LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV060421\
Data File : VV021505.D
Acq On : 04 Jun 2021 14:32
Operator : SY/MD
Sample : M2618-15
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG5W0

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF
Sampling : 1
Start Thrs: 0.2
Stop Thrs : 0

Filtering: 5
Min Area: 0 % of largest Peak
Max Peaks: 100
Peak Location: TOP

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

Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM052721WMA.M
Title : VOC Analysis

Signal : TIC: VV021505.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.108	16	21	33	rVB2	12701	14850	0.73%	0.059%
2	1.307	74	83	94	rVB	182832	192311	9.49%	0.764%
3	1.381	100	106	115	rBV	21980	24236	1.20%	0.096%
4	1.567	156	164	174	rVV	136298	155363	7.67%	0.617%
5	1.632	176	184	194	rVB	71873	90383	4.46%	0.359%
6	1.805	230	238	252	rVB2	43869	64378	3.18%	0.256%
7	1.902	260	268	278	rVB2	14697	19483	0.96%	0.077%
8	2.024	299	306	317	rVV2	16204	25174	1.24%	0.100%
9	2.108	320	332	345	rVV	307012	470388	23.21%	1.869%
10	2.175	345	353	376	rVB	42126	95992	4.74%	0.381%
11	2.500	443	454	456	rVV5	16743	27376	1.35%	0.109%
12	2.532	458	464	472	rVV2	38733	68626	3.39%	0.273%
13	2.580	473	479	494	rVB5	16660	32010	1.58%	0.127%
14	2.763	519	536	554	rVB4	31887	69886	3.45%	0.278%
15	2.953	583	595	609	rBV7	6621	16731	0.83%	0.066%
16	3.043	609	623	645	rVB2	34633	71593	3.53%	0.284%
17	3.278	689	696	711	rVB4	6928	14409	0.71%	0.057%
18	3.583	787	791	805	rVB5	6780	12011	0.59%	0.048%
19	3.767	831	848	868	rVB2	60690	158144	7.80%	0.628%
20	3.882	868	884	912	rBV	145810	409844	20.22%	1.628%
21	4.352	1013	1030	1053	rBV	246003	610707	30.13%	2.426%
22	4.464	1057	1065	1080	rVB3	8982	19952	0.98%	0.079%
23	4.680	1114	1132	1148	rBV3	55384	143883	7.10%	0.572%
24	4.767	1148	1159	1185	rVB6	11630	34347	1.69%	0.136%
25	4.915	1191	1205	1230	rBV6	16615	48449	2.39%	0.192%
26	5.050	1230	1247	1275	rBV2	584744	1527719	75.38%	6.069%
27	5.236	1294	1305	1323	rVB2	41454	111679	5.51%	0.444%
28	5.326	1323	1333	1350	rVB4	12651	28153	1.39%	0.112%
29	5.496	1374	1386	1396	rBV5	15665	33517	1.65%	0.133%
30	5.619	1412	1424	1458	rBV	493243	1014718	50.07%	4.031%
31	5.908	1505	1514	1537	rVB7	7690	22928	1.13%	0.091%
32	6.072	1537	1565	1577	rBV	336556	703149	34.69%	2.793%
33	6.133	1579	1584	1598	rVV5	62624	120022	5.92%	0.477%
34	6.201	1598	1605	1614	rVV4	11074	22530	1.11%	0.089%
35	6.258	1617	1623	1633	rVV4	8232	16111	0.79%	0.064%
36	6.368	1648	1657	1669	rVB6	10858	20877	1.03%	0.083%

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV060421\
Data File : VV021505.D
Acq On : 04 Jun 2021 14:32
Operator : SY/MD
Sample : M2618-15
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG5W0

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_V\Method\SFAMVLM052721WMA.M
Title : VOC Analysis

37	6.657	1731	1747	1764	rVB2	25669	60054	2.96%	0.239%
38	6.779	1771	1785	1798	rBV3	35442	79370	3.92%	0.315%
39	6.844	1799	1805	1817	rVB6	6626	13005	0.64%	0.052%
40	6.988	1836	1850	1872	rBV	208107	374149	18.46%	1.486%
41	7.082	1874	1879	1897	rVB7	7616	16023	0.79%	0.064%
42	7.172	1897	1907	1917	rBV8	8155	15895	0.78%	0.063%
43	7.233	1917	1926	1942	rBV2	35560	86412	4.26%	0.343%
44	7.320	1942	1953	1965	rVV	724100	1252489	61.80%	4.975%
45	7.390	1965	1975	1998	rVB	421556	722358	35.64%	2.869%
46	7.625	2038	2048	2062	rVV	137248	244363	12.06%	0.971%
47	7.696	2062	2070	2087	rVB	70241	134455	6.63%	0.534%
48	7.792	2092	2100	2109	rVB5	7320	13476	0.66%	0.054%
49	7.979	2151	2158	2166	rVB4	7586	10934	0.54%	0.043%
50	8.088	2182	2192	2226	rBV	518097	929397	45.86%	3.692%
51	8.291	2249	2255	2264	rVB2	8528	14101	0.70%	0.056%
52	8.853	2417	2430	2451	rBV	732912	1188507	58.64%	4.721%
53	9.014	2469	2480	2495	rBV	168206	274137	13.53%	1.089%
54	9.136	2506	2518	2535	rBV	1262815	2026670	100.00%	8.050%
55	9.545	2633	2645	2661	rBV	532042	824114	40.66%	3.274%
56	9.712	2690	2697	2712	rVB2	10091	17067	0.84%	0.068%
57	9.815	2718	2729	2738	rBV10	6527	12797	0.63%	0.051%
58	9.934	2754	2766	2780	rBV	20370	33287	1.64%	0.132%
59	10.127	2816	2826	2837	rVB5	7427	12980	0.64%	0.052%
60	10.217	2837	2854	2878	rBV	547741	855924	42.23%	3.400%
61	10.355	2880	2897	2913	rBV	54569	89415	4.41%	0.355%
62	10.448	2914	2926	2945	rBV	590116	1210712	59.74%	4.809%
63	10.541	2945	2955	2966	rVB	251527	371651	18.34%	1.476%
64	10.737	2996	3016	3036	rBV	248850	389252	19.21%	1.546%
65	10.914	3060	3071	3090	rVB	988416	1471234	72.59%	5.844%
66	11.053	3104	3114	3120	rBV7	6706	13349	0.66%	0.053%
67	11.094	3120	3127	3137	rVV7	9991	19725	0.97%	0.078%
68	11.197	3151	3159	3165	rVV	23461	34053	1.68%	0.135%
69	11.249	3165	3175	3192	rVV	851367	1304956	64.39%	5.184%
70	11.335	3192	3202	3215	rVB	277127	419302	20.69%	1.666%
71	11.528	3250	3262	3272	rBV	302541	536848	26.49%	2.133%
72	11.577	3273	3277	3283	rVV2	93951	126414	6.24%	0.502%
73	11.625	3283	3292	3315	rVB2	885688	1516252	74.81%	6.023%
74	11.747	3323	3330	3338	rBV5	9732	14101	0.70%	0.056%
75	11.821	3343	3353	3366	rVB	36802	61365	3.03%	0.244%
76	11.911	3371	3381	3386	rBV	80063	122510	6.04%	0.487%

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV060421\
Data File : VV021505.D
Acq On : 04 Jun 2021 14:32
Operator : SY/MD
Sample : M2618-15
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG5W0

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_V\Method\SFAMVLM052721WMA.M
Title : VOC Analysis

77	11.943	3387	3391	3402	rVB	67640	90053	4.44%	0.358%
78	12.017	3404	3414	3422	rBV	122067	193749	9.56%	0.770%
79	12.117	3433	3445	3467	rVB2	89205	162929	8.04%	0.647%
80	12.316	3498	3507	3518	rVB	33530	55262	2.73%	0.220%
81	12.413	3528	3537	3545	rBV	71388	104971	5.18%	0.417%
82	12.467	3545	3554	3567	rVB	102234	154538	7.63%	0.614%
83	12.551	3571	3580	3586	rBV2	11101	18119	0.89%	0.072%
84	12.724	3625	3634	3649	rVB	97089	154484	7.62%	0.614%
85	12.882	3664	3683	3699	rBV2	172650	305661	15.08%	1.214%
86	12.953	3701	3705	3714	rVV7	8574	11441	0.56%	0.045%
87	13.001	3714	3720	3726	rVV	8517	11136	0.55%	0.044%
88	13.046	3726	3734	3748	rVB	31278	52997	2.61%	0.211%
89	13.168	3760	3772	3778	rBV4	12536	20131	0.99%	0.080%
90	13.216	3779	3787	3795	rVB2	22039	31736	1.57%	0.126%
91	13.271	3796	3804	3811	rBV2	26658	42435	2.09%	0.169%
92	13.368	3828	3834	3842	rVB	14381	17805	0.88%	0.071%
93	13.503	3867	3876	3895	rVB	89674	148613	7.33%	0.590%
94	13.712	3930	3941	3952	rBV7	11655	22694	1.12%	0.090%
95	13.769	3952	3959	3970	rVV3	9842	15753	0.78%	0.063%
96	13.911	3994	4003	4013	rVB3	13893	20771	1.02%	0.083%
97	14.091	4051	4059	4060	rBV2	10893	11814	0.58%	0.047%
98	14.294	4114	4122	4132	rVB4	7657	12143	0.60%	0.048%
99	14.634	4215	4228	4240	rBV	34690	54958	2.71%	0.218%
100	14.818	4276	4285	4297	rVB	22445	35306	1.74%	0.140%

Sum of corrected areas: 25174531

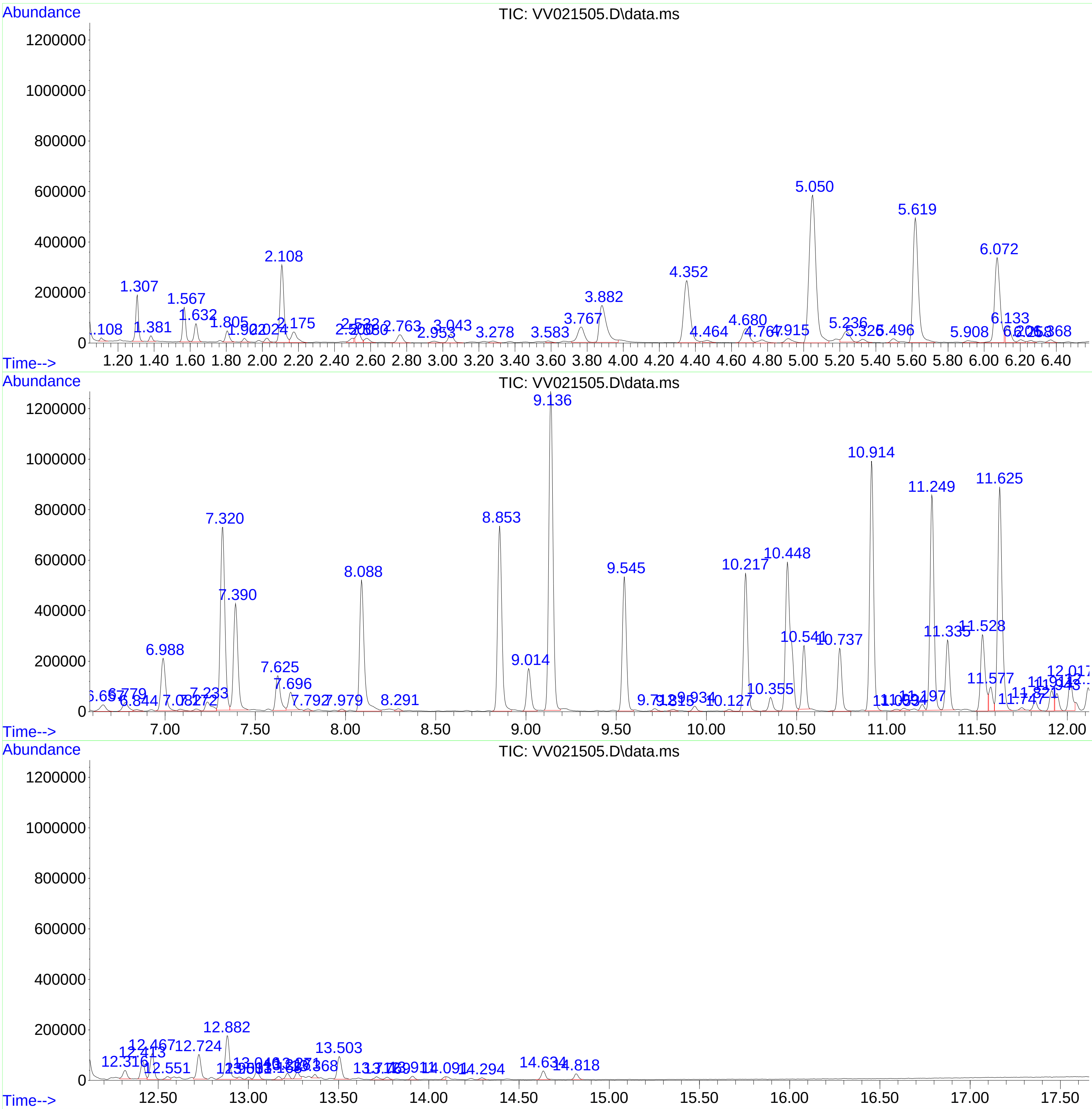
LSC Report - Integrated Chromatogram

```
Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV060421\
Data File : VV021505.D
Acq On    : 04 Jun 2021  14:32
Operator  : SY/MD
Sample    : M2618-15
Misc      : 5.0mL/MSVOA_V/WATER
ALS Vial  : 14   Sample Multiplier: 1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM052721WMA.M
Quant Title  : VOC Analysis

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

Instrument :
MSVOA_V
ClientSampleId :
BG5W0



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Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV060421\
Data File : VV021505.D
Acq On    : 04 Jun 2021  14:32
Operator  : SY/MD
Sample    : M2618-15
Misc      : 5.0mL/MSVOA_V/WATER
ALS Vial  : 14   Sample Multiplier: 1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM052721WMA.M
Quant Title  : VOC Analysis

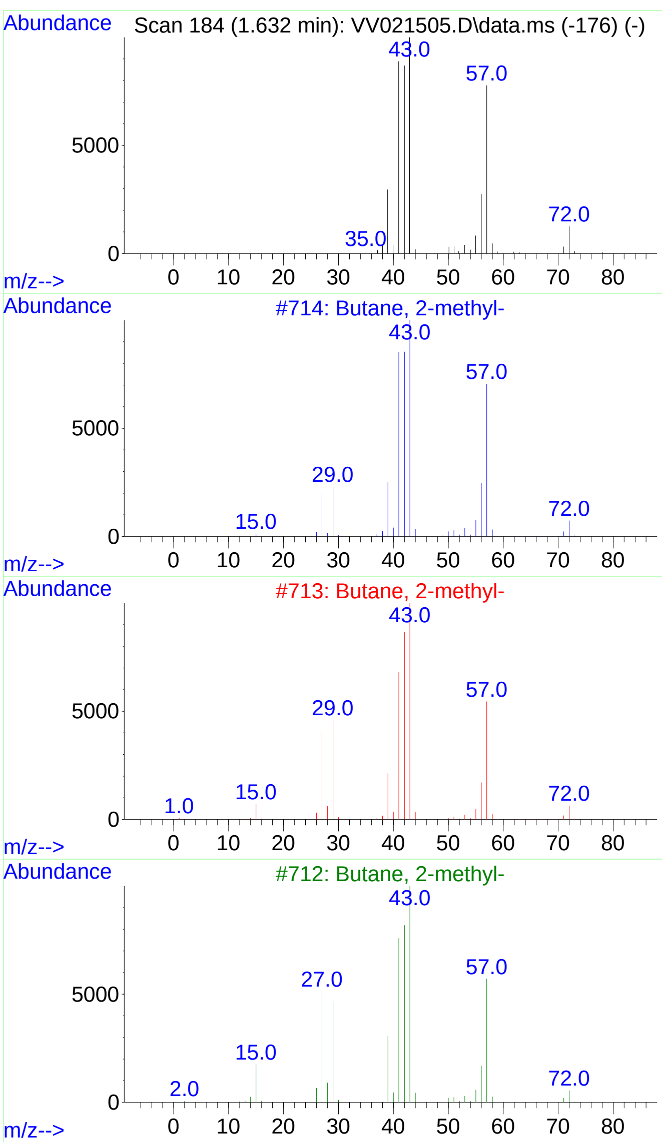
TIC Library    : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P
```

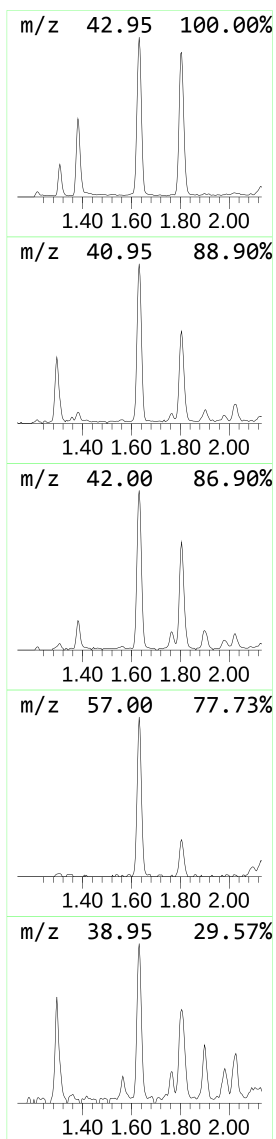
Instrument :
MSVOA_V
ClientSampleId :
BG5W0

```
*********************************************************************
Peak Number  1  (DEL) Alkane: Straight-Chai...   Concentration Rank 16
```

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.632	4.45 ug/L	90383	1,4-Difluorobenzene	5.619

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-	72	C5H12	000078-78-4	91
2	Butane, 2-methyl-	72	C5H12	000078-78-4	91
3	Butane, 2-methyl-	72	C5H12	000078-78-4	83
4	Pentane	72	C5H12	000109-66-0	59
5	Pentane	72	C5H12	000109-66-0	33





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Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV060421\
Data File : VV021505.D
Acq On    : 04 Jun 2021  14:32
Operator  : SY/MD
Sample    : M2618-15
Misc      : 5.0mL/MSVOA_V/WATER
ALS Vial  : 14   Sample Multiplier: 1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM052721WMA.M
Quant Title  : VOC Analysis

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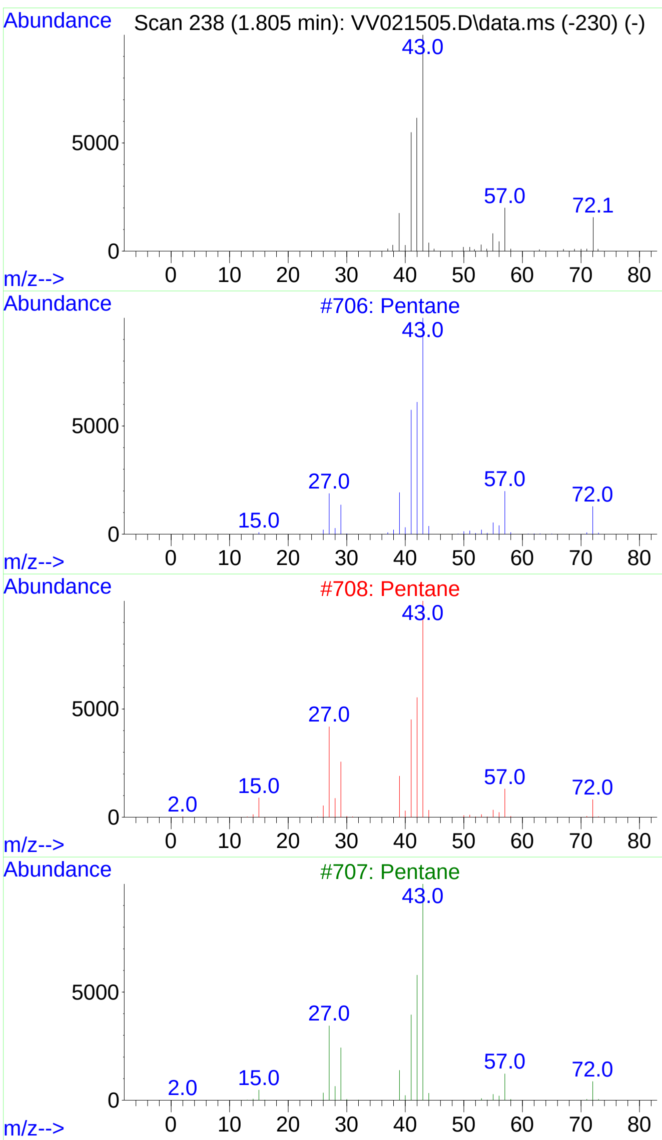
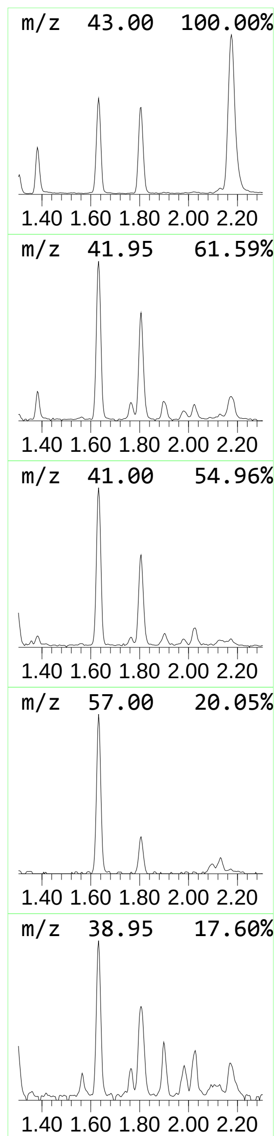

Instrument :
MSVOA_V
ClientSampleId :
BG5W0

****
Peak Number 2 (DEL) Alkane: Straight-Chai...   Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.805	3.17 ug/L	64378	1,4-Difluorobenzene	5.619

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane	72	C5H12	000109-66-0	91
2	Pentane	72	C5H12	000109-66-0	90
3	Pentane	72	C5H12	000109-66-0	86
4	Pentane	72	C5H12	000109-66-0	86
5	Isobutylene epoxide	72	C4H8O	000558-30-5	53





Library Search Compound Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV060421\
Data File : VV021505.D
Acq On : 04 Jun 2021 14:32
Operator : SY/MD
Sample : M2618-15
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG5W0

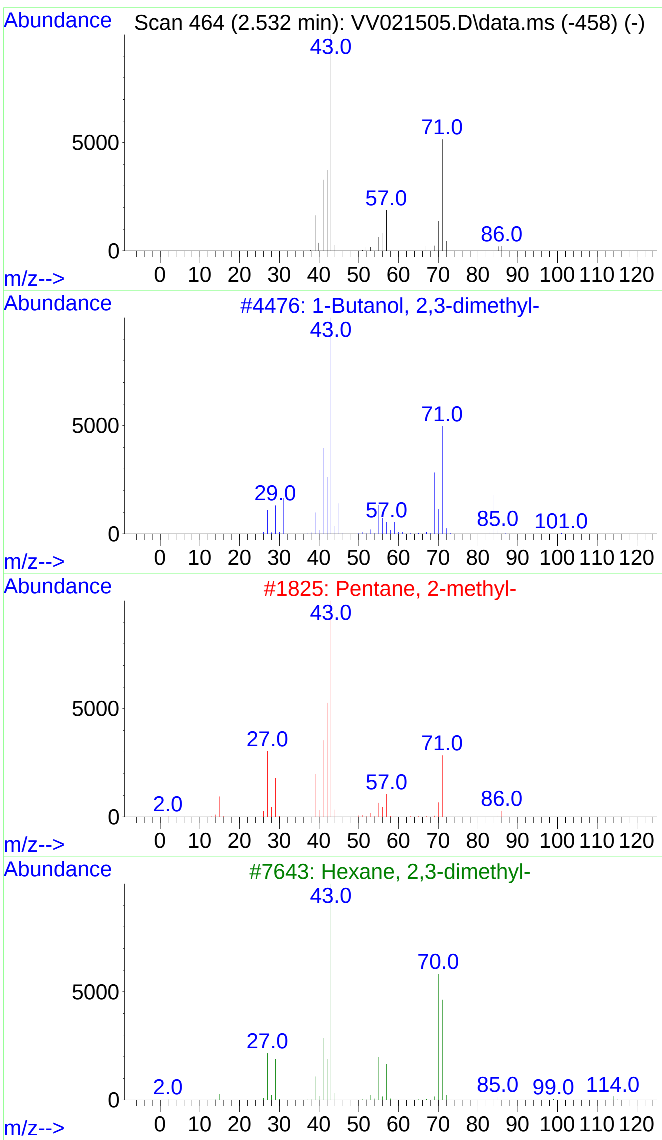
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM052721WMA.M
Quant Title : VOC Analysis

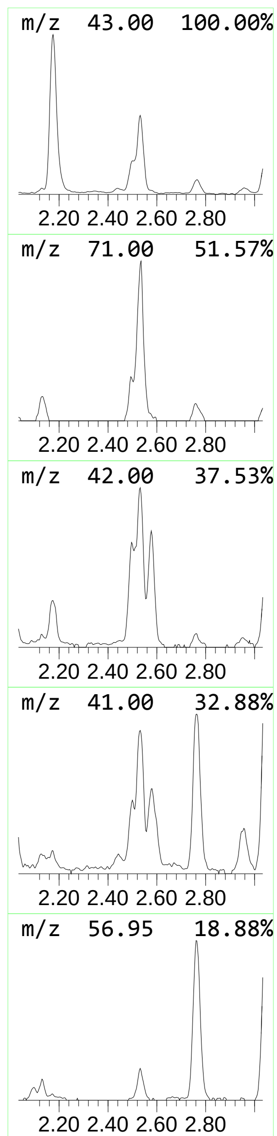
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 1-Butanol, 2,3-dimethyl- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.532	3.38 ug/L	68626	1,4-Difluorobenzene	5.619

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Butanol, 2,3-dimethyl-	102	C6H14O	019550-30-2	78
2	Pentane, 2-methyl-	86	C6H14	000107-83-5	50
3	Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	45
4	Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	43
5	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	33





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV060421\
Data File : VV021505.D
Acq On : 04 Jun 2021 14:32
Operator : SY/MD
Sample : M2618-15
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG5W0

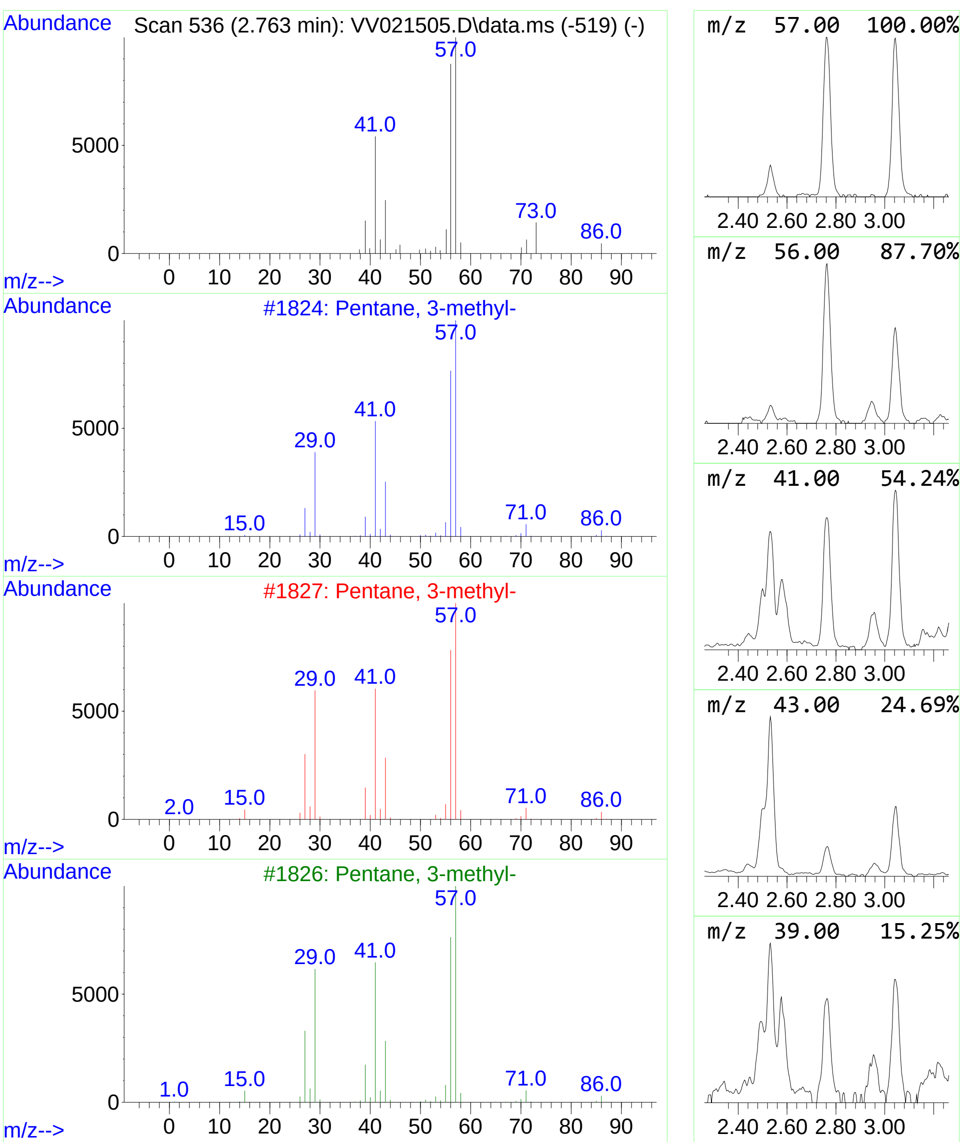
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM052721WMA.M
Quant Title : VOC Analysis

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.763	3.44 ug/L	69886	1,4-Difluorobenzene	5.619

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 3-methyl-	86	C6H14	000096-14-0	87
2	Pentane, 3-methyl-	86	C6H14	000096-14-0	87
3	Pentane, 3-methyl-	86	C6H14	000096-14-0	72
4	Furan, tetrahydro-3-methyl-	86	C5H10O	013423-15-9	59
5	Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	56



Library Search Compound Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV060421\
Data File : VV021505.D
Acq On : 04 Jun 2021 14:32
Operator : SY/MD
Sample : M2618-15
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG5W0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM052721WMA.M
Quant Title : VOC Analysis

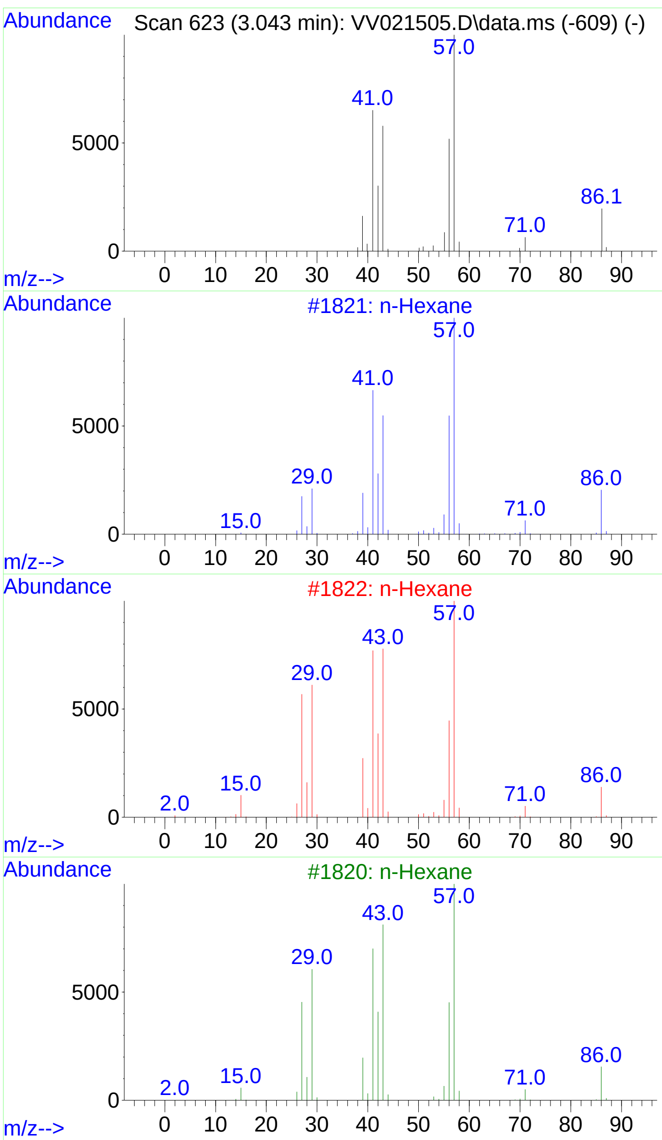
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

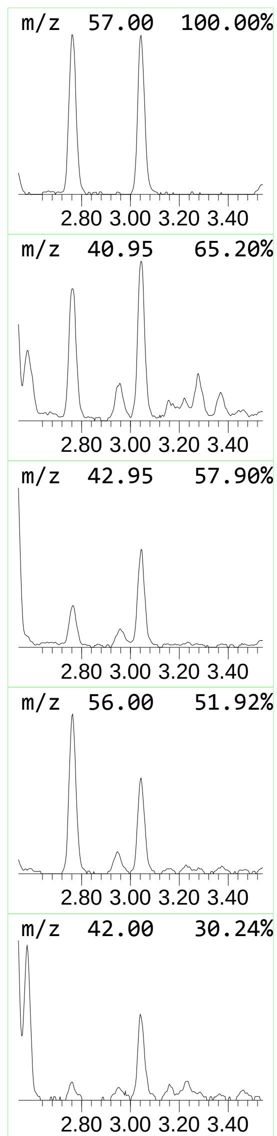
****

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.043	3.53 ug/L	71593	1,4-Difluorobenzene	5.619

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexane	86	C6H14	000110-54-3	91
2	n-Hexane	86	C6H14	000110-54-3	78
3	n-Hexane	86	C6H14	000110-54-3	59
4	Pentane, 3-methyl-	86	C6H14	000096-14-0	50
5	Morpholine	87	C4H9NO	000110-91-8	38





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Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV060421\
Data File : VV021505.D
Acq On    : 04 Jun 2021  14:32
Operator  : SY/MD
Sample    : M2618-15
Misc      : 5.0mL/MSVOA_V/WATER
ALS Vial  : 14   Sample Multiplier: 1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM052721WMA.M
Quant Title  : VOC Analysis

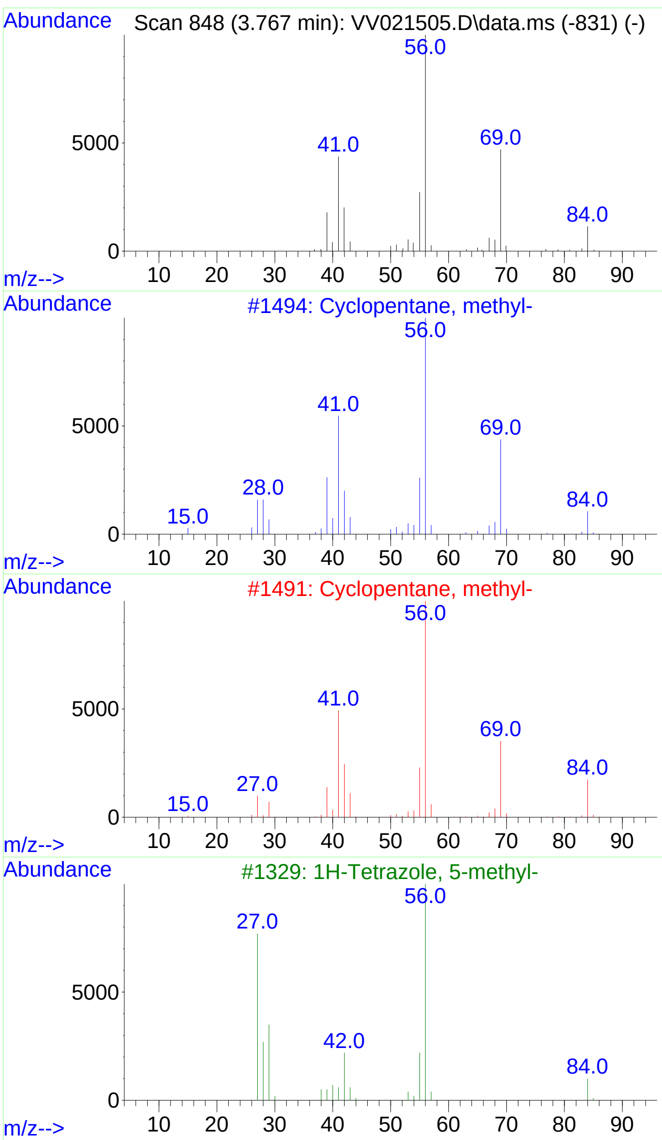
TIC Library    : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P
```

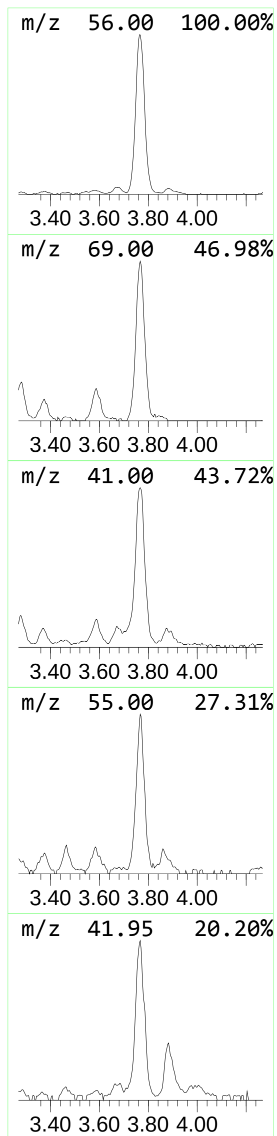
Instrument :
MSVOA_V
ClientSampleId :
BG5W0

```
*********************************************************************
Peak Number  6  (DEL) Alkane: Cyclic3.767       Concentration Rank  7
```

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.767	7.79 ug/L	158144	1,4-Difluorobenzene	5.619

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	90
2	Cyclopentane, methyl-	84	C6H12	000096-37-7	72
3	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	64
4	Cyclopentane, methyl-	84	C6H12	000096-37-7	64
5	Cyclobutane, ethyl-	84	C6H12	004806-61-5	64





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Acq On : 04 Jun 2021 14:32
Operator : SY/MD
Sample : M2618-15
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG5W0

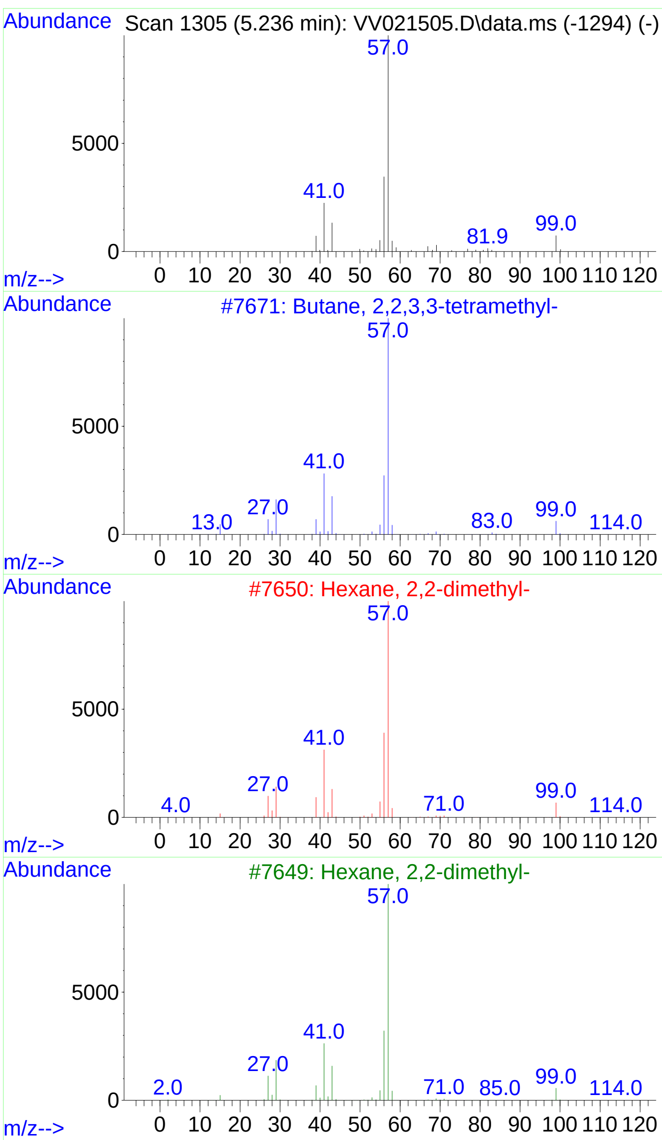
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM052721WMA.M
Quant Title : VOC Analysis

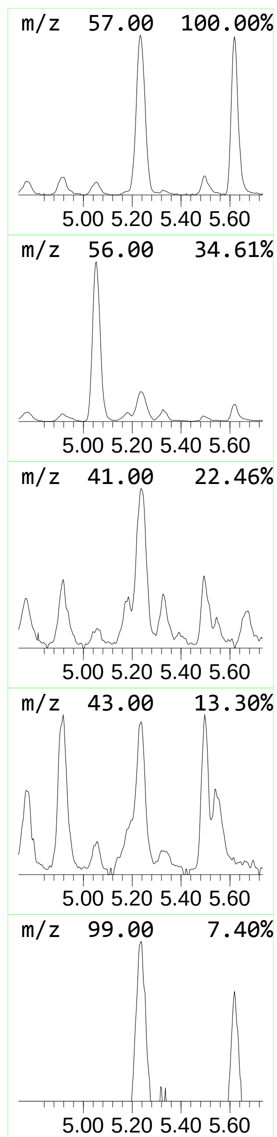
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.236	5.50 ug/L	111679	1,4-Difluorobenzene	5.619

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	72
2	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	64
3	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	64
4	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	56
5	Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	56





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Sample : M2618-15
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG5W0

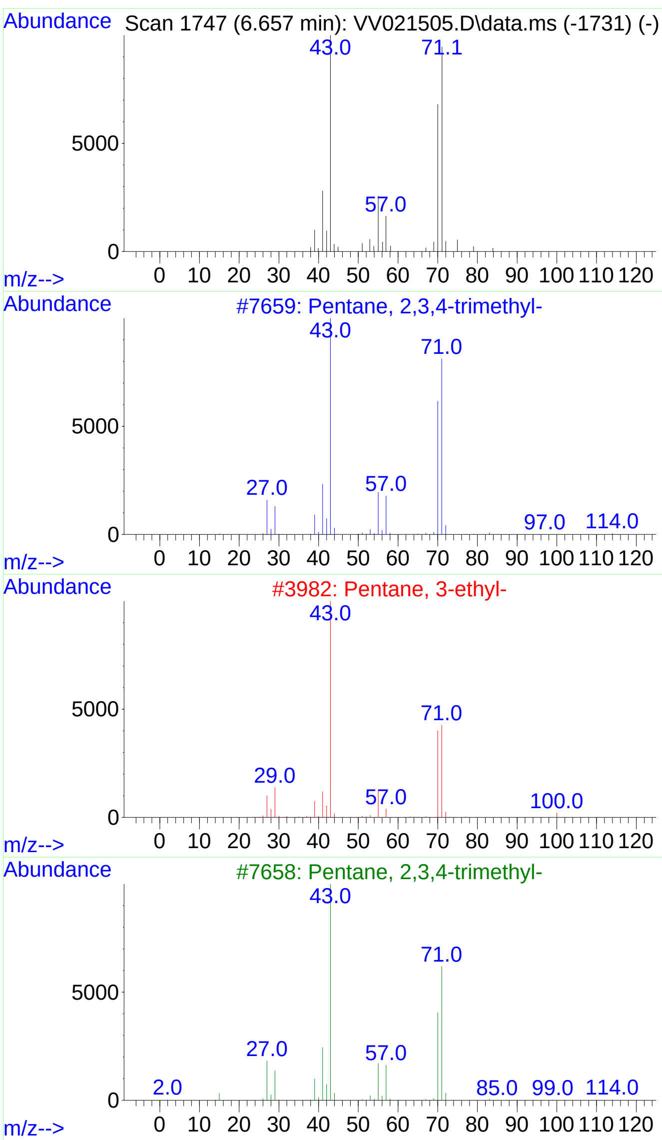
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM052721WMA.M
Quant Title : VOC Analysis

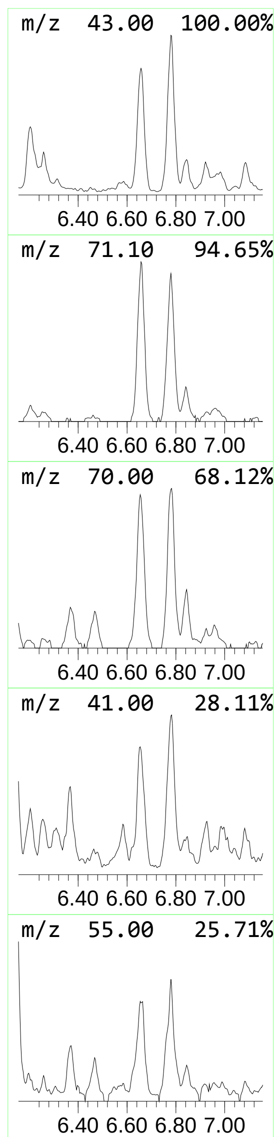
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.657	2.96 ug/L	60054	1,4-Difluorobenzene	5.619

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	78
2	Pentane, 3-ethyl-	100	C7H16	000617-78-7	72
3	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	72
4	Heptane, 3,3,4-trimethyl-	142	C10H22	020278-87-9	56
5	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	53





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ALS Vial  : 14   Sample Multiplier: 1

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Quant Title  : VOC Analysis

TIC Library    : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P
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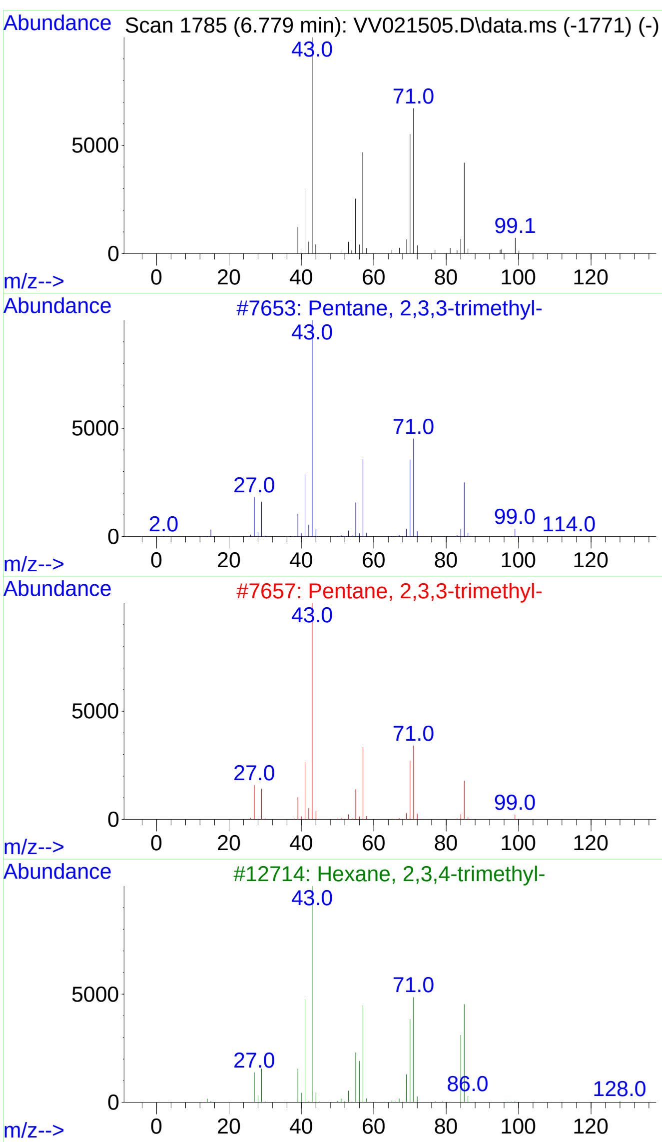
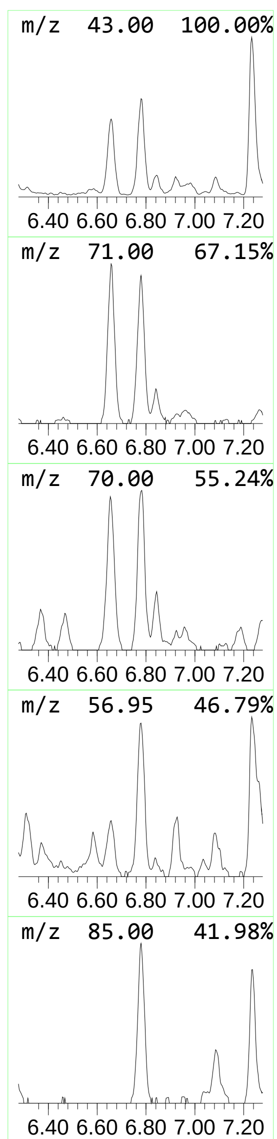
Instrument :
MSVOA_V
ClientSampleId :
BG5W0

```
**********************************************************************
Peak Number  9  (DEL) Alkane: Straight-Chai...  Concentration Rank 18
```

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.779	3.91 ug/L	79370	1,4-Difluorobenzene	5.619

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	90
2	Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	83
3	Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	78
4	Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	59
5	Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	59





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Operator : SY/MD
Sample : M2618-15
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG5W0

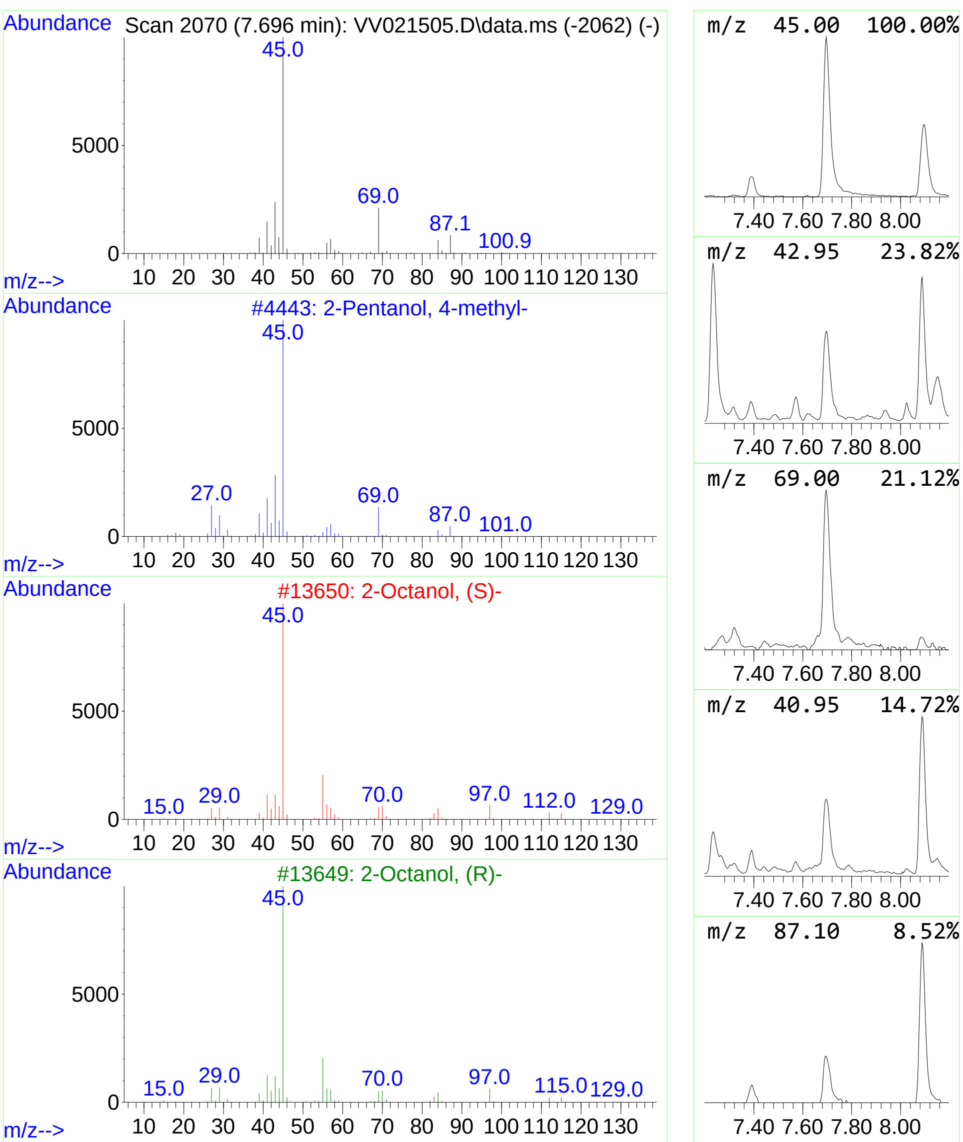
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM052721WMA.M
Quant Title : VOC Analysis

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

****
Peak Number 11 2-Pentanol, 4-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.696	5.66 ug/L	134455	Chlorobenzene-d5	8.853

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	74
2	2-Octanol, (S)-	130	C8H18O	006169-06-8	64
3	2-Octanol, (R)-	130	C8H18O	005978-70-1	64
4	2-Octanol	130	C8H18O	000123-96-6	64
5	2-Hexanol	102	C6H14O	000626-93-7	64



Library Search Compound Report

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Misc      : 5.0mL/MSVOA_V/WATER
ALS Vial  : 14   Sample Multiplier: 1

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Quant Title  : VOC Analysis

TIC Library    : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P
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Instrument :
MSVOA_V
ClientSampleId :
BG5W0

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*********************************************************************
Peak Number 12  Benzene, propyl-              Concentration Rank 22
```
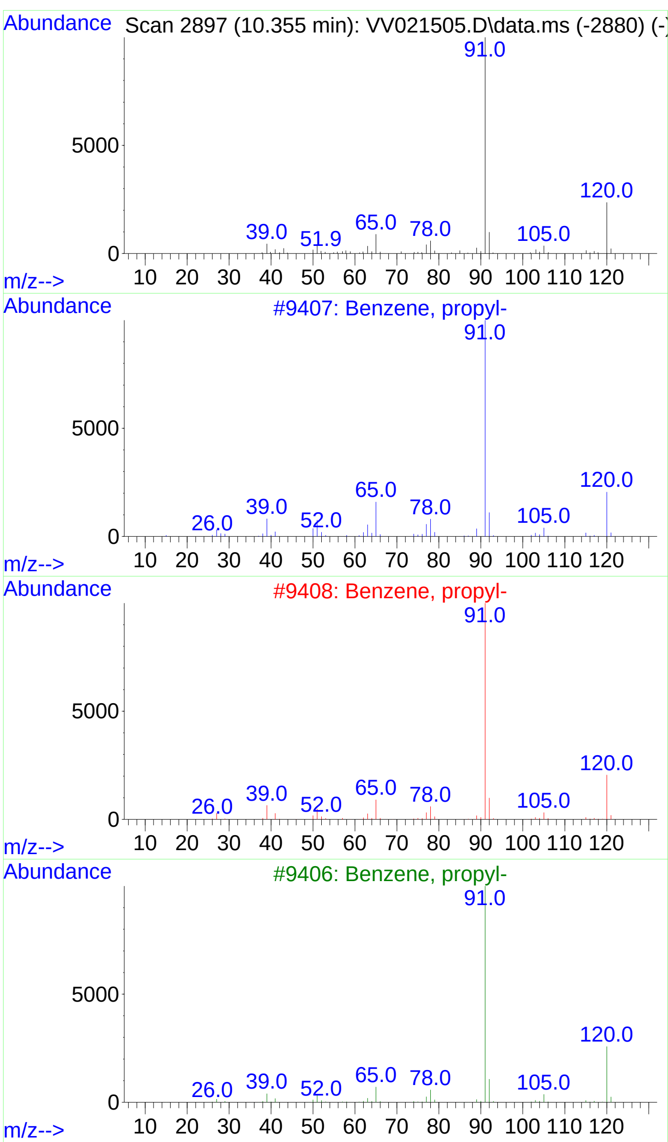
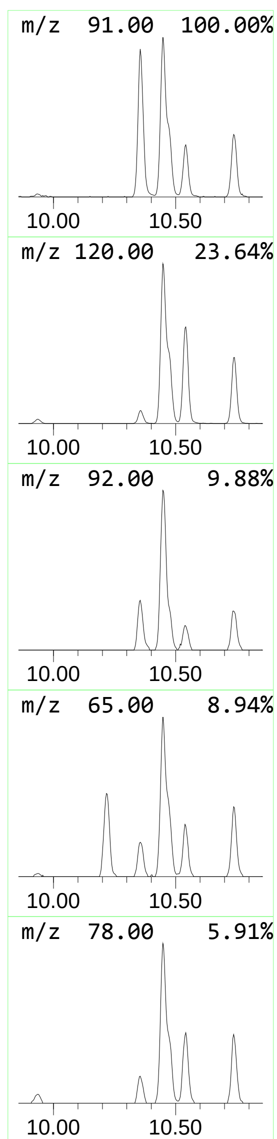

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.355	3.43 ug/L	89415	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, propyl-	120	C9H12	000103-65-1	90
2	Benzene, propyl-	120	C9H12	000103-65-1	87
3	Benzene, propyl-	120	C9H12	000103-65-1	87
4	N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	64
5	Propanenitrile, 3-[(phenylmethyl...	160	C10H12N2	000706-03-6	59





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Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG5W0

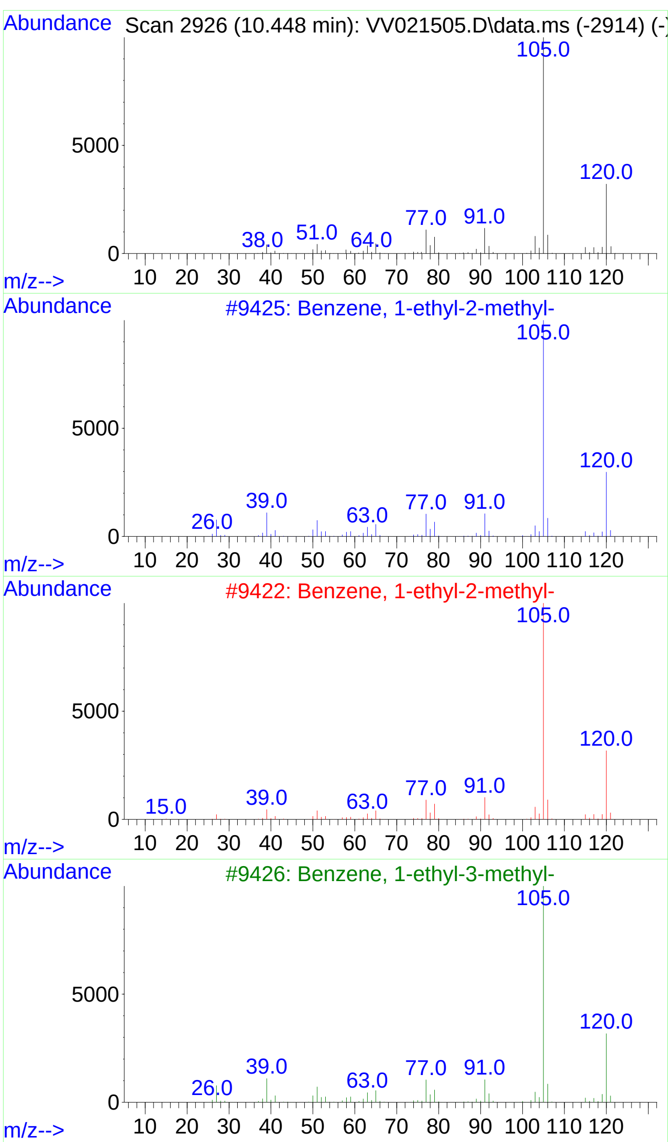
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM052721WMA.M
Quant Title : VOC Analysis

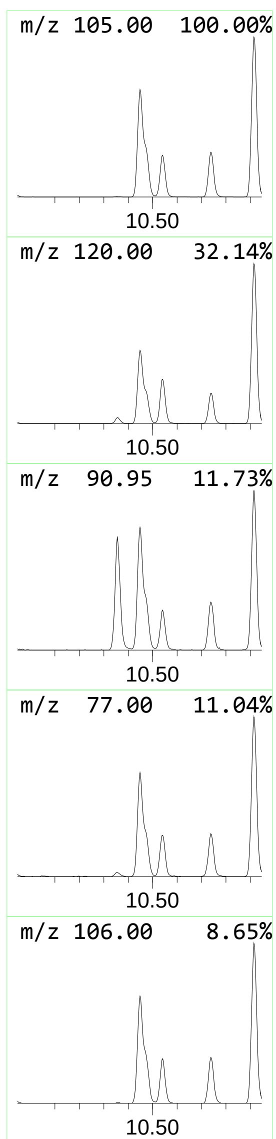
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.448	46.39 ug/L	1210710	1,4-Dichlorobenzene-d4	11.252

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	95
5	Mesitylene	120	C9H12	000108-67-8	91





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Operator : SY/MD
Sample : M2618-15
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG5W0

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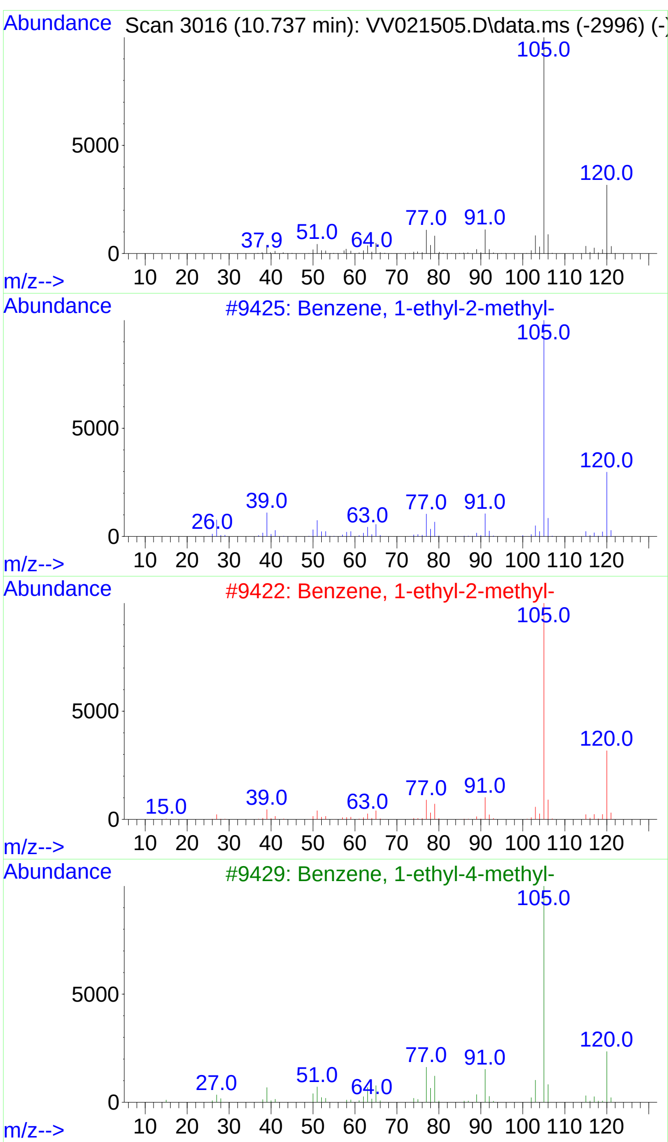
Quant Title : VOC Analysis

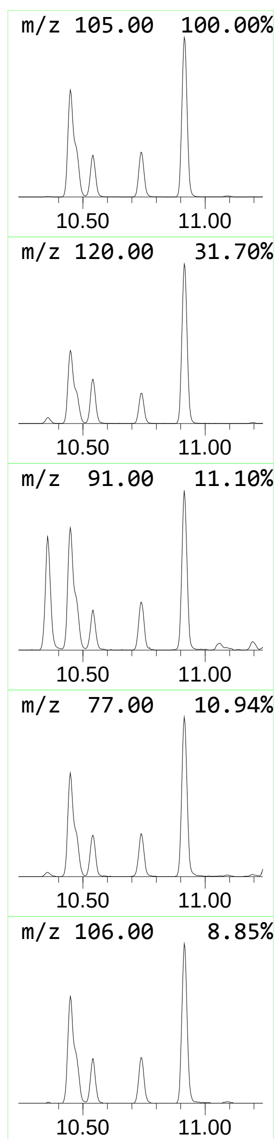
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 Benzene, 1-ethyl-4-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.737	14.91 ug/L	389252	1,4-Dichlorobenzene-d4	11.252

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	93
4	Mesitylene	120	C9H12	000108-67-8	90
5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	90





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Operator : SY/MD
Sample : M2618-15
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG5W0

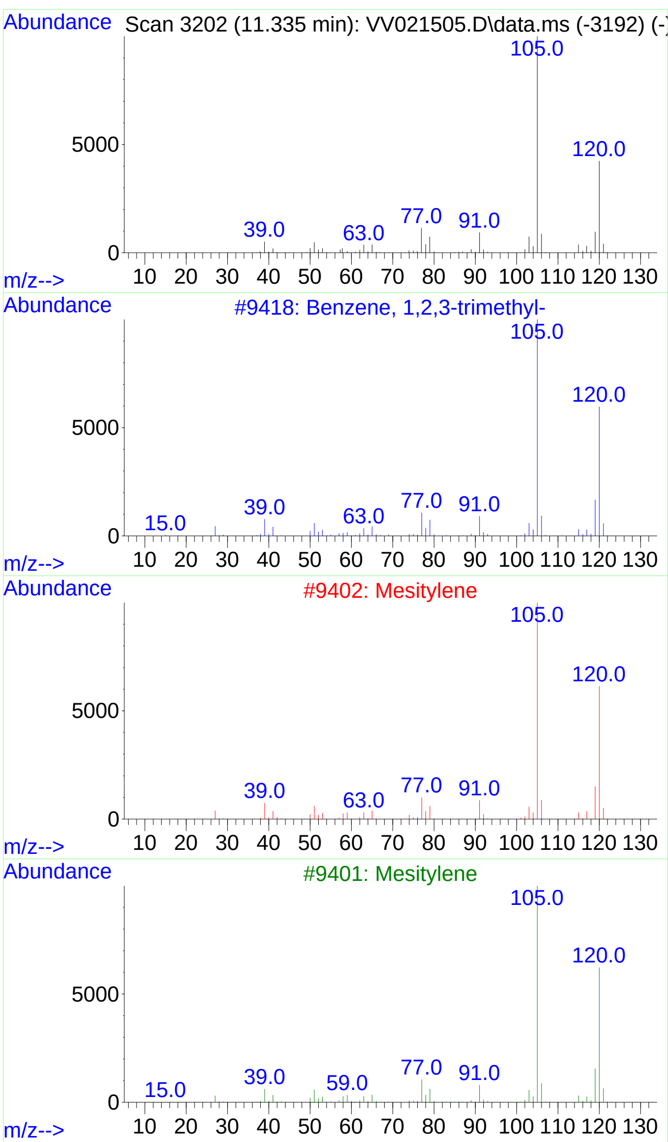
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM052721WMA.M
Quant Title : VOC Analysis

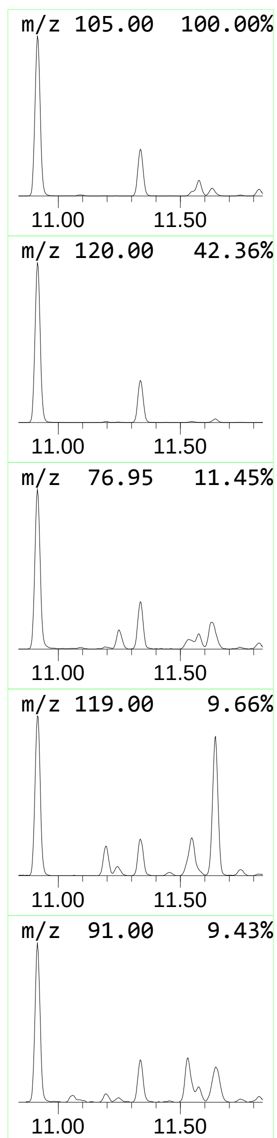
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 Benzene, 1,2,3-trimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.336	16.07 ug/L	419302	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2	Mesitylene	120	C9H12	000108-67-8	93
3	Mesitylene	120	C9H12	000108-67-8	91
4	Mesitylene	120	C9H12	000108-67-8	91
5	Mesitylene	120	C9H12	000108-67-8	91





Instrument :
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ClientSampleId :
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Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 14 Sample Multiplier: 1

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Quant Title : VOC Analysis

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

****

Peak Number 16 Indane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.528	20.57 ug/L	536848	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	95
2	Indane	118	C9H10	000496-11-7	90
3	Indane	118	C9H10	000496-11-7	81
4	Benzene, cyclopropyl-	118	C9H10	000873-49-4	81
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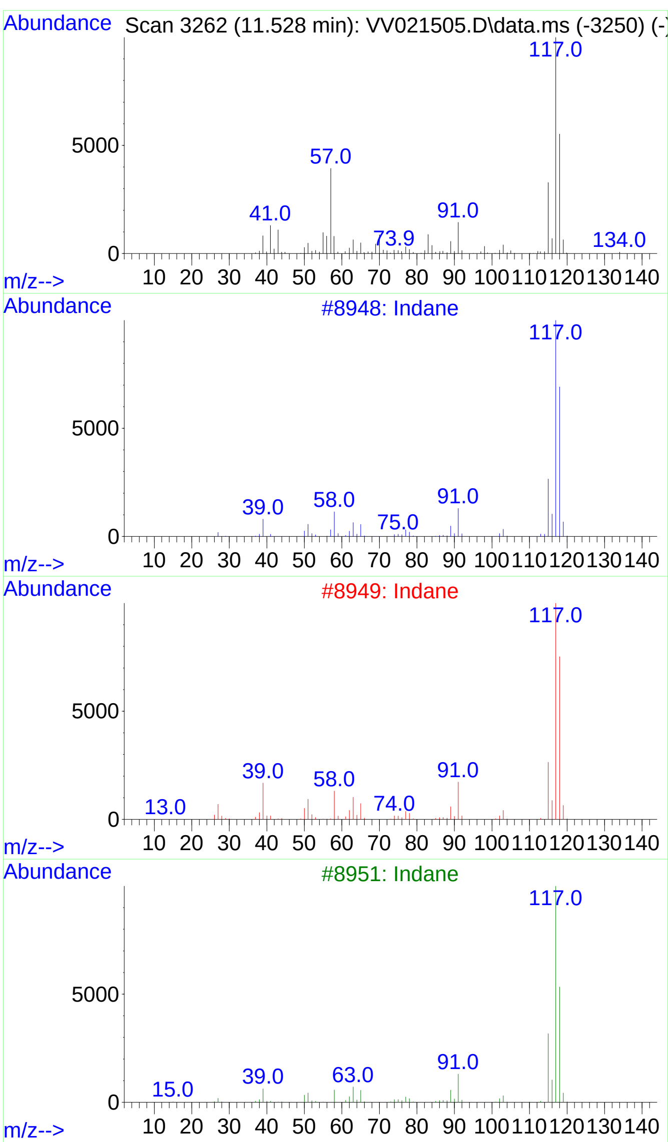

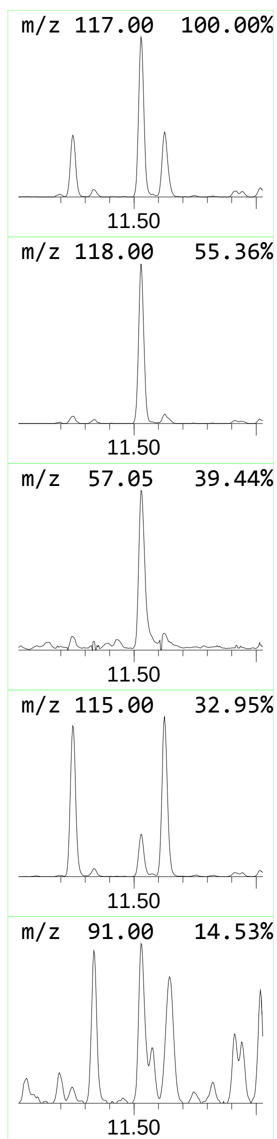

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ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG5W0

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Quant Title : VOC Analysis

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

Peak Number 17 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.911	4.69 ug/L	122510	1,4-Dichlorobenzene-d4	11.252

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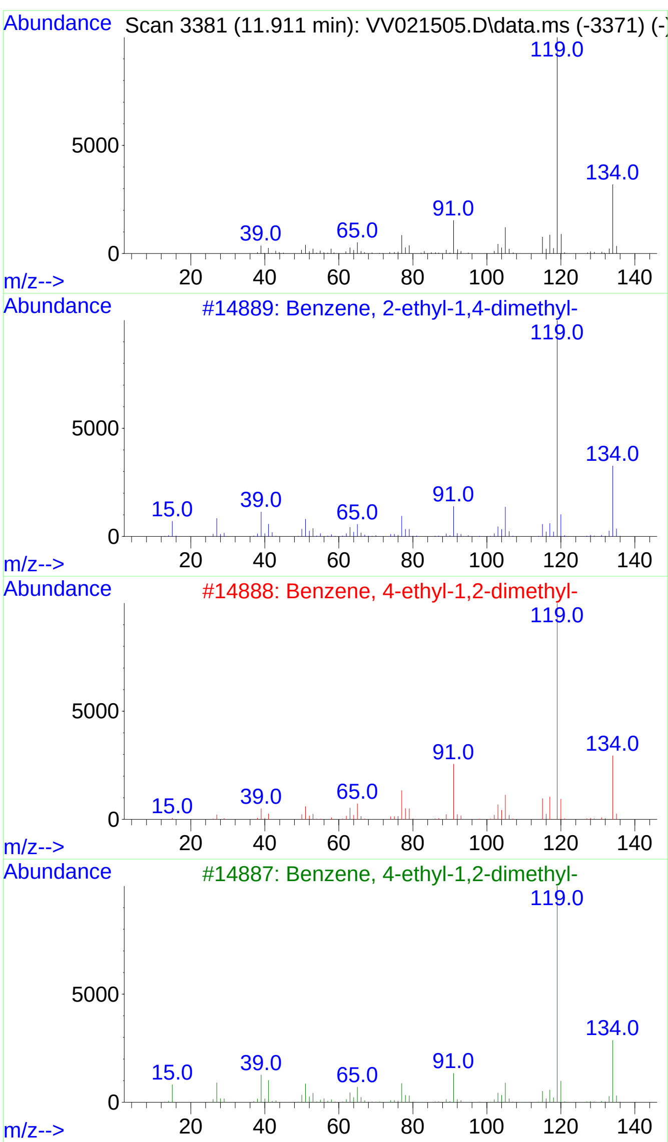

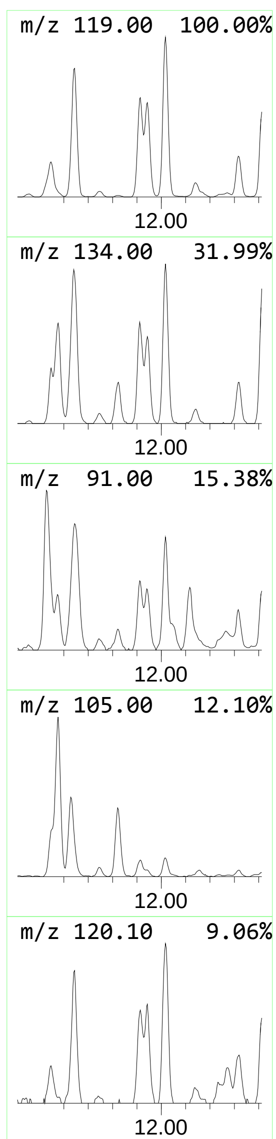

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Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG5W0

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Quant Title : VOC Analysis

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

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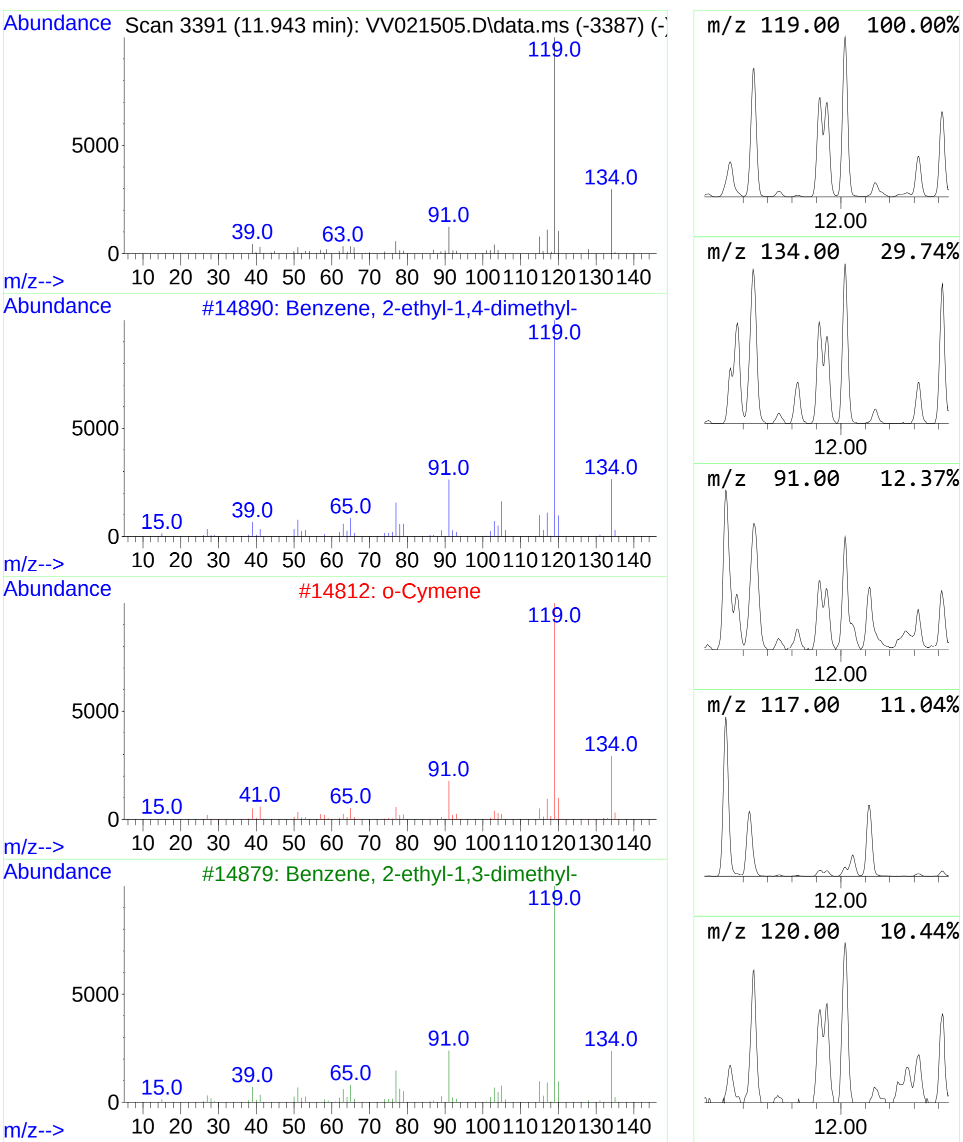
Peak Number 18 o-Cymene Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.943	3.45 ug/L	90053	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91
2	o-Cymene	134	C10H14	000527-84-4	91
3	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	91
4	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	91
5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	90



Library Search Compound Report

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Acq On : 04 Jun 2021 14:32
Operator : SY/MD
Sample : M2618-15
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 14 Sample Multiplier: 1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM052721WMA.M
Quant Title : VOC Analysis

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

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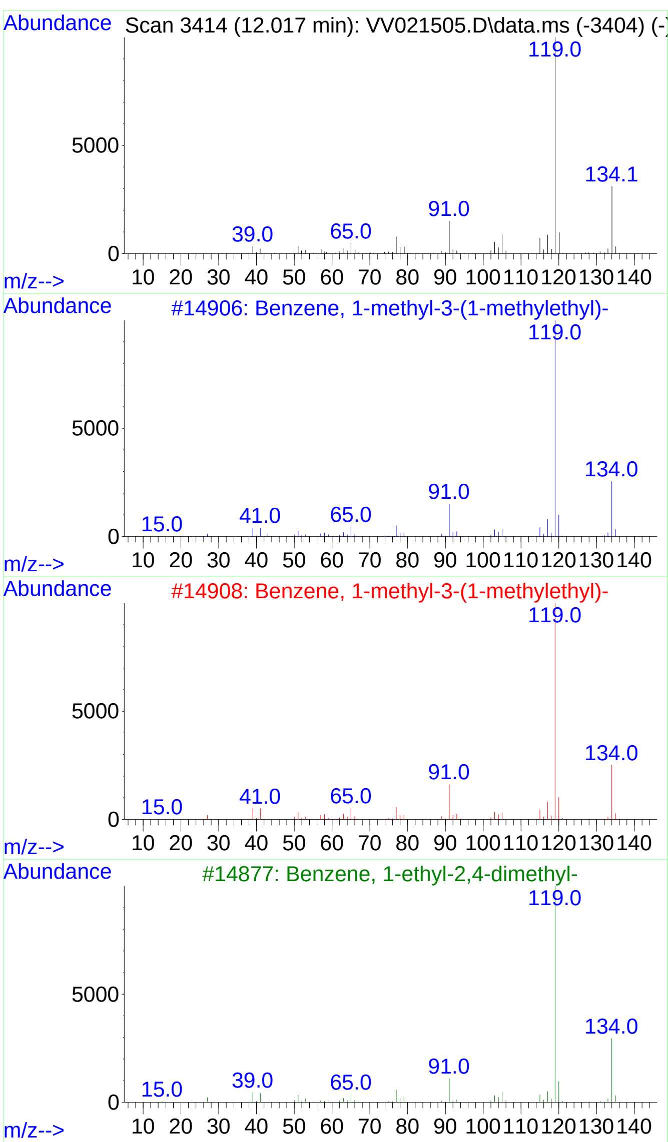
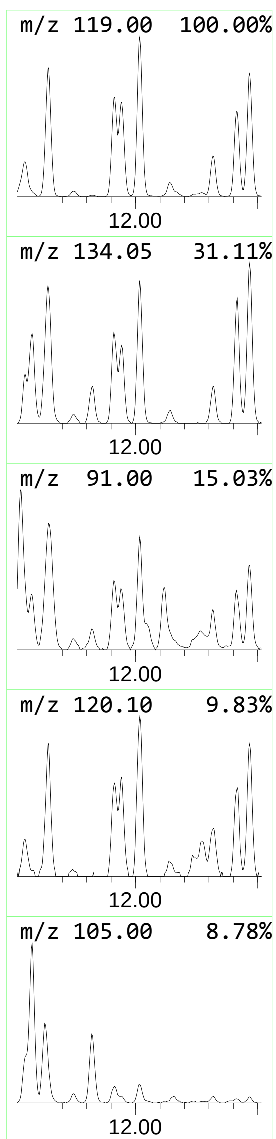
Peak Number 19 Benzene, 1-methyl-3-(1-meth... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.017	7.42 ug/L	193749	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	97
2	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	97
3	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	97
4	o-Cymene	134	C10H14	000527-84-4	97
5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96





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Sample : M2618-15
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG5W0

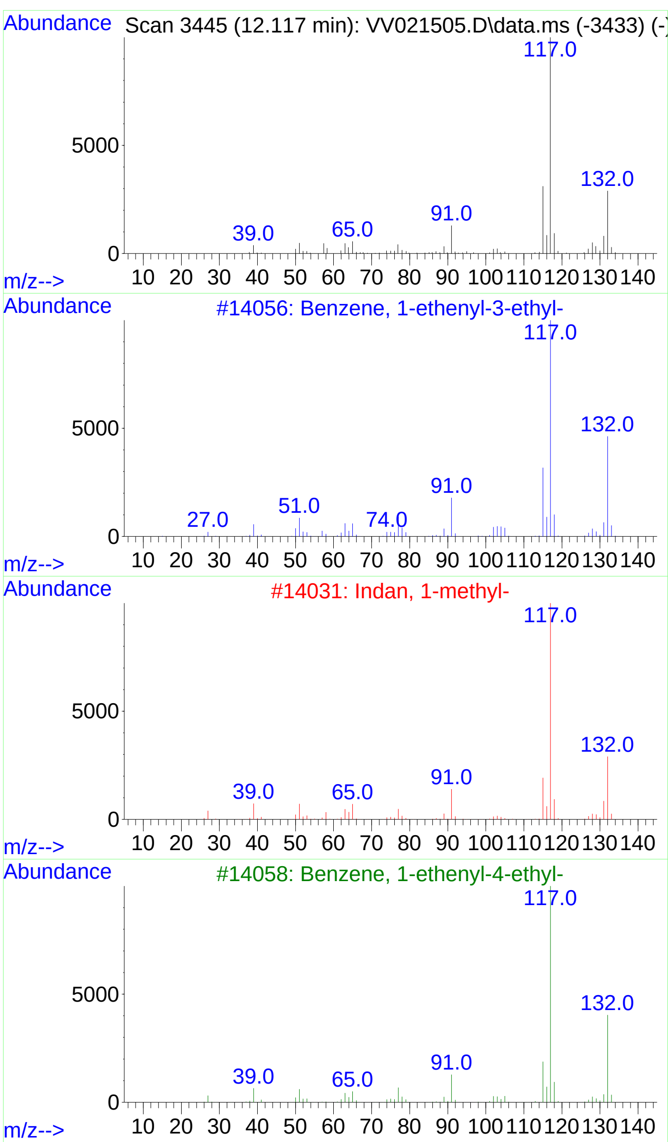
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM052721WMA.M
Quant Title : VOC Analysis

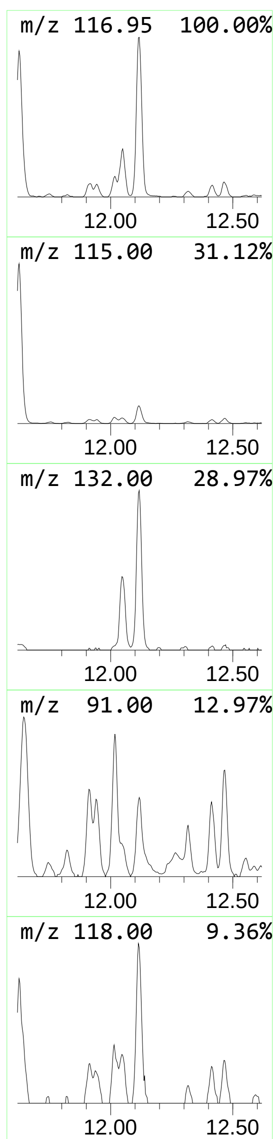
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.117	6.24 ug/L	162929	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	91
2	Indan, 1-methyl-	132	C10H12	000767-58-8	90
3	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87
4	1-Phenyl-1-butene	132	C10H12	000824-90-8	87
5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	87





Library Search Compound Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV060421\
Data File : VV021505.D
Acq On : 04 Jun 2021 14:32
Operator : SY/MD
Sample : M2618-15
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG5W0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM052721WMA.M
Quant Title : VOC Analysis

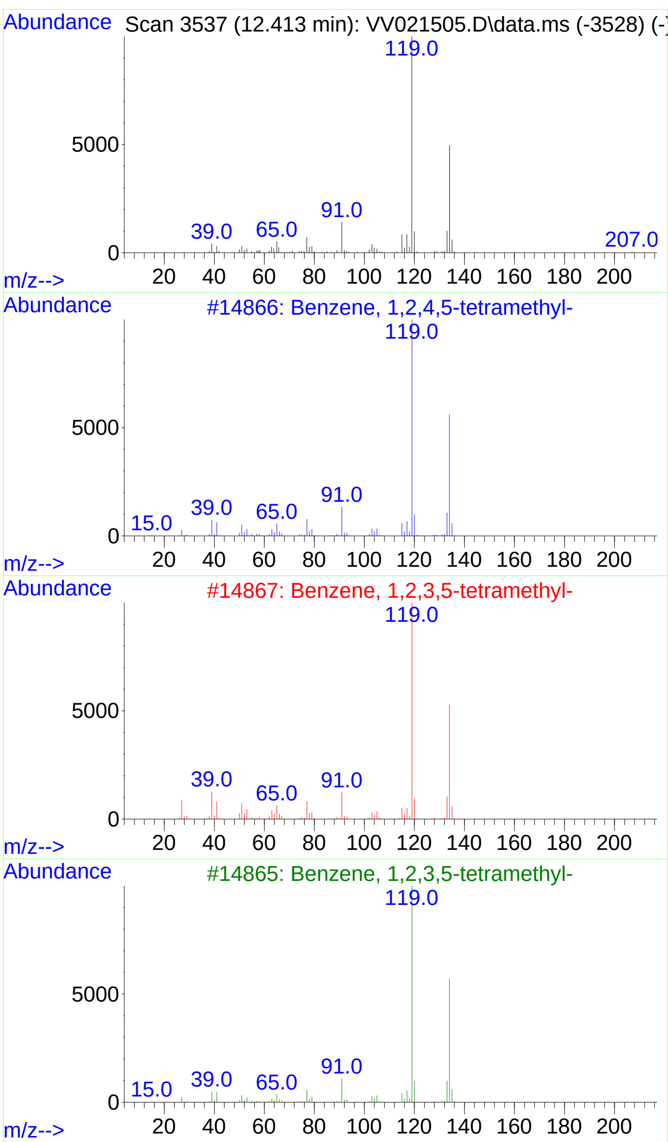
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

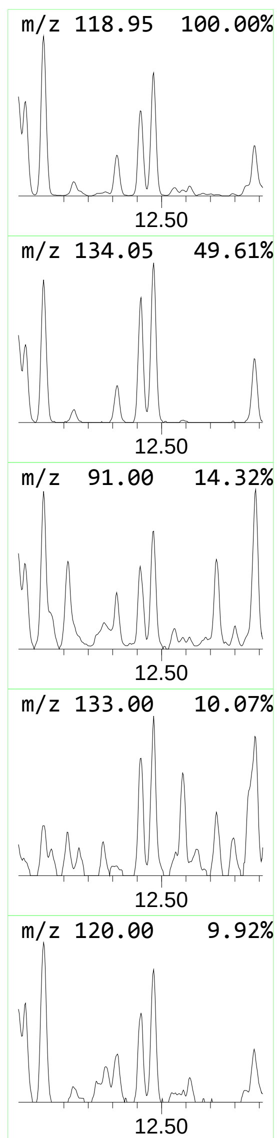
****

Peak Number 21 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.413	4.02 ug/L	104971	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
3	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
4	o-Cymene	134	C10H14	000527-84-4	94
5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94





Library Search Compound Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV060421\
Data File : VV021505.D
Acq On : 04 Jun 2021 14:32
Operator : SY/MD
Sample : M2618-15
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG5W0

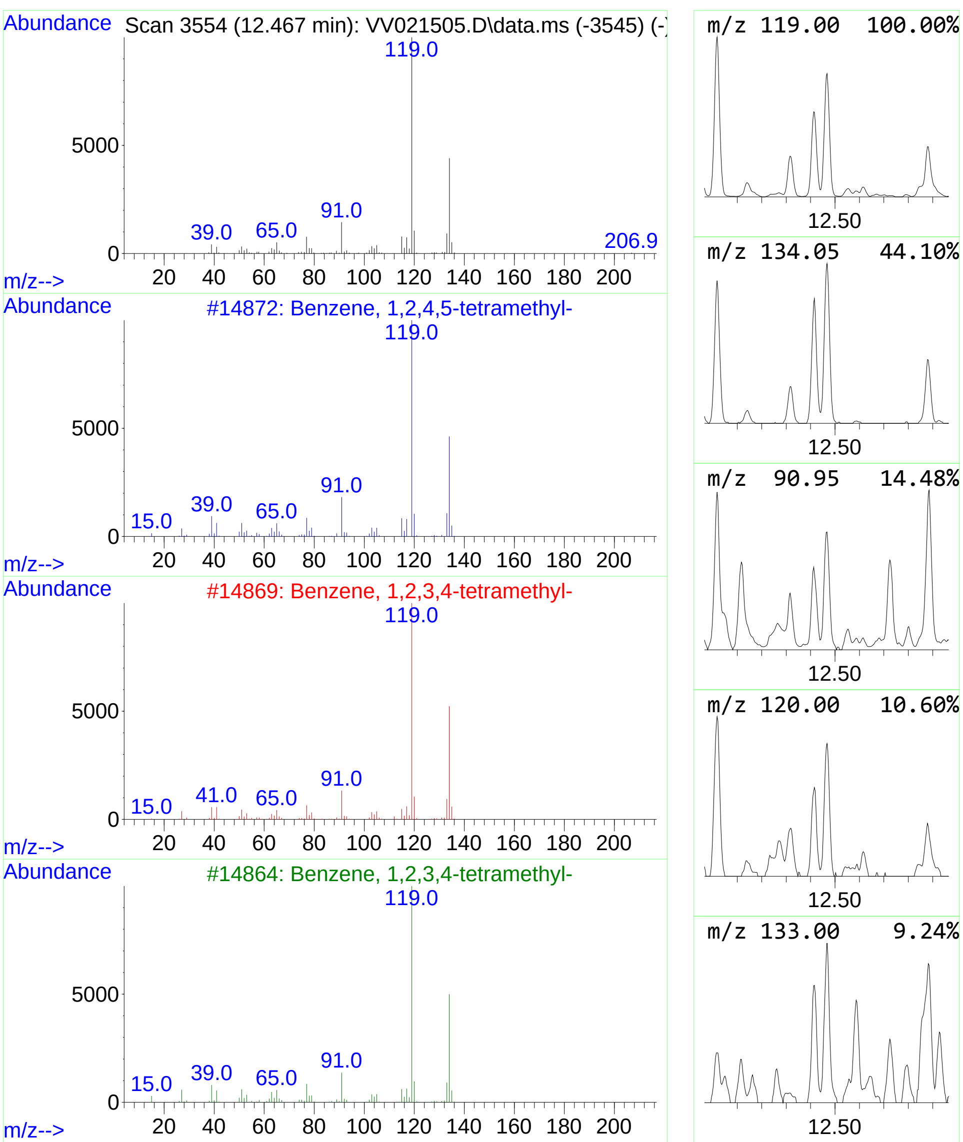
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM052721WMA.M
Quant Title : VOC Analysis

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

****
Peak Number 22 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.467	5.92 ug/L	154538	1,4-Dichlorobenzene-d4	11.252

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV060421\
Data File : VV021505.D
Acq On : 04 Jun 2021 14:32
Operator : SY/MD
Sample : M2618-15
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG5W0

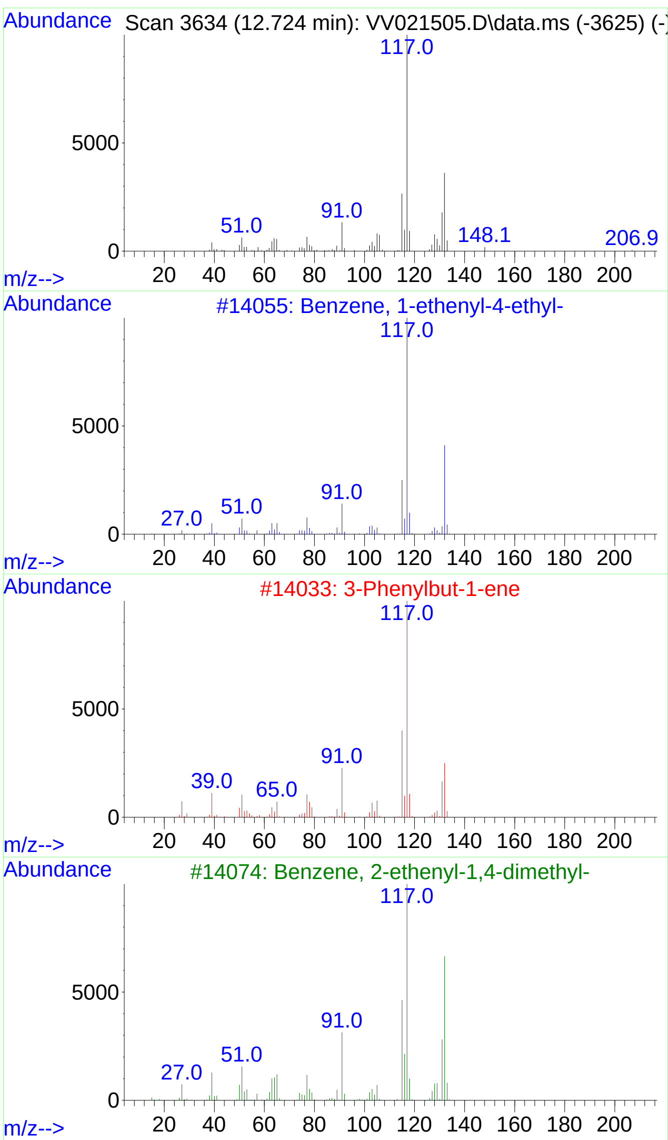
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM052721WMA.M
Quant Title : VOC Analysis

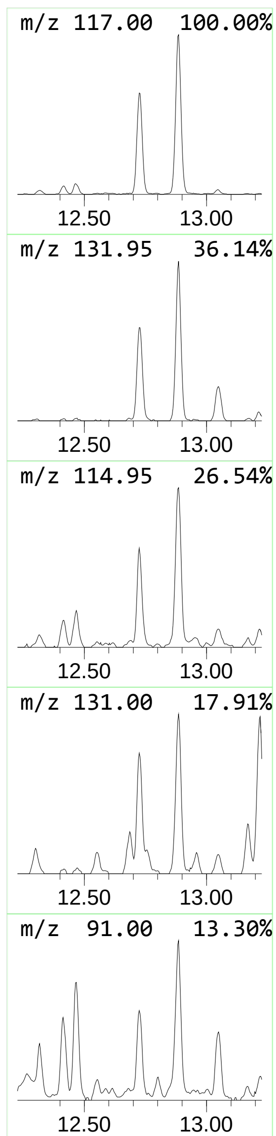
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.725	5.92 ug/L	154484	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
2	3-Phenylbut-1-ene	132	C10H12	000934-10-1	94
3	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92
4	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	91
5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV060421\
Data File : VV021505.D
Acq On : 04 Jun 2021 14:32
Operator : SY/MD
Sample : M2618-15
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG5W0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM052721WMA.M
Quant Title : VOC Analysis

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

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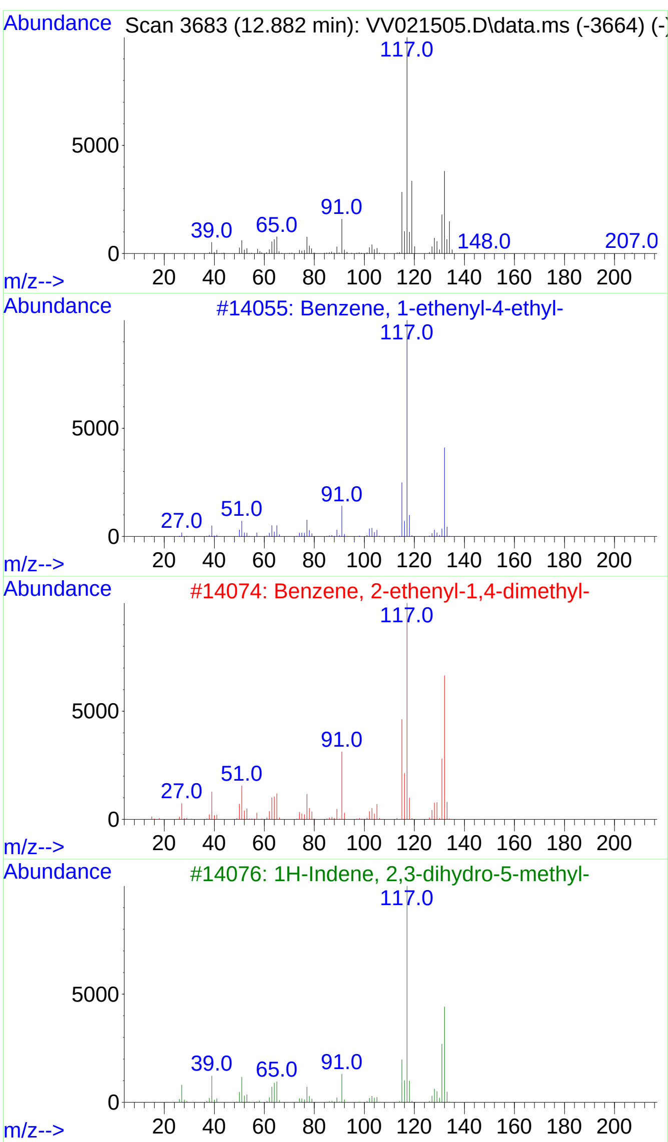
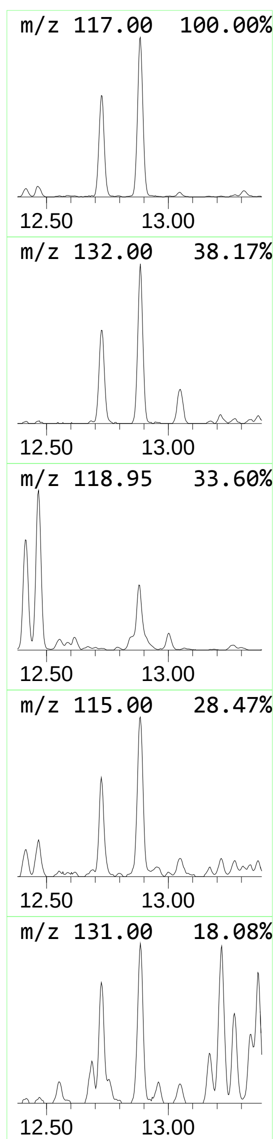
Peak Number 24 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.882	11.71 ug/L	305661	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
2	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	91
3	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	90
4	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	87
5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	87





Library Search Compound Report

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Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV060421\
Data File : VV021505.D
Acq On    : 04 Jun 2021  14:32
Operator  : SY/MD
Sample    : M2618-15
Misc      : 5.0mL/MSVOA_V/WATER
ALS Vial  : 14   Sample Multiplier: 1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM052721WMA.M
Quant Title  : VOC Analysis

TIC Library   : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P
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Instrument : MSVOA_V
ClientSampleId : BG5W0

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*********************************************************************
Peak Number 25  Azulene                        Concentration Rank 12
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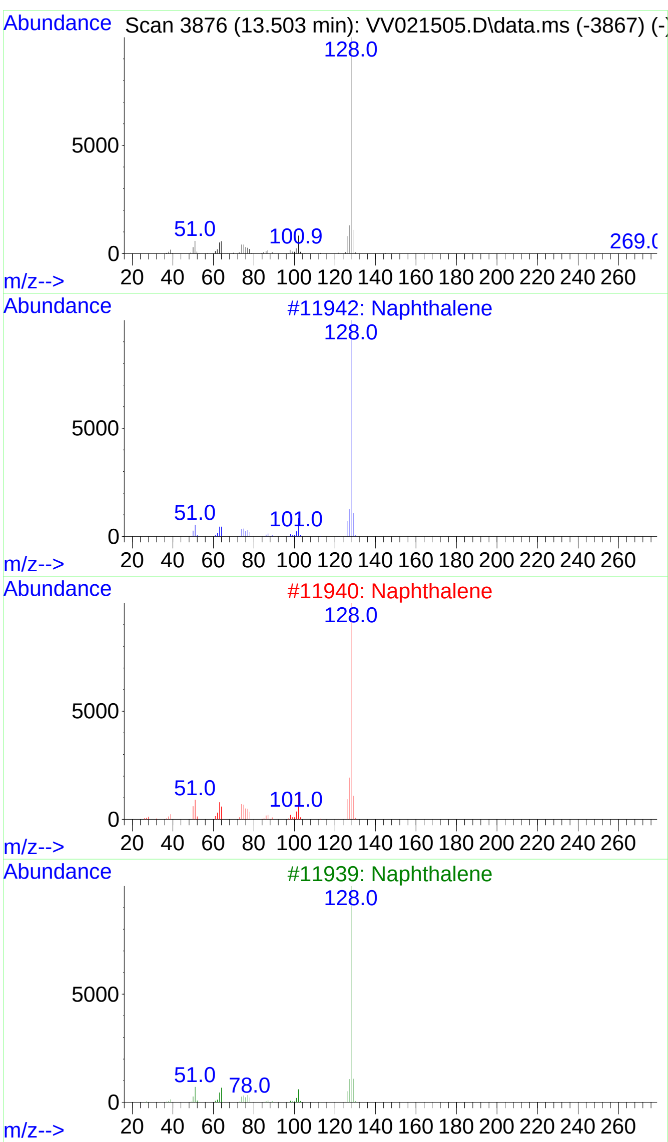
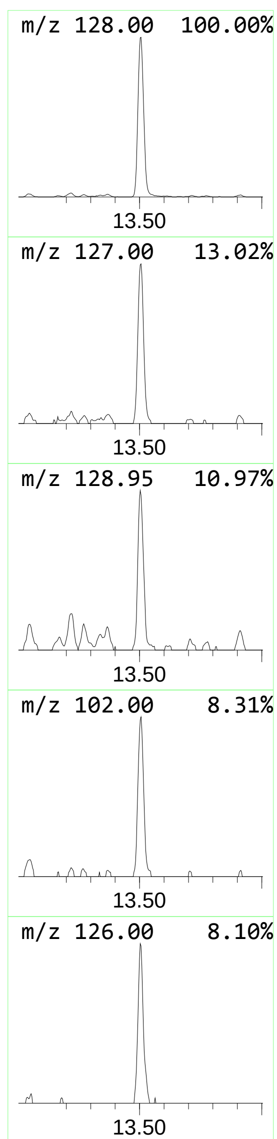
R.T.	EstConc	Area	Relative to ISTD	R.T.
13.503	5.69 ug/L	148613	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene	128	C10H8	000091-20-3	95
2	Naphthalene	128	C10H8	000091-20-3	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_V\Data\VV060421\
Data File : VV021505.D
Acq On : 04 Jun 2021 14:32
Operator : SY/MD
Sample : M2618-15
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG5W0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM052721WMA.M
Quant Title : VOC Analysis

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

TIC Top Hit name	RT	EstConc	Units	Response	Internal Standard #	Internal Standard RT	Internal Standard Resp	Internal Standard Conc
(DEL) Alkane: S...	1.632	4.5	ug/L	90383	1	5.619	1014720	50.0
(DEL) Alkane: S...	1.805	3.2	ug/L	64378	1	5.619	1014720	50.0
1-Butanol, 2,3-...	2.532	3.4	ug/L	68626	1	5.619	1014720	50.0
(DEL) Alkane: S...	2.763	3.4	ug/L	69886	1	5.619	1014720	50.0
(DEL) Alkane: S...	3.043	3.5	ug/L	71593	1	5.619	1014720	50.0
(DEL) Alkane: C...	3.767	7.8	ug/L	158144	1	5.619	1014720	50.0
(DEL) Alkane: S...	5.236	5.5	ug/L	111679	1	5.619	1014720	50.0
(DEL) Alkane: S...	6.657	3.0	ug/L	60054	1	5.619	1014720	50.0
(DEL) Alkane: S...	6.779	3.9	ug/L	79370	1	5.619	1014720	50.0
2-Pentanol, 4-m...	7.696	5.7	ug/L	134455	2	8.853	1188510	50.0
Benzene, propyl-	10.355	3.4	ug/L	89415	3	11.252	1304960	50.0
Benzene, 1-ethy...	10.448	46.4	ug/L	1210710	3	11.252	1304960	50.0
Benzene, 1-ethy...	10.737	14.9	ug/L	389252	3	11.252	1304960	50.0
Benzene, 1,2,3-...	11.336	16.1	ug/L	419302	3	11.252	1304960	50.0
Indane	11.528	20.6	ug/L	536848	3	11.252	1304960	50.0
Benzene, 2-ethy...	11.911	4.7	ug/L	122510	3	11.252	1304960	50.0
o-Cymene	11.943	3.5	ug/L	90053	3	11.252	1304960	50.0
Benzene, 1-meth...	12.017	7.4	ug/L	193749	3	11.252	1304960	50.0
Benzene, 1-ethe...	12.117	6.2	ug/L	162929	3	11.252	1304960	50.0
Benzene, 1,2,4,...	12.413	4.0	ug/L	104971	3	11.252	1304960	50.0
Benzene, 1,2,3,...	12.467	5.9	ug/L	154538	3	11.252	1304960	50.0
Benzene, 1-ethe...	12.725	5.9	ug/L	154484	3	11.252	1304960	50.0
Benzene, 2-ethe...	12.882	11.7	ug/L	305661	3	11.252	1304960	50.0
Azulene	13.503	5.7	ug/L	148613	3	11.252	1304960	50.0