

hbond.lol

64 ALA H	61 PRO O	1.7
64 ALA N	61 PRO O	2.7
65 ALA H	119 PHE O	1.7
65 ALA N	119 PHE O	2.7
73 LEU H	85 LEU O	1.7
73 LEU N	85 LEU O	2.7
74 GLN H	168 LYS O	1.7
74 GLN N	168 LYS O	2.7
82 LEU H	75 CYS O	1.7
82 LEU N	75 CYS O	2.7
85 LEU H	73 LEU O	1.7
85 LEU N	73 LEU O	2.7
93 TYR H	117 LEU O	1.7
93 TYR N	117 LEU O	2.7
94 LEU H	102 ASP OD2	1.7
94 LEU N	102 ASP OD2	2.7
95 PHE H	115 ALA O	1.7
95 PHE N	115 ALA O	2.7
96 GLY H	103 TYR O	1.7
96 GLY N	103 TYR O	2.7
97 ARG H	111 SER O	1.7
97 ARG N	111 SER O	2.7
98 SER H	112 SER O	1.7
98 SER N	112 SER O	2.7
101 CYS H	98 SER O	1.7
101 CYS N	98 SER O	2.7
102 ASP H	94 LEU O	1.7
102 ASP N	94 LEU O	2.7
103 TYR H	94 LEU O	1.7
103 TYR N	94 LEU O	2.7
105 LEU H	96 GLY O	1.7
105 LEU N	96 GLY O	2.7
110 ILE H	107 HIS O	1.7
110 ILE N	107 HIS O	2.7
111 SER H	114 HIS ND1	1.7
111 SER N	114 HIS ND1	2.7
114 HIS H	95 PHE O	1.7
114 HIS N	95 PHE O	2.7
115 ALA H	95 PHE O	1.7
115 ALA N	95 PHE O	2.7
116 VAL H	129 MET O	1.7
116 VAL N	129 MET O	2.7
117 LEU H	93 TYR O	1.7
117 LEU N	93 TYR O	2.7
118 VAL H	127 VAL O	1.7
118 VAL N	127 VAL O	2.7
119 PHE H	91 PRO O	1.7
119 PHE N	91 PRO O	2.7
120 HIS H	125 CYS O	1.7
120 HIS N	125 CYS O	2.7
121 GLY H	63 TRP O	1.7
121 GLY N	63 TRP O	2.7
127 VAL H	118 VAL O	1.7
127 VAL N	118 VAL O	2.7
128 LEU H	150 LEU O	1.7
128 LEU N	150 LEU O	2.7
129 MET H	116 VAL O	1.7
129 MET N	116 VAL O	2.7
131 LEU H	114 HIS O	1.7
131 LEU N	114 HIS O	2.7
138 LYS H	159 GLN O	1.7
138 LYS N	159 GLN O	2.7

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64 ALA H	61 PRO O	2.6
64 ALA N	61 PRO O	3.5
65 ALA H	119 PHE O	2.6
65 ALA N	119 PHE O	3.5
73 LEU H	85 LEU O	2.6
73 LEU N	85 LEU O	3.5
74 GLN H	168 LYS O	2.6
74 GLN N	168 LYS O	3.5
82 LEU H	75 CYS O	2.6
82 LEU N	75 CYS O	3.5
85 LEU H	73 LEU O	2.6
85 LEU N	73 LEU O	3.5
93 TYR H	117 LEU O	2.6
93 TYR N	117 LEU O	3.5
94 LEU H	102 ASP OD2	2.6
94 LEU N	102 ASP OD2	3.5
95 PHE H	115 ALA O	2.6
95 PHE N	115 ALA O	3.5
96 GLY H	103 TYR O	2.6
96 GLY N	103 TYR O	3.5
97 ARG H	111 SER O	2.6
97 ARG N	111 SER O	3.5
98 SER H	112 SER O	2.6
98 SER N	112 SER O	3.5
101 CYS H	98 SER O	2.6
101 CYS N	98 SER O	3.5
102 ASP H	94 LEU O	2.6
102 ASP N	94 LEU O	3.5
103 TYR H	94 LEU O	2.6
103 TYR N	94 LEU O	3.5
105 LEU H	96 GLY O	2.6
105 LEU N	96 GLY O	3.5
110 ILE H	107 HIS O	2.6
110 ILE N	107 HIS O	3.5
111 SER H	114 HIS ND1	2.6
111 SER N	114 HIS ND1	3.5
114 HIS H	95 PHE O	2.6
114 HIS N	95 PHE O	3.5
115 ALA H	95 PHE O	2.6
115 ALA N	95 PHE O	3.5
116 VAL H	129 MET O	2.6
116 VAL N	129 MET O	3.5
117 LEU H	93 TYR O	2.6
117 LEU N	93 TYR O	3.5
118 VAL H	127 VAL O	2.6
118 VAL N	127 VAL O	3.5
119 PHE H	91 PRO O	2.6
119 PHE N	91 PRO O	3.5
120 HIS H	125 CYS O	2.6
120 HIS N	125 CYS O	3.5
121 GLY H	63 TRP O	2.6
121 GLY N	63 TRP O	3.5
127 VAL H	118 VAL O	2.6
127 VAL N	118 VAL O	3.5
128 LEU H	150 LEU O	2.6
128 LEU N	150 LEU O	3.5
129 MET H	116 VAL O	2.6
129 MET N	116 VAL O	3.5
131 LEU H	114 HIS O	2.6
131 LEU N	114 HIS O	3.5
138 LYS H	159 GLN O	2.6
138 LYS N	159 GLN O	3.5

139 LEU H	142 ASN O	1.7	139 LEU H	142 ASN O	2.6
139 LEU N	142 ASN O	2.7	139 LEU N	142 ASN O	3.5
142 ASN H	139 LEU O	1.7	142 ASN H	139 LEU O	2.6
142 ASN N	139 LEU O	2.7	142 ASN N	139 LEU O	3.5
152 ALA H	126 PHE O	1.7	152 ALA H	126 PHE O	2.6
152 ALA N	126 PHE O	2.7	152 ALA N	126 PHE O	3.5
156 SER H	153 PRO O	1.7	156 SER H	153 PRO O	2.6
156 SER N	153 PRO O	2.7	156 SER N	153 PRO O	3.5
158 ILE H	167 TYR O	1.7	158 ILE H	167 TYR O	2.6
158 ILE N	167 TYR O	2.7	158 ILE N	167 TYR O	3.5
159 GLN H	138 LYS O	1.7	159 GLN H	138 LYS O	2.6
159 GLN N	138 LYS O	2.7	159 GLN N	138 LYS O	3.5
167 TYR H	158 ILE O	1.7	167 TYR H	158 ILE O	2.6
167 TYR N	158 ILE O	2.7	167 TYR N	158 ILE O	3.5
168 LYS H	74 GLN O	1.7	168 LYS H	74 GLN O	2.6
168 LYS N	74 GLN O	2.7	168 LYS N	74 GLN O	3.5
169 VAL H	156 SER O	1.7	169 VAL H	156 SER O	2.6
169 VAL N	156 SER O	2.7	169 VAL N	156 SER O	3.5
170 GLN H	72 HIS O	1.7	170 GLN H	72 HIS O	2.6
170 GLN N	72 HIS O	2.7	170 GLN N	72 HIS O	3.5
107 HIS ND1	109 SER OG	2.7	107 HIS ND1	109 SER OG	3.5
109 SER H	107 HIS ND1	2.7	109 SER H	107 HIS ND1	3.5