

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU102618\
 Data File : VU027821.D
 Acq On : 26 Oct 2018 10:27
 Operator : MD/SY
 Sample : J5657-01ME
 Misc : 4.65g/5mL/100uL/5.0mL/MSVOA_U/MEOH
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 A40G4ME

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_U\METHOD\SOMULM102218WMA.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	1.395	62	69	87	rBV	29039	38405	1.06%	0.402%
2	1.598	107	132	133	rBV3	9886	23159	0.64%	0.243%
3	1.643	141	146	157	rVB	27351	35626	0.98%	0.373%
4	2.141	298	301	304	rBV3	817	723	0.02%	0.008%
5	2.234	318	330	344	rBV	56369	89837	2.47%	0.941%
6	2.829	506	515	528	rVB	943	2110	0.06%	0.022%
7	3.276	648	654	659	rBV2	602	728	0.02%	0.008%
8	3.813	815	821	827	rBV	199	348	0.01%	0.004%
9	4.199	927	941	991	rBV	18293	74943	2.06%	0.785%
10	4.643	1064	1079	1108	rVB	45912	117872	3.24%	1.235%
11	4.794	1121	1126	1128	rBV2	360	398	0.01%	0.004%
12	4.961	1168	1178	1180	rBV3	991	1519	0.04%	0.016%
13	5.331	1271	1293	1318	rBV2	104330	300322	8.26%	3.146%
14	5.884	1450	1465	1488	rVV	110654	235893	6.49%	2.471%
15	6.331	1591	1604	1626	rVV2	64890	141537	3.89%	1.483%
16	6.887	1770	1777	1780	rBV3	727	971	0.03%	0.010%
17	7.228	1872	1883	1904	rVB2	36381	73235	2.01%	0.767%
18	7.565	1971	1988	2005	rBV	128183	238072	6.55%	2.494%
19	7.694	2024	2028	2031	rVB5	569	432	0.01%	0.005%
20	7.855	2067	2078	2097	rVB2	24742	50284	1.38%	0.527%
21	8.183	2168	2180	2189	rBV2	13995	33469	0.92%	0.351%
22	8.327	2215	2225	2254	rVB	107175	201450	5.54%	2.110%
23	8.852	2381	2388	2397	rVB2	488	755	0.02%	0.008%
24	9.051	2445	2450	2451	rVV2	569	455	0.01%	0.005%
25	9.093	2451	2463	2483	rVB	174614	309109	8.50%	3.238%
26	9.205	2493	2498	2503	rVB2	290	371	0.01%	0.004%
27	9.253	2505	2513	2521	rBV2	1252	2023	0.06%	0.021%
28	9.379	2543	2552	2566	rBV2	5235	8497	0.23%	0.089%
29	9.784	2670	2678	2689	rBV4	3244	4548	0.13%	0.048%
30	9.958	2725	2732	2740	rVV2	934	1338	0.04%	0.014%
31	10.170	2787	2798	2811	rVB2	5947	9907	0.27%	0.104%
32	10.437	2869	2881	2897	rVB	105611	162426	4.47%	1.701%
33	10.614	2926	2936	2953	rVB2	7758	15751	0.43%	0.165%
34	10.687	2953	2959	2965	rBV2	401	590	0.02%	0.006%

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 ClientSampleId :
 A40G4ME

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

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Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM102218WMA.M
 Title : VOC Analysis

35	10.739	2970	2975	2977	rBV	668	561	0.02%	0.006%
36	10.778	2978	2987	2995	rVV	11432	15901	0.44%	0.167%
37	10.874	3012	3017	3024	rBV2	546	733	0.02%	0.008%
38	11.150	3095	3103	3113	rVB2	10330	14711	0.40%	0.154%
39	11.279	3136	3143	3148	rVB2	706	948	0.03%	0.010%
40	11.324	3152	3157	3161	rVB3	349	414	0.01%	0.004%
41	11.421	3169	3187	3197	rBV	64766	106672	2.93%	1.117%
42	11.507	3197	3214	3229	rVV2	495246	1082499	29.78%	11.339%
43	11.633	3242	3253	3258	rBV4	4126	6730	0.19%	0.070%
44	11.668	3259	3264	3273	rVB3	5326	7098	0.20%	0.074%
45	11.755	3283	3291	3294	rBV5	1186	1883	0.05%	0.020%
46	11.864	3313	3325	3349	rVB2	174237	307645	8.46%	3.222%
47	11.964	3351	3356	3362	rVB4	849	1096	0.03%	0.011%
48	12.025	3367	3375	3383	rVB2	3962	5537	0.15%	0.058%
49	12.150	3407	3414	3417	rBV4	797	1042	0.03%	0.011%
50	12.253	3439	3446	3454	rVV3	5355	7390	0.20%	0.077%
51	12.337	3466	3472	3485	rVB2	3783	5928	0.16%	0.062%
52	12.488	3506	3519	3535	rVV4	4690	11367	0.31%	0.119%
53	12.652	3563	3570	3577	rBV3	2548	3645	0.10%	0.038%
54	12.707	3579	3587	3600	rVV4	3759	6491	0.18%	0.068%
55	12.842	3621	3629	3642	rVB4	12089	21267	0.59%	0.223%
56	12.970	3661	3669	3676	rVB3	1956	2988	0.08%	0.031%
57	13.006	3676	3680	3681	rBV	568	440	0.01%	0.005%
58	13.028	3681	3687	3689	rBV4	972	1059	0.03%	0.011%
59	13.125	3710	3717	3732	rBV5	4799	8329	0.23%	0.087%
60	13.392	3794	3800	3809	rVV5	855	1302	0.04%	0.014%
61	13.507	3827	3836	3849	rVB3	9476	14846	0.41%	0.156%
62	13.642	3872	3878	3881	rBV2	289	349	0.01%	0.004%
63	13.681	3885	3890	3894	rVV2	483	452	0.01%	0.005%
64	13.752	3903	3912	3919	rBV2	2061	3374	0.09%	0.035%
65	13.861	3943	3946	3947	rBV2	722	367	0.01%	0.004%
66	13.993	3980	3987	3994	rVB5	1373	1997	0.05%	0.021%
67	14.112	4015	4024	4030	rBV2	6671	10217	0.28%	0.107%
68	14.228	4053	4060	4061	rBV4	813	773	0.02%	0.008%
69	14.334	4086	4093	4101	rBV6	2341	3343	0.09%	0.035%
70	14.379	4104	4107	4116	rVB3	1270	1452	0.04%	0.015%
71	14.440	4119	4126	4130	rBV3	2704	3627	0.10%	0.038%

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Instrument :
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 ClientSampleId :
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Integration Parameters: LSCINT.P

Integrator: RTE
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 Title : VOC Analysis

72	14.475	4131	4137	4142	rBV4	4284	4927	0.14%	0.052%
73	14.581	4161	4170	4172	rBV3	2780	3326	0.09%	0.035%
74	14.613	4174	4180	4185	rVV5	2690	4435	0.12%	0.046%
75	14.707	4201	4209	4211	rBV5	1852	2503	0.07%	0.026%
76	14.787	4220	4234	4240	rBV	121612	175435	4.83%	1.838%
77	14.832	4240	4248	4260	rVV	218347	331442	9.12%	3.472%
78	14.887	4261	4265	4276	rVB2	33291	44056	1.21%	0.461%
79	14.961	4281	4288	4293	rBV5	2768	3442	0.09%	0.036%
80	15.009	4293	4303	4311	rBV	231743	386381	10.63%	4.047%
81	15.060	4312	4319	4324	rBV	346123	444015	12.21%	4.651%
82	15.147	4338	4346	4350	rBV	33255	45030	1.24%	0.472%
83	15.179	4351	4356	4369	rVB	57568	78409	2.16%	0.821%
84	15.282	4380	4388	4395	rBV3	7862	10643	0.29%	0.111%
85	15.459	4437	4443	4451	rVB4	1299	1913	0.05%	0.020%
86	15.514	4453	4460	4461	rBV4	908	790	0.02%	0.008%
87	15.607	4479	4489	4503	rBV	63559	98432	2.71%	1.031%
88	15.880	4566	4574	4585	rVB2	5485	8676	0.24%	0.091%
89	15.932	4585	4590	4598	rBV4	1385	2007	0.06%	0.021%
90	15.977	4598	4604	4609	rBV4	1761	2046	0.06%	0.021%
91	16.089	4630	4639	4653	rVB4	3454	6233	0.17%	0.065%
92	16.215	4672	4678	4690	rVB5	5365	8318	0.23%	0.087%
93	16.388	4724	4732	4741	rBV3	4321	6542	0.18%	0.069%
94	16.481	4753	4761	4769	rVB3	4491	6182	0.17%	0.065%
95	16.555	4769	4784	4800	rBV	2440686	3635218	100.00%	38.077%
96	16.633	4800	4808	4818	rVV2	10314	18507	0.51%	0.194%
97	16.703	4819	4830	4843	rVV	197982	299751	8.25%	3.140%
98	16.764	4847	4849	4853	rVB3	806	417	0.01%	0.004%
99	17.276	5001	5008	5020	rVB4	13213	20614	0.57%	0.216%
100	17.485	5064	5073	5086	rVB3	30670	55824	1.54%	0.585%

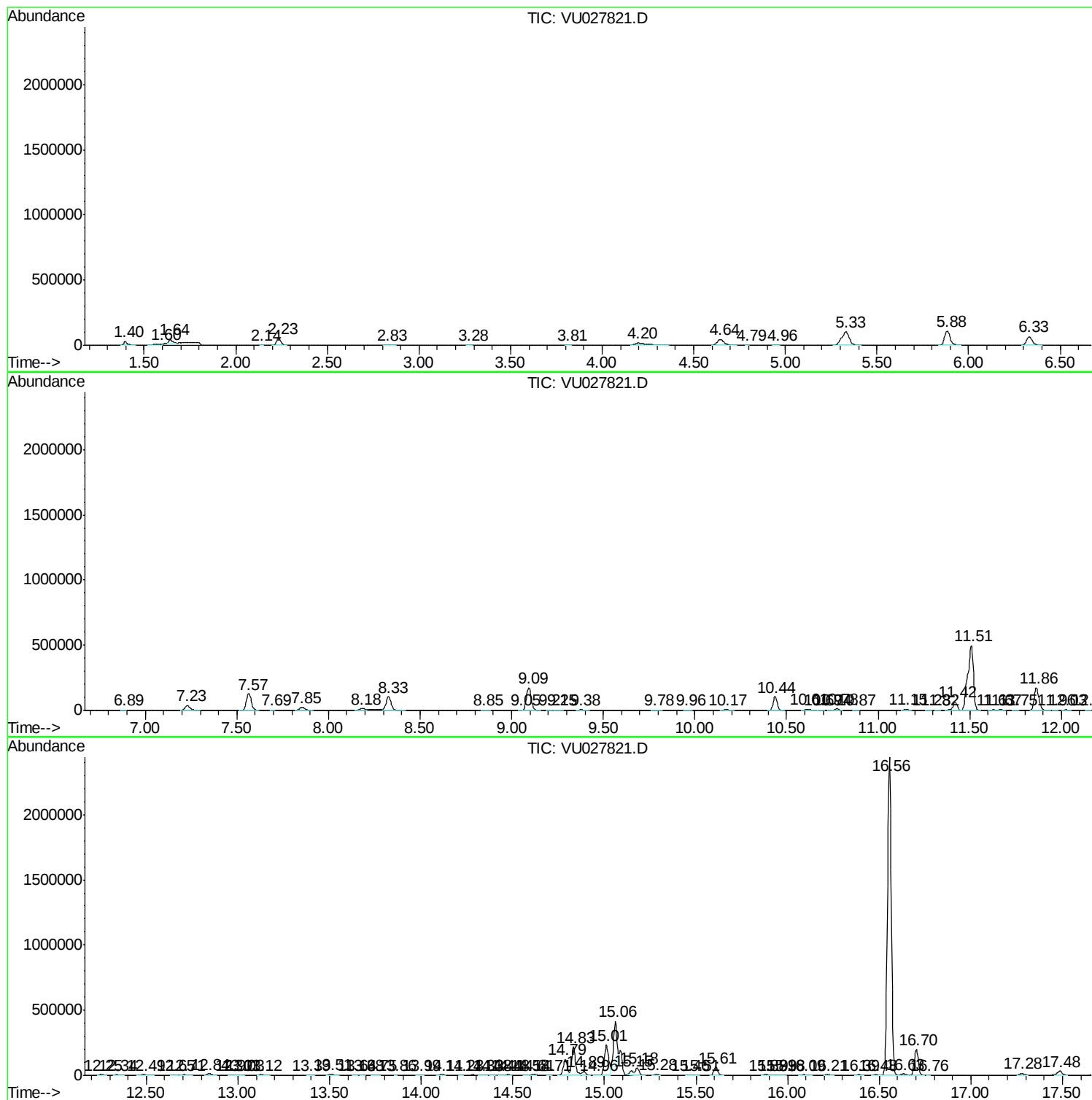
Sum of corrected areas: 9547090

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ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
A40G4ME

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM102218WMA.M
Quant Title : VOC Analysis

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



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Sample : J5657-01ME
Misc : 4.65µ/5mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 5 Sample Multiplier: 1

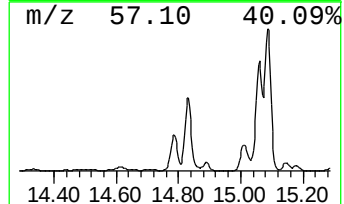
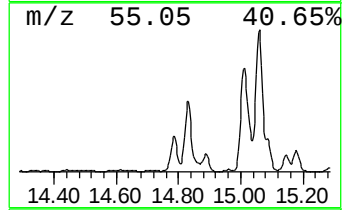
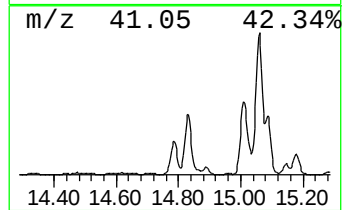
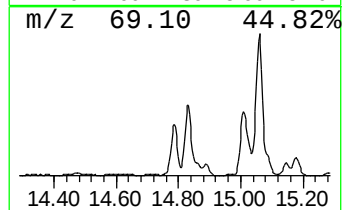
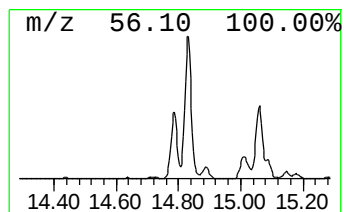
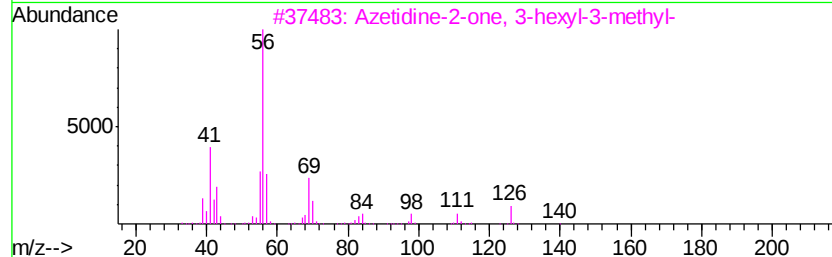
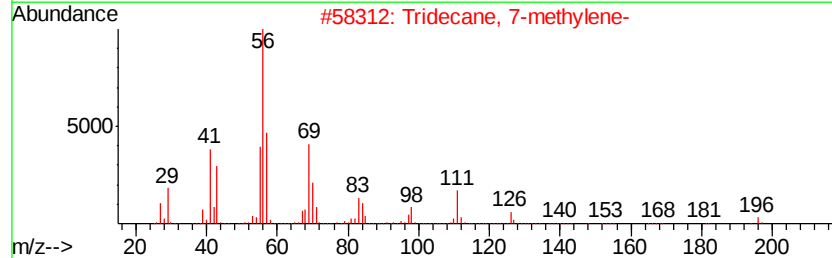
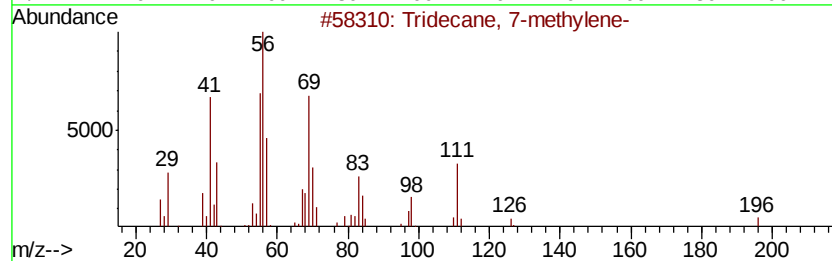
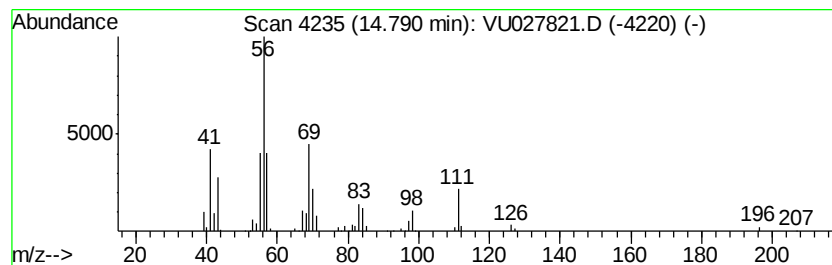
Instrument :
MSVOA_U
ClientSampleId :
A40G4ME

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM102218WMA.M
Quant Title : VOC Analysis

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

Peak Number 3 (DEL) Alkane: Cyclic14.79 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.79	8.10 ug/L	175435	1,4-Dichlorobenzene-d4	11.49	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane, 7-methylene-	196	C14H28	019780-80-4	95
2	Tridecane, 7-methylene-	196	C14H28	019780-80-4	91
3	Azetidine-2-one, 3-hexyl-3-methyl-	169	C10H19NO	105037-97-6	78
4	1-Octene, 7-methyl-	126	C9H18	013151-06-9	68
5	1-Pentene, 2,4-dimethyl-	98	C7H14	002213-32-3	58



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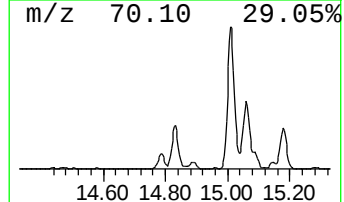
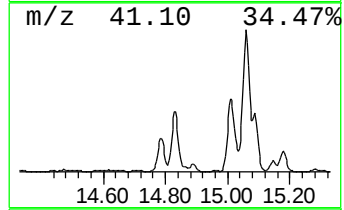
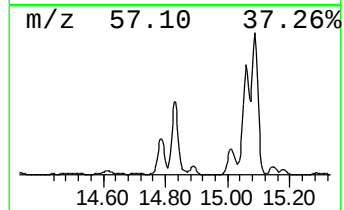
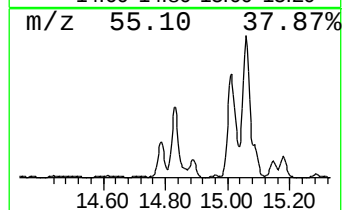
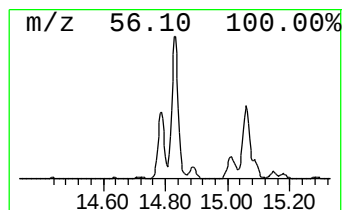
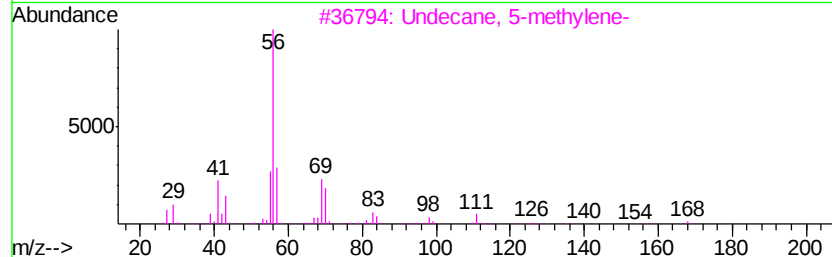
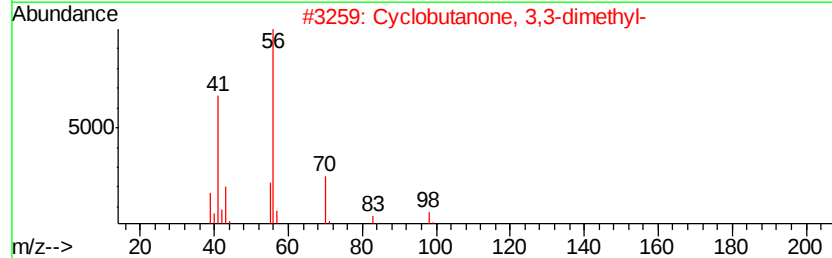
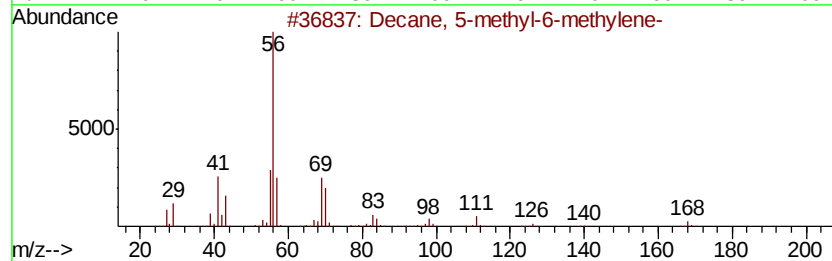
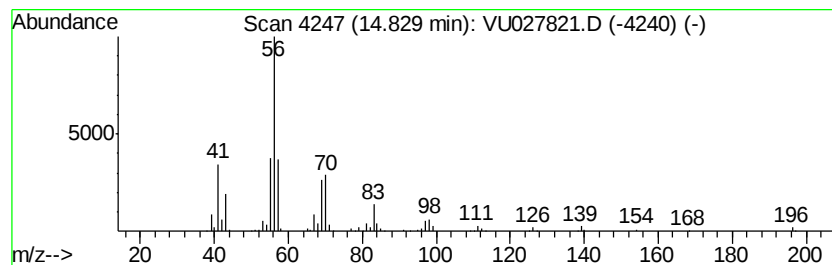
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM102218WMA.M
Quant Title : VOC Analysis

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

Peak Number 4 (DEL) Alkane: Cyclic14.83 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.83	15.31 ug/L	331442	1,4-Dichlorobenzene-d4	11.49

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Decane, 5-methyl-6-methylene-	168	C12H24	075029-95-7	78
2		Cyclobutanone, 3,3-dimethyl-	98	C6H10O	001192-33-2	72
3		Undecane, 5-methylene-	168	C12H24	005698-48-6	72
4		1-Pentene, 2,4-dimethyl-	98	C7H14	002213-32-3	72
5		1-Heptene, 2,6-dimethyl-	126	C9H18	003074-78-0	72



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ClientSampleId :
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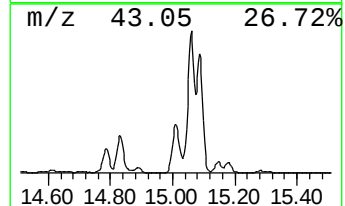
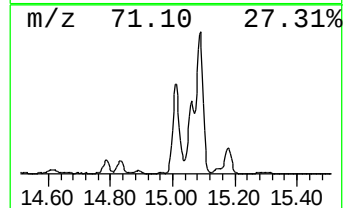
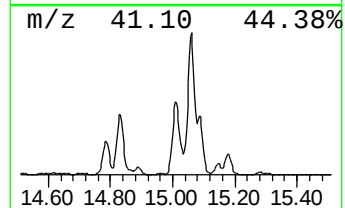
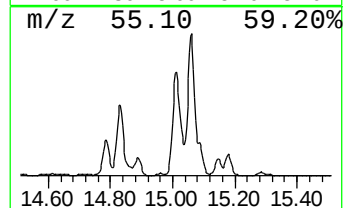
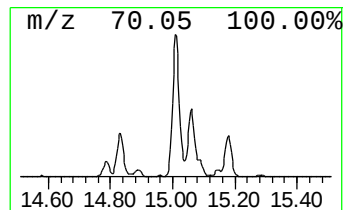
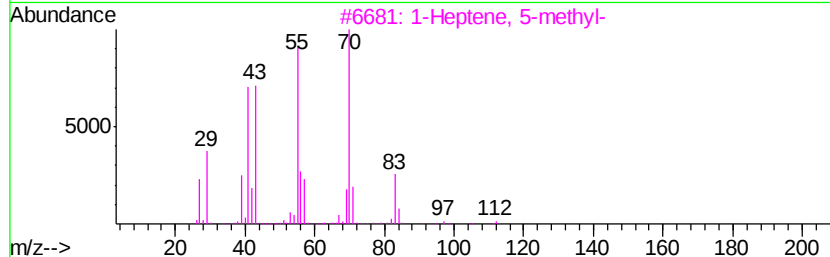
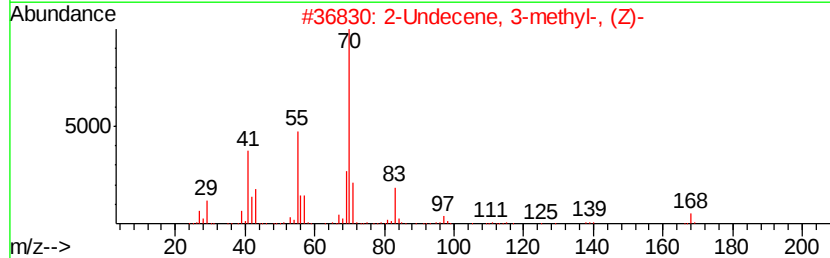
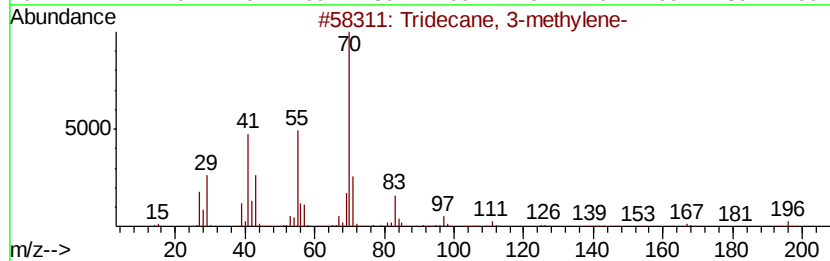
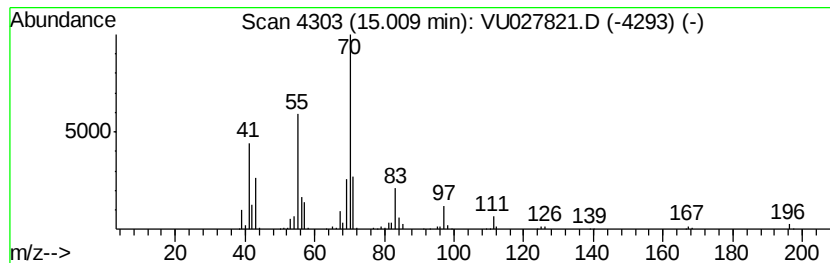
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Quant Title : VOC Analysis

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

Peak Number 5 (DEL) Alkane: Cyclic15.01 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.01	17.85 ug/L	386381	1,4-Dichlorobenzene-d4	11.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane, 3-methylene-	196	C14H28	019780-34-8	87
2	2-Undecene, 3-methyl-, (Z)-	168	C12H24	057024-90-5	78
3	1-Heptene, 5-methyl-	112	C8H16	013151-04-7	46
4	4-Tetradecene, (E)-	196	C14H28	041446-78-0	30
5	5-Tetradecene, (E)-	196	C14H28	041446-66-6	30



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU102618\
Data File : VU027821.D
Acq On : 26 Oct 2018 10:27
Operator : MD/SY
Sample : J5657-01ME
Misc : 4.65u/5mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
A40G4ME

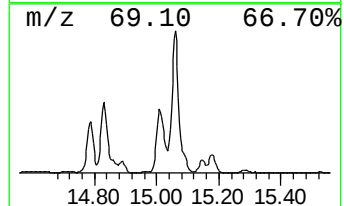
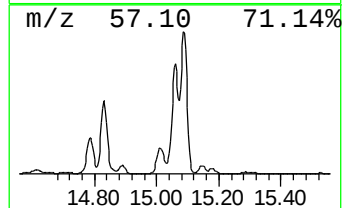
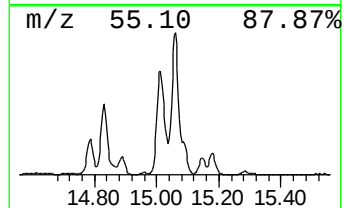
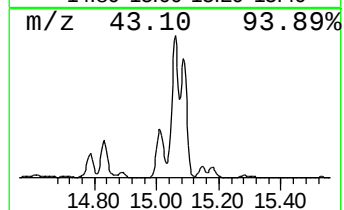
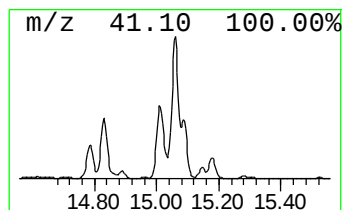
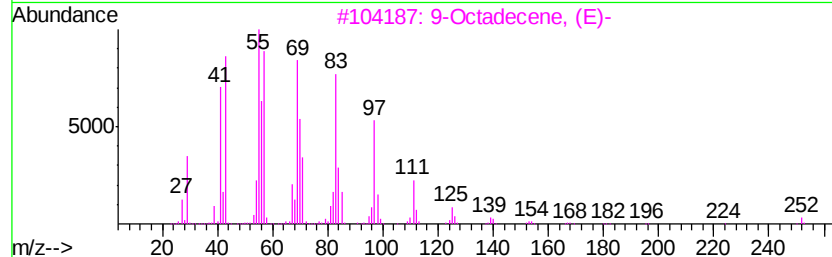
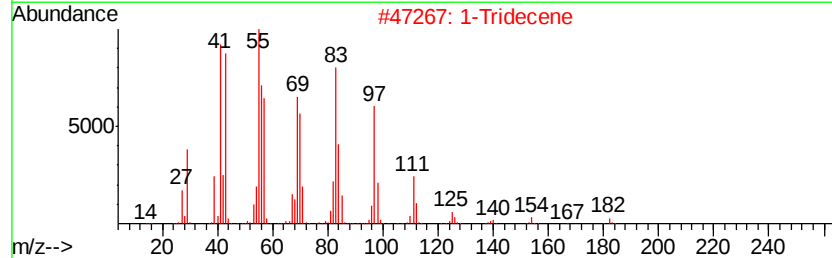
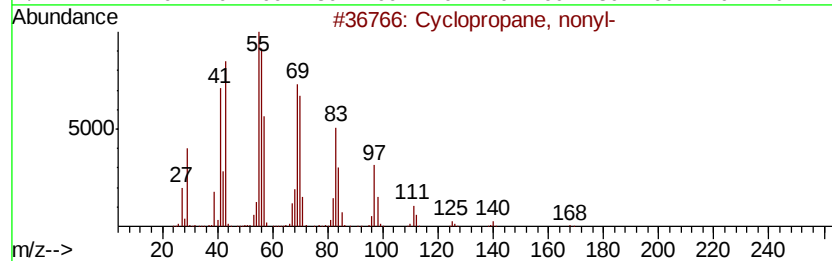
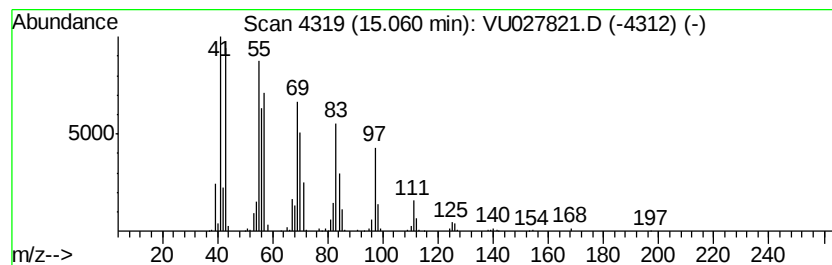
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Quant Title : VOC Analysis

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

Peak Number 6 (DEL) Alkane: Cyclic15.06 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.06	20.51 ug/L	444015	1,4-Dichlorobenzene-d4	11.49

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopropane, nonyl-	168	C12H24	074663-85-7	95
2		1-Tridecene	182	C13H26	002437-56-1	91
3		9-Octadecene, (E)-	252	C18H36	007206-25-9	91
4		Cyclododecane	168	C12H24	000294-62-2	90
5		2-Tetradecene, (E)-	196	C14H28	035953-54-9	90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU102618\
Data File : VU027821.D
Acq On : 26 Oct 2018 10:27
Operator : MD/SY
Sample : J5657-01ME
Misc : 4.65u/5mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
A40G4ME

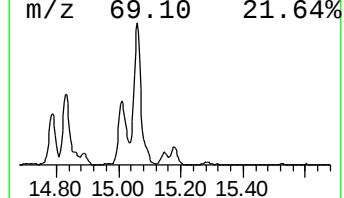
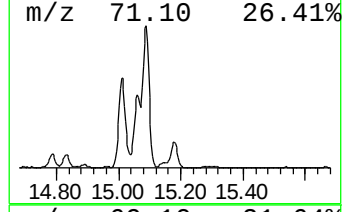
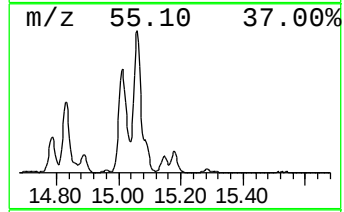
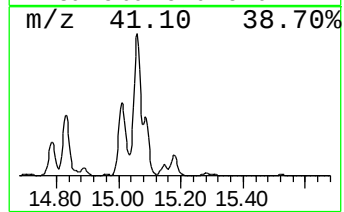
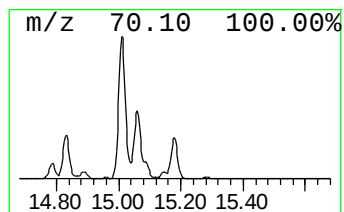
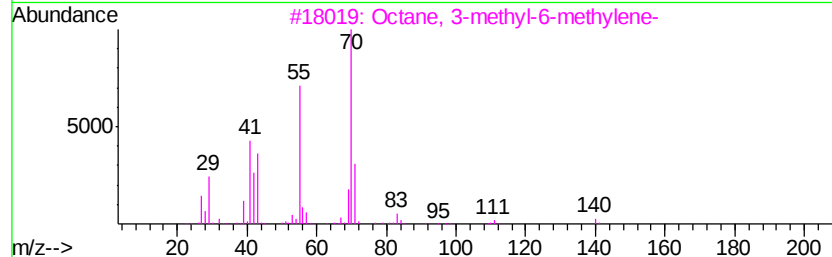
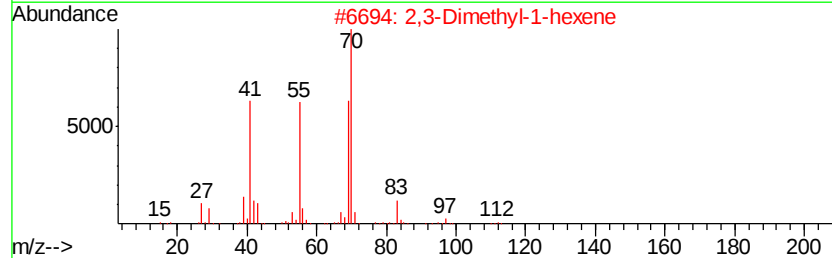
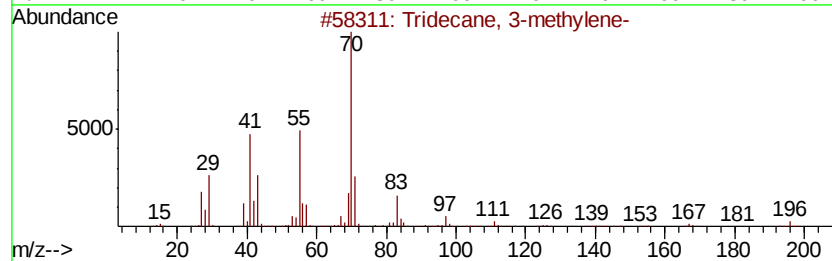
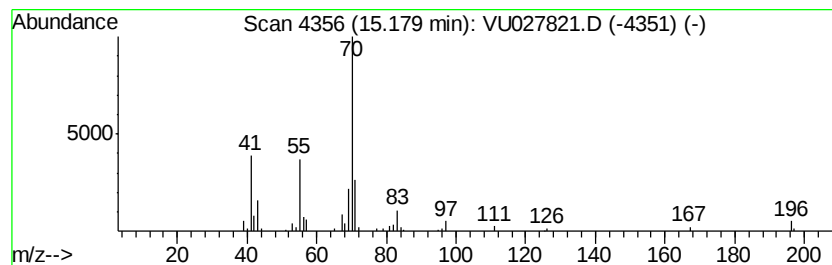
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Quant Title : VOC Analysis

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

Peak Number 7 (DEL) Alkane: Cyclic15.18 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.18	3.62 ug/L	78409	1,4-Dichlorobenzene-d4	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane, 3-methyl-	196	C14H28	019780-34-8	50
2		2,3-Dimethyl-1-hexene	112	C8H16	016746-86-4	50
3		Octane, 3-methyl-6-methylene-	140	C10H20	074630-07-2	50
4		2-Undecene, 3-methyl-, (Z)-	168	C12H24	057024-90-5	39
5		1-Decen-3-one	154	C10H18O	056606-79-2	39



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU102618\
Data File : VU027821.D
Acq On : 26 Oct 2018 10:27
Operator : MD/SY
Sample : J5657-01ME
Misc : 4.65uL/5mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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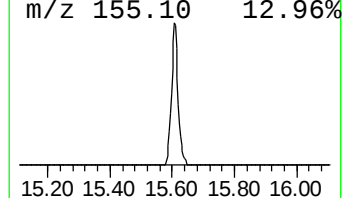
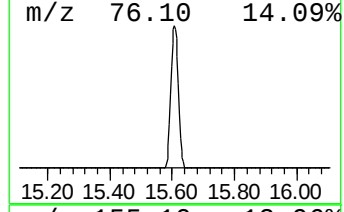
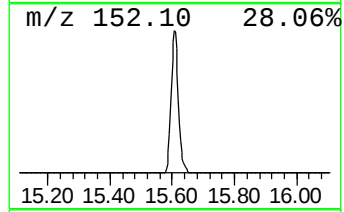
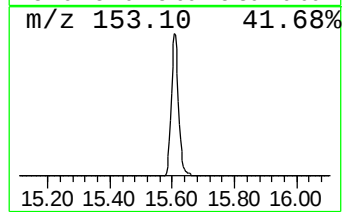
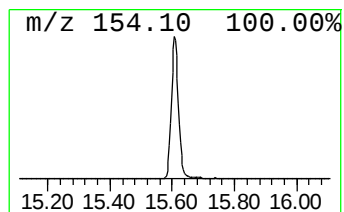
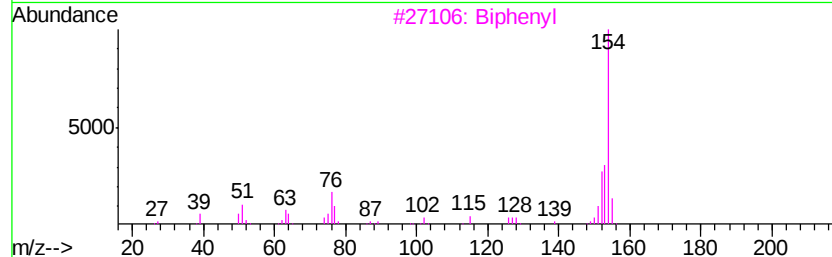
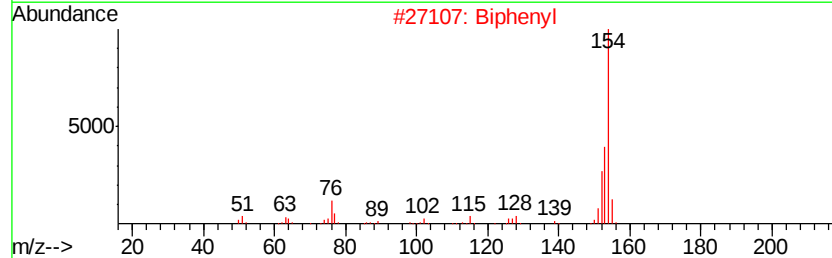
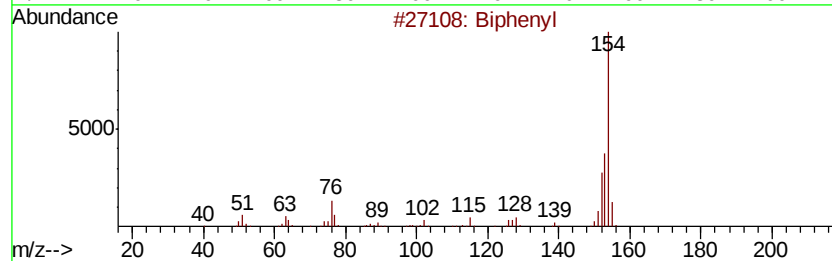
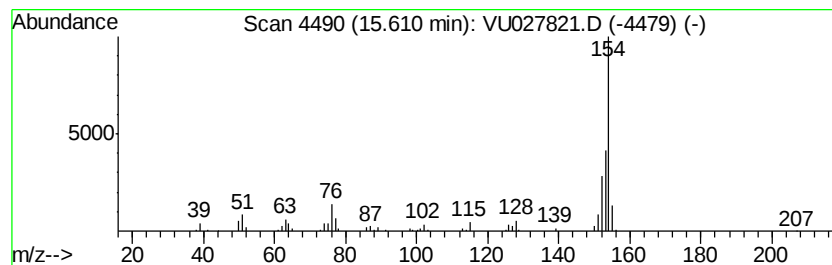
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Quant Title : VOC Analysis

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

Peak Number 8 Biphenyl Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.61	4.55 ug/L	98432	1,4-Dichlorobenzene-d4	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Biphenyl	154	C12H10	000092-52-4	95
2		Biphenyl	154	C12H10	000092-52-4	95
3		Biphenyl	154	C12H10	000092-52-4	94
4		Biphenyl	154	C12H10	000092-52-4	93
5		Biphenyl	154	C12H10	000092-52-4	93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU102618\
Data File : VU027821.D
Acq On : 26 Oct 2018 10:27
Operator : MD/SY
Sample : J5657-01ME
Misc : 4.65uL/5mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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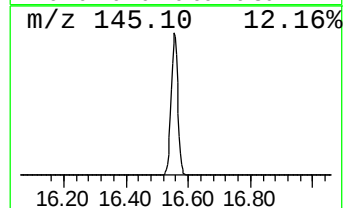
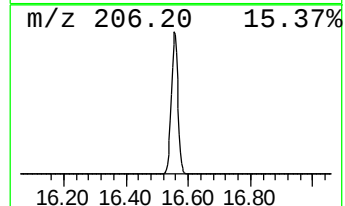
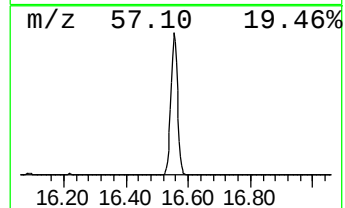
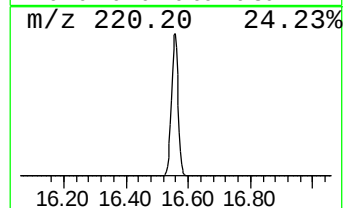
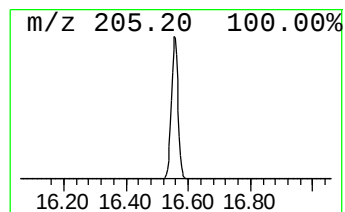
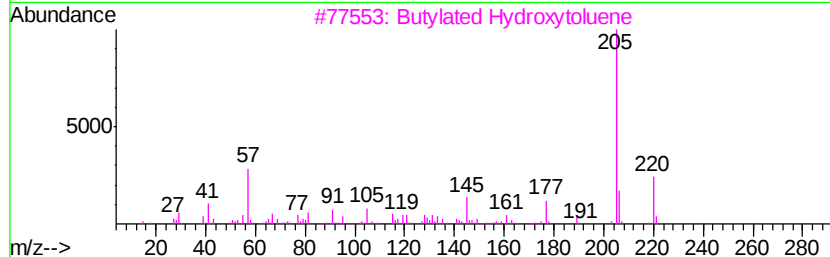
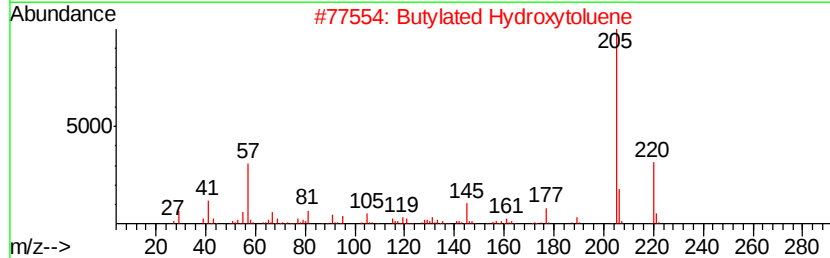
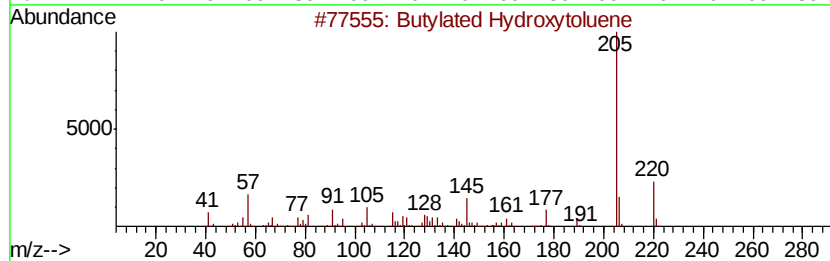
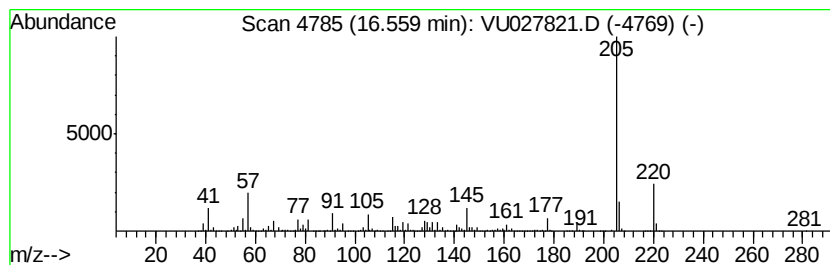
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Quant Title : VOC Analysis

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

Peak Number 9 Butylated Hydroxytoluene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.56	167.91 ug/L	3635220	1,4-Dichlorobenzene-d4	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butylated Hydroxytoluene	220	C15H24O	000128-37-0	98
2		Butylated Hydroxytoluene	220	C15H24O	000128-37-0	96
3		Butylated Hydroxytoluene	220	C15H24O	000128-37-0	95
4		Butylated Hydroxytoluene	220	C15H24O	000128-37-0	94
5		Phenol, 2,4,6-tris(1-methylethyl)-	220	C15H24O	002934-07-8	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU102618\
Data File : VU027821.D
Acq On : 26 Oct 2018 10:27
Operator : MD/SY
Sample : J5657-01ME
Misc : 4.65u/5mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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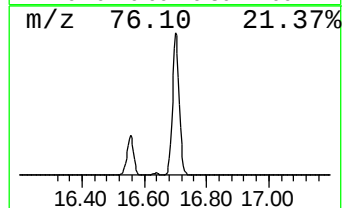
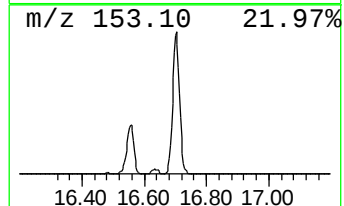
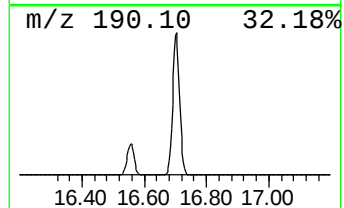
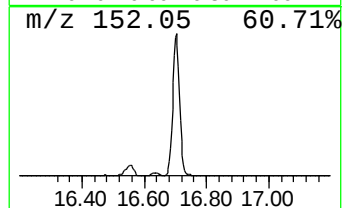
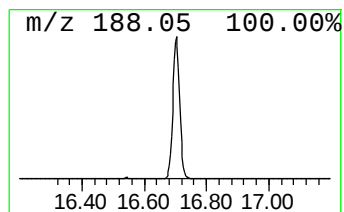
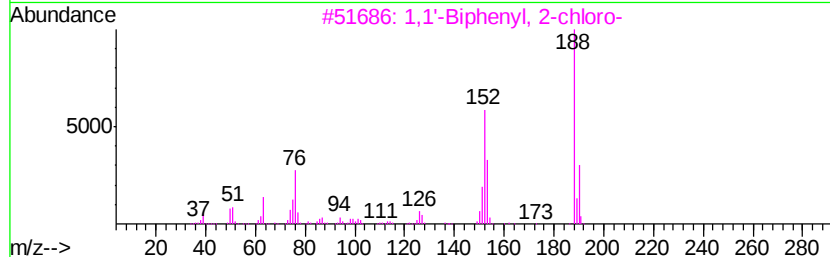
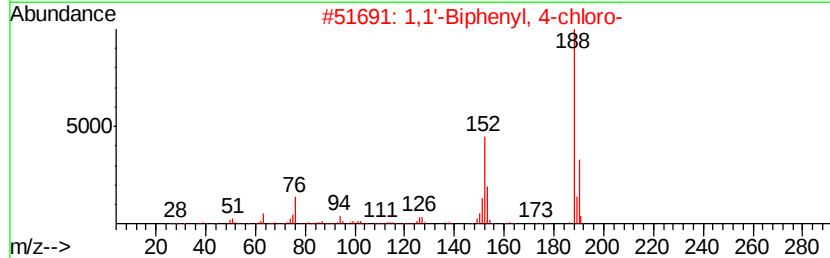
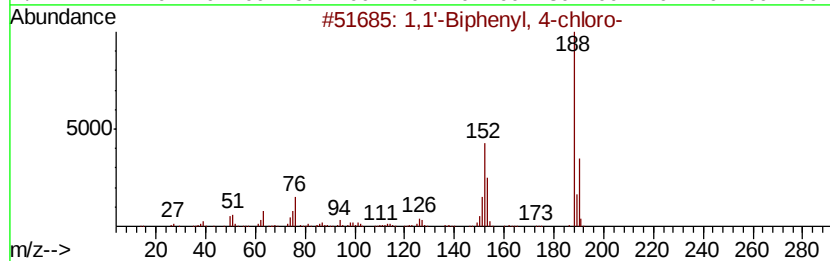
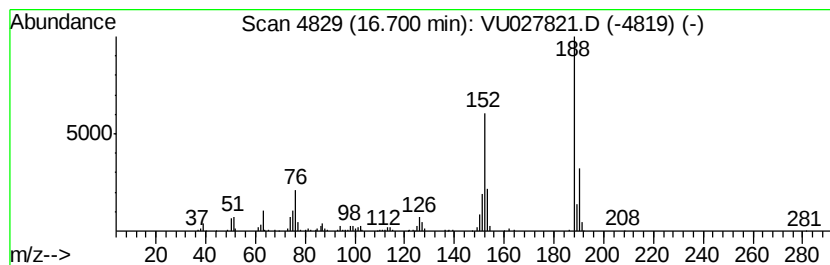
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Quant Title : VOC Analysis

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

Peak Number 10 1,1'-Biphenyl, 4-chloro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.70	13.85 ug/L	299751	1,4-Dichlorobenzene-d4	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 4-chloro-	188	C12H9Cl	002051-62-9	96
2		1,1'-Biphenyl, 4-chloro-	188	C12H9Cl	002051-62-9	94
3		1,1'-Biphenyl, 2-chloro-	188	C12H9Cl	002051-60-7	93
4		1,1'-Biphenyl, 4-chloro-	188	C12H9Cl	002051-62-9	93
5		3-Chlorobiphenyl	188	C12H9Cl	002051-61-8	93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU102618\
Data File : VU027821.D
Acq On : 26 Oct 2018 10:27
Operator : MD/SY
Sample : J5657-01ME
Misc : 4.65µ/5mL/100µL/5.0mL/MSVOA_U/MEOH
ALS Vial : 5 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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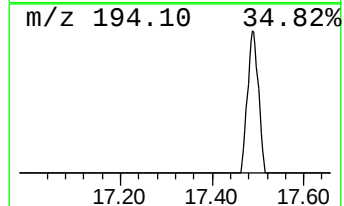
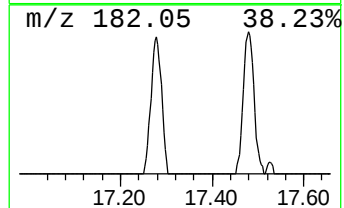
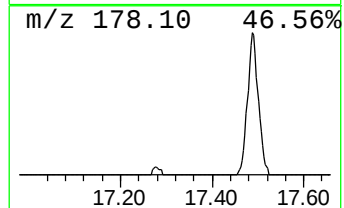
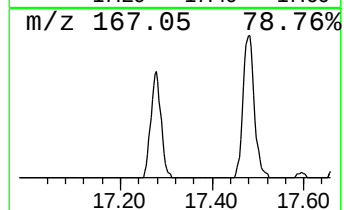
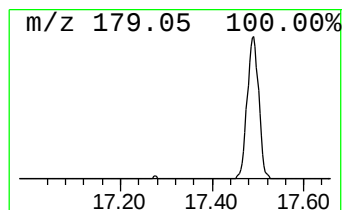
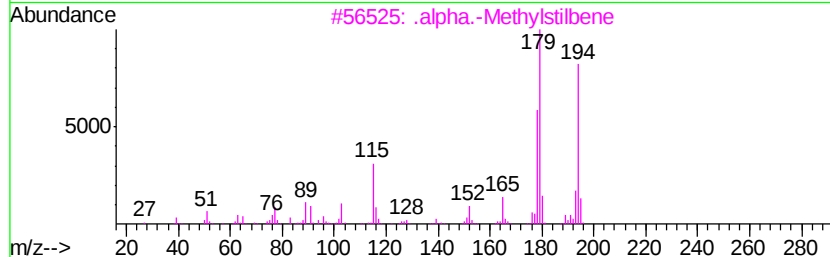
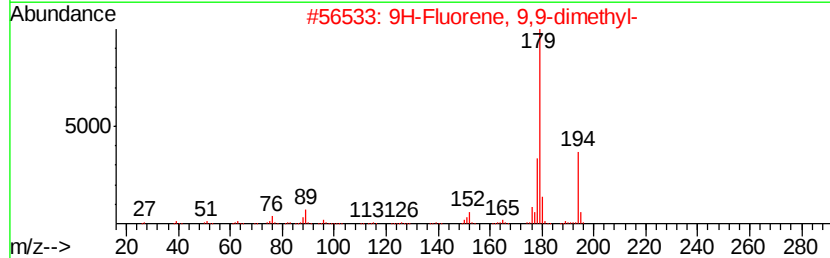
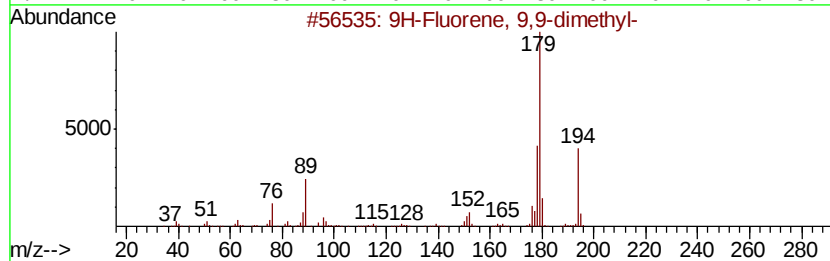
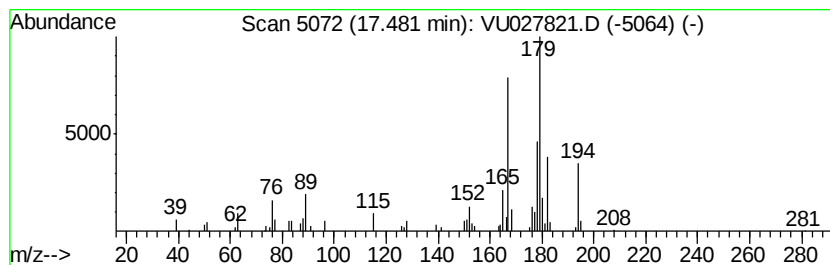
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Quant Title : VOC Analysis

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

Peak Number 11 9H-Fluorene, 9,9-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.48	2.58 ug/L	55824	1,4-Dichlorobenzene-d4	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9H-Fluorene, 9,9-dimethyl-	194	C15H14	004569-45-3	92
2		9H-Fluorene, 9,9-dimethyl-	194	C15H14	004569-45-3	83
3		.alpha.-Methylstilbene	194	C15H14	000779-51-1	70
4		Phenanthrene, 9,10-dihydro-1-met...	194	C15H14	095676-48-5	55
5		Anthracene, 9,10-dihydro-2-methyl-	194	C15H14	000948-67-4	52



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_U\DATA\VU102618\
Data File : VU027821.D
Acq On : 26 Oct 2018 10:27
Operator : MD/SY
Sample : J5657-01ME
Misc : 4.65g/5mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
A40G4ME

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM102218WMA.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
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(DEL) Alkane: Cyc...	14.83	15.3	ug/L	331442	3	11.49	1082500	50.0
(DEL) Alkane: Cyc...	15.01	17.9	ug/L	386381	3	11.49	1082500	50.0
(DEL) Alkane: Cyc...	15.06	20.5	ug/L	444015	3	11.49	1082500	50.0
(DEL) Alkane: Cyc...	15.18	3.6	ug/L	78409	3	11.49	1082500	50.0
Biphenyl	15.61	4.5	ug/L	98432	3	11.49	1082500	50.0
Butylated Hydroxy...	16.56	167.9	ug/L	3635220	3	11.49	1082500	50.0
1,1'-Biphenyl, 4-...	16.70	13.9	ug/L	299751	3	11.49	1082500	50.0
9H-Fluorene, 9,9-...	17.48	2.6	ug/L	55824	3	11.49	1082500	50.0