

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU040221\
 Data File : VU042916.D
 Acq On : 02 Apr 2021 23:16
 Operator : SY/MD
 Sample : M1931-05
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 MW-5

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U032421W.M
 Title : SW846 8260

Signal : TIC: VU042916.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.584	45	76	77	rBV	16801	49081	1.44%	0.084%
2	1.999	194	205	225	rVB	564279	783582	23.07%	1.349%
3	2.208	260	270	286	rVB	156354	223296	6.57%	0.384%
4	2.607	381	394	418	rVB	105066	225557	6.64%	0.388%
5	3.054	517	533	537	rBV2	475090	917162	27.00%	1.579%
6	3.092	538	545	561	rVB	865722	1705128	50.20%	2.935%
7	3.372	611	632	654	rBV	773682	1716092	50.52%	2.954%
8	3.710	718	737	760	rVB2	211674	469128	13.81%	0.807%
9	4.308	903	923	945	rBV5	27361	101090	2.98%	0.174%
10	4.446	947	966	980	rVV	219407	565536	16.65%	0.973%
11	4.543	980	996	1013	rVV	441094	1069988	31.50%	1.842%
12	4.655	1013	1031	1042	rVV4	17241	58365	1.72%	0.100%
13	5.141	1156	1182	1210	rVB3	59156	169339	4.99%	0.291%
14	5.298	1210	1231	1243	rBV	115666	272689	8.03%	0.469%
15	5.385	1243	1258	1271	rBV6	393821	953521	28.07%	1.641%
16	5.472	1272	1285	1308	rVB	645891	1530511	45.06%	2.634%
17	5.600	1308	1325	1336	rBV	250438	552906	16.28%	0.952%
18	5.710	1349	1359	1375	rVB	117316	236932	6.97%	0.408%
19	5.858	1392	1405	1412	rBV	254185	521063	15.34%	0.897%
20	5.912	1413	1422	1438	rVV5	457304	1239213	36.48%	2.133%
21	5.996	1438	1448	1461	rVB	145005	314591	9.26%	0.541%
22	6.141	1479	1493	1514	rBV2	79366	165933	4.88%	0.286%
23	6.256	1515	1529	1538	rVV	249725	489846	14.42%	0.843%
24	6.305	1538	1544	1565	rVV4	102058	250555	7.38%	0.431%
25	6.411	1565	1577	1586	rVV4	25820	51256	1.51%	0.088%
26	6.472	1586	1596	1608	rVB2	28028	49259	1.45%	0.085%
27	6.568	1614	1626	1658	rVB7	32485	102367	3.01%	0.176%
28	6.771	1664	1689	1700	rBV4	409161	1147138	33.77%	1.974%
29	6.874	1712	1721	1731	rVB3	68287	116895	3.44%	0.201%
30	6.932	1731	1739	1747	rBV3	33349	55120	1.62%	0.095%
31	6.989	1747	1757	1769	rVB	104275	188291	5.54%	0.324%
32	7.079	1769	1785	1797	rVB3	48460	104804	3.09%	0.180%
33	7.263	1827	1842	1858	rVB	292099	602284	17.73%	1.037%
34	7.388	1858	1881	1892	rBV	452449	961404	28.30%	1.655%
35	7.626	1939	1955	1961	rVV7	26984	63852	1.88%	0.110%
36	7.774	1990	2001	2013	rVB6	44561	91472	2.69%	0.157%

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Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U032421W.M
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37	7.864	2014	2029	2033	rBV5	95880	191751	5.64%	0.330%
38	7.906	2033	2042	2055	rVB	407122	707325	20.82%	1.217%
39	8.031	2071	2081	2089	rBV6	43131	99258	2.92%	0.171%
40	8.250	2139	2149	2164	rVB4	51158	109395	3.22%	0.188%
41	8.375	2164	2188	2196	rBV3	95425	207255	6.10%	0.357%
42	8.417	2196	2201	2216	rVB2	42622	69323	2.04%	0.119%
43	8.816	2315	2325	2335	rVB3	38693	80552	2.37%	0.139%
44	9.340	2474	2488	2501	rBV9	21997	49763	1.46%	0.086%
45	9.423	2502	2514	2528	rBV	371319	617420	18.18%	1.063%
46	9.517	2538	2543	2552	rVB2	31847	49682	1.46%	0.086%
47	9.578	2552	2562	2580	rVB	120343	200741	5.91%	0.345%
48	9.700	2590	2600	2615	rVB6	30794	77288	2.28%	0.133%
49	10.034	2693	2704	2714	rBV5	19825	44072	1.30%	0.076%
50	10.491	2834	2846	2864	rBV	498158	779611	22.95%	1.342%
51	10.639	2880	2892	2905	rBV	328218	527281	15.52%	0.908%
52	10.912	2957	2977	2996	rBV	2249288	3396881	100.00%	5.846%
53	11.031	2998	3014	3025	rBV	236035	368214	10.84%	0.634%
54	11.095	3025	3034	3044	rVB	38878	57855	1.70%	0.100%
55	11.298	3085	3097	3116	rVB	204577	332120	9.78%	0.572%
56	11.475	3142	3152	3167	rVB	196661	302072	8.89%	0.520%
57	11.616	3182	3196	3201	rBV	256419	400214	11.78%	0.689%
58	11.652	3201	3207	3220	rVB	314627	476645	14.03%	0.820%
59	11.751	3228	3238	3246	rBV	39385	58744	1.73%	0.101%
60	11.819	3246	3259	3269	rBV	446303	713151	20.99%	1.227%
61	11.902	3278	3285	3295	rVB	69409	107611	3.17%	0.185%
62	12.015	3310	3320	3331	rVB	45942	70035	2.06%	0.121%
63	12.105	3334	3348	3363	rBV2	1740176	2631662	77.47%	4.529%
64	12.192	3363	3375	3394	rVV2	634156	1332090	39.22%	2.293%
65	12.308	3399	3411	3423	rVB	246413	378866	11.15%	0.652%
66	12.385	3423	3435	3451	rVB2	470570	875668	25.78%	1.507%
67	12.471	3451	3462	3480	rBV	233684	383498	11.29%	0.660%
68	12.578	3481	3495	3503	rBV	2094686	3142598	92.51%	5.409%
69	12.619	3503	3508	3519	rVB	366751	532087	15.66%	0.916%
70	12.690	3519	3530	3549	rBV2	930241	1801371	53.03%	3.100%
71	12.796	3560	3563	3577	rVB3	27250	48453	1.43%	0.083%
72	12.880	3577	3589	3601	rVB4	143103	275200	8.10%	0.474%
73	12.980	3601	3620	3629	rBV	2095309	3258838	95.94%	5.609%
74	13.034	3629	3637	3651	rVB	590374	898652	26.46%	1.547%
75	13.115	3651	3662	3682	rBV2	327201	557283	16.41%	0.959%
76	13.234	3683	3699	3702	rBV3	111567	213999	6.30%	0.368%

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77	13.259	3702	3707	3712	rVV	121635	179366	5.28%	0.309%
78	13.301	3712	3720	3744	rVB	964308	1520577	44.76%	2.617%
79	13.423	3746	3758	3760	rBV	111608	149210	4.39%	0.257%
80	13.462	3760	3770	3784	rVB3	1815274	3002434	88.39%	5.168%
81	13.568	3795	3803	3815	rVB	256740	390976	11.51%	0.673%
82	13.632	3815	3823	3845	rVB2	185746	317672	9.35%	0.547%
83	13.742	3845	3857	3864	rBV2	202156	347779	10.24%	0.599%
84	13.793	3864	3873	3880	rVV	327951	561948	16.54%	0.967%
85	13.841	3880	3888	3905	rVV5	725812	1452019	42.75%	2.499%
86	13.918	3906	3912	3915	rVV	196641	280882	8.27%	0.483%
87	13.951	3916	3922	3939	rVB2	396368	915011	26.94%	1.575%
88	14.092	3952	3966	3974	rBV	112585	202049	5.95%	0.348%
89	14.198	3992	3999	4012	rVB2	64724	115469	3.40%	0.199%
90	14.295	4012	4029	4039	rBV2	206749	391450	11.52%	0.674%
91	14.352	4039	4047	4057	rVV2	194628	322537	9.50%	0.555%
92	14.494	4079	4091	4104	rBV	299502	469617	13.82%	0.808%
93	14.613	4104	4128	4137	rBV7	30105	90310	2.66%	0.155%
94	14.674	4137	4147	4160	rBV2	244813	463064	13.63%	0.797%
95	14.815	4182	4191	4203	rBV	185308	282636	8.32%	0.486%
96	14.883	4203	4212	4223	rVB2	163841	267282	7.87%	0.460%
97	15.028	4244	4257	4269	rBV4	54575	112293	3.31%	0.193%
98	15.227	4290	4319	4333	rBV	472616	855875	25.20%	1.473%
99	15.417	4366	4378	4389	rVV	524292	826471	24.33%	1.422%
100	16.417	4677	4689	4696	rBV3	27405	45340	1.33%	0.078%

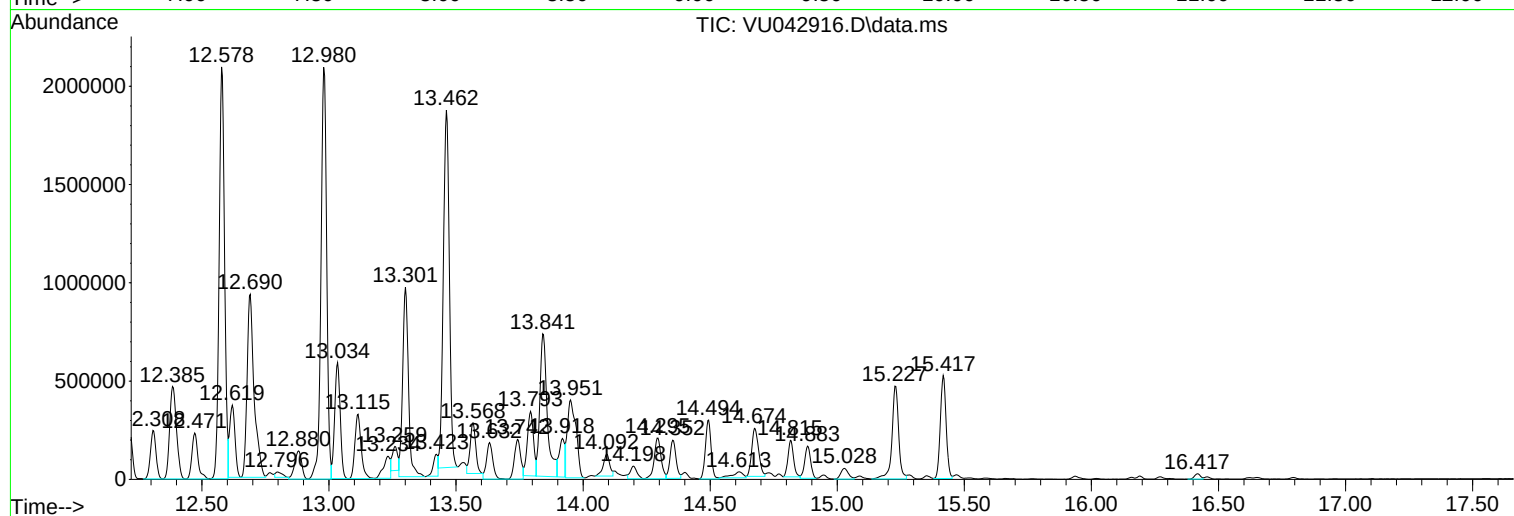
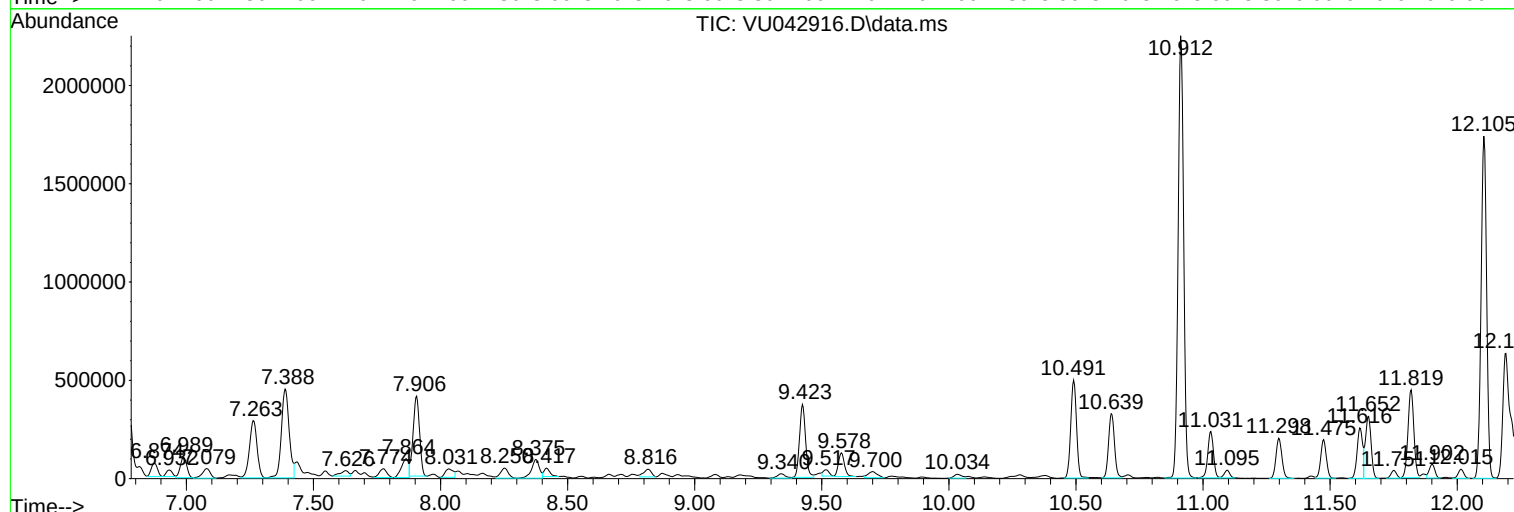
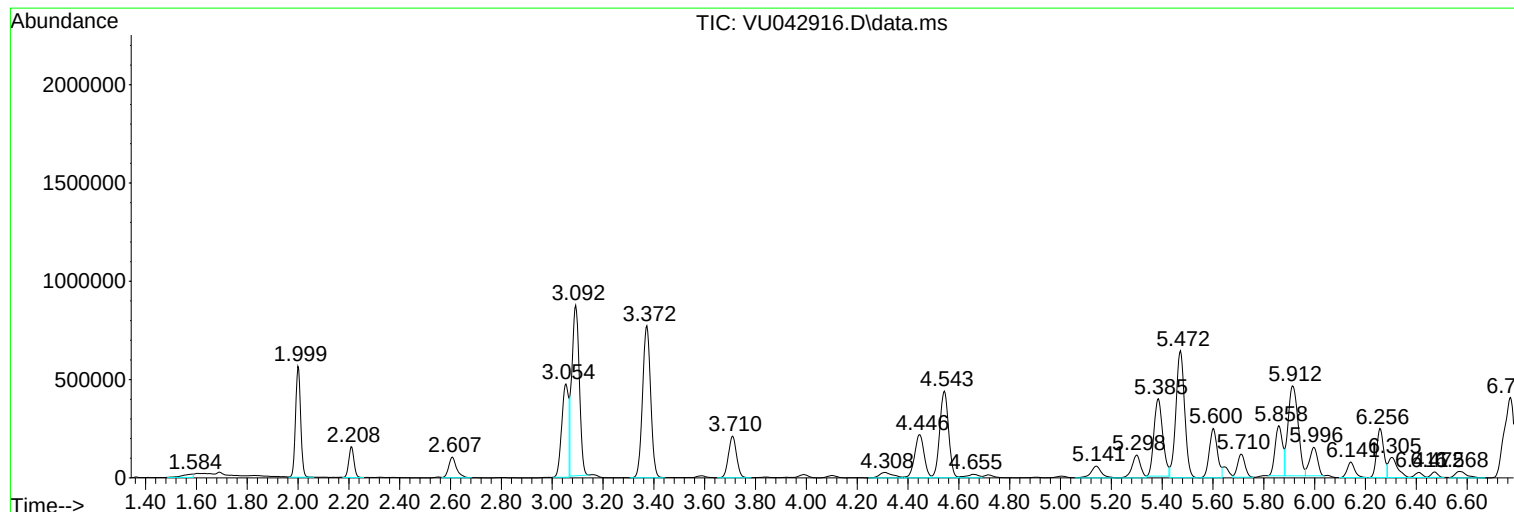
Sum of corrected areas: 58102037

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Instrument :
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ClientSampled :
MW-5

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U032421W.M
Quant Title : SW846 8260

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



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Instrument :
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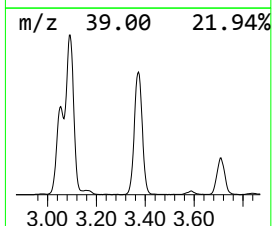
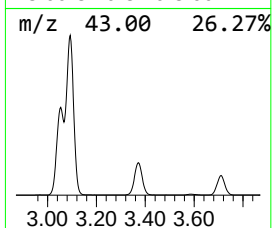
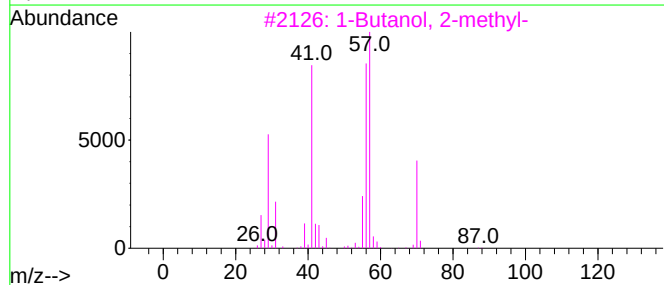
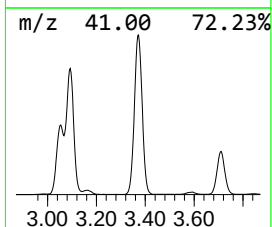
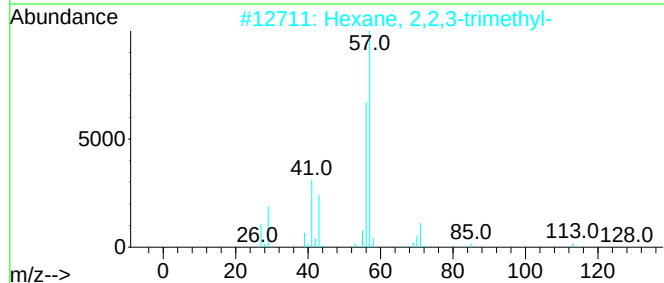
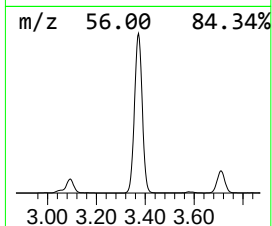
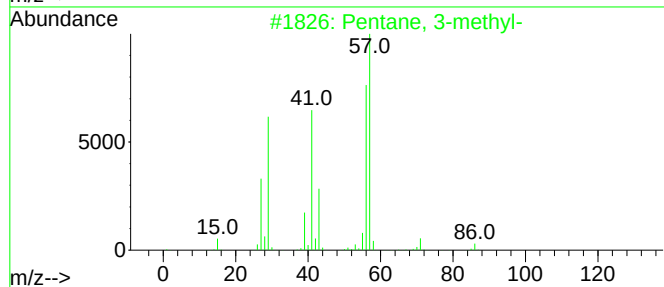
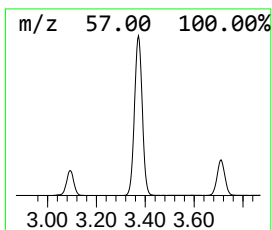
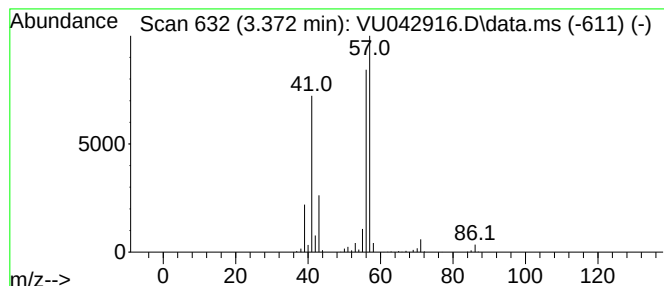
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U032421W.M
Quant Title : SW846 8260

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

Peak Number 1 Pentane, 3-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.372	89.99 ug/l	1716090	Pentafluorobenzene	5.382

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 3-methyl-	86	C6H14	000096-14-0	91
2			Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	43
3			1-Butanol, 2-methyl-	88	C5H12O	000137-32-6	40
4			Butyl isocyanatoacetate	157	C7H11NO3	017046-22-9	40
5			n-Hexane	86	C6H14	000110-54-3	40



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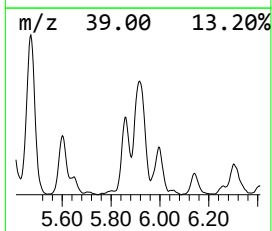
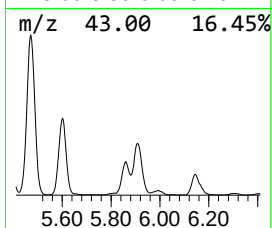
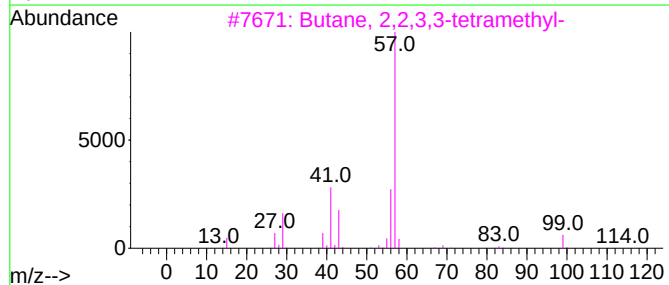
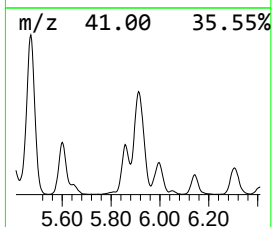
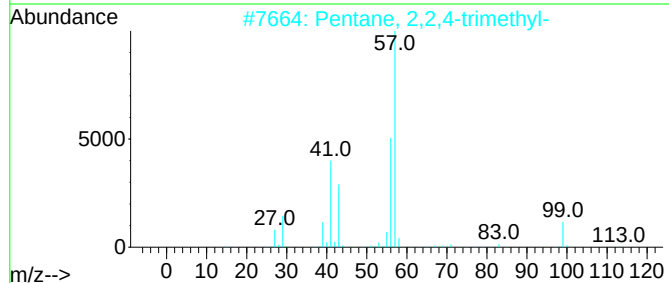
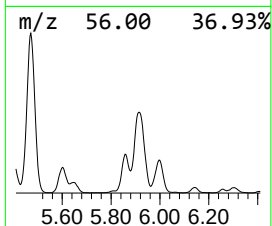
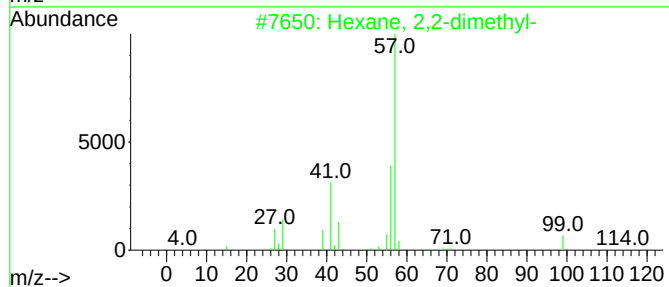
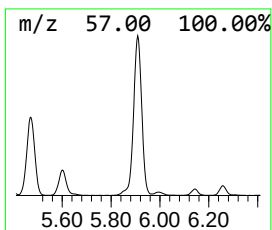
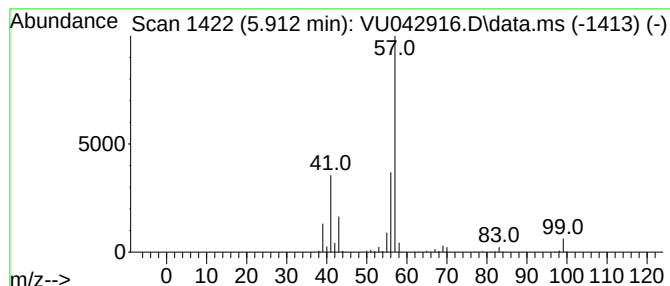
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Quant Title : SW846 8260

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

Peak Number 2 Hexane, 2,2-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.912	126.49 ug/l	1239210	1,4-Difluorobenzene	6.256

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	72
2		Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	72
3		Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	64
4		Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	59
5		Hexane, 2,2,5,5-tetramethyl-	142	C10H22	001071-81-4	59



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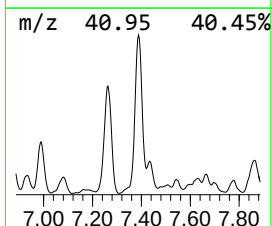
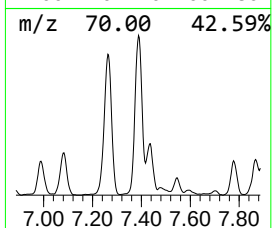
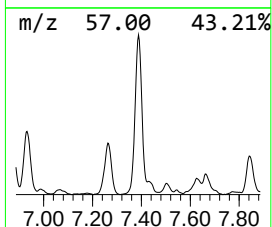
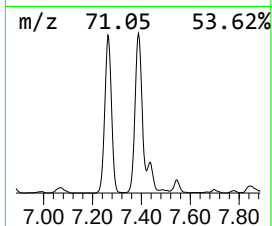
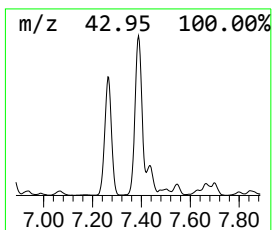
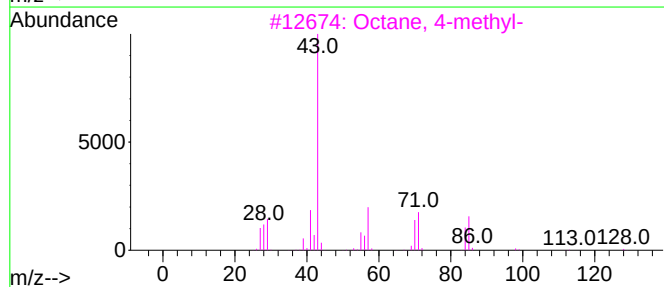
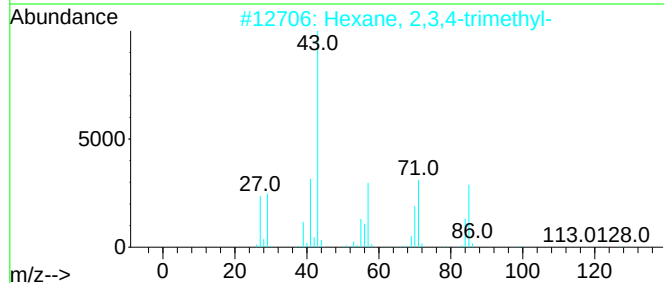
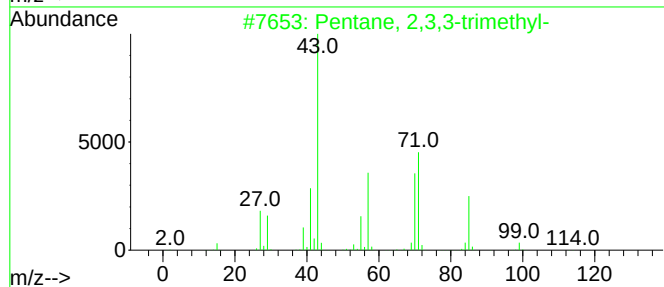
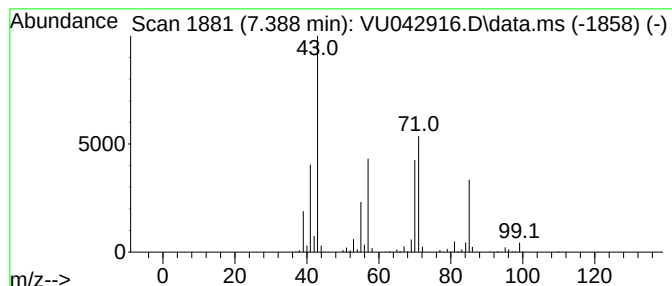
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Quant Title : SW846 8260

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

Peak Number 3 Pentane, 2,3,3-trimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.388	98.13 ug/l	961404	1,4-Difluorobenzene	6.256

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	90
2			Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	78
3			Octane, 4-methyl-	128	C9H20	002216-34-4	78
4			Heptane, 4-methyl-	114	C8H18	000589-53-7	64
5			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU040221\
Data File : VU042916.D
Acq On : 02 Apr 2021 23:16
Operator : SY/MD
Sample : M1931-05
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-5

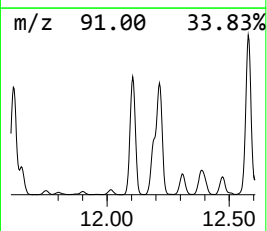
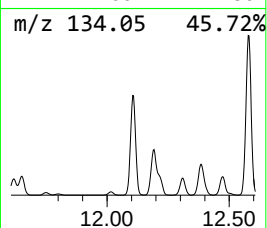
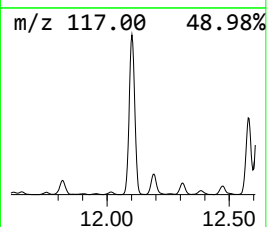
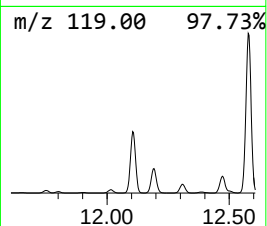
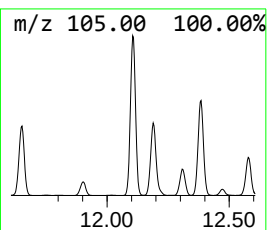
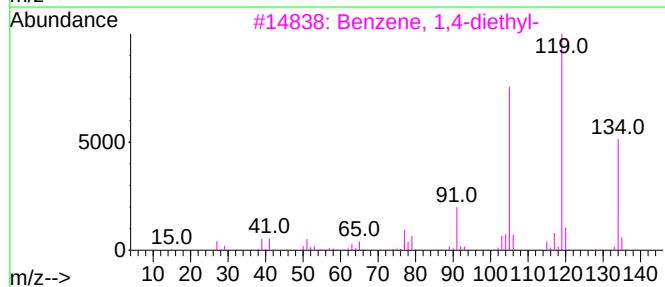
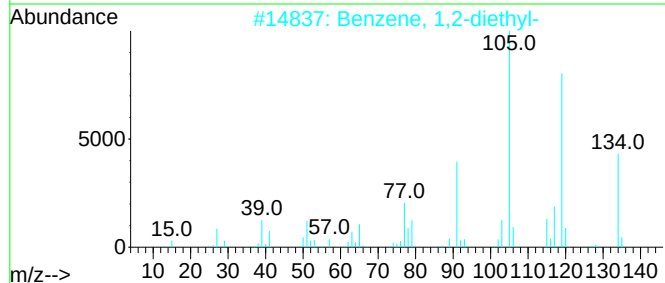
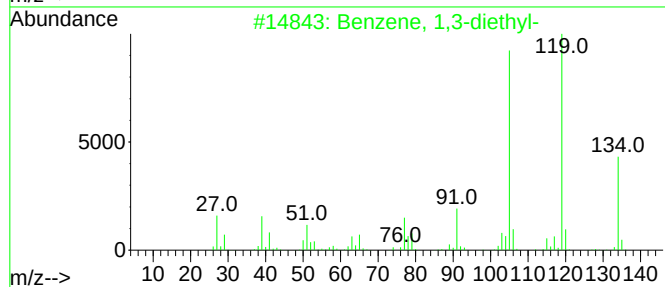
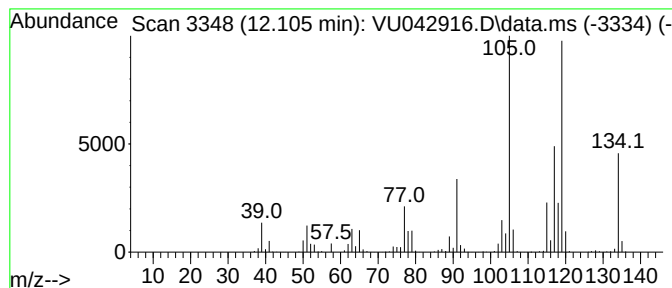
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1,3-diethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.105	184.51 ug/l	2631660	1,4-Dichlorobenzene-d4	11.819

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	92
2			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	89
3			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	76
4			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	53
5			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	49



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU040221\
Data File : VU042916.D
Acq On : 02 Apr 2021 23:16
Operator : SY/MD
Sample : M1931-05
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-5

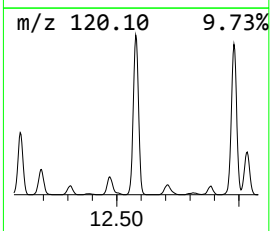
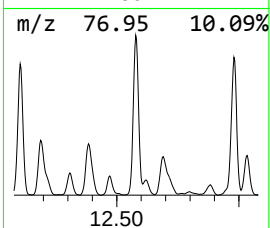
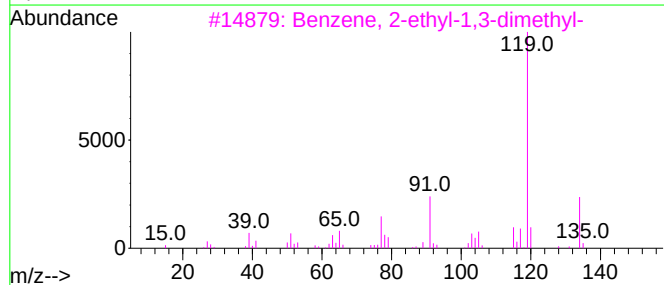
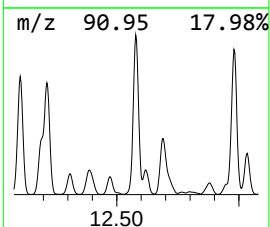
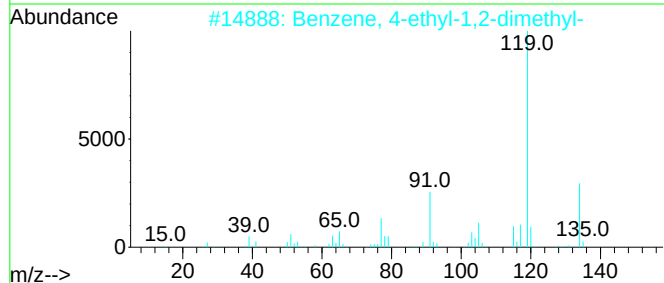
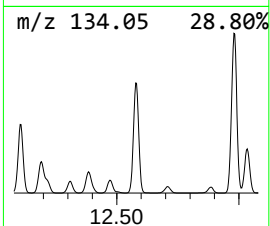
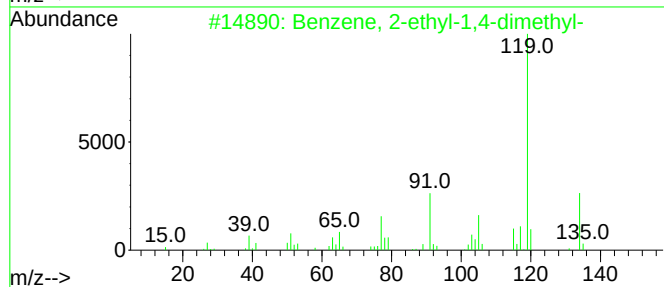
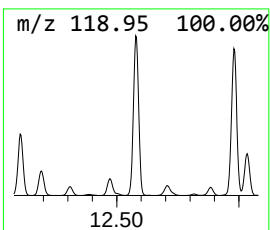
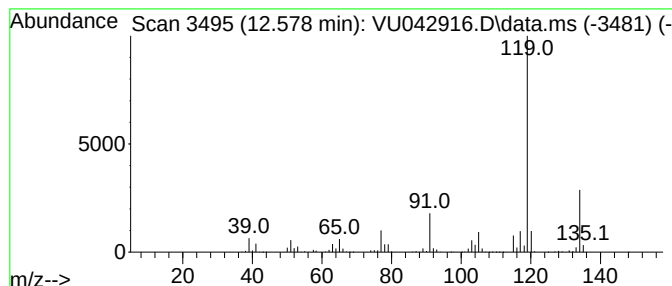
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.578	220.33 ug/l	3142600	1,4-Dichlorobenzene-d4	11.819

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			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			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
4			o-Cymene	134	C10H14	000527-84-4	95
5			p-Cymene	134	C10H14	000099-87-6	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU040221\
Data File : VU042916.D
Acq On : 02 Apr 2021 23:16
Operator : SY/MD
Sample : M1931-05
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-5

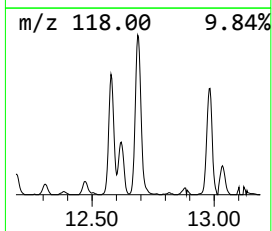
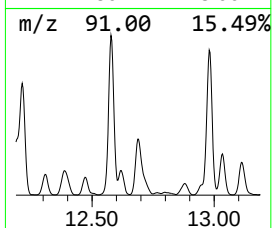
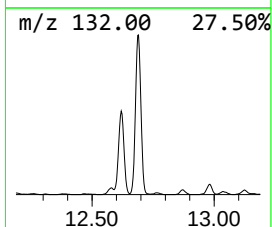
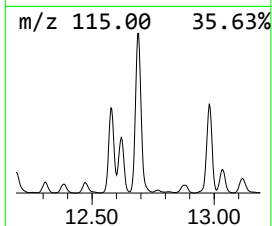
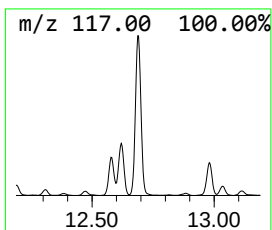
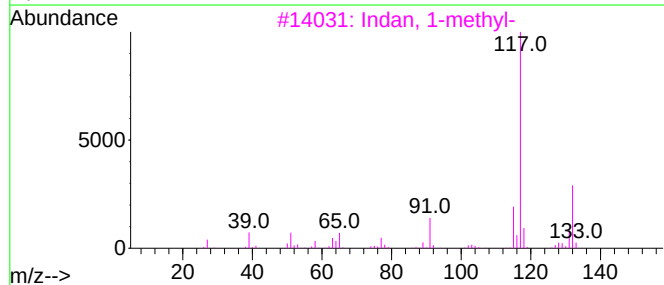
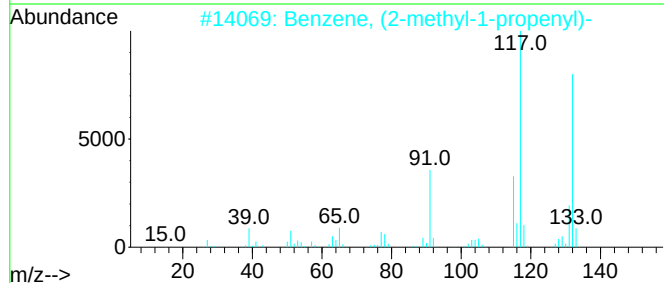
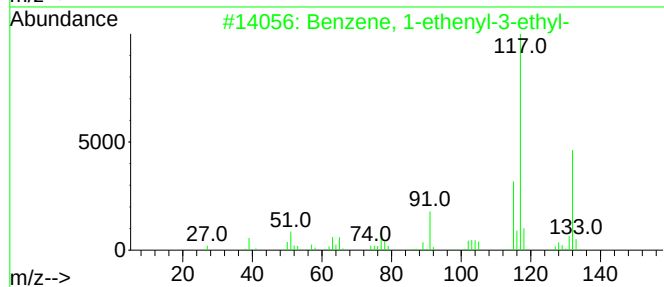
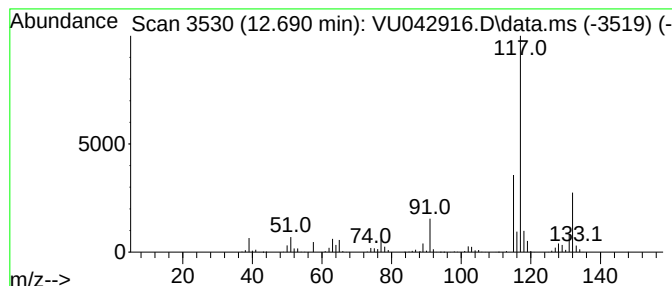
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.690	126.30 ug/l	1801370	1,4-Dichlorobenzene-d4	11.819

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	91
2			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	90
3			Indan, 1-methyl-	132	C10H12	000767-58-8	87
4			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	83
5			3-Phenylbut-1-ene	132	C10H12	000934-10-1	83



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU040221\
Data File : VU042916.D
Acq On : 02 Apr 2021 23:16
Operator : SY/MD
Sample : M1931-05
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-5

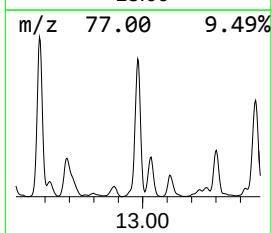
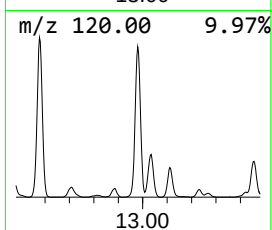
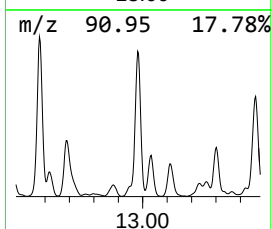
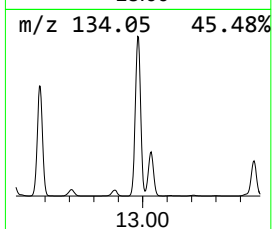
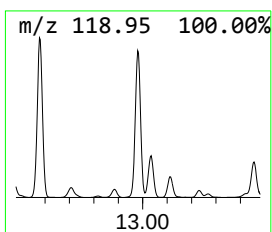
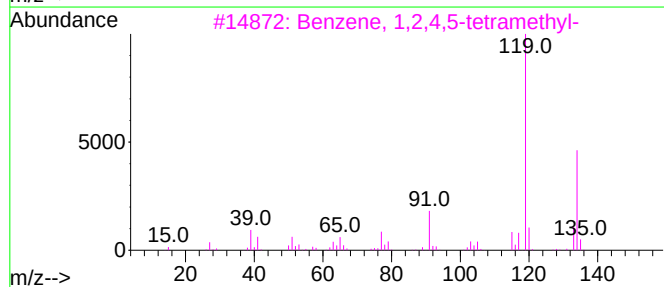
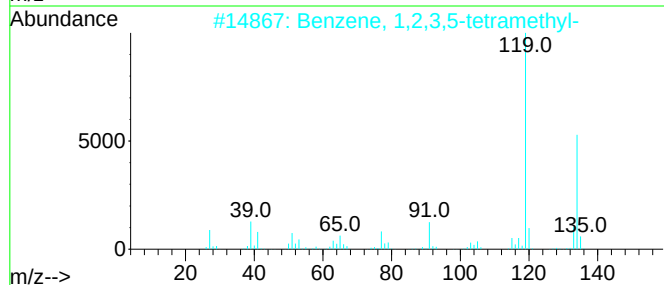
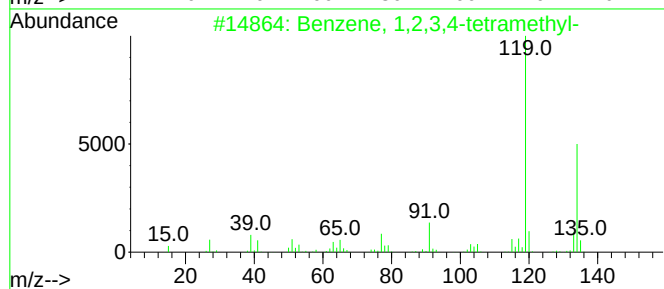
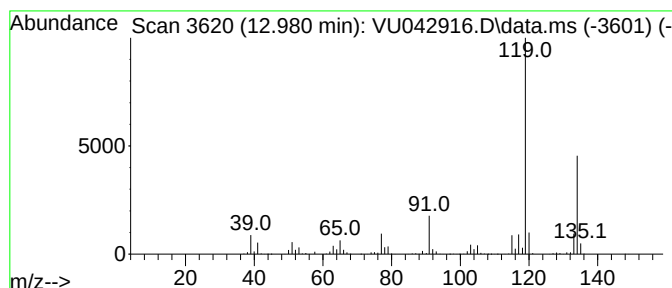
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.980	228.48 ug/l	3258840	1,4-Dichlorobenzene-d4	11.819

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
3			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
4			o-Cymene	134	C10H14	000527-84-4	95
5			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU040221\
Data File : VU042916.D
Acq On : 02 Apr 2021 23:16
Operator : SY/MD
Sample : M1931-05
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampled :
MW-5

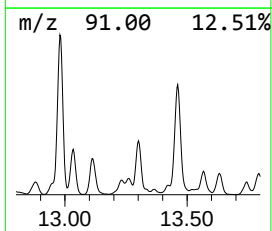
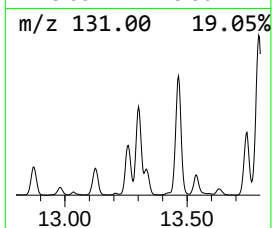
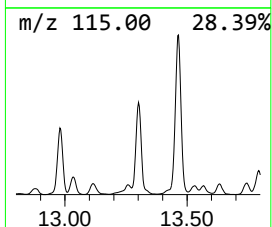
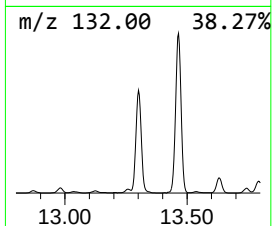
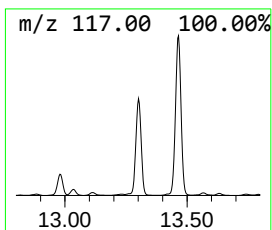
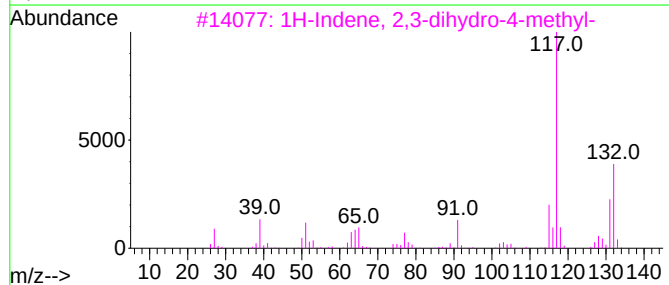
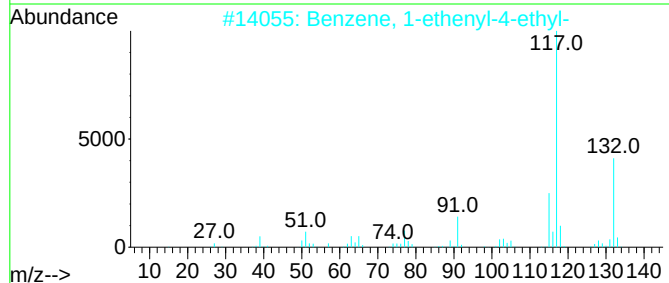
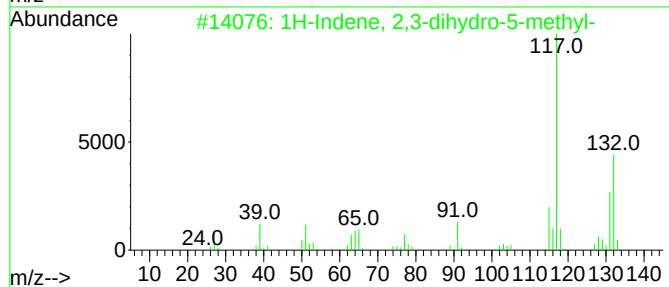
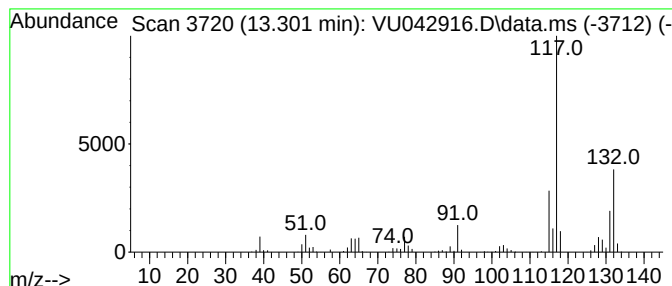
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Quant Title : SW846 8260

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

Peak Number 8 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.301	106.61 ug/l	1520580	1,4-Dichlorobenzene-d4	11.819

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	94	
2	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94	
3	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	93	
4	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92	
5	3-Phenylbut-1-ene	132	C10H12	000934-10-1	90	



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU040221\
Data File : VU042916.D
Acq On : 02 Apr 2021 23:16
Operator : SY/MD
Sample : M1931-05
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-5

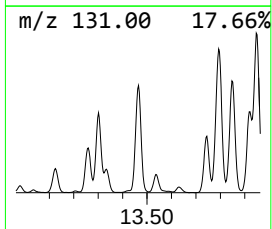
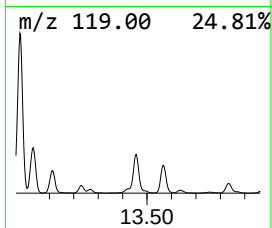
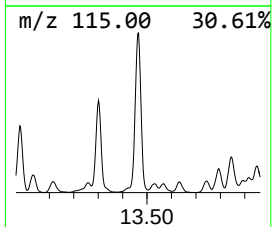
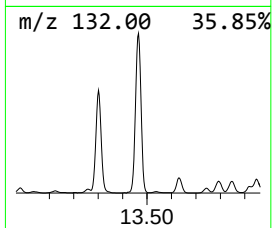
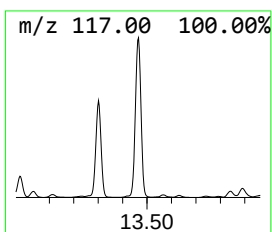
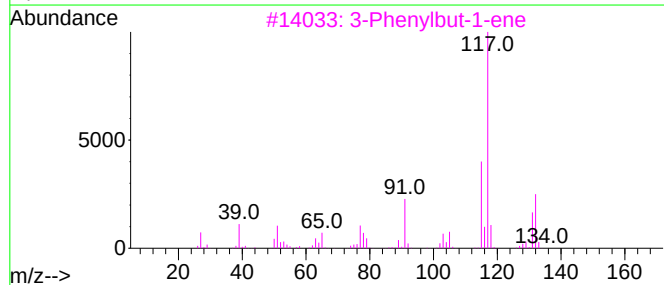
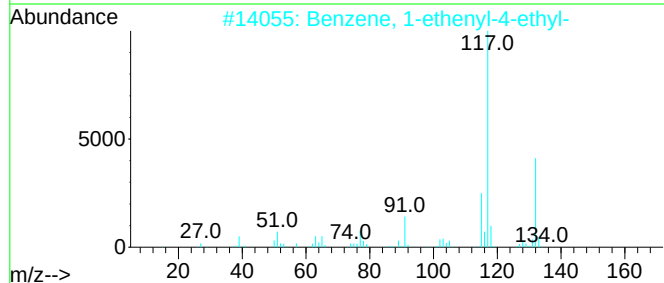
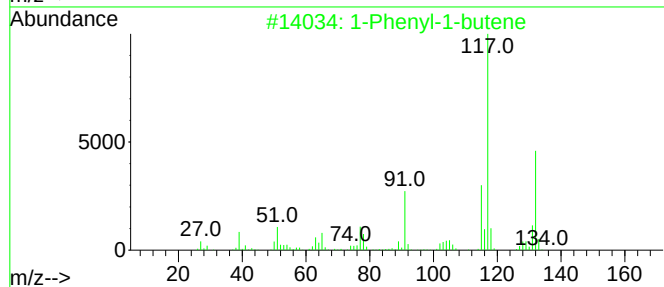
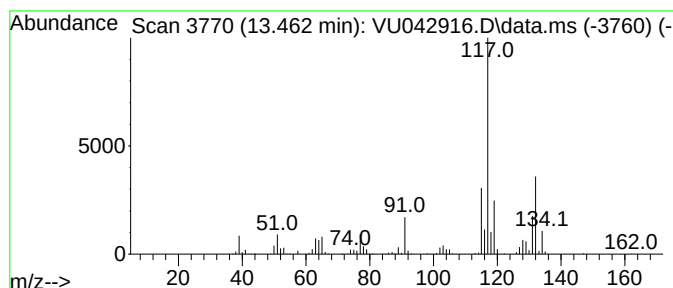
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Quant Title : SW846 8260

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

Peak Number 9 1-Phenyl-1-butene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.462	210.51 ug/l	3002430	1,4-Dichlorobenzene-d4	11.819

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Phenyl-1-butene	132	C10H12	000824-90-8	90
2		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	89
3		3-Phenylbut-1-ene	132	C10H12	000934-10-1	81
4		Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	81
5		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	78



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU040221\
Data File : VU042916.D
Acq On : 02 Apr 2021 23:16
Operator : SY/MD
Sample : M1931-05
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-5

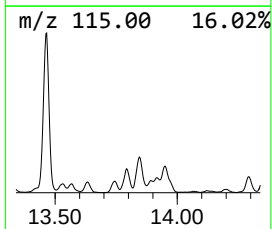
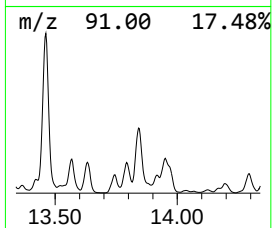
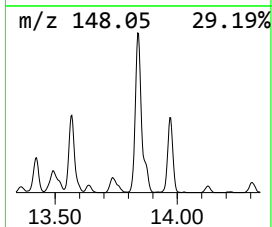
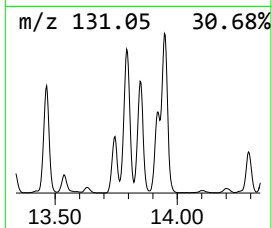
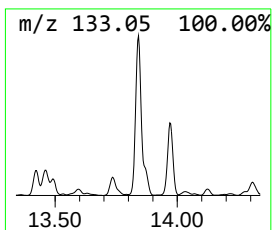
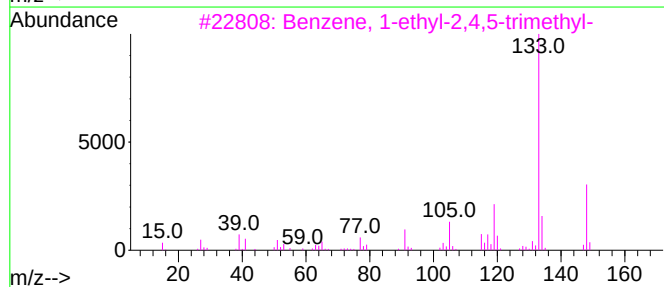
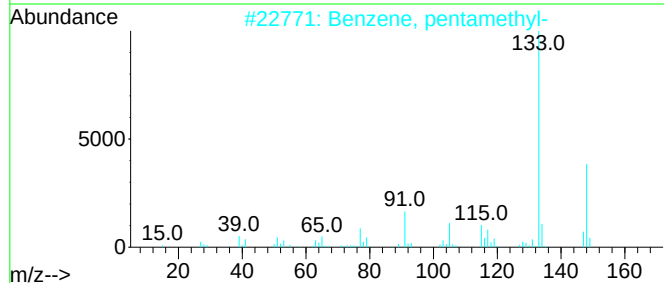
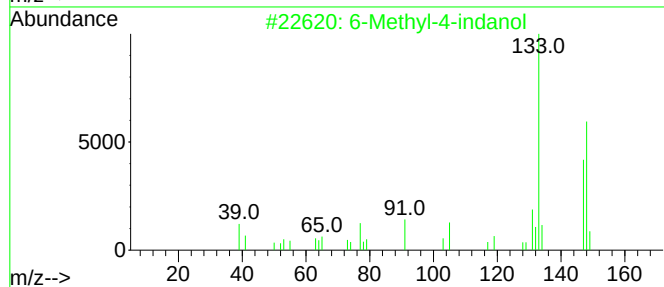
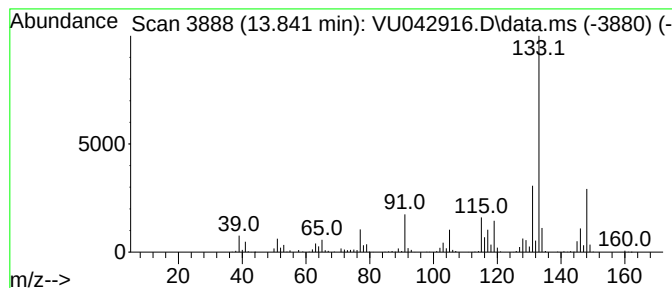
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Quant Title : SW846 8260

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

Peak Number 10 6-Methyl-4-indanol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.841	101.80 ug/l	1452020	1,4-Dichlorobenzene-d4	11.819

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			6-Methyl-4-indanol	148	C10H12O	020294-32-0	80
2			Benzene, pentamethyl-	148	C11H16	000700-12-9	76
3			Benzene, 1-ethyl-2,4,5-trimethyl-	148	C11H16	017851-27-3	74
4			Benzene, 1-ethyl-3-(1-methylethyl)-	148	C11H16	004920-99-4	68
5			2-tert-Butyltoluene	148	C11H16	001074-92-6	68



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU040221\
Data File : VU042916.D
Acq On : 02 Apr 2021 23:16
Operator : SY/MD
Sample : M1931-05
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-5

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U032421W.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, 3-methyl-	3.372	90.0	ug/l	1716090	1	5.382	953521	50.0
Hexane, 2,2-dim...	5.912	126.5	ug/l	1239210	2	6.256	489846	50.0
Pentane, 2,3,3-...	7.388	98.1	ug/l	961404	2	6.256	489846	50.0
Benzene, 1,3-di...	12.105	184.5	ug/l	2631660	4	11.819	713151	50.0
Benzene, 2-ethy...	12.578	220.3	ug/l	3142600	4	11.819	713151	50.0
Benzene, 1-ethe...	12.690	126.3	ug/l	1801370	4	11.819	713151	50.0
Benzene, 1,2,3,...	12.980	228.5	ug/l	3258840	4	11.819	713151	50.0
1H-Indene, 2,3-...	13.301	106.6	ug/l	1520580	4	11.819	713151	50.0
1-Phenyl-1-butene	13.462	210.5	ug/l	3002430	4	11.819	713151	50.0
6-Methyl-4-indanol	13.841	101.8	ug/l	1452020	4	11.819	713151	50.0