

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU100418\
 Data File : VU027399.D
 Acq On : 04 Oct 2018 11:36
 Operator : MD/SY
 Sample : J5084-12 50X
 Misc : 6.42g/5mL/100uL/5.0mL/MSVOA_U/MEOH
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 B-6(8-8.5)

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\82U100118W.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	1.759	173	182	201	rVB	55584	74834	2.80%	0.222%
2	1.948	232	241	259	rVB2	32826	49407	1.85%	0.147%
3	2.746	465	489	501	rBV	144868	369078	13.79%	1.094%
4	3.000	550	568	586	rBV	104502	222159	8.30%	0.659%
5	3.308	647	664	682	rBV2	92626	199753	7.46%	0.592%
6	3.894	819	846	860	rVV5	20009	65070	2.43%	0.193%
7	4.000	861	879	894	rVV	77725	218987	8.18%	0.649%
8	4.099	895	910	927	rVV	94925	260611	9.74%	0.773%
9	4.781	1117	1122	1139	rVV2	22552	44617	1.67%	0.132%
10	4.884	1139	1154	1171	rVV	122160	283988	10.61%	0.842%
11	4.987	1171	1186	1201	rVV3	366711	864714	32.31%	2.564%
12	5.083	1202	1216	1243	rVV2	253637	685187	25.60%	2.032%
13	5.222	1244	1259	1275	rVV	231315	535917	20.02%	1.589%
14	5.312	1275	1287	1309	rVB2	146149	317505	11.86%	0.942%
15	5.485	1328	1341	1345	rBV4	77101	145028	5.42%	0.430%
16	5.540	1346	1358	1374	rVV	533578	1330541	49.71%	3.946%
17	5.623	1375	1384	1396	rVB4	55294	127678	4.77%	0.379%
18	5.784	1420	1434	1445	rBV	153991	311301	11.63%	0.923%
19	5.887	1455	1466	1476	rBV	235736	442091	16.52%	1.311%
20	5.948	1477	1485	1511	rVB3	88300	262607	9.81%	0.779%
21	6.122	1529	1539	1556	rVB2	24927	47545	1.78%	0.141%
22	6.222	1557	1570	1582	rBV4	37004	83410	3.12%	0.247%
23	6.414	1607	1630	1641	rBV3	150202	415950	15.54%	1.233%
24	6.479	1641	1650	1658	rVV	92884	183158	6.84%	0.543%
25	6.537	1658	1668	1678	rVV	117614	234598	8.76%	0.696%
26	6.588	1679	1684	1692	rVV2	23948	41477	1.55%	0.123%
27	6.643	1692	1701	1715	rVB2	49797	95119	3.55%	0.282%
28	6.739	1716	1731	1744	rVB5	35647	80713	3.02%	0.239%
29	6.929	1774	1790	1807	rBV	272775	568572	21.24%	1.686%
30	7.051	1808	1828	1839	rBV2	245341	584387	21.83%	1.733%
31	7.106	1840	1845	1858	rVV2	111876	206599	7.72%	0.613%
32	7.183	1859	1869	1876	rVV	152775	276142	10.32%	0.819%
33	7.221	1877	1881	1897	rVB2	73983	111335	4.16%	0.330%
34	7.344	1910	1919	1938	rVB	167535	338479	12.65%	1.004%

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Integration Parameters: RTEINT.P

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

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Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U100118W.M
 Title : SW846 8260

35	7.440	1939	1949	1963	rVB5	38267	87216	3.26%	0.259%
36	7.527	1963	1976	1980	rBV3	133864	236929	8.85%	0.703%
37	7.566	1980	1988	2001	rVB	498084	864383	32.29%	2.563%
38	7.640	2001	2011	2023	rVB	253663	429011	16.03%	1.272%
39	7.701	2023	2030	2037	rBV3	29884	46748	1.75%	0.139%
40	7.823	2058	2068	2085	rVB2	132748	275522	10.29%	0.817%
41	7.913	2086	2096	2110	rVB4	36690	69951	2.61%	0.207%
42	8.048	2129	2138	2146	rBV2	41454	65557	2.45%	0.194%
43	8.334	2219	2227	2237	rVB4	33624	60358	2.25%	0.179%
44	8.456	2250	2265	2270	rBV	46754	90765	3.39%	0.269%
45	8.755	2344	2358	2369	rBV9	22137	49826	1.86%	0.148%
46	8.913	2398	2407	2419	rVB3	134091	220444	8.24%	0.654%
47	9.032	2434	2444	2453	rBV	106014	174866	6.53%	0.519%
48	9.090	2453	2462	2473	rVV	364531	585077	21.86%	1.735%
49	9.250	2502	2512	2525	rVB	349264	546427	20.41%	1.620%
50	9.376	2538	2551	2565	rBV	1621621	2676690	100.00%	7.938%
51	9.443	2565	2572	2599	rVB	142081	290110	10.84%	0.860%
52	9.717	2647	2657	2665	rBV3	46436	78374	2.93%	0.232%
53	9.781	2665	2677	2701	rVB	545747	905061	33.81%	2.684%
54	9.951	2712	2730	2740	rBV2	48158	103457	3.87%	0.307%
55	10.048	2752	2760	2769	rVB5	29265	47736	1.78%	0.142%
56	10.167	2786	2797	2808	rBV	88752	145735	5.44%	0.432%
57	10.308	2831	2841	2851	rVV	383303	655665	24.50%	1.944%
58	10.360	2851	2857	2871	rVB4	75905	155556	5.81%	0.461%
59	10.456	2879	2887	2895	rBV	47730	65904	2.46%	0.195%
60	10.588	2907	2928	2946	rBV	327393	537777	20.09%	1.595%
61	10.681	2946	2957	2975	rVV	1143821	2398814	89.62%	7.114%
62	10.771	2975	2985	2993	rVV	502593	781870	29.21%	2.319%
63	10.816	2993	2999	3018	rVB	122230	191587	7.16%	0.568%
64	10.974	3037	3048	3067	rVB	389538	622691	23.26%	1.847%
65	11.147	3085	3102	3116	rBV	1690704	2580174	96.39%	7.651%
66	11.292	3132	3147	3152	rBV3	39434	84549	3.16%	0.251%
67	11.324	3152	3157	3169	rVB2	42708	68863	2.57%	0.204%
68	11.427	3182	3189	3197	rVV	55073	81465	3.04%	0.242%
69	11.482	3197	3206	3217	rVV	460566	700633	26.18%	2.078%
70	11.572	3225	3234	3243	rBV	362230	517447	19.33%	1.534%
71	11.774	3283	3297	3302	rBV4	234340	463699	17.32%	1.375%

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 ClientSampleId :
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Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON Filtering: 5
 Sampling : 1 Min Area: 3 % of largest Peak
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 Stop Thrs : 0 Peak Location: TOP

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72	11.810	3303	3308	3316	rVV	289373	398682	14.89%	1.182%
73	11.874	3316	3328	3347	rVB5	399360	831100	31.05%	2.465%
74	11.980	3350	3361	3368	rBV2	50319	84842	3.17%	0.252%
75	12.025	3368	3375	3379	rBV	60154	80969	3.02%	0.240%
76	12.061	3379	3386	3402	rVB3	196474	325976	12.18%	0.967%
77	12.147	3403	3413	3417	rBV	182540	270968	10.12%	0.804%
78	12.176	3418	3422	3432	rVB	178150	235002	8.78%	0.697%
79	12.250	3432	3445	3454	rBV	330410	498603	18.63%	1.479%
80	12.356	3468	3478	3484	rBV	92623	166267	6.21%	0.493%
81	12.552	3530	3539	3549	rVB	64932	101798	3.80%	0.302%
82	12.649	3561	3569	3577	rVB	139745	198187	7.40%	0.588%
83	12.704	3577	3586	3603	rVB	219205	345133	12.89%	1.023%
84	12.787	3603	3612	3617	rBV	50224	79335	2.96%	0.235%
85	12.919	3643	3653	3659	rBV6	29150	54983	2.05%	0.163%
86	12.964	3659	3667	3680	rVB	172655	259597	9.70%	0.770%
87	13.032	3680	3688	3696	rBV2	38879	54493	2.04%	0.162%
88	13.086	3696	3705	3708	rBV3	41563	62811	2.35%	0.186%
89	13.125	3708	3717	3731	rVB2	232563	375841	14.04%	1.115%
90	13.237	3744	3752	3760	rBV	42893	62333	2.33%	0.185%
91	13.289	3760	3768	3789	rVB7	39480	80152	2.99%	0.238%
92	13.456	3812	3820	3828	rBV2	38938	59062	2.21%	0.175%
93	13.507	3828	3836	3844	rBV3	74787	126007	4.71%	0.374%
94	13.742	3899	3909	3919	rVB	216276	324617	12.13%	0.963%
95	13.954	3966	3975	3984	rBV5	23635	47405	1.77%	0.141%
96	14.150	4026	4036	4050	rVB2	45068	80058	2.99%	0.237%
97	14.337	4083	4094	4107	rBV3	36074	78742	2.94%	0.234%
98	14.543	4149	4158	4169	rVB4	22191	42208	1.58%	0.125%
99	14.877	4250	4262	4290	rVB	146979	253436	9.47%	0.752%
100	15.064	4309	4320	4334	rBV	65258	107848	4.03%	0.320%

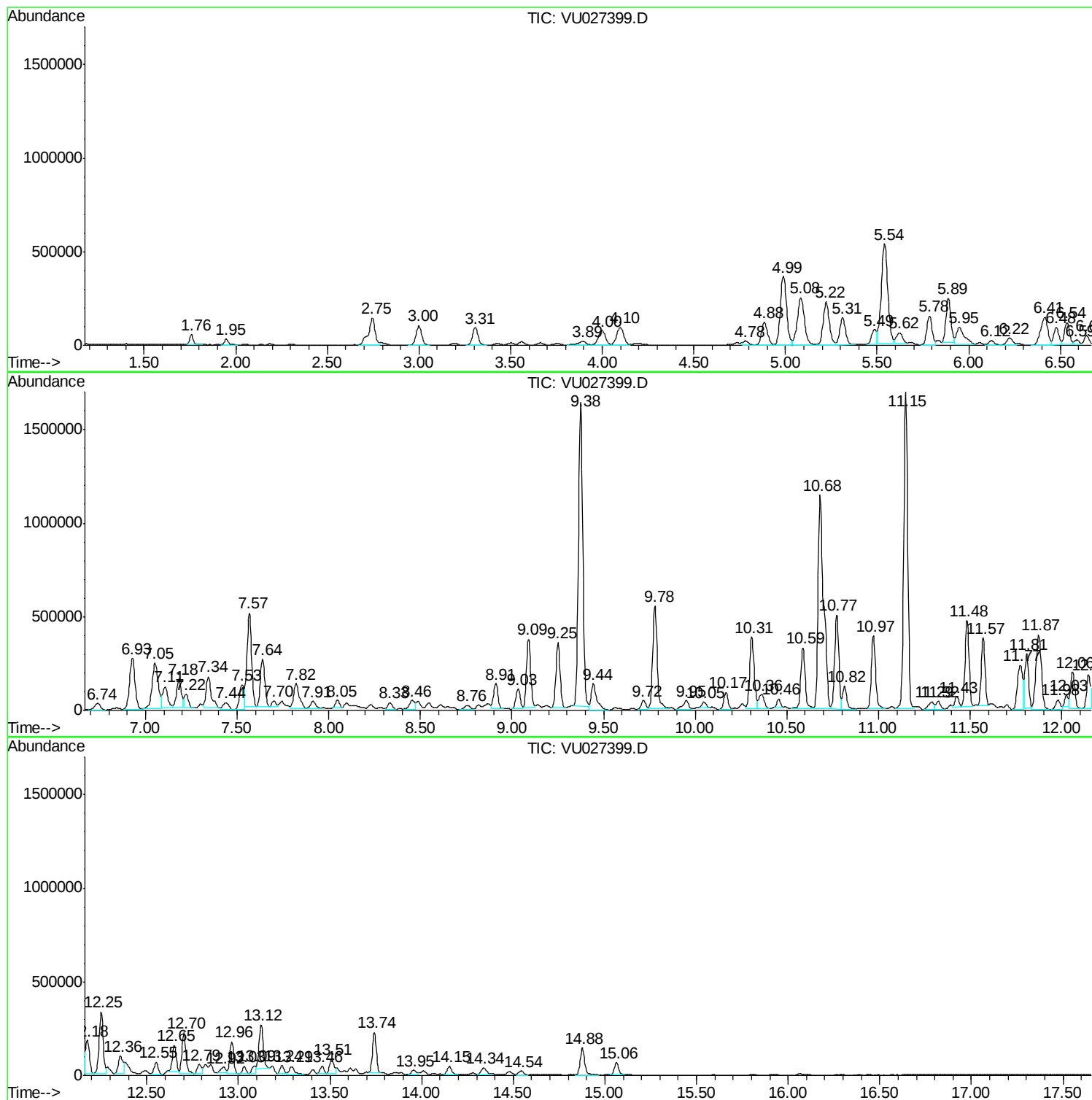
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Instrument :
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B-6(8-8.5)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U100118W.M
Quant Title : SW846 8260

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



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B-6(8-8.5)

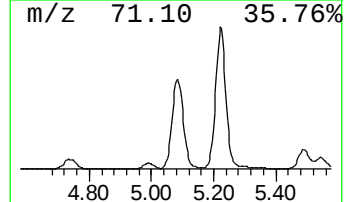
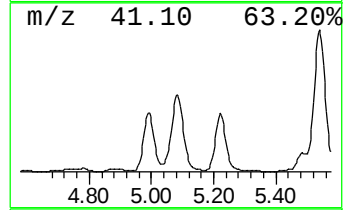
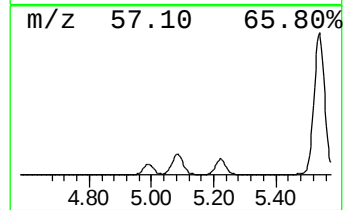
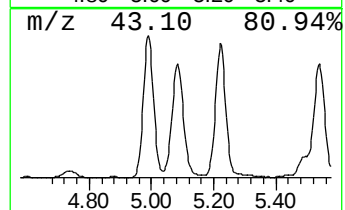
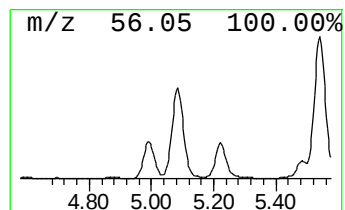
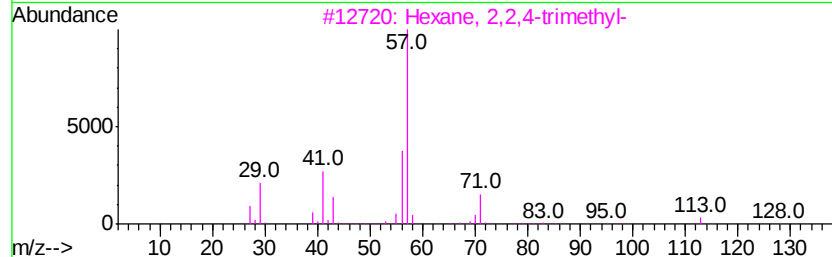
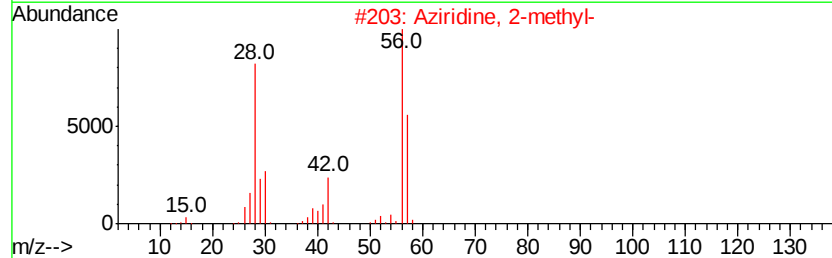
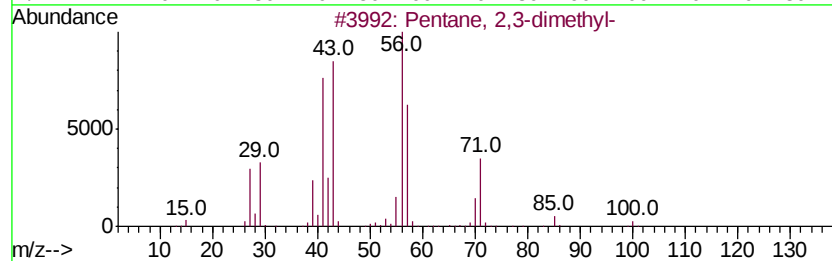
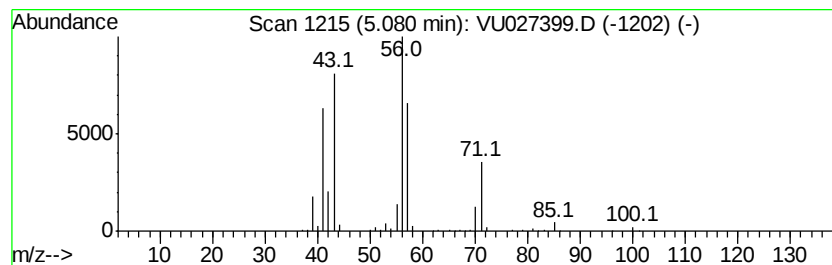
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U100118W.M
Quant Title : SW846 8260

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

Peak Number 1 Pentane, 2,3-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.08	39.62 ug/l	685187	Pentafluorobenzene	4.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	94
2		Aziridine, 2-methyl-	57	C3H7N	000075-55-8	38
3		Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	36
4		Oxalic acid, isobutyl pentyl ester	216	C11H20O4	1000309-37-0	35
5		Hexane, 3-methyl-	100	C7H16	000589-34-4	35



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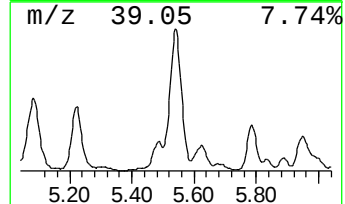
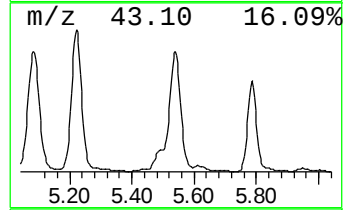
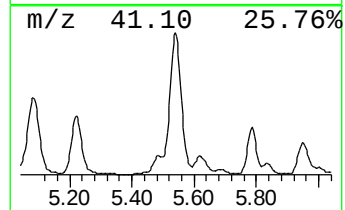
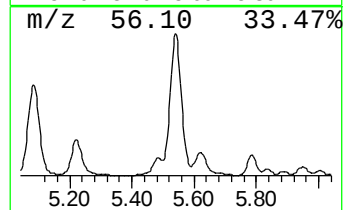
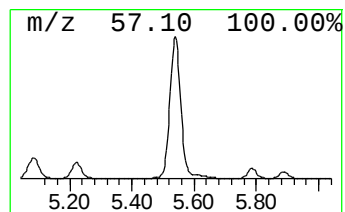
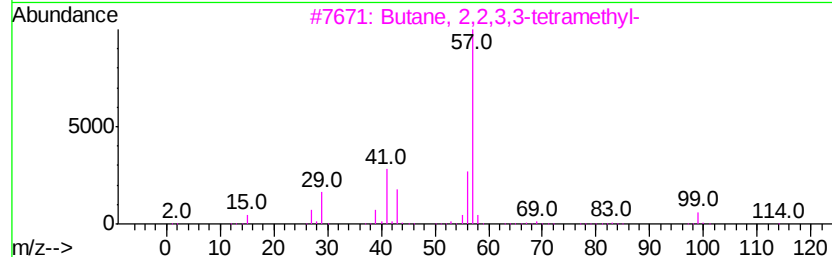
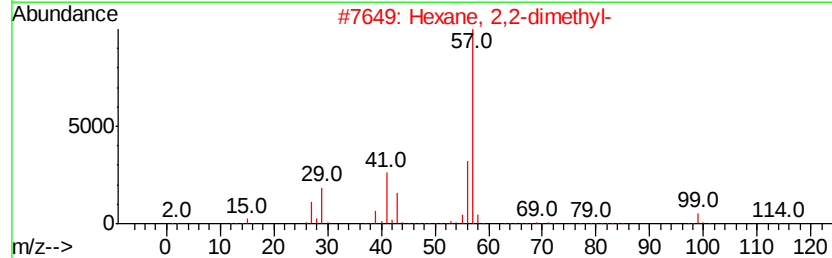
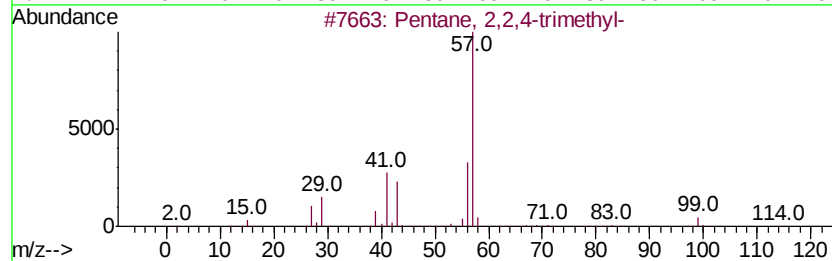
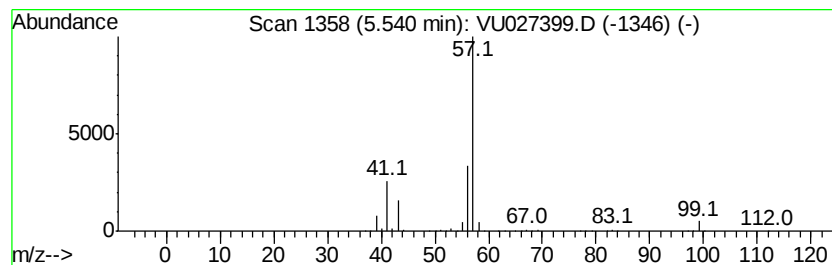
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Quant Title : SW846 8260

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

Peak Number 2 Pentane, 2,2,4-trimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.54	150.48 ug/l	1330540	1,4-Difluorobenzene	5.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	83
2		Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	83
3		Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	83
4		Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	78
5		Heptane, 2,2,4,6,6-pentamethyl-	170	C12H26	013475-82-6	64



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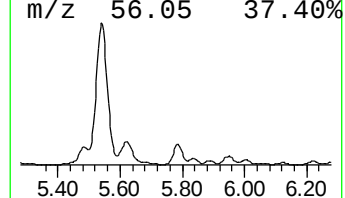
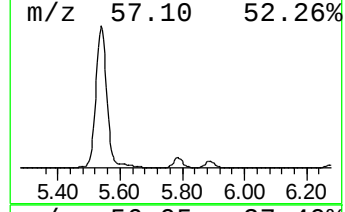
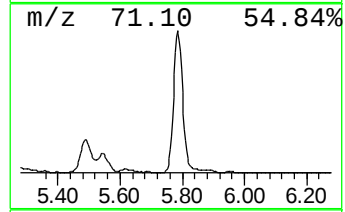
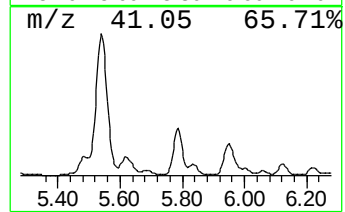
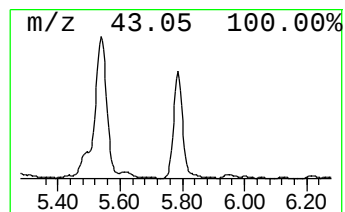
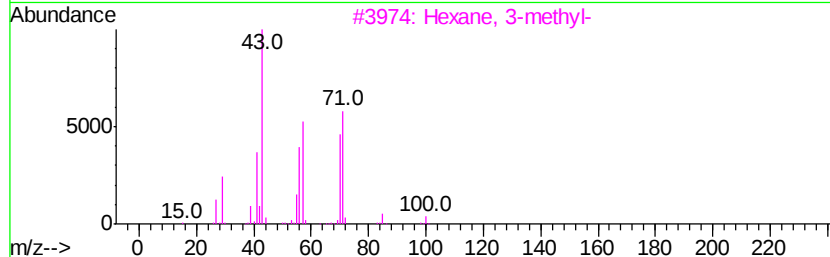
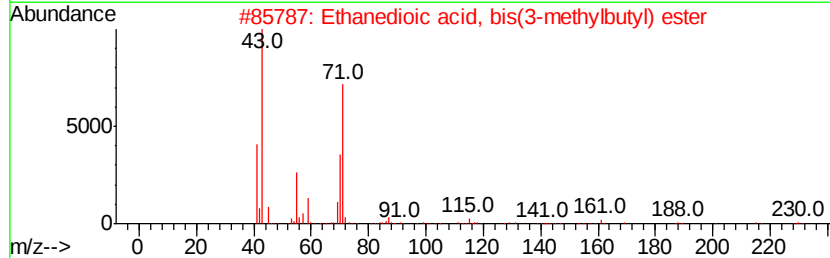
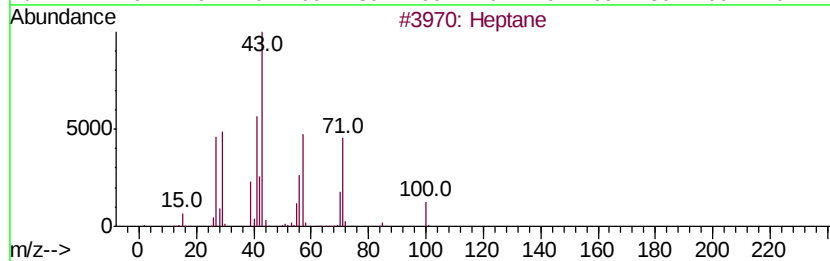
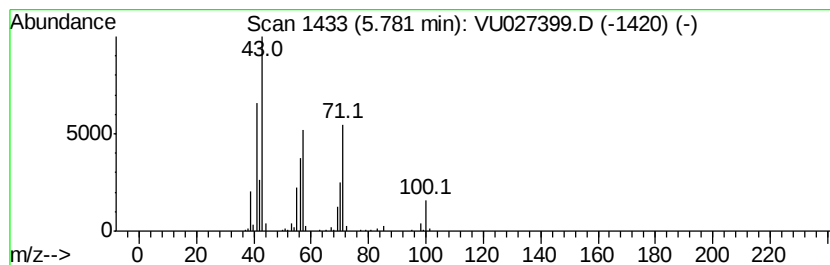
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Quant Title : SW846 8260

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

Peak Number 3 Heptane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.78	35.21 ug/l	311301	1,4-Difluorobenzene	5.89

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane		100	C7H16	000142-82-5	90
2	Ethanedioic acid, bis(3-methylbu...		230	C12H22O4	002051-00-5	50
3	Hexane, 3-methyl-		100	C7H16	000589-34-4	47
4	Hexane, 2,3-dimethyl-		114	C8H18	000584-94-1	38
5	1-Hexene, 3,4,5-trimethyl-		126	C9H18	056728-10-0	38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100418\
Data File : VU027399.D
Acq On : 04 Oct 2018 11:36
Operator : MD/SY
Sample : J5084-12 50X
Misc : 6.42g/5mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
B-6(8-8.5)

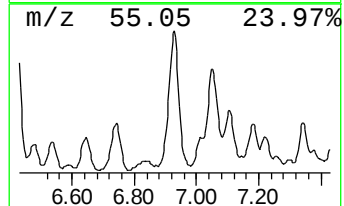
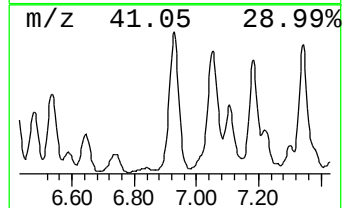
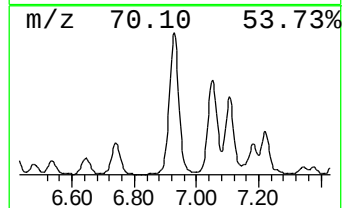
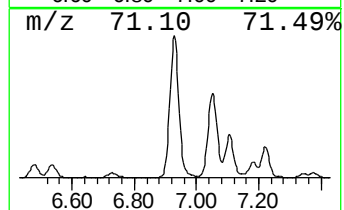
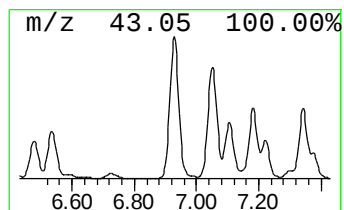
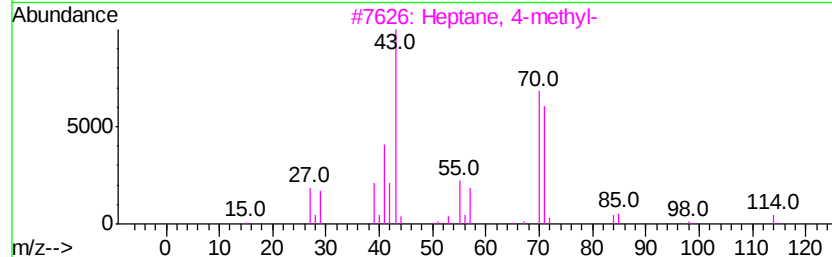
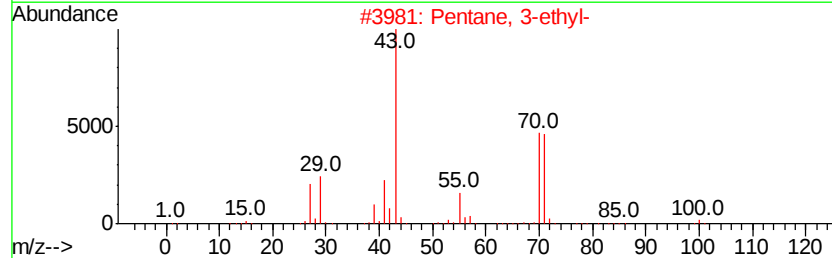
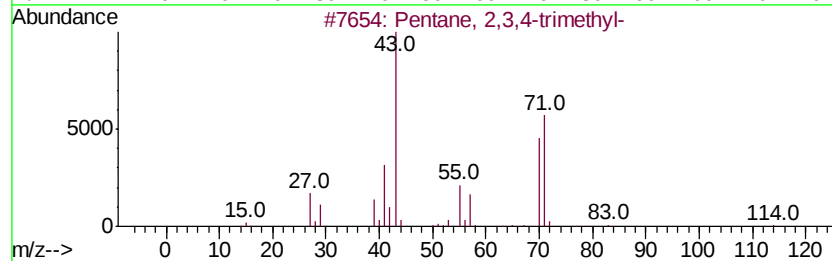
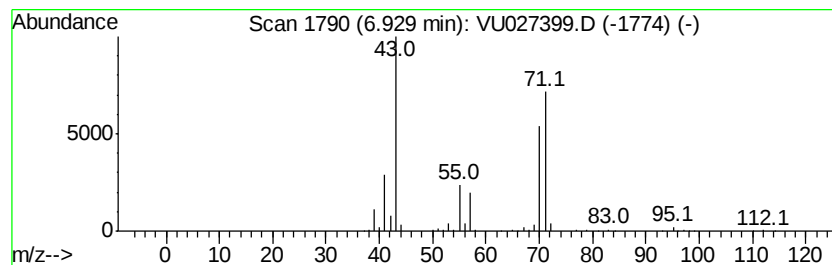
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Quant Title : SW846 8260

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

Peak Number 4 Pentane, 2,3,4-trimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.93	64.30 ug/l	568572	1,4-Difluorobenzene	5.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	91
2		Pentane, 3-ethyl-	100	C7H16	000617-78-7	83
3		Heptane, 4-methyl-	114	C8H18	000589-53-7	83
4		Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	74
5		Decane, 3,3,4-trimethyl-	184	C13H28	049622-18-6	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100418\
Data File : VU027399.D
Acq On : 04 Oct 2018 11:36
Operator : MD/SY
Sample : J5084-12 50X
Misc : 6.42g/5mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
B-6(8-8.5)

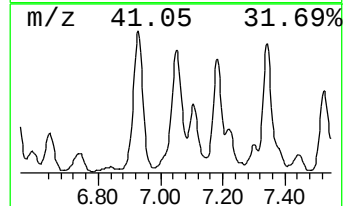
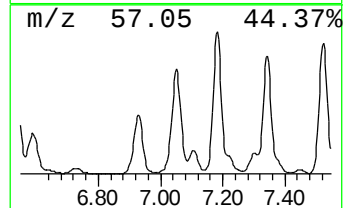
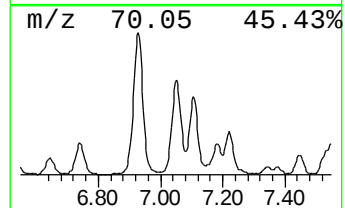
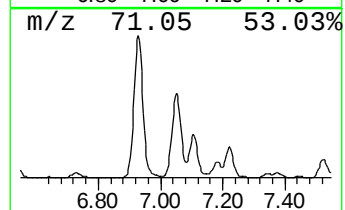
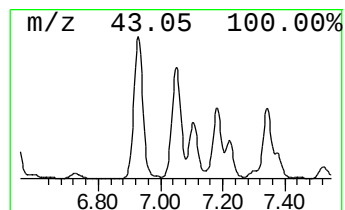
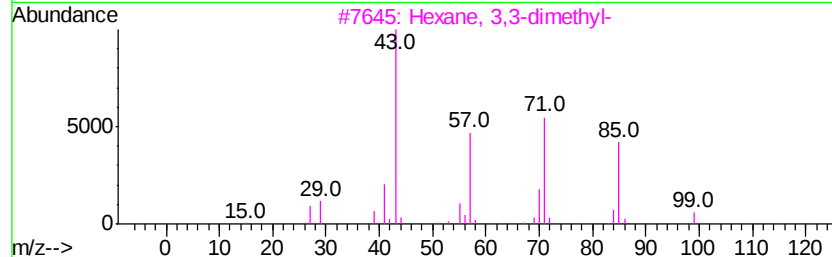
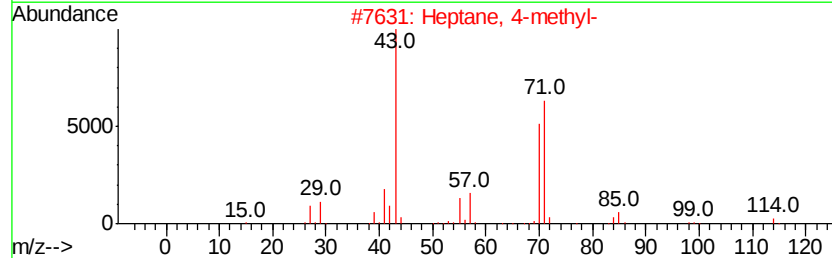
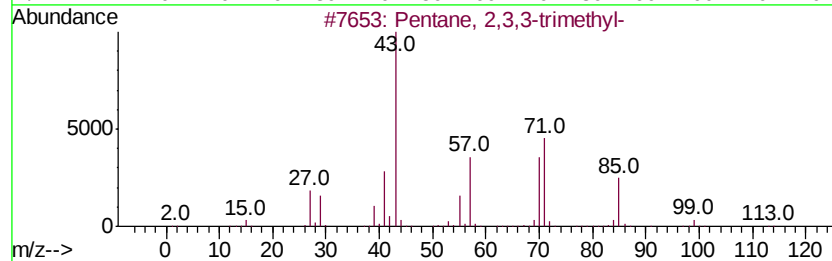
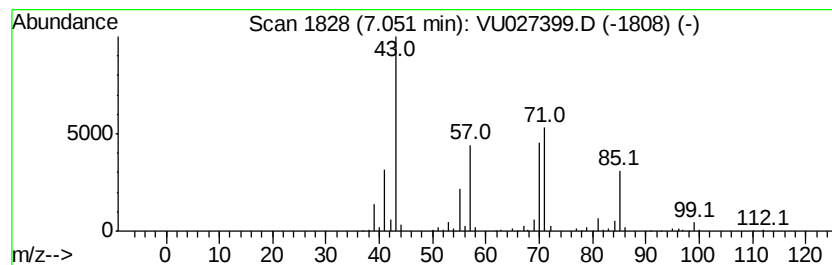
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Quant Title : SW846 8260

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

Peak Number 5 Pentane, 2,3,3-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.05	66.09 ug/l	584387	1,4-Difluorobenzene	5.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	90
2		Heptane, 4-methyl-	114	C8H18	000589-53-7	64
3		Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	59
4		Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	59
5		Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	59



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100418\
Data File : VU027399.D
Acq On : 04 Oct 2018 11:36
Operator : MD/SY
Sample : J5084-12 50X
Misc : 6.42g/5mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
B-6(8-8.5)

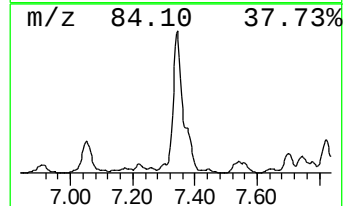
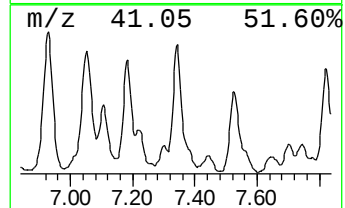
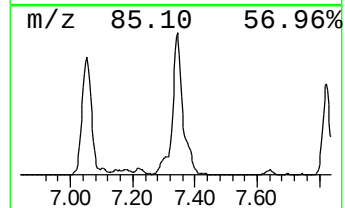
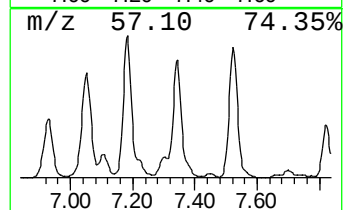
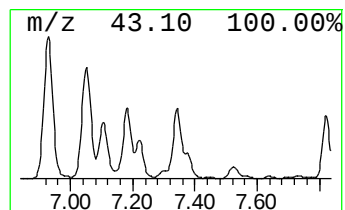
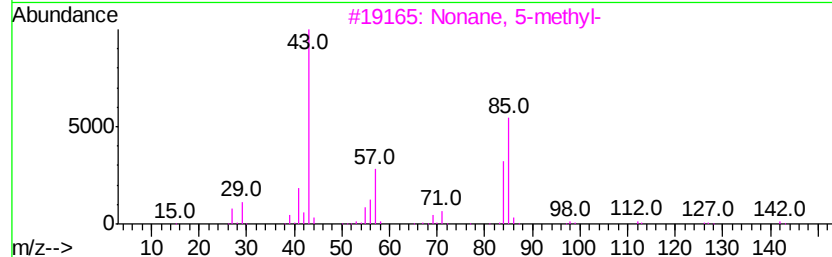
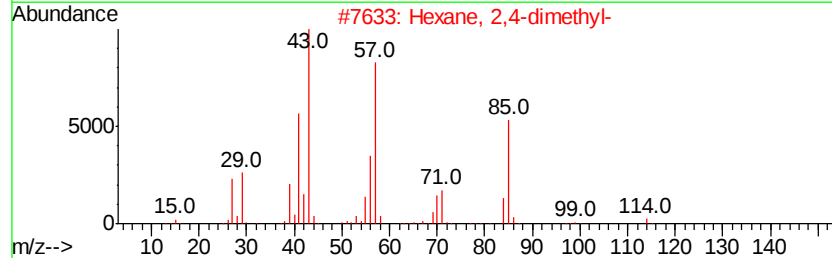
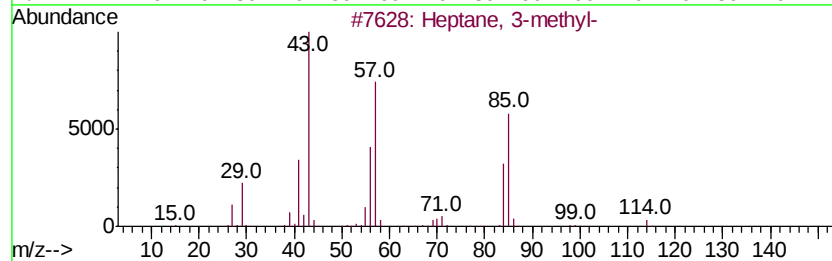
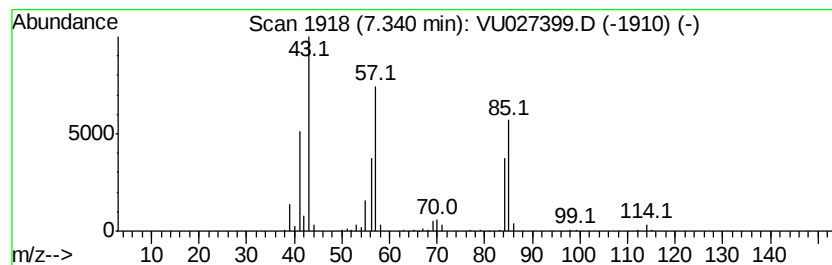
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Quant Title : SW846 8260

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

Peak Number 6 Heptane, 3-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.34	38.28 ug/l	338479	1,4-Difluorobenzene	5.89

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 3-methyl-		114	C8H18	000589-81-1	91
2	Hexane, 2,4-dimethyl-		114	C8H18	000589-43-5	80
3	Nonane, 5-methyl-		142	C10H22	015869-85-9	72
4	Heptane, 3,4,5-trimethyl-		142	C10H22	020278-89-1	72
5	Hexane, 1-(hexyloxy)-2-methyl-		200	C13H28O	074421-17-3	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100418\
Data File : VU027399.D
Acq On : 04 Oct 2018 11:36
Operator : MD/SY
Sample : J5084-12 50X
Misc : 6.42g/5mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 9 Sample Multiplier: 1

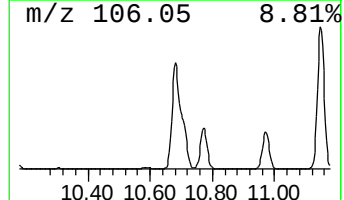
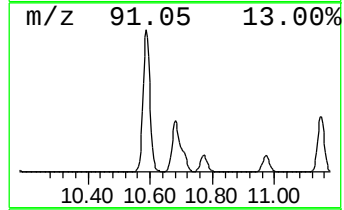
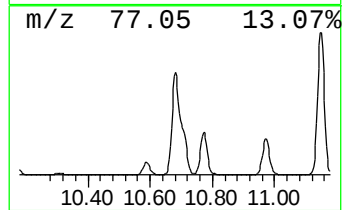
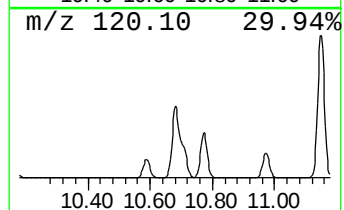
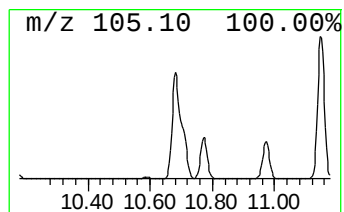
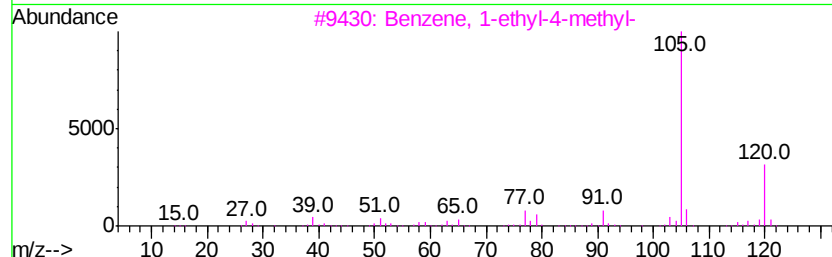
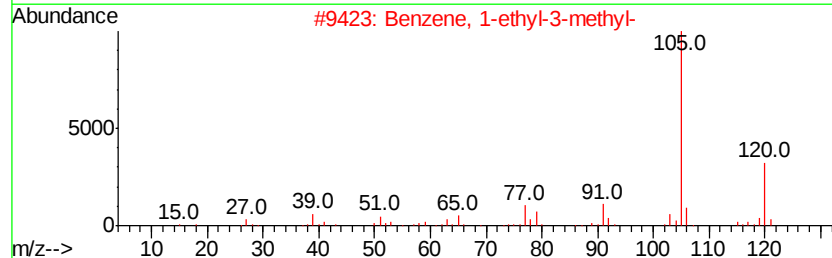
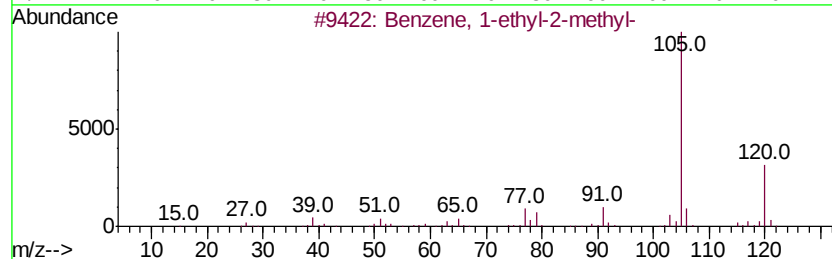
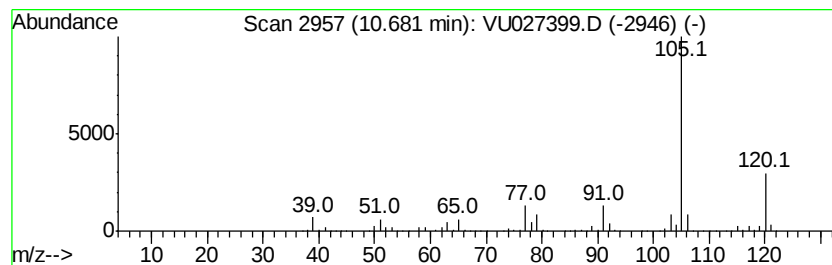
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ClientSampleId :
B-6(8-8.5)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U100118W.M
Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1-ethyl-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.68	171.19 ug/l	2398810	1,4-Dichlorobenzene-d4	11.48
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-ethyl-2-methyl-	120 C9H12	000611-14-3 95
2		Benzene, 1-ethyl-3-methyl-	120 C9H12	000620-14-4 95
3		Benzene, 1-ethyl-4-methyl-	120 C9H12	000622-96-8 91
4		Mesitylene	120 C9H12	000108-67-8 91
5		Benzene, 1,2,4-trimethyl-	120 C9H12	000095-63-6 91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100418\
Data File : VU027399.D
Acq On : 04 Oct 2018 11:36
Operator : MD/SY
Sample : J5084-12 50X
Misc : 6.42g/5mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
B-6(8-8.5)

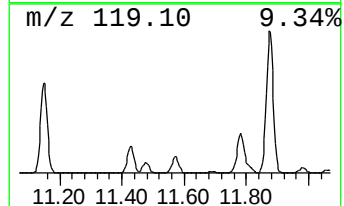
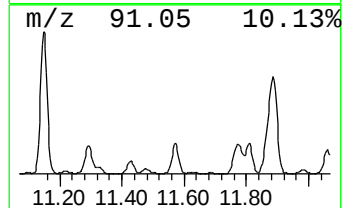
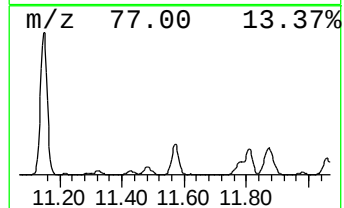
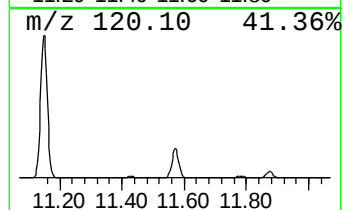
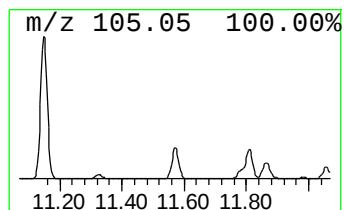
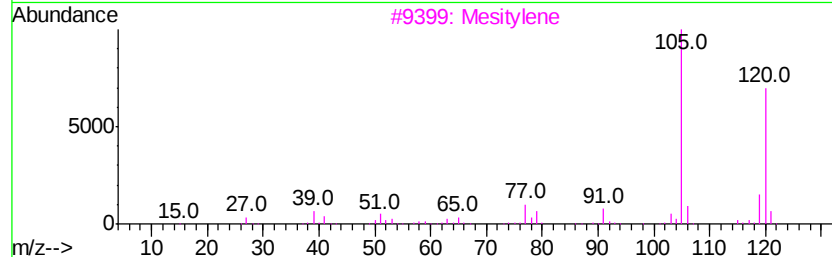
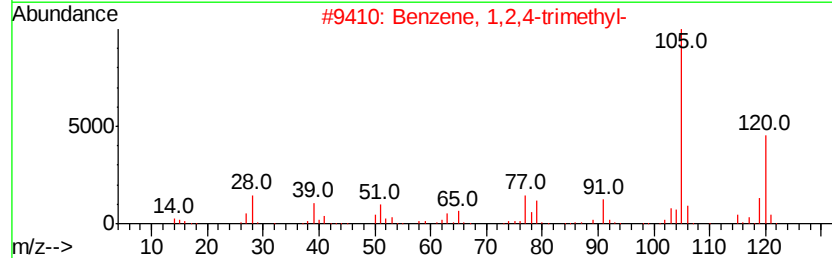
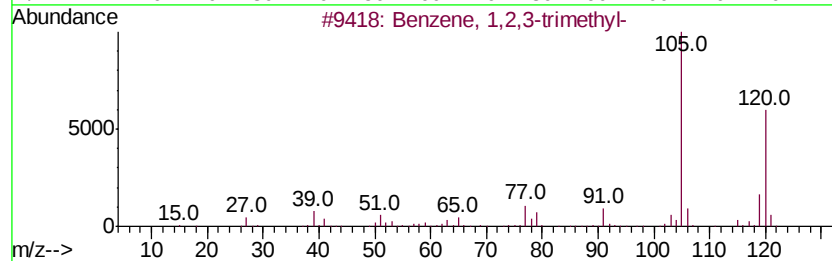
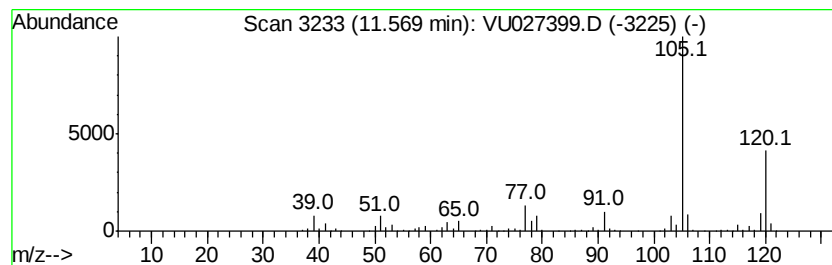
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.57	36.93 ug/l	517447	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
3		Mesitylene	120	C9H12	000108-67-8	91
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90
5		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100418\
Data File : VU027399.D
Acq On : 04 Oct 2018 11:36
Operator : MD/SY
Sample : J5084-12 50X
Misc : 6.42g/5mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
B-6(8-8.5)

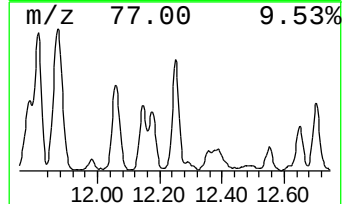
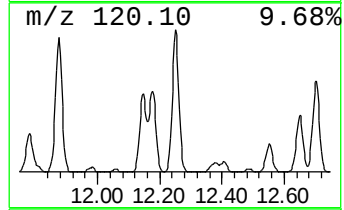
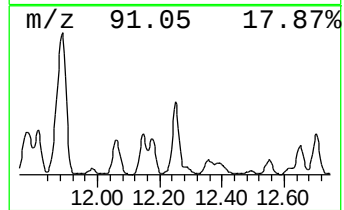
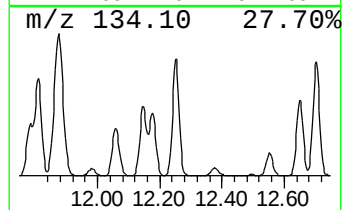
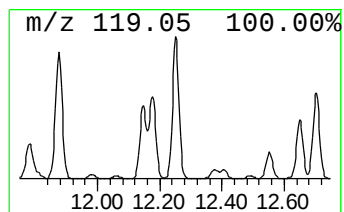
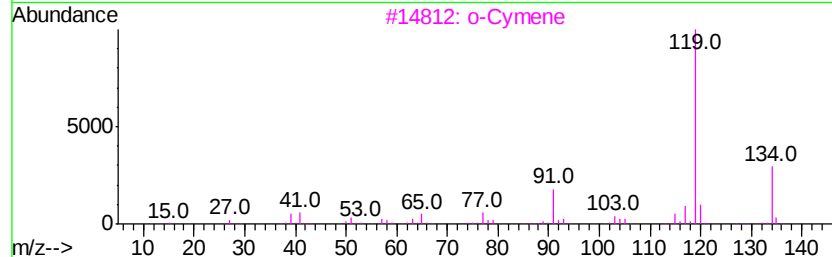
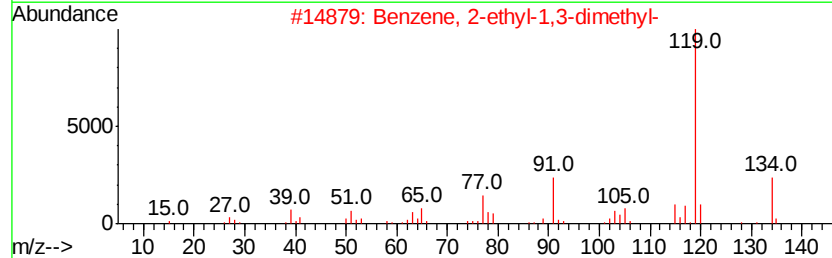
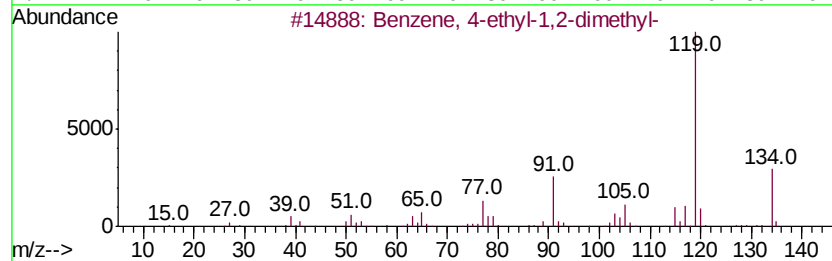
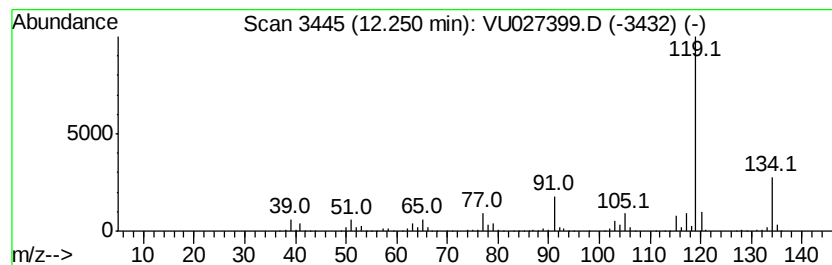
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Quant Title : SW846 8260

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

Peak Number 10 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.25	35.58 ug/l	498603	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
3		o-Cymene	134	C10H14	000527-84-4	95
4		p-Cymene	134	C10H14	000099-87-6	95
5		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU100418\
Data File : VU027399.D
Acq On : 04 Oct 2018 11:36
Operator : MD/SY
Sample : J5084-12 50X
Misc : 6.42g/5mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
B-6(8-8.5)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U100118W.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
Pentane, 2,3-dime...	5.08	39.6	ug/l	685187	1	4.99	864714	50.0
Pentane, 2,2,4-tr...	5.54	150.5	ug/l	1330540	2	5.89	442091	50.0
Heptane	5.78	35.2	ug/l	311301	2	5.89	442091	50.0
Pentane, 2,3,4-tr...	6.93	64.3	ug/l	568572	2	5.89	442091	50.0
Pentane, 2,3,3-tr...	7.05	66.1	ug/l	584387	2	5.89	442091	50.0
Heptane, 3-methyl-	7.34	38.3	ug/l	338479	2	5.89	442091	50.0
Benzene, 1-ethyl-...	10.68	171.2	ug/l	2398810	4	11.48	700633	50.0
Benzene, 1,2,3-tr...	11.57	36.9	ug/l	517447	4	11.48	700633	50.0
Benzene, 4-ethyl-...	12.25	35.6	ug/l	498603	4	11.48	700633	50.0