

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX031619\
 Data File : VX008150.D
 Acq On : 16 Mar 2019 19:20
 Operator : JC/SP
 Sample : K1960-06
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 T3-MW-3-9.0-031419

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\82X031419W.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	1.392	27	39	40	rBV2	14275	41986	3.23%	0.179%
2	1.794	100	105	117	rVB2	24885	39502	3.04%	0.169%
3	2.861	272	280	281	rVV2	21370	43827	3.37%	0.187%
4	2.891	282	285	291	rVB2	28176	51842	3.99%	0.222%
5	3.172	324	331	345	rVB2	29264	78622	6.05%	0.336%
6	4.397	523	532	544	rVB2	44553	128382	9.88%	0.549%
7	4.696	573	581	593	rBV2	21277	61498	4.73%	0.263%
8	5.501	698	713	722	rBV	149054	437169	33.65%	1.869%
9	5.665	730	740	764	rVB	292762	869082	66.89%	3.715%
10	6.068	796	806	818	rBV2	152863	392905	30.24%	1.679%
11	6.257	830	837	845	rBV3	9523	25932	2.00%	0.111%
12	6.354	845	853	861	rVV5	12764	37045	2.85%	0.158%
13	6.452	861	869	879	rVB4	12411	33613	2.59%	0.144%
14	6.860	926	936	946	rBV	396309	966948	74.42%	4.133%
15	7.256	991	1001	1010	rVB3	17727	42242	3.25%	0.181%
16	7.458	1023	1034	1046	rVB4	52200	138029	10.62%	0.590%
17	8.214	1145	1158	1163	rBV	38252	87188	6.71%	0.373%
18	8.567	1211	1216	1224	rVB	22467	40016	3.08%	0.171%
19	8.665	1224	1232	1234	rBV4	11600	21181	1.63%	0.091%
20	8.713	1234	1240	1248	rVB	795143	1244675	95.79%	5.320%
21	9.037	1288	1293	1299	rVB2	12185	18792	1.45%	0.080%
22	9.220	1319	1323	1332	rVB	15504	25054	1.93%	0.107%
23	9.622	1384	1389	1393	rVV2	21460	30009	2.31%	0.128%
24	10.110	1463	1469	1475	rBV	753686	1017204	78.29%	4.348%
25	10.183	1475	1481	1487	rVV	17775	38996	3.00%	0.167%
26	10.335	1502	1506	1507	rBV	16268	20127	1.55%	0.086%
27	10.372	1507	1512	1515	rVV2	31287	62463	4.81%	0.267%
28	10.408	1515	1518	1526	rVB	18559	27424	2.11%	0.117%
29	10.512	1531	1535	1540	rVB	35068	48912	3.76%	0.209%
30	10.658	1546	1559	1564	rBV4	62939	179591	13.82%	0.768%
31	10.707	1564	1567	1575	rVB	19363	30479	2.35%	0.130%
32	10.835	1584	1588	1594	rVB	152999	205276	15.80%	0.877%
33	10.908	1595	1600	1607	rBV3	86165	144184	11.10%	0.616%
34	11.018	1613	1618	1622	rBV	278008	356460	27.43%	1.524%

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35	11.134	1632	1637	1642	rBV	648132	823016	63.34%	3.518%
36	11.183	1642	1645	1649	rVV	76432	96261	7.41%	0.411%
37	11.280	1657	1661	1665	rVB3	68514	91865	7.07%	0.393%
38	11.359	1666	1674	1682	rBV	858109	1049339	80.76%	4.485%
39	11.445	1684	1688	1693	rVB3	27202	44514	3.43%	0.190%
40	11.579	1706	1710	1714	rVB4	16530	24983	1.92%	0.107%
41	11.670	1719	1725	1732	rBV2	117832	173679	13.37%	0.742%
42	11.768	1736	1741	1744	rVV	49566	60919	4.69%	0.260%
43	11.804	1744	1747	1752	rVB5	11880	19626	1.51%	0.084%
44	11.920	1758	1766	1767	rBV	167543	223754	17.22%	0.956%
45	11.945	1767	1770	1775	rVB	297810	357744	27.53%	1.529%
46	12.079	1786	1792	1797	rBV	780275	993597	76.47%	4.247%
47	12.231	1812	1817	1822	rVB	113441	137885	10.61%	0.589%
48	12.298	1823	1828	1834	rBV	1058343	1299318	100.00%	5.554%
49	12.365	1834	1839	1841	rBV2	231940	312864	24.08%	1.337%
50	12.457	1849	1854	1860	rBV	173164	223242	17.18%	0.954%
51	12.524	1860	1865	1869	rVV2	20746	32566	2.51%	0.139%
52	12.664	1880	1888	1891	rVV	494232	640983	49.33%	2.740%
53	12.701	1891	1894	1899	rVV	301424	411402	31.66%	1.758%
54	12.755	1899	1903	1910	rVV2	730529	1237403	95.23%	5.289%
55	12.816	1910	1913	1918	rVV2	47481	72954	5.61%	0.312%
56	12.896	1918	1926	1931	rVB2	163719	264667	20.37%	1.131%
57	12.975	1931	1939	1944	rBV	798468	1037138	79.82%	4.433%
58	13.024	1944	1947	1952	rVB2	35898	50182	3.86%	0.214%
59	13.079	1952	1956	1963	rVB2	89474	133435	10.27%	0.570%
60	13.152	1963	1968	1971	rBV	91927	128385	9.88%	0.549%
61	13.194	1971	1975	1977	rVV2	135909	181358	13.96%	0.775%
62	13.225	1977	1980	1989	rVB	289382	381566	29.37%	1.631%
63	13.310	1989	1994	1996	rBV2	40671	59617	4.59%	0.255%
64	13.347	1996	2000	2006	rVV2	717293	927696	71.40%	3.965%
65	13.402	2006	2009	2011	rVV2	55865	77749	5.98%	0.332%
66	13.426	2011	2013	2015	rVV	65087	66959	5.15%	0.286%
67	13.481	2018	2022	2029	rVB	86326	116515	8.97%	0.498%
68	13.560	2031	2035	2038	rBV2	73664	100221	7.71%	0.428%
69	13.597	2038	2041	2044	rVV	115211	159147	12.25%	0.680%
70	13.639	2044	2048	2052	rVV2	220874	330453	25.43%	1.412%
71	13.694	2052	2057	2059	rVV3	131260	237751	18.30%	1.016%

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72	13.719	2059	2061	2070	rVB2	202593	279848	21.54%	1.196%
73	13.804	2070	2075	2078	rBV	54675	73824	5.68%	0.316%
74	13.853	2078	2083	2088	rVB3	39205	65859	5.07%	0.281%
75	13.908	2088	2092	2097	rVB	200429	242130	18.64%	1.035%
76	13.981	2097	2104	2108	rBV2	166802	256084	19.71%	1.095%
77	14.030	2108	2112	2115	rVV	56603	77849	5.99%	0.333%
78	14.127	2124	2128	2131	rBV	63949	86272	6.64%	0.369%
79	14.164	2131	2134	2138	rVB2	25891	29512	2.27%	0.126%
80	14.267	2146	2151	2161	rVB2	146698	252384	19.42%	1.079%
81	14.377	2165	2169	2173	rVV	67753	82346	6.34%	0.352%
82	14.432	2173	2178	2182	rVB	73216	95751	7.37%	0.409%
83	14.475	2182	2185	2190	rVB	33884	43613	3.36%	0.186%
84	14.536	2190	2195	2200	rBV2	303693	402855	31.01%	1.722%
85	14.670	2208	2217	2223	rBV4	68562	137399	10.57%	0.587%
86	14.731	2223	2227	2231	rVB	60223	77118	5.94%	0.330%
87	14.786	2231	2236	2239	rBV2	17196	22636	1.74%	0.097%
88	14.840	2239	2245	2249	rBV	514247	678376	52.21%	2.900%
89	14.901	2253	2255	2260	rVB4	31589	40378	3.11%	0.173%
90	14.962	2260	2265	2266	rBV2	16169	21192	1.63%	0.091%
91	15.249	2306	2312	2319	rBV	45166	84777	6.52%	0.362%
92	15.322	2319	2324	2329	rVB5	12664	20802	1.60%	0.089%
93	15.444	2339	2344	2346	rBV	26644	39022	3.00%	0.167%
94	15.474	2346	2349	2353	rVV	28112	40256	3.10%	0.172%
95	15.548	2354	2361	2366	rVV	131151	208906	16.08%	0.893%
96	15.596	2366	2369	2375	rVV3	12964	24381	1.88%	0.104%
97	15.688	2376	2384	2388	rVV	144281	247306	19.03%	1.057%
98	15.724	2388	2390	2396	rVB	39335	49630	3.82%	0.212%
99	15.913	2410	2421	2432	rVB	43754	119310	9.18%	0.510%
100	16.066	2441	2446	2451	rBV2	18623	29068	2.24%	0.124%

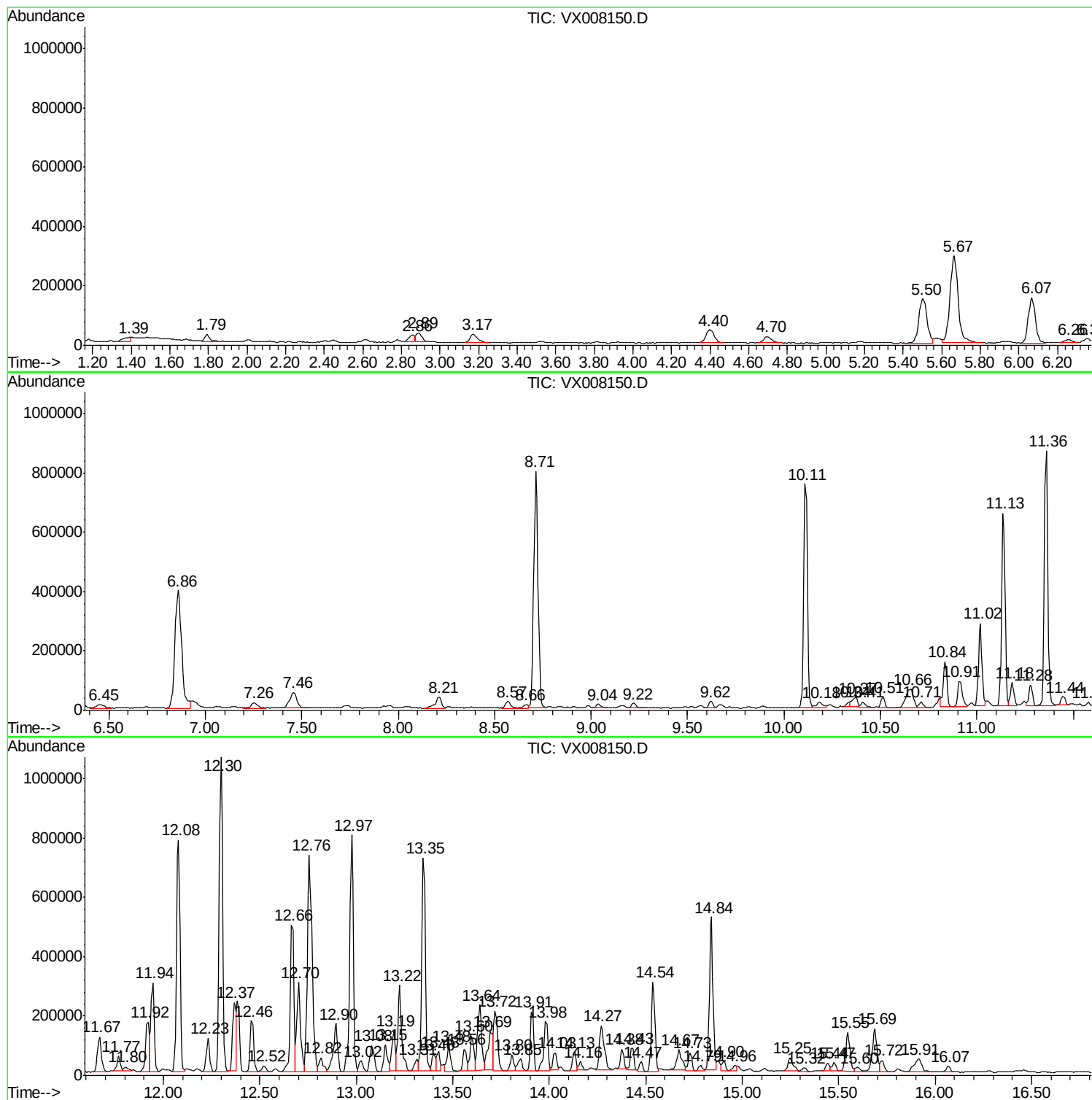
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X031419W.M
Quant Title : SW846 8260

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



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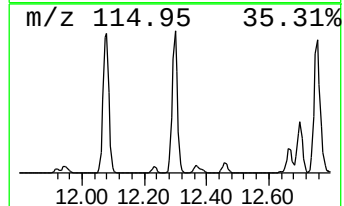
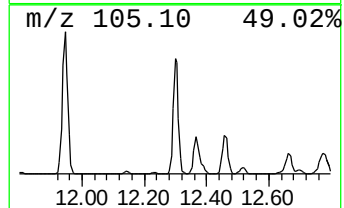
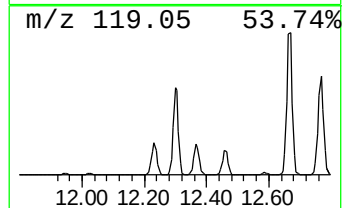
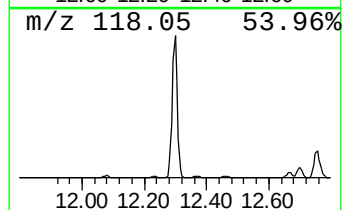
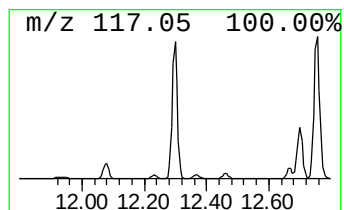
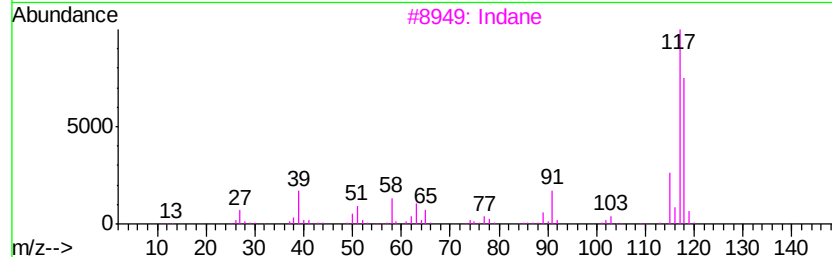
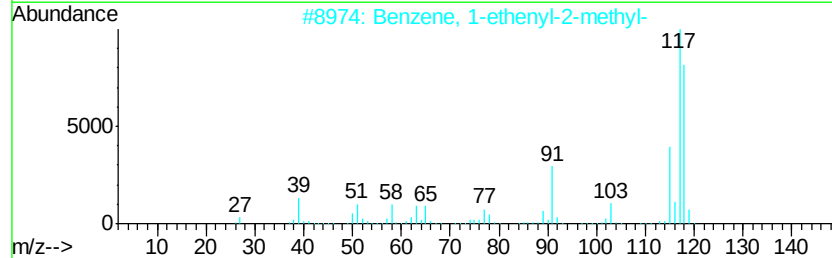
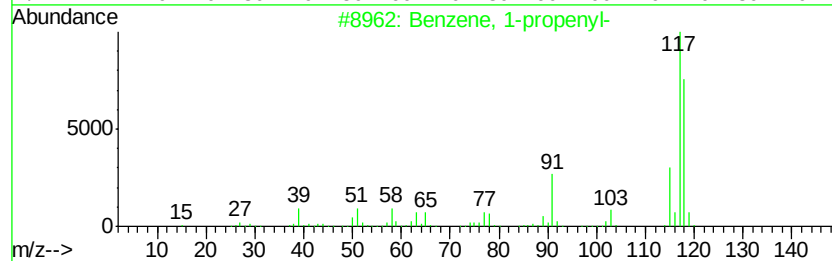
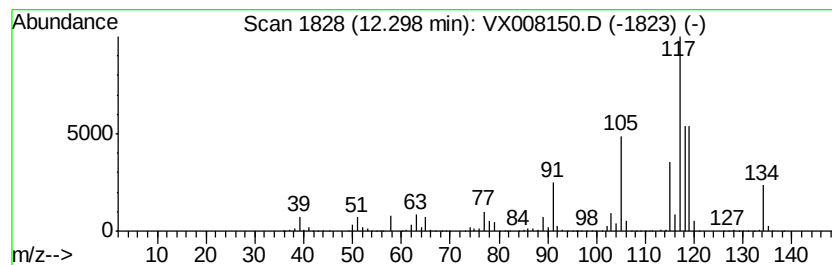
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X031419W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	65.38 ug/l	1299320	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-propenyl-	118 C9H10	000637-50-3 91
2		Benzene, 1-ethenyl-2-methyl-	118 C9H10	000611-15-4 80
3		Indane	118 C9H10	000496-11-7 60
4		Benzene, cyclopropyl-	118 C9H10	000873-49-4 60
5		Benzene, 2-propenyl-	118 C9H10	000300-57-2 60



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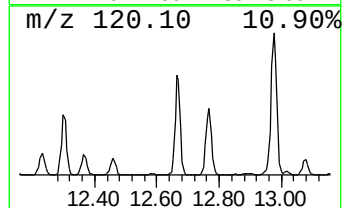
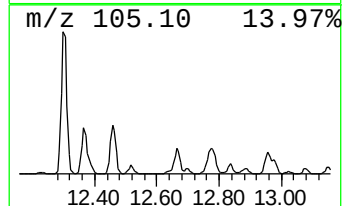
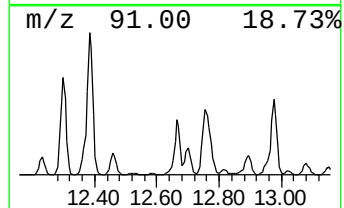
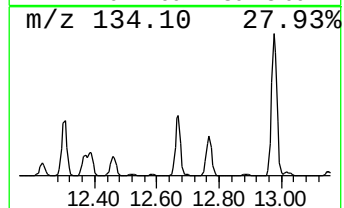
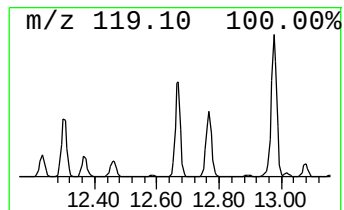
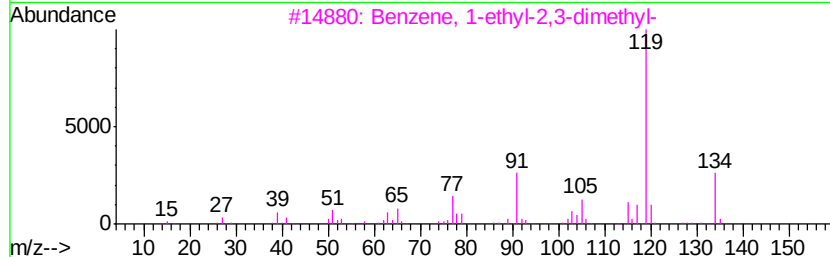
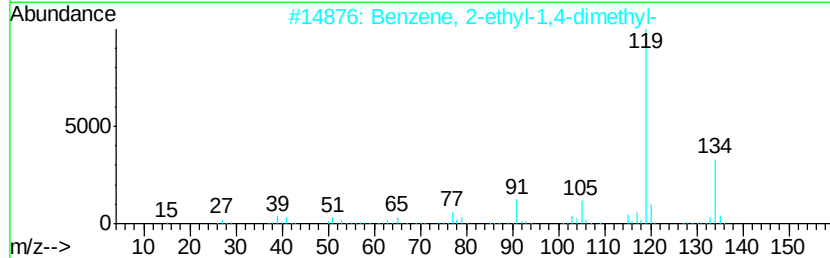
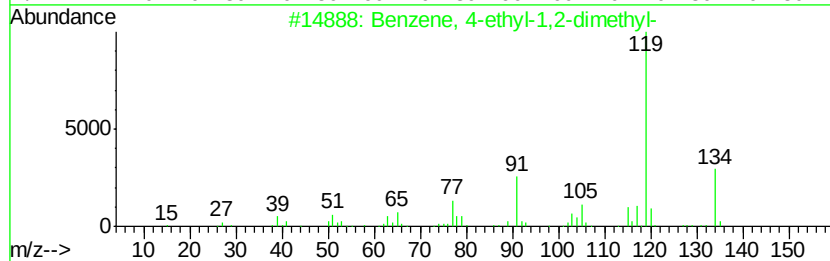
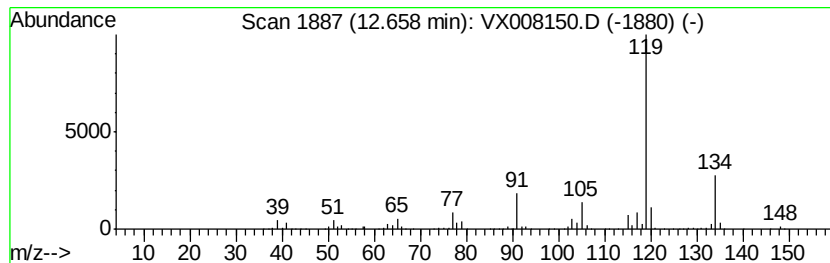
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.		
12.66	32.26 ug/l	640983	1,4-Dichlorobenzene-d4	12.08		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96	
2	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95	
3	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95	
4	o-Cymene	134	C10H14	000527-84-4	95	
5	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94	



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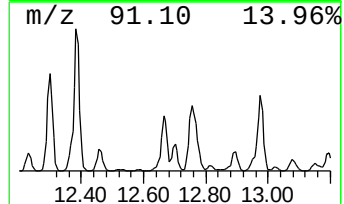
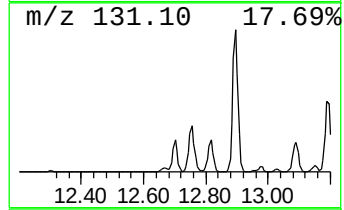
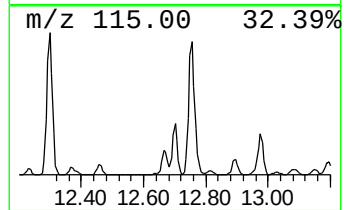
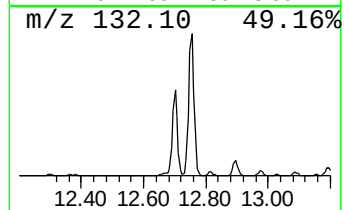
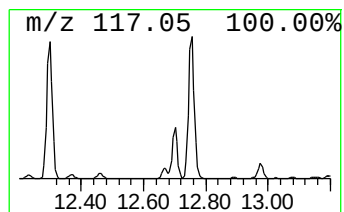
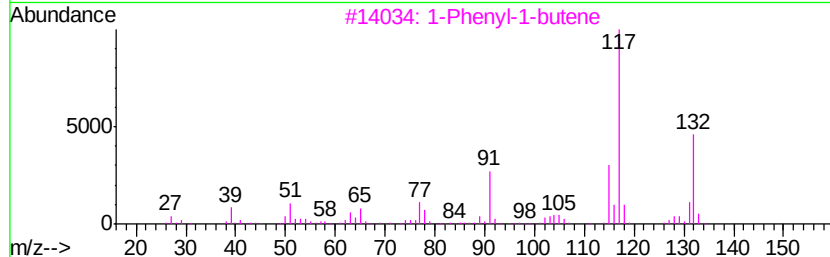
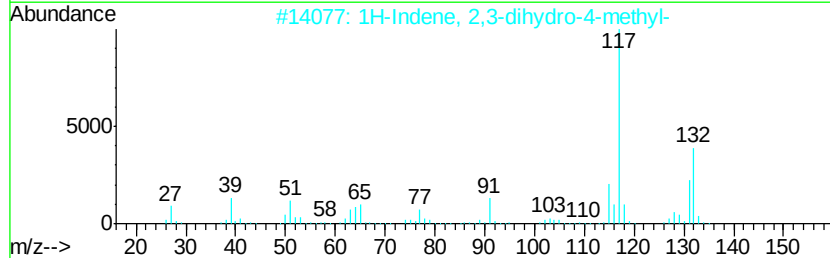
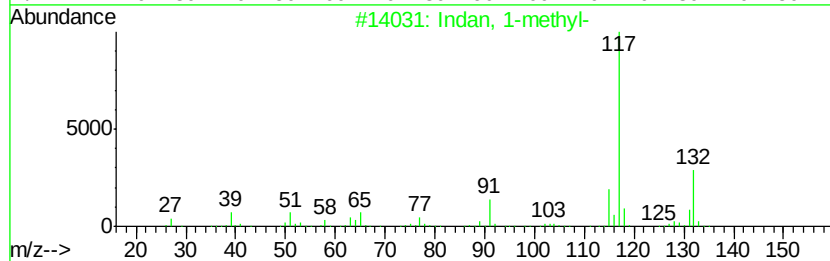
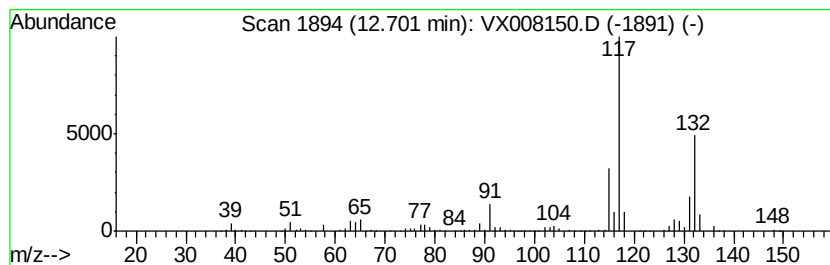
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Peak Number 3 Indan, 1-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.70	20.70 ug/l	411402	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		1H-Indene, 2,3-dihydro-4-methyl-	132 C10H12	000824-22-6 90
3		1-Phenyl-1-butene	132 C10H12	000824-90-8 90
4		Benzene, (2-methyl-2-propenyl)-	132 C10H12	003290-53-7 90
5		(E)-1-Phenyl-1-butene	132 C10H12	001005-64-7 90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX031619\
Data File : VX008150.D
Acq On : 16 Mar 2019 19:20
Operator : JC/SP
Sample : K1960-06
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
T3-MW-3-9.0-031419

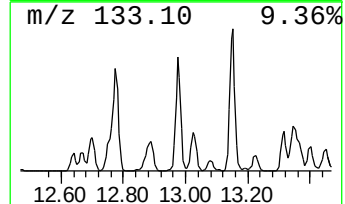
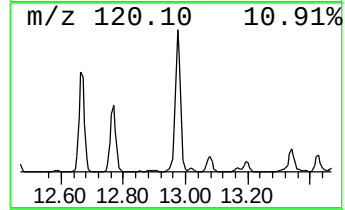
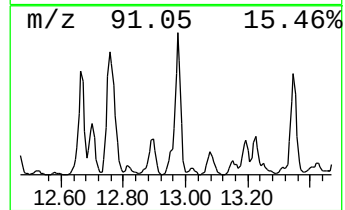
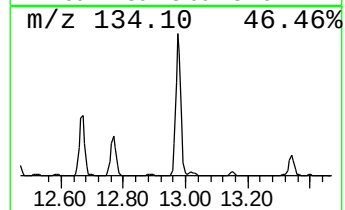
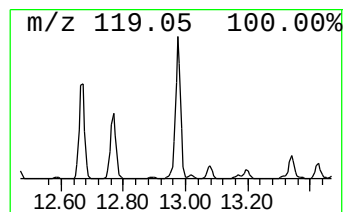
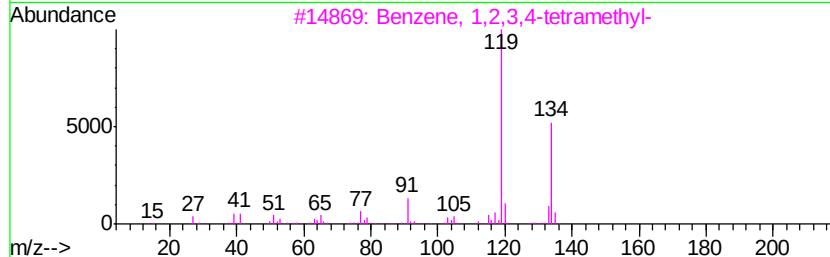
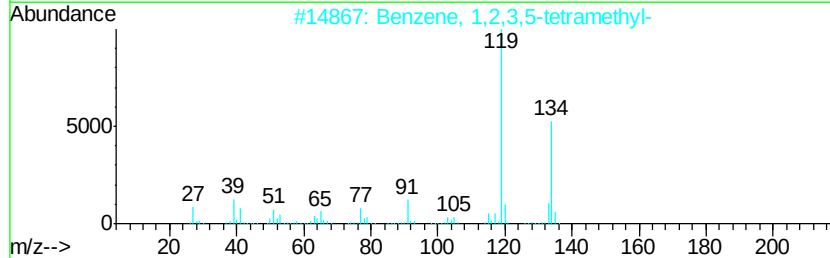
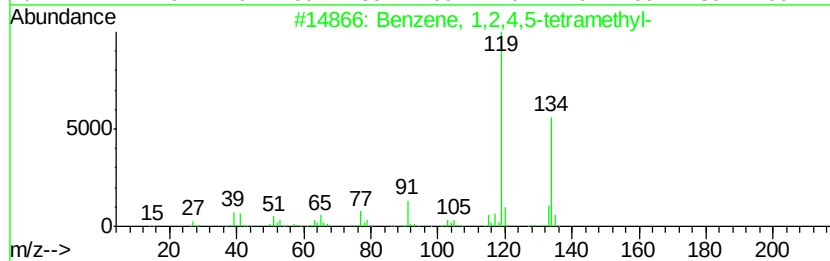
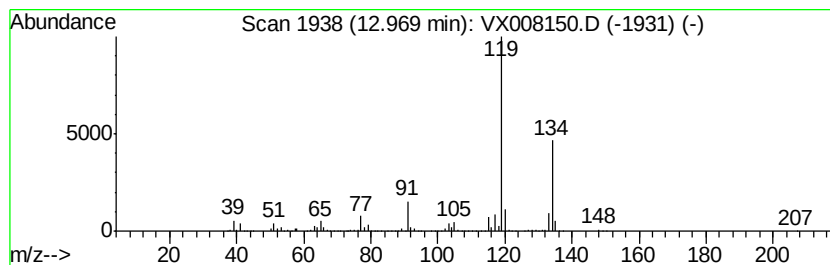
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X031419W.M
Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.97	52.19 ug/l	1037140	1,4-Dichlorobenzene-d4	12.08

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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX031619\
Data File : VX008150.D
Acq On : 16 Mar 2019 19:20
Operator : JC/SP
Sample : K1960-06
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
T3-MW-3-9.0-031419

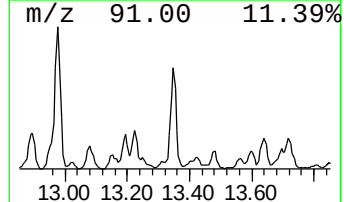
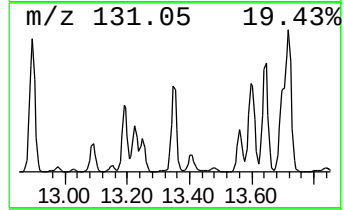
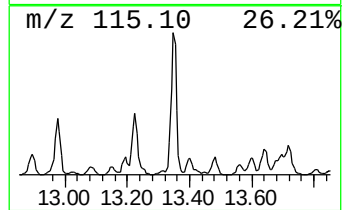
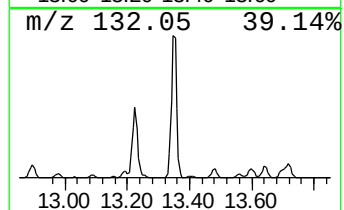
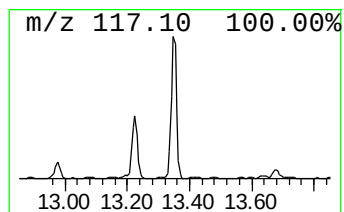
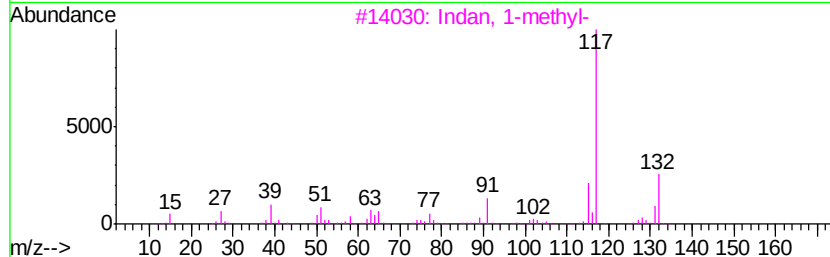
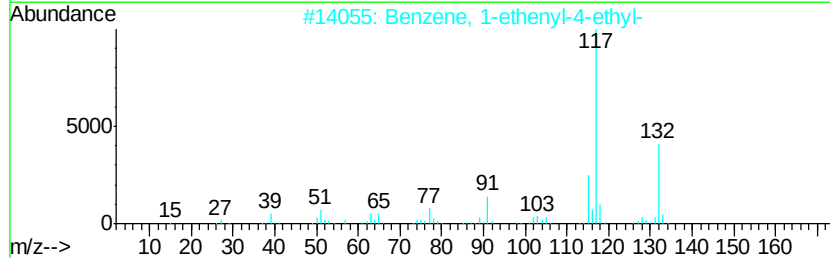
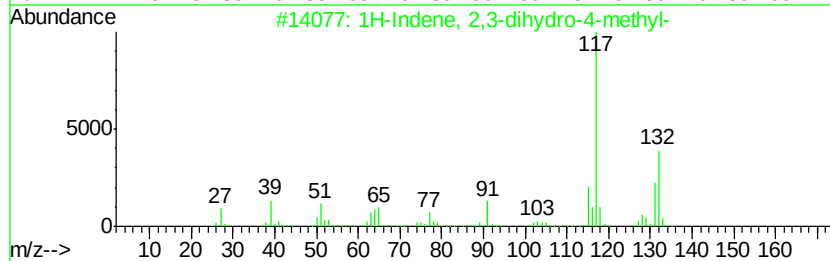
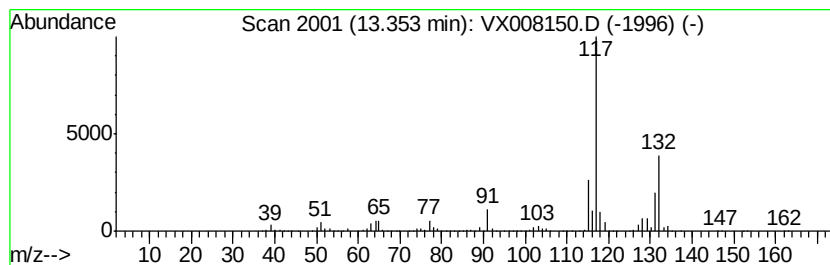
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	46.68 ug/l	927696	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	93
2		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	76
3		Indan, 1-methyl-	132	C10H12	000767-58-8	74
4		Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	72
5		1-Phenyl-1-butene	132	C10H12	000824-90-8	72



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX031619\
Data File : VX008150.D
Acq On : 16 Mar 2019 19:20
Operator : JC/SP
Sample : K1960-06
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 18 Sample Multiplier: 1

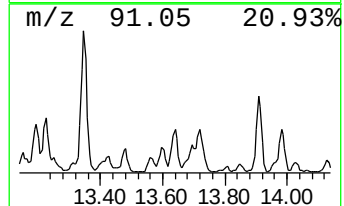
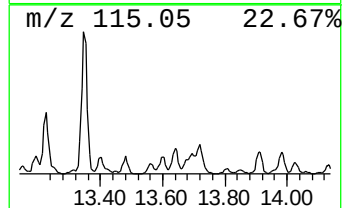
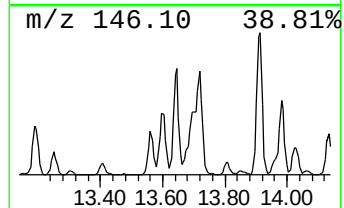
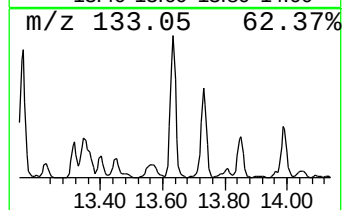
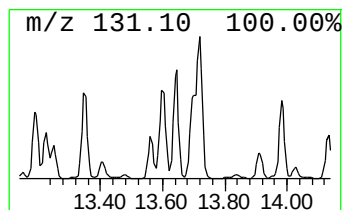
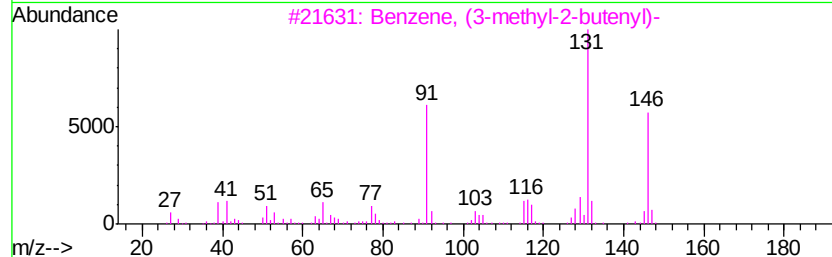
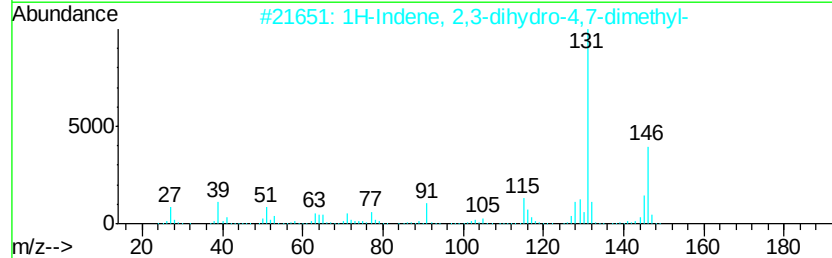
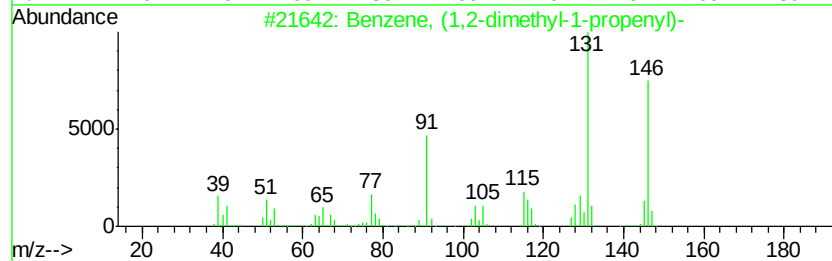
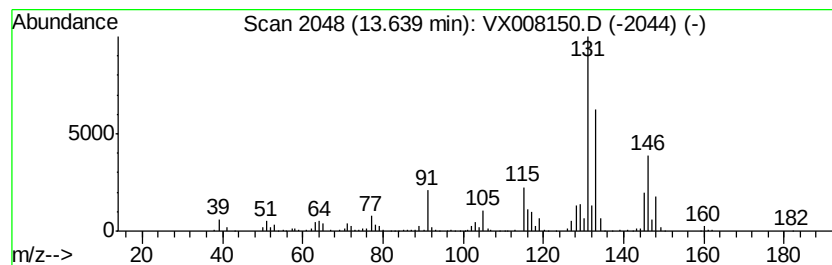
Instrument :
MSVOA_X
ClientSampleId :
T3-MW-3-9.0-031419

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X031419W.M
Quant Title : SW846 8260

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

Peak Number 8 Benzene, (1,2-dimethyl-1-propenyl)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.64	16.63 ug/l	330453	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, (1,2-dimethyl-1-propenyl)-	146 C11H14	000769-57-3 94
2		1H-Indene, 2,3-dihydro-4,7-dimet...	146 C11H14	006682-71-9 94
3		Benzene, (3-methyl-2-butenyl)-	146 C11H14	004489-84-3 92
4		1H-Indene, 2,3-dihydro-1,2-dimet...	146 C11H14	017057-82-8 86
5		Benzene, (1-methyl-1-butenyl)-	146 C11H14	053172-84-2 76



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX031619\
Data File : VX008150.D
Acq On : 16 Mar 2019 19:20
Operator : JC/SP
Sample : K1960-06
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
T3-MW-3-9.0-031419

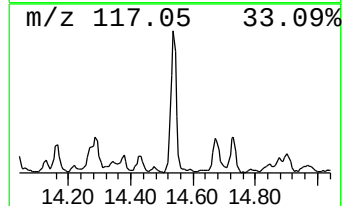
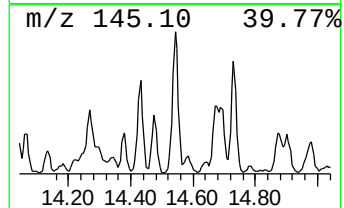
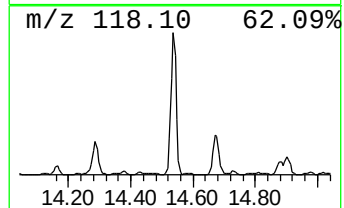
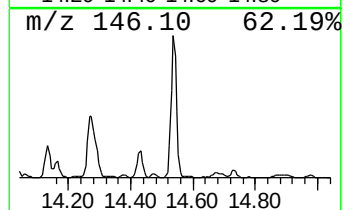
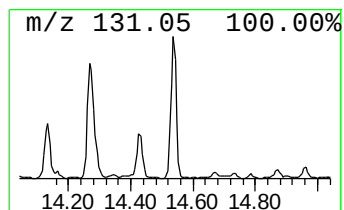
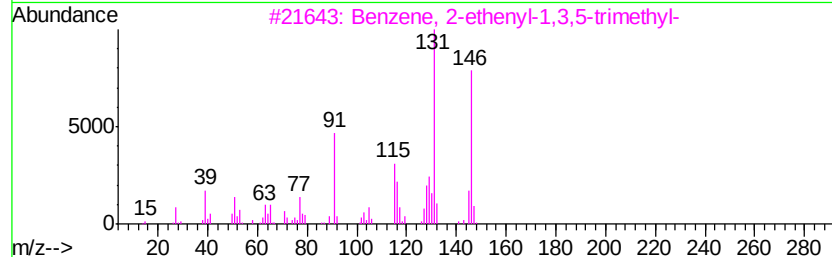
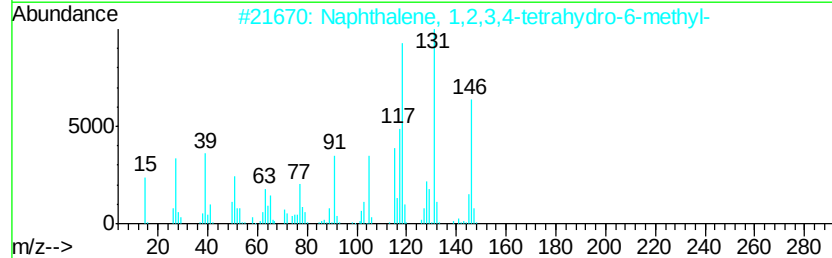
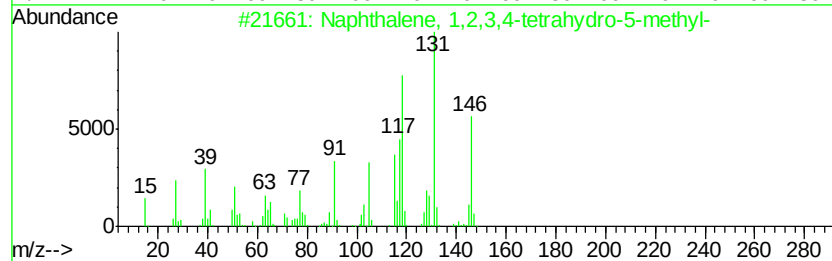
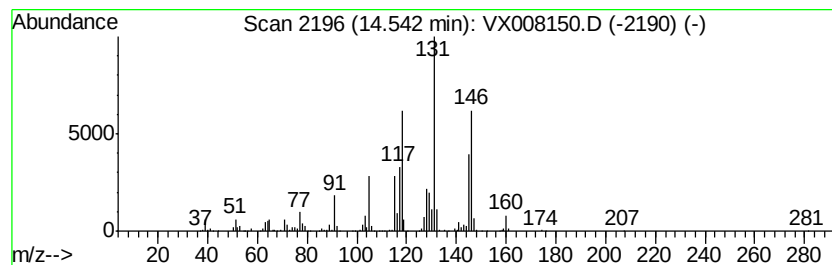
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Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1,2,3,4-tetra... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.54	20.27 ug/l	402855	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	95
2		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	94
3		Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	86
4		1H-1,3,2-Benzodiazaborole, 2-eth...	146	C8H11BN2	074663-81-3	83
5		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	70



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX031619\
Data File : VX008150.D
Acq On : 16 Mar 2019 19:20
Operator : JC/SP
Sample : K1960-06
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 18 Sample Multiplier: 1

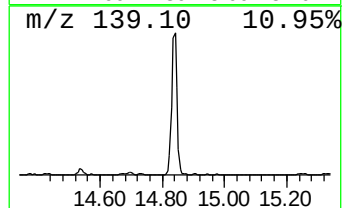
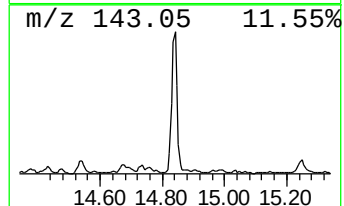
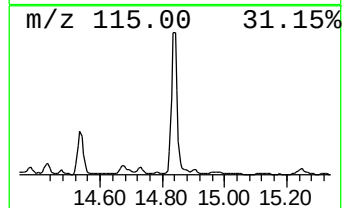
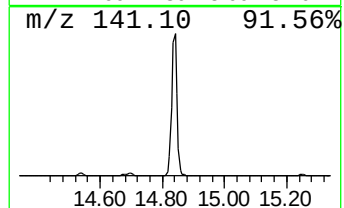
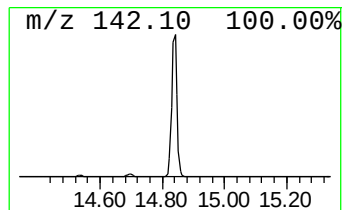
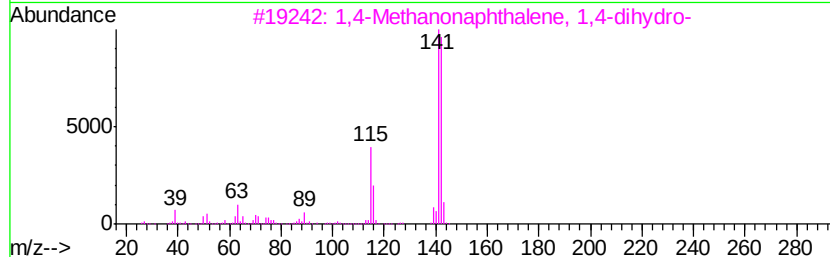
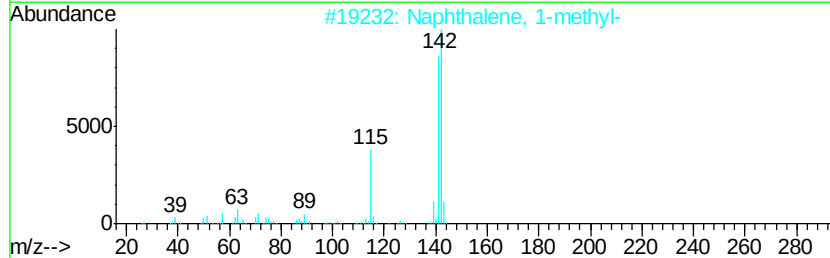
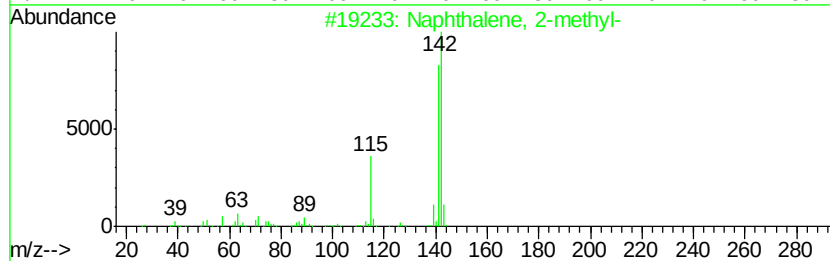
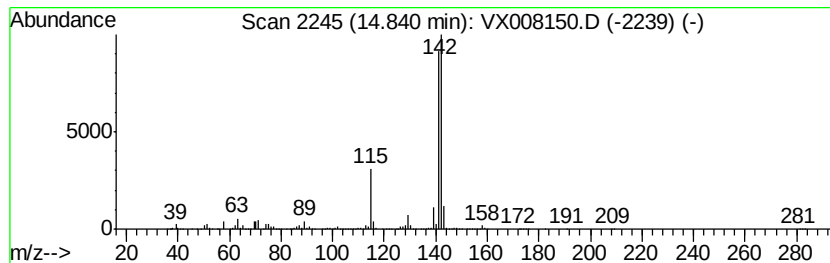
Instrument :
MSVOA_X
ClientSampleId :
T3-MW-3-9.0-031419

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X031419W.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.84	34.14 ug/l	678376	1,4-Dichlorobenzene-d4	12.08	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3	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	59



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX031619\
Data File : VX008150.D
Acq On : 16 Mar 2019 19:20
Operator : JC/SP
Sample : K1960-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
T3-MW-3-9.0-031419

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X031419W.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	65.4	ug/l	1299320	4	12.08	993597	50.0
Benzene, 4-ethyl-...	12.66	32.3	ug/l	640983	4	12.08	993597	50.0
Indan, 1-methyl-	12.70	20.7	ug/l	411402	4	12.08	993597	50.0
Benzene, 1,2,4,5-...	12.97	52.2	ug/l	1037140	4	12.08	993597	50.0
1H-Indene, 2,3-di...	13.35	46.7	ug/l	927696	4	12.08	993597	50.0
Benzene, (1,2-dim...	13.64	16.6	ug/l	330453	4	12.08	993597	50.0
Naphthalene, 1,2,...	14.54	20.3	ug/l	402855	4	12.08	993597	50.0
Naphthalene, 1-me...	14.84	34.1	ug/l	678376	4	12.08	993597	50.0