

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX061920\
 Data File : VX016889.D
 Acq On : 19 Jun 2020 18:53
 Operator : JC/SP
 Sample : L3020-23
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-27

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\82X061820W.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	2.002	143	150	161	rVB3	268296	473409	1.09%	0.145%
2	2.252	185	191	200	rVB	973590	1482907	3.42%	0.455%
3	2.916	296	300	319	rVB2	239559	621046	1.43%	0.190%
4	3.794	432	444	453	rBV	265217	635797	1.47%	0.195%
5	4.172	496	506	516	rBV	182917	479921	1.11%	0.147%
6	4.373	529	539	551	rVB	791284	2303383	5.31%	0.706%
7	5.275	676	687	702	rVB	421345	1288854	2.97%	0.395%
8	5.464	705	718	724	rBV2	170078	542193	1.25%	0.166%
9	5.556	724	733	741	rBV2	705762	2191055	5.05%	0.672%
10	5.903	779	790	803	rBV4	192445	628595	1.45%	0.193%
11	6.275	829	851	855	rBV2	307792	1422395	3.28%	0.436%
12	6.330	856	860	868	rVV3	305515	772120	1.78%	0.237%
13	6.434	868	877	888	rVB	275697	788768	1.82%	0.242%
14	6.678	904	917	931	rVB3	197727	507134	1.17%	0.155%
15	6.836	933	943	947	rBV	358930	918796	2.12%	0.282%
16	6.915	952	956	967	rVB4	376403	1077057	2.48%	0.330%
17	7.232	1000	1008	1017	rBV	314812	699910	1.61%	0.215%
18	7.440	1030	1042	1052	rBV	2507832	5969023	13.76%	1.830%
19	7.714	1081	1087	1098	rVB	255051	500681	1.15%	0.153%
20	8.037	1130	1140	1151	rVB3	206970	652568	1.50%	0.200%
21	8.159	1151	1160	1162	rBV2	357997	687906	1.59%	0.211%
22	8.196	1162	1166	1170	rVV	680578	1154148	2.66%	0.354%
23	8.555	1218	1225	1234	rVB2	725056	1320645	3.04%	0.405%
24	8.653	1234	1241	1244	rBV3	477825	951909	2.19%	0.292%
25	8.696	1244	1248	1254	rVB	822226	1355965	3.13%	0.416%
26	8.763	1254	1259	1264	rBV	304419	478982	1.10%	0.147%
27	8.964	1286	1292	1297	rBV4	226698	442838	1.02%	0.136%
28	9.025	1297	1302	1308	rVV2	271767	488691	1.13%	0.150%
29	9.147	1313	1322	1327	rVV	370876	738260	1.70%	0.226%
30	9.208	1327	1332	1343	rVB2	495088	891977	2.06%	0.273%
31	9.549	1383	1388	1393	rVB3	288252	593332	1.37%	0.182%
32	9.604	1393	1397	1401	rBV	463669	640665	1.48%	0.196%
33	10.098	1474	1478	1482	rVB	978349	1244195	2.87%	0.381%
34	10.244	1495	1502	1507	rBV	27887747	39956100	92.11%	12.250%

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Integration Parameters: RTEINT.P

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

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 Max Peaks: 100
 Peak Location: TOP

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 Title : SW846 8260

35	10.348	1511	1519	1525	rVB	14186689	18559549	42.79%	5.690%
36	11.000	1621	1626	1635	rBV	3047754	3967498	9.15%	1.216%
37	11.122	1641	1646	1650	rBV	900632	1151459	2.65%	0.353%
38	11.347	1672	1683	1689	rBV	8275776	10709475	24.69%	3.283%
39	11.439	1693	1698	1702	rVV	8389077	9983909	23.02%	3.061%
40	11.494	1702	1707	1714	rVB	10083861	11991194	27.64%	3.676%
41	11.652	1728	1733	1742	rBV	2853512	3480629	8.02%	1.067%
42	11.799	1751	1757	1762	rVB	32952669	43377156	100.00%	13.298%
43	11.902	1768	1774	1776	rBV	366538	464853	1.07%	0.143%
44	11.933	1776	1779	1783	rVB	441283	544082	1.25%	0.167%
45	12.061	1795	1800	1806	rVB2	1291909	1815831	4.19%	0.557%
46	12.128	1806	1811	1816	rBV	16693914	21661783	49.94%	6.641%
47	12.286	1831	1837	1843	rBV	19912108	24496654	56.47%	7.510%
48	12.360	1843	1849	1856	rVB3	4651115	7449080	17.17%	2.284%
49	12.445	1858	1863	1868	rVB2	487368	619632	1.43%	0.190%
50	12.500	1868	1872	1879	rVB	1398245	1713379	3.95%	0.525%
51	12.573	1879	1884	1886	rBV	2895379	3796481	8.75%	1.164%
52	12.597	1886	1888	1892	rVB	2053022	2016891	4.65%	0.618%
53	12.652	1892	1897	1900	rBV	6217384	7437712	17.15%	2.280%
54	12.689	1900	1903	1907	rVB	1799865	2147415	4.95%	0.658%
55	12.744	1907	1912	1919	rBV	5595864	7627345	17.58%	2.338%
56	12.890	1931	1936	1940	rVB	1625773	2033085	4.69%	0.623%
57	12.963	1940	1948	1951	rBV	3412306	4095046	9.44%	1.255%
58	13.006	1951	1955	1960	rVB	2971816	3725956	8.59%	1.142%
59	13.067	1960	1965	1966	rBV	373726	491920	1.13%	0.151%
60	13.207	1984	1988	1993	rVB	4559457	5567664	12.84%	1.707%
61	13.298	1998	2003	2004	rBV2	367743	471588	1.09%	0.145%
62	13.335	2004	2009	2014	rVV2	7944759	10712104	24.70%	3.284%
63	13.384	2014	2017	2019	rVV3	430241	572472	1.32%	0.176%
64	13.463	2026	2030	2037	rVB	1343548	1667054	3.84%	0.511%
65	13.542	2038	2043	2046	rBV2	565343	751709	1.73%	0.230%
66	13.585	2046	2050	2053	rVV	1387101	1883023	4.34%	0.577%
67	13.621	2053	2056	2062	rVV3	1113570	2035061	4.69%	0.624%
68	13.707	2062	2070	2075	rVB2	1193874	2383854	5.50%	0.731%
69	13.817	2081	2088	2093	rBV	11216394	14726964	33.95%	4.515%
70	13.908	2098	2103	2107	rVB2	463662	670620	1.55%	0.206%
71	13.963	2107	2112	2116	rBV2	525499	703203	1.62%	0.216%

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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

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72	14.012	2116	2120	2123	rBV	482576	501938	1.16%	0.154%
73	14.115	2132	2137	2141	rBV	895968	1059557	2.44%	0.325%
74	14.256	2155	2160	2170	rBV	720172	1263445	2.91%	0.387%
75	14.414	2182	2186	2190	rVB2	543518	673009	1.55%	0.206%
76	14.676	2225	2229	2236	rVB	5447466	6861443	15.82%	2.104%
77	14.822	2248	2253	2262	rVB	2643784	3417965	7.88%	1.048%

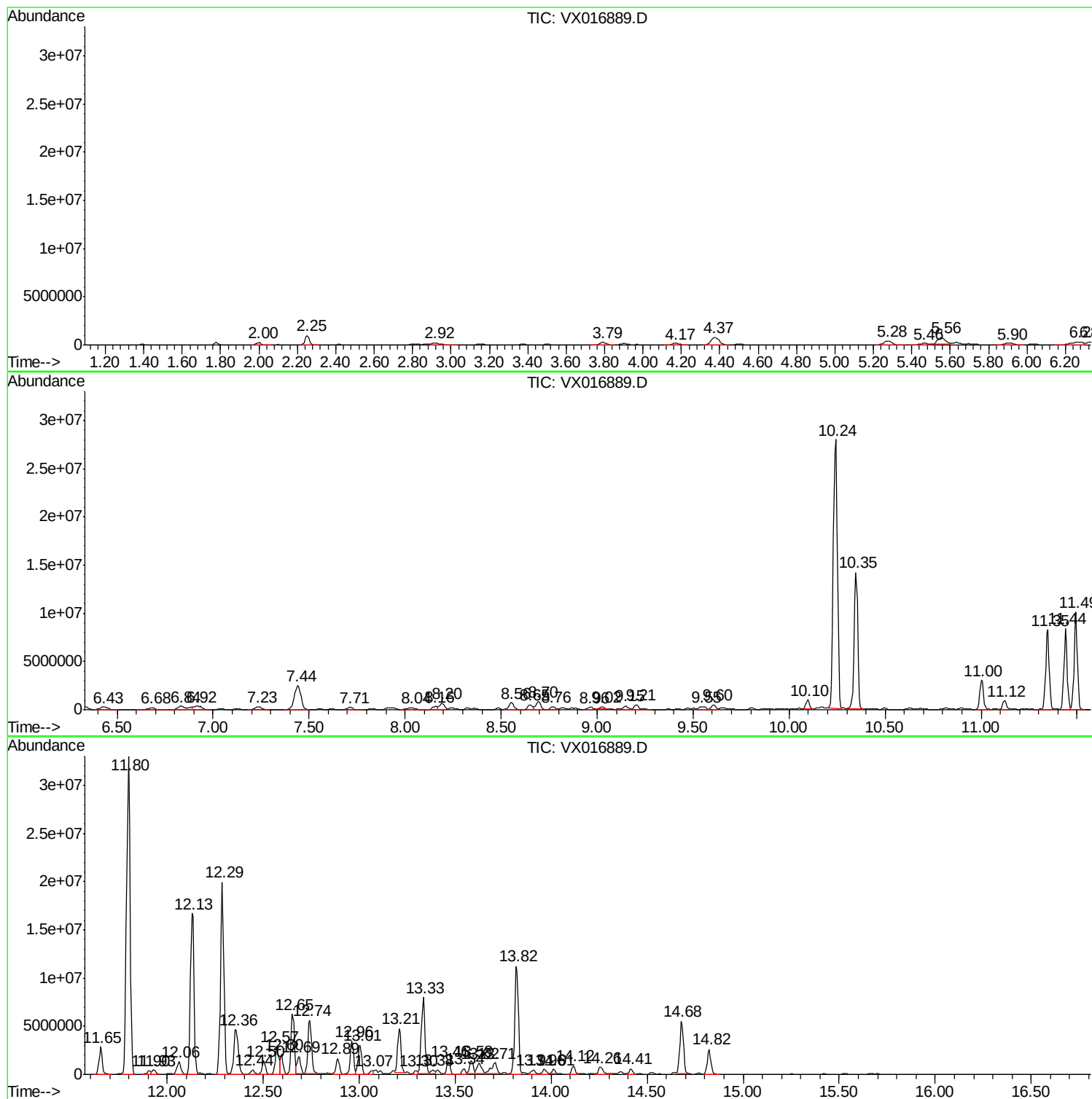
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Instrument :
MSVOA_X
ClientSampleId :
MW-27

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061820W.M
Quant Title : SW846 8260

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



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Sample : L3020-23
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
MW-27

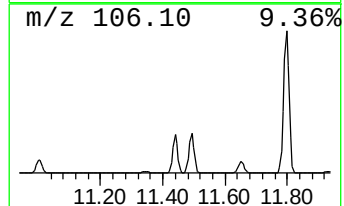
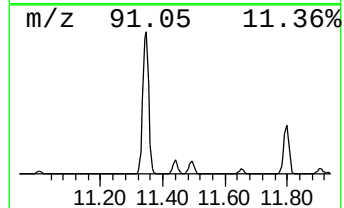
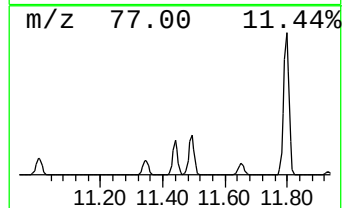
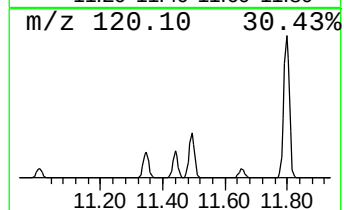
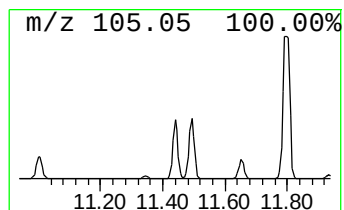
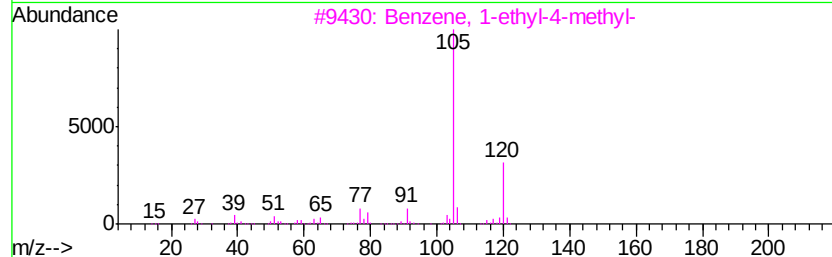
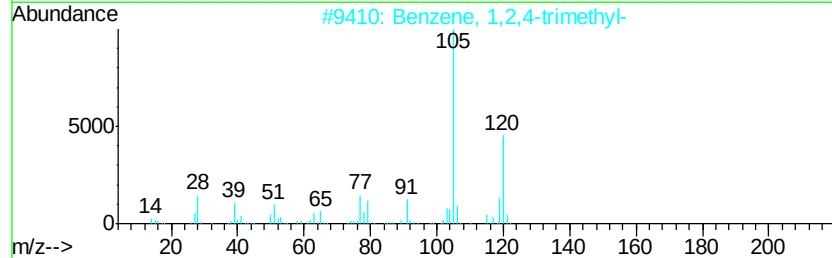
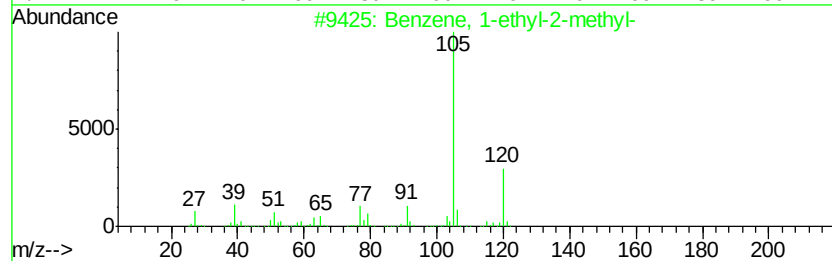
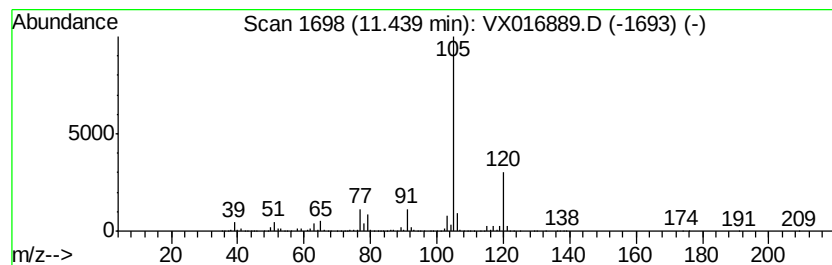
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061820W.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1-ethyl-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.44	274.91 ug/l	9983910	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
3		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5		Mesitylene	120	C9H12	000108-67-8	91



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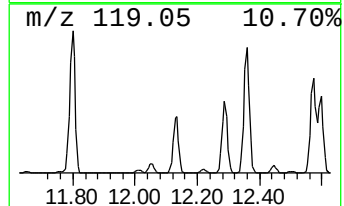
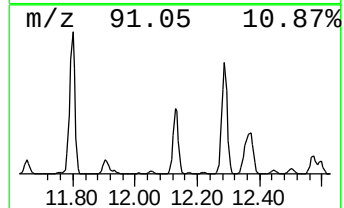
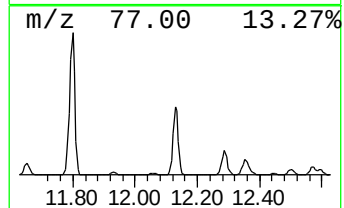
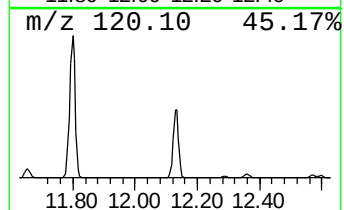
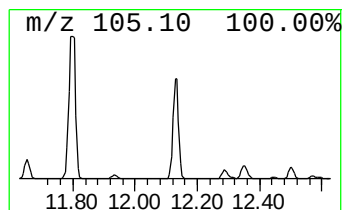
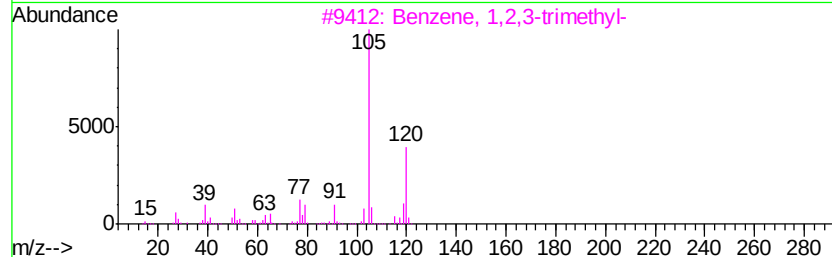
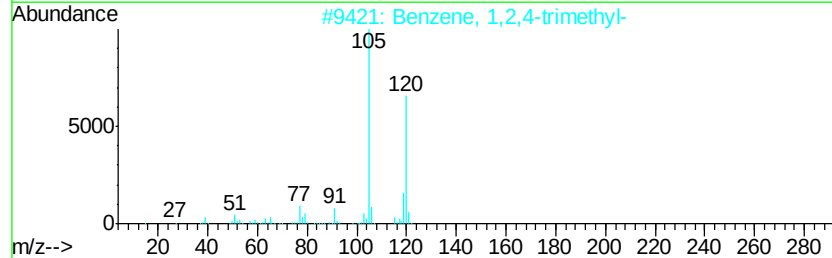
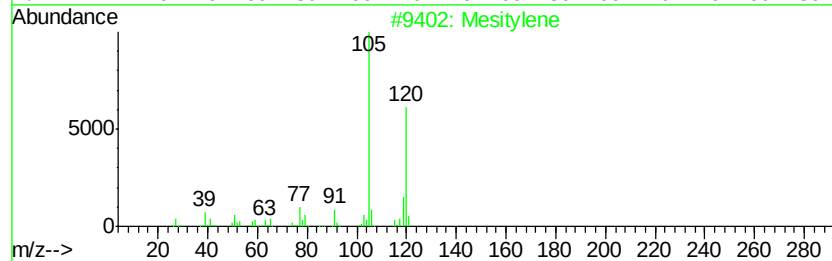
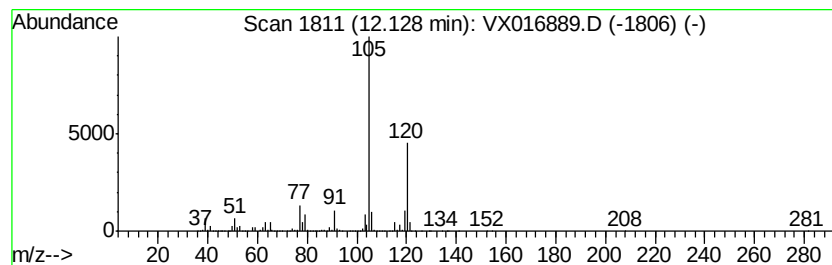
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061820W.M
Quant Title : SW846 8260

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

Peak Number 2 Benzene, 1-ethyl-3-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.13	596.47 ug/l	21661800	1,4-Dichlorobenzene-d4	12.06

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Mesitylene	120	C9H12	000108-67-8	97
2		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
3		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



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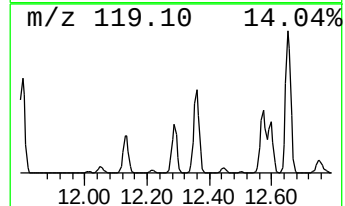
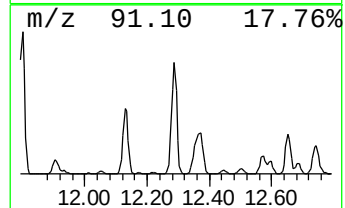
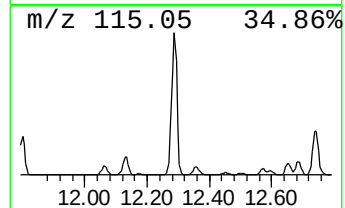
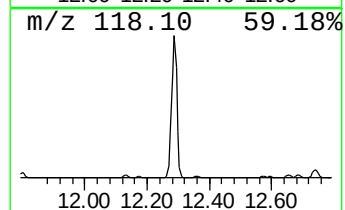
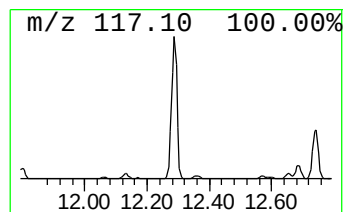
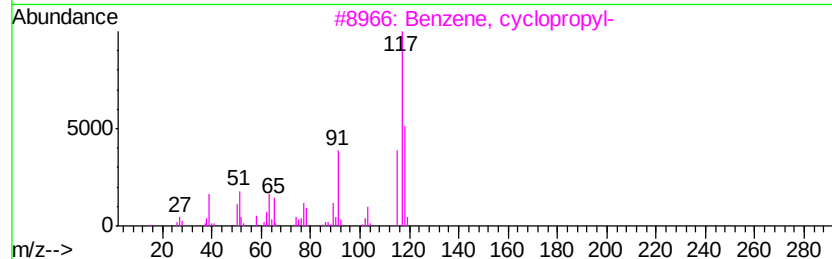
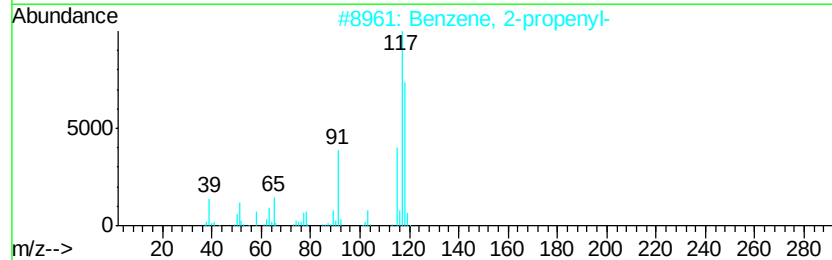
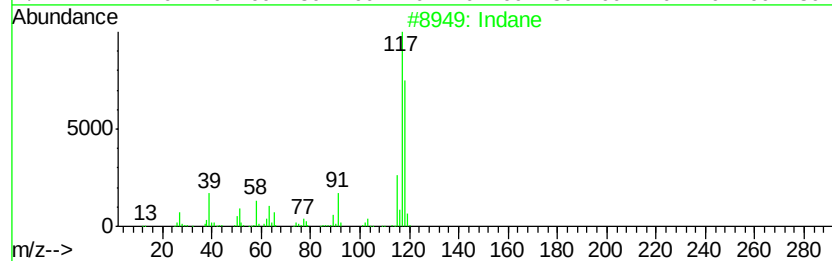
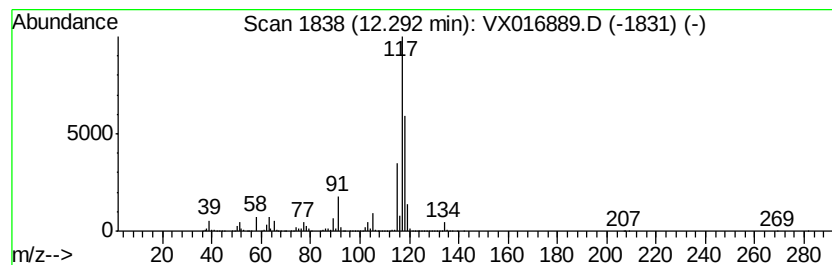
Instrument :
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Quant Title : SW846 8260

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.
12.29	674.53 ug/l	24496700	1,4-Dichlorobenzene-d4	12.06
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 76
3	Benzene, cyclopropyl-		118 C9H10	000873-49-4 76
4	Benzene, 1-ethenyl-2-methyl-		118 C9H10	000611-15-4 55
5	Benzene, 1-ethenyl-4-methyl-		118 C9H10	000622-97-9 49



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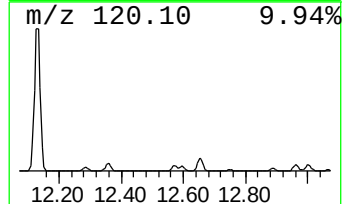
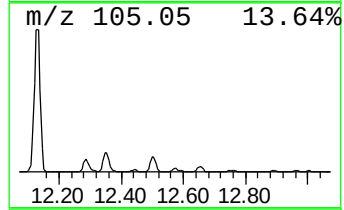
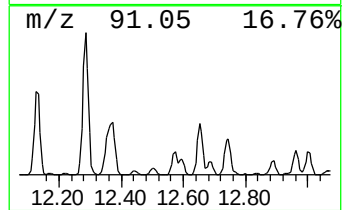
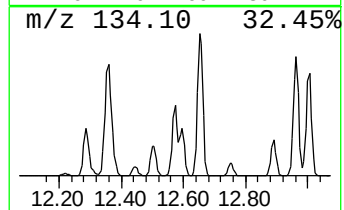
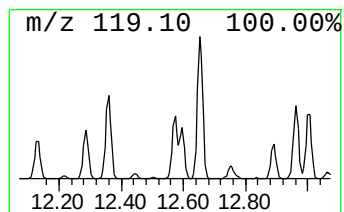
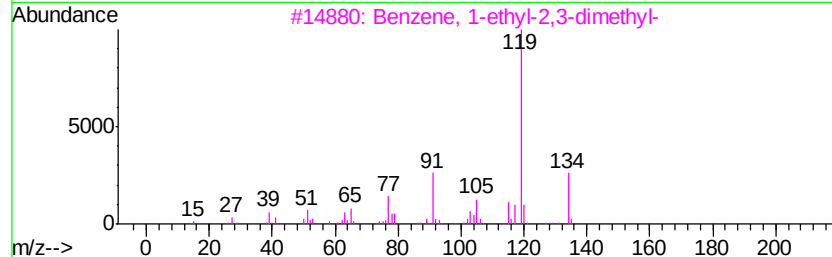
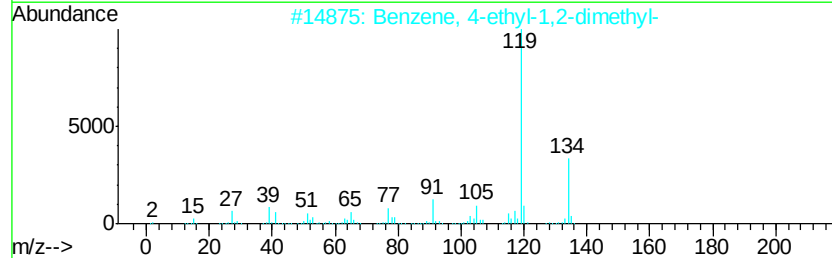
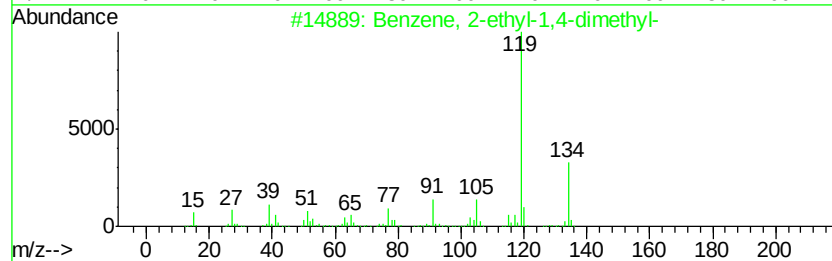
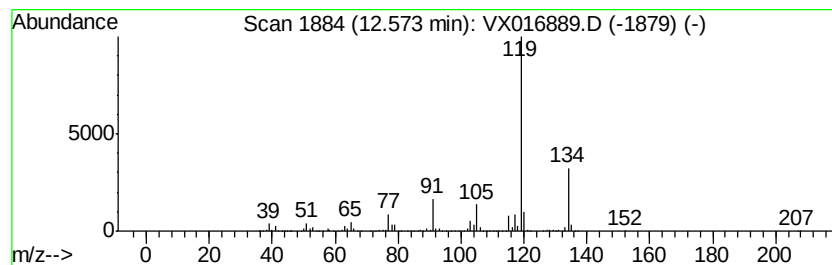
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.57	104.54 ug/l	3796480	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	97
2		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
3		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
4		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94
5		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX061920\
Data File : VX016889.D
Acq On : 19 Jun 2020 18:53
Operator : JC/SP
Sample : L3020-23
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-27

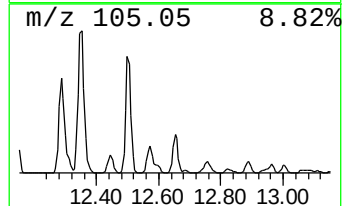
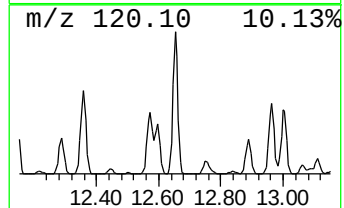
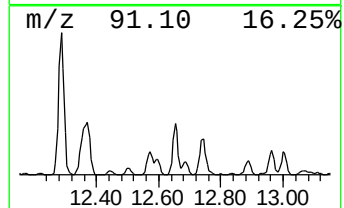
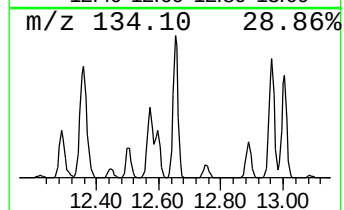
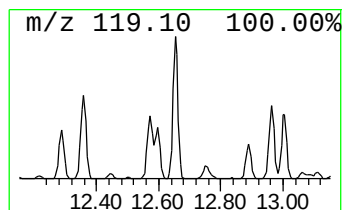
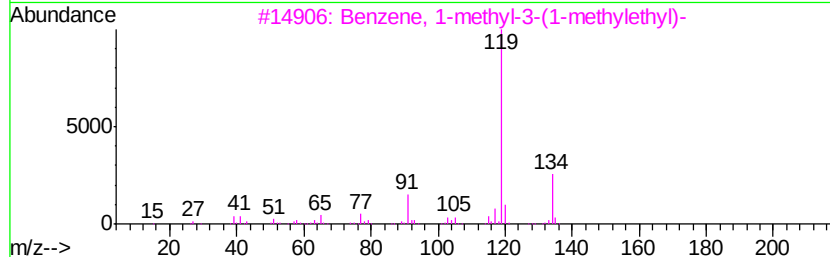
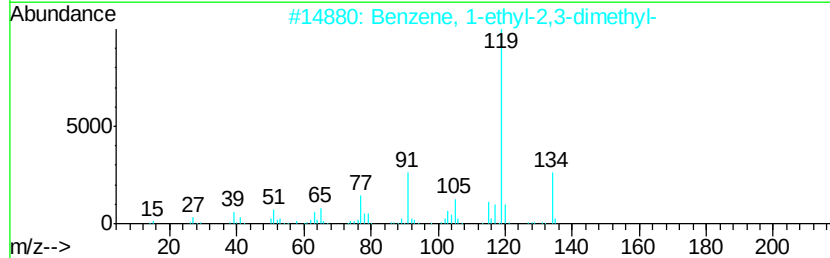
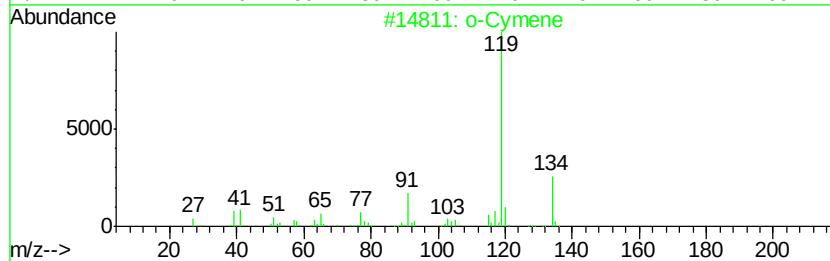
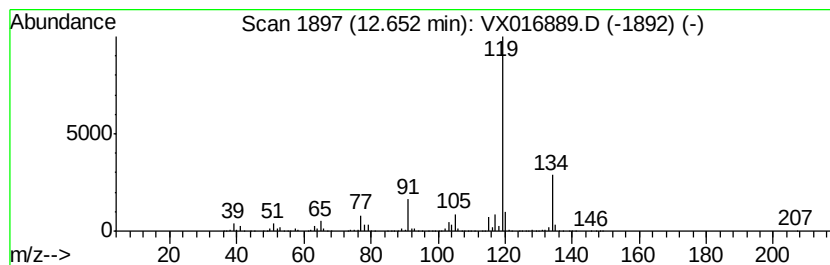
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Quant Title : SW846 8260

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

Peak Number 5 o-Cymene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.65	204.80 ug/l	7437710	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	o-Cymene	134	C10H14	000527-84-4	97
2		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
3		Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	95
4		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
5		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX061920\
Data File : VX016889.D
Acq On : 19 Jun 2020 18:53
Operator : JC/SP
Sample : L3020-23
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-27

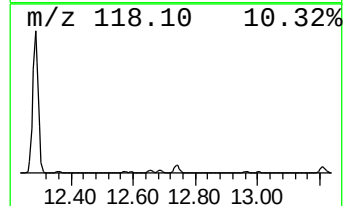
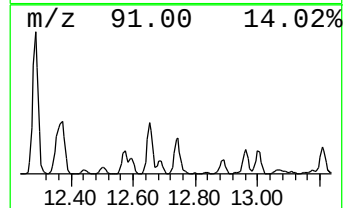
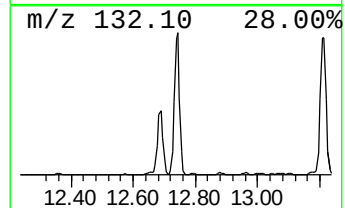
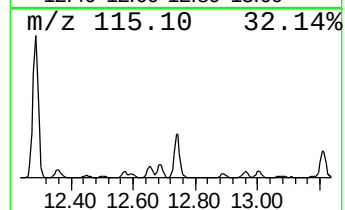
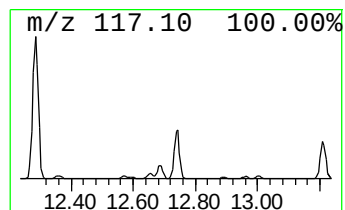
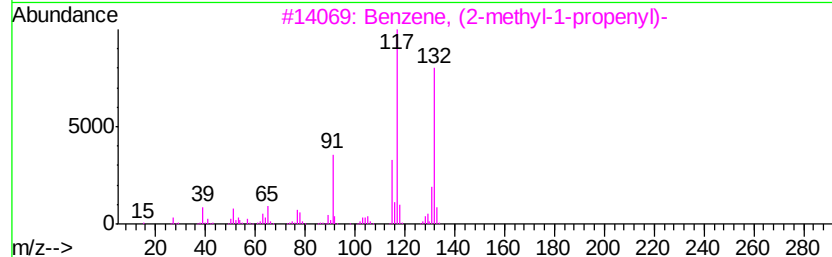
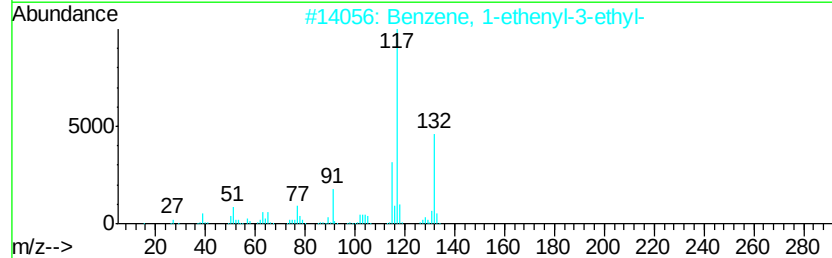
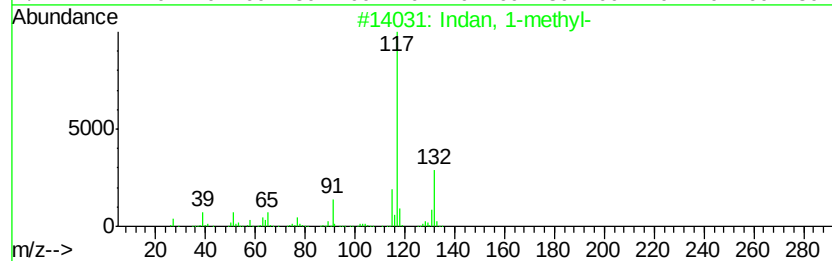
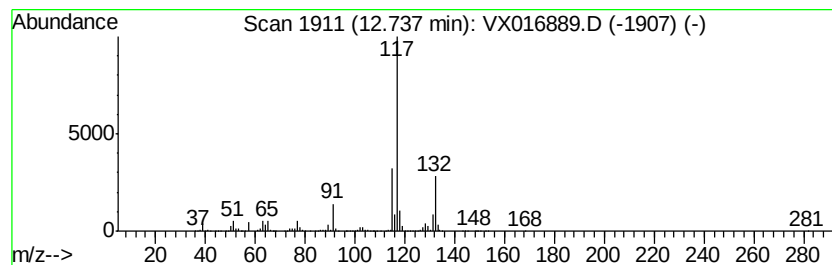
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Quant Title : SW846 8260

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

Peak Number 6 Indan, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.74	210.02 ug/l	7627350	1,4-Dichlorobenzene-d4	12.06

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indan, 1-methyl-		132	C10H12	000767-58-8	93
2	Benzene, 1-ethenyl-3-ethyl-		132	C10H12	007525-62-4	90
3	Benzene, (2-methyl-1-propenyl)-		132	C10H12	000768-49-0	90
4	Benzene, (2-methylcyclopropyl)-		132	C10H12	1000327-39-0	87
5	Benzene, (2-methyl-2-propenyl)-		132	C10H12	003290-53-7	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX061920\
Data File : VX016889.D
Acq On : 19 Jun 2020 18:53
Operator : JC/SP
Sample : L3020-23
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-27

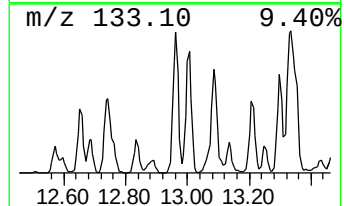
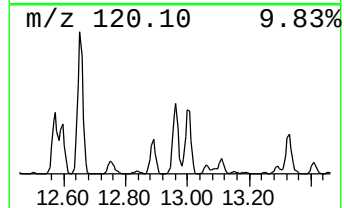
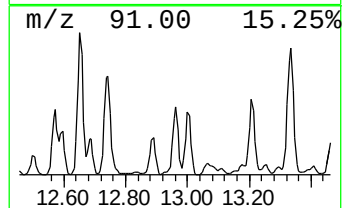
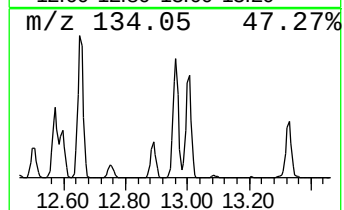
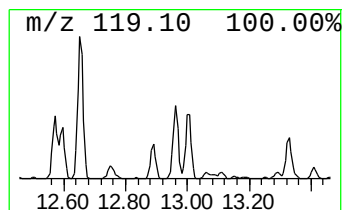
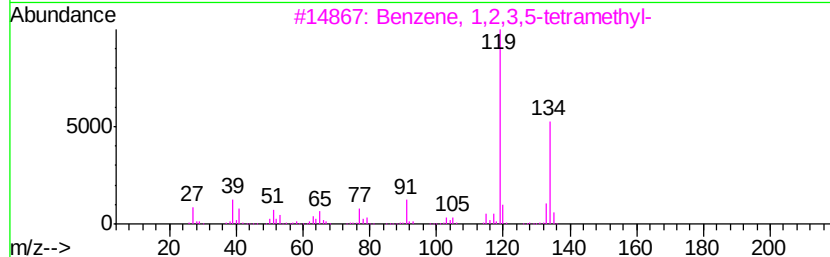
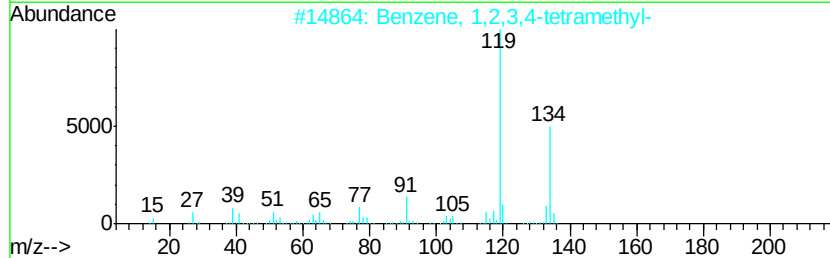
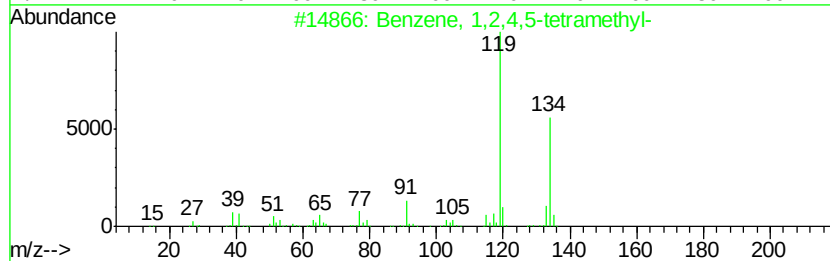
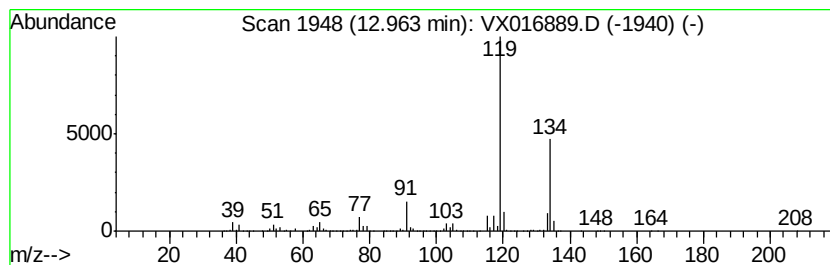
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.96	112.76 ug/l	4095050	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	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\VX061920\
Data File : VX016889.D
Acq On : 19 Jun 2020 18:53
Operator : JC/SP
Sample : L3020-23
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-27

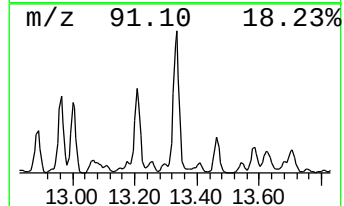
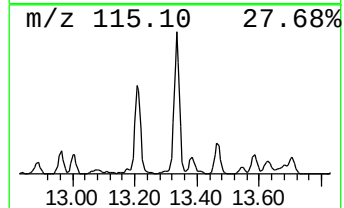
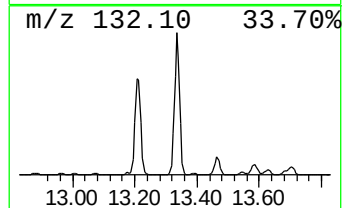
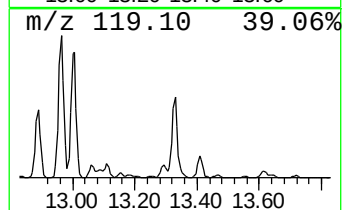
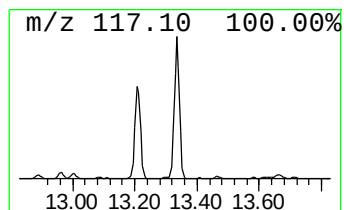
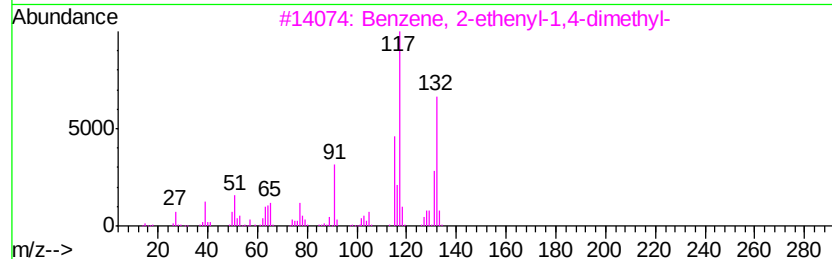
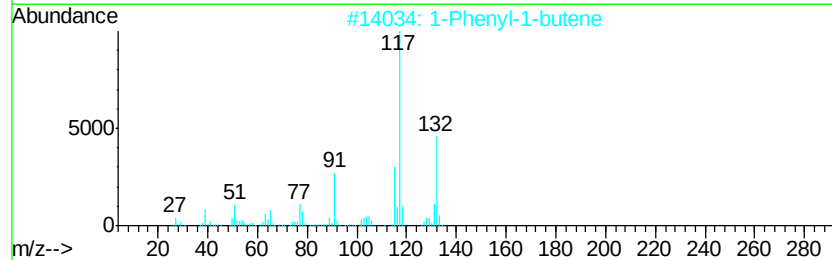
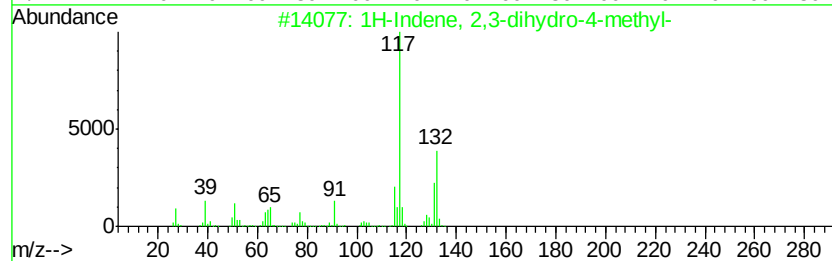
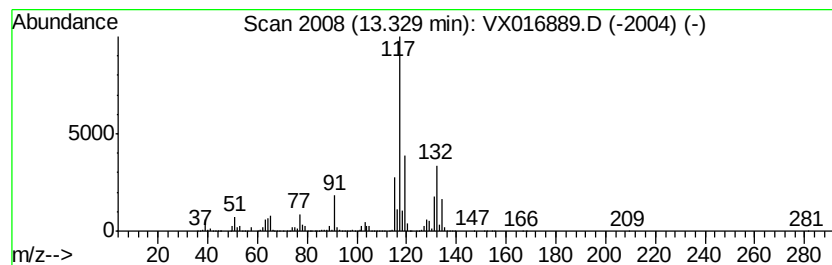
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.33	294.96 ug/l	10712100	1,4-Dichlorobenzene-d4	12.06

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90
2		1-Phenyl-1-butene	132	C10H12	000824-90-8	83
3		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	83
4		Indan, 1-methyl-	132	C10H12	000767-58-8	81
5		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	81



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX061920\
Data File : VX016889.D
Acq On : 19 Jun 2020 18:53
Operator : JC/SP
Sample : L3020-23
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-27

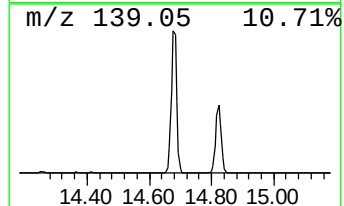
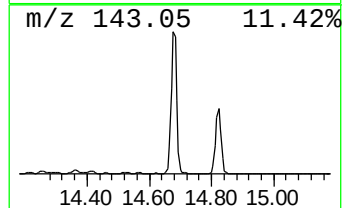
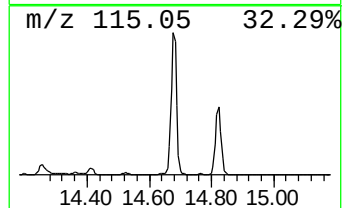
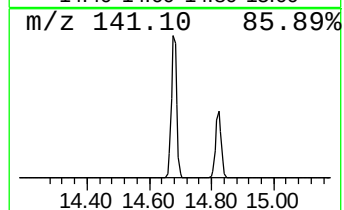
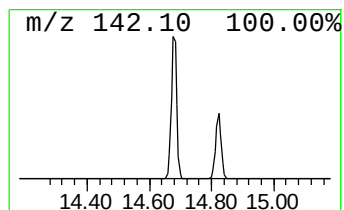
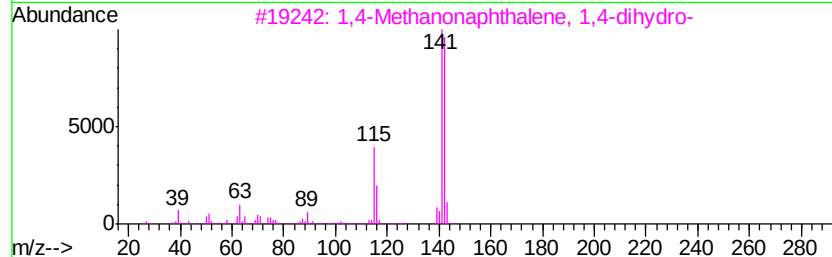
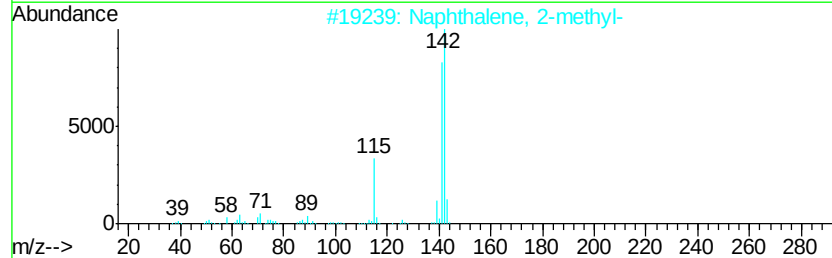
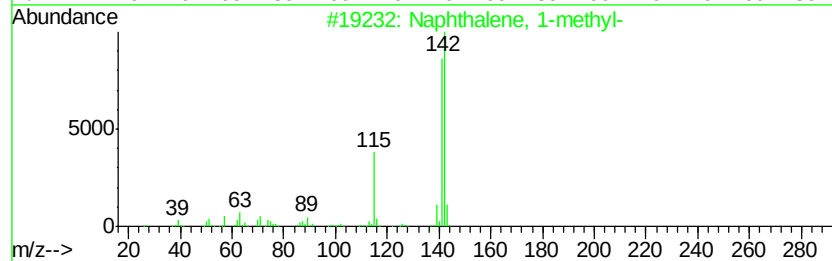
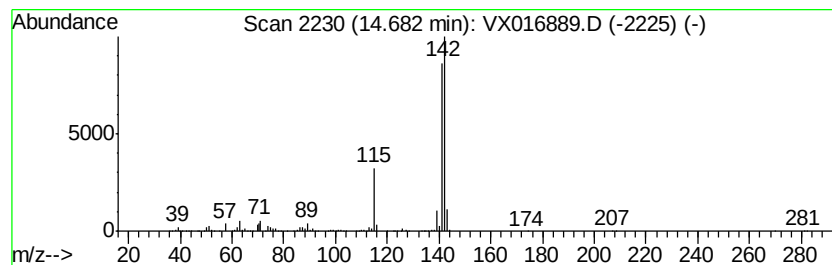
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Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.68	188.93 ug/l	6861440	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3		1,4-Methanonaphthalene, 1,4-dihv...	142	C11H10	004453-90-1	91
4		Benzocycloheptatriene	142	C11H10	000264-09-5	91
5		1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	49



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX061920\
Data File : VX016889.D
Acq On : 19 Jun 2020 18:53
Operator : JC/SP
Sample : L3020-23
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-27

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061820W.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
Benzene, 1-ethyl-...	11.44	274.9	ug/l	9983910	4	12.06	1815830	50.0
Benzene, 1-ethyl-...	12.13	596.5	ug/l	21661800	4	12.06	1815830	50.0
Indane	12.29	674.5	ug/l	24496700	4	12.06	1815830	50.0
Benzene, 2-ethyl-...	12.57	104.5	ug/l	3796480	4	12.06	1815830	50.0
o-Cymene	12.65	204.8	ug/l	7437710	4	12.06	1815830	50.0
Indan, 1-methyl-	12.74	210.0	ug/l	7627350	4	12.06	1815830	50.0
Benzene, 1,2,4,5-...	12.96	112.8	ug/l	4095050	4	12.06	1815830	50.0
1H-Indene, 2,3-di...	13.33	295.0	ug/l	10712100	4	12.06	1815830	50.0
Naphthalene, 1-me...	14.68	188.9	ug/l	6861440	4	12.06	1815830	50.0