LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU100721\
Data File : VU045231.D
Acq On : 08 Oct 2021 00:39
Operator : SY/MD
Sample : M4138-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F3

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\SFAMULM092321WMA.M
Title : VOC Analysis

Signal : TIC: VU045231.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.360	3	6	17	rVB3	40457	37543	0.02%	0.010%
2	1.475	31	42	43	rBV2	407581	501913	0.31%	0.139%
3	1.597	75	80	92	rVB	179148	195459	0.12%	0.054%
4	1.842	150	156	160	rBV2	36471	55166	0.03%	0.015%
5	1.893	166	172	196	rVB	146618	239061	0.15%	0.066%
6	2.562	367	380	395	rBV	247425	426096	0.26%	0.118%
7	2.633	395	402	410	rVV2	12182	20139	0.01%	0.006%
8	2.684	411	418	431	rVB	13197	23057	0.01%	0.006%
9	2.781	433	448	463	rBV2	28350	70035	0.04%	0.019%
10	3.048	521	531	533	rBV7	6323	10517	0.01%	0.003%
11	3.199	569	578	593	rVB3	5440	13354	0.01%	0.004%
12	3.359	615	628	645	rVB2	18450	42267	0.03%	0.012%
13	3.697	721	733	752	rVB3	18099	43340	0.03%	0.012%
14	3.871	770	787	807	rBV2	122739	293250	0.18%	0.081%
15	4.533	978	993	1009	rBV2	26580	69246	0.04%	0.019%
16	4.629	1009	1023	1038	rVV	143096	345542	0.21%	0.096%
17	4.710	1038	1048	1069	rVB	83675	190903	0.12%	0.053%
18	4.813	1077	1080	1097	rVB3	2098	3249	0.00%	0.001%
19	4.990	1120	1135	1148	rBV	203530	467364	0.29%	0.130%
20	5.067	1148	1159	1195	rVB	227839	518339	0.32%	0.144%
21	5.392	1243	1260	1274	rVV3	31972	74784	0.05%	0.021%
22	5.459	1276	1281	1292	rVB6	2743	5353	0.00%	0.001%
23	5.591	1313	1322	1323	rBV5	3168	3701	0.00%	0.001%
24	5.639	1332	1337	1345	rVB5	2449	3250	0.00%	0.001%
25	5.729	1345	1365	1374	rVV2	473039	1278227	0.79%	0.355%
26	5.777	1375	1380	1394	rVV	186913	331951	0.21%	0.092%
27	5.851	1395	1403	1414	rVB4	9354	17982	0.01%	0.005%
28	5.925	1414	1426	1435	rBV5	8343	16887	0.01%	0.005%
29	5.990	1436	1446	1466	rVB6	15349	33880	0.02%	0.009%
30	6.137	1480	1492	1498	rBV6	2163	3990	0.00%	0.001%
31	6.253	1511	1528	1557	rBV	383426	726050	0.45%	0.202%
32	6.533	1604	1615	1621	rBV3	8605	17441	0.01%	0.005%
33	6.694	1651	1665	1677	rVV	304839	594326	0.37%	0.165%
34	6.764	1677	1687	1708	rVB	129807	269973	0.17%	0.075%
35	6.983	1742	1755	1767	rVB4	12245	23856	0.01%	0.007%
36	7.076	1767	1784	1788	rBV6	1191	2600	0.00%	0.001%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU100721\
Data File : VU045231.D
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Operator : SY/MD
Sample : M4138-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F3

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\SFAMULM092321WMA.M
Title : VOC Analysis

37	7.121	1791	1798	1799	rBV3	3091	3163	0.00%	0.001%
38	7.176	1808	1815	1827	rVB8	1985	4084	0.00%	0.001%
39	7.244	1827	1836	1844	rVB8	3101	5376	0.00%	0.001%
40	7.568	1923	1937	1954	rBV	168635	297600	0.18%	0.083%
41	7.800	1993	2009	2022	rBV	132084	239626	0.15%	0.067%
42	7.903	2022	2041	2050	rBV	607957	1044673	0.65%	0.290%
43	7.977	2050	2064	2098	rVB	16288989	27684404	17.21%	7.691%
44	8.182	2118	2128	2136	rBV	112664	179350	0.11%	0.050%
45	8.234	2136	2144	2179	rVB	108729	233088	0.14%	0.065%
46	8.375	2179	2188	2197	rVB6	4490	8213	0.01%	0.002%
47	8.559	2239	2245	2252	rVB7	1891	2350	0.00%	0.001%
48	8.639	2252	2270	2300	rBV	465112	825928	0.51%	0.229%
49	8.813	2318	2324	2332	rVB6	4590	7795	0.00%	0.002%
50	8.870	2332	2342	2348	rBV3	5540	9044	0.01%	0.003%
51	9.105	2409	2415	2427	rVB5	1262	2367	0.00%	0.001%
52	9.218	2438	2450	2458	rBV5	5259	9480	0.01%	0.003%
53	9.337	2476	2487	2492	rBV6	2607	4639	0.00%	0.001%
54	9.382	2492	2501	2502	rBV2	7306	8681	0.01%	0.002%
55	9.424	2502	2514	2528	rVV	578970	949299	0.59%	0.264%
56	9.494	2528	2536	2547	rVV3	25348	57227	0.04%	0.016%
57	9.587	2547	2565	2586	rVV2	40613390	74130553	46.08%	20.595%
58	9.726	2587	2608	2643	rVB3	60127921	160884233	100.00%	44.697%
59	9.954	2672	2679	2691	rVB3	7020	11886	0.01%	0.003%
60	10.022	2691	2700	2710	rVB7	6794	12329	0.01%	0.003%
61	10.115	2711	2729	2759	rBV2	39077869	67365371	41.87%	18.716%
62	10.279	2774	2780	2789	rVB7	5047	8112	0.01%	0.002%
63	10.366	2793	2807	2818	rBV5	16749	30374	0.02%	0.008%
64	10.488	2832	2845	2864	rBV	1130138	1755362	1.09%	0.488%
65	10.568	2867	2870	2879	rVB6	2330	2943	0.00%	0.001%
66	10.658	2888	2898	2917	rVB10	6117	16629	0.01%	0.005%
67	10.761	2917	2930	2946	rBV	441495	707152	0.44%	0.196%
68	10.909	2964	2976	2990	rBV	489033	749452	0.47%	0.208%
69	11.002	2990	3005	3022	rVV	1530639	3264587	2.03%	0.907%
70	11.092	3022	3033	3047	rVB	759882	1174985	0.73%	0.326%
71	11.237	3066	3078	3085	rBV	51184	82623	0.05%	0.023%
72	11.295	3085	3096	3108	rVV	689456	1059344	0.66%	0.294%
73	11.411	3123	3132	3139	rBV7	7934	16380	0.01%	0.005%
74	11.472	3139	3151	3167	rVB	2524602	3784996	2.35%	1.052%
75	11.645	3191	3205	3216	rVB5	22337	45634	0.03%	0.013%
76	11.748	3224	3237	3245	rBV2	51051	86986	0.05%	0.024%

LSC Area Percent Report

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Acq On : 08 Oct 2021 00:39
Operator : SY/MD
Sample : M4138-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F3

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\SFAMULM092321WMA.M
Title : VOC Analysis

77	11.816	3245	3258	3271	rVB	619753	995806	0.62%	0.277%
78	11.899	3271	3284	3298	rBV	899859	1410885	0.88%	0.392%
79	12.018	3312	3321	3329	rBV4	21375	35317	0.02%	0.010%
80	12.099	3329	3346	3364	rVV	369276	642849	0.40%	0.179%
81	12.195	3364	3376	3397	rVB	673297	1106847	0.69%	0.308%
82	12.308	3401	3411	3418	rBV2	13277	21291	0.01%	0.006%
83	12.375	3419	3432	3451	rVB4	46149	99706	0.06%	0.028%
84	12.488	3451	3467	3485	rBV4	180008	493960	0.31%	0.137%
85	12.575	3486	3494	3502	rVV	77289	125632	0.08%	0.035%
86	12.616	3502	3507	3518	rVV2	25757	39641	0.02%	0.011%
87	12.687	3518	3529	3548	rVB2	74235	148949	0.09%	0.041%
88	12.819	3548	3570	3580	rBV	59714	137359	0.09%	0.038%
89	12.880	3580	3589	3600	rVB2	39983	66963	0.04%	0.019%
90	12.973	3601	3618	3626	rBV2	50513	132671	0.08%	0.037%
91	13.028	3626	3635	3646	rVB	79907	134354	0.08%	0.037%
92	13.298	3710	3719	3733	rVB	26328	40868	0.03%	0.011%
93	13.455	3747	3768	3779	rVV4	107331	193720	0.12%	0.054%
94	13.520	3780	3788	3799	rVB3	11389	17725	0.01%	0.005%
95	13.626	3812	3821	3831	rBV2	26797	41335	0.03%	0.011%
96	13.742	3842	3857	3868	rBV2	19172	33210	0.02%	0.009%
97	13.845	3882	3889	3900	rVB9	3149	5403	0.00%	0.002%
98	14.089	3951	3965	3983	rVV	172579	275650	0.17%	0.077%
99	15.230	4312	4320	4330	rBV4	1299	2316	0.00%	0.001%
100	15.417	4371	4378	4386	rBV4	1431	2375	0.00%	0.001%

Sum of corrected areas: 359941241

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU100721\
Data File : VU045231.D
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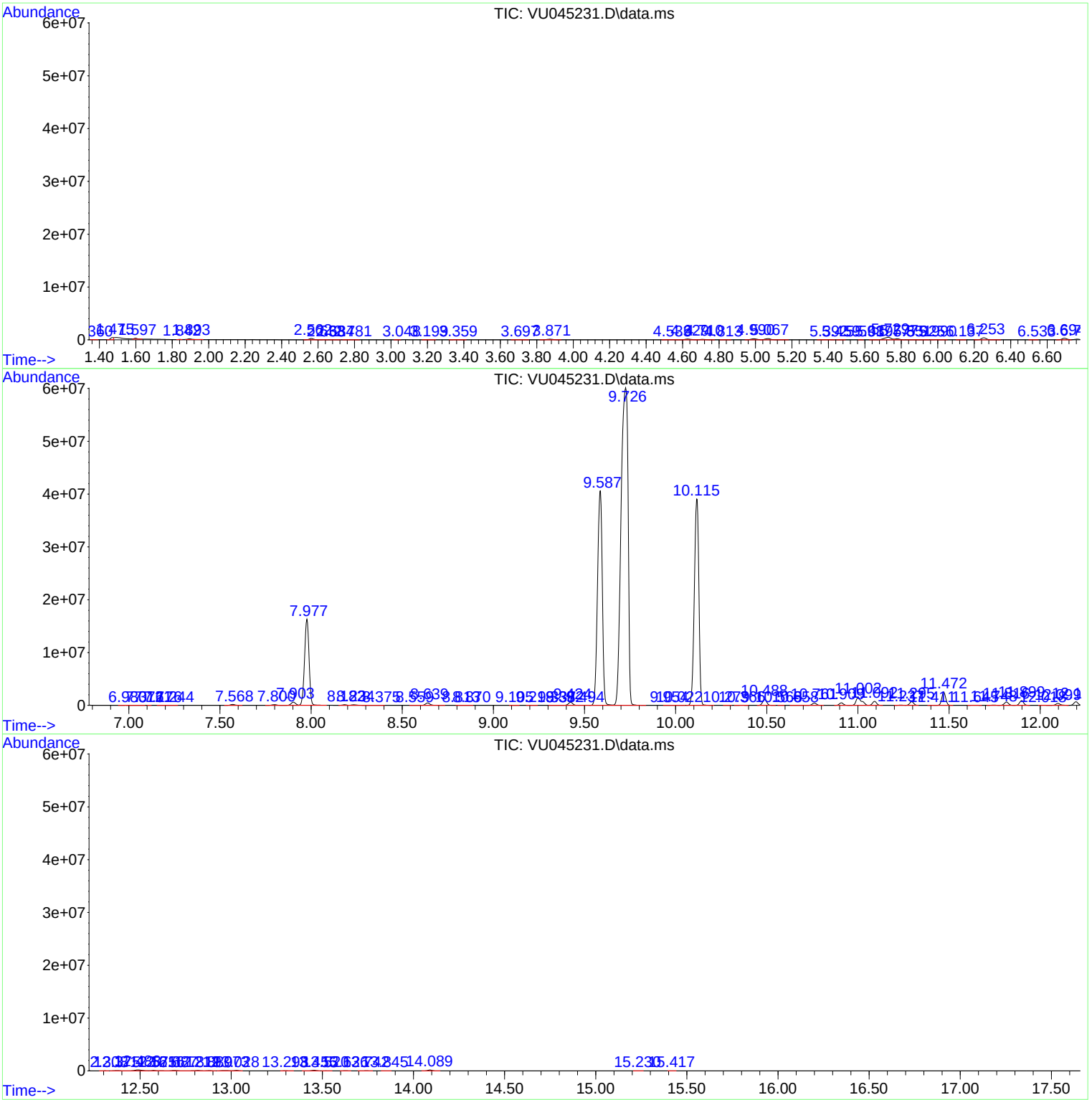
Operator : SY/MD
Sample : M4138-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW7F3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092321WMA.M
Quant Title : VOC Analysis

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU100721\
Data File : VU045231.D
Acq On : 08 Oct 2021 00:39
Operator : SY/MD
Sample : M4138-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F3

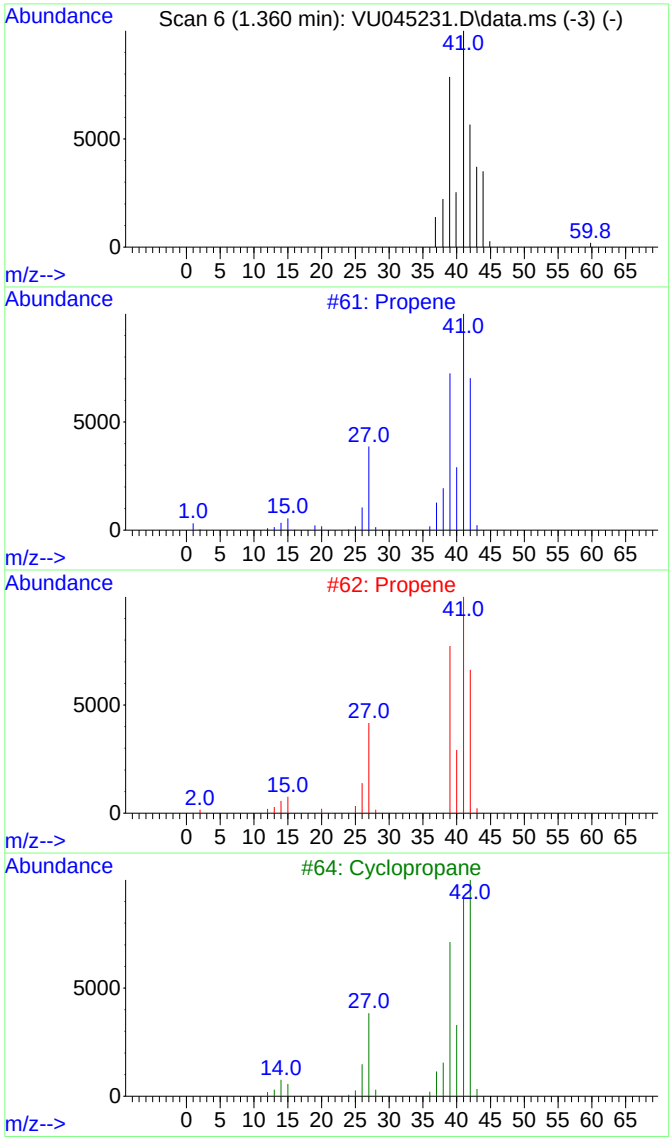
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092321WMA.M
Quant Title : VOC Analysis

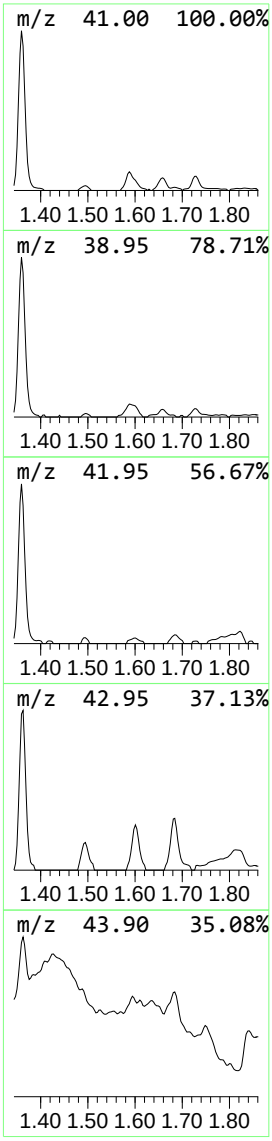
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.360	2.59 ug/L	37543	1,4-Difluorobenzene	6.253

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propene	42	C3H6	000115-07-1	90
2	Propene	42	C3H6	000115-07-1	42
3	Cyclopropane	42	C3H6	000075-19-4	9
4	Cyclopropene	40	C3H4	002781-85-3	9
5	Cyclopropane	42	C3H6	000075-19-4	9





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU100721\
Data File : VU045231.D
Acq On : 08 Oct 2021 00:39
Operator : SY/MD
Sample : M4138-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092321WMA.M
Quant Title : VOC Analysis

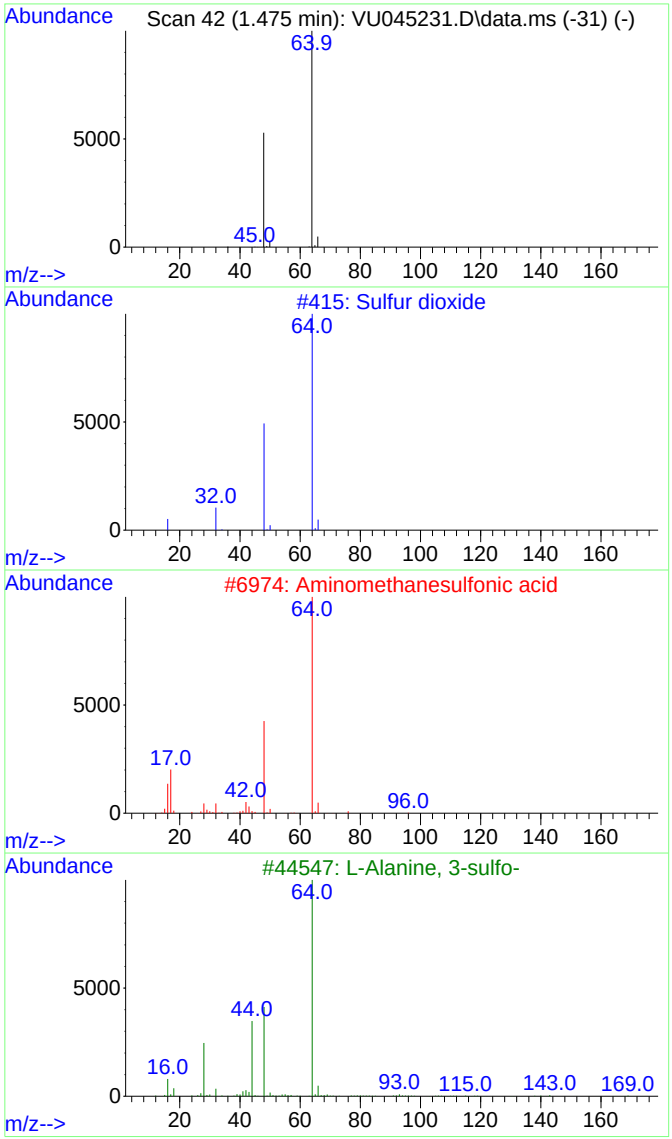
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

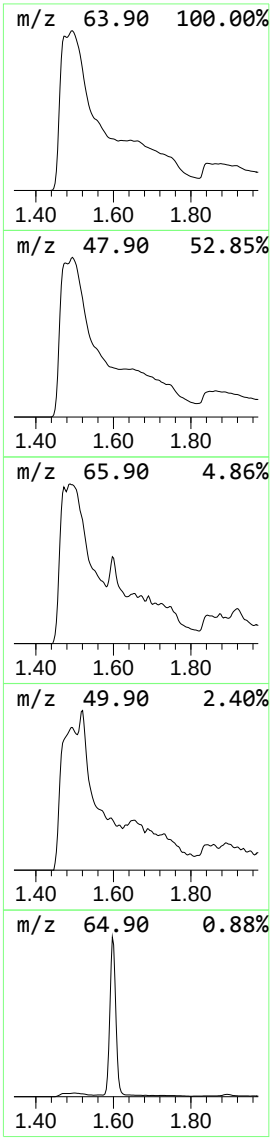
****

Peak Number 2 Sulfur dioxide Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.475	34.56 ug/L	501913	1,4-Difluorobenzene	6.253

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Sulfur dioxide	64	O2S	007446-09-5	83
2	Aminomethanesulfonic acid	111	CH5NO3S	013881-91-9	74
3	L-Alanine, 3-sulfo-	169	C3H7NO5S	000498-40-8	9
4	Aminomethanesulfonic acid	111	CH5NO3S	013881-91-9	9
5	Aminomethanesulfonic acid	111	CH5NO3S	013881-91-9	9





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU100721\
Data File : VU045231.D
Acq On : 08 Oct 2021 00:39
Operator : SY/MD
Sample : M4138-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F3

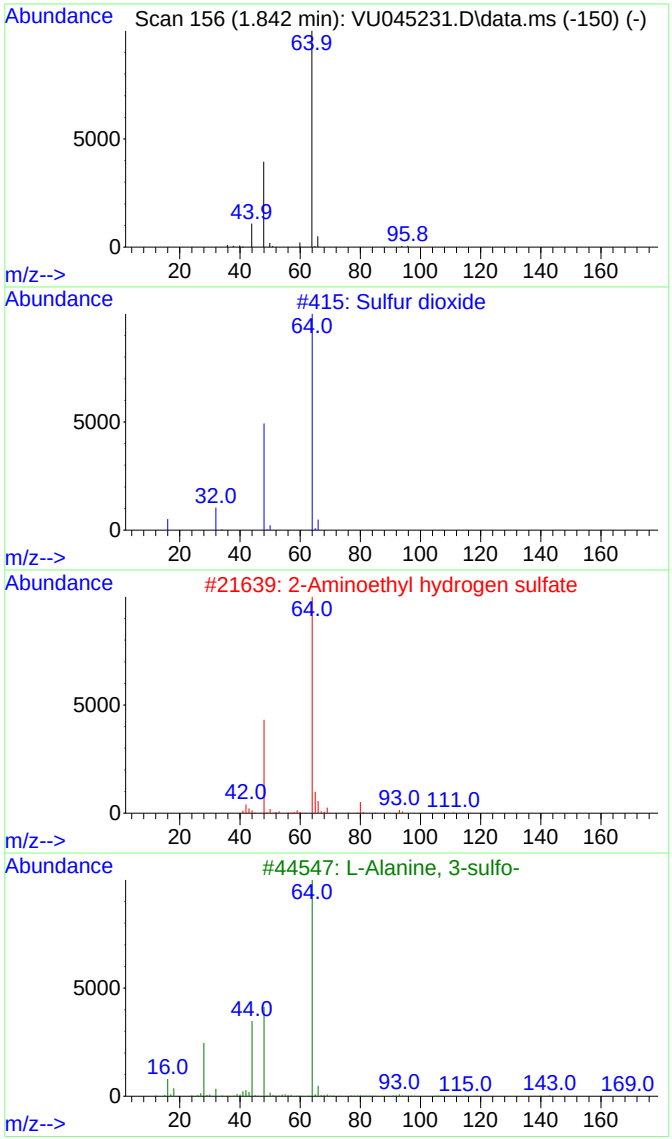
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092321WMA.M
Quant Title : VOC Analysis

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

****
Peak Number 3 2-Aminoethyl hydrogen sulfate Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.842	3.80 ug/L	55166	1,4-Difluorobenzene	6.253

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Sulfur dioxide	64	O2S	007446-09-5	74
2	2-Aminoethyl hydrogen sulfate	141	C2H7NO4S	000926-39-6	64
3	L-Alanine, 3-sulfo-	169	C3H7NO5S	000498-40-8	64
4	(N-(-2-Acetamido))-2-aminoethane...	182	C4H10N2O4S	007365-82-4	64
5	Aminomethanesulfonic acid	111	CH5NO3S	013881-91-9	9



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU100721\
Data File : VU045231.D
Acq On : 08 Oct 2021 00:39
Operator : SY/MD
Sample : M4138-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW7F3

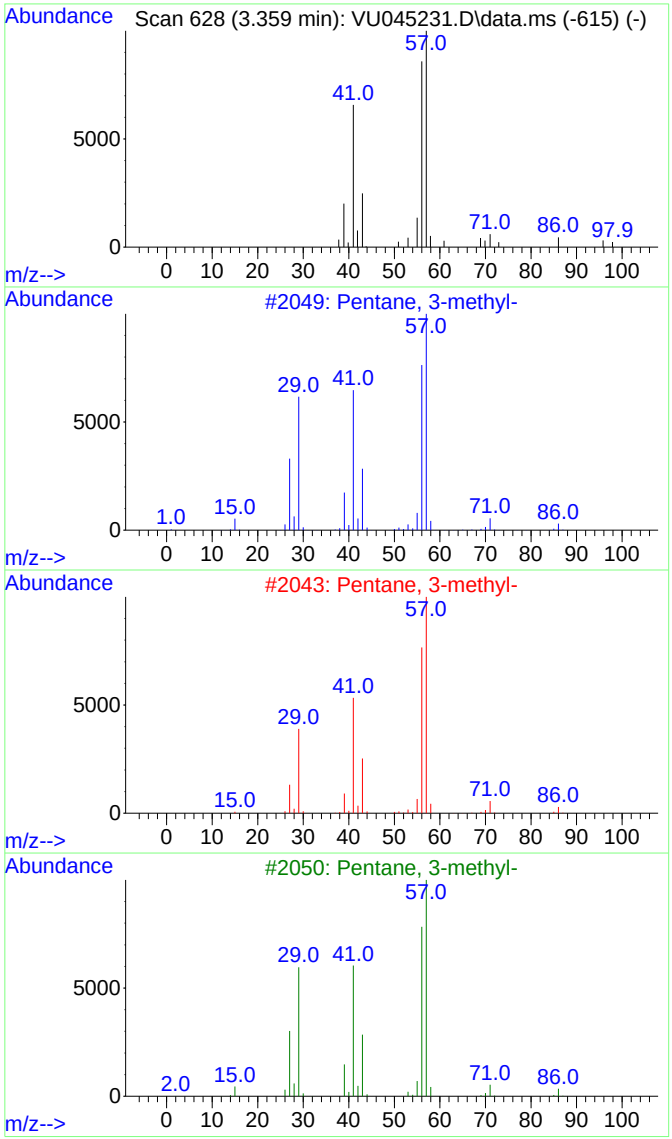
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092321WMA.M
Quant Title : VOC Analysis

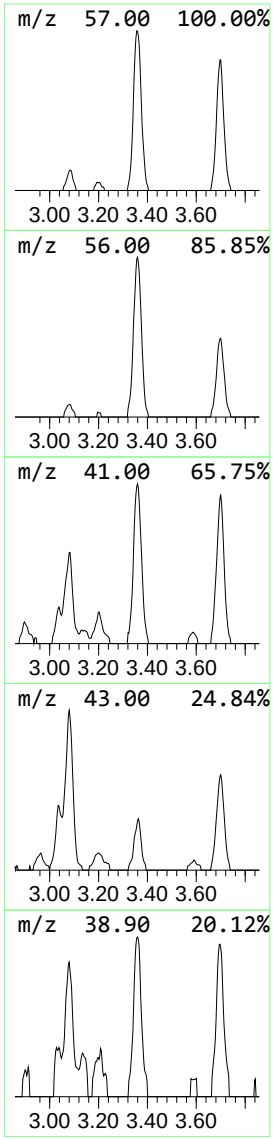
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.359	2.91 ug/L	42267	1,4-Difluorobenzene	6.253

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 3-methyl-	86	C6H14	000096-14-0	83
2	Pentane, 3-methyl-	86	C6H14	000096-14-0	80
3	Pentane, 3-methyl-	86	C6H14	000096-14-0	80
4	Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	64
5	Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	56





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU100721\
Data File : VU045231.D
Acq On : 08 Oct 2021 00:39
Operator : SY/MD
Sample : M4138-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F3

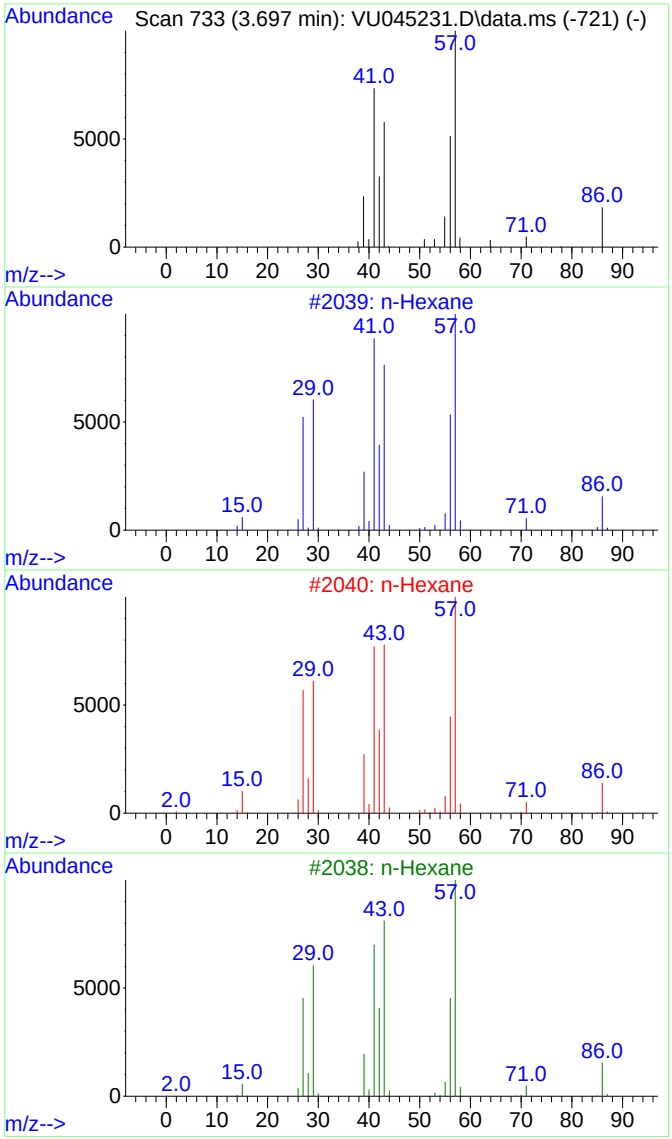
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092321WMA.M
Quant Title : VOC Analysis

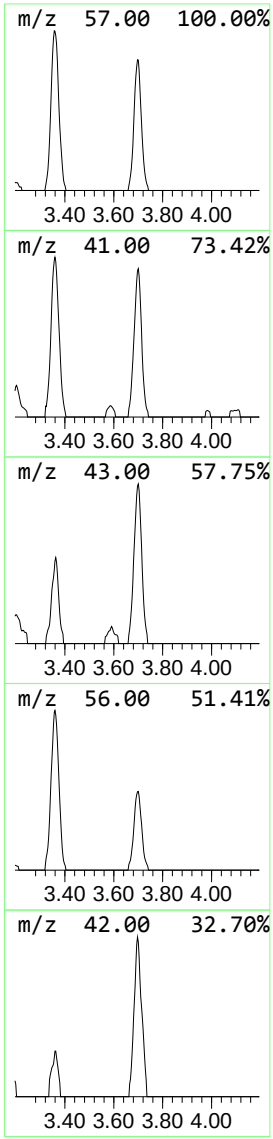
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.697	2.98 ug/L	43340	1,4-Difluorobenzene	6.253

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexane	86	C6H14	000110-54-3	78
2	n-Hexane	86	C6H14	000110-54-3	72
3	n-Hexane	86	C6H14	000110-54-3	72
4	2-Ethyl-oxetane	86	C5H10O	1010386-40-2	72
5	n-Hexane	86	C6H14	000110-54-3	53





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU100721\
Data File : VU045231.D
Acq On : 08 Oct 2021 00:39
Operator : SY/MD
Sample : M4138-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F3

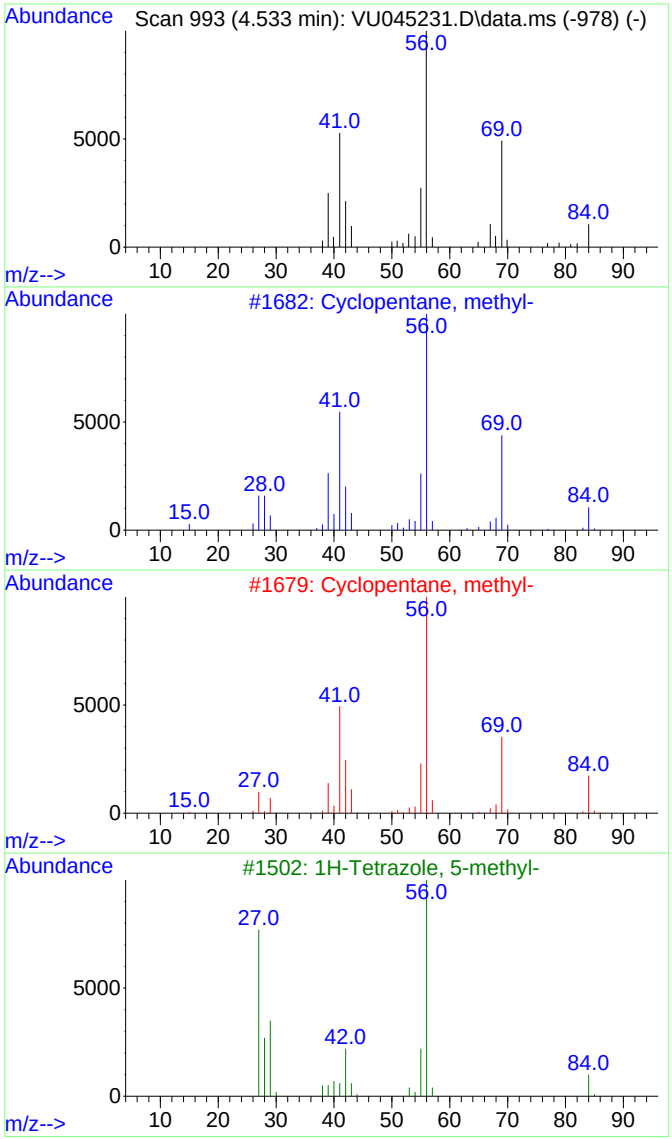
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092321WMA.M
Quant Title : VOC Analysis

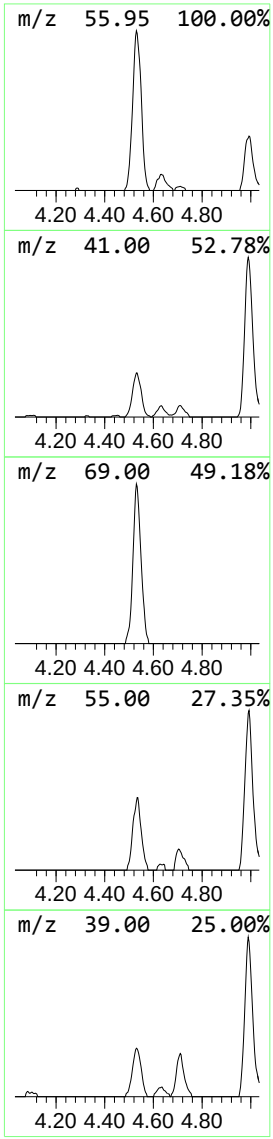
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 (DEL) Alkane: Cyclic4.533 Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.533	4.77 ug/L	69246	1,4-Difluorobenzene	6.253

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	91
2	Cyclopentane, methyl-	84	C6H12	000096-37-7	80
3	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	72
4	Cyclopentane, methyl-	84	C6H12	000096-37-7	64
5	Cyclobutane, ethyl-	84	C6H12	004806-61-5	64





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Acq On : 08 Oct 2021 00:39
Operator : SY/MD
Sample : M4138-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW7F3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092321WMA.M
Quant Title : VOC Analysis

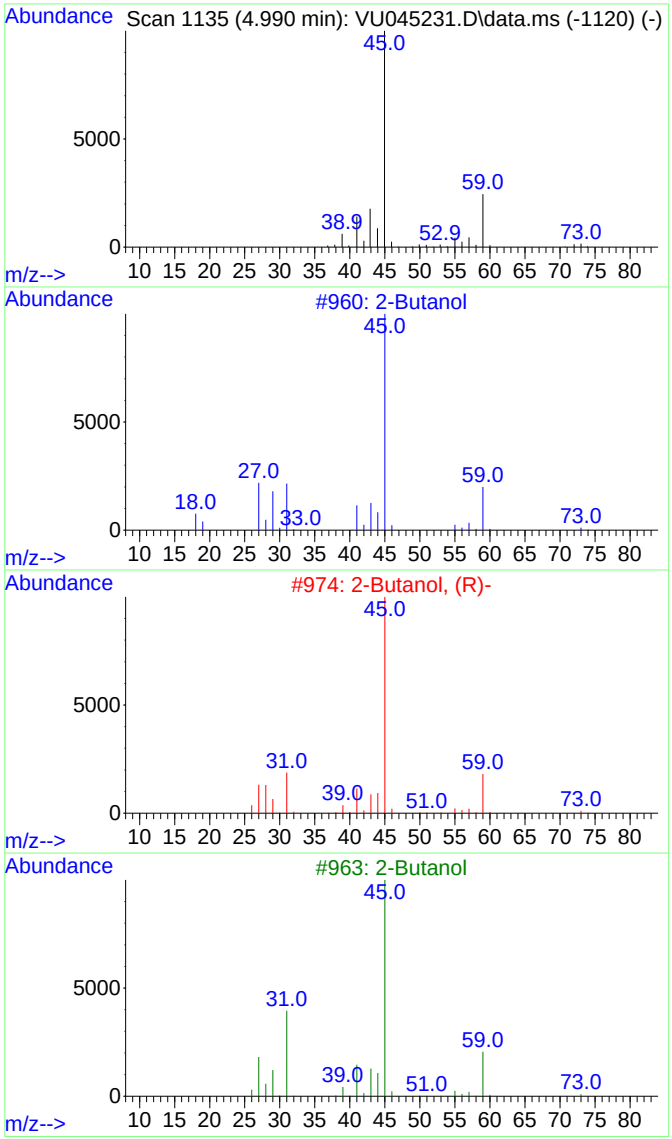
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

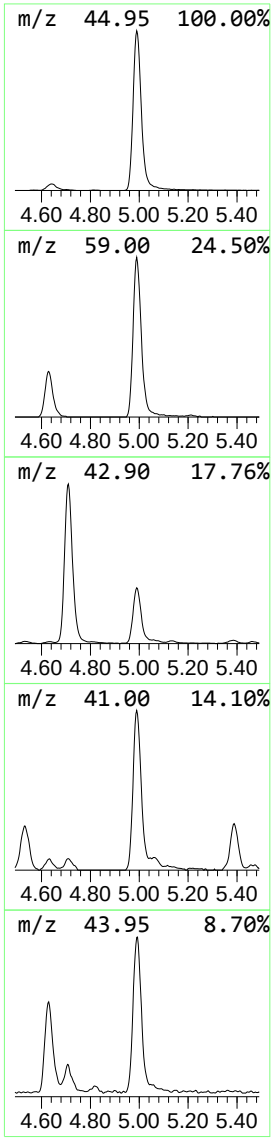
****

Peak Number 7 2-Butanol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.990	32.19 ug/L	467364	1,4-Difluorobenzene	6.253

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butanol	74	C4H10O	000078-92-2	83
2	2-Butanol, (R)-	74	C4H10O	014898-79-4	83
3	2-Butanol	74	C4H10O	000078-92-2	78
4	2-Butanol	74	C4H10O	000078-92-2	74
5	2-Butanol, (R)-	74	C4H10O	014898-79-4	72





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Acq On : 08 Oct 2021 00:39
Operator : SY/MD
Sample : M4138-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F3

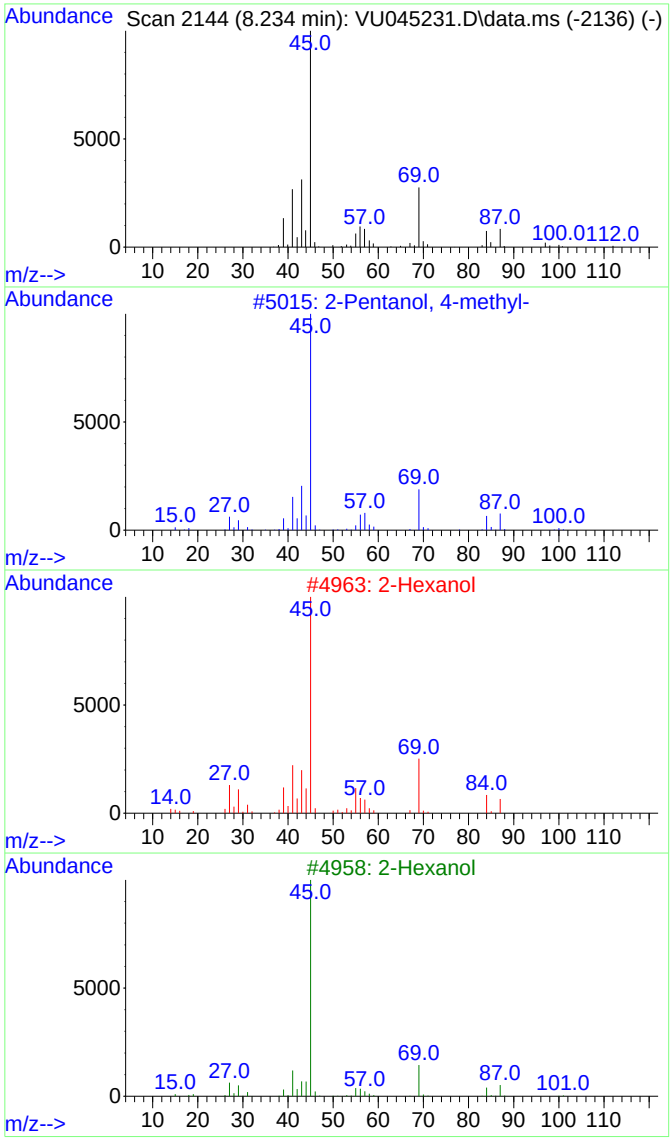
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092321WMA.M
Quant Title : VOC Analysis

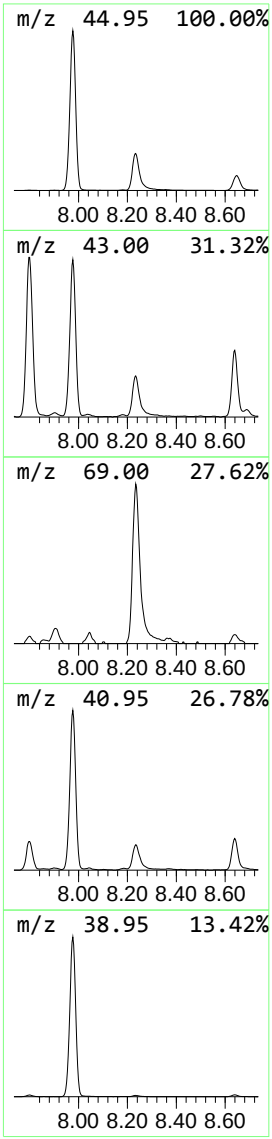
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 2-Pentanol, 4-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.234	12.28 ug/L	233088	Chlorobenzene-d5	9.424

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	90
2	2-Hexanol	102	C6H14O	000626-93-7	83
3	2-Hexanol	102	C6H14O	000626-93-7	78
4	2-Nonanol	144	C9H20O	000628-99-9	78
5	2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	72





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Acq On : 08 Oct 2021 00:39
Operator : SY/MD
Sample : M4138-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092321WMA.M
Quant Title : VOC Analysis

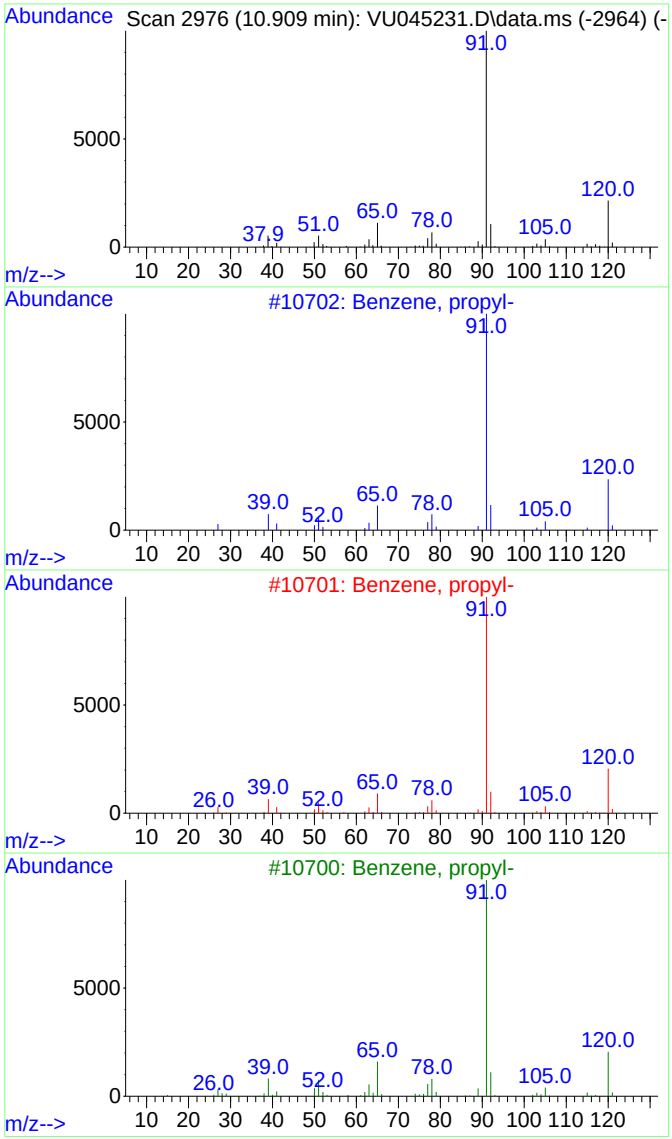
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

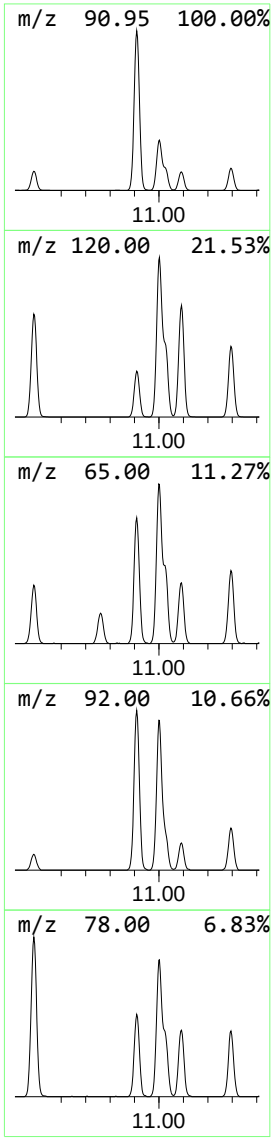
****

Peak Number 10 Benzene, propyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.909	37.63 ug/L	749452	1,4-Dichlorobenzene-d4	11.816

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	90
3	Benzene, propyl-	120	C9H12	000103-65-1	90
4	Benzene, propyl-	120	C9H12	000103-65-1	86
5	Phenethylamine, N-benzyl-p-chloro-	245	C15H16ClN	013622-43-0	72





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Acq On : 08 Oct 2021 00:39
Operator : SY/MD
Sample : M4138-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092321WMA.M
Quant Title : VOC Analysis

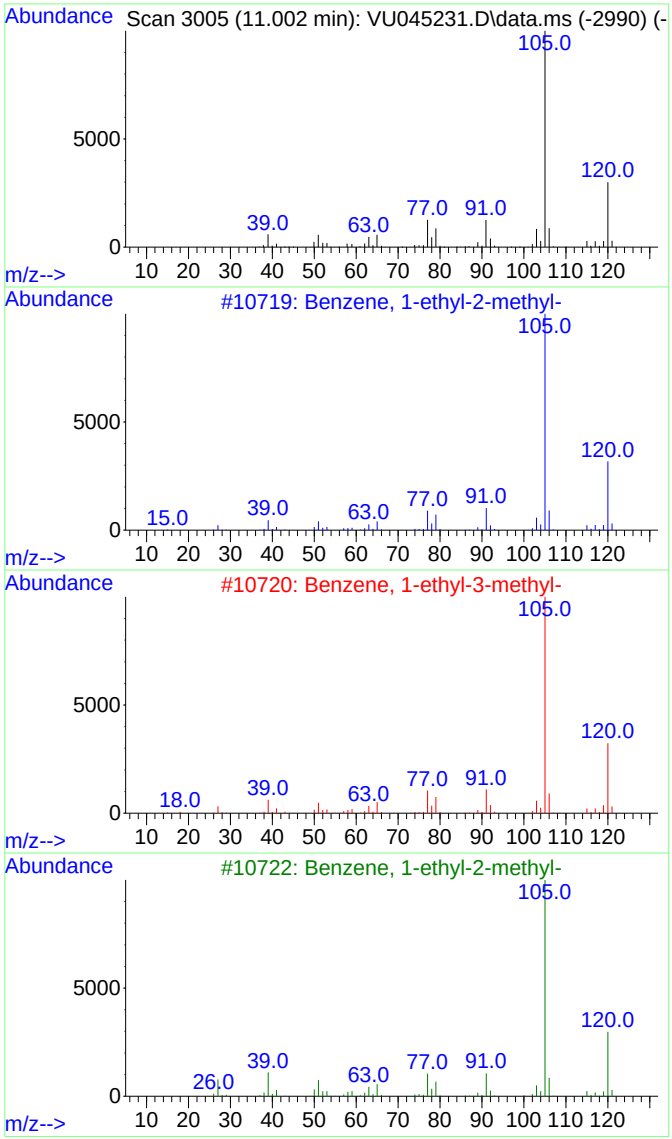
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

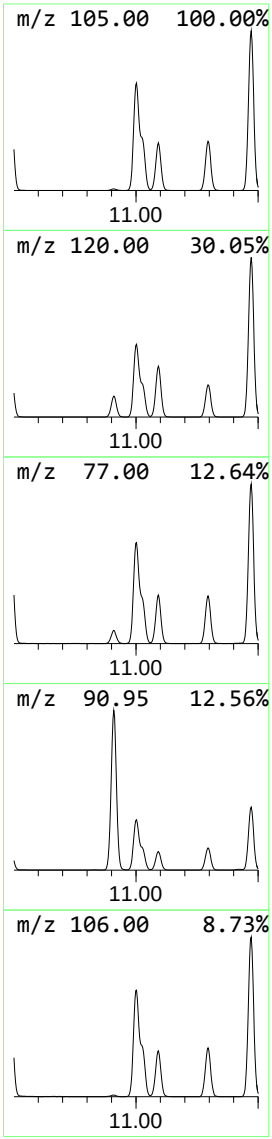
****

Peak Number 11 Benzene, 1-ethyl-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.002	163.92 ug/L	3264590	1,4-Dichlorobenzene-d4	11.816

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





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Acq On : 08 Oct 2021 00:39
Operator : SY/MD
Sample : M4138-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F3

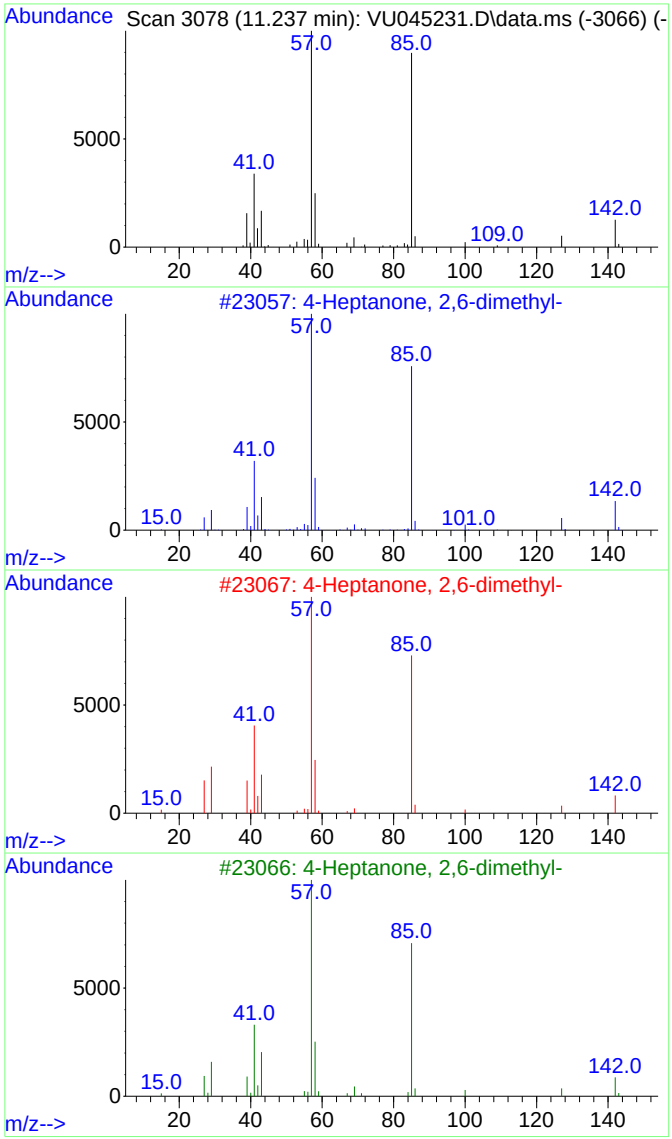
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092321WMA.M
Quant Title : VOC Analysis

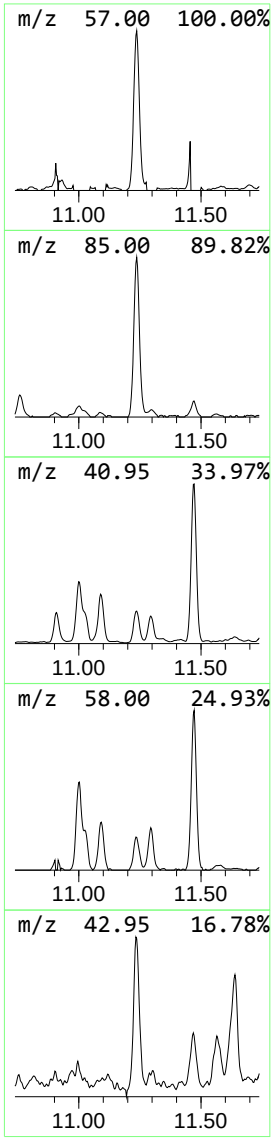
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 4-Heptanone, 2,6-dimethyl- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.237	4.15 ug/L	82623	1,4-Dichlorobenzene-d4	11.816

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	4-Octanone, 2-methyl-	142	C9H18O	007492-38-8	72





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Acq On : 08 Oct 2021 00:39
Operator : SY/MD
Sample : M4138-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092321WMA.M
Quant Title : VOC Analysis

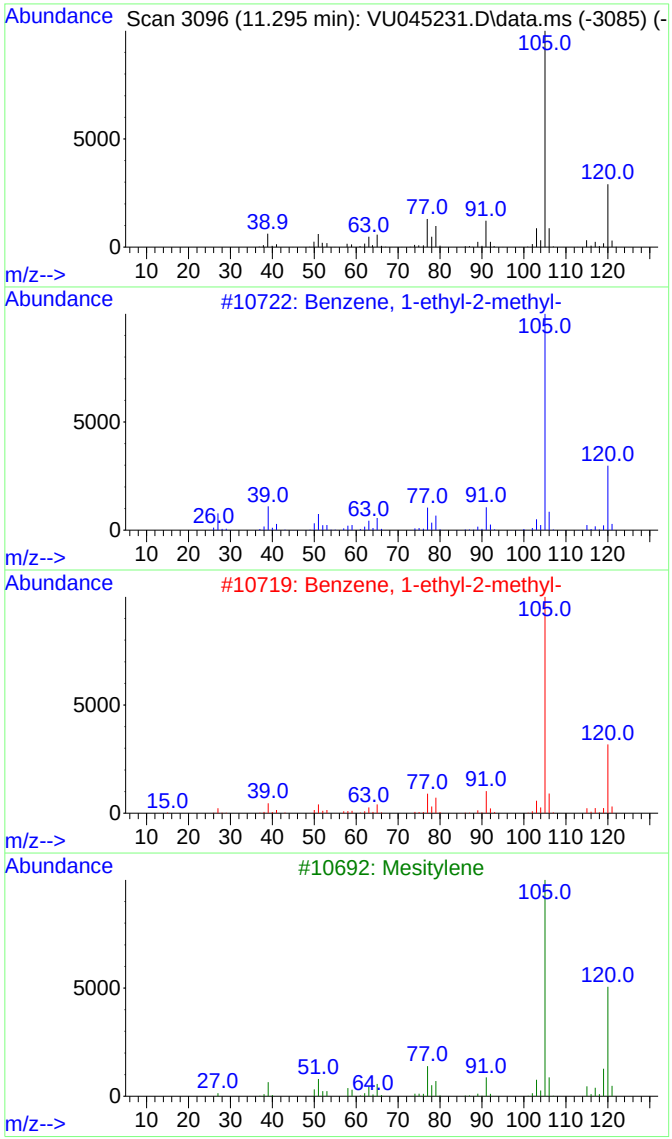
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

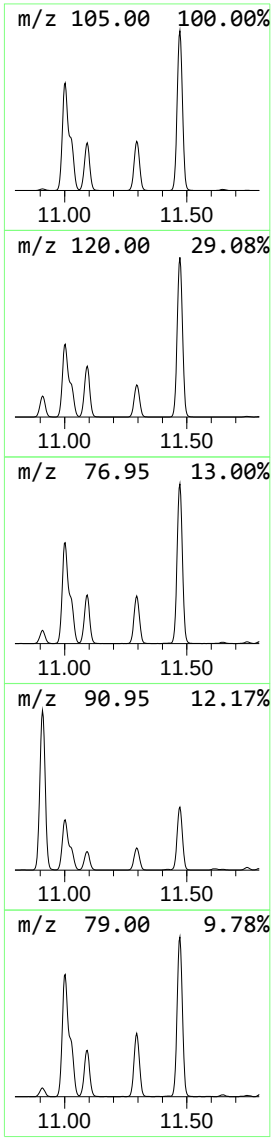
****

Peak Number 13 Benzene, 1-ethyl-4-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.295	53.19 ug/L	1059340	1,4-Dichlorobenzene-d4	11.816

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-2-methyl-	120	C9H12	000611-14-3	94
3	Mesitylene	120	C9H12	000108-67-8	91
4	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91





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Acq On : 08 Oct 2021 00:39
Operator : SY/MD
Sample : M4138-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW7F3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092321WMA.M
Quant Title : VOC Analysis

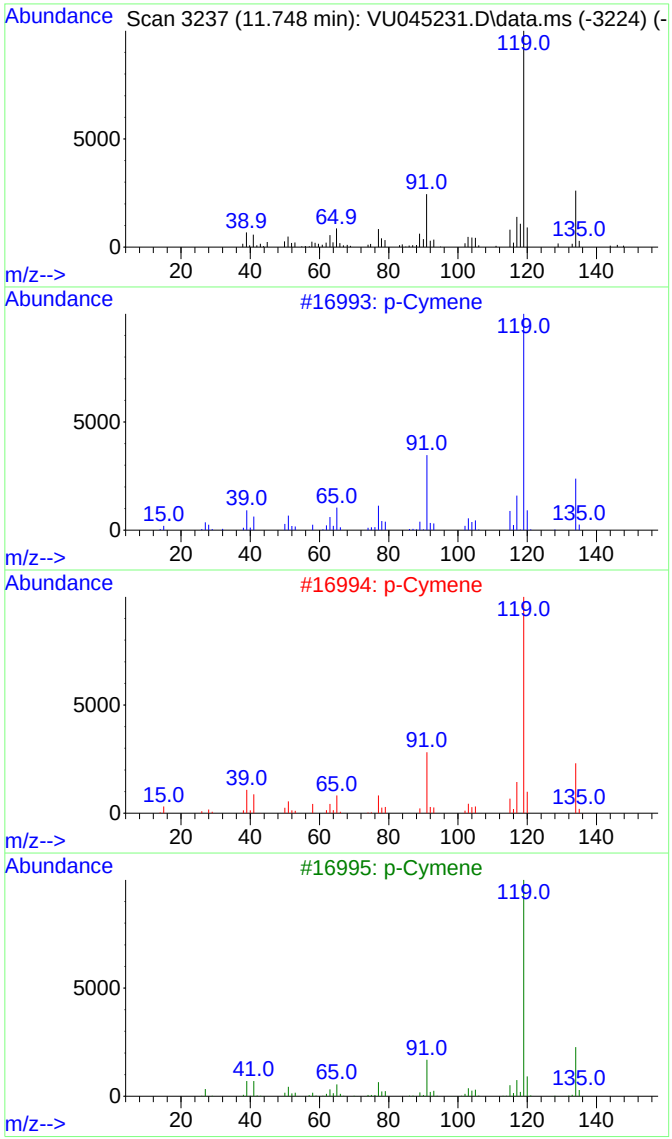
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

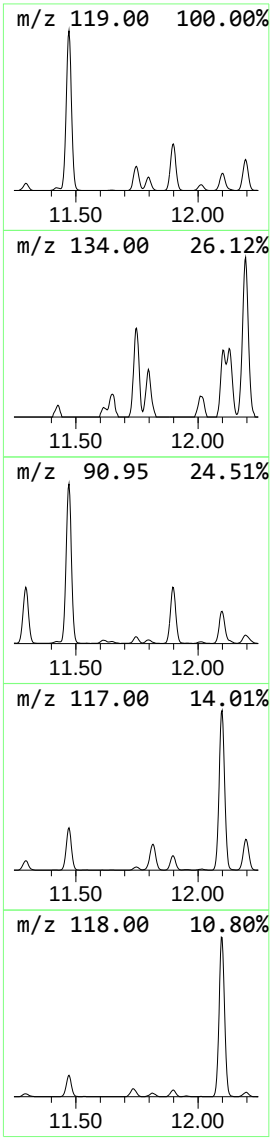
****

Peak Number 14 p-Cymene Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.748	4.37 ug/L	86986	1,4-Dichlorobenzene-d4	11.816

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Cymene	134	C10H14	000099-87-6	95
2	p-Cymene	134	C10H14	000099-87-6	94
3	p-Cymene	134	C10H14	000099-87-6	93
4	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	87
5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	87





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Acq On : 08 Oct 2021 00:39
Operator : SY/MD
Sample : M4138-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F3

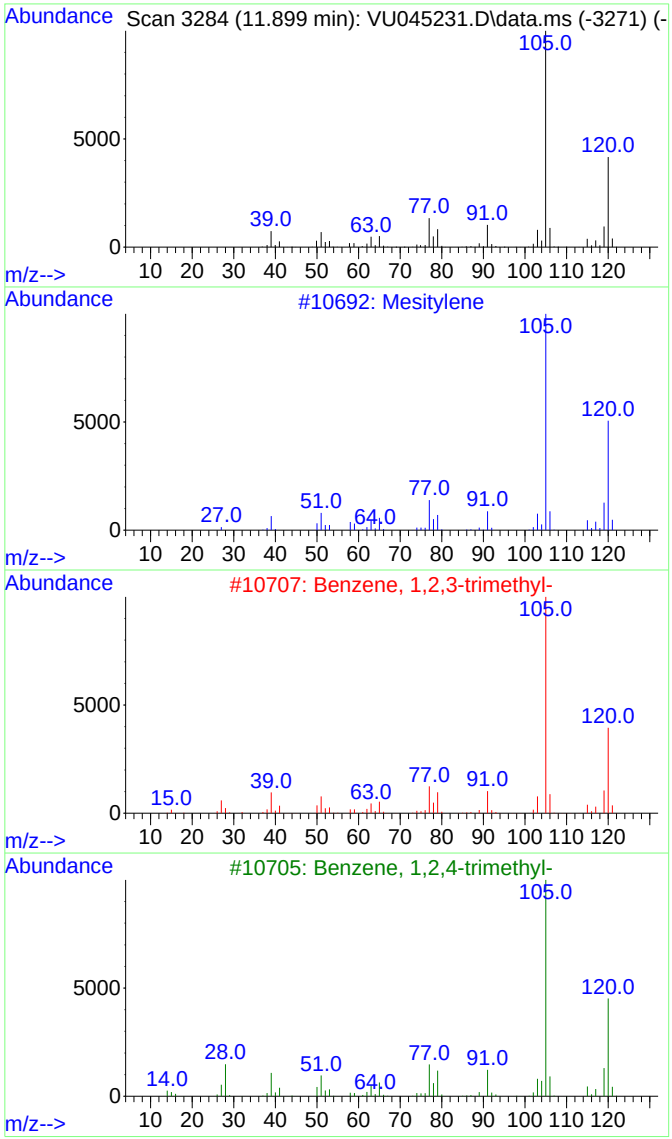
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092321WMA.M
Quant Title : VOC Analysis

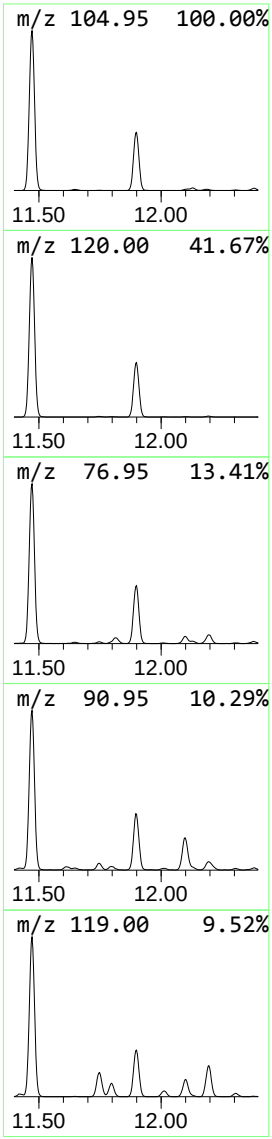
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 Benzene, 1,2,3-trimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.899	70.84 ug/L	1410890	1,4-Dichlorobenzene-d4	11.816

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Mesitylene	120	C9H12	000108-67-8	97
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
5	Mesitylene	120	C9H12	000108-67-8	91





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Data File : VU045231.D
Acq On : 08 Oct 2021 00:39
Operator : SY/MD
Sample : M4138-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092321WMA.M
Quant Title : VOC Analysis

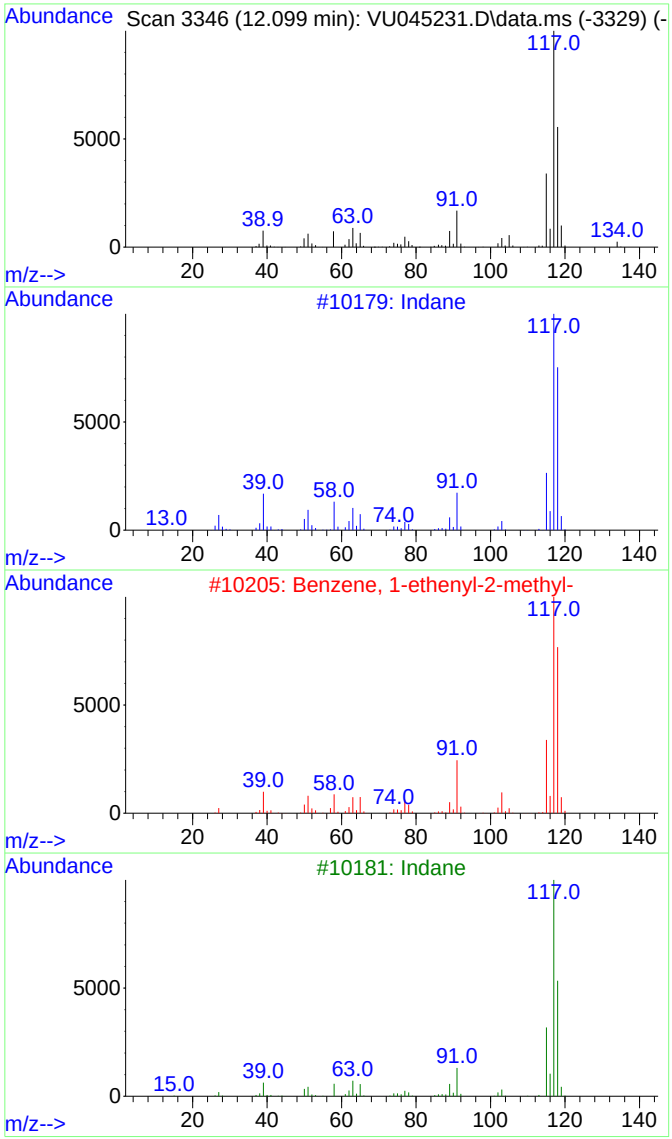
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

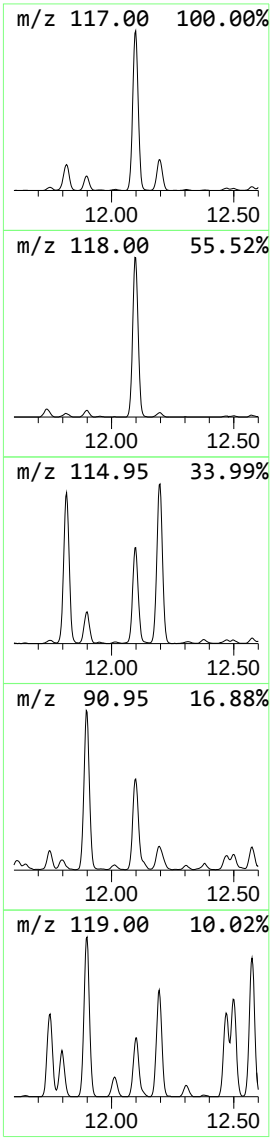
****

Peak Number 16 Indane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.099	32.28 ug/L	642849	1,4-Dichlorobenzene-d4	11.816

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	91
2	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	90
3	Indane	118	C9H10	000496-11-7	87
4	Benzene, cyclopropyl-	118	C9H10	000873-49-4	81
5	Benzene, 2-propenyl-	118	C9H10	000300-57-2	81





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Acq On : 08 Oct 2021 00:39
Operator : SY/MD
Sample : M4138-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F3

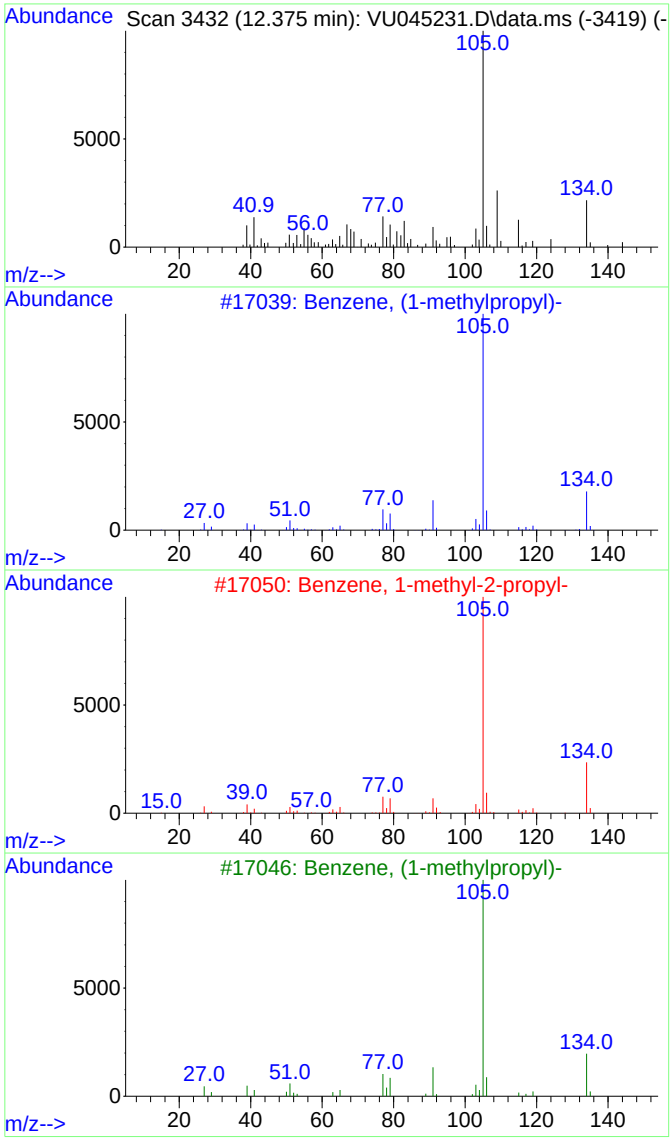
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092321WMA.M
Quant Title : VOC Analysis

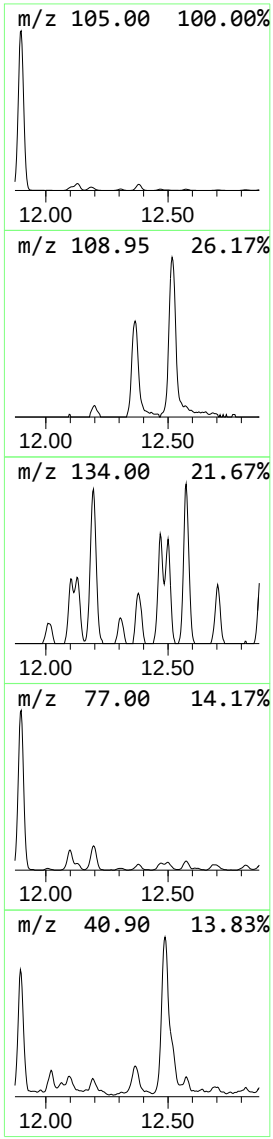
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 Benzene, (1-methylpropyl)- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.375	5.01 ug/L	99706	1,4-Dichlorobenzene-d4	11.816

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	60
2	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	60
3	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	55
4	Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	55
5	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	55





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Data File : VU045231.D
Acq On : 08 Oct 2021 00:39
Operator : SY/MD
Sample : M4138-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F3

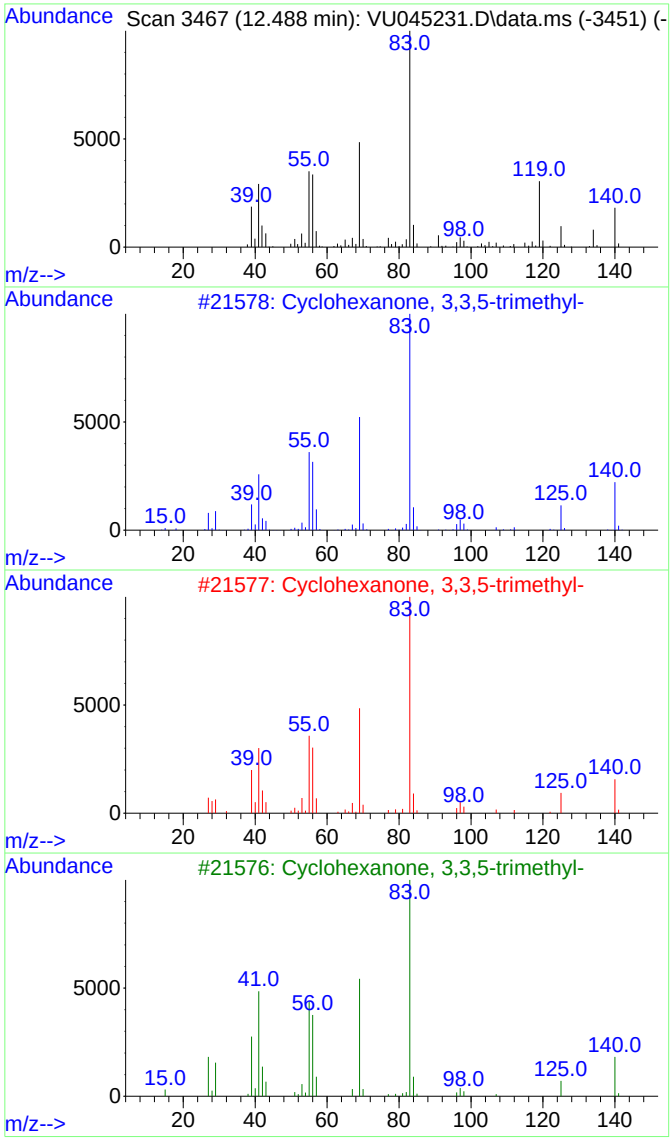
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092321WMA.M
Quant Title : VOC Analysis

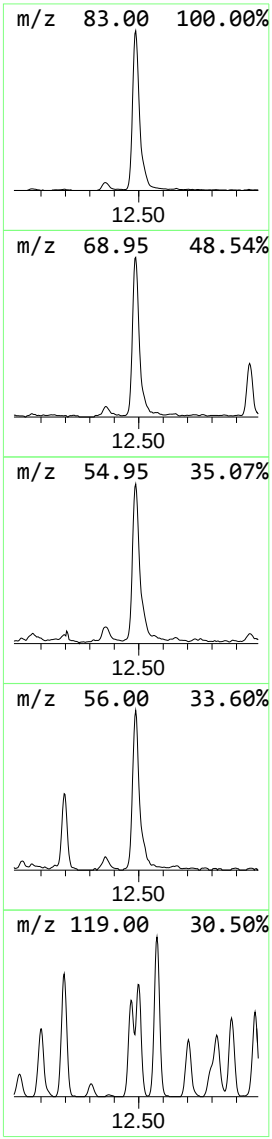
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 18 Cyclohexanone, 3,3,5-trimet... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.488	24.80 ug/L	493960	1,4-Dichlorobenzene-d4	11.816

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	95
3	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	94
4	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	87
5	Cyclopentanone, 2,4,4-trimethyl-	126	C8H14O	004694-12-6	53





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU100721\
Data File : VU045231.D
Acq On : 08 Oct 2021 00:39
Operator : SY/MD
Sample : M4138-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F3

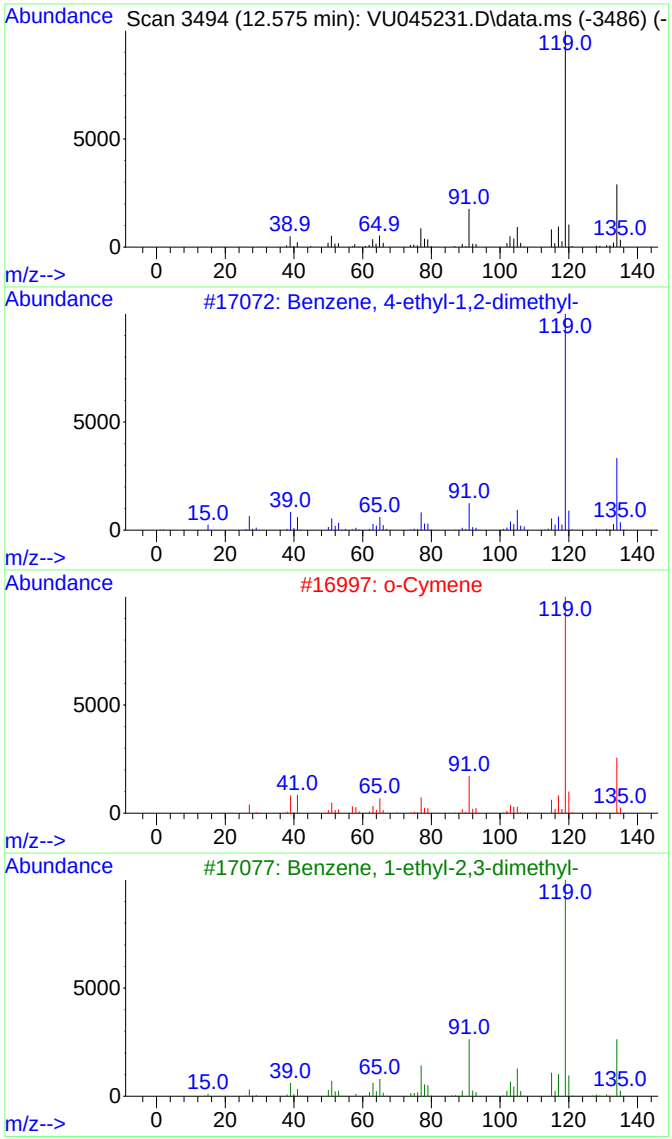
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092321WMA.M
Quant Title : VOC Analysis

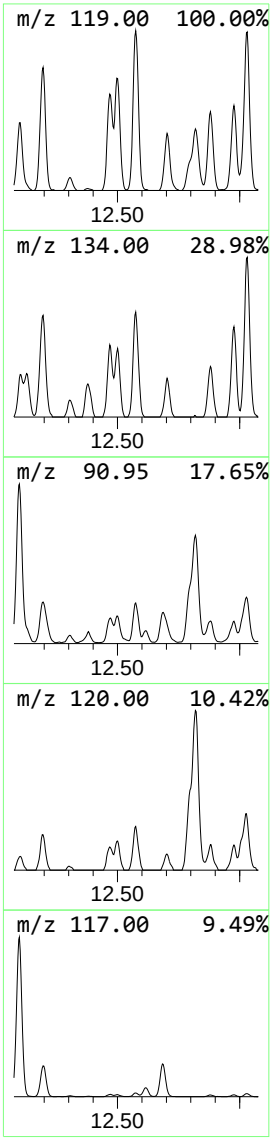
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.575	6.31 ug/L	125632	1,4-Dichlorobenzene-d4	11.816

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
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	94
5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU100721\
Data File : VU045231.D
Acq On : 08 Oct 2021 00:39
Operator : SY/MD
Sample : M4138-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW7F3

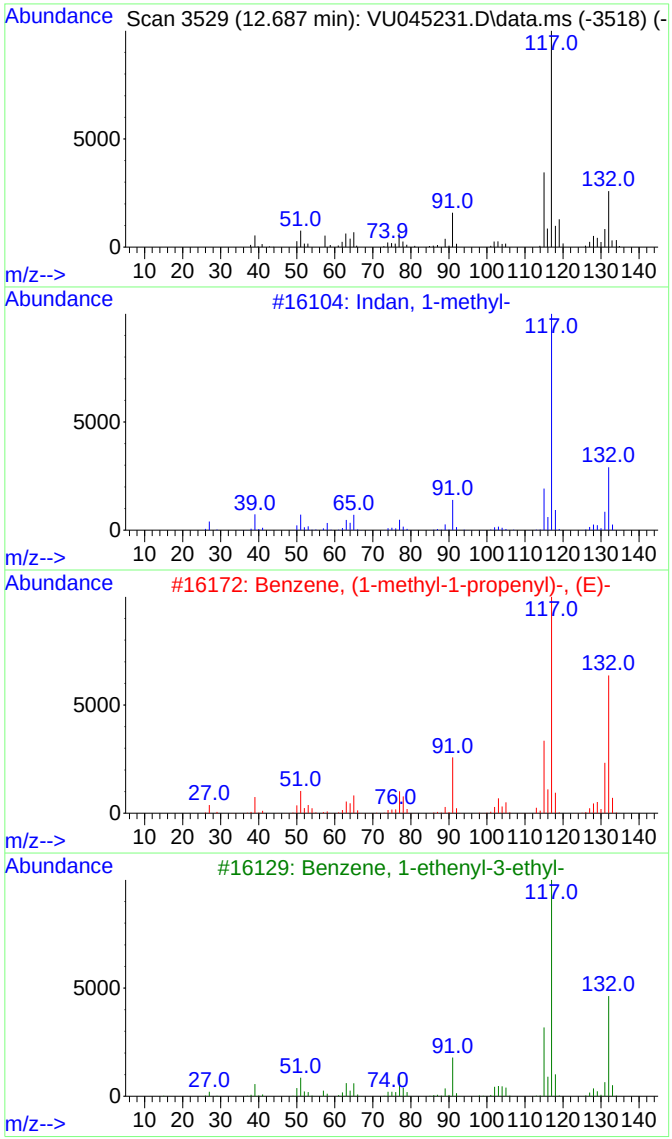
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092321WMA.M
Quant Title : VOC Analysis

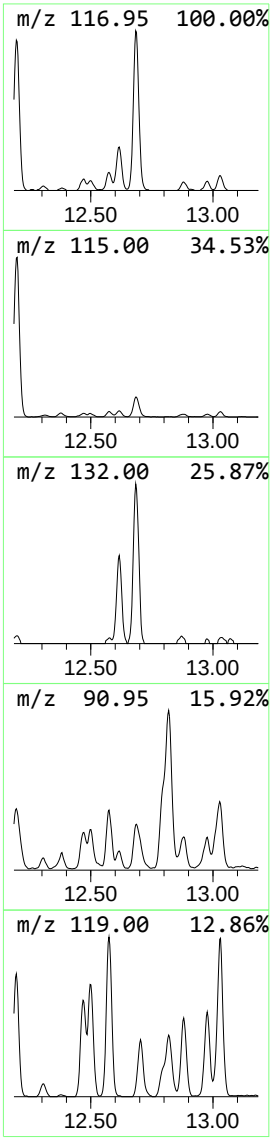
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 Indan, 1-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.687	7.48 ug/L	148949	1,4-Dichlorobenzene-d4	11.816

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indan, 1-methyl-	132	C10H12	000767-58-8	93
2	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	87
3	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	86
4	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	83
5	1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	81





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Data File : VU045231.D
Acq On : 08 Oct 2021 00:39
Operator : SY/MD
Sample : M4138-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F3

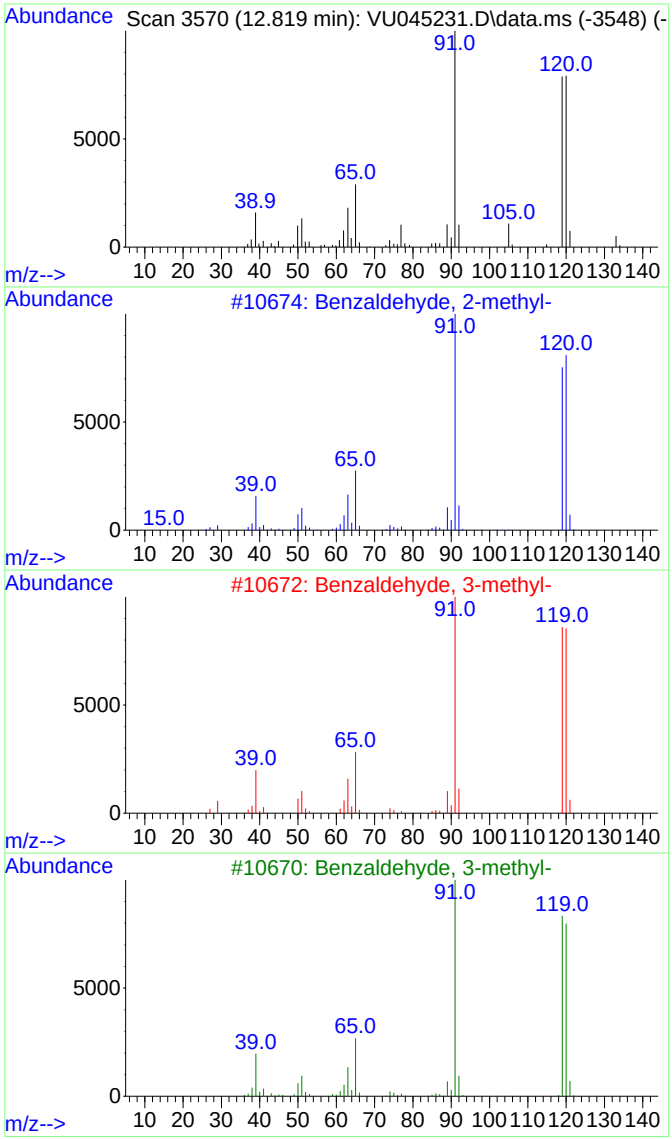
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092321WMA.M
Quant Title : VOC Analysis

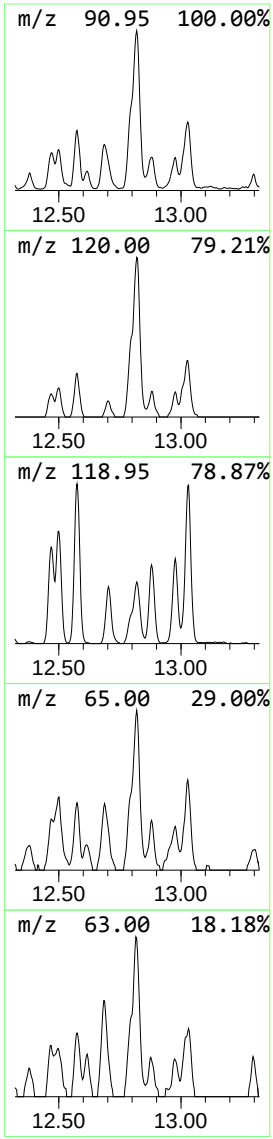
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 21 Benzaldehyde, 2-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.819	6.90 ug/L	137359	1,4-Dichlorobenzene-d4	11.816

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzaldehyde, 2-methyl-	120	C8H8O	000529-20-4	97
2	Benzaldehyde, 3-methyl-	120	C8H8O	000620-23-5	97
3	Benzaldehyde, 3-methyl-	120	C8H8O	000620-23-5	96
4	Benzaldehyde, 2-methyl-	120	C8H8O	000529-20-4	96
5	Benzaldehyde, 2-methyl-	120	C8H8O	000529-20-4	96





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU100721\
Data File : VU045231.D
Acq On : 08 Oct 2021 00:39
Operator : SY/MD
Sample : M4138-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F3

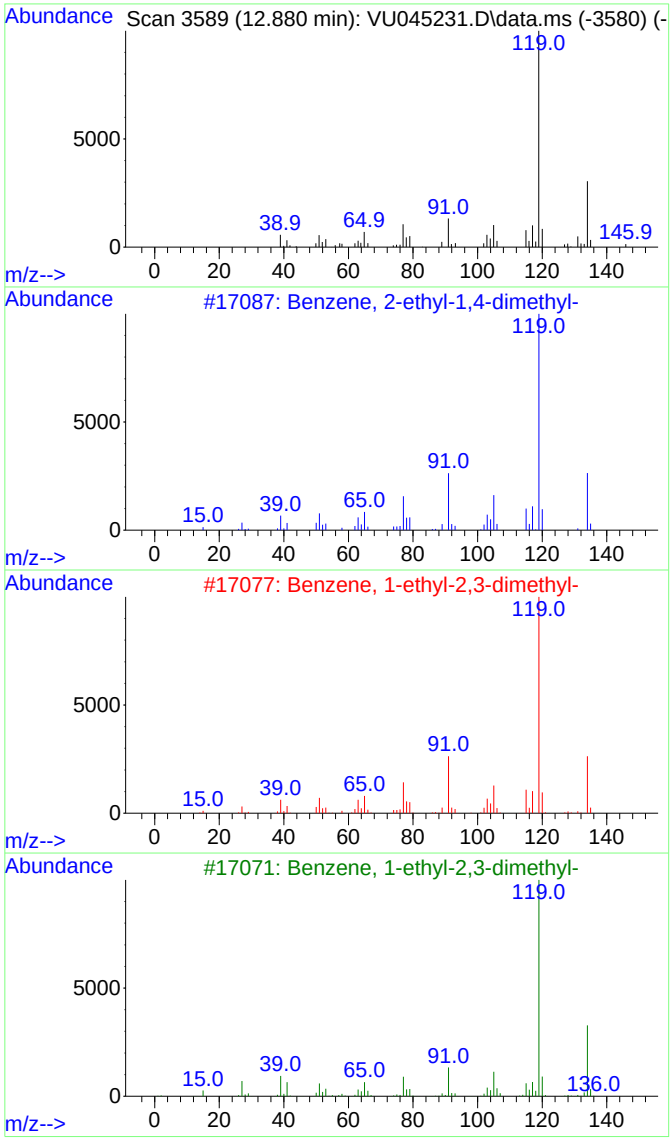
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092321WMA.M
Quant Title : VOC Analysis

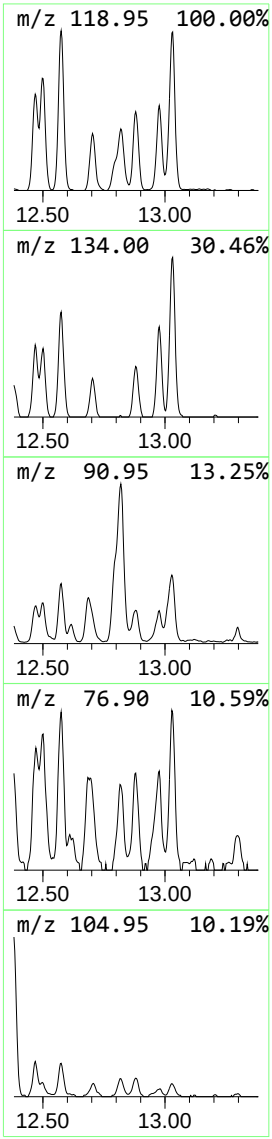
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 22 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.880	3.36 ug/L	66963	1,4-Dichlorobenzene-d4	11.816

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
3	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	94
4	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	94
5	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	93





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU100721\
Data File : VU045231.D
Acq On : 08 Oct 2021 00:39
Operator : SY/MD
Sample : M4138-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F3

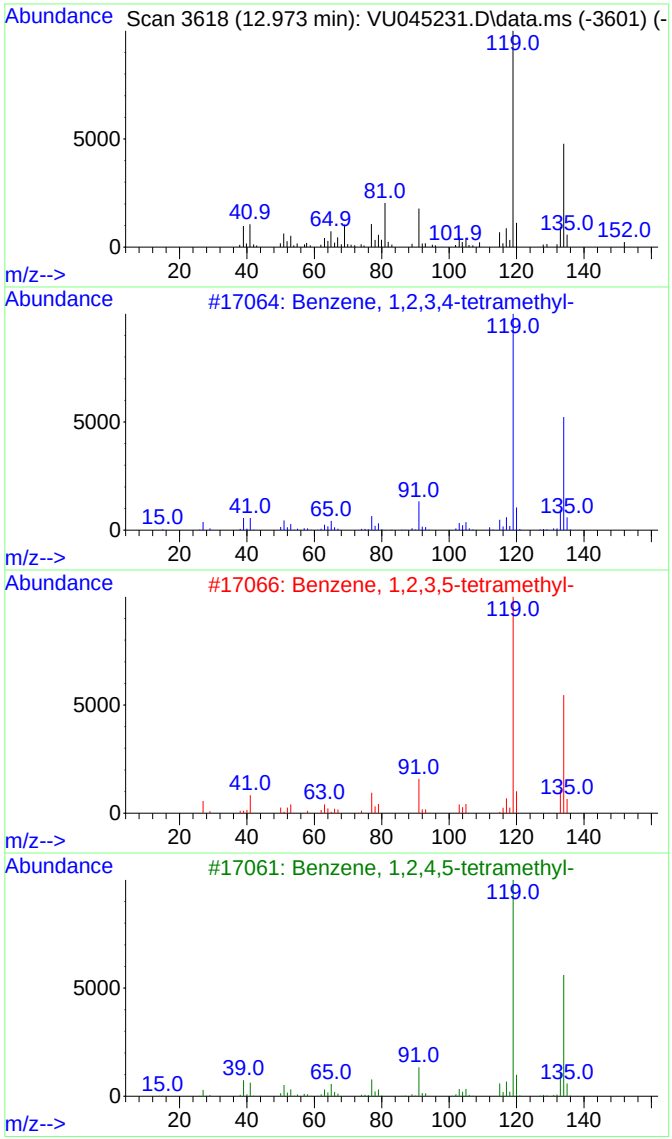
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092321WMA.M
Quant Title : VOC Analysis

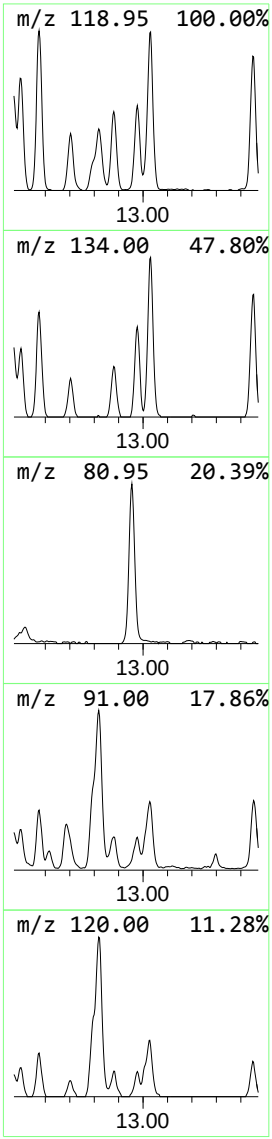
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.973	6.66 ug/L	132671	1,4-Dichlorobenzene-d4	11.816

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
4	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
5	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU100721\
Data File : VU045231.D
Acq On : 08 Oct 2021 00:39
Operator : SY/MD
Sample : M4138-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F3

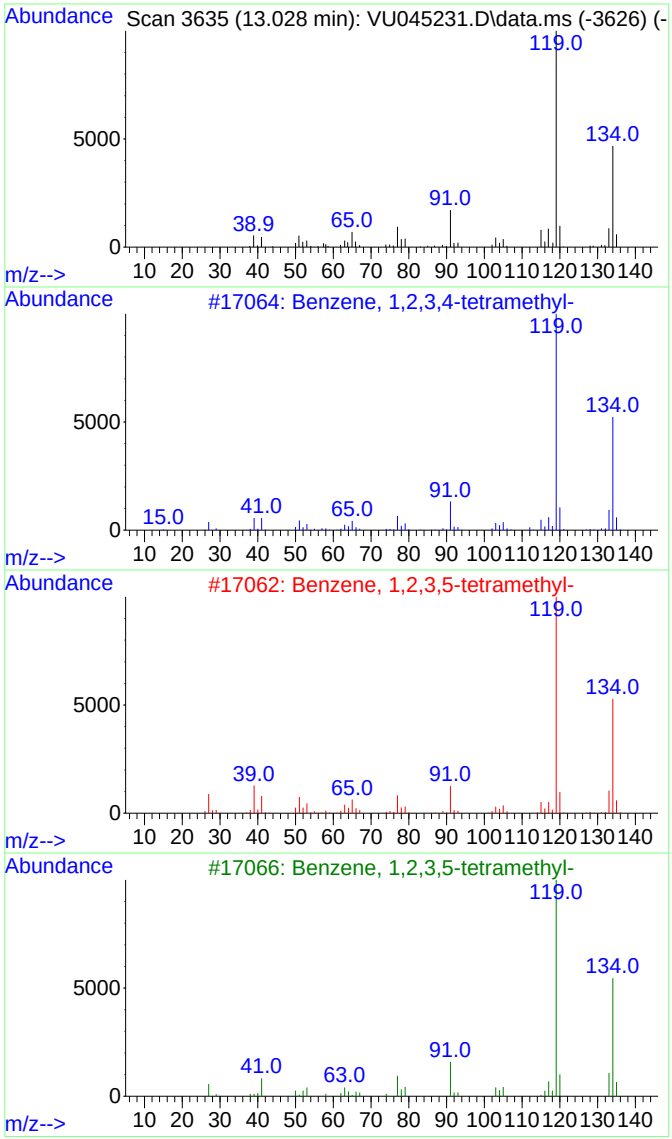
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092321WMA.M
Quant Title : VOC Analysis

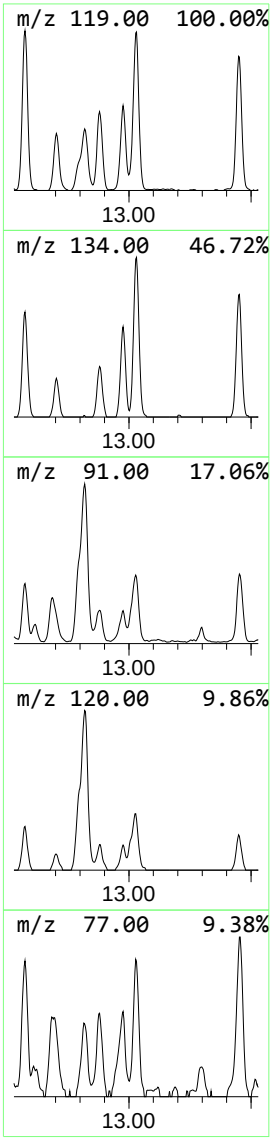
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.028	6.75 ug/L	134354	1,4-Dichlorobenzene-d4	11.816

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
3	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
4	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	95





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU100721\
Data File : VU045231.D
Acq On : 08 Oct 2021 00:39
Operator : SY/MD
Sample : M4138-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092321WMA.M
Quant Title : VOC Analysis

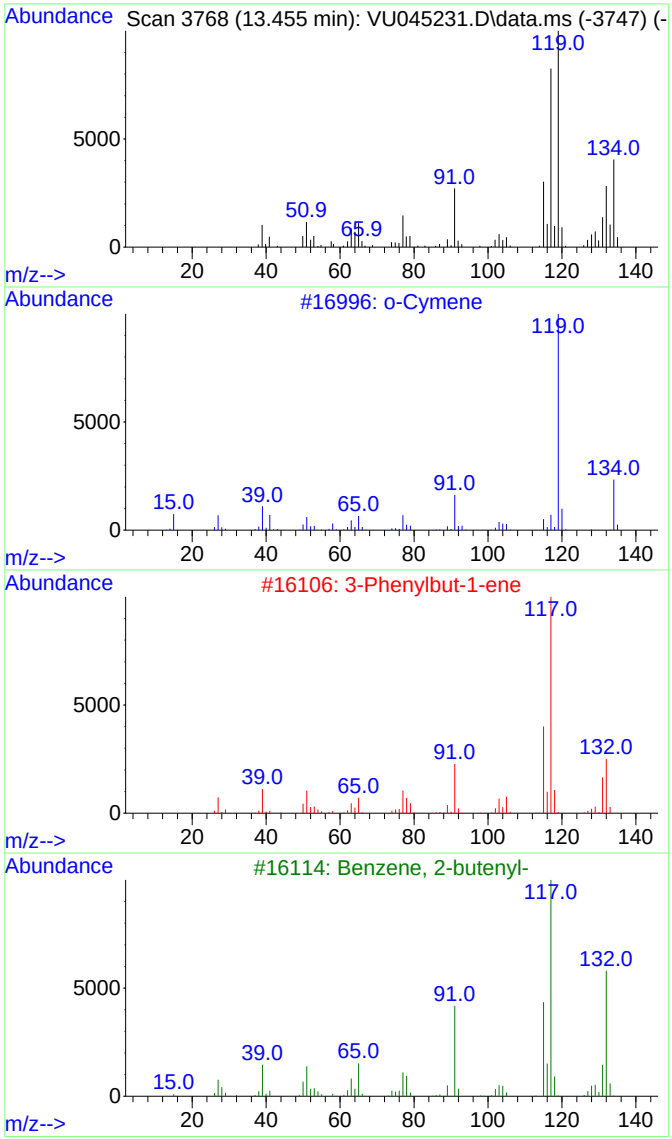
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

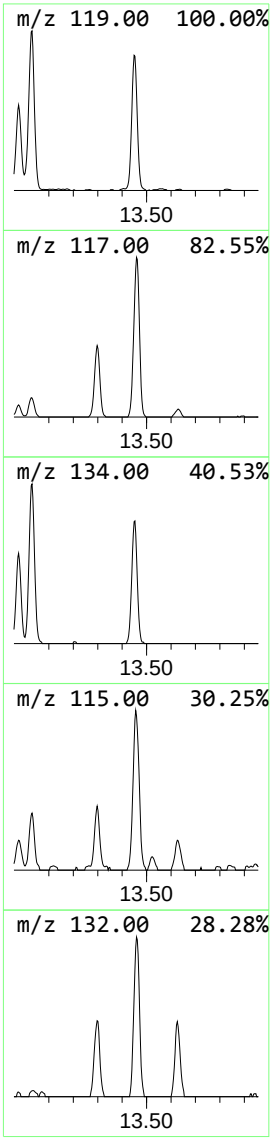
****

Peak Number 25 o-Cymene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.456	9.73 ug/L	193720	1,4-Dichlorobenzene-d4	11.816

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	50
2	3-Phenylbut-1-ene	132	C10H12	000934-10-1	50
3	Benzene, 2-butenyl-	132	C10H12	001560-06-1	50
4	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	50
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	50





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU100721\
Data File : VU045231.D
Acq On : 08 Oct 2021 00:39
Operator : SY/MD
Sample : M4138-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092321WMA.M
Quant Title : VOC Analysis

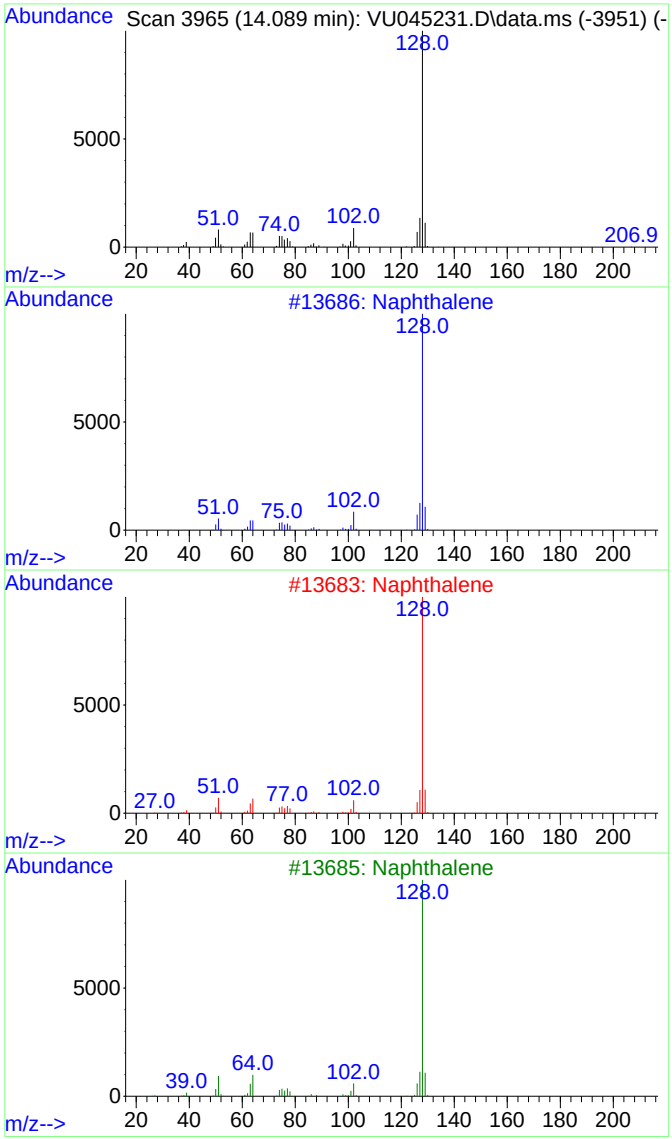
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

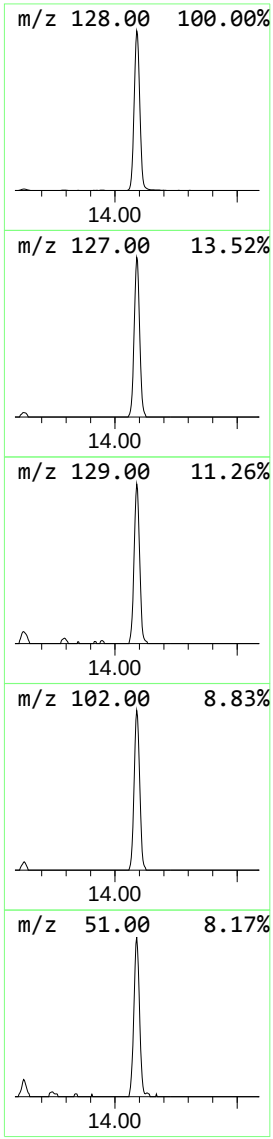
****

Peak Number 26 Azulene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.089	13.84 ug/L	275650	1,4-Dichlorobenzene-d4	11.816

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene	128	C10H8	000091-20-3	95
2	Naphthalene	128	C10H8	000091-20-3	95
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\VU100721\
Data File : VU045231.D
Acq On : 08 Oct 2021 00:39
Operator : SY/MD
Sample : M4138-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092321WMA.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
(DEL) Alkane: S...	1.360	2.6	ug/L	37543	1	6.253	726050	50.0
Sulfur dioxide	1.475	34.6	ug/L	501913	1	6.253	726050	50.0
2-Aminoethyl hy...	1.842	3.8	ug/L	55166	1	6.253	726050	50.0
(DEL) Alkane: S...	3.359	2.9	ug/L	42267	1	6.253	726050	50.0
(DEL) Alkane: S...	3.697	3.0	ug/L	43340	1	6.253	726050	50.0
(DEL) Alkane: C...	4.533	4.8	ug/L	69246	1	6.253	726050	50.0
2-Butanol	4.990	32.2	ug/L	467364	1	6.253	726050	50.0
2-Pentanol, 4-m...	8.234	12.3	ug/L	233088	2	9.424	949299	50.0
Benzene, propyl-	10.909	37.6	ug/L	749452	3	11.816	995806	50.0
Benzene, 1-ethy...	11.002	163.9	ug/L	3264590	3	11.816	995806	50.0
4-Heptanone, 2,...	11.237	4.2	ug/L	82623	3	11.816	995806	50.0
Benzene, 1-ethy...	11.295	53.2	ug/L	1059340	3	11.816	995806	50.0
p-Cymene	11.748	4.4	ug/L	86986	3	11.816	995806	50.0
Benzene, 1,2,3-...	11.899	70.8	ug/L	1410890	3	11.816	995806	50.0
Indane	12.099	32.3	ug/L	642849	3	11.816	995806	50.0
Benzene, (1-met...	12.375	5.0	ug/L	99706	3	11.816	995806	50.0
Cyclohexanone, ...	12.488	24.8	ug/L	493960	3	11.816	995806	50.0
Benzene, 4-ethy...	12.575	6.3	ug/L	125632	3	11.816	995806	50.0
Indan, 1-methyl-	12.687	7.5	ug/L	148949	3	11.816	995806	50.0
Benzaldehyde, 2...	12.819	6.9	ug/L	137359	3	11.816	995806	50.0
Benzene, 2-ethy...	12.880	3.4	ug/L	66963	3	11.816	995806	50.0
Benzene, 1,2,3,...	12.973	6.7	ug/L	132671	3	11.816	995806	50.0
Benzene, 1,2,3,...	13.028	6.8	ug/L	134354	3	11.816	995806	50.0
o-Cymene	13.456	9.7	ug/L	193720	3	11.816	995806	50.0
Azulene	14.089	13.8	ug/L	275650	3	11.816	995806	50.0