

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX021119\
 Data File : VX007463.D
 Acq On : 11 Feb 2019 13:06
 Operator : JC/SP
 Sample : K1433-02
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 KY001MW030-190207

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	2.849	270	278	282	rBV2	48630	114752	1.94%	0.234%
2	3.178	323	332	335	rBV3	78261	206082	3.49%	0.420%
3	4.306	508	517	526	rBV3	43571	121669	2.06%	0.248%
4	5.507	703	714	722	rBV2	296933	875203	14.81%	1.784%
5	5.665	731	740	746	rBV	466743	1270265	21.49%	2.590%
6	5.933	773	784	797	rBV5	55015	190572	3.22%	0.389%
7	6.068	797	806	816	rBV	300235	768787	13.01%	1.567%
8	6.257	828	837	842	rBV3	50702	157209	2.66%	0.321%
9	6.348	842	852	863	rVV	458099	1470053	24.87%	2.997%
10	6.458	863	870	880	rVB2	75972	212102	3.59%	0.432%
11	6.860	926	936	946	rBV	749250	1821769	30.83%	3.714%
12	6.939	946	949	959	rVB3	53321	125742	2.13%	0.256%
13	7.256	993	1001	1008	rBV	62185	131155	2.22%	0.267%
14	7.445	1021	1032	1043	rBV4	140504	423747	7.17%	0.864%
15	7.573	1045	1053	1058	rBV2	41055	81241	1.37%	0.166%
16	7.634	1058	1063	1069	rBV2	61871	123885	2.10%	0.253%
17	7.848	1091	1098	1105	rVB2	37407	74737	1.26%	0.152%
18	7.951	1105	1115	1122	rBV6	25133	82934	1.40%	0.169%
19	8.055	1122	1132	1141	rVB	295925	621922	10.52%	1.268%
20	8.177	1141	1152	1161	rBV2	321123	747262	12.64%	1.524%
21	8.256	1161	1165	1175	rVV3	63202	131027	2.22%	0.267%
22	8.372	1175	1184	1191	rVB7	27240	61609	1.04%	0.126%
23	8.573	1212	1217	1226	rVB3	88777	169884	2.87%	0.346%
24	8.671	1226	1233	1235	rBV3	77271	139018	2.35%	0.283%
25	8.713	1235	1240	1249	rVB	1581708	2454582	41.53%	5.004%
26	8.982	1280	1284	1289	rVV	38767	61733	1.04%	0.126%
27	9.165	1305	1314	1318	rVB3	65449	122150	2.07%	0.249%
28	9.219	1318	1323	1328	rBV	155374	232678	3.94%	0.474%
29	9.561	1375	1379	1385	rVB3	42549	85292	1.44%	0.174%
30	9.823	1416	1422	1426	rBV4	36775	62325	1.05%	0.127%
31	10.110	1464	1469	1475	rBV	1579426	2063613	34.92%	4.207%
32	11.018	1611	1618	1624	rBV	2431098	3010776	50.94%	6.138%
33	11.134	1632	1637	1643	rBV	1374625	1717340	29.06%	3.501%
34	11.359	1669	1674	1683	rVB	2309042	2782153	47.08%	5.672%

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Integrator: RTE
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35	11.664	1719	1724	1734	rVB	140076	191809	3.25%	0.391%
36	11.768	1734	1741	1745	rBV	94177	121560	2.06%	0.248%
37	11.920	1760	1766	1768	rBV	230167	320418	5.42%	0.653%
38	11.945	1768	1770	1779	rVB	283769	308551	5.22%	0.629%
39	12.079	1786	1792	1798	rBV	1685161	2154979	36.46%	4.394%
40	12.231	1812	1817	1822	rVB	353313	416005	7.04%	0.848%
41	12.298	1822	1828	1835	rBV	4965698	5909968	100.00%	12.049%
42	12.365	1835	1839	1849	rVB2	315724	620110	10.49%	1.264%
43	12.457	1849	1854	1860	rBV	240850	305476	5.17%	0.623%
44	12.585	1869	1875	1881	rBV	42804	60050	1.02%	0.122%
45	12.664	1881	1888	1892	rBV	1433863	1848742	31.28%	3.769%
46	12.701	1892	1894	1898	rVV	491987	517793	8.76%	1.056%
47	12.755	1898	1903	1910	rVV2	1809271	2443781	41.35%	4.982%
48	12.896	1919	1926	1931	rVB	87495	130389	2.21%	0.266%
49	12.975	1931	1939	1944	rBV	1666152	1985944	33.60%	4.049%
50	13.079	1952	1956	1964	rVB2	152065	238564	4.04%	0.486%
51	13.170	1964	1971	1973	rBV2	194885	263209	4.45%	0.537%
52	13.225	1976	1980	1987	rVB	1001165	1231058	20.83%	2.510%
53	13.292	1987	1991	1993	rBV	84997	110385	1.87%	0.225%
54	13.347	1995	2000	2005	rVB2	1893859	2464110	41.69%	5.024%
55	13.402	2005	2009	2012	rBV2	126306	164313	2.78%	0.335%
56	13.481	2018	2022	2028	rVB	204148	258615	4.38%	0.527%
57	13.560	2030	2035	2038	rBV2	126013	187563	3.17%	0.382%
58	13.597	2038	2041	2044	rVV	356949	488477	8.27%	0.996%
59	13.639	2044	2048	2052	rVV3	480988	751868	12.72%	1.533%
60	13.719	2052	2061	2067	rVB2	497243	1032351	17.47%	2.105%
61	13.853	2078	2083	2089	rVB3	36930	59977	1.01%	0.122%
62	13.981	2096	2104	2108	rBV2	152199	219947	3.72%	0.448%
63	14.024	2108	2111	2115	rVB	90897	113155	1.91%	0.231%
64	14.133	2124	2129	2134	rBV	208188	255946	4.33%	0.522%
65	14.273	2147	2152	2157	rBV	193888	320165	5.42%	0.653%
66	14.377	2164	2169	2174	rBV	185370	249475	4.22%	0.509%
67	14.426	2174	2177	2182	rVB	122347	154879	2.62%	0.316%
68	14.542	2191	2196	2200	rBV2	43573	70295	1.19%	0.143%
69	14.694	2217	2221	2225	rVB	130991	153342	2.59%	0.313%
70	14.840	2239	2245	2255	rBV	142242	210078	3.55%	0.428%

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Peak separation: 5

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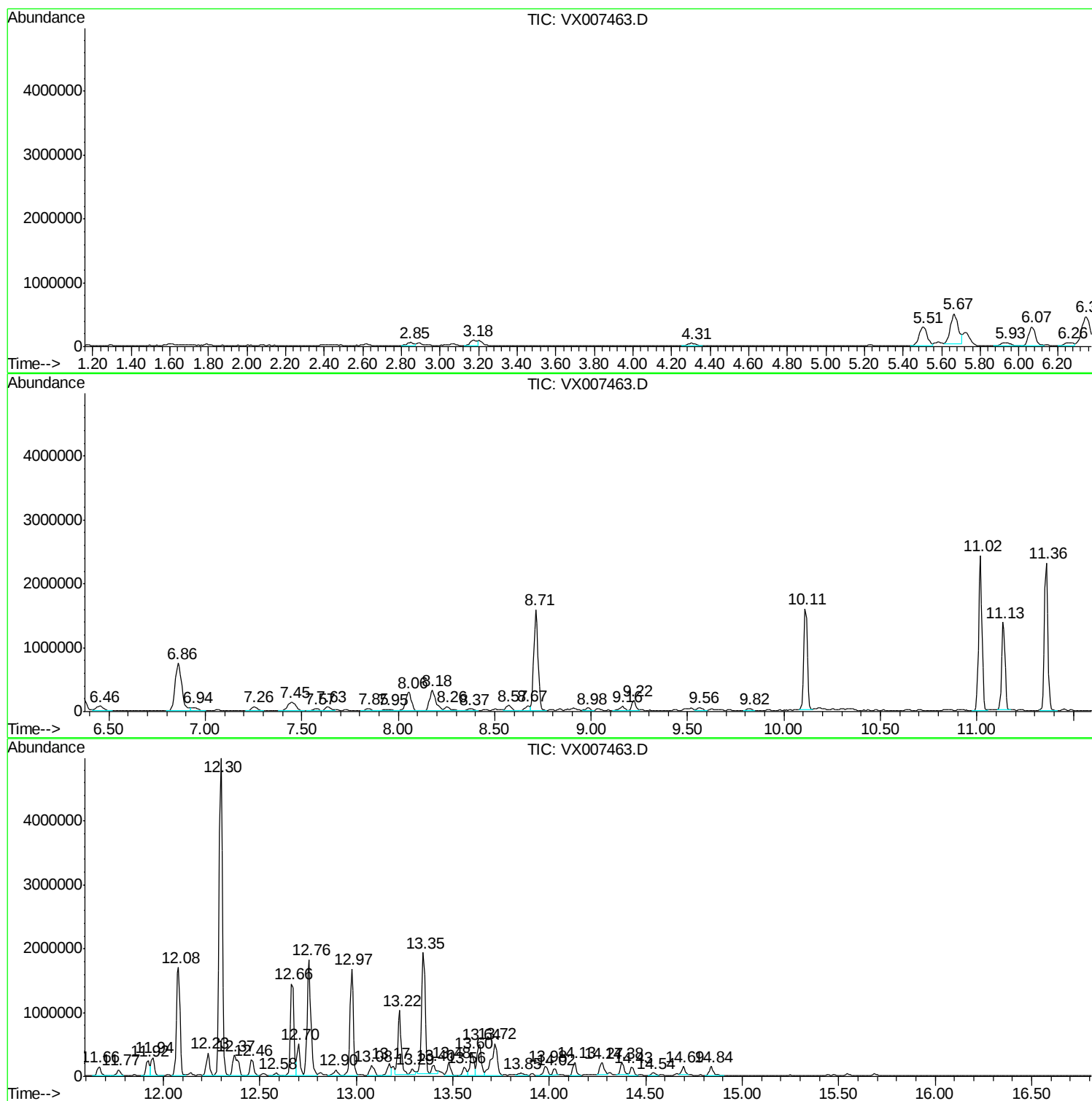
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X020119W.M
Quant Title : SW846 8260

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



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Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
KY001MW030-190207

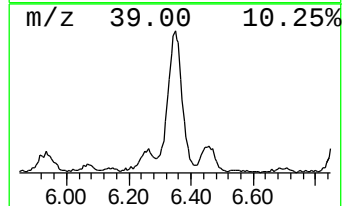
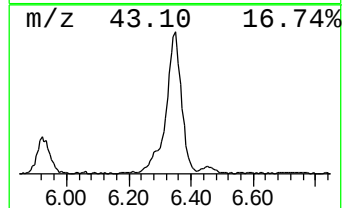
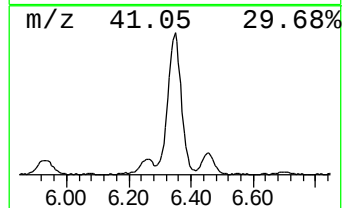
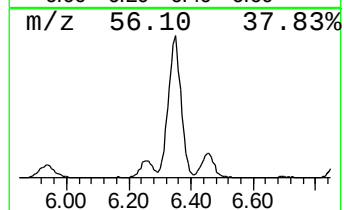
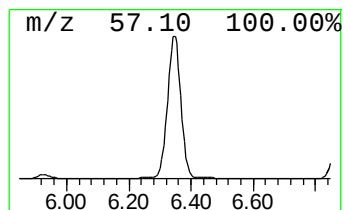
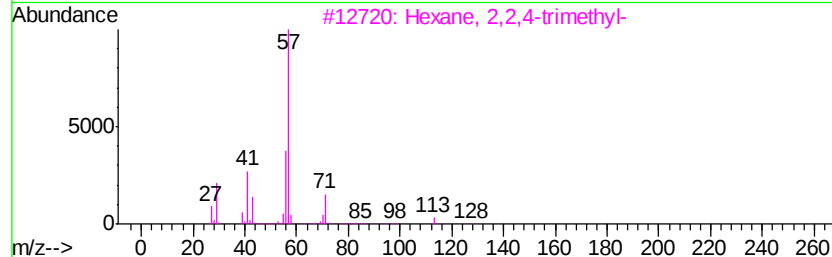
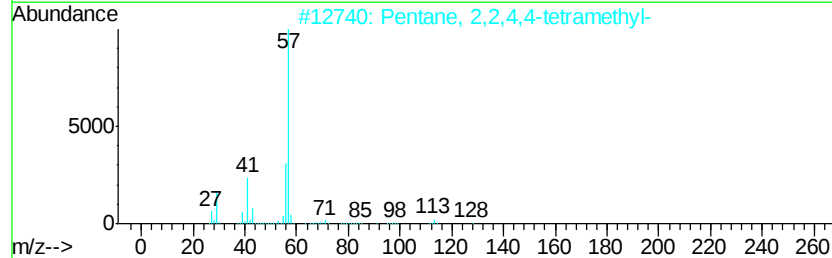
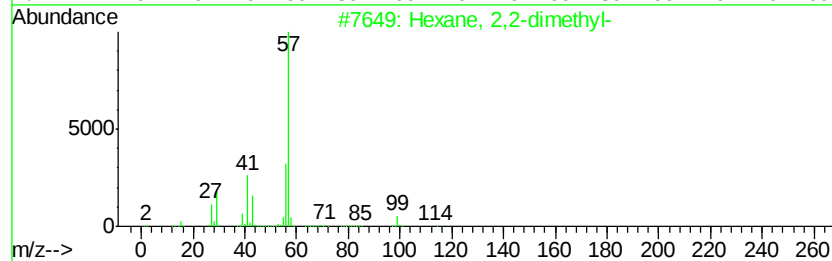
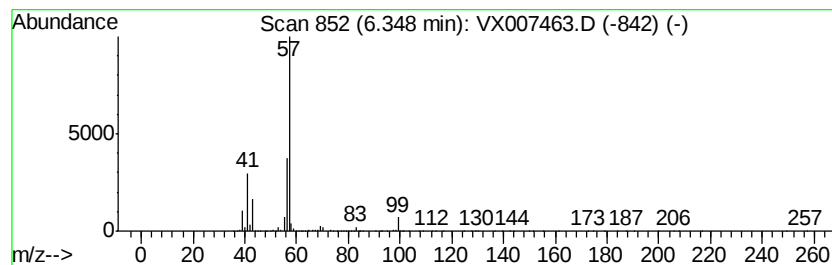
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X020119W.M
Quant Title : SW846 8260

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

Peak Number 1 Hexane, 2,2-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.35	40.35 ug/l	1470050	1,4-Difluorobenzene	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	64
2		Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	59
3		Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	59
4		Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	56
5		Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	50



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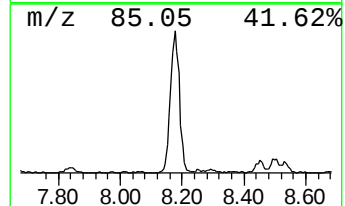
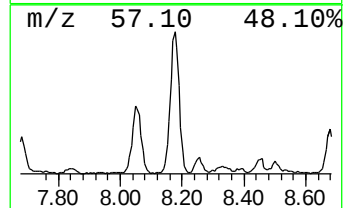
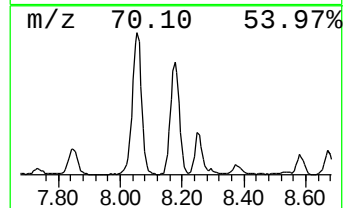
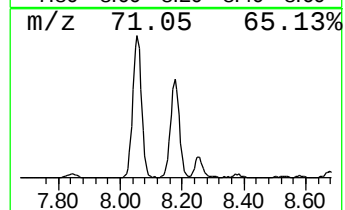
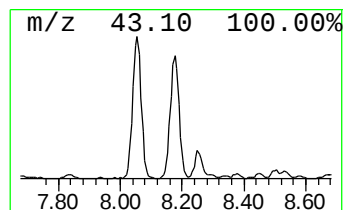
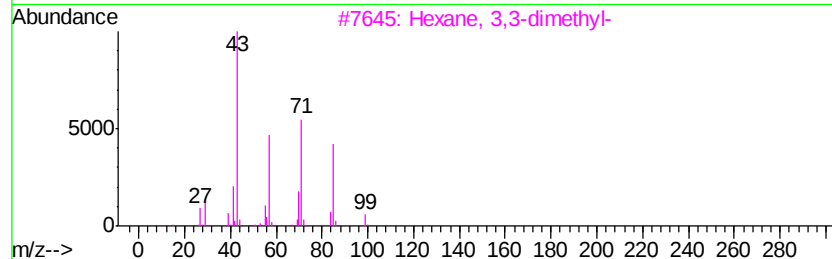
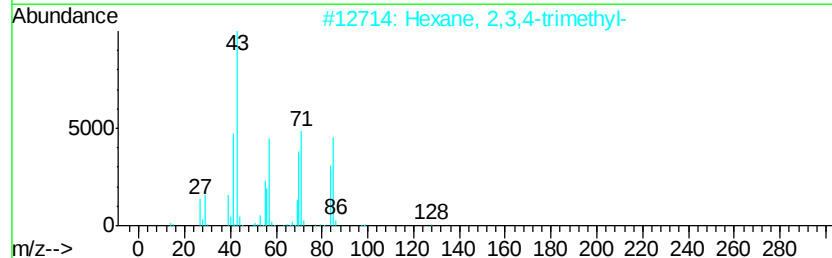
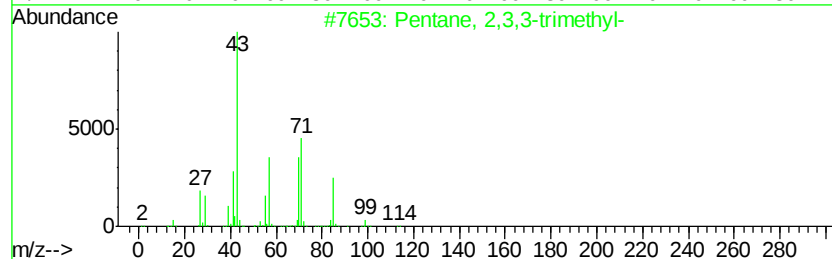
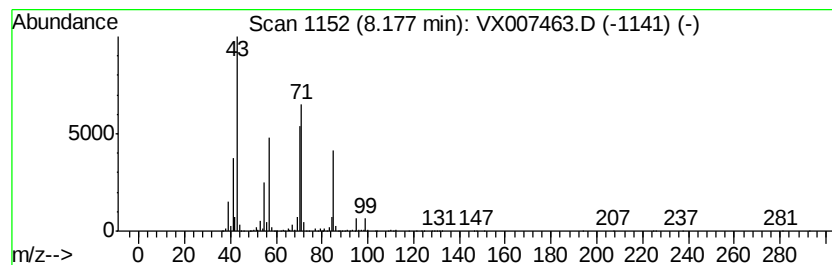
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X020119W.M
 Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.18	20.51 ug/l	747262	1,4-Difluorobenzene	6.86

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			Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	64
4			Sulfurous acid, hexyl pentyl ester	236	C11H24O3S	1000309-14-1	59
5			Propanoic acid, 2-methyl-, 2-eth...	200	C12H24O2	035061-61-1	53



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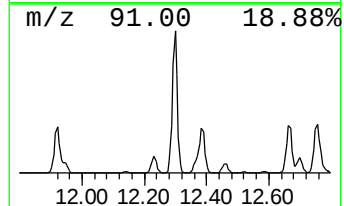
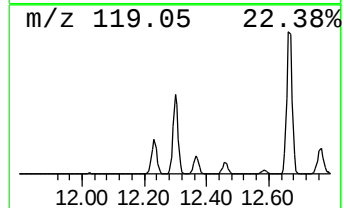
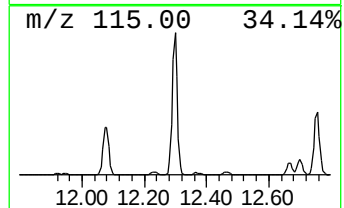
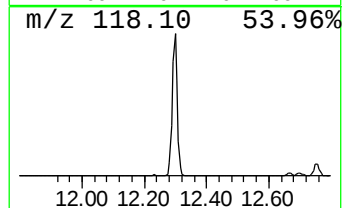
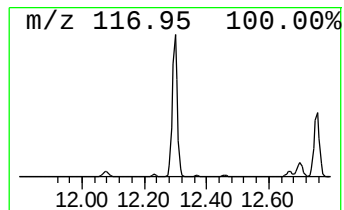
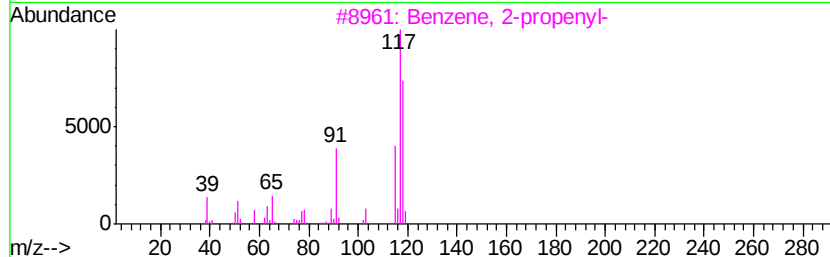
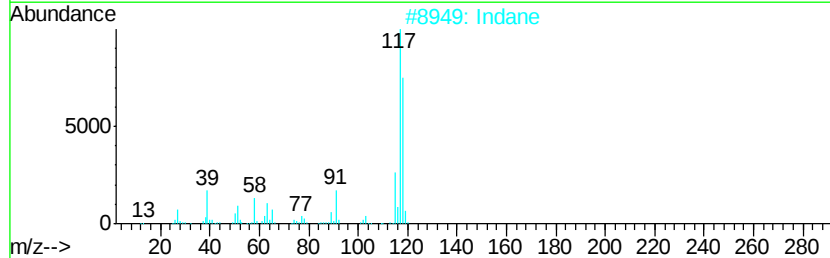
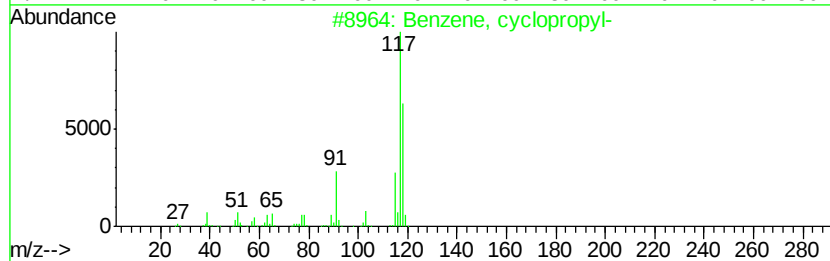
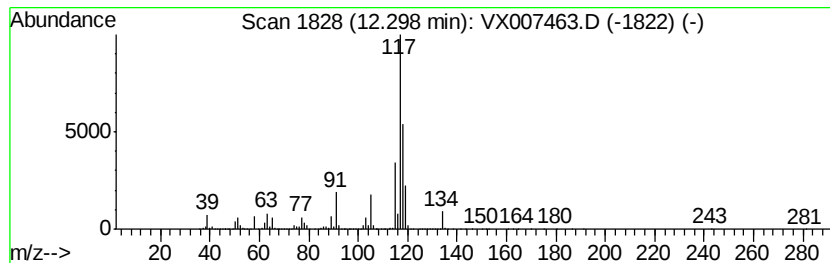
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Quant Title : SW846 8260

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

Peak Number 3 Benzene, cyclopropyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	137.12 ug/l	5909970	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, cvclopropyl-	118	C9H10	000873-49-4	76
2		Indane	118	C9H10	000496-11-7	76
3		Benzene, 2-propenyl-	118	C9H10	000300-57-2	76
4		Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	55
5		Benzene, 1-propenyl-	118	C9H10	000637-50-3	50



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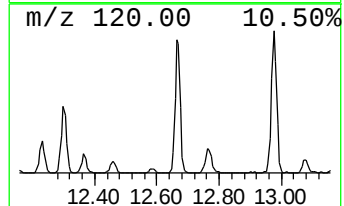
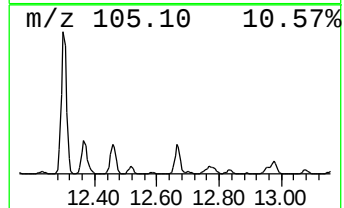
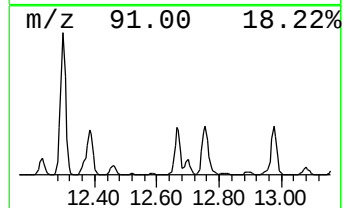
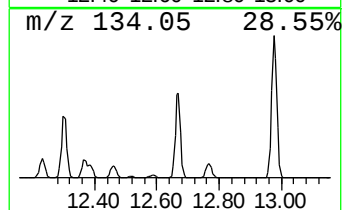
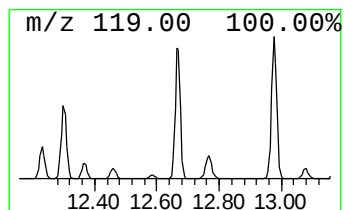
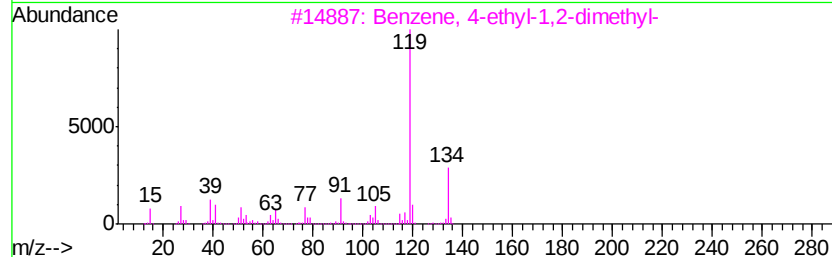
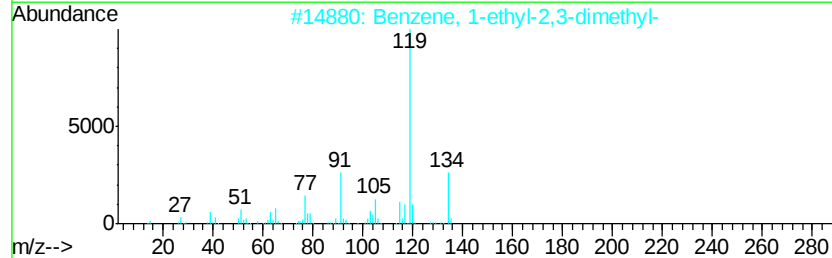
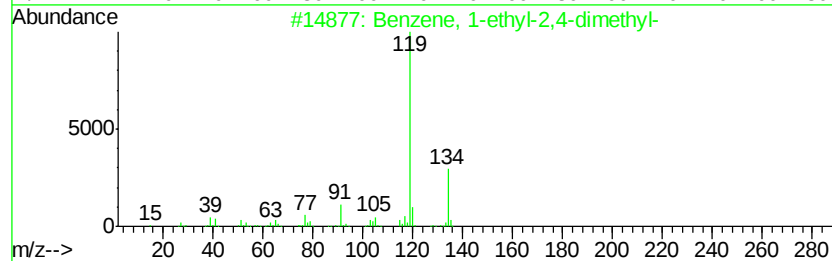
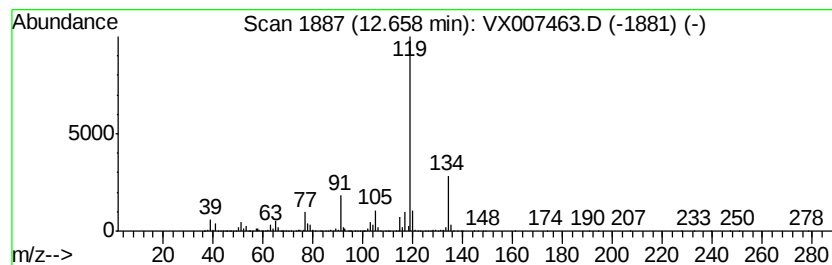
Instrument :
MSVOA_X
ClientSampleId :
KY001MW030-190207

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.66	42.89 ug/l	1848740	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Benzene, 1-ethyl-2,4-dimethyl-	134 C10H14	000874-41-9	97
2	Benzene, 1-ethyl-2,3-dimethyl-	134 C10H14	000933-98-2	96
3	Benzene, 4-ethyl-1,2-dimethyl-	134 C10H14	000934-80-5	95
4	Benzene, 2-ethyl-1,4-dimethyl-	134 C10H14	001758-88-9	95
5	Benzene, 2-ethyl-1,3-dimethyl-	134 C10H14	002870-04-4	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX021119\
Data File : VX007463.D
Acq On : 11 Feb 2019 13:06
Operator : JC/SP
Sample : K1433-02
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
KY001MW030-190207

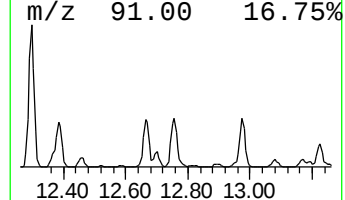
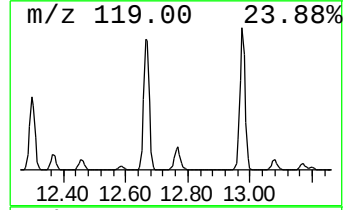
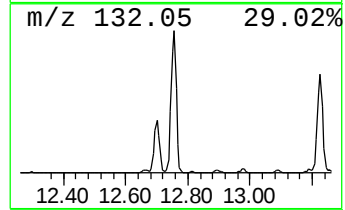
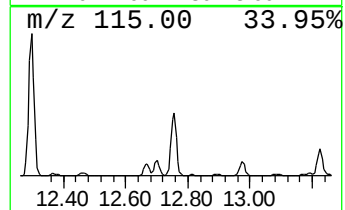
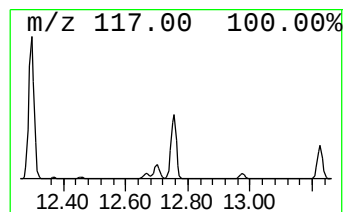
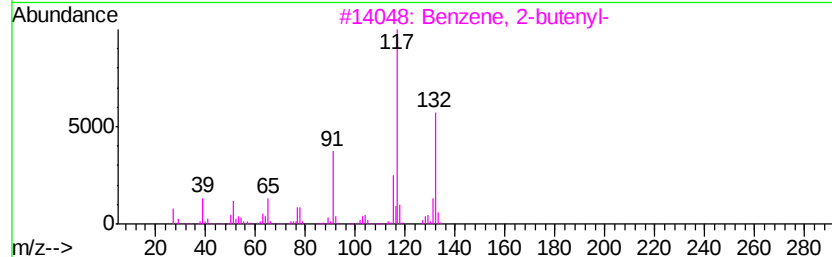
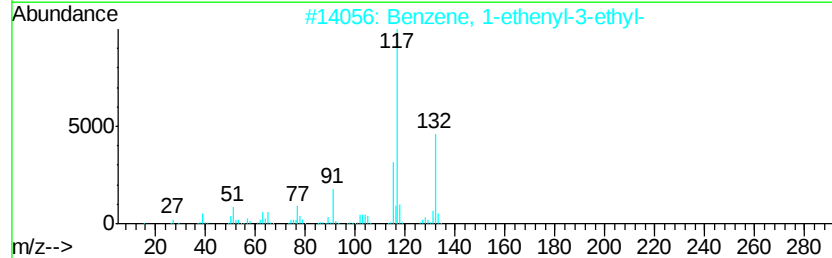
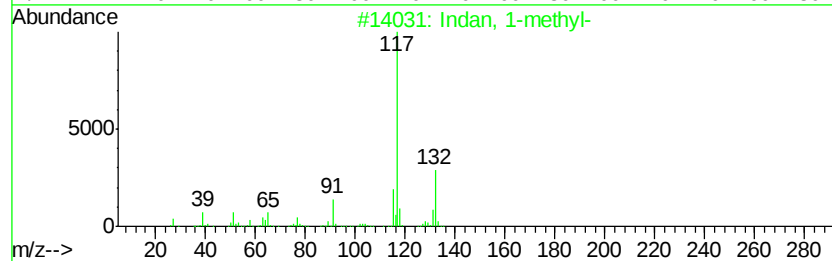
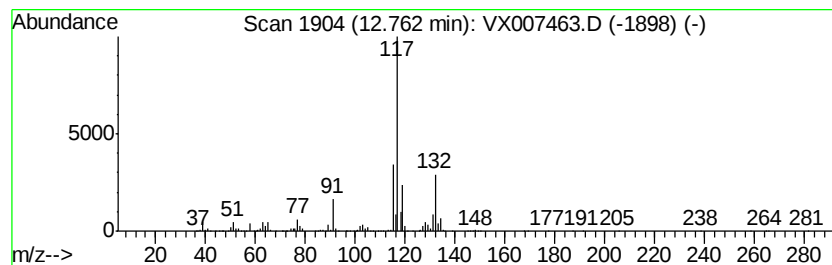
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Quant Title : SW846 8260

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

Peak Number 5 Indan, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	56.70 ug/l	2443780	1,4-Dichlorobenzene-d4	12.08

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indan, 1-methyl-		132	C10H12	000767-58-8	90
2	Benzene, 1-ethenyl-3-ethyl-		132	C10H12	007525-62-4	87
3	Benzene, 2-butenyl-		132	C10H12	001560-06-1	83
4	Benzene, (2-methylcyclopropyl)-		132	C10H12	1000327-39-0	83
5	Benzene, (1-methyl-1-propenyl)-, ...		132	C10H12	000768-00-3	83



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX021119\
Data File : VX007463.D
Acq On : 11 Feb 2019 13:06
Operator : JC/SP
Sample : K1433-02
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
KY001MW030-190207

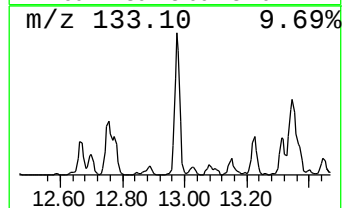
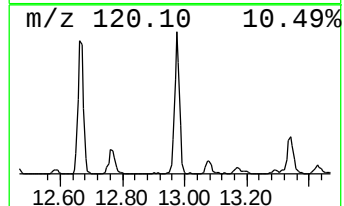
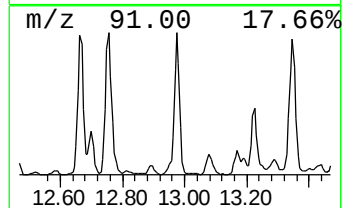
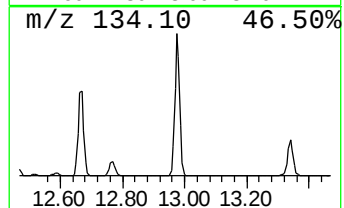
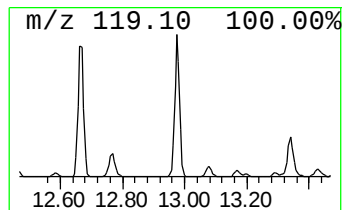
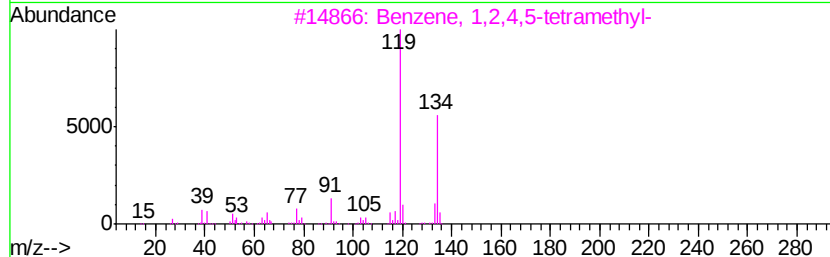
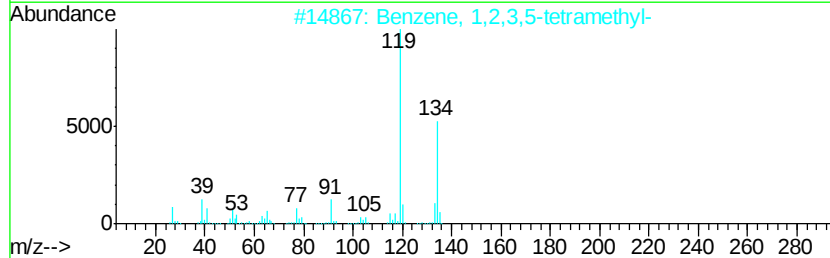
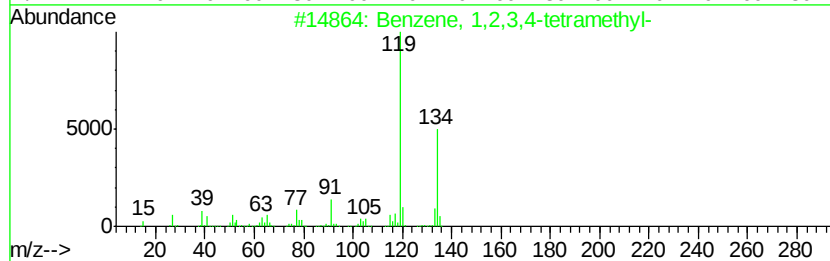
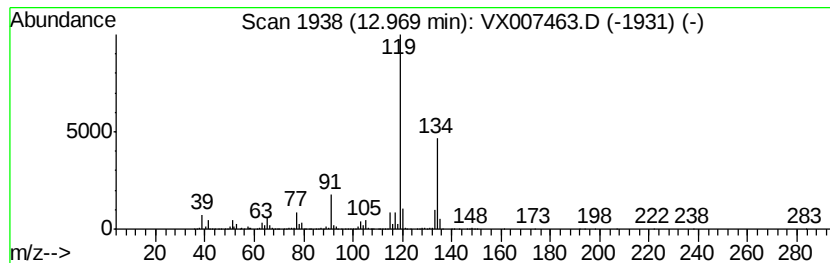
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.97	46.08 ug/l	1985940	1,4-Dichlorobenzene-d4	12.08

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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX021119\
Data File : VX007463.D
Acq On : 11 Feb 2019 13:06
Operator : JC/SP
Sample : K1433-02
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
KY001MW030-190207

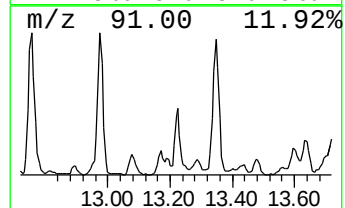
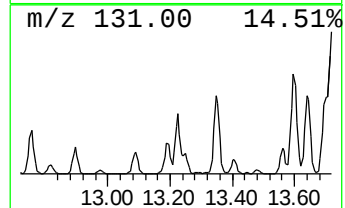
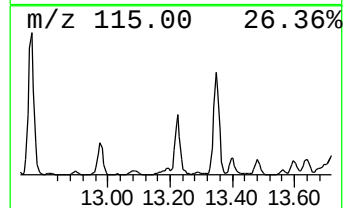
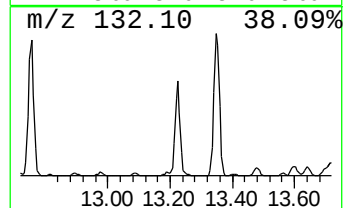
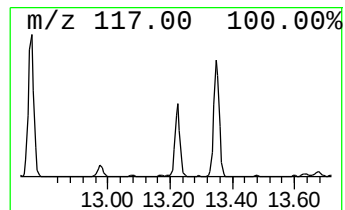
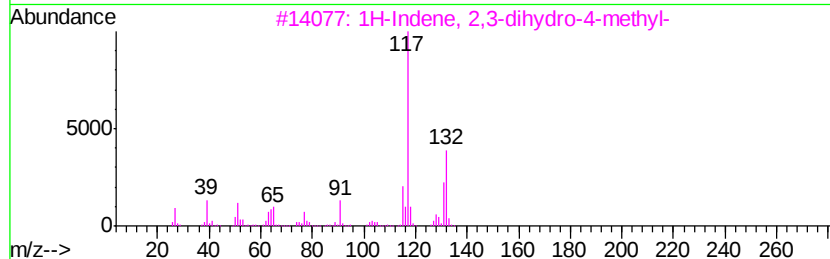
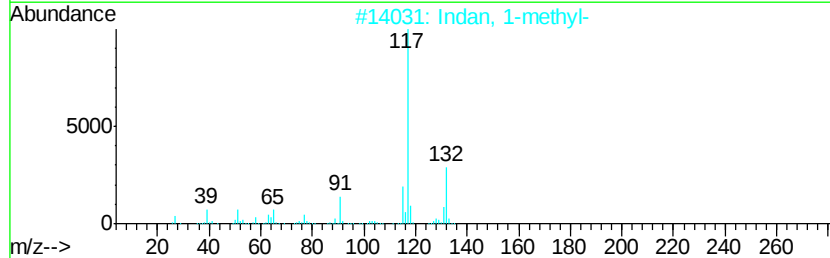
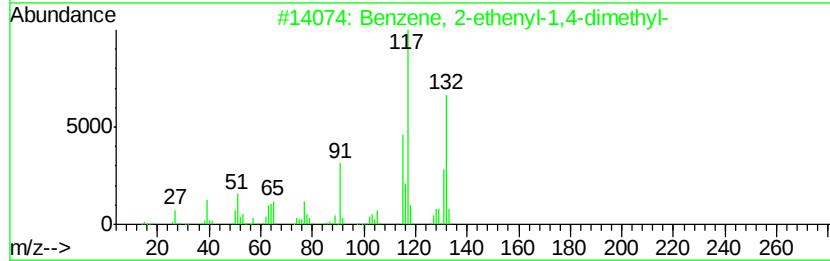
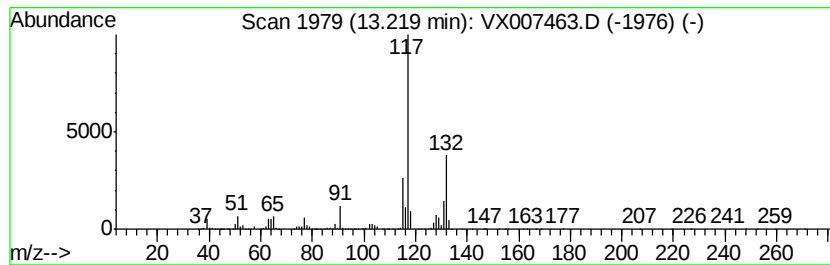
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.22	28.56 ug/l	1231060	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	92
2		Indan, 1-methyl-	132	C10H12	000767-58-8	90
3		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90
4		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	90
5		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX021119\
Data File : VX007463.D
Acq On : 11 Feb 2019 13:06
Operator : JC/SP
Sample : K1433-02
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
KY001MW030-190207

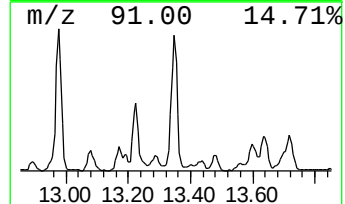
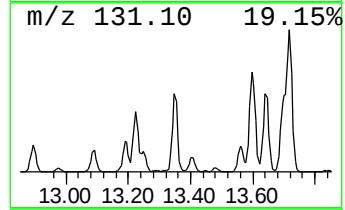
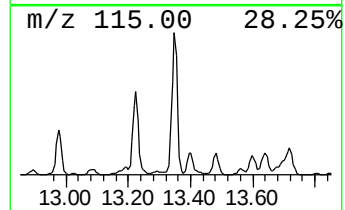
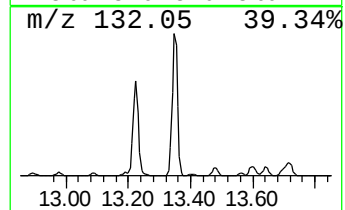
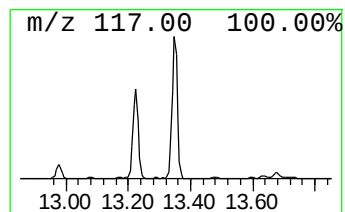
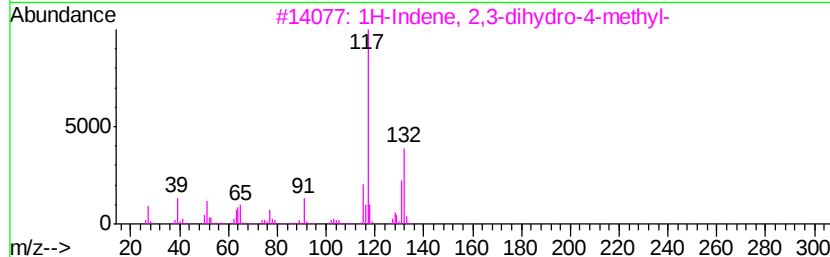
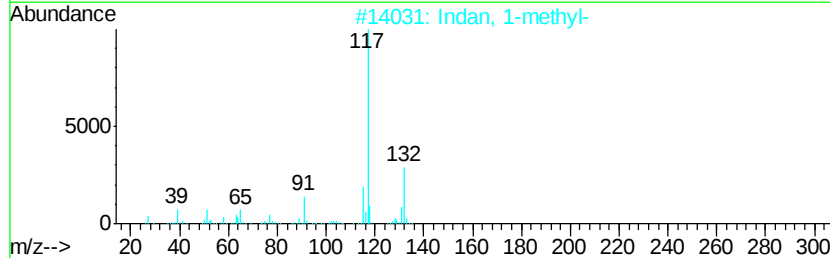
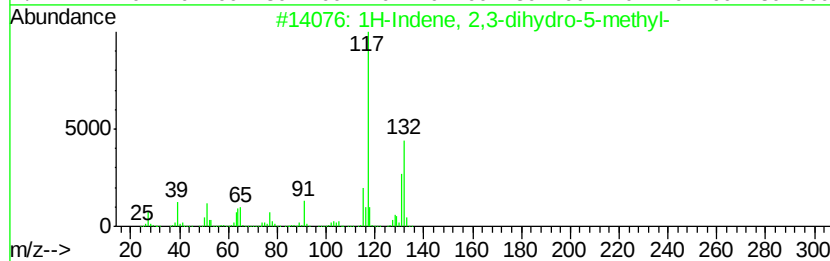
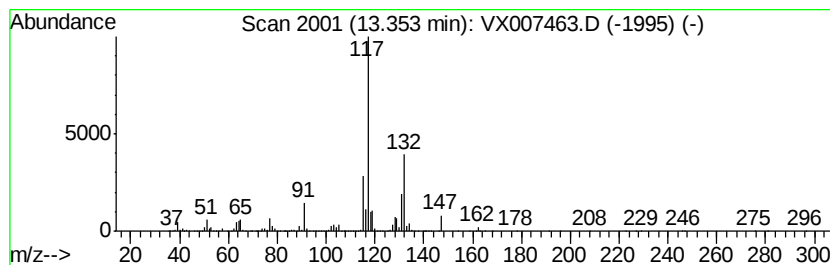
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	57.17 ug/l	2464110	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	87
2		Indan, 1-methyl-	132	C10H12	000767-58-8	76
3		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	76
4		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	74
5		1-Phenyl-1-butene	132	C10H12	000824-90-8	74



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX021119\
Data File : VX007463.D
Acq On : 11 Feb 2019 13:06
Operator : JC/SP
Sample : K1433-02
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
KY001MW030-190207

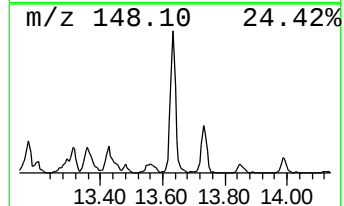
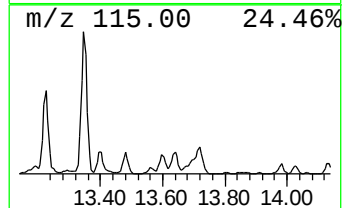
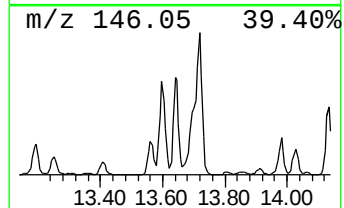
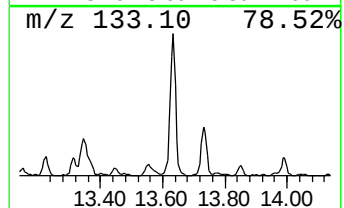
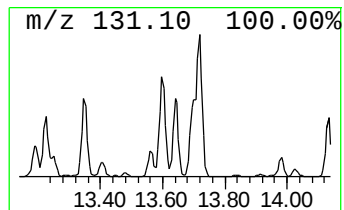
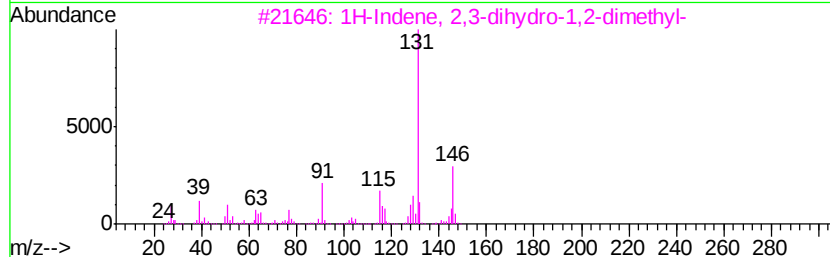
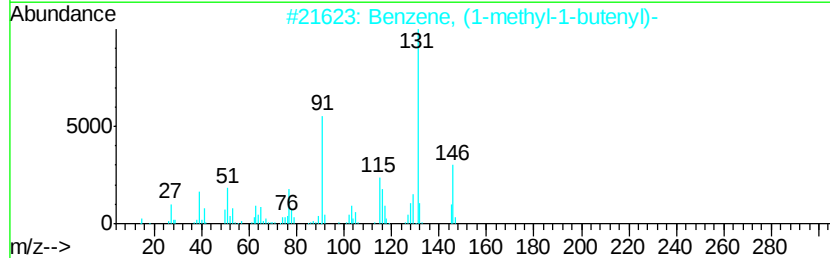
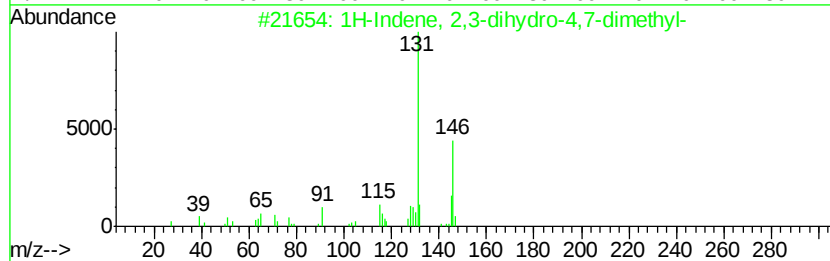
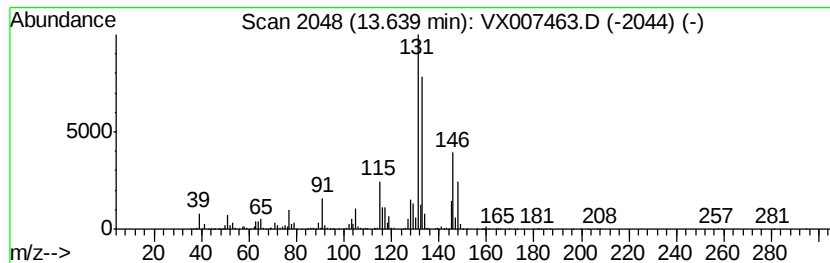
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Quant Title : SW846 8260

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

Peak Number 9 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.64	17.44 ug/l	751868	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	91
2		Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	83
3		1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	83
4		Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	78
5		2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	78



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX021119\
Data File : VX007463.D
Acq On : 11 Feb 2019 13:06
Operator : JC/SP
Sample : K1433-02
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
KY001MW030-190207

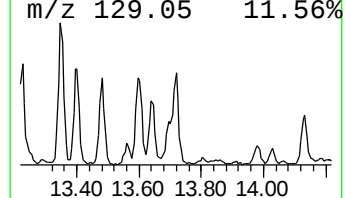
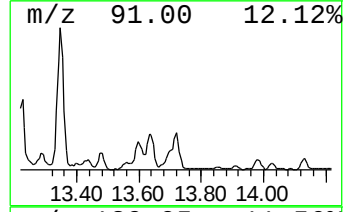
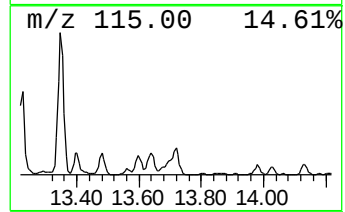
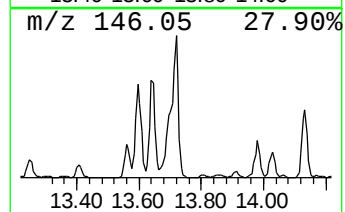
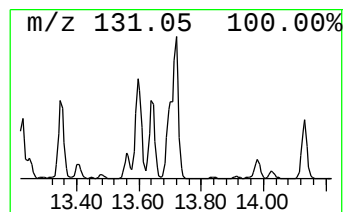
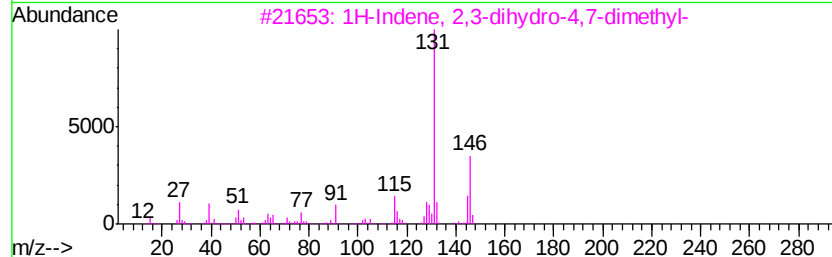
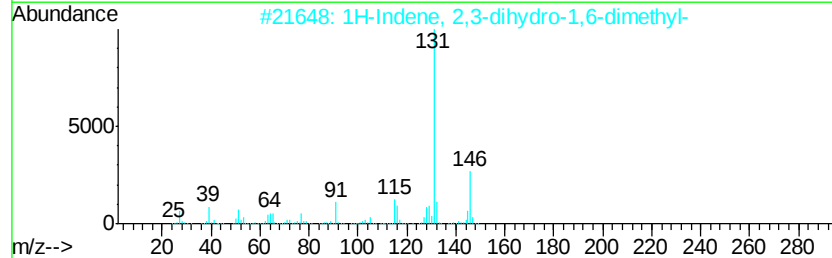
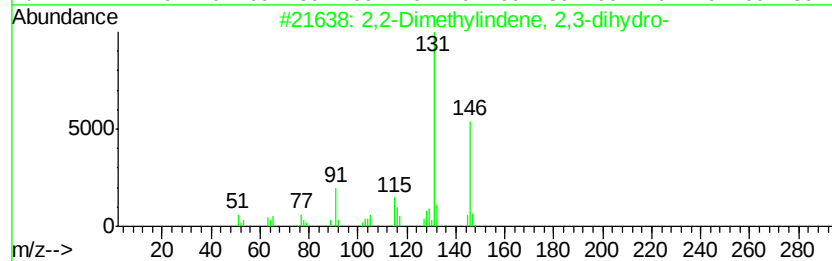
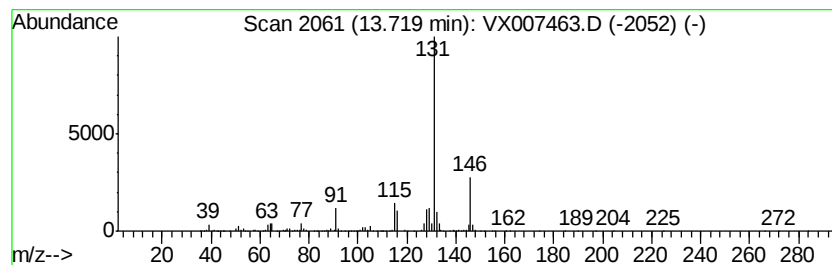
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Quant Title : SW846 8260

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

Peak Number 10 2,2-Dimethylindene, 2,3-dih... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.72	23.95 ug/l	1032350	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	94
2		1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	91
3		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	91
4		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	91
5		1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	91



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX021119\
Data File : VX007463.D
Acq On : 11 Feb 2019 13:06
Operator : JC/SP
Sample : K1433-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
KY001MW030-190207

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
Hexane, 2,2-dimet...	6.35	40.4	ug/l	1470050	2	6.86	1821770	50.0
Pentane, 2,3,3-tr...	8.18	20.5	ug/l	747262	2	6.86	1821770	50.0
Benzene, cyclopro...	12.30	137.1	ug/l	5909970	4	12.08	2154980	50.0
Benzene, 1-ethyl-...	12.66	42.9	ug/l	1848740	4	12.08	2154980	50.0
Indan, 1-methyl-	12.76	56.7	ug/l	2443780	4	12.08	2154980	50.0
Benzene, 1,2,3,4-...	12.97	46.1	ug/l	1985940	4	12.08	2154980	50.0
Benzene, 2-etheny...	13.22	28.6	ug/l	1231060	4	12.08	2154980	50.0
1H-Indene, 2,3-di...	13.35	57.2	ug/l	2464110	4	12.08	2154980	50.0
1H-Indene, 2,3-di...	13.64	17.4	ug/l	751868	4	12.08	2154980	50.0
2,2-Dimethylinden...	13.72	23.9	ug/l	1032350	4	12.08	2154980	50.0