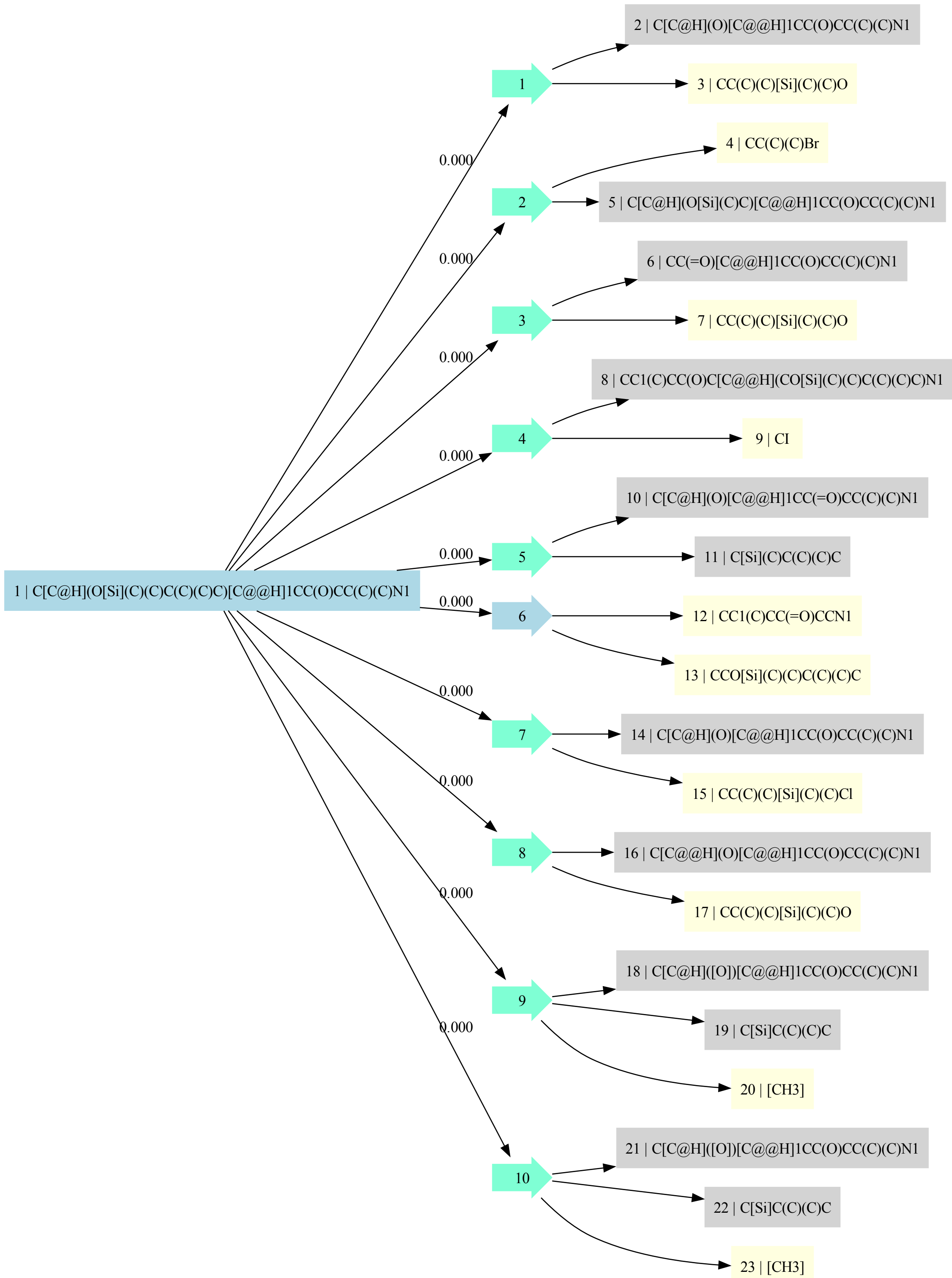
1 | C[C@H](O[Si](C)(C)C(C)(C)C)[C@@H]1CC(O)CC(C)(C)N1
0.000
0.000
0.000
0.000
0.000
0.000
0.000
0.000
0.000
0.000
1
2
3
4
5
6
7
8
9
10
2 | C[C@H](O)[C@@H]1CC(O)CC(C)(C)N1
3 | CC(C)(C)[Si](C)(C)O
4 | CC(C)(C)Br
5 | C[C@H](O[Si](C)C)[C@@H]1CC(O)CC(C)(C)N1
6 | CC(=O)[C@@H]1CC(O)CC(C)(C)N1
7 | CC(C)(C)[Si](C)(C)O
8 | CC1(C)CC(O)C[C@@H](CO[Si](C)(C)C(C)(C)C)N1
9 | CI
10 | C[C@H](O)[C@@H]1CC(=O)CC(C)(C)N1
11 | C[Si](C)C(C)(C)C
12 | CC1(C)CC(=O)CCN1
13 | CCO[Si](C)(C)C(C)(C)C
14 | C[C@H](O)[C@@H]1CC(O)CC(C)(C)N1
15 | CC(C)(C)[Si](C)(C)Cl
16 | C[C@@H](O)[C@@H]1CC(O)CC(C)(C)N1
17 | CC(C)(C)[Si](C)(C)O
18 | C[C@H]([O])[C@@H]1CC(O)CC(C)(C)N1
19 | C[Si]C(C)(C)C
20 | [CH3]
21 | C[C@H]([O])[C@@H]1CC(O)CC(C)(C)N1
22 | C[Si]C(C)(C)C
23 | [CH3]