

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX122018\
 Data File : VX006702.D
 Acq On : 21 Dec 2018 03:48
 Operator : JC/MD
 Sample : J6517-01
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 T3-MW-2-9.0-121918

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\82X121818W.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	2.891	280	285	305	rVB	131005	313159	1.24%	0.163%
2	3.172	323	331	346	rVB2	129875	339595	1.34%	0.177%
3	4.403	523	533	543	rVB	245258	706155	2.79%	0.368%
4	5.507	703	714	721	rBV	343825	998236	3.95%	0.520%
5	5.665	732	740	767	rVB	623504	2254743	8.91%	1.174%
6	5.933	773	784	796	rVB5	99568	311470	1.23%	0.162%
7	6.067	796	806	816	rBV	356381	928374	3.67%	0.483%
8	6.256	823	837	844	rBV3	112748	349665	1.38%	0.182%
9	6.354	845	853	861	rVV3	191457	558061	2.21%	0.291%
10	6.458	861	870	881	rVB2	124309	354992	1.40%	0.185%
11	6.860	925	936	943	rBV	926322	2277428	9.00%	1.186%
12	6.939	944	949	961	rVB4	247014	702705	2.78%	0.366%
13	7.256	989	1001	1010	rVB	205303	466226	1.84%	0.243%
14	7.457	1022	1034	1046	rBV4	495836	1429974	5.65%	0.744%
15	7.732	1073	1079	1089	rVB	150886	293024	1.16%	0.153%
16	7.957	1105	1116	1123	rBV5	89762	309192	1.22%	0.161%
17	8.055	1123	1132	1143	rVB3	135461	359798	1.42%	0.187%
18	8.177	1143	1152	1154	rBV	223818	437901	1.73%	0.228%
19	8.213	1154	1158	1163	rVV	565400	987832	3.90%	0.514%
20	8.573	1211	1217	1227	rVB2	302440	551172	2.18%	0.287%
21	8.713	1235	1240	1248	rVB	1938517	3026989	11.96%	1.576%
22	8.988	1280	1285	1289	rBV	162137	253739	1.00%	0.132%
23	9.164	1305	1314	1319	rBV2	141646	288782	1.14%	0.150%
24	9.219	1319	1323	1334	rVB	481816	756215	2.99%	0.394%
25	9.561	1375	1379	1385	rVB4	129277	259395	1.03%	0.135%
26	10.115	1464	1470	1474	rBV	1860384	2551956	10.09%	1.329%
27	11.018	1613	1618	1627	rBV	2331596	2878366	11.38%	1.498%
28	11.140	1633	1638	1642	rBV	1611457	2049154	8.10%	1.067%
29	11.359	1669	1674	1684	rVB	8772440	10357060	40.94%	5.392%
30	11.670	1719	1725	1733	rBV	900971	1171592	4.63%	0.610%
31	11.920	1760	1766	1768	rBV	436681	612082	2.42%	0.319%
32	11.944	1768	1770	1777	rVB	766987	872141	3.45%	0.454%
33	12.079	1787	1792	1798	rBV	2160170	2617946	10.35%	1.363%
34	12.231	1813	1817	1823	rVB	551255	655289	2.59%	0.341%

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35	12.298	1823	1828	1835	rBV	10378995	13146891	51.96%	6.844%
36	12.371	1835	1840	1847	rBV2	2383478	3959130	15.65%	2.061%
37	12.463	1850	1855	1860	rVB	1412075	1767633	6.99%	0.920%
38	12.524	1860	1865	1871	rBV3	373288	596378	2.36%	0.310%
39	12.585	1871	1875	1882	rBV	1894270	2363005	9.34%	1.230%
40	12.670	1882	1889	1892	rBV	14914238	16930130	66.92%	8.814%
41	12.700	1892	1894	1899	rVV	2915955	3290842	13.01%	1.713%
42	12.755	1899	1903	1911	rVB2	9610277	13789210	54.50%	7.178%
43	12.895	1922	1926	1932	rVB2	628470	889292	3.51%	0.463%
44	12.981	1932	1940	1943	rBV	8607058	10418710	41.18%	5.424%
45	13.017	1943	1946	1952	rVB	5283553	6327259	25.01%	3.294%
46	13.084	1952	1957	1963	rBV3	566521	1312278	5.19%	0.683%
47	13.194	1972	1975	1977	rBV2	513450	647673	2.56%	0.337%
48	13.225	1977	1980	1990	rVB	7466939	9530058	37.67%	4.961%
49	13.316	1990	1995	1996	rBV	622722	877366	3.47%	0.457%
50	13.347	1996	2000	2006	rVV2	17637329	25300583	100.00%	13.171%
51	13.401	2006	2009	2011	rVV2	487730	609871	2.41%	0.317%
52	13.426	2011	2013	2018	rVB	396067	440602	1.74%	0.229%
53	13.481	2018	2022	2028	rVB	726960	932979	3.69%	0.486%
54	13.560	2028	2035	2038	rBV2	675312	963244	3.81%	0.501%
55	13.603	2038	2042	2045	rVV	1798313	2474699	9.78%	1.288%
56	13.639	2045	2048	2054	rVV2	1910960	3608947	14.26%	1.879%
57	13.700	2054	2058	2059	rVV	1021971	1331907	5.26%	0.693%
58	13.725	2059	2062	2068	rVB2	1804722	2681197	10.60%	1.396%
59	13.853	2079	2083	2088	rVB2	342830	426996	1.69%	0.222%
60	13.914	2088	2093	2097	rVB	423736	530838	2.10%	0.276%
61	13.987	2097	2105	2109	rBV2	967965	1413482	5.59%	0.736%
62	14.029	2109	2112	2116	rVB	755345	857874	3.39%	0.447%
63	14.133	2124	2129	2133	rBV	1220798	1520864	6.01%	0.792%
64	14.273	2148	2152	2158	rBV2	1322001	2470544	9.76%	1.286%
65	14.377	2165	2169	2175	rBV	699121	1054335	4.17%	0.549%
66	14.432	2175	2178	2182	rVB	795594	924048	3.65%	0.481%
67	14.541	2190	2196	2200	rBV2	815821	1104696	4.37%	0.575%
68	14.657	2209	2215	2217	rBV	248050	360970	1.43%	0.188%
69	14.694	2217	2221	2226	rVB	2686074	3509049	13.87%	1.827%
70	14.840	2240	2245	2254	rBV	6843078	8532808	33.73%	4.442%
71	15.553	2355	2362	2366	rBV	271927	446217	1.76%	0.232%

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Integration Parameters: RTEINT.P

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Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
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Stop Thrs : 0 Peak Location: TOP

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72	15.688	2376	2384	2388	rBV	412670	729612	2.88%	0.380%
73	15.730	2388	2391	2396	rVB	280784	396714	1.57%	0.207%
74	15.919	2412	2422	2433	rVB2	120386	342477	1.35%	0.178%

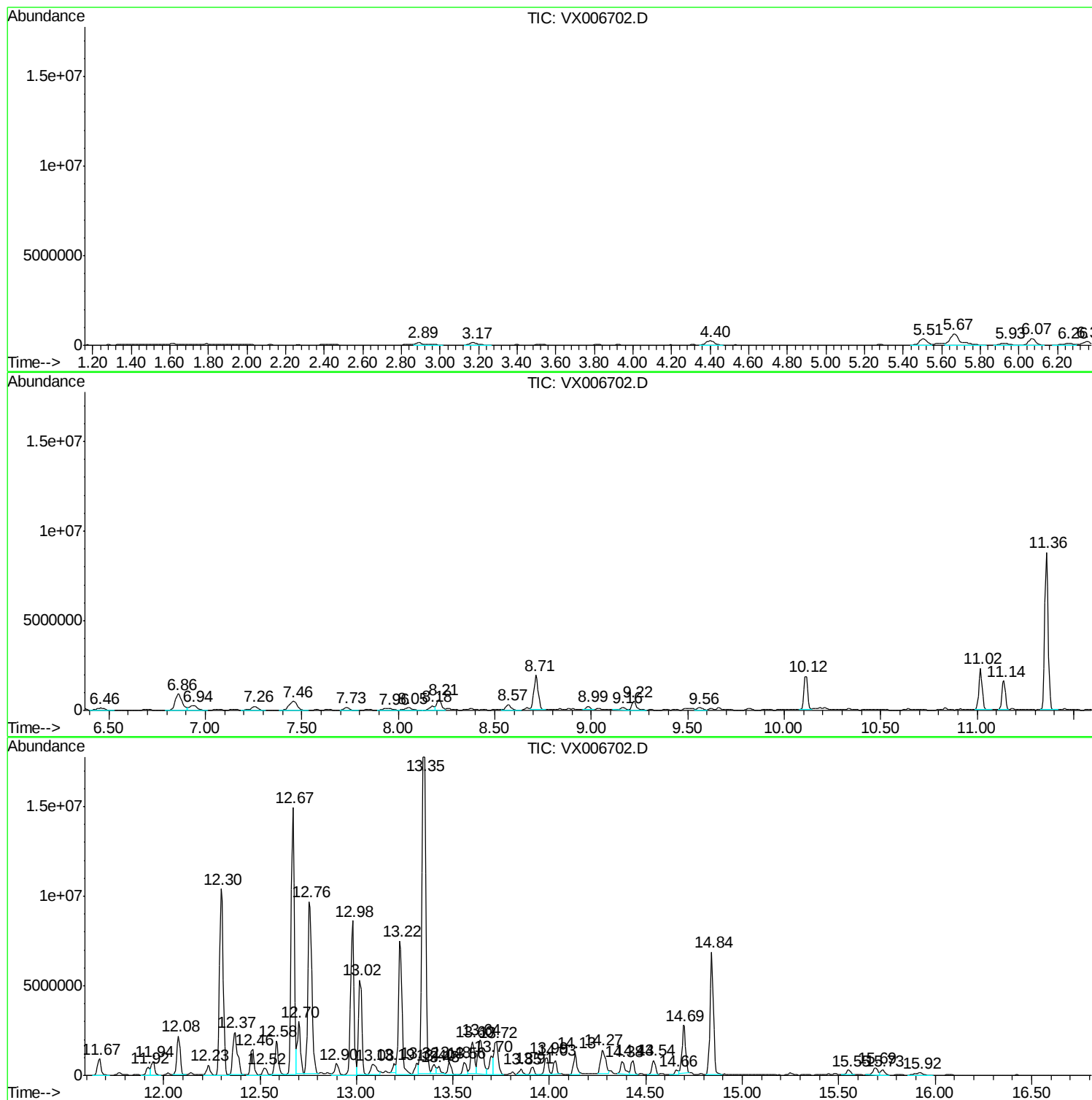
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X121818W.M
Quant Title : SW846 8260

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



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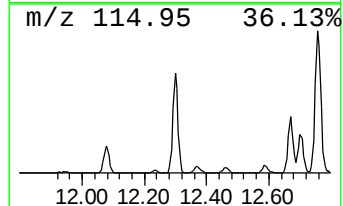
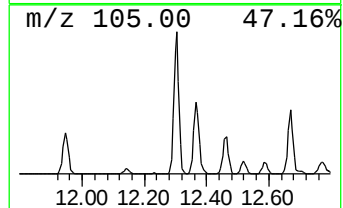
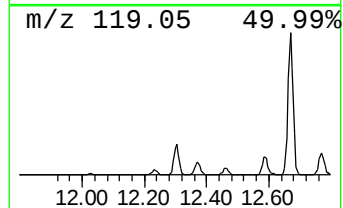
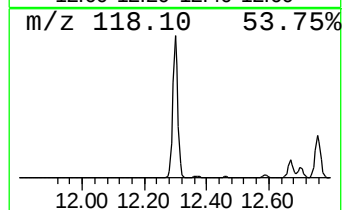
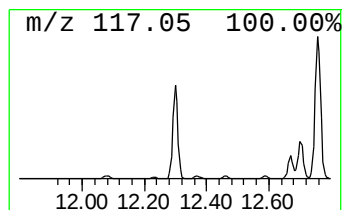
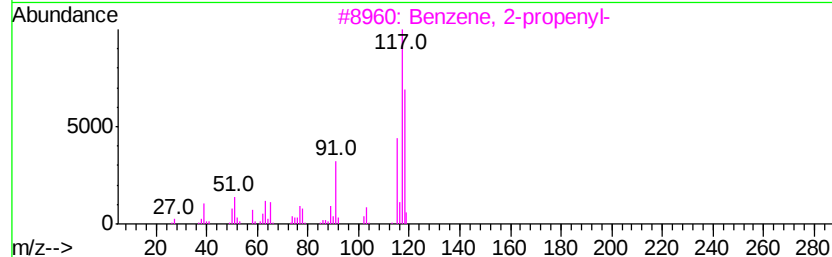
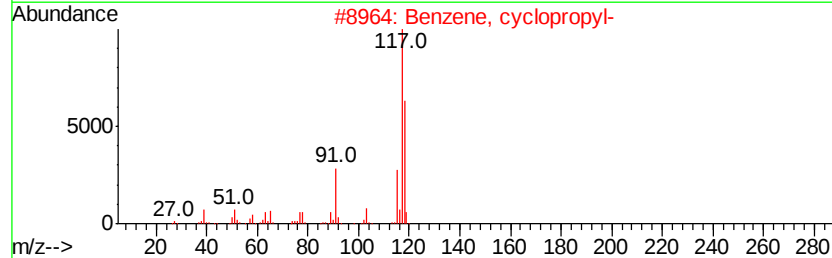
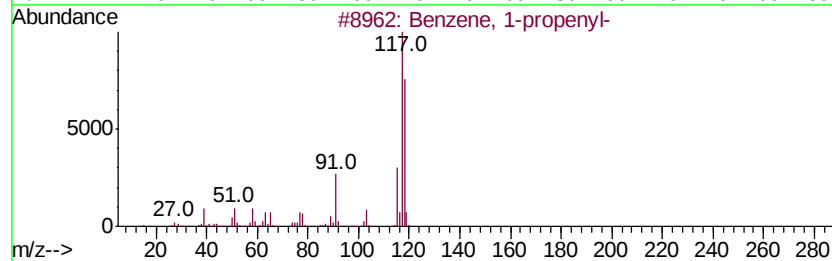
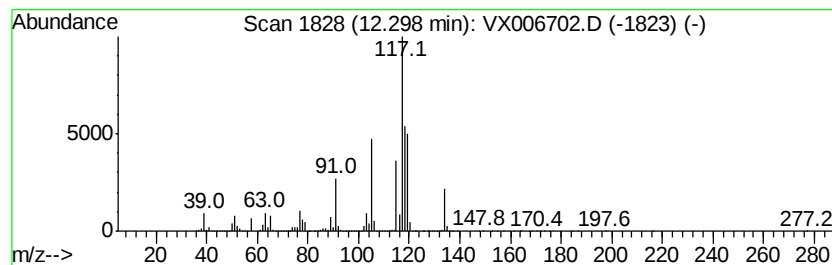
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X121818W.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1-propenyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	251.09 ug/l	13146900	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-propenyl-	118	C9H10	000637-50-3	91
2		Benzene, cyclopropyl-	118	C9H10	000873-49-4	90
3		Benzene, 2-propenyl-	118	C9H10	000300-57-2	60
4		Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	60
5		Indane	118	C9H10	000496-11-7	55



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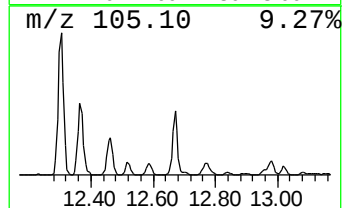
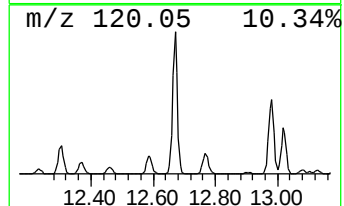
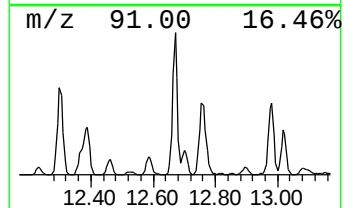
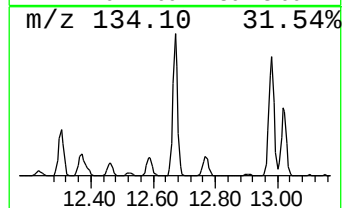
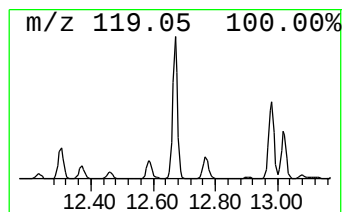
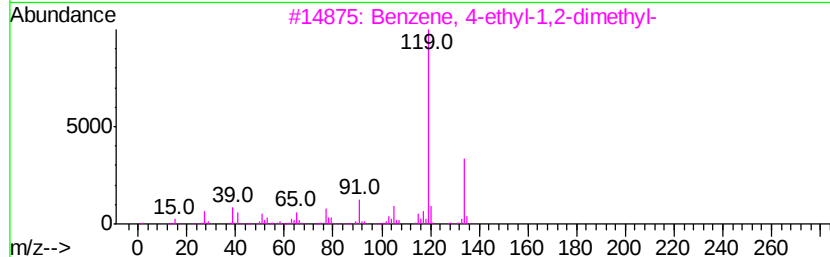
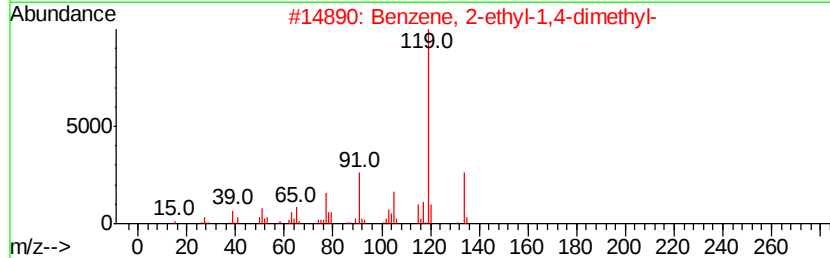
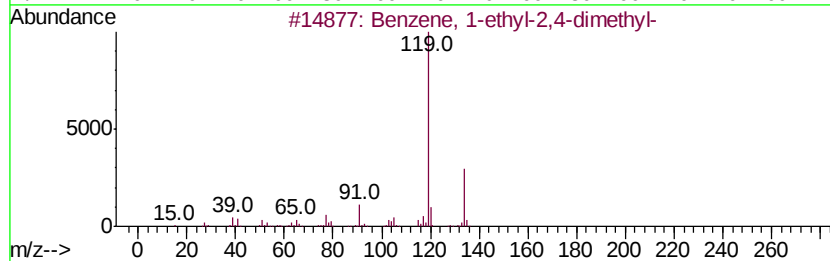
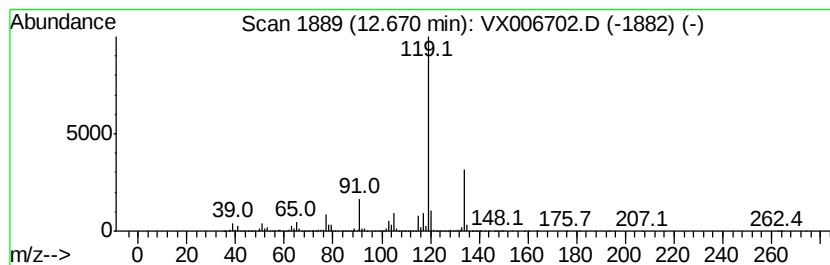
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.67	323.35 ug/l	16930100	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		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
3		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
4		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94



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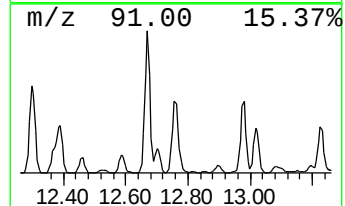
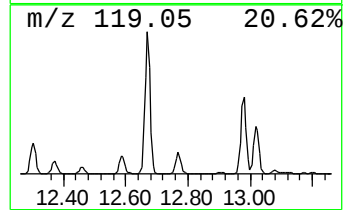
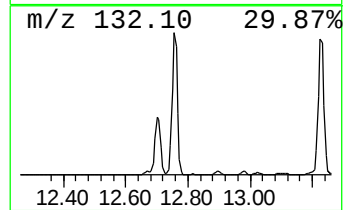
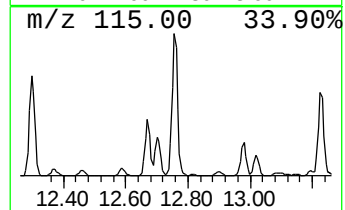
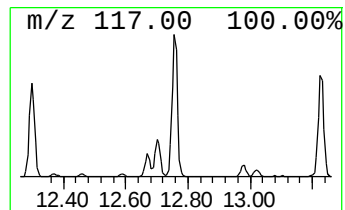
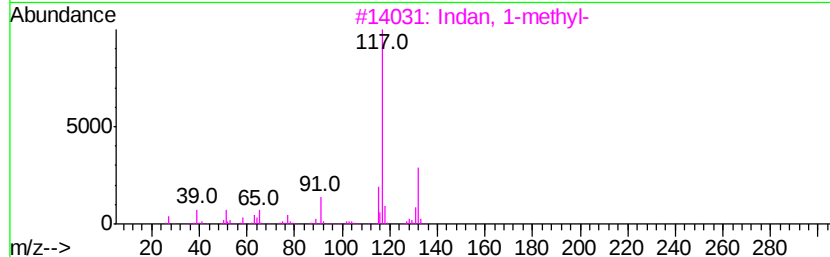
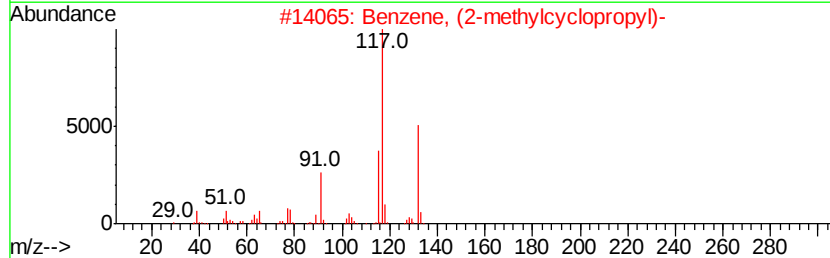
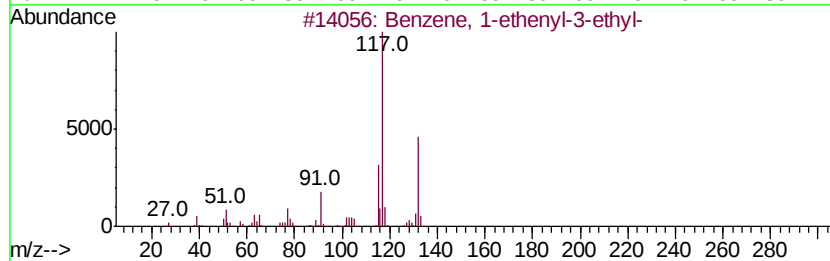
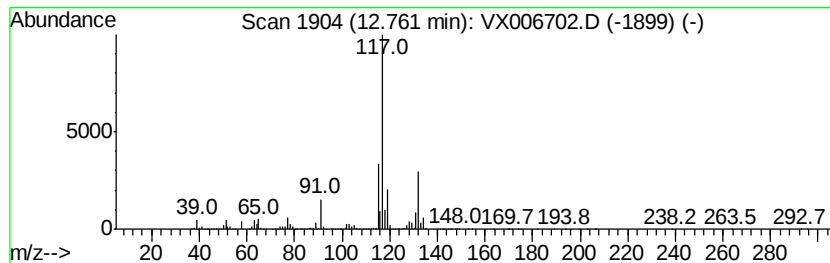
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	263.36 ug/l	13789200	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-methylcyclopropyl)-	132	C10H12	1000327-39-0	87
3		Indan, 1-methyl-	132	C10H12	000767-58-8	87
4		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	86
5		3-Phenylbut-1-ene	132	C10H12	000934-10-1	83



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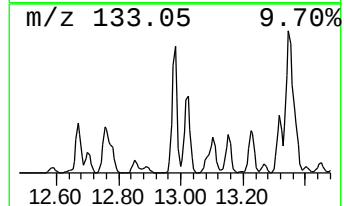
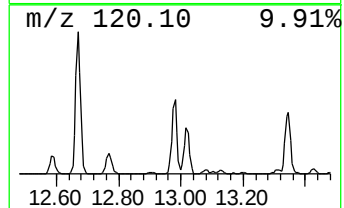
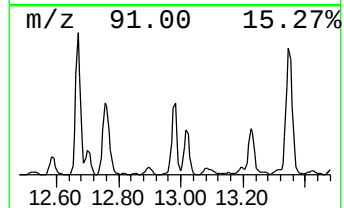
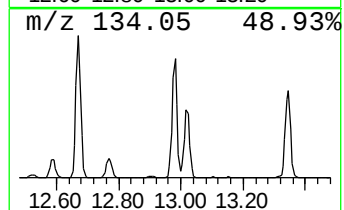
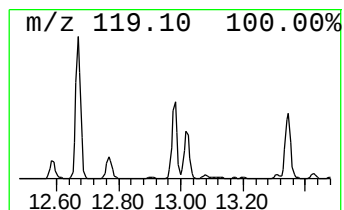
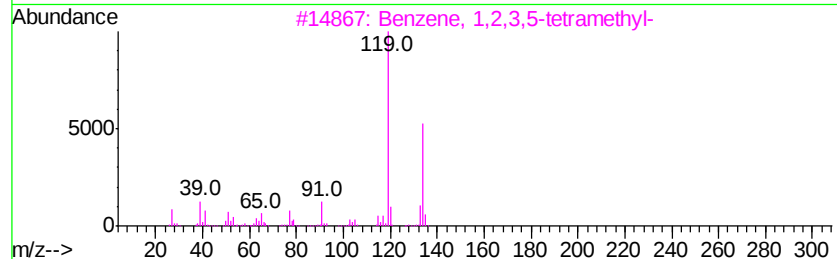
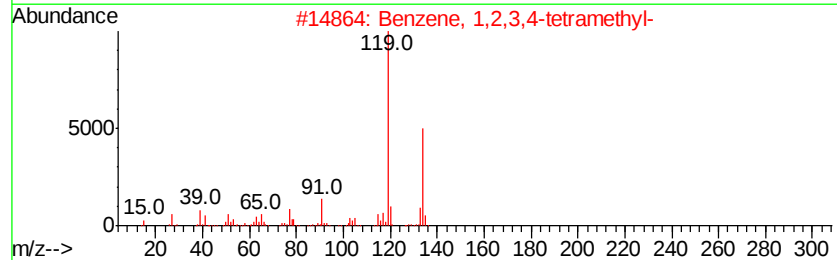
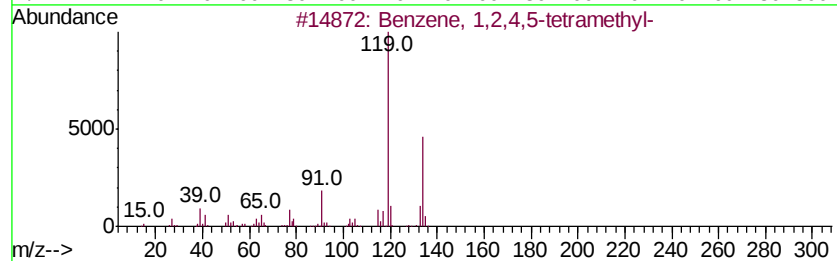
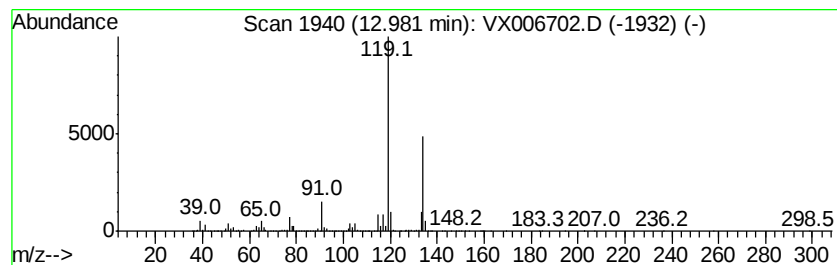
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.98	198.99 ug/l	10418700	1,4-Dichlorobenzene-d4	12.08

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, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX122018\
Data File : VX006702.D
Acq On : 21 Dec 2018 03:48
Operator : JC/MD
Sample : J6517-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 44 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
T3-MW-2-9.0-121918

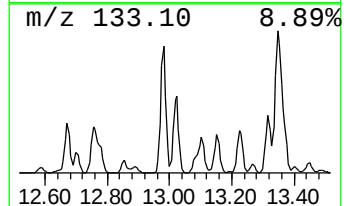
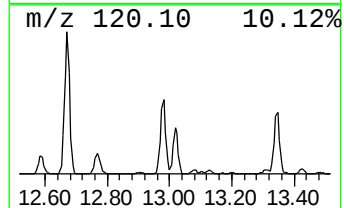
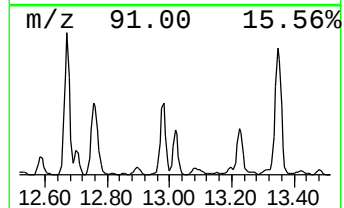
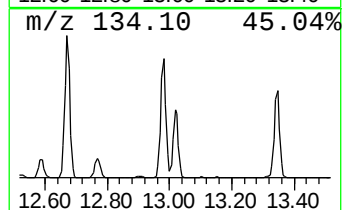
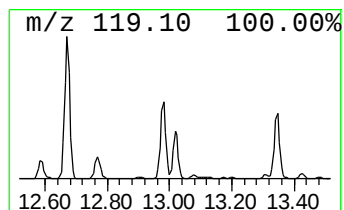
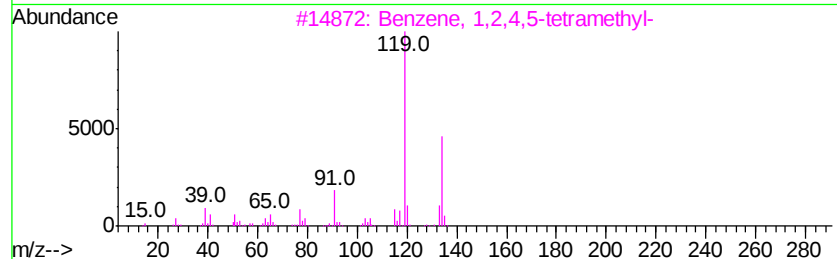
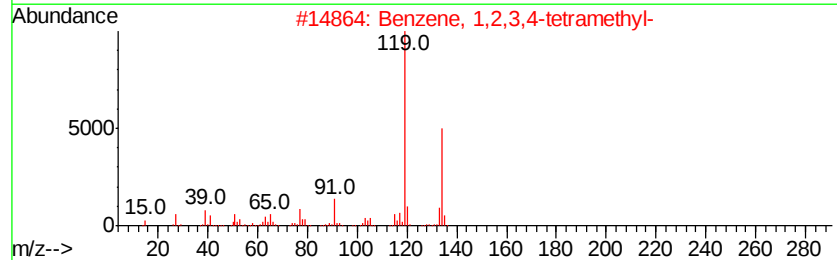
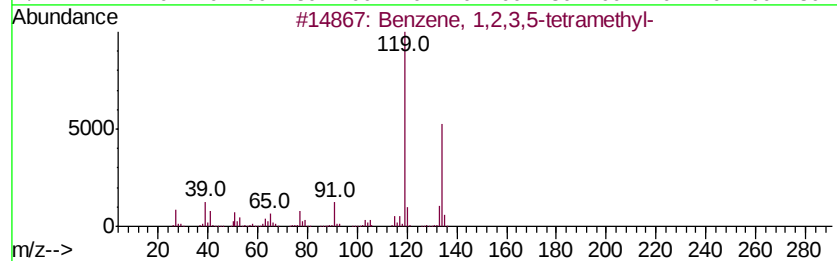
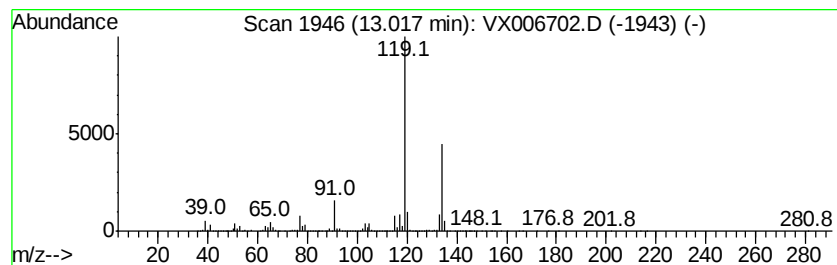
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.02	120.84 ug/l	6327260	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, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX122018\
Data File : VX006702.D
Acq On : 21 Dec 2018 03:48
Operator : JC/MD
Sample : J6517-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 44 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
T3-MW-2-9.0-121918

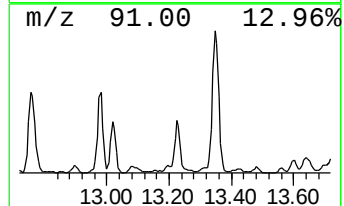
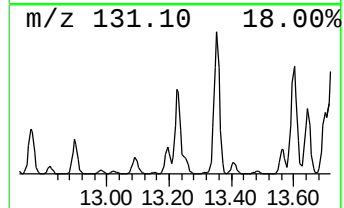
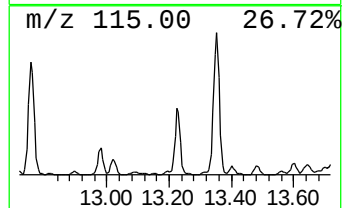
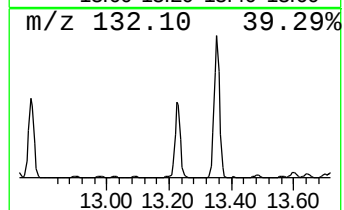
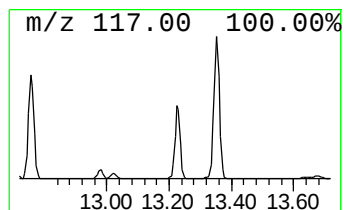
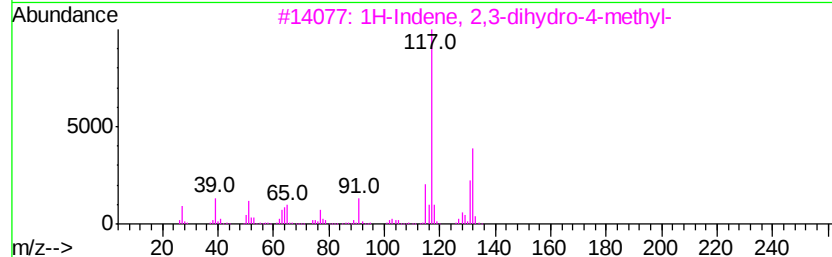
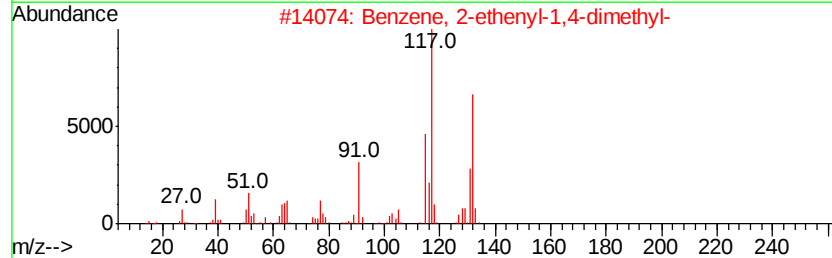
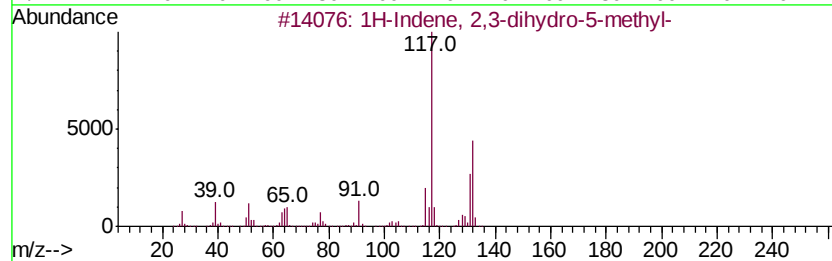
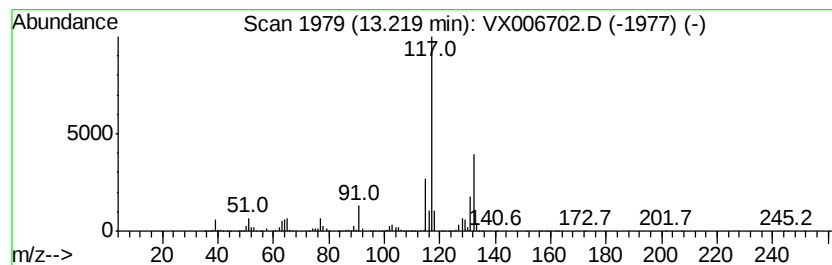
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Quant Title : SW846 8260

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

Peak Number 6 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.22	182.01 ug/l	9530060	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	93
2		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92
3		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90
4		Indan, 1-methyl-	132	C10H12	000767-58-8	90
5		3-Phenylbut-1-ene	132	C10H12	000934-10-1	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX122018\
Data File : VX006702.D
Acq On : 21 Dec 2018 03:48
Operator : JC/MD
Sample : J6517-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 44 Sample Multiplier: 1

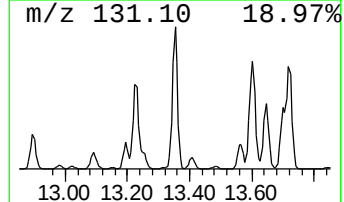
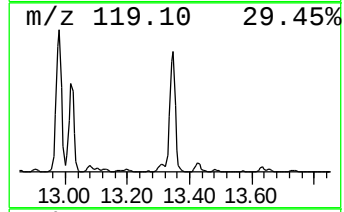
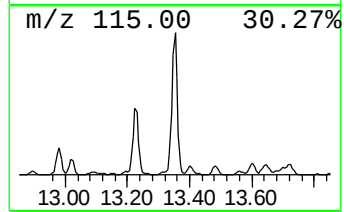
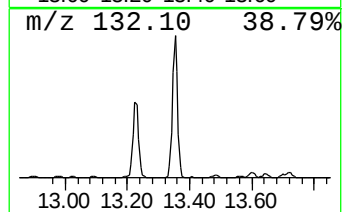
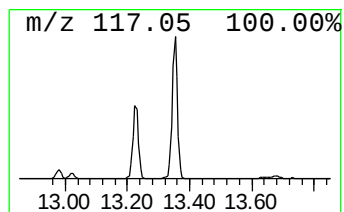
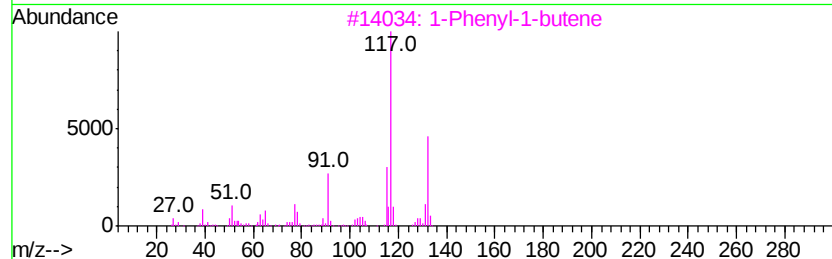
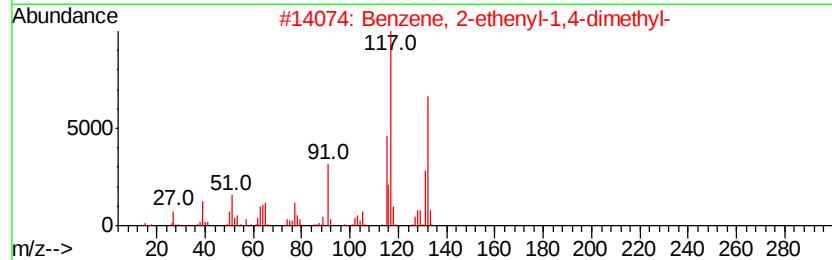
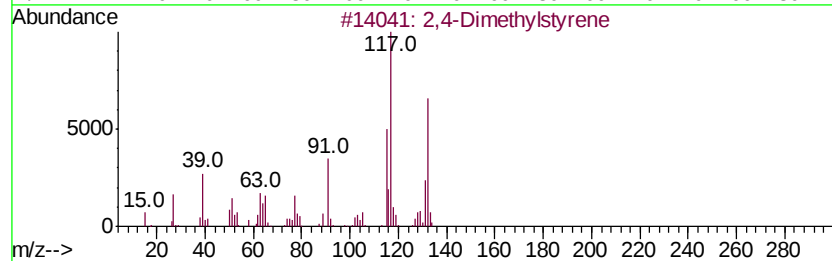
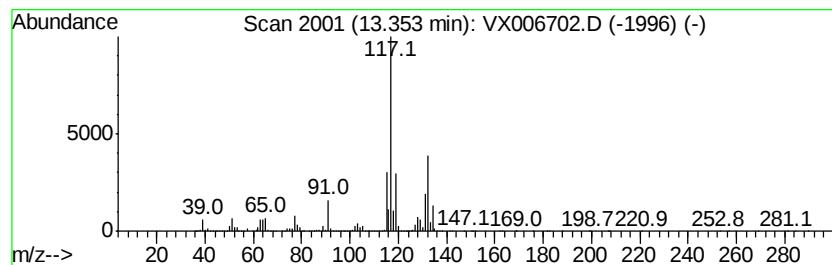
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MSVOA_X
ClientSampleId :
T3-MW-2-9.0-121918

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X121818W.M
Quant Title : SW846 8260

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

Peak Number 7 2,4-Dimethylstyrene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	483.21 ug/l	25300600	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	2,4-Dimethylstvrene	132 C10H12	002234-20-0	95
2	Benzene, 2-ethenyl-1,4-dimethyl-	132 C10H12	002039-89-6	95
3	1-Phenyl-1-butene	132 C10H12	000824-90-8	93
4	Benzene, 1-ethenyl-3-ethyl-	132 C10H12	007525-62-4	92
5	1H-Indene, 2,3-dihydro-4-methyl-	132 C10H12	000824-22-6	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX122018\
Data File : VX006702.D
Acq On : 21 Dec 2018 03:48
Operator : JC/MD
Sample : J6517-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 44 Sample Multiplier: 1

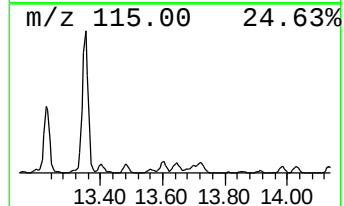
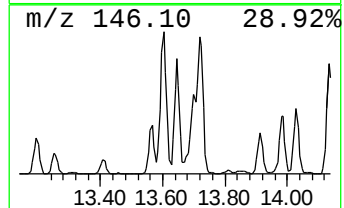
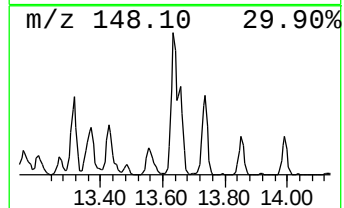
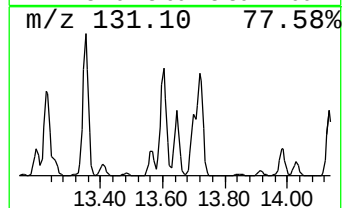
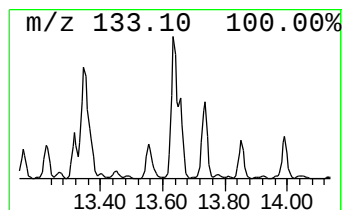
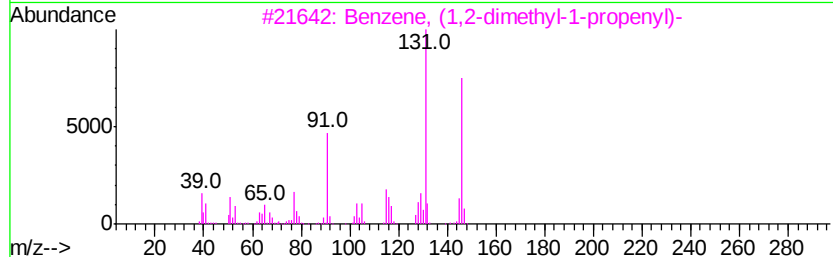
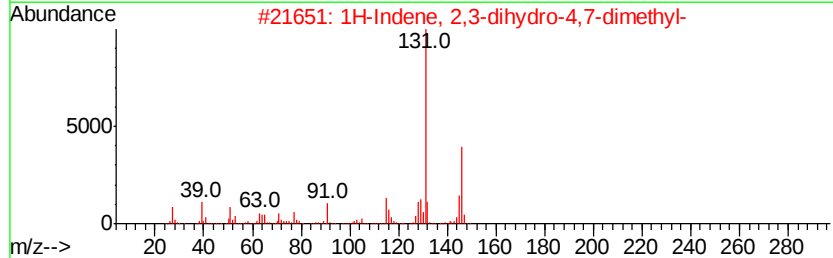
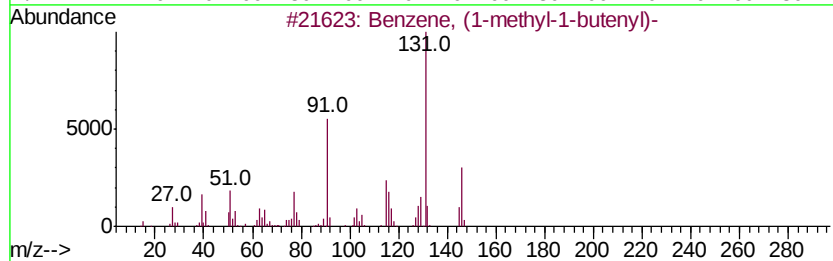
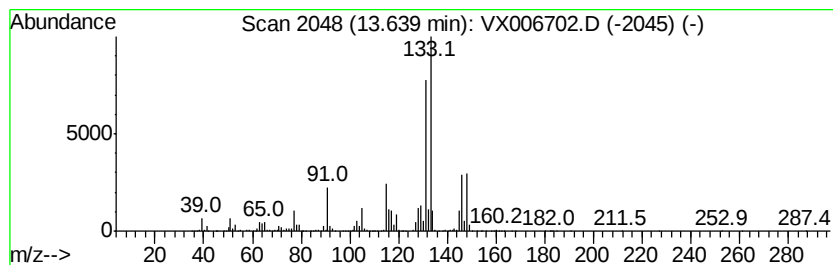
Instrument :
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ClientSampleId :
T3-MW-2-9.0-121918

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X121818W.M
Quant Title : SW846 8260

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

Peak Number 8 Benzene, (1-methyl-1-butenyl)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.64	68.93 ug/l	3608950	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, (1-methyl-1-butenyl)-	146 C11H14	053172-84-2 89
2		1H-Indene, 2,3-dihydro-4,7-dimet...	146 C11H14	006682-71-9 86
3		Benzene, (1,2-dimethyl-1-propenyl)-	146 C11H14	000769-57-3 70
4		Benzene, (2-methyl-1-butenyl)-	146 C11H14	056253-64-6 70
5		1H-Indene, 2,3-dihydro-1,2-dimet...	146 C11H14	017057-82-8 70



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX122018\
Data File : VX006702.D
Acq On : 21 Dec 2018 03:48
Operator : JC/MD
Sample : J6517-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 44 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
T3-MW-2-9.0-121918

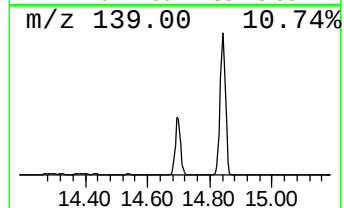
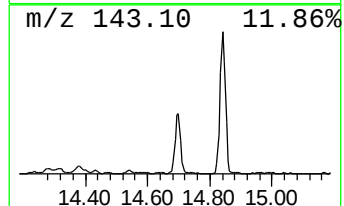
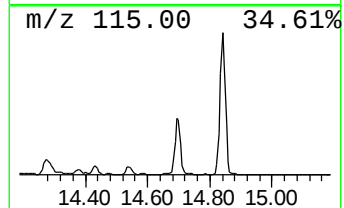
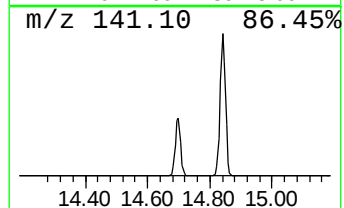
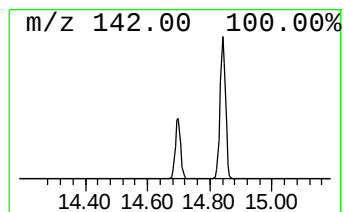
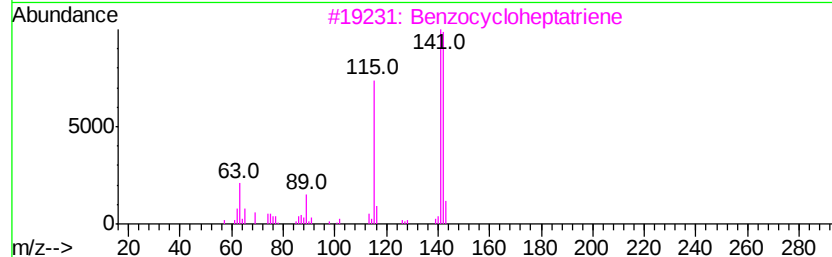
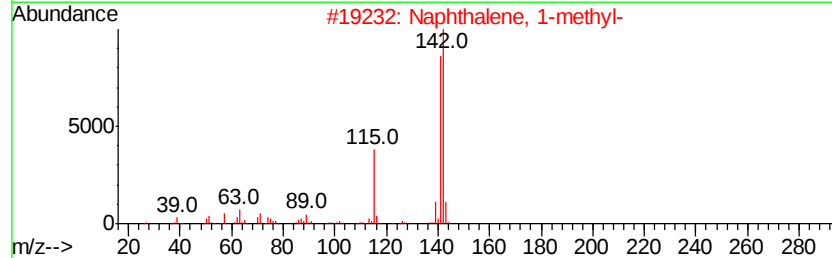
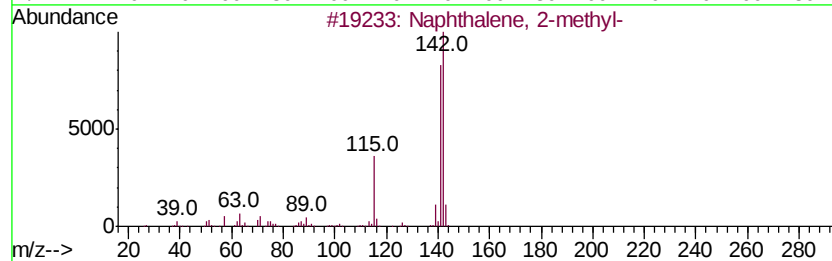
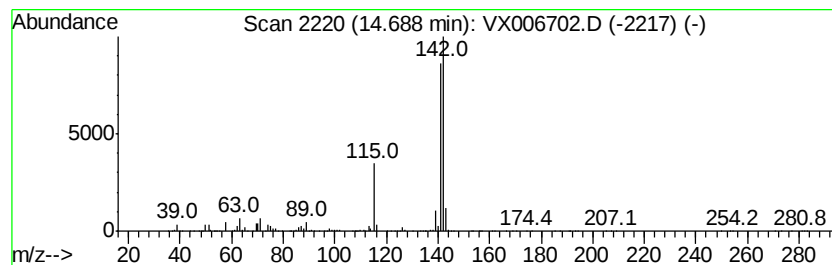
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Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.69	67.02 ug/l	3509050	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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	91
5		1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	64



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX122018\
Data File : VX006702.D
Acq On : 21 Dec 2018 03:48
Operator : JC/MD
Sample : J6517-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 44 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
T3-MW-2-9.0-121918

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X121818W.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-propenyl-	12.30	251.1	ug/l	13146900	4	12.08	2617950	50.0
Benzene, 1-ethyl-...	12.67	323.4	ug/l	16930100	4	12.08	2617950	50.0
Benzene, 1-etheny...	12.76	263.4	ug/l	13789200	4	12.08	2617950	50.0
Benzene, 1,2,4,5-...	12.98	199.0	ug/l	10418700	4	12.08	2617950	50.0
Benzene, 1,2,3,5-...	13.02	120.8	ug/l	6327260	4	12.08	2617950	50.0
1H-Indene, 2,3-di...	13.22	182.0	ug/l	9530060	4	12.08	2617950	50.0
2,4-Dimethylstyrene	13.35	483.2	ug/l	25300600	4	12.08	2617950	50.0
Benzene, (1-methy...	13.64	68.9	ug/l	3608950	4	12.08	2617950	50.0
Naphthalene, 1-me...	14.69	67.0	ug/l	3509050	4	12.08	2617950	50.0