LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU100721\
Data File : VU045227.D
Acq On : 07 Oct 2021 23:04
Operator : SY/MD
Sample : M4138-08
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F8

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: VU045227.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.359	3	6	18	rVB3	37141	37407	0.02%	0.011%
2	1.498	33	49	74	rBV2	458795	2226712	1.42%	0.637%
3	1.597	76	80	92	rVB	194232	271557	0.17%	0.078%
4	1.845	150	157	164	rBV2	43851	98596	0.06%	0.028%
5	1.896	166	173	196	rVB	151359	271905	0.17%	0.078%
6	2.469	346	351	367	rVB5	2007	4760	0.00%	0.001%
7	2.562	367	380	395	rBV	256994	439361	0.28%	0.126%
8	2.636	396	403	411	rVV	10349	15269	0.01%	0.004%
9	2.684	411	418	434	rVB2	12219	21228	0.01%	0.006%
10	2.784	434	449	464	rBV2	27344	71423	0.05%	0.020%
11	3.076	536	540	554	rVB2	11156	18938	0.01%	0.005%
12	3.198	571	578	594	rVB3	5579	11736	0.01%	0.003%
13	3.359	616	628	645	rVB	18124	41838	0.03%	0.012%
14	3.697	720	733	751	rVB3	17293	42936	0.03%	0.012%
15	3.870	770	787	809	rBV	117536	279242	0.18%	0.080%
16	4.533	976	993	1009	rBV2	26608	68428	0.04%	0.020%
17	4.629	1009	1023	1038	rVV	146808	353427	0.23%	0.101%
18	4.710	1038	1048	1071	rVB	78819	183308	0.12%	0.052%
19	4.989	1120	1135	1147	rBV	191147	436991	0.28%	0.125%
20	5.067	1148	1159	1189	rVB	232609	525601	0.34%	0.150%
21	5.388	1242	1259	1274	rBV2	30989	71916	0.05%	0.021%
22	5.462	1274	1282	1294	rVV6	3045	6746	0.00%	0.002%
23	5.591	1313	1322	1325	rBV6	3610	5187	0.00%	0.001%
24	5.639	1331	1337	1344	rVB5	3084	4350	0.00%	0.001%
25	5.729	1344	1365	1374	rVV2	486834	1309722	0.84%	0.375%
26	5.777	1375	1380	1394	rVV	187689	332668	0.21%	0.095%
27	5.851	1395	1403	1415	rVV4	9179	18606	0.01%	0.005%
28	5.925	1415	1426	1436	rVV5	8339	18109	0.01%	0.005%
29	5.993	1436	1447	1464	rVB4	14685	33133	0.02%	0.009%
30	6.131	1480	1490	1500	rBV6	2151	4465	0.00%	0.001%
31	6.253	1511	1528	1552	rBV	386424	731201	0.47%	0.209%
32	6.533	1600	1615	1622	rBV3	8748	19079	0.01%	0.005%
33	6.697	1650	1666	1677	rVV	307330	605081	0.39%	0.173%
34	6.764	1677	1687	1709	rVV	124690	264088	0.17%	0.076%
35	6.983	1741	1755	1768	rVV4	11736	22766	0.01%	0.007%
36	7.121	1790	1798	1809	rVV2	3488	7228	0.00%	0.002%

LSC Area Percent Report

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Data File : VU045227.D
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Operator : SY/MD
Sample : M4138-08
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F8

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	7.182	1812	1817	1827	rVV5	2246	3765	0.00%	0.001%
38	7.243	1827	1836	1845	rVB10	2858	5500	0.00%	0.002%
39	7.568	1924	1937	1954	rBV	170377	302365	0.19%	0.086%
40	7.697	1967	1977	1983	rBV7	1306	2379	0.00%	0.001%
41	7.800	1992	2009	2021	rBV	125982	227057	0.14%	0.065%
42	7.902	2028	2041	2050	rBV	615318	1049819	0.67%	0.300%
43	7.976	2050	2064	2098	rVB	15440519	26110460	16.67%	7.467%
44	8.182	2118	2128	2136	rBV	116933	186277	0.12%	0.053%
45	8.234	2136	2144	2173	rVB	103172	220557	0.14%	0.063%
46	8.372	2180	2187	2196	rVB5	4885	7863	0.01%	0.002%
47	8.420	2197	2202	2211	rVB8	1961	3275	0.00%	0.001%
48	8.639	2251	2270	2301	rBV	471462	834985	0.53%	0.239%
49	8.812	2311	2324	2332	rVB8	4746	9714	0.01%	0.003%
50	8.870	2332	2342	2349	rBV2	5552	9392	0.01%	0.003%
51	9.217	2437	2450	2459	rBV5	5226	9532	0.01%	0.003%
52	9.333	2478	2486	2492	rBV4	2385	3894	0.00%	0.001%
53	9.378	2492	2500	2502	rBV3	6562	8038	0.01%	0.002%
54	9.423	2502	2514	2528	rVV	574822	957754	0.61%	0.274%
55	9.497	2529	2537	2547	rVV6	26359	61058	0.04%	0.017%
56	9.584	2547	2564	2586	rVV2	39991605	71294341	45.51%	20.389%
57	9.726	2587	2608	2643	rVB3	59407104	156669534	100.00%	44.804%
58	9.954	2672	2679	2687	rVB5	6167	8977	0.01%	0.003%
59	10.021	2691	2700	2710	rVB5	6746	12274	0.01%	0.004%
60	10.115	2710	2729	2761	rBV2	38336832	64350270	41.07%	18.403%
61	10.237	2762	2767	2772	rVV6	3473	5030	0.00%	0.001%
62	10.275	2775	2779	2792	rVB9	5173	8032	0.01%	0.002%
63	10.365	2795	2807	2821	rBV4	16481	31294	0.02%	0.009%
64	10.488	2832	2845	2864	rBV	1084839	1677276	1.07%	0.480%
65	10.568	2865	2870	2875	rVB4	1917	2623	0.00%	0.001%
66	10.655	2886	2897	2899	rBV7	5612	7433	0.00%	0.002%
67	10.761	2917	2930	2946	rBV	452333	725953	0.46%	0.208%
68	10.909	2964	2976	2990	rBV	468224	720585	0.46%	0.206%
69	11.002	2991	3005	3023	rVV	1485371	3128312	2.00%	0.895%
70	11.092	3023	3033	3047	rVB	742473	1130439	0.72%	0.323%
71	11.237	3064	3078	3085	rBV	48471	80200	0.05%	0.023%
72	11.295	3085	3096	3120	rVV	673965	1056055	0.67%	0.302%
73	11.410	3122	3132	3139	rVV7	8186	18405	0.01%	0.005%
74	11.471	3139	3151	3174	rVB	2388932	3660649	2.34%	1.047%
75	11.645	3188	3205	3216	rVB5	21337	42387	0.03%	0.012%
76	11.745	3224	3236	3245	rBV2	49734	81625	0.05%	0.023%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU100721\
Data File : VU045227.D
Acq On : 07 Oct 2021 23:04
Operator : SY/MD
Sample : M4138-08
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F8

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	11.815	3245	3258	3271	rVB	619900	994052	0.63%	0.284%
78	11.899	3271	3284	3297	rBV	877184	1353592	0.86%	0.387%
79	12.018	3312	3321	3329	rBV4	21805	36424	0.02%	0.010%
80	12.098	3332	3346	3364	rVV2	353339	618756	0.39%	0.177%
81	12.195	3364	3376	3395	rVB	683342	1113321	0.71%	0.318%
82	12.304	3400	3410	3418	rBV3	12617	20509	0.01%	0.006%
83	12.375	3418	3432	3450	rVB4	45266	97938	0.06%	0.028%
84	12.487	3450	3467	3486	rBV3	182677	483853	0.31%	0.138%
85	12.574	3486	3494	3502	rVV	75920	121211	0.08%	0.035%
86	12.616	3502	3507	3518	rVV	24661	38993	0.02%	0.011%
87	12.687	3518	3529	3547	rVB3	69934	140433	0.09%	0.040%
88	12.819	3553	3570	3580	rBV	67108	149464	0.10%	0.043%
89	12.880	3580	3589	3601	rVB3	39707	65184	0.04%	0.019%
90	12.973	3601	3618	3626	rBV2	49229	128980	0.08%	0.037%
91	13.028	3626	3635	3648	rVB2	79836	130255	0.08%	0.037%
92	13.256	3696	3706	3709	rBV6	1516	2213	0.00%	0.001%
93	13.298	3709	3719	3740	rVB	25487	41342	0.03%	0.012%
94	13.455	3745	3768	3779	rBV3	101128	183194	0.12%	0.052%
95	13.520	3780	3788	3798	rVB3	10412	16880	0.01%	0.005%
96	13.629	3811	3822	3834	rVB2	25416	40377	0.03%	0.012%
97	13.741	3847	3857	3868	rVV2	17396	29855	0.02%	0.009%
98	13.790	3868	3872	3878	rVB5	1975	2126	0.00%	0.001%
99	13.841	3880	3888	3901	rVB9	3474	7253	0.00%	0.002%
100	14.089	3953	3965	3982	rVV	161441	260874	0.17%	0.075%

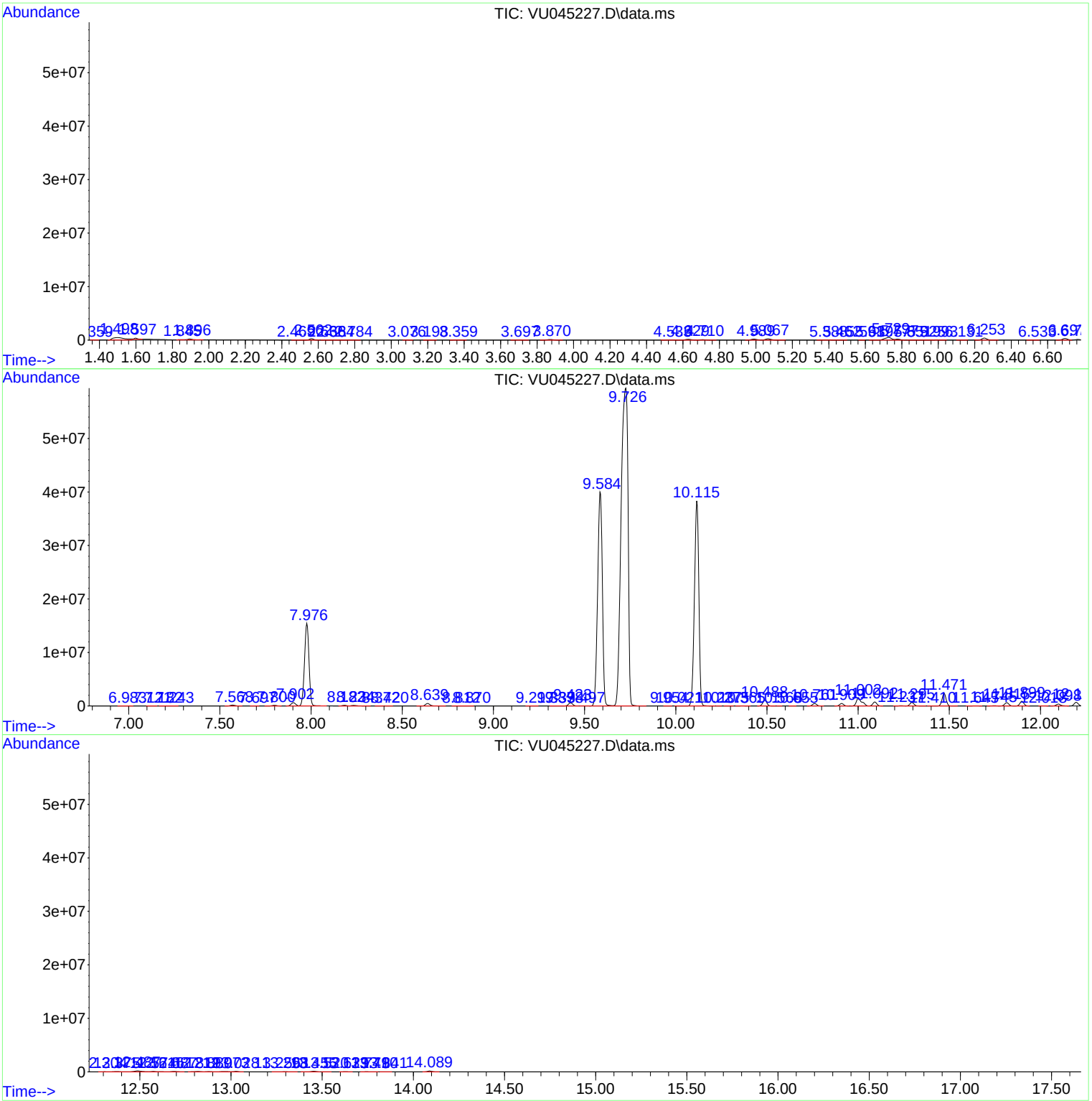
Sum of corrected areas: 349677465

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU100721\
Data File : VU045227.D
Acq On : 07 Oct 2021 23:04
Operator : SY/MD
Sample : M4138-08
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW7F8

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 : VU045227.D
Acq On : 07 Oct 2021 23:04
Operator : SY/MD
Sample : M4138-08
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F8

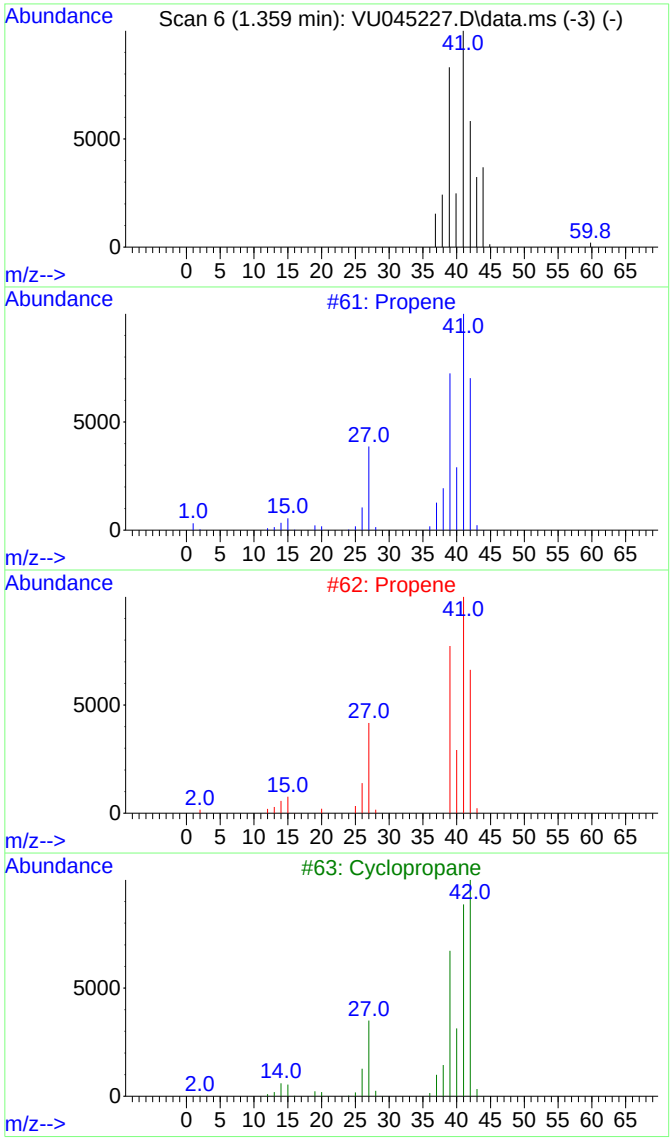
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092321WMA.M
Quant Title : VOC Analysis

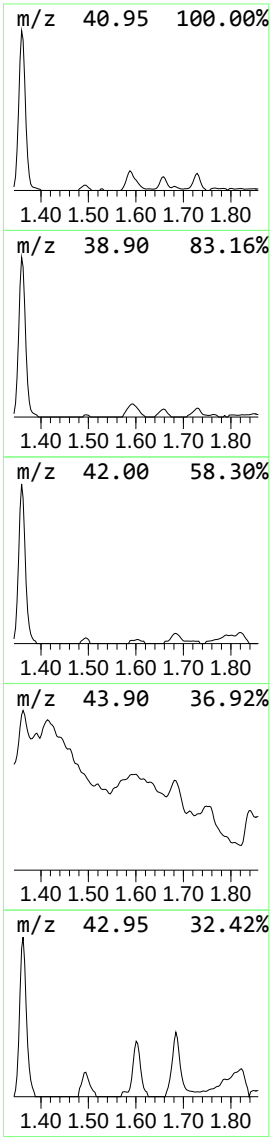
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.359	2.56 ug/L	37407	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	Cyclopropane	42	C3H6	000075-19-4	9
5	Cyclopropene	40	C3H4	002781-85-3	4





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU100721\
Data File : VU045227.D
Acq On : 07 Oct 2021 23:04
Operator : SY/MD
Sample : M4138-08
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F8

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

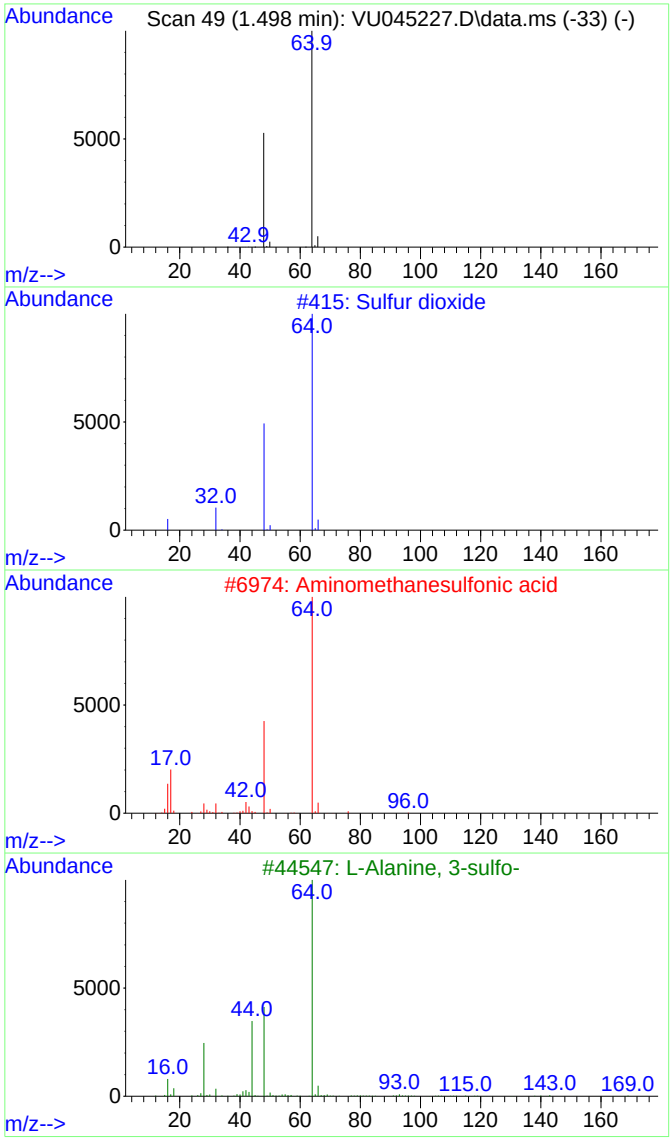
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

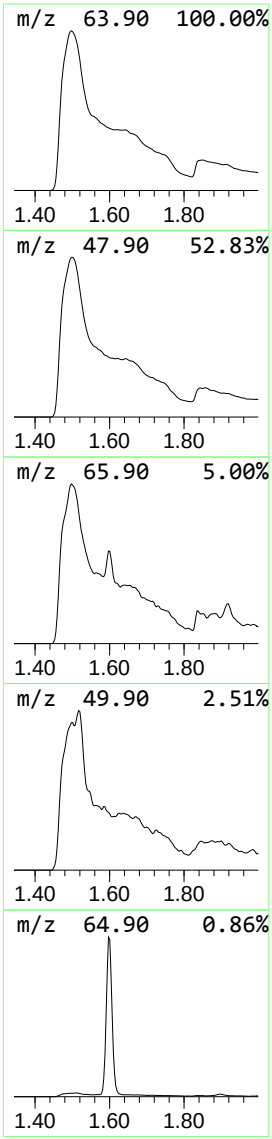
****

Peak Number 2 Sulfur dioxide Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.498	152.26 ug/L	2226710	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	Ethene, 1,1-difluoro-	64	C2H2F2	000075-38-7	4





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU100721\
Data File : VU045227.D
Acq On : 07 Oct 2021 23:04
Operator : SY/MD
Sample : M4138-08
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F8

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

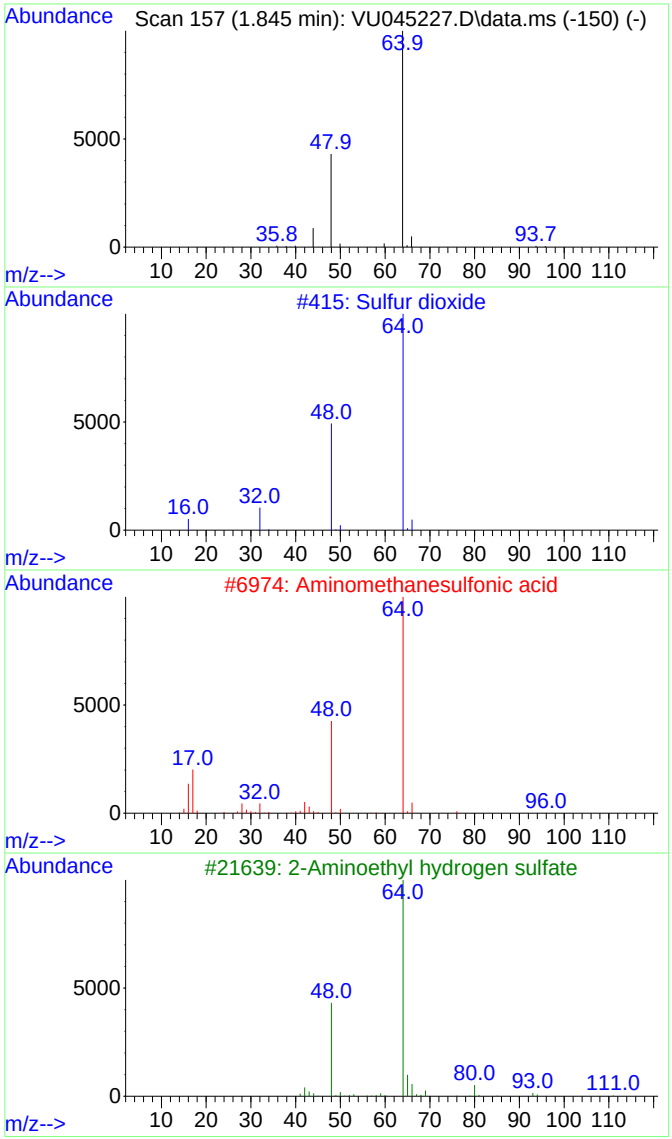
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

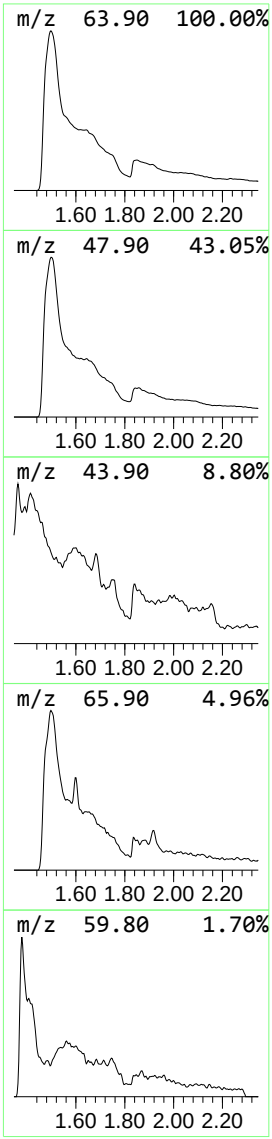
****

Peak Number 3 Aminomethanesulfonic acid Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.845	6.74 ug/L	98596	1,4-Difluorobenzene	6.253

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Sulfur dioxide	64	O2S	007446-09-5	90
2	Aminomethanesulfonic acid	111	CH5NO3S	013881-91-9	64
3	2-Aminoethyl hydrogen sulfate	141	C2H7NO4S	000926-39-6	64
4	L-Alanine, 3-sulfo-	169	C3H7NO5S	000498-40-8	64
5	(N-(-2-Acetamido))-2-aminoethane...	182	C4H10N2O4S	007365-82-4	64





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU100721\
Data File : VU045227.D
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Operator : SY/MD
Sample : M4138-08
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F8

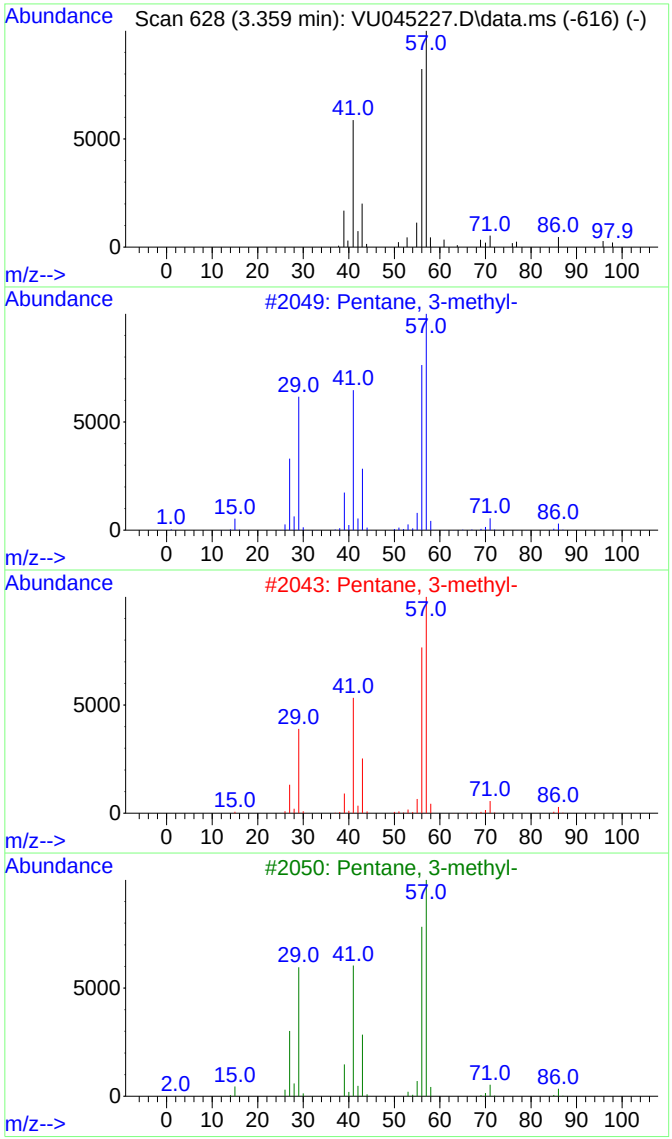
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092321WMA.M
Quant Title : VOC Analysis

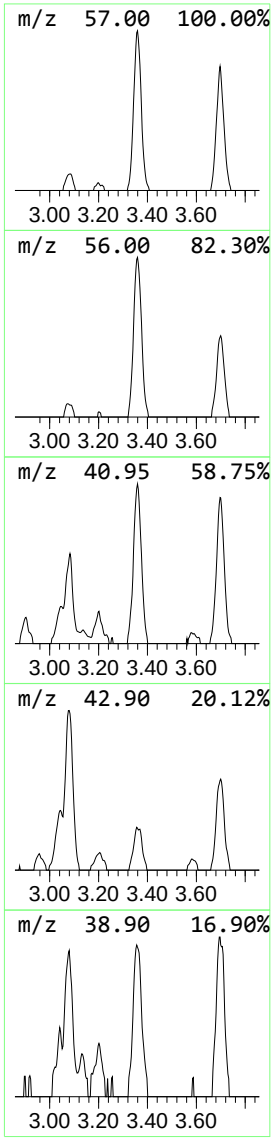
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.359	2.86 ug/L	41838	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	Sulphuric acid dibutyl ester	210	C8H18O4S	000625-22-9	50
5	Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	50





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU100721\
Data File : VU045227.D
Acq On : 07 Oct 2021 23:04
Operator : SY/MD
Sample : M4138-08
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW7F8

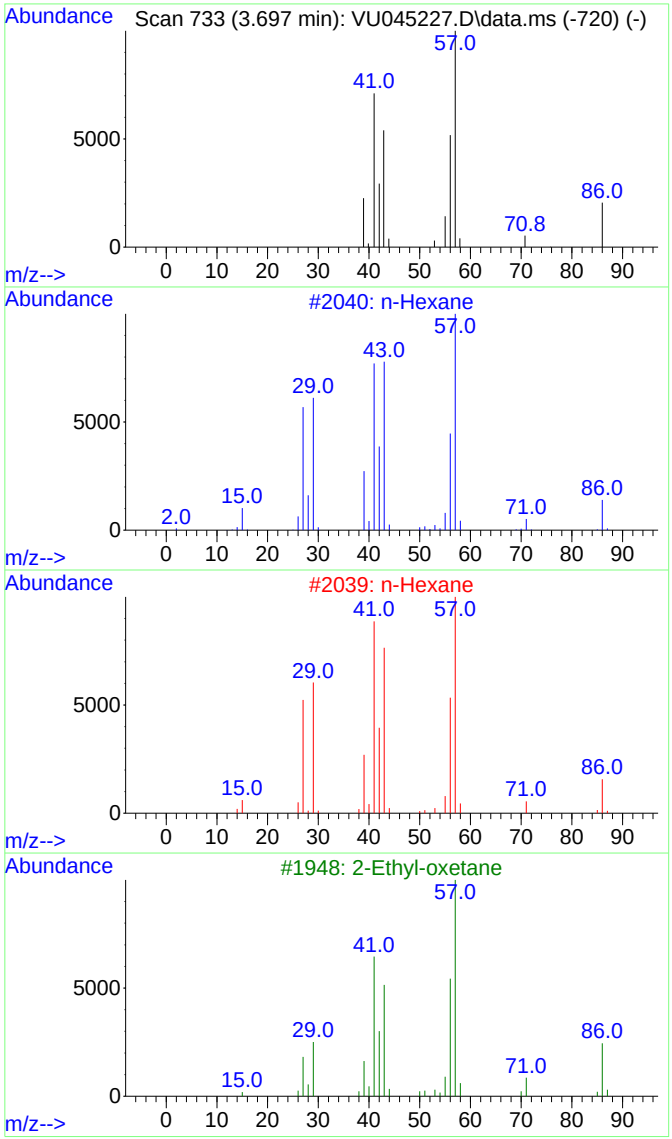
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092321WMA.M
Quant Title : VOC Analysis

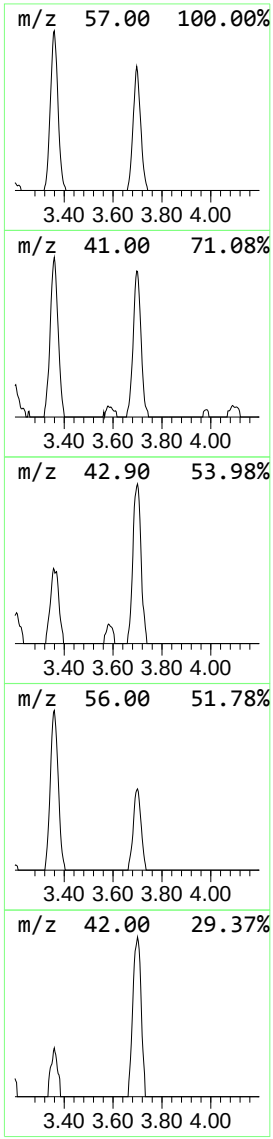
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.697	2.94 ug/L	42936	1,4-Difluorobenzene	6.253

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexane	86	C6H14	000110-54-3	72
2	n-Hexane	86	C6H14	000110-54-3	72
3	2-Ethyl-oxetane	86	C5H10O	1010386-40-2	72
4	n-Hexane	86	C6H14	000110-54-3	56
5	n-Hexane	86	C6H14	000110-54-3	50





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Data File : VU045227.D
Acq On : 07 Oct 2021 23:04
Operator : SY/MD
Sample : M4138-08
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F8

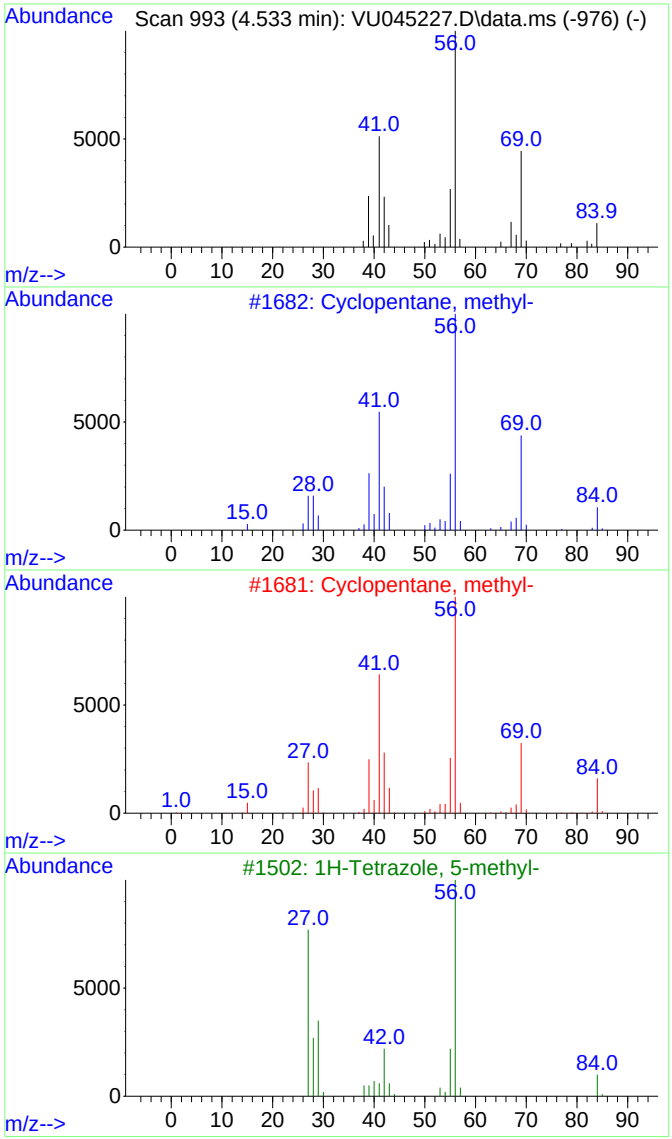
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092321WMA.M
Quant Title : VOC Analysis

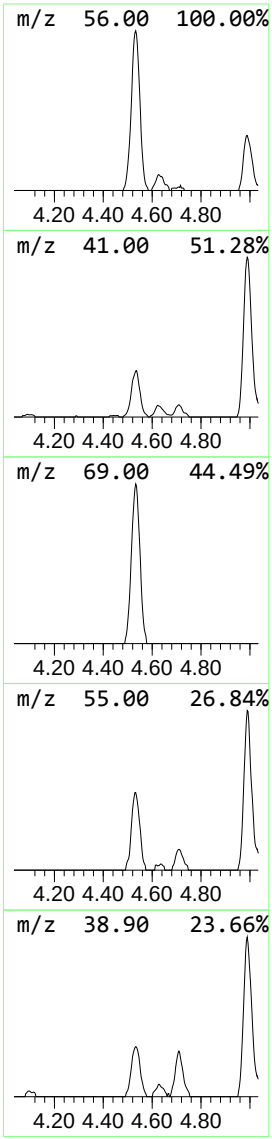
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 (DEL) Alkane: Cyclic4.533 Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.533	4.68 ug/L	68428	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	86
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 : 07 Oct 2021 23:04
Operator : SY/MD
Sample : M4138-08
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F8

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

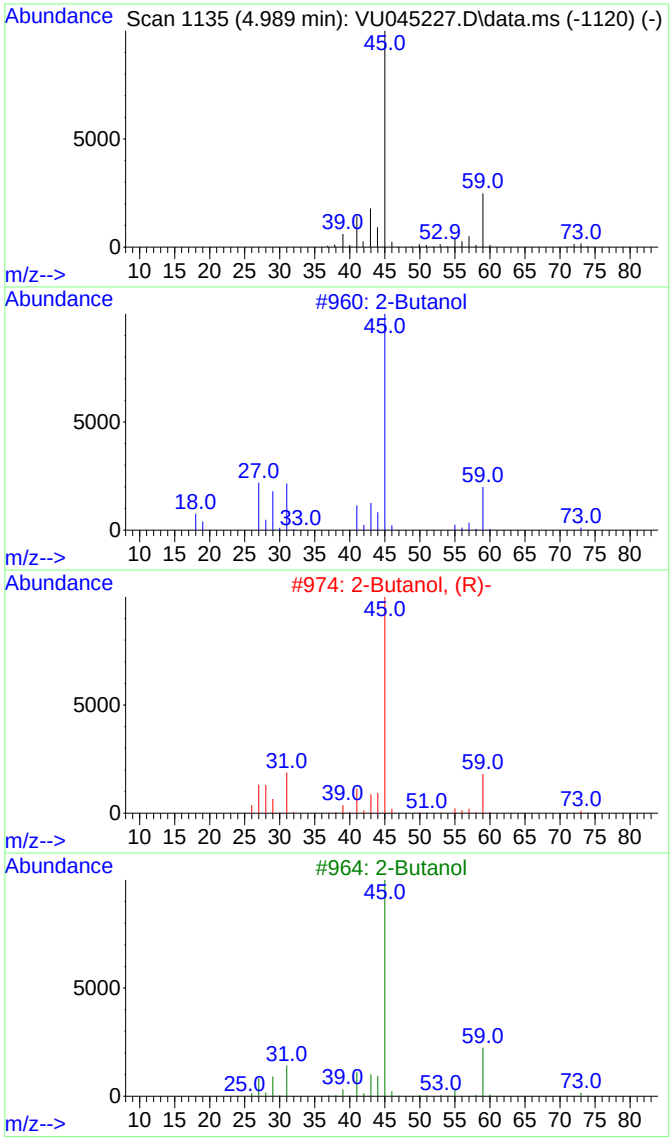
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

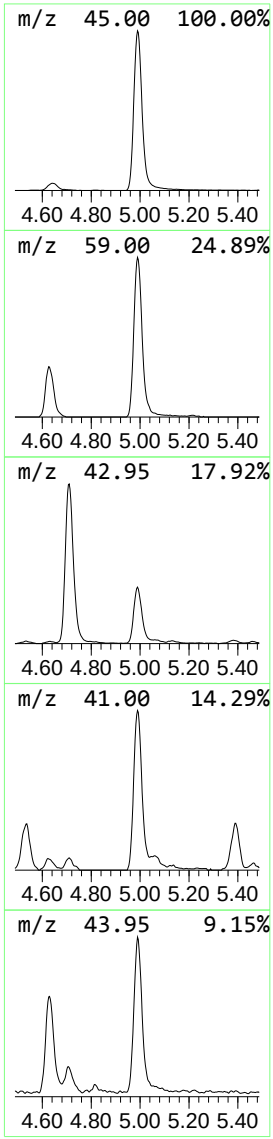
****

Peak Number 7 2-Butanol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.989	29.88 ug/L	436991	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 : 07 Oct 2021 23:04
Operator : SY/MD
Sample : M4138-08
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW7F8

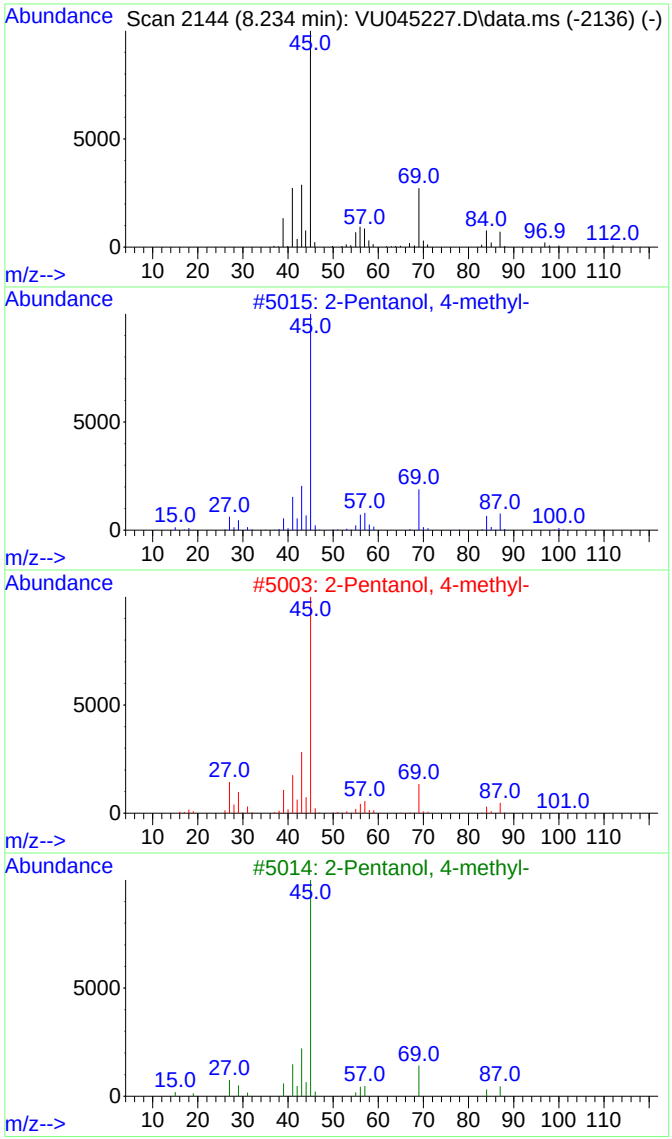
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092321WMA.M
Quant Title : VOC Analysis

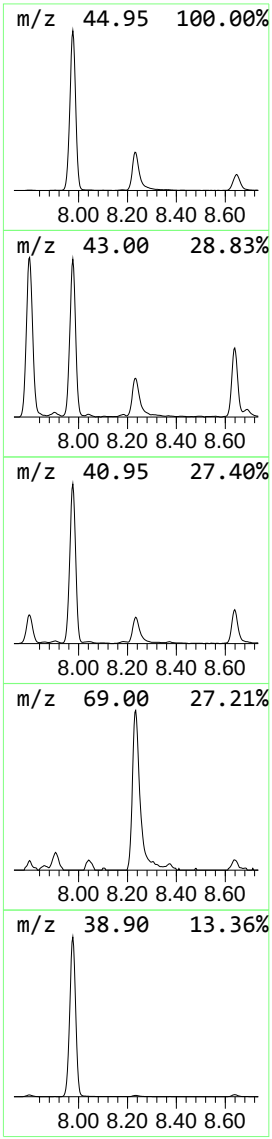
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.234	11.51 ug/L	220557	Chlorobenzene-d5	9.423

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	83
2	2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	83
3	2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	78
4	2-Hexanol	102	C6H14O	000626-93-7	74
5	2-Hexanol	102	C6H14O	000626-93-7	72





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Acq On : 07 Oct 2021 23:04
Operator : SY/MD
Sample : M4138-08
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW7F8

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

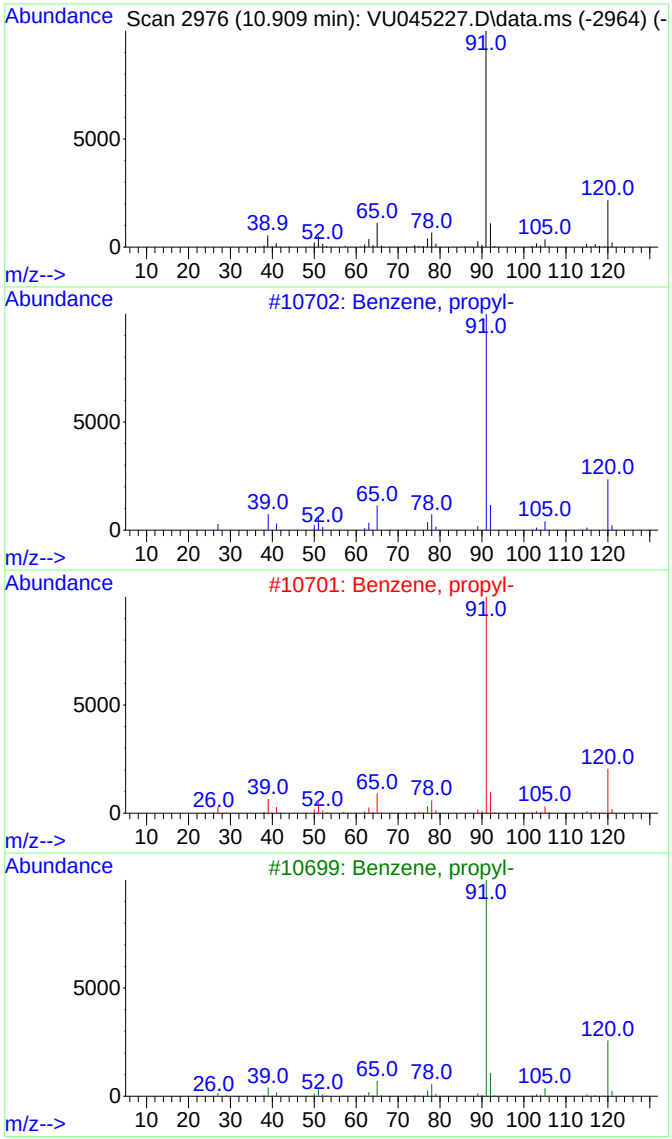
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

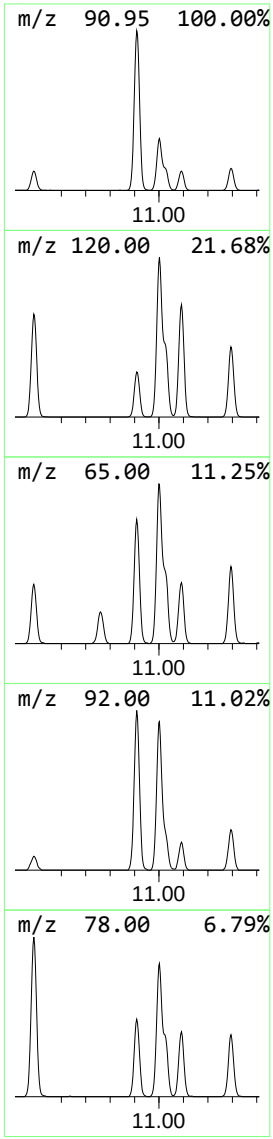
****

Peak Number 10 Benzene, propyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.909	36.24 ug/L	720585	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	86
4	Phenethylamine, N-benzyl-p-chloro-	245	C15H16ClN	013622-43-0	72
5	1,2-Ethanediamine, N,N'-bis(phen...	240	C16H20N2	000140-28-3	64





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Operator : SY/MD
Sample : M4138-08
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F8

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

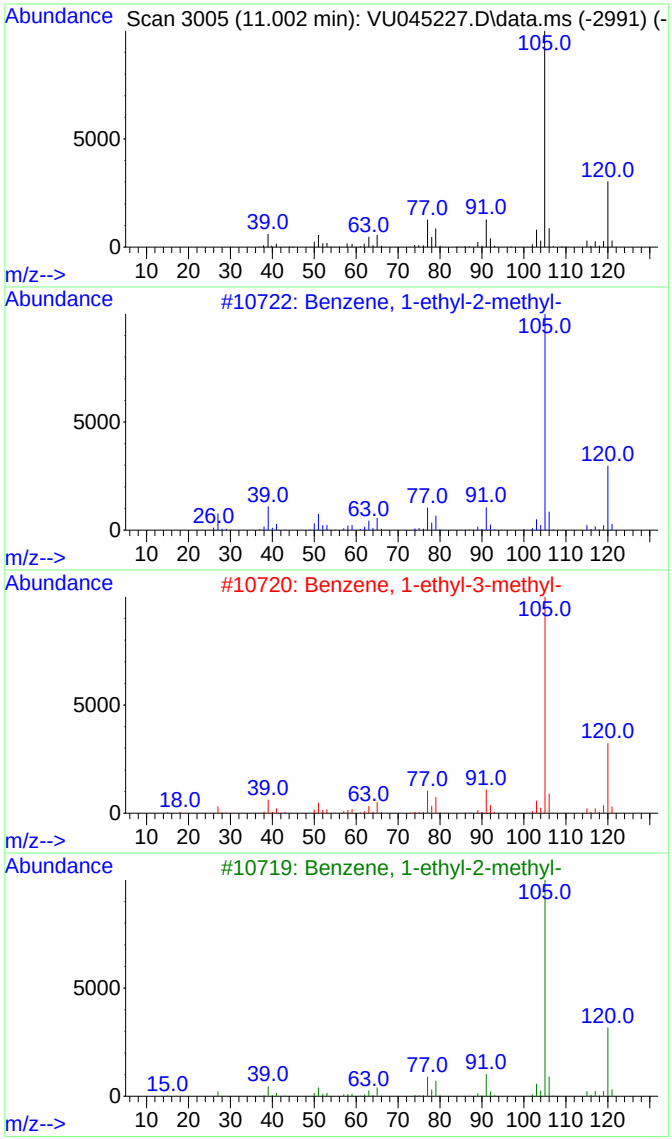
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

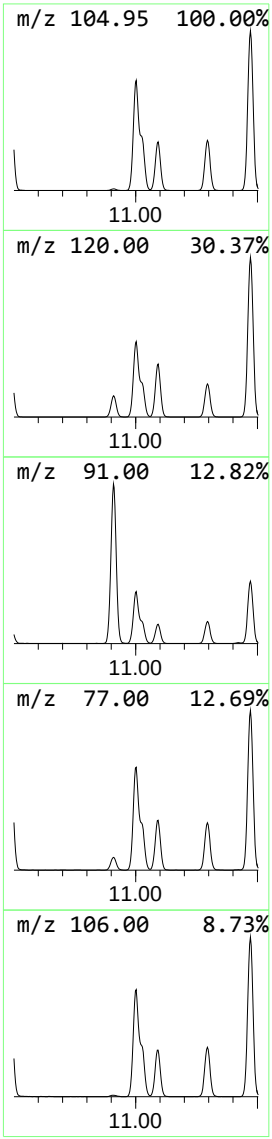
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.002	157.35 ug/L	3128310	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	Mesitylene	120	C9H12	000108-67-8	91





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Operator : SY/MD
Sample : M4138-08
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F8

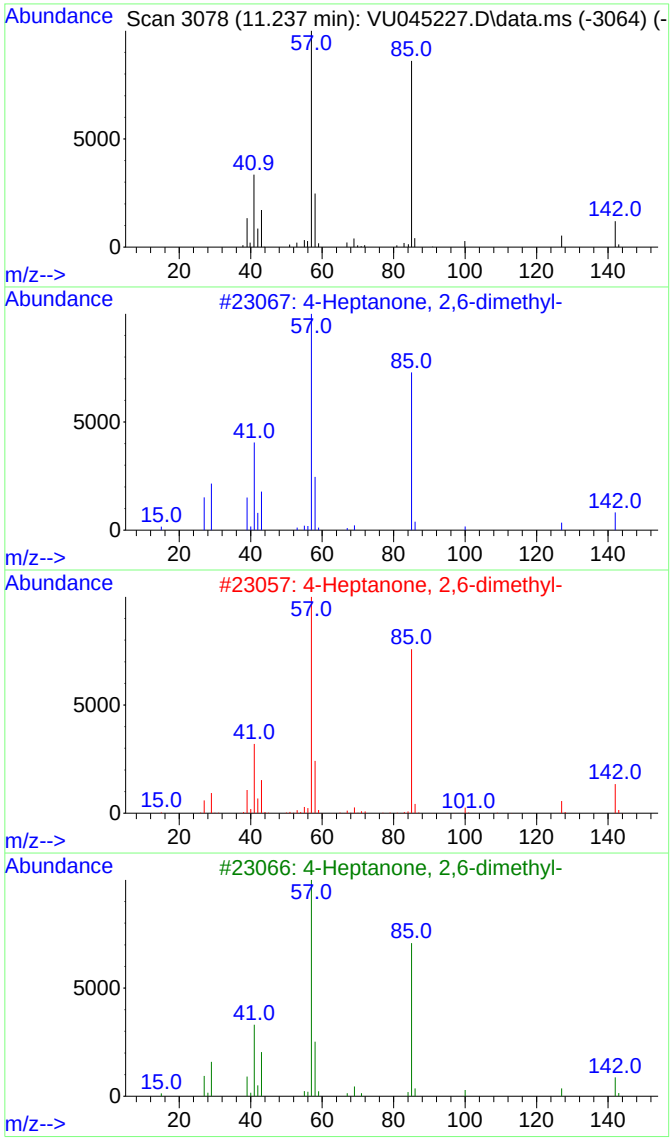
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092321WMA.M
Quant Title : VOC Analysis

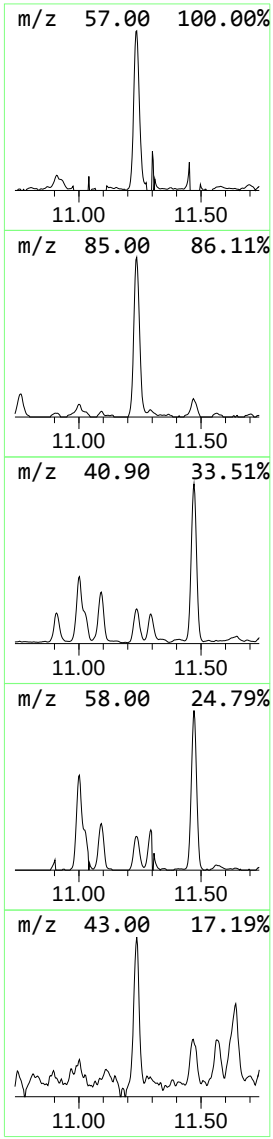
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 4-Heptanone, 2,6-dimethyl- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.237	4.03 ug/L	80200	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	94
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	64





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Acq On : 07 Oct 2021 23:04
Operator : SY/MD
Sample : M4138-08
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F8

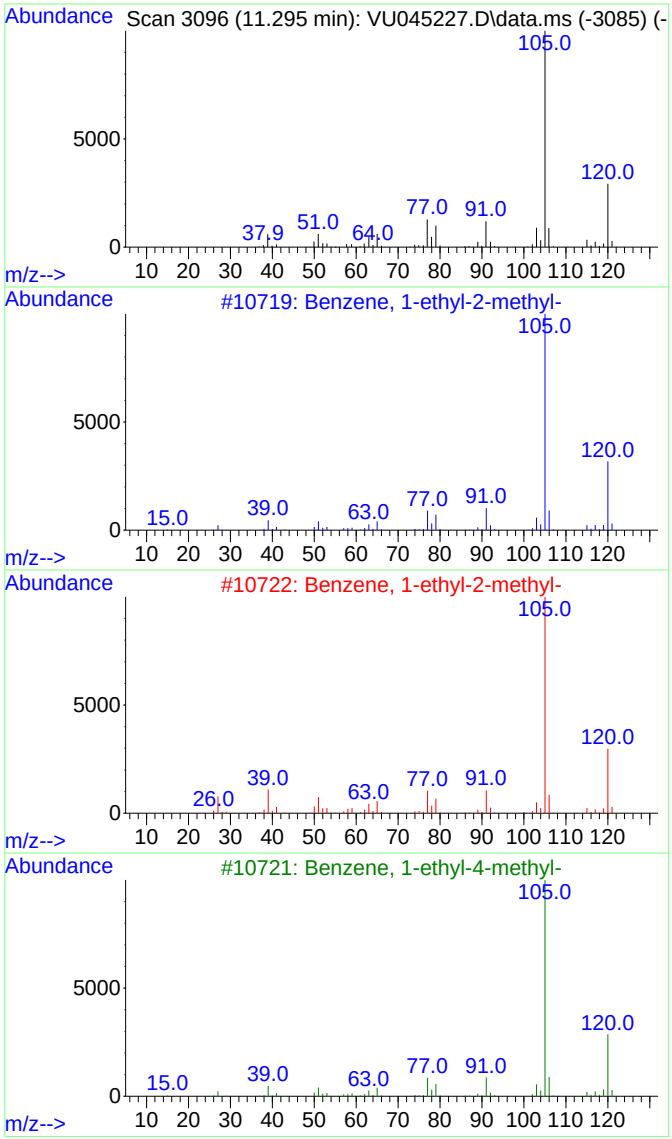
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092321WMA.M
Quant Title : VOC Analysis

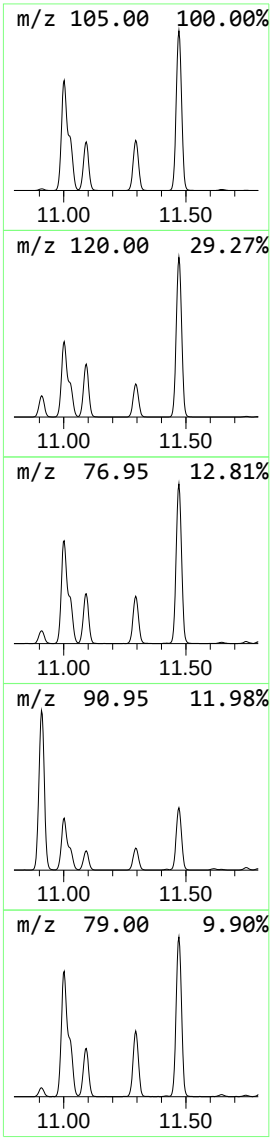
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.295	53.12 ug/L	1056060	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	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91





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Operator : SY/MD
Sample : M4138-08
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F8

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

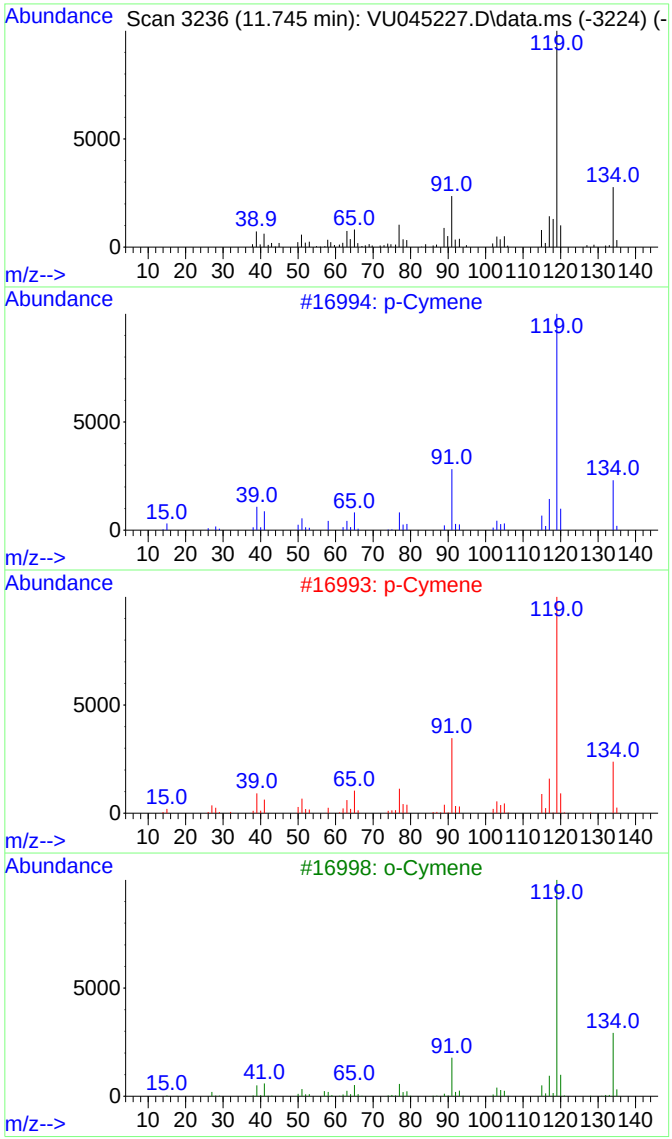
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

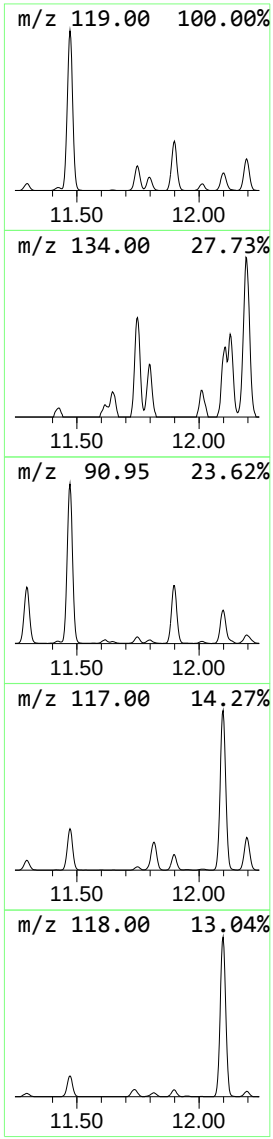
****

Peak Number 14 p-Cymene Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.745	4.11 ug/L	81625	1,4-Dichlorobenzene-d4	11.816

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Cymene	134	C10H14	000099-87-6	94
2	p-Cymene	134	C10H14	000099-87-6	94
3	o-Cymene	134	C10H14	000527-84-4	93
4	o-Cymene	134	C10H14	000527-84-4	93
5	o-Cymene	134	C10H14	000527-84-4	93





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Operator : SY/MD
Sample : M4138-08
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F8

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

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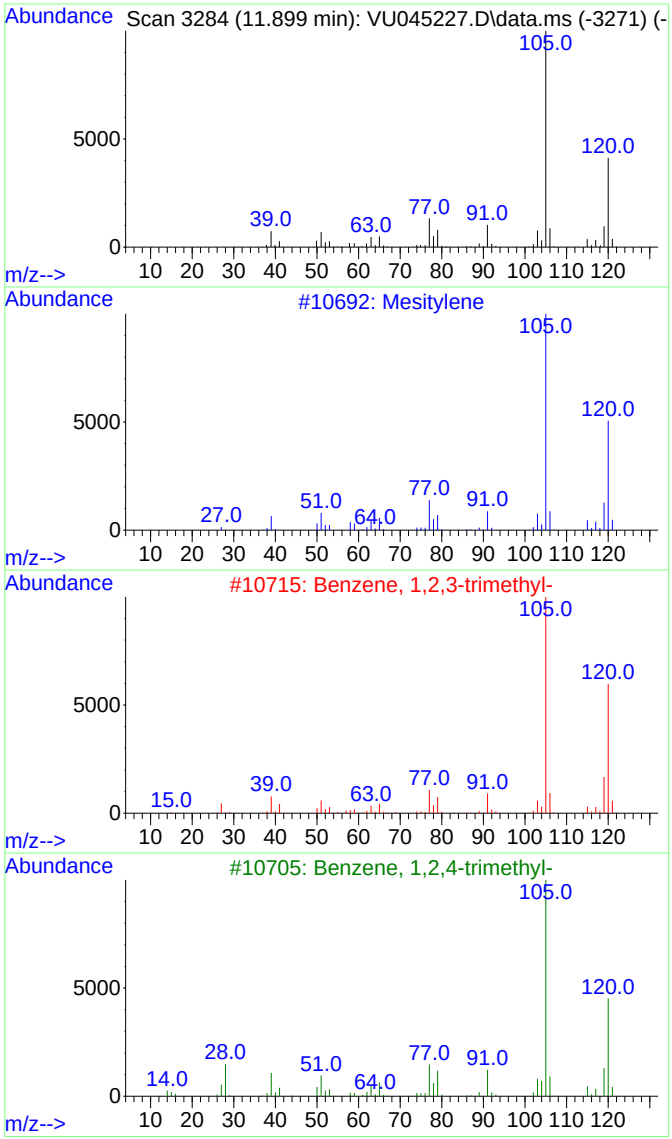
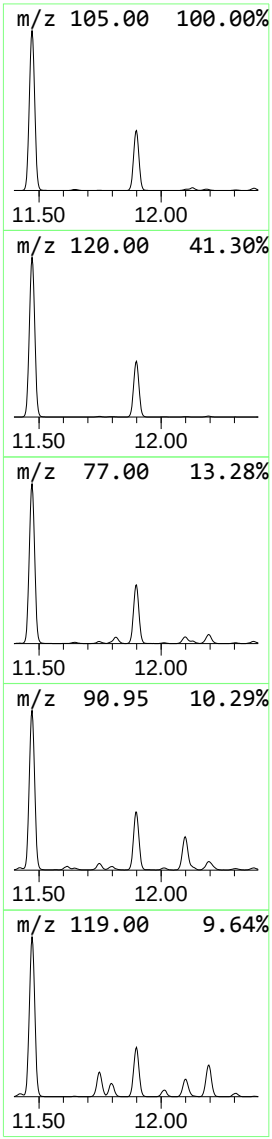
Peak Number 15 Benzene, 1,2,3-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.899	68.08 ug/L	1353590	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-ethyl-2-methyl-	120	C9H12	000611-14-3	93
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91





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Sample : M4138-08
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F8

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

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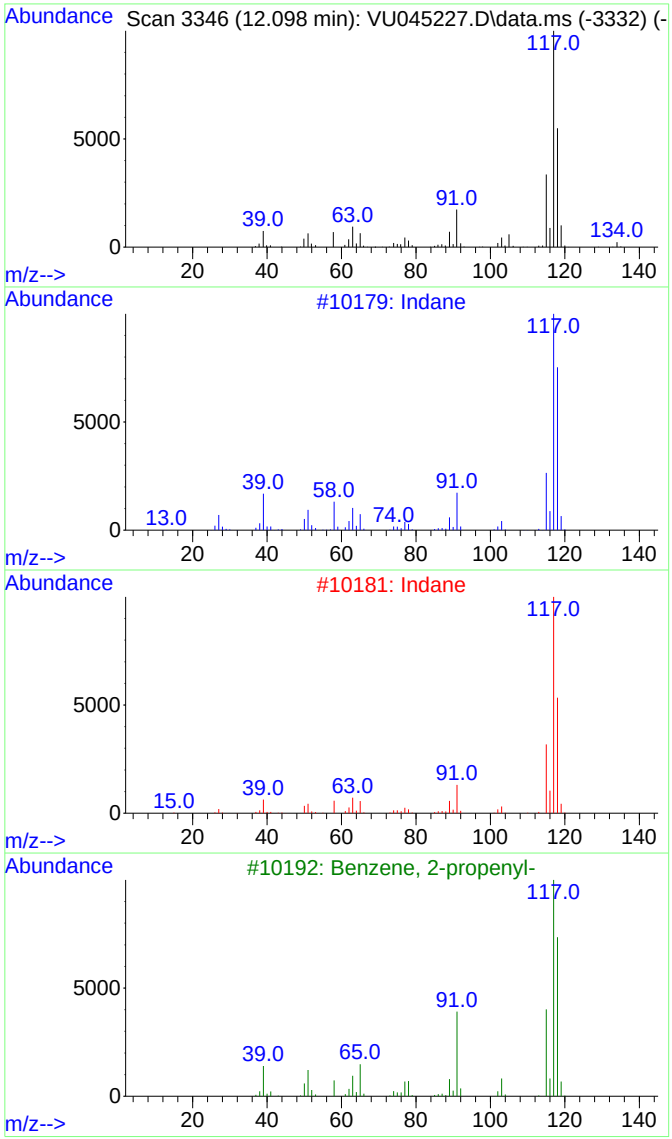
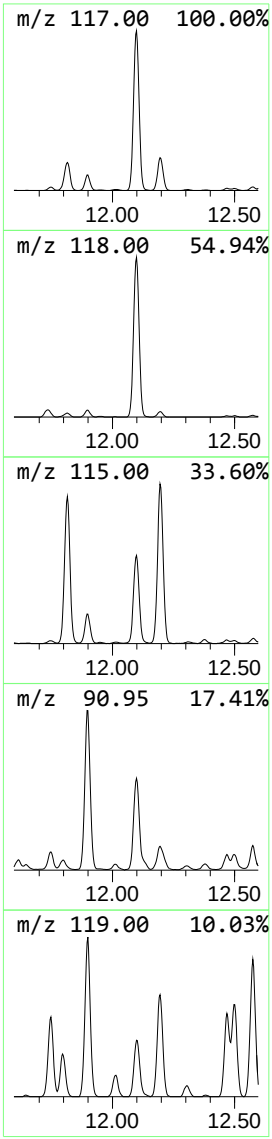
Peak Number 16 Indane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.098	31.12 ug/L	618756	1,4-Dichlorobenzene-d4	11.816

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	87
2	Indane	118	C9H10	000496-11-7	87
3	Benzene, 2-propenyl-	118	C9H10	000300-57-2	81
4	Benzene, cyclopropyl-	118	C9H10	000873-49-4	81
5	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	76





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Sample : M4138-08
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW7F8

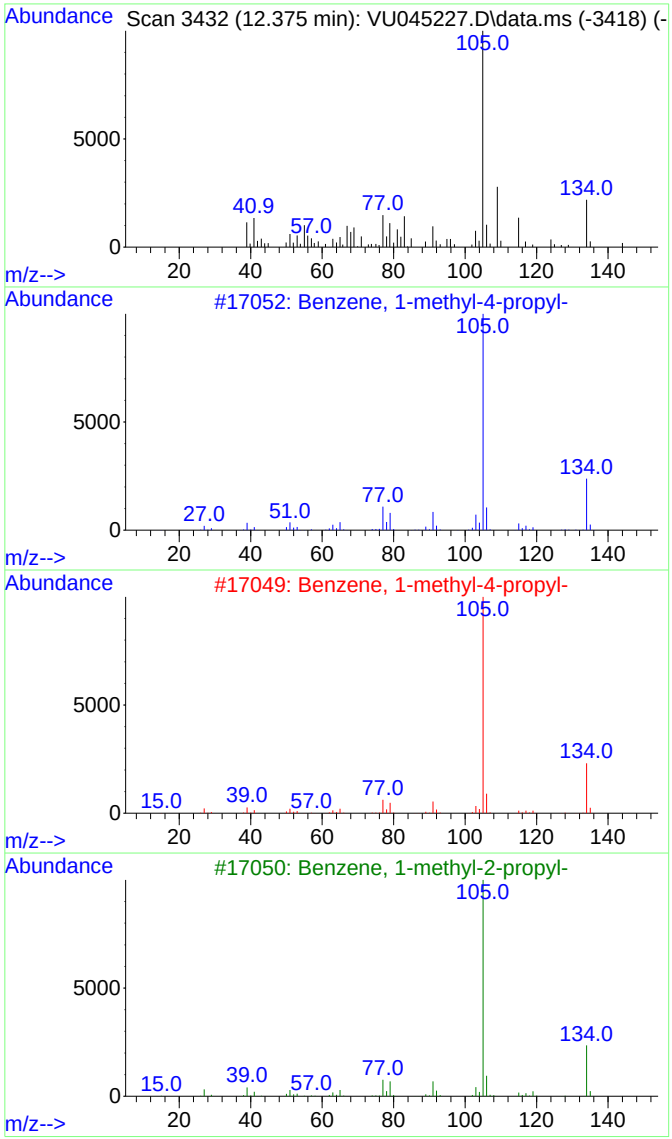
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092321WMA.M
Quant Title : VOC Analysis

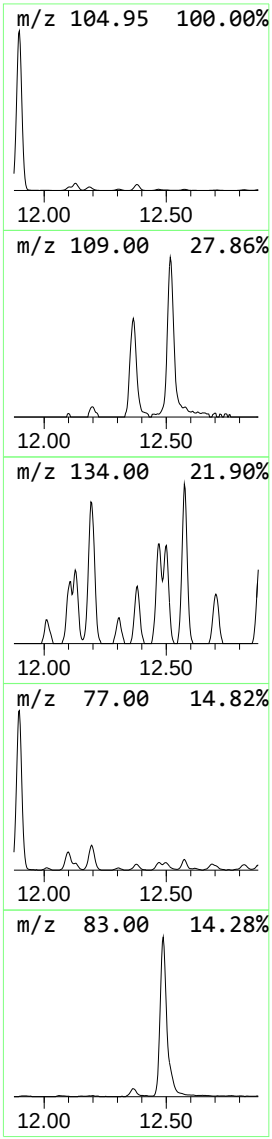
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 Benzene, 1-methyl-4-propyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.375	4.93 ug/L	97938	1,4-Dichlorobenzene-d4	11.816

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	90
2	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	60
3	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	60
4	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	60
5	Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	60





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Data File : VU045227.D
Acq On : 07 Oct 2021 23:04
Operator : SY/MD
Sample : M4138-08
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F8

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

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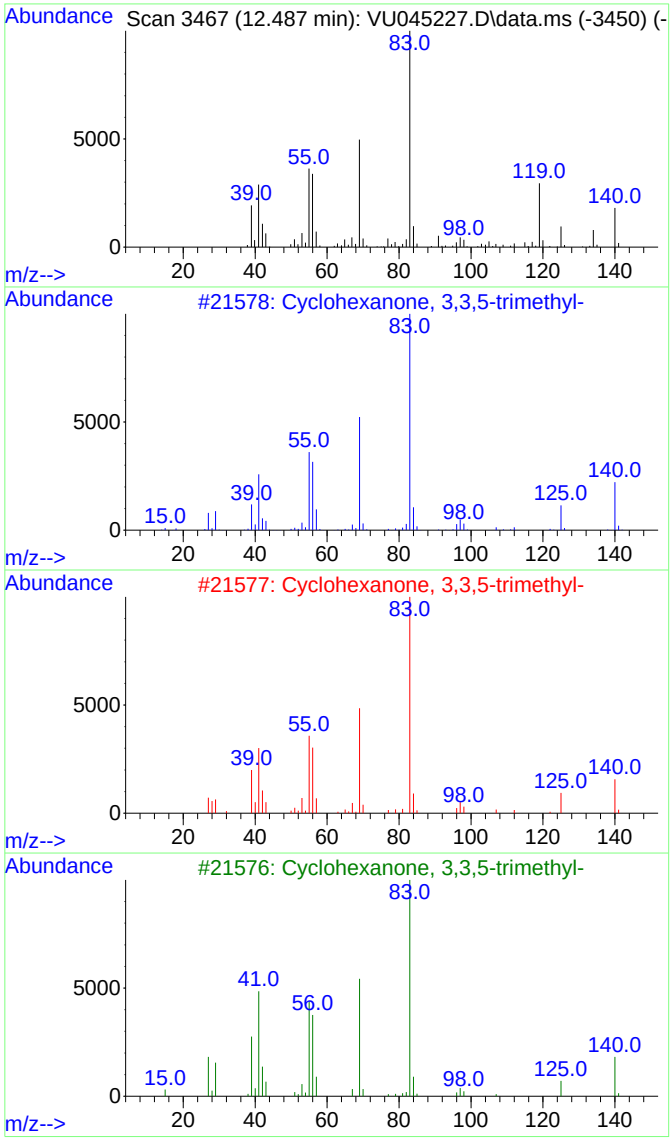
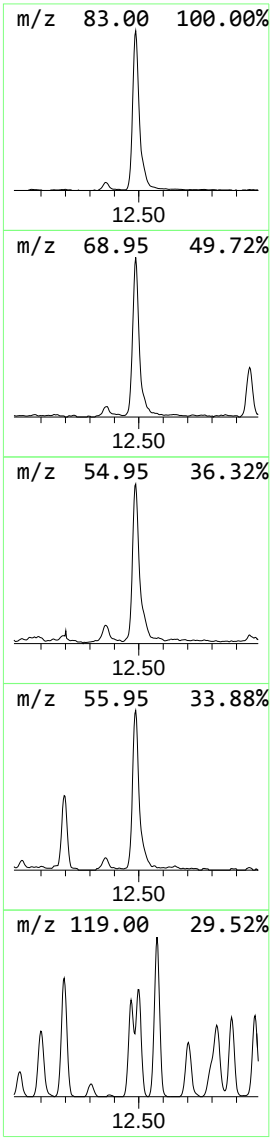
Peak Number 18 Cyclohexanone, 3,3,5-trimet... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.488	24.34 ug/L	483853	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	1,3-Cyclohexanedione, 5,5-dimethyl-	140	C8H12O2	000126-81-8	47





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU100721\
Data File : VU045227.D
Acq On : 07 Oct 2021 23:04
Operator : SY/MD
Sample : M4138-08
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F8

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

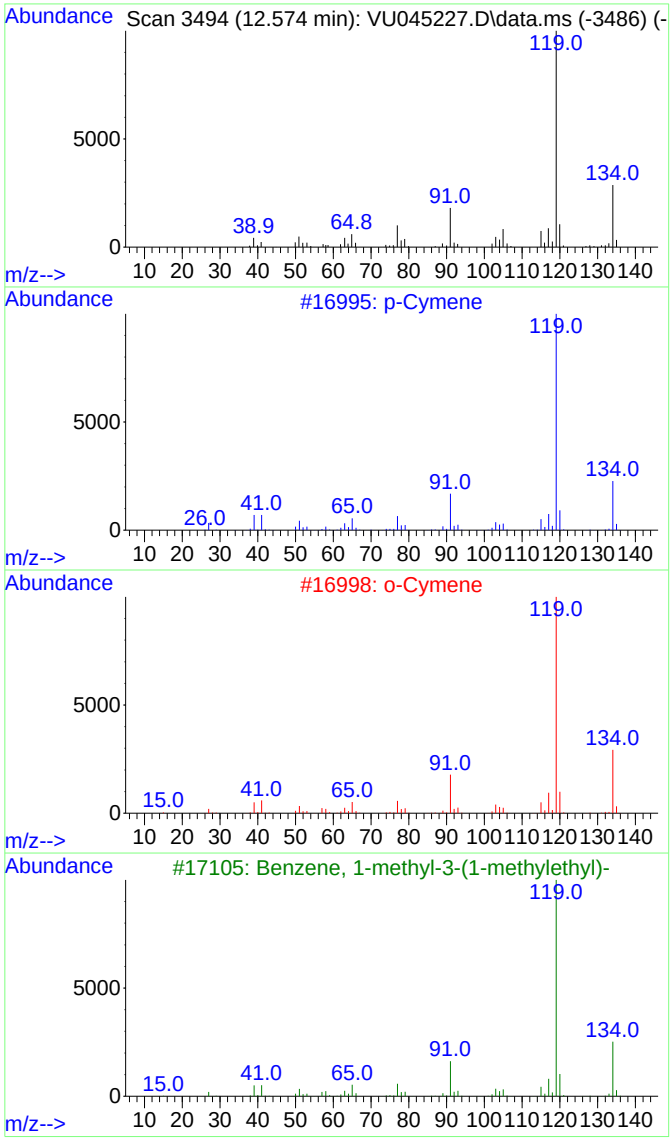
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

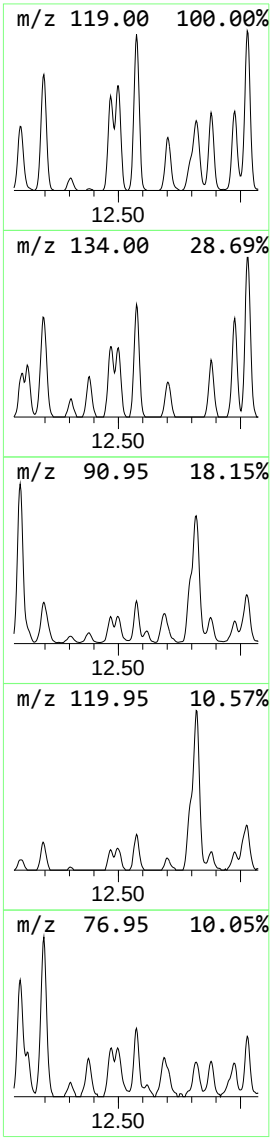
****

Peak Number 19 o-Cymene Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.574	6.10 ug/L	121211	1,4-Dichlorobenzene-d4	11.816

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





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Data File : VU045227.D
Acq On : 07 Oct 2021 23:04
Operator : SY/MD
Sample : M4138-08
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F8

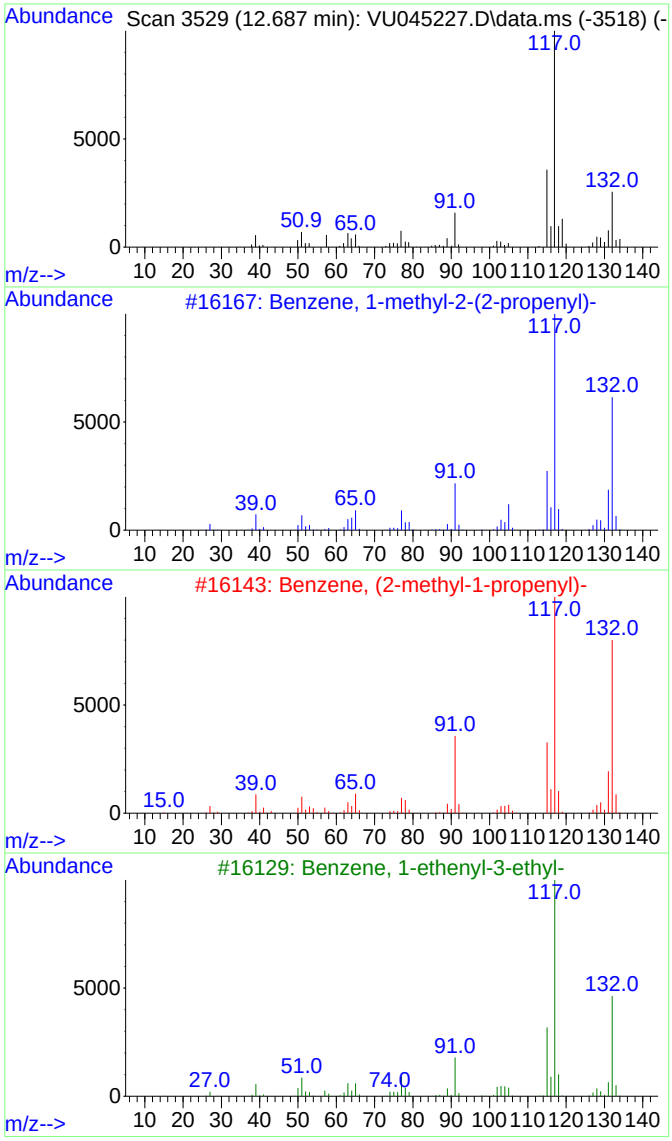
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092321WMA.M
Quant Title : VOC Analysis

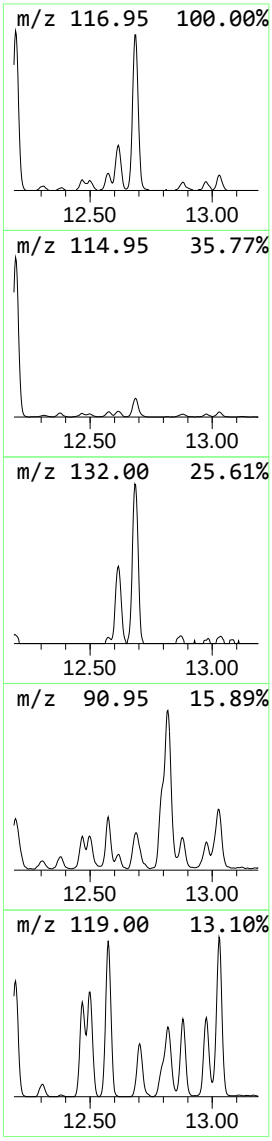
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 Benzene, 1-methyl-2-(2-prop... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.687	7.06 ug/L	140433	1,4-Dichlorobenzene-d4	11.816

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	87
2	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	86
3	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	83
4	Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	81
5	Indan, 1-methyl-	132	C10H12	000767-58-8	81





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Acq On : 07 Oct 2021 23:04
Operator : SY/MD
Sample : M4138-08
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW7F8

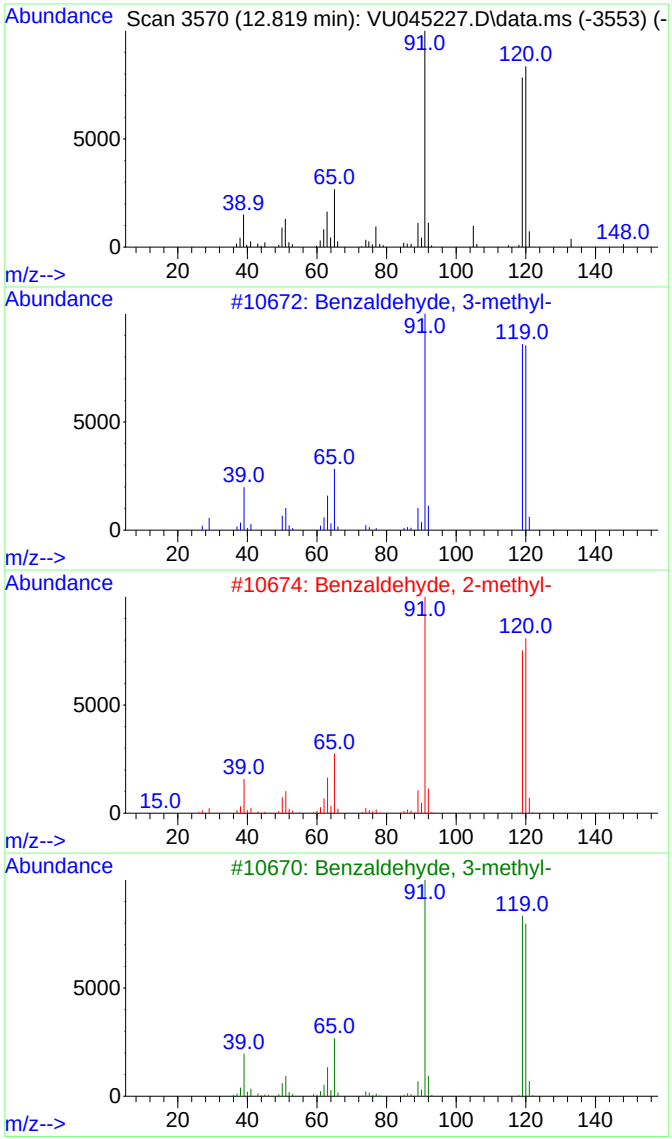
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092321WMA.M
Quant Title : VOC Analysis

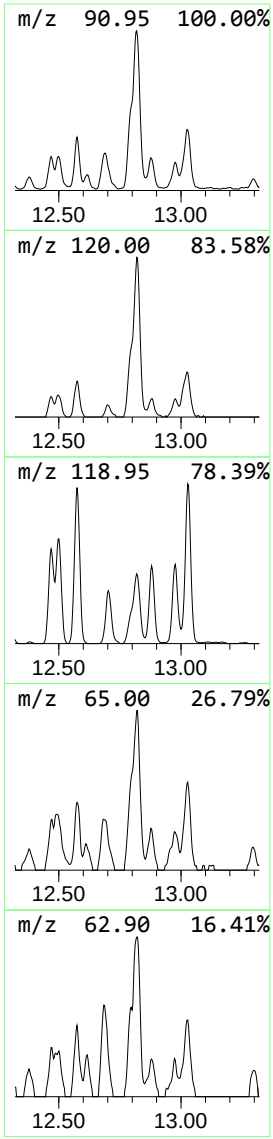
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 21 Benzaldehyde, 3-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.819	7.52 ug/L	149464	1,4-Dichlorobenzene-d4	11.816

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzaldehyde, 3-methyl-	120	C8H8O	000620-23-5	97
2	Benzaldehyde, 2-methyl-	120	C8H8O	000529-20-4	97
3	Benzaldehyde, 3-methyl-	120	C8H8O	000620-23-5	96
4	Benzaldehyde, 4-methyl-	120	C8H8O	000104-87-0	96
5	Benzaldehyde, 4-methyl-	120	C8H8O	000104-87-0	96





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Data File : VU045227.D
Acq On : 07 Oct 2021 23:04
Operator : SY/MD
Sample : M4138-08
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F8

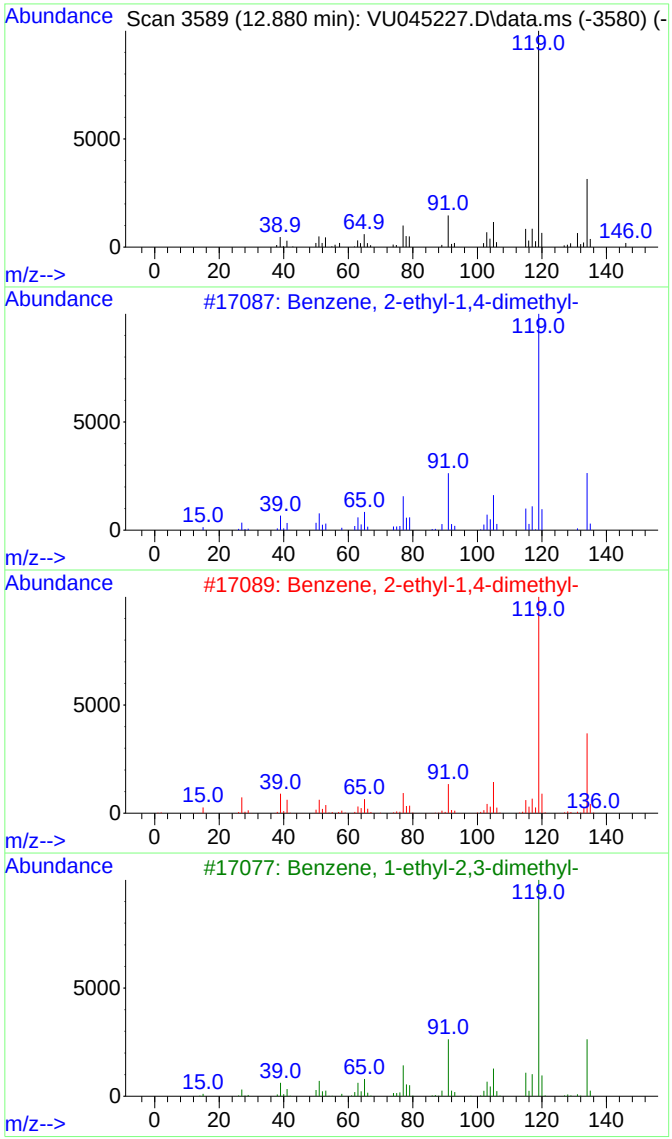
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092321WMA.M
Quant Title : VOC Analysis

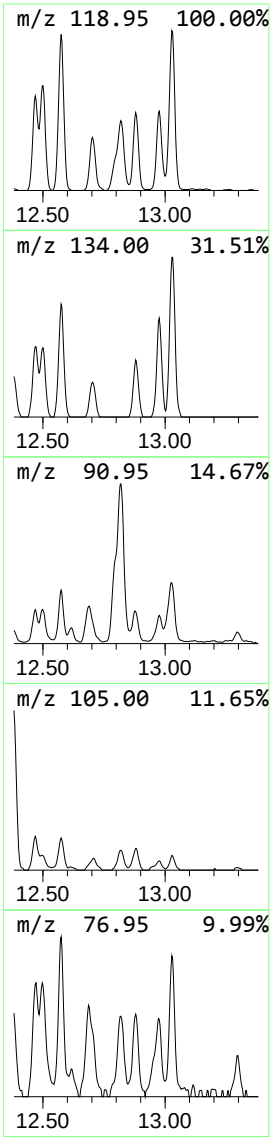
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.880	3.28 ug/L	65184	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, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
3	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
4	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94
5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	93





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU100721\
Data File : VU045227.D
Acq On : 07 Oct 2021 23:04
Operator : SY/MD
Sample : M4138-08
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F8

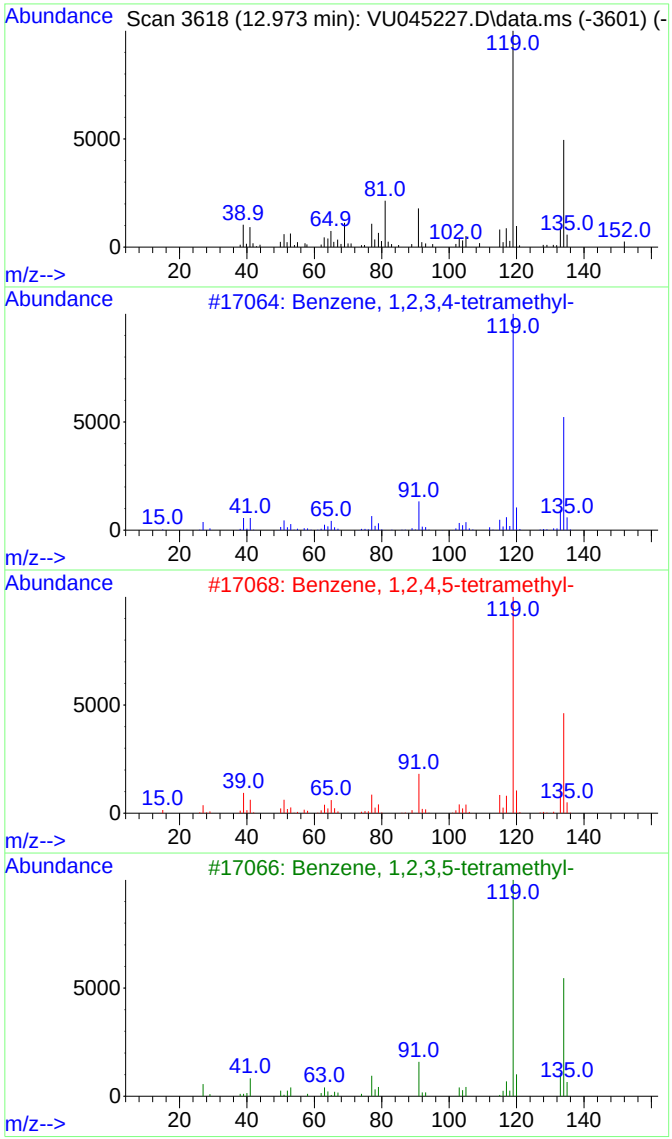
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092321WMA.M
Quant Title : VOC Analysis

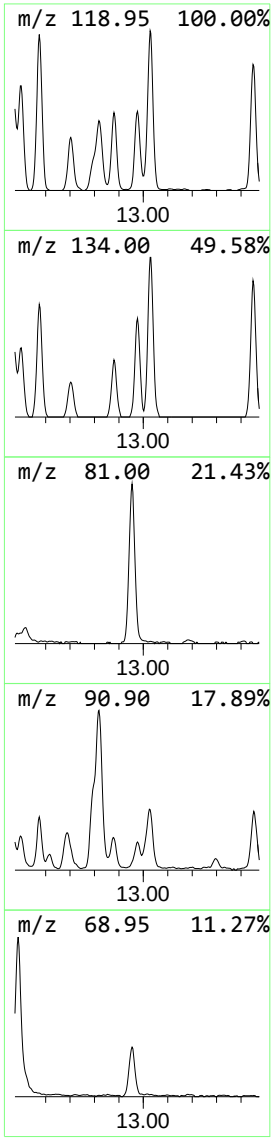
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.973	6.49 ug/L	128980	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, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
3	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
4	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	96
5	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95





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Data File : VU045227.D
Acq On : 07 Oct 2021 23:04
Operator : SY/MD
Sample : M4138-08
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F8

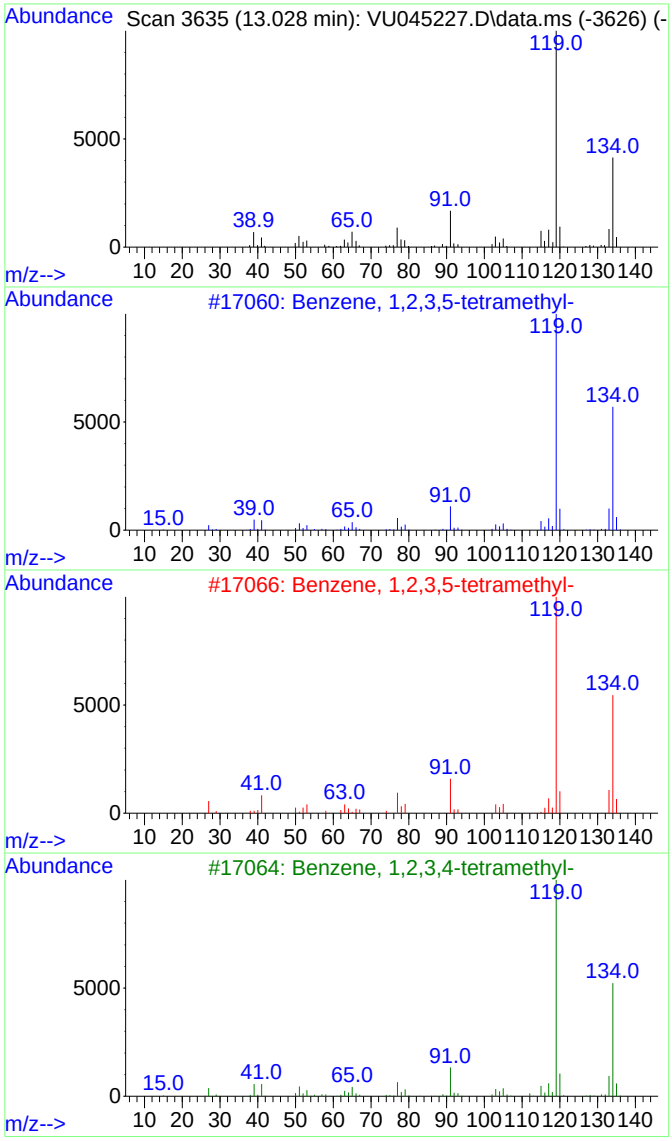
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092321WMA.M
Quant Title : VOC Analysis

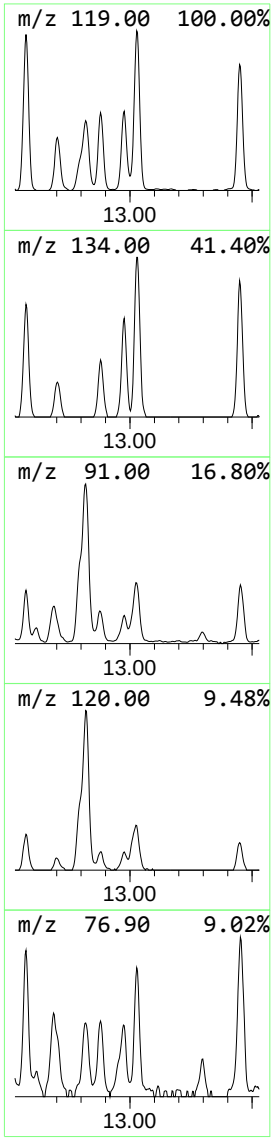
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.028	6.55 ug/L	130255	1,4-Dichlorobenzene-d4	11.816

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





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Data File : VU045227.D
Acq On : 07 Oct 2021 23:04
Operator : SY/MD
Sample : M4138-08
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F8

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

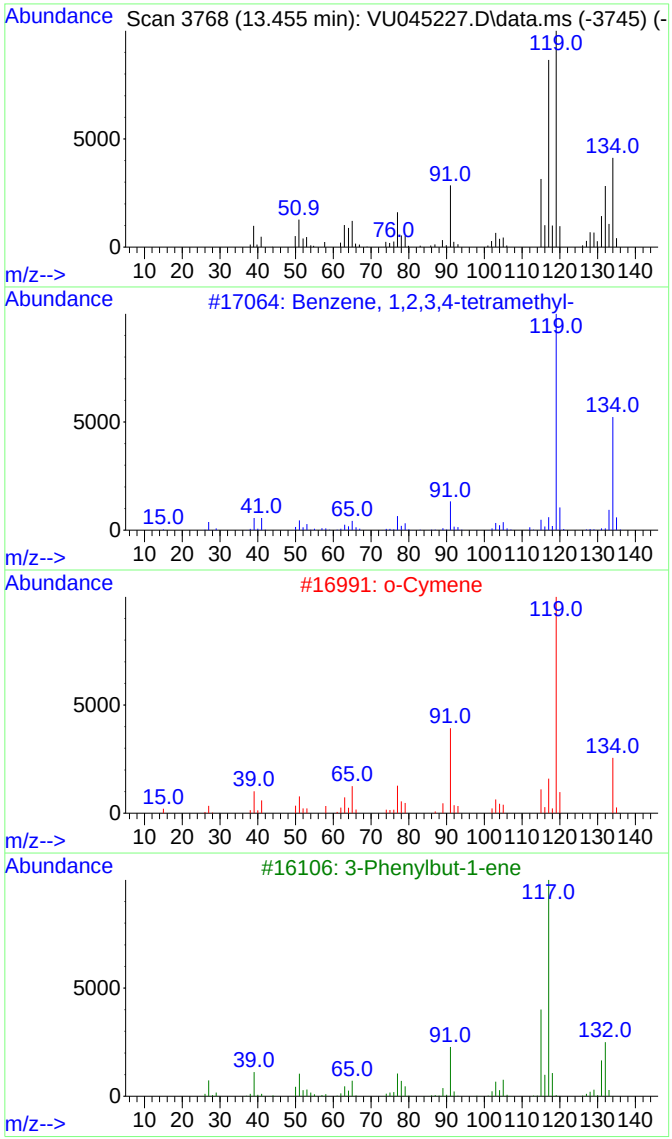
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

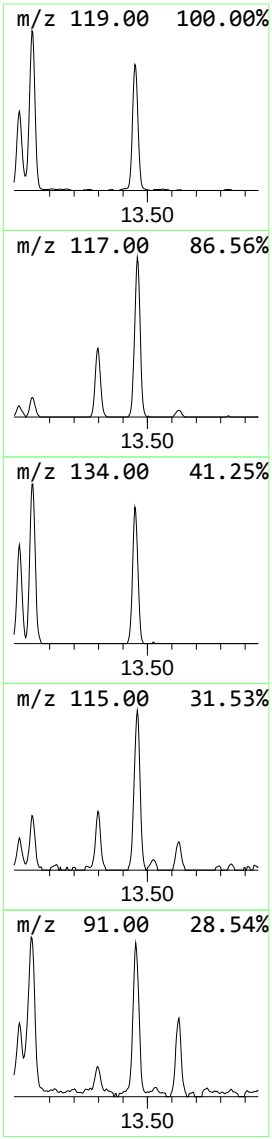
****

Peak Number 25 3-Phenylbut-1-ene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.455	9.21 ug/L	183194	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	50
2	o-Cymene	134	C10H14	000527-84-4	50
3	3-Phenylbut-1-ene	132	C10H12	000934-10-1	50
4	1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	50
5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	50





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU100721\
Data File : VU045227.D
Acq On : 07 Oct 2021 23:04
Operator : SY/MD
Sample : M4138-08
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F8

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

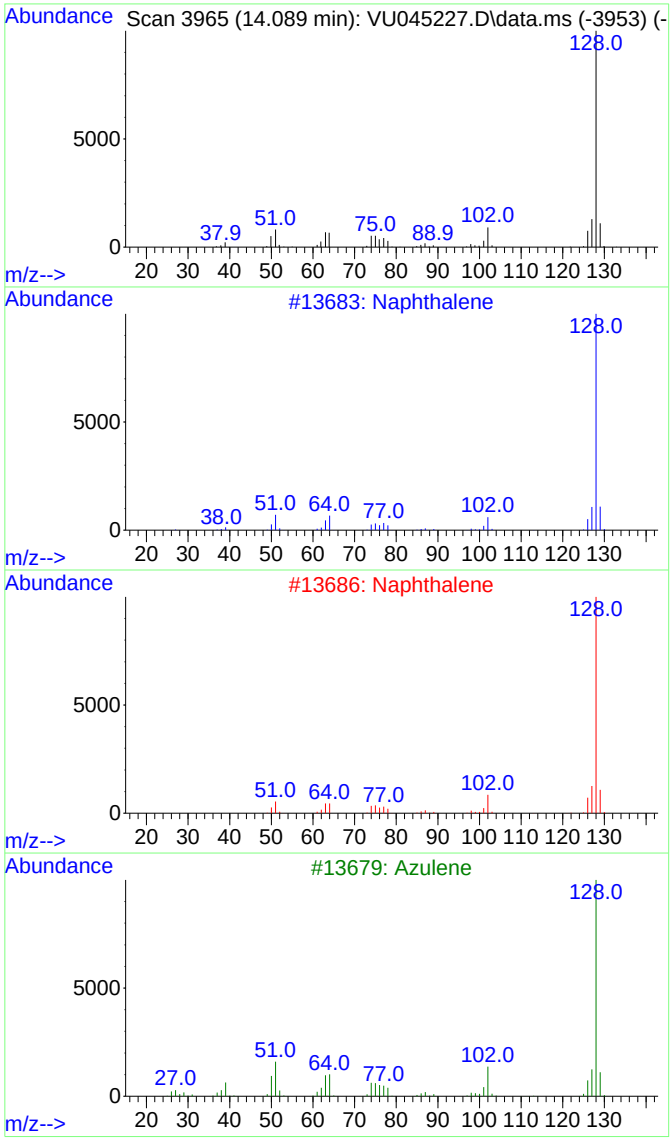
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

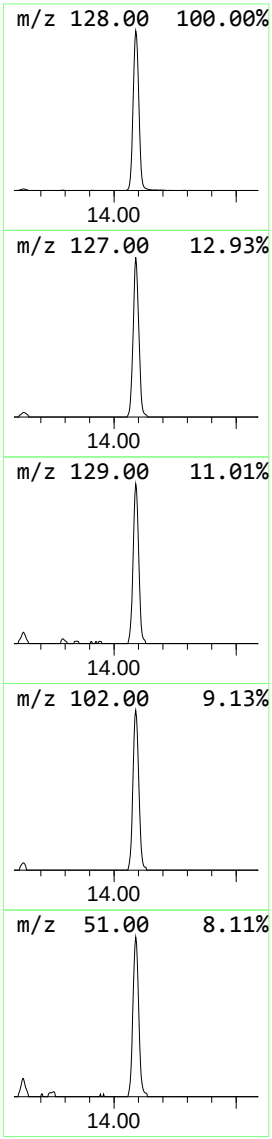
****

Peak Number 26 Azulene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.089	13.12 ug/L	260874	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	Azulene	128	C10H8	000275-51-4	94
4	Naphthalene	128	C10H8	000091-20-3	93
5	Azulene	128	C10H8	000275-51-4	91





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU100721\
Data File : VU045227.D
Acq On : 07 Oct 2021 23:04
Operator : SY/MD
Sample : M4138-08
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : EW7F8

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...	1.359	2.6	ug/L	37407	1	6.253	731201	50.0
Sulfur dioxide	1.498	152.3	ug/L	2226710	1	6.253	731201	50.0
Aminomethanesul...	1.845	6.7	ug/L	98596	1	6.253	731201	50.0
(DEL) Alkane: S...	3.359	2.9	ug/L	41838	1	6.253	731201	50.0
(DEL) Alkane: S...	3.697	2.9	ug/L	42936	1	6.253	731201	50.0
(DEL) Alkane: C...	4.533	4.7	ug/L	68428	1	6.253	731201	50.0
2-Butanol	4.989	29.9	ug/L	436991	1	6.253	731201	50.0
2-Pentanol, 4-m...	8.234	11.5	ug/L	220557	2	9.423	957754	50.0
Benzene, propyl-	10.909	36.2	ug/L	720585	3	11.816	994052	50.0
Benzene, 1-ethy...	11.002	157.3	ug/L	3128310	3	11.816	994052	50.0
4-Heptanone, 2,...	11.237	4.0	ug/L	80200	3	11.816	994052	50.0
Benzene, 1-ethy...	11.295	53.1	ug/L	1056060	3	11.816	994052	50.0
p-Cymene	11.745	4.1	ug/L	81625	3	11.816	994052	50.0
Benzene, 1,2,3-...	11.899	68.1	ug/L	1353590	3	11.816	994052	50.0
Indane	12.098	31.1	ug/L	618756	3	11.816	994052	50.0
Benzene, 1-meth...	12.375	4.9	ug/L	97938	3	11.816	994052	50.0
Cyclohexanone, ...	12.488	24.3	ug/L	483853	3	11.816	994052	50.0
o-Cymene	12.574	6.1	ug/L	121211	3	11.816	994052	50.0
Benzene, 1-meth...	12.687	7.1	ug/L	140433	3	11.816	994052	50.0
Benzaldehyde, 3...	12.819	7.5	ug/L	149464	3	11.816	994052	50.0
Benzene, 2-ethy...	12.880	3.3	ug/L	65184	3	11.816	994052	50.0
Benzene, 1,2,3,...	12.973	6.5	ug/L	128980	3	11.816	994052	50.0
Benzene, 1,2,3,...	13.028	6.5	ug/L	130255	3	11.816	994052	50.0
3-Phenylbut-1-ene	13.455	9.2	ug/L	183194	3	11.816	994052	50.0
Azulene	14.089	13.1	ug/L	260874	3	11.816	994052	50.0