

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX021919\
 Data File : VX007615.D
 Acq On : 19 Feb 2019 16:04
 Operator : JC/SP
 Sample : K1528-01
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-1

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_X\METHOD\82X020119W.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	1.300	21	24	38	rBV	6776246	8159409	3.96%	0.526%
2	1.404	38	41	59	rVB2	57275686	73448181	35.64%	4.737%
3	1.794	99	105	126	rVV2	65845256	102442131	49.71%	6.607%
4	1.952	126	131	135	rVV	3176033	4583410	2.22%	0.296%
5	2.001	135	139	153	rVV3	44765255	76979615	37.35%	4.965%
6	2.123	153	159	167	rVV	11928962	17088496	8.29%	1.102%
7	2.221	169	175	178	rVV	5687834	8670126	4.21%	0.559%
8	2.269	178	183	196	rVV	21748552	33115233	16.07%	2.136%
9	2.397	196	204	216	rVB	1015363	2078720	1.01%	0.134%
10	2.842	255	277	280	rBV3	2804590	9272143	4.50%	0.598%
11	2.891	280	285	288	rVV	9686296	19540029	9.48%	1.260%
12	2.934	288	292	320	rVV	18092308	41999314	20.38%	2.709%
13	3.172	321	331	350	rVB	6199652	13836339	6.71%	0.892%
14	3.397	357	368	379	rBV3	1897347	4786564	2.32%	0.309%
15	3.519	379	388	401	rVB	3959580	8820989	4.28%	0.569%
16	3.666	401	412	418	rBV	1293130	3366136	1.63%	0.217%
17	3.751	418	426	431	rBV	1456138	3569868	1.73%	0.230%
18	3.812	431	436	446	rVB	2785721	6658762	3.23%	0.429%
19	3.922	446	454	460	rBV	1654498	4274708	2.07%	0.276%
20	4.196	489	499	510	rVB	2604155	6983435	3.39%	0.450%
21	4.403	522	533	545	rVV	10985452	31179128	15.13%	2.011%
22	4.525	545	553	566	rVB	804815	2364189	1.15%	0.152%
23	5.305	656	681	700	rBV	4743127	14796027	7.18%	0.954%
24	5.580	717	726	736	rBV2	3486958	11181161	5.43%	0.721%
25	5.927	771	783	796	rBV3	994468	3071980	1.49%	0.198%
26	6.269	821	839	840	rBV	845098	2316674	1.12%	0.149%
27	6.458	862	870	880	rVB	796213	2212071	1.07%	0.143%
28	6.945	940	950	957	rVB6	1399643	5520269	2.68%	0.356%
29	7.256	990	1001	1011	rBV	1251886	2945338	1.43%	0.190%
30	7.464	1022	1035	1046	rBV4	3208681	8272303	4.01%	0.534%
31	8.213	1145	1158	1163	rBV	2108538	3949309	1.92%	0.255%
32	8.573	1210	1217	1227	rVB	1428637	2486887	1.21%	0.160%
33	8.720	1235	1241	1246	rVB	1741196	2956732	1.43%	0.191%
34	8.805	1246	1255	1260	rBV3	76647354	147048726	71.35%	9.484%

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35	10.116	1465	1470	1474	rBV	1833216	2486320	1.21%	0.160%
36	10.262	1487	1494	1500	rVB3	81572658	130194834	63.17%	8.397%
37	10.378	1503	1513	1518	rBV2	114787493	206089347	100.00%	13.292%
38	10.707	1560	1567	1572	rVB	38615855	52547899	25.50%	3.389%
39	11.018	1613	1618	1626	rVB	5943863	7139143	3.46%	0.460%
40	11.140	1631	1638	1643	rBV	1670153	2175521	1.06%	0.140%
41	11.359	1668	1674	1681	rBV	14313631	17773488	8.62%	1.146%
42	11.439	1681	1687	1694	rVV2	50100852	93954149	45.59%	6.060%
43	11.512	1694	1699	1705	rVB	22831385	28412791	13.79%	1.833%
44	11.670	1719	1725	1735	rVB	23141291	27781073	13.48%	1.792%
45	11.817	1742	1749	1753	rBV2	63887261	84970495	41.23%	5.480%
46	12.079	1787	1792	1797	rVB2	2473024	3415227	1.66%	0.220%
47	12.146	1797	1803	1807	rBV	13783991	17198625	8.35%	1.109%
48	12.298	1823	1828	1836	rBV2	32889047	46636570	22.63%	3.008%
49	12.371	1836	1840	1849	rVB3	7628374	11113282	5.39%	0.717%
50	12.463	1849	1855	1860	rBV2	1550973	2278897	1.11%	0.147%
51	12.518	1860	1864	1870	rVB	1958279	2301480	1.12%	0.148%
52	12.585	1870	1875	1877	rBV	5036181	6162196	2.99%	0.397%
53	12.609	1877	1879	1884	rVB	4450426	4698512	2.28%	0.303%
54	12.670	1884	1889	1892	rBV	7113504	8959191	4.35%	0.578%
55	12.700	1892	1894	1898	rVB	2630694	2686861	1.30%	0.173%
56	12.755	1898	1903	1911	rBV	7667613	10187332	4.94%	0.657%
57	12.902	1922	1927	1932	rVB	1947902	2370239	1.15%	0.153%
58	12.975	1932	1939	1943	rBV	4362428	5357384	2.60%	0.346%
59	13.018	1943	1946	1951	rVB	6454390	7202251	3.49%	0.465%
60	13.225	1975	1980	1984	rVB	7539825	9014530	4.37%	0.581%
61	13.347	1996	2000	2005	rVB2	13037576	17071609	8.28%	1.101%
62	13.481	2017	2022	2029	rVB	3817483	4829336	2.34%	0.311%
63	13.834	2072	2080	2089	rBV	26623822	33660389	16.33%	2.171%
64	14.694	2217	2221	2231	rVB	7036626	8484809	4.12%	0.547%
65	14.840	2239	2245	2253	rVB	2524216	3312285	1.61%	0.214%

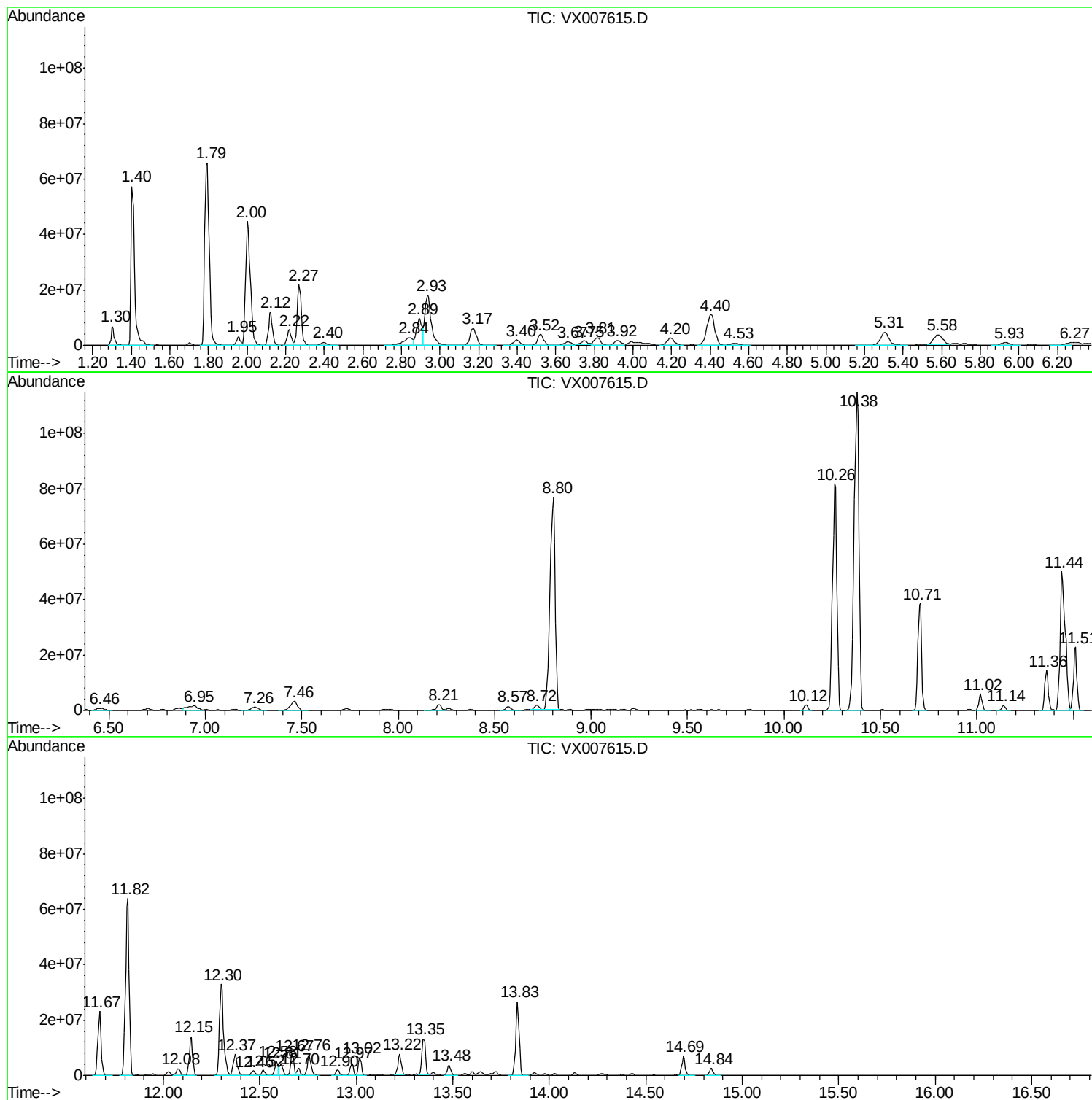
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Quant Title : SW846 8260

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



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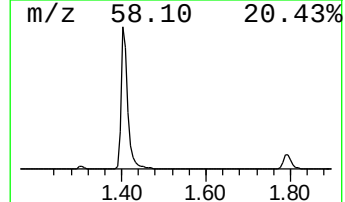
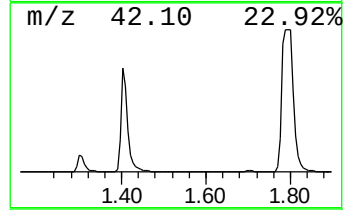
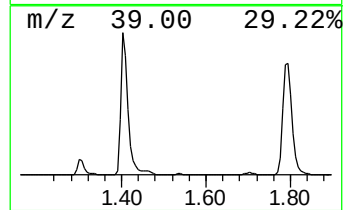
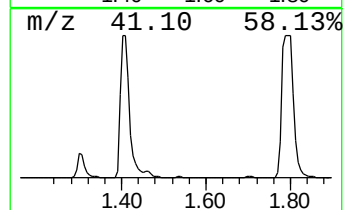
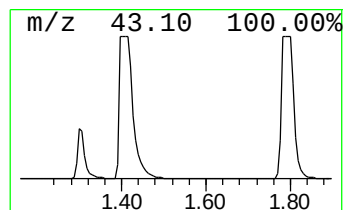
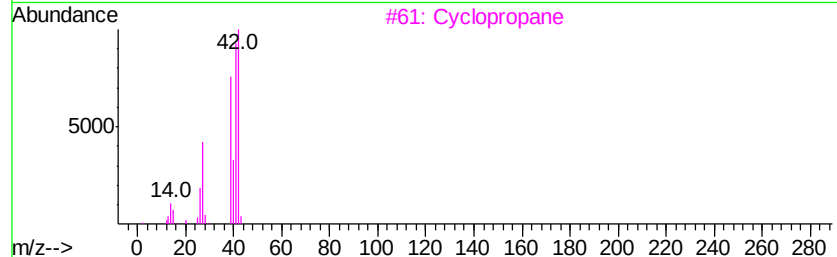
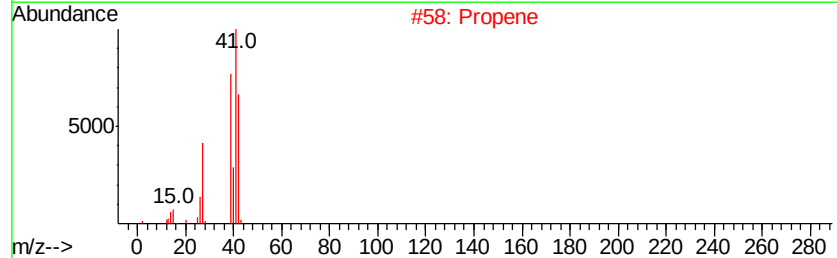
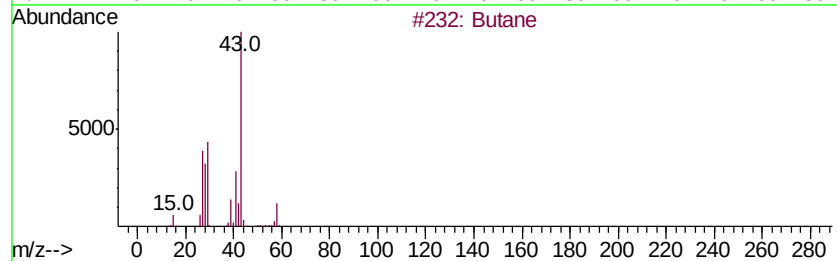
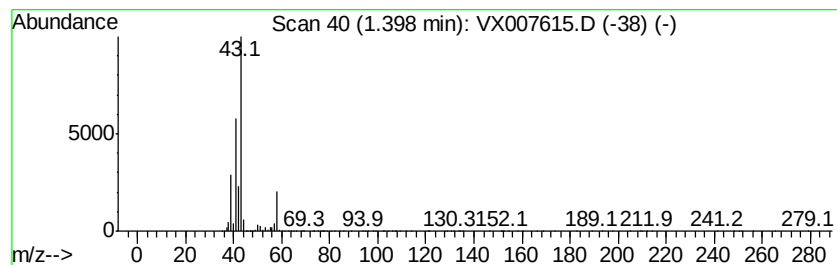
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X020119W.M
Quant Title : SW846 8260

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

Peak Number 1 Butane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.40	328.45 ug/l	73448200	Pentafluorobenzene	5.67

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane		58	C4H10	000106-97-8	78
2	Propene		42	C3H6	000115-07-1	9
3	Cyclopropane		42	C3H6	000075-19-4	9
4	Oxirane, trimethyl-		86	C5H10O	005076-19-7	9
5	Trimethylene oxide		58	C3H6O	000503-30-0	5



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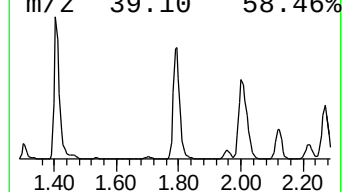
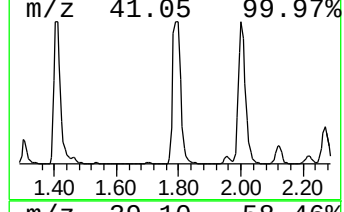
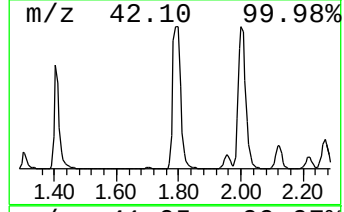
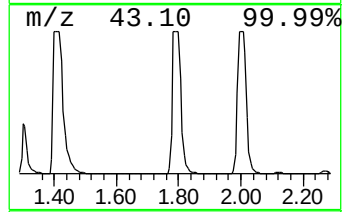
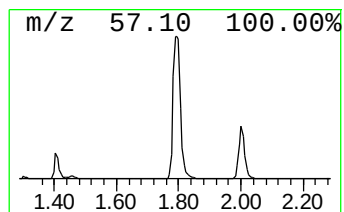
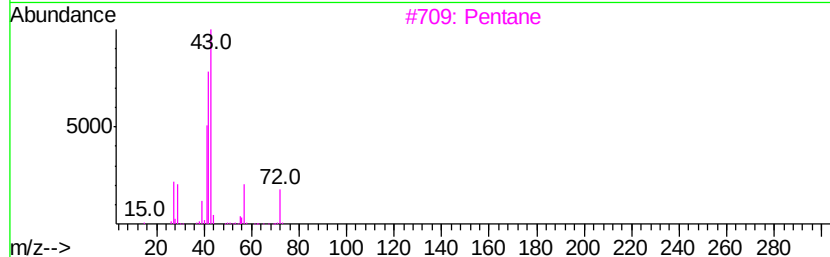
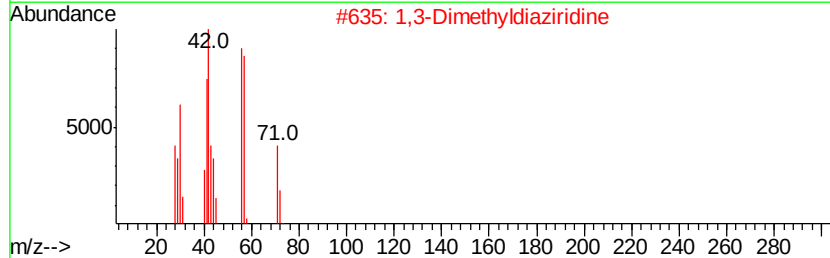
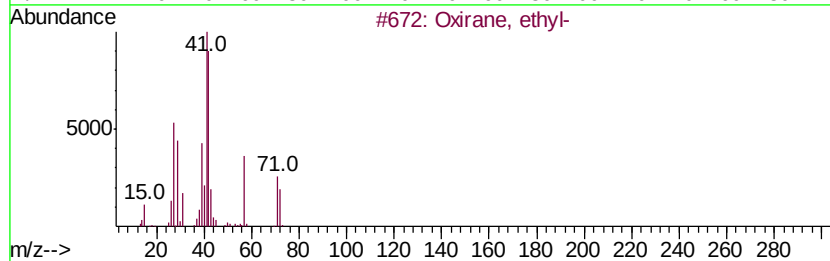
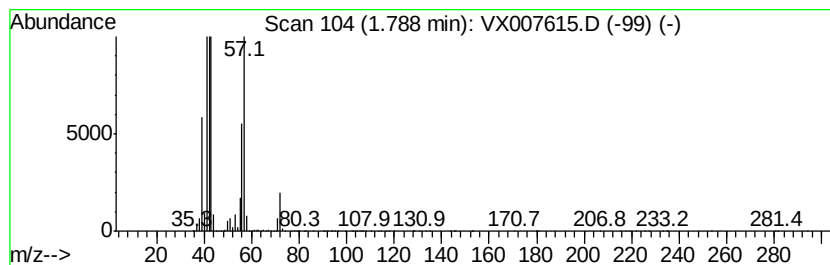
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Oxirane, ethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.79	458.10 ug/l	102442000	Pentafluorobenzene	5.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Oxirane, ethyl-	72	C4H8O	000106-88-7	50
2		1,3-Dimethyldiaziridine	72	C3H8N2	026177-36-6	38
3		Pentane	72	C5H12	000109-66-0	37
4		Diaziridine,3,3-dimethyl-	72	C3H8N2	004901-76-2	37
5		Isobutylene epoxide	72	C4H8O	000558-30-5	32



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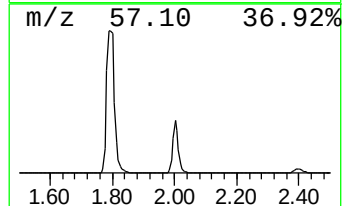
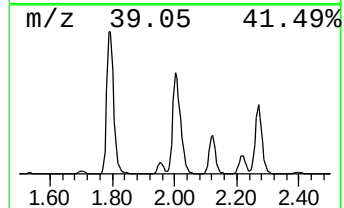
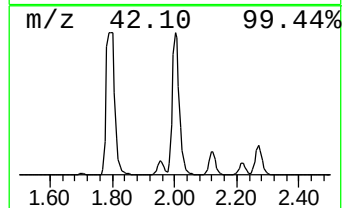
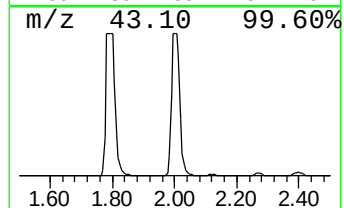
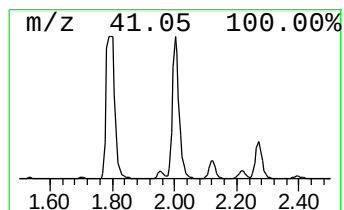
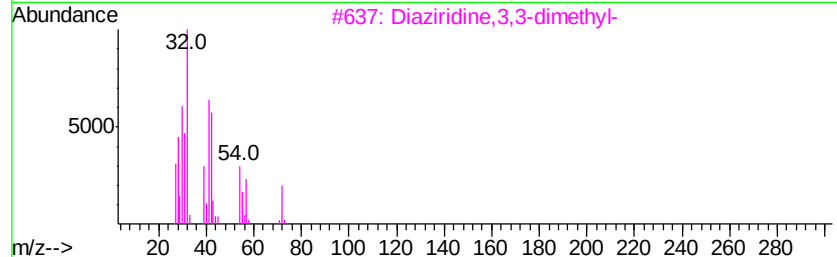
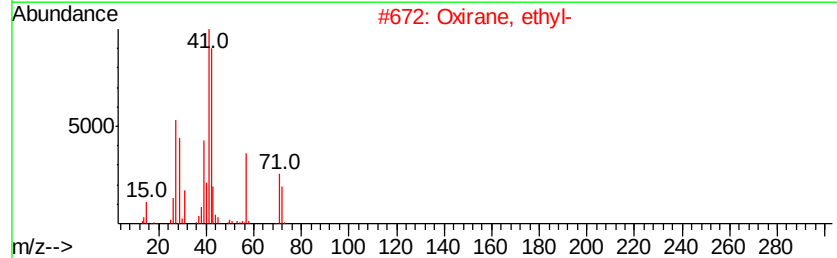
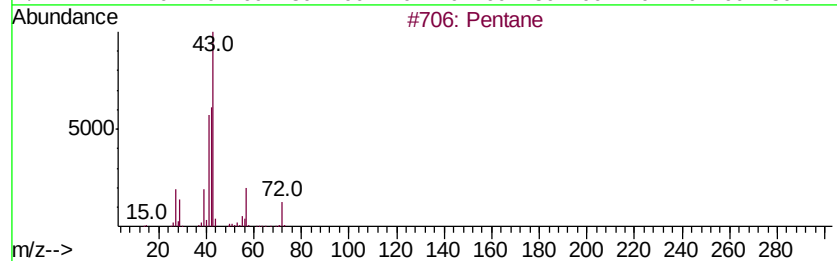
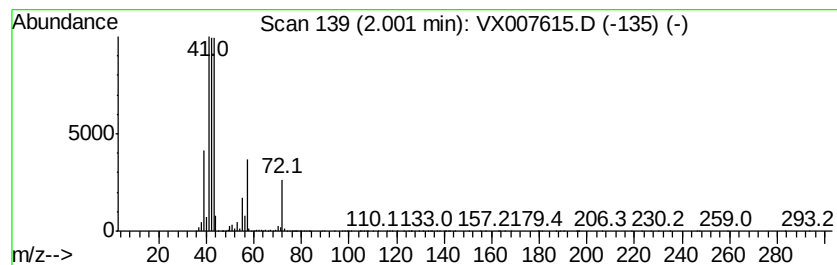
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 Pentane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.00	344.24 ug/l	76979600	Pentafluorobenzene	5.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane	72	C5H12	000109-66-0	83
2		Oxirane, ethyl-	72	C4H8O	000106-88-7	64
3		Diaziridine,3,3-dimethyl-	72	C3H8N2	004901-76-2	56
4		Isobutylene epoxide	72	C4H8O	000558-30-5	49
5		Butane, 2-methyl-	72	C5H12	000078-78-4	40



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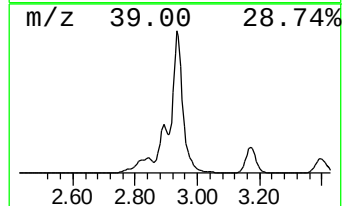
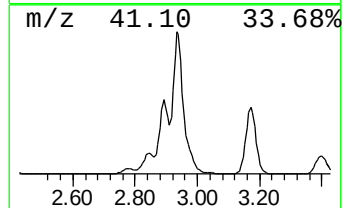
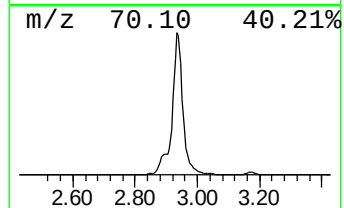
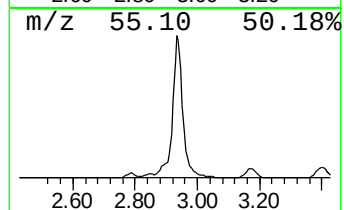
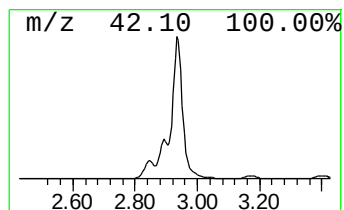
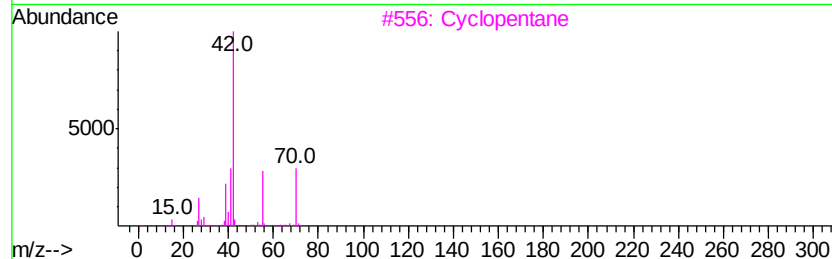
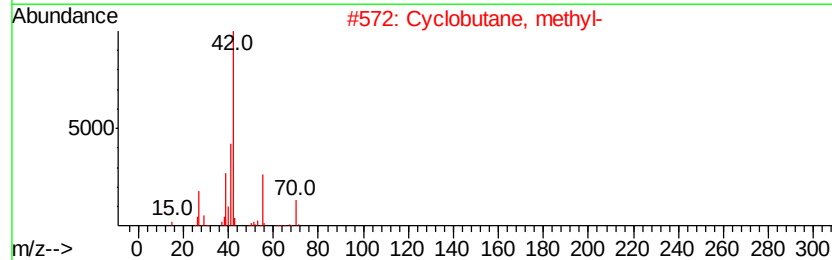
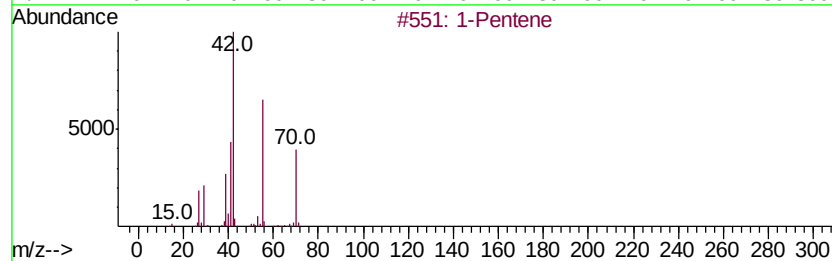
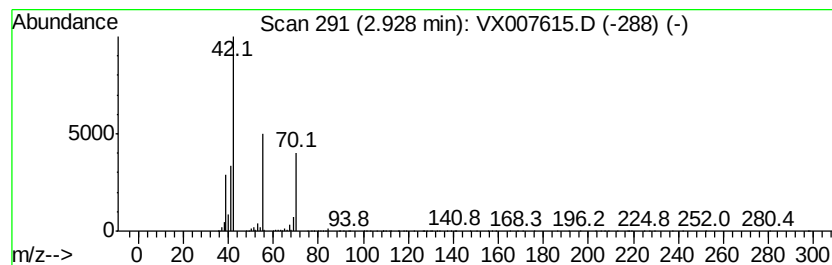
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Quant Title : SW846 8260

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

Peak Number 4 1-Pentene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.93	187.81 ug/l	41999300	Pentafluorobenzene	5.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Pentene	70	C5H10	000109-67-1	86
2		Cyclobutane, methyl-	70	C5H10	000598-61-8	86
3		Cyclopentane	70	C5H10	000287-92-3	78
4		2-Pentene, (Z)-	70	C5H10	000627-20-3	47
5		2-Pentene	70	C5H10	000109-68-2	47



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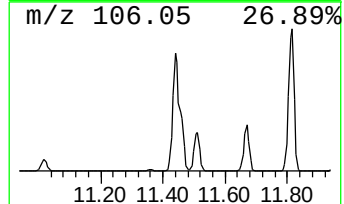
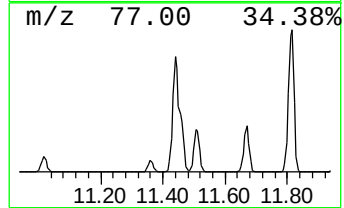
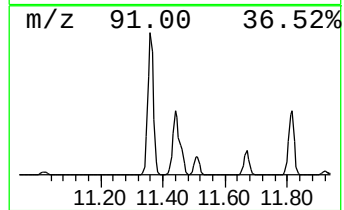
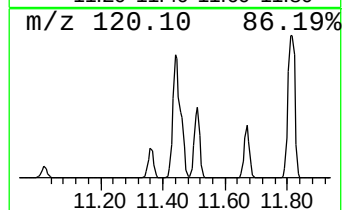
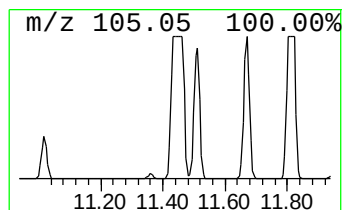
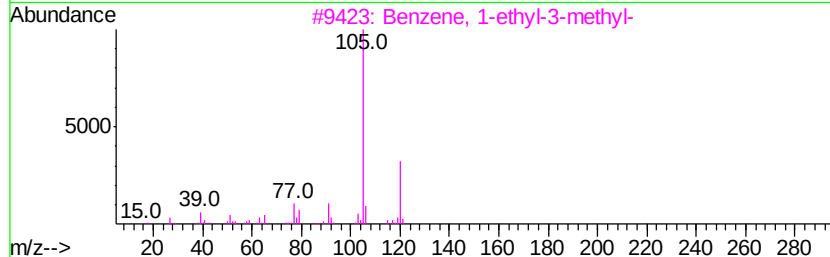
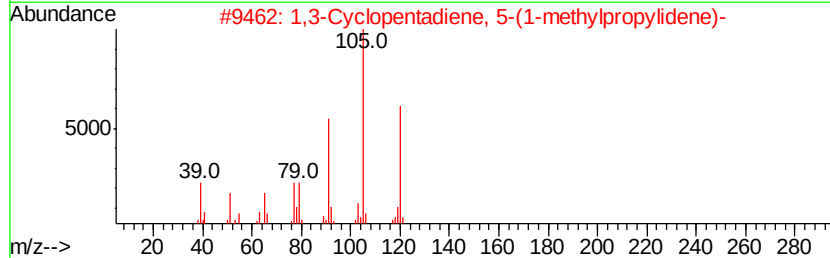
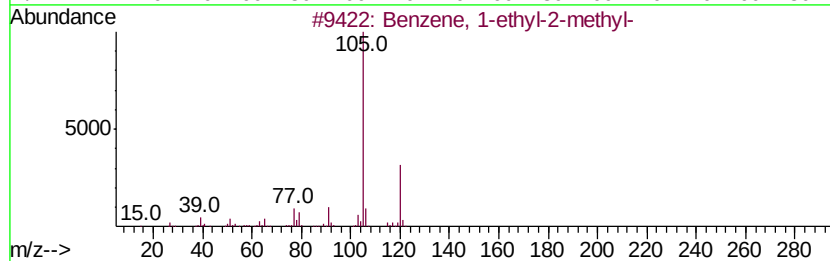
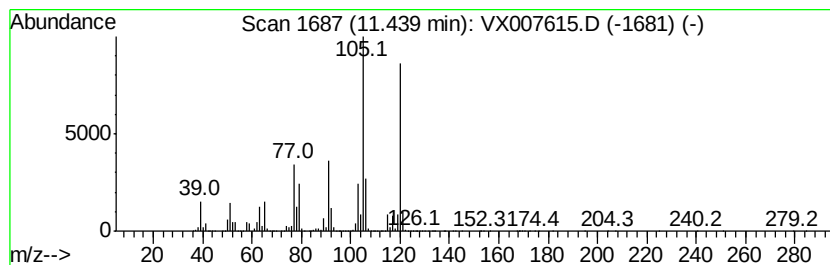
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1-ethyl-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.44	1375.52 ug/l	93954100	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	87
2		1,3-Cyclopentadiene, 5-(1-methyl...	120	C9H12	003141-02-4	87
3		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	87
4		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	87
5		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	83



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX021919\
Data File : VX007615.D
Acq On : 19 Feb 2019 16:04
Operator : JC/SP
Sample : K1528-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 16 Sample Multiplier: 1

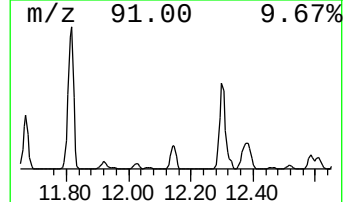
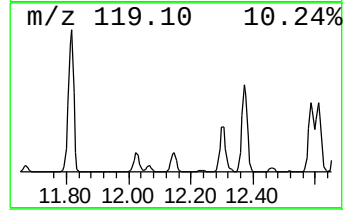
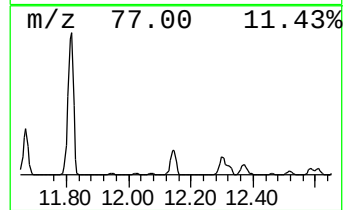
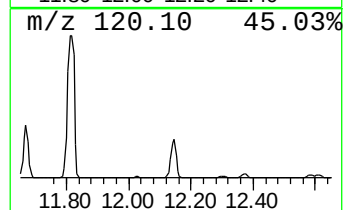
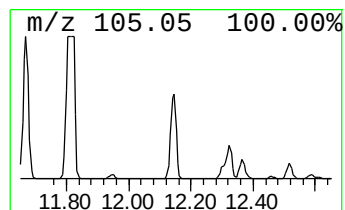
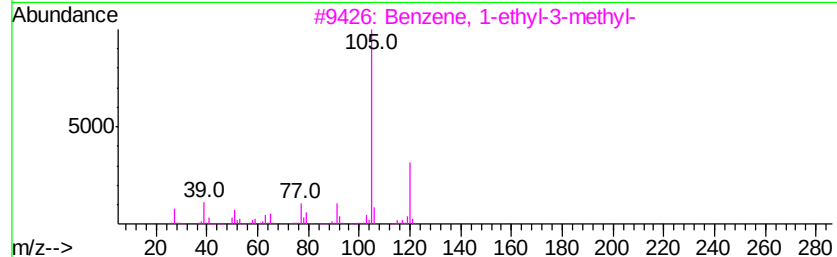
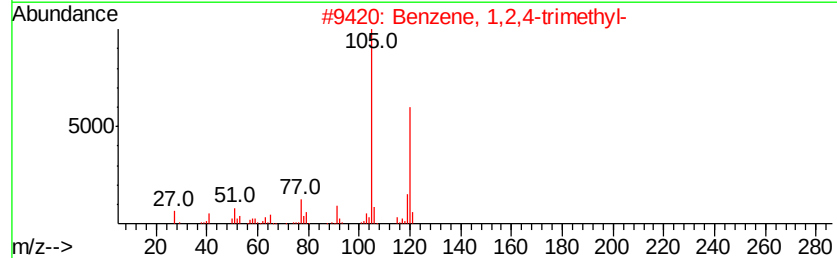
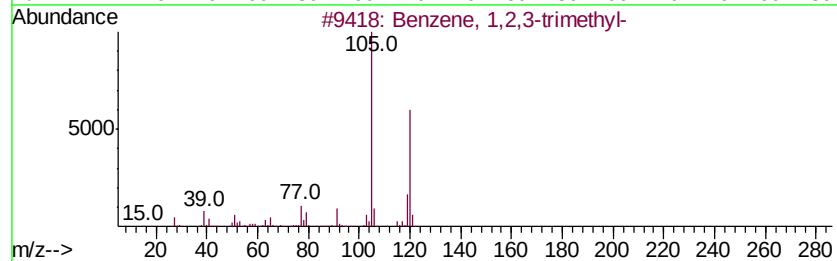
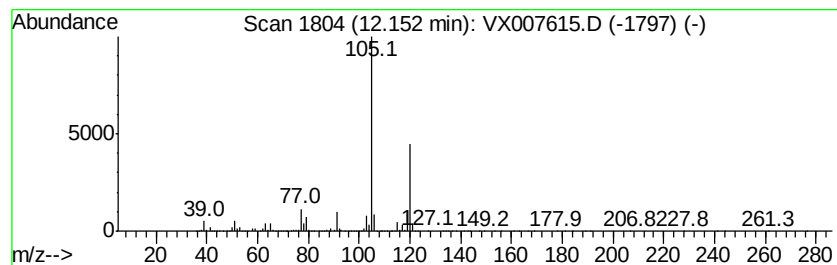
Instrument :
MSVOA_X
ClientSampleId :
MW-1

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X020119W.M
Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1,2,3-trimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.15	251.79 ug/l	17198600	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1,2,3-trimethyl-	120 C9H12	000526-73-8 95
2		Benzene, 1,2,4-trimethyl-	120 C9H12	000095-63-6 94
3		Benzene, 1-ethyl-3-methyl-	120 C9H12	000620-14-4 91
4		Mesitylene	120 C9H12	000108-67-8 91
5		Benzene, 1-ethyl-2-methyl-	120 C9H12	000611-14-3 90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX021919\
Data File : VX007615.D
Acq On : 19 Feb 2019 16:04
Operator : JC/SP
Sample : K1528-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-1

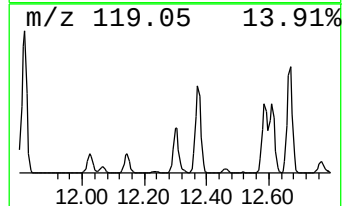
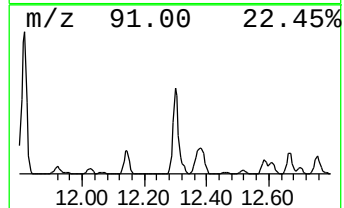
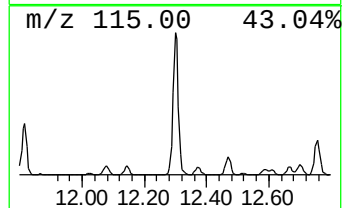
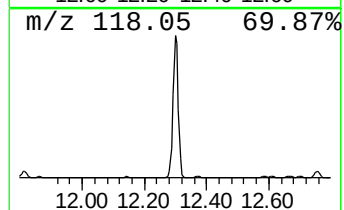
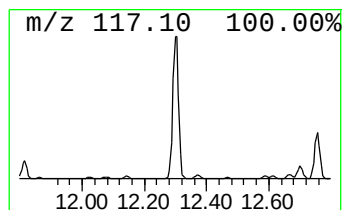
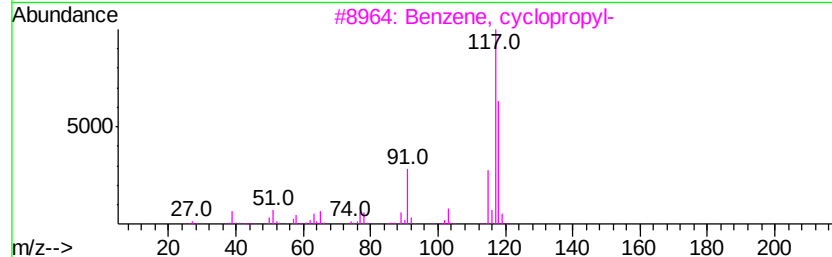
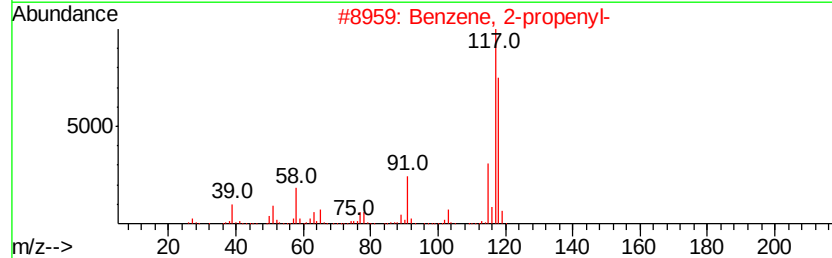
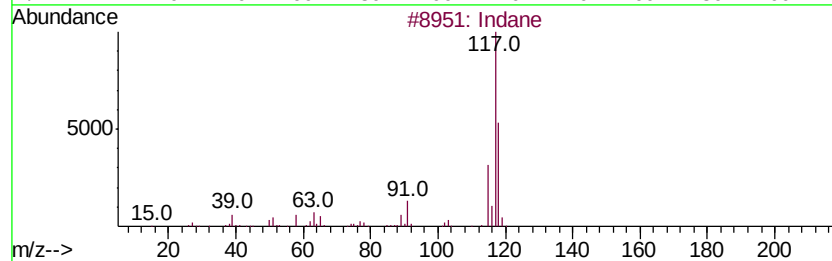
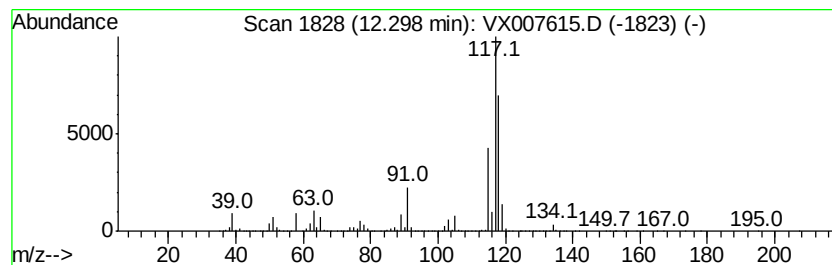
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X020119W.M
Quant Title : SW846 8260

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

Peak Number 8 Indane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	682.77 ug/l	46636600	1,4-Dichlorobenzene-d4	12.08

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane		118	C9H10	000496-11-7	76
2	Benzene, 2-propenyl-		118	C9H10	000300-57-2	70
3	Benzene, cyclopropyl-		118	C9H10	000873-49-4	70
4	Benzene, 1-ethenyl-2-methyl-		118	C9H10	000611-15-4	62
5	Benzene, 1-propenyl-		118	C9H10	000637-50-3	62



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX021919\
Data File : VX007615.D
Acq On : 19 Feb 2019 16:04
Operator : JC/SP
Sample : K1528-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-1

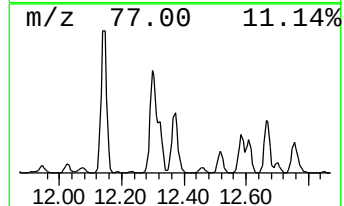
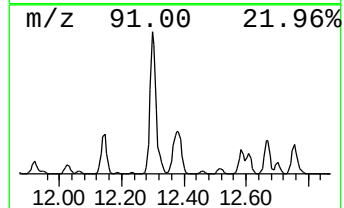
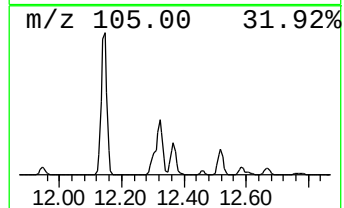
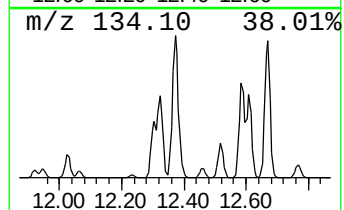
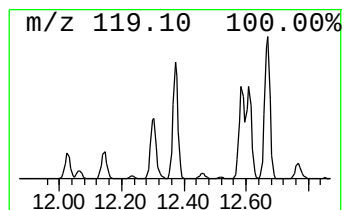
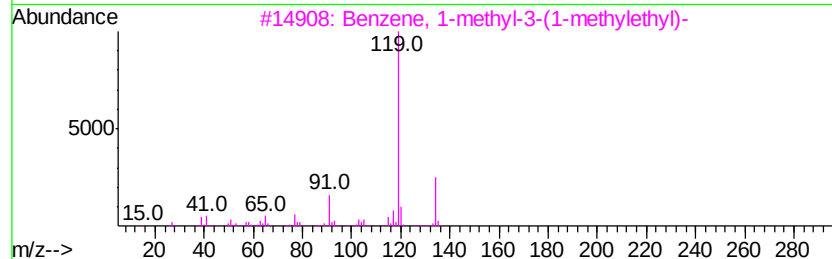
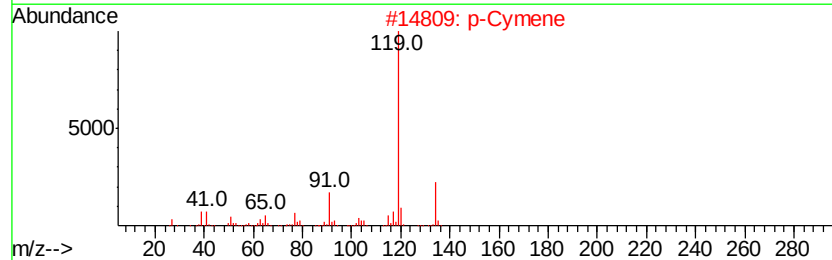
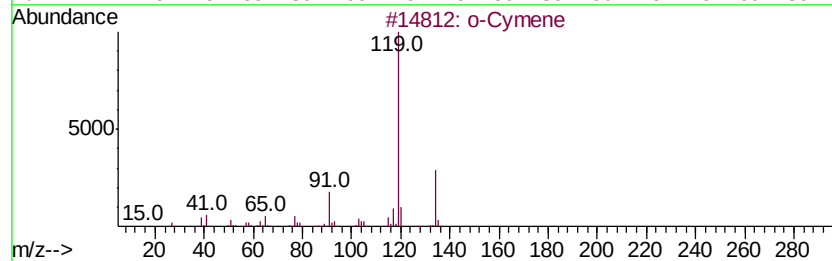
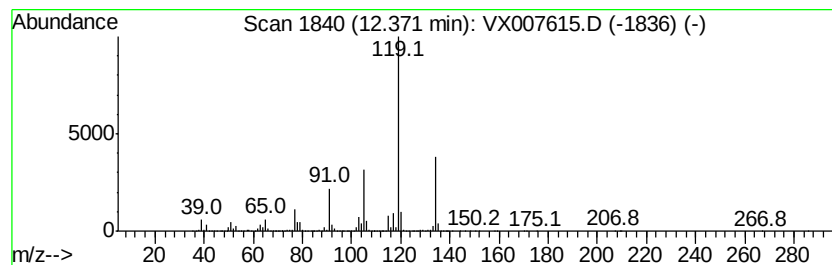
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Quant Title : SW846 8260

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

Peak Number 9 o-Cymene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.37	162.70 ug/l	11113300	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	o-Cymene	134	C10H14	000527-84-4	95
2		p-Cymene	134	C10H14	000099-87-6	94
3		Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	94
4		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	93
5		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX021919\
Data File : VX007615.D
Acq On : 19 Feb 2019 16:04
Operator : JC/SP
Sample : K1528-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-1

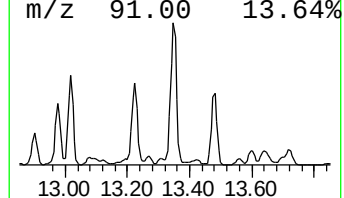
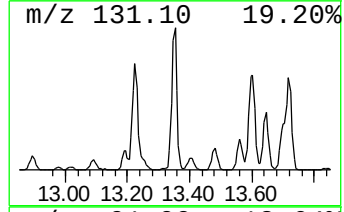
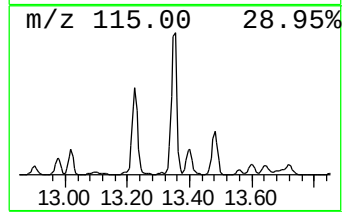
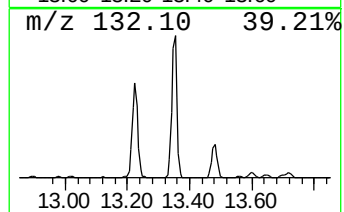
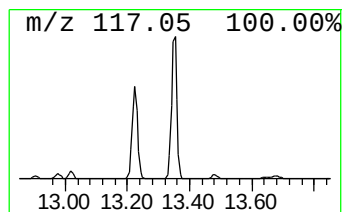
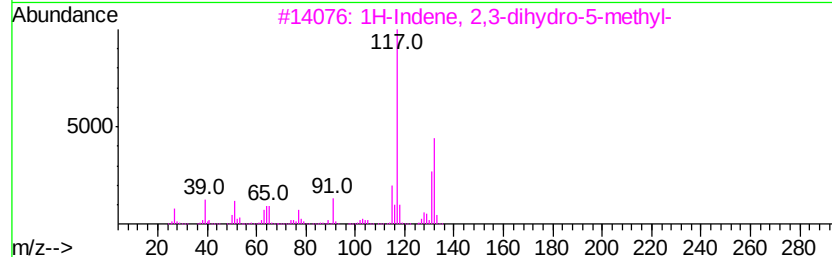
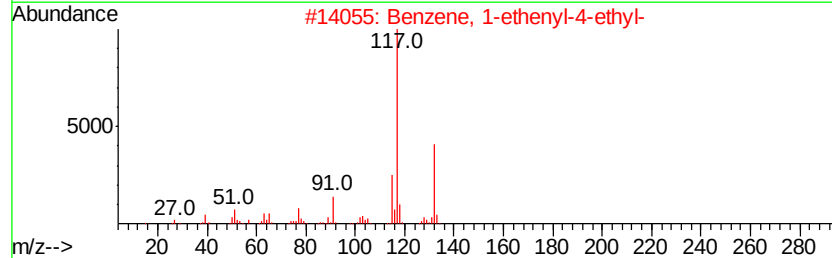
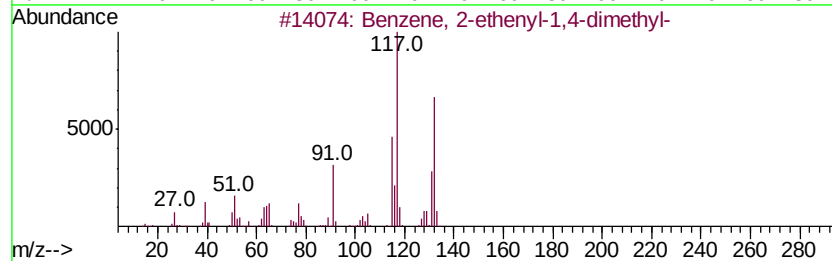
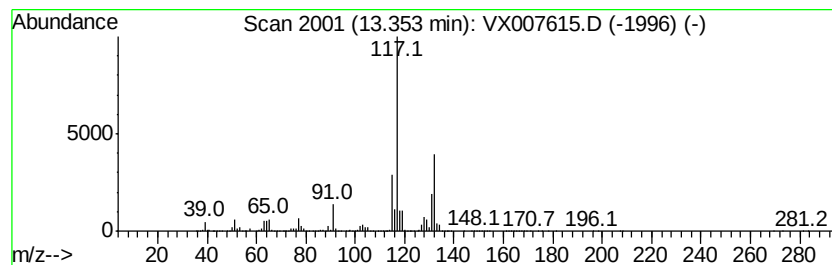
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X020119W.M
Quant Title : SW846 8260

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

Peak Number 10 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	249.93 ug/l	17071600	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
2		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
3		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	87
4		1-Phenyl-1-butene	132	C10H12	000824-90-8	87
5		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX021919\
Data File : VX007615.D
Acq On : 19 Feb 2019 16:04
Operator : JC/SP
Sample : K1528-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-1

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X020119W.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
Butane	1.40	328.4	ug/l	73448200	1	5.67	11181200	50.0
Oxirane, ethyl-	1.79	458.1	ug/l	102442000	1	5.67	11181200	50.0
Pentane	2.00	344.2	ug/l	76979600	1	5.67	11181200	50.0
1-Pentene	2.93	187.8	ug/l	41999300	1	5.67	11181200	50.0
Benzene, 1-ethyl-...	11.44	1375.5	ug/l	93954100	4	12.08	3415230	50.0
Benzene, 1,2,3-tr...	12.15	251.8	ug/l	17198600	4	12.08	3415230	50.0
Indane	12.30	682.8	ug/l	46636600	4	12.08	3415230	50.0
o-Cymene	12.37	162.7	ug/l	11113300	4	12.08	3415230	50.0
Benzene, 2-etheny...	13.35	249.9	ug/l	17071600	4	12.08	3415230	50.0