

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV053019\
 Data File : VV011176.D
 Acq On : 31 May 2019 12:43
 Operator : SY/MD
 Sample : K3017-21
 Misc : 4.78G/10ML/MSVOA_V/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 A51F2

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\SOM2VLM053019S.M
 Title : VOC Analysis

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.110	3	6	21	rVB2	29793	30058	2.31%	0.241%
2	1.200	30	34	54	rVB4	23227	46506	3.57%	0.373%
3	1.309	61	68	77	rVB	196715	193575	14.87%	1.554%
4	1.380	84	90	99	rVB	97543	98322	7.55%	0.790%
5	1.557	138	145	153	rVB	143378	171939	13.21%	1.381%
6	2.113	308	318	329	rVV2	319580	464730	35.71%	3.732%
7	2.184	331	340	352	rVB	158541	232012	17.83%	1.863%
8	2.450	414	423	434	rVB	54010	84990	6.53%	0.682%
9	2.730	508	510	516	rBV7	1401	1366	0.10%	0.011%
10	2.984	580	589	600	rBV2	9274	15950	1.23%	0.128%
11	3.058	605	612	625	rVV5	8840	16527	1.27%	0.133%
12	3.196	645	655	666	rBV6	4989	11951	0.92%	0.096%
13	3.470	737	740	743	rBV4	1125	942	0.07%	0.008%
14	3.528	756	758	760	rBV2	1613	864	0.07%	0.007%
15	3.743	811	825	840	rBV2	23874	56461	4.34%	0.453%
16	3.820	847	849	854	rVB4	1213	890	0.07%	0.007%
17	3.862	854	862	865	rBV8	2043	2886	0.22%	0.023%
18	3.929	872	883	901	rBV	51580	131059	10.07%	1.052%
19	4.142	947	949	957	rVV6	1303	975	0.07%	0.008%
20	4.286	990	994	997	rBV4	1142	843	0.06%	0.007%
21	4.389	1008	1026	1048	rBV	206382	508887	39.10%	4.086%
22	4.714	1125	1127	1136	rVB9	1850	1838	0.14%	0.015%
23	4.801	1151	1154	1158	rVB4	1446	1085	0.08%	0.009%
24	5.000	1211	1216	1217	rBV4	1199	969	0.07%	0.008%
25	5.087	1225	1243	1271	rBV2	511670	1271102	97.66%	10.207%
26	5.225	1284	1286	1289	rBV3	1341	891	0.07%	0.007%
27	5.286	1303	1305	1308	rVB4	1438	884	0.07%	0.007%
28	5.370	1318	1331	1346	rBV2	15631	37820	2.91%	0.304%
29	5.531	1376	1381	1383	rBV6	3336	3733	0.29%	0.030%
30	5.656	1406	1420	1445	rBV	549679	1094949	84.13%	8.793%
31	5.833	1473	1475	1479	rVB5	1765	1168	0.09%	0.009%
32	5.952	1499	1512	1527	rBV8	14422	32661	2.51%	0.262%
33	6.113	1547	1562	1581	rVV	292868	587746	45.16%	4.720%
34	6.225	1594	1597	1600	rBV4	2301	2000	0.15%	0.016%

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35	6.325	1615	1628	1644	rBV	124723	242115	18.60%	1.944%
36	7.026	1826	1846	1868	rBV	173246	317277	24.38%	2.548%
37	7.180	1888	1894	1895	rBV5	4229	3937	0.30%	0.032%
38	7.357	1929	1949	1964	rBV	643333	1163704	89.41%	9.345%
39	7.608	2021	2027	2033	rVB9	3346	4750	0.36%	0.038%
40	7.659	2033	2043	2061	rVV2	111523	190786	14.66%	1.532%
41	7.749	2064	2071	2083	rVB9	6986	12998	1.00%	0.104%
42	7.871	2106	2109	2115	rBV6	973	920	0.07%	0.007%
43	7.974	2131	2141	2151	rBV5	8682	20386	1.57%	0.164%
44	8.132	2179	2190	2208	rBV	202420	364128	27.98%	2.924%
45	8.280	2224	2236	2258	rBV	719595	1180958	90.74%	9.483%
46	8.781	2389	2392	2397	rVB6	1292	1063	0.08%	0.009%
47	8.894	2407	2427	2450	rBV	774489	1301526	100.00%	10.452%
48	9.106	2485	2493	2495	rBV7	2799	3172	0.24%	0.025%
49	9.148	2504	2506	2508	rBV2	1458	868	0.07%	0.007%
50	9.177	2509	2515	2526	rVV2	9646	16867	1.30%	0.135%
51	9.260	2530	2541	2563	rVV2	13845	44357	3.41%	0.356%
52	9.354	2567	2570	2580	rVB7	5603	7465	0.57%	0.060%
53	9.469	2601	2606	2608	rBV4	1206	1118	0.09%	0.009%
54	9.527	2614	2624	2635	rBV4	13093	29312	2.25%	0.235%
55	9.749	2687	2693	2697	rVV2	4527	5369	0.41%	0.043%
56	9.794	2699	2707	2716	rVV2	28800	48026	3.69%	0.386%
57	9.849	2716	2724	2740	rVB2	52469	85342	6.56%	0.685%
58	10.177	2819	2826	2828	rVV8	2331	2677	0.21%	0.021%
59	10.225	2831	2841	2845	rVV2	80912	126833	9.74%	1.018%
60	10.260	2846	2852	2865	rVB	243971	375826	28.88%	3.018%
61	10.421	2893	2902	2915	rVB4	12854	23606	1.81%	0.190%
62	10.566	2935	2947	2958	rBV10	10936	26745	2.05%	0.215%
63	10.624	2958	2965	2982	rVB3	19754	36634	2.81%	0.294%
64	10.759	2999	3007	3015	rBV5	6569	9847	0.76%	0.079%
65	10.801	3017	3020	3021	rBV2	3173	1781	0.14%	0.014%
66	10.987	3059	3078	3088	rBV3	46997	86210	6.62%	0.692%
67	11.042	3091	3095	3101	rVB6	4293	5228	0.40%	0.042%
68	11.132	3120	3123	3126	rBV4	1726	1053	0.08%	0.008%
69	11.196	3134	3143	3150	rBV2	39854	66667	5.12%	0.535%
70	11.292	3162	3173	3190	rVB	506712	819823	62.99%	6.583%
71	11.575	3254	3261	3271	rBV8	7305	12757	0.98%	0.102%

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 Title : VOC Analysis

72	11.672	3279	3291	3303	rBV	345266	555756	42.70%	4.463%
73	11.823	3333	3338	3340	rBV5	2496	2126	0.16%	0.017%
74	12.055	3405	3410	3425	rVB9	11732	19237	1.48%	0.154%
75	12.296	3476	3485	3495	rBV2	7278	13277	1.02%	0.107%
76	12.392	3506	3515	3530	rVV3	25646	44131	3.39%	0.354%
77	12.588	3571	3576	3578	rBV4	2331	1920	0.15%	0.015%
78	12.717	3610	3616	3623	rBV5	7344	9375	0.72%	0.075%
79	12.788	3629	3638	3646	rVB5	5285	9122	0.70%	0.073%
80	12.929	3676	3682	3690	rVB5	4246	5503	0.42%	0.044%
81	13.048	3717	3719	3725	rVB7	1325	860	0.07%	0.007%
82	13.119	3739	3741	3747	rVB5	1897	1460	0.11%	0.012%
83	13.286	3787	3793	3798	rVB9	4057	4744	0.36%	0.038%
84	13.482	3849	3854	3855	rBV4	2286	2078	0.16%	0.017%
85	13.633	3896	3901	3902	rBV4	2614	2299	0.18%	0.018%
86	13.768	3939	3943	3945	rBV4	1640	1086	0.08%	0.009%
87	14.016	4016	4020	4022	rBV4	1736	1392	0.11%	0.011%
88	14.132	4054	4056	4062	rVB6	1650	1405	0.11%	0.011%
89	14.263	4094	4097	4100	rBV4	1751	1129	0.09%	0.009%
90	14.279	4100	4102	4108	rVB6	1704	1341	0.10%	0.011%
91	14.485	4164	4166	4170	rVB5	1789	1243	0.10%	0.010%
92	14.588	4193	4198	4200	rVV4	1503	1436	0.11%	0.012%
93	14.617	4203	4207	4211	rVB7	1475	1146	0.09%	0.009%
94	14.652	4211	4218	4220	rBV8	1152	1207	0.09%	0.010%
95	14.913	4296	4299	4300	rBV3	1336	932	0.07%	0.007%
96	15.061	4342	4345	4348	rBV4	1393	889	0.07%	0.007%
97	15.431	4454	4460	4470	rVB10	4882	8165	0.63%	0.066%
98	15.556	4492	4499	4504	rBV10	3621	4894	0.38%	0.039%
99	15.961	4621	4625	4629	rVB6	2056	1655	0.13%	0.013%
100	16.414	4764	4766	4770	rBV4	2566	2436	0.19%	0.020%

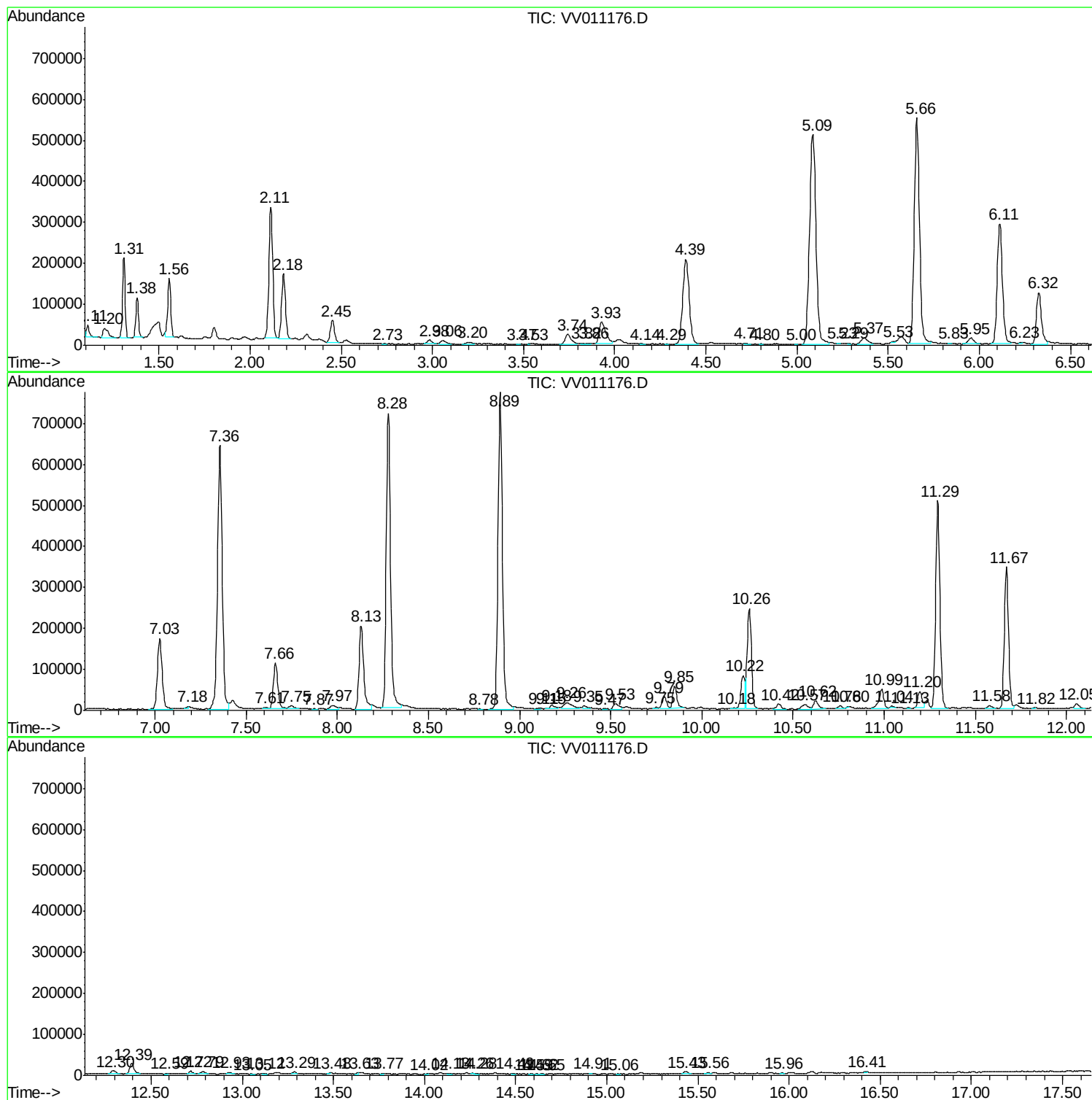
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_V\METHOD\SOM2VLM053019S.M
Quant Title : VOC Analysis

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



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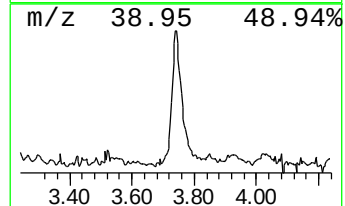
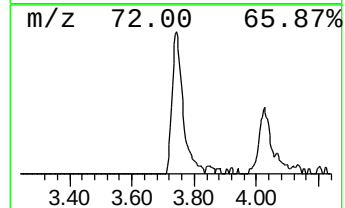
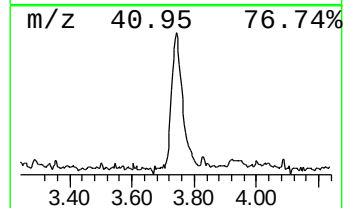
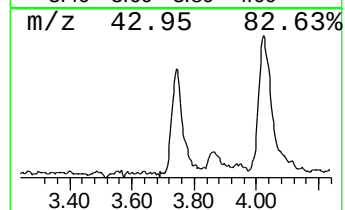
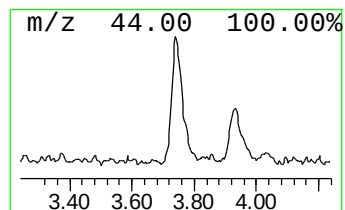
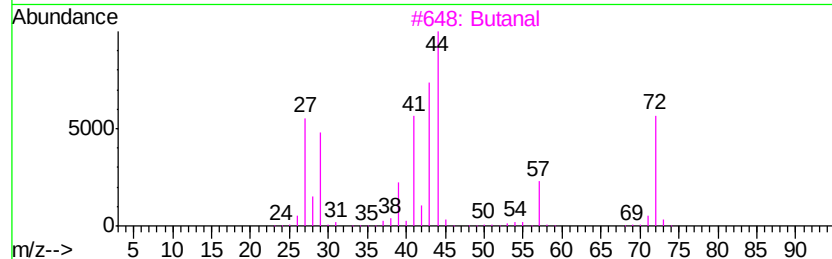
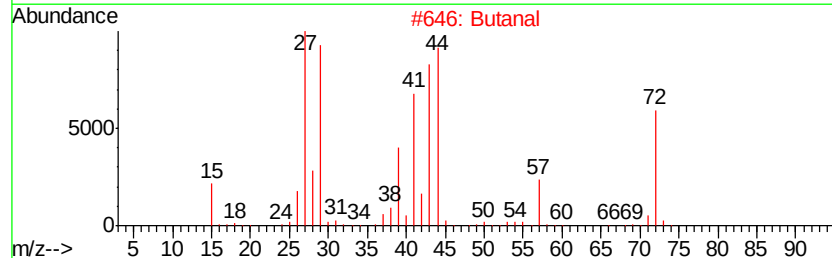
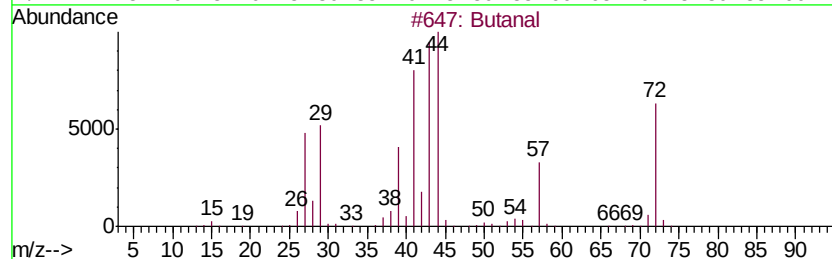
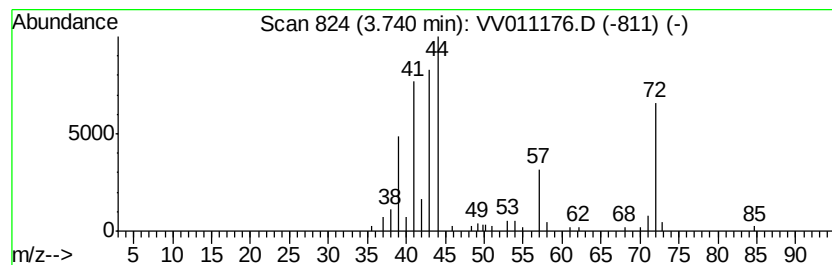
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_V\METHOD\SOM2VLM053019S.M
Quant Title : VOC Analysis

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

Peak Number 2 Butanal Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.74	1.29 ug/L	56461	1,4-Difluorobenzene	5.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butanal	72	C4H8O	000123-72-8	94
2		Butanal	72	C4H8O	000123-72-8	91
3		Butanal	72	C4H8O	000123-72-8	64
4		Butanal	72	C4H8O	000123-72-8	64
5		Ethene, ethoxy-	72	C4H8O	000109-92-2	58



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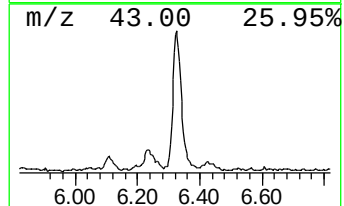
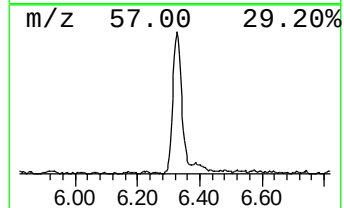
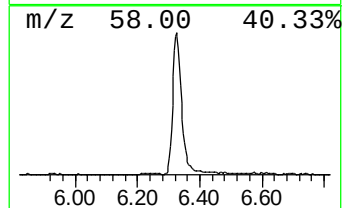
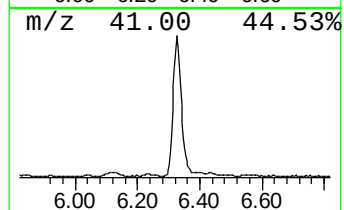
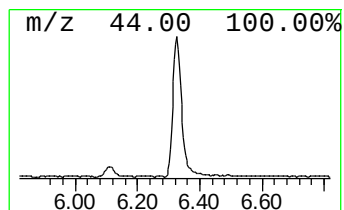
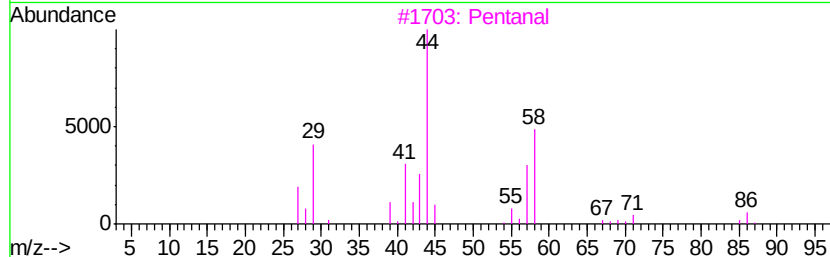
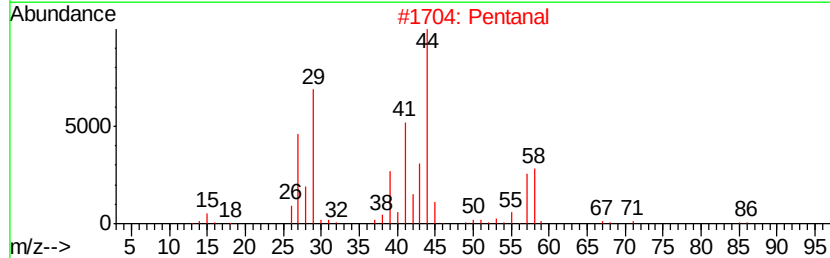
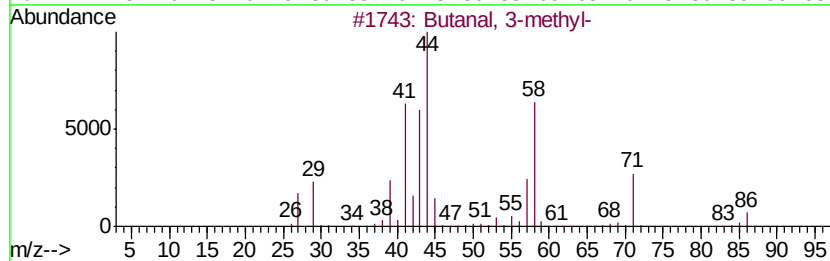
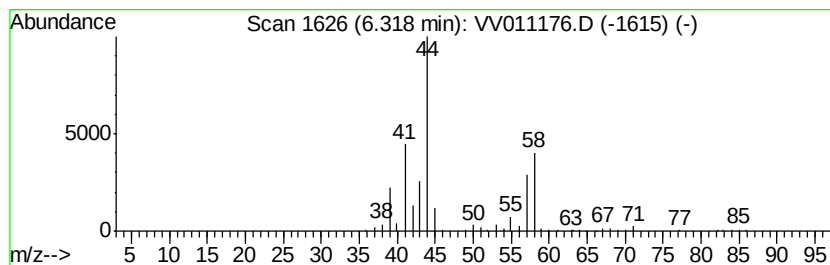
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Butanal, 3-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.32	5.53 ug/L	242115	1,4-Difluorobenzene	5.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butanal, 3-methyl-	86	C5H10O	000590-86-3	83
2		Pentanal	86	C5H10O	000110-62-3	80
3		Pentanal	86	C5H10O	000110-62-3	78
4		Butanal, 3-methyl-	86	C5H10O	000590-86-3	64
5		Pentanal	86	C5H10O	000110-62-3	58



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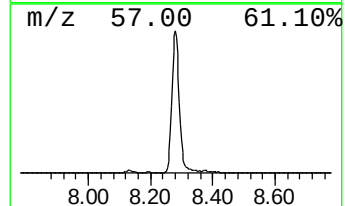
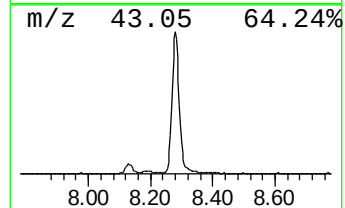
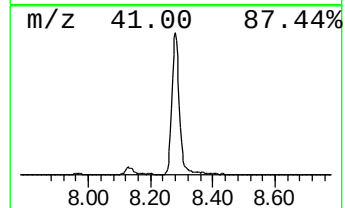
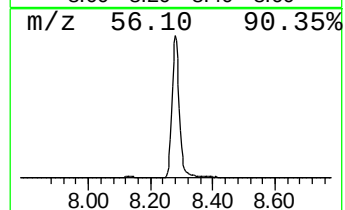
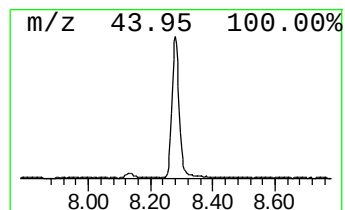
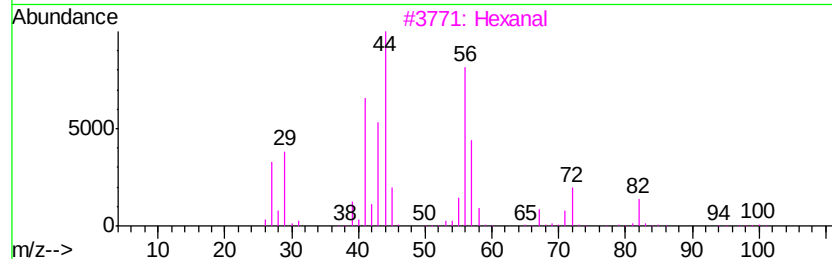
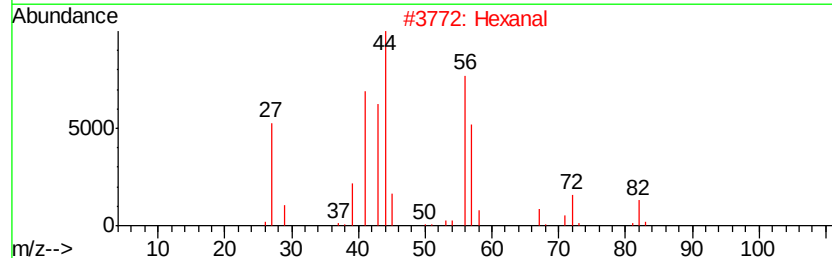
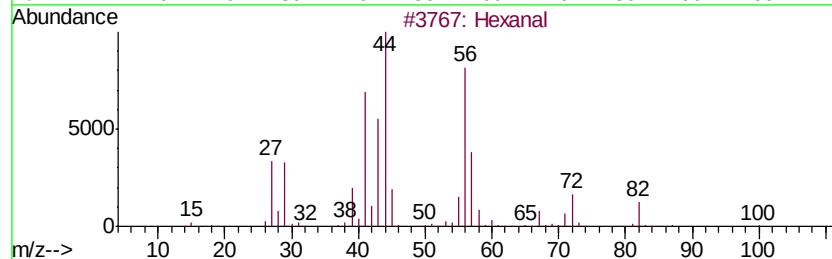
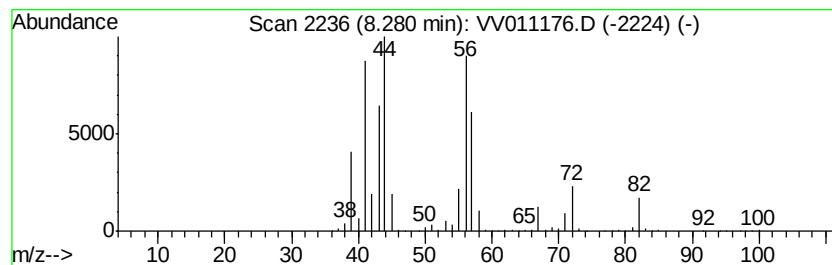
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Peak Number 5 Hexanal Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.28	22.68 ug/L	1180960	Chlorobenzene-d5	8.89

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanal		100	C6H12O	000066-25-1	81
2	Hexanal		100	C6H12O	000066-25-1	72
3	Hexanal		100	C6H12O	000066-25-1	72
4	Hexanal		100	C6H12O	000066-25-1	50
5	Glutaraldehyde		100	C5H8O2	000111-30-8	30



Data Path : Z:\V0ASRV\HPCHEM1\MSV0A_V\DATA\VV053019\
Data File : VV011176.D
Acq On : 31 May 2019 12:43
Operator : SY/MD
Sample : K3017-21
Misc : 4.78G/10ML/MSV0A_V/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
A51F2

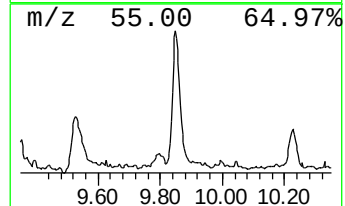
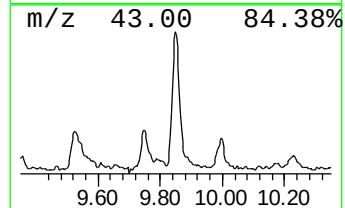
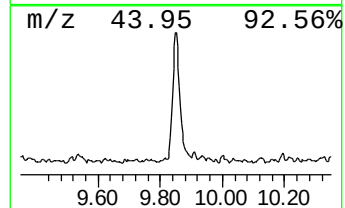
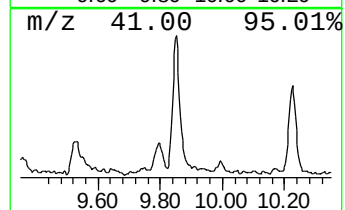
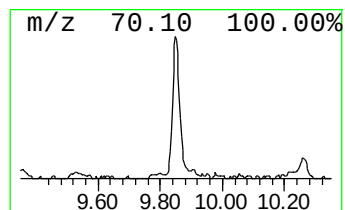
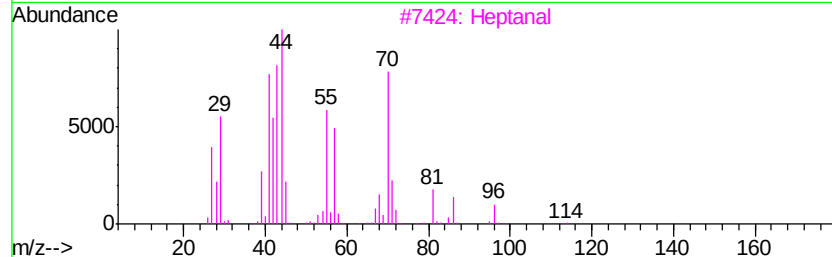
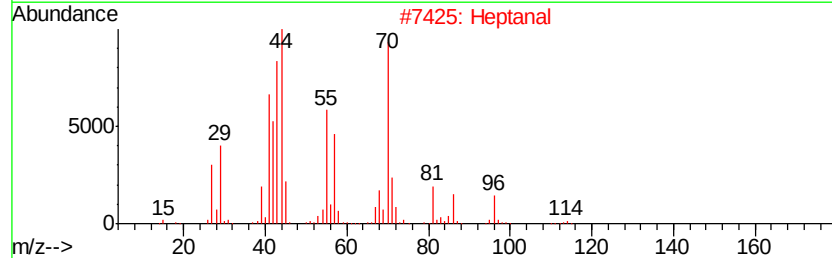
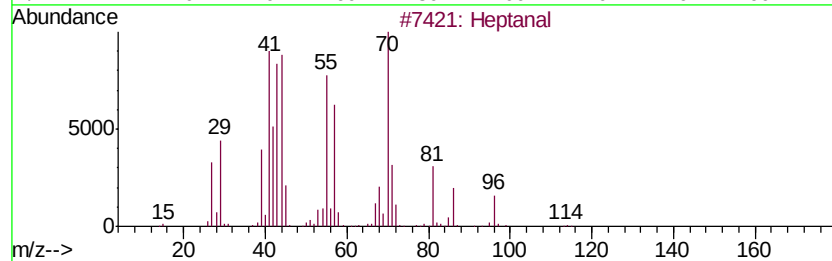
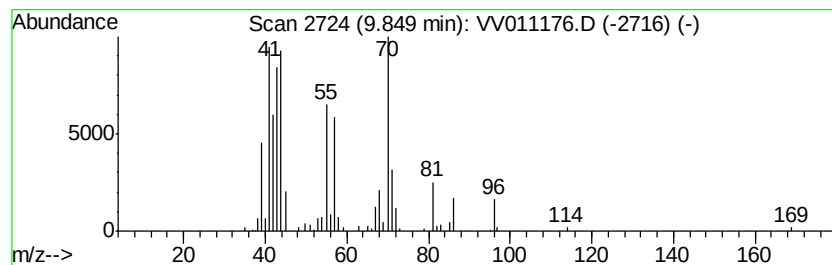
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Quant Title : VOC Analysis

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

Peak Number 6 Heptanal Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.85	1.64 ug/L	85342	Chlorobenzene-d5	8.89

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptanal		114	C7H14O	000111-71-7	98
2	Heptanal		114	C7H14O	000111-71-7	97
3	Heptanal		114	C7H14O	000111-71-7	90
4	Heptanal		114	C7H14O	000111-71-7	43
5	Heptanal		114	C7H14O	000111-71-7	43



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_V\DATA\VV053019\
Data File : VV011176.D
Acq On : 31 May 2019 12:43
Operator : SY/MD
Sample : K3017-21
Misc : 4.78G/10ML/MSVOA_V/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
A51F2

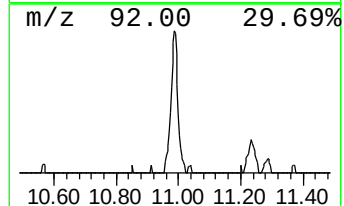
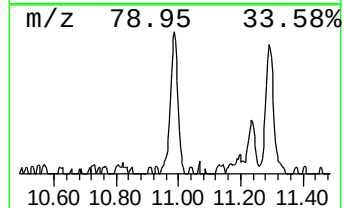
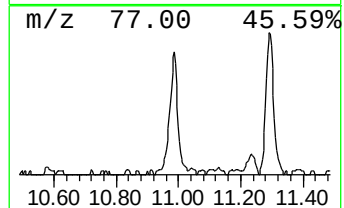
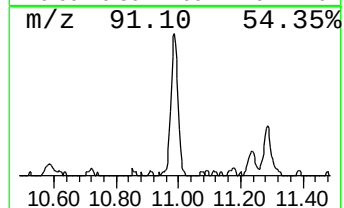
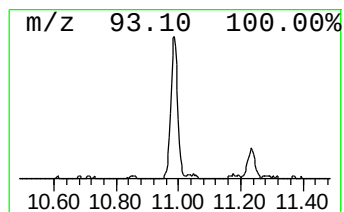
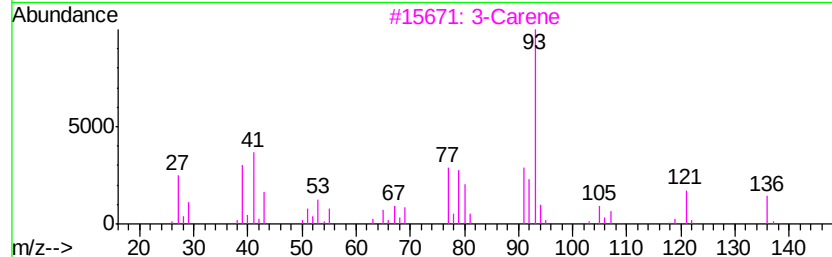
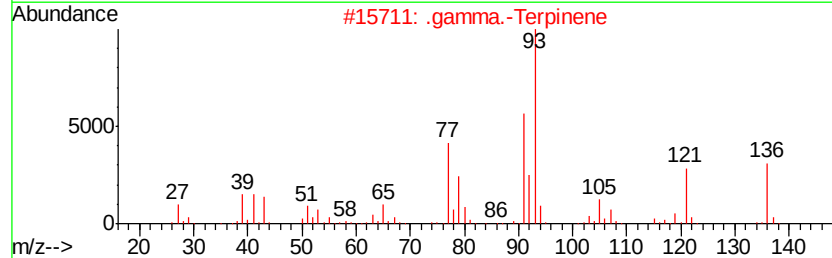
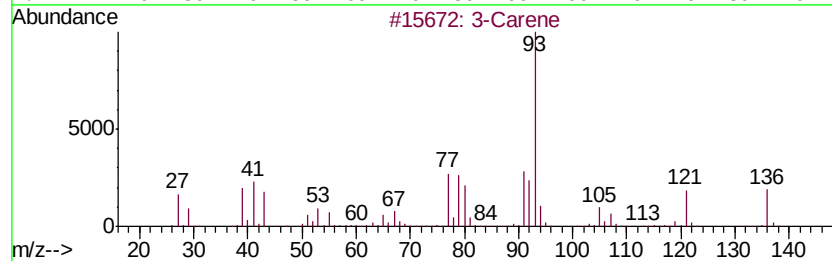
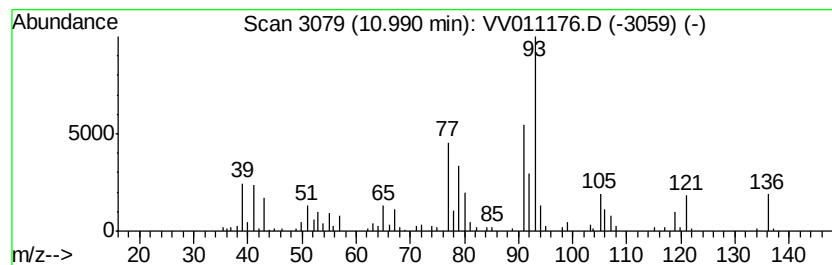
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Quant Title : VOC Analysis

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

Peak Number 7 3-Carene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.99	2.63 ug/L	86210	1,4-Dichlorobenzene-d4	11.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Carene	136	C10H16	013466-78-9	93
2		.gamma.-Terpinene	136	C10H16	000099-85-4	93
3		3-Carene	136	C10H16	013466-78-9	89
4		(1S)-2,6,6-Trimethylbicyclo[3.1....	136	C10H16	007785-26-4	87
5		1,3,7-Octatriene, 3,7-dimethyl-	136	C10H16	000502-99-8	83



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_V\DATA\VV053019\
Data File : VV011176.D
Acq On : 31 May 2019 12:43
Operator : SY/MD
Sample : K3017-21
Misc : 4.78G/10ML/MSVOA_V/SOIL
ALS Vial : 1 Sample Multiplier: 1

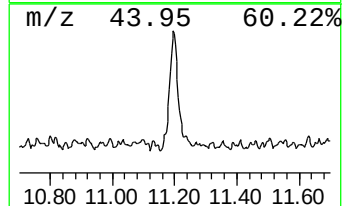
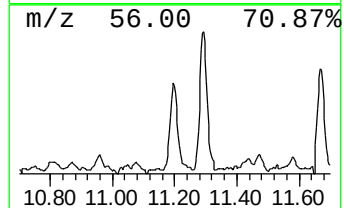
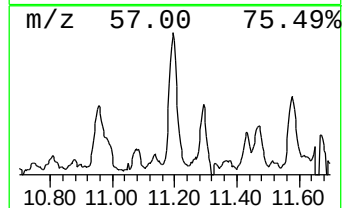
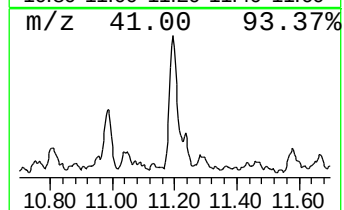
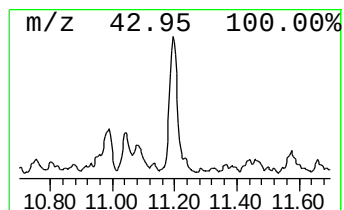
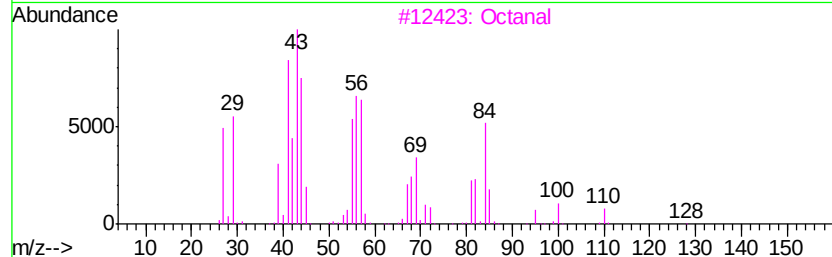
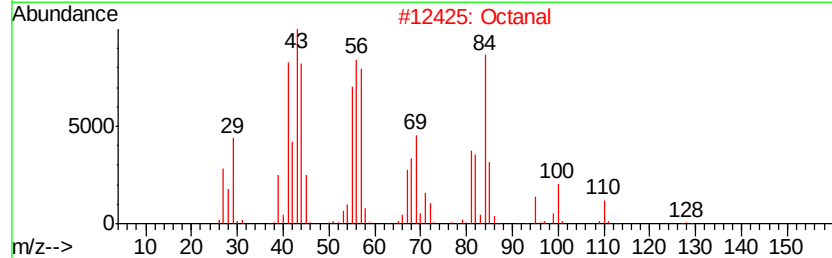
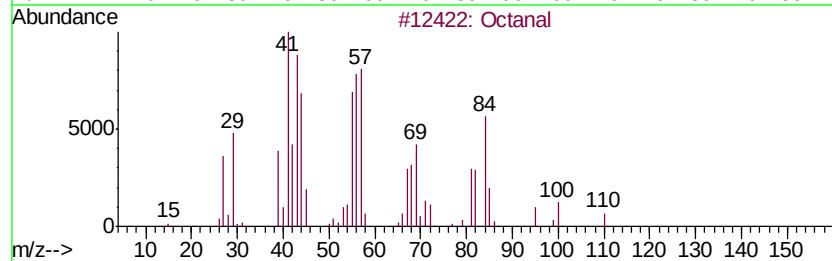
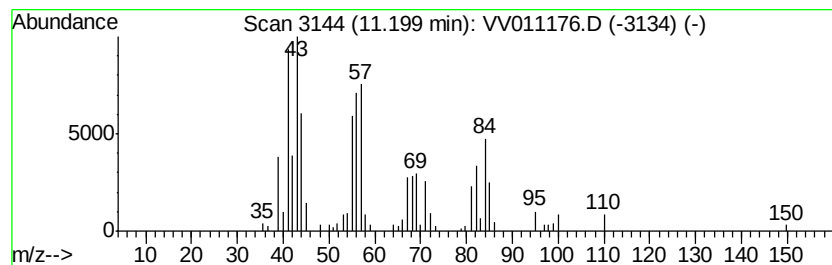
Instrument :
MSVOA_V
ClientSampleId :
A51F2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_V\METHOD\SOM2VLM053019S.M
Quant Title : VOC Analysis

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

Peak Number 8 Octanal Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.20	2.03 ug/L	66667	1,4-Dichlorobenzene-d4	11.29
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Octanal		128 C8H16O	000124-13-0 80
2	Octanal		128 C8H16O	000124-13-0 80
3	Octanal		128 C8H16O	000124-13-0 64
4	Octanal		128 C8H16O	000124-13-0 50
5	Cyclopentanol, 2-methyl-, trans-		100 C6H12O	025144-04-1 42



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_V\DATA\VV053019\
Data File : VV011176.D
Acq On : 31 May 2019 12:43
Operator : SY/MD
Sample : K3017-21
Misc : 4.78G/10ML/MSVOA_V/SOIL
ALS Vial : 1 Sample Multiplier: 1

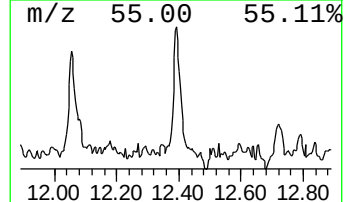
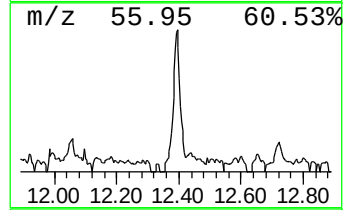
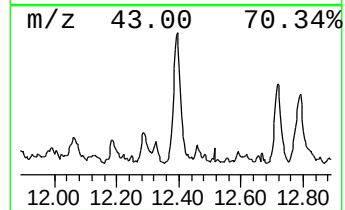
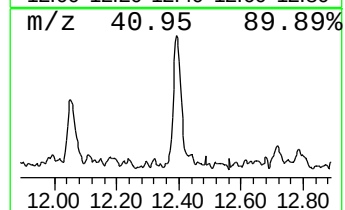
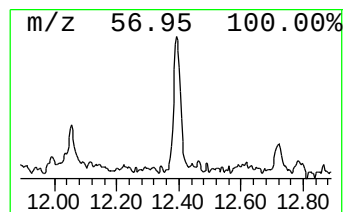
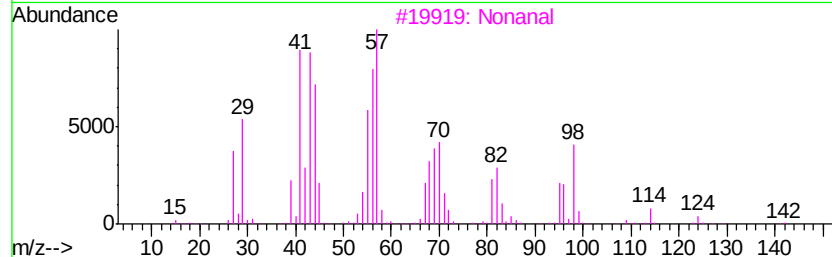
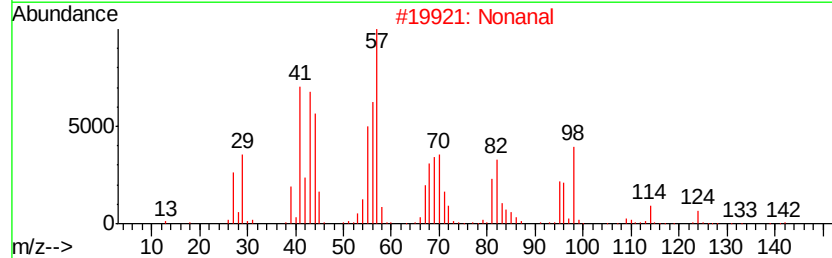
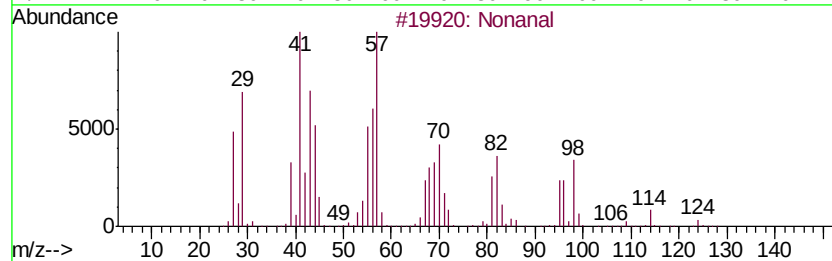
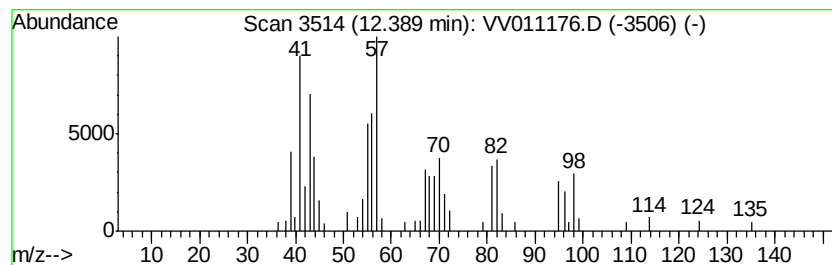
Instrument :
MSVOA_V
ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_V\METHOD\SOM2VLM053019S.M
Quant Title : VOC Analysis

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

Peak Number 9 Nonanal Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.39	1.35 ug/L	44131	1,4-Dichlorobenzene-d4	11.29
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Nonanal		142 C9H18O	000124-19-6 91
2	Nonanal		142 C9H18O	000124-19-6 83
3	Nonanal		142 C9H18O	000124-19-6 78
4	1,3-Cyclohexanediamine		114 C6H14N2	003385-21-5 46
5	Cyclohexanol, 4-methyl-, cis-		114 C7H14O	007731-28-4 25



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV053019\
Data File : VV011176.D
Acq On : 31 May 2019 12:43
Operator : SY/MD
Sample : K3017-21
Misc : 4.78G/10ML/MSVOA_V/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
A51F2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOM2VLM053019S.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--			
					#	RT	Resp	Conc
Butanal	3.74	1.3	ug/L	56461	1	5.66	1094950	25.0
Butanal, 3-methyl-	6.32	5.5	ug/L	242115	1	5.66	1094950	25.0
Hexanal	8.28	22.7	ug/L	1180960	2	8.89	1301530	25.0
Heptanal	9.85	1.6	ug/L	85342	2	8.89	1301530	25.0
3-Carene	10.99	2.6	ug/L	86210	3	11.29	819823	25.0
Octanal	11.20	2.0	ug/L	66667	3	11.29	819823	25.0
Nonanal	12.39	1.4	ug/L	44131	3	11.29	819823	25.0