

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX021119\
 Data File : VX007468.D
 Acq On : 11 Feb 2019 15:19
 Operator : JC/SP
 Sample : K1433-10
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 KY001MW031-190207

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	2.848	267	278	280	rBV2	298074	652855	1.02%	0.230%
2	2.891	280	285	289	rVV	650778	1432440	2.23%	0.505%
3	2.934	289	292	304	rVB2	469706	937683	1.46%	0.330%
4	3.202	322	336	353	rVB	5037598	13226213	20.62%	4.661%
5	4.397	523	532	545	rVB	1082319	3120012	4.86%	1.099%
6	5.506	701	714	719	rBV	337696	1010648	1.58%	0.356%
7	5.579	719	726	734	rVV3	644018	2178112	3.40%	0.768%
8	5.665	734	740	745	rVV	566110	1638661	2.55%	0.577%
9	5.726	746	750	768	rVB2	433204	1273090	1.98%	0.449%
10	5.927	773	783	797	rBV5	256940	806335	1.26%	0.284%
11	6.067	797	806	810	rBV	305237	753358	1.17%	0.265%
12	6.146	810	819	830	rVB	5622274	14329902	22.34%	5.050%
13	6.348	847	852	861	rVB2	685403	1966490	3.07%	0.693%
14	6.445	861	868	882	rVB3	417052	1273754	1.99%	0.449%
15	6.860	926	936	945	rBV2	825841	2419790	3.77%	0.853%
16	7.256	992	1001	1011	rBV2	297375	660272	1.03%	0.233%
17	7.451	1023	1033	1047	rVB4	543096	1607534	2.51%	0.566%
18	7.981	1113	1120	1126	rVB3	352731	783720	1.22%	0.276%
19	8.055	1126	1132	1143	rVB	311821	689886	1.08%	0.243%
20	8.177	1144	1152	1155	rBV	417686	854602	1.33%	0.301%
21	8.579	1211	1218	1225	rVB2	512974	974992	1.52%	0.344%
22	8.713	1235	1240	1246	rVB	1704586	2640309	4.12%	0.930%
23	8.780	1246	1251	1257	rVB2	816594	1225706	1.91%	0.432%
24	8.908	1269	1272	1280	rVB	433742	655268	1.02%	0.231%
25	9.518	1363	1372	1376	rBV3	369910	681733	1.06%	0.240%
26	9.853	1424	1427	1434	rVB	484082	695810	1.08%	0.245%
27	10.085	1459	1465	1467	rVV2	896509	1598498	2.49%	0.563%
28	10.109	1467	1469	1474	rVV	1673219	2241815	3.50%	0.790%
29	10.164	1474	1478	1481	rVV2	619171	1003680	1.56%	0.354%
30	10.207	1481	1485	1487	rVV	606347	847421	1.32%	0.299%
31	10.255	1487	1493	1498	rVV	46559449	64139761	100.00%	22.601%
32	10.298	1498	1500	1506	rVV3	781312	1524469	2.38%	0.537%
33	10.359	1506	1510	1515	rVB	3481117	4526171	7.06%	1.595%
34	10.432	1515	1522	1525	rBV	605629	865037	1.35%	0.305%

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35	10.701	1561	1566	1575	rVB	4360881	5670999	8.84%	1.998%
36	10.847	1586	1590	1596	rVB4	360984	658179	1.03%	0.232%
37	11.018	1611	1618	1623	rVB	7584854	9754598	15.21%	3.437%
38	11.139	1633	1638	1643	rVV3	1823665	2599652	4.05%	0.916%
39	11.219	1646	1651	1659	rVB2	542719	1101171	1.72%	0.388%
40	11.359	1670	1674	1684	rVB	8135416	9726237	15.16%	3.427%
41	11.450	1684	1689	1693	rVV	1306784	1698956	2.65%	0.599%
42	11.505	1693	1698	1704	rVB	1135121	1682672	2.62%	0.593%
43	11.670	1720	1725	1731	rBV	12588427	15445168	24.08%	5.443%
44	11.804	1743	1747	1752	rVB	2276456	2777104	4.33%	0.979%
45	12.078	1786	1792	1797	rVB	2078932	2685124	4.19%	0.946%
46	12.139	1797	1802	1806	rBV	768375	1050749	1.64%	0.370%
47	12.231	1813	1817	1822	rVB	1825050	2186973	3.41%	0.771%
48	12.298	1822	1828	1833	rBV	26314696	31924604	49.77%	11.250%
49	12.365	1836	1839	1846	rVB2	864418	1454821	2.27%	0.513%
50	12.468	1851	1856	1860	rVV	2082754	2575148	4.01%	0.907%
51	12.517	1860	1864	1870	rVB2	685286	1005284	1.57%	0.354%
52	12.584	1870	1875	1882	rBV	553125	736364	1.15%	0.259%
53	12.670	1884	1889	1892	rVV	2067056	2612333	4.07%	0.921%
54	12.700	1892	1894	1898	rVV	1102761	1151246	1.79%	0.406%
55	12.755	1898	1903	1911	rVB2	3631521	4891847	7.63%	1.724%
56	12.974	1932	1939	1943	rBV	2834684	3464795	5.40%	1.221%
57	13.224	1976	1980	1990	rVB	2746399	3294914	5.14%	1.161%
58	13.346	1996	2000	2005	rVB2	5350807	6918729	10.79%	2.438%
59	13.401	2005	2009	2017	rVV2	1272704	1658322	2.59%	0.584%
60	13.480	2017	2022	2028	rVB	1534883	1889873	2.95%	0.666%
61	13.602	2038	2042	2045	rBV	520285	688476	1.07%	0.243%
62	13.645	2045	2049	2054	rVB4	521695	979629	1.53%	0.345%
63	13.718	2059	2061	2066	rVB2	563044	698519	1.09%	0.246%
64	13.834	2071	2080	2088	rVB	16149781	19832103	30.92%	6.988%
65	13.926	2089	2095	2100	rBV	1280659	1649517	2.57%	0.581%
66	14.694	2218	2221	2232	rVB	1063994	1433290	2.23%	0.505%
67	14.840	2240	2245	2251	rBV2	1975743	2653216	4.14%	0.935%

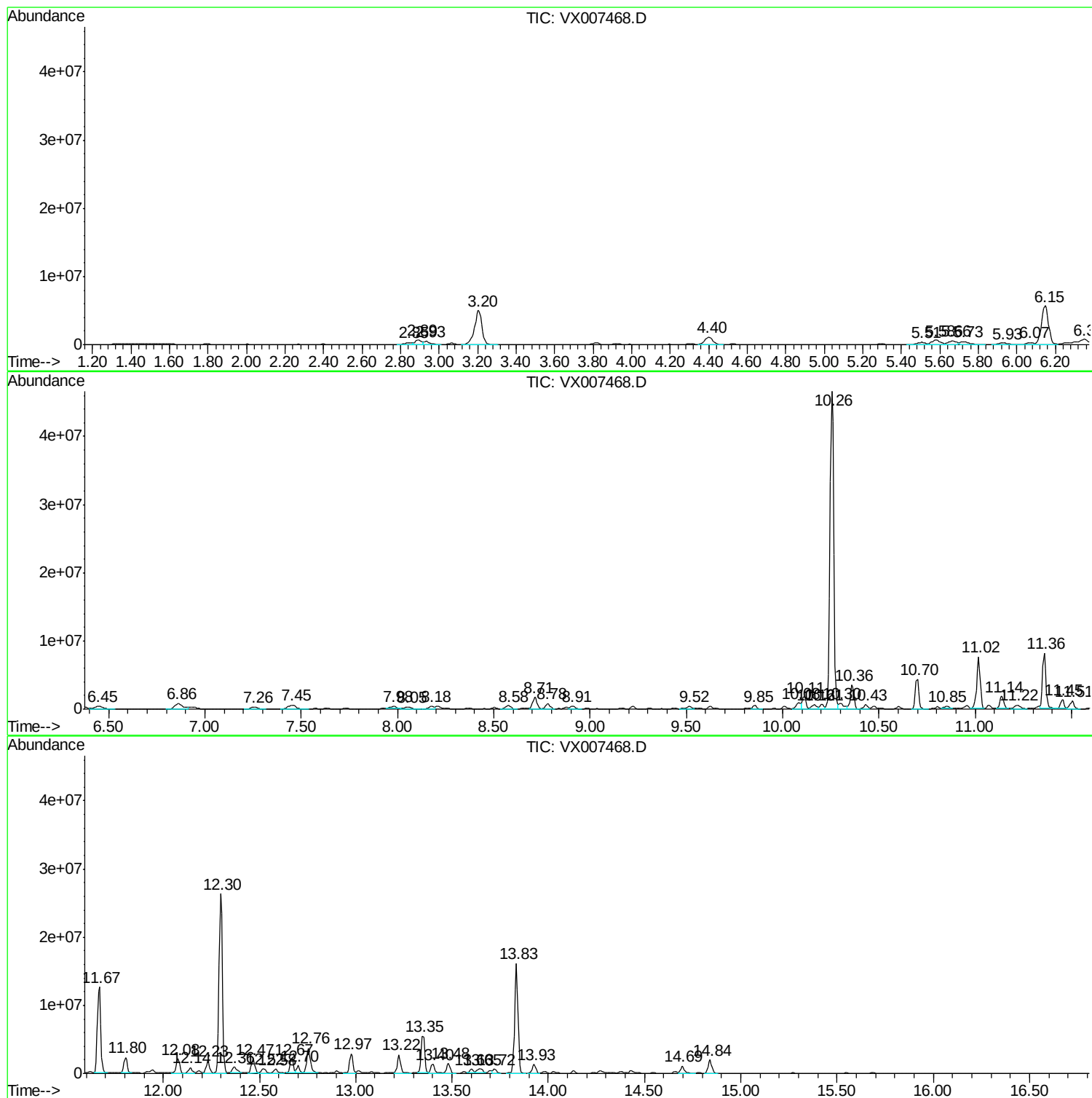
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Quant Title : SW846 8260

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



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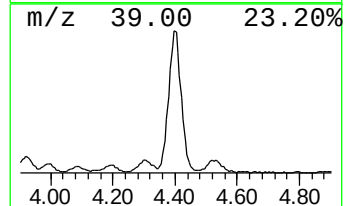
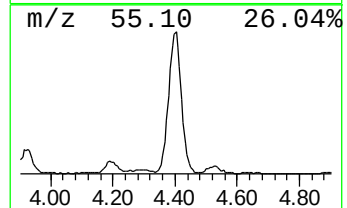
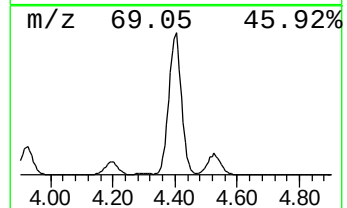
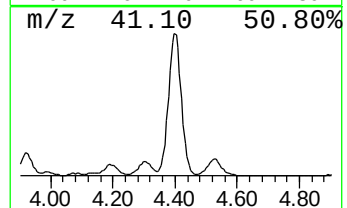
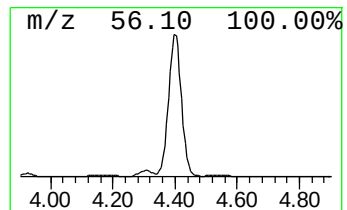
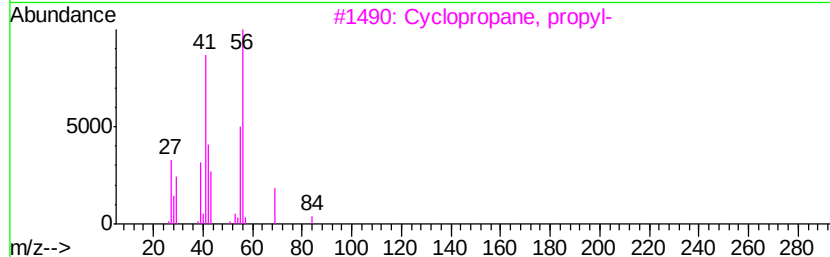
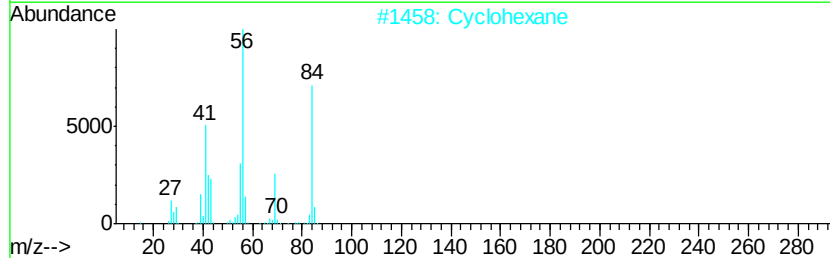
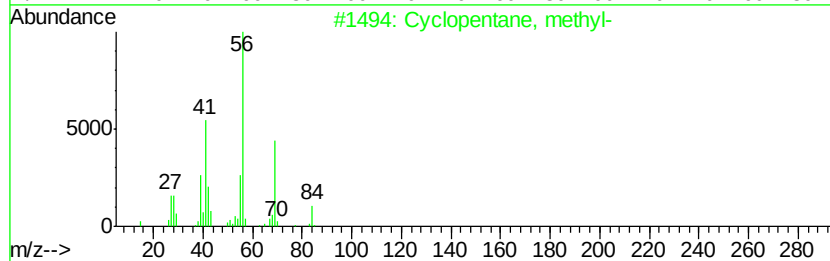
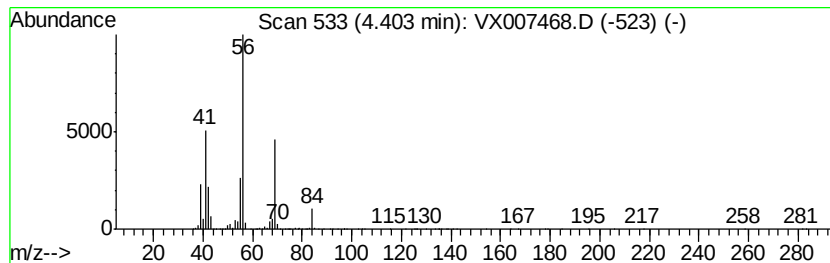
Instrument :
MSVOA_X
ClientSampleId :
KY001MW031-190207

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 Cyclopentane, methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
4.40	95.20 ug/l	3120010	Pentafluorobenzene	5.66	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	94
2	Cyclohexane	84	C6H12	000110-82-7	83
3	Cyclopropane, propyl-	84	C6H12	002415-72-7	78
4	Cyclobutane, ethyl-	84	C6H12	004806-61-5	72
5	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	64



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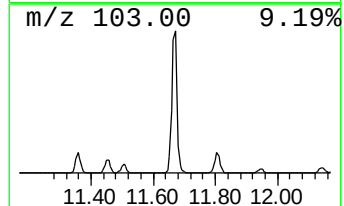
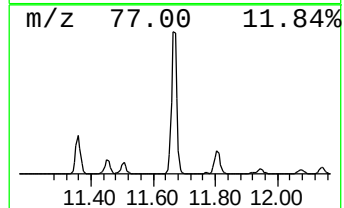
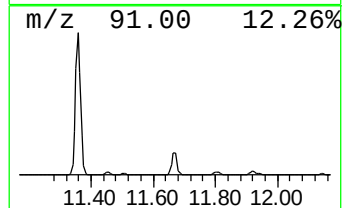
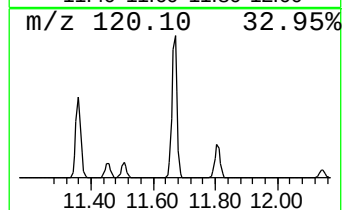
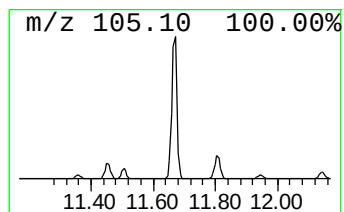
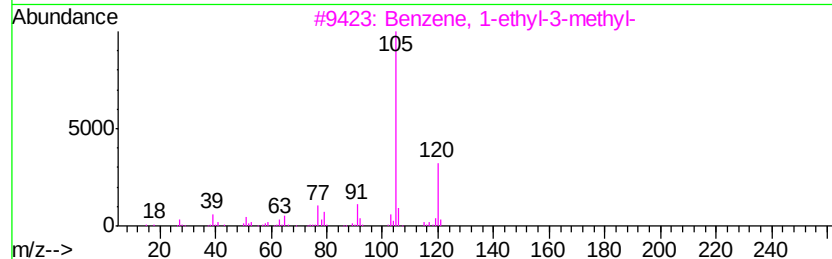
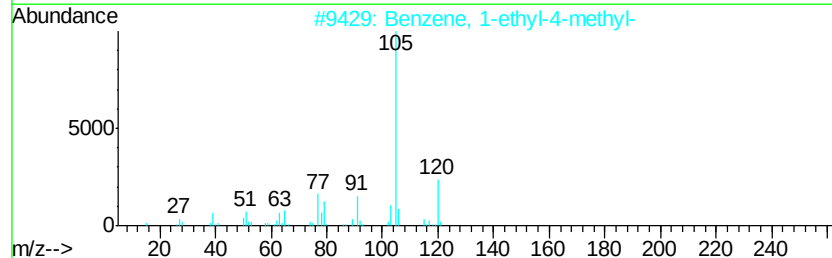
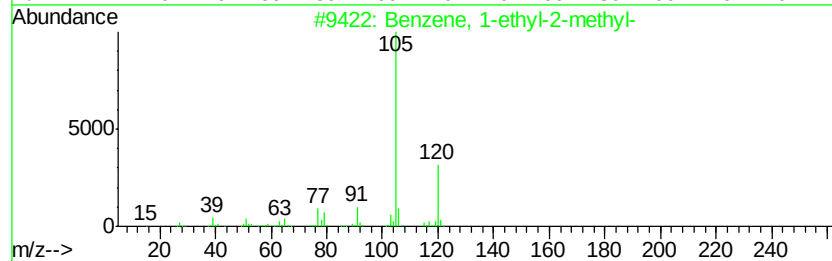
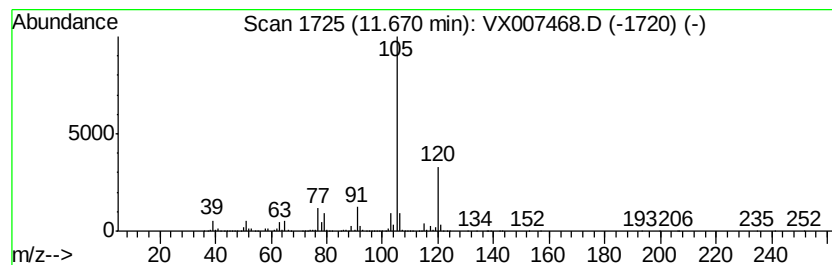
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.67	287.61 ug/l	15445200	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	94
2		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
3		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5		Mesitylene	120	C9H12	000108-67-8	91



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ALS Vial : 13 Sample Multiplier: 1

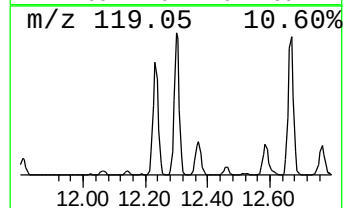
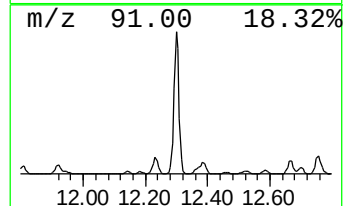
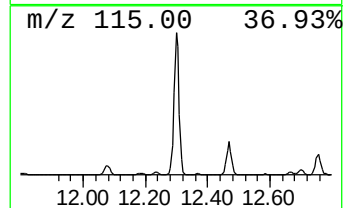
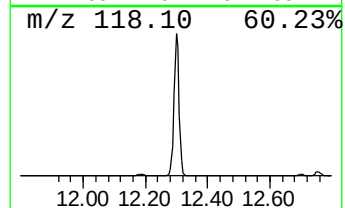
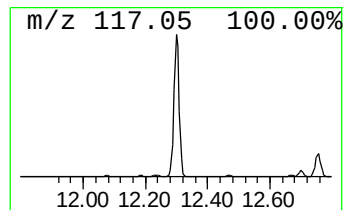
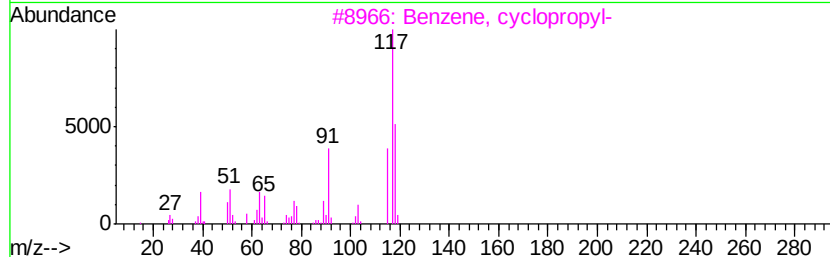
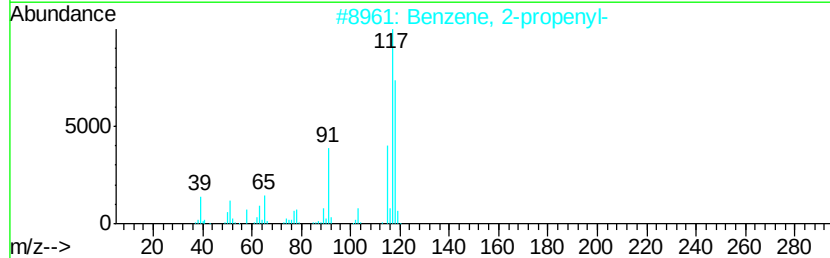
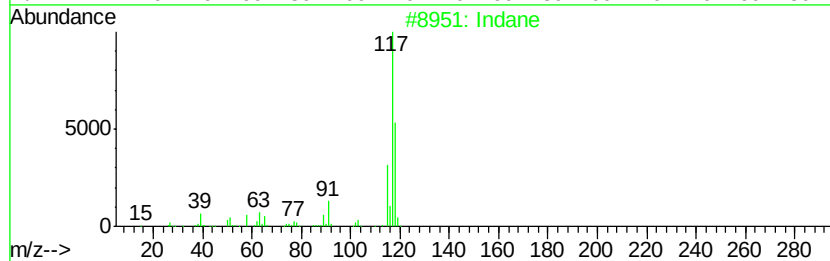
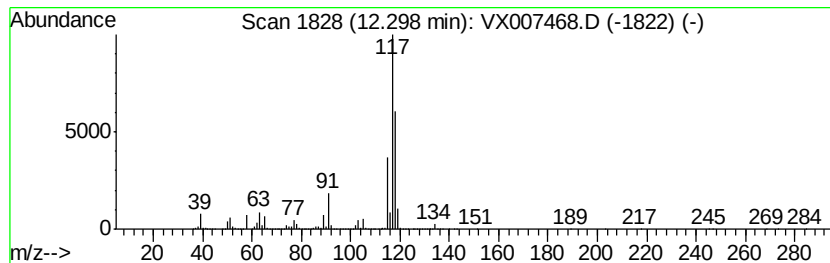
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	594.47 ug/l	31924600	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Indane	118 C9H10	000496-11-7	90
2	Benzene, 2-propenyl-	118 C9H10	000300-57-2	87
3	Benzene, cyclopropyl-	118 C9H10	000873-49-4	81
4	Benzene, 1-ethenyl-2-methyl-	118 C9H10	000611-15-4	74
5	Benzeneethanol, .beta.-ethenyl-....	162 C11H14O	040596-14-3	50



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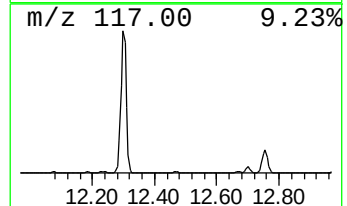
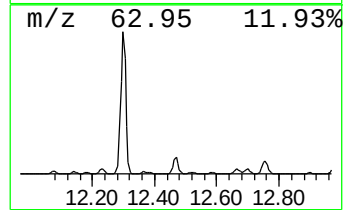
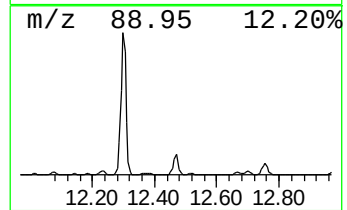
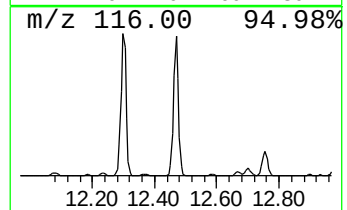
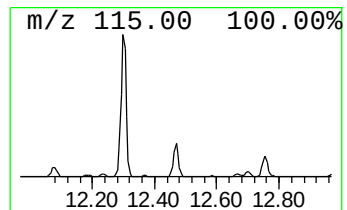
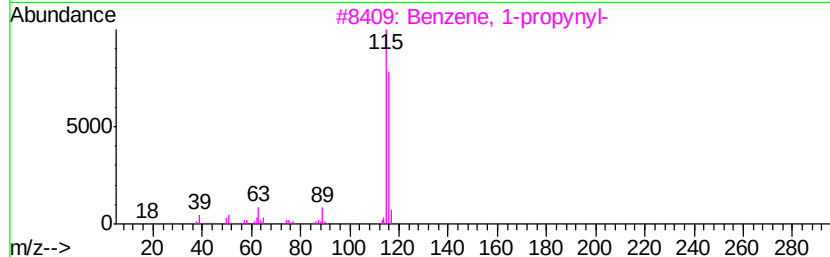
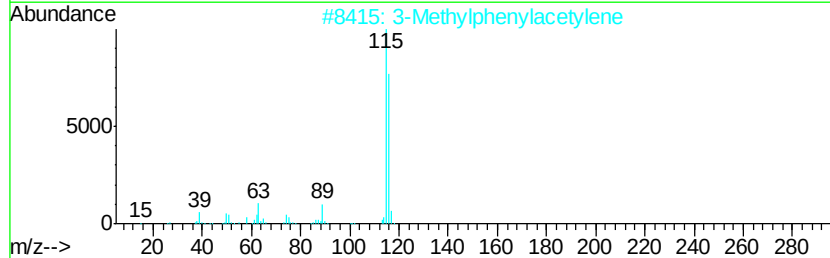
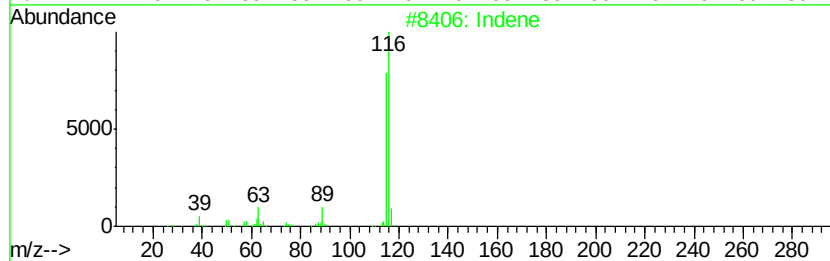
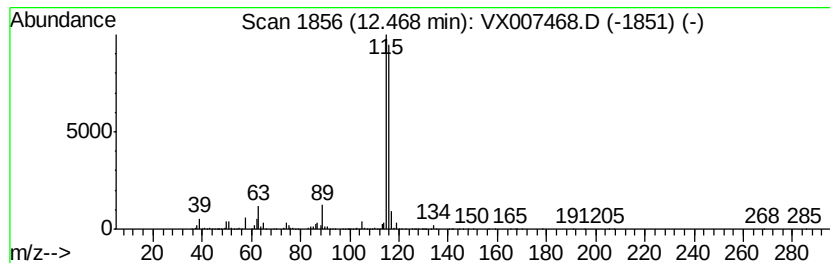
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Peak Number 4 Indene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.47	47.95 ug/l	2575150	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indene	116	C9H8	000095-13-6	97
2		3-Methylphenylacetylene	116	C9H8	000766-82-5	91
3		Benzene, 1-propynyl-	116	C9H8	000673-32-5	91
4		Benzene, 1-ethynyl-4-methyl-	116	C9H8	000766-97-2	90
5		Benzene, 1,2-propadienyl-	116	C9H8	002327-99-3	90



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ClientSampleId :
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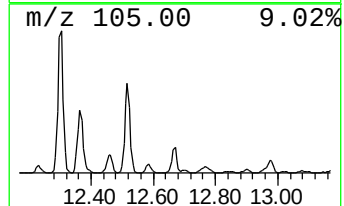
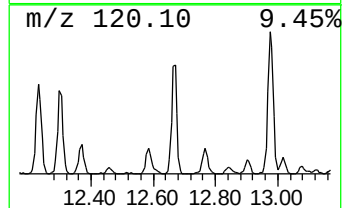
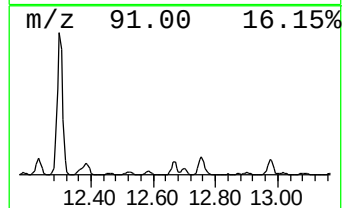
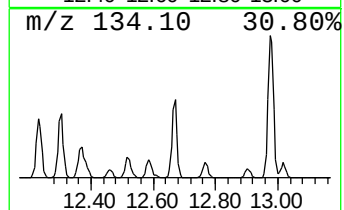
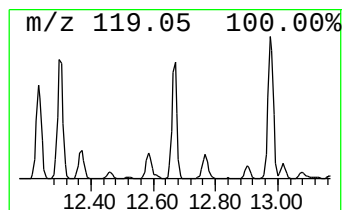
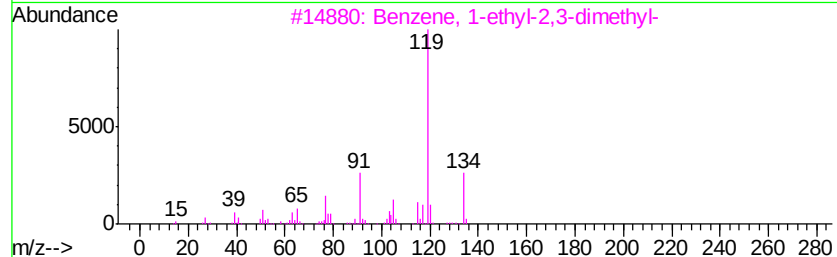
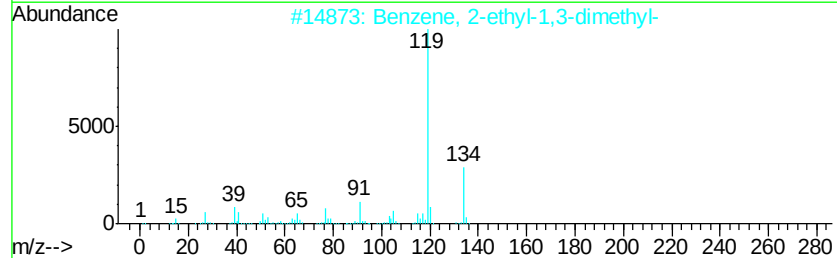
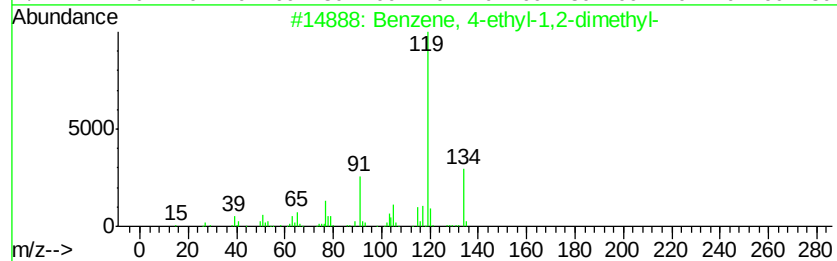
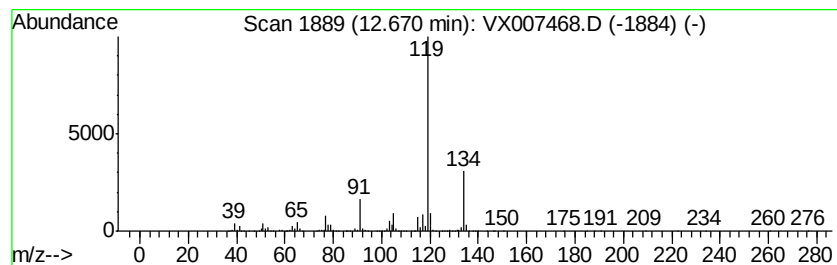
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.67	48.64 ug/l	2612330	1,4-Dichlorobenzene-d4	12.08

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		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
4		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
5		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX021119\
Data File : VX007468.D
Acq On : 11 Feb 2019 15:19
Operator : JC/SP
Sample : K1433-10
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
KY001MW031-190207

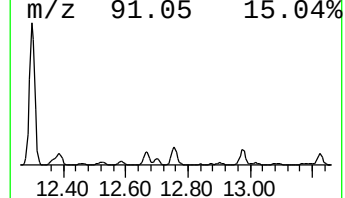
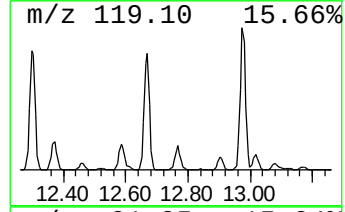
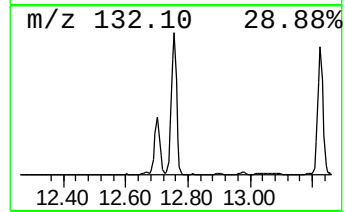
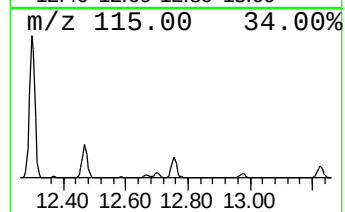
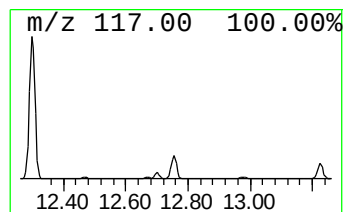
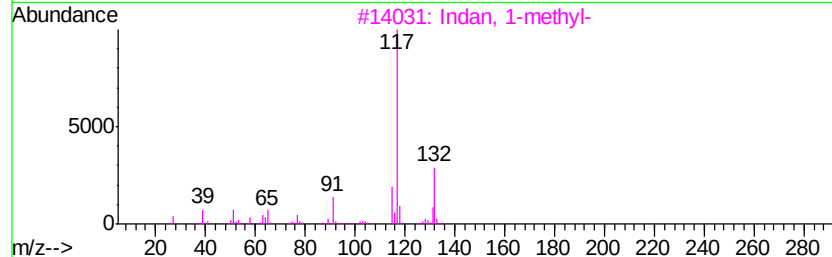
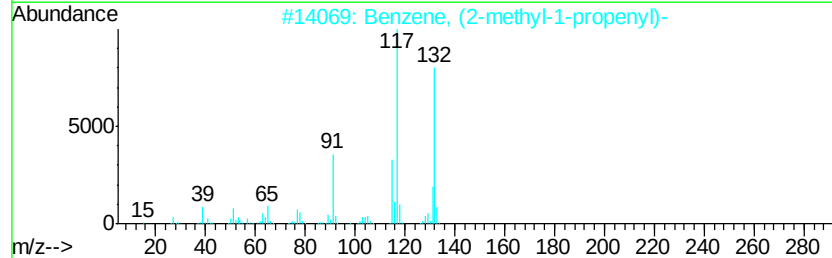
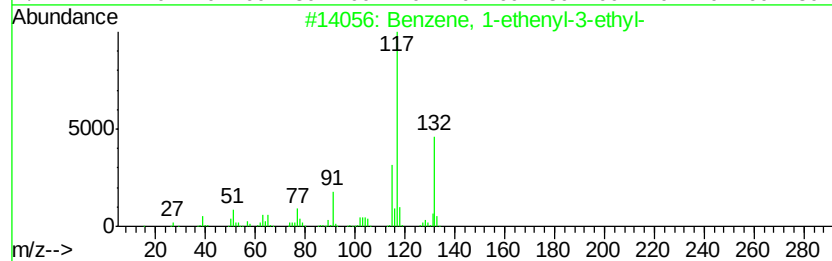
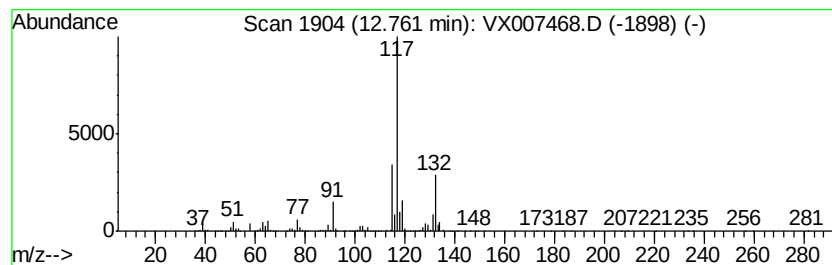
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	91.09 ug/l	4891850	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	90
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	87
5		Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX021119\
Data File : VX007468.D
Acq On : 11 Feb 2019 15:19
Operator : JC/SP
Sample : K1433-10
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
KY001MW031-190207

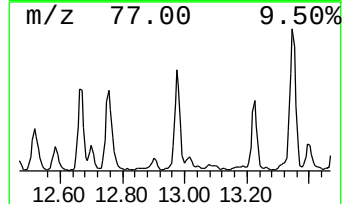
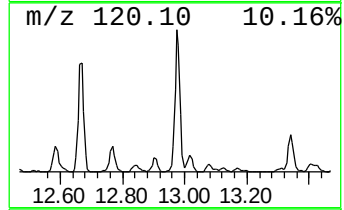
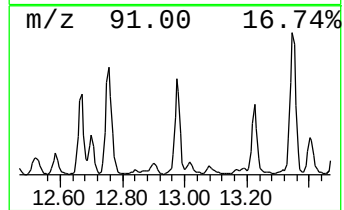
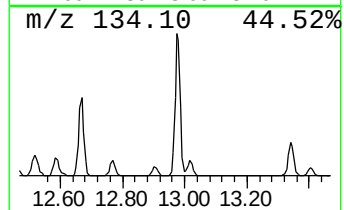
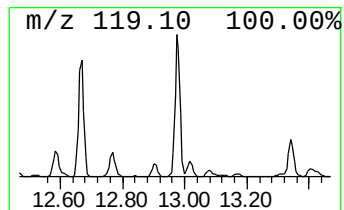
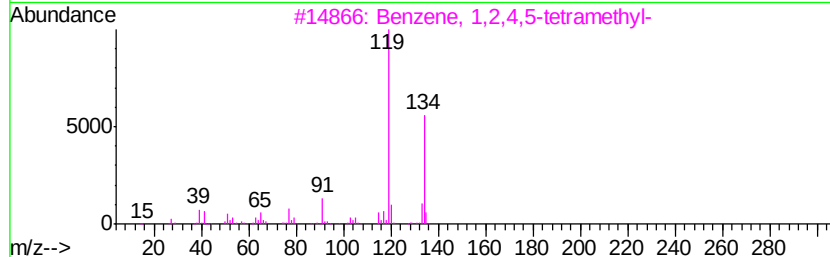
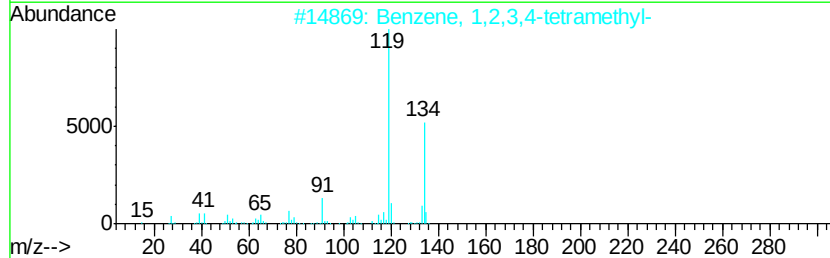
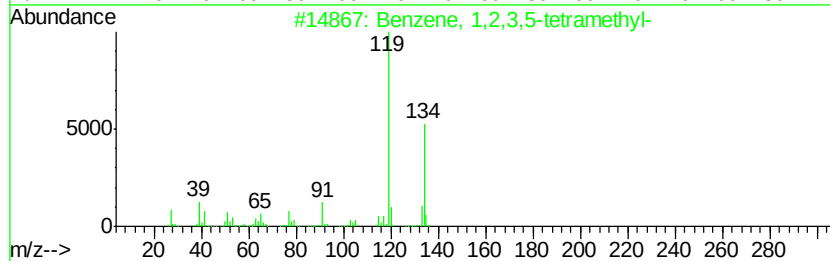
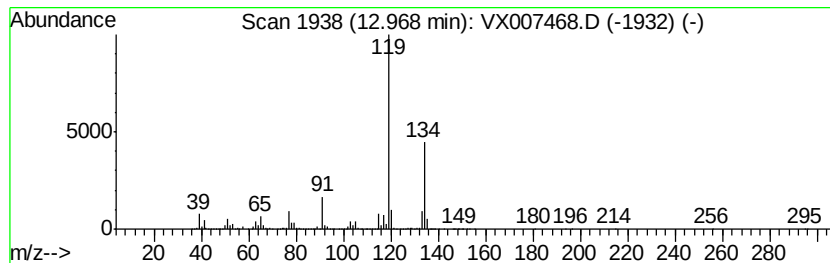
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.97	64.52 ug/l	3464800	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
2		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
4		o-Cymene	134	C10H14	000527-84-4	95
5		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX021119\
Data File : VX007468.D
Acq On : 11 Feb 2019 15:19
Operator : JC/SP
Sample : K1433-10
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

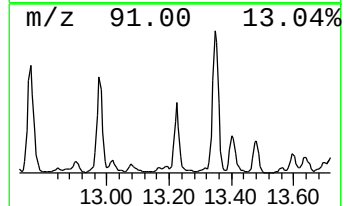
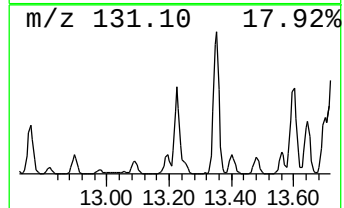
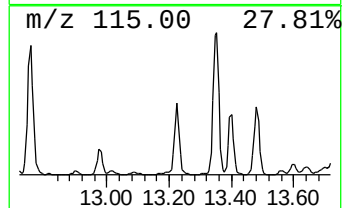
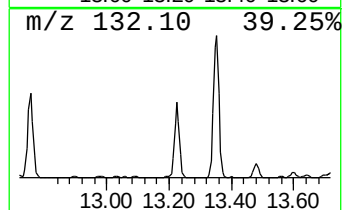
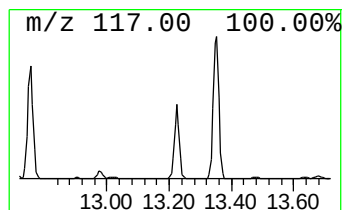
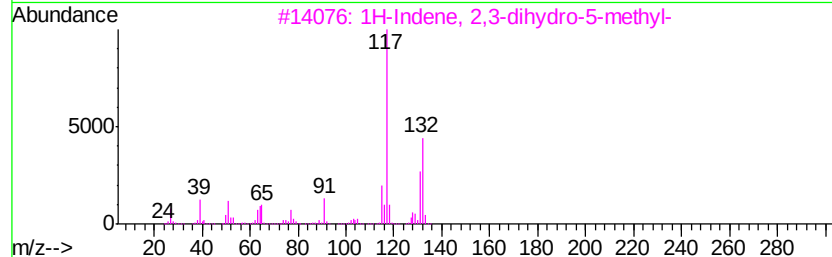
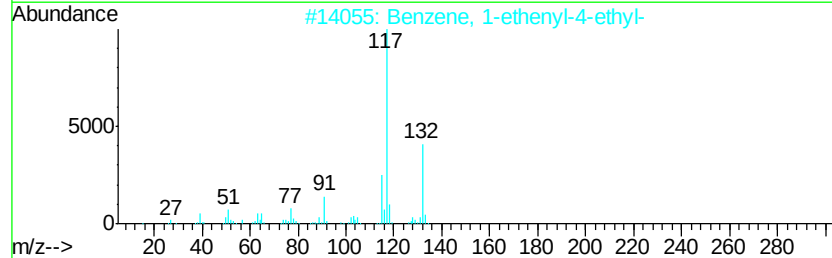
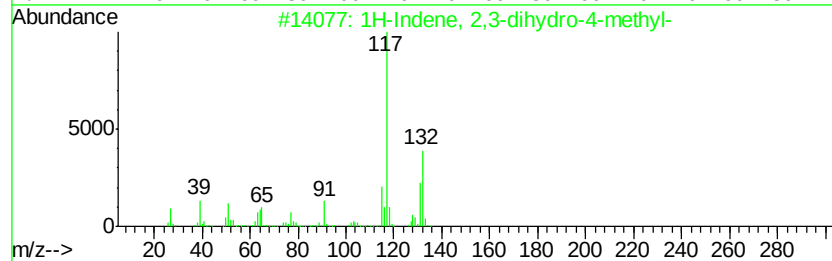
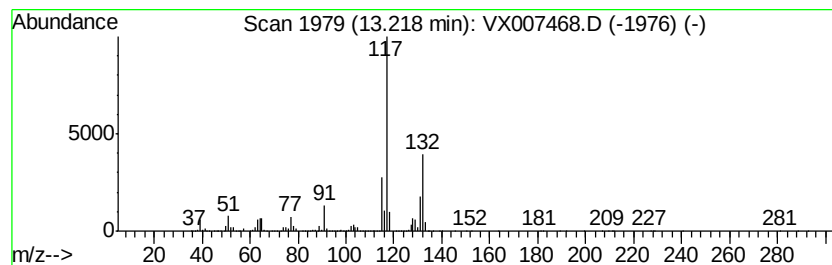
Instrument :
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ClientSampleId :
KY001MW031-190207

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 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.22	61.36 ug/l	3294910	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		1H-Indene, 2,3-dihydro-4-methyl-	132 C10H12	000824-22-6 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 93
4		Benzene, 2-ethenyl-1,4-dimethyl-	132 C10H12	002039-89-6 92
5		Benzene, 1-ethenyl-3-ethyl-	132 C10H12	007525-62-4 90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX021119\
Data File : VX007468.D
Acq On : 11 Feb 2019 15:19
Operator : JC/SP
Sample : K1433-10
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
KY001MW031-190207

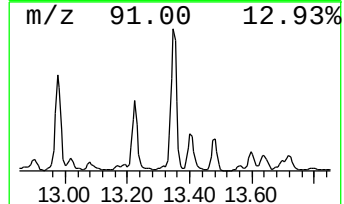
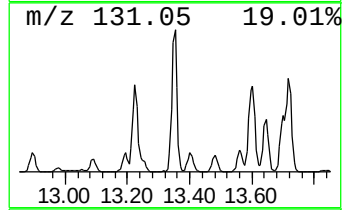
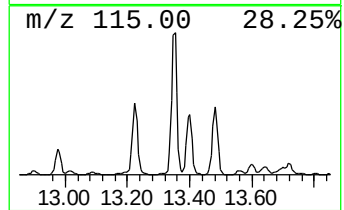
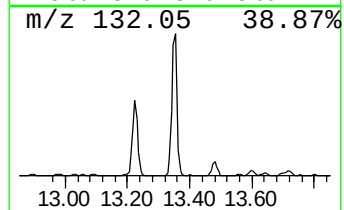
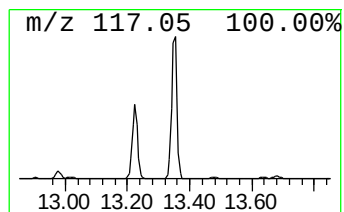
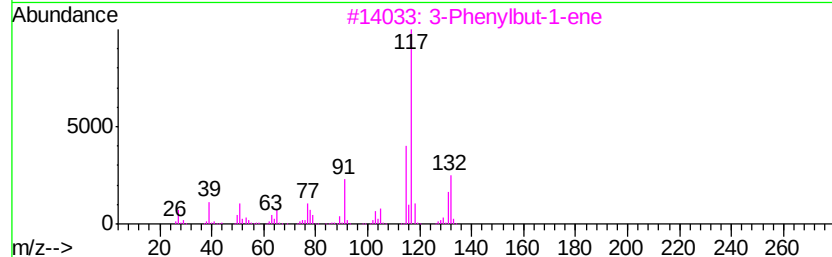
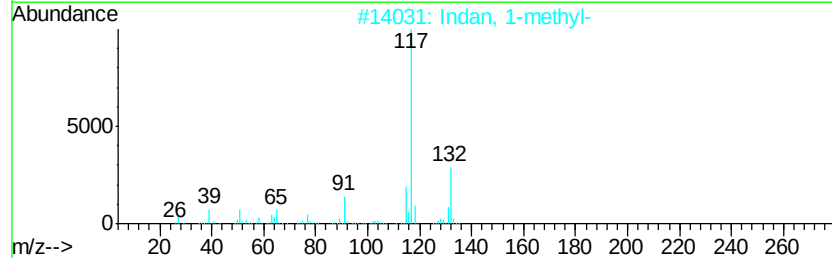
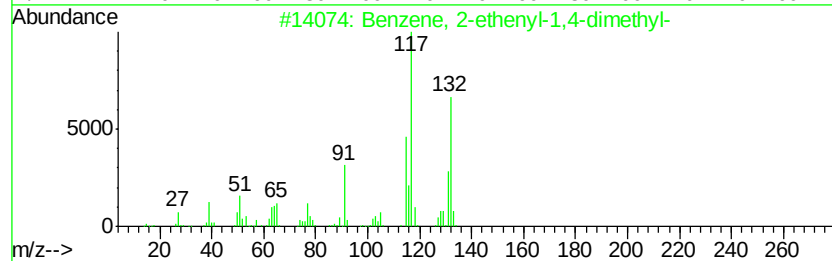
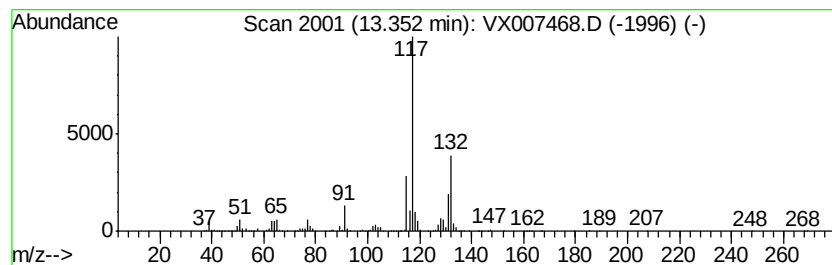
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X020119W.M
Quant Title : SW846 8260

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

Peak Number 9 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	128.83 ug/l	6918730	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	96
2		Indan, 1-methyl-	132	C10H12	000767-58-8	93
3		3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
4		Benzene, 2-butenyl-	132	C10H12	001560-06-1	90
5		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX021119\
Data File : VX007468.D
Acq On : 11 Feb 2019 15:19
Operator : JC/SP
Sample : K1433-10
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

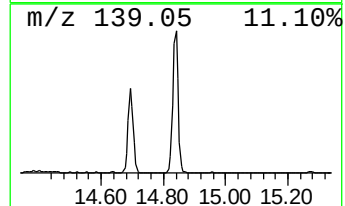
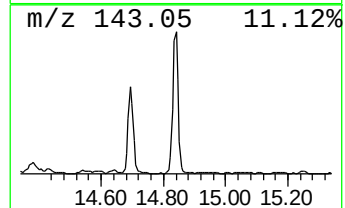
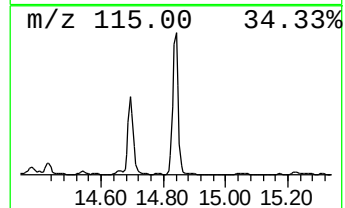
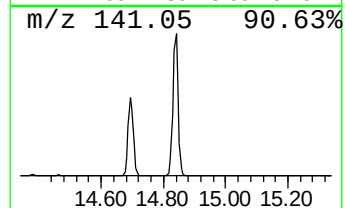
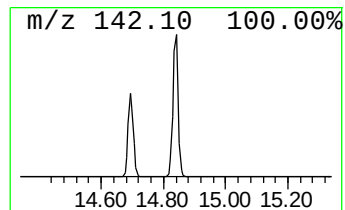
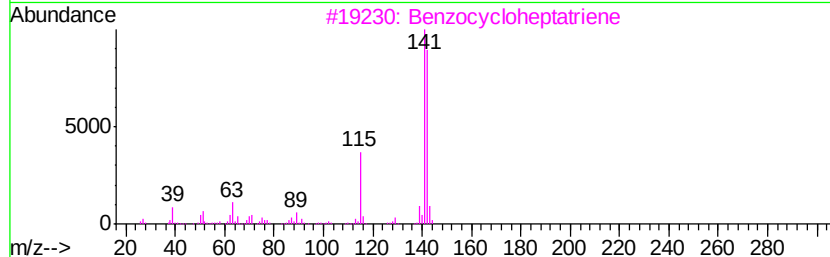
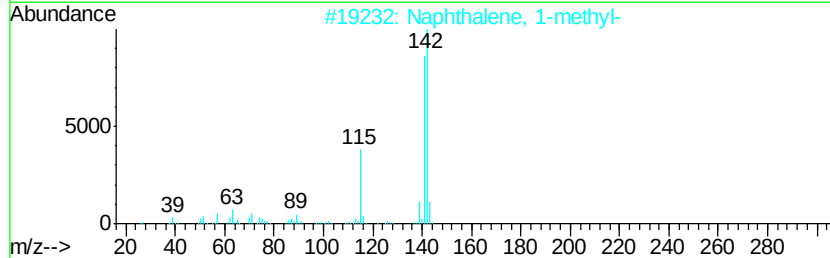
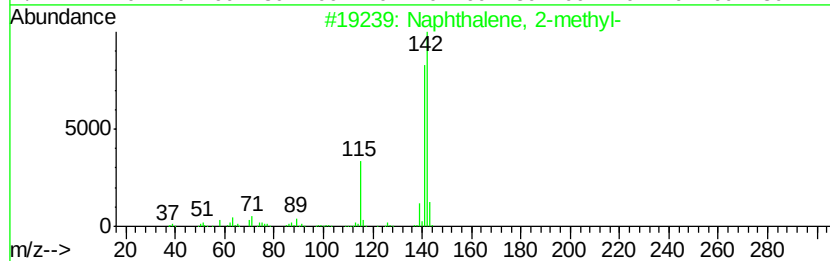
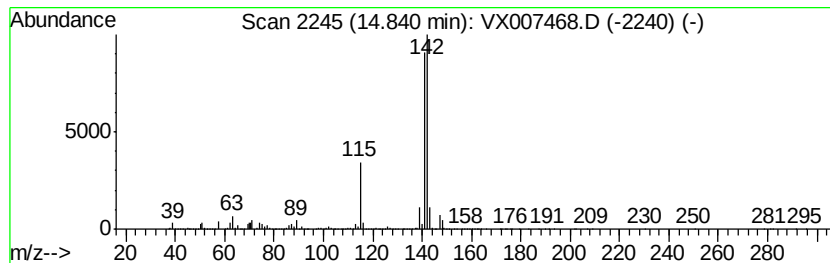
Instrument :
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X020119W.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.84	49.41 ug/l	2653220	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 2-methyl-	142 C11H10	000091-57-6 96
2		Naphthalene, 1-methyl-	142 C11H10	000090-12-0 96
3		Benzocycloheptatriene	142 C11H10	000264-09-5 91
4		1,4-Methanonaphthalene, 1,4-dihy...	142 C11H10	004453-90-1 90
5		1H-Indene, 1-ethylidene-	142 C11H10	002471-83-2 80



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX021119\
Data File : VX007468.D
Acq On : 11 Feb 2019 15:19
Operator : JC/SP
Sample : K1433-10
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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Quant Method : Z:\VOASRV\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
Cyclopentane, met...	4.40	95.2	ug/l	3120010	1	5.66	1638660	50.0
Benzene, 1-ethyl-...	11.67	287.6	ug/l	15445200	4	12.08	2685120	50.0
Indane	12.30	594.5	ug/l	31924600	4	12.08	2685120	50.0
Indene	12.47	48.0	ug/l	2575150	4	12.08	2685120	50.0
Benzene, 4-ethyl-...	12.67	48.6	ug/l	2612330	4	12.08	2685120	50.0
Benzene, 1-etheny...	12.76	91.1	ug/l	4891850	4	12.08	2685120	50.0
Benzene, 1,2,3,5-...	12.97	64.5	ug/l	3464800	4	12.08	2685120	50.0
1H-Indene, 2,3-di...	13.22	61.4	ug/l	3294910	4	12.08	2685120	50.0
Benzene, 2-etheny...	13.35	128.8	ug/l	6918730	4	12.08	2685120	50.0
Naphthalene, 1-me...	14.84	49.4	ug/l	2653220	4	12.08	2685120	50.0