL:HG22	1.90	0.53
1:B:65:LEU:HB3	1:B:98:LEU:HD21	1.91	0.53
1:A:762:ARG:NH2	1:A:833:LEU:HD11	2.24	0.53
1:A:718:ILE:HD12	1:A:718:ILE:N	2.23	0.53
1:B:771:GLU:OE2	1:B:792:LEU:HD13	2.07	0.53
1:B:894:PRO:HB2	1:B:959:LEU:C	2.29	0.53
1:B:963:ASP:OD2	1:B:964:LEU:N	2.41	0.53
1:A:249:LEU:O	1:A:252:LYS:N	2.39	0.53
1:A:306:ALA:O	1:A:308:PRO:HD3	2.08	0.53
1:B:762:ARG:NH2	1:B:833:LEU:CD1	2.71	0.53
1:B:161:ALA:HA	1:B:210:SER:HB2	1.91	0.53
1:A:315:ILE:HD12	1:A:760:PHE:CE1	2.44	0.53
1:B:735:LEU:HD11	1:B:743:ILE:CD1	2.17	0.53
1:A:847:ALA:HA	1:A:973:ILE:HG21	1.90	0.53
1:B:725:ALA:O	1:B:729:THR:HG23	2.09	0.52
1:B:248:PRO:HD2	1:B:341:THR:HG21	1.91	0.52
1:A:782:GLY:CA	1:A:871:THR:HA	2.39	0.52
1:B:23:GLY:HA2	1:B:150:ILE:CD1	2.39	0.52
1:A:116:ILE:HG13	1:A:334:ARG:HD2	1.91	0.52
1:A:286:GLY:HA2	1:A:290:ARG:HB3	1.90	0.52
1:B:964:LEU:C	1:B:964:LEU:HD23	2.30	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:633:ILE:HD11	1:A:652:ALA:CB	2.39	0.52
1:A:311:LEU:HD21	1:A:761:ILE:HG23	1.90	0.52
1:B:743:ILE:HD12	1:B:743:ILE:N	2.25	0.52
1:B:847:ALA:HA	1:B:973:ILE:CG2	2.39	0.52
1:B:856:PHE:CZ	1:B:896:PRO:HG3	2.45	0.52
1:B:897:MET:HE1	1:B:958:LYS:HE3	1.90	0.52
1:B:633:ILE:HD11	1:B:652:ALA:HB1	1.91	0.52
1:B:662:PRO:HG2	1:B:665:GLU:CG	2.40	0.52
1:A:754:TYR:O	1:A:757:MET:N	2.37	0.52
1:B:545:ILE:O	1:B:549:ILE:HG12	2.09	0.52
1:B:96:LEU:C	1:B:96:LEU:HD23	2.30	0.52
1:A:978:ILE:O	1:A:982:GLU:HB2	2.09	0.52
1:B:789:PRO:C	1:B:791:GLN:H	2.13	0.52
1:B:865:VAL:HG12	1:B:866:THR:O	2.08	0.52
1:B:56:GLN:HG3	1:B:106:VAL:CG2	2.39	0.52
1:B:112:ALA:C	1:B:116:ILE:HD12	2.30	0.52
1:A:75:LEU:CD1	1:A:300:VAL:HB	2.38	0.52
1:B:600:LEU:HD13	1:B:600:LEU:O	2.09	0.52
1:B:315:ILE:HD11	1:B:757:MET:HE3	1.92	0.52
1:A:248:PRO:CB	1:A:341:THR:HG23	2.40	0.52
1:A:116:ILE:CD1	1:A:332:ILE:HG21	2.40	0.52
1:A:42:PRO:HD3	1:A:228:VAL:HG12	1.92	0.52
1:B:291:GLY:O	1:B:292:ALA:HB2	2.09	0.52
1:B:988:ALA:HA	1:B:992:LEU:HB2	1.91	0.52
1:A:859:ALA:O	1:A:861:ASP:N	2.43	0.52
1:B:847:ALA:HB1	1:B:903:VAL:CG1	2.39	0.52
1:A:51:GLU:O	1:A:54:ILE:HB	2.10	0.52
1:A:489:ARG:H	1:A:489:ARG:HD3	1.75	0.52
1:A:611:ILE:CD1	1:A:639:ILE:HD12	2.40	0.52
1:A:651:ARG:HG3	1:A:651:ARG:NH1	2.23	0.52
1:B:94:ILE:HD13	1:B:98:LEU:HG	1.92	0.52
1:A:189:LYS:HD2	1:A:205:LYS:O	2.10	0.52
1:B:249:LEU:O	1:B:249:LEU:HD12	2.10	0.52
1:A:912:ALA:HB1	1:A:933:LEU:HD11	1.91	0.52
1:B:735:LEU:HD13	1:B:743:ILE:HD11	1.86	0.51
1:B:48:SER:OG	1:B:51:GLU:HG2	2.09	0.51
1:A:369:ILE:N	1:A:369:ILE:HD12	2.25	0.51
1:A:680:GLU:O	1:A:683:HIS:HB2	2.10	0.51
1:B:329:LYS:O	1:B:330:ASN:CB	2.59	0.51
1:B:458:GLU:OE2	1:B:460:ARG:HG2	2.10	0.51
1:A:737:ASP:O	1:A:738:ASP:HB2	2.10	0.51
1:B:111:ASN:O	1:B:114:ASN:HB2	2.10	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:249:LEU:HG	1:B:250:GLN:N	2.25	0.51
1:A:747:VAL:HG12	1:A:817:MET:HE1	1.91	0.51
1:B:782:GLY:HA3	1:B:871:THR:HB	1.92	0.51
1:B:792:LEU:C	1:B:795:VAL:HG23	2.31	0.51
1:A:792:LEU:C	1:A:795:VAL:HG23	2.31	0.51
1:A:44:GLU:N	1:A:44:GLU:CD	2.64	0.51
1:B:341:THR:HB	1:B:716:ILE:HD11	1.91	0.51
1:B:762:ARG:NH2	1:B:833:LEU:HD11	2.25	0.51
1:B:801:GLY:C	1:B:803:PRO:HD2	2.31	0.51
1:A:383:SER:O	1:A:384:ILE:HD13	2.10	0.51
1:A:611:ILE:HD11	1:A:639:ILE:HD12	1.92	0.51
1:B:311:LEU:CD1	1:B:761:ILE:HD13	2.35	0.51
1:A:94:ILE:HG23	1:A:95:LEU:N	2.26	0.51
1:A:964:LEU:HD23	1:A:964:LEU:C	2.31	0.51
1:B:866:THR:HG23	1:B:869:GLN:HE21	1.76	0.51
1:A:66:LEU:O	1:A:69:ALA:HB3	2.11	0.51
1:A:94:ILE:HG23	1:A:95:LEU:H	1.76	0.51
1:B:663:LEU:N	1:B:663:LEU:CD1	2.74	0.51
1:B:44:GLU:N	1:B:44:GLU:CD	2.63	0.51
1:B:463:SER:OG	1:B:466:GLU:HG3	2.10	0.51
1:A:60:LEU:HD23	1:A:257:GLY:HA3	1.93	0.51
1:A:60:LEU:N	1:A:63:ARG:HH21	2.09	0.51
1:A:810:ASN:OD1	1:A:930:ASN:ND2	2.43	0.51
1:B:369:ILE:HD11	1:B:593:PHE:CD1	2.46	0.51
1:A:600:LEU:N	1:A:600:LEU:HD12	2.26	0.51
1:A:463:SER:OG	1:A:466:GLU:HG3	2.11	0.51
1:A:633:ILE:HD11	1:A:652:ALA:HB1	1.93	0.51
1:B:810:ASN:OD1	1:B:930:ASN:ND2	2.44	0.51
1:B:748:GLU:N	1:B:817:MET:HE1	2.26	0.51
1:B:355:THR:HG23	1:B:720:MET:CG	2.40	0.51
1:A:600:LEU:CD1	1:A:600:LEU:O	2.58	0.51
1:A:855:TRP:CA	1:A:859:ALA:CB	2.74	0.51
1:A:65:LEU:HG	1:A:94:ILE:CD1	2.40	0.51
1:A:91:PRO:HA	1:A:94:ILE:HG22	1.91	0.50
1:B:41:LEU:HB3	1:B:44:GLU:OE2	2.11	0.50
1:A:988:ALA:HA	1:A:992:LEU:HB2	1.92	0.50
1:B:749:GLU:OE2	1:B:753:ILE:HD11	2.10	0.50
1:A:907:ILE:HD11	1:A:974:SER:CA	2.41	0.50
1:A:93:VAL:O	1:A:97:ILE:HG12	2.12	0.50
1:B:235:ILE:HG21	1:B:705:VAL:HG12	1.93	0.50
1:A:222:ILE:HG23	1:A:222:ILE:O	2.11	0.50
1:B:939:LEU:O	1:B:943:LEU:HG	2.12	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:783:LEU:HG	1:B:870:LEU:HD23	1.94	0.50
1:A:104:VAL:O	1:A:107:TRP:HB3	2.12	0.50
1:A:914:ASN:C	1:A:916:LEU:N	2.64	0.50
1:B:369:ILE:HD12	1:B:369:ILE:N	2.26	0.50
1:B:238:GLN:O	1:B:242:THR:HG23	2.10	0.50
1:A:813:ASP:OD2	1:A:813:ASP:N	2.33	0.50
1:A:762:ARG:HH22	1:A:918:GLU:CA	2.25	0.50
1:B:802:LEU:HD11	1:B:939:LEU:HG	1.89	0.50
1:A:963:ASP:H	1:A:966:GLN:NE2	2.08	0.50
1:B:855:TRP:HD1	1:B:966:GLN:OE1	1.94	0.50
1:A:116:ILE:HD11	1:A:332:ILE:HG21	1.92	0.50
1:B:172:THR:CG2	1:B:489:ARG:HD2	2.41	0.50
1:A:291:GLY:O	1:A:292:ALA:HB2	2.11	0.50
1:B:237:ASP:O	1:B:238:GLN:C	2.48	0.50
1:A:252:LYS:HA	1:A:252:LYS:HE2	1.94	0.50
1:B:60:LEU:CD2	1:B:257:GLY:HA3	2.42	0.50
1:B:222:ILE:HG23	1:B:222:ILE:O	2.10	0.50
1:B:103:ILE:HG23	1:B:104:VAL:N	2.27	0.50
1:A:771:GLU:OE2	1:A:792:LEU:HD13	2.12	0.50
1:A:307:ILE:O	1:A:309:GLU:N	2.45	0.50
1:B:116:ILE:HG13	1:B:332:ILE:HG21	1.93	0.50
1:B:320:ALA:O	1:B:324:ARG:HG2	2.12	0.50
1:B:338:SER:HB2	1:B:716:ILE:HD12	1.94	0.50
1:B:947:ILE:HD11	1:B:957:PHE:CZ	2.46	0.50
1:A:82:GLU:HG3	1:A:83:GLU:N	2.26	0.50
1:B:893:ALA:O	1:B:896:PRO:HD2	2.12	0.50
1:A:489:ARG:CD	1:A:489:ARG:N	2.75	0.50
1:B:236:ARG:O	1:B:239:MET:HB3	2.11	0.50
1:B:352:LYS:HE2	1:B:627:ASP:HB3	1.93	0.50
1:B:350:SER:OG	1:B:356:LEU:HD13	2.12	0.50
1:A:969:MET:HE2	1:A:973:ILE:HG12	1.94	0.50
1:B:782:GLY:C	1:B:871:THR:HA	2.32	0.50
1:B:334:ARG:HD3	1:B:731:SER:O	2.12	0.50
1:B:505:ARG:HB3	1:B:508:VAL:CG2	2.42	0.50
1:A:762:ARG:HH22	1:A:918:GLU:HG3	1.77	0.49
1:A:311:LEU:HB3	1:A:312:PRO:HD3	1.95	0.49
1:B:757:MET:HG2	1:B:760:PHE:HZ	1.71	0.49
1:A:775:ILE:H	1:A:775:ILE:HD13	1.77	0.49
1:B:465:VAL:HG23	1:B:466:GLU:N	2.27	0.49
1:A:505:ARG:HB3	1:A:508:VAL:CG2	2.42	0.49
1:A:624:ILE:CG2	1:A:684:LYS:HG2	2.42	0.49
1:B:743:ILE:H	1:B:743:ILE:HD12	1.77	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:787:LEU:HD23	1:B:791:GLN:OE1	2.12	0.49
1:B:907:ILE:HD11	1:B:974:SER:CA	2.40	0.49
1:A:326:MET:O	1:A:331:ALA:HB3	2.12	0.49
1:A:261:SER:O	1:A:264:ILE:HB	2.11	0.49
1:B:600:LEU:CD1	1:B:600:LEU:C	2.80	0.49
1:A:127:GLY:HA3	1:A:145:ILE:HD11	1.94	0.49
1:B:306:ALA:O	1:B:308:PRO:HD3	2.12	0.49
1:B:340:GLU:HA	1:B:750:GLY:CA	2.42	0.49
1:A:947:ILE:HD11	1:A:957:PHE:CZ	2.46	0.49
1:A:369:ILE:HD11	1:A:593:PHE:CD1	2.47	0.49
1:B:352:LYS:HE2	1:B:627:ASP:CB	2.42	0.49
1:B:25:THR:HA	1:B:132:ALA:HB3	1.94	0.49
1:A:784:PRO:HD3	1:A:870:LEU:HD12	1.94	0.49
1:A:342:LEU:HD22	1:A:747:VAL:CG2	2.42	0.49
1:B:383:SER:CA	1:B:384:ILE:HD12	2.42	0.49
1:B:518:PRO:C	1:B:522:ILE:HD13	2.33	0.49
1:A:458:GLU:OE2	1:A:460:ARG:HG2	2.12	0.49
1:A:320:ALA:O	1:A:324:ARG:HG2	2.11	0.49
1:B:978:ILE:O	1:B:982:GLU:HB2	2.12	0.49
1:A:939:LEU:O	1:A:943:LEU:HG	2.12	0.49
1:A:957:PHE:O	1:A:959:LEU:HG	2.13	0.49
1:B:953:LEU:HB2	1:B:954:PRO:HD3	1.95	0.49
1:B:633:ILE:HD12	1:B:642:PHE:CE1	2.47	0.49
1:B:654:THR:OG1	1:B:657:GLU:HG3	2.12	0.49
1:A:108:GLN:NE2	1:A:108:GLN:HA	2.26	0.49
1:B:402:ILE:HD12	1:B:403:ARG:C	2.32	0.49
1:A:782:GLY:C	1:A:871:THR:HA	2.32	0.49
1:A:847:ALA:HB1	1:A:903:VAL:CG1	2.42	0.49
1:B:244:GLN:O	1:B:246:LYS:N	2.45	0.49
1:B:489:ARG:H	1:B:489:ARG:HD3	1.75	0.49
1:A:832:TRP:NE1	1:A:988:ALA:HB2	2.27	0.49
1:B:782:GLY:CA	1:B:871:THR:HA	2.43	0.49
1:B:140:ILE:HD11	1:B:145:ILE:HD11	1.95	0.49
1:B:749:GLU:C	1:B:751:ARG:H	2.16	0.49
1:B:795:VAL:HA	1:B:799:THR:CB	2.43	0.49
1:B:267:ILE:HG21	1:B:302:LEU:CD1	2.42	0.49
1:A:757:MET:HE1	1:A:761:ILE:CD1	2.32	0.49
1:B:333:VAL:HG12	1:B:335:SER:O	2.13	0.49
1:A:856:PHE:CZ	1:A:896:PRO:HG3	2.48	0.49
1:A:715:GLU:C	1:A:716:ILE:HD12	2.32	0.49
1:B:55:GLU:O	1:B:58:GLU:HB2	2.13	0.49
1:A:179:ILE:HG22	1:A:179:ILE:O	2.13	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:762:ARG:HH22	1:B:918:GLU:CA	2.26	0.48
1:A:787:LEU:CD1	1:A:900:ALA:HB1	2.40	0.48
1:A:855:TRP:HD1	1:A:966:GLN:OE1	1.96	0.48
1:B:622:ILE:O	1:B:624:ILE:HD12	2.13	0.48
1:B:763:TYR:CE1	1:B:912:ALA:HB2	2.48	0.48
1:A:320:ALA:O	1:A:323:THR:HB	2.13	0.48
1:A:179:ILE:HG23	1:A:724:THR:CG2	2.43	0.48
1:A:705:VAL:HG22	1:A:726:VAL:HG21	1.94	0.48
1:A:758:LYS:O	1:A:762:ARG:HG3	2.13	0.48
1:B:758:LYS:O	1:B:762:ARG:HG3	2.14	0.48
1:A:48:SER:OG	1:A:51:GLU:HG2	2.12	0.48
1:B:358:THR:OG1	1:B:602:PRO:HG2	2.13	0.48
1:A:708:ALA:HB3	1:A:709:PRO:CD	2.39	0.48
1:A:305:ALA:C	1:A:768:ASN:HD22	2.17	0.48
1:A:79:GLU:HG2	1:A:84:THR:HB	1.94	0.48
1:A:795:VAL:HA	1:A:799:THR:CB	2.43	0.48
1:B:855:TRP:HZ3	1:B:856:PHE:CE2	2.32	0.48
1:A:48:SER:H	1:A:51:GLU:HG2	1.79	0.48
1:A:697:ILE:N	1:A:697:ILE:HD12	2.29	0.48
1:A:833:LEU:CA	1:A:836:ARG:HB2	2.39	0.48
1:B:611:ILE:CD1	1:B:621:VAL:HG21	2.41	0.48
1:A:931:ILE:HD12	1:A:931:ILE:N	2.28	0.48
1:A:314:VAL:HG11	1:A:804:ALA:HB1	1.96	0.48
1:B:832:TRP:NE1	1:B:988:ALA:HB2	2.28	0.48
1:B:799:THR:HG21	1:B:905:VAL:HG22	1.95	0.48
1:A:878:GLU:HG3	1:A:879:ASP:N	2.23	0.48
1:B:48:SER:H	1:B:51:GLU:HG2	1.79	0.48
1:B:363:VAL:CG1	1:B:448:LEU:HD13	2.44	0.48
1:B:914:ASN:C	1:B:916:LEU:N	2.65	0.48
1:A:235:ILE:HG21	1:A:705:VAL:HG12	1.94	0.48
1:B:267:ILE:HG21	1:B:302:LEU:HD13	1.94	0.48
1:B:46:GLY:O	1:B:47:LYS:C	2.52	0.48
1:A:833:LEU:HD11	1:A:837:TYR:HE1	1.79	0.48
1:A:947:ILE:HD11	1:A:957:PHE:CE2	2.49	0.48
1:B:787:LEU:HD11	1:B:900:ALA:CB	2.37	0.48
1:A:42:PRO:O	1:A:43:ALA:O	2.31	0.48
1:B:813:ASP:HB2	1:B:815:ASP:OD1	2.14	0.48
1:B:762:ARG:HH22	1:B:918:GLU:CB	2.27	0.48
1:A:787:LEU:HD23	1:A:791:GLN:OE1	2.13	0.48
1:B:489:ARG:CD	1:B:489:ARG:N	2.75	0.48
1:B:663:LEU:N	1:B:663:LEU:HD12	2.29	0.48
1:A:359:ASN:O	1:A:359:ASN:ND2	2.46	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:756:ASN:O	1:B:759:GLN:CB	2.55	0.47
1:A:855:TRP:O	1:A:862:GLY:HA3	2.14	0.47
1:B:784:PRO:HD3	1:B:870:LEU:HD12	1.94	0.47
1:B:852:ALA:O	1:B:855:TRP:HB3	2.14	0.47
1:B:899:MET:O	1:B:903:VAL:HG23	2.14	0.47
1:B:84:THR:HA	1:B:87:ALA:HB3	1.95	0.47
1:A:387:SER:HB2	1:A:602:PRO:HB2	1.96	0.47
1:A:413:LEU:C	1:A:413:LEU:HD23	2.35	0.47
1:B:757:MET:HE2	1:B:761:ILE:HD11	1.95	0.47
1:A:247:THR:CG2	1:A:250:GLN:HB3	2.44	0.47
1:B:857:MET:HE3	1:B:867:TYR:HA	1.96	0.47
1:B:880:HIS:HB3	1:B:881:PRO:HD3	1.95	0.47
1:B:150:ILE:CD1	1:B:150:ILE:N	2.78	0.47
1:A:763:TYR:CE1	1:A:912:ALA:HB2	2.49	0.47
1:B:848:THR:O	1:B:852:ALA:CB	2.62	0.47
1:B:854:TRP:CE3	1:B:855:TRP:N	2.82	0.47
1:B:783:LEU:HG	1:B:870:LEU:CD2	2.43	0.47
1:B:899:MET:HE1	1:B:970:VAL:CG2	2.30	0.47
1:B:947:ILE:HD11	1:B:957:PHE:CE2	2.49	0.47
1:B:898:THR:HG1	1:B:959:LEU:HA	1.78	0.47
1:B:48:SER:H	1:B:51:GLU:CG	2.27	0.47
1:A:356:LEU:CD1	1:A:740:PHE:CE1	2.97	0.47
1:B:633:ILE:HD12	1:B:642:PHE:HE1	1.79	0.47
1:A:840:ILE:HD13	1:A:840:ILE:N	2.29	0.47
1:B:51:GLU:O	1:B:54:ILE:HB	2.14	0.47
1:A:73:PHE:O	1:A:77:TRP:HB2	2.15	0.47
1:A:247:THR:CB	1:A:250:GLN:HB3	2.38	0.47
1:A:893:ALA:O	1:A:896:PRO:HD2	2.14	0.47
1:B:782:GLY:O	1:B:783:LEU:O	2.33	0.47
1:B:917:SER:CB	1:B:920:GLN:HB2	2.44	0.47
1:A:813:ASP:HB2	1:A:815:ASP:OD1	2.14	0.47
1:A:84:THR:HA	1:A:87:ALA:HB3	1.97	0.47
1:A:340:GLU:OE2	1:A:344:CYS:SG	2.69	0.47
1:B:931:ILE:N	1:B:931:ILE:HD12	2.27	0.47
1:B:639:ILE:HD11	1:B:641:ILE:CD1	2.43	0.47
1:B:735:LEU:HD22	1:B:742:THR:HB	1.97	0.47
1:A:895:GLU:H	1:A:895:GLU:CD	2.17	0.47
1:B:898:THR:OG1	1:B:959:LEU:HA	2.15	0.47
1:B:880:HIS:N	1:B:881:PRO:HD2	2.30	0.47
1:A:545:ILE:CD1	1:A:545:ILE:N	2.77	0.47
1:B:315:ILE:CD1	1:B:757:MET:CE	2.93	0.47
1:A:782:GLY:CA	1:A:871:THR:HB	2.45	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:787:LEU:HD21	1:B:901:LEU:CA	2.44	0.47
1:A:971:LEU:O	1:A:975:LEU:HD23	2.14	0.47
1:B:18:VAL:CG2	1:B:150:ILE:HD11	2.45	0.47
1:A:762:ARG:HH22	1:A:918:GLU:CB	2.28	0.47
1:A:783:LEU:H	1:A:783:LEU:CD1	2.12	0.47
1:B:969:MET:O	1:B:969:MET:HE2	2.15	0.47
1:A:465:VAL:HG23	1:A:466:GLU:N	2.29	0.47
1:B:71:ILE:N	1:B:71:ILE:HD13	2.29	0.47
1:B:173:LEU:O	1:B:188:ILE:HA	2.15	0.47
1:A:367:PHE:C	1:A:367:PHE:CD2	2.88	0.47
1:A:899:MET:O	1:A:903:VAL:HG23	2.16	0.46
1:B:858:TYR:N	1:B:858:TYR:CD1	2.83	0.46
1:B:957:PHE:O	1:B:959:LEU:HG	2.15	0.46
1:B:413:LEU:HD23	1:B:413:LEU:C	2.35	0.46
1:B:737:ASP:C	1:B:739:ASN:N	2.69	0.46
1:A:86:THR:HA	1:A:89:VAL:HG23	1.97	0.46
1:B:60:LEU:O	1:B:64:ILE:HG12	2.15	0.46
1:B:10:GLU:CD	1:B:10:GLU:H	2.13	0.46
1:B:308:PRO:HB2	1:B:764:LEU:HD23	1.96	0.46
1:A:787:LEU:HD21	1:A:901:LEU:CA	2.45	0.46
1:A:899:MET:CE	1:A:970:VAL:CG2	2.88	0.46
1:B:895:GLU:HB2	1:B:896:PRO:HD3	1.96	0.46
1:B:878:GLU:HG3	1:B:879:ASP:N	2.23	0.46
1:B:553:GLY:O	1:B:554:THR:HG23	2.15	0.46
1:B:840:ILE:HD13	1:B:911:ASN:HD21	1.80	0.46
1:B:249:LEU:HD12	1:B:249:LEU:C	2.36	0.46
1:B:747:VAL:HG12	1:B:817:MET:HE1	1.97	0.46
1:B:322:GLY:HA2	1:B:325:ARG:HE	1.78	0.46
1:A:654:THR:OG1	1:A:657:GLU:HG3	2.16	0.46
1:A:10:GLU:CD	1:A:10:GLU:H	2.12	0.46
1:B:155:VAL:HG13	1:B:155:VAL:O	2.15	0.46
1:B:315:ILE:HD13	1:B:760:PHE:CE1	2.50	0.46
1:B:817:MET:C	1:B:819:ARG:N	2.66	0.46
1:A:855:TRP:HZ3	1:A:856:PHE:CE2	2.33	0.46
1:A:715:GLU:CB	1:A:716:ILE:HD12	2.44	0.46
1:B:549:ILE:CD1	1:B:596:VAL:HG21	2.25	0.46
1:B:522:ILE:N	1:B:522:ILE:HD12	2.31	0.46
1:A:759:GLN:HE22	1:A:762:ARG:NH1	2.13	0.46
1:B:757:MET:HA	1:B:760:PHE:CE2	2.51	0.46
1:A:783:LEU:HG	1:A:870:LEU:CD2	2.46	0.46
1:B:878:GLU:C	1:B:880:HIS:H	2.19	0.46
1:A:48:SER:H	1:A:51:GLU:CG	2.28	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:261:SER:O	1:A:264:ILE:CG1	2.62	0.46
1:A:762:ARG:HB3	1:A:837:TYR:CE1	2.49	0.46
1:B:311:LEU:CB	1:B:312:PRO:CD	2.88	0.46
1:B:751:ARG:O	1:B:755:ASN:OD1	2.34	0.46
1:B:367:PHE:C	1:B:367:PHE:CD2	2.88	0.46
1:B:859:ALA:O	1:B:861:ASP:N	2.48	0.46
1:A:716:ILE:N	1:A:716:ILE:HD12	2.30	0.46
1:A:868:HIS:O	1:A:872:HIS:ND1	2.46	0.46
1:B:39:ASN:O	1:B:40:GLU:HB3	2.15	0.46
1:B:326:MET:HE3	1:B:331:ALA:CB	2.45	0.46
1:B:737:ASP:OD2	1:B:739:ASN:HB2	2.16	0.46
1:B:633:ILE:CD1	1:B:642:PHE:HE1	2.27	0.46
1:A:633:ILE:HD12	1:A:642:PHE:CE1	2.50	0.46
1:B:108:GLN:HE21	1:B:108:GLN:HA	1.80	0.46
1:B:249:LEU:HB2	1:B:340:GLU:OE1	2.16	0.46
1:B:765:ILE:O	1:B:766:SER:C	2.53	0.46
1:A:917:SER:CB	1:A:920:GLN:HB2	2.46	0.46
1:A:854:TRP:CE3	1:A:855:TRP:N	2.84	0.46
1:A:716:ILE:HG22	1:A:718:ILE:HD11	1.98	0.46
1:B:86:THR:HA	1:B:89:VAL:CG2	2.44	0.46
1:A:443:THR:O	1:A:447:THR:HG23	2.16	0.46
1:B:0:ACE:H3	1:B:36:TYR:CE2	2.51	0.46
1:A:299:ALA:HA	1:A:302:LEU:HD11	1.98	0.46
1:A:75:LEU:CD2	1:A:297:LYS:HA	2.46	0.46
1:A:858:TYR:N	1:A:858:TYR:CD1	2.83	0.46
1:B:782:GLY:CA	1:B:871:THR:HB	2.46	0.46
1:A:53:VAL:HG22	1:A:106:VAL:HG21	1.97	0.46
1:B:827:PRO:HB2	1:B:828:LEU:H	1.50	0.45
1:A:947:ILE:HD11	1:A:957:PHE:CE1	2.52	0.45
1:B:788:ILE:HD11	1:B:958:LYS:HB2	1.98	0.45
1:B:491:ARG:HD2	1:B:588:GLU:OE2	2.15	0.45
1:A:790:VAL:O	1:A:790:VAL:CG1	2.61	0.45
1:A:832:TRP:HZ2	1:A:984:LEU:O	1.99	0.45
1:B:757:MET:HA	1:B:760:PHE:CZ	2.51	0.45
1:B:762:ARG:HB3	1:B:837:TYR:CE1	2.50	0.45
1:B:823:SER:HB3	1:B:826:GLU:CB	2.27	0.45
1:B:845:GLY:C	1:B:847:ALA:H	2.18	0.45
1:B:868:HIS:O	1:B:872:HIS:ND1	2.47	0.45
1:A:304:VAL:CG1	1:A:793:LEU:HD21	2.45	0.45
1:A:852:ALA:O	1:A:855:TRP:HB3	2.16	0.45
1:B:898:THR:OG1	1:B:960:LYS:N	2.45	0.45
1:A:880:HIS:HB3	1:A:881:PRO:HD3	1.97	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:652:ALA:HA	1:A:675:CYS:O	2.16	0.45
1:B:259:GLN:HE21	1:B:259:GLN:HB3	1.58	0.45
1:B:882:HIS:HB3	1:B:883:PHE:CD1	2.52	0.45
1:A:91:PRO:HA	1:A:94:ILE:CG2	2.46	0.45
1:B:97:ILE:O	1:B:100:ALA:HB3	2.17	0.45
1:B:62:VAL:HG12	1:B:98:LEU:HD22	1.99	0.45
1:A:79:GLU:CG	1:A:84:THR:HB	2.47	0.45
1:A:338:SER:O	1:A:339:VAL:C	2.55	0.45
1:B:299:ALA:O	1:B:302:LEU:HD12	2.17	0.45
1:B:349:CYS:HB3	1:B:624:ILE:HD11	1.95	0.45
1:A:261:SER:HA	1:A:264:ILE:HG12	1.97	0.45
1:A:757:MET:HA	1:A:760:PHE:CZ	2.52	0.45
1:B:767:SER:O	1:B:771:GLU:CG	2.63	0.45
1:B:947:ILE:HD11	1:B:957:PHE:CE1	2.51	0.45
1:B:108:GLN:NE2	1:B:108:GLN:HA	2.32	0.45
1:A:335:SER:OG	1:A:337:PRO:HD2	2.17	0.45
1:A:765:ILE:O	1:A:766:SER:C	2.55	0.45
1:B:311:LEU:HD11	1:B:761:ILE:HG23	1.98	0.45
1:B:342:LEU:O	1:B:747:VAL:HG22	2.17	0.45
1:B:833:LEU:CA	1:B:836:ARG:HB2	2.41	0.45
1:A:898:THR:CA	1:A:959:LEU:CD2	2.88	0.45
1:A:854:TRP:HH2	1:A:966:GLN:HG2	1.81	0.45
1:B:895:GLU:CD	1:B:895:GLU:H	2.20	0.45
1:B:971:LEU:O	1:B:975:LEU:HD23	2.16	0.45
1:A:639:ILE:HD11	1:A:641:ILE:CD1	2.47	0.45
1:A:757:MET:HA	1:A:760:PHE:CE2	2.51	0.45
1:B:810:ASN:HA	1:B:930:ASN:ND2	2.32	0.45
1:A:849:VAL:O	1:A:853:ALA:HB2	2.12	0.45
1:A:817:MET:C	1:A:819:ARG:N	2.66	0.45
1:A:39:ASN:O	1:A:40:GLU:CB	2.62	0.45
1:A:983:ILE:O	1:A:987:ILE:HG12	2.16	0.45
1:A:86:THR:HG22	1:A:89:VAL:HG21	1.99	0.45
1:B:352:LYS:HB2	1:B:625:THR:HB	1.99	0.45
1:A:190:HIS:O	1:A:206:ASN:HA	2.17	0.45
1:B:885:GLY:O	1:B:886:LEU:HG	2.17	0.45
1:A:898:THR:OG1	1:A:960:LYS:N	2.47	0.45
1:A:799:THR:HG21	1:A:905:VAL:HG22	1.99	0.45
1:A:975:LEU:H	1:A:975:LEU:CD2	2.30	0.45
1:A:914:ASN:HD21	1:A:922:LEU:HD11	1.80	0.45
1:A:633:ILE:CD1	1:A:642:PHE:HE1	2.29	0.45
1:B:963:ASP:H	1:B:966:GLN:NE2	2.11	0.44
1:B:322:GLY:N	1:B:325:ARG:HH21	2.15	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:633:ILE:HD12	1:A:642:PHE:HE1	1.82	0.44
1:B:193:PRO:O	1:B:195:PRO:HD3	2.18	0.44
1:A:663:LEU:H	1:A:663:LEU:CD1	2.30	0.44
1:A:308:PRO:CB	1:A:764:LEU:HD23	2.46	0.44
1:A:782:GLY:HA3	1:A:871:THR:CB	2.46	0.44
1:B:794:TRP:O	1:B:794:TRP:CD1	2.70	0.44
1:B:671:ARG:HD3	1:B:694:TYR:CZ	2.52	0.44
1:A:765:ILE:HD13	1:A:765:ILE:C	2.38	0.44
1:A:857:MET:CE	1:A:867:TYR:HA	2.48	0.44
1:B:774:CYS:O	1:B:778:THR:HG23	2.17	0.44
1:A:48:SER:O	1:A:52:LEU:HG	2.18	0.44
1:B:680:GLU:O	1:B:683:HIS:HB2	2.16	0.44
1:B:403:ARG:HH11	1:B:456:ASN:HD21	1.65	0.44
1:A:553:GLY:O	1:A:554:THR:HG23	2.17	0.44
1:B:786:ALA:C	1:B:897:MET:HG2	2.37	0.44
1:B:897:MET:CE	1:B:958:LYS:HE3	2.47	0.44
1:A:491:ARG:HD2	1:A:588:GLU:OE2	2.17	0.44
1:A:271:VAL:O	1:A:273:LEU:N	2.50	0.44
1:A:300:VAL:HA	1:A:303:ALA:HB2	2.00	0.44
1:A:767:SER:O	1:A:771:GLU:CG	2.63	0.44
1:A:248:PRO:HB2	1:A:341:THR:HG23	1.99	0.44
1:B:794:TRP:C	1:B:794:TRP:CD1	2.90	0.44
1:A:557:ASP:HB3	1:A:559:LEU:HG	1.98	0.44
1:A:338:SER:HA	1:A:341:THR:CB	2.46	0.44
1:A:155:VAL:HG13	1:A:155:VAL:O	2.17	0.44
1:A:853:ALA:O	1:A:856:PHE:N	2.51	0.44
1:B:899:MET:HE2	1:B:966:GLN:O	2.15	0.44
1:B:914:ASN:HD21	1:B:922:LEU:HD11	1.82	0.44
1:B:443:THR:O	1:B:447:THR:HG23	2.18	0.44
1:A:762:ARG:CZ	1:A:833:LEU:HD11	2.48	0.44
1:A:810:ASN:HA	1:A:930:ASN:ND2	2.33	0.44
1:B:833:LEU:HD11	1:B:837:TYR:HE1	1.83	0.44
1:A:786:ALA:C	1:A:897:MET:HG2	2.38	0.44
1:A:878:GLU:C	1:A:880:HIS:H	2.20	0.44
1:B:663:LEU:H	1:B:663:LEU:HD13	1.82	0.44
1:A:624:ILE:CD1	1:A:676:PHE:CD1	3.01	0.44
1:B:85:ILE:HD13	1:B:85:ILE:O	2.17	0.44
1:B:336:LEU:O	1:B:339:VAL:HG23	2.18	0.44
1:B:759:GLN:O	1:B:760:PHE:C	2.55	0.44
1:A:748:GLU:N	1:A:817:MET:HE1	2.33	0.43
1:A:307:ILE:N	1:A:307:ILE:HD13	2.33	0.43
1:B:755:ASN:HD21	1:B:816:ILE:HD11	1.82	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:895:GLU:N	1:A:896:PRO:HD2	2.33	0.43
1:B:81:GLY:HA2	1:B:84:THR:HG22	2.00	0.43
1:A:261:SER:O	1:A:264:ILE:CB	2.66	0.43
1:B:633:ILE:HD13	1:B:675:CYS:SG	2.59	0.43
1:B:982:GLU:OE1	1:B:982:GLU:HA	2.18	0.43
1:B:873:PHE:HB2	1:B:891:PHE:CD2	2.53	0.43
1:B:804:ALA:O	1:B:805:THR:C	2.56	0.43
1:B:975:LEU:H	1:B:975:LEU:CD2	2.31	0.43
1:B:524:ARG:HH21	1:B:588:GLU:HB2	1.83	0.43
1:A:926:PRO:HA	1:A:927:PRO:HD3	1.95	0.43
1:A:293:ILE:H	1:A:293:ILE:HG13	1.59	0.43
1:A:762:ARG:HH21	1:A:833:LEU:HD13	1.83	0.43
1:B:802:LEU:H	1:B:802:LEU:HG	1.49	0.43
1:A:895:GLU:HB2	1:A:896:PRO:HD3	2.00	0.43
1:A:342:LEU:O	1:A:747:VAL:HG22	2.18	0.43
1:B:907:ILE:CD1	1:B:974:SER:HB3	2.49	0.43
1:A:102:ALA:O	1:A:106:VAL:CG2	2.64	0.43
1:A:794:TRP:C	1:A:794:TRP:CD1	2.91	0.43
1:B:759:GLN:HE22	1:B:762:ARG:NH1	2.15	0.43
1:B:617:ALA:O	1:B:820:PRO:HA	2.19	0.43
1:B:624:ILE:CG2	1:B:684:LYS:HG2	2.49	0.43
1:B:70:CYS:O	1:B:73:PHE:HB3	2.18	0.43
1:A:873:PHE:HB2	1:A:891:PHE:CD2	2.53	0.43
1:A:299:ALA:O	1:A:303:ALA:HB2	2.18	0.43
1:B:825:LYS:HA	1:B:825:LYS:HE2	2.00	0.43
1:B:897:MET:HE3	1:B:958:LYS:O	2.18	0.43
1:B:49:LEU:HD12	1:B:110:ARG:HG3	1.98	0.43
1:B:522:ILE:HD11	1:B:563:ALA:CB	2.48	0.43
1:B:38:HIS:HD2	1:B:143:ARG:NH1	2.17	0.43
1:A:625:THR:HA	3:A:998:BEF:F1	2.08	0.43
1:A:267:ILE:HD12	1:A:302:LEU:HD22	2.00	0.43
1:A:897:MET:HE3	1:A:958:LYS:O	2.19	0.43
1:A:898:THR:OG1	1:A:959:LEU:HA	2.18	0.43
1:B:367:PHE:HE2	1:B:545:ILE:HD12	1.83	0.43
1:B:48:SER:O	1:B:52:LEU:HG	2.19	0.43
1:A:680:GLU:HB3	1:A:681:PRO:CD	2.48	0.43
1:B:255:GLU:C	1:B:257:GLY:H	2.22	0.43
1:A:670:CYS:HG	1:A:690:TYR:HD1	1.66	0.43
1:A:783:LEU:HG	1:A:870:LEU:HD23	1.99	0.43
1:A:845:GLY:C	1:A:847:ALA:H	2.20	0.43
1:B:895:GLU:N	1:B:896:PRO:HD2	2.33	0.43
1:A:975:LEU:N	1:A:976:PRO:HD2	2.34	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:55:GLU:O	1:A:58:GLU:HB2	2.19	0.43
1:B:519:GLU:CD	1:B:519:GLU:N	2.63	0.43
1:A:811:PRO:HG3	1:A:929:VAL:O	2.18	0.43
1:A:606:GLU:N	1:A:606:GLU:CD	2.71	0.43
1:A:624:ILE:HD13	1:A:676:PHE:HD1	1.84	0.43
1:A:852:ALA:CA	1:A:899:MET:HB3	2.48	0.43
1:A:103:ILE:HG23	1:A:104:VAL:H	1.83	0.43
1:B:214:ILE:H	1:B:214:ILE:CD1	2.31	0.43
1:A:524:ARG:HH21	1:A:588:GLU:HB2	1.84	0.43
1:A:604:ARG:NE	1:A:720:MET:HE2	2.34	0.43
1:A:347:VAL:HG12	1:A:348:ILE:N	2.34	0.43
1:A:825:LYS:HE2	1:A:825:LYS:HA	2.00	0.42
1:B:342:LEU:HA	1:B:716:ILE:HG13	2.00	0.42
1:B:172:THR:CG2	1:B:489:ARG:CD	2.96	0.42
1:B:119:LEU:HD12	1:B:119:LEU:HA	1.75	0.42
1:A:774:CYS:O	1:A:778:THR:HG23	2.19	0.42
1:B:811:PRO:HG3	1:B:929:VAL:O	2.18	0.42
1:B:763:TYR:HE1	1:B:908:GLU:O	2.02	0.42
1:A:982:GLU:HA	1:A:982:GLU:OE1	2.19	0.42
1:B:851:ALA:HB3	1:B:903:VAL:HG21	2.00	0.42
1:B:894:PRO:HB3	1:B:958:LYS:C	2.39	0.42
1:A:314:VAL:CG1	1:A:804:ALA:HB1	2.49	0.42
1:B:100:ALA:O	1:B:103:ILE:HG22	2.19	0.42
1:B:78:PHE:HE2	1:B:293:ILE:HG23	1.83	0.42
1:A:823:SER:HB3	1:A:826:GLU:CB	2.26	0.42
1:A:848:THR:O	1:A:852:ALA:CB	2.68	0.42
1:B:62:VAL:HG13	1:B:98:LEU:HD22	2.00	0.42
1:B:28:GLN:HG2	1:B:31:ARG:HH22	1.85	0.42
1:A:236:ARG:C	1:A:236:ARG:HD3	2.39	0.42
1:B:315:ILE:HD13	1:B:760:PHE:HE1	1.85	0.42
1:B:315:ILE:HD11	1:B:761:ILE:HD11	2.00	0.42
1:B:832:TRP:HZ2	1:B:984:LEU:O	2.03	0.42
1:A:342:LEU:HD22	1:A:747:VAL:HG22	2.00	0.42
1:B:60:LEU:HD23	1:B:257:GLY:CA	2.50	0.42
1:B:391:PRO:HB3	1:B:450:GLU:HB3	2.01	0.42
1:A:884:GLU:HG2	1:A:884:GLU:H	1.63	0.42
1:A:880:HIS:N	1:A:881:PRO:HD2	2.34	0.42
1:A:882:HIS:HB3	1:A:883:PHE:CD1	2.54	0.42
1:A:975:LEU:HD22	1:A:975:LEU:N	2.35	0.42
1:A:112:ALA:O	1:A:113:GLU:C	2.58	0.42
1:A:624:ILE:HD13	1:A:676:PHE:HB2	2.01	0.42
1:A:28:GLN:HG2	1:A:31:ARG:HH22	1.85	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:260:LEU:O	1:B:263:VAL:HB	2.19	0.42
1:B:753:ILE:O	1:B:757:MET:HB2	2.19	0.42
1:A:899:MET:HE2	1:A:966:GLN:O	2.19	0.42
1:A:103:ILE:HG23	1:A:104:VAL:N	2.35	0.42
1:A:914:ASN:O	1:A:916:LEU:N	2.53	0.42
1:B:402:ILE:HD12	1:B:402:ILE:C	2.40	0.42
1:B:258:GLU:O	1:B:259:GLN:C	2.57	0.42
1:A:827:PRO:HB2	1:A:828:LEU:H	1.49	0.42
1:B:311:LEU:HD21	1:B:761:ILE:HD12	2.01	0.42
1:B:314:VAL:CG1	1:B:804:ALA:HB1	2.45	0.42
1:B:795:VAL:HA	1:B:799:THR:HB	2.02	0.42
1:B:899:MET:CE	1:B:970:VAL:CG2	2.89	0.42
1:B:962:LEU:HB3	1:B:966:GLN:HB2	2.01	0.42
1:A:120:LYS:CE	1:A:723:GLY:O	2.68	0.42
1:B:950:VAL:O	1:B:954:PRO:CD	2.68	0.42
1:A:387:SER:O	1:A:602:PRO:HG2	2.20	0.42
1:A:286:GLY:HA3	1:A:290:ARG:HB3	1.99	0.42
1:A:577:VAL:HG12	1:A:579:ASP:OD2	2.20	0.42
1:B:231:GLU:O	1:B:235:ILE:HG12	2.18	0.42
1:B:680:GLU:HB3	1:B:681:PRO:CD	2.49	0.42
1:A:834:PHE:O	1:A:838:MET:HB2	2.20	0.42
1:B:338:SER:O	1:B:339:VAL:C	2.57	0.42
1:A:899:MET:SD	1:A:962:LEU:CD2	3.08	0.42
1:B:851:ALA:O	1:B:854:TRP:HB3	2.19	0.42
1:B:577:VAL:HG12	1:B:579:ASP:OD2	2.19	0.42
1:B:235:ILE:HD11	1:B:681:PRO:HG2	2.02	0.42
1:B:82:GLU:C	1:B:83:GLU:HG3	2.40	0.42
1:A:181:THR:CB	1:A:183:GLU:HG3	2.48	0.42
1:A:232:ILE:HD13	1:A:232:ILE:C	2.40	0.42
1:A:174:ARG:NH2	1:A:188:ILE:HD11	2.35	0.42
1:B:59:ASP:OD2	1:B:62:VAL:HG23	2.19	0.42
1:B:2:GLU:O	1:B:3:ALA:CB	2.66	0.42
1:B:600:LEU:N	1:B:600:LEU:HD12	2.35	0.42
1:A:624:ILE:HG22	1:A:684:LYS:HG2	2.00	0.42
1:A:265:SER:O	1:A:268:CYS:HB2	2.20	0.42
1:A:193:PRO:O	1:A:195:PRO:HD3	2.20	0.42
1:B:751:ARG:NH1	1:B:819:ARG:O	2.52	0.41
1:A:962:LEU:HB3	1:A:966:GLN:HB2	2.02	0.41
1:B:606:GLU:CD	1:B:606:GLU:N	2.72	0.41
1:B:606:GLU:HG3	1:B:739:ASN:OD1	2.20	0.41
1:B:290:ARG:HE	1:B:291:GLY:N	2.18	0.41
1:A:624:ILE:HD13	1:A:676:PHE:CD1	2.55	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:770:GLY:HA2	1:B:841:GLY:O	2.20	0.41
1:B:336:LEU:O	1:B:339:VAL:CG2	2.68	0.41
1:B:549:ILE:CD1	1:B:596:VAL:HG11	2.50	0.41
1:B:975:LEU:N	1:B:976:PRO:HD2	2.35	0.41
1:A:2:GLU:O	1:A:3:ALA:CB	2.65	0.41
1:A:311:LEU:HD11	1:A:761:ILE:HG12	2.02	0.41
1:B:336:LEU:HB2	1:B:337:PRO:HD3	2.02	0.41
1:A:901:LEU:O	1:A:905:VAL:HG23	2.20	0.41
1:A:868:HIS:NE2	1:A:883:PHE:HZ	2.18	0.41
1:A:112:ALA:O	1:A:114:ASN:N	2.53	0.41
1:A:931:ILE:CD1	1:A:931:ILE:H	2.31	0.41
1:A:290:ARG:HG3	1:A:291:GLY:N	2.34	0.41
1:B:60:LEU:HD21	1:B:257:GLY:HA3	2.03	0.41
1:B:584:PHE:CD2	1:B:584:PHE:N	2.86	0.41
1:B:1:MET:O	1:B:1:MET:HG2	2.20	0.41
1:A:311:LEU:HD22	1:A:761:ILE:HG23	2.02	0.41
1:B:749:GLU:C	1:B:751:ARG:N	2.74	0.41
1:B:762:ARG:CZ	1:B:833:LEU:HD11	2.51	0.41
1:B:784:PRO:HG2	1:B:856:PHE:CE1	2.54	0.41
1:B:854:TRP:HH2	1:B:966:GLN:HG2	1.84	0.41
1:B:895:GLU:OE2	1:B:960:LYS:HD2	2.21	0.41
1:B:255:GLU:C	1:B:257:GLY:N	2.74	0.41
1:A:231:GLU:O	1:A:235:ILE:HG12	2.20	0.41
1:A:434:TYR:HE2	1:A:468:ALA:HB2	1.85	0.41
1:A:762:ARG:NH2	1:A:833:LEU:HD13	2.36	0.41
1:A:762:ARG:O	1:A:765:ILE:HG22	2.20	0.41
1:B:762:ARG:HH21	1:B:833:LEU:HD13	1.85	0.41
1:A:897:MET:CE	1:A:958:LYS:HE3	2.50	0.41
1:A:90:GLU:OE1	1:A:793:LEU:HD11	2.21	0.41
1:B:42:PRO:O	1:B:43:ALA:O	2.37	0.41
1:B:556:ARG:NE	1:B:644:GLU:CG	2.83	0.41
1:A:763:TYR:HE1	1:A:908:GLU:O	2.03	0.41
1:B:804:ALA:O	1:B:807:LEU:N	2.52	0.41
1:A:969:MET:O	1:A:969:MET:HE2	2.20	0.41
1:B:868:HIS:NE2	1:B:883:PHE:HZ	2.18	0.41
1:B:975:LEU:N	1:B:975:LEU:HD22	2.36	0.41
1:A:70:CYS:O	1:A:73:PHE:HB3	2.21	0.41
1:A:697:ILE:H	1:A:697:ILE:HD12	1.85	0.41
1:B:686:LYS:HE3	1:B:686:LYS:HB2	1.90	0.41
1:A:584:PHE:N	1:A:584:PHE:CD2	2.87	0.41
1:A:851:ALA:HB3	1:A:903:VAL:HG21	2.01	0.41
1:A:895:GLU:OE2	1:A:960:LYS:CD	2.69	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:103:ILE:HD13	1:A:103:ILE:C	2.41	0.41
1:B:352:LYS:HG2	1:B:353:THR:N	2.35	0.41
1:A:305:ALA:O	1:A:768:ASN:HA	2.20	0.41
1:B:309:GLU:H	1:B:309:GLU:CD	2.23	0.41
1:A:775:ILE:HD13	1:A:775:ILE:N	2.31	0.41
1:B:964:LEU:O	1:B:964:LEU:HD23	2.20	0.41
1:B:926:PRO:HA	1:B:927:PRO:HD3	1.95	0.41
1:A:325:ARG:NH2	1:A:753:ILE:HD11	2.36	0.41
1:B:116:ILE:CD1	1:B:334:ARG:HG2	2.51	0.41
1:A:663:LEU:CD1	1:A:663:LEU:N	2.83	0.41
1:A:794:TRP:O	1:A:794:TRP:CD1	2.73	0.41
1:B:390:ALA:HA	1:B:391:PRO:HD3	1.82	0.41
1:A:96:LEU:C	1:A:96:LEU:HD23	2.41	0.41
1:B:338:SER:HA	1:B:341:THR:CB	2.50	0.41
1:B:740:PHE:HA	1:B:743:ILE:HD13	2.01	0.41
1:B:947:ILE:HD11	1:B:957:PHE:CD2	2.56	0.41
1:B:853:ALA:O	1:B:856:PHE:N	2.54	0.41
1:B:969:MET:HE2	1:B:973:ILE:HG12	2.03	0.41
1:B:880:HIS:N	1:B:881:PRO:CD	2.83	0.41
1:B:624:ILE:HG22	1:B:684:LYS:HG2	2.03	0.41
1:A:907:ILE:CD1	1:A:974:SER:HB3	2.51	0.41
1:A:512:MET:HB2	1:A:567:ARG:HB3	2.03	0.41
1:B:652:ALA:HA	1:B:675:CYS:O	2.20	0.41
1:A:885:GLY:O	1:A:886:LEU:HG	2.21	0.41
1:A:797:LEU:HA	1:A:797:LEU:HD23	1.64	0.41
1:B:895:GLU:OE2	1:B:960:LYS:CD	2.68	0.41
1:B:261:SER:O	1:B:264:ILE:HB	2.20	0.41
1:A:416:ILE:HD11	1:A:566:THR:HG22	2.03	0.41
1:B:369:ILE:HD11	1:B:593:PHE:HD1	1.85	0.41
1:B:290:ARG:HE	1:B:291:GLY:CA	2.34	0.41
1:A:390:ALA:HA	1:A:391:PRO:HD3	1.83	0.41
1:B:305:ALA:HB3	1:B:772:VAL:HG22	2.03	0.41
1:B:315:ILE:HG22	1:B:316:THR:N	2.36	0.40
1:A:895:GLU:OE2	1:A:960:LYS:HD2	2.21	0.40
1:B:369:ILE:HG12	1:B:528:VAL:HG11	2.03	0.40
1:A:606:GLU:HG3	1:A:739:ASN:OD1	2.22	0.40
1:B:112:ALA:O	1:B:113:GLU:C	2.58	0.40
1:B:345:THR:HG23	1:B:697:ILE:O	2.21	0.40
1:B:304:VAL:C	1:B:306:ALA:H	2.25	0.40
1:B:740:PHE:O	1:B:743:ILE:CD1	2.60	0.40
1:B:556:ARG:NE	1:B:644:GLU:HG2	2.37	0.40
1:A:363:VAL:HG11	1:A:448:LEU:HD22	2.03	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:269:VAL:O	1:B:272:TRP:HE3	2.05	0.40
1:A:843:TYR:HH	1:A:976:PRO:HG2	1.84	0.40
1:B:914:ASN:O	1:B:916:LEU:N	2.55	0.40
1:B:357:THR:HB	1:B:602:PRO:O	2.21	0.40
1:A:75:LEU:HD11	1:A:300:VAL:HG23	2.04	0.40
1:B:762:ARG:NH2	1:B:833:LEU:HD13	2.36	0.40
1:B:740:PHE:C	1:B:743:ILE:HD13	2.40	0.40
1:A:795:VAL:HA	1:A:799:THR:HB	2.03	0.40
1:A:947:ILE:HD11	1:A:957:PHE:CD2	2.57	0.40
1:B:898:THR:HG21	1:B:960:LYS:O	2.22	0.40
1:B:236:ARG:HD3	1:B:236:ARG:C	2.41	0.40
1:A:604:ARG:HB2	1:A:607:VAL:HG23	2.03	0.40
1:A:684:LYS:NZ	1:A:707:ASP:OD1	2.53	0.40
1:B:797:LEU:HA	1:B:797:LEU:HD23	1.62	0.40
1:B:416:ILE:HD11	1:B:566:THR:HG22	2.03	0.40
1:B:103:ILE:CG2	1:B:104:VAL:N	2.84	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	993/995 (100%)	883 (89%)	84 (8%)	26 (3%)	8	62
1	B	993/995 (100%)	876 (88%)	91 (9%)	26 (3%)	8	62
All	All	1986/1990 (100%)	1759 (89%)	175 (9%)	52 (3%)	8	62

All (52) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	47	LYS
1	A	245	ASP
1	A	292	ALA
1	A	783	LEU

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Mol	Chain	Res	Type
1	A	818	ASP
1	A	827	PRO
1	A	828	LEU
1	B	47	LYS
1	B	245	ASP
1	B	292	ALA
1	B	783	LEU
1	B	818	ASP
1	B	827	PRO
1	B	828	LEU
1	B	863	PRO
1	A	43	ALA
1	A	293	ILE
1	A	309	GLU
1	A	790	VAL
1	A	860	GLU
1	A	863	PRO
1	A	878	GLU
1	B	43	ALA
1	B	293	ILE
1	B	790	VAL
1	B	860	GLU
1	B	878	GLU
1	A	762	ARG
1	A	830	SER
1	B	309	GLU
1	B	330	ASN
1	B	762	ARG
1	B	830	SER
1	A	272	TRP
1	A	287	SER
1	A	754	TYR
1	A	915	SER
1	B	42	PRO
1	B	831	GLY
1	A	294	TYR
1	A	951	ASP
1	B	294	TYR
1	B	915	SER
1	A	831	GLY
1	B	352	LYS
1	B	951	ASP

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Mol	Chain	Res	Type
1	A	789	PRO
1	A	601	ASP
1	B	46	GLY
1	B	789	PRO
1	A	950	VAL
1	B	950	VAL

### 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	840/840 (100%)	757 (90%)	83 (10%)	11	53
1	B	840/840 (100%)	761 (91%)	79 (9%)	13	56
All	All	1680/1680 (100%)	1518 (90%)	162 (10%)	12	55

All (162) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	GLU
1	A	56	GLN
1	A	59	ASP
1	A	79	GLU
1	A	86	THR
1	A	88	PHE
1	A	90	GLU
1	A	92	PHE
1	A	94	ILE
1	A	103	ILE
1	A	110	ARG
1	A	114	ASN
1	A	125	GLU
1	A	134	ARG
1	A	149	ASP
1	A	164	ARG
1	A	171	THR
1	A	232	ILE
1	A	239	MET

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Mol	Chain	Res	Type
1	A	247	THR
1	A	249	LEU
1	A	250	GLN
1	A	251	GLN
1	A	252	LYS
1	A	255	GLU
1	A	256	PHE
1	A	259	GLN
1	A	284	HIS
1	A	289	ILE
1	A	296	PHE
1	A	298	ILE
1	A	302	LEU
1	A	307	ILE
1	A	334	ARG
1	A	338	SER
1	A	342	LEU
1	A	353	THR
1	A	355	THR
1	A	356	LEU
1	A	359	ASN
1	A	367	PHE
1	A	402	ILE
1	A	403	ARG
1	A	421	ASN
1	A	426	ASP
1	A	439	GLU
1	A	447	THR
1	A	484	THR
1	A	488	SER
1	A	489	ARG
1	A	519	GLU
1	A	534	ARG
1	A	545	ILE
1	A	566	THR
1	A	572	LYS
1	A	573	ARG
1	A	575	GLU
1	A	600	LEU
1	A	691	LEU
1	A	713	LYS
1	A	726	VAL

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Mol	Chain	Res	Type
1	A	757	MET
1	A	763	TYR
1	A	765	ILE
1	A	775	ILE
1	A	793	LEU
1	A	795	VAL
1	A	797	LEU
1	A	802	LEU
1	A	813	ASP
1	A	816	ILE
1	A	817	MET
1	A	829	ILE
1	A	840	ILE
1	A	849	VAL
1	A	854	TRP
1	A	882	HIS
1	A	911	ASN
1	A	925	MET
1	A	951	ASP
1	A	955	MET
1	A	956	ILE
1	A	993	GLU
1	B	10	GLU
1	B	78	PHE
1	B	79	GLU
1	B	85	ILE
1	B	86	THR
1	B	88	PHE
1	B	90	GLU
1	B	92	PHE
1	B	94	ILE
1	B	95	LEU
1	B	101	ASN
1	B	103	ILE
1	B	109	GLU
1	B	110	ARG
1	B	114	ASN
1	B	125	GLU
1	B	134	ARG
1	B	140	ILE
1	B	149	ASP
1	B	164	ARG

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Mol	Chain	Res	Type
1	B	172	THR
1	B	175	VAL
1	B	246	LYS
1	B	250	GLN
1	B	251	GLN
1	B	252	LYS
1	B	255	GLU
1	B	256	PHE
1	B	259	GLN
1	B	265	SER
1	B	272	TRP
1	B	281	ASP
1	B	296	PHE
1	B	300	VAL
1	B	302	LEU
1	B	307	ILE
1	B	338	SER
1	B	342	LEU
1	B	355	THR
1	B	361	MET
1	B	384	ILE
1	B	403	ARG
1	B	421	ASN
1	B	426	ASP
1	B	439	GLU
1	B	447	THR
1	B	484	THR
1	B	488	SER
1	B	489	ARG
1	B	519	GLU
1	B	534	ARG
1	B	566	THR
1	B	572	LYS
1	B	573	ARG
1	B	575	GLU
1	B	600	LEU
1	B	691	LEU
1	B	713	LYS
1	B	726	VAL
1	B	755	ASN
1	B	757	MET
1	B	763	TYR

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Mol	Chain	Res	Type
1	B	783	LEU
1	B	793	LEU
1	B	795	VAL
1	B	797	LEU
1	B	802	LEU
1	B	813	ASP
1	B	816	ILE
1	B	817	MET
1	B	829	ILE
1	B	849	VAL
1	B	854	TRP
1	B	882	HIS
1	B	911	ASN
1	B	925	MET
1	B	951	ASP
1	B	955	MET
1	B	993	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (36) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	108	GLN
1	A	111	ASN
1	A	114	ASN
1	A	177	GLN
1	A	251	GLN
1	A	275	ASN
1	A	359	ASN
1	A	406	GLN
1	A	421	ASN
1	A	461	ASN
1	A	510	ASN
1	A	759	GLN
1	A	869	GLN
1	A	875	GLN
1	A	911	ASN
1	A	914	ASN
1	A	966	GLN
1	B	38	HIS
1	B	108	GLN
1	B	111	ASN
1	B	114	ASN

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Mol	Chain	Res	Type
1	B	250	GLN
1	B	251	GLN
1	B	259	GLN
1	B	359	ASN
1	B	406	GLN
1	B	421	ASN
1	B	456	ASN
1	B	461	ASN
1	B	510	ASN
1	B	759	GLN
1	B	869	GLN
1	B	875	GLN
1	B	911	ASN
1	B	914	ASN
1	B	966	GLN

### 5.3.3 RNA ⓘ

There are no RNA chains in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. 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  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	BEF	A	998	1	0,3,3	0.00	-	0,3,3	0.00	-
3	BEF	B	1098	1	0,3,3	0.00	-	0,3,3	0.00	-

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	BEF	A	998	1	-	0/0/0/0	0/0/0/0
3	BEF	B	1098	1	-	0/0/0/0	0/0/0/0

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.

## 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

### 6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

### 6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

### 6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.