LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120821\
Data File : VU046192.D
Acq On : 08 Dec 2021 21:50
Operator : SY/MD
Sample : M4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5N1

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\SFAMULM112921WMA.M
Title : VOC Analysis

Signal : TIC: VU046192.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.607	73	83	110	rBV	3334237	3724984	18.39%	2.513%
2	1.938	173	186	203	rBV	3010746	4062485	20.05%	2.741%
3	2.578	372	385	402	rBV3	211233	417525	2.06%	0.282%
4	2.809	449	457	469	rBV	7548	14986	0.07%	0.010%
5	3.051	522	532	542	rBV2	20036	39033	0.19%	0.026%
6	3.359	615	628	653	rVB	238330	473595	2.34%	0.319%
7	3.707	723	736	750	rVB6	3893	8838	0.04%	0.006%
8	3.877	767	789	819	rBV	5729624	13683941	67.55%	9.231%
9	4.671	1013	1036	1064	rBV	7367870	16750921	82.69%	11.300%
10	5.067	1146	1159	1182	rVB	135940	303174	1.50%	0.205%
11	5.321	1219	1238	1256	rBV	1528516	3418105	16.87%	2.306%
12	5.597	1317	1324	1334	rVB6	1572	2684	0.01%	0.002%
13	5.639	1334	1337	1344	rVB6	830	922	0.00%	0.001%
14	5.729	1344	1365	1375	rBV2	270000	744397	3.67%	0.502%
15	6.157	1484	1498	1513	rBV2	3998	10445	0.05%	0.007%
16	6.250	1513	1527	1543	rBV	245271	462438	2.28%	0.312%
17	6.546	1606	1619	1636	rBV	110033	204083	1.01%	0.138%
18	6.694	1651	1665	1678	rBV	171926	334062	1.65%	0.225%
19	6.764	1679	1687	1699	rVB3	21454	43319	0.21%	0.029%
20	6.925	1726	1737	1739	rBV3	1343	1890	0.01%	0.001%
21	6.983	1749	1755	1756	rBV4	1521	1423	0.01%	0.001%
22	7.256	1835	1840	1846	rVB2	1531	1764	0.01%	0.001%
23	7.568	1922	1937	1954	rVV	102762	181299	0.89%	0.122%
24	7.687	1964	1974	1979	rBV2	1104	1613	0.01%	0.001%
25	7.793	1995	2007	2020	rBV2	28734	49954	0.25%	0.034%
26	7.899	2020	2040	2050	rBV	350555	604280	2.98%	0.408%
27	7.970	2051	2062	2077	rVV	1391773	2328715	11.50%	1.571%
28	8.179	2116	2127	2139	rBV	70686	117669	0.58%	0.079%
29	8.240	2140	2146	2160	rVB6	3879	8154	0.04%	0.006%
30	8.401	2183	2196	2209	rVB2	30170	49741	0.25%	0.034%
31	8.475	2209	2219	2234	rVB3	4180	9118	0.05%	0.006%
32	8.555	2234	2244	2254	rVB2	19710	32613	0.16%	0.022%
33	8.636	2254	2269	2308	rBV	302294	556551	2.75%	0.375%
34	8.867	2335	2341	2354	rVB7	3012	4872	0.02%	0.003%
35	9.095	2406	2412	2414	rBV2	1173	1012	0.00%	0.001%
36	9.275	2455	2468	2477	rBV3	3832	6487	0.03%	0.004%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120821\
Data File : VU046192.D
Acq On : 08 Dec 2021 21:50
Operator : SY/MD
Sample : M4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5N1

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\SFAMULM112921WMA.M
Title : VOC Analysis

37	9.349	2484	2491	2498	rVV5	1432	2471	0.01%	0.002%
38	9.417	2500	2512	2533	rVV	354757	593277	2.93%	0.400%
39	9.510	2536	2541	2547	rVV5	4165	6018	0.03%	0.004%
40	9.571	2547	2560	2584	rVV	3159091	4986893	24.62%	3.364%
41	9.697	2584	2599	2633	rVB	7224993	11991218	59.19%	8.089%
42	9.918	2660	2668	2677	rVB4	6018	9765	0.05%	0.007%
43	10.102	2710	2725	2752	rVB	3927829	6299930	31.10%	4.250%
44	10.234	2761	2766	2772	rBV3	627	1020	0.01%	0.001%
45	10.275	2772	2779	2789	rVB5	2478	4084	0.02%	0.003%
46	10.346	2794	2801	2807	rBV4	2737	3477	0.02%	0.002%
47	10.484	2832	2844	2867	rVB	413397	639533	3.16%	0.431%
48	10.578	2867	2873	2874	rBV2	1100	1073	0.01%	0.001%
49	10.600	2874	2880	2887	rBV2	3392	4250	0.02%	0.003%
50	10.645	2887	2894	2900	rBV2	5611	7553	0.04%	0.005%
51	10.758	2917	2929	2942	rBV	287346	467077	2.31%	0.315%
52	10.906	2959	2975	2990	rBV	1592457	2444719	12.07%	1.649%
53	10.999	2990	3004	3022	rVV	5229825	11671540	57.61%	7.874%
54	11.089	3022	3032	3049	rVB	2856461	4383498	21.64%	2.957%
55	11.234	3066	3077	3083	rBV	73580	111821	0.55%	0.075%
56	11.295	3083	3096	3117	rVB	3197847	5108638	25.22%	3.446%
57	11.420	3123	3135	3137	rBV5	8603	11373	0.06%	0.008%
58	11.471	3137	3151	3167	rVV	13020637	20257935	100.00%	13.666%
59	11.558	3171	3178	3185	rVV3	8564	15074	0.07%	0.010%
60	11.613	3186	3195	3198	rVV	30626	44487	0.22%	0.030%
61	11.645	3199	3205	3218	rVB	65987	102119	0.50%	0.069%
62	11.745	3223	3236	3244	rBV	181143	279133	1.38%	0.188%
63	11.812	3244	3257	3270	rVB2	431796	768239	3.79%	0.518%
64	11.896	3270	3283	3298	rBV	4174429	6458430	31.88%	4.357%
65	12.012	3312	3319	3329	rVB2	27046	40507	0.20%	0.027%
66	12.095	3331	3345	3364	rBV2	1735794	3116791	15.39%	2.103%
67	12.192	3364	3375	3396	rVB5	815280	1699781	8.39%	1.147%
68	12.304	3400	3410	3420	rBV3	51730	88615	0.44%	0.060%
69	12.378	3421	3433	3444	rVB	156187	258193	1.27%	0.174%
70	12.484	3447	3466	3485	rBV	7527099	13139666	64.86%	8.864%
71	12.571	3485	3493	3502	rVB	395486	563258	2.78%	0.380%
72	12.684	3518	3528	3547	rVB3	159510	342771	1.69%	0.231%
73	12.809	3563	3567	3575	rVB4	7588	10295	0.05%	0.007%
74	12.877	3577	3588	3601	rBV	154941	239482	1.18%	0.162%
75	12.973	3602	3618	3626	rBV2	252813	437849	2.16%	0.295%
76	13.028	3626	3635	3649	rVB	393499	604787	2.99%	0.408%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120821\
Data File : VU046192.D
Acq On : 08 Dec 2021 21:50
Operator : SY/MD
Sample : M4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5N1

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\SFAMULM112921WMA.M
Title : VOC Analysis

77	13.108	3652	3660	3664	rBV4	7990	11780	0.06%	0.008%
78	13.295	3709	3718	3729	rVB	120129	179277	0.88%	0.121%
79	13.352	3729	3736	3745	rVB5	6041	8843	0.04%	0.006%
80	13.452	3756	3767	3779	rVB3	404496	674921	3.33%	0.455%
81	13.516	3780	3787	3796	rVB3	26563	40601	0.20%	0.027%
82	13.626	3809	3821	3836	rVB2	96607	161732	0.80%	0.109%
83	13.738	3844	3856	3865	rBV3	52801	89902	0.44%	0.061%
84	13.835	3878	3886	3903	rVB7	18880	43961	0.22%	0.030%
85	13.912	3903	3910	3914	rBV3	4226	6239	0.03%	0.004%
86	14.005	3934	3939	3950	rVB5	6604	10589	0.05%	0.007%
87	14.086	3950	3964	3980	rBV	619603	974600	4.81%	0.657%
88	14.291	4016	4028	4036	rBV7	5204	9414	0.05%	0.006%
89	14.346	4038	4045	4057	rVB7	2382	4648	0.02%	0.003%
90	14.484	4081	4088	4095	rBV3	3614	4858	0.02%	0.003%
91	14.574	4110	4116	4126	rVB3	2957	4783	0.02%	0.003%
92	14.677	4136	4148	4151	rBV5	5040	9316	0.05%	0.006%
93	14.745	4163	4169	4170	rBV4	1036	902	0.00%	0.001%
94	14.809	4182	4189	4196	rBV2	3615	4842	0.02%	0.003%
95	14.876	4203	4210	4218	rVB6	2927	3905	0.02%	0.003%
96	15.018	4247	4254	4259	rVV6	2221	3119	0.02%	0.002%
97	15.221	4306	4317	4331	rVB	40051	61056	0.30%	0.041%
98	15.410	4365	4376	4386	rVB	26374	42253	0.21%	0.029%
99	16.044	4567	4573	4581	rBV3	693	1132	0.01%	0.001%
100	16.770	4793	4799	4808	rVB6	5492	6679	0.03%	0.005%

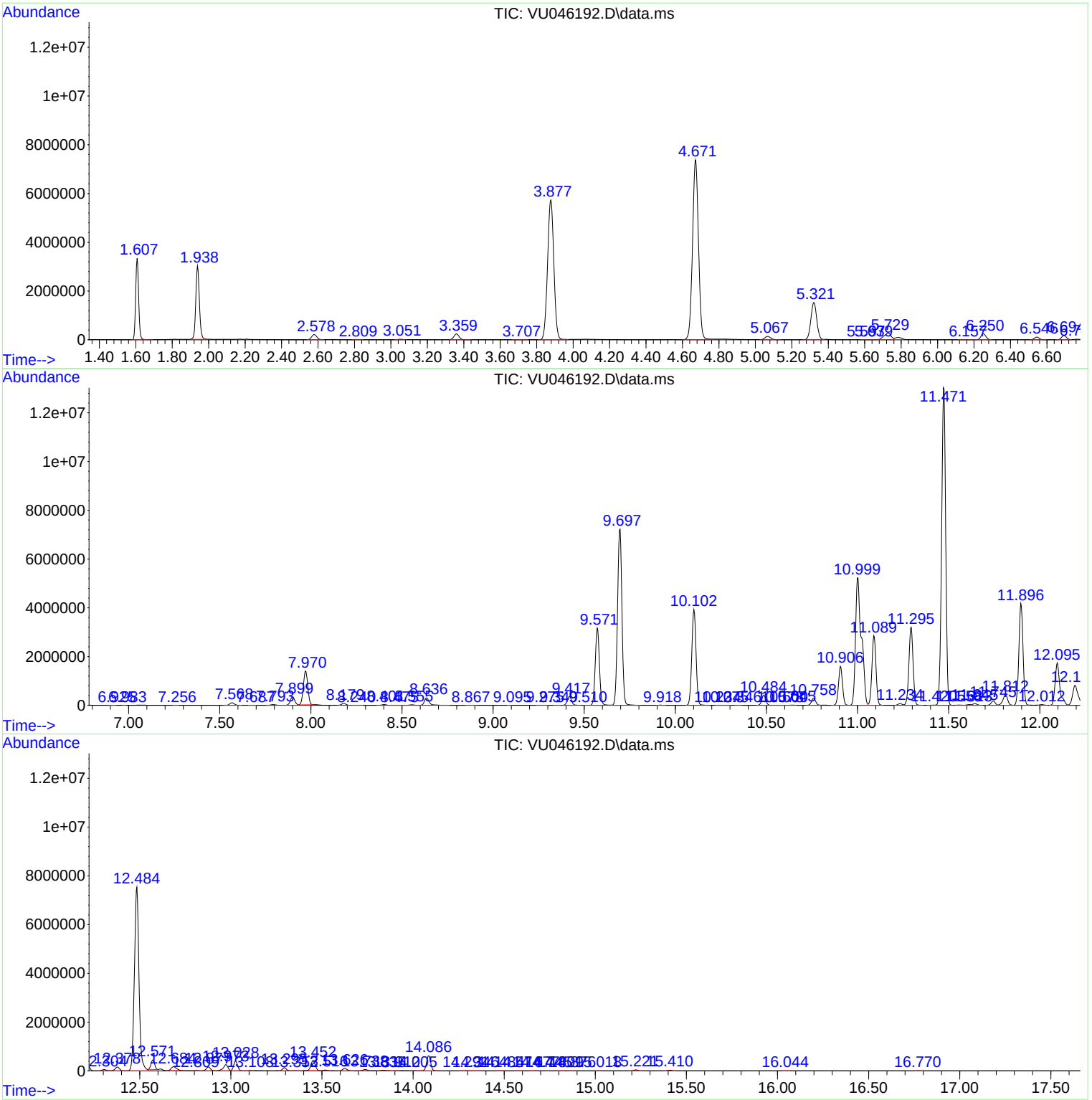
Sum of corrected areas: 148233084

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120821\
Data File : VU046192.D
Acq On : 08 Dec 2021 21:50
Operator : SY/MD
Sample : M4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5N1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120821\
Data File : VU046192.D
Acq On : 08 Dec 2021 21:50
Operator : SY/MD
Sample : M4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5N1

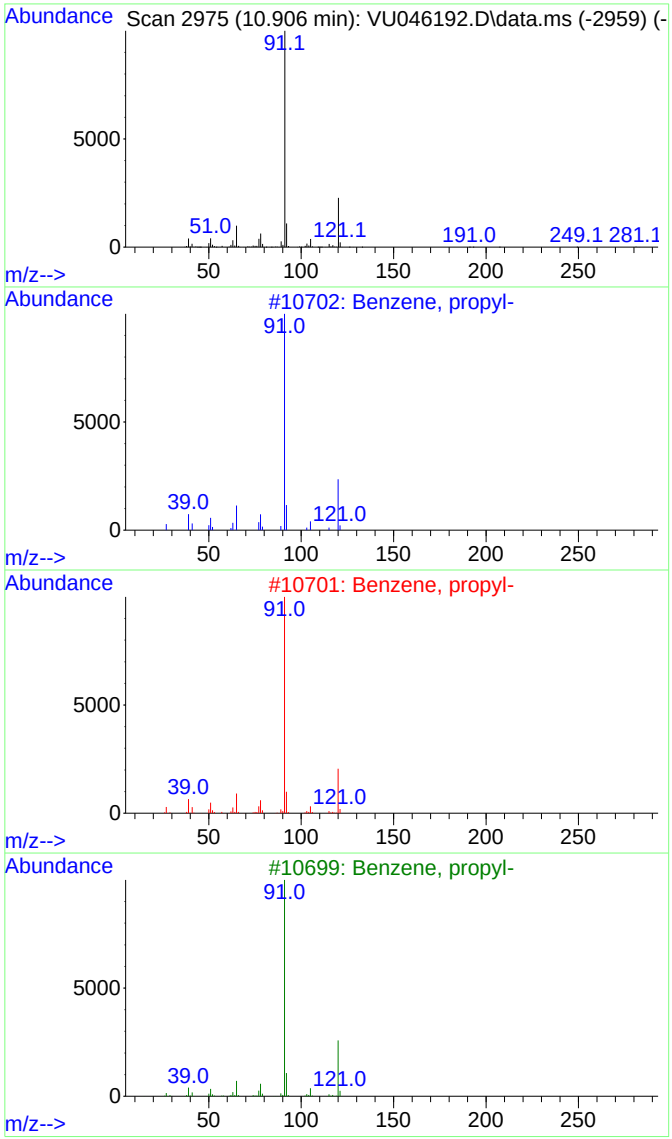
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
Quant Title : VOC Analysis

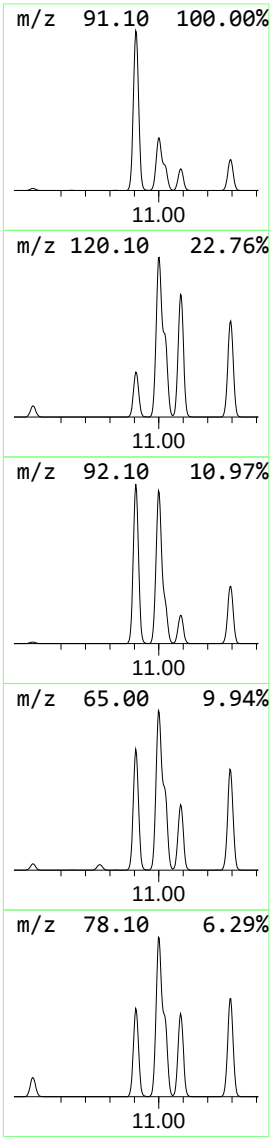
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Benzene, propyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.906	159.11 ug/L	2444720	1,4-Dichlorobenzene-d4	11.812

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, propyl-	120	C9H12	000103-65-1	94
2	Benzene, propyl-	120	C9H12	000103-65-1	90
3	Benzene, propyl-	120	C9H12	000103-65-1	90
4	Benzene, propyl-	120	C9H12	000103-65-1	86
5	N-Isobutyl(phenyl)methanesulfona...	227	C11H17NO2S	144615-94-1	72





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120821\
Data File : VU046192.D
Acq On : 08 Dec 2021 21:50
Operator : SY/MD
Sample : M4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5N1

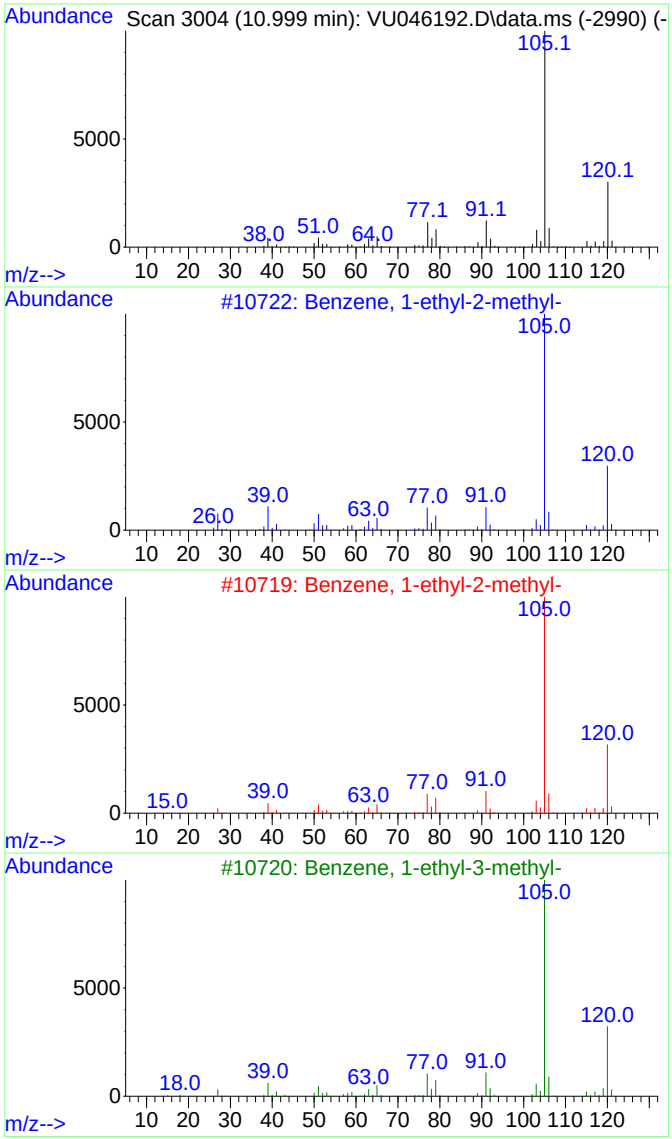
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
Quant Title : VOC Analysis

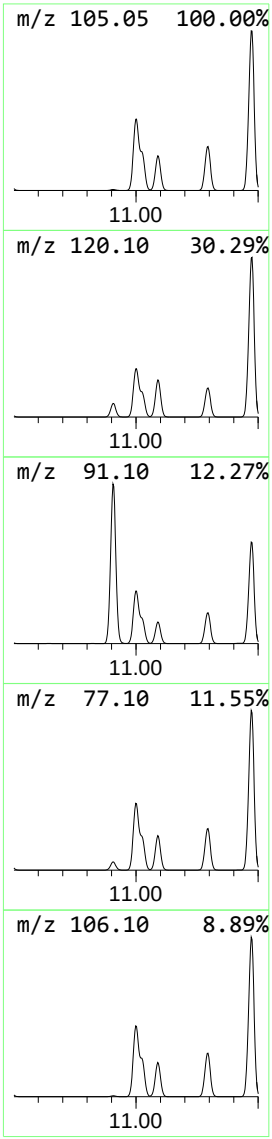
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.999	759.63 ug/L	11671500	1,4-Dichlorobenzene-d4	11.812

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-2-methyl-	120	C9H12	000611-14-3	95
3	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
5	Mesitylene	120	C9H12	000108-67-8	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120821\
Data File : VU046192.D
Acq On : 08 Dec 2021 21:50
Operator : SY/MD
Sample : M4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5N1

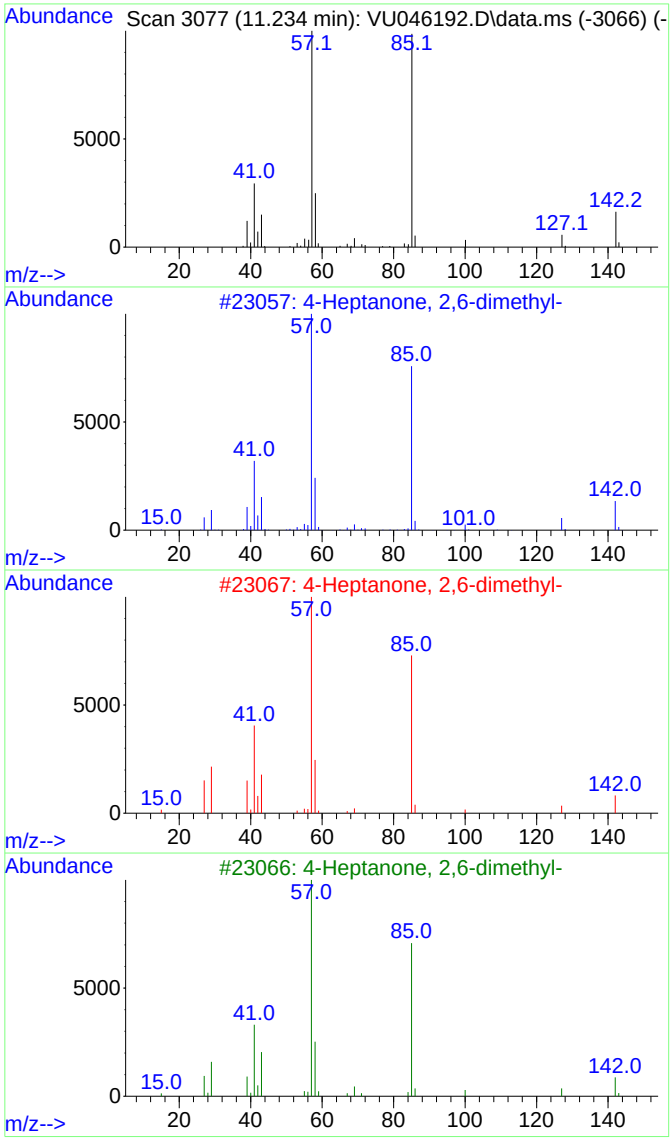
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
Quant Title : VOC Analysis

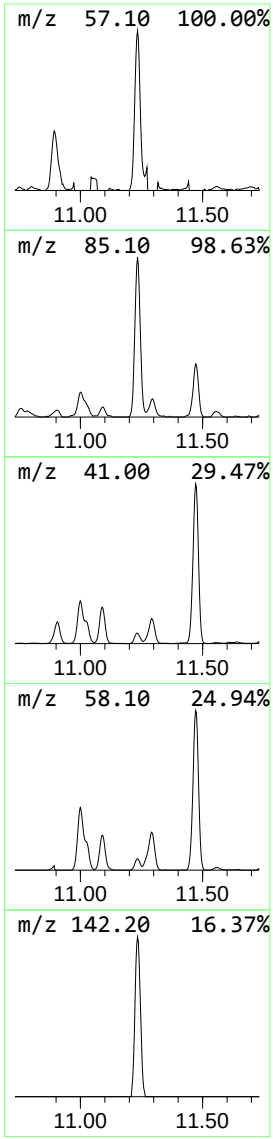
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 4-Heptanone, 2,6-dimethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.233	7.28 ug/L	111821	1,4-Dichlorobenzene-d4	11.812

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Heptanone, 2,6-dimethyl-	142	C9H18O	000108-83-8	94
2	4-Heptanone, 2,6-dimethyl-	142	C9H18O	000108-83-8	91
3	4-Heptanone, 2,6-dimethyl-	142	C9H18O	000108-83-8	91
4	4-Heptanone, 2,6-dimethyl-	142	C9H18O	000108-83-8	90
5	2,4-Hexanedione, 5,5-dimethyl-	142	C8H14O2	007307-04-2	59





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120821\
Data File : VU046192.D
Acq On : 08 Dec 2021 21:50
Operator : SY/MD
Sample : M4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5N1

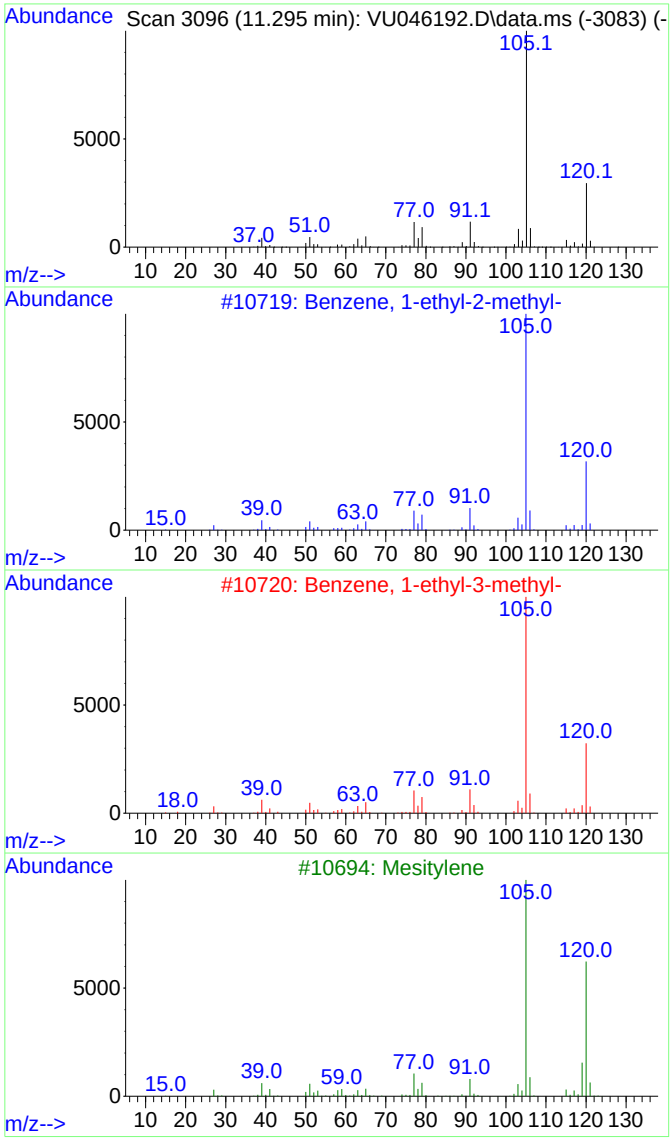
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
Quant Title : VOC Analysis

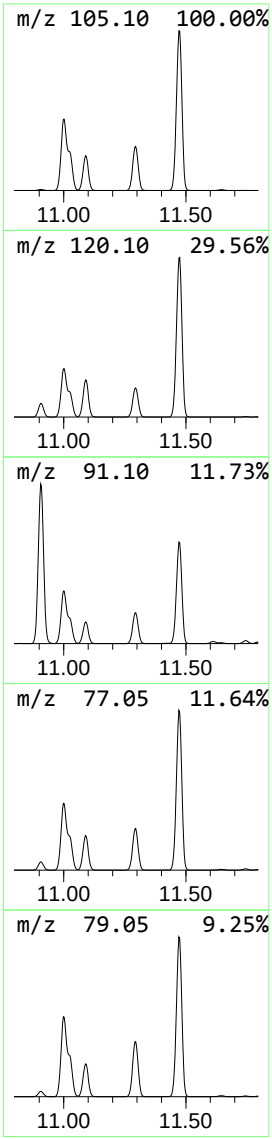
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzene, 1-ethyl-3-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.295	332.49 ug/L	5108640	1,4-Dichlorobenzene-d4	11.812

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
3	Mesitylene	120	C9H12	000108-67-8	91
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120821\
Data File : VU046192.D
Acq On : 08 Dec 2021 21:50
Operator : SY/MD
Sample : M4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5N1

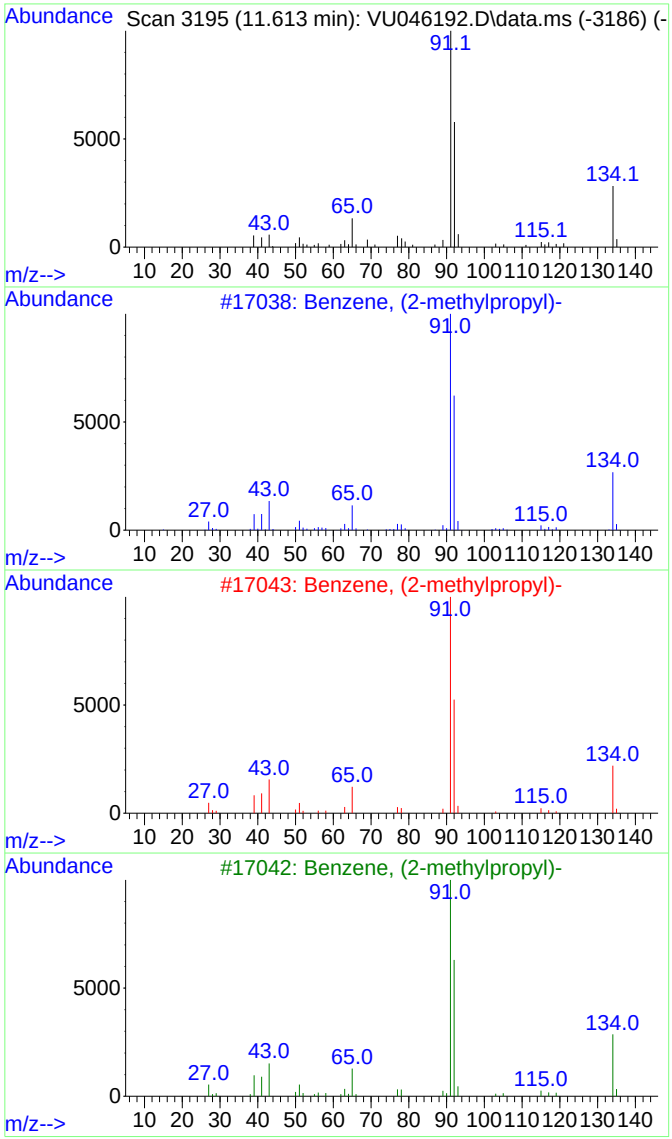
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
Quant Title : VOC Analysis

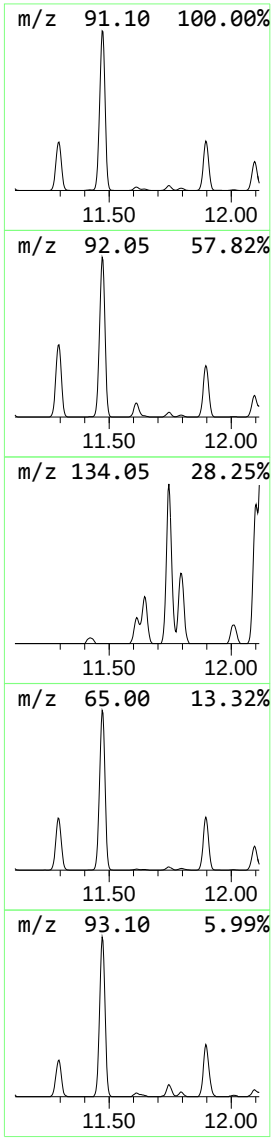
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzene, (2-methylpropyl)- Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.613	2.90 ug/L	44487	1,4-Dichlorobenzene-d4	11.812

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	90
2	Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	90
3	Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	90
4	Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	90
5	Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	90





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120821\
Data File : VU046192.D
Acq On : 08 Dec 2021 21:50
Operator : SY/MD
Sample : M4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5N1

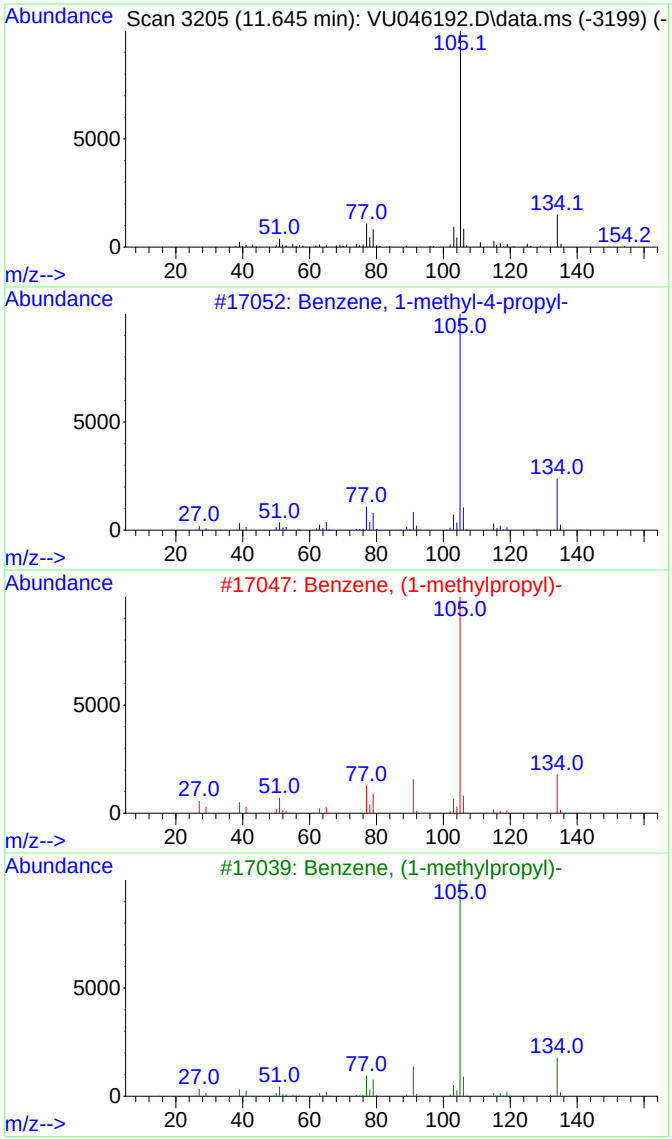
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
Quant Title : VOC Analysis

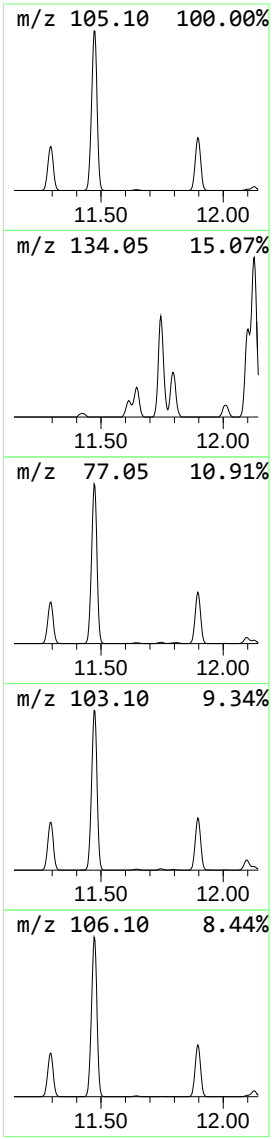
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzene, 1-methyl-4-propyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.645	6.65 ug/L	102119	1,4-Dichlorobenzene-d4	11.812

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	87
3	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	86
4	Oxirane, 2-methyl-2-phenyl-	134	C9H10O	002085-88-3	80
5	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	72





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120821\
Data File : VU046192.D
Acq On : 08 Dec 2021 21:50
Operator : SY/MD
Sample : M4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5N1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
Quant Title : VOC Analysis

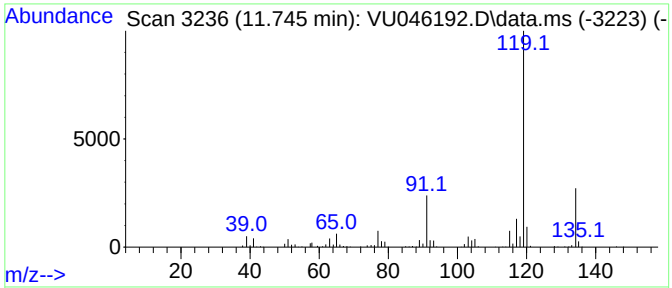
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

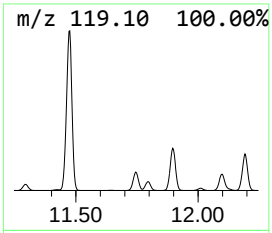
****

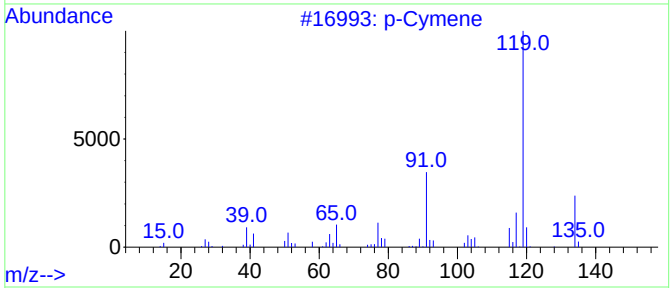
Peak Number 8 p-Cymene Concentration Rank 14

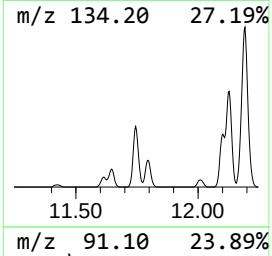
R.T.	EstConc	Area	Relative to ISTD	R.T.
11.745	18.17 ug/L	279133	1,4-Dichlorobenzene-d4	11.812

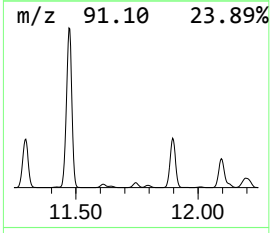
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Cymene	134	C10H14	000099-87-6	95
2	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94
3	1,3,5-Cycloheptatriene, 3,7,7-tr...	134	C10H14	003479-89-8	94
4	p-Cymene	134	C10H14	000099-87-6	93
5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	91

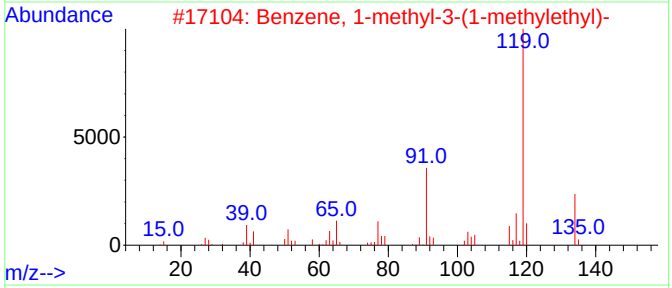


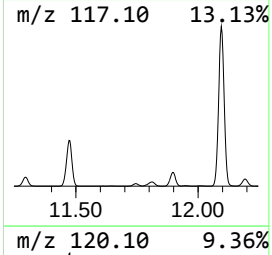


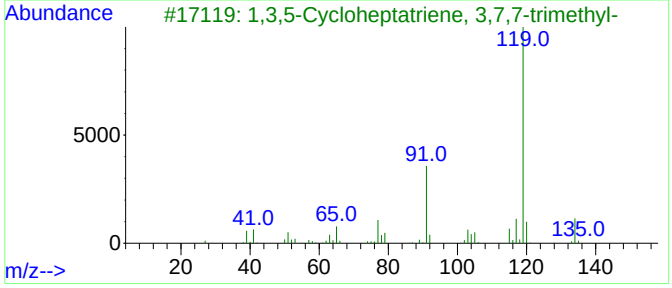


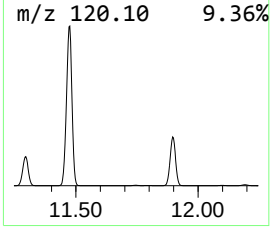












Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120821\
Data File : VU046192.D
Acq On : 08 Dec 2021 21:50
Operator : SY/MD
Sample : M4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5N1

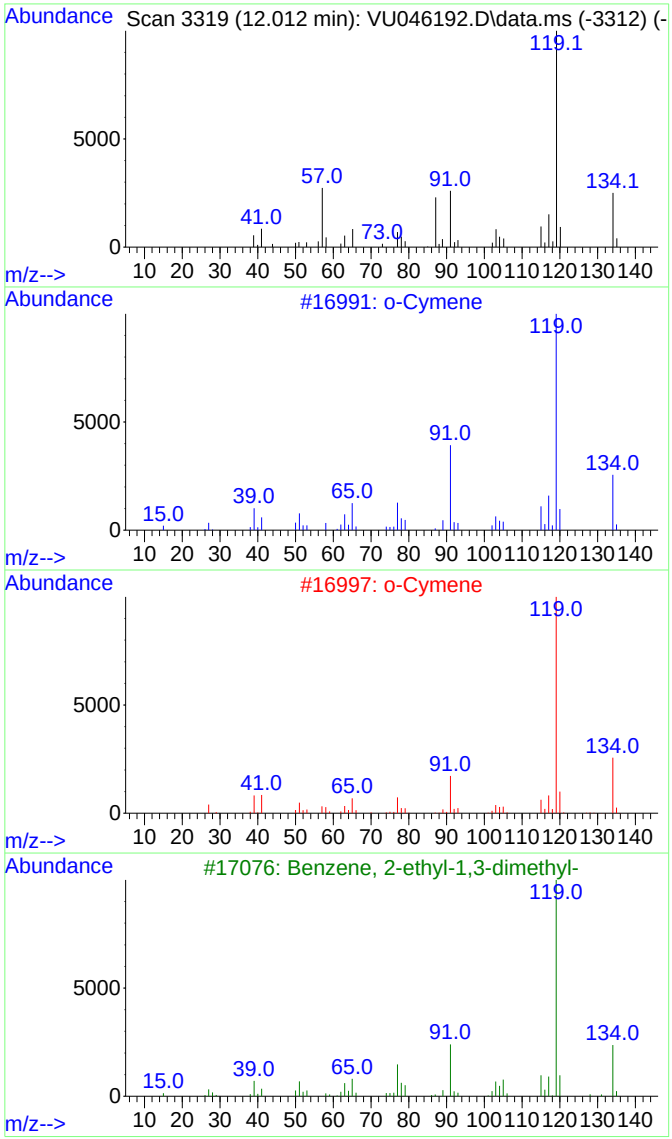
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
Quant Title : VOC Analysis

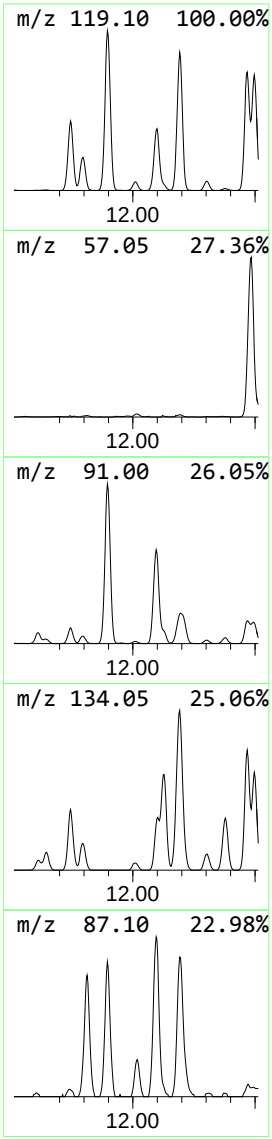
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 o-Cymene Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.012	2.64 ug/L	40507	1,4-Dichlorobenzene-d4	11.812

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	91
2	o-Cymene	134	C10H14	000527-84-4	87
3	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	87
4	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	87
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	87





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120821\
Data File : VU046192.D
Acq On : 08 Dec 2021 21:50
Operator : SY/MD
Sample : M4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5N1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
Quant Title : VOC Analysis

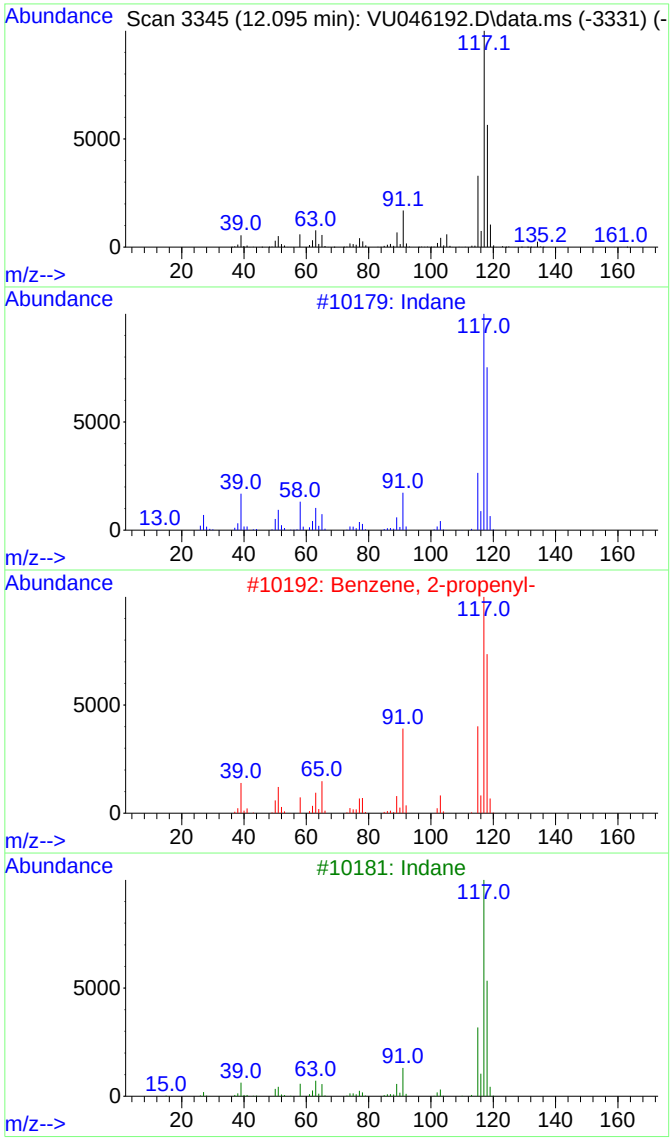
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

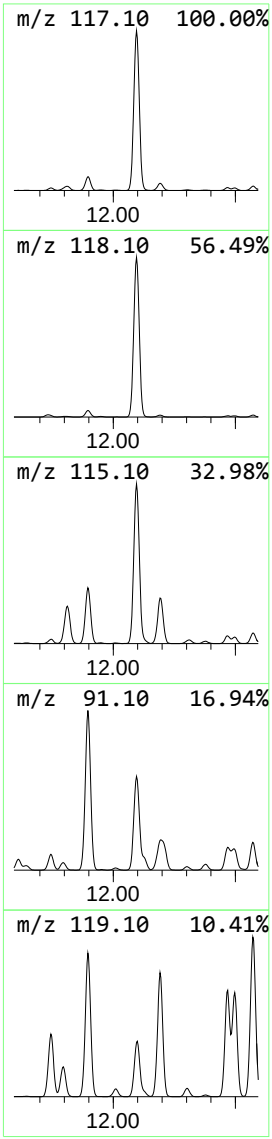
****

Peak Number 11 Indane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.095	202.85 ug/L	3116790	1,4-Dichlorobenzene-d4	11.812

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	91
2	Benzene, 2-propenyl-	118	C9H10	000300-57-2	83
3	Indane	118	C9H10	000496-11-7	81
4	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	81
5	Benzene, cyclopropyl-	118	C9H10	000873-49-4	81





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120821\
Data File : VU046192.D
Acq On : 08 Dec 2021 21:50
Operator : SY/MD
Sample : M4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5N1

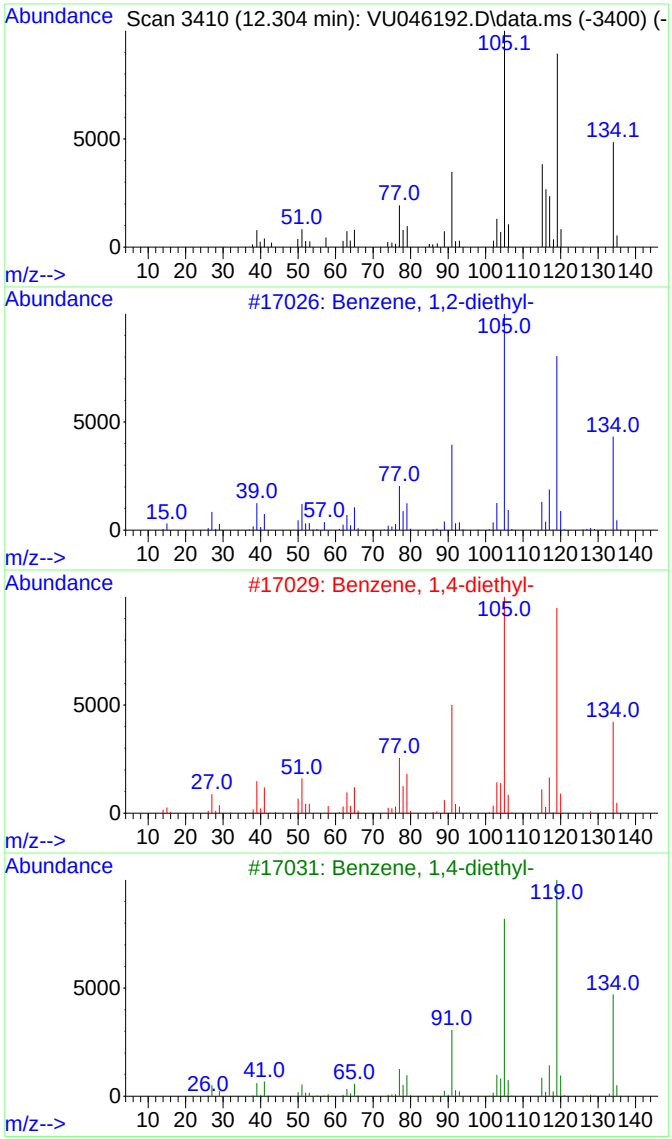
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
Quant Title : VOC Analysis

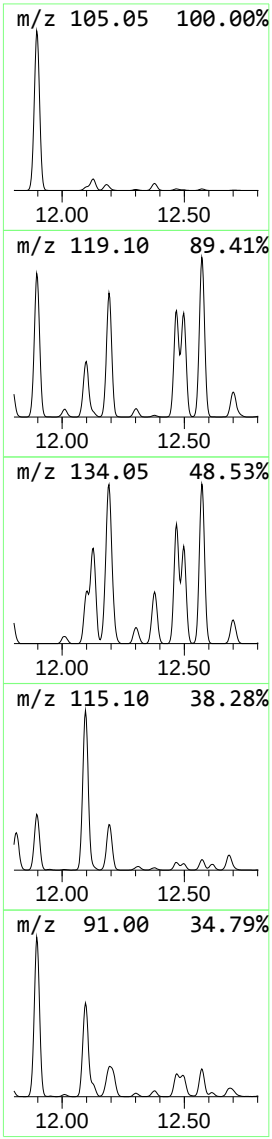
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 Benzene, 1,2-diethyl- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.304	5.77 ug/L	88615	1,4-Dichlorobenzene-d4	11.812

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	93
2	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	76
3	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	70
4	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	70
5	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	70





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120821\
Data File : VU046192.D
Acq On : 08 Dec 2021 21:50
Operator : SY/MD
Sample : M4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5N1

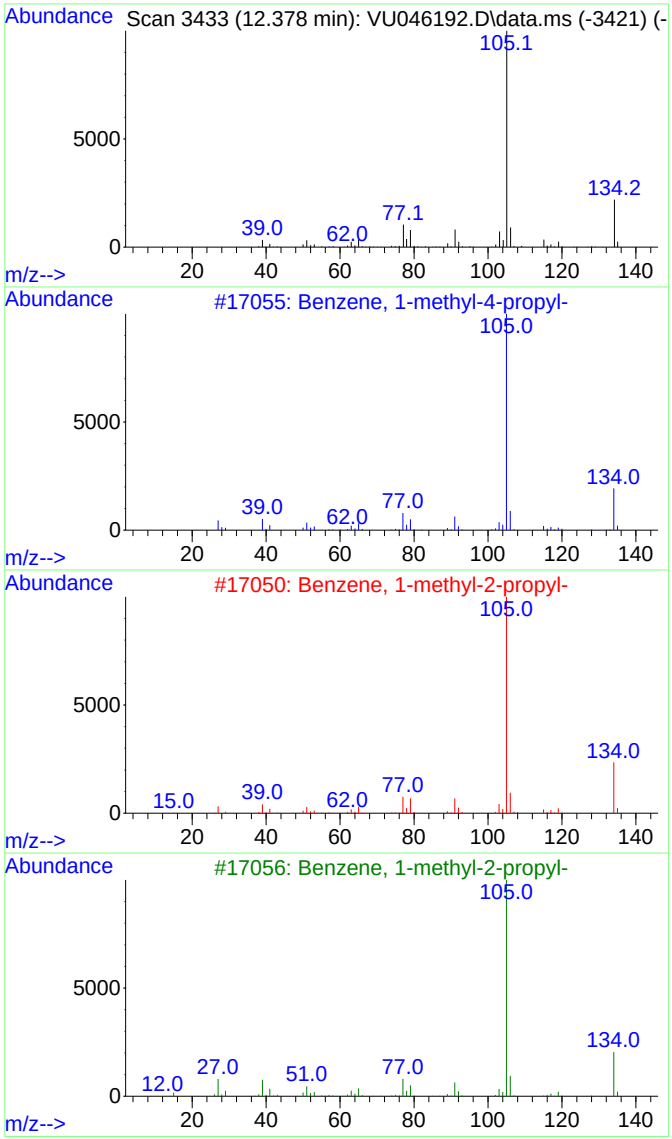
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
Quant Title : VOC Analysis

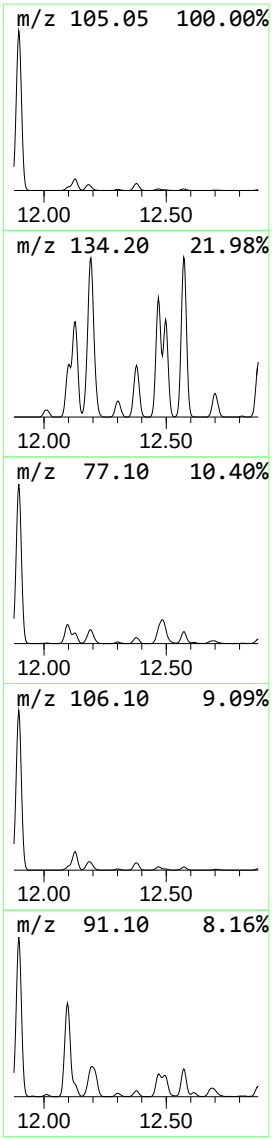
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Benzene, 1-methyl-2-propyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.378	16.80 ug/L	258193	1,4-Dichlorobenzene-d4	11.812

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
2	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	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





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120821\
Data File : VU046192.D
Acq On : 08 Dec 2021 21:50
Operator : SY/MD
Sample : M4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5N1

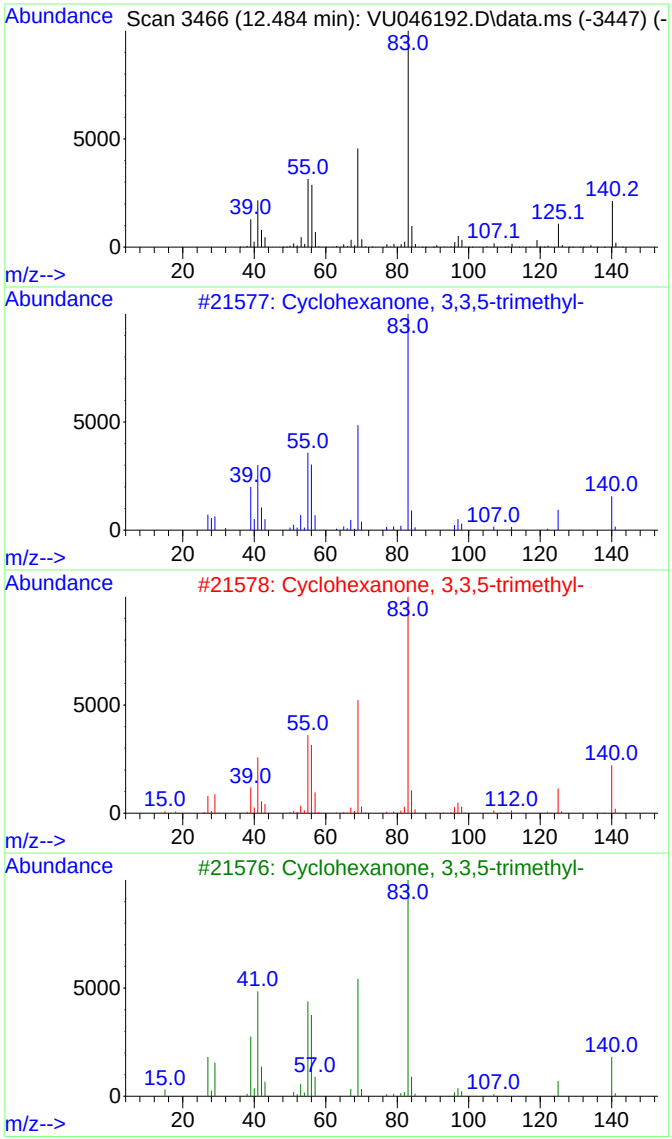
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
Quant Title : VOC Analysis

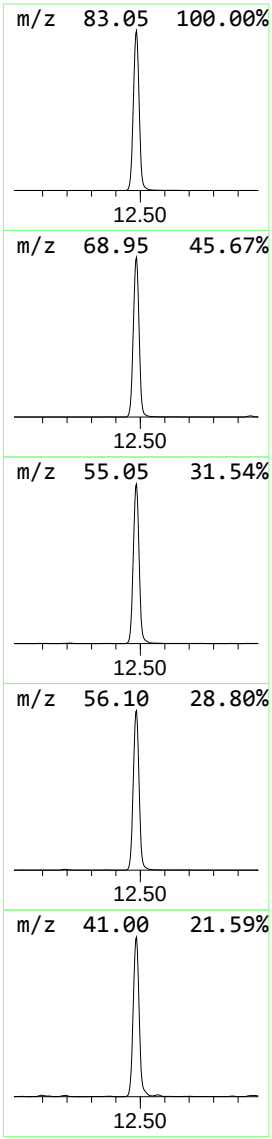
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 Cyclohexanone, 3,3,5-trimet... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.484	855.18 ug/L	13139700	1,4-Dichlorobenzene-d4	11.812

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	96
2	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	96
3	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	91
4	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	91
5	1H-1,2,4-Triazole, 3-(2-methylpr...	125	C6H11N3	040515-29-5	50





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120821\
Data File : VU046192.D
Acq On : 08 Dec 2021 21:50
Operator : SY/MD
Sample : M4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5N1

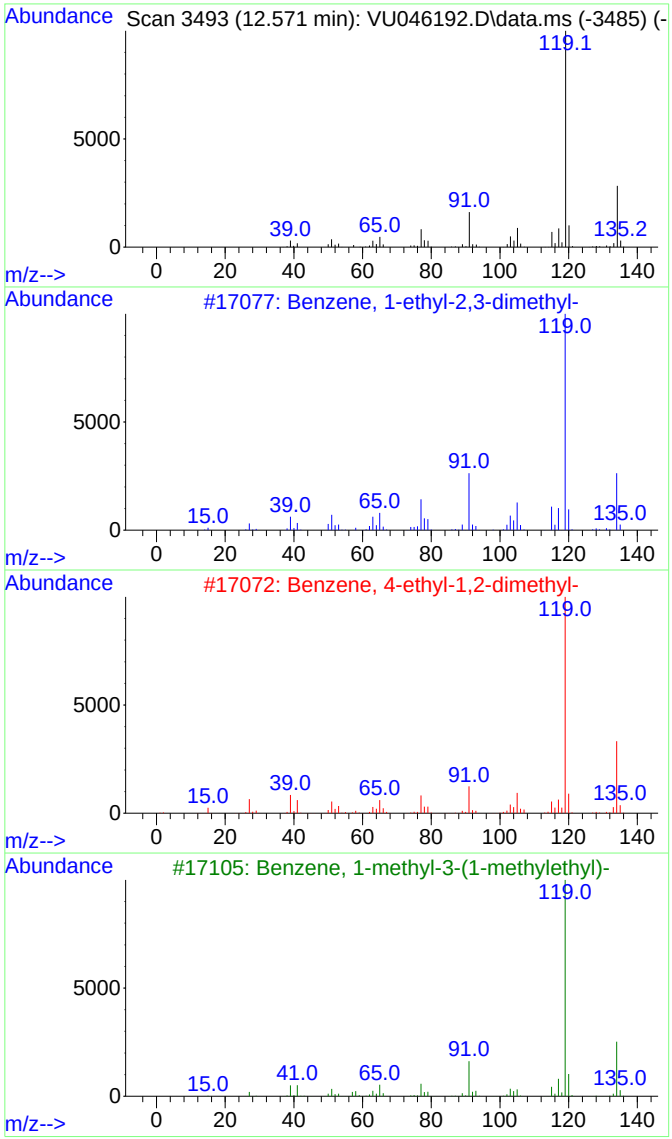
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
Quant Title : VOC Analysis

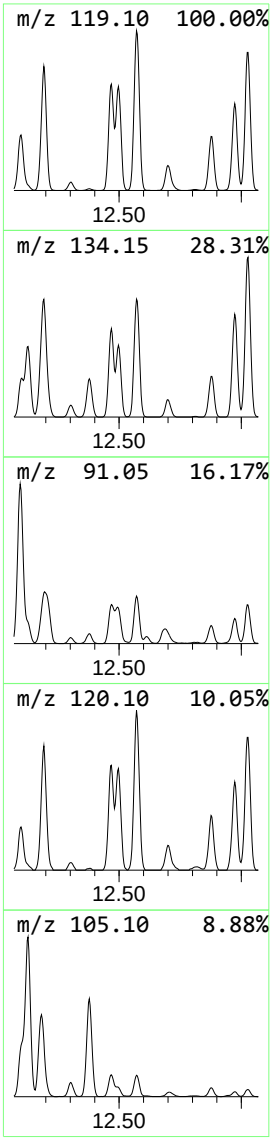
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.571	36.66 ug/L	563258	1,4-Dichlorobenzene-d4	11.812

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
2	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
3	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
4	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
5	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120821\
Data File : VU046192.D
Acq On : 08 Dec 2021 21:50
Operator : SY/MD
Sample : M4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5N1

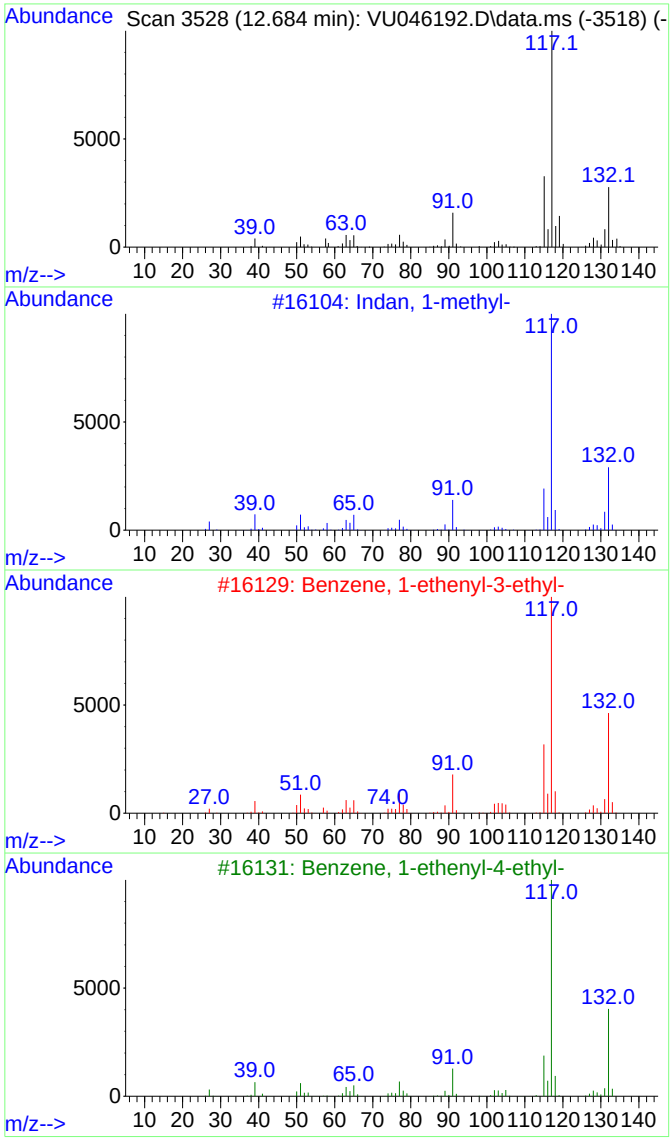
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
Quant Title : VOC Analysis

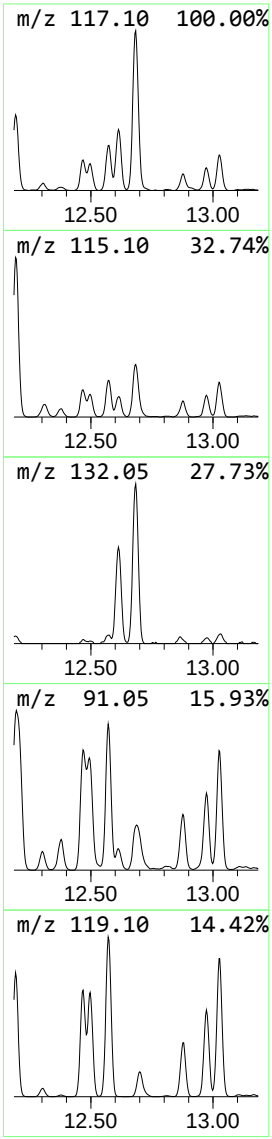
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 Indan, 1-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.684	22.31 ug/L	342771	1,4-Dichlorobenzene-d4	11.812

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indan, 1-methyl-	132	C10H12	000767-58-8	94
2	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
3	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	81
4	Indan, 1-methyl-	132	C10H12	000767-58-8	76
5	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	74





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120821\
Data File : VU046192.D
Acq On : 08 Dec 2021 21:50
Operator : SY/MD
Sample : M4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5N1

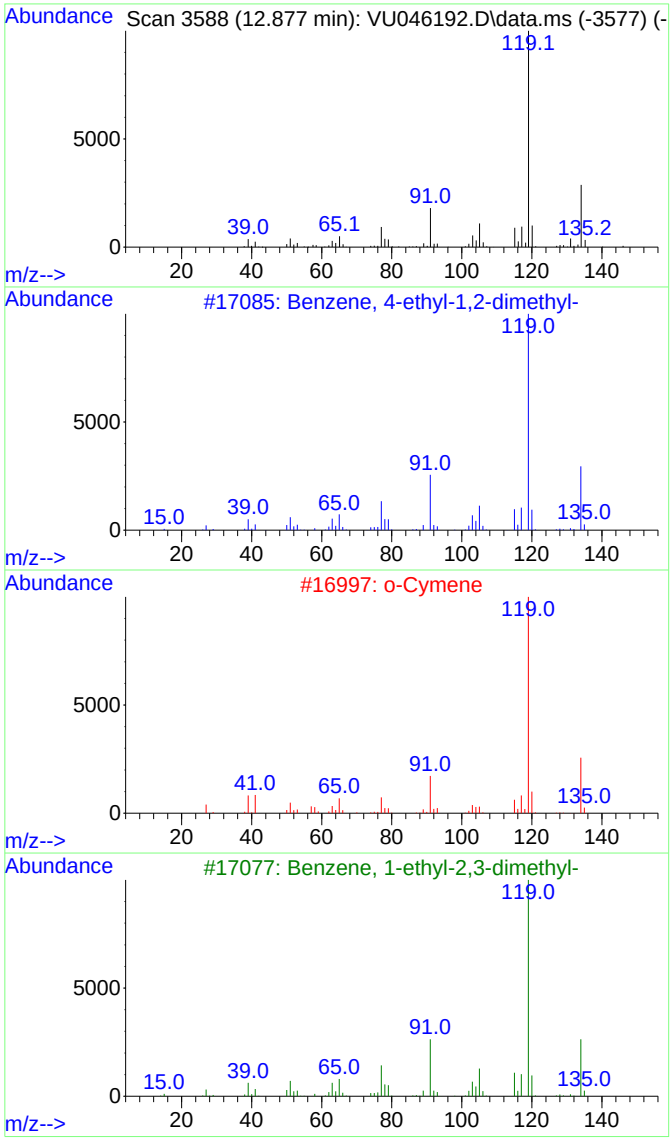
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
Quant Title : VOC Analysis

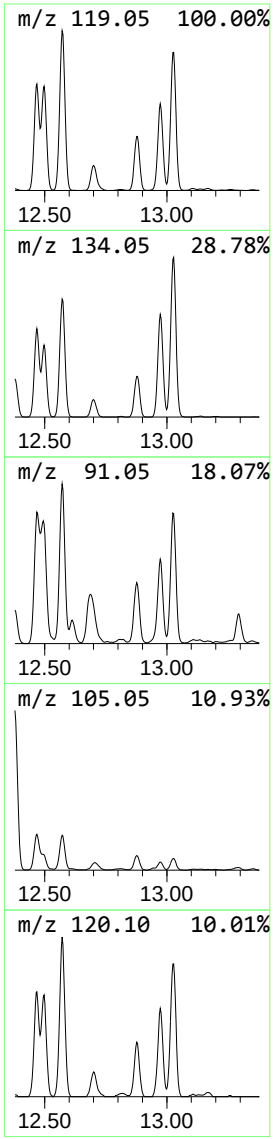
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.877	15.59 ug/L	239482	1,4-Dichlorobenzene-d4	11.812

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2	o-Cymene	134	C10H14	000527-84-4	95
3	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
4	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120821\
Data File : VU046192.D
Acq On : 08 Dec 2021 21:50
Operator : SY/MD
Sample : M4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5N1

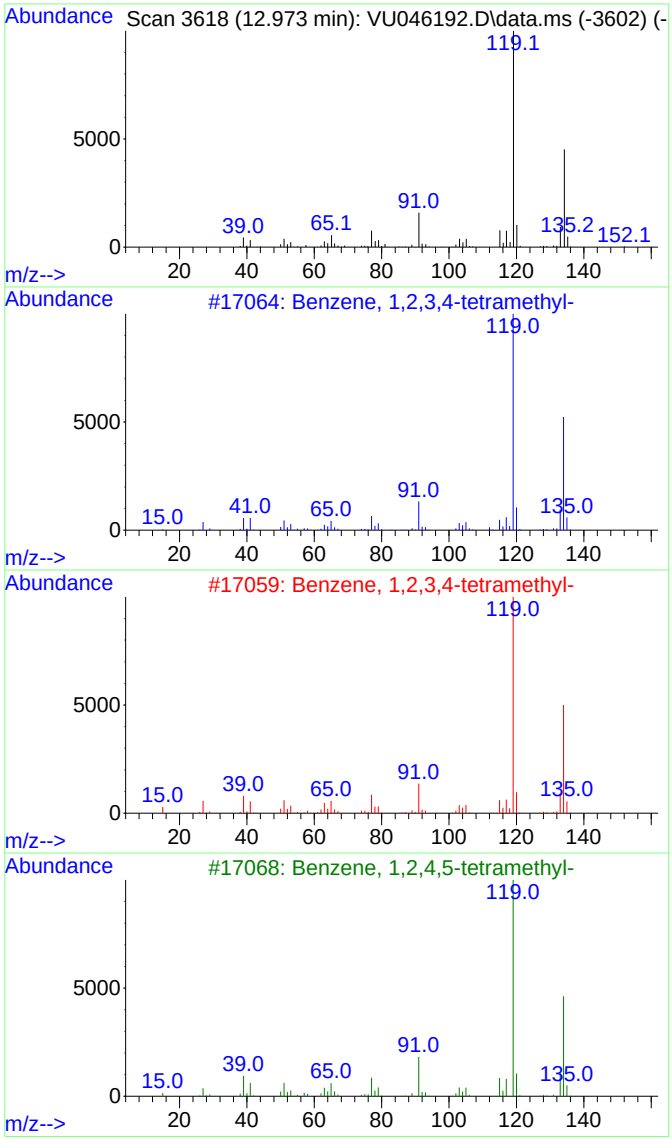
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
Quant Title : VOC Analysis

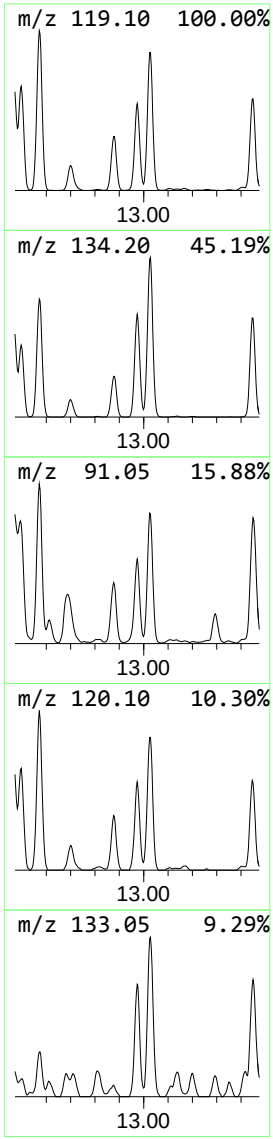
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 18 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.973	28.50 ug/L	437849	1,4-Dichlorobenzene-d4	11.812

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
4	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
5	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120821\
Data File : VU046192.D
Acq On : 08 Dec 2021 21:50
Operator : SY/MD
Sample : M4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5N1

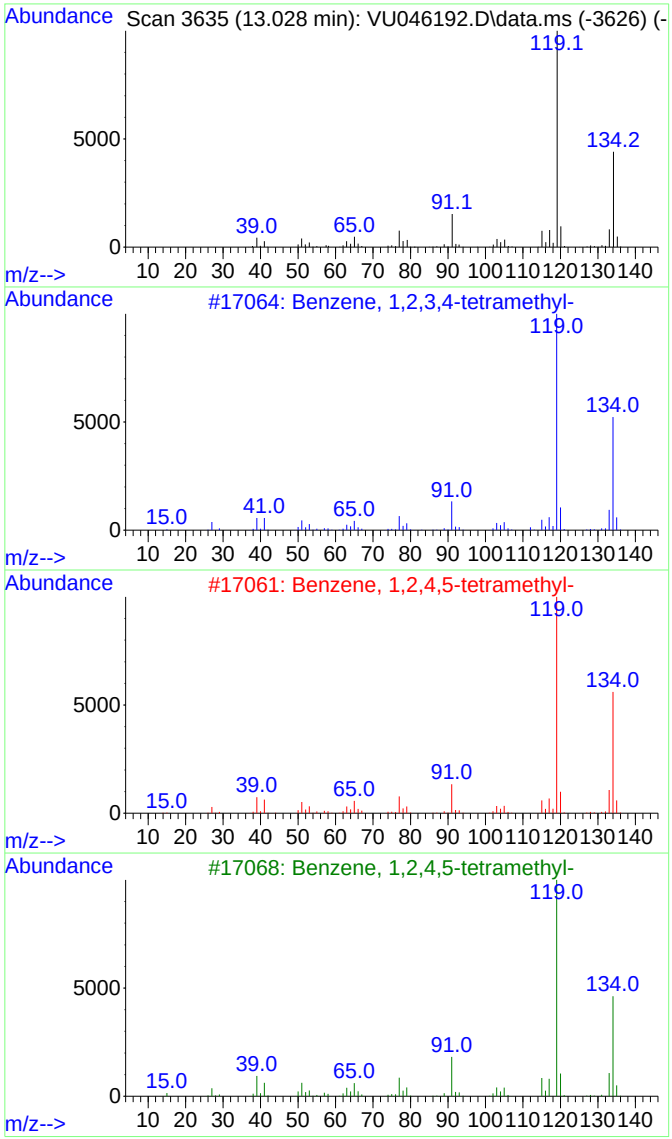
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
Quant Title : VOC Analysis

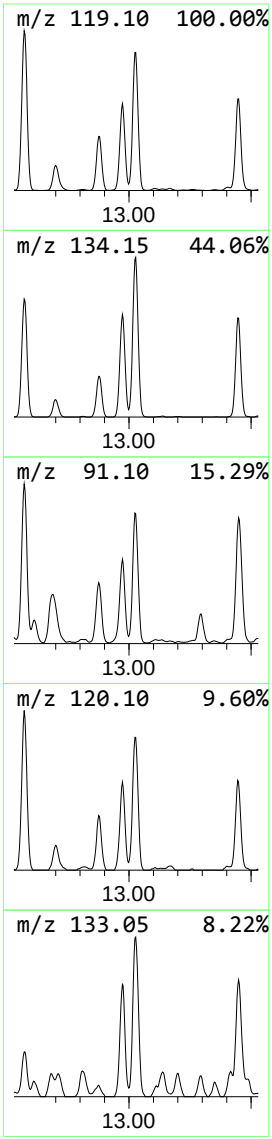
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.028	39.36 ug/L	604787	1,4-Dichlorobenzene-d4	11.812

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
4	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
5	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120821\
Data File : VU046192.D
Acq On : 08 Dec 2021 21:50
Operator : SY/MD
Sample : M4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5N1

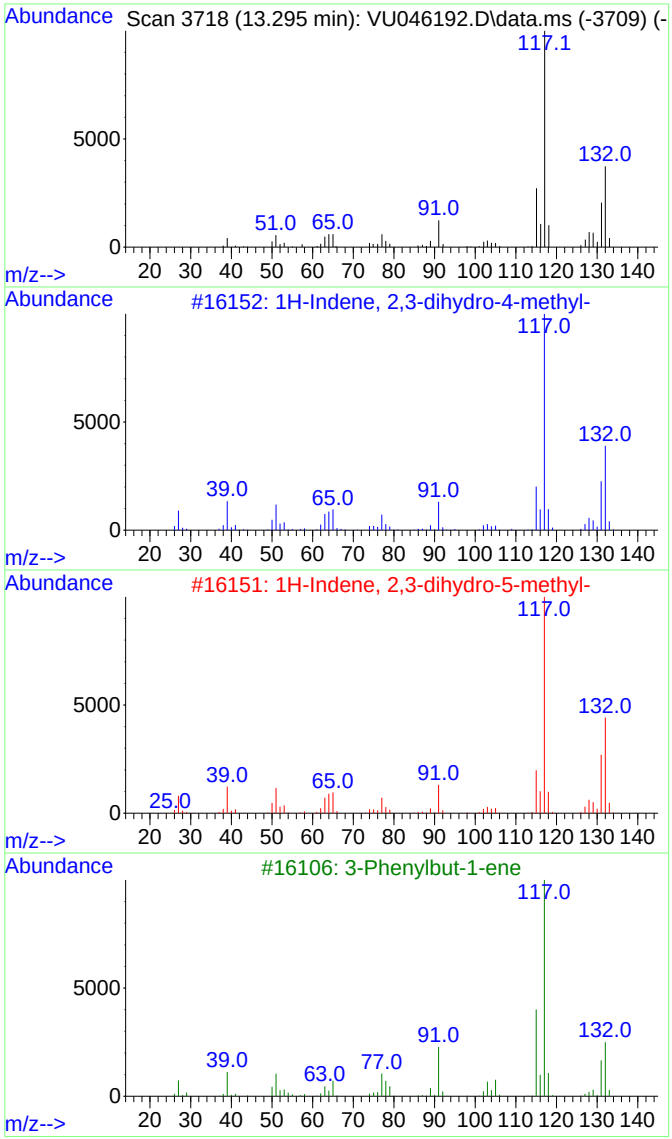
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
Quant Title : VOC Analysis

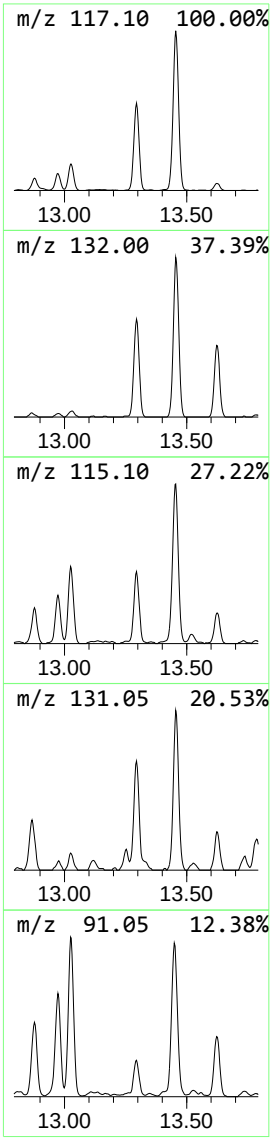
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.294	11.67 ug/L	179277	1,4-Dichlorobenzene-d4	11.812

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	95
2	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	91
3	3-Phenylbut-1-ene	132	C10H12	000934-10-1	91
4	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	87
5	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	83





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120821\
Data File : VU046192.D
Acq On : 08 Dec 2021 21:50
Operator : SY/MD
Sample : M4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5N1

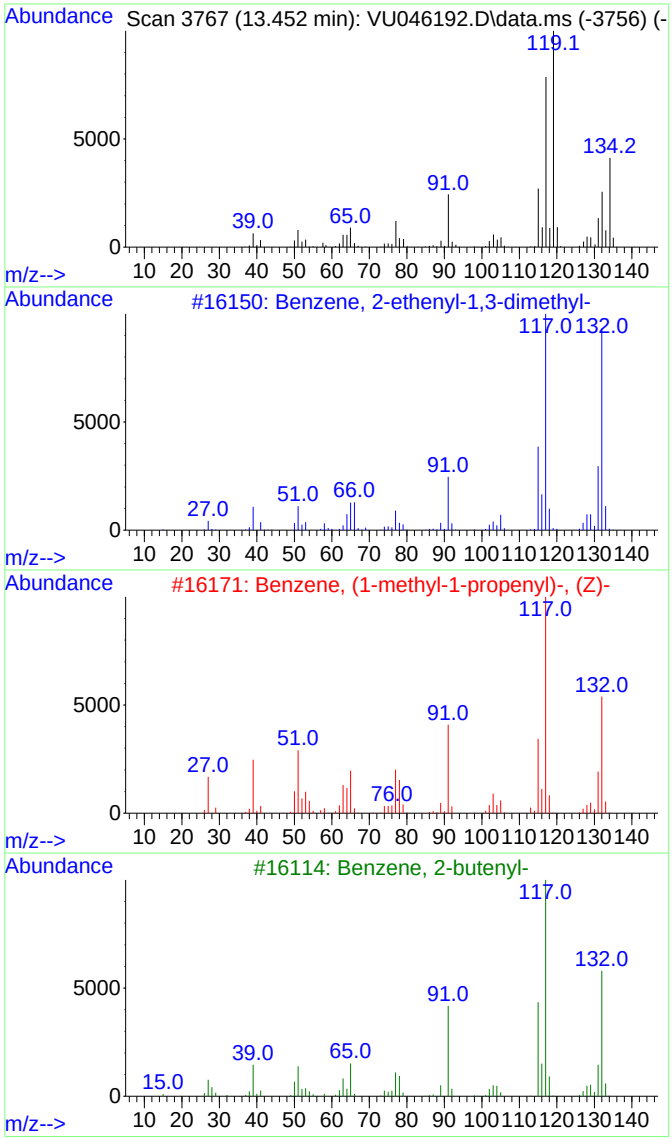
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
Quant Title : VOC Analysis

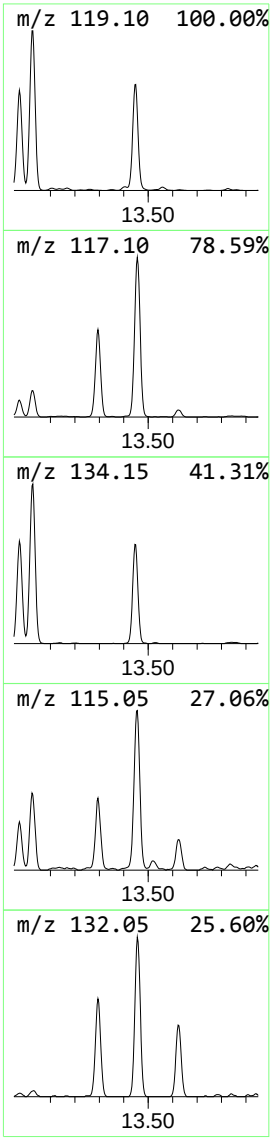
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 21 Benzene, 2-ethenyl-1,3-dime... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.452	43.93 ug/L	674921	1,4-Dichlorobenzene-d4	11.812

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	64
2	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	64
3	Benzene, 2-butenyl-	132	C10H12	001560-06-1	64
4	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	55
5	p-Cymene	134	C10H14	000099-87-6	55





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120821\
Data File : VU046192.D
Acq On : 08 Dec 2021 21:50
Operator : SY/MD
Sample : M4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5N1

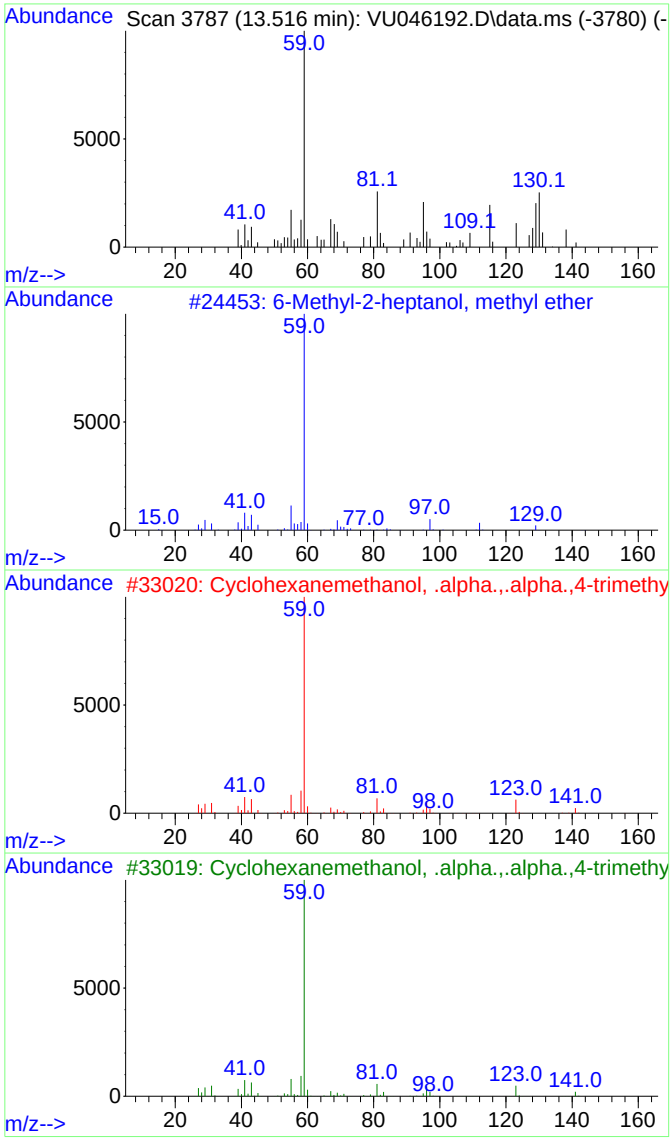
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
Quant Title : VOC Analysis

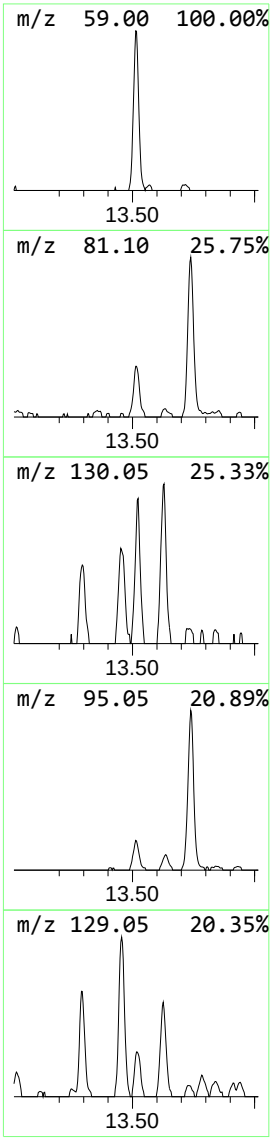
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 22 6-Methyl-2-heptanol, methyl... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.516	2.64 ug/L	40601	1,4-Dichlorobenzene-d4	11.812

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	6-Methyl-2-heptanol, methyl ether	144	C9H20O	1000365-52-0	27
2	Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	005114-00-1	25
3	Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	007322-63-6	25
4	7-Octen-2-ol, 2,6-dimethyl-	156	C10H20O	018479-58-8	23
5	Cyclopentanemethanol, .alpha.,.a...	128	C8H16O	001462-06-2	23





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120821\
Data File : VU046192.D
Acq On : 08 Dec 2021 21:50
Operator : SY/MD
Sample : M4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5N1

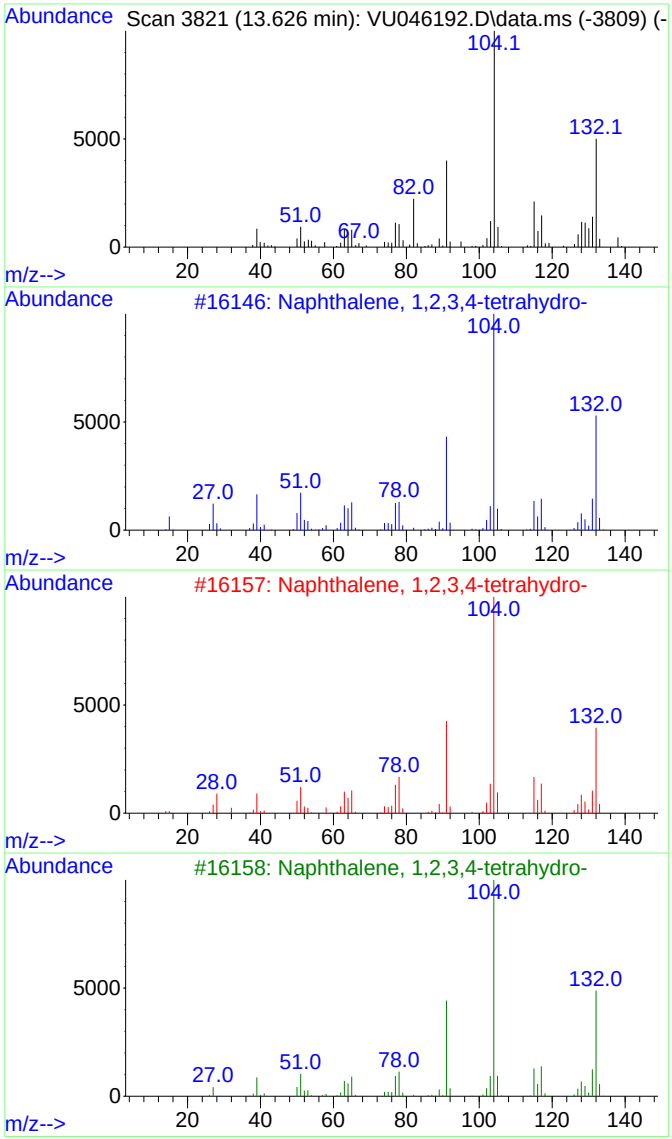
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
Quant Title : VOC Analysis

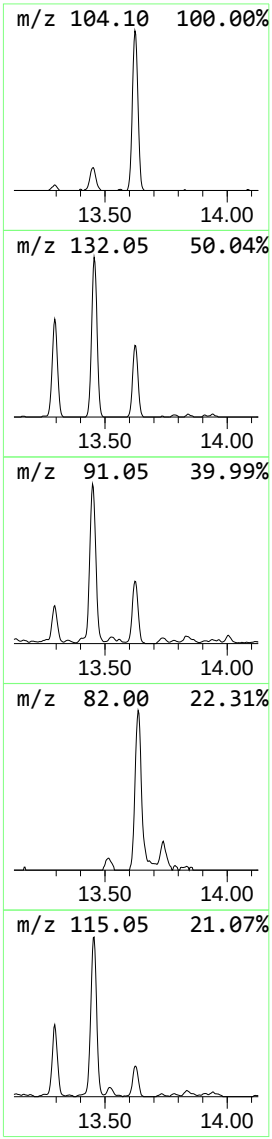
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.626	10.53 ug/L	161732	1,4-Dichlorobenzene-d4	11.812

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	94
2	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	93
3	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	92
4	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	70
5	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	68





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120821\
Data File : VU046192.D
Acq On : 08 Dec 2021 21:50
Operator : SY/MD
Sample : M4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5N1

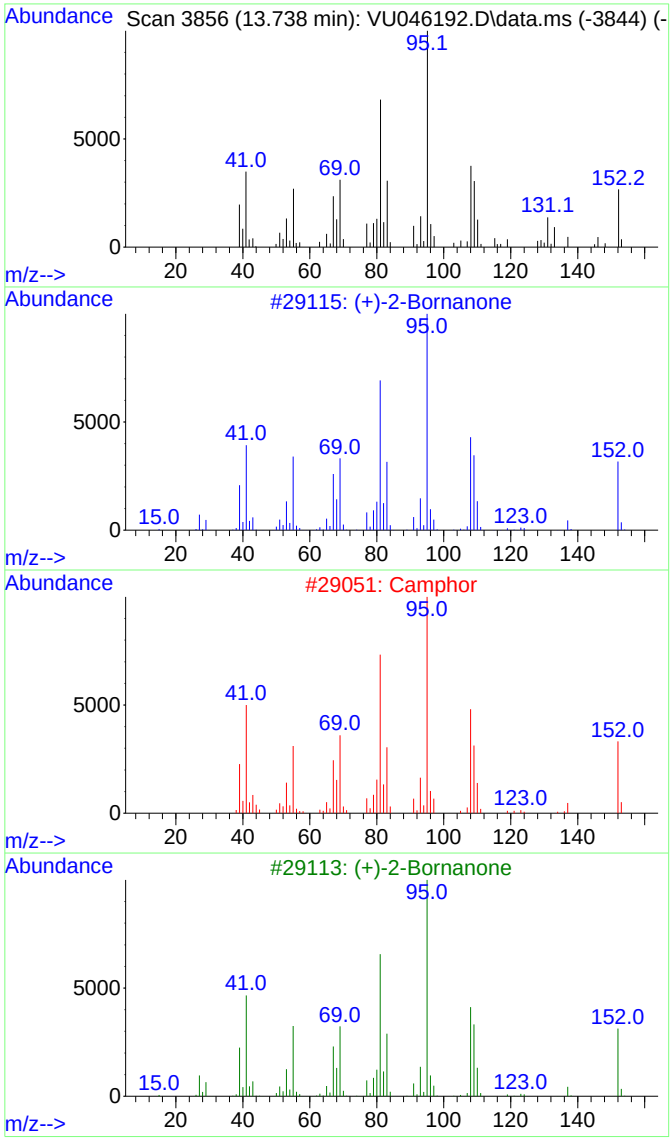
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
Quant Title : VOC Analysis

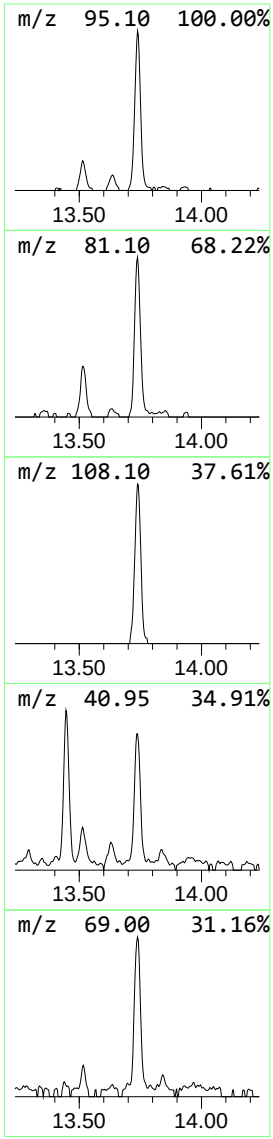
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 (+)-2-Bornanone Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.738	5.85 ug/L	89902	1,4-Dichlorobenzene-d4	11.812

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(+)-2-Bornanone	152	C10H16O	000464-49-3	98
2	Camphor	152	C10H16O	000076-22-2	98
3	(+)-2-Bornanone	152	C10H16O	000464-49-3	98
4	Camphor	152	C10H16O	000076-22-2	98
5	(+)-2-Bornanone	152	C10H16O	000464-49-3	97





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120821\
Data File : VU046192.D
Acq On : 08 Dec 2021 21:50
Operator : SY/MD
Sample : M4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5N1

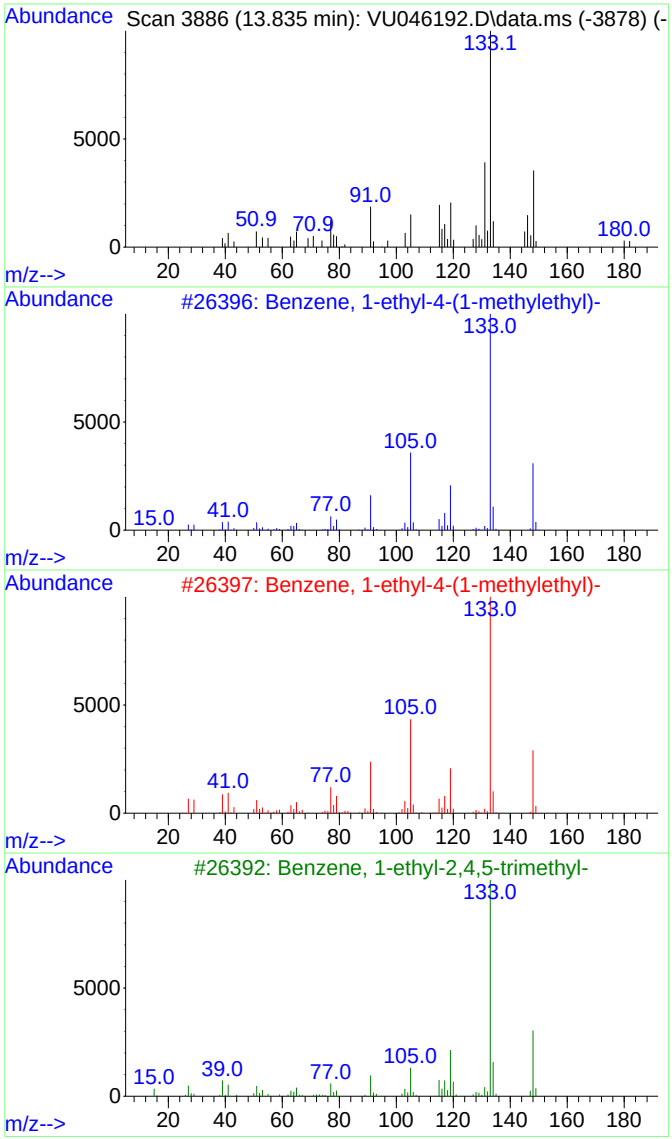
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
Quant Title : VOC Analysis

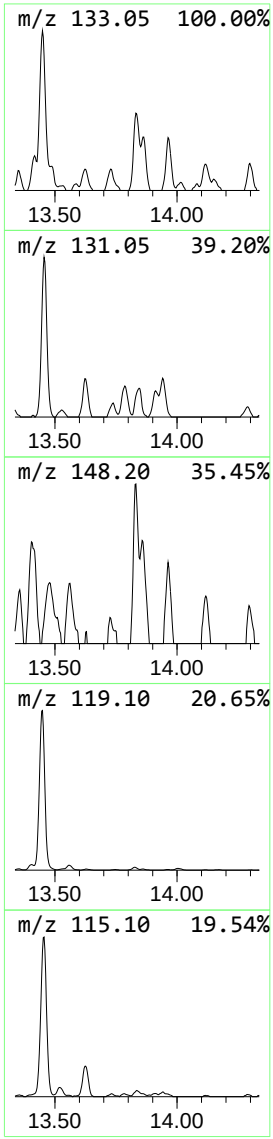
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 25 Benzene, 1-ethyl-4-(1-methy... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.835	2.86 ug/L	43961	1,4-Dichlorobenzene-d4	11.812

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	64
2	Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	58
3	Benzene, 1-ethyl-2,4,5-trimethyl-	148	C11H16	017851-27-3	52
4	Benzene, 1-ethyl-3-(1-methylethyl)-	148	C11H16	004920-99-4	50
5	Benzene, 1,3-dimethyl-5-(1-methy...	148	C11H16	004706-90-5	49





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120821\
Data File : VU046192.D
Acq On : 08 Dec 2021 21:50
Operator : SY/MD
Sample : M4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5N1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
Quant Title : VOC Analysis

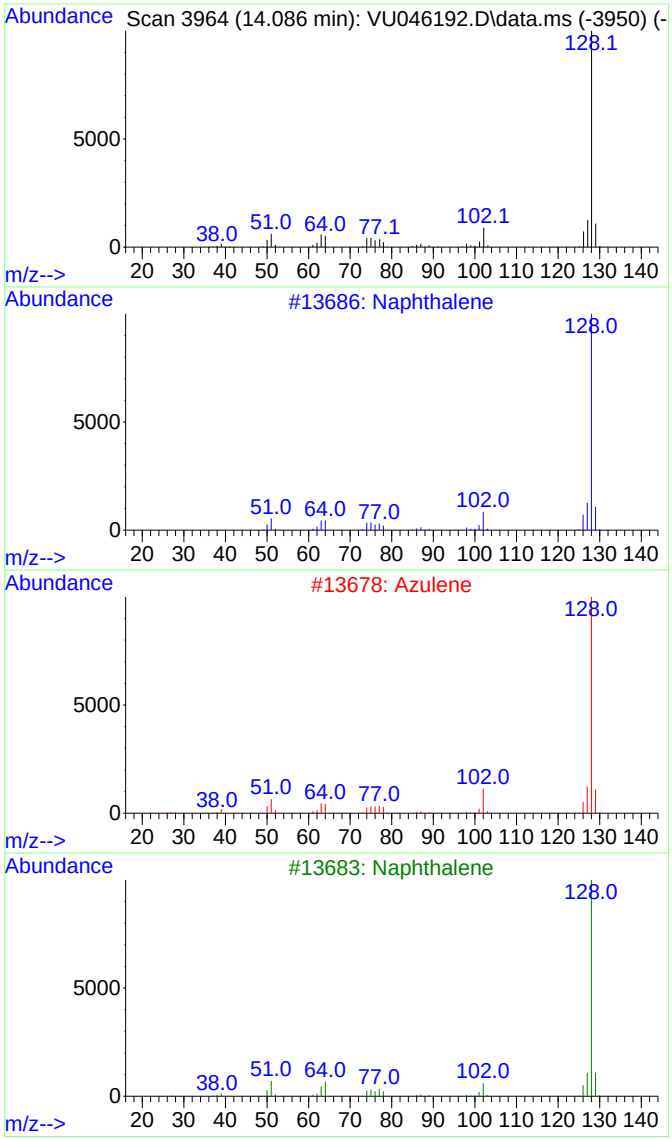
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

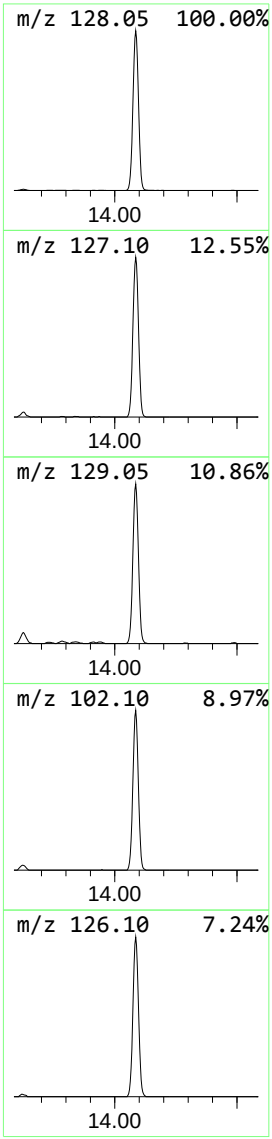
****

Peak Number 26 Azulene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.086	63.43 ug/L	974600	1,4-Dichlorobenzene-d4	11.812

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene	128	C10H8	000091-20-3	95
2	Azulene	128	C10H8	000275-51-4	94
3	Naphthalene	128	C10H8	000091-20-3	93
4	Azulene	128	C10H8	000275-51-4	91
5	Azulene	128	C10H8	000275-51-4	91





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120821\
Data File : VU046192.D
Acq On : 08 Dec 2021 21:50
Operator : SY/MD
Sample : M4983-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5N1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Benzene, propyl-	10.906	159.1	ug/L	2444720	3	11.812	768239	50.0
Benzene, 1-ethy...	10.999	759.6	ug/L	11671500	3	11.812	768239	50.0
4-Heptanone, 2,...	11.233	7.3	ug/L	111821	3	11.812	768239	50.0
Benzene, 1-ethy...	11.295	332.5	ug/L	5108640	3	11.812	768239	50.0
Benzene, (2-met...	11.613	2.9	ug/L	44487	3	11.812	768239	50.0
Benzene, 1-meth...	11.645	6.7	ug/L	102119	3	11.812	768239	50.0
p-Cymene	11.745	18.2	ug/L	279133	3	11.812	768239	50.0
o-Cymene	12.012	2.6	ug/L	40507	3	11.812	768239	50.0
Indane	12.095	202.8	ug/L	3116790	3	11.812	768239	50.0
Benzene, 1,2-di...	12.304	5.8	ug/L	88615	3	11.812	768239	50.0
Benzene, 1-meth...	12.378	16.8	ug/L	258193	3	11.812	768239	50.0
Cyclohexanone, ...	12.484	855.2	ug/L	13139700	3	11.812	768239	50.0
Benzene, 1-ethy...	12.571	36.7	ug/L	563258	3	11.812	768239	50.0
Indan, 1-methyl-	12.684	22.3	ug/L	342771	3	11.812	768239	50.0
Benzene, 4-ethy...	12.877	15.6	ug/L	239482	3	11.812	768239	50.0
Benzene, 1,2,3,...	12.973	28.5	ug/L	437849	3	11.812	768239	50.0
Benzene, 1,2,4,...	13.028	39.4	ug/L	604787	3	11.812	768239	50.0
1H-Indene, 2,3-...	13.294	11.7	ug/L	179277	3	11.812	768239	50.0
Benzene, 2-ethe...	13.452	43.9	ug/L	674921	3	11.812	768239	50.0
6-Methyl-2-hept...	13.516	2.6	ug/L	40601	3	11.812	768239	50.0
Naphthalene, 1,...	13.626	10.5	ug/L	161732	3	11.812	768239	50.0
(+)-2-Bornanone	13.738	5.8	ug/L	89902	3	11.812	768239	50.0
Benzene, 1-ethy...	13.835	2.9	ug/L	43961	3	11.812	768239	50.0
Azulene	14.086	63.4	ug/L	974600	3	11.812	768239	50.0