

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX101618\
 Data File : VX005362.D
 Acq On : 16 Oct 2018 11:26
 Operator : JC/MD
 Sample : J5521-14
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TWP-04

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\82X101218W.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	1.636	67	79	84	rBV7	13969	36172	2.74%	0.230%
2	2.446	206	212	219	rVB	29217	49050	3.72%	0.312%
3	2.617	234	240	245	rBV2	11232	19973	1.51%	0.127%
4	4.397	518	532	541	rBV4	18195	49834	3.78%	0.317%
5	4.702	574	582	587	rBV	6586	17925	1.36%	0.114%
6	5.501	701	713	723	rBV2	155839	454936	34.47%	2.896%
7	5.665	730	740	757	rVB	273664	782294	59.27%	4.980%
8	6.068	794	806	814	rBV	172651	451381	34.20%	2.873%
9	6.141	814	818	829	rVB2	21597	54193	4.11%	0.345%
10	6.860	926	936	947	rBV	401410	969890	73.48%	6.174%
11	7.458	1024	1034	1045	rBV5	14922	40052	3.03%	0.255%
12	8.567	1211	1216	1225	rVB5	10176	18879	1.43%	0.120%
13	8.714	1233	1240	1249	rBV	828588	1319886	100.00%	8.401%
14	10.116	1464	1470	1484	rBV	785108	1134956	85.99%	7.224%
15	10.250	1487	1492	1499	rBV	47874	66665	5.05%	0.424%
16	10.360	1504	1510	1516	rBV	52544	76555	5.80%	0.487%
17	10.701	1561	1566	1573	rVB	24506	36781	2.79%	0.234%
18	11.018	1612	1618	1625	rBV	109346	143195	10.85%	0.911%
19	11.140	1632	1638	1651	rBV	715717	944559	71.56%	6.012%
20	11.359	1667	1674	1681	rBV	284254	355373	26.92%	2.262%
21	11.433	1681	1686	1688	rBV	148685	194084	14.70%	1.235%
22	11.506	1694	1698	1706	rVB	117869	147241	11.16%	0.937%
23	11.670	1719	1725	1736	rVV	178920	240735	18.24%	1.532%
24	11.804	1742	1747	1753	rBV	485897	613826	46.51%	3.907%
25	11.920	1763	1766	1768	rBV	20412	28009	2.12%	0.178%
26	11.945	1768	1770	1774	rVB	37251	40368	3.06%	0.257%
27	12.024	1780	1783	1786	rVB	19162	22563	1.71%	0.144%
28	12.079	1786	1792	1798	rVV	994467	1214913	92.05%	7.733%
29	12.140	1798	1802	1807	rVB	197609	248907	18.86%	1.584%
30	12.298	1821	1828	1836	rBV	933364	1204456	91.25%	7.667%
31	12.371	1836	1840	1850	rVB3	117943	201441	15.26%	1.282%
32	12.463	1850	1855	1860	rBV2	41854	56200	4.26%	0.358%
33	12.518	1860	1864	1870	rVB	42414	58142	4.41%	0.370%
34	12.585	1870	1875	1877	rBV	58596	74944	5.68%	0.477%

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Integrator: RTE
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35	12.609	1877	1879	1883	rVV	40896	48439	3.67%	0.308%
36	12.670	1884	1889	1892	rVV	382332	455830	34.54%	2.902%
37	12.701	1892	1894	1898	rVV	65681	70125	5.31%	0.446%
38	12.756	1898	1903	1911	rVV2	218047	318073	24.10%	2.025%
39	12.908	1923	1928	1932	rBV2	49674	72688	5.51%	0.463%
40	12.975	1935	1939	1943	rVV	203370	254695	19.30%	1.621%
41	13.018	1943	1946	1952	rVB	62035	75315	5.71%	0.479%
42	13.079	1952	1956	1961	rBV3	18896	40001	3.03%	0.255%
43	13.225	1976	1980	1984	rVB	187219	216379	16.39%	1.377%
44	13.347	1995	2000	2005	rBV2	444687	621175	47.06%	3.954%
45	13.402	2006	2009	2012	rVV3	23529	38027	2.88%	0.242%
46	13.481	2019	2022	2031	rVV5	38727	58794	4.45%	0.374%
47	13.560	2032	2035	2038	rVV3	14614	19830	1.50%	0.126%
48	13.603	2038	2042	2045	rVV	33784	45959	3.48%	0.293%
49	13.640	2045	2048	2053	rVB4	39511	54635	4.14%	0.348%
50	13.719	2058	2061	2066	rVB2	35108	57483	4.36%	0.366%
51	13.835	2076	2080	2086	rBV	139639	177514	13.45%	1.130%
52	13.896	2087	2090	2098	rVB4	40915	66095	5.01%	0.421%
53	13.981	2099	2104	2108	rVB3	25392	41699	3.16%	0.265%
54	14.030	2108	2112	2116	rBV	16743	23462	1.78%	0.149%
55	14.133	2125	2129	2133	rBV	39248	55551	4.21%	0.354%
56	14.274	2148	2152	2158	rVB2	38597	64525	4.89%	0.411%
57	14.377	2166	2169	2173	rVB	17297	22305	1.69%	0.142%
58	14.432	2174	2178	2183	rVB	28426	38554	2.92%	0.245%
59	14.542	2192	2196	2199	rBV2	17950	24955	1.89%	0.159%
60	14.670	2212	2217	2218	rBV2	51526	61245	4.64%	0.390%
61	14.694	2218	2221	2231	rVB	190697	259918	19.69%	1.654%
62	14.786	2232	2236	2240	rBV6	12541	21885	1.66%	0.139%
63	14.841	2240	2245	2251	rBV	225516	284678	21.57%	1.812%
64	15.273	2313	2316	2320	rVB	43862	56794	4.30%	0.362%
65	15.444	2340	2344	2347	rBV2	26442	36061	2.73%	0.230%
66	15.487	2347	2351	2356	rVV4	18127	33828	2.56%	0.215%
67	15.548	2356	2361	2374	rVB	67910	121818	9.23%	0.775%
68	15.688	2379	2384	2387	rVV	99011	159357	12.07%	1.014%
69	15.725	2387	2390	2399	rVB	56998	97952	7.42%	0.623%
70	15.816	2403	2405	2412	rVV7	10153	15075	1.14%	0.096%
71	15.914	2413	2421	2431	rVB3	26530	74281	5.63%	0.473%

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72	16.072	2443	2447	2457	rVB	21154	44594	3.38%	0.284%
73	16.170	2458	2463	2470	rBV3	22240	37294	2.83%	0.237%
74	16.688	2544	2548	2553	rVB2	19494	36682	2.78%	0.233%
75	16.755	2554	2559	2564	rBV3	17323	38194	2.89%	0.243%

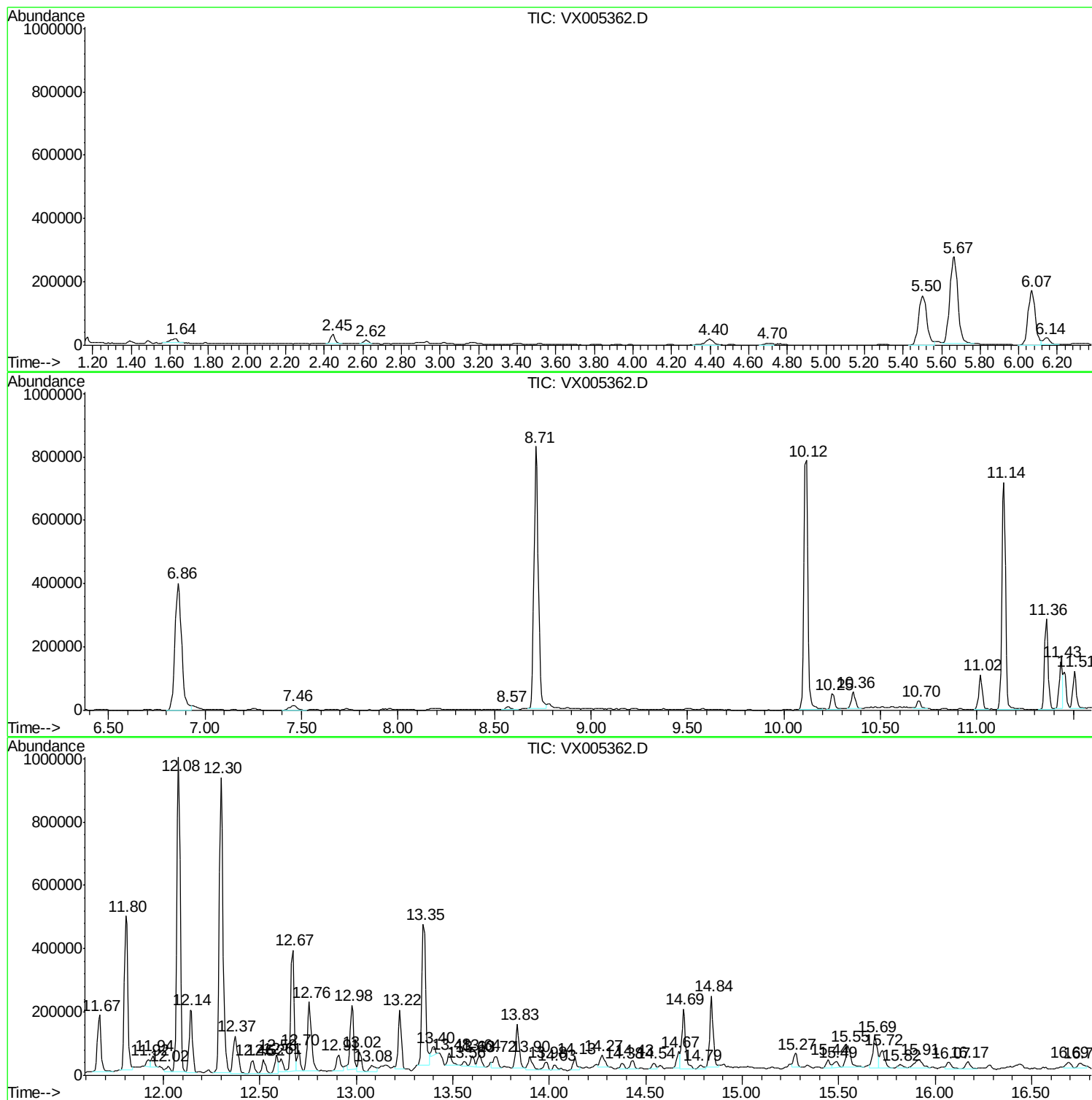
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X101218W.M
Quant Title : SW846 8260

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



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Instrument :
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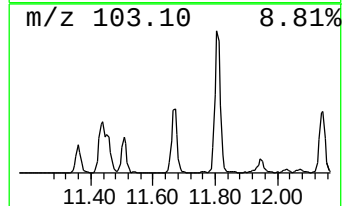
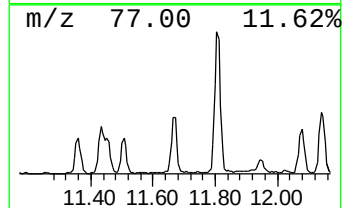
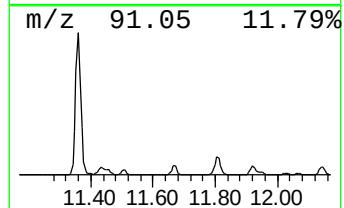
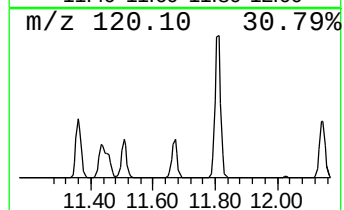
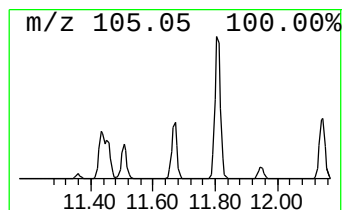
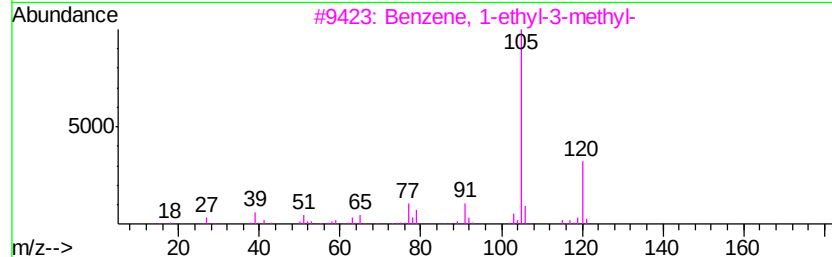
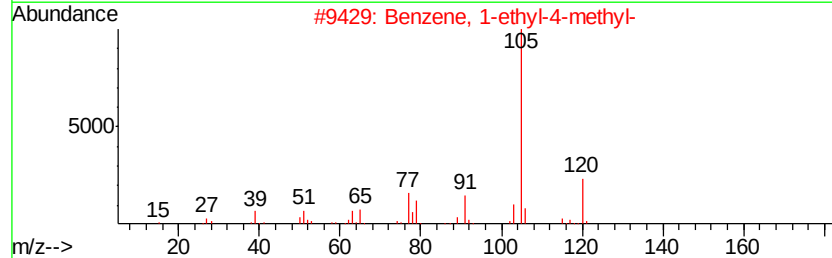
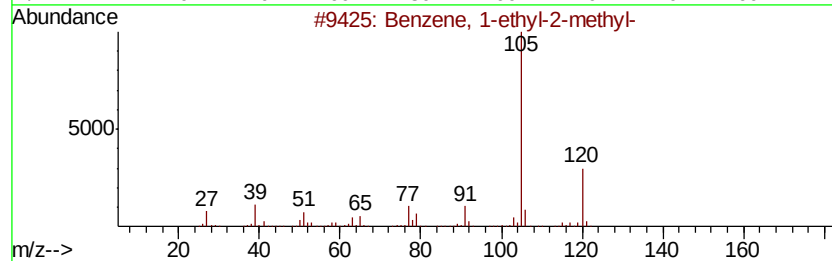
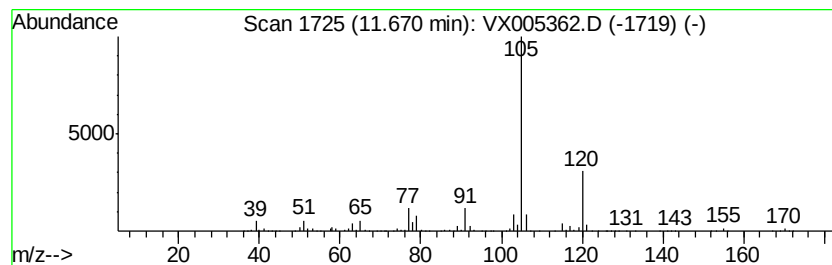
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X101218W.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1-ethyl-2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.67	9.91 ug/l	240735	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	90
4		Mesitylene	120	C9H12	000108-67-8	90
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	90



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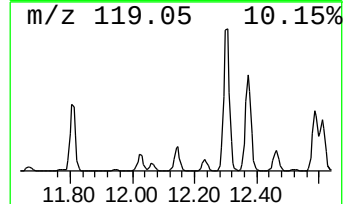
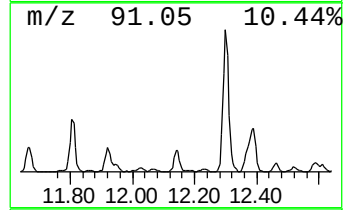
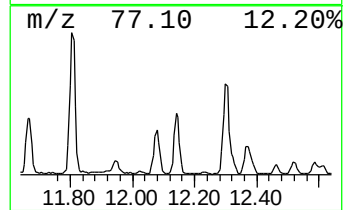
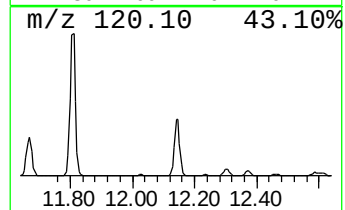
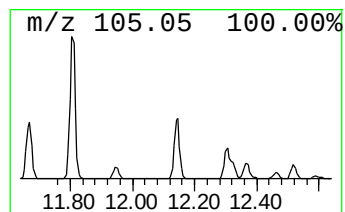
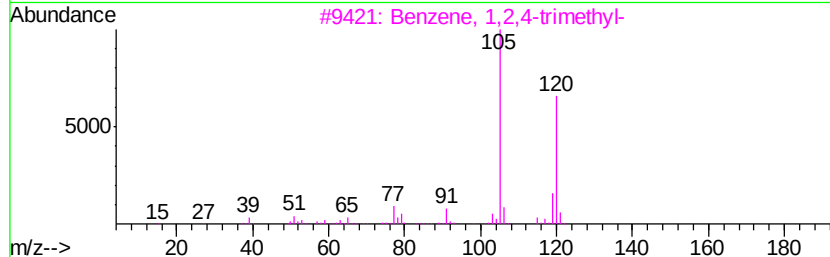
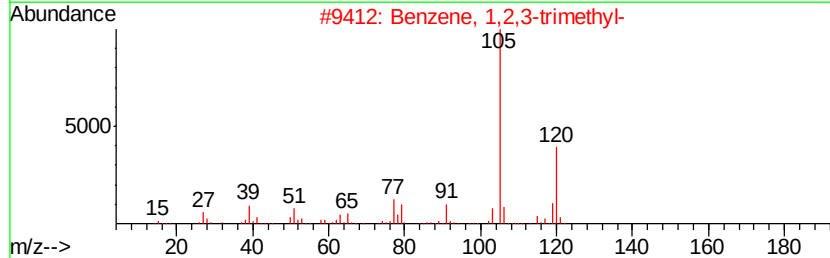
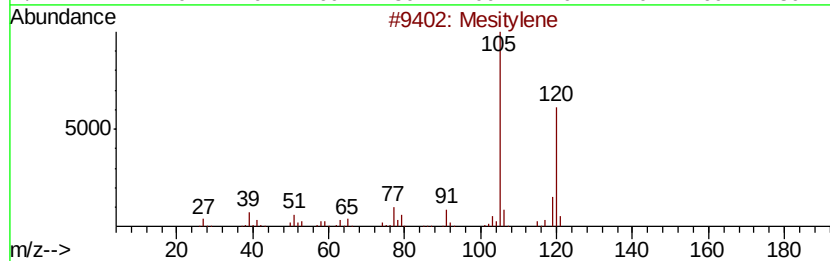
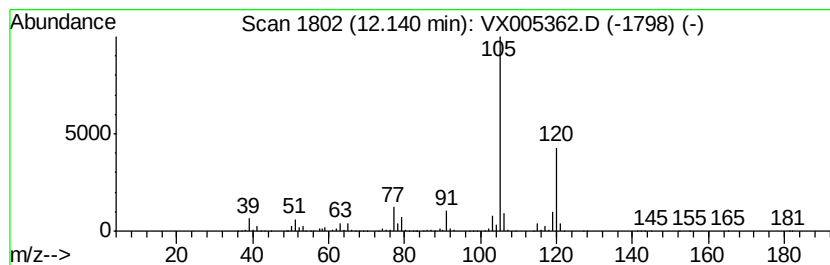
Instrument :
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ClientSampled :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X101218W.M
Quant Title : SW846 8260

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

Peak Number 2 Mesitylene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.14	10.24 ug/l	248907	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Mesitylene	120 C9H12	000108-67-8	97
2	Benzene, 1,2,3-trimethyl-	120 C9H12	000526-73-8	95
3	Benzene, 1,2,4-trimethyl-	120 C9H12	000095-63-6	94
4	Benzene, 1-ethyl-3-methyl-	120 C9H12	000620-14-4	91
5	Benzene, 1-ethyl-2-methyl-	120 C9H12	000611-14-3	90



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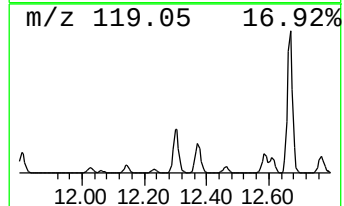
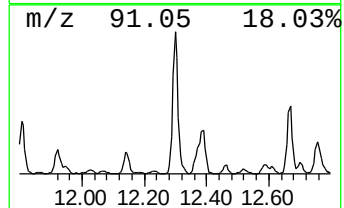
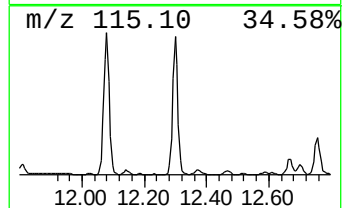
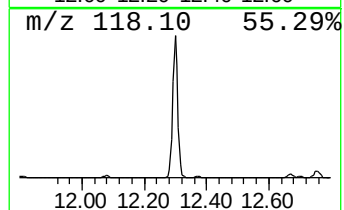
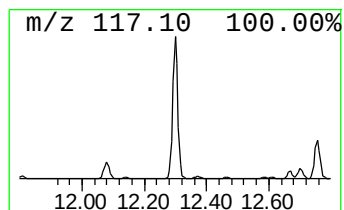
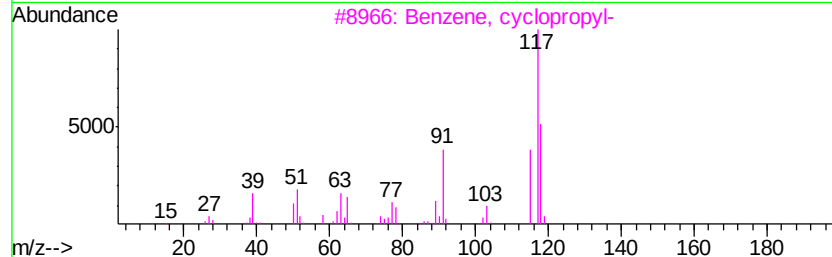
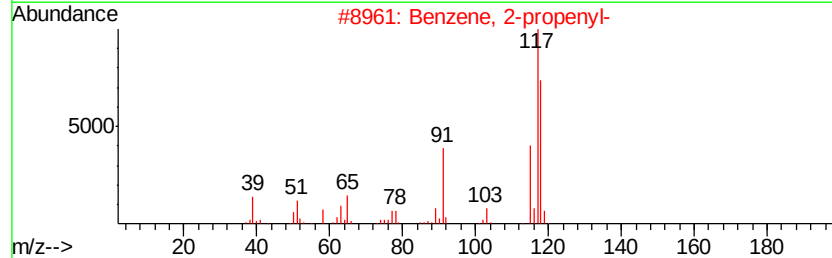
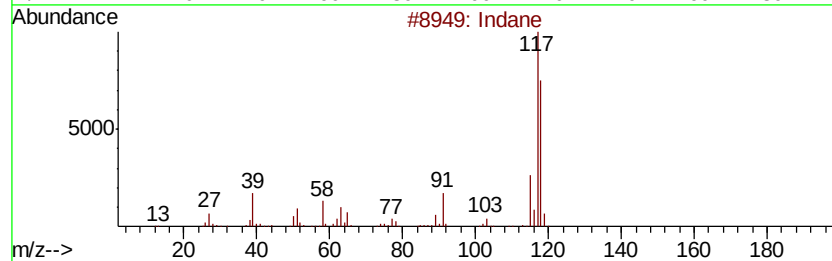
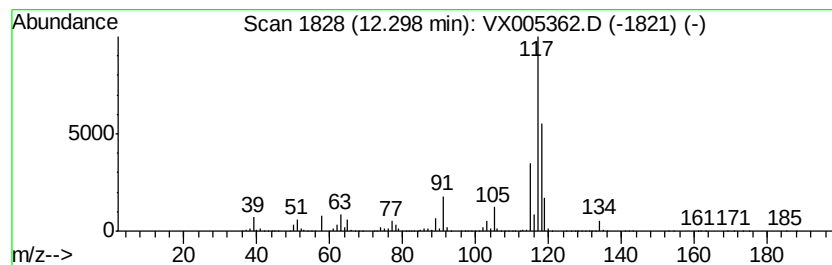
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	49.57 ug/l	1204460	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 76
3	Benzene, cyclopropyl-		118 C9H10	000873-49-4 76
4	Benzene, 1-propenyl-		118 C9H10	000637-50-3 68
5	Deltacyclene		118 C9H10	007785-10-6 58



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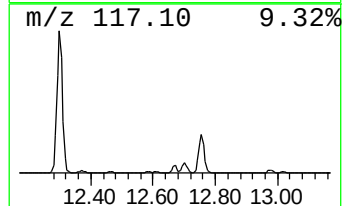
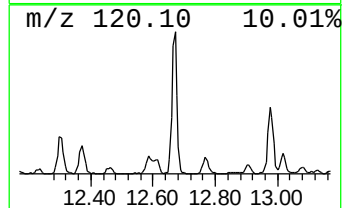
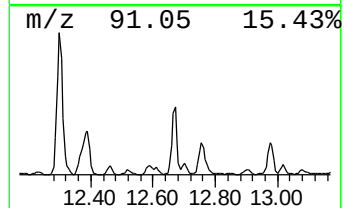
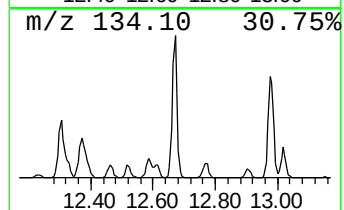
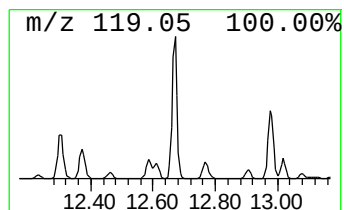
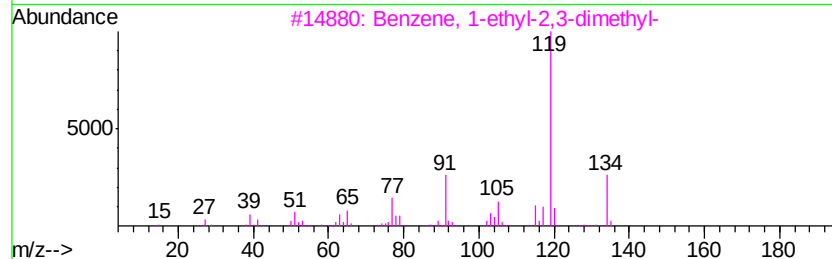
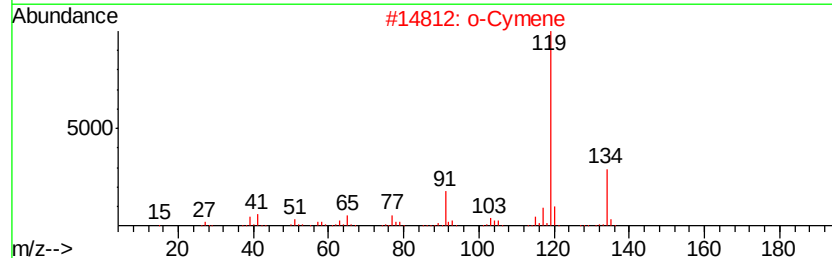
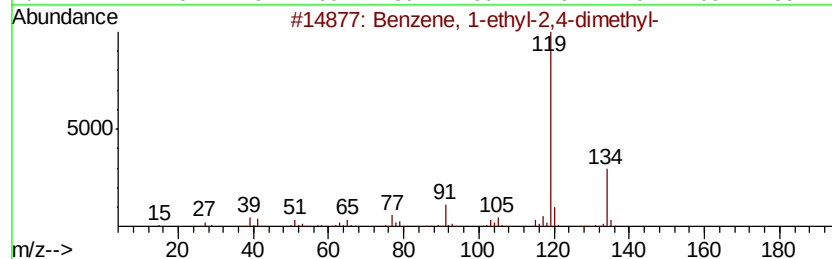
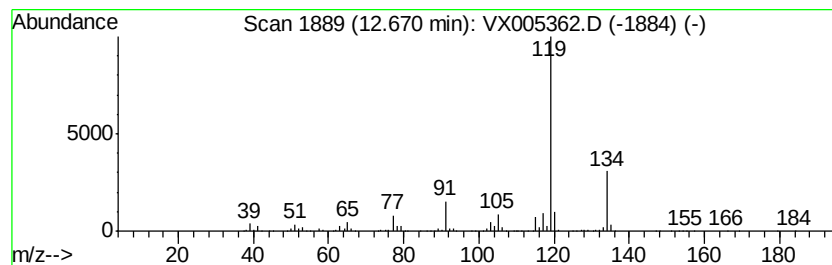
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.67	18.76 ug/l	455830	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	97
2		o-Cymene	134	C10H14	000527-84-4	97
3		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
4		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
5		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX101618\
Data File : VX005362.D
Acq On : 16 Oct 2018 11:26
Operator : JC/MD
Sample : J5521-14
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 9 Sample Multiplier: 1

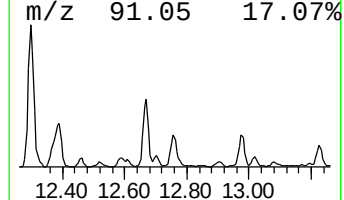
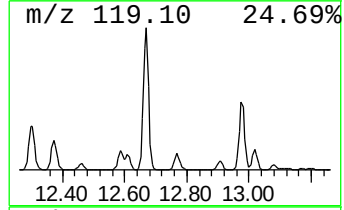
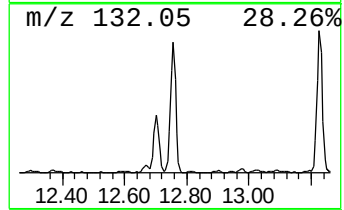
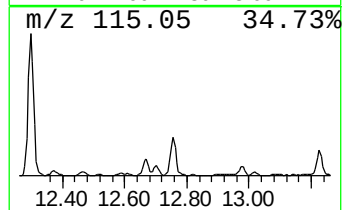
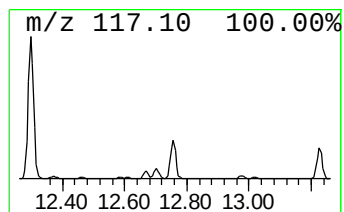
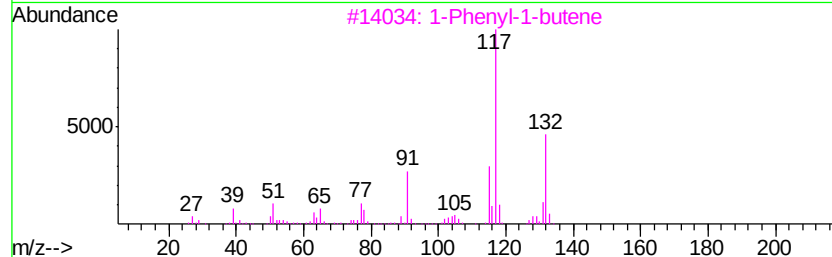
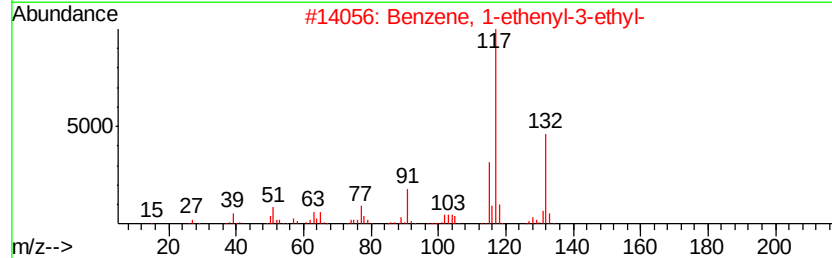
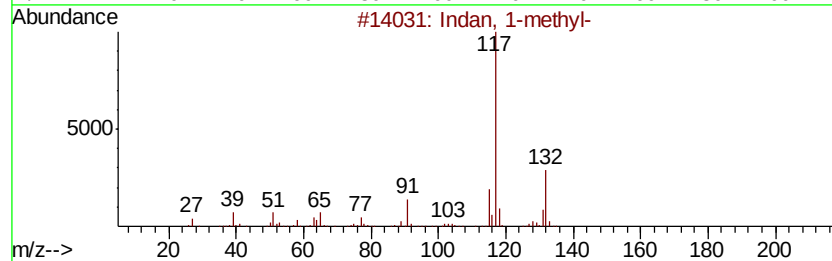
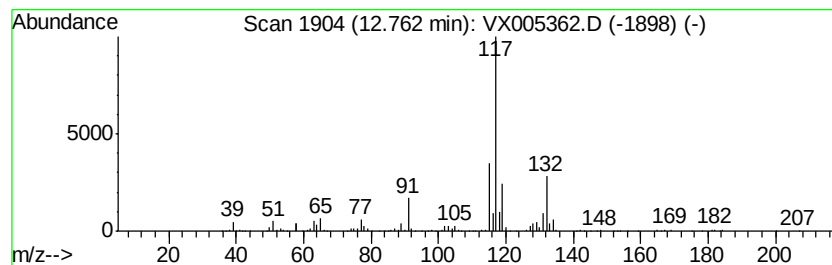
Instrument :
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ClientSampleId :
TWP-04

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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 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	13.09 ug/l	318073	1,4-Dichlorobenzene-d4	12.08
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	87
3	1-Phenyl-1-butene	132 C10H12	000824-90-8	83
4	Benzene, (2-methylcyclopropyl)-	132 C10H12	1000327-39-0	80
5	Benzene, 2-butenyl-	132 C10H12	001560-06-1	80



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX101618\
Data File : VX005362.D
Acq On : 16 Oct 2018 11:26
Operator : JC/MD
Sample : J5521-14
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 9 Sample Multiplier: 1

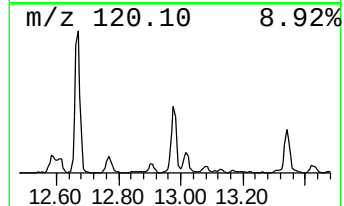
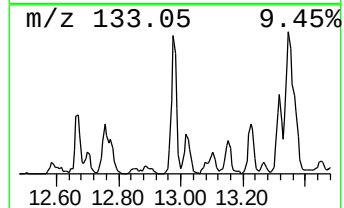
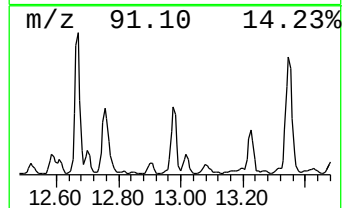
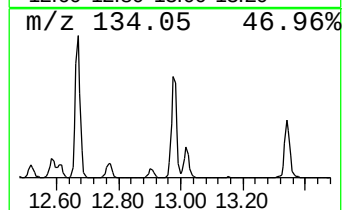
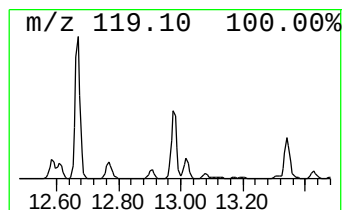
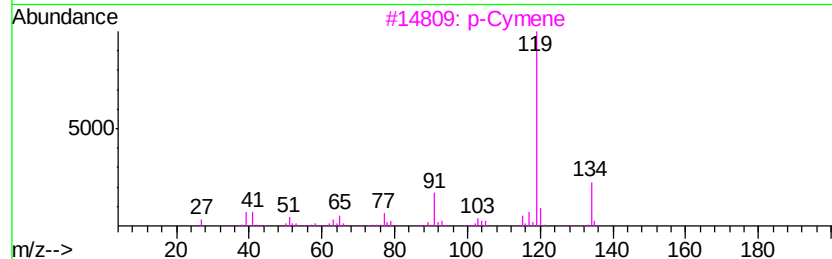
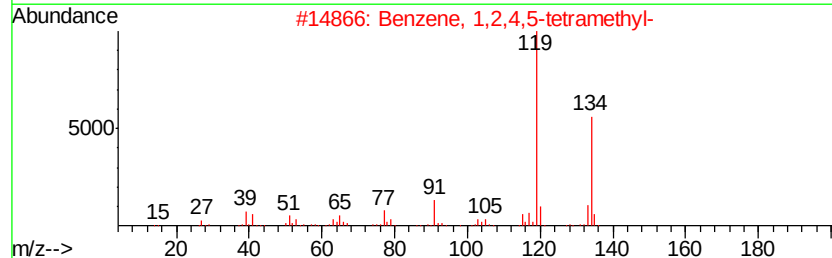
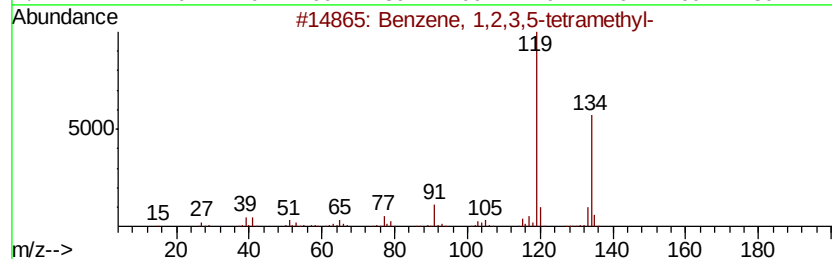
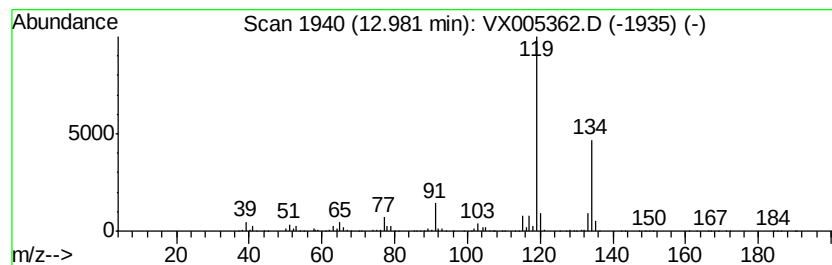
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ClientSampleId :
TWP-04

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X101218W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.98	10.48 ug/l	254695	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1,2,3,5-tetramethyl-	134 C10H14	000527-53-7 97
2		Benzene, 1,2,4,5-tetramethyl-	134 C10H14	000095-93-2 97
3		p-Cymene	134 C10H14	000099-87-6 93
4		o-Cymene	134 C10H14	000527-84-4 93
5		Benzene, 1,2,3,4-tetramethyl-	134 C10H14	000488-23-3 93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX101618\
Data File : VX005362.D
Acq On : 16 Oct 2018 11:26
Operator : JC/MD
Sample : J5521-14
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TWP-04

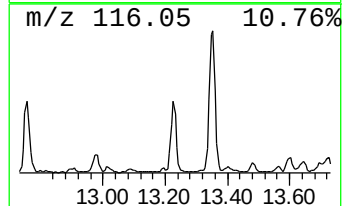
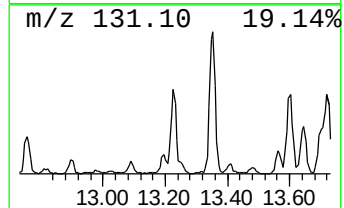
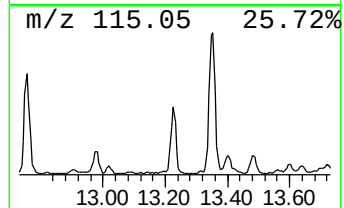
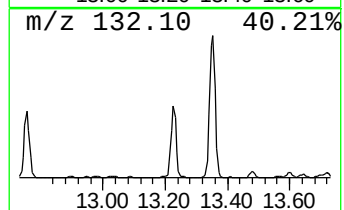
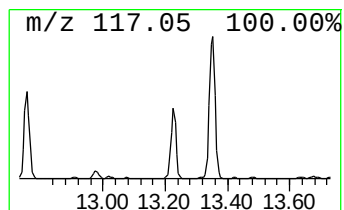
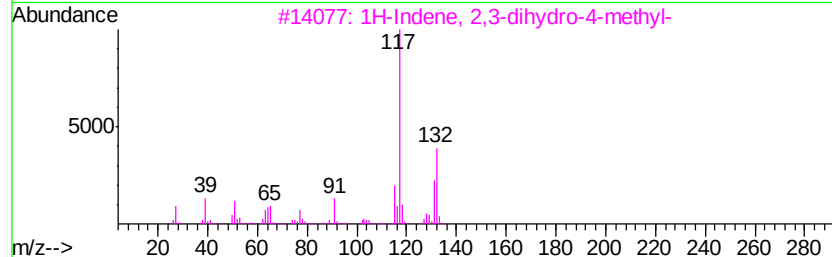
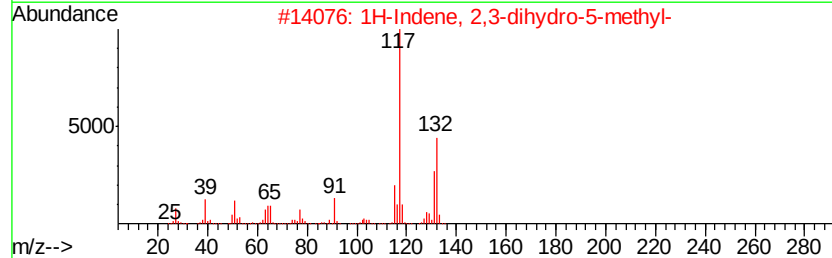
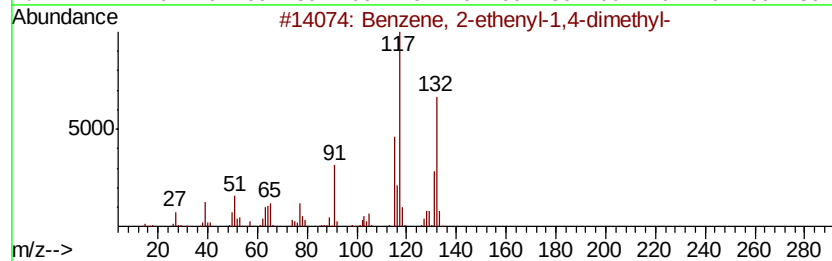
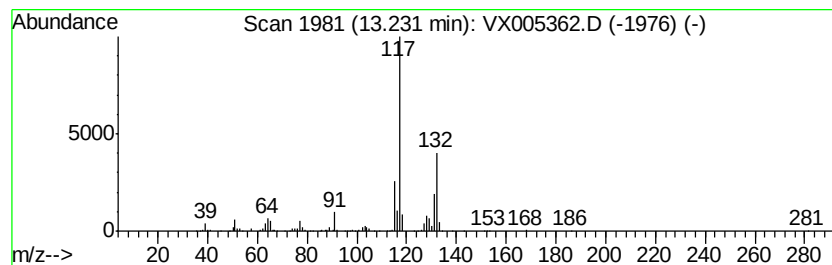
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.23	8.91 ug/l	216379	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		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	90
3		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90
4		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	90
5		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX101618\
Data File : VX005362.D
Acq On : 16 Oct 2018 11:26
Operator : JC/MD
Sample : J5521-14
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TWP-04

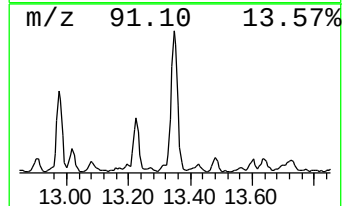
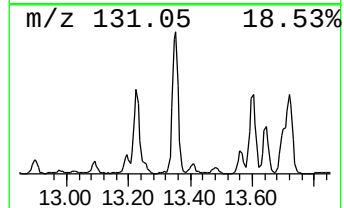
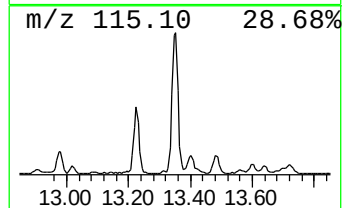
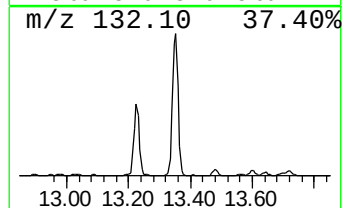
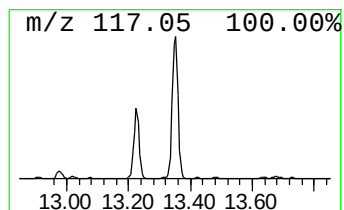
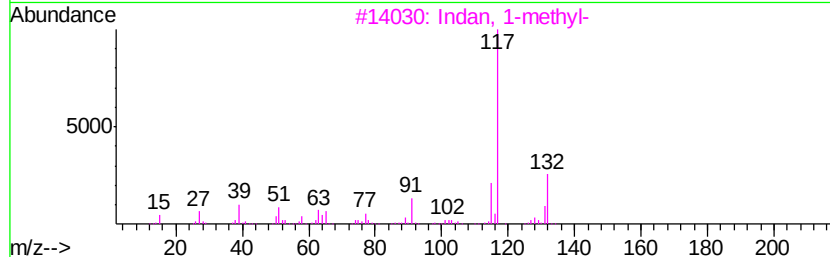
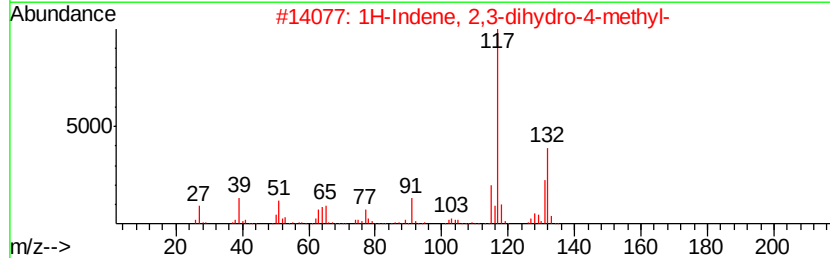
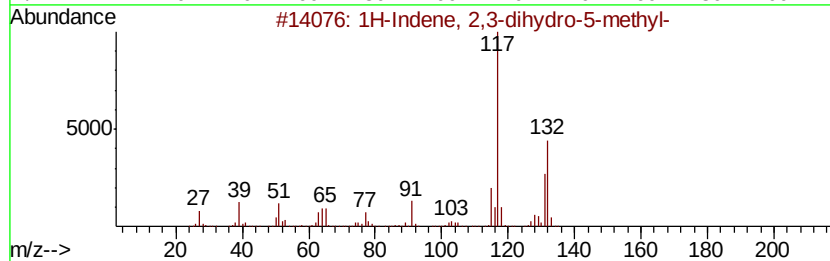
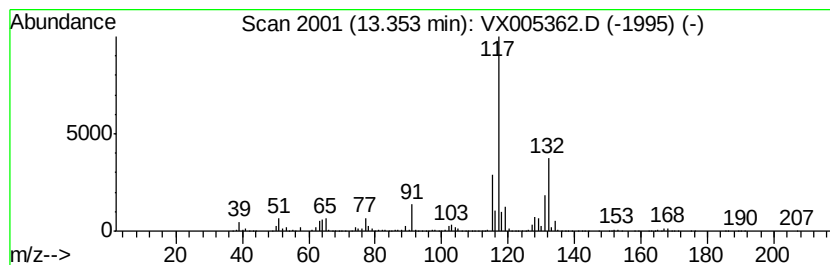
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Quant Title : SW846 8260

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	25.56 ug/l	621175	1,4-Dichlorobenzene-d4	12.08

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	89
2			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	76
3			Indan, 1-methyl-	132	C10H12	000767-58-8	76
4			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	74
5			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	72



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX101618\
Data File : VX005362.D
Acq On : 16 Oct 2018 11:26
Operator : JC/MD
Sample : J5521-14
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 9 Sample Multiplier: 1

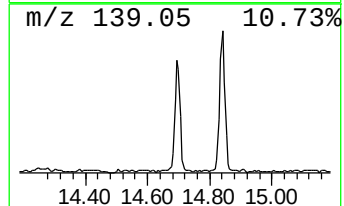
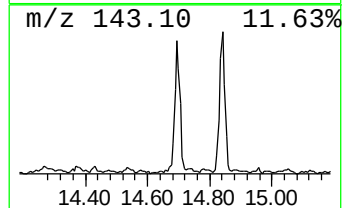
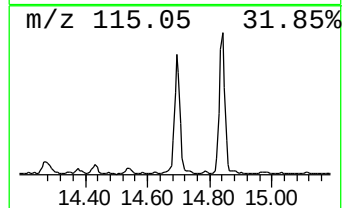
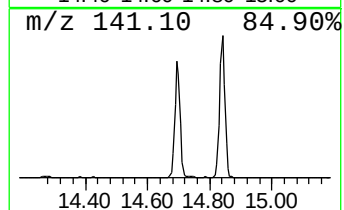
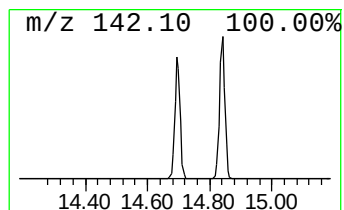
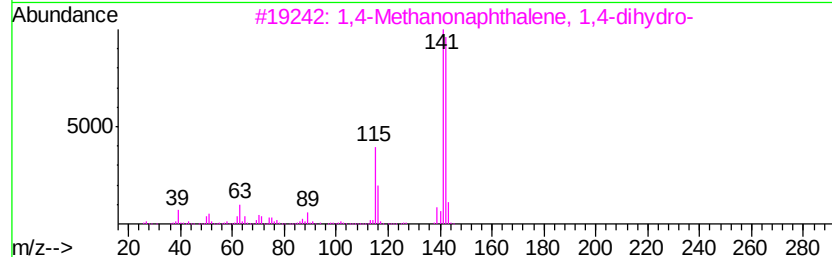
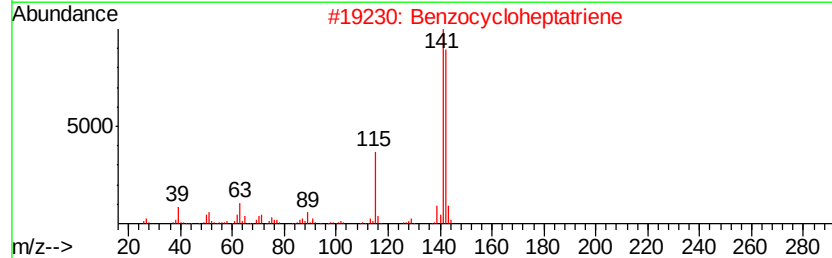
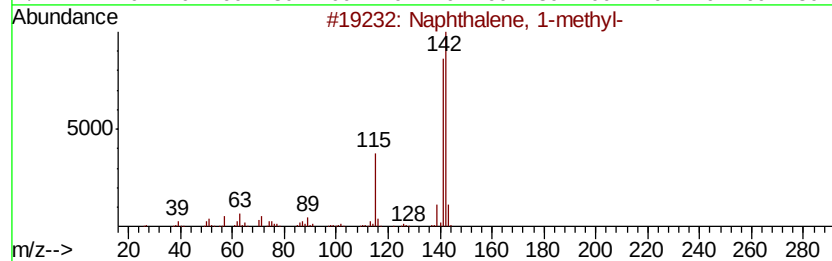
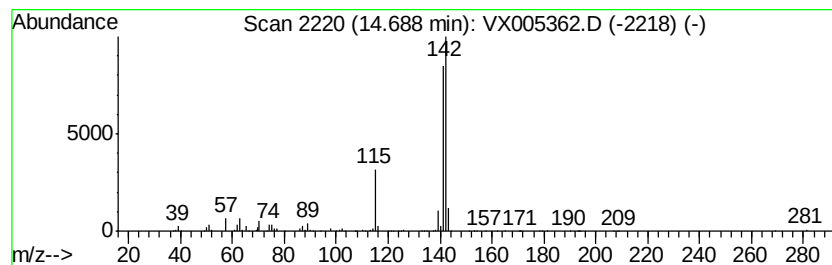
Instrument :
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ClientSampleId :
TWP-04

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X101218W.M
Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.69	10.70 ug/l	259918	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1-methyl-	142 C11H10	000090-12-0 94
2		Benzocycloheptatriene	142 C11H10	000264-09-5 91
3		1,4-Methanonaphthalene, 1,4-dihydro-	142 C11H10	004453-90-1 91
4		Naphthalene, 2-methyl-	142 C11H10	000091-57-6 90
5		Naphthalene, 1-(2-hydroxypropyl)	186 C13H14O	027653-13-0 83



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX101618\
Data File : VX005362.D
Acq On : 16 Oct 2018 11:26
Operator : JC/MD
Sample : J5521-14
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
TWP-04

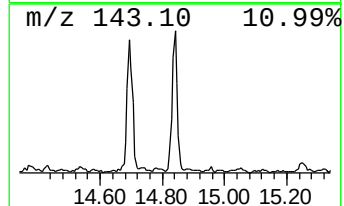
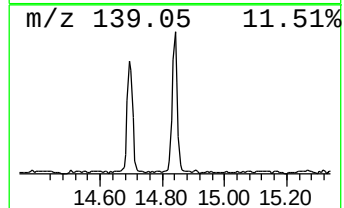
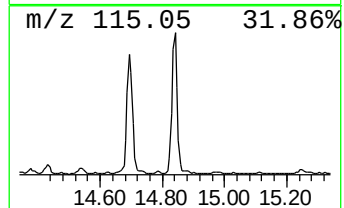
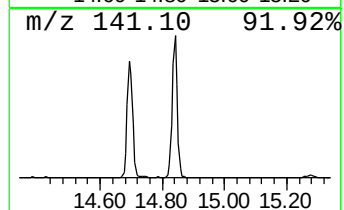
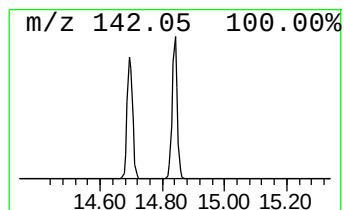
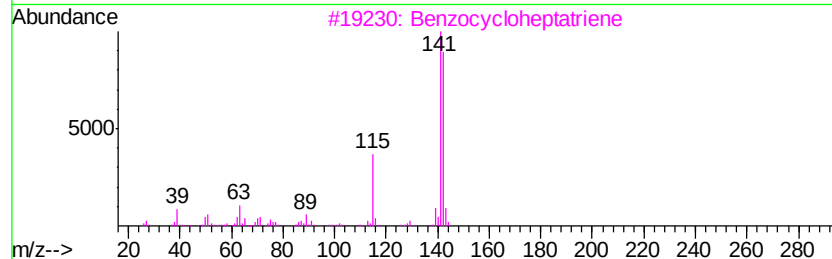
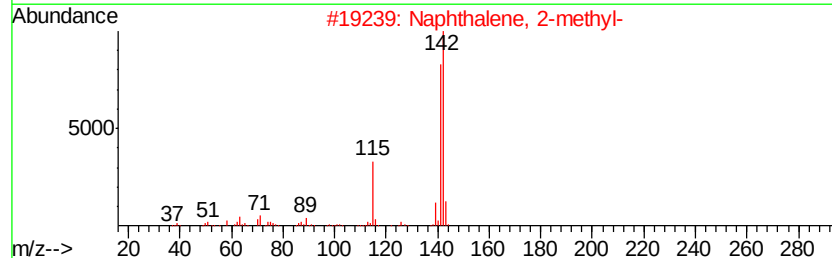
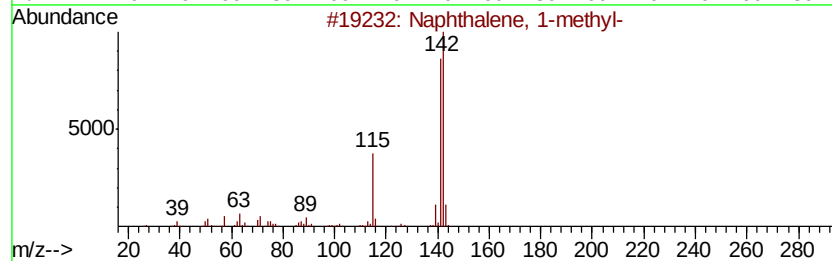
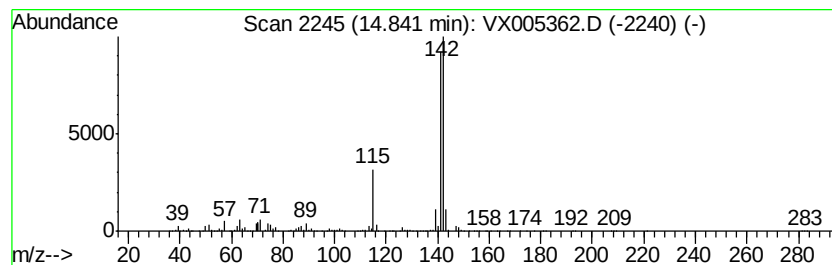
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Quant Title : SW846 8260

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

Peak Number 10 Benzocycloheptatriene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.84	11.72 ug/l	284678	1,4-Dichlorobenzene-d4	12.08

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		Benzocycloheptatriene	142	C11H10	000264-09-5	90
4		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	90
5		1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	59



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX101618\
Data File : VX005362.D
Acq On : 16 Oct 2018 11:26
Operator : JC/MD
Sample : J5521-14
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TWP-04

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X101218W.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-...	11.67	9.9	ug/l	240735	4	12.08	1214910	50.0
Mesitylene	12.14	10.2	ug/l	248907	4	12.08	1214910	50.0
Indane	12.30	49.6	ug/l	1204460	4	12.08	1214910	50.0
Benzene, 1-ethyl-...	12.67	18.8	ug/l	455830	4	12.08	1214910	50.0
Indan, 1-methyl-	12.76	13.1	ug/l	318073	4	12.08	1214910	50.0
Benzene, 1,2,3,5-...	12.98	10.5	ug/l	254695	4	12.08	1214910	50.0
Benzene, 2-etheny...	13.23	8.9	ug/l	216379	4	12.08	1214910	50.0
1H-Indene, 2,3-di...	13.35	25.6	ug/l	621175	4	12.08	1214910	50.0
Naphthalene, 1-me...	14.69	10.7	ug/l	259918	4	12.08	1214910	50.0
Benzocycloheptatr...	14.84	11.7	ug/l	284678	4	12.08	1214910	50.0