



Full wwPDB X-ray Structure Validation Report i

Feb 28, 2014 – 11:27 AM GMT

PDB ID : 3DBR
Title : Structural Dissection of a Gating Mechanism Preventing Misactivation of Ubiquitin by NEDD8's E1 (APPBP1-UBA3Arg190Gln-NEDD8Ala72Arg)
Authors : Souphron, J.; Schulman, B.A.
Deposited on : 2008-06-02
Resolution : 3.05 Å(reported)

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

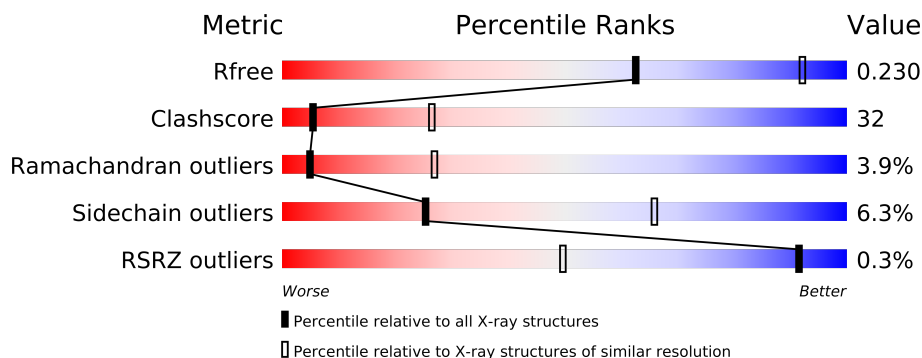
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.15 2013
Xtriage (Phenix)	:	dev-1323
EDS	:	stable22639
Percentile statistics	:	21963
Refmac	:	5.8.0049
CCP4	:	6.3.0 (Settle)
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 3.05 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	66092	2079 (3.12-3.00)
Clashscore	79885	2629 (3.12-3.00)
Ramachandran outliers	78287	2536 (3.12-3.00)
Sidechain outliers	78261	2539 (3.12-3.00)
RSRZ outliers	66119	2081 (3.12-3.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	531	
1	C	531	
1	E	531	
1	G	531	
2	B	434	
2	D	434	
2	F	434	
2	H	434	
3	I	88	
3	J	88	
3	K	88	
3	L	88	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 32424 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 NEDD8-activating enzyme E1 regulatory subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	517	Total	C	N	O	S	0	0	0
			4104	2600	699	790	15			
1	C	516	Total	C	N	O	S	0	0	0
			4096	2596	698	787	15			
1	E	515	Total	C	N	O	S	0	0	0
			4093	2597	696	785	15			
1	G	508	Total	C	N	O	S	0	0	0
			4038	2563	686	774	15			

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	EXPRESSION TAG	UNP Q13564
A	0	SER	-	EXPRESSION TAG	UNP Q13564
A	?	-	ASN	DELETION	UNP Q13564
A	?	-	GLU	DELETION	UNP Q13564
A	?	-	ASN	DELETION	UNP Q13564
A	?	-	GLY	DELETION	UNP Q13564
A	?	-	ALA	DELETION	UNP Q13564
C	-1	GLY	-	EXPRESSION TAG	UNP Q13564
C	0	SER	-	EXPRESSION TAG	UNP Q13564
C	?	-	ASN	DELETION	UNP Q13564
C	?	-	GLU	DELETION	UNP Q13564
C	?	-	ASN	DELETION	UNP Q13564
C	?	-	GLY	DELETION	UNP Q13564
C	?	-	ALA	DELETION	UNP Q13564
E	-1	GLY	-	EXPRESSION TAG	UNP Q13564
E	0	SER	-	EXPRESSION TAG	UNP Q13564
E	?	-	ASN	DELETION	UNP Q13564
E	?	-	GLU	DELETION	UNP Q13564
E	?	-	ASN	DELETION	UNP Q13564
E	?	-	GLY	DELETION	UNP Q13564
E	?	-	ALA	DELETION	UNP Q13564

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
G	-1	GLY	-	EXPRESSION TAG	UNP Q13564
G	0	SER	-	EXPRESSION TAG	UNP Q13564
G	?	-	ASN	DELETION	UNP Q13564
G	?	-	GLU	DELETION	UNP Q13564
G	?	-	ASN	DELETION	UNP Q13564
G	?	-	GLY	DELETION	UNP Q13564
G	?	-	ALA	DELETION	UNP Q13564

- Molecule 2 is a protein called NEDD8-activating enzyme E1 catalytic subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	432	Total	C	N	O	S	0	0	0
			3402	2174	575	636	17			
2	D	431	Total	C	N	O	S	0	0	0
			3399	2174	575	633	17			
2	F	431	Total	C	N	O	S	0	0	0
			3395	2171	574	633	17			
2	H	431	Total	C	N	O	S	0	0	0
			3387	2164	573	633	17			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	9	MET	-	EXPRESSION TAG	UNP Q8TBC4
B	10	LYS	-	EXPRESSION TAG	UNP Q8TBC4
B	11	LEU	-	EXPRESSION TAG	UNP Q8TBC4
B	190	GLN	ARG	ENGINEERED	UNP Q8TBC4
B	216	ALA	CYS	ENGINEERED	UNP Q8TBC4
D	9	MET	-	EXPRESSION TAG	UNP Q8TBC4
D	10	LYS	-	EXPRESSION TAG	UNP Q8TBC4
D	11	LEU	-	EXPRESSION TAG	UNP Q8TBC4
D	190	GLN	ARG	ENGINEERED	UNP Q8TBC4
D	216	ALA	CYS	ENGINEERED	UNP Q8TBC4
F	9	MET	-	EXPRESSION TAG	UNP Q8TBC4
F	10	LYS	-	EXPRESSION TAG	UNP Q8TBC4
F	11	LEU	-	EXPRESSION TAG	UNP Q8TBC4
F	190	GLN	ARG	ENGINEERED	UNP Q8TBC4
F	216	ALA	CYS	ENGINEERED	UNP Q8TBC4
H	9	MET	-	EXPRESSION TAG	UNP Q8TBC4
H	10	LYS	-	EXPRESSION TAG	UNP Q8TBC4
H	11	LEU	-	EXPRESSION TAG	UNP Q8TBC4
H	190	GLN	ARG	ENGINEERED	UNP Q8TBC4

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
H	216	ALA	CYS	ENGINEERED	UNP Q8TBC4

- Molecule 3 is a protein called NEDD8.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	I	85	Total	C	N	O	S	0	0	0
			666	414	122	128	2			
3	J	77	Total	C	N	O	S	0	0	0
			612	384	108	118	2			
3	K	77	Total	C	N	O	S	0	0	0
			612	384	108	118	2			
3	L	78	Total	C	N	O	S	0	0	0
			616	386	109	119	2			

There are 52 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
I	89	GLY	-	EXPRESSION TAG	UNP Q15843
I	90	SER	-	EXPRESSION TAG	UNP Q15843
I	91	ARG	-	EXPRESSION TAG	UNP Q15843
I	92	ARG	-	EXPRESSION TAG	UNP Q15843
I	93	ALA	-	EXPRESSION TAG	UNP Q15843
I	94	SER	-	EXPRESSION TAG	UNP Q15843
I	95	VAL	-	EXPRESSION TAG	UNP Q15843
I	96	GLY	-	EXPRESSION TAG	UNP Q15843
I	97	SER	-	EXPRESSION TAG	UNP Q15843
I	98	GLY	-	EXPRESSION TAG	UNP Q15843
I	99	GLY	-	EXPRESSION TAG	UNP Q15843
I	100	SER	-	EXPRESSION TAG	UNP Q15843
I	172	ARG	ALA	ENGINEERED	UNP Q15843
J	89	GLY	-	EXPRESSION TAG	UNP Q15843
J	90	SER	-	EXPRESSION TAG	UNP Q15843
J	91	ARG	-	EXPRESSION TAG	UNP Q15843
J	92	ARG	-	EXPRESSION TAG	UNP Q15843
J	93	ALA	-	EXPRESSION TAG	UNP Q15843
J	94	SER	-	EXPRESSION TAG	UNP Q15843
J	95	VAL	-	EXPRESSION TAG	UNP Q15843
J	96	GLY	-	EXPRESSION TAG	UNP Q15843
J	97	SER	-	EXPRESSION TAG	UNP Q15843
J	98	GLY	-	EXPRESSION TAG	UNP Q15843
J	99	GLY	-	EXPRESSION TAG	UNP Q15843
J	100	SER	-	EXPRESSION TAG	UNP Q15843

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
J	172	ARG	ALA	ENGINEERED	UNP Q15843
K	89	GLY	-	EXPRESSION TAG	UNP Q15843
K	90	SER	-	EXPRESSION TAG	UNP Q15843
K	91	ARG	-	EXPRESSION TAG	UNP Q15843
K	92	ARG	-	EXPRESSION TAG	UNP Q15843
K	93	ALA	-	EXPRESSION TAG	UNP Q15843
K	94	SER	-	EXPRESSION TAG	UNP Q15843
K	95	VAL	-	EXPRESSION TAG	UNP Q15843
K	96	GLY	-	EXPRESSION TAG	UNP Q15843
K	97	SER	-	EXPRESSION TAG	UNP Q15843
K	98	GLY	-	EXPRESSION TAG	UNP Q15843
K	99	GLY	-	EXPRESSION TAG	UNP Q15843
K	100	SER	-	EXPRESSION TAG	UNP Q15843
K	172	ARG	ALA	ENGINEERED	UNP Q15843
L	89	GLY	-	EXPRESSION TAG	UNP Q15843
L	90	SER	-	EXPRESSION TAG	UNP Q15843
L	91	ARG	-	EXPRESSION TAG	UNP Q15843
L	92	ARG	-	EXPRESSION TAG	UNP Q15843
L	93	ALA	-	EXPRESSION TAG	UNP Q15843
L	94	SER	-	EXPRESSION TAG	UNP Q15843
L	95	VAL	-	EXPRESSION TAG	UNP Q15843
L	96	GLY	-	EXPRESSION TAG	UNP Q15843
L	97	SER	-	EXPRESSION TAG	UNP Q15843
L	98	GLY	-	EXPRESSION TAG	UNP Q15843
L	99	GLY	-	EXPRESSION TAG	UNP Q15843
L	100	SER	-	EXPRESSION TAG	UNP Q15843
L	172	ARG	ALA	ENGINEERED	UNP Q15843

- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

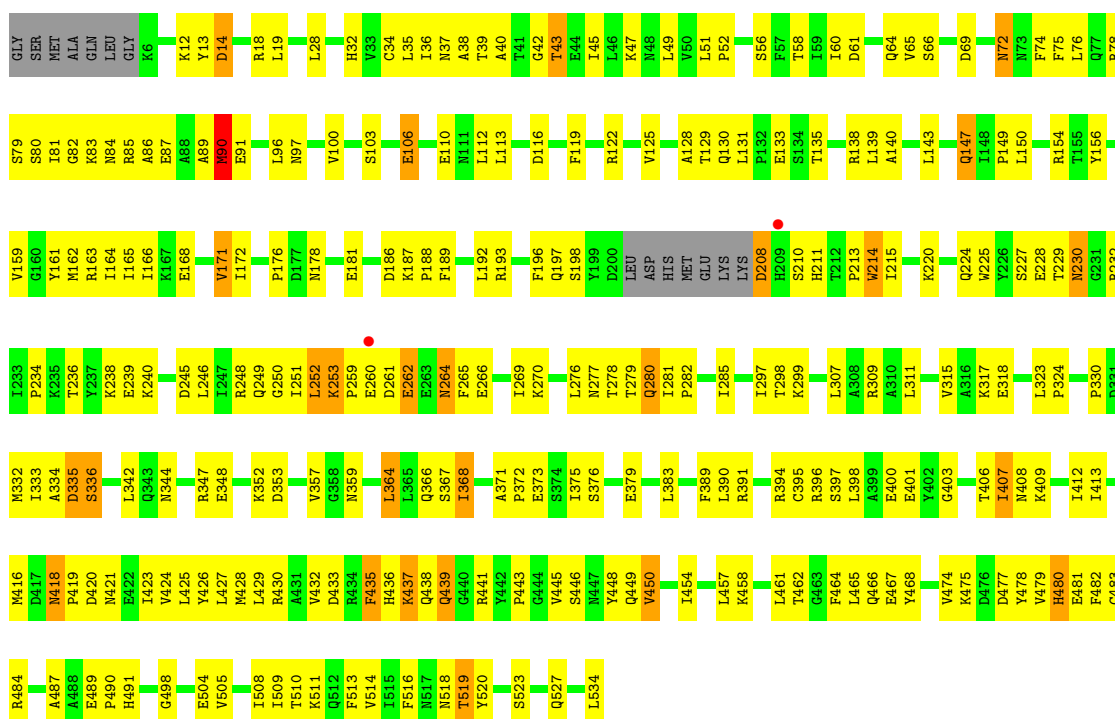
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	H	1	Total Zn 1 1	0	0
4	B	1	Total Zn 1 1	0	0
4	D	1	Total Zn 1 1	0	0
4	F	1	Total Zn 1 1	0	0

3 Residue-property plots

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

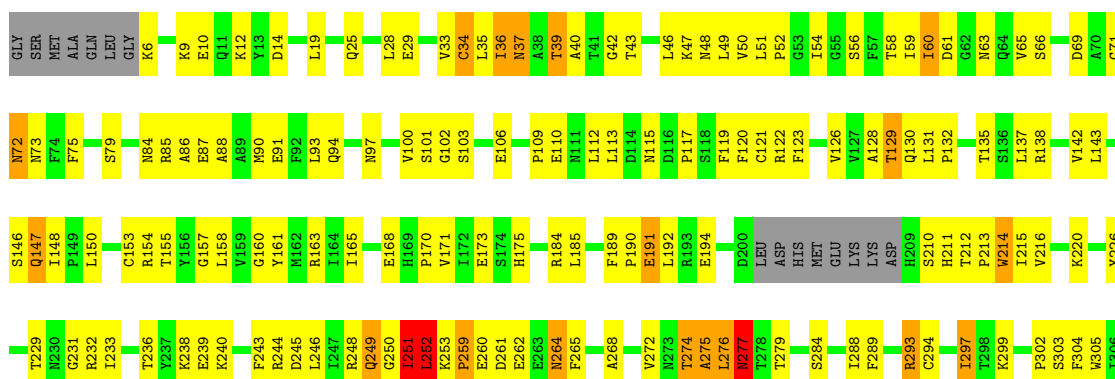
- Molecule 1: NEDD8-activating enzyme E1 regulatory subunit

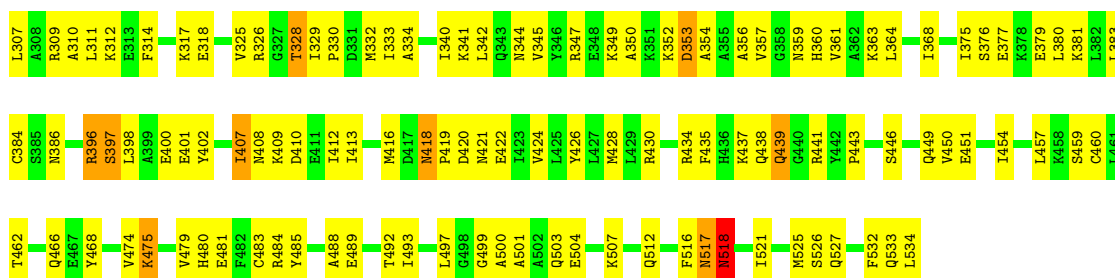
Chain A:



- Molecule 1: NEDD8-activating enzyme E1 regulatory subunit

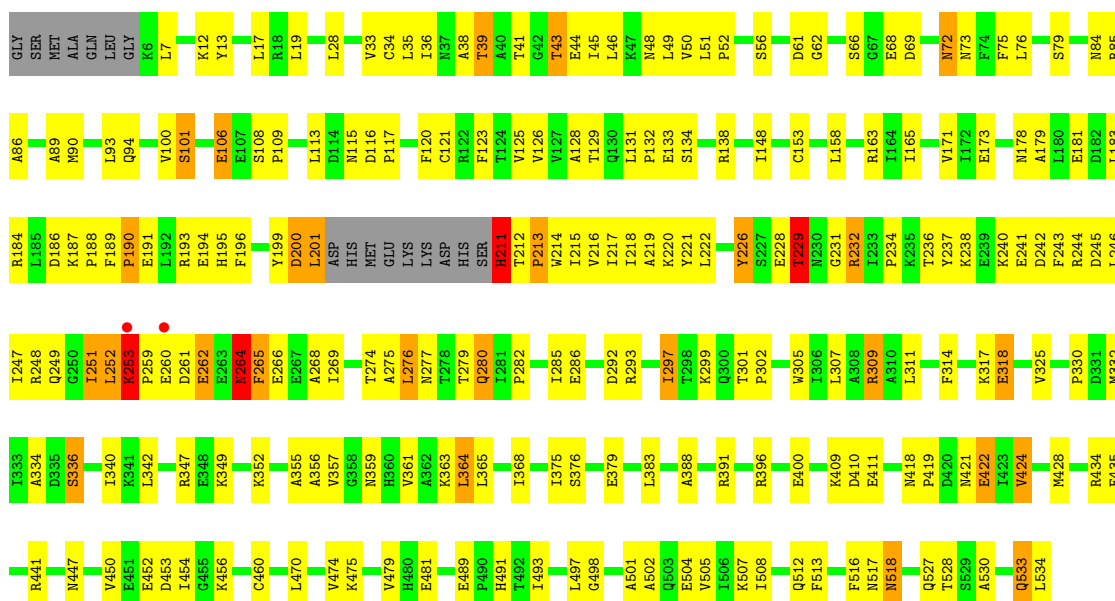
Chain C:





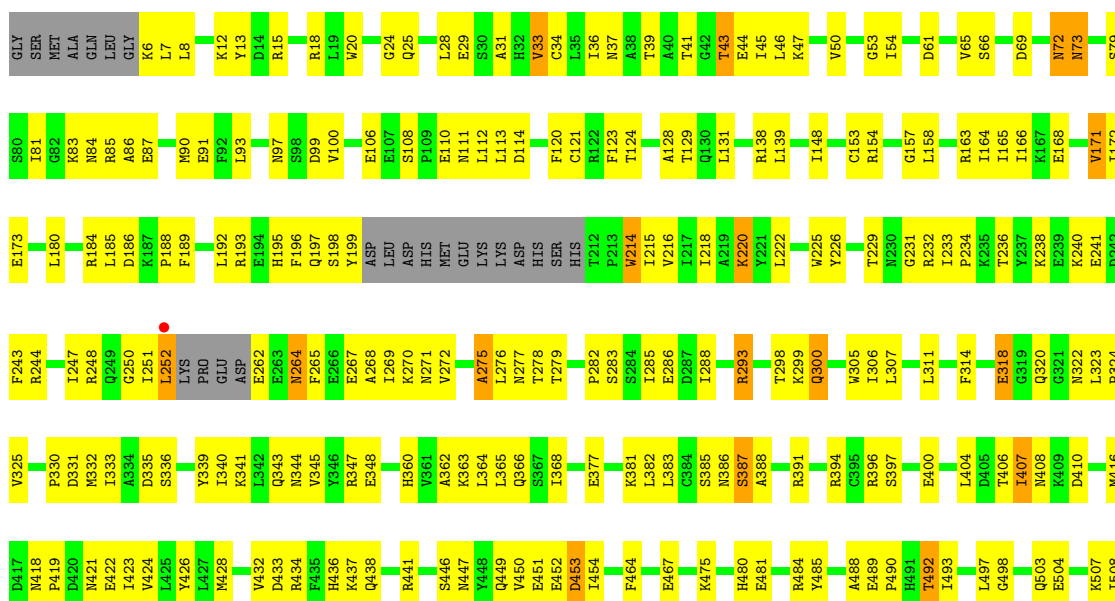
• Molecule 1: NEDD8-activating enzyme E1 regulatory subunit

Chain E:



• Molecule 1: NEDD8-activating enzyme E1 regulatory subunit

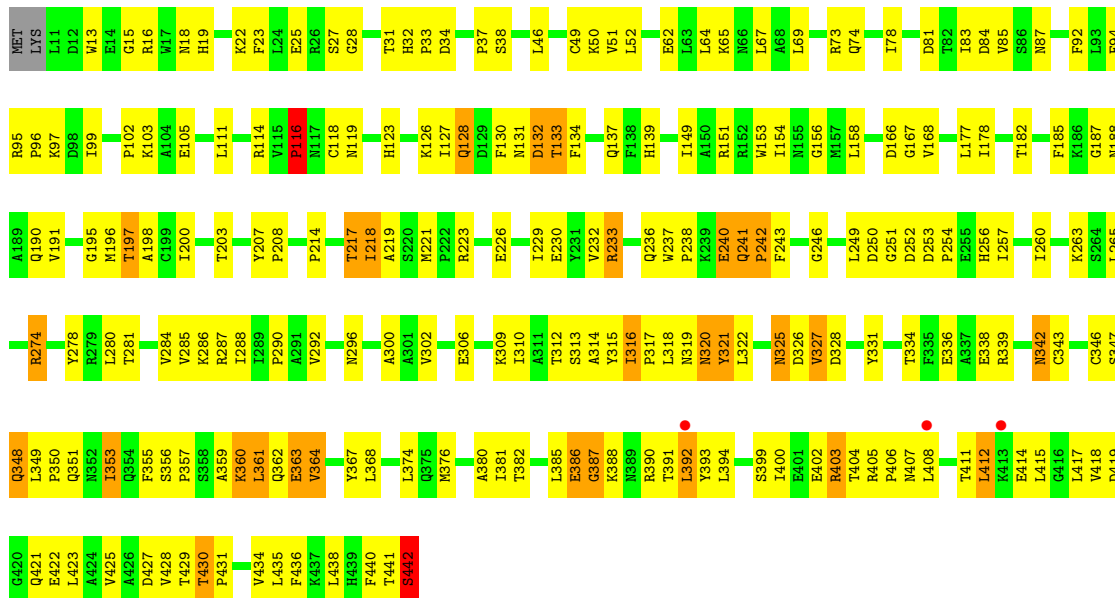
Chain G:





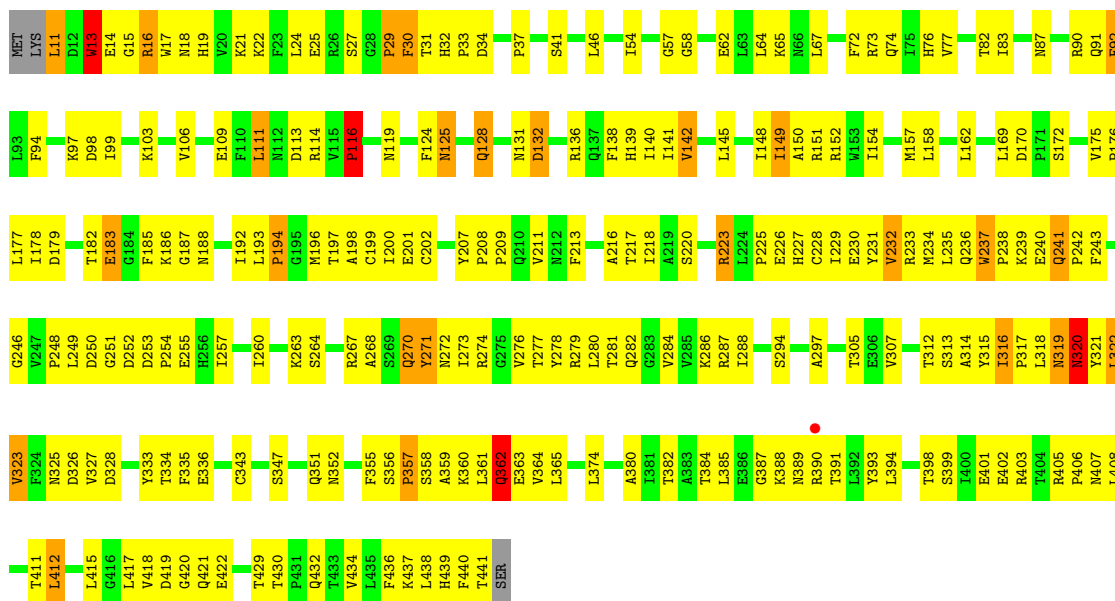
• Molecule 2: NEDD8-activating enzyme E1 catalytic subunit

Chain B:



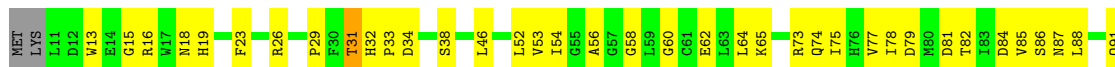
• Molecule 2: NEDD8-activating enzyme E1 catalytic subunit

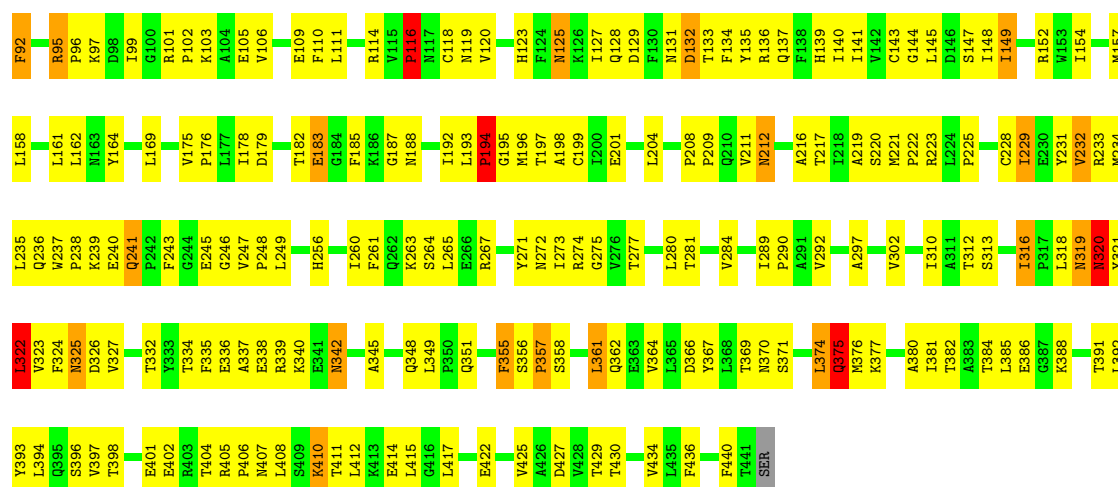
Chain D:



• Molecule 2: NEDD8-activating enzyme E1 catalytic subunit

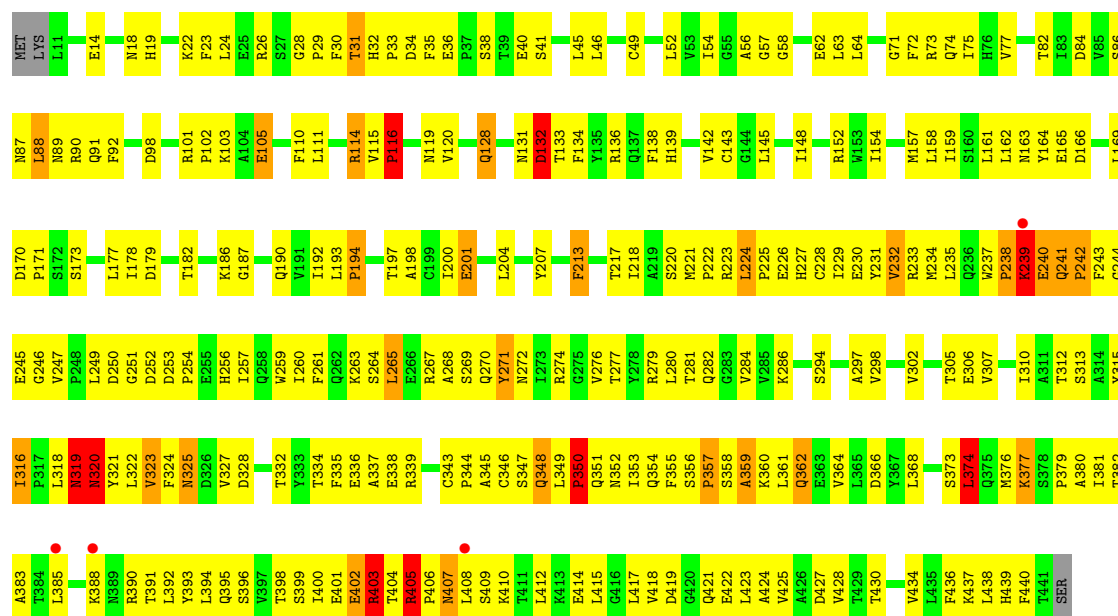
Chain F:





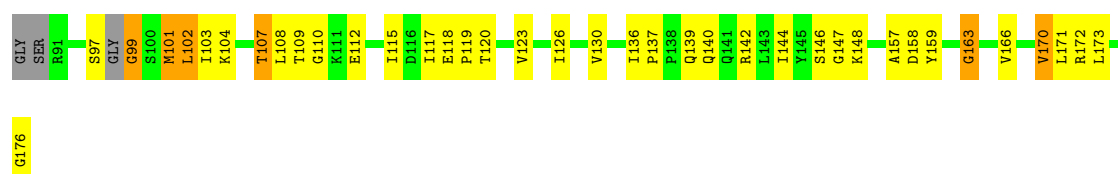
• Molecule 2: NEDD8-activating enzyme E1 catalytic subunit

Chain H:



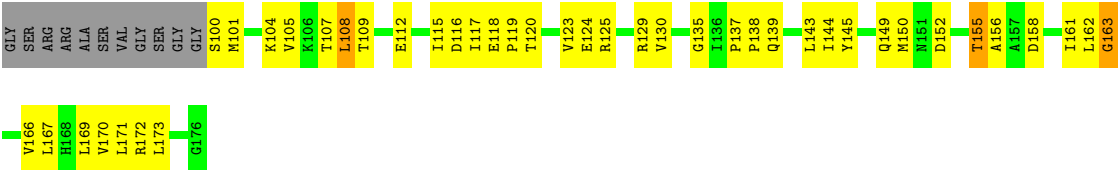
• Molecule 3: NEDD8

Chain I:



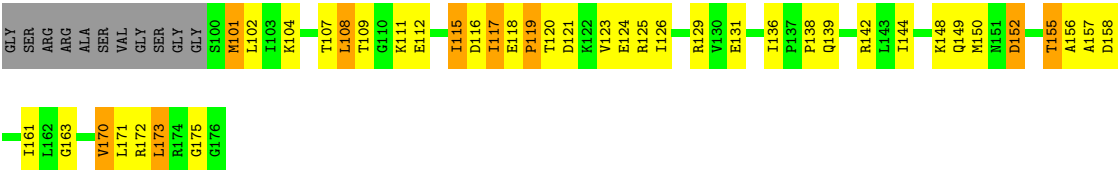
• Molecule 3: NEDD8

Chain J:



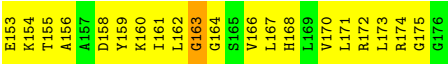
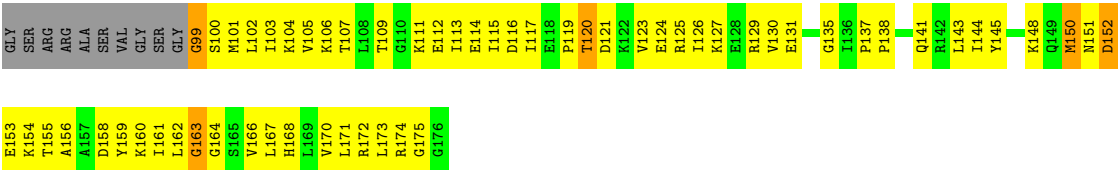
• Molecule 3: NEDD8

Chain K:



• Molecule 3: NEDD8

Chain L:



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	134.32Å 198.53Å 208.79Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 3.05 48.29 – 3.07	Depositor EDS
% Data completeness (in resolution range)	93.7 (50.00-3.05) 95.3 (48.29-3.07)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.90 (at 3.07Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.230 , 0.280 0.233 , 0.230	Depositor DCC
R_{free} test set	4982 reflections (4.98%)	DCC
Wilson B-factor (Å ²)	67.5	Xtriage
Anisotropy	0.524	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 41.6	EDS
Estimated twinning fraction	0.010 for -h,l,k	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 100052 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	32424	wwPDB-VP
Average B, all atoms (Å ²)	80.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.84% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

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

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	0.43	0/4184	0.68	1/5660 (0.0%)
1	C	0.45	0/4176	0.73	9/5649 (0.2%)
1	E	0.48	1/4172 (0.0%)	0.71	2/5643 (0.0%)
1	G	0.41	0/4114	0.67	1/5563 (0.0%)
2	B	0.50	1/3480 (0.0%)	0.75	4/4736 (0.1%)
2	D	0.44	0/3477	0.72	2/4732 (0.0%)
2	F	0.46	0/3473	0.73	3/4728 (0.1%)
2	H	0.50	0/3465	0.85	10/4718 (0.2%)
3	I	0.55	2/670 (0.3%)	0.77	0/891
3	J	0.43	0/617	0.82	1/823 (0.1%)
3	K	0.46	0/617	0.78	0/823
3	L	0.61	1/621 (0.2%)	0.98	5/828 (0.6%)
All	All	0.46	5/33066 (0.0%)	0.74	38/44794 (0.1%)

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
2	F	0	1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	442	SER	C-O	8.06	1.38	1.23
1	E	211	HIS	CG-CD2	-6.33	1.25	1.35
3	I	99	GLY	C-O	-5.79	1.14	1.23
3	I	99	GLY	CA-C	5.25	1.60	1.51
3	L	99	GLY	N-CA	5.09	1.53	1.46

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	320	ASN	N-CA-C	12.33	144.30	111.00
2	B	442	SER	N-CA-C	-9.77	84.61	111.00
2	D	320	ASN	N-CA-C	9.75	137.32	111.00
3	L	100	SER	N-CA-C	9.41	136.40	111.00
2	H	271	TYR	N-CA-C	-8.89	86.99	111.00
2	H	403	ARG	C-N-CA	-8.72	99.91	121.70
2	B	320	ASN	N-CA-C	8.68	134.43	111.00
1	C	518	ASN	N-CA-C	8.31	133.44	111.00
2	H	240	GLU	N-CA-C	-7.99	89.43	111.00
1	C	277	ASN	N-CA-CB	7.19	123.55	110.60
1	E	252	LEU	CA-CB-CG	-7.04	99.11	115.30
2	F	320	ASN	N-CA-C	7.00	129.91	111.00
3	L	101	MET	N-CA-C	-6.75	92.78	111.00
1	C	276	LEU	N-CA-C	6.74	129.20	111.00
2	H	240	GLU	C-N-CA	-6.69	104.97	121.70
2	F	375	GLN	N-CA-C	6.57	128.75	111.00
3	L	100	SER	CA-C-N	-6.55	102.79	117.20
2	D	319	ASN	N-CA-C	-6.54	93.35	111.00
2	H	239	LYS	N-CA-C	6.17	127.65	111.00
1	C	252	LEU	N-CA-C	6.14	127.58	111.00
2	B	386	GLU	N-CA-C	-5.95	94.94	111.00
1	C	210	SER	N-CA-C	-5.85	95.20	111.00
2	H	359	ALA	N-CA-C	5.79	126.62	111.00
2	B	240	GLU	N-CA-C	-5.74	95.50	111.00
3	J	100	SER	CA-C-N	-5.64	104.80	117.20
1	C	211	HIS	CB-CA-C	5.58	121.57	110.40
1	C	129	THR	CA-CB-CG2	-5.57	104.60	112.40
2	H	241	GLN	C-N-CD	-5.57	108.35	120.60
2	F	322	LEU	CA-CB-CG	5.50	127.96	115.30
2	H	319	ASN	C-N-CA	-5.43	108.12	121.70
1	A	519	THR	N-CA-C	5.40	125.58	111.00
3	L	120	THR	N-CA-C	5.29	125.29	111.00
3	L	100	SER	O-C-N	5.25	131.09	122.70
1	C	260	GLU	N-CA-C	-5.19	96.98	111.00
2	H	321	TYR	N-CA-CB	-5.12	101.38	110.60
1	E	232	ARG	N-CA-C	5.11	124.80	111.00
1	G	252	LEU	CA-CB-CG	-5.08	103.62	115.30
1	C	129	THR	CA-C-N	5.07	128.35	117.20

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	F	319	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	4104	0	4052	264	0
1	C	4096	0	4048	255	0
1	E	4093	0	4059	220	0
1	G	4038	0	4009	220	0
2	B	3402	0	3380	253	0
2	D	3399	0	3386	241	0
2	F	3395	0	3375	269	0
2	H	3387	0	3354	311	0
3	I	666	0	703	49	0
3	J	612	0	648	46	0
3	K	612	0	648	47	0
3	L	616	0	651	74	0
4	B	1	0	0	0	0
4	D	1	0	0	0	0
4	F	1	0	0	0	0
4	H	1	0	0	0	0
All	All	32424	0	32313	2093	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 32.

All (2093) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:252:LEU:O	1:C:259:PRO:HD2	1.17	1.33
1:A:252:LEU:O	1:A:259:PRO:HD2	1.25	1.30
1:C:517:ASN:ND2	1:C:518:ASN:H	1.31	1.28
2:F:371:SER:CB	2:F:374:LEU:HD11	1.64	1.26
1:C:517:ASN:HD22	1:C:518:ASN:N	1.35	1.24
2:F:371:SER:HB3	2:F:374:LEU:HD11	1.20	1.12
1:A:259:PRO:HB2	1:A:262:GLU:OE1	1.51	1.11

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:K:123:VAL:HB	3:K:152:ASP:HA	1.29	1.11
1:G:333:ILE:HG22	2:H:223:ARG:NH2	1.66	1.10
2:F:149:ILE:H	2:F:149:ILE:HD12	1.06	1.10
1:A:36:ILE:HD12	1:A:60:ILE:HB	1.25	1.10
3:J:123:VAL:HB	3:J:152:ASP:HA	1.33	1.09
2:D:149:ILE:HD12	2:D:149:ILE:H	1.11	1.09
2:B:338:GLU:HG3	3:I:148:LYS:HG2	1.30	1.09
2:F:325:ASN:HD21	2:F:327:VAL:HG23	0.98	1.08
2:D:207:TYR:CE1	3:J:172:ARG:HD3	1.89	1.07
1:E:226:TYR:CD1	1:E:231:GLY:HA2	1.89	1.07
1:A:208:ASP:N	1:A:211:HIS:HB2	1.69	1.07
2:B:357:PRO:HD2	2:B:442:SER:O	1.56	1.05
1:C:35:LEU:HD11	1:C:129:THR:CG2	1.87	1.05
1:G:333:ILE:HG22	2:H:223:ARG:HH22	1.13	1.04
1:A:518:ASN:HD22	1:A:534:LEU:N	1.56	1.03
2:H:362:GLN:HG2	2:H:408:LEU:HB3	1.36	1.03
1:C:35:LEU:HD11	1:C:129:THR:HG21	1.42	1.01
2:H:403:ARG:HG2	2:H:403:ARG:O	1.54	1.01
2:H:232:VAL:HG21	2:H:264:SER:HA	1.39	1.00
1:C:252:LEU:O	1:C:259:PRO:CD	2.10	1.00
1:E:226:TYR:HD1	1:E:231:GLY:HA2	1.23	1.00
1:A:236:THR:HB	1:A:239:GLU:HG3	1.43	0.99
2:F:325:ASN:ND2	2:F:327:VAL:HG23	1.77	0.98
1:G:489:GLU:H	2:H:19:HIS:HD2	1.00	0.98
2:F:380:ALA:CB	2:F:394:LEU:HD12	1.93	0.98
2:H:380:ALA:CB	2:H:394:LEU:HD12	1.93	0.98
1:C:130:GLN:OE1	1:C:155:THR:HB	1.63	0.97
1:A:518:ASN:ND2	1:A:534:LEU:N	2.12	0.97
2:D:402:GLU:HA	2:D:405:ARG:HH12	1.26	0.97
3:K:155:THR:HG22	3:K:158:ASP:OD1	1.63	0.97
1:E:184:ARG:HH12	1:E:325:VAL:HG22	1.27	0.97
2:F:362:GLN:HG2	2:F:408:LEU:HD22	1.45	0.96
1:C:259:PRO:HB2	1:C:262:GLU:OE1	1.66	0.96
2:H:128:GLN:HE21	2:H:128:GLN:H	1.12	0.96
2:F:361:LEU:HB3	2:F:408:LEU:HD23	1.48	0.96
2:B:386:GLU:O	2:B:388:LYS:N	1.99	0.95
1:A:518:ASN:ND2	1:A:534:LEU:H	1.64	0.95
1:A:252:LEU:C	1:A:259:PRO:HD2	1.87	0.95
2:B:207:TYR:CE2	3:I:172:ARG:HG2	2.02	0.95
1:E:347:ARG:HH22	2:F:274:ARG:HD2	1.30	0.94
1:E:211:HIS:CE1	2:F:221:MET:SD	2.60	0.94
2:F:376:MET:HB3	2:F:427:ASP:OD1	1.67	0.94

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:F:374:LEU:H	2:F:374:LEU:HD12	1.33	0.94
1:G:489:GLU:H	2:H:19:HIS:CD2	1.85	0.94
1:G:293:ARG:HH11	1:G:293:ARG:HG2	1.29	0.94
2:F:371:SER:HB3	2:F:374:LEU:CD1	1.98	0.93
3:I:99:GLY:O	3:I:118:GLU:CB	2.16	0.93
1:A:252:LEU:O	1:A:259:PRO:CD	2.16	0.93
2:D:257:ILE:HD13	2:D:282:GLN:HG2	1.50	0.93
3:I:107:THR:HG22	3:I:109:THR:H	1.31	0.92
2:D:149:ILE:CD1	2:D:149:ILE:H	1.81	0.92
2:B:214:PRO:HG2	2:B:217:THR:HG23	1.52	0.92
1:C:35:LEU:CD1	1:C:129:THR:CG2	2.48	0.92
1:E:201:LEU:HD23	1:E:201:LEU:N	1.85	0.91
2:F:64:LEU:HD11	2:F:77:VAL:HG21	1.51	0.91
2:F:362:GLN:HE21	2:F:408:LEU:HD13	1.35	0.90
2:B:236:GLN:HE22	2:B:263:LYS:HD2	1.35	0.90
3:L:107:THR:HG23	3:L:109:THR:H	1.35	0.90
2:F:342:ASN:HD22	2:F:342:ASN:H	1.18	0.90
2:F:149:ILE:H	2:F:149:ILE:CD1	1.82	0.90
2:F:380:ALA:HB1	2:F:394:LEU:HA	1.54	0.90
2:B:342:ASN:HD22	2:B:342:ASN:H	1.18	0.89
2:H:223:ARG:HB2	2:H:227:HIS:CE1	2.08	0.89
2:F:74:GLN:HE22	2:F:119:ASN:HD22	1.15	0.89
1:G:121:CYS:HA	1:G:148:ILE:HD11	1.53	0.89
2:D:380:ALA:HB2	2:D:394:LEU:HD12	1.53	0.89
3:K:118:GLU:HG2	3:K:121:ASP:OD1	1.73	0.88
1:G:180:LEU:H	1:G:180:LEU:HD12	1.39	0.88
2:B:380:ALA:HB2	2:B:394:LEU:HD12	1.56	0.88
2:H:128:GLN:NE2	2:H:128:GLN:H	1.72	0.87
3:L:120:THR:HA	3:L:155:THR:OG1	1.73	0.87
2:F:131:ASN:HB3	2:H:131:ASN:HD22	1.35	0.87
2:F:340:LYS:HB3	2:F:342:ASN:HD21	1.38	0.87
2:B:380:ALA:CB	2:B:394:LEU:HD12	2.04	0.87
2:B:327:VAL:HG23	2:B:328:ASP:H	1.39	0.87
2:D:232:VAL:HG11	2:D:263:LYS:HB2	1.56	0.86
2:F:380:ALA:HB1	2:F:394:LEU:HD12	1.56	0.86
3:J:161:ILE:HD11	3:J:167:LEU:HD21	1.56	0.86
1:E:211:HIS:CD2	1:E:211:HIS:N	2.39	0.86
2:H:58:GLY:H	2:H:91:GLN:HG2	1.40	0.86
1:E:184:ARG:NH1	1:E:325:VAL:HG22	1.89	0.86
2:H:318:LEU:HD11	2:H:334:THR:CG2	2.06	0.86
2:F:325:ASN:HD21	2:F:327:VAL:CG2	1.87	0.85
2:B:343:CYS:O	2:B:347:SER:HB3	1.76	0.85

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:36:ILE:HB	1:C:128:ALA:HA	1.58	0.85
1:E:297:ILE:H	1:E:297:ILE:HD12	1.41	0.85
1:A:307:LEU:HB3	1:A:383:LEU:HD22	1.57	0.85
1:E:212:THR:OG1	1:E:217:ILE:HD11	1.76	0.85
3:L:170:VAL:HG12	3:L:171:LEU:H	1.40	0.85
2:H:380:ALA:HB2	2:H:394:LEU:HD12	1.59	0.85
1:E:307:LEU:HD13	1:E:383:LEU:HD22	1.58	0.85
1:G:421:ASN:O	1:G:424:VAL:HG23	1.76	0.85
2:B:218:ILE:HD11	2:B:230:GLU:HG2	1.59	0.85
1:C:37:ASN:O	1:C:129:THR:OG1	1.94	0.84
2:F:134:PHE:O	2:F:137:GLN:HG2	1.77	0.84
1:E:51:LEU:HD11	2:F:92:PHE:HB3	1.58	0.84
2:H:271:TYR:CD2	2:H:271:TYR:N	2.42	0.84
2:B:355:PHE:O	2:B:440:PHE:HA	1.76	0.84
2:F:411:THR:H	2:F:414:GLU:HB3	1.40	0.84
3:K:125:ARG:HD3	3:K:129:ARG:HH21	1.41	0.83
2:B:361:LEU:HD23	2:B:408:LEU:HD23	1.60	0.83
2:D:141:ILE:HD12	2:D:158:LEU:HD21	1.58	0.83
2:F:149:ILE:N	2:F:149:ILE:HD12	1.90	0.83
2:F:411:THR:H	2:F:414:GLU:CB	1.91	0.83
2:F:371:SER:OG	2:F:374:LEU:HD11	1.79	0.83
1:A:229:THR:CG2	1:A:232:ARG:HB3	2.08	0.83
2:H:318:LEU:HG	2:H:319:ASN:O	1.79	0.83
1:G:518:ASN:HB3	1:G:533:GLN:HA	1.61	0.83
1:E:248:ARG:O	1:E:251:ILE:CG1	2.27	0.83
3:L:123:VAL:HG11	3:L:150:MET:HB3	1.60	0.82
2:B:316:ILE:H	2:B:316:ILE:HD12	1.42	0.82
2:D:380:ALA:CB	2:D:394:LEU:HD12	2.08	0.82
2:B:342:ASN:H	2:B:342:ASN:ND2	1.75	0.82
2:H:318:LEU:HD11	2:H:334:THR:HG23	1.61	0.81
1:G:518:ASN:CB	1:G:533:GLN:HA	2.10	0.81
1:C:35:LEU:CD1	1:C:129:THR:HG21	2.08	0.81
1:C:299:LYS:HA	1:C:368:ILE:HG23	1.63	0.81
1:A:418:ASN:ND2	1:A:420:ASP:H	1.79	0.81
2:D:149:ILE:HD12	2:D:149:ILE:N	1.94	0.81
1:A:224:GLN:NE2	1:A:246:LEU:HD11	1.96	0.81
1:C:35:LEU:CD1	1:C:129:THR:HG23	2.10	0.80
3:J:101:MET:SD	3:J:162:LEU:HA	2.21	0.80
1:C:527:GLN:HB2	2:D:318:LEU:HD13	1.64	0.80
2:B:13:TRP:CZ3	2:B:116:PRO:HG2	2.16	0.80
2:B:240:GLU:O	2:B:242:PRO:CD	2.30	0.80
1:E:252:LEU:C	1:E:259:PRO:HD2	2.01	0.80

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:211:HIS:HD2	1:E:211:HIS:N	1.76	0.79
2:F:74:GLN:NE2	2:F:119:ASN:HD22	1.79	0.79
2:D:318:LEU:HG	2:D:319:ASN:O	1.82	0.79
3:J:118:GLU:HG3	3:J:120:THR:HG22	1.62	0.79
2:H:277:THR:HG23	2:H:280:LEU:H	1.47	0.79
3:I:99:GLY:O	3:I:118:GLU:HB2	1.81	0.79
1:C:229:THR:CG2	1:C:232:ARG:HB2	2.12	0.79
3:J:170:VAL:HG22	3:J:171:LEU:H	1.48	0.79
3:I:99:GLY:O	3:I:118:GLU:HB3	1.81	0.79
1:E:201:LEU:N	1:E:201:LEU:CD2	2.45	0.79
2:F:412:LEU:HD13	2:F:440:PHE:HE2	1.49	0.78
2:H:231:TYR:HA	2:H:235:LEU:HD23	1.64	0.78
1:G:6:LYS:HD2	1:G:6:LYS:N	1.97	0.78
2:H:380:ALA:HB1	2:H:394:LEU:HD12	1.64	0.78
1:A:462:THR:O	1:A:466:GLN:HG3	1.83	0.78
1:C:518:ASN:HB2	1:C:533:GLN:HA	1.63	0.78
1:G:360:HIS:O	1:G:364:LEU:HD13	1.83	0.78
1:C:297:ILE:H	1:C:297:ILE:HD13	1.48	0.78
1:E:262:GLU:HA	1:E:262:GLU:OE2	1.83	0.78
1:A:259:PRO:CB	1:A:262:GLU:OE1	2.31	0.78
1:A:125:VAL:HG12	1:A:149:PRO:HB2	1.65	0.78
2:F:87:ASN:ND2	2:F:103:LYS:HE2	1.99	0.78
1:E:186:ASP:OD2	1:E:279:THR:HB	1.84	0.78
2:H:241:GLN:OE1	2:H:245:GLU:HA	1.83	0.77
2:D:402:GLU:HA	2:D:405:ARG:NH1	1.98	0.77
1:A:418:ASN:HD22	1:A:420:ASP:H	1.29	0.77
2:B:240:GLU:O	2:B:242:PRO:HD3	1.84	0.77
2:H:351:GLN:H	2:H:436:PHE:HA	1.47	0.77
2:F:277:THR:HG23	2:F:280:LEU:H	1.47	0.77
1:G:333:ILE:HA	2:H:223:ARG:CZ	2.15	0.77
1:C:35:LEU:HD11	1:C:129:THR:HG23	1.65	0.77
1:A:317:LYS:HB3	1:A:318:GLU:OE1	1.85	0.77
2:B:356:SER:HB3	2:B:441:THR:OG1	1.84	0.77
1:G:489:GLU:N	2:H:19:HIS:HD2	1.81	0.77
2:F:340:LYS:HB3	2:F:342:ASN:ND2	2.00	0.77
2:H:162:LEU:HA	2:H:173:SER:OG	1.83	0.77
2:H:403:ARG:O	2:H:403:ARG:CG	2.33	0.77
1:G:447:ASN:HD22	2:H:26:ARG:HH21	1.30	0.77
2:H:316:ILE:HD12	2:H:316:ILE:H	1.48	0.77
1:A:28:LEU:HD12	1:A:509:ILE:HG21	1.67	0.76
3:L:105:VAL:HG23	3:L:167:LEU:HB2	1.67	0.76
1:G:226:TYR:OH	1:G:233:ILE:HG22	1.85	0.76

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:163:ARG:CZ	1:C:518:ASN:OD1	2.34	0.76
1:C:418:ASN:ND2	1:C:420:ASP:H	1.81	0.76
1:C:51:LEU:HD11	2:D:92:PHE:HB3	1.68	0.76
1:E:248:ARG:O	1:E:251:ILE:HG13	1.84	0.76
2:H:164:TYR:CE2	2:H:169:LEU:HB2	2.21	0.76
1:G:267:GLU:OE2	1:G:333:ILE:HG12	1.85	0.76
3:L:120:THR:CA	3:L:155:THR:OG1	2.34	0.76
2:H:217:THR:HG21	2:H:223:ARG:HH21	1.50	0.75
2:B:357:PRO:CD	2:B:442:SER:O	2.34	0.75
1:G:41:THR:O	1:G:45:ILE:HG13	1.85	0.75
1:A:75:PHE:O	1:A:76:LEU:HD23	1.85	0.75
1:C:129:THR:O	1:C:153:CYS:O	2.03	0.75
1:E:253:LYS:HA	1:E:253:LYS:HE3	1.69	0.75
2:D:380:ALA:CB	2:D:394:LEU:CD1	2.64	0.75
1:E:489:GLU:H	2:F:19:HIS:CD2	2.04	0.75
2:F:62:GLU:HG2	2:F:297:ALA:HA	1.68	0.74
1:A:229:THR:HG21	1:A:232:ARG:HB3	1.68	0.74
1:E:409:LYS:HG2	1:E:470:LEU:HD21	1.68	0.74
1:A:112:LEU:H	1:A:112:LEU:HD22	1.50	0.74
1:C:259:PRO:HB2	1:C:262:GLU:CD	2.07	0.74
2:B:342:ASN:HD22	2:B:342:ASN:N	1.79	0.74
1:G:341:LYS:O	1:G:345:VAL:HG23	1.87	0.74
1:E:61:ASP:HB3	1:E:86:ALA:HB2	1.70	0.74
2:H:62:GLU:HG2	2:H:297:ALA:HA	1.69	0.74
2:B:81:ASP:HB2	2:B:103:LYS:HD2	1.68	0.74
1:C:115:ASN:O	1:C:117:PRO:HD3	1.88	0.74
1:C:252:LEU:C	1:C:259:PRO:HD2	2.06	0.74
1:C:396:ARG:HD2	1:C:534:LEU:O	1.88	0.74
1:A:34:CYS:SG	1:A:36:ILE:HD11	2.28	0.74
2:B:214:PRO:HG2	2:B:217:THR:CG2	2.18	0.74
2:B:419:ASP:OD1	2:B:440:PHE:HD1	1.71	0.74
1:G:293:ARG:NH1	1:G:293:ARG:HG2	2.00	0.73
2:B:229:ILE:HD13	2:B:281:THR:HA	1.69	0.73
2:B:95:ARG:HH11	2:B:95:ARG:HG3	1.52	0.73
1:C:229:THR:HG21	1:C:232:ARG:HB2	1.69	0.73
2:H:240:GLU:O	2:H:242:PRO:HD3	1.88	0.73
2:H:339:ARG:HH21	2:H:346:CYS:HB3	1.52	0.73
2:B:13:TRP:HZ3	2:B:116:PRO:HG2	1.53	0.73
1:C:397:SER:OG	1:C:400:GLU:HG3	1.89	0.73
1:C:84:ASN:HB3	1:C:87:GLU:HB3	1.71	0.73
2:F:316:ILE:H	2:F:316:ILE:HD12	1.52	0.73
1:A:36:ILE:CD1	1:A:60:ILE:HB	2.11	0.73

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:H:252:ASP:O	2:H:254:PRO:HD3	1.88	0.73
1:C:143:LEU:HD22	1:C:148:ILE:HB	1.70	0.73
2:H:356:SER:O	2:H:358:SER:N	2.21	0.73
2:H:223:ARG:HB2	2:H:227:HIS:HE1	1.52	0.73
1:A:236:THR:CB	1:A:239:GLU:HG3	2.19	0.73
2:F:382:THR:HG22	2:F:391:THR:HA	1.70	0.73
1:E:307:LEU:HB3	1:E:383:LEU:HD22	1.71	0.73
2:F:217:THR:HG21	2:F:223:ARG:NH2	2.03	0.73
2:F:348:GLN:O	2:F:349:LEU:HD12	1.89	0.73
2:F:199:CYS:SG	2:F:201:GLU:HB2	2.29	0.73
2:B:200:ILE:HD13	3:I:172:ARG:HH22	1.54	0.73
1:A:61:ASP:HB3	1:A:86:ALA:HB2	1.69	0.73
1:C:130:GLN:HG2	1:C:154:ARG:HA	1.71	0.72
2:D:223:ARG:HB2	2:D:223:ARG:HH11	1.52	0.72
2:F:371:SER:O	2:F:374:LEU:HD12	1.90	0.72
2:H:404:THR:HG22	2:H:407:ASN:HD22	1.54	0.72
2:B:318:LEU:HD11	2:B:334:THR:CG2	2.19	0.72
1:A:236:THR:HG22	1:A:238:LYS:H	1.55	0.72
2:B:353:ILE:HG23	2:B:355:PHE:HD1	1.54	0.72
2:H:38:SER:HB3	2:H:41:SER:OG	1.90	0.72
2:F:58:GLY:H	2:F:91:GLN:HG2	1.54	0.72
2:F:351:GLN:HB2	2:F:436:PHE:CD2	2.24	0.72
1:G:45:ILE:HG12	1:G:498:GLY:HA2	1.72	0.72
2:D:162:LEU:HD22	2:D:169:LEU:HD11	1.71	0.72
1:E:347:ARG:NH2	2:F:274:ARG:HD2	2.04	0.71
1:A:489:GLU:H	2:B:19:HIS:CD2	2.08	0.71
2:D:16:ARG:HD3	2:D:17:TRP:NE1	2.05	0.71
1:C:259:PRO:C	1:C:261:ASP:H	1.93	0.71
1:A:259:PRO:C	1:A:261:ASP:H	1.94	0.71
1:E:211:HIS:HE1	2:F:221:MET:SD	2.13	0.71
2:H:224:LEU:HG	2:H:227:HIS:CE1	2.26	0.71
2:B:217:THR:HA	2:B:221:MET:HG3	1.70	0.71
1:C:409:LYS:HE2	1:C:468:TYR:O	1.91	0.71
2:B:253:ASP:HB3	2:B:256:HIS:HB2	1.70	0.71
3:L:170:VAL:HG12	3:L:171:LEU:N	2.05	0.71
2:H:207:TYR:CE1	3:L:172:ARG:HD3	2.25	0.71
1:A:266:GLU:O	1:A:270:LYS:HG3	1.91	0.71
1:G:446:SER:HB2	1:G:449:GLN:HG3	1.73	0.71
2:H:325:ASN:ND2	2:H:327:VAL:HG23	2.06	0.71
1:C:251:ILE:HG22	1:C:262:GLU:OE1	1.90	0.70
2:H:235:LEU:O	2:H:238:PRO:HD2	1.91	0.70
1:A:397:SER:OG	1:A:400:GLU:HG3	1.90	0.70

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:217:THR:HB	2:B:223:ARG:NH2	2.06	0.70
2:D:232:VAL:CG1	2:D:263:LYS:HB2	2.20	0.70
1:G:196:PHE:O	1:G:220:LYS:HE3	1.91	0.70
2:F:374:LEU:N	2:F:374:LEU:HD12	1.98	0.70
1:E:184:ARG:HH11	1:E:184:ARG:HG3	1.55	0.70
2:F:411:THR:OG1	2:F:414:GLU:HB2	1.91	0.70
2:B:325:ASN:HD21	2:B:327:VAL:HG22	1.56	0.70
2:B:240:GLU:HG3	2:B:240:GLU:O	1.91	0.70
2:H:239:LYS:O	2:H:239:LYS:HG3	1.91	0.70
1:A:264:ASN:N	1:A:264:ASN:HD22	1.89	0.70
2:H:318:LEU:C	2:H:319:ASN:O	2.24	0.70
2:H:74:GLN:HE22	2:H:119:ASN:HD22	1.39	0.70
1:A:299:LYS:HG2	1:A:368:ILE:O	1.90	0.70
2:B:74:GLN:HE22	2:B:119:ASN:HD22	1.40	0.70
3:K:116:ASP:O	3:K:117:ILE:HD12	1.92	0.70
2:F:241:GLN:HG2	2:F:245:GLU:HA	1.72	0.70
1:C:248:ARG:O	1:C:251:ILE:HG13	1.92	0.70
2:H:234:MET:O	2:H:235:LEU:HD13	1.92	0.70
1:A:297:ILE:N	1:A:297:ILE:HD12	2.06	0.70
2:F:32:HIS:CD2	2:F:34:ASP:H	2.10	0.70
2:D:325:ASN:HD21	2:D:327:VAL:HG22	1.56	0.70
2:H:164:TYR:CD2	2:H:169:LEU:HB2	2.27	0.69
2:H:398:THR:O	2:H:402:GLU:HG3	1.92	0.69
2:H:157:MET:O	2:H:157:MET:HE3	1.92	0.69
2:H:187:GLY:HA2	3:L:173:LEU:HD13	1.74	0.69
2:H:148:ILE:HG12	3:L:174:ARG:HG2	1.73	0.69
2:B:64:LEU:HB3	2:B:111:LEU:CD1	2.22	0.69
2:H:294:SER:O	2:H:298:VAL:HG23	1.92	0.69
1:A:84:ASN:OD1	1:A:106:GLU:HG2	1.92	0.69
1:A:518:ASN:CG	1:A:519:THR:H	1.94	0.69
2:D:318:LEU:HD11	2:D:334:THR:CG2	2.22	0.69
1:A:36:ILE:O	1:A:60:ILE:O	2.08	0.69
2:D:141:ILE:CD1	2:D:158:LEU:HD21	2.22	0.69
1:C:317:LYS:HB3	1:C:318:GLU:OE1	1.93	0.69
2:D:380:ALA:HB1	2:D:394:LEU:CD1	2.21	0.69
1:G:507:LYS:HG2	1:G:513:PHE:HB2	1.73	0.69
2:F:237:TRP:HB3	2:F:238:PRO:HD3	1.75	0.69
2:B:356:SER:HB3	2:B:442:SER:H	1.55	0.69
2:B:380:ALA:CB	2:B:394:LEU:CD1	2.71	0.69
1:C:146:SER:O	1:C:147:GLN:HB2	1.90	0.69
1:G:185:LEU:HD12	1:G:275:ALA:HB1	1.73	0.69
2:B:361:LEU:HB3	2:B:408:LEU:HA	1.74	0.69

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:297:ILE:H	1:E:297:ILE:CD1	2.03	0.69
1:A:527:GLN:HB2	2:B:318:LEU:HD13	1.75	0.69
2:F:125:ASN:HD22	2:F:125:ASN:N	1.91	0.69
2:B:203:THR:OG1	3:I:172:ARG:NH1	2.25	0.68
2:F:131:ASN:HD22	2:H:131:ASN:HB3	1.58	0.68
2:B:64:LEU:HB3	2:B:111:LEU:HD11	1.74	0.68
1:G:33:VAL:CG2	1:G:54:ILE:HG12	2.22	0.68
1:C:274:THR:O	1:C:276:LEU:N	2.26	0.68
1:A:154:ARG:HB3	1:A:161:TYR:HB3	1.76	0.68
1:C:163:ARG:NH1	1:C:165:ILE:HG12	2.07	0.68
2:D:17:TRP:O	2:D:21:LYS:HB2	1.92	0.68
3:K:104:LYS:HD2	3:K:112:GLU:OE2	1.91	0.68
2:H:355:PHE:O	2:H:440:PHE:HA	1.93	0.68
2:F:355:PHE:CE1	2:F:364:VAL:HG22	2.28	0.68
2:F:217:THR:HG21	2:F:223:ARG:HH22	1.56	0.68
2:D:125:ASN:HD22	2:D:125:ASN:N	1.92	0.68
1:A:61:ASP:OD2	1:A:85:ARG:HB3	1.92	0.68
2:D:74:GLN:NE2	2:D:119:ASN:HD22	1.92	0.68
2:F:54:ILE:HG22	2:F:145:LEU:HD21	1.75	0.68
2:H:271:TYR:HD2	2:H:271:TYR:N	1.90	0.68
2:H:232:VAL:HG21	2:H:264:SER:CA	2.21	0.68
1:C:46:LEU:O	1:C:50:VAL:HG23	1.94	0.68
1:G:248:ARG:O	1:G:251:ILE:HG13	1.93	0.68
2:D:322:LEU:HD12	2:D:323:VAL:N	2.08	0.68
1:E:309:ARG:HH11	1:E:309:ARG:HG3	1.59	0.68
2:B:182:THR:HG21	2:B:296:ASN:OD1	1.94	0.68
2:F:362:GLN:NE2	2:F:408:LEU:HD13	2.08	0.68
1:A:87:GLU:O	1:A:91:GLU:HG3	1.94	0.68
2:H:192:ILE:HG22	2:H:194:PRO:HD3	1.76	0.68
1:C:163:ARG:HD2	1:C:518:ASN:O	1.93	0.68
1:E:259:PRO:C	1:E:261:ASP:H	1.96	0.67
1:C:42:GLY:HA2	1:C:129:THR:HG21	1.77	0.67
2:H:228:CYS:O	2:H:232:VAL:HG23	1.94	0.67
2:F:357:PRO:HA	2:F:412:LEU:HB2	1.76	0.67
2:F:229:ILE:HD13	2:F:281:THR:HA	1.75	0.67
2:H:357:PRO:HA	2:H:440:PHE:CE2	2.29	0.67
2:D:62:GLU:HG2	2:D:297:ALA:HA	1.77	0.67
1:C:248:ARG:C	1:C:250:GLY:H	1.98	0.67
2:H:362:GLN:OE1	2:H:408:LEU:HD13	1.93	0.67
2:D:322:LEU:HD12	2:D:322:LEU:C	2.14	0.67
2:H:361:LEU:O	2:H:364:VAL:HG22	1.95	0.67
1:A:446:SER:HB2	1:A:449:GLN:NE2	2.10	0.67

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:226:TYR:CE1	1:C:231:GLY:HA2	2.29	0.67
2:H:64:LEU:HD21	2:H:77:VAL:CG2	2.25	0.67
2:F:132:ASP:HB3	2:F:157:MET:HE1	1.74	0.67
1:C:421:ASN:O	1:C:424:VAL:HG22	1.95	0.67
2:H:325:ASN:HD21	2:H:327:VAL:HG23	1.58	0.67
2:F:380:ALA:CB	2:F:394:LEU:HA	2.25	0.67
1:E:297:ILE:HB	1:E:368:ILE:HD11	1.75	0.67
1:G:518:ASN:HD22	1:G:534:LEU:H	1.43	0.67
2:B:400:ILE:H	2:B:400:ILE:HD12	1.58	0.67
1:A:461:LEU:O	1:A:465:LEU:HG	1.95	0.66
2:H:253:ASP:HB3	2:H:256:HIS:HB2	1.77	0.66
1:C:264:ASN:HD22	1:C:264:ASN:H	1.44	0.66
1:G:333:ILE:HG22	2:H:223:ARG:CZ	2.25	0.66
2:H:128:GLN:HE21	2:H:128:GLN:N	1.91	0.66
2:F:220:SER:C	2:F:222:PRO:HD3	2.15	0.66
2:B:361:LEU:HB2	2:B:407:ASN:O	1.95	0.66
2:D:380:ALA:HB1	2:D:394:LEU:HD13	1.78	0.66
2:F:398:THR:O	2:F:401:GLU:HG2	1.96	0.66
2:F:362:GLN:HE21	2:F:408:LEU:CD1	2.06	0.66
3:L:117:ILE:HD13	3:L:126:ILE:HG12	1.76	0.66
1:E:307:LEU:HB3	1:E:383:LEU:CD2	2.25	0.66
1:A:446:SER:HB2	1:A:449:GLN:HE21	1.59	0.66
1:E:422:GLU:HG3	1:E:530:ALA:HB3	1.78	0.66
2:D:411:THR:O	2:D:415:LEU:HG	1.96	0.66
2:H:348:GLN:HB3	2:H:349:LEU:HD12	1.77	0.66
2:D:412:LEU:HD22	2:D:417:LEU:HD12	1.77	0.66
2:D:192:ILE:CD1	2:D:200:ILE:HG13	2.26	0.66
2:D:418:VAL:HG22	2:D:421:GLN:NE2	2.10	0.66
1:G:112:LEU:C	1:G:114:ASP:H	1.96	0.66
1:A:125:VAL:HG21	1:A:505:VAL:HG13	1.77	0.66
2:F:241:GLN:HE21	2:F:246:GLY:H	1.42	0.66
1:A:430:ARG:NH1	1:A:430:ARG:HB3	2.10	0.66
2:D:231:TYR:C	2:D:233:ARG:H	1.98	0.66
1:G:61:ASP:OD2	1:G:85:ARG:HD2	1.95	0.66
2:H:412:LEU:HD22	2:H:440:PHE:CE2	2.31	0.66
1:G:307:LEU:HB3	1:G:383:LEU:CD2	2.25	0.66
2:B:407:ASN:ND2	2:B:415:LEU:HD22	2.10	0.65
3:K:156:ALA:HA	3:K:161:ILE:HD12	1.78	0.65
3:L:123:VAL:HA	3:L:126:ILE:HD12	1.78	0.65
1:A:61:ASP:OD2	1:A:85:ARG:HD2	1.95	0.65
2:D:360:LYS:HA	2:D:411:THR:HA	1.77	0.65
2:B:385:LEU:HD12	2:B:390:ARG:HD3	1.78	0.65

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:58:THR:HA	1:A:103:SER:O	1.96	0.65
2:B:400:ILE:HD12	2:B:400:ILE:N	2.12	0.65
2:B:427:ASP:HB3	2:B:429:THR:HG22	1.77	0.65
2:D:200:ILE:HG12	3:J:172:ARG:HH22	1.61	0.65
1:C:407:ILE:HG13	1:C:408:ASN:N	2.12	0.65
2:F:125:ASN:ND2	2:F:125:ASN:N	2.41	0.65
1:C:43:THR:HG21	1:C:73:ASN:OD1	1.97	0.65
2:H:128:GLN:N	2:H:128:GLN:NE2	2.42	0.65
3:L:107:THR:CG2	3:L:111:LYS:HB3	2.26	0.65
3:K:124:GLU:HB2	3:K:152:ASP:O	1.96	0.65
3:K:125:ARG:HD3	3:K:129:ARG:NH2	2.09	0.65
1:G:226:TYR:HB3	1:G:231:GLY:HA2	1.79	0.65
3:I:104:LYS:HD3	3:I:112:GLU:OE2	1.96	0.65
2:B:200:ILE:CD1	3:I:172:ARG:HH22	2.09	0.65
2:F:342:ASN:ND2	2:F:342:ASN:H	1.92	0.65
1:A:220:LYS:HE2	1:A:220:LYS:HA	1.78	0.65
3:I:101:MET:N	3:I:117:ILE:O	2.23	0.65
3:K:117:ILE:HG12	3:K:126:ILE:HG12	1.77	0.65
2:F:154:ILE:O	2:F:158:LEU:HD23	1.97	0.65
2:H:354:GLN:HE22	2:H:439:HIS:HB3	1.61	0.65
1:A:311:LEU:O	1:A:315:VAL:HG23	1.97	0.65
2:H:405:ARG:HB3	2:H:406:PRO:HD3	1.77	0.65
3:K:172:ARG:C	3:K:173:LEU:HG	2.17	0.65
1:A:518:ASN:HD21	1:A:534:LEU:HB2	1.60	0.64
2:B:380:ALA:HB1	2:B:394:LEU:CD1	2.28	0.64
2:D:343:CYS:O	2:D:347:SER:HB3	1.98	0.64
2:B:318:LEU:HG	2:B:319:ASN:O	1.96	0.64
1:A:150:LEU:HD23	1:A:165:ILE:HD12	1.79	0.64
1:A:248:ARG:HA	1:A:251:ILE:HD11	1.79	0.64
1:E:45:ILE:HG12	1:E:498:GLY:HA2	1.79	0.64
1:C:357:VAL:O	1:C:361:VAL:HG23	1.97	0.64
1:A:229:THR:O	1:A:230:ASN:HB2	1.98	0.64
1:A:252:LEU:HD22	1:A:253:LYS:N	2.12	0.64
2:B:102:PRO:HG2	2:B:105:GLU:HB3	1.78	0.64
2:B:425:VAL:HB	2:B:434:VAL:HG13	1.77	0.64
1:A:430:ARG:HB3	1:A:430:ARG:HH11	1.62	0.64
3:L:99:GLY:O	3:L:116:ASP:HB3	1.97	0.64
2:H:32:HIS:HD2	2:H:34:ASP:H	1.43	0.64
3:L:125:ARG:HD3	3:L:129:ARG:NH2	2.12	0.64
1:C:163:ARG:NH2	1:C:518:ASN:OD1	2.31	0.64
1:G:527:GLN:HB2	2:H:318:LEU:HD13	1.79	0.64
2:B:302:VAL:HG11	2:B:322:LEU:HD23	1.79	0.64

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:D:357:PRO:HG3	2:D:440:PHE:CG	2.33	0.64
2:D:29:PRO:HG2	2:D:30:PHE:H	1.63	0.64
1:E:211:HIS:HB2	1:E:334:ALA:HA	1.79	0.64
2:F:201:GLU:HG3	2:F:345:ALA:HB2	1.80	0.64
2:D:132:ASP:HB3	2:D:157:MET:HE1	1.80	0.64
2:F:412:LEU:HD23	2:F:417:LEU:HD22	1.78	0.64
2:B:233:ARG:HH11	2:B:233:ARG:HB3	1.63	0.64
2:F:16:ARG:HH22	2:F:116:PRO:HB2	1.63	0.64
3:J:161:ILE:CD1	3:J:167:LEU:HD21	2.26	0.64
2:B:188:ASN:OD1	3:I:173:LEU:HD12	1.98	0.64
2:F:371:SER:CB	2:F:374:LEU:CD1	2.58	0.63
2:B:200:ILE:CD1	3:I:172:ARG:NH2	2.61	0.63
1:C:84:ASN:HB3	1:C:87:GLU:CB	2.27	0.63
2:D:237:TRP:HB3	2:D:238:PRO:HD3	1.79	0.63
1:C:396:ARG:HG3	1:C:518:ASN:HD21	1.62	0.63
2:H:35:PHE:HD1	2:H:313:SER:O	1.80	0.63
2:H:344:PRO:HG3	2:H:374:LEU:HD13	1.80	0.63
3:L:107:THR:HG21	3:L:111:LYS:HB3	1.79	0.63
1:C:51:LEU:HB2	1:C:52:PRO:HD3	1.81	0.63
3:L:120:THR:HA	3:L:155:THR:HG1	1.63	0.63
2:H:132:ASP:OD1	2:H:132:ASP:N	2.30	0.63
1:G:215:ILE:H	1:G:332:MET:HE1	1.64	0.63
2:H:270:GLN:O	2:H:272:ASN:OD1	2.16	0.63
2:H:318:LEU:HD11	2:H:334:THR:HG21	1.80	0.63
2:B:430:THR:HG22	2:B:431:PRO:HD2	1.79	0.63
2:D:268:ALA:HB2	2:D:276:VAL:HG21	1.80	0.63
1:C:259:PRO:C	1:C:261:ASP:N	2.49	0.63
2:H:238:PRO:O	2:H:241:GLN:HG3	1.98	0.63
1:C:274:THR:C	1:C:276:LEU:H	2.01	0.63
1:C:517:ASN:ND2	1:C:518:ASN:N	2.10	0.63
2:F:384:THR:O	2:F:385:LEU:HD23	1.99	0.63
2:H:178:ILE:HD11	2:H:310:ILE:CD1	2.29	0.63
3:J:155:THR:HG23	3:J:158:ASP:OD1	1.99	0.63
2:H:187:GLY:CA	3:L:173:LEU:HD13	2.29	0.63
3:I:170:VAL:HG13	3:I:171:LEU:N	2.13	0.62
2:D:380:ALA:HB2	2:D:394:LEU:CD1	2.26	0.62
2:B:96:PRO:O	2:B:99:ILE:HG13	1.99	0.62
1:A:409:LYS:HE2	1:A:468:TYR:O	1.99	0.62
1:C:33:VAL:HG22	1:C:34:CYS:N	2.12	0.62
1:A:458:LYS:O	1:A:458:LYS:HD3	1.99	0.62
2:D:207:TYR:CE1	3:J:172:ARG:CD	2.76	0.62
2:H:404:THR:HG22	2:H:407:ASN:ND2	2.14	0.62

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:178:ASN:ND2	3:I:136:ILE:HG12	2.14	0.62
1:C:503:GLN:O	1:C:507:LYS:HG3	1.99	0.62
2:D:250:ASP:OD2	2:D:253:ASP:HB2	1.99	0.62
1:E:309:ARG:HB3	1:E:364:LEU:HD21	1.81	0.62
1:E:113:LEU:O	1:E:117:PRO:HG3	1.99	0.62
2:F:267:ARG:HG2	2:F:267:ARG:HH11	1.64	0.62
2:D:232:VAL:HG12	2:D:232:VAL:O	1.99	0.62
1:G:158:LEU:HD12	1:G:493:ILE:HD11	1.80	0.62
1:C:434:ARG:HD3	1:C:460:CYS:HB3	1.80	0.62
1:G:311:LEU:O	1:G:314:PHE:HB3	2.00	0.62
2:F:46:LEU:HD22	2:F:73:ARG:CZ	2.30	0.62
1:G:397:SER:OG	1:G:400:GLU:HG3	1.98	0.62
1:G:264:ASN:HB3	1:G:333:ILE:HG13	1.80	0.62
2:B:356:SER:CB	2:B:441:THR:OG1	2.47	0.62
3:I:107:THR:HG22	3:I:109:THR:N	2.09	0.62
1:C:61:ASP:OD1	1:C:63:ASN:HB2	1.99	0.62
1:E:36:ILE:HG21	1:E:131:LEU:HD21	1.82	0.62
1:E:243:PHE:O	1:E:247:ILE:HG13	1.99	0.62
1:G:333:ILE:N	2:H:223:ARG:NH1	2.46	0.62
2:B:442:SER:OG	2:B:442:SER:O	2.12	0.62
1:G:332:MET:HG2	1:G:339:TYR:HE1	1.65	0.62
2:F:183:GLU:HG3	2:F:289:ILE:HG21	1.80	0.62
2:F:342:ASN:HD22	2:F:342:ASN:N	1.85	0.62
2:H:58:GLY:N	2:H:91:GLN:HG2	2.14	0.62
1:E:253:LYS:N	1:E:259:PRO:HD2	2.14	0.62
2:H:237:TRP:HB3	2:H:238:PRO:HD3	1.81	0.62
2:H:391:THR:O	2:H:404:THR:HG21	1.99	0.61
3:K:155:THR:HG23	3:K:157:ALA:H	1.65	0.61
3:L:123:VAL:HA	3:L:126:ILE:CD1	2.29	0.61
1:A:112:LEU:H	1:A:112:LEU:CD2	2.13	0.61
1:C:294:CYS:O	1:C:309:ARG:HD3	2.00	0.61
2:H:233:ARG:NH1	2:H:234:MET:HB2	2.15	0.61
1:E:214:TRP:O	1:E:218:ILE:HG13	1.99	0.61
2:B:74:GLN:NE2	2:B:119:ASN:HD22	1.97	0.61
2:F:31:THR:HG23	2:F:32:HIS:O	2.00	0.61
2:F:232:VAL:HG12	2:F:232:VAL:O	1.99	0.61
2:H:237:TRP:CE3	2:H:242:PRO:HG2	2.35	0.61
1:G:199:TYR:CD2	1:G:216:VAL:HG11	2.35	0.61
1:G:185:LEU:CD1	1:G:275:ALA:HB1	2.30	0.61
2:F:192:ILE:HG22	2:F:194:PRO:HD3	1.82	0.61
1:A:234:PRO:HB2	1:A:276:LEU:CD1	2.30	0.61
2:H:234:MET:C	2:H:235:LEU:HD22	2.20	0.61

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:D:157:MET:HE3	2:D:157:MET:O	2.00	0.61
1:E:447:ASN:ND2	2:F:26:ARG:HH21	1.97	0.61
2:D:390:ARG:HG2	2:D:391:THR:N	2.14	0.61
1:G:241:GLU:O	1:G:244:ARG:HB2	1.99	0.61
1:G:50:VAL:HG13	1:G:100:VAL:HG21	1.82	0.61
2:B:415:LEU:HB2	2:B:417:LEU:CD2	2.30	0.61
2:F:316:ILE:H	2:F:316:ILE:CD1	2.09	0.61
2:F:139:HIS:C	2:F:140:ILE:HD12	2.20	0.61
3:J:104:LYS:HB2	3:J:166:VAL:HG12	1.81	0.61
1:A:450:VAL:O	1:A:454:ILE:HG13	1.99	0.61
2:B:316:ILE:CD1	2:B:316:ILE:H	2.08	0.61
1:A:366:GLN:C	1:A:368:ILE:H	2.04	0.61
1:G:61:ASP:HB3	1:G:86:ALA:HB2	1.83	0.61
1:C:58:THR:HA	1:C:103:SER:O	2.01	0.61
2:H:320:ASN:HD22	2:H:337:ALA:HB3	1.66	0.61
1:C:492:THR:HB	2:D:305:THR:OG1	2.01	0.61
1:G:72:ASN:HD22	1:G:72:ASN:H	1.48	0.61
2:F:362:GLN:CG	2:F:408:LEU:HD22	2.27	0.61
2:D:15:GLY:HA2	2:D:18:ASN:ND2	2.16	0.61
2:H:405:ARG:HH11	2:H:405:ARG:CB	2.13	0.61
2:B:368:LEU:HD21	2:B:436:PHE:CE2	2.35	0.61
2:H:92:PHE:H	2:H:92:PHE:HD2	1.48	0.61
2:H:58:GLY:H	2:H:91:GLN:CG	2.14	0.61
1:E:252:LEU:C	1:E:259:PRO:CD	2.69	0.61
1:A:423:ILE:O	1:A:426:TYR:HB3	2.01	0.61
2:D:192:ILE:HD11	2:D:200:ILE:HG13	1.83	0.61
3:L:150:MET:CE	3:L:167:LEU:HD13	2.31	0.61
1:A:281:ILE:HD11	1:A:315:VAL:HG11	1.82	0.61
2:B:131:ASN:HB3	2:D:131:ASN:HD22	1.64	0.61
2:D:185:PHE:HB3	2:D:326:ASP:HB2	1.83	0.61
1:A:36:ILE:HD12	1:A:60:ILE:CB	2.16	0.61
2:H:405:ARG:HB3	2:H:405:ARG:HH11	1.65	0.61
2:H:419:ASP:OD2	2:H:440:PHE:HB2	2.01	0.61
2:H:32:HIS:CG	2:H:33:PRO:HD2	2.36	0.61
1:G:192:LEU:O	1:G:195:HIS:HB3	2.01	0.61
1:G:214:TRP:HB2	1:G:268:ALA:HB2	1.82	0.60
1:G:336:SER:HB2	2:H:221:MET:HA	1.83	0.60
1:C:376:SER:OG	1:C:379:GLU:HG3	2.01	0.60
1:C:501:ALA:O	1:C:504:GLU:N	2.34	0.60
1:A:168:GLU:HG3	1:A:394:ARG:HE	1.66	0.60
1:C:430:ARG:NH1	1:C:430:ARG:HB3	2.16	0.60
2:H:217:THR:HG21	2:H:223:ARG:HE	1.65	0.60

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:233:ARG:HD2	2:B:285:VAL:HG13	1.82	0.60
2:B:353:ILE:HG23	2:B:355:PHE:CD1	2.35	0.60
3:J:101:MET:HB2	3:J:117:ILE:O	1.99	0.60
2:B:374:LEU:O	2:B:376:MET:HG3	2.01	0.60
2:B:49:CYS:HA	2:B:139:HIS:HD2	1.65	0.60
1:G:229:THR:HG21	1:G:232:ARG:HH21	1.65	0.60
2:B:218:ILE:CD1	2:B:230:GLU:HG2	2.29	0.60
2:H:349:LEU:HB3	2:H:350:PRO:HD2	1.83	0.60
2:H:178:ILE:N	2:H:178:ILE:HD12	2.16	0.60
1:G:306:ILE:HD13	1:G:365:LEU:CD2	2.31	0.60
1:C:438:GLN:HG2	1:C:438:GLN:O	2.01	0.60
2:F:178:ILE:HD11	2:F:310:ILE:HD12	1.82	0.60
1:C:6:LYS:O	1:C:9:LYS:HG2	2.01	0.60
1:A:407:ILE:HD13	1:A:408:ASN:N	2.16	0.60
2:H:390:ARG:NH2	2:H:407:ASN:HD21	1.99	0.60
2:B:412:LEU:O	2:B:417:LEU:HB2	2.02	0.60
1:E:489:GLU:H	2:F:19:HIS:HD2	1.46	0.60
2:F:223:ARG:HG3	2:F:223:ARG:HH11	1.67	0.60
1:G:306:ILE:HD13	1:G:365:LEU:HD23	1.83	0.60
1:E:240:LYS:HE2	1:E:276:LEU:HD12	1.84	0.60
3:L:124:GLU:HB2	3:L:152:ASP:O	2.00	0.60
2:D:236:GLN:HE22	2:D:263:LYS:HD2	1.67	0.60
1:G:307:LEU:HB3	1:G:383:LEU:HD22	1.84	0.60
1:C:284:SER:O	1:C:288:ILE:HG13	2.01	0.60
1:A:147:GLN:N	1:A:147:GLN:HE21	1.99	0.60
2:H:400:ILE:O	2:H:404:THR:N	2.35	0.60
2:F:376:MET:HE1	2:F:427:ASP:HB2	1.82	0.60
1:E:365:LEU:HD11	1:E:375:ILE:HD13	1.83	0.60
2:F:384:THR:C	2:F:385:LEU:HD23	2.22	0.60
2:D:199:CYS:O	2:D:202:CYS:HB2	2.02	0.60
1:A:437:LYS:HZ3	1:A:437:LYS:HA	1.65	0.60
2:D:325:ASN:ND2	2:D:327:VAL:HG22	2.17	0.60
2:H:158:LEU:CD1	2:H:177:LEU:HD22	2.32	0.60
3:J:144:ILE:HD13	3:J:149:GLN:HA	1.84	0.60
2:F:425:VAL:HB	2:F:434:VAL:HG13	1.82	0.60
1:G:265:PHE:O	1:G:269:ILE:HG13	2.02	0.60
2:H:237:TRP:CD2	2:H:242:PRO:HG2	2.37	0.60
2:F:361:LEU:HD12	2:F:407:ASN:O	2.02	0.60
2:F:131:ASN:HB3	2:H:131:ASN:ND2	2.14	0.60
2:D:158:LEU:HD22	2:D:175:VAL:HB	1.84	0.59
1:E:219:ALA:O	1:E:222:LEU:HB3	2.02	0.59
3:K:144:ILE:HD11	3:K:170:VAL:HG11	1.83	0.59

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:G:184:ARG:HD3	1:G:279:THR:OG1	2.02	0.59
1:C:126:VAL:O	1:C:150:LEU:HD12	2.02	0.59
1:C:360:HIS:O	1:C:364:LEU:HD23	2.02	0.59
2:H:131:ASN:O	2:H:133:THR:N	2.35	0.59
1:C:407:ILE:HG13	1:C:408:ASN:H	1.66	0.59
2:H:90:ARG:HG2	2:H:90:ARG:HH11	1.68	0.59
1:G:36:ILE:HB	1:G:128:ALA:HA	1.85	0.59
2:F:410:LYS:HB3	2:F:414:GLU:HG2	1.84	0.59
1:C:113:LEU:HD13	1:C:138:ARG:HG2	1.84	0.59
1:E:447:ASN:ND2	2:F:26:ARG:HE	2.00	0.59
1:E:39:THR:O	1:E:43:THR:HB	2.01	0.59
3:L:117:ILE:CD1	3:L:126:ILE:HG23	2.32	0.59
2:F:241:GLN:CG	2:F:245:GLU:HA	2.32	0.59
1:C:61:ASP:CB	1:C:86:ALA:HB2	2.33	0.59
2:B:158:LEU:HD12	2:B:177:LEU:HB2	1.85	0.59
1:C:277:ASN:O	1:C:277:ASN:ND2	2.36	0.59
3:I:103:ILE:CD1	3:I:117:ILE:HD12	2.33	0.59
1:G:422:GLU:OE1	1:G:422:GLU:N	2.33	0.59
1:A:47:LYS:HG3	1:A:75:PHE:HZ	1.67	0.59
1:C:113:LEU:O	1:C:117:PRO:HG3	2.03	0.59
1:G:492:THR:HB	2:H:305:THR:OG1	2.02	0.59
2:D:200:ILE:HG12	3:J:172:ARG:NH2	2.17	0.59
3:L:117:ILE:HG23	3:L:121:ASP:HB2	1.82	0.59
2:B:385:LEU:CD1	2:B:390:ARG:HD3	2.32	0.59
1:C:158:LEU:HA	1:C:493:ILE:HG13	1.82	0.59
2:H:335:PHE:HE2	3:L:170:VAL:HG11	1.67	0.59
2:F:15:GLY:HA2	2:F:18:ASN:HD21	1.67	0.59
1:E:226:TYR:CD1	1:E:231:GLY:CA	2.78	0.59
1:C:42:GLY:CA	1:C:129:THR:HG21	2.32	0.59
2:F:411:THR:O	2:F:414:GLU:HB3	2.02	0.59
2:F:219:ALA:HA	2:F:267:ARG:HH22	1.67	0.59
2:B:286:LYS:HB2	2:B:288:ILE:HG13	1.85	0.59
1:E:13:TYR:O	1:E:17:LEU:HG	2.03	0.59
3:K:149:GLN:HG3	3:K:149:GLN:O	2.03	0.59
3:I:117:ILE:HG22	3:I:118:GLU:N	2.17	0.59
2:D:237:TRP:CE3	2:D:242:PRO:HG2	2.38	0.59
2:F:249:LEU:CD1	2:F:260:ILE:HD11	2.33	0.59
1:G:264:ASN:HA	1:G:333:ILE:HD11	1.85	0.58
2:H:224:LEU:H	2:H:227:HIS:CE1	2.20	0.58
2:H:368:LEU:HD13	2:H:425:VAL:HG11	1.84	0.58
1:C:441:ARG:HB3	2:H:270:GLN:OE1	2.03	0.58
1:E:248:ARG:O	1:E:251:ILE:HG12	2.00	0.58

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:340:ILE:HG21	2:F:272:ASN:ND2	2.18	0.58
1:C:173:GLU:HA	1:C:386:ASN:ND2	2.18	0.58
1:C:262:GLU:HG2	1:C:265:PHE:CD1	2.38	0.58
2:H:356:SER:C	2:H:358:SER:N	2.57	0.58
2:D:124:PHE:C	2:D:125:ASN:HD22	2.05	0.58
1:E:282:PRO:HB2	1:E:285:ILE:HD13	1.84	0.58
2:D:128:GLN:NE2	2:D:128:GLN:H	2.01	0.58
2:F:222:PRO:O	2:F:273:ILE:HD11	2.01	0.58
2:B:393:TYR:CE2	2:B:408:LEU:HD21	2.39	0.58
2:D:384:THR:HA	2:D:389:ASN:HB3	1.83	0.58
1:E:236:THR:HG22	1:E:238:LYS:H	1.67	0.58
1:E:507:LYS:HG2	1:E:513:PHE:HB2	1.84	0.58
1:E:200:ASP:C	1:E:201:LEU:HD23	2.24	0.58
2:B:380:ALA:HB1	2:B:394:LEU:HD12	1.84	0.58
1:A:307:LEU:HB3	1:A:383:LEU:CD2	2.32	0.58
2:F:225:PRO:O	2:F:229:ILE:HG13	2.03	0.58
2:H:35:PHE:CD1	2:H:313:SER:O	2.57	0.58
3:L:106:LYS:HG2	3:L:112:GLU:HG2	1.84	0.58
2:D:430:THR:OG1	2:D:432:GLN:HG2	2.02	0.58
1:A:518:ASN:OD1	1:A:519:THR:N	2.35	0.58
2:H:249:LEU:CD1	2:H:260:ILE:HD11	2.34	0.58
3:L:126:ILE:O	3:L:130:VAL:HG23	2.04	0.58
1:C:229:THR:HG22	1:C:232:ARG:HD2	1.85	0.58
2:H:177:LEU:O	2:H:177:LEU:HG	2.03	0.58
3:J:105:VAL:CG1	3:J:115:ILE:HD11	2.34	0.58
2:F:366:ASP:HB3	2:F:370:ASN:HD22	1.68	0.58
2:F:221:MET:N	2:F:222:PRO:HD3	2.19	0.58
1:E:41:THR:O	1:E:45:ILE:HG13	2.04	0.58
3:L:103:ILE:O	3:L:103:ILE:HD12	2.04	0.58
1:A:19:LEU:HD21	2:B:290:PRO:HB2	1.86	0.58
2:B:15:GLY:HA2	2:B:18:ASN:OD1	2.03	0.58
2:D:320:ASN:HB2	2:D:336:GLU:HA	1.86	0.58
1:A:49:LEU:O	1:A:52:PRO:HG2	2.03	0.58
2:B:196:MET:O	2:B:339:ARG:HD2	2.03	0.58
2:F:325:ASN:HD22	2:F:326:ASP:N	2.01	0.58
2:H:238:PRO:O	2:H:241:GLN:CG	2.51	0.58
1:C:236:THR:HB	1:C:239:GLU:HG3	1.86	0.58
1:C:130:GLN:OE1	1:C:155:THR:CB	2.48	0.58
1:C:418:ASN:C	1:C:418:ASN:HD22	2.07	0.58
2:D:13:TRP:HZ3	2:D:116:PRO:HG2	1.67	0.58
1:A:518:ASN:HD22	1:A:534:LEU:H	1.29	0.58
1:E:184:ARG:HD3	1:E:279:THR:OG1	2.04	0.58

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:H:84:ASP:H	2:H:87:ASN:ND2	2.02	0.58
1:G:226:TYR:HE2	1:G:233:ILE:HA	1.67	0.58
2:B:153:TRP:CE2	2:B:431:PRO:HG3	2.39	0.58
1:A:436:HIS:HA	1:A:441:ARG:O	2.04	0.58
3:K:107:THR:HG22	3:K:109:THR:H	1.69	0.58
2:F:239:LYS:O	2:F:240:GLU:HG2	2.04	0.57
2:D:193:LEU:H	2:D:197:THR:HB	1.68	0.57
2:F:410:LYS:CB	2:F:414:GLU:HG2	2.33	0.57
2:B:325:ASN:ND2	2:B:327:VAL:HG22	2.19	0.57
3:K:115:ILE:HG23	3:K:117:ILE:HD13	1.85	0.57
2:H:32:HIS:CD2	2:H:34:ASP:H	2.22	0.57
1:G:84:ASN:OD1	1:G:106:GLU:HG3	2.04	0.57
1:G:488:ALA:C	1:G:490:PRO:HD3	2.25	0.57
1:A:418:ASN:C	1:A:418:ASN:HD22	2.08	0.57
1:C:72:ASN:OD1	1:C:489:GLU:HG2	2.04	0.57
1:C:88:ALA:O	1:C:91:GLU:HB2	2.05	0.57
3:L:102:LEU:HD11	3:L:114:GLU:HB3	1.85	0.57
1:G:65:VAL:HB	1:G:81:ILE:HA	1.86	0.57
2:H:318:LEU:O	2:H:319:ASN:O	2.22	0.57
2:H:270:GLN:C	2:H:272:ASN:OD1	2.43	0.57
2:F:371:SER:O	2:F:374:LEU:CD1	2.52	0.57
2:B:327:VAL:HG23	2:B:328:ASP:N	2.17	0.57
2:F:236:GLN:HE22	2:F:263:LYS:HB3	1.70	0.57
2:F:187:GLY:CA	3:K:173:LEU:HD13	2.35	0.57
2:F:152:ARG:CZ	2:F:429:THR:HG21	2.35	0.57
1:A:224:GLN:HE21	1:A:246:LEU:HD11	1.70	0.57
2:F:187:GLY:HA2	3:K:173:LEU:HD13	1.87	0.57
1:G:20:TRP:O	1:G:24:GLY:HA3	2.04	0.57
2:F:381:ILE:H	2:F:381:ILE:HD12	1.70	0.57
1:C:438:GLN:O	1:C:439:GLN:HG2	2.04	0.57
2:F:178:ILE:HD11	2:F:310:ILE:CD1	2.34	0.57
1:A:484:ARG:HH12	2:B:31:THR:HG22	1.68	0.57
3:K:118:GLU:HG3	3:K:120:THR:HG22	1.86	0.57
2:B:353:ILE:HD11	2:B:367:TYR:CE2	2.40	0.57
2:D:335:PHE:HE2	3:J:170:VAL:HG21	1.70	0.57
1:G:33:VAL:HG22	1:G:54:ILE:HG12	1.86	0.57
2:H:359:ALA:O	2:H:412:LEU:HD12	2.05	0.57
1:C:97:ASN:HB3	1:C:100:VAL:HG23	1.87	0.57
3:L:102:LEU:HD12	3:L:115:ILE:O	2.04	0.57
2:H:201:GLU:HB3	2:H:343:CYS:SG	2.44	0.57
2:D:316:ILE:H	2:D:316:ILE:HD12	1.70	0.57
1:C:163:ARG:CD	1:C:518:ASN:O	2.52	0.57

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:F:411:THR:N	2:F:414:GLU:HB3	2.16	0.57
1:G:6:LYS:C	1:G:8:LEU:H	2.08	0.57
1:C:418:ASN:HD22	1:C:419:PRO:N	2.02	0.57
2:B:318:LEU:HD11	2:B:334:THR:HG21	1.87	0.57
1:G:275:ALA:O	1:G:278:THR:HG23	2.05	0.57
1:C:189:PHE:CE1	1:C:192:LEU:HB2	2.39	0.57
1:G:518:ASN:ND2	1:G:534:LEU:H	2.03	0.56
2:B:385:LEU:HD12	2:B:390:ARG:CD	2.34	0.56
2:F:385:LEU:HD11	2:F:392:LEU:HD21	1.86	0.56
1:A:156:TYR:CE1	1:A:487:ALA:HA	2.40	0.56
1:A:259:PRO:C	1:A:261:ASP:N	2.59	0.56
2:H:376:MET:HB3	2:H:427:ASP:OD1	2.04	0.56
2:D:252:ASP:O	2:D:254:PRO:HD3	2.05	0.56
2:F:222:PRO:HG2	2:F:267:ARG:NH2	2.19	0.56
1:A:527:GLN:HG3	2:B:306:GLU:OE1	2.05	0.56
2:F:236:GLN:O	2:F:240:GLU:HG3	2.05	0.56
1:G:215:ILE:O	1:G:218:ILE:HG22	2.05	0.56
1:A:441:ARG:HH11	1:A:441:ARG:HG2	1.71	0.56
1:G:434:ARG:NH1	1:G:464:PHE:HA	2.19	0.56
2:H:422:GLU:C	2:H:423:LEU:HD12	2.25	0.56
2:B:427:ASP:OD2	2:B:428:VAL:N	2.38	0.56
1:A:484:ARG:NH1	2:B:31:THR:HG22	2.21	0.56
2:F:96:PRO:O	2:F:99:ILE:HG13	2.05	0.56
2:D:216:ALA:O	2:D:220:SER:HB2	2.06	0.56
2:D:352:ASN:HB3	2:D:439:HIS:CD2	2.39	0.56
2:H:383:ALA:HB3	2:H:392:LEU:HD21	1.87	0.56
1:E:215:ILE:HG13	1:E:332:MET:SD	2.46	0.56
1:C:59:ILE:HG22	1:C:60:ILE:H	1.69	0.56
1:A:371:ALA:C	1:A:373:GLU:H	2.08	0.56
1:C:307:LEU:HD22	1:C:383:LEU:HD22	1.88	0.56
1:C:430:ARG:HH11	1:C:430:ARG:HB3	1.70	0.56
1:G:331:ASP:OD1	2:H:223:ARG:HB3	2.04	0.56
2:F:164:TYR:CZ	2:F:348:GLN:HG3	2.41	0.56
2:H:46:LEU:HD23	2:H:71:GLY:O	2.05	0.56
1:G:299:LYS:HA	1:G:368:ILE:CG2	2.36	0.56
2:H:229:ILE:HD13	2:H:281:THR:HA	1.88	0.56
1:E:193:ARG:HG3	1:E:193:ARG:HH11	1.70	0.56
2:D:148:ILE:HG12	2:D:207:TYR:CE2	2.40	0.56
3:L:155:THR:HG23	3:L:158:ASP:H	1.70	0.56
1:E:218:ILE:O	1:E:222:LEU:HB2	2.06	0.56
1:E:264:ASN:HD22	1:E:265:PHE:N	2.04	0.56
2:F:147:SER:HB2	2:F:149:ILE:HD13	1.87	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:H:356:SER:C	2:H:358:SER:H	2.08	0.56
1:C:435:PHE:O	1:C:439:GLN:HB2	2.05	0.56
1:A:147:GLN:CA	1:A:147:GLN:HE21	2.18	0.56
3:K:107:THR:HG22	3:K:109:THR:N	2.20	0.56
2:H:324:PHE:HD1	2:H:332:THR:HG22	1.70	0.56
1:G:407:ILE:HG13	1:G:408:ASN:N	2.21	0.56
2:D:158:LEU:HD13	2:D:177:LEU:HB2	1.86	0.56
3:J:116:ASP:O	3:J:117:ILE:HD13	2.06	0.56
1:C:297:ILE:CD1	1:C:297:ILE:H	2.16	0.56
2:B:425:VAL:HB	2:B:434:VAL:CG1	2.35	0.56
2:F:16:ARG:NH2	2:F:116:PRO:HB2	2.21	0.56
1:E:518:ASN:HB2	1:E:533:GLN:HA	1.87	0.56
2:F:312:THR:O	2:F:313:SER:HB2	2.05	0.56
1:E:299:LYS:HE2	1:E:368:ILE:O	2.06	0.55
1:C:233:ILE:N	1:C:233:ILE:HD12	2.21	0.55
2:D:57:GLY:C	2:D:91:GLN:HG2	2.26	0.55
2:F:256:HIS:O	2:F:260:ILE:HG13	2.06	0.55
2:B:198:ALA:H	2:B:320:ASN:HD21	1.54	0.55
2:B:252:ASP:O	2:B:254:PRO:HD3	2.06	0.55
1:C:426:TYR:O	1:C:430:ARG:HG2	2.06	0.55
2:F:381:ILE:HG22	2:F:382:THR:N	2.20	0.55
1:A:66:SER:O	1:A:81:ILE:HG23	2.06	0.55
1:A:264:ASN:H	1:A:264:ASN:HD22	1.53	0.55
1:E:447:ASN:HD22	2:F:26:ARG:NH2	2.04	0.55
2:H:52:LEU:HD22	2:H:138:PHE:CE1	2.42	0.55
1:E:158:LEU:HD12	1:E:158:LEU:N	2.21	0.55
1:G:7:LEU:O	1:G:7:LEU:HD23	2.06	0.55
2:B:356:SER:N	2:B:441:THR:OG1	2.38	0.55
2:F:412:LEU:HD13	2:F:440:PHE:CE2	2.35	0.55
2:F:348:GLN:HG2	2:F:348:GLN:O	2.06	0.55
2:F:351:GLN:H	2:F:436:PHE:HA	1.70	0.55
1:C:481:GLU:OE2	2:D:29:PRO:HD2	2.05	0.55
1:C:61:ASP:HB2	1:C:86:ALA:HB2	1.88	0.55
2:B:197:THR:CG2	2:B:198:ALA:N	2.70	0.55
1:C:259:PRO:CB	1:C:262:GLU:OE1	2.47	0.55
1:E:212:THR:O	1:E:213:PRO:C	2.44	0.55
1:E:234:PRO:HG2	1:E:243:PHE:CD1	2.42	0.55
2:F:15:GLY:HA2	2:F:18:ASN:ND2	2.21	0.55
1:A:19:LEU:HD21	2:B:290:PRO:CB	2.36	0.55
1:C:119:PHE:O	1:C:122:ARG:HG2	2.07	0.55
1:G:18:ARG:HH12	2:H:279:ARG:HG3	1.71	0.55
1:E:84:ASN:OD1	1:E:106:GLU:HG2	2.06	0.55

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:F:320:ASN:ND2	2:F:337:ALA:O	2.40	0.55
2:F:234:MET:O	2:F:235:LEU:HD23	2.06	0.55
1:E:357:VAL:O	1:E:361:VAL:HG23	2.06	0.55
2:D:230:GLU:HG3	2:D:230:GLU:O	2.07	0.55
1:A:74:PHE:CD1	2:B:65:LYS:HG3	2.42	0.55
1:G:267:GLU:O	1:G:270:LYS:HG2	2.06	0.55
1:G:180:LEU:H	1:G:180:LEU:CD1	2.17	0.55
3:L:113:ILE:HG13	3:L:113:ILE:O	2.06	0.55
1:A:112:LEU:N	1:A:112:LEU:HD22	2.19	0.55
1:G:164:ILE:HD11	1:G:166:ILE:HD11	1.89	0.55
1:G:299:LYS:HG2	1:G:368:ILE:HG22	1.89	0.55
1:A:14:ASP:O	1:A:18:ARG:HG3	2.07	0.55
2:F:425:VAL:HB	2:F:434:VAL:CG1	2.37	0.55
3:J:170:VAL:HG22	3:J:171:LEU:N	2.18	0.55
1:E:396:ARG:NH2	1:E:400:GLU:HB3	2.22	0.55
1:A:193:ARG:O	1:A:197:GLN:HG3	2.07	0.55
2:B:435:LEU:C	2:B:436:PHE:HD1	2.10	0.55
2:D:240:GLU:O	2:D:240:GLU:HG3	2.07	0.55
2:H:217:THR:HG21	2:H:223:ARG:NH2	2.19	0.55
3:J:123:VAL:CB	3:J:152:ASP:HA	2.22	0.55
2:B:353:ILE:HD11	2:B:367:TYR:HE2	1.71	0.55
1:A:47:LYS:HG3	1:A:75:PHE:CZ	2.42	0.55
1:G:279:THR:HG22	1:G:322:ASN:ND2	2.22	0.55
1:E:173:GLU:O	1:E:512:GLN:O	2.25	0.55
2:H:228:CYS:SG	2:H:268:ALA:HB2	2.47	0.54
2:D:74:GLN:HE22	2:D:119:ASN:HD22	1.52	0.54
1:A:418:ASN:HD22	1:A:419:PRO:N	2.05	0.54
1:C:412:ILE:O	1:C:416:MET:HB2	2.07	0.54
2:F:236:GLN:HE22	2:F:263:LYS:HD2	1.72	0.54
1:G:248:ARG:O	1:G:251:ILE:CG1	2.55	0.54
1:G:87:GLU:O	1:G:91:GLU:HG3	2.07	0.54
1:G:516:PHE:HD2	1:G:516:PHE:H	1.54	0.54
1:E:252:LEU:O	1:E:259:PRO:N	2.40	0.54
2:F:87:ASN:HD21	2:F:103:LYS:HE2	1.70	0.54
1:E:215:ILE:HD12	1:E:332:MET:HE1	1.89	0.54
2:H:200:ILE:HD13	3:L:172:ARG:HH22	1.72	0.54
2:H:412:LEU:HD23	2:H:438:LEU:HD21	1.88	0.54
2:D:229:ILE:HD13	2:D:281:THR:HA	1.89	0.54
1:C:243:PHE:O	1:C:246:LEU:HB3	2.08	0.54
1:C:154:ARG:HB3	1:C:161:TYR:HB3	1.88	0.54
3:K:123:VAL:CB	3:K:152:ASP:HA	2.21	0.54
1:G:518:ASN:HB3	1:G:532:PHE:O	2.07	0.54

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:232:ARG:C	1:C:233:ILE:HD12	2.28	0.54
2:B:182:THR:HG22	3:I:176:GLY:HA2	1.90	0.54
1:G:37:ASN:HB3	1:G:39:THR:HG23	1.89	0.54
2:F:318:LEU:HD11	2:F:334:THR:CG2	2.37	0.54
1:E:332:MET:O	2:F:223:ARG:NE	2.40	0.54
1:E:171:VAL:CG1	1:E:391:ARG:HB2	2.37	0.54
1:A:461:LEU:HD21	1:A:474:VAL:HG11	1.89	0.54
1:G:215:ILE:H	1:G:332:MET:CE	2.20	0.54
1:C:34:CYS:HB2	1:C:123:PHE:CD2	2.42	0.54
1:A:178:ASN:CG	3:I:136:ILE:HG12	2.28	0.54
1:A:45:ILE:HG12	1:A:498:GLY:HA2	1.90	0.54
1:C:248:ARG:O	1:C:250:GLY:N	2.41	0.54
1:A:78:ARG:HA	1:A:81:ILE:HG13	1.90	0.54
1:G:434:ARG:O	1:G:438:GLN:HG2	2.08	0.54
2:H:229:ILE:HD12	2:H:284:VAL:HG21	1.88	0.54
2:D:32:HIS:ND1	2:D:33:PRO:HD2	2.23	0.54
3:L:143:LEU:C	3:L:144:ILE:HD12	2.28	0.54
1:E:302:PRO:HG2	1:E:305:TRP:HD1	1.72	0.54
2:F:415:LEU:HB2	2:F:417:LEU:CD1	2.38	0.54
1:E:260:GLU:O	1:E:261:ASP:OD1	2.25	0.54
1:A:96:LEU:HD23	2:B:95:ARG:HH21	1.73	0.54
1:E:220:LYS:C	1:E:222:LEU:H	2.11	0.54
2:H:38:SER:C	2:H:40:GLU:H	2.10	0.54
2:H:132:ASP:HB3	2:H:157:MET:HE1	1.90	0.54
2:F:54:ILE:CG2	2:F:145:LEU:HD21	2.37	0.54
1:G:307:LEU:HB3	1:G:383:LEU:HD21	1.89	0.54
2:H:259:TRP:CE2	2:H:263:LYS:HG3	2.42	0.54
1:A:396:ARG:HD3	1:A:534:LEU:O	2.08	0.54
2:H:231:TYR:O	2:H:233:ARG:N	2.41	0.54
1:C:274:THR:C	1:C:276:LEU:N	2.61	0.54
2:D:357:PRO:HG3	2:D:440:PHE:CD2	2.43	0.54
2:B:226:GLU:HG3	2:B:280:LEU:HD11	1.90	0.54
1:G:269:ILE:O	1:G:269:ILE:HG22	2.08	0.54
2:B:359:ALA:O	2:B:412:LEU:HD23	2.08	0.54
1:E:163:ARG:NH1	1:E:518:ASN:HD21	2.05	0.54
1:A:12:LYS:HG3	1:A:13:TYR:CD2	2.43	0.54
1:A:214:TRP:CZ3	1:A:332:MET:HB3	2.43	0.54
3:L:117:ILE:CG2	3:L:121:ASP:HB2	2.38	0.53
1:A:78:ARG:O	1:A:81:ILE:HG13	2.08	0.53
2:B:62:GLU:HG2	2:B:300:ALA:HB3	1.89	0.53
2:D:362:GLN:NE2	2:D:365:LEU:HD23	2.23	0.53
1:C:130:GLN:CD	1:C:155:THR:H	2.12	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:K:119:PRO:O	3:K:157:ALA:HB2	2.08	0.53
2:B:419:ASP:HA	2:B:440:PHE:HE1	1.72	0.53
2:B:241:GLN:OE1	2:B:246:GLY:N	2.40	0.53
2:D:415:LEU:CD1	2:D:417:LEU:HG	2.38	0.53
2:D:415:LEU:HD11	2:D:417:LEU:HG	1.90	0.53
2:B:351:GLN:HB3	2:B:436:PHE:CG	2.43	0.53
2:B:46:LEU:HD22	2:B:73:ARG:CZ	2.38	0.53
2:D:183:GLU:HB3	3:J:173:LEU:HD22	1.89	0.53
2:H:224:LEU:N	2:H:227:HIS:ND1	2.55	0.53
3:J:118:GLU:HG3	3:J:120:THR:CG2	2.37	0.53
2:H:207:TYR:OH	3:L:172:ARG:NE	2.40	0.53
2:F:125:ASN:HB2	2:F:129:ASP:OD2	2.07	0.53
2:H:344:PRO:HG3	2:H:374:LEU:HD22	1.90	0.53
1:E:158:LEU:HD11	2:F:23:PHE:CE2	2.43	0.53
1:A:214:TRP:CD1	1:A:214:TRP:C	2.82	0.53
1:A:162:MET:HB3	1:A:520:TYR:HB3	1.90	0.53
2:B:74:GLN:HE22	2:B:119:ASN:ND2	2.06	0.53
2:F:322:LEU:HD13	2:F:323:VAL:N	2.23	0.53
1:A:282:PRO:HB2	1:A:285:ILE:HD13	1.91	0.53
2:H:402:GLU:HA	2:H:405:ARG:HH12	1.73	0.53
2:F:356:SER:C	2:F:358:SER:H	2.12	0.53
1:E:246:LEU:O	1:E:246:LEU:HD23	2.08	0.53
1:G:333:ILE:CA	2:H:223:ARG:NH1	2.71	0.53
2:F:391:THR:HG21	2:F:394:LEU:HB2	1.90	0.53
1:C:297:ILE:HD13	1:C:297:ILE:N	2.19	0.53
1:C:59:ILE:O	1:C:60:ILE:HG13	2.08	0.53
2:B:49:CYS:HA	2:B:139:HIS:CD2	2.43	0.53
1:A:437:LYS:HA	1:A:437:LYS:NZ	2.24	0.53
2:B:195:GLY:O	2:B:339:ARG:NH1	2.41	0.53
1:E:171:VAL:HG12	1:E:391:ARG:HB2	1.89	0.53
2:B:361:LEU:O	2:B:364:VAL:HG23	2.09	0.53
1:E:68:GLU:HG2	2:F:18:ASN:HD22	1.73	0.53
2:B:251:GLY:O	2:B:257:ILE:HD11	2.09	0.53
1:C:341:LYS:O	1:C:345:VAL:HG23	2.08	0.53
1:E:447:ASN:HD22	2:F:26:ARG:HH21	1.55	0.53
1:C:120:PHE:C	1:C:122:ARG:H	2.12	0.53
1:E:527:GLN:HB2	2:F:318:LEU:HD13	1.89	0.53
1:A:172:ILE:HA	1:A:390:LEU:HD23	1.91	0.53
2:B:190:GLN:HG3	2:B:191:VAL:N	2.23	0.53
1:A:72:ASN:C	1:A:72:ASN:ND2	2.62	0.53
1:C:36:ILE:CB	1:C:128:ALA:HA	2.36	0.53
2:B:241:GLN:O	2:B:242:PRO:C	2.47	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:130:GLN:OE1	1:A:154:ARG:HA	2.09	0.53
2:D:398:THR:O	2:D:402:GLU:HG3	2.09	0.53
2:B:207:TYR:O	3:I:142:ARG:NH1	2.42	0.53
3:L:126:ILE:HD11	3:L:156:ALA:HB2	1.91	0.53
1:C:59:ILE:HG22	1:C:60:ILE:N	2.23	0.53
1:E:113:LEU:HB3	1:E:138:ARG:NH2	2.23	0.53
1:A:72:ASN:HD22	1:A:72:ASN:C	2.11	0.53
2:B:178:ILE:HD11	2:B:310:ILE:CD1	2.39	0.53
1:C:132:PRO:HD2	1:C:135:THR:HB	1.90	0.53
1:G:404:LEU:HD21	1:G:467:GLU:HG2	1.89	0.53
1:A:518:ASN:ND2	1:A:534:LEU:CA	2.72	0.52
2:H:380:ALA:HB1	2:H:394:LEU:CD1	2.35	0.52
2:H:382:THR:OG1	2:H:424:ALA:HB3	2.08	0.52
2:D:405:ARG:HH11	2:D:405:ARG:HB3	1.73	0.52
2:H:234:MET:O	2:H:234:MET:HG2	2.08	0.52
2:H:243:PHE:O	2:H:247:VAL:HB	2.09	0.52
2:B:360:LYS:O	2:B:363:GLU:HB2	2.09	0.52
2:B:407:ASN:HB3	2:B:415:LEU:HD11	1.91	0.52
2:H:348:GLN:CA	2:H:348:GLN:HE21	2.21	0.52
1:G:226:TYR:CZ	1:G:233:ILE:HG22	2.44	0.52
2:B:422:GLU:HG3	2:B:436:PHE:O	2.10	0.52
2:H:338:GLU:CD	3:L:148:LYS:HE3	2.29	0.52
2:B:51:VAL:HG11	2:B:67:LEU:HD13	1.89	0.52
3:I:102:LEU:O	3:I:102:LEU:HD23	2.09	0.52
3:K:123:VAL:HB	3:K:152:ASP:CA	2.21	0.52
1:G:518:ASN:HB2	1:G:533:GLN:HA	1.89	0.52
2:H:412:LEU:HD22	2:H:440:PHE:HE2	1.73	0.52
1:G:299:LYS:HE2	1:G:368:ILE:O	2.09	0.52
1:E:194:GLU:O	1:E:196:PHE:N	2.42	0.52
2:D:246:GLY:O	2:D:248:PRO:HD3	2.09	0.52
2:F:195:GLY:O	2:F:196:MET:HG2	2.10	0.52
1:A:248:ARG:C	1:A:250:GLY:H	2.12	0.52
1:G:244:ARG:HH11	1:G:244:ARG:HG3	1.74	0.52
2:H:343:CYS:O	2:H:347:SER:HB3	2.10	0.52
3:K:102:LEU:O	3:K:163:GLY:O	2.28	0.52
3:J:105:VAL:HG11	3:J:115:ILE:HD11	1.90	0.52
2:H:102:PRO:HG2	2:H:105:GLU:HB2	1.91	0.52
1:A:119:PHE:O	1:A:122:ARG:HG2	2.10	0.52
1:E:307:LEU:HD21	1:E:375:ILE:HG21	1.91	0.52
2:H:231:TYR:C	2:H:233:ARG:H	2.14	0.52
2:B:434:VAL:HG22	2:B:436:PHE:HE1	1.75	0.52
2:B:83:ILE:HD13	2:B:94:PHE:HB3	1.91	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:F:183:GLU:HG3	2:F:289:ILE:CG2	2.39	0.52
1:E:376:SER:OG	1:E:379:GLU:HG3	2.09	0.52
1:G:25:GLN:O	1:G:29:GLU:HG3	2.10	0.52
1:G:377:GLU:OE2	1:G:381:LYS:HE2	2.09	0.52
2:H:381:ILE:N	2:H:381:ILE:HD12	2.25	0.52
2:F:324:PHE:HD1	2:F:332:THR:HG22	1.74	0.52
2:H:385:LEU:HD12	2:H:390:ARG:HD3	1.92	0.52
3:I:115:ILE:HD11	3:I:126:ILE:HG23	1.90	0.52
2:B:419:ASP:CA	2:B:440:PHE:HE1	2.23	0.52
1:G:333:ILE:HA	2:H:223:ARG:NH1	2.24	0.52
1:A:516:PHE:CD1	1:A:518:ASN:O	2.63	0.52
2:D:318:LEU:HD11	2:D:334:THR:HG21	1.89	0.52
2:D:231:TYR:CD1	2:D:235:LEU:HD12	2.45	0.52
1:C:66:SER:H	1:C:69:ASP:HB2	1.74	0.52
2:B:232:VAL:CG1	2:B:260:ILE:HG23	2.40	0.52
1:A:309:ARG:HG3	1:A:364:LEU:HD21	1.92	0.52
2:D:152:ARG:HB3	2:D:429:THR:CG2	2.39	0.52
2:D:152:ARG:CZ	2:D:429:THR:HG21	2.40	0.52
2:D:327:VAL:HG23	2:D:328:ASP:H	1.75	0.52
1:E:309:ARG:NH1	1:E:309:ARG:HG3	2.24	0.52
3:K:142:ARG:HB2	3:K:170:VAL:HG12	1.92	0.52
2:D:362:GLN:HG2	2:D:408:LEU:HB3	1.92	0.52
1:E:450:VAL:O	1:E:454:ILE:HG13	2.10	0.52
1:A:128:ALA:HB1	1:A:131:LEU:CD1	2.40	0.52
1:C:521:ILE:HD12	1:C:532:PHE:HE1	1.75	0.52
2:H:148:ILE:HD11	3:L:174:ARG:HD3	1.91	0.51
1:C:226:TYR:CZ	1:C:231:GLY:HA2	2.45	0.51
3:K:107:THR:C	3:K:109:THR:H	2.13	0.51
1:C:462:THR:O	1:C:466:GLN:HB2	2.09	0.51
1:G:331:ASP:OD1	2:H:224:LEU:HD21	2.10	0.51
2:B:13:TRP:CH2	2:B:116:PRO:HG2	2.44	0.51
1:C:49:LEU:C	1:C:52:PRO:HD2	2.30	0.51
1:C:33:VAL:CG2	1:C:34:CYS:N	2.74	0.51
1:E:121:CYS:HA	1:E:148:ILE:HD11	1.92	0.51
1:G:298:THR:OG1	1:G:300:GLN:HG2	2.09	0.51
1:E:317:LYS:HD3	1:E:318:GLU:OE1	2.10	0.51
1:A:516:PHE:CE1	1:A:518:ASN:O	2.64	0.51
2:F:228:CYS:SG	2:F:273:ILE:HD12	2.50	0.51
2:H:350:PRO:HB2	2:H:437:LYS:HG3	1.91	0.51
2:B:132:ASP:O	2:B:133:THR:C	2.49	0.51
1:A:35:LEU:O	1:A:36:ILE:HD13	2.11	0.51
2:F:64:LEU:HD21	2:F:77:VAL:CG2	2.41	0.51

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:397:SER:HG	1:C:400:GLU:HG3	1.74	0.51
2:B:318:LEU:HD11	2:B:334:THR:HG23	1.93	0.51
2:B:64:LEU:CB	2:B:111:LEU:HD11	2.41	0.51
1:A:448:TYR:CE1	1:A:449:GLN:HG3	2.45	0.51
1:E:43:THR:HG21	1:E:73:ASN:OD1	2.10	0.51
1:E:72:ASN:HD22	1:E:73:ASN:N	2.08	0.51
2:B:309:LYS:HG2	2:B:315:TYR:HB2	1.92	0.51
1:A:432:VAL:HG21	1:A:482:PHE:HD2	1.75	0.51
2:D:223:ARG:NH1	2:D:223:ARG:HB2	2.24	0.51
2:H:355:PHE:O	2:H:440:PHE:HD2	1.94	0.51
1:A:12:LYS:HE3	1:A:13:TYR:CE2	2.46	0.51
1:E:317:LYS:HB3	1:E:318:GLU:OE1	2.10	0.51
2:B:32:HIS:ND1	2:B:33:PRO:HD2	2.26	0.51
1:C:426:TYR:CE2	1:C:534:LEU:HD13	2.45	0.51
1:G:112:LEU:C	1:G:114:ASP:N	2.64	0.51
2:H:90:ARG:HG2	2:H:90:ARG:NH1	2.24	0.51
2:D:152:ARG:HB3	2:D:429:THR:HG23	1.90	0.51
2:H:24:LEU:HD21	2:H:312:THR:HG21	1.91	0.51
1:G:168:GLU:HG3	1:G:394:ARG:HE	1.74	0.51
2:H:350:PRO:HB3	2:H:436:PHE:C	2.31	0.51
1:C:48:ASN:O	1:C:52:PRO:HG2	2.11	0.51
1:E:265:PHE:O	1:E:269:ILE:HG13	2.10	0.51
3:J:108:LEU:HD22	3:J:169:LEU:O	2.10	0.51
2:D:270:GLN:C	2:D:272:ASN:H	2.13	0.51
2:B:441:THR:C	2:B:442:SER:OXT	2.44	0.51
1:E:184:ARG:NH1	1:E:184:ARG:HG3	2.23	0.51
1:A:229:THR:HG22	1:A:232:ARG:HB3	1.88	0.51
2:H:316:ILE:CD1	2:H:316:ILE:H	2.14	0.51
1:G:199:TYR:CE2	1:G:216:VAL:HG11	2.45	0.51
2:H:136:ARG:HG2	2:H:161:LEU:HD22	1.93	0.51
2:F:366:ASP:HB3	2:F:370:ASN:ND2	2.26	0.51
1:A:445:VAL:O	2:B:22:LYS:HE2	2.11	0.51
1:C:446:SER:HB2	1:C:449:GLN:HB3	1.93	0.51
1:C:185:LEU:HD12	1:C:275:ALA:HB1	1.92	0.51
2:H:380:ALA:CB	2:H:394:LEU:CD1	2.78	0.51
2:H:339:ARG:HE	2:H:346:CYS:HB3	1.75	0.51
2:H:339:ARG:NH2	2:H:346:CYS:HB3	2.25	0.51
2:D:25:GLU:O	2:D:37:PRO:HB2	2.11	0.51
2:B:187:GLY:HA2	3:I:173:LEU:HD13	1.93	0.51
1:E:116:ASP:N	1:E:117:PRO:HD3	2.26	0.51
1:C:236:THR:HG22	1:C:238:LYS:H	1.76	0.51
1:G:408:ASN:OD1	1:G:408:ASN:O	2.29	0.51

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:84:ASP:OD1	2:B:85:VAL:N	2.44	0.51
2:B:27:SER:HB3	2:B:37:PRO:HG3	1.93	0.51
1:A:176:PRO:HG2	1:A:389:PHE:CD2	2.45	0.51
1:G:447:ASN:HD22	2:H:26:ARG:NH2	2.05	0.51
1:C:113:LEU:HD22	1:C:142:VAL:HG21	1.91	0.51
2:H:74:GLN:NE2	2:H:119:ASN:HD22	2.07	0.51
2:H:418:VAL:HG22	2:H:421:GLN:OE1	2.11	0.51
1:A:248:ARG:HA	1:A:251:ILE:CD1	2.39	0.51
2:H:320:ASN:HB2	2:H:337:ALA:N	2.26	0.51
1:A:215:ILE:HD12	1:A:342:LEU:HD21	1.93	0.51
1:C:344:ASN:HD22	1:C:347:ARG:HH12	1.58	0.51
2:H:350:PRO:HB3	2:H:436:PHE:O	2.11	0.50
1:G:330:PRO:O	1:G:339:TYR:OH	2.24	0.50
2:B:422:GLU:O	2:B:423:LEU:HG	2.10	0.50
1:G:72:ASN:HD22	1:G:72:ASN:N	2.06	0.50
1:A:65:VAL:HB	1:A:82:GLY:H	1.75	0.50
1:E:61:ASP:CB	1:E:86:ALA:HB2	2.40	0.50
2:B:95:ARG:HH11	2:B:95:ARG:CG	2.22	0.50
1:G:184:ARG:NH1	1:G:325:VAL:HG22	2.26	0.50
2:F:197:THR:HG23	2:F:320:ASN:OD1	2.12	0.50
1:C:216:VAL:O	1:C:220:LYS:HG2	2.11	0.50
1:C:259:PRO:CB	1:C:262:GLU:OE2	2.59	0.50
1:G:214:TRP:CZ3	1:G:332:MET:HB3	2.46	0.50
2:H:64:LEU:HD21	2:H:77:VAL:HG21	1.92	0.50
2:H:320:ASN:HB2	2:H:337:ALA:H	1.77	0.50
2:F:320:ASN:HD22	2:F:337:ALA:N	2.09	0.50
1:E:108:SER:HB2	1:E:109:PRO:HD2	1.92	0.50
1:E:187:LYS:HD2	1:E:279:THR:HG21	1.93	0.50
1:A:112:LEU:O	1:A:116:ASP:O	2.29	0.50
1:G:396:ARG:NH2	1:G:406:THR:O	2.44	0.50
1:C:268:ALA:O	1:C:272:VAL:HG23	2.12	0.50
1:E:441:ARG:NH2	1:E:453:ASP:OD1	2.44	0.50
1:C:264:ASN:ND2	1:C:265:PHE:N	2.60	0.50
2:B:237:TRP:HB3	2:B:238:PRO:HD3	1.93	0.50
2:H:270:GLN:C	2:H:271:TYR:HD2	2.14	0.50
2:F:280:LEU:O	2:F:284:VAL:HG23	2.11	0.50
2:D:16:ARG:HD3	2:D:17:TRP:HE1	1.75	0.50
2:F:236:GLN:NE2	2:F:263:LYS:HB3	2.26	0.50
2:H:250:ASP:HB3	2:H:253:ASP:HB2	1.94	0.50
2:F:232:VAL:HG21	2:F:264:SER:HA	1.93	0.50
1:E:264:ASN:ND2	1:E:265:PHE:N	2.60	0.50
1:E:66:SER:H	1:E:69:ASP:HB2	1.75	0.50

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:G:226:TYR:O	1:G:231:GLY:N	2.44	0.50
2:D:314:ALA:O	2:D:315:TYR:CD1	2.65	0.50
2:H:322:LEU:HD12	2:H:323:VAL:O	2.12	0.50
2:H:220:SER:O	2:H:222:PRO:HD3	2.11	0.50
1:G:450:VAL:O	1:G:454:ILE:HG13	2.11	0.50
2:F:338:GLU:OE1	3:K:148:LYS:HE3	2.12	0.50
1:E:375:ILE:HD12	1:E:375:ILE:N	2.27	0.50
1:A:227:SER:C	1:A:229:THR:H	2.15	0.50
2:D:24:LEU:CD2	2:D:312:THR:HB	2.41	0.50
1:C:215:ILE:HD12	1:C:342:LEU:HD21	1.93	0.50
2:D:142:VAL:HG12	2:D:178:ILE:HB	1.93	0.50
1:C:325:VAL:HG21	1:C:349:LYS:HG2	1.94	0.50
1:A:491:HIS:CD2	2:B:69:LEU:HD12	2.46	0.50
2:H:393:TYR:HA	2:H:404:THR:HB	1.93	0.50
2:F:376:MET:CE	2:F:427:ASP:HB2	2.42	0.50
2:F:401:GLU:HG3	2:F:402:GLU:N	2.27	0.50
2:B:376:MET:HB3	2:B:427:ASP:OD1	2.12	0.50
1:E:447:ASN:HD21	2:F:26:ARG:HE	1.59	0.50
2:D:83:ILE:HD12	2:D:98:ASP:HB3	1.92	0.50
2:D:277:THR:HG22	2:D:278:TYR:N	2.27	0.50
3:I:103:ILE:HD11	3:I:117:ILE:HD12	1.94	0.49
3:L:120:THR:CB	3:L:155:THR:OG1	2.60	0.49
2:H:231:TYR:C	2:H:233:ARG:N	2.66	0.49
3:I:104:LYS:O	3:I:166:VAL:HA	2.12	0.49
2:B:430:THR:HG22	2:B:431:PRO:CD	2.41	0.49
1:C:34:CYS:SG	1:C:60:ILE:HD12	2.51	0.49
1:E:19:LEU:HG	2:F:290:PRO:HB2	1.94	0.49
2:H:407:ASN:C	2:H:409:SER:H	2.15	0.49
2:B:415:LEU:HB2	2:B:417:LEU:HD21	1.93	0.49
1:G:218:ILE:O	1:G:222:LEU:HB2	2.12	0.49
1:E:309:ARG:HB3	1:E:364:LEU:CD2	2.42	0.49
2:B:391:THR:HG21	2:B:400:ILE:HG21	1.94	0.49
2:B:197:THR:HG23	2:B:198:ALA:N	2.27	0.49
1:C:454:ILE:HD13	1:C:480:HIS:ND1	2.27	0.49
1:E:199:TYR:CD2	1:E:216:VAL:HG11	2.46	0.49
3:L:105:VAL:O	3:L:105:VAL:HG13	2.11	0.49
2:H:240:GLU:O	2:H:242:PRO:CD	2.59	0.49
2:B:319:ASN:CG	2:B:336:GLU:HG3	2.32	0.49
2:D:352:ASN:HB3	2:D:439:HIS:HD2	1.78	0.49
2:D:152:ARG:NE	2:D:429:THR:HG21	2.27	0.49
2:H:322:LEU:HD12	2:H:323:VAL:N	2.26	0.49
2:B:284:VAL:O	2:B:287:ARG:NH1	2.45	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:G:43:THR:HG21	1:G:73:ASN:OD1	2.11	0.49
2:F:342:ASN:ND2	2:F:342:ASN:N	2.57	0.49
1:C:409:LYS:O	1:C:413:ILE:HG13	2.11	0.49
2:H:344:PRO:CD	2:H:374:LEU:HD22	2.42	0.49
2:B:197:THR:HG23	2:B:320:ASN:HD21	1.78	0.49
1:G:34:CYS:HB2	1:G:123:PHE:CG	2.48	0.49
1:G:288:ILE:HG23	1:G:305:TRP:CZ3	2.47	0.49
3:I:171:LEU:HD12	3:I:172:ARG:H	1.76	0.49
2:H:233:ARG:HH12	2:H:234:MET:HB2	1.76	0.49
2:D:234:MET:O	2:D:235:LEU:HD23	2.13	0.49
2:F:188:ASN:OD1	3:K:173:LEU:HD12	2.12	0.49
1:E:72:ASN:ND2	1:E:72:ASN:C	2.65	0.49
2:B:37:PRO:O	2:B:38:SER:HB2	2.11	0.49
1:G:12:LYS:HE3	1:G:13:TYR:CZ	2.48	0.49
1:C:36:ILE:HG22	1:C:37:ASN:N	2.27	0.49
3:I:126:ILE:O	3:I:130:VAL:HG23	2.11	0.49
2:D:232:VAL:HA	2:D:236:GLN:HB3	1.94	0.49
2:F:275:GLY:O	2:F:280:LEU:HD23	2.13	0.49
2:D:18:ASN:O	2:D:22:LYS:HG3	2.13	0.49
2:H:357:PRO:HG3	2:H:440:PHE:CG	2.48	0.49
1:A:133:GLU:HG3	1:A:433:ASP:HB3	1.95	0.49
3:J:104:LYS:HD3	3:J:112:GLU:OE2	2.12	0.49
2:D:434:VAL:HG13	2:D:436:PHE:HE1	1.77	0.49
1:G:12:LYS:HG3	2:H:88:LEU:HB2	1.94	0.49
2:B:348:GLN:HA	2:B:348:GLN:OE1	2.13	0.49
1:C:35:LEU:HD12	1:C:129:THR:HG23	1.94	0.49
1:G:518:ASN:ND2	1:G:534:LEU:N	2.61	0.49
2:F:401:GLU:O	2:F:405:ARG:N	2.44	0.49
1:C:157:GLY:HA3	1:C:485:TYR:CD1	2.47	0.49
3:L:162:LEU:O	3:L:164:GLY:N	2.45	0.49
2:H:18:ASN:O	2:H:22:LYS:HG2	2.12	0.49
2:F:102:PRO:HG2	2:F:105:GLU:CB	2.43	0.49
2:H:45:LEU:HD11	2:H:72:PHE:CE2	2.48	0.49
3:I:118:GLU:O	3:I:120:THR:N	2.46	0.49
3:K:118:GLU:HG2	3:K:121:ASP:CG	2.33	0.49
2:H:401:GLU:HG3	2:H:402:GLU:N	2.28	0.49
2:F:132:ASP:HB3	2:F:157:MET:CE	2.42	0.49
1:C:131:LEU:HB3	1:C:132:PRO:CD	2.43	0.49
1:G:34:CYS:HB2	1:G:123:PHE:CD2	2.47	0.49
1:C:137:LEU:HA	1:C:398:LEU:HD23	1.95	0.49
1:E:336:SER:OG	2:F:221:MET:HA	2.12	0.49
1:A:409:LYS:O	1:A:413:ILE:HG13	2.12	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:D:286:LYS:HB2	2:D:288:ILE:HG13	1.94	0.49
3:J:107:THR:HG23	3:J:109:THR:H	1.77	0.49
1:E:297:ILE:N	1:E:297:ILE:HD12	2.19	0.49
2:B:418:VAL:HG22	2:B:421:GLN:HG3	1.95	0.49
2:H:148:ILE:O	2:H:152:ARG:HG3	2.13	0.49
1:A:164:ILE:HD11	1:A:166:ILE:HD11	1.95	0.49
1:C:220:LYS:HE3	1:C:220:LYS:HA	1.95	0.49
2:D:178:ILE:HD12	2:D:307:VAL:HG22	1.94	0.49
1:A:64:GLN:HA	1:A:83:LYS:O	2.13	0.49
1:E:307:LEU:CD1	1:E:383:LEU:HD22	2.39	0.48
1:C:229:THR:HG22	1:C:232:ARG:HB2	1.91	0.48
2:H:349:LEU:HB3	2:H:350:PRO:CD	2.42	0.48
2:H:170:ASP:O	2:H:173:SER:HB3	2.12	0.48
1:G:164:ILE:HD13	1:G:515:ILE:HD12	1.95	0.48
2:D:268:ALA:CB	2:D:276:VAL:HG21	2.43	0.48
2:H:98:ASP:OD1	2:H:101:ARG:NH1	2.46	0.48
1:A:425:LEU:HD11	1:A:523:SER:HB2	1.95	0.48
1:A:323:LEU:HB3	1:A:324:PRO:HD2	1.95	0.48
3:J:124:GLU:HB3	3:J:152:ASP:O	2.12	0.48
3:K:118:GLU:O	3:K:120:THR:N	2.46	0.48
2:B:353:ILE:CG2	2:B:355:PHE:HD1	2.23	0.48
2:D:231:TYR:C	2:D:233:ARG:N	2.64	0.48
2:H:320:ASN:HB2	2:H:336:GLU:HA	1.95	0.48
1:E:12:LYS:HE3	1:E:13:TYR:CZ	2.47	0.48
1:E:158:LEU:HA	1:E:493:ILE:HG13	1.95	0.48
2:F:197:THR:CG2	2:F:198:ALA:N	2.75	0.48
2:F:322:LEU:CD1	2:F:322:LEU:C	2.82	0.48
1:A:348:GLU:HG3	1:E:115:ASN:ND2	2.27	0.48
1:G:165:ILE:HG22	1:G:165:ILE:O	2.11	0.48
2:D:207:TYR:CD1	3:J:172:ARG:HD3	2.46	0.48
1:A:211:HIS:HB3	1:A:335:ASP:HB3	1.96	0.48
2:B:353:ILE:HG22	2:B:438:LEU:HB2	1.95	0.48
1:C:526:SER:O	1:C:527:GLN:HB2	2.13	0.48
1:A:297:ILE:HG22	1:A:298:THR:N	2.29	0.48
1:C:19:LEU:HD22	2:D:185:PHE:CE1	2.48	0.48
1:E:527:GLN:CB	2:F:318:LEU:HD13	2.42	0.48
2:D:225:PRO:HG3	2:D:274:ARG:O	2.13	0.48
1:C:293:ARG:HB3	1:C:305:TRP:CD2	2.48	0.48
1:A:35:LEU:C	1:A:36:ILE:HD13	2.33	0.48
1:A:401:GLU:O	1:A:407:ILE:HG12	2.13	0.48
3:I:107:THR:CG2	3:I:108:LEU:N	2.76	0.48
1:E:220:LYS:C	1:E:222:LEU:N	2.66	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:G:218:ILE:HD11	1:G:272:VAL:HG22	1.95	0.48
2:D:208:PRO:HB2	2:D:209:PRO:HD2	1.95	0.48
1:C:240:LYS:O	1:C:244:ARG:HG3	2.13	0.48
2:H:56:ALA:O	2:H:103:LYS:HD3	2.12	0.48
2:B:417:LEU:CD1	2:B:421:GLN:HE22	2.26	0.48
1:E:259:PRO:CB	1:E:261:ASP:O	2.62	0.48
2:B:400:ILE:H	2:B:400:ILE:CD1	2.27	0.48
2:D:419:ASP:HB2	2:D:440:PHE:CD1	2.49	0.48
1:G:108:SER:HB3	1:G:111:ASN:HB3	1.94	0.48
1:C:481:GLU:HG2	1:C:525:MET:SD	2.53	0.48
2:H:52:LEU:HB2	2:H:138:PHE:CD2	2.48	0.48
1:C:521:ILE:HD12	1:C:532:PHE:CE1	2.49	0.48
1:C:190:PRO:O	1:C:194:GLU:HG3	2.13	0.48
1:A:181:GLU:O	1:A:278:THR:HB	2.13	0.48
1:E:34:CYS:HB2	1:E:123:PHE:CD2	2.48	0.48
1:C:354:ALA:HA	1:C:384:CYS:SG	2.52	0.48
1:A:401:GLU:OE2	1:A:534:LEU:HB2	2.13	0.48
1:E:48:ASN:HB2	1:E:502:ALA:CB	2.43	0.48
2:H:158:LEU:HD13	2:H:177:LEU:HD22	1.96	0.48
2:H:322:LEU:HD12	2:H:322:LEU:C	2.34	0.48
1:A:236:THR:O	1:A:240:LYS:HG3	2.13	0.48
2:H:382:THR:HA	2:H:392:LEU:HG	1.96	0.48
1:E:183:LEU:O	1:E:184:ARG:HB2	2.14	0.48
3:L:170:VAL:CG1	3:L:171:LEU:H	2.18	0.48
2:D:249:LEU:HD11	2:D:260:ILE:HD11	1.95	0.48
2:B:198:ALA:HB3	2:B:320:ASN:ND2	2.28	0.48
2:D:87:ASN:ND2	2:D:103:LYS:HE2	2.29	0.48
1:E:481:GLU:OE1	2:F:29:PRO:HD2	2.13	0.48
2:F:231:TYR:CG	2:F:267:ARG:NH1	2.81	0.48
1:C:50:VAL:HG13	1:C:100:VAL:HG21	1.95	0.48
1:E:422:GLU:HG3	1:E:530:ALA:CB	2.42	0.48
1:E:435:PHE:HE1	1:E:456:LYS:HB2	1.78	0.48
2:B:405:ARG:HB3	2:B:406:PRO:HD3	1.95	0.48
2:D:361:LEU:O	2:D:363:GLU:N	2.47	0.48
1:A:47:LYS:O	1:A:47:LYS:HD2	2.14	0.48
3:K:115:ILE:CG2	3:K:117:ILE:HD13	2.43	0.48
2:H:89:ASN:OD1	2:H:90:ARG:NH1	2.47	0.48
2:H:377:LYS:HD3	2:H:428:VAL:HG21	1.96	0.48
2:D:393:TYR:CZ	2:D:408:LEU:HD11	2.49	0.48
3:L:161:ILE:O	3:L:162:LEU:HD23	2.14	0.48
1:A:186:ASP:C	1:A:188:PRO:HD3	2.35	0.48
2:H:414:GLU:HG2	2:H:414:GLU:O	2.14	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:527:GLN:NE2	2:B:302:VAL:HG13	2.28	0.48
2:B:391:THR:HG22	2:B:391:THR:O	2.13	0.48
1:E:491:HIS:NE2	2:F:65:LYS:HE3	2.29	0.48
2:D:13:TRP:C	2:D:13:TRP:CD1	2.87	0.48
2:F:102:PRO:HG2	2:F:105:GLU:HB3	1.95	0.48
2:H:251:GLY:O	2:H:286:LYS:HD2	2.13	0.48
1:E:187:LYS:HD2	1:E:279:THR:CG2	2.44	0.47
2:B:208:PRO:HG3	3:I:171:LEU:HD11	1.95	0.47
1:C:418:ASN:HD22	1:C:419:PRO:CD	2.27	0.47
2:D:73:ARG:HG3	2:D:73:ARG:NH1	2.29	0.47
1:E:422:GLU:H	1:E:422:GLU:CD	2.15	0.47
2:F:158:LEU:HD12	2:F:175:VAL:HB	1.96	0.47
3:J:115:ILE:HG23	3:J:129:ARG:HB3	1.95	0.47
1:G:18:ARG:NH1	2:H:279:ARG:HG3	2.28	0.47
1:E:34:CYS:HB2	1:E:123:PHE:CE2	2.49	0.47
1:C:175:HIS:HD2	1:C:512:GLN:O	1.97	0.47
2:H:197:THR:HG22	2:H:198:ALA:N	2.29	0.47
1:C:264:ASN:HD22	1:C:264:ASN:N	2.10	0.47
2:H:249:LEU:HD11	2:H:260:ILE:HD11	1.95	0.47
1:G:120:PHE:O	1:G:148:ILE:HD13	2.14	0.47
2:H:57:GLY:HA2	2:H:91:GLN:HG2	1.96	0.47
2:B:187:GLY:CA	3:I:173:LEU:HD13	2.44	0.47
1:G:36:ILE:HG21	1:G:131:LEU:HD11	1.96	0.47
1:G:43:THR:HG22	1:G:44:GLU:N	2.29	0.47
1:E:75:PHE:O	1:E:76:LEU:HD23	2.14	0.47
1:A:40:ALA:O	1:A:43:THR:HG22	2.14	0.47
1:C:168:GLU:HG2	1:C:170:PRO:HD3	1.96	0.47
1:C:457:LEU:C	1:C:459:SER:H	2.16	0.47
2:B:403:ARG:HH11	2:B:403:ARG:HG3	1.80	0.47
1:G:225:TRP:CE2	1:G:234:PRO:HG3	2.49	0.47
1:E:19:LEU:HD22	2:F:185:PHE:CE1	2.48	0.47
2:D:73:ARG:HG3	2:D:73:ARG:HH11	1.78	0.47
2:B:351:GLN:H	2:B:436:PHE:HA	1.78	0.47
1:E:43:THR:HG22	1:E:44:GLU:N	2.29	0.47
1:E:236:THR:HG22	1:E:237:TYR:N	2.28	0.47
1:E:163:ARG:HH11	1:E:165:ILE:HG12	1.79	0.47
1:C:25:GLN:O	1:C:29:GLU:HG3	2.14	0.47
2:H:267:ARG:O	2:H:267:ARG:HD3	2.14	0.47
2:F:231:TYR:C	2:F:233:ARG:H	2.16	0.47
2:H:190:GLN:OE1	3:L:172:ARG:NH1	2.47	0.47
2:H:423:LEU:N	2:H:423:LEU:HD12	2.29	0.47
2:D:182:THR:CG2	2:D:183:GLU:N	2.77	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:435:PHE:O	1:A:438:GLN:N	2.48	0.47
1:A:32:HIS:HD2	1:A:56:SER:O	1.97	0.47
1:C:130:GLN:NE2	1:C:155:THR:O	2.47	0.47
3:J:143:LEU:HB3	3:J:150:MET:HE1	1.97	0.47
3:J:156:ALA:O	3:J:161:ILE:HB	2.14	0.47
2:B:421:GLN:HB2	2:B:438:LEU:HD21	1.96	0.47
2:D:322:LEU:C	2:D:322:LEU:CD1	2.82	0.47
2:H:402:GLU:HA	2:H:405:ARG:NH1	2.30	0.47
2:H:412:LEU:O	2:H:417:LEU:HB2	2.15	0.47
2:H:354:GLN:NE2	2:H:354:GLN:HA	2.29	0.47
2:B:274:ARG:HD3	1:E:62:GLY:O	2.14	0.47
1:C:248:ARG:C	1:C:250:GLY:N	2.65	0.47
2:B:81:ASP:CB	2:B:103:LYS:HD2	2.42	0.47
2:D:225:PRO:HB2	2:D:280:LEU:HD21	1.97	0.47
1:C:484:ARG:HD2	2:D:27:SER:O	2.13	0.47
1:C:163:ARG:NE	1:C:518:ASN:OD1	2.46	0.47
1:A:516:PHE:H	1:A:516:PHE:HD2	1.62	0.47
2:F:380:ALA:HB3	2:F:394:LEU:HD12	1.90	0.47
2:F:425:VAL:O	2:F:434:VAL:HG12	2.15	0.47
2:F:361:LEU:O	2:F:364:VAL:N	2.42	0.47
3:L:123:VAL:HA	3:L:126:ILE:CG1	2.45	0.47
3:L:155:THR:O	3:L:158:ASP:HB2	2.14	0.47
3:J:145:TYR:HB2	3:J:167:LEU:CD2	2.44	0.47
2:H:87:ASN:HB3	2:H:91:GLN:NE2	2.29	0.47
2:H:351:GLN:HB3	2:H:436:PHE:CD2	2.50	0.47
2:H:159:ILE:HG23	2:H:162:LEU:HD12	1.96	0.47
2:F:422:GLU:HA	2:F:436:PHE:O	2.15	0.47
2:F:123:HIS:HB3	2:F:125:ASN:HD21	1.79	0.47
1:A:154:ARG:HH11	1:A:154:ARG:HG2	1.79	0.47
2:H:412:LEU:CD2	2:H:438:LEU:HD21	2.44	0.47
2:H:419:ASP:HB2	2:H:440:PHE:HD1	1.79	0.47
1:G:158:LEU:CD2	2:H:23:PHE:CE2	2.98	0.47
1:G:184:ARG:NH2	1:G:323:LEU:O	2.47	0.47
2:D:197:THR:CG2	2:D:198:ALA:N	2.77	0.47
1:E:194:GLU:C	1:E:196:PHE:H	2.18	0.47
2:D:152:ARG:NH2	2:D:374:LEU:O	2.48	0.47
2:D:280:LEU:O	2:D:284:VAL:HG23	2.15	0.47
1:G:97:ASN:OD1	1:G:99:ASP:HB2	2.13	0.47
1:C:326:ARG:HG2	1:C:328:THR:HG23	1.96	0.47
2:B:166:ASP:O	2:B:168:VAL:N	2.45	0.47
2:F:136:ARG:NH1	2:F:161:LEU:HD22	2.30	0.47
1:A:424:VAL:HA	1:A:427:LEU:HD12	1.97	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:252:LEU:C	1:A:252:LEU:HD22	2.35	0.47
3:I:123:VAL:O	3:I:126:ILE:N	2.45	0.47
2:H:207:TYR:CE1	3:L:172:ARG:CD	2.97	0.47
2:F:141:ILE:HD12	2:F:158:LEU:HD11	1.97	0.47
2:D:382:THR:HG22	2:D:391:THR:HA	1.96	0.47
2:H:306:GLU:O	2:H:307:VAL:C	2.53	0.47
1:A:467:GLU:HG2	1:A:467:GLU:O	2.15	0.47
1:A:262:GLU:HG2	1:A:265:PHE:HD1	1.79	0.47
1:A:265:PHE:O	1:A:269:ILE:HG13	2.15	0.47
2:B:321:TYR:OH	3:I:172:ARG:CG	2.63	0.47
2:B:218:ILE:HG22	2:B:219:ALA:N	2.29	0.47
2:F:225:PRO:HG2	2:F:275:GLY:HA3	1.97	0.47
1:C:138:ARG:O	1:C:142:VAL:HG23	2.15	0.47
2:B:427:ASP:CB	2:B:429:THR:HG22	2.45	0.47
2:D:312:THR:O	2:D:313:SER:HB3	2.14	0.47
1:A:128:ALA:HB1	1:A:131:LEU:HD12	1.97	0.47
1:C:214:TRP:C	1:C:214:TRP:CD1	2.88	0.47
2:D:209:PRO:HG3	3:J:139:GLN:HG2	1.96	0.47
1:C:75:PHE:CD2	1:C:93:LEU:HD23	2.50	0.47
2:H:63:LEU:HD22	2:H:142:VAL:HG11	1.97	0.47
1:A:262:GLU:HG2	1:A:265:PHE:CD1	2.50	0.47
1:C:401:GLU:OE1	1:C:426:TYR:OH	2.29	0.47
1:A:34:CYS:SG	1:A:36:ILE:CD1	3.02	0.47
1:G:180:LEU:HD12	1:G:180:LEU:N	2.19	0.47
1:A:489:GLU:O	1:A:489:GLU:HG2	2.14	0.47
2:F:129:ASP:HB3	2:H:134:PHE:HB2	1.95	0.47
1:E:447:ASN:ND2	2:F:26:ARG:NH2	2.62	0.47
1:C:307:LEU:HB3	1:C:383:LEU:HD22	1.97	0.47
1:C:449:GLN:O	1:C:449:GLN:HG2	2.15	0.47
1:A:504:GLU:O	1:A:508:ILE:HG13	2.15	0.47
2:F:386:GLU:CB	2:F:388:LYS:HE2	2.45	0.47
2:D:192:ILE:O	2:D:194:PRO:HD3	2.15	0.46
2:F:355:PHE:O	2:F:440:PHE:HA	2.16	0.46
1:G:510:THR:OG1	1:G:512:GLN:HB2	2.16	0.46
1:E:72:ASN:HD22	1:E:72:ASN:C	2.19	0.46
3:L:104:LYS:HB2	3:L:166:VAL:HG22	1.98	0.46
1:A:371:ALA:O	1:A:373:GLU:N	2.47	0.46
2:B:126:LYS:O	2:B:128:GLN:N	2.48	0.46
2:B:312:THR:O	2:B:314:ALA:N	2.48	0.46
1:G:433:ASP:O	1:G:436:HIS:HB3	2.15	0.46
2:B:380:ALA:HB1	2:B:394:LEU:HD13	1.94	0.46
2:H:87:ASN:HB3	2:H:91:GLN:CD	2.35	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:D:412:LEU:HD22	2:D:417:LEU:CD1	2.44	0.46
1:C:131:LEU:HB3	1:C:132:PRO:HD2	1.97	0.46
1:G:480:HIS:HB2	2:H:29:PRO:HG2	1.98	0.46
2:F:381:ILE:HG13	2:F:425:VAL:HG22	1.96	0.46
2:F:348:GLN:C	2:F:349:LEU:HD12	2.34	0.46
2:F:335:PHE:HE2	3:K:170:VAL:HG21	1.80	0.46
1:C:184:ARG:HD3	1:C:279:THR:OG1	2.14	0.46
3:J:150:MET:HB2	3:J:150:MET:HE2	1.57	0.46
1:A:518:ASN:HD21	1:A:534:LEU:CB	2.26	0.46
1:E:347:ARG:HH22	2:F:274:ARG:CD	2.15	0.46
1:A:78:ARG:C	1:A:80:SER:H	2.19	0.46
2:B:16:ARG:HH22	2:B:116:PRO:HB2	1.80	0.46
1:G:447:ASN:HA	2:H:26:ARG:HH21	1.80	0.46
1:A:366:GLN:O	1:A:368:ILE:N	2.47	0.46
1:C:481:GLU:OE1	2:D:315:TYR:OH	2.32	0.46
1:C:489:GLU:O	1:C:489:GLU:HG2	2.16	0.46
1:C:72:ASN:HB2	2:D:19:HIS:ND1	2.31	0.46
2:D:187:GLY:CA	3:J:173:LEU:HD12	2.46	0.46
2:B:27:SER:HB3	2:B:37:PRO:CG	2.46	0.46
1:G:124:THR:HG21	1:G:509:ILE:HG12	1.96	0.46
2:H:257:ILE:HD13	2:H:282:GLN:HG2	1.98	0.46
3:L:151:ASN:O	3:L:153:GLU:N	2.48	0.46
1:C:428:MET:HE1	1:C:479:VAL:HA	1.97	0.46
2:F:357:PRO:HD3	2:F:440:PHE:CG	2.51	0.46
2:B:207:TYR:HE2	3:I:172:ARG:HG2	1.72	0.46
3:K:121:ASP:HB3	3:K:125:ARG:HD2	1.98	0.46
1:G:504:GLU:OE2	1:G:507:LYS:NZ	2.42	0.46
3:J:105:VAL:HG11	3:J:130:VAL:HG22	1.97	0.46
2:F:312:THR:O	2:F:313:SER:CB	2.63	0.46
2:F:356:SER:O	2:F:358:SER:N	2.48	0.46
2:F:106:VAL:O	2:F:109:GLU:HB3	2.16	0.46
2:B:418:VAL:CG2	2:B:421:GLN:HG3	2.46	0.46
2:B:240:GLU:CG	2:B:240:GLU:O	2.62	0.46
2:H:233:ARG:HG2	2:H:234:MET:N	2.31	0.46
1:C:488:ALA:HB2	2:D:22:LYS:HD2	1.98	0.46
1:C:146:SER:O	1:C:147:GLN:CB	2.61	0.46
1:G:386:ASN:O	1:G:387:SER:C	2.54	0.46
2:H:158:LEU:HD13	2:H:177:LEU:HB2	1.97	0.46
1:C:137:LEU:HD21	1:C:402:TYR:CG	2.51	0.46
2:H:213:PHE:O	2:H:218:ILE:HD11	2.16	0.46
2:D:67:LEU:O	2:D:72:PHE:HB2	2.15	0.46
1:C:259:PRO:CB	1:C:262:GLU:CD	2.83	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:264:ASN:ND2	1:C:265:PHE:H	2.12	0.46
2:D:405:ARG:N	2:D:406:PRO:CD	2.79	0.46
1:A:130:GLN:HA	1:A:154:ARG:HD2	1.98	0.46
2:D:419:ASP:CG	2:D:420:GLY:N	2.69	0.46
1:E:44:GLU:OE1	2:F:65:LYS:NZ	2.48	0.46
1:C:214:TRP:CD1	1:C:215:ILE:N	2.84	0.46
3:I:142:ARG:HA	3:I:142:ARG:HD3	1.82	0.46
2:D:125:ASN:ND2	2:D:125:ASN:N	2.60	0.46
1:G:248:ARG:O	1:G:251:ILE:CD1	2.64	0.46
3:J:135:GLY:O	3:J:137:PRO:HD3	2.16	0.46
2:F:85:VAL:HG13	2:F:86:SER:N	2.31	0.46
1:A:110:GLU:OE1	1:A:135:THR:OG1	2.31	0.46
2:H:163:ASN:O	2:H:163:ASN:OD1	2.33	0.46
1:G:396:ARG:HH21	1:G:400:GLU:HB3	1.80	0.46
2:F:366:ASP:HA	2:F:369:THR:HB	1.98	0.46
2:B:28:GLY:O	2:B:31:THR:HG23	2.16	0.46
2:D:438:LEU:HD12	2:D:439:HIS:H	1.81	0.46
1:E:301:THR:HG23	1:E:305:TRP:HB2	1.98	0.46
1:G:347:ARG:HH12	2:H:274:ARG:HD2	1.81	0.46
1:G:157:GLY:HA3	1:G:485:TYR:CG	2.51	0.46
2:B:349:LEU:HD23	2:B:349:LEU:N	2.30	0.46
2:F:412:LEU:HB3	2:F:440:PHE:CZ	2.51	0.46
3:L:170:VAL:CG1	3:L:171:LEU:N	2.76	0.46
2:D:228:CYS:O	2:D:231:TYR:HB3	2.16	0.46
2:H:373:SER:OG	2:H:374:LEU:HD23	2.16	0.46
1:A:324:PRO:HB3	1:A:353:ASP:HB3	1.98	0.46
1:E:178:ASN:O	1:E:179:ALA:HB2	2.15	0.46
3:J:125:ARG:HH11	3:J:125:ARG:HG3	1.81	0.46
1:E:49:LEU:O	1:E:52:PRO:HG2	2.14	0.45
1:A:78:ARG:HA	1:A:81:ILE:CG1	2.46	0.45
2:H:357:PRO:HD3	2:H:440:PHE:HB3	1.96	0.45
2:H:417:LEU:HD23	2:H:421:GLN:NE2	2.31	0.45
1:C:119:PHE:HE1	1:C:122:ARG:NH1	2.14	0.45
1:A:348:GLU:O	1:A:352:LYS:HG3	2.16	0.45
2:H:75:ILE:O	2:H:120:VAL:HA	2.15	0.45
1:A:490:PRO:HG3	2:B:23:PHE:HE1	1.81	0.45
1:G:173:GLU:CG	1:G:382:LEU:HD11	2.46	0.45
1:E:132:PRO:O	1:E:133:GLU:C	2.54	0.45
1:G:28:LEU:HD12	1:G:28:LEU:HA	1.77	0.45
2:F:393:TYR:CE2	2:F:408:LEU:HD11	2.52	0.45
2:F:131:ASN:O	2:F:133:THR:N	2.49	0.45
1:A:38:ALA:O	1:A:85:ARG:HD3	2.16	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:G:197:GLN:C	1:G:199:TYR:H	2.19	0.45
2:H:74:GLN:HE22	2:H:119:ASN:ND2	2.12	0.45
2:F:32:HIS:CG	2:F:33:PRO:HD2	2.51	0.45
1:A:164:ILE:HG23	1:A:164:ILE:O	2.15	0.45
1:C:310:ALA:CB	1:C:361:VAL:HG22	2.47	0.45
2:H:158:LEU:HD12	2:H:177:LEU:HD22	1.98	0.45
2:D:76:HIS:CD2	2:D:138:PHE:HE2	2.34	0.45
1:E:181:GLU:HG3	1:E:330:PRO:HD3	1.96	0.45
3:I:139:GLN:O	3:I:142:ARG:NH2	2.36	0.45
2:B:217:THR:HG21	2:B:223:ARG:HH22	1.80	0.45
2:D:158:LEU:CD1	2:D:177:LEU:HB2	2.46	0.45
2:F:316:ILE:N	2:F:316:ILE:HD12	2.27	0.45
2:B:130:PHE:HE2	2:D:125:ASN:OD1	1.99	0.45
2:F:13:TRP:CZ3	2:F:116:PRO:HG2	2.50	0.45
2:H:381:ILE:H	2:H:381:ILE:HD12	1.81	0.45
1:C:443:PRO:HB2	1:C:483:CYS:SG	2.56	0.45
1:E:46:LEU:HD23	1:E:93:LEU:HD13	1.97	0.45
1:C:94:GLN:HE22	1:C:102:GLY:H	1.64	0.45
1:E:125:VAL:HG12	1:E:126:VAL:N	2.31	0.45
1:C:28:LEU:HD12	1:C:54:ILE:HD12	1.97	0.45
2:D:257:ILE:HG21	2:D:282:GLN:HG2	1.97	0.45
1:A:474:VAL:O	1:A:475:LYS:C	2.53	0.45
1:C:318:GLU:HG3	1:C:356:ALA:HB1	1.98	0.45
1:G:504:GLU:O	1:G:508:ILE:HG13	2.17	0.45
1:G:214:TRP:C	1:G:214:TRP:CD1	2.90	0.45
1:G:158:LEU:HA	1:G:493:ILE:HG13	1.99	0.45
1:E:236:THR:C	1:E:238:LYS:H	2.20	0.45
3:L:103:ILE:HG22	3:L:163:GLY:O	2.17	0.45
1:C:236:THR:CB	1:C:239:GLU:HG3	2.45	0.45
1:E:527:GLN:OE1	2:F:302:VAL:HG13	2.16	0.45
2:F:204:LEU:HD21	2:F:375:GLN:CG	2.46	0.45
1:G:262:GLU:HG3	1:G:264:ASN:ND2	2.32	0.45
2:F:325:ASN:HD22	2:F:326:ASP:H	1.65	0.45
3:L:107:THR:HG22	3:L:111:LYS:HB3	1.97	0.45
2:B:421:GLN:O	2:B:438:LEU:HD23	2.17	0.45
2:F:87:ASN:HD22	2:F:103:LYS:HE2	1.78	0.45
1:A:430:ARG:CZ	1:A:464:PHE:CE1	3.00	0.45
3:K:156:ALA:CA	3:K:161:ILE:HD12	2.45	0.45
1:E:245:ASP:C	1:E:247:ILE:H	2.19	0.45
2:B:405:ARG:HH11	2:B:405:ARG:HG3	1.82	0.45
1:A:344:ASN:HA	1:A:347:ARG:HB2	1.97	0.45
1:G:172:ILE:HD12	1:G:172:ILE:N	2.32	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:L:107:THR:HG22	3:L:111:LYS:N	2.32	0.45
1:C:110:GLU:O	1:C:113:LEU:N	2.49	0.45
1:A:489:GLU:H	2:B:19:HIS:HD2	1.61	0.45
2:D:58:GLY:O	2:D:62:GLU:HB2	2.17	0.45
2:F:158:LEU:CD1	2:F:175:VAL:HB	2.47	0.45
3:K:107:THR:CG2	3:K:108:LEU:N	2.79	0.45
2:D:229:ILE:HA	2:D:264:SER:OG	2.16	0.45
2:D:270:GLN:O	2:D:272:ASN:N	2.50	0.45
2:F:101:ARG:HB3	2:F:102:PRO:HD2	1.98	0.45
1:E:286:GLU:HA	1:E:286:GLU:OE1	2.16	0.45
1:A:163:ARG:HG2	1:A:163:ARG:HH11	1.82	0.45
2:D:136:ARG:HG2	2:D:136:ARG:NH1	2.31	0.45
2:D:11:LEU:N	2:D:11:LEU:HD23	2.32	0.45
3:J:162:LEU:O	3:J:163:GLY:C	2.55	0.45
1:E:220:LYS:O	1:E:222:LEU:N	2.50	0.45
2:H:178:ILE:HD11	2:H:310:ILE:HD13	1.96	0.45
2:B:25:GLU:O	2:B:37:PRO:HB3	2.16	0.45
1:C:47:LYS:HZ1	2:D:65:LYS:HE2	1.82	0.45
2:D:145:LEU:HD12	2:D:151:ARG:HG3	1.99	0.45
2:D:54:ILE:HG22	2:D:145:LEU:HD21	1.97	0.45
1:G:340:ILE:HD13	1:G:340:ILE:HA	1.86	0.45
1:A:252:LEU:C	1:A:259:PRO:CD	2.72	0.45
1:A:36:ILE:HD11	1:A:60:ILE:HD12	1.99	0.45
3:I:117:ILE:CG2	3:I:118:GLU:N	2.79	0.45
2:B:342:ASN:ND2	2:B:342:ASN:N	2.43	0.45
2:B:64:LEU:HB3	2:B:111:LEU:HD13	1.97	0.45
2:B:226:GLU:N	2:B:226:GLU:OE1	2.44	0.45
2:H:110:PHE:CE2	2:H:114:ARG:NH1	2.85	0.45
1:A:513:PHE:N	1:A:513:PHE:CD1	2.85	0.45
2:F:185:PHE:HB3	2:F:326:ASP:HB3	1.98	0.45
2:B:356:SER:CA	2:B:441:THR:OG1	2.65	0.45
2:B:217:THR:CG2	2:B:223:ARG:HH22	2.30	0.45
1:A:418:ASN:HB3	1:A:421:ASN:HB2	1.99	0.45
2:D:188:ASN:HA	2:D:322:LEU:O	2.17	0.45
2:B:83:ILE:HB	2:B:99:ILE:HA	1.99	0.45
1:C:500:ALA:O	1:C:504:GLU:HG2	2.17	0.45
2:D:151:ARG:HG2	2:D:179:ASP:OD2	2.17	0.45
2:B:185:PHE:HE2	2:B:292:VAL:HG21	1.82	0.45
1:A:113:LEU:O	1:A:138:ARG:NH2	2.50	0.45
1:C:262:GLU:HB3	1:C:265:PHE:HB2	1.99	0.45
2:D:207:TYR:HE1	3:J:172:ARG:HD3	1.64	0.45
2:H:232:VAL:HG12	2:H:232:VAL:O	2.17	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:252:LEU:O	1:E:259:PRO:CD	2.65	0.45
1:C:407:ILE:HG23	1:C:409:LYS:HG3	1.99	0.45
2:F:32:HIS:CD2	2:F:34:ASP:HB2	2.52	0.45
1:G:271:ASN:O	1:G:275:ALA:N	2.47	0.45
2:D:58:GLY:N	2:D:91:GLN:HG2	2.32	0.45
2:B:32:HIS:CD2	2:B:34:ASP:HB2	2.52	0.45
1:E:178:ASN:ND2	3:K:136:ILE:HG12	2.32	0.45
1:C:90:MET:O	1:C:94:GLN:HB2	2.17	0.45
2:D:64:LEU:HB3	2:D:111:LEU:CD1	2.47	0.45
2:D:356:SER:O	2:D:358:SER:N	2.50	0.45
2:D:149:ILE:CD1	2:D:149:ILE:N	2.62	0.44
1:E:51:LEU:N	1:E:52:PRO:HD2	2.32	0.44
1:C:418:ASN:C	1:C:418:ASN:ND2	2.69	0.44
2:H:402:GLU:HG3	2:H:402:GLU:H	1.51	0.44
2:B:350:PRO:HB2	2:B:436:PHE:C	2.38	0.44
1:E:36:ILE:HB	1:E:128:ALA:HA	1.99	0.44
2:H:225:PRO:HG3	2:H:274:ARG:O	2.17	0.44
1:G:113:LEU:HD11	1:G:139:LEU:HB2	1.99	0.44
3:L:138:PRO:HA	3:L:141:GLN:NE2	2.31	0.44
3:L:123:VAL:HA	3:L:126:ILE:HG13	1.98	0.44
2:B:419:ASP:HA	2:B:440:PHE:CE1	2.52	0.44
1:A:28:LEU:CD1	1:A:509:ILE:HG21	2.44	0.44
2:B:87:ASN:ND2	2:B:103:LYS:HE3	2.33	0.44
2:F:208:PRO:HG3	3:K:171:LEU:HD11	2.00	0.44
1:A:49:LEU:C	1:A:52:PRO:HD2	2.38	0.44
1:G:428:MET:O	1:G:432:VAL:HG23	2.17	0.44
1:A:376:SER:OG	1:A:379:GLU:HG3	2.18	0.44
2:D:226:GLU:N	2:D:226:GLU:OE2	2.49	0.44
1:A:259:PRO:CG	1:A:262:GLU:OE1	2.65	0.44
1:A:229:THR:O	1:A:229:THR:HG22	2.17	0.44
1:A:419:PRO:HB2	1:A:475:LYS:HE3	2.00	0.44
1:E:259:PRO:HB3	1:E:261:ASP:O	2.17	0.44
2:H:230:GLU:HA	2:H:230:GLU:OE1	2.17	0.44
2:H:230:GLU:O	2:H:234:MET:HB3	2.17	0.44
2:H:344:PRO:HD3	2:H:374:LEU:HD22	1.98	0.44
1:C:504:GLU:HG3	1:C:516:PHE:CE2	2.51	0.44
1:E:229:THR:O	1:E:232:ARG:HG3	2.18	0.44
2:H:391:THR:O	2:H:392:LEU:C	2.55	0.44
3:L:155:THR:O	3:L:158:ASP:N	2.50	0.44
1:A:78:ARG:NH2	2:B:13:TRP:HB3	2.32	0.44
2:H:132:ASP:O	2:H:136:ARG:HG3	2.18	0.44
3:K:170:VAL:HG13	3:K:171:LEU:N	2.33	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:439:GLN:HE21	1:A:441:ARG:HH12	1.65	0.44
1:A:347:ARG:HH12	2:B:274:ARG:HH22	1.65	0.44
1:G:363:LYS:O	1:G:366:GLN:HB2	2.16	0.44
1:G:320:GLN:OE1	1:G:320:GLN:HA	2.17	0.44
2:D:405:ARG:NH1	2:D:405:ARG:CB	2.80	0.44
2:D:251:GLY:O	2:D:257:ILE:HD11	2.17	0.44
2:H:269:SER:O	2:H:272:ASN:OD1	2.36	0.44
1:G:534:LEU:HD23	1:G:534:LEU:N	2.32	0.44
2:H:245:GLU:CG	2:H:246:GLY:N	2.80	0.44
2:D:74:GLN:HE22	2:D:119:ASN:ND2	2.15	0.44
2:B:399:SER:HB3	2:B:400:ILE:HD12	1.99	0.44
2:F:261:PHE:O	2:F:264:SER:HB2	2.17	0.44
1:E:194:GLU:C	1:E:196:PHE:N	2.71	0.44
3:K:102:LEU:HD23	3:K:102:LEU:C	2.37	0.44
2:F:355:PHE:CZ	2:F:364:VAL:HG22	2.51	0.44
1:G:426:TYR:CE2	1:G:534:LEU:HD13	2.53	0.44
1:E:61:ASP:OD2	1:E:85:ARG:HD2	2.18	0.44
2:H:186:LYS:HG2	2:H:325:ASN:ND2	2.32	0.44
1:A:409:LYS:HB3	1:A:413:ILE:HD11	1.98	0.44
1:A:412:ILE:HG21	1:A:427:LEU:HD21	2.00	0.44
2:F:84:ASP:OD1	2:F:85:VAL:N	2.49	0.44
1:E:46:LEU:O	1:E:50:VAL:HG23	2.18	0.44
2:H:54:ILE:HG22	2:H:145:LEU:HD21	2.00	0.44
1:A:60:ILE:HG22	1:A:60:ILE:O	2.16	0.44
1:A:210:SER:O	1:A:211:HIS:HD2	2.01	0.44
1:E:49:LEU:C	1:E:52:PRO:HD2	2.37	0.44
1:C:368:ILE:O	1:C:368:ILE:HG22	2.17	0.44
2:D:237:TRP:CZ2	2:D:249:LEU:HA	2.53	0.44
1:A:441:ARG:NH1	1:A:441:ARG:HG2	2.32	0.44
1:C:333:ILE:O	1:C:334:ALA:HB2	2.16	0.44
1:G:283:SER:O	1:G:286:GLU:HB2	2.18	0.44
2:H:380:ALA:O	2:H:425:VAL:HA	2.17	0.44
2:B:386:GLU:O	2:B:387:GLY:C	2.54	0.44
2:D:327:VAL:HG23	2:D:328:ASP:N	2.33	0.44
1:G:222:LEU:HD13	1:G:243:PHE:CZ	2.53	0.44
1:A:430:ARG:CZ	1:A:464:PHE:HE1	2.31	0.44
2:F:192:ILE:O	2:F:193:LEU:HD23	2.18	0.44
2:B:237:TRP:CE2	2:B:249:LEU:HB2	2.52	0.44
2:H:213:PHE:CD1	2:H:213:PHE:N	2.79	0.44
1:E:50:VAL:HG13	1:E:100:VAL:HG21	1.99	0.44
2:B:362:GLN:OE1	2:B:362:GLN:HA	2.17	0.44
1:G:293:ARG:NH1	1:G:293:ARG:CG	2.77	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:F:53:VAL:HB	2:F:77:VAL:HG22	1.99	0.44
2:H:84:ASP:OD1	2:H:86:SER:N	2.44	0.44
2:H:136:ARG:CG	2:H:161:LEU:HD22	2.48	0.44
2:H:353:ILE:HG22	2:H:355:PHE:CD1	2.53	0.44
2:D:73:ARG:O	2:D:119:ASN:HB3	2.18	0.44
1:C:349:LYS:O	1:C:352:LYS:N	2.50	0.44
2:D:356:SER:C	2:D:358:SER:H	2.21	0.44
1:G:46:LEU:HD23	1:G:93:LEU:HD13	1.99	0.44
1:E:359:ASN:HB3	1:E:363:LYS:HE3	1.99	0.44
2:H:410:LYS:CB	2:H:415:LEU:HD21	2.48	0.44
2:H:379:PRO:HD2	2:H:395:GLN:NE2	2.33	0.44
3:L:105:VAL:HA	3:L:167:LEU:O	2.18	0.43
2:D:141:ILE:HD12	2:D:158:LEU:HD11	2.00	0.43
2:D:223:ARG:HH12	2:D:227:HIS:CE1	2.36	0.43
2:F:240:GLU:O	2:F:241:GLN:C	2.56	0.43
1:G:214:TRP:CE2	1:G:271:ASN:ND2	2.86	0.43
2:D:29:PRO:C	2:D:31:THR:H	2.22	0.43
1:G:386:ASN:O	1:G:388:ALA:N	2.51	0.43
1:C:61:ASP:HB3	1:C:86:ALA:HB2	2.00	0.43
3:L:104:LYS:HG2	3:L:114:GLU:HG2	2.00	0.43
1:G:416:MET:C	1:G:418:ASN:H	2.22	0.43
2:B:402:GLU:C	2:B:404:THR:H	2.21	0.43
2:H:165:GLU:O	2:H:166:ASP:HB2	2.18	0.43
2:F:56:ALA:HA	2:F:60:GLY:HA3	1.99	0.43
1:G:226:TYR:CB	1:G:231:GLY:HA2	2.46	0.43
1:C:34:CYS:HB2	1:C:123:PHE:CE2	2.54	0.43
3:K:170:VAL:HG22	3:K:171:LEU:H	1.83	0.43
3:L:159:TYR:O	3:L:160:LYS:HB2	2.18	0.43
1:C:65:VAL:HG22	1:C:85:ARG:HA	1.99	0.43
1:E:505:VAL:O	1:E:508:ILE:HB	2.18	0.43
2:F:377:LYS:HD2	2:F:377:LYS:O	2.18	0.43
1:C:56:SER:HB2	1:C:101:SER:O	2.17	0.43
2:F:367:TYR:O	2:F:371:SER:HB3	2.18	0.43
3:L:120:THR:O	3:L:120:THR:OG1	2.31	0.43
1:C:416:MET:HE1	1:C:424:VAL:HG12	2.01	0.43
2:B:123:HIS:CE1	2:D:125:ASN:HD21	2.36	0.43
2:D:412:LEU:CD1	2:D:440:PHE:HE2	2.31	0.43
1:E:314:PHE:HB2	1:E:357:VAL:HG22	1.99	0.43
1:A:214:TRP:CD1	1:A:215:ILE:N	2.87	0.43
1:A:215:ILE:HG13	1:A:332:MET:SD	2.58	0.43
2:D:408:LEU:HD23	2:D:408:LEU:HA	1.86	0.43
1:C:293:ARG:HH11	1:C:293:ARG:HG3	1.84	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:F:162:LEU:HD22	2:F:169:LEU:HD11	2.00	0.43
1:A:457:LEU:HD23	1:A:479:VAL:O	2.18	0.43
2:H:265:LEU:HD23	2:H:276:VAL:CG1	2.48	0.43
1:C:437:LYS:HD2	1:C:437:LYS:HA	1.90	0.43
1:E:411:GLU:OE1	1:E:411:GLU:HA	2.18	0.43
1:G:264:ASN:N	1:G:264:ASN:HD22	2.16	0.43
1:G:264:ASN:CA	1:G:333:ILE:HD11	2.48	0.43
3:L:145:TYR:HB2	3:L:167:LEU:HD22	2.00	0.43
1:A:366:GLN:C	1:A:368:ILE:N	2.70	0.43
1:G:164:ILE:HD11	1:G:508:ILE:HD11	2.00	0.43
2:F:139:HIS:O	2:F:176:PRO:HD2	2.19	0.43
2:B:226:GLU:CG	2:B:280:LEU:HD11	2.48	0.43
1:C:349:LYS:O	1:C:350:ALA:C	2.57	0.43
2:D:98:ASP:O	2:D:99:ILE:C	2.56	0.43
1:A:159:VAL:HG21	1:A:425:LEU:HD22	2.00	0.43
2:D:106:VAL:O	2:D:109:GLU:HG2	2.18	0.43
1:A:189:PHE:CE1	1:A:192:LEU:HB2	2.54	0.43
1:G:47:LYS:C	1:G:47:LYS:HD2	2.39	0.43
1:A:311:LEU:HD12	1:A:311:LEU:HA	1.89	0.43
2:D:237:TRP:O	2:D:241:GLN:HG2	2.18	0.43
2:F:249:LEU:HD13	2:F:260:ILE:HD11	1.99	0.43
1:G:516:PHE:N	1:G:516:PHE:CD2	2.86	0.43
2:B:312:THR:C	2:B:314:ALA:N	2.72	0.43
3:J:123:VAL:HG13	3:J:150:MET:HE3	1.99	0.43
2:H:393:TYR:CG	2:H:394:LEU:N	2.86	0.43
2:B:418:VAL:HG22	2:B:421:GLN:OE1	2.19	0.43
2:D:419:ASP:CG	2:D:420:GLY:H	2.21	0.43
2:B:351:GLN:N	2:B:436:PHE:HA	2.33	0.43
1:A:429:LEU:HD23	1:A:429:LEU:HA	1.85	0.43
2:D:361:LEU:C	2:D:363:GLU:N	2.70	0.43
2:F:412:LEU:O	2:F:417:LEU:HB2	2.18	0.43
3:L:120:THR:OG1	3:L:155:THR:OG1	2.37	0.43
2:D:124:PHE:C	2:D:125:ASN:ND2	2.72	0.43
2:B:391:THR:O	2:B:392:LEU:C	2.56	0.43
2:F:401:GLU:O	2:F:405:ARG:HB2	2.19	0.43
2:B:249:LEU:HD12	2:B:250:ASP:N	2.33	0.43
1:E:28:LEU:HD12	1:E:28:LEU:HA	1.81	0.43
2:B:407:ASN:HD22	2:B:415:LEU:HD13	1.83	0.43
2:H:348:GLN:C	2:H:348:GLN:HE21	2.22	0.43
2:B:253:ASP:O	2:B:256:HIS:HB2	2.19	0.43
2:H:92:PHE:N	2:H:92:PHE:CD2	2.80	0.43
1:E:76:LEU:HD21	1:E:89:ALA:HB2	1.99	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:G:129:THR:HA	1:G:153:CYS:O	2.19	0.43
2:H:398:THR:O	2:H:401:GLU:HB3	2.19	0.43
1:E:447:ASN:ND2	2:F:26:ARG:NE	2.65	0.43
2:H:46:LEU:O	2:H:73:ARG:HB3	2.19	0.43
1:E:396:ARG:HH21	1:E:400:GLU:HB3	1.82	0.43
2:B:166:ASP:C	2:B:168:VAL:H	2.22	0.43
1:G:157:GLY:HA3	1:G:485:TYR:CD1	2.54	0.43
2:B:331:TYR:CE1	3:I:110:GLY:HA2	2.54	0.43
1:E:421:ASN:O	1:E:424:VAL:HG22	2.19	0.43
2:D:243:PHE:CD1	2:D:243:PHE:N	2.87	0.43
2:F:357:PRO:HB3	2:F:440:PHE:CZ	2.54	0.43
2:F:231:TYR:O	2:F:233:ARG:N	2.52	0.43
1:A:225:TRP:O	1:A:229:THR:HB	2.18	0.43
2:H:233:ARG:HG2	2:H:233:ARG:HH11	1.84	0.43
1:E:214:TRP:CH2	1:E:332:MET:HG2	2.54	0.43
2:H:200:ILE:HD13	3:L:172:ARG:NH2	2.34	0.43
2:D:325:ASN:OD1	2:D:327:VAL:HG22	2.19	0.43
1:G:31:ALA:HB3	1:G:54:ILE:HD11	2.01	0.43
2:B:385:LEU:HA	2:B:385:LEU:HD23	1.53	0.43
1:C:435:PHE:CE2	1:C:439:GLN:HG3	2.54	0.43
1:A:364:LEU:HD12	1:A:364:LEU:HA	1.80	0.43
2:B:128:GLN:HB2	2:B:128:GLN:HE21	1.63	0.43
2:D:76:HIS:CD2	2:D:138:PHE:CE2	3.07	0.43
2:D:139:HIS:O	2:D:176:PRO:HD2	2.19	0.43
1:E:56:SER:HB3	1:E:101:SER:HB2	2.00	0.43
3:I:157:ALA:C	3:I:159:TYR:H	2.22	0.43
2:D:422:GLU:HA	2:D:437:LYS:HA	2.01	0.43
1:G:110:GLU:OE2	1:G:138:ARG:NH1	2.52	0.43
2:B:356:SER:HA	2:B:441:THR:H	1.84	0.42
1:A:518:ASN:HD21	1:A:534:LEU:CA	2.31	0.42
3:I:170:VAL:CG1	3:I:171:LEU:N	2.82	0.42
2:B:217:THR:CB	2:B:223:ARG:NH2	2.79	0.42
2:H:272:ASN:OD1	2:H:272:ASN:N	2.49	0.42
2:D:90:ARG:HG2	2:D:90:ARG:HH11	1.84	0.42
2:F:405:ARG:N	2:F:406:PRO:CD	2.82	0.42
1:A:430:ARG:O	1:A:433:ASP:HB2	2.19	0.42
1:A:140:ALA:CB	1:A:398:LEU:HD23	2.48	0.42
2:B:156:GLY:HA3	2:B:430:THR:CG2	2.49	0.42
2:D:390:ARG:CG	2:D:391:THR:N	2.80	0.42
2:F:256:HIS:CD2	2:F:256:HIS:N	2.87	0.42
1:A:481:GLU:O	1:A:484:ARG:HB3	2.18	0.42
2:D:351:GLN:O	2:D:436:PHE:HA	2.19	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:181:GLU:OE1	1:A:330:PRO:HD3	2.19	0.42
1:E:120:PHE:HA	1:E:123:PHE:CD1	2.54	0.42
1:A:139:LEU:HD12	1:A:143:LEU:HG	2.01	0.42
1:E:190:PRO:O	1:E:191:GLU:C	2.58	0.42
1:A:279:THR:O	1:A:280:GLN:HB2	2.17	0.42
1:G:171:VAL:HG12	1:G:391:ARG:O	2.19	0.42
1:C:474:VAL:O	1:C:475:LYS:C	2.57	0.42
1:A:260:GLU:O	1:A:261:ASP:OD1	2.37	0.42
1:A:403:GLY:HA3	1:A:406:THR:OG1	2.19	0.42
2:D:405:ARG:HB3	2:D:406:PRO:HD3	2.00	0.42
1:E:279:THR:O	1:E:280:GLN:HB2	2.20	0.42
1:G:447:ASN:ND2	2:H:26:ARG:HE	2.17	0.42
1:C:409:LYS:O	1:C:410:ASP:C	2.56	0.42
1:A:248:ARG:O	1:A:251:ILE:HG13	2.20	0.42
1:G:193:ARG:C	1:G:195:HIS:N	2.73	0.42
2:F:148:ILE:O	2:F:152:ARG:HG3	2.20	0.42
1:C:25:GLN:HE21	1:C:29:GLU:CG	2.31	0.42
1:A:171:VAL:HG13	1:A:391:ARG:HB2	2.01	0.42
2:H:226:GLU:N	2:H:226:GLU:OE1	2.53	0.42
2:B:151:ARG:HG3	2:B:151:ARG:HH11	1.83	0.42
2:D:321:TYR:OH	3:J:172:ARG:HA	2.20	0.42
1:A:396:ARG:NH2	1:A:406:THR:O	2.47	0.42
1:A:518:ASN:CG	1:A:519:THR:N	2.65	0.42
1:A:238:LYS:HD3	1:A:238:LYS:O	2.19	0.42
2:B:221:MET:HE2	2:B:221:MET:HB3	1.91	0.42
1:E:48:ASN:O	1:E:52:PRO:HG3	2.19	0.42
2:B:415:LEU:HB2	2:B:417:LEU:HD23	2.01	0.42
2:B:417:LEU:HD22	2:B:417:LEU:N	2.35	0.42
2:D:286:LYS:O	2:D:287:ARG:C	2.57	0.42
2:D:399:SER:HB2	2:D:403:ARG:HH12	1.83	0.42
1:A:225:TRP:CE2	1:A:234:PRO:HG3	2.54	0.42
1:G:318:GLU:OE1	1:G:360:HIS:NE2	2.52	0.42
1:C:71:GLY:HA3	2:D:16:ARG:O	2.19	0.42
1:C:446:SER:HB2	1:C:449:GLN:CB	2.50	0.42
1:C:354:ALA:O	1:C:380:LEU:HD21	2.18	0.42
2:F:204:LEU:HD21	2:F:375:GLN:HG2	2.01	0.42
1:G:514:VAL:HB	2:H:328:ASP:OD1	2.20	0.42
1:G:236:THR:O	1:G:240:LYS:HG3	2.19	0.42
1:C:377:GLU:O	1:C:381:LYS:HG2	2.19	0.42
1:G:484:ARG:NH1	2:H:315:TYR:CZ	2.88	0.42
2:H:260:ILE:O	2:H:264:SER:HB2	2.20	0.42
2:B:230:GLU:CD	2:B:233:ARG:HH12	2.22	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:299:LYS:HA	1:C:368:ILE:CG2	2.43	0.42
1:A:165:ILE:HD13	1:A:398:LEU:HB2	2.01	0.42
2:D:32:HIS:HB3	2:D:313:SER:O	2.19	0.42
3:J:137:PRO:HA	3:J:138:PRO:HD3	1.89	0.42
1:C:160:GLY:HA3	1:C:497:LEU:HD11	2.01	0.42
3:I:103:ILE:HD11	3:I:117:ILE:CD1	2.49	0.42
2:H:169:LEU:HD12	2:H:170:ASP:N	2.34	0.42
2:D:217:THR:CG2	2:D:223:ARG:HH22	2.33	0.42
2:H:423:LEU:HD11	2:H:438:LEU:HB2	2.02	0.42
1:E:504:GLU:O	1:E:508:ILE:HG13	2.20	0.42
1:E:352:LYS:O	1:E:355:ALA:N	2.52	0.42
1:A:89:ALA:O	1:A:90:MET:C	2.58	0.42
1:A:236:THR:HB	1:A:239:GLU:H	1.85	0.42
2:H:425:VAL:O	2:H:434:VAL:HG12	2.19	0.42
1:E:211:HIS:HE1	2:F:221:MET:CE	2.33	0.42
3:L:154:LYS:HB3	3:L:158:ASP:OD1	2.19	0.42
1:E:259:PRO:HB2	1:E:261:ASP:O	2.20	0.42
1:A:318:GLU:N	1:A:318:GLU:OE1	2.50	0.42
1:C:416:MET:CE	1:C:419:PRO:HA	2.50	0.42
1:G:214:TRP:H	1:G:332:MET:HE2	1.85	0.42
1:G:53:GLY:O	1:G:54:ILE:C	2.58	0.42
2:D:316:ILE:HA	2:D:317:PRO:HD3	1.84	0.42
1:G:299:LYS:HA	1:G:368:ILE:HG21	2.01	0.42
3:I:102:LEU:CG	3:I:102:LEU:O	2.67	0.42
3:J:169:LEU:HD23	3:J:169:LEU:HA	1.87	0.42
1:C:252:LEU:HD23	1:C:252:LEU:HA	1.77	0.42
1:A:307:LEU:HD21	1:A:375:ILE:HG21	2.02	0.42
2:H:356:SER:OG	2:H:358:SER:HB3	2.19	0.42
2:F:217:THR:CG2	2:F:223:ARG:NH2	2.79	0.42
2:B:351:GLN:HB3	2:B:436:PHE:CD1	2.54	0.42
1:G:311:LEU:HD21	1:G:387:SER:HB2	2.02	0.42
3:K:107:THR:HB	3:K:111:LYS:HB3	2.02	0.42
2:D:316:ILE:CD1	2:D:316:ILE:H	2.32	0.42
2:D:359:ALA:HB1	2:D:363:GLU:CD	2.40	0.42
1:G:225:TRP:CD1	1:G:225:TRP:C	2.92	0.42
3:I:137:PRO:O	3:I:140:GLN:HG2	2.20	0.42
2:F:127:ILE:O	2:F:135:TYR:OH	2.38	0.42
1:E:428:MET:CE	1:E:479:VAL:HA	2.50	0.42
2:H:30:PHE:N	2:H:30:PHE:CD1	2.87	0.42
2:F:212:ASN:HA	2:F:212:ASN:HD22	1.58	0.42
2:B:357:PRO:O	2:B:411:THR:HB	2.20	0.42
2:F:393:TYR:HD1	2:F:404:THR:OG1	2.03	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:417:LEU:HD12	2:B:421:GLN:HE22	1.84	0.42
1:A:96:LEU:HD23	2:B:95:ARG:NH2	2.34	0.42
1:E:214:TRP:CZ3	1:E:332:MET:HG2	2.54	0.42
1:G:185:LEU:O	1:G:188:PRO:HD3	2.19	0.42
2:D:46:LEU:HB3	2:D:73:ARG:NH1	2.34	0.42
2:D:412:LEU:HD12	2:D:440:PHE:HE2	1.84	0.42
2:B:381:ILE:CG2	2:B:423:LEU:HD22	2.50	0.42
2:H:178:ILE:N	2:H:178:ILE:CD1	2.82	0.42
1:G:323:LEU:HB3	1:G:324:PRO:HD2	2.02	0.42
1:C:189:PHE:O	1:C:192:LEU:N	2.52	0.42
1:E:396:ARG:HD3	1:E:534:LEU:O	2.20	0.42
1:C:446:SER:HB2	1:C:449:GLN:HE21	1.84	0.42
1:A:457:LEU:HD22	1:A:483:CYS:SG	2.60	0.42
2:B:52:LEU:HD11	2:B:78:ILE:HG13	2.02	0.42
1:G:282:PRO:HG2	1:G:285:ILE:HD13	2.02	0.42
1:C:359:ASN:O	1:C:363:LYS:HG3	2.20	0.42
2:B:441:THR:O	2:B:442:SER:OXT	2.38	0.42
2:H:318:LEU:O	2:H:319:ASN:C	2.55	0.42
2:B:419:ASP:OD1	2:B:440:PHE:CD1	2.62	0.42
1:A:78:ARG:HB2	1:A:81:ILE:HD11	2.02	0.42
1:E:215:ILE:HD12	1:E:342:LEU:HD21	2.01	0.42
2:D:237:TRP:C	2:D:239:LYS:H	2.23	0.42
2:B:178:ILE:N	2:B:178:ILE:HD12	2.34	0.42
1:G:362:ALA:O	1:G:366:GLN:HG3	2.20	0.42
2:F:52:LEU:HD11	2:F:78:ILE:HG13	2.01	0.42
2:H:193:LEU:HD23	2:H:193:LEU:HA	1.74	0.42
1:G:481:GLU:HA	1:G:481:GLU:OE1	2.20	0.42
3:I:107:THR:CG2	3:I:109:THR:HG23	2.50	0.41
1:C:143:LEU:HD22	1:C:148:ILE:CB	2.45	0.41
1:A:84:ASN:HB3	1:A:87:GLU:HB3	2.02	0.41
2:F:154:ILE:HG13	2:F:154:ILE:O	2.20	0.41
1:C:157:GLY:HA3	1:C:485:TYR:CG	2.55	0.41
1:E:240:LYS:HG2	1:E:276:LEU:CD1	2.49	0.41
1:E:43:THR:CG2	1:E:44:GLU:N	2.83	0.41
2:D:64:LEU:HB3	2:D:111:LEU:HD13	2.02	0.41
1:A:42:GLY:HA2	1:A:129:THR:HG21	2.02	0.41
1:C:340:ILE:CD1	2:D:271:TYR:O	2.68	0.41
3:L:135:GLY:O	3:L:137:PRO:HD3	2.20	0.41
2:F:243:PHE:O	2:F:247:VAL:HG21	2.19	0.41
3:K:123:VAL:HG13	3:K:150:MET:HE2	2.02	0.41
1:G:264:ASN:HD22	1:G:265:PHE:N	2.18	0.41
1:E:187:LYS:N	1:E:188:PRO:CD	2.82	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:L:123:VAL:HG23	3:L:154:LYS:O	2.20	0.41
1:A:514:VAL:HB	2:B:328:ASP:HA	2.01	0.41
1:G:196:PHE:O	1:G:199:TYR:HB2	2.20	0.41
2:H:148:ILE:CG1	3:L:174:ARG:HG2	2.46	0.41
1:C:39:THR:O	1:C:43:THR:HB	2.19	0.41
1:G:238:LYS:O	1:G:241:GLU:HB3	2.20	0.41
1:A:426:TYR:O	1:A:429:LEU:HB2	2.19	0.41
1:E:282:PRO:HG2	1:E:388:ALA:HB1	2.03	0.41
3:L:106:LYS:CG	3:L:112:GLU:HG2	2.49	0.41
2:B:339:ARG:NH2	2:B:346:CYS:O	2.44	0.41
2:F:320:ASN:HB2	2:F:336:GLU:HA	2.02	0.41
1:A:416:MET:CE	1:A:424:VAL:HG22	2.50	0.41
3:L:151:ASN:C	3:L:153:GLU:H	2.23	0.41
2:F:143:CYS:HB2	2:F:179:ASP:HA	2.01	0.41
1:C:329:ILE:HB	1:C:330:PRO:HD2	2.02	0.41
2:F:75:ILE:O	2:F:120:VAL:HA	2.19	0.41
2:F:357:PRO:HD3	2:F:440:PHE:CD2	2.54	0.41
2:B:207:TYR:CE2	3:I:172:ARG:CG	2.91	0.41
2:B:325:ASN:HD22	2:B:326:ASP:N	2.18	0.41
2:D:90:ARG:HG3	2:D:91:GLN:HG3	2.02	0.41
2:B:132:ASP:O	2:B:134:PHE:N	2.53	0.41
2:D:277:THR:HG22	2:D:279:ARG:H	1.85	0.41
2:H:217:THR:HG21	2:H:223:ARG:NE	2.31	0.41
1:G:333:ILE:CG2	2:H:223:ARG:NH1	2.83	0.41
2:F:241:GLN:CB	2:F:245:GLU:HA	2.51	0.41
1:E:364:LEU:HA	1:E:364:LEU:HD12	1.84	0.41
2:H:427:ASP:OD2	2:H:428:VAL:N	2.49	0.41
2:H:220:SER:O	2:H:222:PRO:CD	2.68	0.41
1:A:186:ASP:OD2	1:A:187:LYS:N	2.53	0.41
1:C:212:THR:HA	1:C:213:PRO:HD3	1.94	0.41
3:L:155:THR:HG22	3:L:158:ASP:CG	2.41	0.41
1:E:311:LEU:HD22	1:E:383:LEU:HD11	2.02	0.41
2:D:418:VAL:CG2	2:D:421:GLN:NE2	2.80	0.41
2:H:354:GLN:NE2	2:H:439:HIS:HB3	2.34	0.41
2:D:268:ALA:O	2:D:273:ILE:N	2.53	0.41
3:J:125:ARG:HG3	3:J:125:ARG:NH1	2.35	0.41
1:C:47:LYS:NZ	2:D:65:LYS:HE2	2.36	0.41
1:G:344:ASN:O	1:G:348:GLU:HB2	2.20	0.41
1:C:314:PHE:HZ	1:C:353:ASP:OD2	2.04	0.41
1:C:245:ASP:O	1:C:249:GLN:HG3	2.21	0.41
1:A:335:ASP:O	1:A:336:SER:C	2.59	0.41
2:H:249:LEU:CD2	2:H:260:ILE:HD11	2.50	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:I:117:ILE:HD13	3:I:126:ILE:HG12	2.03	0.41
1:A:276:LEU:HD23	1:A:276:LEU:HA	1.84	0.41
1:A:428:MET:SD	1:A:461:LEU:HD22	2.61	0.41
2:H:349:LEU:O	2:H:350:PRO:O	2.38	0.41
1:A:154:ARG:NH1	1:A:154:ARG:HG2	2.36	0.41
1:C:40:ALA:O	1:C:43:THR:HG22	2.21	0.41
2:B:434:VAL:HG22	2:B:436:PHE:CE1	2.55	0.41
3:L:125:ARG:C	3:L:127:LYS:N	2.72	0.41
2:H:35:PHE:O	2:H:36:GLU:HG2	2.19	0.41
2:F:271:TYR:O	2:F:272:ASN:OD1	2.37	0.41
1:E:501:ALA:O	1:E:505:VAL:HG23	2.20	0.41
1:G:282:PRO:HG2	1:G:285:ILE:HB	2.02	0.41
1:E:419:PRO:HB2	1:E:475:LYS:HE3	2.03	0.41
2:F:79:ASP:OD2	2:F:81:ASP:HB2	2.20	0.41
1:E:241:GLU:OE1	1:E:244:ARG:HD2	2.21	0.41
1:G:497:LEU:HD12	1:G:497:LEU:HA	1.92	0.41
1:G:333:ILE:CG2	2:H:223:ARG:HH12	2.34	0.41
1:G:333:ILE:HD13	1:G:333:ILE:HG21	1.74	0.41
1:A:297:ILE:N	1:A:297:ILE:CD1	2.76	0.41
2:D:228:CYS:SG	2:D:267:ARG:HG2	2.61	0.41
1:G:311:LEU:HA	1:G:311:LEU:HD12	1.83	0.41
2:F:261:PHE:CZ	2:F:265:LEU:HD11	2.55	0.41
2:B:50:LYS:H	2:B:139:HIS:CD2	2.38	0.41
1:E:318:GLU:HG3	1:E:356:ALA:HB1	2.02	0.41
2:F:136:ARG:HH11	2:F:136:ARG:HG3	1.85	0.41
1:A:37:ASN:HB2	1:A:129:THR:O	2.21	0.41
2:D:385:LEU:O	2:D:387:GLY:N	2.52	0.41
1:C:499:GLY:O	2:D:294:SER:HB2	2.20	0.41
2:H:204:LEU:O	2:H:204:LEU:HG	2.20	0.41
1:G:15:ARG:HD3	1:G:15:ARG:HA	1.92	0.41
1:A:261:ASP:O	1:A:262:GLU:CG	2.69	0.41
2:D:333:TYR:CE2	2:D:335:PHE:HB3	2.56	0.41
1:G:248:ARG:O	1:G:251:ILE:HD12	2.21	0.41
2:D:231:TYR:HD1	2:D:235:LEU:HD12	1.85	0.41
1:C:481:GLU:HG3	1:C:485:TYR:CE2	2.56	0.41
1:E:518:ASN:CB	1:E:533:GLN:HA	2.50	0.41
3:I:102:LEU:O	3:I:163:GLY:O	2.39	0.41
2:B:309:LYS:HD3	2:B:317:PRO:HA	2.03	0.41
1:A:353:ASP:O	1:A:357:VAL:HG23	2.20	0.41
2:F:247:VAL:HA	2:F:248:PRO:HD3	1.81	0.41
1:C:289:PHE:HB3	1:C:312:LYS:HD2	2.03	0.41
2:F:209:PRO:HG2	3:K:139:GLN:O	2.20	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:97:ASN:HB3	1:A:100:VAL:HG23	2.02	0.41
1:G:66:SER:N	1:G:69:ASP:HB2	2.35	0.41
1:A:359:ASN:HD22	1:A:359:ASN:HA	1.60	0.41
3:L:131:GLU:HG2	3:L:131:GLU:O	2.19	0.41
1:C:163:ARG:HH12	1:C:165:ILE:HG12	1.81	0.41
2:H:385:LEU:CD1	2:H:390:ARG:HD3	2.51	0.41
2:F:410:LYS:HD3	2:F:415:LEU:HG	2.03	0.41
2:F:74:GLN:NE2	2:F:74:GLN:HA	2.36	0.41
3:L:120:THR:OG1	3:L:155:THR:CB	2.68	0.41
1:G:422:GLU:O	1:G:424:VAL:N	2.54	0.41
2:D:141:ILE:HD12	2:D:158:LEU:CD2	2.41	0.41
2:B:16:ARG:NH2	2:B:116:PRO:HB2	2.36	0.41
2:H:171:PRO:C	2:H:173:SER:H	2.24	0.41
1:A:165:ILE:C	1:A:166:ILE:HG13	2.40	0.41
1:E:45:ILE:CG1	1:E:498:GLY:HA2	2.49	0.41
1:C:158:LEU:HD12	1:C:525:MET:HG2	2.01	0.41
1:E:35:LEU:HD12	1:E:36:ILE:N	2.36	0.41
1:G:72:ASN:C	1:G:72:ASN:ND2	2.74	0.41
1:E:274:THR:O	1:E:276:LEU:N	2.54	0.41
3:L:106:LYS:HB2	3:L:168:HIS:CD2	2.55	0.41
1:C:120:PHE:O	1:C:122:ARG:N	2.54	0.41
2:D:186:LYS:C	3:J:173:LEU:HD13	2.41	0.41
1:C:293:ARG:HB3	1:C:305:TRP:CE3	2.55	0.41
1:E:241:GLU:O	1:E:244:ARG:HB2	2.21	0.41
2:D:213:PHE:HB2	2:D:218:ILE:HD11	2.02	0.41
2:D:405:ARG:NH1	2:D:405:ARG:HB3	2.35	0.41
2:D:140:ILE:HG22	2:D:141:ILE:N	2.35	0.41
2:H:233:ARG:HG2	2:H:233:ARG:NH1	2.35	0.41
2:H:241:GLN:HA	2:H:242:PRO:HD3	1.57	0.41
2:D:15:GLY:O	2:D:17:TRP:N	2.54	0.41
3:K:101:MET:HE3	3:K:117:ILE:HG22	2.02	0.41
1:G:512:GLN:C	1:G:513:PHE:CD1	2.94	0.41
2:H:64:LEU:HD11	2:H:77:VAL:HG21	2.03	0.41
2:H:352:ASN:HD22	2:H:439:HIS:CD2	2.39	0.41
2:F:321:TYR:OH	3:K:172:ARG:HG3	2.21	0.41
2:D:249:LEU:CD1	2:D:260:ILE:HD11	2.50	0.41
2:D:253:ASP:OD1	2:D:255:GLU:HB2	2.20	0.41
1:E:245:ASP:C	1:E:247:ILE:N	2.74	0.41
1:C:9:LYS:HB3	1:C:9:LYS:HE2	1.90	0.41
2:D:32:HIS:CE1	2:D:33:PRO:HD2	2.56	0.41
2:B:257:ILE:HG22	2:B:278:TYR:CE1	2.56	0.41
3:I:102:LEU:O	3:I:102:LEU:HG	2.21	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:D:356:SER:C	2:D:358:SER:N	2.75	0.41
1:G:286:GLU:OE1	1:G:286:GLU:HA	2.20	0.41
1:E:474:VAL:O	1:E:475:LYS:C	2.59	0.41
1:A:510:THR:O	1:A:511:LYS:HB2	2.21	0.41
1:A:213:PRO:HG3	1:A:334:ALA:CB	2.50	0.41
3:K:131:GLU:CD	3:K:138:PRO:HD3	2.41	0.41
2:D:150:ALA:O	2:D:154:ILE:HG22	2.21	0.41
2:H:115:VAL:HA	2:H:116:PRO:HD2	1.88	0.41
1:G:437:LYS:HD2	1:G:437:LYS:HA	1.90	0.41
2:F:219:ALA:HB2	2:F:231:TYR:CD1	2.56	0.40
1:A:140:ALA:HB3	1:A:398:LEU:HD23	2.03	0.40
1:A:371:ALA:C	1:A:373:GLU:N	2.75	0.40
1:E:528:THR:HG22	2:F:334:THR:OG1	2.20	0.40
2:F:356:SER:C	2:F:358:SER:N	2.75	0.40
2:B:265:LEU:HD13	1:E:66:SER:HA	2.02	0.40
1:A:424:VAL:HG11	1:A:478:TYR:CD1	2.56	0.40
2:F:95:ARG:HB2	2:F:97:LYS:HG2	2.03	0.40
1:C:302:PRO:O	1:C:304:PHE:N	2.54	0.40
1:E:201:LEU:HG	1:E:201:LEU:H	1.34	0.40
2:D:323:VAL:HG21	3:J:170:VAL:HG23	2.03	0.40
1:A:430:ARG:NH2	1:A:464:PHE:HE1	2.18	0.40
1:A:248:ARG:HH11	1:A:248:ARG:HG3	1.87	0.40
1:C:9:LYS:HA	1:C:12:LYS:HB3	2.03	0.40
2:F:195:GLY:HA2	2:F:339:ARG:NH2	2.35	0.40
1:C:190:PRO:HG2	1:C:191:GLU:OE2	2.21	0.40
1:E:241:GLU:OE1	1:E:241:GLU:HA	2.21	0.40
1:E:434:ARG:HD3	1:E:460:CYS:O	2.21	0.40
2:H:143:CYS:HB3	2:H:179:ASP:OD2	2.20	0.40
1:E:129:THR:HA	1:E:153:CYS:O	2.21	0.40
1:C:261:ASP:O	1:C:262:GLU:OE2	2.39	0.40
1:C:161:TYR:OH	1:C:163:ARG:HB2	2.21	0.40
2:B:355:PHE:CE2	2:B:363:GLU:O	2.74	0.40
2:B:87:ASN:ND2	2:B:103:LYS:CE	2.85	0.40
2:D:15:GLY:O	2:D:16:ARG:C	2.59	0.40
1:G:243:PHE:CE2	1:G:247:ILE:HD11	2.56	0.40
1:G:111:ASN:O	1:G:114:ASP:HB3	2.22	0.40
1:G:158:LEU:HD13	1:G:158:LEU:HA	1.83	0.40
2:F:65:LYS:HE2	2:F:65:LYS:HB3	1.91	0.40
1:A:51:LEU:N	1:A:52:PRO:HD2	2.36	0.40
1:G:66:SER:H	1:G:69:ASP:HB2	1.86	0.40
2:H:28:GLY:O	2:H:31:THR:HG22	2.22	0.40
1:C:419:PRO:O	1:C:424:VAL:HG11	2.21	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:361:VAL:CG1	1:C:375:ILE:HD12	2.51	0.40
1:E:236:THR:C	1:E:238:LYS:N	2.74	0.40
1:G:65:VAL:HG23	1:G:83:LYS:O	2.21	0.40
2:D:351:GLN:HG2	2:D:352:ASN:N	2.36	0.40
2:D:270:GLN:C	2:D:272:ASN:N	2.75	0.40
2:H:302:VAL:HG11	2:H:322:LEU:HD23	2.04	0.40
1:C:215:ILE:HG13	1:C:332:MET:CE	2.52	0.40
1:E:33:VAL:HG12	1:E:34:CYS:N	2.36	0.40
1:E:266:GLU:C	1:E:268:ALA:H	2.23	0.40
1:E:189:PHE:CE1	1:E:349:LYS:HB2	2.56	0.40
1:G:441:ARG:NH2	1:G:453:ASP:OD2	2.55	0.40
1:E:90:MET:O	1:E:94:GLN:HB2	2.20	0.40
2:D:170:ASP:OD2	2:D:172:SER:HB2	2.21	0.40
1:E:251:ILE:HG13	1:E:251:ILE:H	1.44	0.40
2:F:351:GLN:HB2	2:F:436:PHE:HD2	1.77	0.40
2:D:21:LYS:O	2:D:25:GLU:HG3	2.22	0.40
2:D:325:ASN:HD21	2:D:327:VAL:CG2	2.29	0.40
1:G:251:ILE:C	1:G:252:LEU:O	2.58	0.40
1:C:309:ARG:HG3	1:C:309:ARG:HH11	1.86	0.40
1:A:454:ILE:HG21	1:A:480:HIS:CE1	2.56	0.40
2:F:335:PHE:CZ	3:K:144:ILE:HD13	2.57	0.40
1:A:215:ILE:HD11	1:A:332:MET:CE	2.51	0.40
1:E:229:THR:HG22	1:E:232:ARG:HB2	2.04	0.40
1:A:192:LEU:HD11	1:A:196:PHE:CZ	2.57	0.40
2:D:355:PHE:CZ	2:D:364:VAL:HA	2.57	0.40
1:E:7:LEU:HD23	1:E:7:LEU:O	2.20	0.40
2:B:243:PHE:N	2:B:243:PHE:CD1	2.90	0.40
2:D:94:PHE:N	2:D:94:PHE:CD2	2.87	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	513/531 (97%)	433 (84%)	64 (12%)	16 (3%)	7 34

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	512/531 (96%)	432 (84%)	59 (12%)	21 (4%)	4	25
1	E	511/531 (96%)	447 (88%)	46 (9%)	18 (4%)	6	30
1	G	502/531 (94%)	434 (86%)	54 (11%)	14 (3%)	8	37
2	B	430/434 (99%)	362 (84%)	56 (13%)	12 (3%)	8	37
2	D	429/434 (99%)	347 (81%)	62 (14%)	20 (5%)	4	22
2	F	429/434 (99%)	362 (84%)	51 (12%)	16 (4%)	5	29
2	H	429/434 (99%)	344 (80%)	60 (14%)	25 (6%)	3	16
3	I	81/88 (92%)	71 (88%)	5 (6%)	5 (6%)	2	14
3	J	75/88 (85%)	66 (88%)	7 (9%)	2 (3%)	8	38
3	K	75/88 (85%)	63 (84%)	8 (11%)	4 (5%)	3	19
3	L	76/88 (86%)	60 (79%)	12 (16%)	4 (5%)	3	19
All	All	4062/4212 (96%)	3421 (84%)	484 (12%)	157 (4%)	5	27

All (157) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	253	LYS
2	B	116	PRO
2	B	127	ILE
2	B	241	GLN
2	B	387	GLY
1	C	253	LYS
2	D	13	TRP
2	D	16	ARG
2	D	41	SER
2	D	92	PHE
2	D	116	PRO
2	D	320	ASN
1	E	253	LYS
1	E	275	ALA
1	E	276	LEU
2	F	116	PRO
2	F	132	ASP
2	F	375	GLN
1	G	276	LEU
2	H	116	PRO
2	H	132	ASP
2	H	242	PRO

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
2	H	320	ASN
2	H	350	PRO
2	H	377	LYS
2	H	396	SER
3	L	163	GLY
1	A	228	GLU
1	A	336	SER
1	A	367	SER
1	A	450	VAL
3	I	147	GLY
3	I	163	GLY
1	C	36	ILE
1	C	79	SER
1	C	249	GLN
1	C	275	ALA
1	C	451	GLU
2	D	270	GLN
2	D	271	TYR
1	E	195	HIS
1	E	200	ASP
1	E	280	GLN
2	F	320	ASN
1	G	171	VAL
1	G	335	ASP
1	G	451	GLU
1	G	453	ASP
1	G	475	LYS
2	H	88	LEU
2	H	244	GLY
2	H	319	ASN
2	H	362	GLN
2	H	399	SER
2	H	407	ASN
3	L	119	PRO
1	A	90	MET
1	A	106	GLU
1	A	280	GLN
2	B	92	PHE
2	B	133	THR
2	B	313	SER
2	B	403	ARG
3	I	119	PRO

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
3	I	158	ASP
1	C	121	CYS
1	C	252	LEU
1	C	475	LYS
1	C	518	ASN
2	D	29	PRO
2	D	30	PHE
2	D	237	TRP
2	D	362	GLN
1	E	134	SER
1	E	229	THR
1	E	249	GLN
1	E	264	ASN
1	E	318	GLU
1	E	336	SER
2	F	183	GLU
2	F	232	VAL
2	F	396	SER
2	F	410	LYS
3	K	119	PRO
3	K	152	ASP
1	G	452	GLU
2	H	194	PRO
2	H	232	VAL
2	H	239	LYS
2	H	265	LEU
2	H	345	ALA
2	H	357	PRO
3	L	152	ASP
1	A	79	SER
1	A	198	SER
1	A	230	ASN
1	A	435	PHE
2	B	361	LEU
3	I	146	SER
1	C	72	ASN
1	C	147	GLN
1	C	251	ILE
2	D	34	ASP
2	D	183	GLU
2	D	412	LEU
3	J	163	GLY

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	E	38	ALA
1	E	106	GLU
2	F	88	LEU
2	F	194	PRO
2	F	211	VAL
2	F	216	ALA
2	F	357	PRO
1	G	198	SER
1	G	275	ALA
1	G	387	SER
2	H	31	THR
2	H	374	LEU
2	H	388	LYS
1	A	249	GLN
2	B	242	PRO
1	C	106	GLU
1	C	303	SER
1	C	397	SER
2	D	241	GLN
1	E	190	PRO
2	F	92	PHE
3	K	108	LEU
3	K	175	GLY
1	G	250	GLY
2	H	261	PHE
1	A	372	PRO
2	B	348	GLN
1	C	259	PRO
1	E	221	TYR
1	E	228	GLU
1	G	407	ILE
1	C	109	PRO
2	D	357	PRO
2	F	144	GLY
1	C	407	ILE
2	D	232	VAL
1	E	213	PRO
1	G	423	ILE
2	H	405	ARG
1	C	450	VAL
2	D	211	VAL
1	G	419	PRO

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
2	H	238	PRO
1	A	333	ILE
1	A	443	PRO
2	B	167	GLY
1	C	60	ILE
2	D	194	PRO
3	L	175	GLY
3	J	119	PRO
2	F	229	ILE

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	449/462 (97%)	424 (94%)	25 (6%)	30	71
1	C	448/462 (97%)	425 (95%)	23 (5%)	33	75
1	E	448/462 (97%)	417 (93%)	31 (7%)	22	62
1	G	442/462 (96%)	419 (95%)	23 (5%)	32	74
2	B	378/382 (99%)	349 (92%)	29 (8%)	18	56
2	D	378/382 (99%)	352 (93%)	26 (7%)	22	62
2	F	377/382 (99%)	350 (93%)	27 (7%)	21	59
2	H	375/382 (98%)	348 (93%)	27 (7%)	21	59
3	I	73/74 (99%)	67 (92%)	6 (8%)	17	52
3	J	68/74 (92%)	66 (97%)	2 (3%)	55	89
3	K	68/74 (92%)	62 (91%)	6 (9%)	14	48
3	L	68/74 (92%)	67 (98%)	1 (2%)	76	95
All	All	3572/3672 (97%)	3346 (94%)	226 (6%)	25	66

All (226) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	ASP
1	A	39	THR

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	43	THR
1	A	69	ASP
1	A	72	ASN
1	A	90	MET
1	A	147	GLN
1	A	171	VAL
1	A	208	ASP
1	A	214	TRP
1	A	245	ASP
1	A	252	LEU
1	A	262	GLU
1	A	264	ASN
1	A	277	ASN
1	A	335	ASP
1	A	364	LEU
1	A	368	ILE
1	A	395	CYS
1	A	407	ILE
1	A	418	ASN
1	A	437	LYS
1	A	439	GLN
1	A	477	ASP
1	A	480	HIS
2	B	97	LYS
2	B	114	ARG
2	B	116	PRO
2	B	118	CYS
2	B	128	GLN
2	B	132	ASP
2	B	137	GLN
2	B	149	ILE
2	B	154	ILE
2	B	197	THR
2	B	217	THR
2	B	218	ILE
2	B	233	ARG
2	B	274	ARG
2	B	316	ILE
2	B	321	TYR
2	B	325	ASN
2	B	327	VAL
2	B	342	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
2	B	353	ILE
2	B	360	LYS
2	B	363	GLU
2	B	364	VAL
2	B	382	THR
2	B	392	LEU
2	B	412	LEU
2	B	414	GLU
2	B	430	THR
2	B	442	SER
3	I	97	SER
3	I	101	MET
3	I	102	LEU
3	I	107	THR
3	I	144	ILE
3	I	170	VAL
1	C	10	GLU
1	C	14	ASP
1	C	34	CYS
1	C	37	ASN
1	C	39	THR
1	C	112	LEU
1	C	171	VAL
1	C	191	GLU
1	C	214	TRP
1	C	251	ILE
1	C	264	ASN
1	C	274	THR
1	C	277	ASN
1	C	293	ARG
1	C	297	ILE
1	C	311	LEU
1	C	328	THR
1	C	353	ASP
1	C	396	ARG
1	C	418	ASN
1	C	422	GLU
1	C	439	GLN
1	C	517	ASN
2	D	11	LEU
2	D	13	TRP
2	D	14	GLU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
2	D	77	VAL
2	D	82	THR
2	D	97	LYS
2	D	111	LEU
2	D	113	ASP
2	D	114	ARG
2	D	116	PRO
2	D	125	ASN
2	D	128	GLN
2	D	132	ASP
2	D	142	VAL
2	D	149	ILE
2	D	196	MET
2	D	201	GLU
2	D	223	ARG
2	D	316	ILE
2	D	322	LEU
2	D	323	VAL
2	D	362	GLN
2	D	388	LYS
2	D	401	GLU
2	D	407	ASN
2	D	441	THR
3	J	108	LEU
3	J	155	THR
1	E	39	THR
1	E	43	THR
1	E	72	ASN
1	E	79	SER
1	E	101	SER
1	E	201	LEU
1	E	211	HIS
1	E	226	TYR
1	E	229	THR
1	E	242	ASP
1	E	251	ILE
1	E	253	LYS
1	E	262	GLU
1	E	264	ASN
1	E	265	PHE
1	E	277	ASN
1	E	292	ASP

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	E	293	ARG
1	E	297	ILE
1	E	309	ARG
1	E	364	LEU
1	E	410	ASP
1	E	418	ASN
1	E	422	GLU
1	E	424	VAL
1	E	452	GLU
1	E	497	LEU
1	E	516	PHE
1	E	517	ASN
1	E	518	ASN
1	E	533	GLN
2	F	31	THR
2	F	38	SER
2	F	82	THR
2	F	95	ARG
2	F	110	PHE
2	F	111	LEU
2	F	114	ARG
2	F	116	PRO
2	F	118	CYS
2	F	125	ASN
2	F	128	GLN
2	F	149	ILE
2	F	182	THR
2	F	194	PRO
2	F	212	ASN
2	F	241	GLN
2	F	292	VAL
2	F	316	ILE
2	F	319	ASN
2	F	322	LEU
2	F	325	ASN
2	F	342	ASN
2	F	355	PHE
2	F	361	LEU
2	F	374	LEU
2	F	397	VAL
2	F	430	THR
3	K	101	MET

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
3	K	115	ILE
3	K	117	ILE
3	K	155	THR
3	K	170	VAL
3	K	173	LEU
1	G	33	VAL
1	G	43	THR
1	G	72	ASN
1	G	73	ASN
1	G	79	SER
1	G	90	MET
1	G	154	ARG
1	G	163	ARG
1	G	186	ASP
1	G	189	PHE
1	G	214	TRP
1	G	220	LYS
1	G	264	ASN
1	G	277	ASN
1	G	293	ARG
1	G	300	GLN
1	G	318	GLU
1	G	343	GLN
1	G	385	SER
1	G	410	ASP
1	G	492	THR
1	G	503	GLN
1	G	516	PHE
2	H	14	GLU
2	H	49	CYS
2	H	82	THR
2	H	105	GLU
2	H	111	LEU
2	H	114	ARG
2	H	116	PRO
2	H	128	GLN
2	H	132	ASP
2	H	139	HIS
2	H	154	ILE
2	H	182	THR
2	H	201	GLU
2	H	213	PHE

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
2	H	224	LEU
2	H	316	ILE
2	H	323	VAL
2	H	325	ASN
2	H	348	GLN
2	H	350	PRO
2	H	360	LYS
2	H	366	ASP
2	H	374	LEU
2	H	402	GLU
2	H	403	ARG
2	H	405	ARG
2	H	430	THR
3	L	150	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (138) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	32	HIS
1	A	37	ASN
1	A	72	ASN
1	A	111	ASN
1	A	147	GLN
1	A	169	HIS
1	A	178	ASN
1	A	197	GLN
1	A	211	HIS
1	A	224	GLN
1	A	264	ASN
1	A	277	ASN
1	A	320	GLN
1	A	359	ASN
1	A	418	ASN
1	A	439	GLN
1	A	480	HIS
1	A	512	GLN
1	A	518	ASN
2	B	19	HIS
2	B	66	ASN
2	B	74	GLN
2	B	137	GLN
2	B	210	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
2	B	236	GLN
2	B	319	ASN
2	B	325	ASN
2	B	342	ASN
2	B	351	GLN
2	B	389	ASN
2	B	395	GLN
2	B	407	ASN
3	I	140	GLN
1	C	37	ASN
1	C	197	GLN
1	C	224	GLN
1	C	230	ASN
1	C	264	ASN
1	C	320	GLN
1	C	322	ASN
1	C	344	ASN
1	C	359	ASN
1	C	386	ASN
1	C	418	ASN
1	C	436	HIS
1	C	439	GLN
1	C	447	ASN
1	C	512	GLN
1	C	517	ASN
2	D	18	ASN
2	D	74	GLN
2	D	125	ASN
2	D	128	GLN
2	D	131	ASN
2	D	139	HIS
2	D	210	GLN
2	D	236	GLN
2	D	270	GLN
2	D	325	ASN
2	D	348	GLN
2	D	362	GLN
2	D	421	GLN
2	D	432	GLN
1	E	72	ASN
1	E	115	ASN
1	E	178	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	E	197	GLN
1	E	211	HIS
1	E	264	ASN
1	E	271	ASN
1	E	277	ASN
1	E	359	ASN
1	E	370	GLN
1	E	418	ASN
1	E	436	HIS
1	E	439	GLN
1	E	447	ASN
1	E	466	GLN
1	E	518	ASN
2	F	18	ASN
2	F	19	HIS
2	F	32	HIS
2	F	74	GLN
2	F	125	ASN
2	F	128	GLN
2	F	131	ASN
2	F	212	ASN
2	F	236	GLN
2	F	241	GLN
2	F	319	ASN
2	F	320	ASN
2	F	325	ASN
2	F	342	ASN
2	F	351	GLN
2	F	362	GLN
2	F	370	ASN
2	F	395	GLN
2	F	407	ASN
1	G	37	ASN
1	G	64	GLN
1	G	72	ASN
1	G	197	GLN
1	G	230	ASN
1	G	264	ASN
1	G	271	ASN
1	G	277	ASN
1	G	300	GLN
1	G	322	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	G	343	GLN
1	G	344	ASN
1	G	359	ASN
1	G	366	GLN
1	G	386	ASN
1	G	438	GLN
1	G	439	GLN
1	G	447	ASN
1	G	518	ASN
1	G	527	GLN
2	H	18	ASN
2	H	19	HIS
2	H	32	HIS
2	H	74	GLN
2	H	87	ASN
2	H	125	ASN
2	H	128	GLN
2	H	131	ASN
2	H	163	ASN
2	H	236	GLN
2	H	258	GLN
2	H	325	ASN
2	H	348	GLN
2	H	351	GLN
2	H	352	ASN
2	H	354	GLN
2	H	362	GLN
2	H	395	GLN
2	H	407	ASN
3	L	141	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, 4 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.

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 ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	517/531 (97%)	-0.12	2 (0%) 90 42	46, 76, 107, 130	0
1	C	516/531 (97%)	-0.08	0 100 100	53, 82, 116, 140	0
1	E	515/531 (96%)	-0.10	2 (0%) 90 42	36, 62, 128, 141	0
1	G	508/531 (95%)	-0.08	1 (0%) 93 56	44, 71, 135, 147	0
2	B	432/434 (99%)	-0.07	3 (0%) 84 30	46, 68, 122, 131	0
2	D	431/434 (99%)	0.02	1 (0%) 93 56	47, 85, 118, 131	0
2	F	431/434 (99%)	-0.01	0 100 100	41, 70, 121, 134	0
2	H	431/434 (99%)	-0.02	4 (0%) 81 24	45, 79, 127, 140	0
3	I	85/88 (96%)	-0.07	0 100 100	54, 77, 99, 110	0
3	J	77/88 (87%)	0.12	0 100 100	59, 90, 106, 116	0
3	K	77/88 (87%)	-0.10	0 100 100	56, 81, 99, 104	0
3	L	78/88 (88%)	0.05	0 100 100	70, 105, 125, 129	0
All	All	4098/4212 (97%)	-0.06	13 (0%) 91 50	36, 76, 122, 147	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	209	HIS	3.3
1	G	252	LEU	2.7
1	A	260	GLU	2.4
1	E	253	LYS	2.3
2	D	390	ARG	2.3
1	E	260	GLU	2.3
2	H	408	LEU	2.2
2	B	413	LYS	2.2
2	H	385	LEU	2.1
2	B	408	LEU	2.1
2	B	392	LEU	2.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
2	H	388	LYS	2.1
2	H	239	LYS	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(Å ²)	Q<0.9
4	ZN	D	3	1/1	0.13	-0.56	67,67,67,67	0
4	ZN	H	2	1/1	0.16	-0.65	84,84,84,84	0
4	ZN	B	1	1/1	0.14	-0.65	70,70,70,70	0
4	ZN	F	4	1/1	0.14	-0.72	67,67,67,67	0

6.5 Other polymers ⓘ

There are no such residues in this entry.