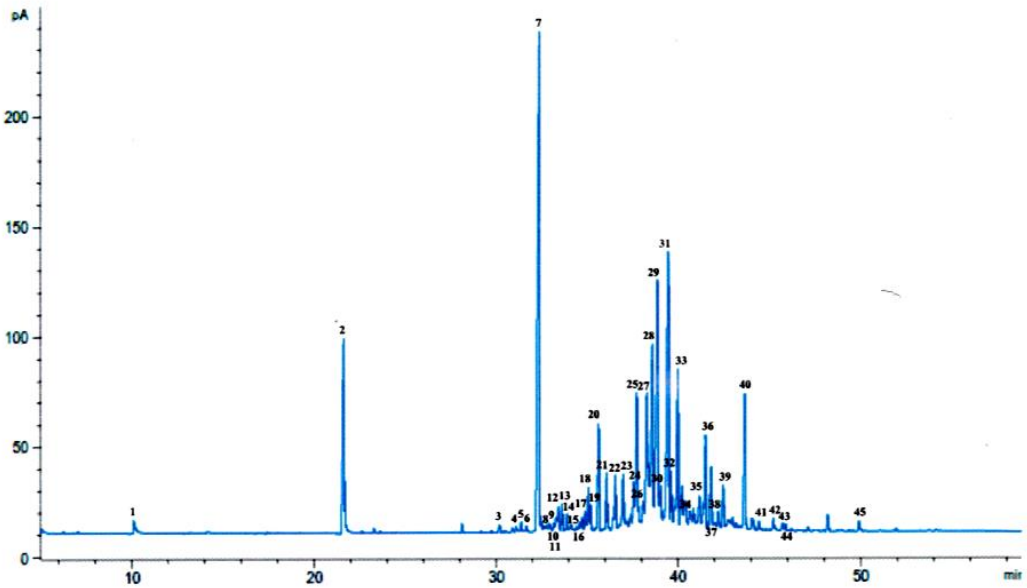
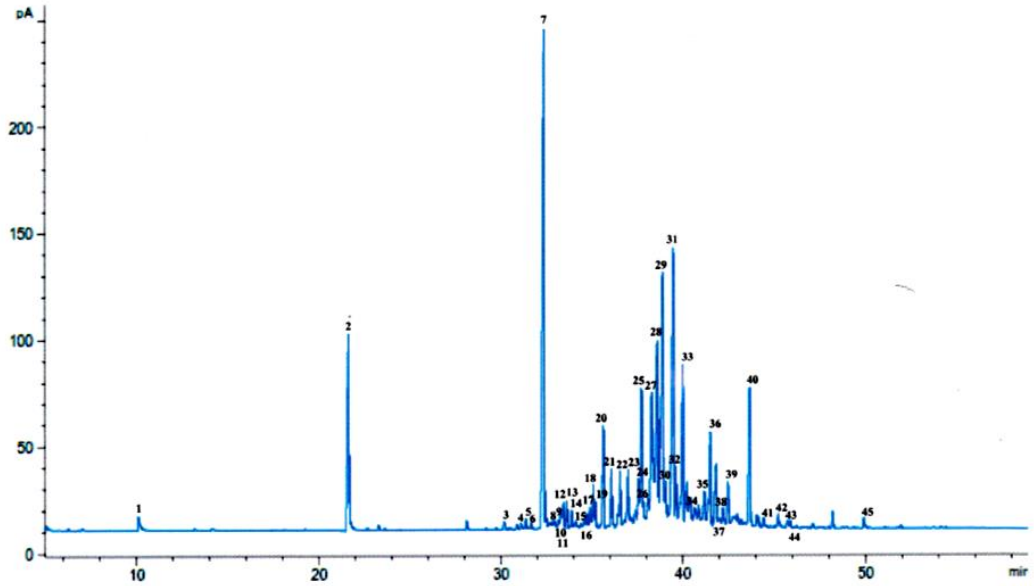
A



No.	Ret. time, t_r (min)	Peak area (%)	Compounds	Ident. mode
1	10.066	0.54	Benzaldehyde	FID,MS
2	21.585	6.36	4-Phenyl-2-butanone	FID,MS
3	30.152	0.20	α-Gurjunene	FID
4	31.071	0.14	β-Caryophyllene	FID
5	31.344	0.19	Aromadendrene	FID
6	31.653	0.13	α-Caryophyllene	FID,MS
7	32.362	17.80	*allo*-Aromadendrene	FID
8	32.750	0.13	γ-Selinene	FID,MS
9	32.879	0.15	γ-Muurolene	FID
10	32.989	0.12	β-Vetispirene	FID,MS
11	33.284	0.23	Valencene	FID,MS
12	33.428	0.69	Dihydro-β-agarofuran	FID,MS
13	33.609	0.60	δ-Guaiene	FID,MS
14	33.880	0.40	γ-Cadinene	FID
15	34.591	0.20	Kessane	FID,MS
16	34.715	0.28	Selina-3,7(11)-diene	FID
17	34.888	0.43	α-Calacorene	FID
18	35.060	1.04	Dehydro-aromadendrene	FID
19	35.185	0.52	α-Agarofuran	FID
20	35.621	2.62	*cis*-Nerolidol	FID
21	36.064	1.41	*nor*-Ketoagarofuran	FID
22	36.538	1.32	Epoxybulnesene	FID
23	36.973	1.56	Caryophyllene oxide	FID
24	37.559	1.11	Guaiol	FID
25	37.714	3.77	10-*epi*-γ-Eudesmol	FID,MS
26	38.071	0.41	γ-Eudesmol	FID,MS
27	38.275	3.95	Agarospirol	FID,MS
28	38.591	6.04	τ-Muurolol	FID
29	38.878	9.04	α-Eudesmol	FID
30	39.048	0.76	Jinkoh-eremol	FID
31	39.505	10.27	Kusunol	FID
32	39.635	1.07	Bulnesol	FID
33	40.016	4.53	Dehydrojinkoh-eremol	FID
34	40.653	0.60	α-Bisabolol	FID
35	41.180	1.12	Rotundone	FID
36	41.529	2.35	Valerenol	FID
37	42.050	0.15	Selina-4,11-dien-14-oic acid	FID
38	42.211	0.45	α-Costol	FID,MS
39	42.482	1.15	Selina-3,11-dien-9-al	FID
40	43.678	3.99	Dehydrofukinone	FID,MS
41	44.434	0.24	Hexadecanal	FID
42	45.230	0.41	Guaia-1(10),11-dien-15-oic acid	FID
43	45.752	0.24	*oxo*-Agarospirol	FID
44	45.904	0.18	Dihydrocolumellarin	FID,MS
45	49.919	0.29	1,5-Diphenyl-3-pentanone	FID,MS

FID	Identification by linear retention indices, determined relative to the retention times on DB1 column of homologous series of C_7-C_{20} n-alkanes having difference in I value of ±10 between calculated Kovats retention index (I_{cal}) and reference Kovats retention index (I_{ref}).
MS	Identification by comparison of mass spectrum generated from sample analysis with library of National Institute of Standards and Technology (NIST) with acceptance similarity ≥ 80%

B

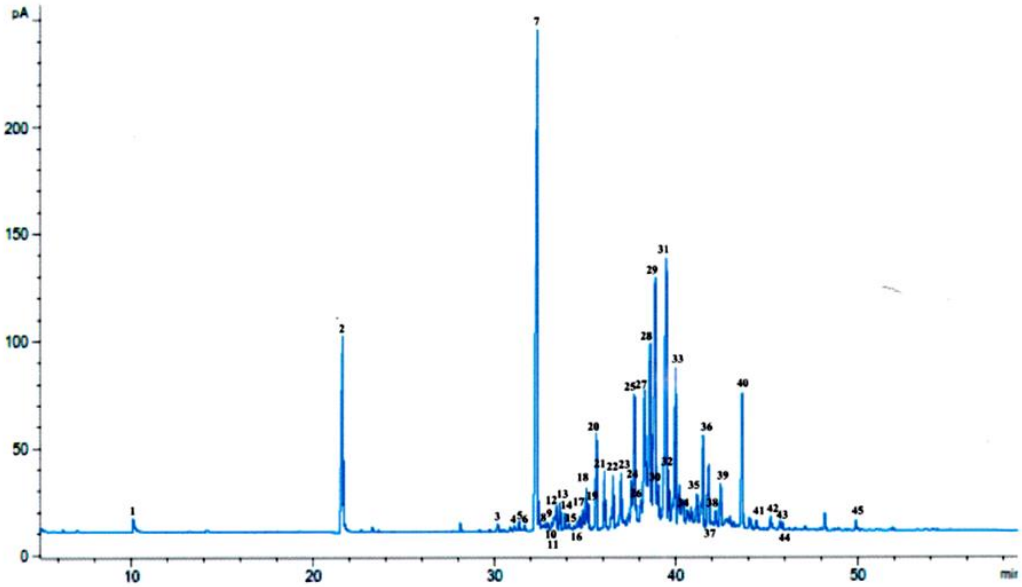


No.	Ret. time, t_r (min)	Peak area (%)	Compounds	Ident. mode
1	10.070	0.55	Benzaldehyde	FID,MS
2	21.594	6.39	4-Phenyl-2-butanone	FID,MS
3	30.152	0.20	α-Gurjunene	FID
4	31.073	0.15	β-Caryophyllene	FID
5	31.346	0.20	Aromadendrene	FID
6	31.652	0.13	α-Caryophyllene	FID,MS
7	32.368	17.90	*allo*-Aromadendrene	FID
8	32.752	0.14	γ-Selinene	FID,MS
9	32.882	0.15	γ-Muurolene	FID
10	32.991	0.12	β-Vetispirene	FID,MS
11	33.285	0.23	Valencene	FID,MS
12	33.431	0.69	Dihydro-β-agarofuran	FID,MS
13	33.612	0.60	δ-Guaiene	FID,MS
14	33.882	0.39	γ-Cadinene	FID
15	34.591	0.20	Kessane	FID,MS
16	34.718	0.27	Selina-3,7(11)-diene	FID
17	34.889	0.44	α-Calacorene	FID
18	35.061	1.02	Dehydro-aromadendrene	FID
19	35.187	0.52	α-Agarofuran	FID
20	35.623	2.47	*cis*-Nerolidol	FID
21	36.067	1.47	*nor*-Ketoagarofuran	FID
22	36.539	1.30	Epoxybulnesene	FID
23	36.975	1.57	Caryophyllene oxide	FID
24	37.563	1.12	Guaiol	FID
25	37.717	3.78	10-*epi*-γ-Eudesmol	FID,MS
26	38.077	0.39	γ-Eudesmol	FID,MS
27	38.276	3.96	Agarospirol	FID,MS
28	38.596	6.02	τ-Muurolol	FID
29	38.886	9.04	α-Eudesmol	FID
30	39.050	0.75	Jinkoh-eremol	FID
31	39.513	10.20	Kusunol	FID
32	39.643	1.07	Bulnesol	FID
33	40.022	4.55	Dehydrojinkoh-eremol	FID
34	40.653	0.60	α-Bisabolol	FID
35	41.183	1.10	Rotundone	FID
36	41.532	2.34	Valerenol	FID
37	42.053	0.16	Selina-4,11-dien-14-oic acid	FID
38	42.214	0.45	α-Costol	FID,MS
39	42.486	1.15	Selina-3,11-dien-9-al	FID
40	43.681	4.02	Dehydrofukinone	FID,MS
41	44.435	0.20	Hexadecanal	FID
42	45.230	0.45	Guaia-1(10),11-dien-15-oic acid	FID
43	45.752	0.23	*oxo*-Agarospirol	FID
44	45.902	0.18	Dihydrocolumellarin	FID,MS
45	49.917	0.28	1,5-Diphenyl-3-pentanone	FID,MS

FID Identification by linear retention indices, determined relative to the retention times on DB1 column of homologous series of C_7-C_{20} n-alkanes having difference in I value of ±10 between calculated Kovats retention index (I_{cal}) and reference Kovats retention index (I_{ref}).

MS Identification by comparison of mass spectrum generated from sample analysis with library of National Institute of Standards and Technology (NIST) with acceptance similarity ≥ 80%

C

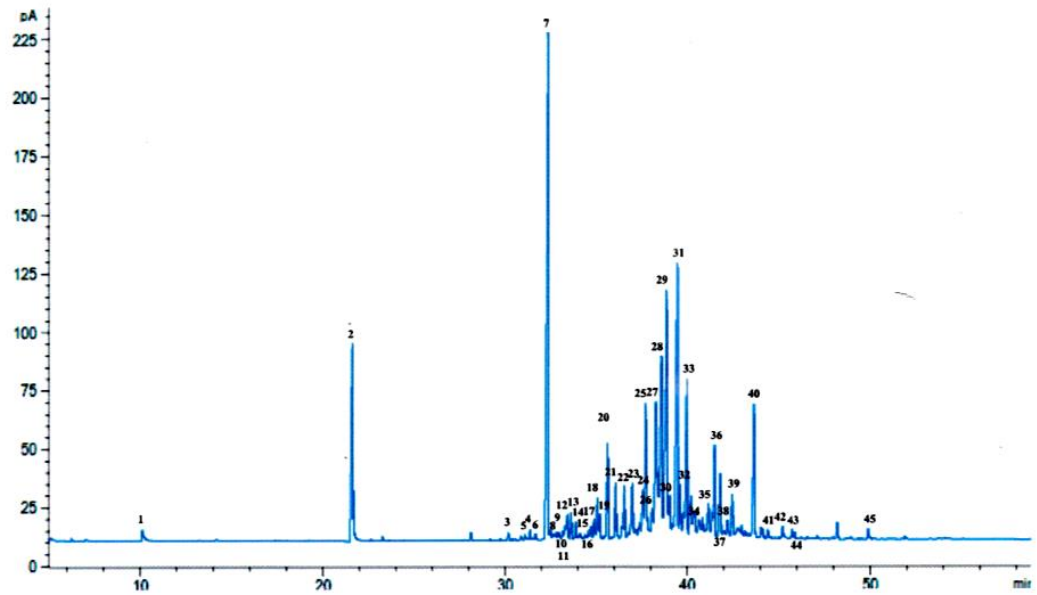


No.	Ret. time, t_r (min)	Peak area (%)	Compounds	Ident. mode
1	10.070	0.54	Benzaldehyde	FID,MS
2	21.592	6.49	4-Phenyl-2-butanone	FID,MS
3	30.153	0.20	α-Gurjunene	FID
4	31.073	0.14	β-Caryophyllene	FID
5	31.346	0.19	Aromadendrene	FID
6	31.653	0.13	α-Caryophyllene	FID,MS
7	32.367	18.04	*allo*-Aromadendrene	FID
8	32.754	0.14	γ-Selinene	FID,MS
9	32.882	0.15	γ-Muurolene	FID
10	32.990	0.12	β-Vetispirene	FID,MS
11	33.285	0.23	Valencene	FID,MS
12	33.430	0.70	Dihydro-β-agarofuran	FID,MS
13	33.611	0.60	δ-Guaiene	FID,MS
14	33.881	0.42	γ-Cadinene	FID
15	34.592	0.18	Kessane	FID,MS
16	34.717	0.27	Selina-3,7(11)-diene	FID
17	34.889	0.44	α-Calacorene	FID
18	35.063	1.00	Dehydro-aromadendrene	FID
19	35.187	0.53	α-Agarofuran	FID
20	35.620	2.37	*cis*-Nerolidol	FID
21	36.066	1.42	*nor*-Ketoagarofuran	FID
22	36.539	1.28	Epoxybulnesene	FID
23	36.976	1.58	Caryophyllene oxide	FID
24	37.560	1.12	Guaiol	FID
25	37.715	3.72	10-*epi*-γ-Eudesmol	FID,MS
26	38.074	0.41	γ-Eudesmol	FID,MS
27	38.278	3.97	Agarospirol	FID,MS
28	38.594	6.11	τ-Muurolol	FID
29	38.886	9.14	α-Eudesmol	FID
30	39.050	0.75	Jinkoh-eremol	FID
31	39.508	10.20	Kusunol	FID
32	39.641	1.07	Bulnesol	FID
33	40.020	4.57	Dehydrojinkoh-eremol	FID
34	40.654	0.61	α-Bisabolol	FID
35	41.183	1.11	Rotundone	FID
36	41.531	2.35	Valerenol	FID
37	42.054	0.16	Selina-4,11-dien-14-oic acid	FID
38	42.212	0.45	α-Costol	FID,MS
39	42.484	1.16	Selina-3,11-dien-9-al	FID
40	43.680	4.05	Dehydrofukinone	FID,MS
41	44.435	0.20	Hexadecanal	FID
42	45.232	0.46	Guaia-1(10),11-dien-15-oic acid	FID
43	45.756	0.28	*oxo*-Agarospirol	FID
44	45.904	0.18	Dihydrocolumellarin	FID,MS
45	49.920	0.30	1,5-Diphenyl-3-pentanone	FID,MS

FID Identification by linear retention indices, determined relative to the retention times on DB1 column of homologous series of C_7-C_{20} n-alkanes having difference in I value of ±10 between calculated Kovats retention index (I_{cal}) and reference Kovats retention index (I_{ref}).

MS Identification by comparison of mass spectrum generated from sample analysis with library of National Institute of Standards and Technology (NIST) with acceptance similarity ≥ 80%

D

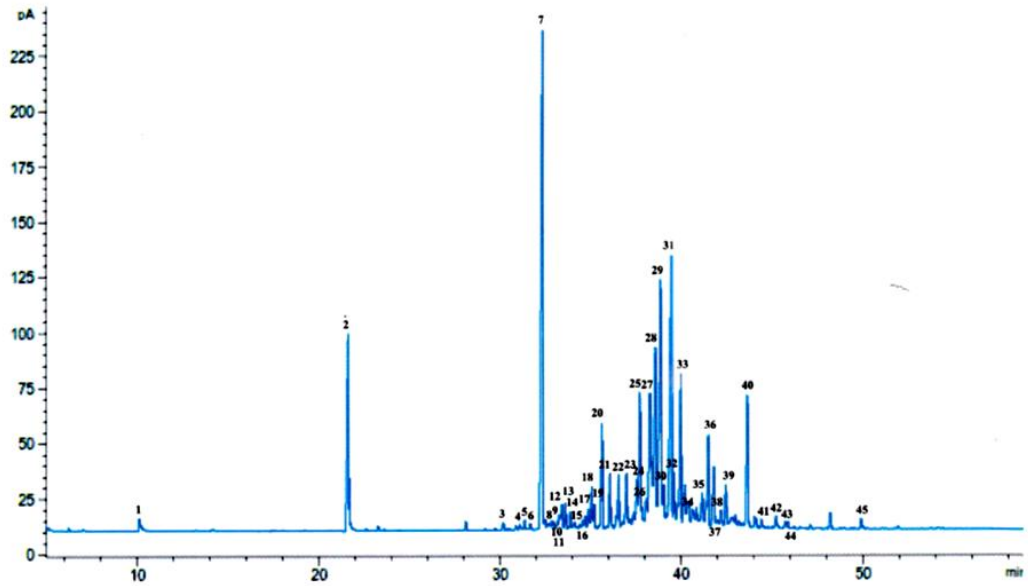


No.	Ret. time, t_r (min)	Peak area (%)	Compounds	Ident. mode
1	10.079	0.48	Benzaldehyde	FID,MS
2	21.584	6.49	4-Phenyl-2-butanone	FID,MS
3	30.152	0.20	α-Gurjunene	FID
4	31.072	0.14	β-Caryophyllene	FID
5	31.345	0.19	Aromadendrene	FID
6	31.650	0.13	α-Caryophyllene	FID,MS
7	32.354	18.03	*allo*-Aromadendrene	FID
8	32.751	0.14	γ-Selinene	FID,MS
9	32.881	0.15	γ-Muurolene	FID
10	32.989	0.12	β-Vetispirene	FID,MS
11	33.284	0.23	Valencene	FID,MS
12	33.428	0.70	Dihydro-β-agarofuran	FID,MS
13	33.610	0.60	δ-Guaiene	FID,MS
14	33.881	0.36	γ-Cadinene	FID
15	34.590	0.18	Kessane	FID,MS
16	34.714	0.27	Selina-3,7(11)-diene	FID
17	34.887	0.43	α-Calacorene	FID
18	35.061	1.00	Dehydro-aromadendrene	FID
19	35.184	0.52	α-Agarofuran	FID
20	35.614	2.40	*cis*-Nerolidol	FID
21	36.061	1.41	*nor*-Ketoagarofuran	FID
22	36.534	1.29	Epoxybulnesene	FID
23	36.970	1.57	Caryophyllene oxide	FID
24	37.551	1.11	Guaiol	FID
25	37.707	3.78	10-*epi*-γ-Eudesmol	FID,MS
26	38.065	0.45	γ-Eudesmol	FID,MS
27	38.266	4.00	Agarospirol	FID,MS
28	38.578	6.15	τ-Muurolol	FID
29	38.866	9.18	α-Eudesmol	FID
30	39.038	0.75	Jinkoh-eremol	FID
31	39.489	10.09	Kusunol	FID
32	39.623	1.06	Bulnesol	FID
33	40.005	4.53	Dehydrojinkoh-eremol	FID
34	40.646	0.53	α-Bisabolol	FID
35	41.174	0.92	Rotundone	FID
36	41.521	2.28	Valerenol	FID
37	42.046	0.10	Selina-4,11-dien-14-oic acid	FID
38	42.208	0.45	α-Costol	FID,MS
39	42.477	1.13	Selina-3,11-dien-9-al	FID
40	43.668	3.94	Dehydrofukinone	FID,MS
41	44.430	0.18	Hexadecanal	FID
42	45.226	0.45	Guaia-1(10),11-dien-15-oic acid	FID
43	45.753	0.27	*oxo*-Agarospirol	FID
44	45.902	0.18	Dihydrocolumellarin	FID,MS
45	49.920	0.31	1,5-Diphenyl-3-pentanone	FID,MS

FID Identification by linear retention indices, determined relative to the retention times on DB1 column of homologous series of C_7-C_{20} n-alkanes having difference in I value of ±10 between calculated Kovats retention index (I_{cal}) and reference Kovats retention index (I_{ref}).

MS Identification by comparison of mass spectrum generated from sample analysis with library of National Institute of Standards and Technology (NIST) with acceptance similarity ≥ 80%

E



No.	Ret. time, t_r (min)	Peak area (%)	Compounds	Ident. mode
1	10.071	0.54	Benzaldehyde	FID,MS
2	21.585	6.53	4-Phenyl-2-butanone	FID,MS
3	30.148	0.20	α-Gurjunene	FID
4	31.068	0.15	β-Caryophyllene	FID
5	31.340	0.19	Aromadendrene	FID
6	31.648	0.14	α-Caryophyllene	FID,MS
7	32.355	18.12	*allo*-Aromadendrene	FID
8	32.749	0.14	γ-Selinene	FID,MS
9	32.877	0.15	γ-Muurolene	FID
10	32.984	0.13	β-Vetispirene	FID,MS
11	33.279	0.23	Valencene	FID,MS
12	33.425	0.70	Dihydro-β-agarofuran	FID,MS
13	33.606	0.60	δ-Guaiene	FID,MS
14	33.877	0.40	γ-Cadinene	FID
15	34.585	0.20	Kessane	FID,MS
16	34.710	0.28	Selina-3,7(11)-diene	FID
17	34.883	0.43	α-Calacorene	FID
18	35.057	1.03	Dehydro-aromadendrene	FID
19	35.180	0.52	α-Agarofuran	FID
20	35.615	2.56	*cis*-Nerolidol	FID
21	36.059	1.41	*nor*-Ketoagarofuran	FID
22	36.532	1.31	Epoxybulnesene	FID
23	36.969	1.68	Caryophyllene oxide	FID
24	37.549	1.09	Guaiol	FID
25	37.707	3.66	10-*epi*-γ-Eudesmol	FID,MS
26	38.067	0.43	γ-Eudesmol	FID,MS
27	38.268	3.97	Agarospirol	FID,MS
28	38.580	6.08	τ-Muurolol	FID
29	38.870	9.10	α-Eudesmol	FID
30	39.039	0.76	Jinkoh-eremol	FID
31	39.493	10.17	Kusunol	FID
32	39.626	1.06	Bulnesol	FID
33	40.006	4.49	Dehydrojinkoh-eremol	FID
34	40.643	0.60	α-Bisabolol	FID
35	41.173	1.10	Rotundone	FID
36	41.521	2.31	Valerenol	FID
37	42.042	0.08	Selina-4,11-dien-14-oic acid	FID
38	42.205	0.44	α-Costol	FID,MS
39	42.475	1.13	Selina-3,11-dien-9-al	FID
40	43.668	3.98	Dehydrofukinone	FID,MS
41	44.428	0.18	Hexadecanal	FID
42	45.223	0.43	Guaia-1(10),11-dien-15-oic acid	FID
43	45.744	0.22	*oxo*-Agarospirol	FID
44	45.897	0.18	Dihydrocolumellarin	FID,MS
45	49.912	0.30	1,5-Diphenyl-3-pentanone	FID,MS

FID Identification by linear retention indices, determined relative to the retention times on DB1 column of homologous series of C_7-C_{20} n-alkanes having difference in I value of ±10 between calculated Kovats retention index (I_{cal}) and reference Kovats retention index (I_{ref}).

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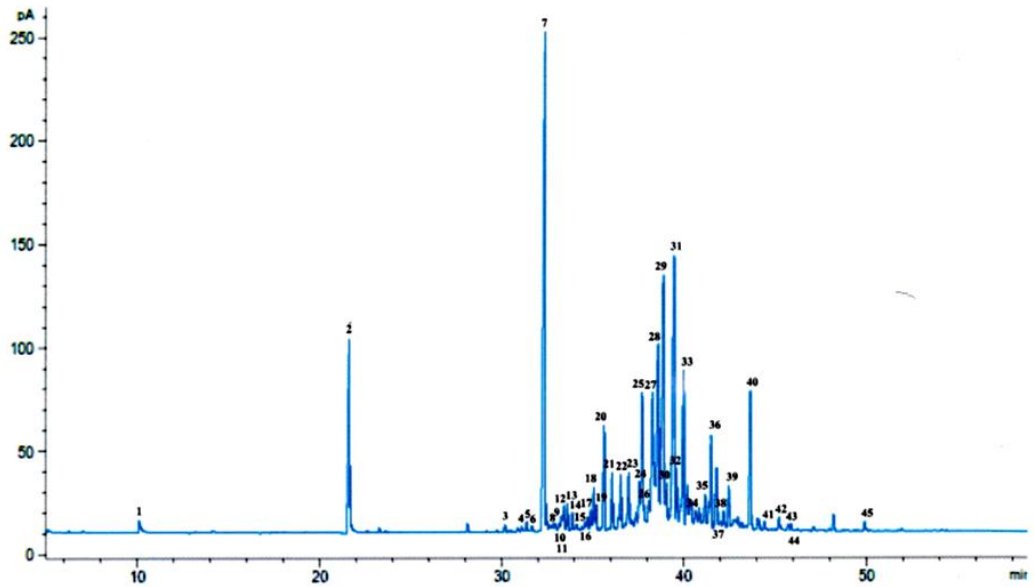
F



No.	Ret. time, t_r (min)	Peak area (%)	Compounds	Ident. mode
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4	31.070	0.15	β-Caryophyllene	FID
5	31.342	0.20	Aromadendrene	FID
6	31.650	0.13	α-Caryophyllene	FID,MS
7	32.366	17.91	*allo*-Aromadendrene	FID
8	32.748	0.14	γ-Selinene	FID,MS
9	32.877	0.15	γ-Muurolene	FID
10	32.987	0.12	β-Vetispirene	FID,MS
11	33.281	0.24	Valencene	FID,MS
12	33.428	0.70	Dihydro-β-agarofuran	FID,MS
13	33.608	0.60	δ-Guaiene	FID,MS
14	33.877	0.42	γ-Cadinene	FID
15	34.586	0.20	Kessane	FID,MS
16	34.714	0.28	Selina-3,7(11)-diene	FID
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18	35.059	1.03	Dehydro-aromadendrene	FID
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29	38.886	9.00	α-Eudesmol	FID
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31	39.513	10.23	Kusunol	FID
32	39.643	1.08	Bulnesol	FID
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34	40.650	0.54	α-Bisabolol	FID
35	41.183	0.86	Rotundone	FID
36	41.529	2.31	Valerenol	FID
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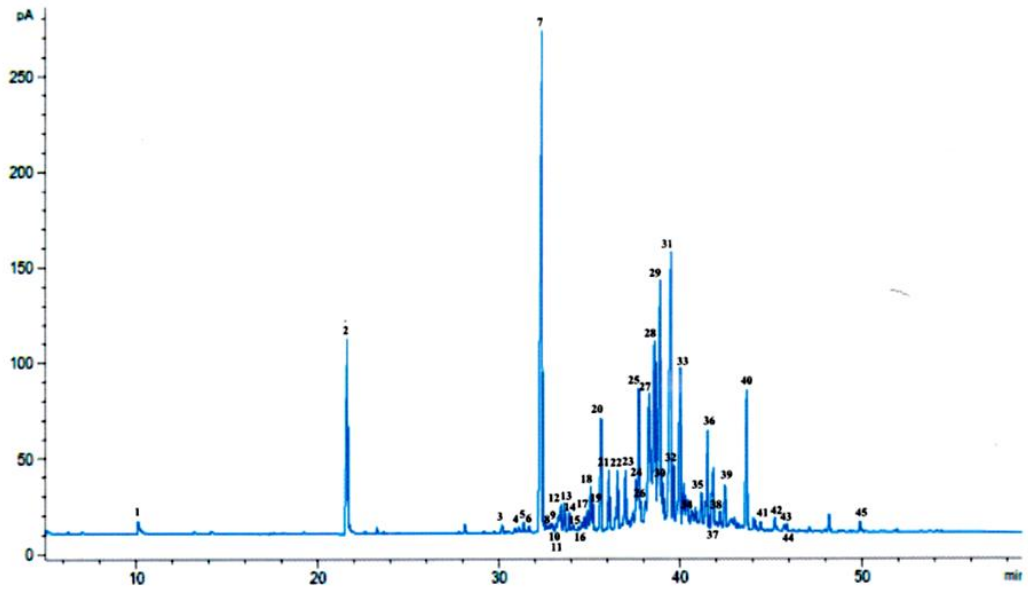
38	42.211	0.44	α-Costol	FID,MS
39	42.481	1.14	Selina-3,11-dien-9-al	FID
40	43.679	4.09	Dehydrofukinone	FID,MS
41	44.431	0.20	Hexadecanal	FID
42	45.225	0.44	Guaia-1(10),11-dien-15-oic acid	FID
43	45.746	0.21	*oxo*-Agarospirol	FID
44	45.898	0.17	Dihydrocolumellarin	FID,MS
45	49.914	0.28	1,5-Diphenyl-3-pentanone	FID,MS

FID Identification by linear retention indices, determined relative to the retention times on DB1 column of homologous series of C_7-C_{20} n-alkanes having difference in I value of ±10 between calculated Kovats retention index (I_{cal}) and reference Kovats retention index (I_{ref}).

MS Identification by comparison of mass spectrum generated from sample analysis with library of National Institute of Standards and Technology (NIST) with acceptance similarity ≥ 80%

G



No.	Ret. time, t_r (min)	Peak area (%)	Compounds	Ident. mode
1	10.067	0.49	Benzaldehyde	FID,MS
2	21.601	6.39	4-Phenyl-2-butanone	FID,MS
3	30.149	0.20	α-Gurjunene	FID
4	31.068	0.15	β-Caryophyllene	FID
5	31.342	0.20	Aromadendrene	FID
6	31.650	0.13	α-Caryophyllene	FID,MS
7	32.380	18.16	*allo*-Aromadendrene	FID
8	32.750	0.14	γ-Selinene	FID,MS
9	32.880	0.15	γ-Muurolene	FID
10	32.988	0.13	β-Vetispirene	FID,MS
11	33.283	0.24	Valencene	FID,MS
12	33.431	0.70	Dihydro-β-agarofuran	FID,MS
13	33.610	0.60	δ-Guaiene	FID,MS
14	33.880	0.42	γ-Cadinene	FID
15	34.588	0.20	Kessane	FID,MS
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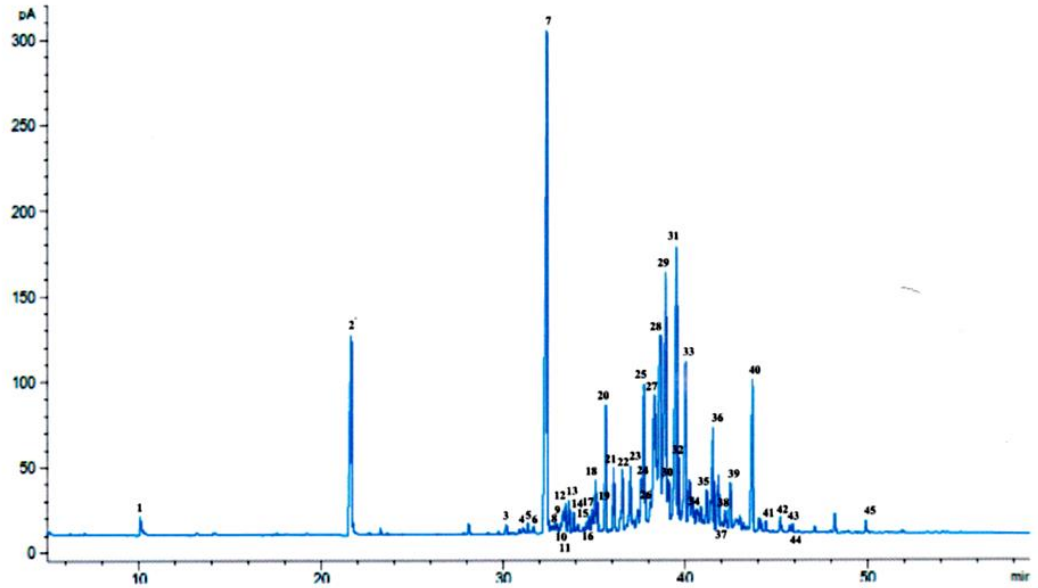
18	35.062	1.05	Dehydro-aromadendrene	FID
19	35.187	0.53	α-Agarofuran	FID
20	35.627	2.66	*cis*-Nerolidol	FID
21	36.068	1.50	*nor*-Ketoagarofuran	FID
22	36.541	1.34	Epoxybulnesene	FID
23	36.976	1.60	Caryophyllene oxide	FID
24	37.568	1.15	Guaiol	FID
25	37.722	3.76	10-*epi*-γ-Eudesmol	FID,MS
26	38.084	0.33	γ-Eudesmol	FID,MS
27	38.291	3.95	Agarospirol	FID,MS
28	38.602	5.91	τ-Muurolol	FID
29	38.902	8.92	α-Eudesmol	FID
30	39.065	0.74	Jinkoh-eremol	FID
31	39.537	10.34	Kusunol	FID
32	39.660	1.08	Bulnesol	FID
33	40.037	4.85	Dehydrojinkoh-eremol	FID
34	40.656	0.33	α-Bisabolol	FID
35	41.191	0.88	Rotundone	FID
36	41.540	2.34	Valerenol	FID
37	42.057	0.16	Selina-4,11-dien-14-oic acid	FID
38	42.215	0.44	α-Costol	FID,MS
39	42.487	1.15	Selina-3,11-dien-9-al	FID
40	43.692	4.14	Dehydrofukinone	FID,MS
41	44.436	0.26	Hexadecanal	FID
42	45.229	0.44	Guaia-1(10),11-dien-15-oic acid	FID
43	45.750	0.21	*oxo*-Agarospirol	FID
44	45.901	0.18	Dihydrocolumellarin	FID,MS
45	49.915	0.28	1,5-Diphenyl-3-pentanone	FID,MS

FID Identification by linear retention indices, determined relative to the retention times on DB1 column of homologous series of C_7-C_{20} n-alkanes having difference in I value of ±10 between calculated Kovats retention index (I_{cal}) and reference Kovats retention index (I_{ref}).

MS Identification by comparison of mass spectrum generated from sample analysis with library of National Institute of Standards and Technology (NIST) with acceptance similarity ≥ 80%

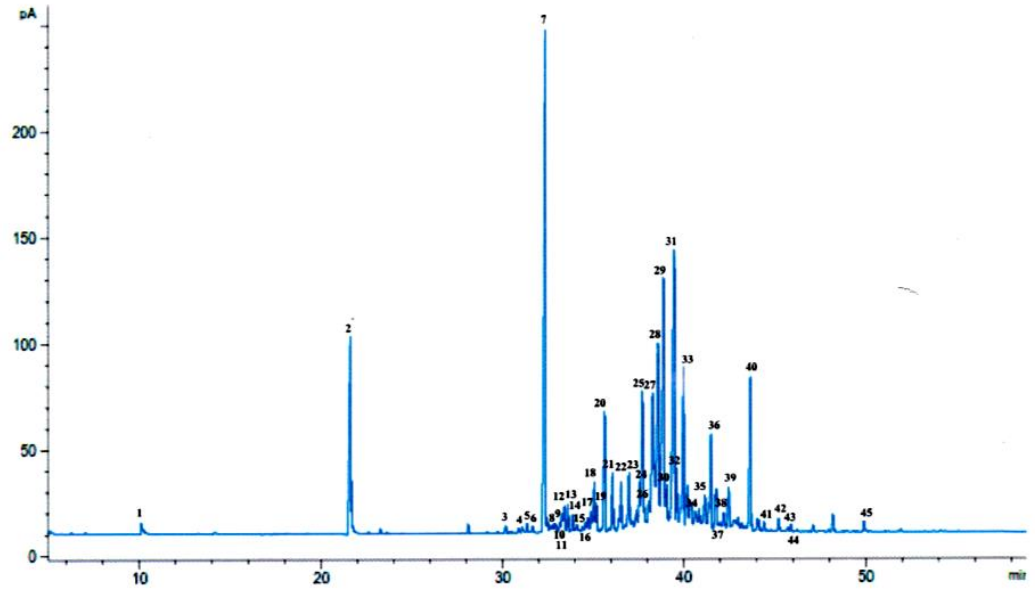
H



No.	Ret. time, t_r (min)	Peak area (%)	Compounds	Ident. mode
1	10.057	0.55	Benzaldehyde	FID,MS
2	21.618	6.13	4-Phenyl-2-butanone	FID,MS
3	30.149	0.20	α-Gurjunene	FID
4	31.071	0.19	β-Caryophyllene	FID
5	31.342	0.23	Aromadendrene	FID
6	31.650	0.25	α-Caryophyllene	FID,MS
7	32.402	17.55	*allo*-Aromadendrene	FID
8	32.753	0.19	γ-Selinene	FID,MS
9	32.884	0.17	γ-Muurolene	FID
10	32.991	0.13	β-Vetispirene	FID,MS
11	33.287	0.28	Valencene	FID,MS
12	33.436	0.65	Dihydro-β-agarofuran	FID,MS
13	33.613	0.58	δ-Guaiene	FID,MS
14	33.882	0.40	γ-Cadinene	FID
15	34.592	0.19	Kessane	FID,MS
16	34.721	0.27	Selina-3,7(11)-diene	FID
17	34.891	0.42	α-Calacorene	FID
18	35.068	1.10	Dehydro-aromadendrene	FID
19	35.194	0.51	α-Agarofuran	FID
20	35.641	2.75	*cis*-Nerolidol	FID
21	36.078	1.40	*nor*-Ketoagarofuran	FID
22	36.549	1.23	Epoxybulnesene	FID
23	36.983	1.52	Caryophyllene oxide	FID
24	37.587	1.11	Guaiol	FID
25	37.737	3.63	10-*epi*-γ-Eudesmol	FID,MS
26	38.091	0.41	γ-Eudesmol	FID,MS
27	38.314	5.99	Agarospirol	FID,MS
28	38.631	6.78	τ-Muurolol	FID
29	38.935	9.45	α-Eudesmol	FID
30	39.087	0.98	Jinkoh-eremol	FID
31	39.575	9.97	Kusunol	FID
32	39.692	1.02	Bulnesol	FID
33	40.066	4.52	Dehydrojinkoh-eremol	FID
34	40.669	0.32	α-Bisabolol	FID
35	41.207	0.77	Rotundone	FID
36	41.557	2.21	Valerenol	FID
37	42.072	0.14	Selina-4,11-dien-14-oic acid	FID
38	42.226	0.42	α-Costol	FID,MS
39	42.501	1.10	Selina-3,11-dien-9-al	FID
40	43.717	4.29	Dehydrofukinone	FID,MS
41	44.443	0.18	Hexadecanal	FID
42	45.235	0.40	Guaia-1(10),11-dien-15-oic acid	FID
43	45.757	0.17	*oxo*-Agarospirol	FID
44	45.907	0.17	Dihydrocolumellarin	FID,MS
45	49.915	0.29	1,5-Diphenyl-3-pentanone	FID,MS

FID Identification by linear retention indices, determined relative to the retention times on DB1 column of homologous series of C_7-C_{20} n-alkanes having difference in I value of ±10 between calculated Kovats retention index (I_{cal}) and reference Kovats retention index (I_{ref}).

MS Identification by comparison of mass spectrum generated from sample analysis with library of National Institute of Standards and Technology (NIST) with acceptance similarity ≥ 80%

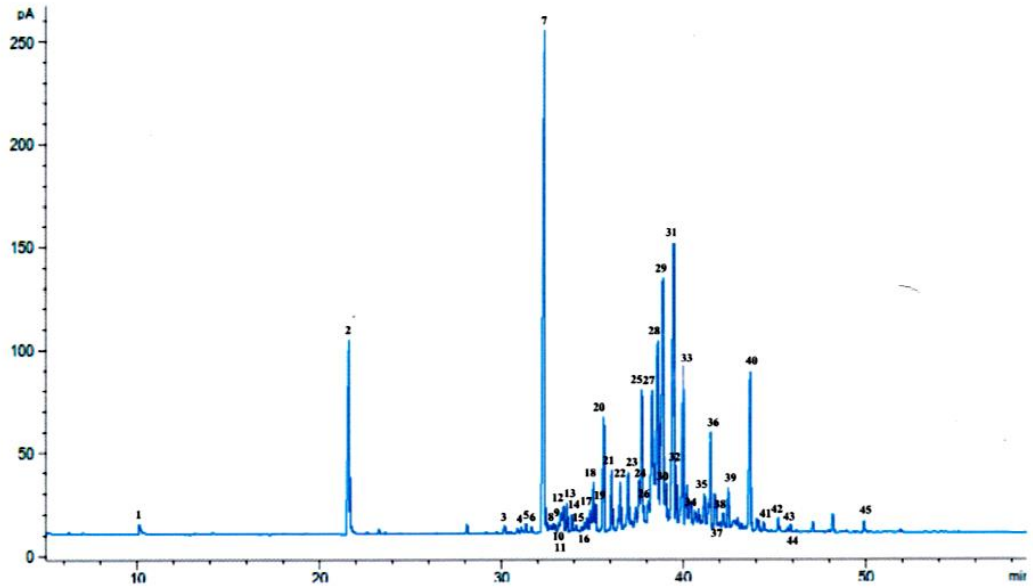


No.	Ret. time, t_r (min)	Peak area (%)	Compounds	Ident. mode
1	10.071	0.49	Benzaldehyde	FID,MS
2	21.586	6.25	4-Phenyl-2-butanone	FID,MS
3	30.146	0.20	α-Gurjunene	FID
4	31.066	0.14	β-Caryophyllene	FID
5	31.338	0.17	Aromadendrene	FID
6	31.646	0.13	α-Caryophyllene	FID,MS
7	32.361	17.70	*allo*-Aromadendrene	FID
8	32.745	0.15	γ-Selinene	FID,MS
9	32.875	0.16	γ-Muurolene	FID
10	32.983	0.13	β-Vetispirene	FID,MS
11	33.278	0.30	Valencene	FID,MS
12	33.425	0.67	Dihydro-β-agarofuran	FID,MS
13	33.605	0.60	δ-Guaiene	FID,MS
14	33.875	0.40	γ-Cadinene	FID
15	34.585	0.16	Kessane	FID,MS
16	34.713	0.28	Selina-3,7(11)-diene	FID
17	34.884	0.42	α-Calacorene	FID
18	35.058	1.26	Dehydro-aromadendrene	FID
19	35.182	0.53	α-Agarofuran	FID
20	35.623	2.89	*cis*-Nerolidol	FID
21	36.060	1.44	*nor*-Ketoagarofuran	FID
22	36.531	1.14	Epoxybulnesene	FID
23	36.969	1.56	Caryophyllene oxide	FID
24	37.554	1.09	Guaiol	FID
25	37.710	3.73	10-*epi*-γ-Eudesmol	FID,MS
26	38.070	0.38	γ-Eudesmol	FID,MS
27	38.275	3.88	Agarospirol	FID,MS
28	38.579	5.98	τ-Muurolol	FID
29	38.879	8.80	α-Eudesmol	FID
30	39.045	0.90	Jinkoh-eremol	FID
31	39.506	10.36	Kusunol	FID
32	39.638	1.12	Bulnesol	FID
33	40.015	4.52	Dehydrojinkoh-eremol	FID
34	40.645	0.58	α-Bisabolol	FID
35	41.177	1.09	Rotundone	FID
36	41.525	2.31	Valerenol	FID
37	42.042	0.16	Selina-4,11-dien-14-oic acid	FID
38	42.205	0.43	α-Costol	FID,MS
39	42.478	1.10	Selina-3,11-dien-9-al	FID
40	43.681	4.50	Dehydrofukinone	FID,MS
41	44.426	0.20	Hexadecanal	FID
42	45.222	0.40	Guaia-1(10),11-dien-15-oic acid	FID
43	45.741	0.12	*oxo*-Agarospirol	FID
44	45.895	0.17	Dihydrocolumellarin	FID,MS
45	49.910	0.29	1,5-Diphenyl-3-pentanone	FID,MS

FID Identification by linear retention indices, determined relative to the retention times on DB1 column of homologous series of C_7-C_{20} n-alkanes having difference in I value of ±10 between calculated Kovats retention index (I_{cal}) and reference Kovats retention index (I_{ref}).

MS Identification by comparison of mass spectrum generated from sample analysis with library of National Institute of Standards and Technology (NIST) with acceptance similarity ≥ 80%

J



No.	Ret. time, t_r (min)	Peak area (%)	Compounds	Ident. mode
1	10.075	0.43	Benzaldehyde	FID,MS
2	21.589	6.13	4-Phenyl-2-butanone	FID,MS
3	30.144	0.20	α-Gurjunene	FID
4	31.065	0.14	β-Caryophyllene	FID
5	31.339	0.17	Aromadendrene	FID
6	31.646	0.13	α-Caryophyllene	FID,MS
7	32.366	17.68	*allo*-Aromadendrene	FID
8	32.742	0.15	γ-Selinene	FID,MS
9	32.875	0.16	γ-Muurolene	FID
10	32.983	0.13	β-Vetispirene	FID,MS
11	33.279	0.30	Valencene	FID,MS
12	33.426	0.67	Dihydro-β-agarofuran	FID,MS
13	33.605	0.60	δ-Guaiene	FID,MS
14	33.875	0.34	γ-Cadinene	FID
15	34.586	0.19	Kessane	FID,MS
16	34.712	0.28	Selina-3,7(11)-diene	FID
17	34.885	0.44	α-Calacorene	FID
18	35.059	1.22	Dehydro-aromadendrene	FID
19	35.181	0.54	α-Agarofuran	FID
20	35.620	2.68	*cis*-Nerolidol	FID
21	36.059	1.45	*nor*-Ketoagarofuran	FID
22	36.530	1.09	Epoxybulnesene	FID
23	36.970	1.58	Caryophyllene oxide	FID
24	37.559	1.10	Guaiol	FID
25	37.713	3.73	10-*epi*-γ-Eudesmol	FID,MS
26	38.068	0.36	γ-Eudesmol	FID,MS
27	38.278	3.91	Agarospirol	FID,MS
28	38.594	5.93	τ-Muurolol	FID
29	38.886	8.98	α-Eudesmol	FID
30	39.050	0.74	Jinkoh-eremol	FID
31	39.516	10.34	Kusunol	FID
32	39.644	1.13	Bulnesol	FID
33	40.020	4.80	Dehydrojinkoh-eremol	FID
34	40.644	0.59	α-Bisabolol	FID
35	41.179	1.10	Rotundone	FID
36	41.527	2.29	Valerenol	FID
37	42.044	0.17	Selina-4,11-dien-14-oic acid	FID
38	42.212	0.43	α-Costol	FID,MS
39	42.480	1.09	Selina-3,11-dien-9-al	FID
40	43.693	5.09	Dehydrofukinone	FID,MS
41	44.430	0.20	Hexadecanal	FID
42	45.224	0.43	Guaia-1(10),11-dien-15-oic acid	FID
43	45.742	0.14	*oxo*-Agarospirol	FID
44	45.895	0.17	Dihydrocolumellarin	FID,MS
45	49.910	0.30	1,5-Diphenyl-3-pentanone	FID,MS

FID Identification by linear retention indices, determined relative to the retention times on DB1 column of homologous series of C_7-C_{20} n-alkanes having difference in I value of ±10 between calculated Kovats retention index (I_{cal}) and reference Kovats retention index (I_{ref}).

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