LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092519\
Data File : VU034814.D
Acq On : 25 Sep 2019 19:57
Operator : JC/SP
Sample : K4983-03
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BFDF4

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM091919WMA.M
Title : VOC Analysis

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.177	22	27	32	rBV2	9373	9099	0.33%	0.023%
2	1.392	86	94	108	rBV	314836	340958	12.36%	0.854%
3	1.466	110	117	126	rBV	43281	49837	1.81%	0.125%
4	1.543	133	141	150	rVB2	32969	44629	1.62%	0.112%
5	1.672	172	181	196	rBV	259934	338836	12.28%	0.849%
6	2.257	352	363	373	rBV	460549	707440	25.65%	1.773%
7	2.305	373	378	392	rVB	72236	114296	4.14%	0.286%
8	2.685	483	496	515	rVV	292079	525765	19.06%	1.318%
9	2.775	515	524	533	rBV6	8683	16516	0.60%	0.041%
10	3.630	780	790	796	rBV6	4500	8353	0.30%	0.021%
11	3.968	888	895	912	rVB4	9181	21979	0.80%	0.055%
12	4.071	912	927	939	rVB7	9232	23422	0.85%	0.059%
13	4.151	939	952	981	rBV	226238	633949	22.98%	1.589%
14	4.624	1078	1099	1124	rBV2	345287	845586	30.65%	2.119%
15	4.762	1131	1142	1154	rVB7	9441	21981	0.80%	0.055%
16	4.971	1196	1207	1218	rVB7	13003	30495	1.11%	0.076%
17	5.315	1289	1314	1326	rBV2	753878	2214007	80.26%	5.548%
18	5.527	1366	1380	1398	rVV2	73661	166484	6.04%	0.417%
19	5.768	1444	1455	1469	rVB8	5649	14116	0.51%	0.035%
20	5.865	1469	1485	1501	rBV	523472	1050440	38.08%	2.632%
21	6.157	1565	1576	1600	rVB	36008	99776	3.62%	0.250%
22	6.309	1608	1623	1640	rBV	506071	1034118	37.49%	2.592%
23	6.402	1642	1652	1678	rVB4	30117	77765	2.82%	0.195%
24	6.563	1695	1702	1705	rBV	12635	15936	0.58%	0.040%
25	6.588	1707	1710	1726	rVB3	19524	34091	1.24%	0.085%
26	6.685	1728	1740	1770	rBV	1503940	2758396	100.00%	6.913%
27	7.048	1841	1853	1862	rVB5	9854	17661	0.64%	0.044%
28	7.206	1889	1902	1924	rBV2	292089	539944	19.57%	1.353%
29	7.315	1929	1936	1945	rVB5	13684	22926	0.83%	0.057%
30	7.399	1951	1962	1970	rBV	196130	331492	12.02%	0.831%
31	7.440	1971	1975	1990	rVB2	56450	89237	3.24%	0.224%
32	7.543	1994	2007	2019	rBV	1032926	1821926	66.05%	4.566%
33	7.614	2019	2029	2045	rVB	619395	1078052	39.08%	2.702%
34	7.826	2085	2095	2111	rBV2	243950	437230	15.85%	1.096%

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Operator : JC/SP
Sample : K4983-03
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BFDF4

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM091919WMA.M
Title : VOC Analysis

35	7.897	2111	2117	2138	rVB4	26596	70379	2.55%	0.176%
36	8.135	2178	2191	2204	rBV	289478	486211	17.63%	1.218%
37	8.212	2206	2215	2219	rBV4	33995	53036	1.92%	0.133%
38	8.241	2220	2224	2229	rVV4	24778	30276	1.10%	0.076%
39	8.289	2229	2239	2262	rVV	853402	1480874	53.69%	3.711%
40	8.395	2262	2272	2285	rBV	475934	739921	26.82%	1.854%
41	8.514	2302	2309	2320	rVB2	23761	41442	1.50%	0.104%
42	8.884	2411	2424	2449	rBV	575493	925757	33.56%	2.320%
43	9.067	2456	2481	2508	rBV	800002	1409076	51.08%	3.531%
44	9.228	2518	2531	2550	rBV	331960	549002	19.90%	1.376%
45	9.350	2555	2569	2585	rBV	1419153	2355865	85.41%	5.904%
46	9.588	2635	2643	2654	rVV2	18405	27487	1.00%	0.069%
47	9.755	2683	2695	2720	rVB	1171009	1937262	70.23%	4.855%
48	9.916	2736	2745	2758	rVV2	12074	28280	1.03%	0.071%
49	10.016	2767	2776	2785	rVV2	13102	27899	1.01%	0.070%
50	10.144	2806	2816	2827	rBV3	53609	88706	3.22%	0.222%
51	10.299	2858	2864	2872	rBV3	5911	9457	0.34%	0.024%
52	10.408	2886	2898	2914	rVV	669683	1077735	39.07%	2.701%
53	10.466	2914	2916	2925	rVB9	5925	6352	0.23%	0.016%
54	10.566	2936	2947	2961	rBV	103890	176519	6.40%	0.442%
55	10.659	2964	2976	2994	rVV	454896	1031572	37.40%	2.585%
56	10.749	2994	3004	3028	rVB	325702	534954	19.39%	1.341%
57	10.897	3041	3050	3056	rBV2	27583	45366	1.64%	0.114%
58	10.948	3056	3066	3085	rVB	351108	574515	20.83%	1.440%
59	11.125	3109	3121	3137	rVV	917203	1419248	51.45%	3.557%
60	11.263	3154	3164	3169	rBV8	8075	15449	0.56%	0.039%
61	11.305	3171	3177	3183	rVB4	9454	14073	0.51%	0.035%
62	11.405	3199	3208	3214	rBV2	43365	62251	2.26%	0.156%
63	11.459	3214	3225	3241	rVV	790331	1298317	47.07%	3.254%
64	11.546	3242	3252	3266	rVB	590355	900751	32.65%	2.257%
65	11.742	3301	3313	3322	rBV	339844	591402	21.44%	1.482%
66	11.836	3332	3342	3369	rVB2	1065535	2072998	75.15%	5.195%
67	11.958	3372	3380	3395	rBV3	34102	58095	2.11%	0.146%
68	12.032	3396	3403	3415	rVB	44124	74769	2.71%	0.187%
69	12.125	3419	3432	3436	rBV2	71058	115672	4.19%	0.290%
70	12.154	3437	3441	3451	rVB	65312	87538	3.17%	0.219%
71	12.228	3455	3464	3471	rBV2	134342	202774	7.35%	0.508%

LSC Area Percent Report

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Sample : K4983-03
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BFDF4

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM091919WMA.M
Title : VOC Analysis

72	12.331	3486	3496	3518	rVB2	135108	262663	9.52%	0.658%
73	12.472	3529	3540	3544	rBV2	12020	21188	0.77%	0.053%
74	12.530	3548	3558	3568	rVV4	60006	106022	3.84%	0.266%
75	12.594	3568	3578	3583	rVV2	79190	134935	4.89%	0.338%
76	12.627	3583	3588	3596	rVV	103083	158550	5.75%	0.397%
77	12.681	3596	3605	3619	rVB	171292	264019	9.57%	0.662%
78	12.768	3624	3632	3637	rBV8	15413	23409	0.85%	0.059%
79	12.942	3678	3686	3702	rVB	115415	177584	6.44%	0.445%
80	13.099	3717	3735	3750	rBV2	315497	550451	19.96%	1.379%
81	13.164	3750	3755	3766	rVV6	27434	45816	1.66%	0.115%
82	13.218	3769	3772	3778	rVV	10222	12469	0.45%	0.031%
83	13.270	3778	3788	3802	rVB4	74894	135721	4.92%	0.340%
84	13.376	3811	3821	3831	rBV7	71409	128662	4.66%	0.322%
85	13.434	3831	3839	3848	rVB3	37210	56717	2.06%	0.142%
86	13.488	3848	3856	3862	rBV4	52396	79843	2.89%	0.200%
87	13.559	3873	3878	3881	rBV3	11399	14740	0.53%	0.037%
88	13.588	3881	3887	3909	rVB5	53608	114866	4.16%	0.288%
89	13.720	3917	3928	3951	rBV	387488	682849	24.76%	1.711%
90	13.842	3956	3966	3983	rVV3	27011	62406	2.26%	0.156%
91	13.929	3983	3993	4003	rVV6	20280	44117	1.60%	0.111%
92	13.990	4007	4012	4021	rVB4	18443	27865	1.01%	0.070%
93	14.135	4047	4057	4071	rVV4	34345	66206	2.40%	0.166%
94	14.311	4103	4112	4125	rBV7	33474	79029	2.87%	0.198%
95	14.459	4150	4158	4166	rVV7	13800	25237	0.91%	0.063%
96	14.520	4170	4177	4186	rVB6	20852	35856	1.30%	0.090%
97	14.655	4213	4219	4231	rVB6	11498	20396	0.74%	0.051%
98	14.807	4258	4266	4271	rBV3	11429	16420	0.60%	0.041%
99	14.858	4272	4282	4308	rVV	101563	199832	7.24%	0.501%
100	15.045	4329	4340	4352	rBV	109036	198080	7.18%	0.496%

Sum of corrected areas: 39903966

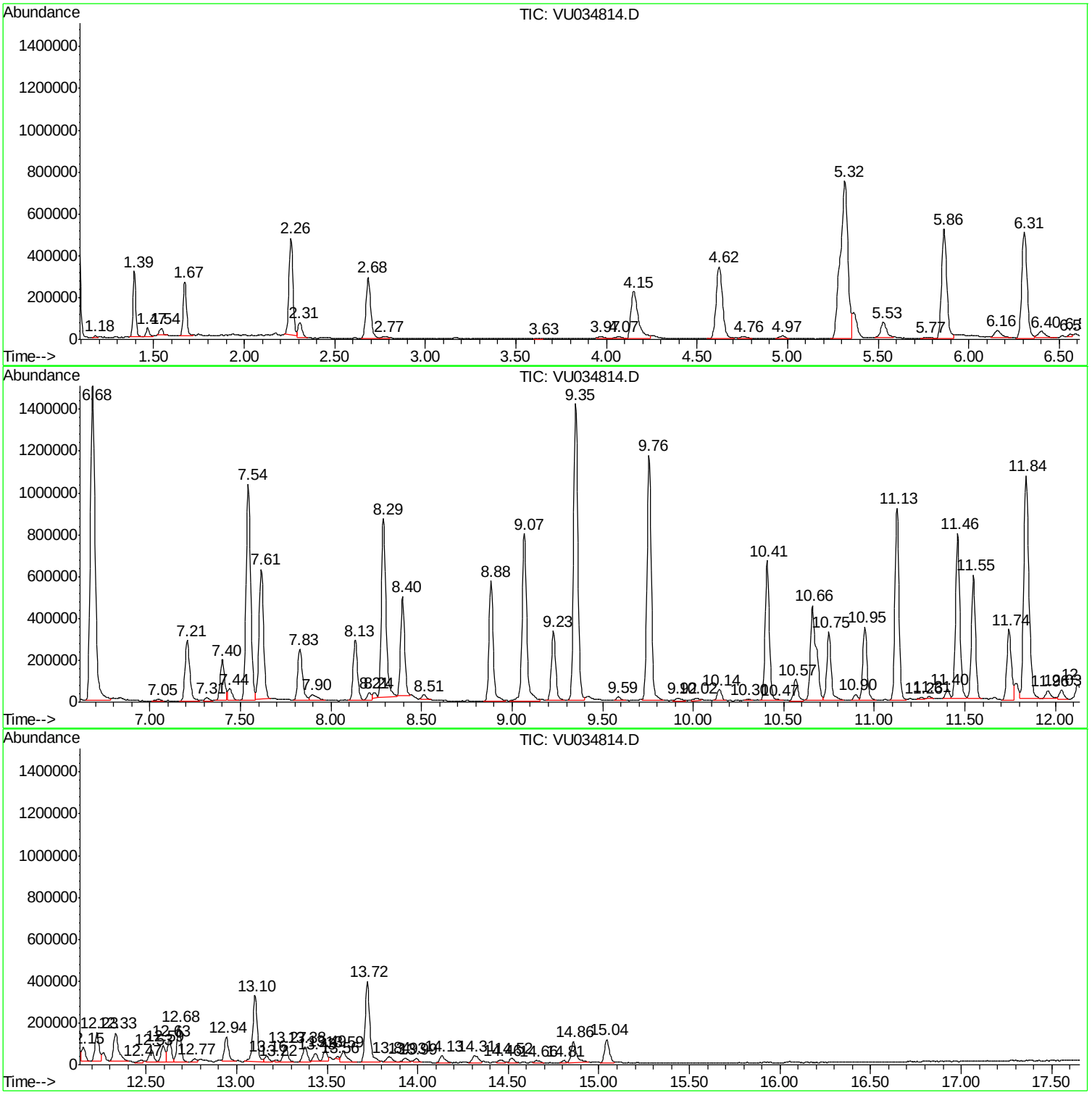
LSC Report - Integrated Chromatogram

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092519\
Data File : VU034814.D
Acq On : 25 Sep 2019 19:57
Operator : JC/SP
Sample : K4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BFDF4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM091919WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092519\
Data File : VU034814.D
Acq On : 25 Sep 2019 19:57
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ALS Vial : 25 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BFDF4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM091919WMA.M
Quant Title : VOC Analysis

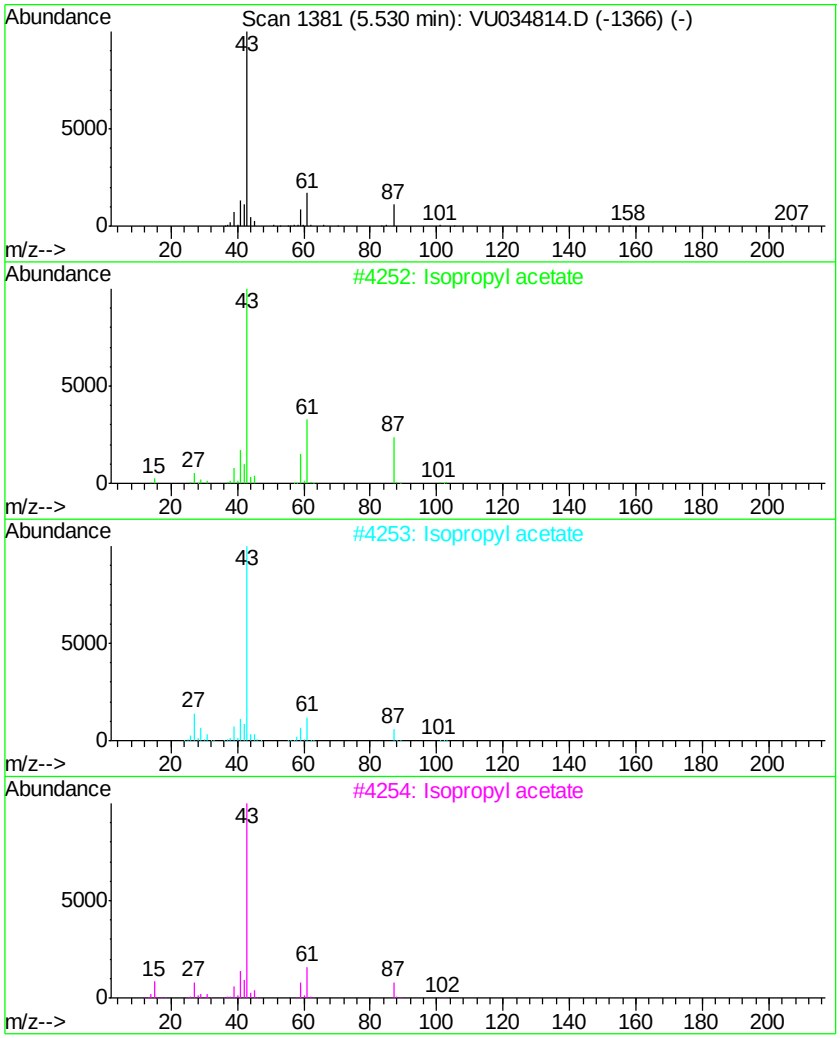
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

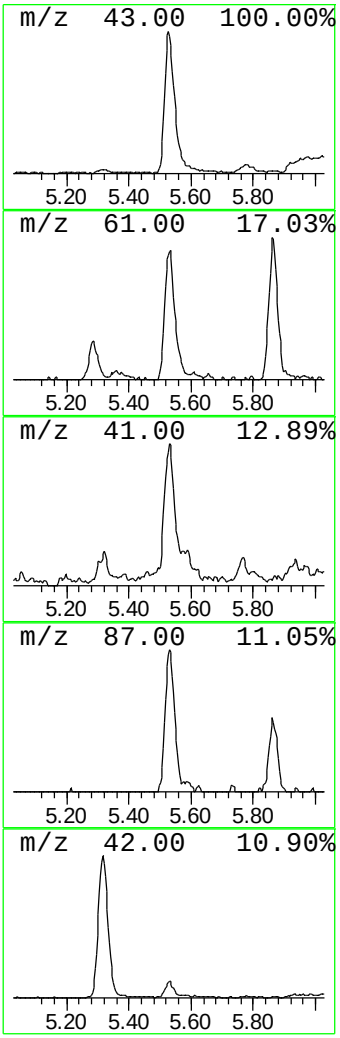
****

Peak Number 1 Isopropyl acetate Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.53	7.92 ug/L	166484	1,4-Difluorobenzene	5.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Isopropyl acetate	102	C5H10O2	000108-21-4	72
2	Isopropyl acetate	102	C5H10O2	000108-21-4	56
3	Isopropyl acetate	102	C5H10O2	000108-21-4	56
4	Isopropyl acetate	102	C5H10O2	000108-21-4	50
5	Acetaldehyde, O-ethyloxime-	87	C4H9NO	1000288-34-6	9





Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092519\
Data File : VU034814.D
Acq On : 25 Sep 2019 19:57
Operator : JC/SP
Sample : K4983-03
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BFDF4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM091919WMA.M
Quant Title : VOC Analysis

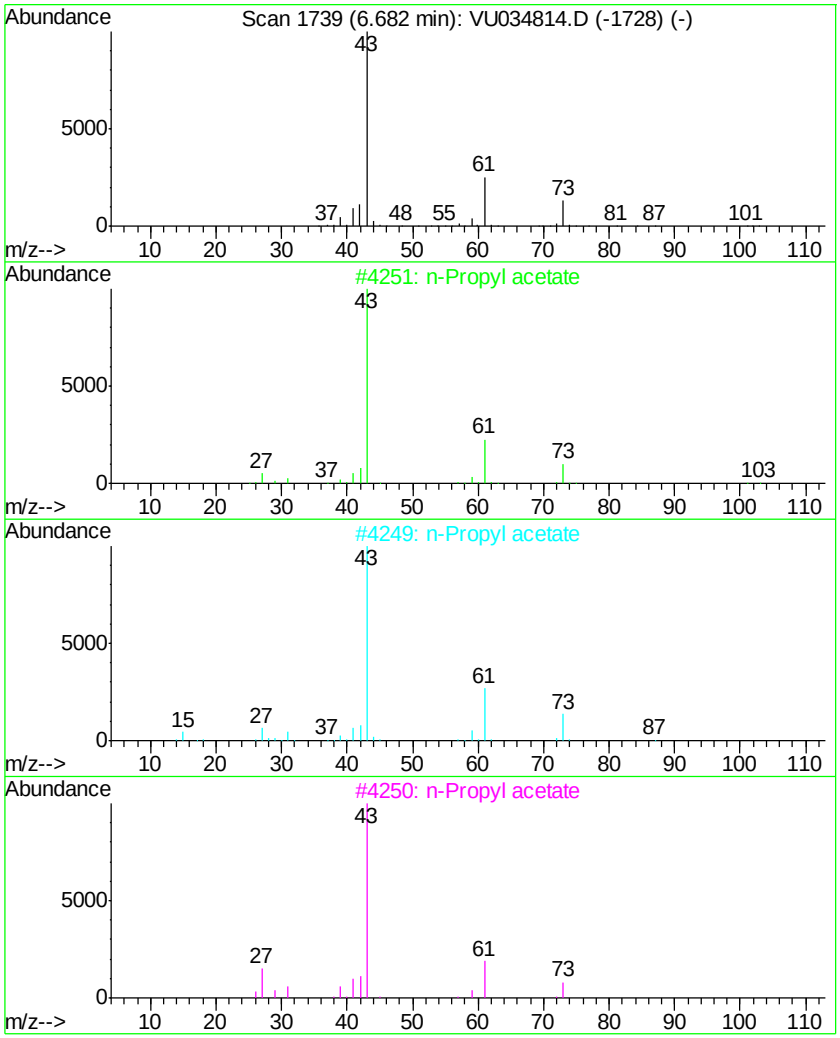
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

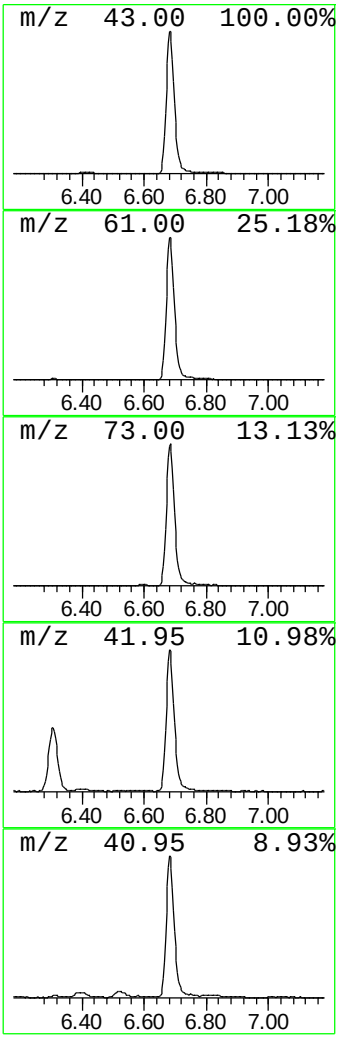
****

Peak Number 2 n-Propyl acetate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.68	131.30 ug/L	2758400	1,4-Difluorobenzene	5.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Propyl acetate	102	C5H10O2	000109-60-4	83
2	n-Propyl acetate	102	C5H10O2	000109-60-4	78
3	n-Propyl acetate	102	C5H10O2	000109-60-4	78
4	Isopropyl acetate	102	C5H10O2	000108-21-4	36
5	Isopropyl acetate	102	C5H10O2	000108-21-4	9





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ClientSampleId : BFDF4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM091919WMA.M
Quant Title : VOC Analysis

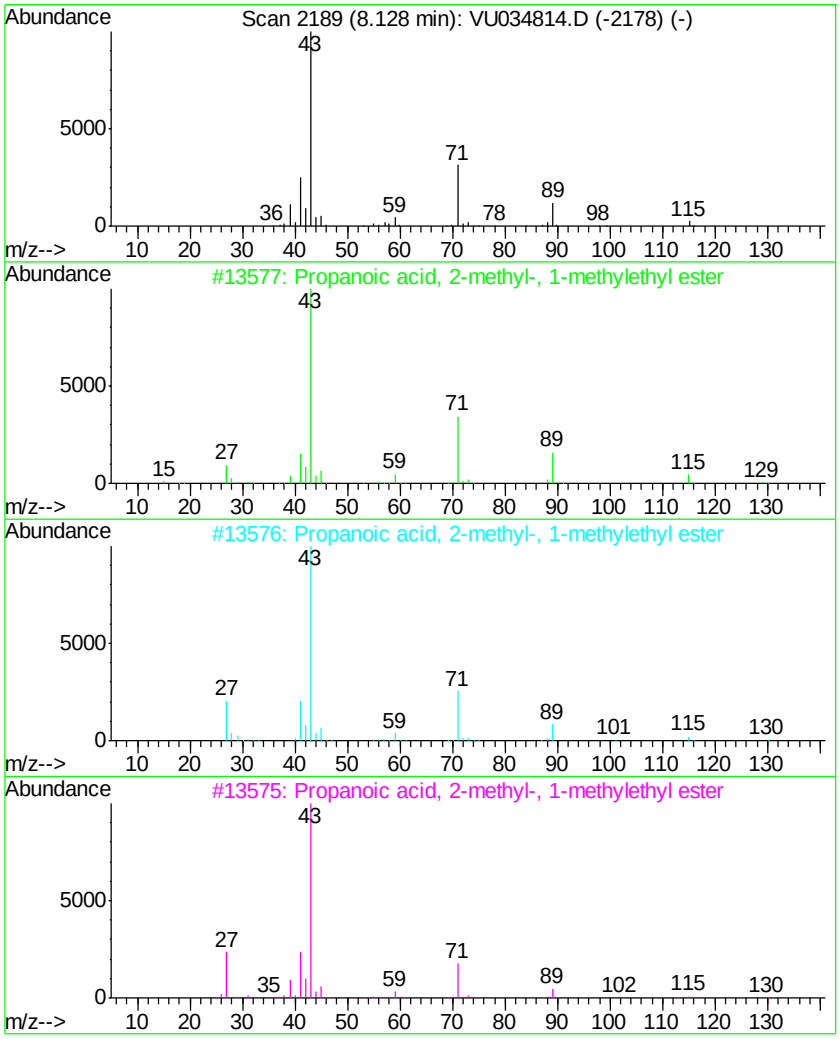
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

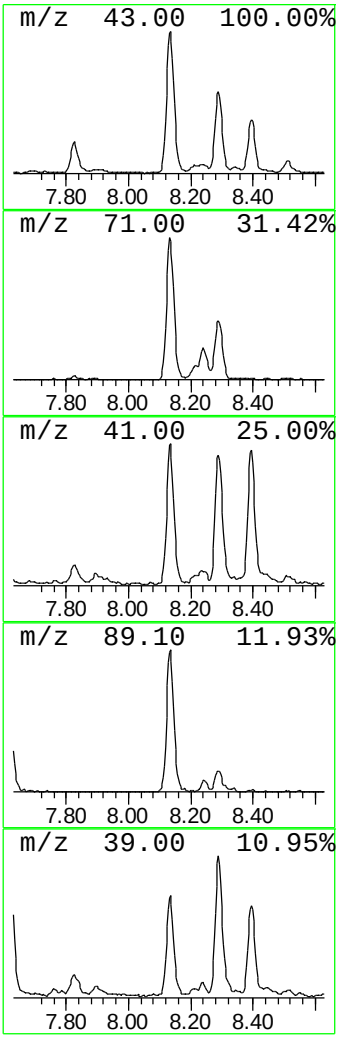
****

Peak Number 4 Propanoic acid, 2-methyl-, ... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.13	17.25 ug/L	486211	Chlorobenzene-d5	9.07

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	83
2	Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	83
3	Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	64
4	Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	59
5	Propanoic acid, 2-methyl-, propy...	130	C7H14O2	000644-49-5	45





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ALS Vial : 25 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BFDF4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM091919WMA.M
Quant Title : VOC Analysis

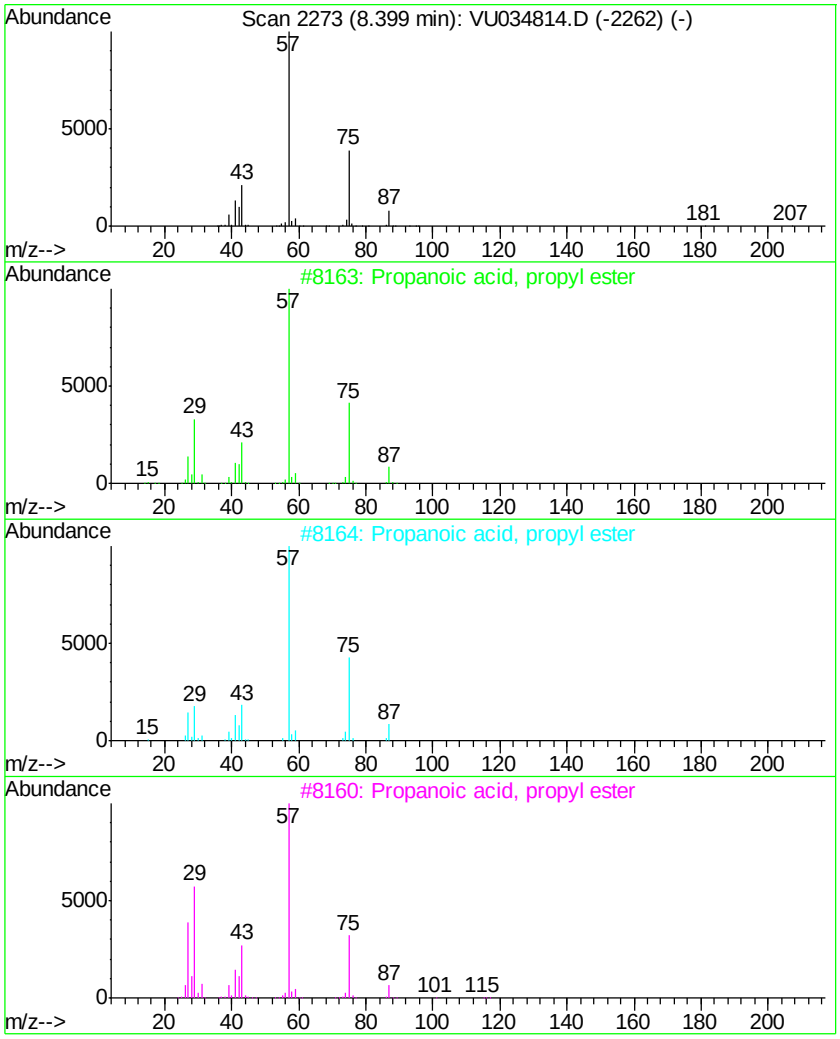
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

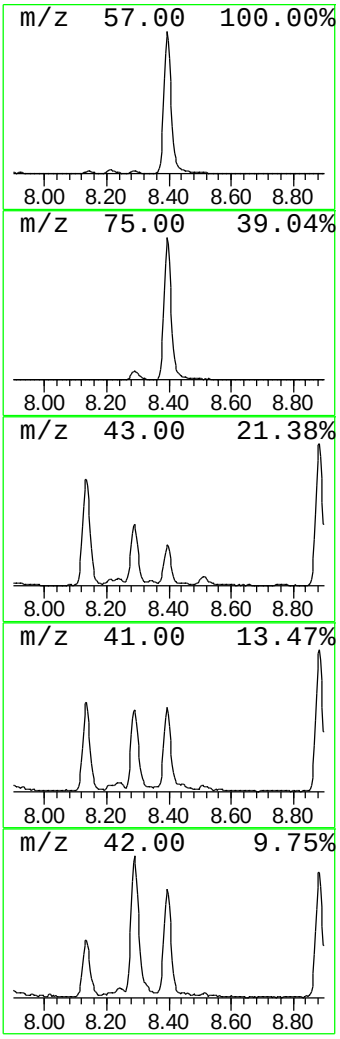
****

Peak Number 5 Propanoic acid, propyl ester Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.40	26.26 ug/L	739921	Chlorobenzene-d5	9.07

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	83
2	Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	83
3	Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	83
4	Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	74
5	Propanoic acid, butyl ester	130	C7H14O2	000590-01-2	38





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Operator : JC/SP
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Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BFDF4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM091919WMA.M
Quant Title : VOC Analysis

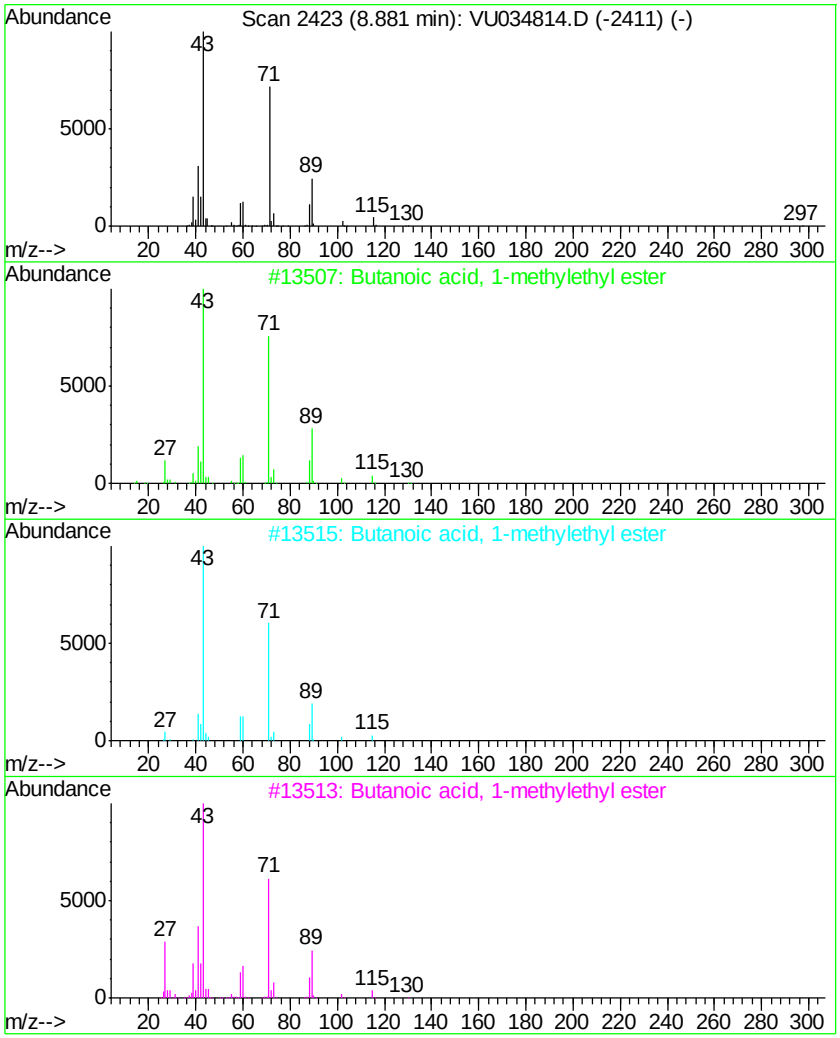
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

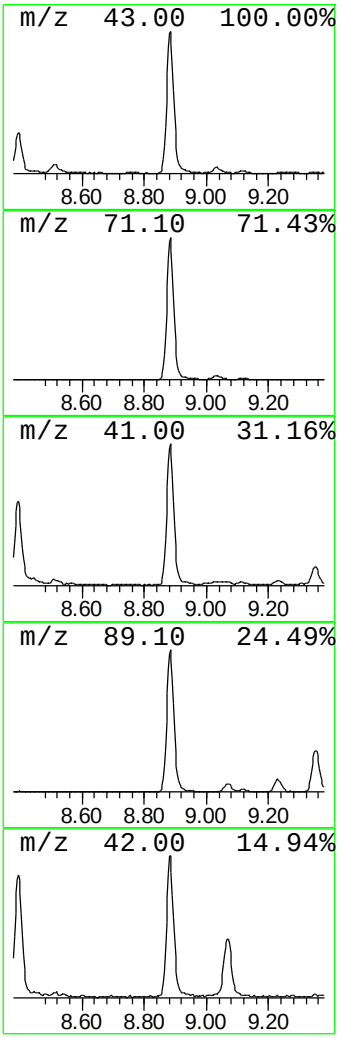
****

Peak Number 6 Butanoic acid, 1-methylethy... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.88	32.85 ug/L	925757	Chlorobenzene-d5	9.07

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid, 1-methylethyl ester	130	C7H14O2	000638-11-9	95
2	Butanoic acid, 1-methylethyl ester	130	C7H14O2	000638-11-9	90
3	Butanoic acid, 1-methylethyl ester	130	C7H14O2	000638-11-9	90
4	Butanoic acid, 1-methylethyl ester	130	C7H14O2	000638-11-9	86
5	Propanoic acid, 2-methyl-, propy...	130	C7H14O2	000644-49-5	59





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Data File : VU034814.D
Acq On : 25 Sep 2019 19:57
Operator : JC/SP
Sample : K4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BFDF4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM091919WMA.M
Quant Title : VOC Analysis

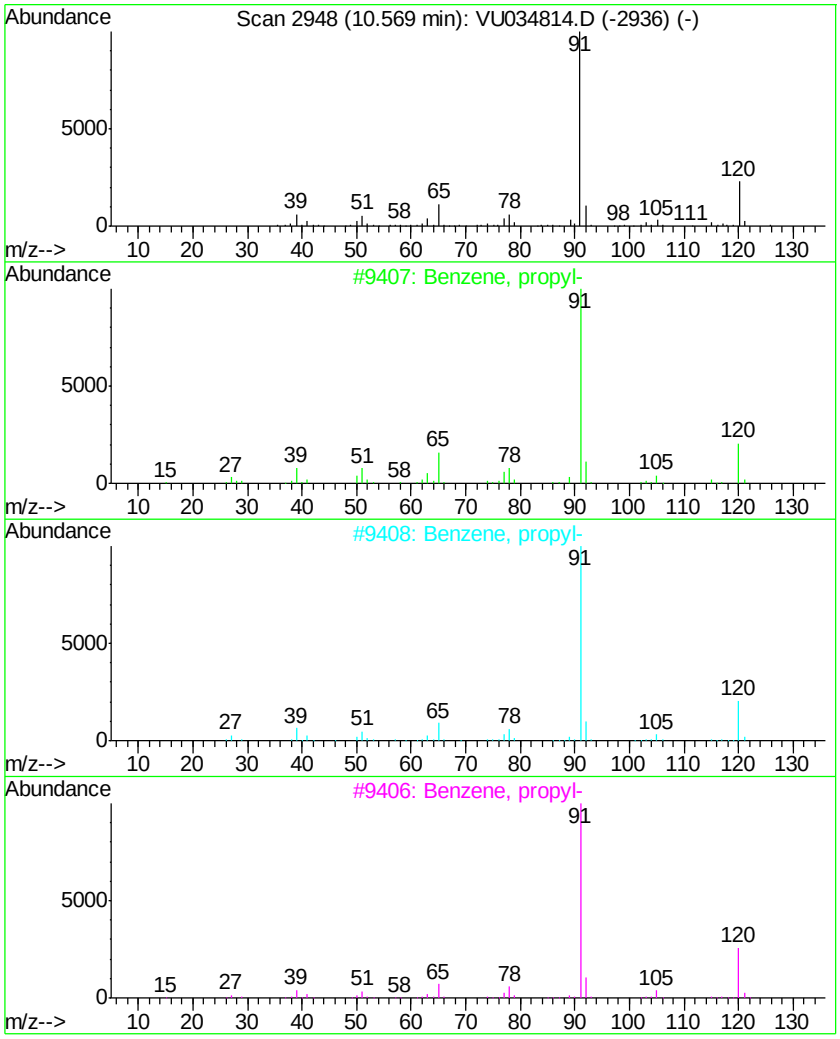
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

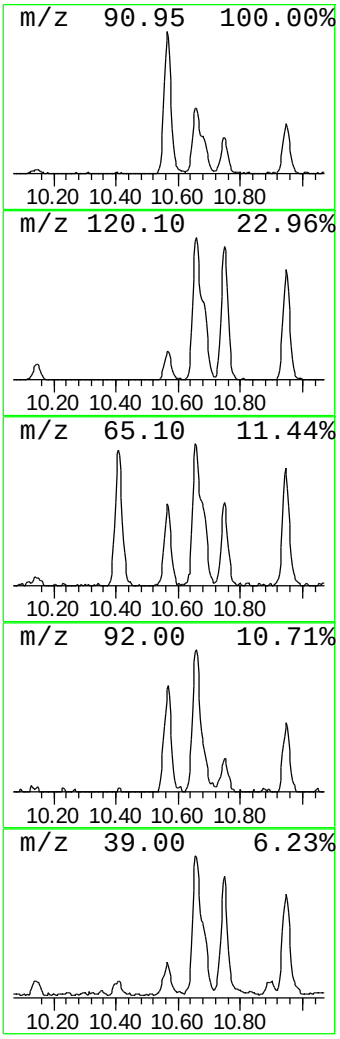
****

Peak Number 7 Benzene, propyl- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.57	6.80 ug/L	176519	1,4-Dichlorobenzene-d4	11.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, propyl-	120	C9H12	000103-65-1	93
2	Benzene, propyl-	120	C9H12	000103-65-1	87
3	Benzene, propyl-	120	C9H12	000103-65-1	80
4	Phenethylamine, N-benzyl-p-chloro-	245	C15H16ClN	013622-43-0	78
5	1,2-Ethanediamine, N,N'-bis(phen...	240	C16H20N2	000140-28-3	78





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Data File : VU034814.D
Acq On : 25 Sep 2019 19:57
Operator : JC/SP
Sample : K4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BFDF4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM091919WMA.M
Quant Title : VOC Analysis

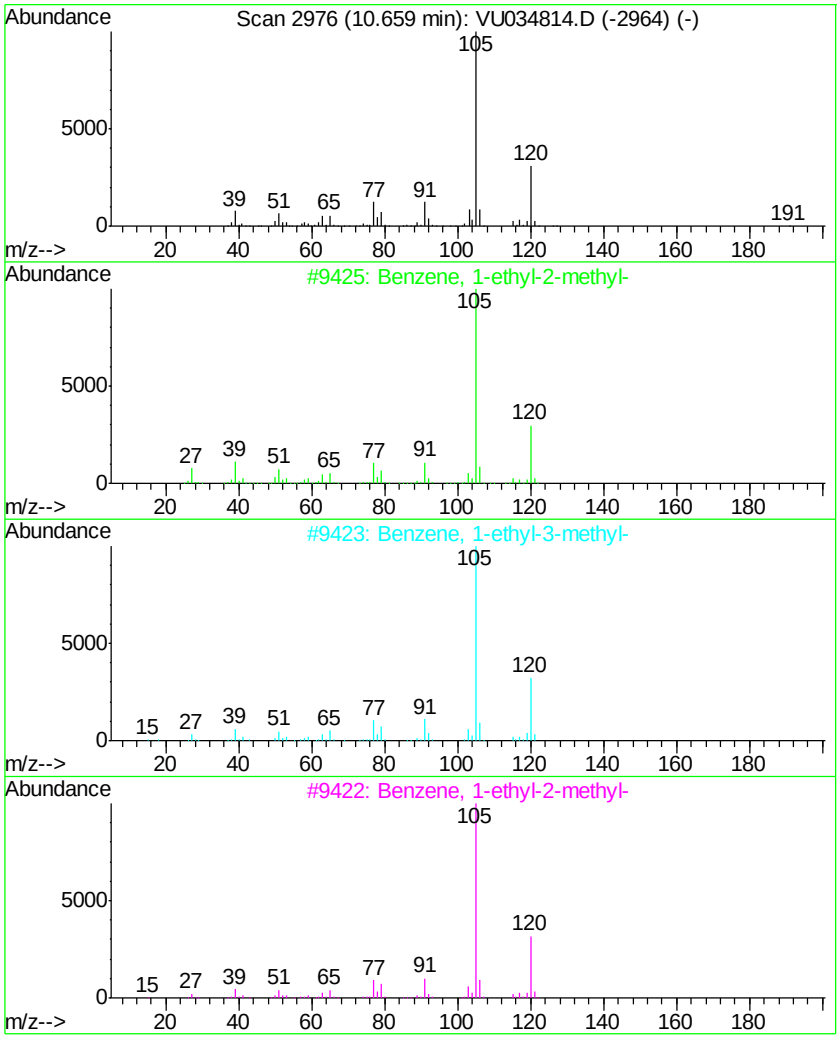
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

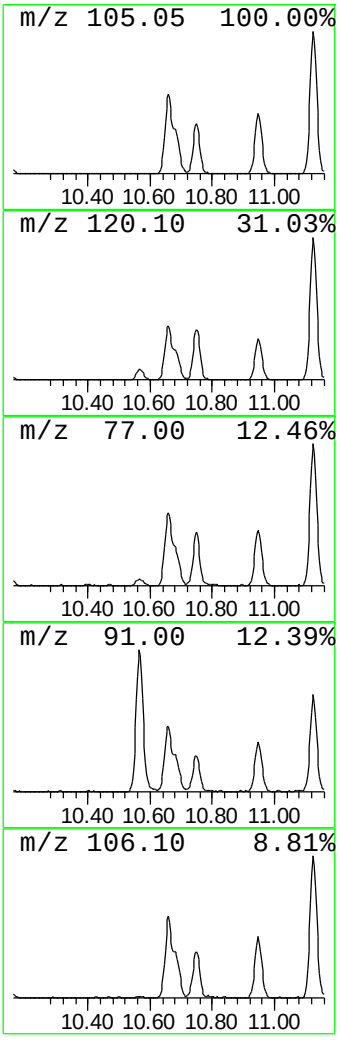
****

Peak Number 8 Benzene, 1-ethyl-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.66	39.73 ug/L	1031570	1,4-Dichlorobenzene-d4	11.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
3	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092519\
Data File : VU034814.D
Acq On : 25 Sep 2019 19:57
Operator : JC/SP
Sample : K4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BFDF4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM091919WMA.M
Quant Title : VOC Analysis

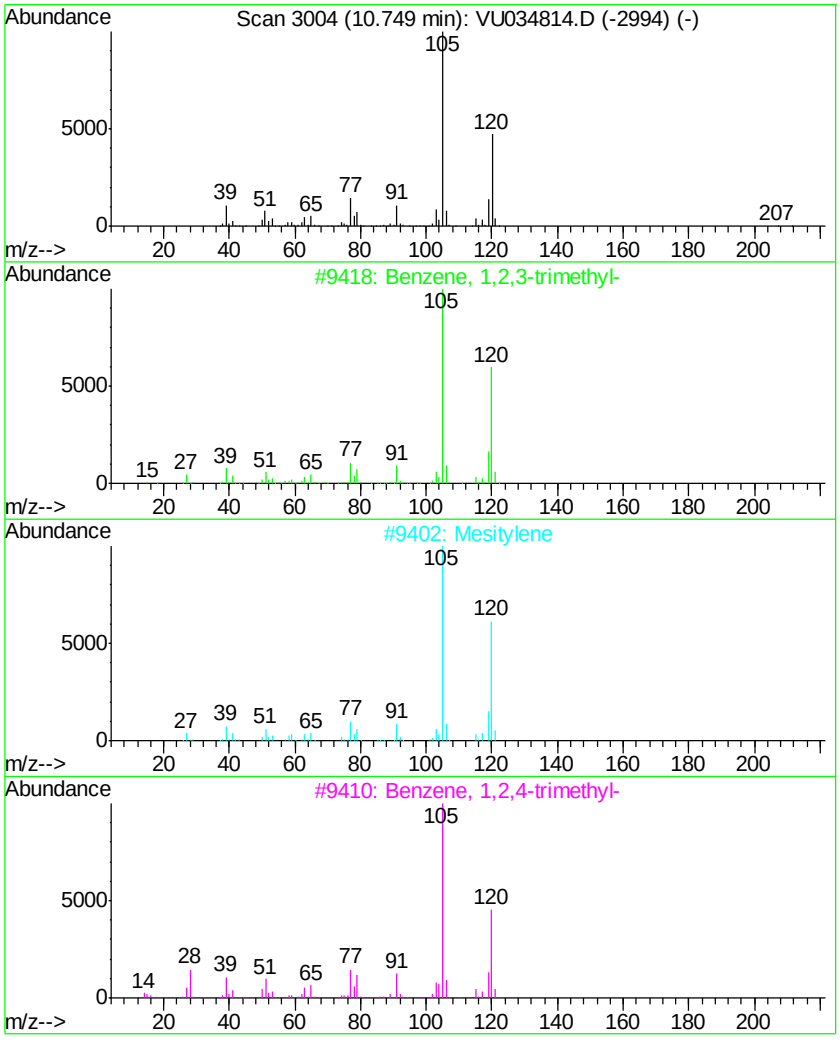
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

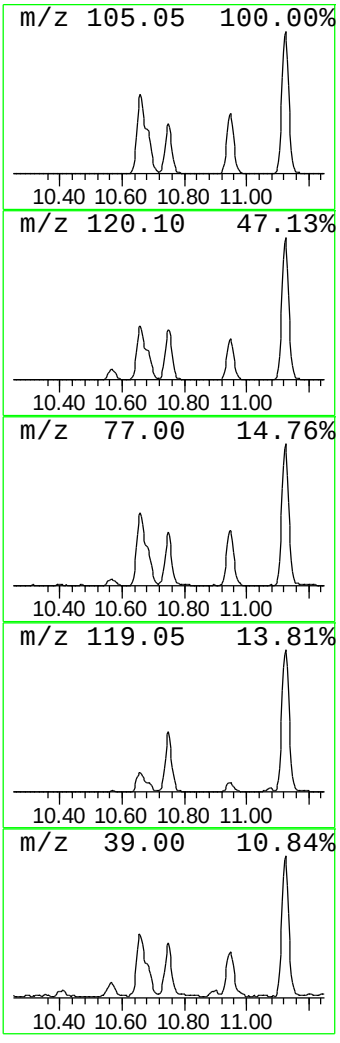
****

Peak Number 9 Benzene, 1,2,3-trimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.75	20.60 ug/L	534954	1,4-Dichlorobenzene-d4	11.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2	Mesitylene	120	C9H12	000108-67-8	94
3	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4	Mesitylene	120	C9H12	000108-67-8	91
5	Mesitylene	120	C9H12	000108-67-8	91





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092519\
Data File : VU034814.D
Acq On : 25 Sep 2019 19:57
Operator : JC/SP
Sample : K4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BFDF4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM091919WMA.M
Quant Title : VOC Analysis

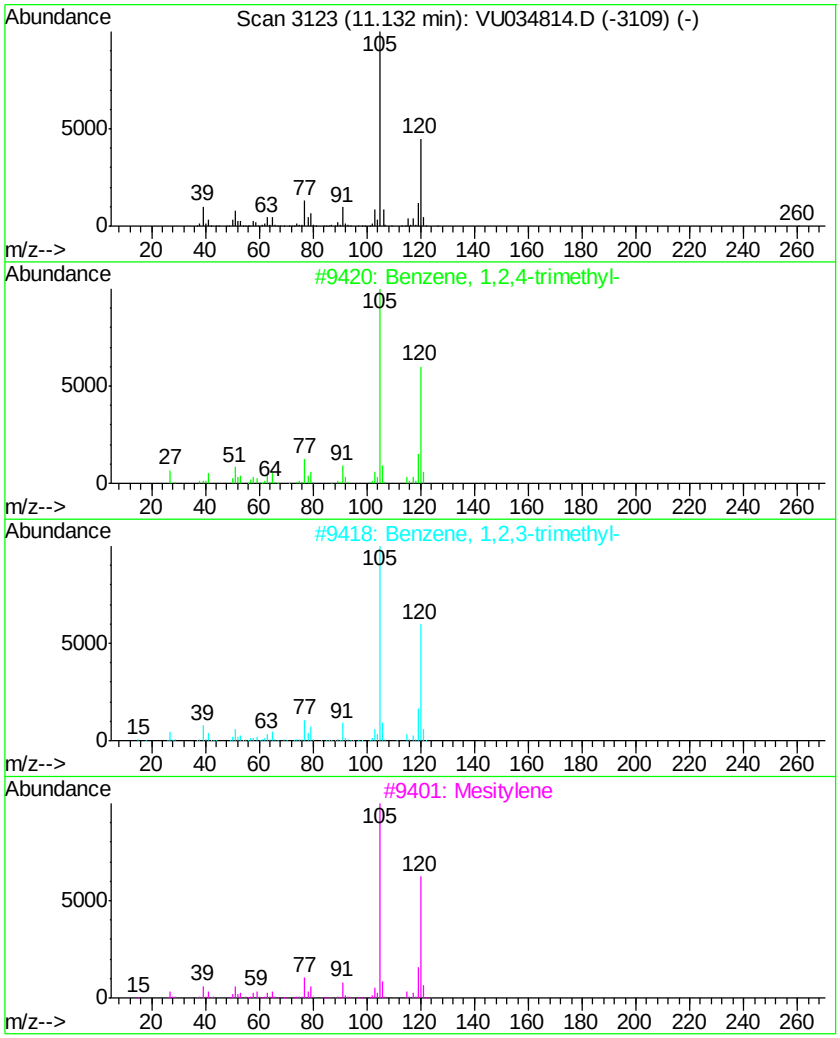
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

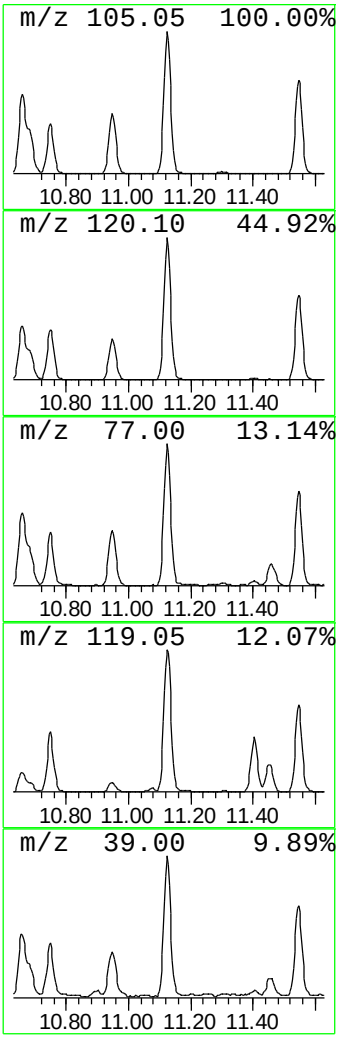
****

Peak Number 11 Benzene, 1,2,4-trimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.13	54.66 ug/L	1419250	1,4-Dichlorobenzene-d4	11.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	95
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3	Mesitylene	120	C9H12	000108-67-8	94
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94





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Data File : VU034814.D
Acq On : 25 Sep 2019 19:57
Operator : JC/SP
Sample : K4983-03
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BFDF4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM091919WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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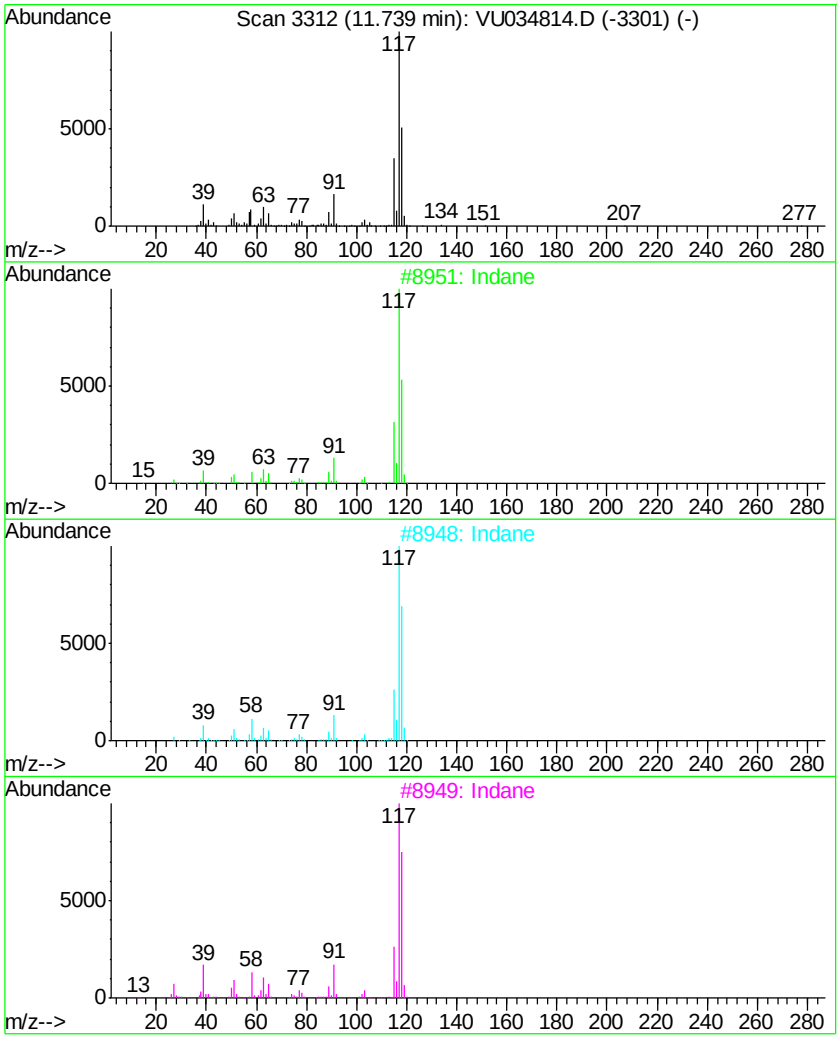
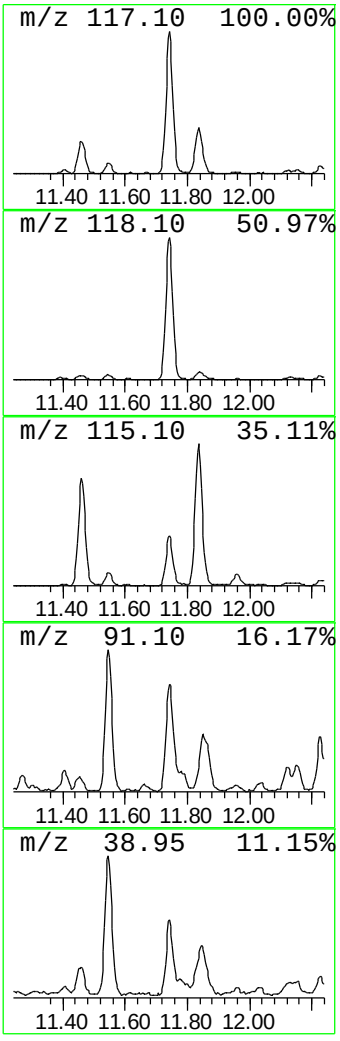
Peak Number 13 Indane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.74	22.78 ug/L	591402	1,4-Dichlorobenzene-d4	11.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	93
2	Indane	118	C9H10	000496-11-7	93
3	Indane	118	C9H10	000496-11-7	92
4	(Z)-1-Phenylpropene	118	C9H10	000766-90-5	81
5	Deltacyclene	118	C9H10	007785-10-6	68





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Data File : VU034814.D
Acq On : 25 Sep 2019 19:57
Operator : JC/SP
Sample : K4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BFDF4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM091919WMA.M
Quant Title : VOC Analysis

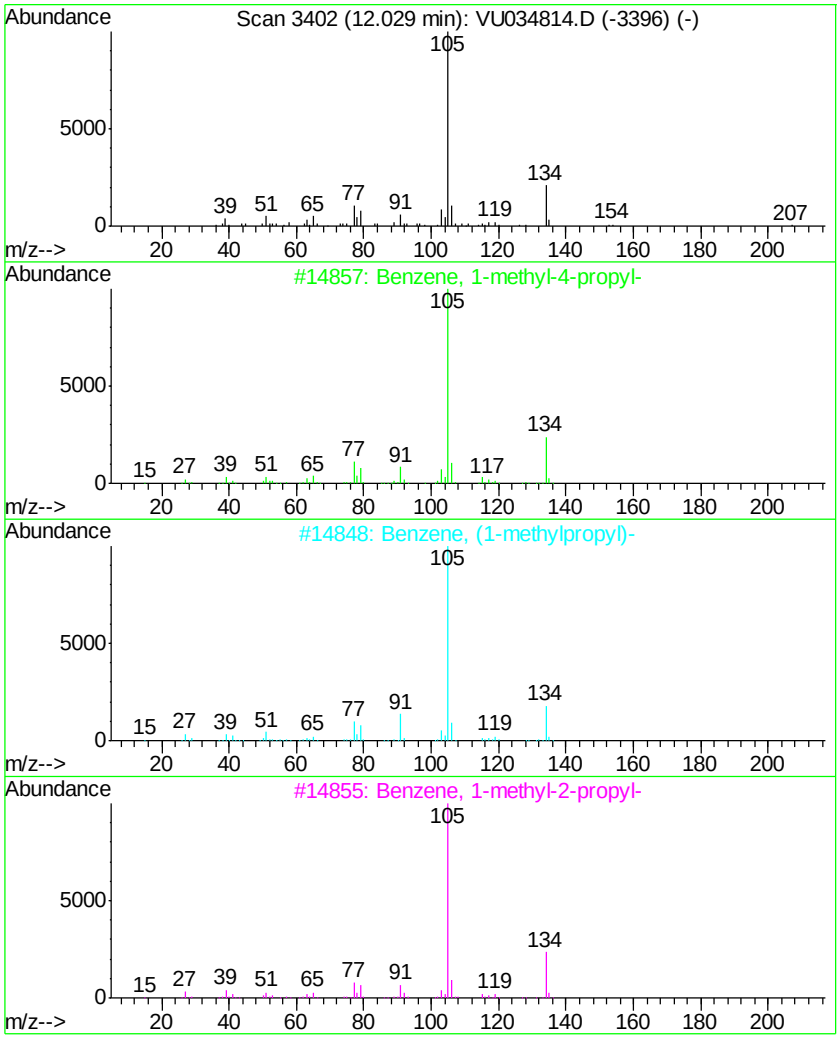
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

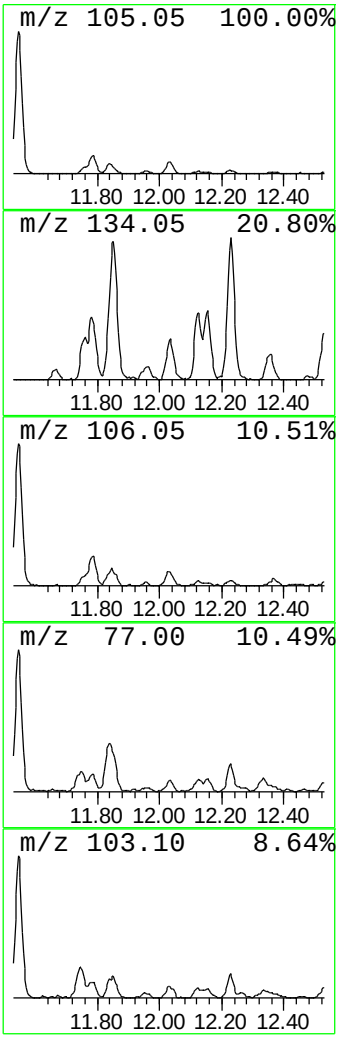
****

Peak Number 14 Benzene, 1-methyl-4-propyl- Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.03	2.88 ug/L	74769	1,4-Dichlorobenzene-d4	11.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
2	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	91
3	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	90
4	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	90
5	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	90





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Data File : VU034814.D
Acq On : 25 Sep 2019 19:57
Operator : JC/SP
Sample : K4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BFDF4

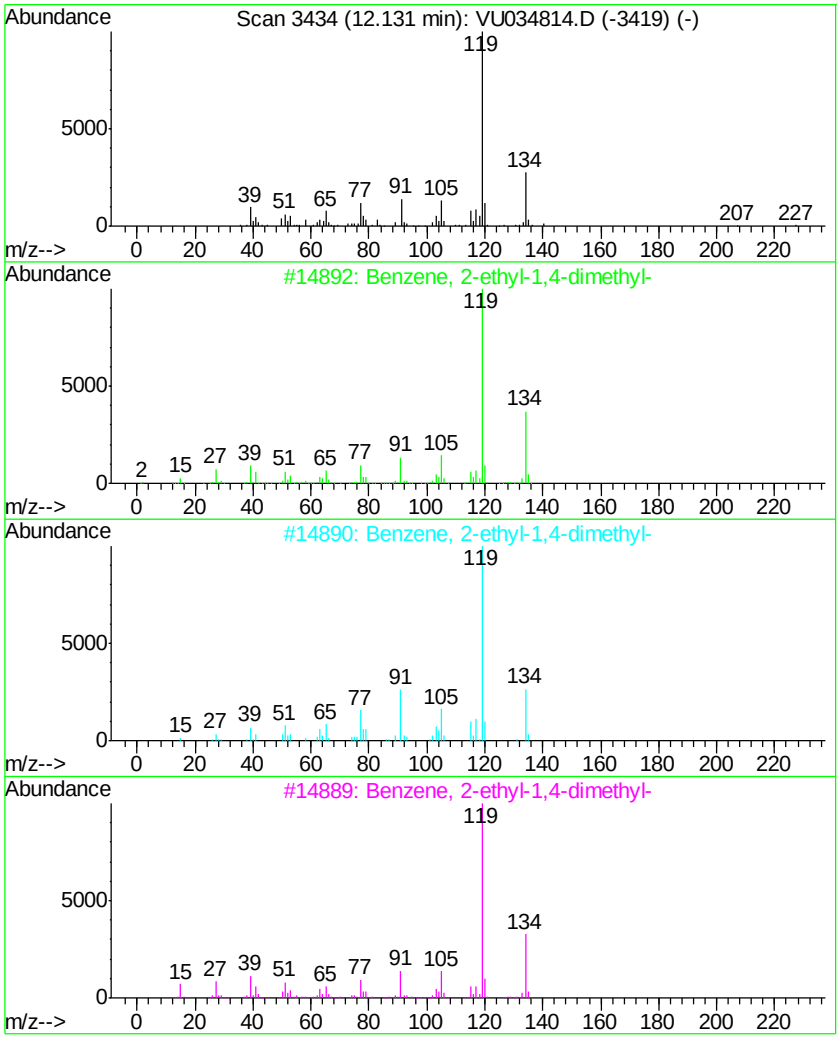
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM091919WMA.M
Quant Title : VOC Analysis

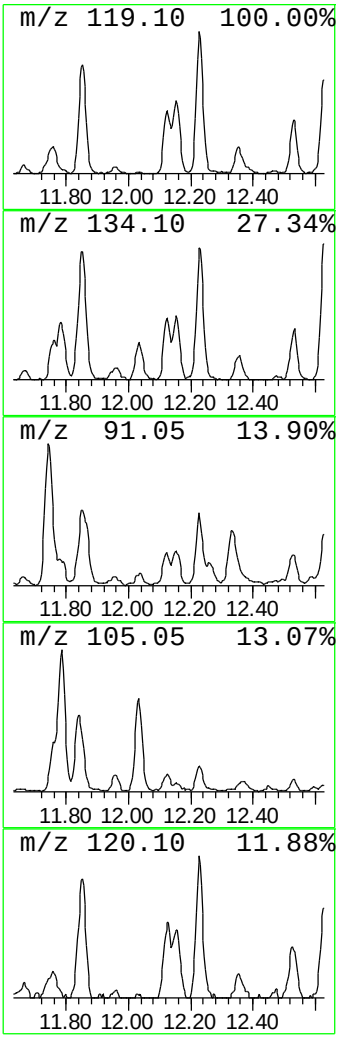
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.13	4.45 ug/L	115672	1,4-Dichlorobenzene-d4	11.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	97
2	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
3	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
4	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95





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Data File : VU034814.D
Acq On : 25 Sep 2019 19:57
Operator : JC/SP
Sample : K4983-03
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BFDF4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM091919WMA.M
Quant Title : VOC Analysis

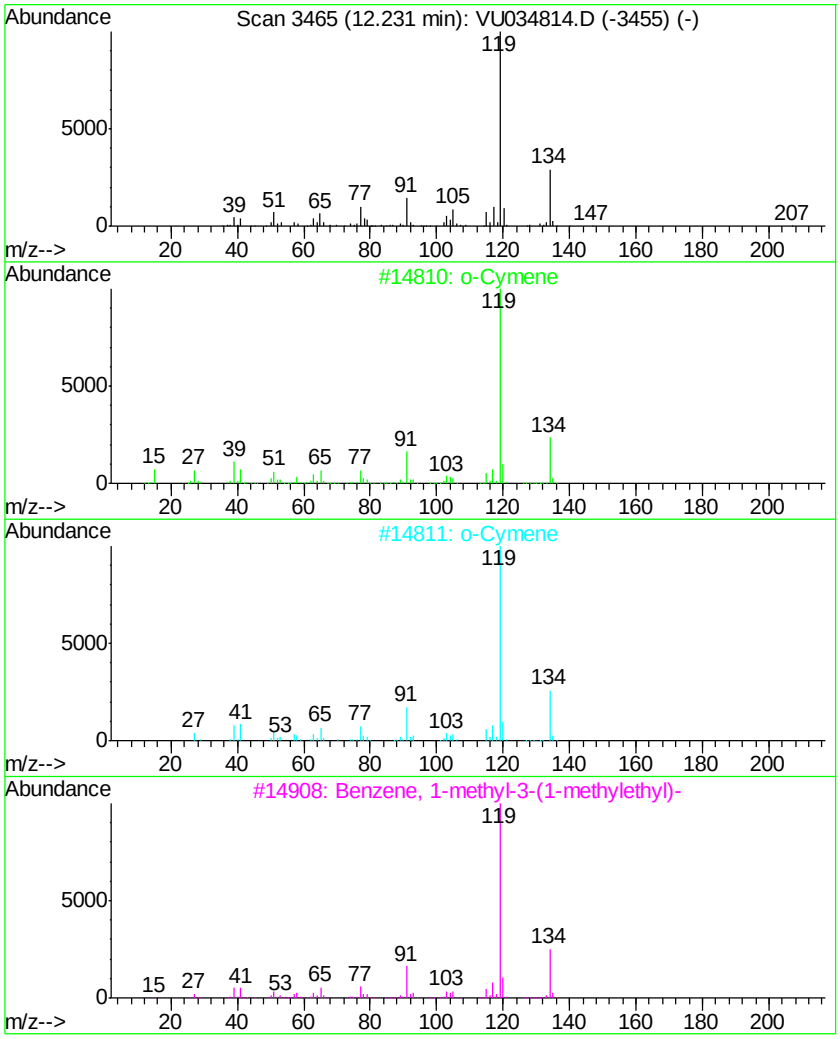
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

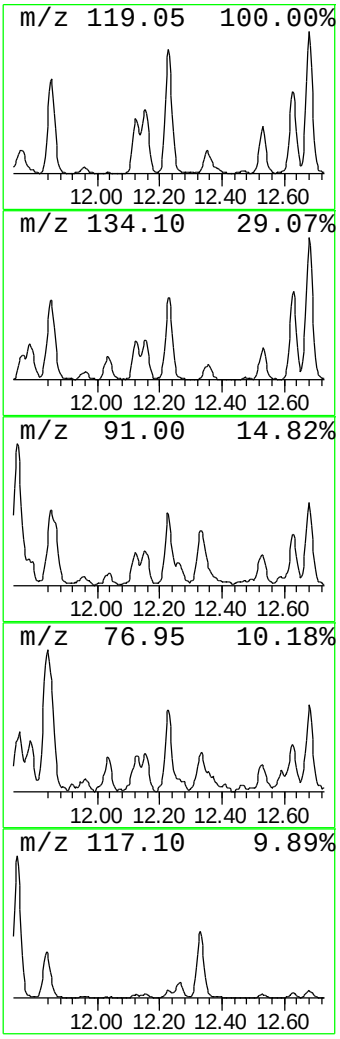
****

Peak Number 17 o-Cymene Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.23	7.81 ug/L	202774	1,4-Dichlorobenzene-d4	11.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	97
2	o-Cymene	134	C10H14	000527-84-4	97
3	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
4	p-Cymene	134	C10H14	000099-87-6	95
5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95





Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092519\
Data File : VU034814.D
Acq On : 25 Sep 2019 19:57
Operator : JC/SP
Sample : K4983-03
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BFDF4

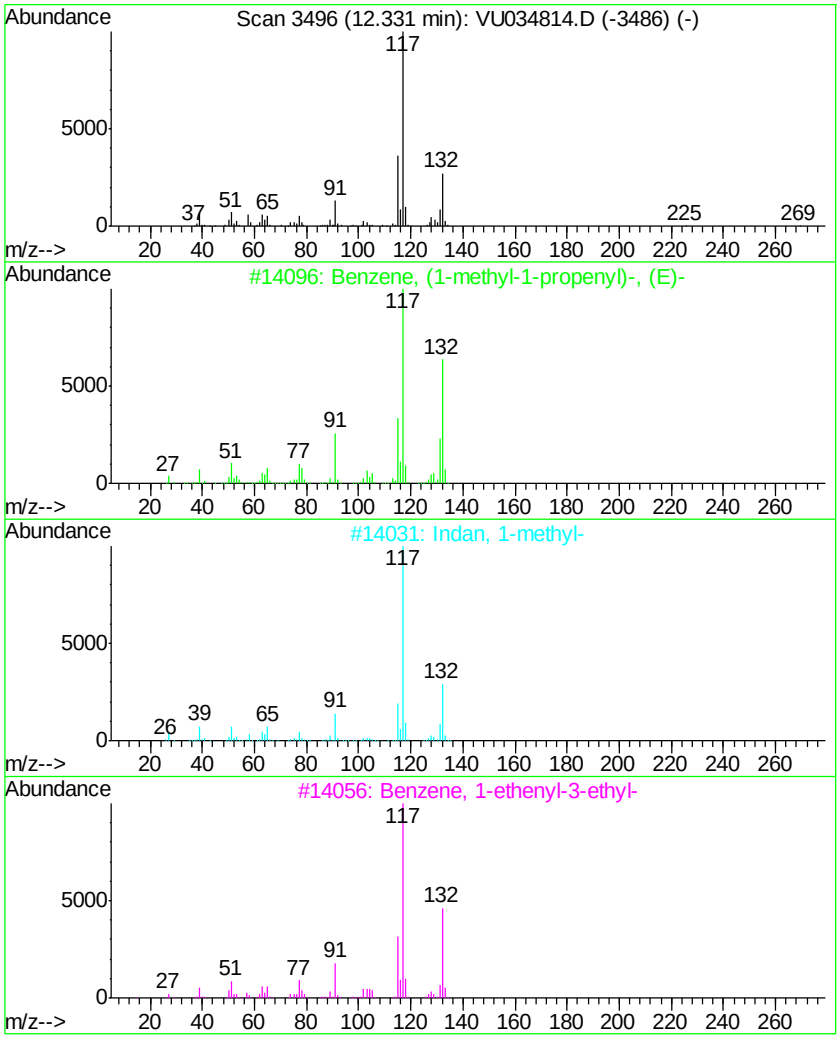
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM091919WMA.M
Quant Title : VOC Analysis

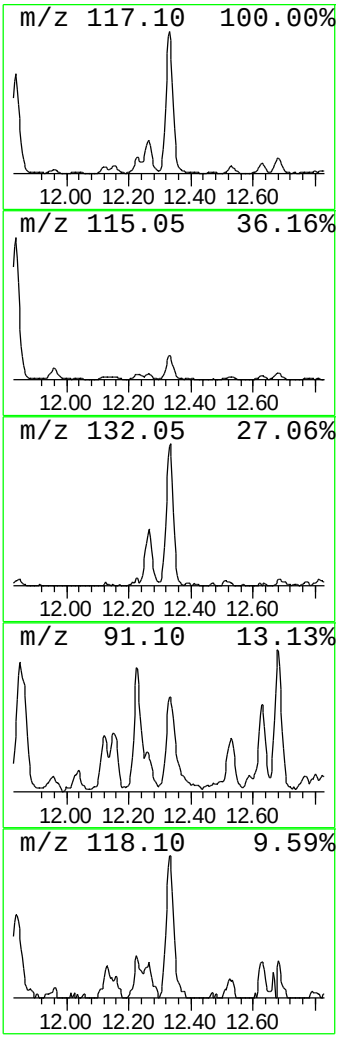
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 18 Benzene, (1-methyl-1-propen... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.33	10.12 ug/L	262663	1,4-Dichlorobenzene-d4	11.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	87
2	Indan, 1-methyl-	132	C10H12	000767-58-8	87
3	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	86
4	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	83
5	Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	80





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092519\
Data File : VU034814.D
Acq On : 25 Sep 2019 19:57
Operator : JC/SP
Sample : K4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BFDF4

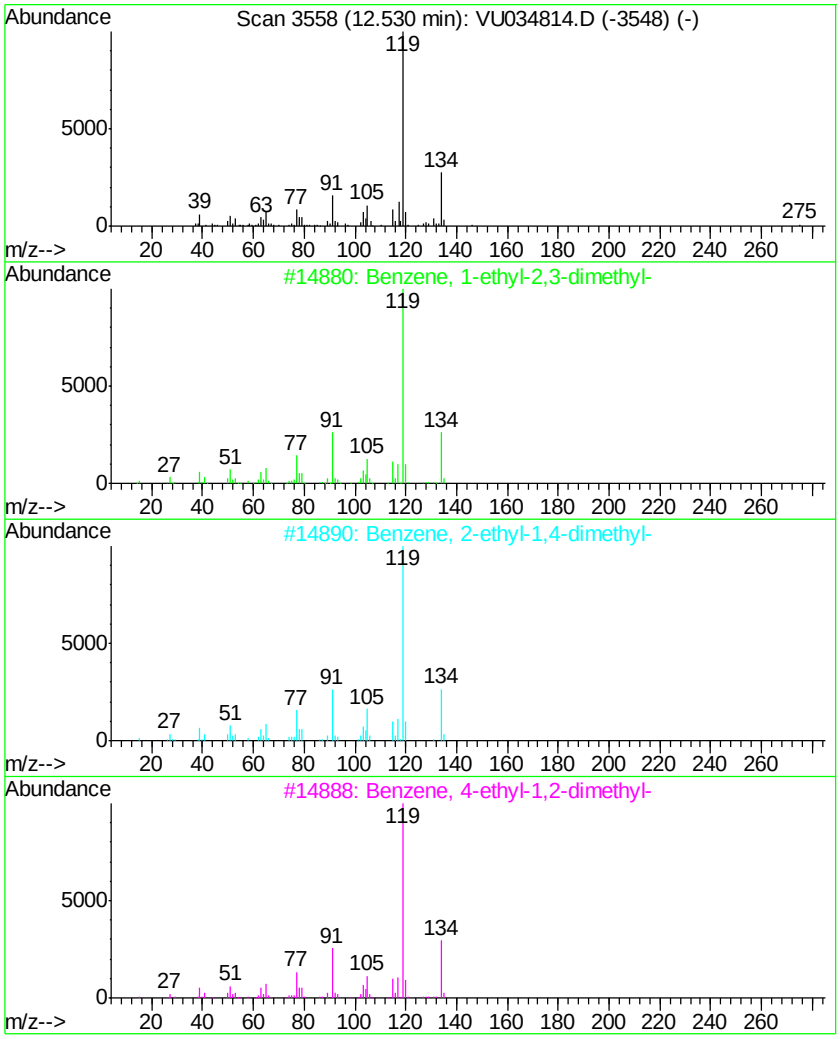
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM091919WMA.M
Quant Title : VOC Analysis

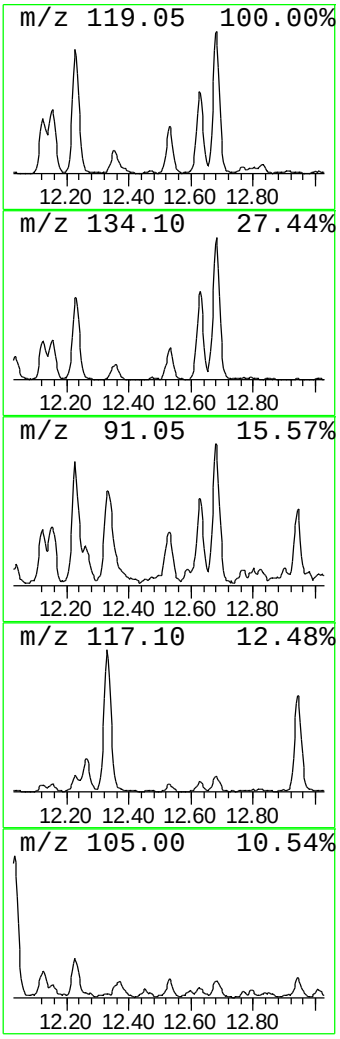
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.53	4.08 ug/L	106022	1,4-Dichlorobenzene-d4	11.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
2	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
3	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	93
4	1,4-Cyclohexadiene, 3-ethenyl-1,...	134	C10H14	062338-57-2	90
5	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	90





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Data File : VU034814.D
Acq On : 25 Sep 2019 19:57
Operator : JC/SP
Sample : K4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BFDF4

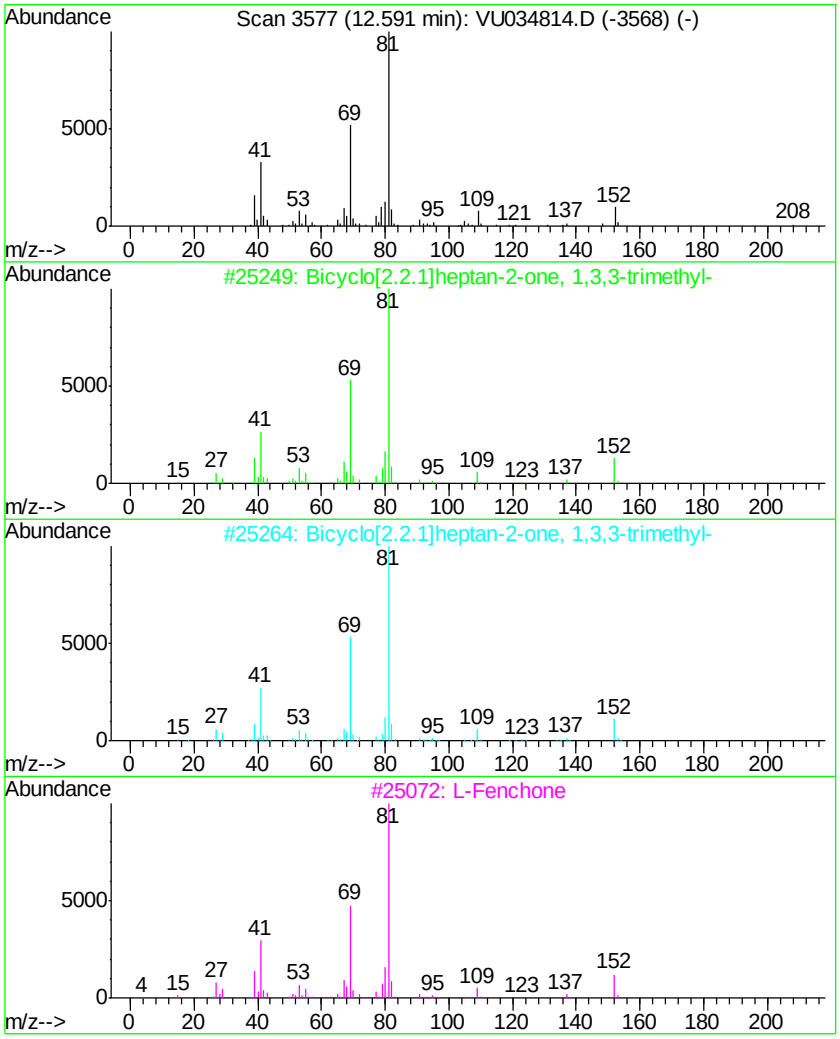
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM091919WMA.M
Quant Title : VOC Analysis

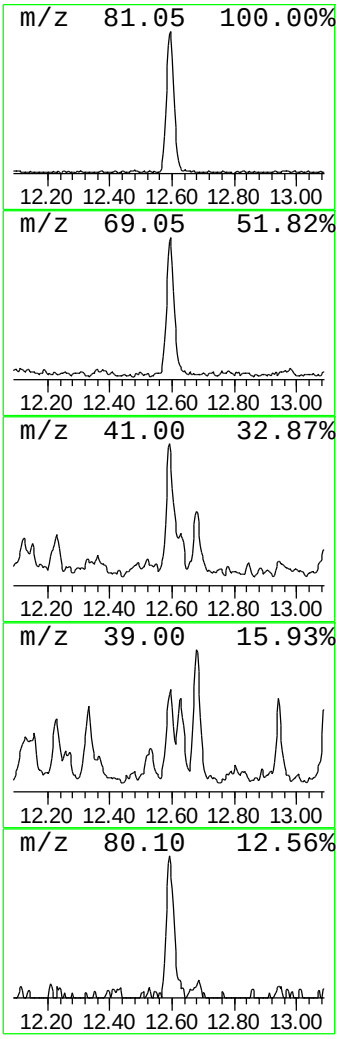
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 Bicyclo[2.2.1]heptan-2-one,... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.59	5.20 ug/L	134935	1,4-Dichlorobenzene-d4	11.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[2.2.1]heptan-2-one, 1,3,...	152	C10H16O	001195-79-5	95
2	Bicyclo[2.2.1]heptan-2-one, 1,3,...	152	C10H16O	001195-79-5	90
3	L-Fenchone	152	C10H16O	007787-20-4	90
4	Bicyclo[2.2.1]heptan-2-one, 1,3,...	152	C10H16O	001195-79-5	90
5	Bicyclo[2.2.1]heptan-2-one, 1,3,...	152	C10H16O	001195-79-5	80





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092519\
Data File : VU034814.D
Acq On : 25 Sep 2019 19:57
Operator : JC/SP
Sample : K4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BFDF4

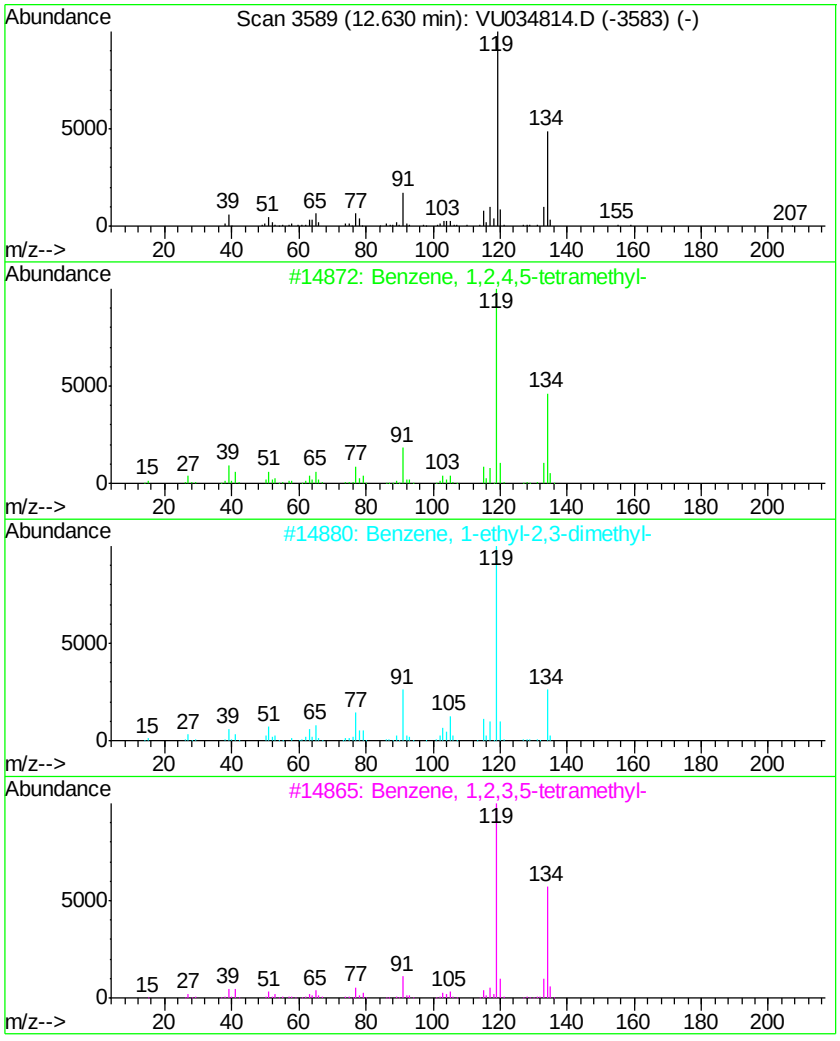
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM091919WMA.M
Quant Title : VOC Analysis

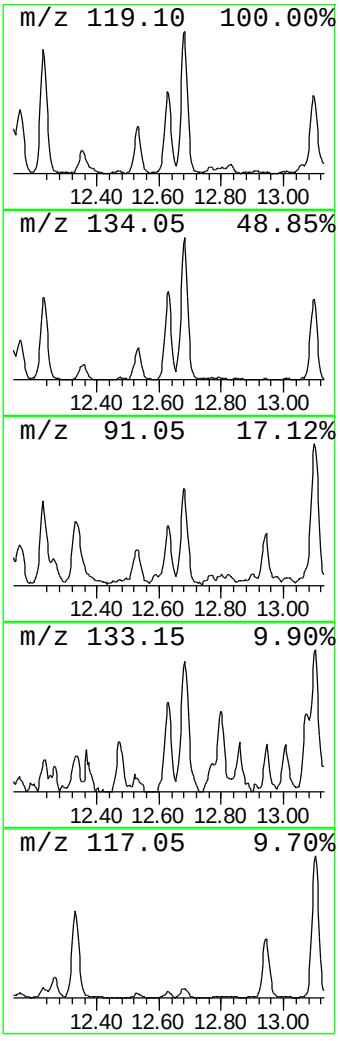
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 21 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.63	6.11 ug/L	158550	1,4-Dichlorobenzene-d4	11.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	94
2	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	93
3	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	91
4	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	91
5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	91





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092519\
Data File : VU034814.D
Acq On : 25 Sep 2019 19:57
Operator : JC/SP
Sample : K4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BFDF4

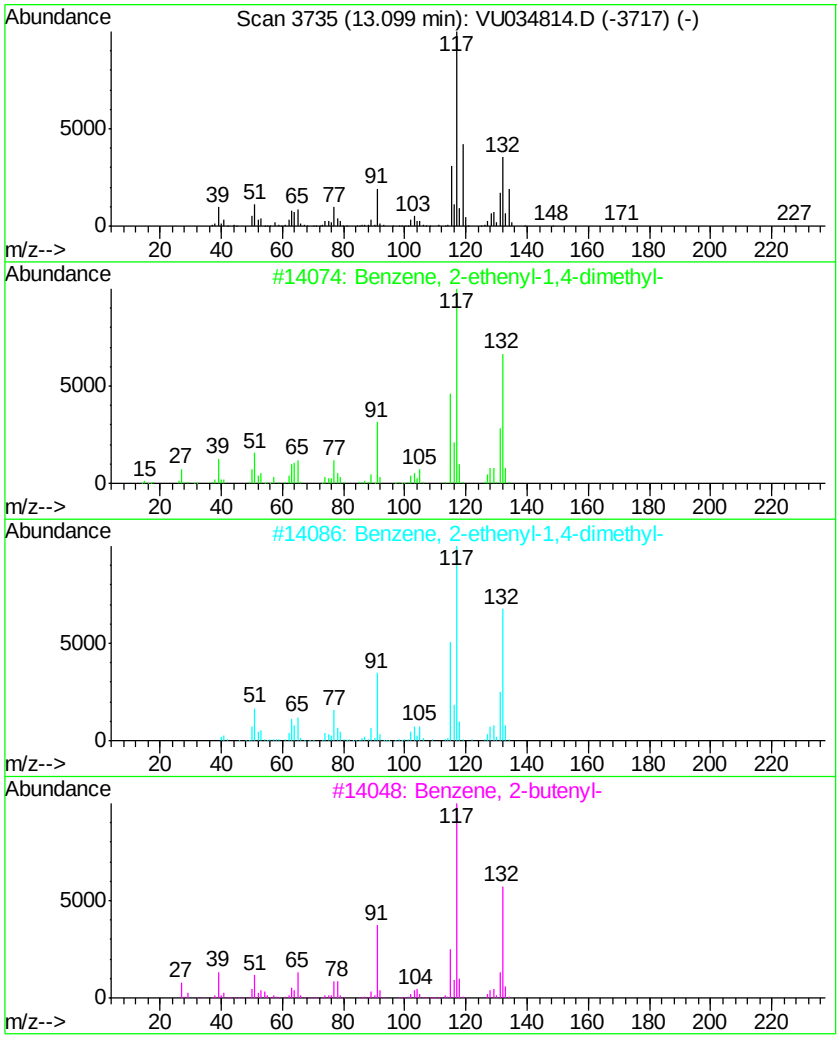
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM091919WMA.M
Quant Title : VOC Analysis

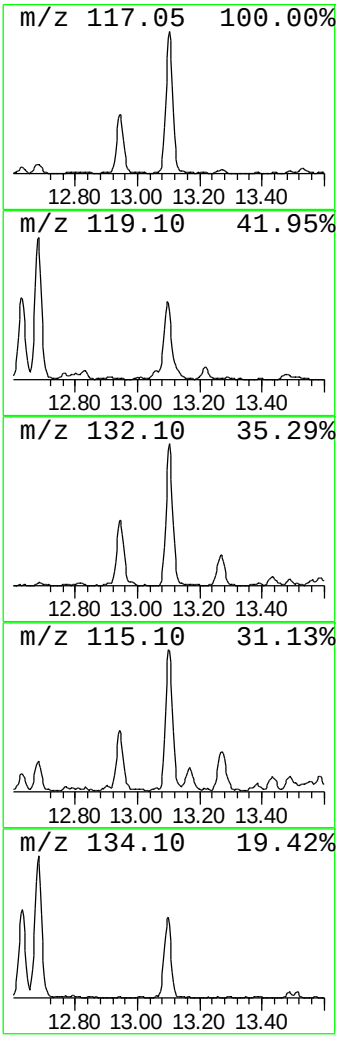
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.10	21.20 ug/L	550451	1,4-Dichlorobenzene-d4	11.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
2	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	87
3	Benzene, 2-butenyl-	132	C10H12	001560-06-1	70
4	3-Phenylbut-1-ene	132	C10H12	000934-10-1	70
5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	64





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Data File : VU034814.D
Acq On : 25 Sep 2019 19:57
Operator : JC/SP
Sample : K4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BFDF4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM091919WMA.M
Quant Title : VOC Analysis

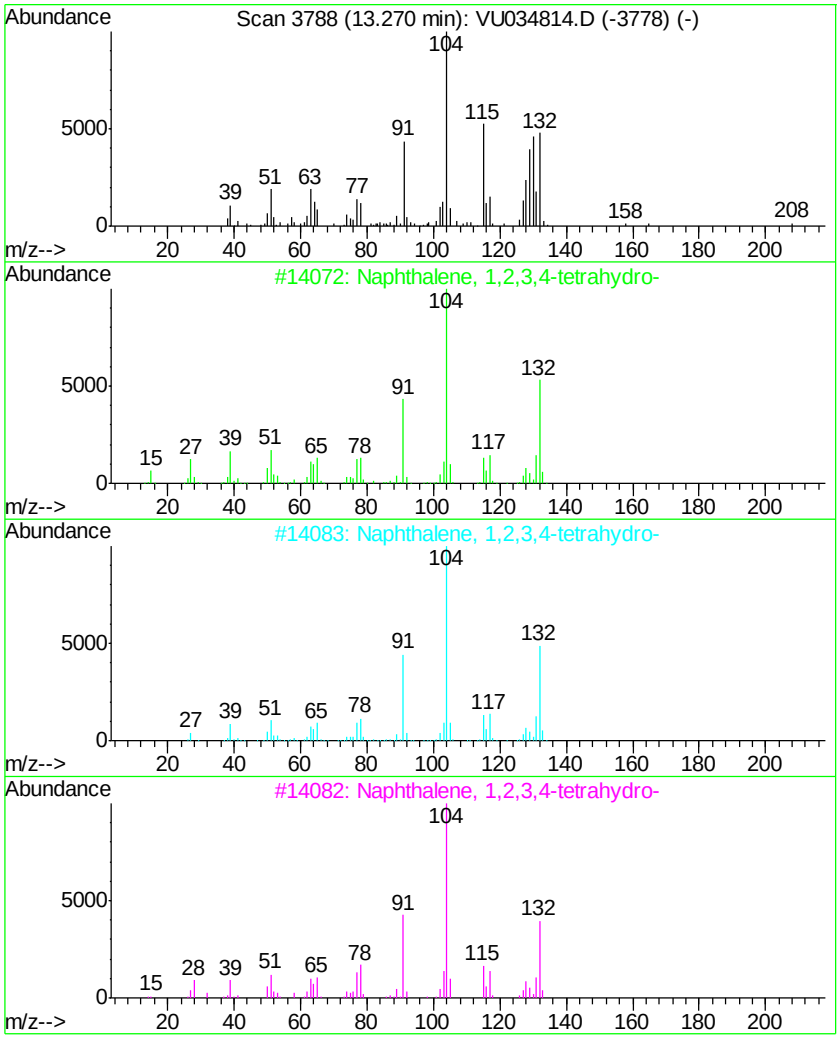
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

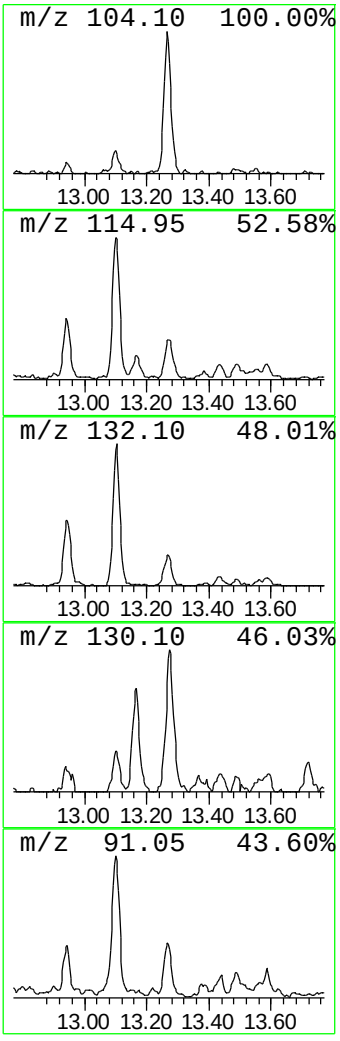
****

Peak Number 25 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.27	5.23 ug/L	135721	1,4-Dichlorobenzene-d4	11.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	70
2	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	55
3	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	50
4	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	50
5	Benzene,1-methyl-1,2-propadienyl-	130	C10H10	022433-39-2	42





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Data File : VU034814.D
Acq On : 25 Sep 2019 19:57
Operator : JC/SP
Sample : K4983-03
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BFDF4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM091919WMA.M
Quant Title : VOC Analysis

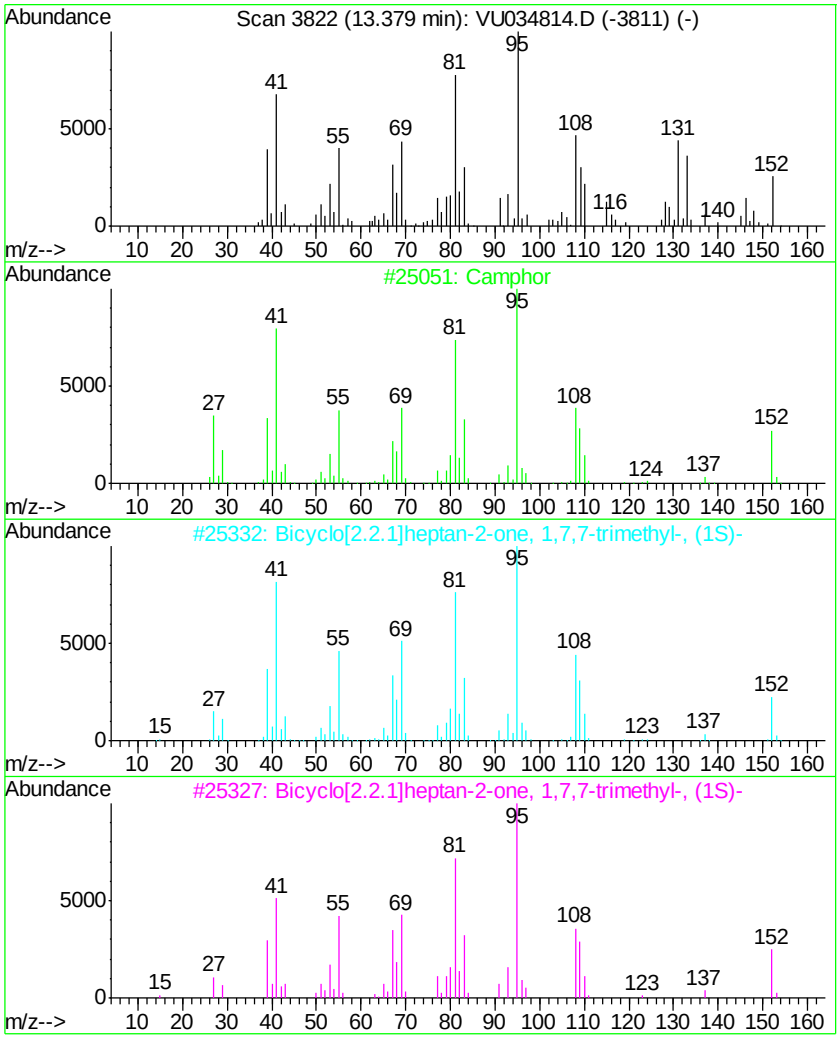
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

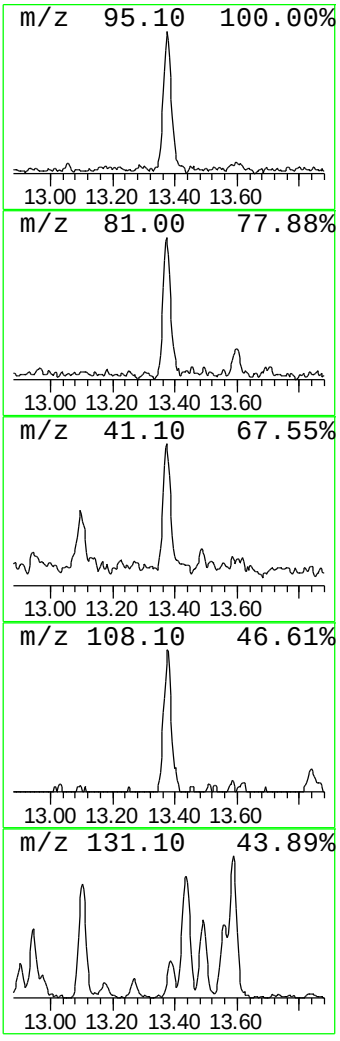
****

Peak Number 26 Camphor Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.38	4.95 ug/L	128662	1,4-Dichlorobenzene-d4	11.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Camphor	152	C10H16O	000076-22-2	93
2	Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-48-2	93
3	Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-48-2	89
4	Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-48-2	89
5	(+)-2-Bornanone	152	C10H16O	000464-49-3	89





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092519\
Data File : VU034814.D
Acq On : 25 Sep 2019 19:57
Operator : JC/SP
Sample : K4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BFDF4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM091919WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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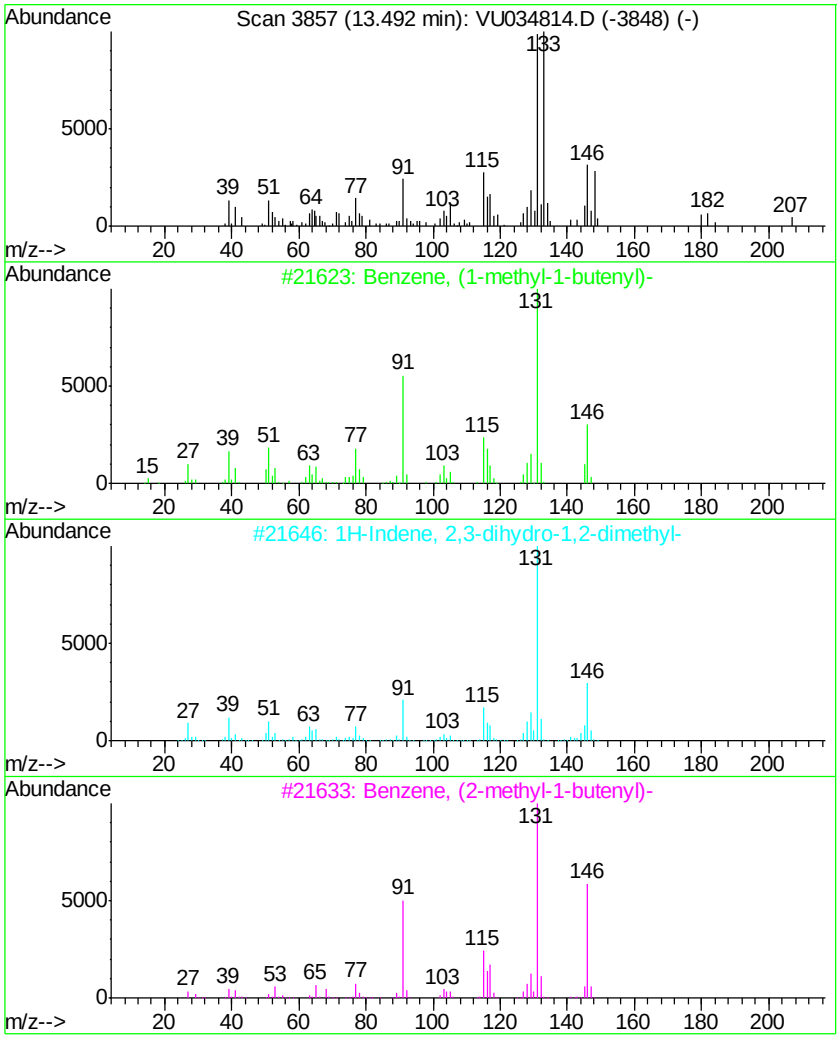
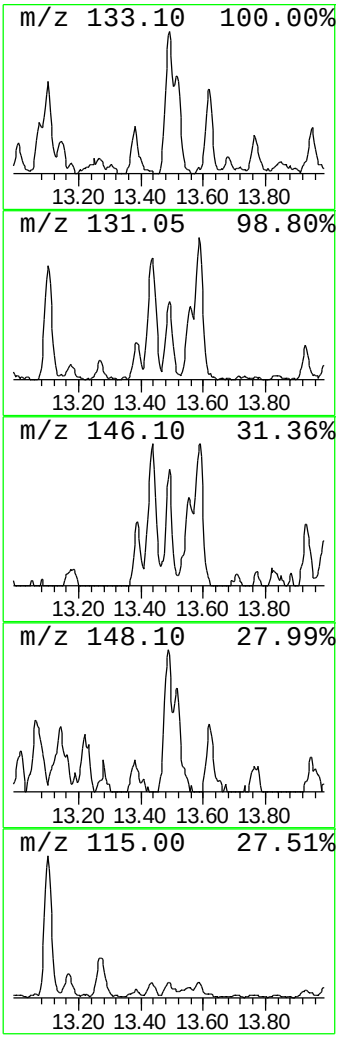
Peak Number 27 Benzene, (1-methyl-1-butenyl)- Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.49	3.07 ug/L	79843	1,4-Dichlorobenzene-d4	11.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	91
2	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	87
3	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	86
4	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	86
5	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	70





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092519\
Data File : VU034814.D
Acq On : 25 Sep 2019 19:57
Operator : JC/SP
Sample : K4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BFDF4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM091919WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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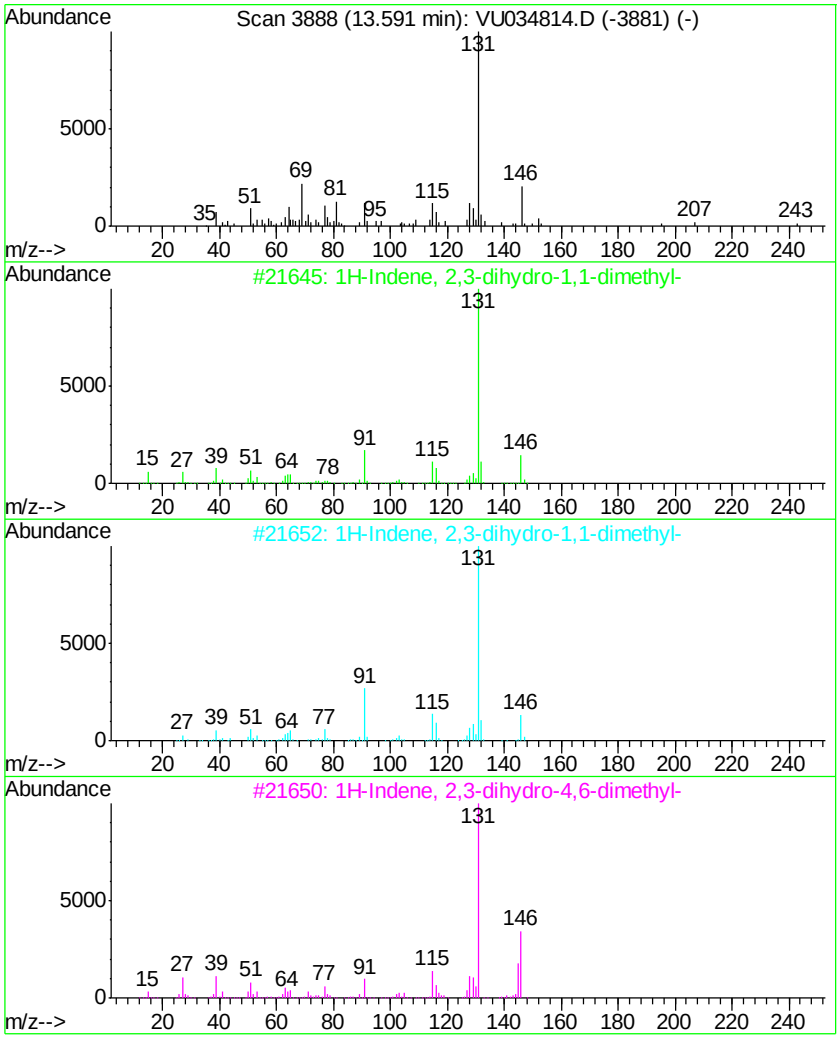
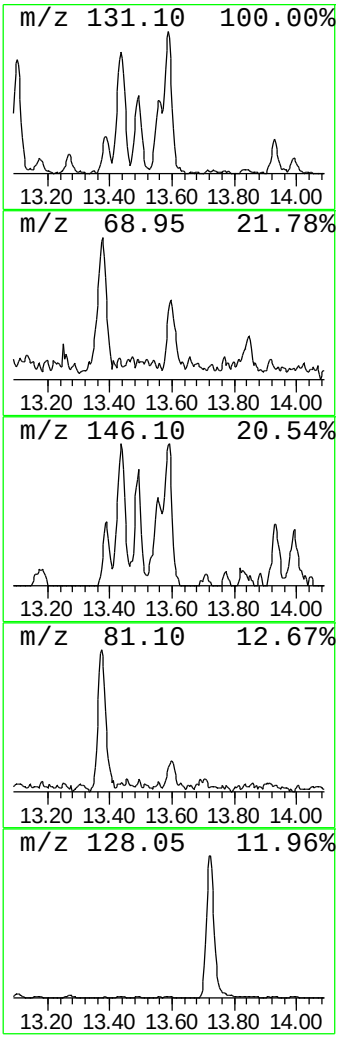
Peak Number 28 1H-Indene, 2,3-dihydro-1,1-... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.59	4.42 ug/L	114866	1,4-Dichlorobenzene-d4	11.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	80
2	1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	80
3	1H-Indene, 2,3-dihydro-4,6-dimet...	146	C11H14	001685-82-1	72
4	Benzene, 1-(chloromethyl)-4-(2-p...	166	C10H11Cl	036875-10-2	64
5	Cyclopropane, 2-bromo-1-methyl-1...	210	C10H11Br	055091-64-0	64





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092519\
Data File : VU034814.D
Acq On : 25 Sep 2019 19:57
Operator : JC/SP
Sample : K4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BFDF4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM091919WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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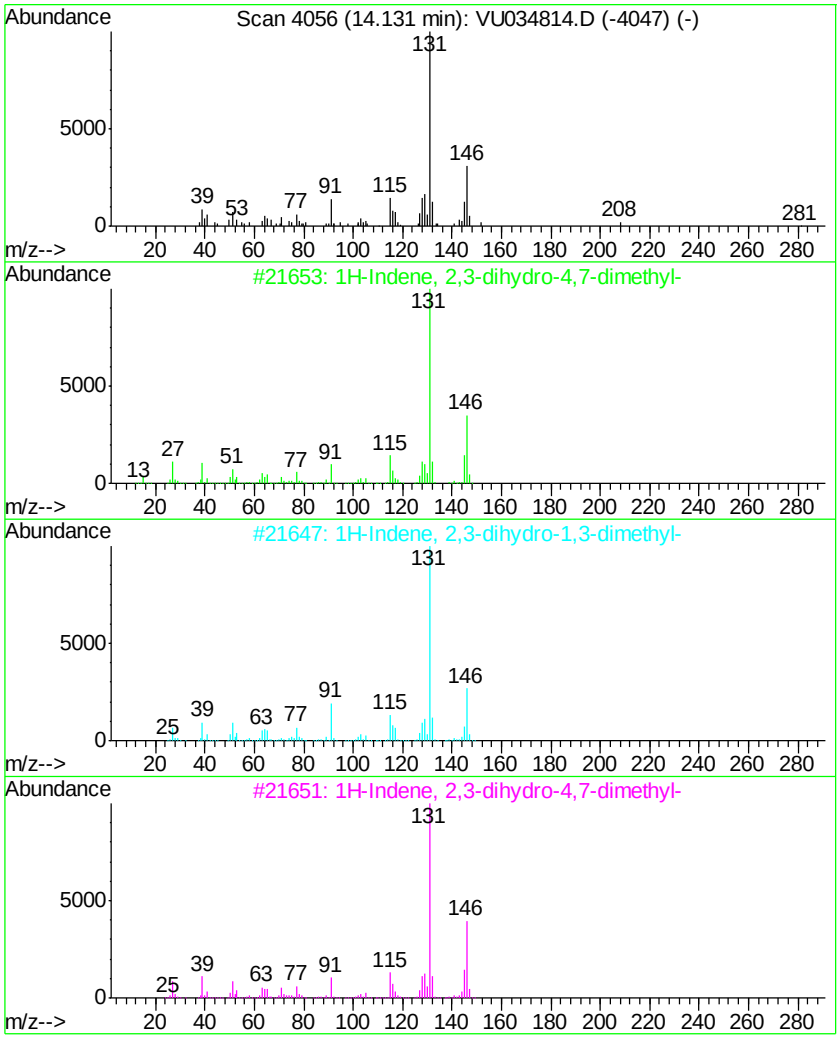
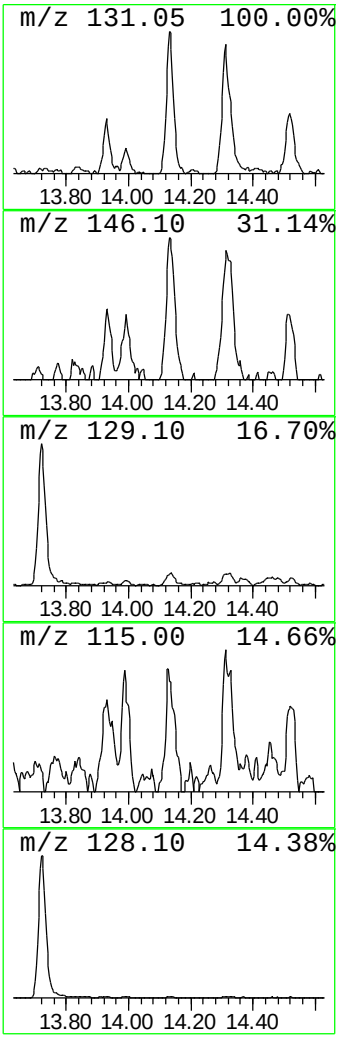
Peak Number 30 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.13	2.55 ug/L	66206	1,4-Dichlorobenzene-d4	11.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	96
2	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	95
3	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	94
4	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	91
5	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	91





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092519\
Data File : VU034814.D
Acq On : 25 Sep 2019 19:57
Operator : JC/SP
Sample : K4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BFDF4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM091919WMA.M
Quant Title : VOC Analysis

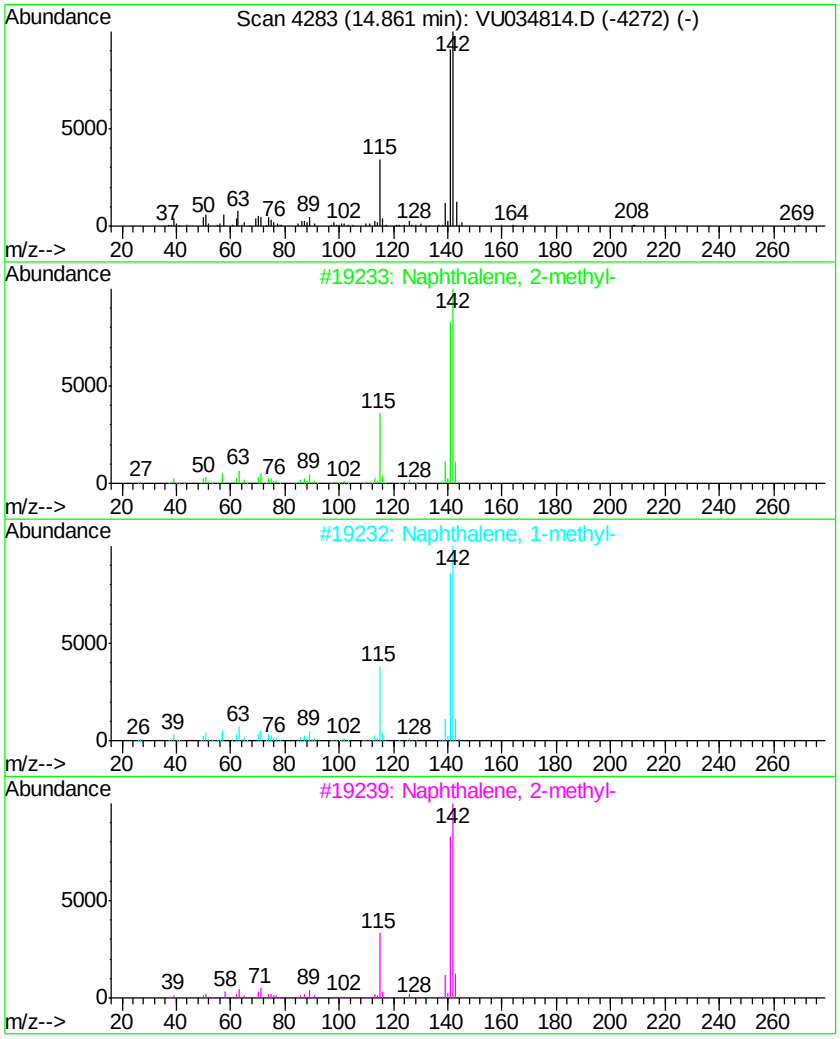
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

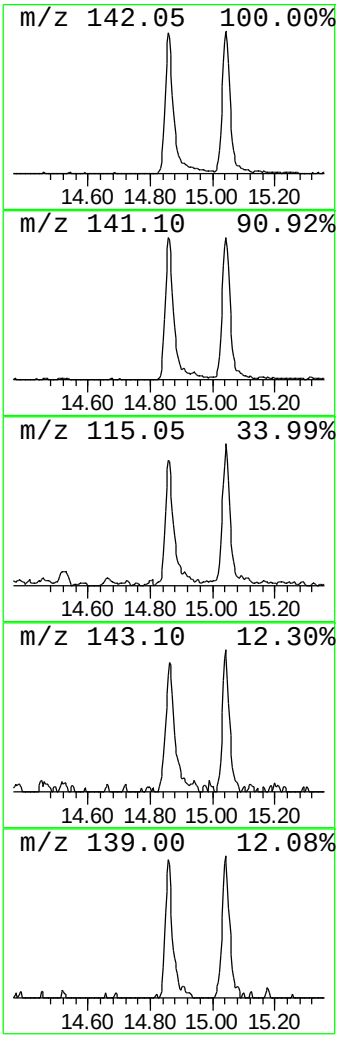
****

Peak Number 32 Naphthalene, 1-methyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.86	7.70 ug/L	199832	1,4-Dichlorobenzene-d4	11.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	95
4	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	94
5	Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	94





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092519\
Data File : VU034814.D
Acq On : 25 Sep 2019 19:57
Operator : JC/SP
Sample : K4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BFDF4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM091919WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Isopropyl acetate	5.53	7.9	ug/L	166484	1	5.86	1050440	50.0
n-Propyl acetate	6.68	131.3	ug/L	2758400	1	5.86	1050440	50.0
Propanoic acid, 2...	8.13	17.3	ug/L	486211	2	9.07	1409080	50.0
Propanoic acid, p...	8.40	26.3	ug/L	739921	2	9.07	1409080	50.0
Butanoic acid, 1-...	8.88	32.9	ug/L	925757	2	9.07	1409080	50.0
Benzene, propyl-	10.57	6.8	ug/L	176519	3	11.46	1298320	50.0
Benzene, 1-ethyl-...	10.66	39.7	ug/L	1031570	3	11.46	1298320	50.0
Benzene, 1,2,3-tr...	10.75	20.6	ug/L	534954	3	11.46	1298320	50.0
Benzene, 1,2,4-tr...	11.13	54.7	ug/L	1419250	3	11.46	1298320	50.0
Indane	11.74	22.8	ug/L	591402	3	11.46	1298320	50.0
Benzene, 1-methyl...	12.03	2.9	ug/L	74769	3	11.46	1298320	50.0
Benzene, 2-ethyl-...	12.13	4.5	ug/L	115672	3	11.46	1298320	50.0
o-Cymene	12.23	7.8	ug/L	202774	3	11.46	1298320	50.0
Benzene, (1-methy...	12.33	10.1	ug/L	262663	3	11.46	1298320	50.0
Benzene, 1-ethyl-...	12.53	4.1	ug/L	106022	3	11.46	1298320	50.0
Bicyclo[2.2.1]hep...	12.59	5.2	ug/L	134935	3	11.46	1298320	50.0
Benzene, 1,2,4,5-...	12.63	6.1	ug/L	158550	3	11.46	1298320	50.0
Benzene, 2-etheny...	13.10	21.2	ug/L	550451	3	11.46	1298320	50.0
Naphthalene, 1,2,...	13.27	5.2	ug/L	135721	3	11.46	1298320	50.0
Camphor	13.38	5.0	ug/L	128662	3	11.46	1298320	50.0
Benzene, (1-methy...	13.49	3.1	ug/L	79843	3	11.46	1298320	50.0
1H-Indene, 2,3-di...	13.59	4.4	ug/L	114866	3	11.46	1298320	50.0
1H-Indene, 2,3-di...	14.13	2.5	ug/L	66206	3	11.46	1298320	50.0
Naphthalene, 1-me...	14.86	7.7	ug/L	199832	3	11.46	1298320	50.0