

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX082719\
 Data File : VX011865.D
 Acq On : 27 Aug 2019 21:09
 Operator : JC/SP
 Sample : K4566-03
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-1

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\82X082219W.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	1.782	99	103	110	rBV	86390	120646	2.82%	0.215%
2	2.599	230	237	249	rVB	61028	112391	2.62%	0.200%
3	2.842	269	277	280	rBV	27326	59430	1.39%	0.106%
4	2.928	287	291	298	rVB2	24115	48579	1.13%	0.087%
5	3.166	322	330	339	rBV	61016	139782	3.26%	0.249%
6	4.397	522	532	542	rVB2	61463	175305	4.09%	0.313%
7	5.494	699	712	720	rBV	203453	594582	13.88%	1.060%
8	5.574	720	725	729	rVV	39615	104372	2.44%	0.186%
9	5.653	729	738	759	rVB	368333	1117547	26.08%	1.993%
10	5.927	772	783	794	rVB4	13808	47560	1.11%	0.085%
11	6.055	794	804	816	rBV	214855	557551	13.01%	0.994%
12	6.250	825	836	845	rBV6	19485	73896	1.72%	0.132%
13	6.348	845	852	862	rVV4	26283	79242	1.85%	0.141%
14	6.854	925	935	948	rBV	524891	1261646	29.44%	2.250%
15	7.458	1022	1034	1042	rBV3	72509	186184	4.34%	0.332%
16	8.665	1225	1232	1234	rBV2	36347	62118	1.45%	0.111%
17	8.713	1234	1240	1250	rVB	1041440	1666837	38.90%	2.972%
18	9.037	1288	1293	1299	rVB	38464	62695	1.46%	0.112%
19	9.158	1305	1313	1319	rBV	29054	49060	1.14%	0.087%
20	9.622	1384	1389	1393	rVV	38456	51039	1.19%	0.091%
21	10.109	1463	1469	1476	rBV	1096679	1435639	33.50%	2.560%
22	10.640	1547	1556	1563	rBV3	22694	60994	1.42%	0.109%
23	10.835	1576	1588	1595	rBV2	38990	71738	1.67%	0.128%
24	10.908	1595	1600	1608	rBV3	30950	50644	1.18%	0.090%
25	11.012	1612	1617	1623	rBV	275891	365609	8.53%	0.652%
26	11.134	1631	1637	1642	rBV	983813	1187854	27.72%	2.118%
27	11.353	1668	1673	1681	rVV	549756	697846	16.29%	1.244%
28	11.670	1719	1725	1732	rVB3	34051	56068	1.31%	0.100%
29	11.914	1758	1765	1767	rBV	80460	116379	2.72%	0.208%
30	11.945	1767	1770	1775	rVB	262088	317856	7.42%	0.567%
31	12.073	1786	1791	1797	rBV	1170832	1442323	33.66%	2.572%
32	12.231	1811	1817	1821	rVB	82020	108172	2.52%	0.193%
33	12.298	1822	1828	1833	rVV	1737821	2179350	50.86%	3.886%
34	12.365	1834	1839	1847	rVB2	250970	443659	10.35%	0.791%

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35	12.457	1848	1854	1860	rBV	245557	301852	7.04%	0.538%
36	12.664	1880	1888	1891	rBV	745235	918276	21.43%	1.637%
37	12.694	1891	1893	1898	rVV	411517	514774	12.01%	0.918%
38	12.749	1898	1902	1910	rVV2	1004637	1522249	35.52%	2.714%
39	12.810	1910	1912	1919	rVV3	49004	99775	2.33%	0.178%
40	12.889	1919	1925	1930	rVB2	174586	267162	6.23%	0.476%
41	12.975	1930	1939	1944	rBV	1148334	1457210	34.01%	2.598%
42	13.024	1944	1947	1951	rVB2	39070	50775	1.18%	0.091%
43	13.078	1951	1956	1963	rVB3	207055	306572	7.15%	0.547%
44	13.146	1963	1967	1971	rBV2	116473	195942	4.57%	0.349%
45	13.188	1971	1974	1977	rVV2	331397	454577	10.61%	0.811%
46	13.219	1977	1979	1989	rVV	538594	777108	18.14%	1.386%
47	13.310	1989	1994	1995	rVV2	124664	156107	3.64%	0.278%
48	13.347	1995	2000	2005	rVV2	2404798	3461350	80.78%	6.172%
49	13.402	2005	2009	2010	rVV2	121657	153992	3.59%	0.275%
50	13.420	2010	2012	2015	rVV	163831	198101	4.62%	0.353%
51	13.475	2018	2021	2029	rVB2	108241	158076	3.69%	0.282%
52	13.560	2030	2035	2037	rBV	251958	322952	7.54%	0.576%
53	13.597	2037	2041	2044	rVV	524324	747036	17.43%	1.332%
54	13.633	2044	2047	2051	rVV2	729209	1107990	25.86%	1.976%
55	13.694	2051	2057	2058	rVV2	443648	707508	16.51%	1.262%
56	13.719	2058	2061	2069	rVV2	787929	1187496	27.71%	2.117%
57	13.804	2069	2075	2078	rVV	135864	175178	4.09%	0.312%
58	13.847	2078	2082	2087	rVV2	159544	235143	5.49%	0.419%
59	13.908	2087	2092	2097	rVB	720726	876710	20.46%	1.563%
60	13.981	2097	2104	2108	rBV2	866171	1215886	28.37%	2.168%
61	14.023	2108	2111	2115	rVV	372530	489167	11.42%	0.872%
62	14.060	2115	2117	2123	rVV2	81844	105246	2.46%	0.188%
63	14.127	2123	2128	2132	rVV	643858	827594	19.31%	1.476%
64	14.158	2132	2133	2140	rVV3	91661	145399	3.39%	0.259%
65	14.212	2140	2142	2146	rVV2	62072	96834	2.26%	0.173%
66	14.267	2146	2151	2161	rVV2	818442	1538046	35.89%	2.742%
67	14.371	2165	2168	2173	rVV	430239	580953	13.56%	1.036%
68	14.426	2173	2177	2181	rVV	687300	878278	20.50%	1.566%
69	14.475	2181	2185	2190	rVB	124875	170596	3.98%	0.304%
70	14.536	2190	2195	2200	rBV2	1161041	1515991	35.38%	2.703%
71	14.578	2200	2202	2208	rVV2	71042	108678	2.54%	0.194%

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72	14.694	2208	2221	2224	rVV2	1369990	2541454	59.31%	4.532%
73	14.725	2224	2226	2231	rVV	343936	437223	10.20%	0.780%
74	14.779	2231	2235	2239	rVV2	155757	212018	4.95%	0.378%
75	14.834	2239	2244	2248	rVV	3443891	4285076	100.00%	7.641%
76	14.871	2248	2250	2253	rVV2	237552	318212	7.43%	0.567%
77	14.901	2253	2255	2259	rVV2	179836	212121	4.95%	0.378%
78	14.956	2259	2264	2271	rVV4	161727	357382	8.34%	0.637%
79	15.023	2271	2275	2281	rVB4	39988	94869	2.21%	0.169%
80	15.109	2282	2289	2293	rBV2	82370	125623	2.93%	0.224%
81	15.243	2305	2311	2319	rBV	455677	760276	17.74%	1.356%
82	15.316	2319	2323	2330	rVB2	111285	189375	4.42%	0.338%
83	15.438	2337	2343	2346	rBV2	208319	334855	7.81%	0.597%
84	15.468	2346	2348	2353	rVV	336422	440228	10.27%	0.785%
85	15.541	2353	2360	2365	rVV	570130	955316	22.29%	1.703%
86	15.584	2365	2367	2373	rVV4	139733	230284	5.37%	0.411%
87	15.682	2373	2383	2387	rVV	978792	1634957	38.15%	2.915%
88	15.718	2387	2389	2397	rVB	512696	681114	15.90%	1.214%
89	15.804	2397	2403	2410	rBV3	85087	193246	4.51%	0.345%
90	15.883	2410	2416	2418	rVV	294622	488135	11.39%	0.870%
91	15.913	2418	2421	2431	rVB	418187	729844	17.03%	1.301%
92	16.060	2439	2445	2457	rVB	446671	776360	18.12%	1.384%
93	16.243	2471	2475	2477	rBV2	29800	47570	1.11%	0.085%
94	16.273	2477	2480	2486	rVB2	60265	97004	2.26%	0.173%
95	16.340	2486	2491	2494	rBV	46519	82230	1.92%	0.147%
96	16.413	2499	2503	2505	rBV	37959	62254	1.45%	0.111%
97	16.560	2522	2527	2532	rBV	60082	101477	2.37%	0.181%
98	16.645	2532	2541	2543	rBV5	45172	125224	2.92%	0.223%
99	16.681	2543	2547	2552	rVV2	92140	169215	3.95%	0.302%
100	16.742	2552	2557	2564	rVV2	136957	268612	6.27%	0.479%

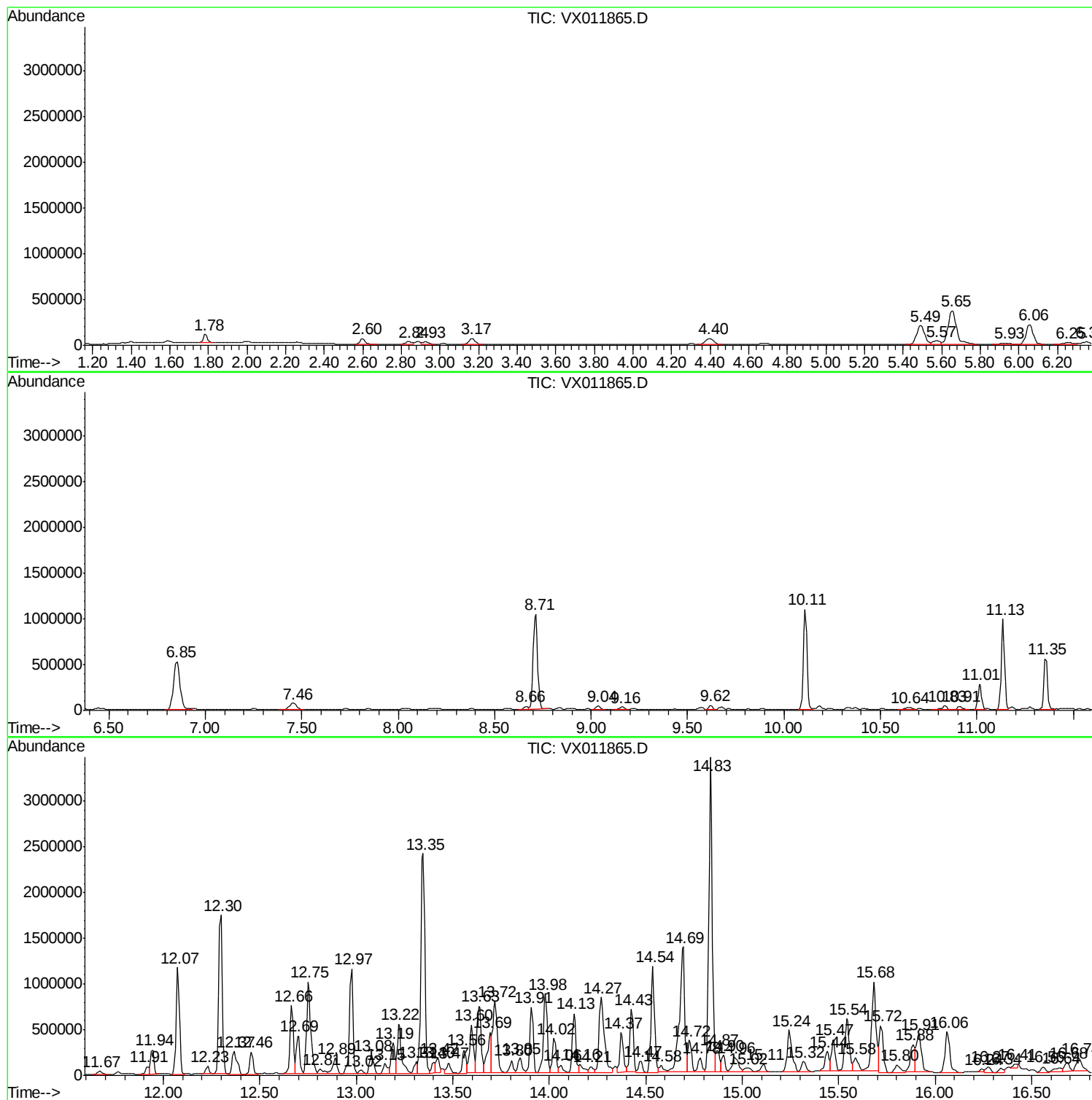
Sum of corrected areas: 56082726

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X082219W.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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ClientSampled :
MW-1

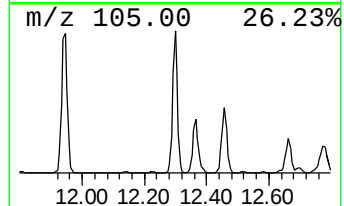
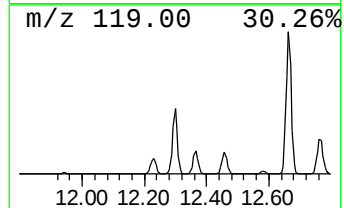
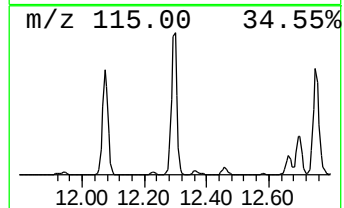
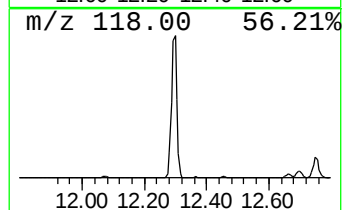
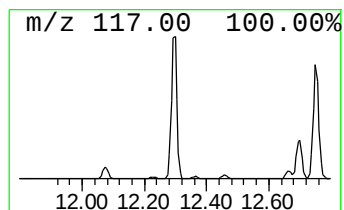
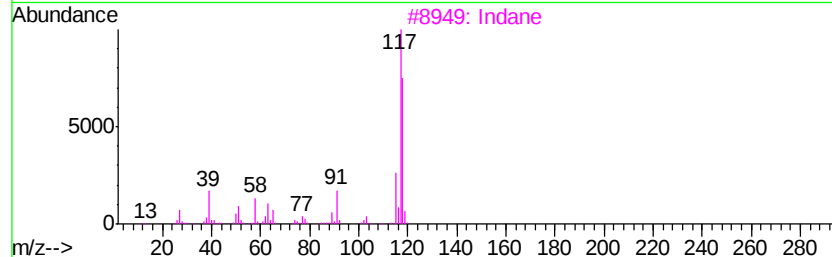
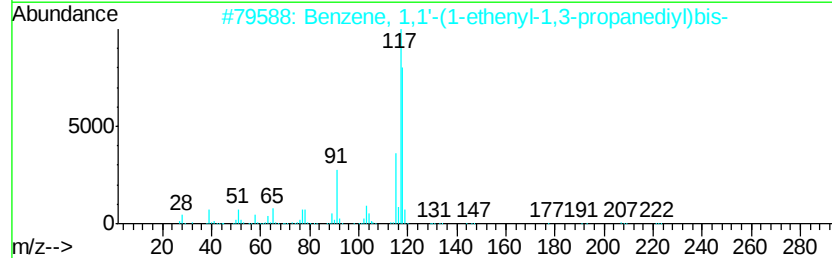
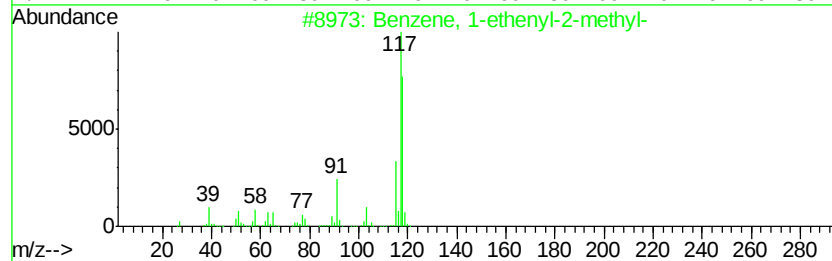
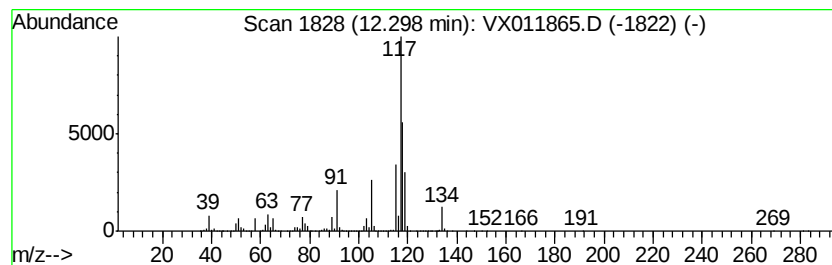
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X082219W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	75.55 ug/l	2179350	1,4-Dichlorobenzene-d4	12.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	91
2		Benzene, 1,1'-(1-ethenyl-1,3-pro...	222	C17H18	061141-97-7	64
3		Indane	118	C9H10	000496-11-7	64
4		Benzene, cyclopropyl-	118	C9H10	000873-49-4	64
5		Benzene, 2-propenyl-	118	C9H10	000300-57-2	64



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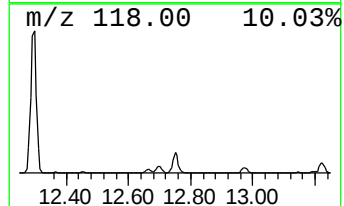
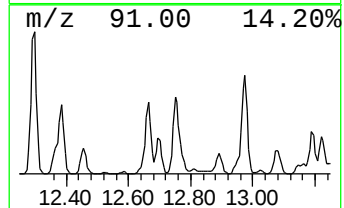
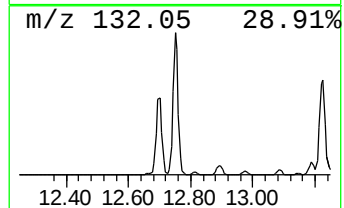
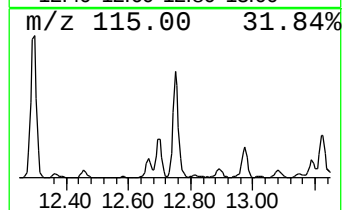
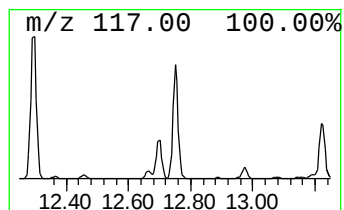
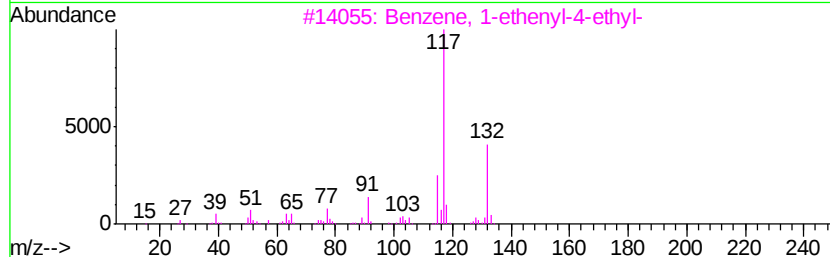
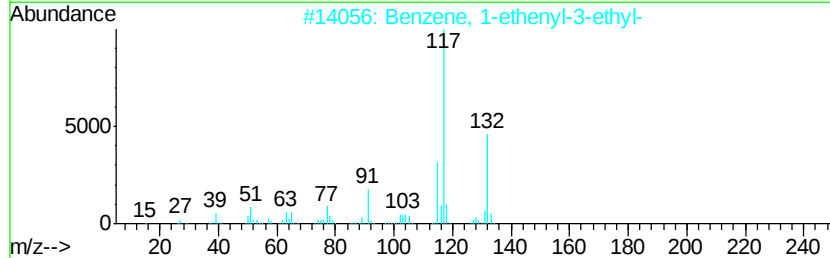
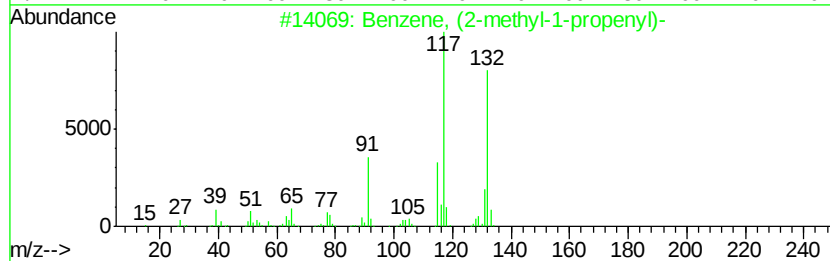
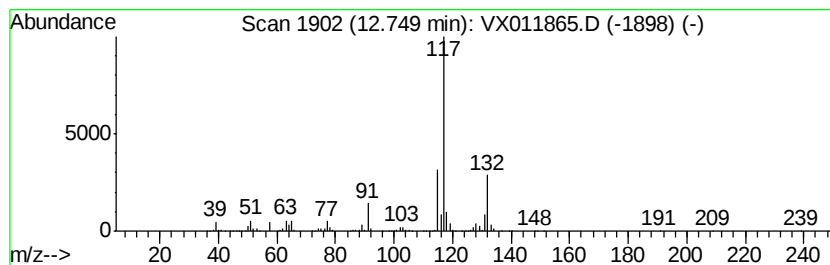
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzene, (2-methyl-1-propen... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.75	52.77 ug/l	1522250	1,4-Dichlorobenzene-d4	12.07

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	90
2		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	90
3		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87
4		Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	87
5		Indan, 1-methyl-	132	C10H12	000767-58-8	87



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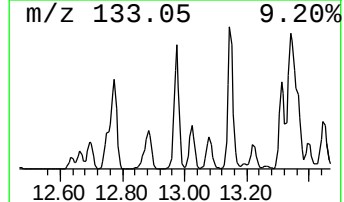
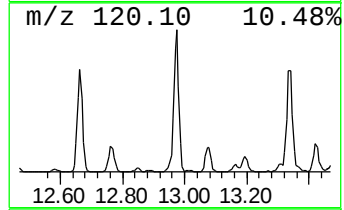
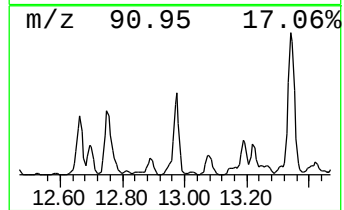
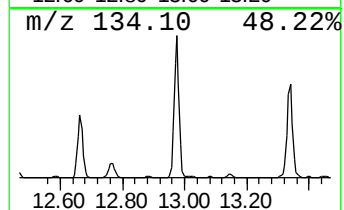
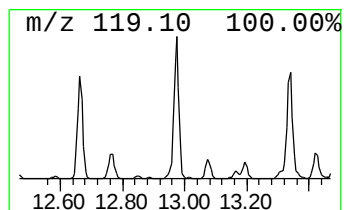
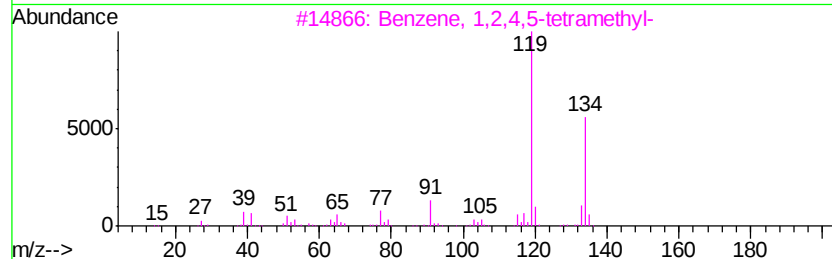
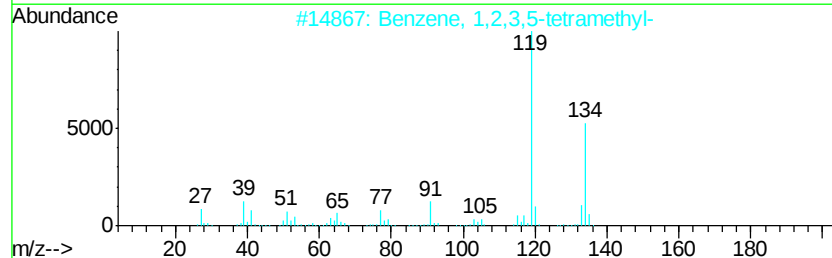
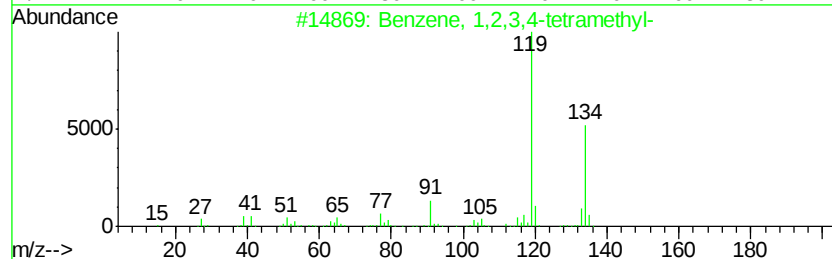
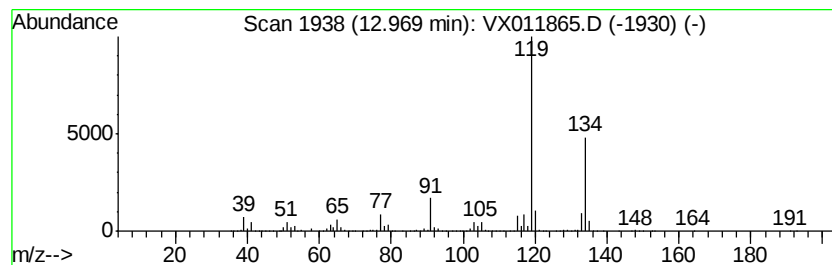
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.97	50.52 ug/l	1457210	1,4-Dichlorobenzene-d4	12.07

Hit# of	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
4	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94
5	o-Cymene	134	C10H14	000527-84-4	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX082719\
Data File : VX011865.D
Acq On : 27 Aug 2019 21:09
Operator : JC/SP
Sample : K4566-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-1

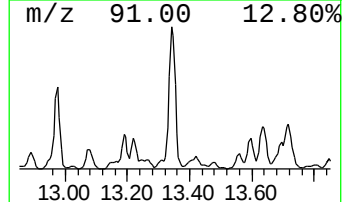
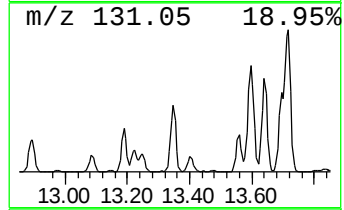
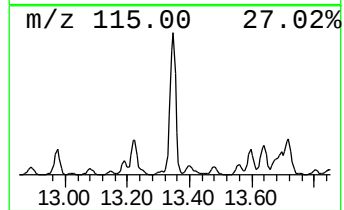
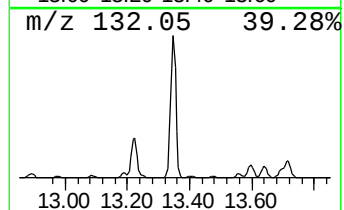
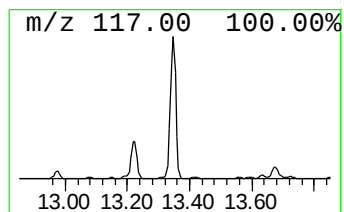
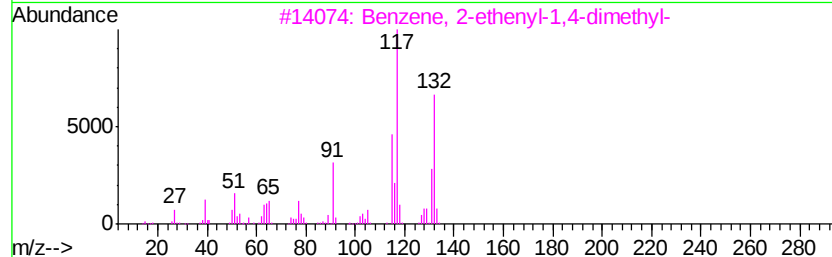
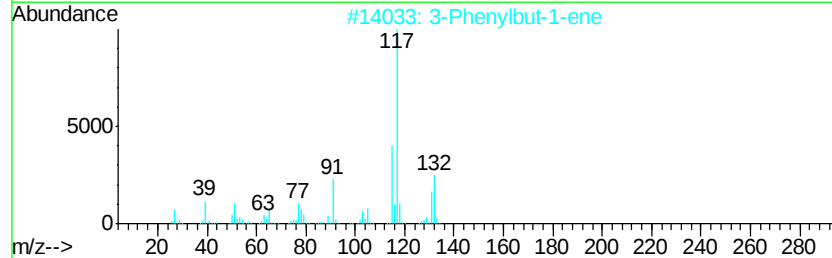
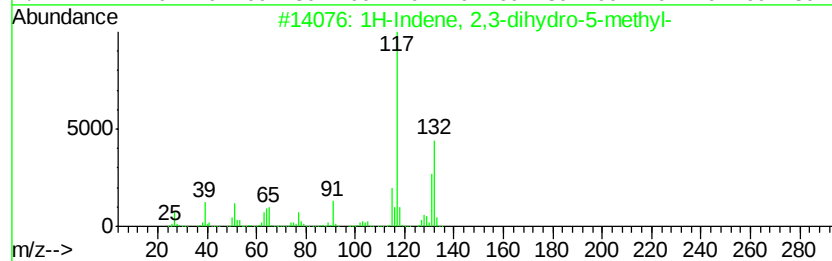
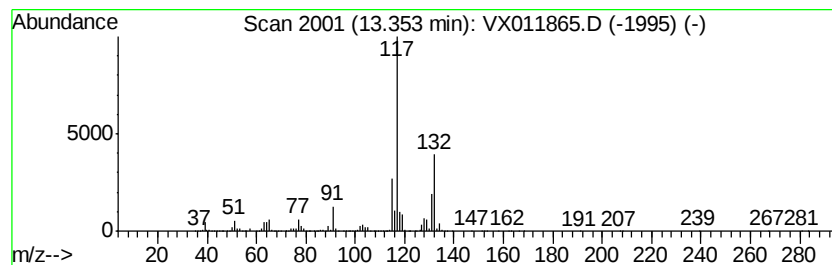
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Quant Title : SW846 8260

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

Peak Number 4 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	119.99 ug/l	3461350	1,4-Dichlorobenzene-d4	12.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	81
2		3-Phenylbut-1-ene	132	C10H12	000934-10-1	81
3		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	78
4		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	76
5		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	76



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX082719\
Data File : VX011865.D
Acq On : 27 Aug 2019 21:09
Operator : JC/SP
Sample : K4566-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-1

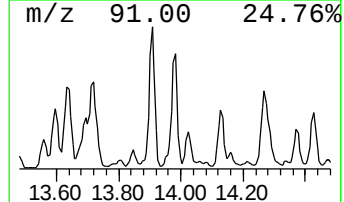
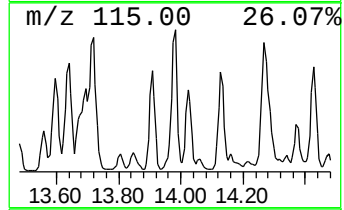
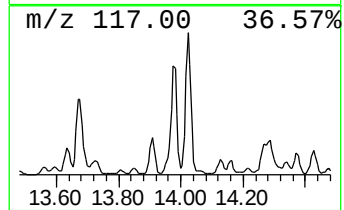
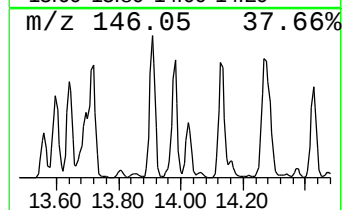
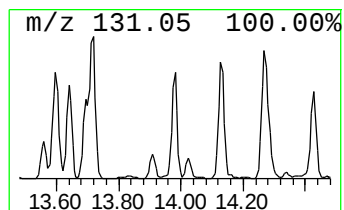
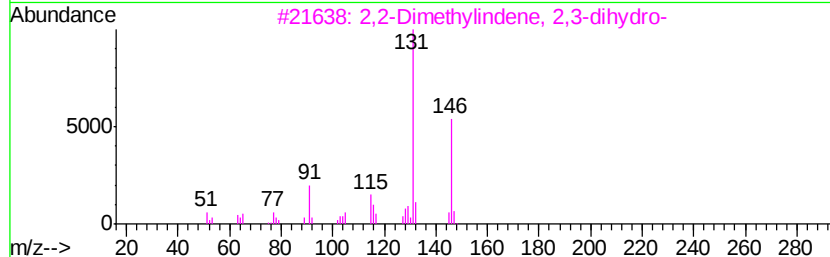
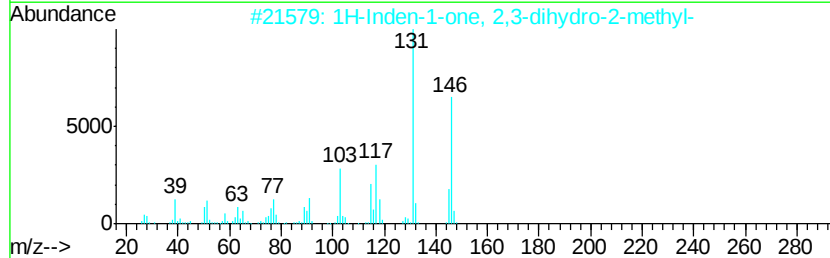
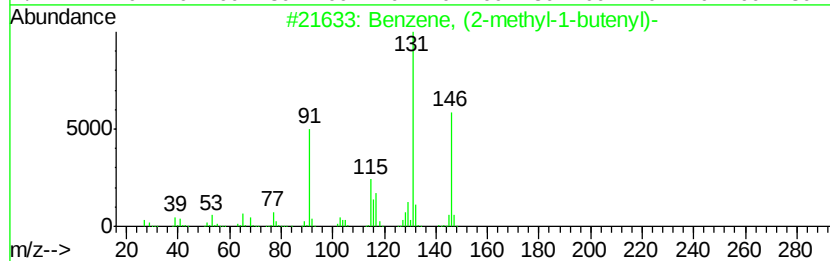
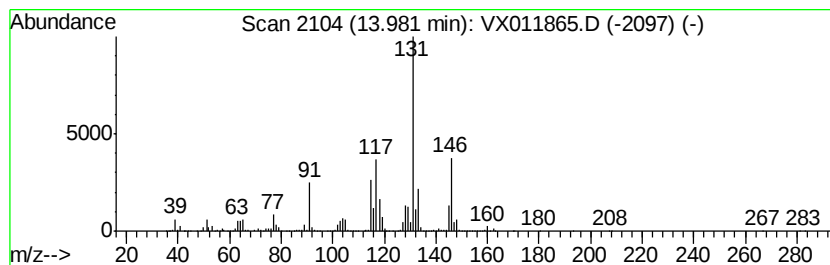
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, (2-methyl-1-butenyl)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.98	42.15 ug/l	1215890	1,4-Dichlorobenzene-d4	12.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	94
2		1H-Inden-1-one, 2,3-dihydro-2-methyl-	146	C10H10O	017496-14-9	90
3		2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	89
4		Naphthalene, 1,2,3,4-tetrahydro-	146	C11H14	001559-81-5	81
5		1H-Indene, 2,3-dihydro-4,7-dimethyl-	146	C11H14	006682-71-9	76



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX082719\
Data File : VX011865.D
Acq On : 27 Aug 2019 21:09
Operator : JC/SP
Sample : K4566-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-1

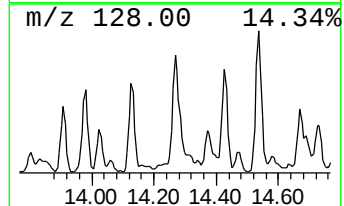
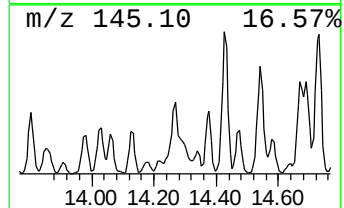
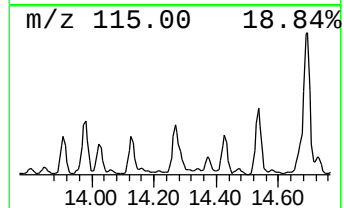
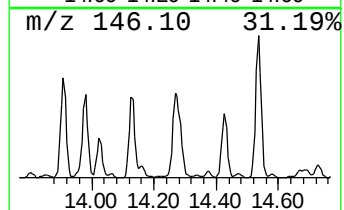
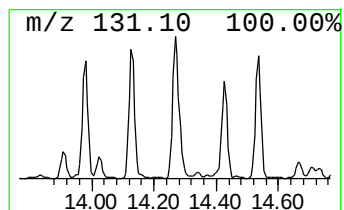
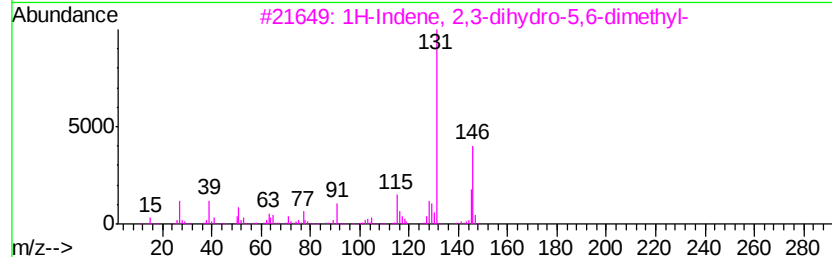
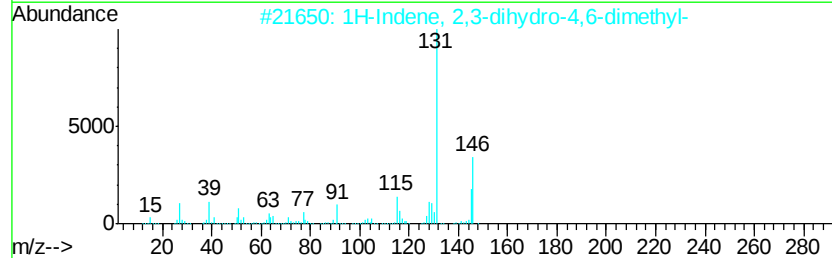
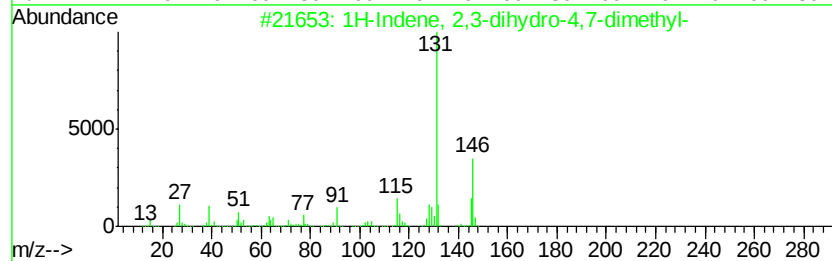
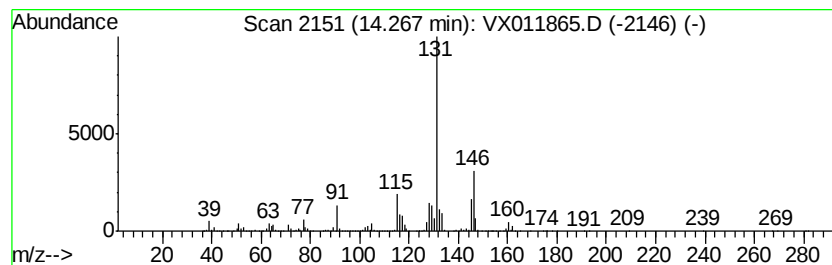
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Quant Title : SW846 8260

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

Peak Number 6 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.27	53.32 ug/l	1538050	1,4-Dichlorobenzene-d4	12.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	96
2		1H-Indene, 2,3-dihydro-4,6-dimet...	146	C11H14	001685-82-1	94
3		1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	91
4		Benzene, 1-methyl-3-(1-methyl-2-...	146	C11H14	052161-57-6	90
5		1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX082719\
Data File : VX011865.D
Acq On : 27 Aug 2019 21:09
Operator : JC/SP
Sample : K4566-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-1

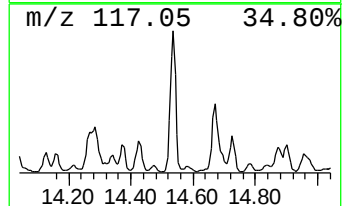
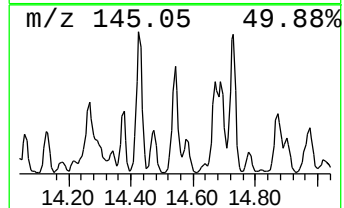
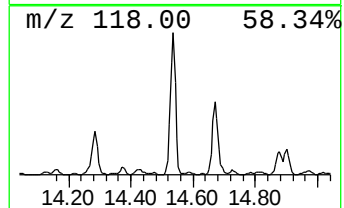
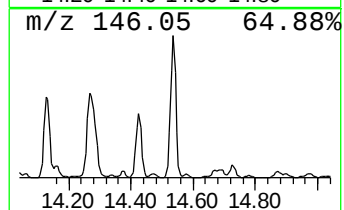
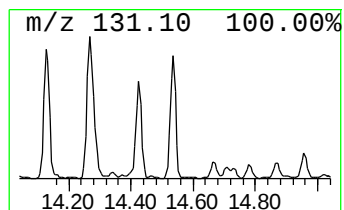
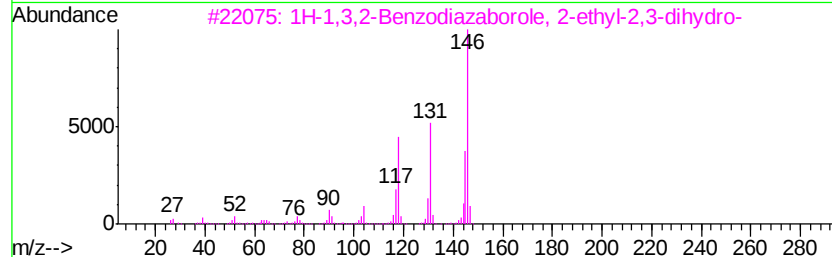
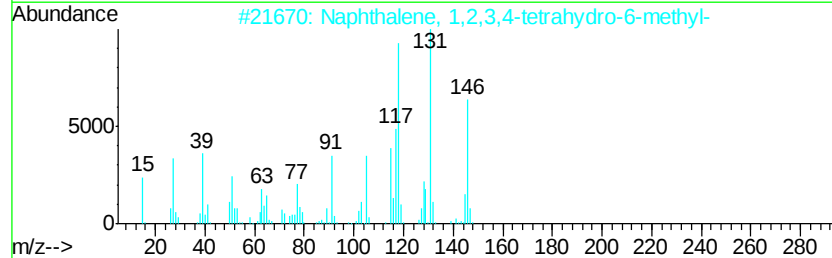
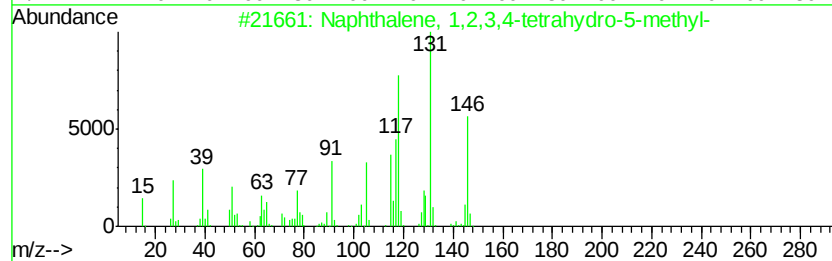
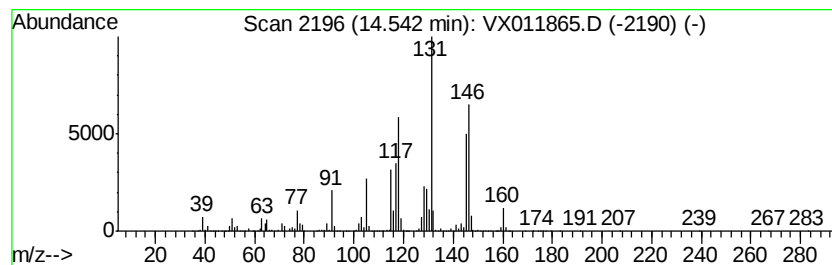
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Quant Title : SW846 8260

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

Peak Number 7 Naphthalene, 1,2,3,4-tetra... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.54	52.55 ug/l	1515990	1,4-Dichlorobenzene-d4	12.07

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	95
3		1H-1,3,2-Benzodiazaborole, 2-eth...	146	C8H11BN2	074663-81-3	90
4		Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	64
5		Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	64



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX082719\
Data File : VX011865.D
Acq On : 27 Aug 2019 21:09
Operator : JC/SP
Sample : K4566-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-1

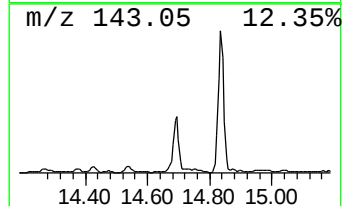
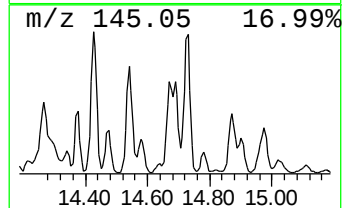
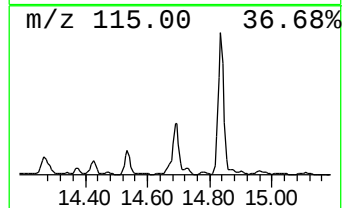
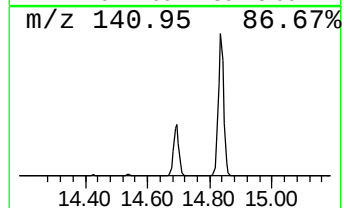
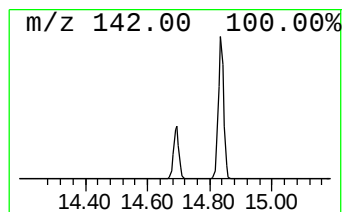
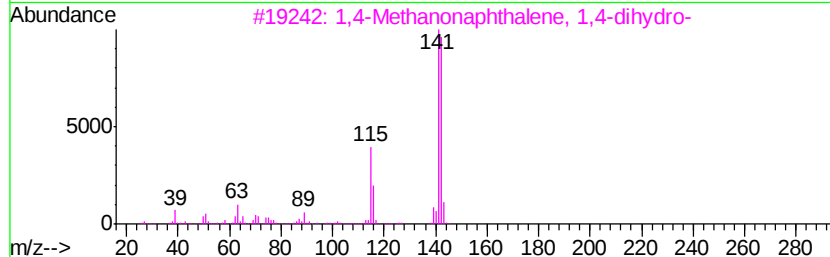
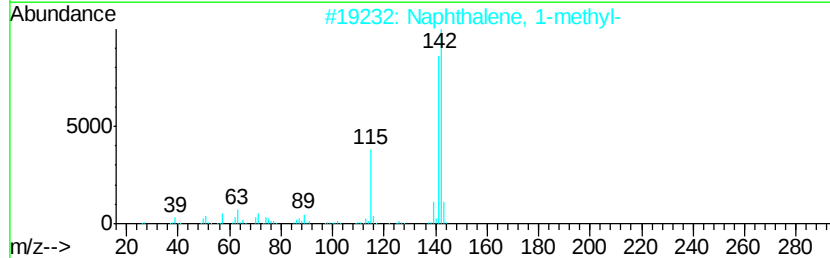
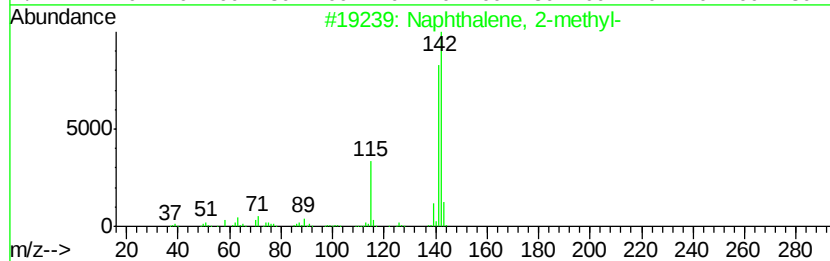
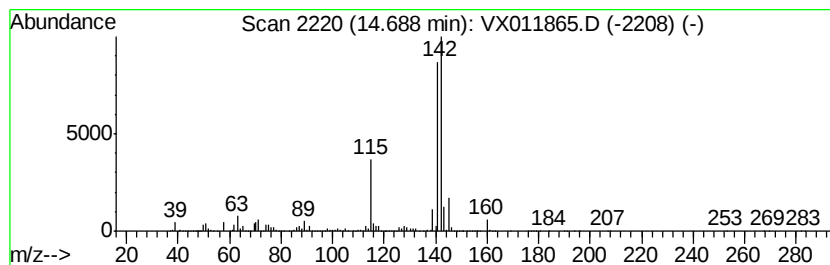
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Quant Title : SW846 8260

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

Peak Number 8 Naphthalene, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.69	88.10 ug/l	2541450	1,4-Dichlorobenzene-d4	12.07

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	90
4		Benzocycloheptatriene	142	C11H10	000264-09-5	90
5		Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	59



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX082719\
Data File : VX011865.D
Acq On : 27 Aug 2019 21:09
Operator : JC/SP
Sample : K4566-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
MW-1

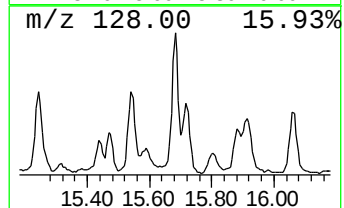
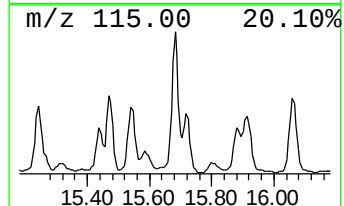
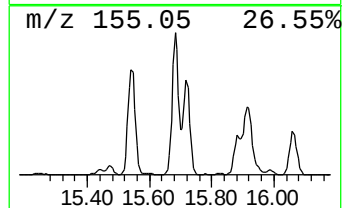
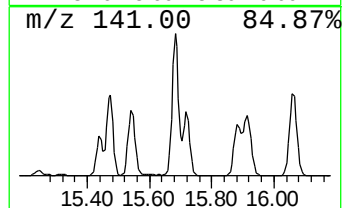
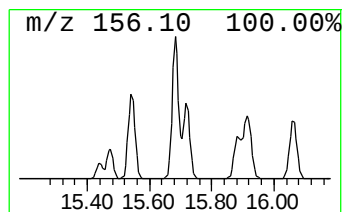
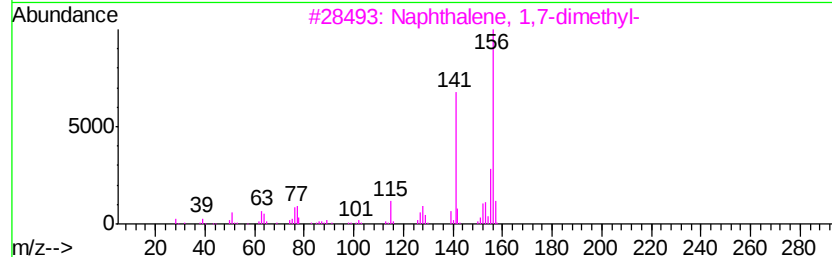
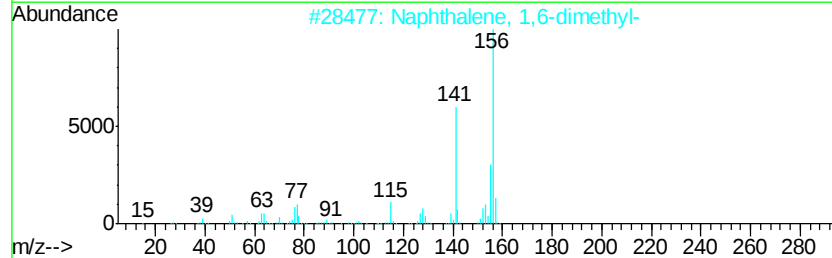
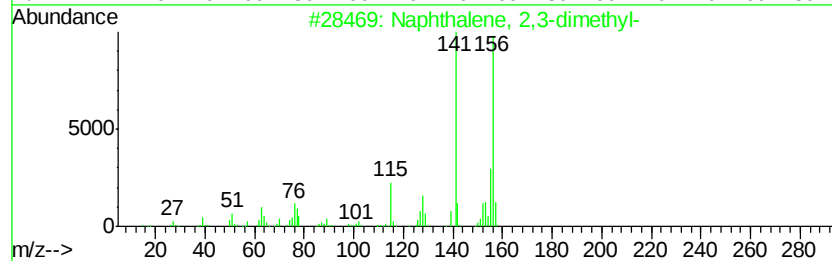
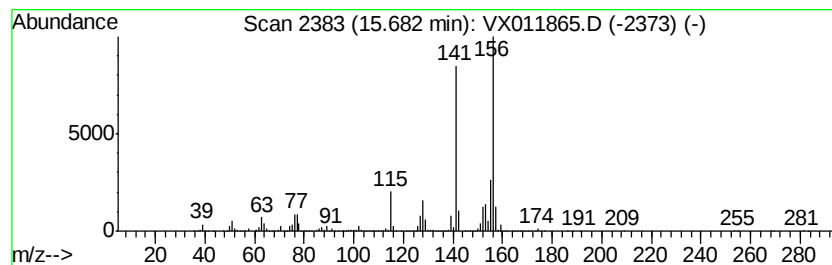
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Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 2,3-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.68	56.68 ug/l	1634960	1,4-Dichlorobenzene-d4	12.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
2		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
3		Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
4		Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	97
5		Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	96



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX082719\
Data File : VX011865.D
Acq On : 27 Aug 2019 21:09
Operator : JC/SP
Sample : K4566-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-1

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X082219W.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-etheny...	12.30	75.5	ug/l	2179350	4	12.07	1442320	50.0
Benzene, (2-methy...	12.75	52.8	ug/l	1522250	4	12.07	1442320	50.0
Benzene, 1,2,3,4-...	12.97	50.5	ug/l	1457210	4	12.07	1442320	50.0
1H-Indene, 2,3-di...	13.35	120.0	ug/l	3461350	4	12.07	1442320	50.0
Benzene, (2-methy...	13.98	42.1	ug/l	1215890	4	12.07	1442320	50.0
1H-Indene, 2,3-di...	14.27	53.3	ug/l	1538050	4	12.07	1442320	50.0
Naphthalene, 1,2,...	14.54	52.5	ug/l	1515990	4	12.07	1442320	50.0
Naphthalene, 1-me...	14.69	88.1	ug/l	2541450	4	12.07	1442320	50.0
Naphthalene, 2,3-...	15.68	56.7	ug/l	1634960	4	12.07	1442320	50.0