

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU062818\
 Data File : VU025025.D
 Acq On : 28 Jun 2018 17:06
 Operator : MD/SY
 Sample : J3707-26 20X
 Misc : 5.80g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 WP021SB489-31-33-180622

Integration Parameters: RTEINT.P

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

Filtering: 5
 Min Area: 3 % 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\82U061318W.M
 Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.088	136	155	180	rBV	1171018	1342182	51.15%	8.279%
2	1.580	291	308	316	rBV6	12082	34545	1.32%	0.213%
3	4.230	1107	1132	1158	rBV2	218426	545338	20.78%	3.364%
4	4.886	1315	1336	1353	rBV	191678	435268	16.59%	2.685%
5	4.985	1353	1367	1386	rVV	293064	656742	25.03%	4.051%
6	5.313	1453	1469	1488	rBV2	206627	432995	16.50%	2.671%
7	5.542	1525	1540	1556	rVB3	34399	82370	3.14%	0.508%
8	5.889	1631	1648	1677	rBV	396037	782395	29.81%	4.826%
9	6.191	1726	1742	1787	rBV	1337338	2624200	100.00%	16.187%
10	6.416	1797	1812	1824	rBV2	19189	39565	1.51%	0.244%
11	6.931	1955	1972	1984	rBV2	20465	42093	1.60%	0.260%
12	7.053	1996	2010	2020	rBV3	23224	45321	1.73%	0.280%
13	7.567	2141	2170	2188	rBV	730677	1303658	49.68%	8.042%
14	8.915	2580	2589	2602	rVB2	17499	29785	1.14%	0.184%
15	9.091	2633	2644	2671	rVB	572646	940057	35.82%	5.799%
16	9.448	2746	2755	2776	rVB	84691	137156	5.23%	0.846%
17	9.722	2827	2840	2857	rBV6	21068	44629	1.70%	0.275%
18	9.818	2857	2870	2879	rBV4	15994	30075	1.15%	0.186%
19	9.953	2895	2912	2923	rBV	46610	83494	3.18%	0.515%
20	10.046	2933	2941	2949	rBV3	29455	46060	1.76%	0.284%
21	10.091	2950	2955	2967	rVB2	20994	34362	1.31%	0.212%
22	10.259	2988	3007	3013	rBV4	29423	69169	2.64%	0.427%
23	10.310	3013	3023	3033	rVV	603652	1022899	38.98%	6.310%
24	10.355	3034	3037	3051	rVB2	76229	108726	4.14%	0.671%
25	10.458	3061	3069	3076	rBV	59290	84326	3.21%	0.520%
26	10.689	3133	3141	3154	rVV3	33261	81106	3.09%	0.500%
27	10.773	3154	3167	3172	rVV3	42928	102159	3.89%	0.630%
28	10.815	3172	3180	3193	rVB	569848	828423	31.57%	5.110%
29	10.972	3222	3229	3235	rBV3	18862	30544	1.16%	0.188%
30	11.014	3236	3242	3249	rVB2	28859	39394	1.50%	0.243%
31	11.078	3254	3262	3267	rBV3	31649	45610	1.74%	0.281%
32	11.117	3267	3274	3280	rVV	114139	158183	6.03%	0.976%
33	11.152	3280	3285	3293	rVB	67951	93610	3.57%	0.577%
34	11.271	3313	3322	3331	rBV3	39749	82902	3.16%	0.511%

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Integration Parameters: RTEINT.P

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35	11.400	3353	3362	3369	rBV3	51893	91345	3.48%	0.563%
36	11.487	3379	3389	3399	rBV	668327	1047911	39.93%	6.464%
37	11.532	3400	3403	3410	rVV	57389	68619	2.61%	0.423%
38	11.574	3410	3416	3422	rVV	57728	75542	2.88%	0.466%
39	11.612	3422	3428	3439	rVB	86988	118557	4.52%	0.731%
40	11.702	3450	3456	3470	rVB	68988	109090	4.16%	0.673%
41	11.812	3473	3490	3498	rBV2	58761	129660	4.94%	0.800%
42	11.879	3498	3511	3525	rVB5	111321	264729	10.09%	1.633%
43	12.027	3548	3557	3565	rBV	446101	622945	23.74%	3.843%
44	12.152	3587	3596	3599	rBV	53433	77598	2.96%	0.479%
45	12.181	3600	3605	3613	rVV	71998	105115	4.01%	0.648%
46	12.252	3614	3627	3636	rVB2	86991	156475	5.96%	0.965%
47	12.365	3653	3662	3666	rBV4	28850	49432	1.88%	0.305%
48	12.500	3693	3704	3712	rBV7	27215	54976	2.09%	0.339%
49	12.551	3714	3720	3731	rVB3	37564	61833	2.36%	0.381%
50	12.615	3731	3740	3744	rBV5	32974	50876	1.94%	0.314%
51	12.651	3745	3751	3760	rVB2	71598	110483	4.21%	0.682%
52	12.705	3760	3768	3778	rBV	96017	154480	5.89%	0.953%
53	12.792	3787	3795	3799	rBV2	20277	29602	1.13%	0.183%
54	12.966	3842	3849	3859	rVB2	28717	44798	1.71%	0.276%
55	13.127	3891	3899	3915	rVB3	119078	194840	7.42%	1.202%
56	13.509	4010	4018	4025	rBV4	18259	31169	1.19%	0.192%
57	13.747	4084	4092	4101	rVB	19505	29286	1.12%	0.181%
58	13.889	4129	4136	4148	rVB2	23548	38426	1.46%	0.237%
59	14.146	4208	4216	4224	rBV2	24806	34300	1.31%	0.212%

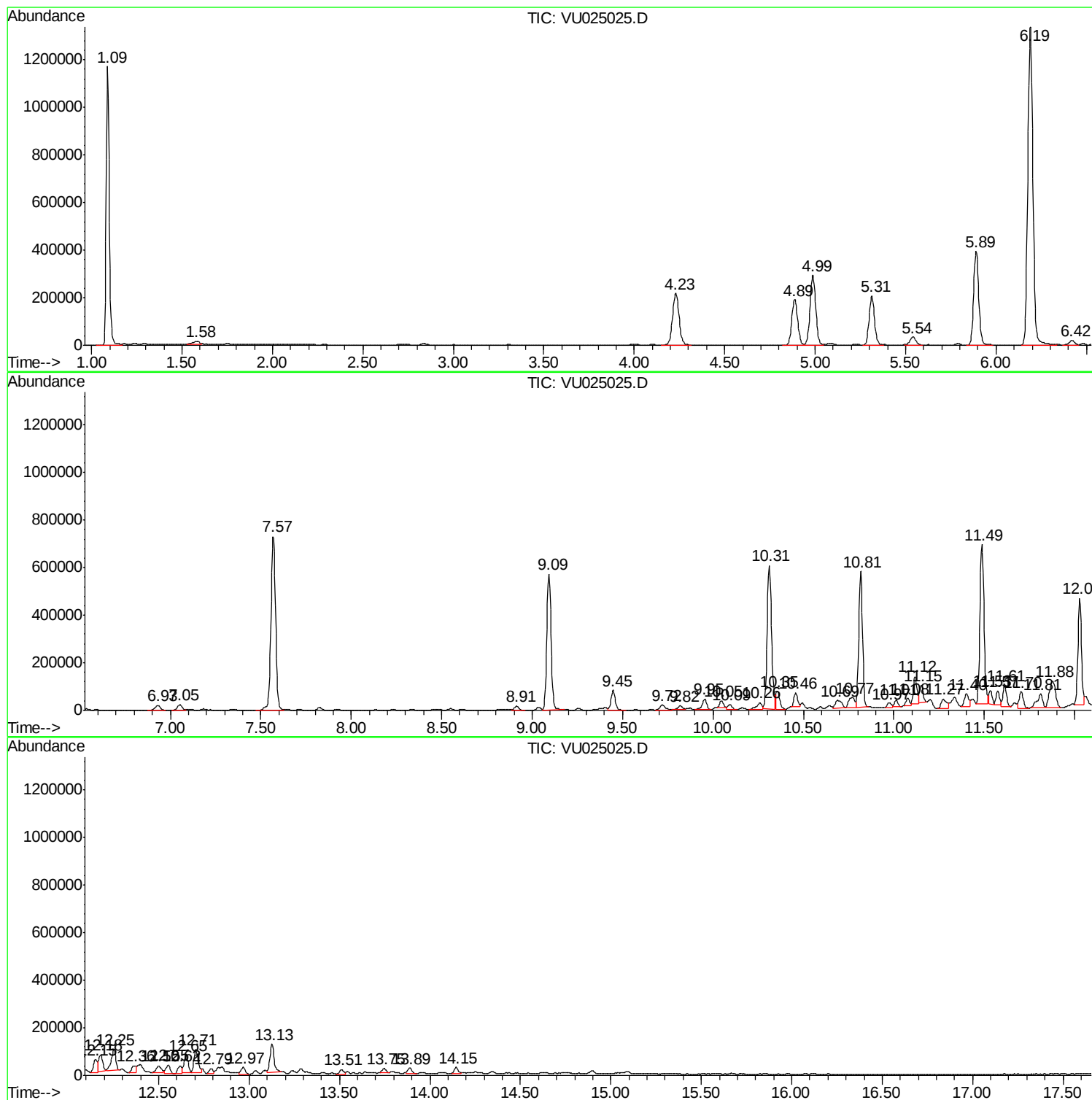
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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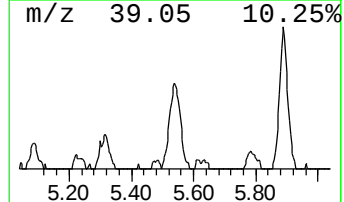
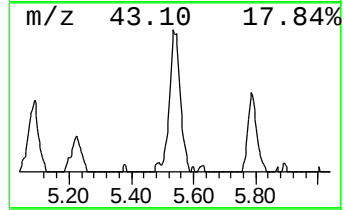
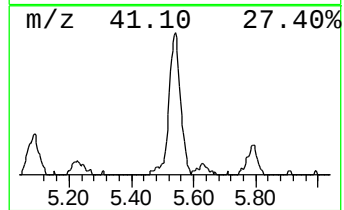
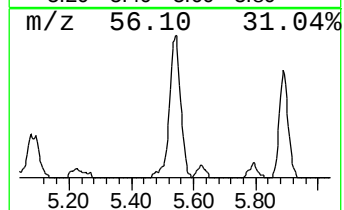
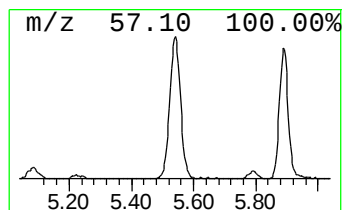
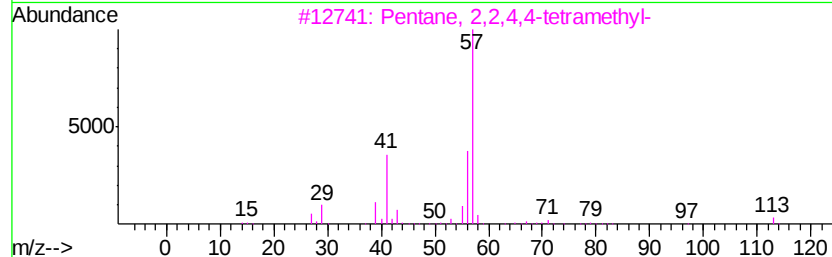
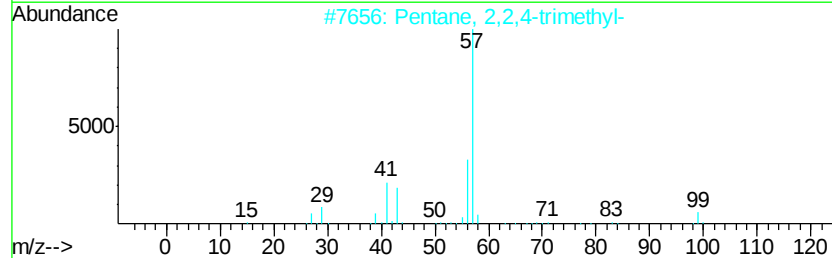
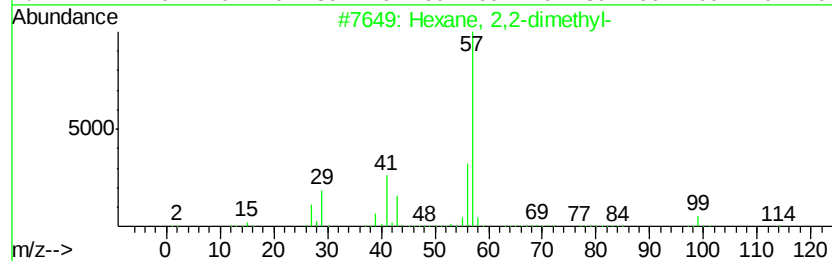
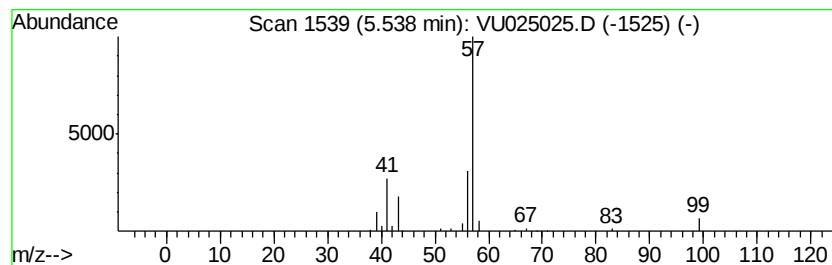
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U061318W.M
Quant Title : SW846 8260

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

Peak Number 2 Hexane, 2,2-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.54	5.26 ug/l	82370	1,4-Difluorobenzene	5.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	72
2		Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	72
3		Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	64
4		Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	56
5		Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	50



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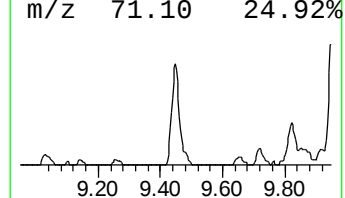
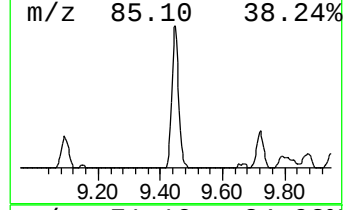
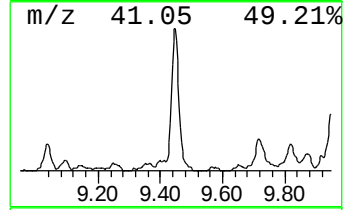
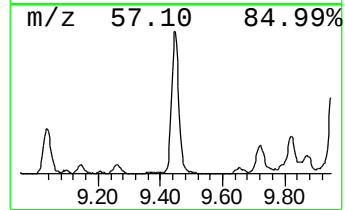
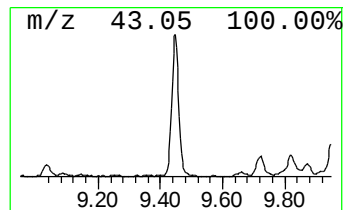
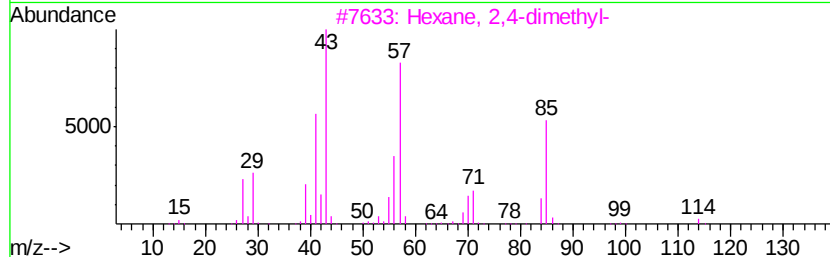
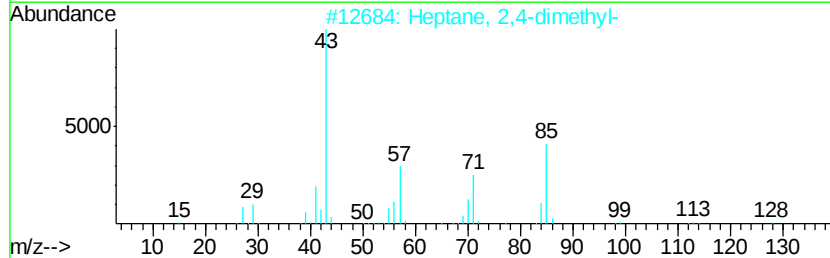
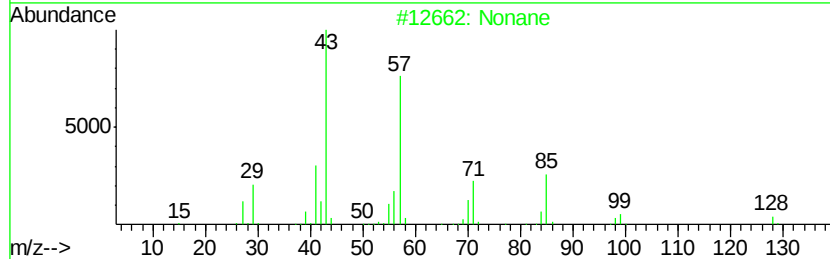
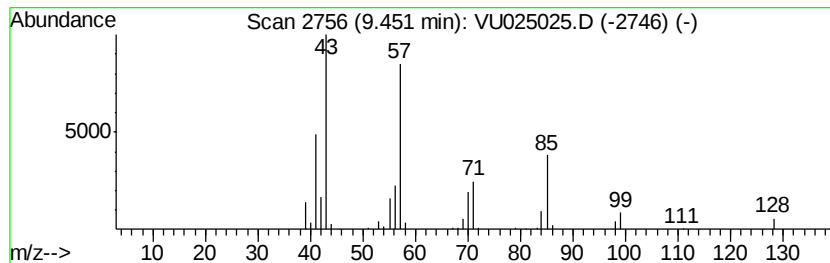
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 Nonane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.		
9.45	7.30 ug/l	137156	Chlorobenzene-d5	9.09		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonane		128	C9H20	000111-84-2	90
2	Heptane, 2,4-dimethyl-		128	C9H20	002213-23-2	53
3	Hexane, 2,4-dimethyl-		114	C8H18	000589-43-5	53
4	Decane, 2,5,6-trimethyl-		184	C13H28	062108-23-0	50
5	Sulfurous acid, hexyl heptyl ester		264	C13H28O3S	1000309-12-9	50



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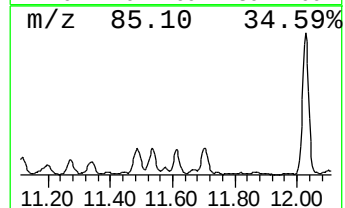
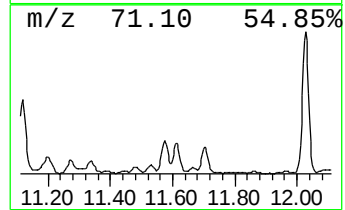
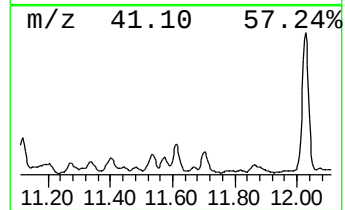
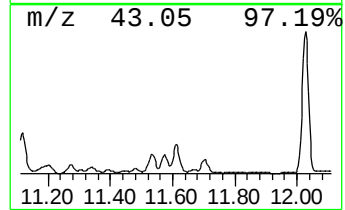
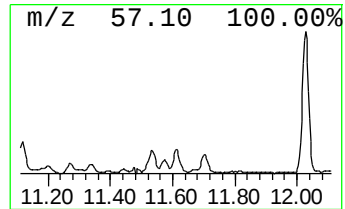
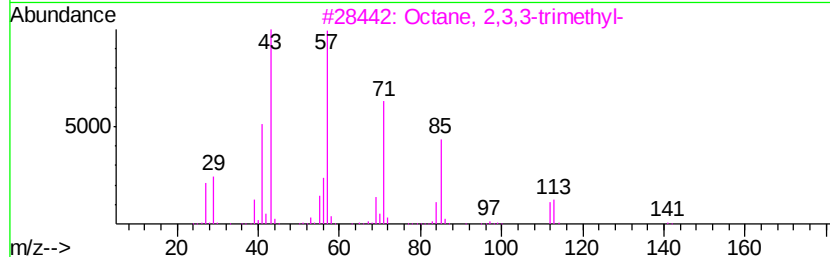
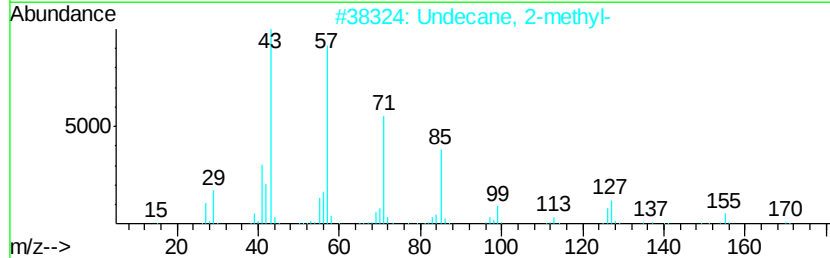
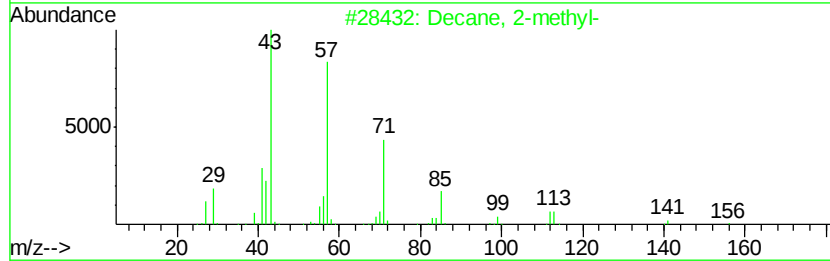
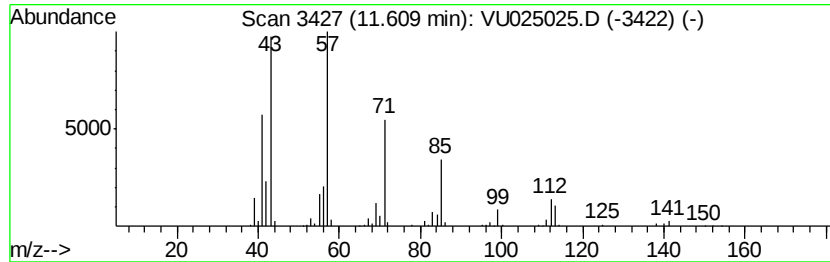
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Decane, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.61	5.66 ug/l	118557	1,4-Dichlorobenzene-d4	11.49	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane, 2-methyl-	156	C11H24	006975-98-0	87
2	Undecane, 2-methyl-	170	C12H26	007045-71-8	72
3	Octane, 2,3,3-trimethyl-	156	C11H24	062016-30-2	72
4	Tridecane	184	C13H28	000629-50-5	64
5	1-Iodo-2-methylnonane	268	C10H21I	1000101-47-9	64



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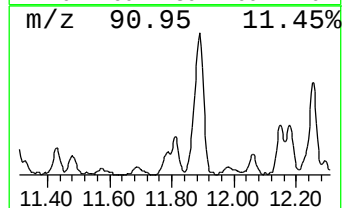
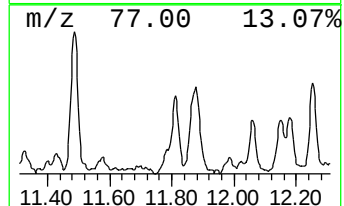
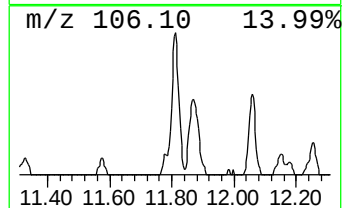
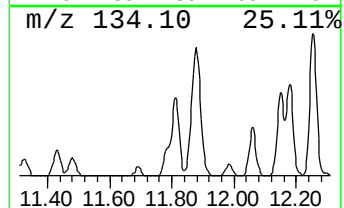
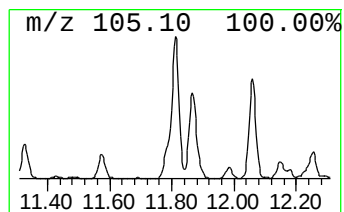
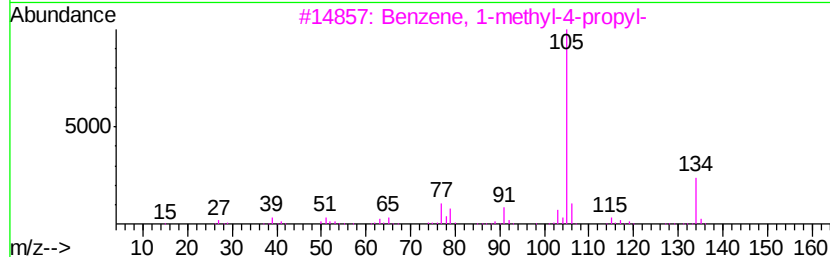
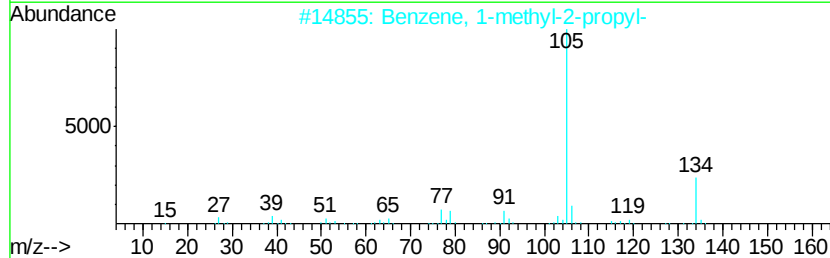
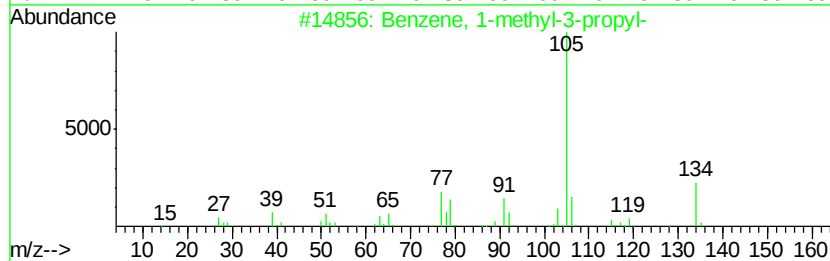
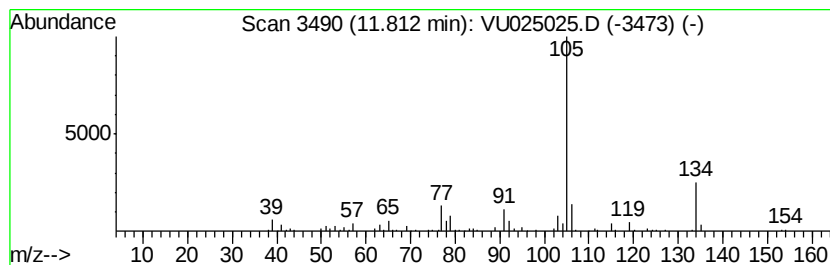
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, 1-methyl-3-propyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.81	6.19 ug/l	129660	1,4-Dichlorobenzene-d4	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	94
2		Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	91
3		Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
4		Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	72
5		2-Tolyloxirane	134	C9H10O	002783-26-8	72



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Misc : 5.80uL/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 19 Sample Multiplier: 1

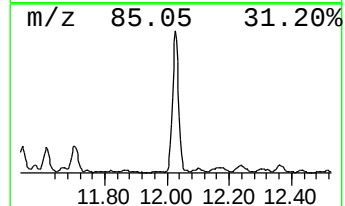
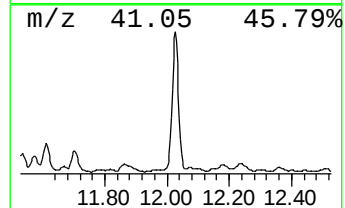
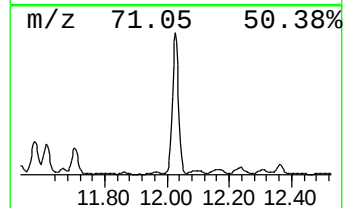
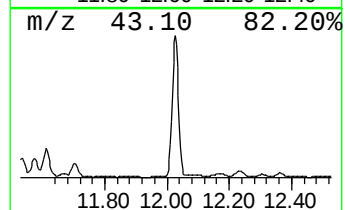
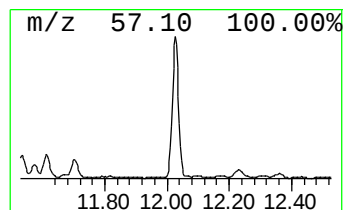
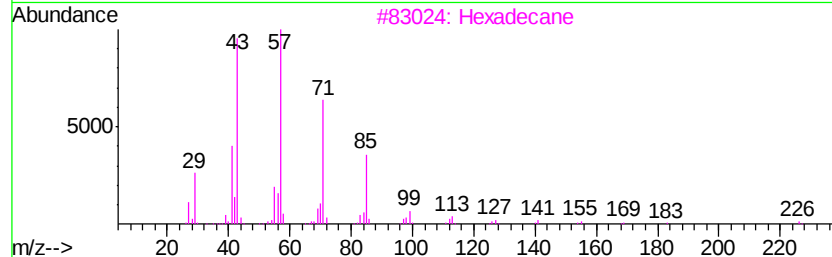
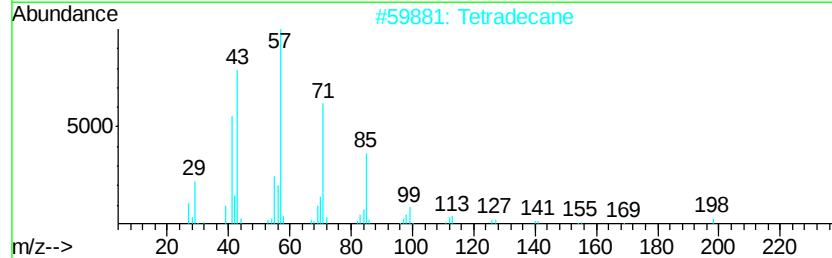
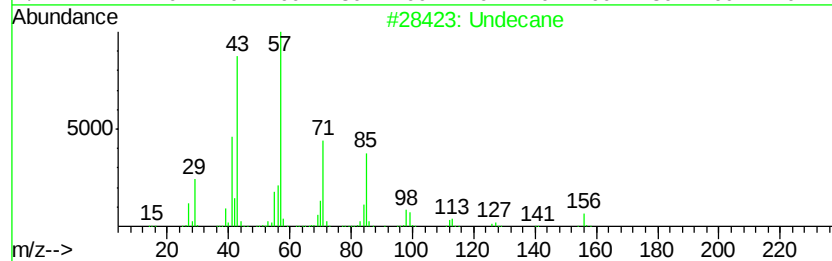
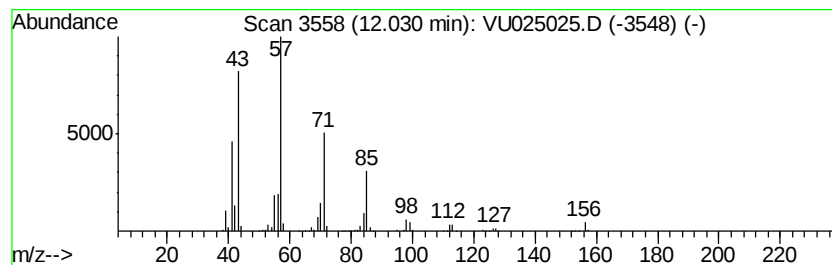
Instrument :
MSVOA_U
ClientSampleId :
WP021SB489-31-33-180622

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U061318W.M
Quant Title : SW846 8260

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

Peak Number 6 Undecane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.03	29.72 ug/l	622945	1,4-Dichlorobenzene-d4	11.49
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Undecane	156 C11H24	001120-21-4	94
2	Tetradecane	198 C14H30	000629-59-4	86
3	Hexadecane	226 C16H34	000544-76-3	80
4	Tridecane	184 C13H28	000629-50-5	78
5	Nonane	128 C9H20	000111-84-2	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU062818\
Data File : VU025025.D
Acq On : 28 Jun 2018 17:06
Operator : MD/SY
Sample : J3707-26 20X
Misc : 5.80g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
WP021SB489-31-33-180622

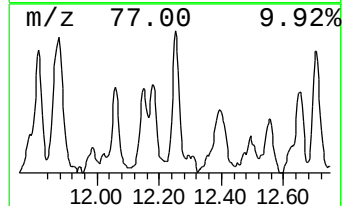
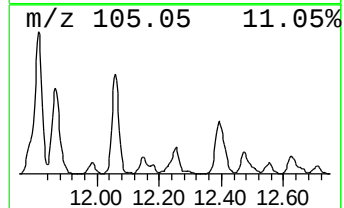
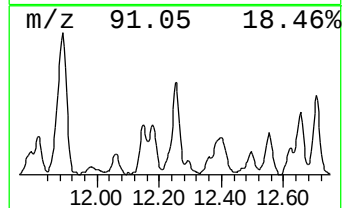
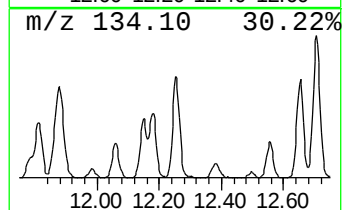
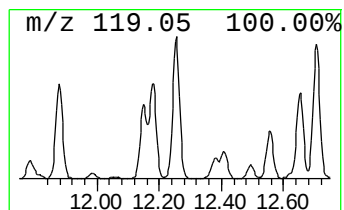
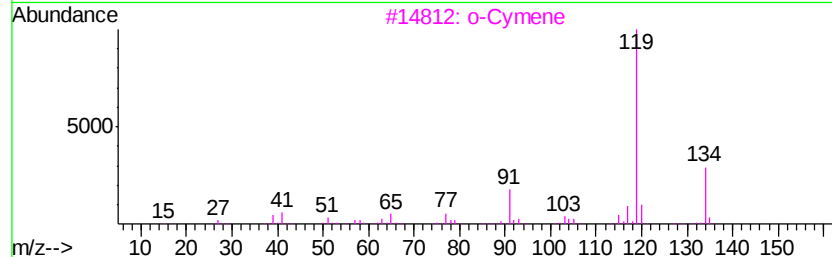
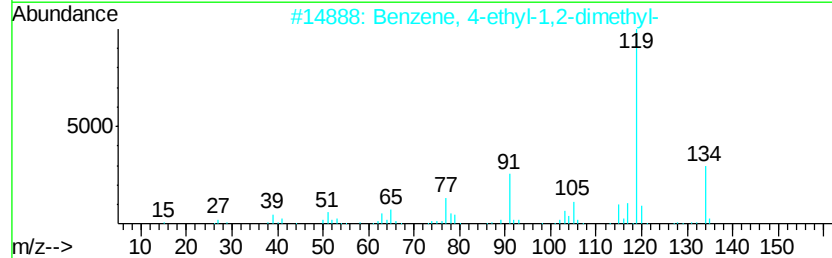
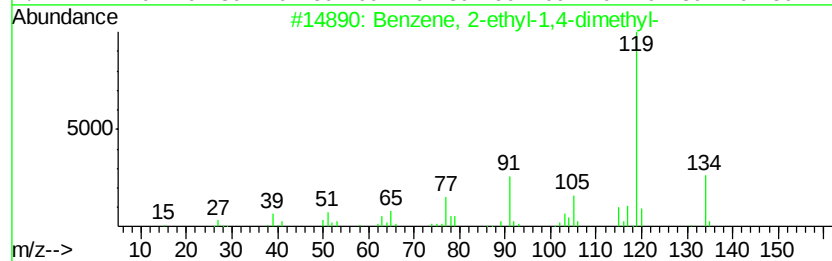
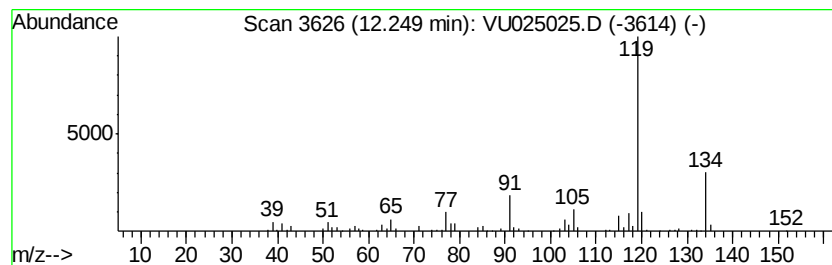
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.25	7.47 ug/l	156475	1,4-Dichlorobenzene-d4	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
3		o-Cymene	134	C10H14	000527-84-4	95
4		p-Cymene	134	C10H14	000099-87-6	95
5		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU062818\
Data File : VU025025.D
Acq On : 28 Jun 2018 17:06
Operator : MD/SY
Sample : J3707-26 20X
Misc : 5.80u/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
WP021SB489-31-33-180622

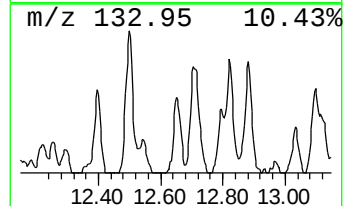
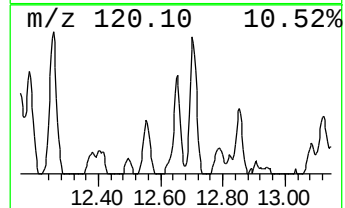
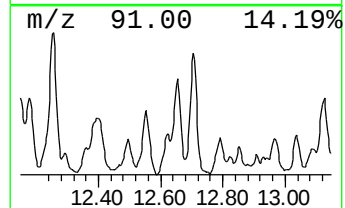
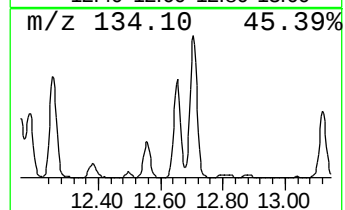
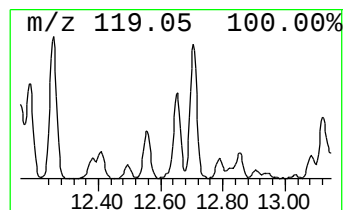
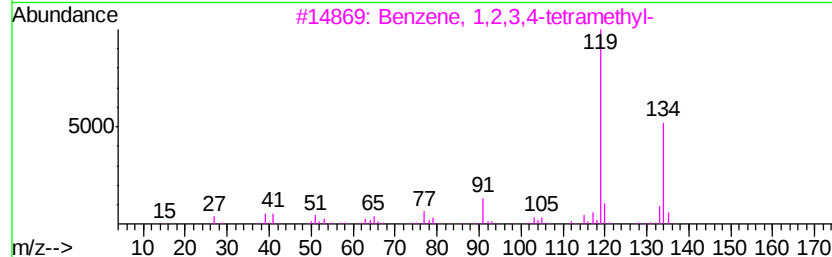
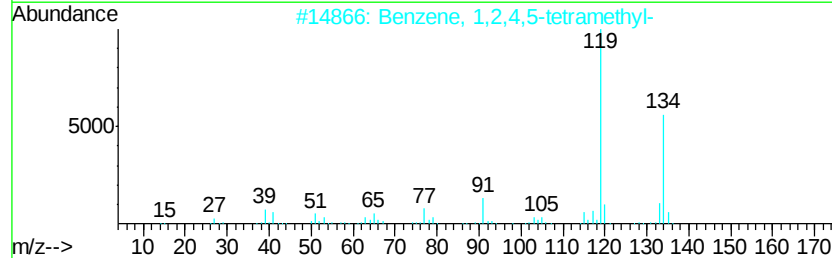
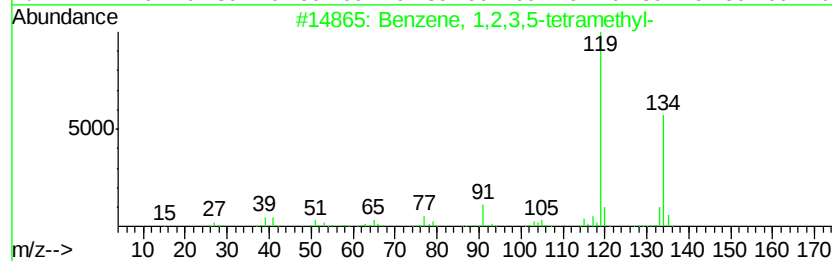
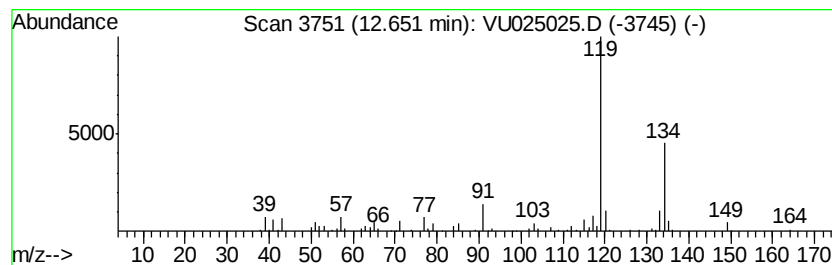
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.65	5.27 ug/l	110483	1,4-Dichlorobenzene-d4	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
2		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
3		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
4		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	93
5		o-Cymene	134	C10H14	000527-84-4	93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU062818\
Data File : VU025025.D
Acq On : 28 Jun 2018 17:06
Operator : MD/SY
Sample : J3707-26 20X
Misc : 5.80µ/5.0mL/100µL/5.0mL/MSVOA_U/MEOH
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
WP021SB489-31-33-180622

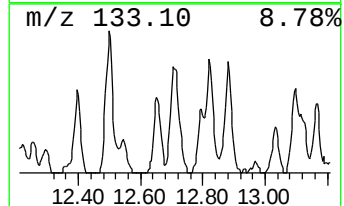
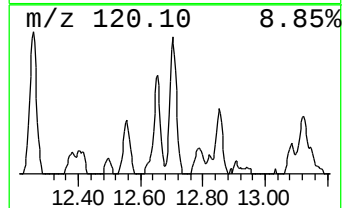
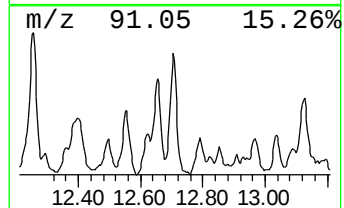
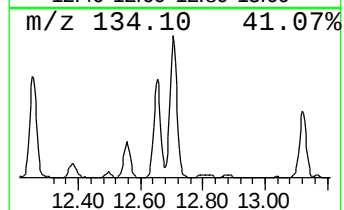
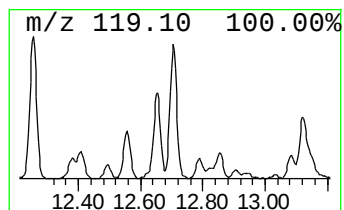
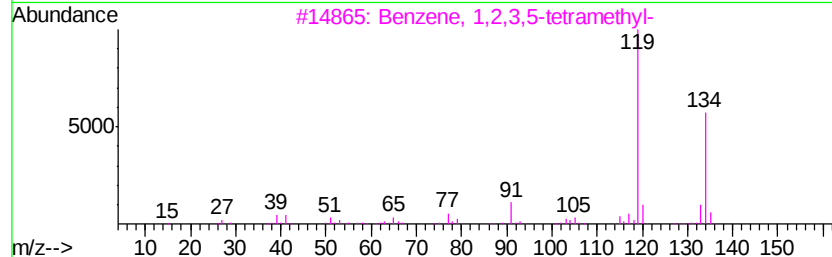
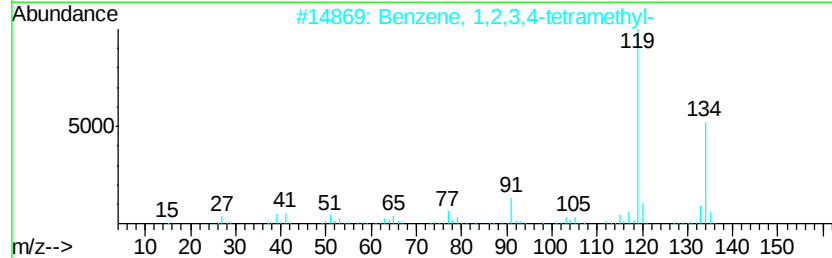
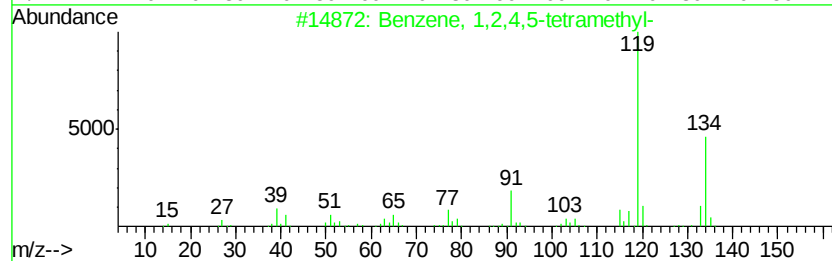
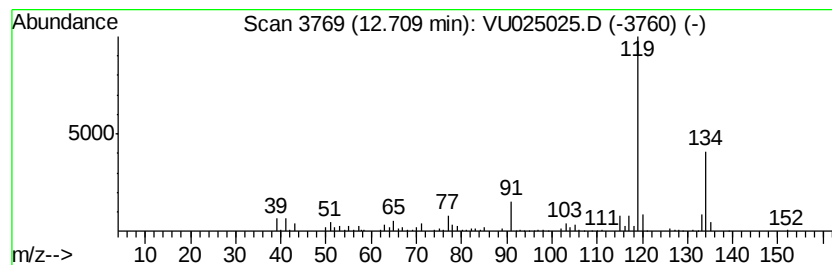
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.71	7.37 ug/l	154480	1,4-Dichlorobenzene-d4	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
4		o-Cymene	134	C10H14	000527-84-4	94
5		p-Cymene	134	C10H14	000099-87-6	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU062818\
Data File : VU025025.D
Acq On : 28 Jun 2018 17:06
Operator : MD/SY
Sample : J3707-26 20X
Misc : 5.80µ/5.0mL/100µL/5.0mL/MSVOA_U/MEOH
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
WP021SB489-31-33-180622

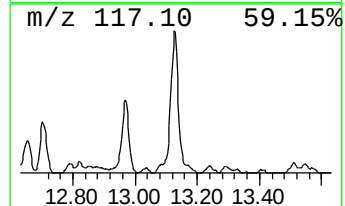
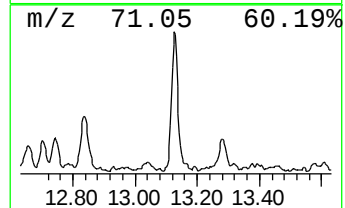
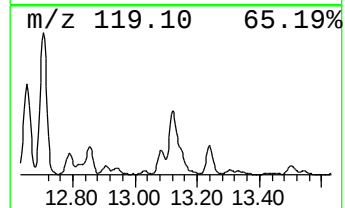
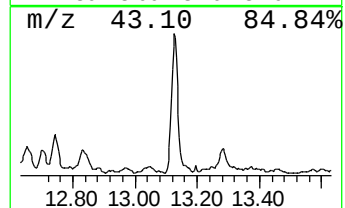
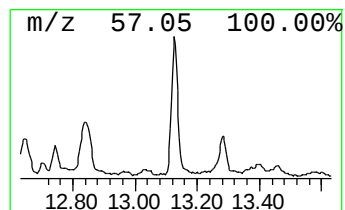
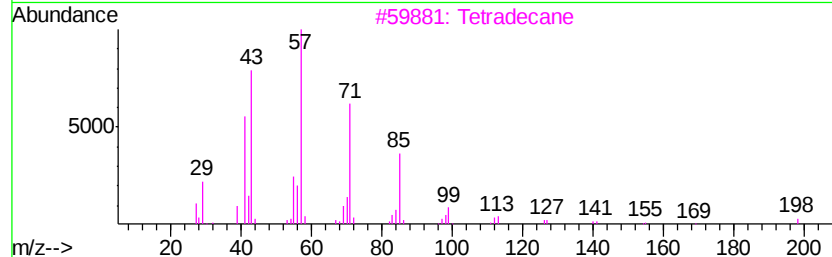
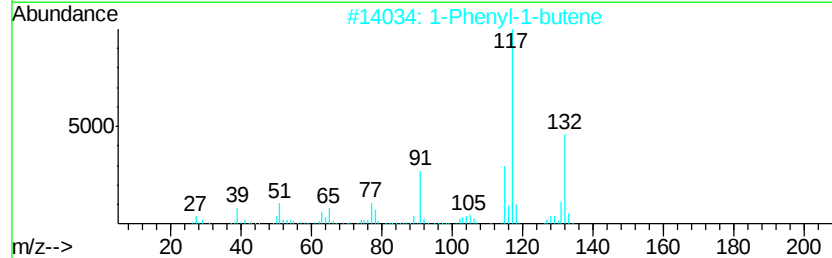
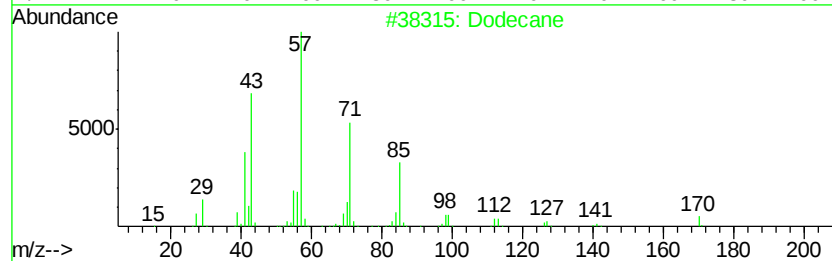
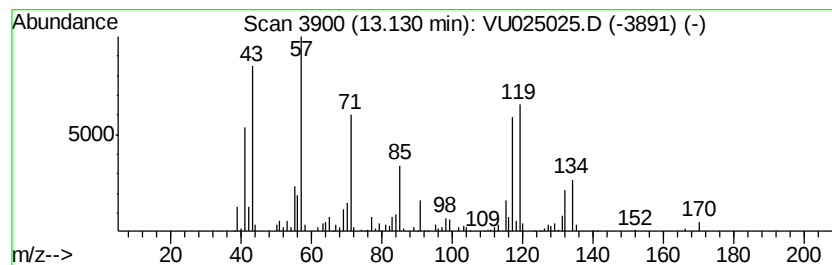
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U061318W.M
Quant Title : SW846 8260

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

Peak Number 10 Dodecane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.13	9.30 ug/l	194840	1,4-Dichlorobenzene-d4	11.49

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane	170	C12H26	000112-40-3	74
2			1-Phenyl-1-butene	132	C10H12	000824-90-8	55
3			Tetradecane	198	C14H30	000629-59-4	30
4			p-Cymene	134	C10H14	000099-87-6	25
5			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	25



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_U\DATA\VU062818\
Data File : VU025025.D
Acq On : 28 Jun 2018 17:06
Operator : MD/SY
Sample : J3707-26 20X
Misc : 5.80g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
WP021SB489-31-33-180622

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U061318W.M
Quant Title : SW846 8260

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Hexane, 2,2-dimet...	5.54	5.3	ug/l	82370	2	5.89	782395	50.0
Nonane	9.45	7.3	ug/l	137156	3	9.09	940057	50.0
Decane, 2-methyl-	11.61	5.7	ug/l	118557	4	11.49	1047910	50.0
Benzene, 1-methyl...	11.81	6.2	ug/l	129660	4	11.49	1047910	50.0
Undecane	12.03	29.7	ug/l	622945	4	11.49	1047910	50.0
Benzene, 2-ethyl-...	12.25	7.5	ug/l	156475	4	11.49	1047910	50.0
Benzene, 1,2,3,5-...	12.65	5.3	ug/l	110483	4	11.49	1047910	50.0
Benzene, 1,2,4,5-...	12.71	7.4	ug/l	154480	4	11.49	1047910	50.0
Dodecane	13.13	9.3	ug/l	194840	4	11.49	1047910	50.0