

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX031523\
 Data File : VX034511.D
 Acq On : 15 Mar 2023 16:22
 Operator : JC/MD
 Sample : 01892-01 10X
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BOOM-030923

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X031023W.M
 Title : SW846 8260

Signal : TIC: VX034511.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.569	70	79	84	rBV3	9682	22964	1.19%	0.070%
2	5.385	693	705	721	rBV	144482	424351	22.05%	1.298%
3	5.550	721	732	747	rVB	245700	703869	36.57%	2.153%
4	5.952	787	798	808	rBV	168692	454623	23.62%	1.391%
5	6.037	809	812	825	rVB2	9852	25711	1.34%	0.079%
6	6.757	916	930	945	rBV	394791	956994	49.73%	2.927%
7	8.647	1233	1240	1248	rBV	857700	1323475	68.77%	4.049%
8	8.720	1248	1252	1259	rVB	28439	46396	2.41%	0.142%
9	10.055	1465	1471	1487	rVB	813791	1127241	58.57%	3.448%
10	10.195	1487	1494	1500	rBV	30317	39960	2.08%	0.122%
11	10.299	1503	1511	1517	rVV	104616	148111	7.70%	0.453%
12	10.640	1562	1567	1577	rVB	79519	112432	5.84%	0.344%
13	10.963	1614	1620	1624	rBV	15758	20134	1.05%	0.062%
14	11.079	1634	1639	1652	rBV	770756	968793	50.34%	2.964%
15	11.305	1671	1676	1683	rBV	46823	61407	3.19%	0.188%
16	11.378	1683	1688	1695	rVV	191249	338841	17.61%	1.037%
17	11.451	1695	1700	1702	rVV	106600	142572	7.41%	0.436%
18	11.476	1702	1704	1711	rVB	68452	81750	4.25%	0.250%
19	11.610	1722	1726	1734	rVB	125988	178868	9.29%	0.547%
20	11.713	1740	1743	1745	rBV	19951	22740	1.18%	0.070%
21	11.750	1745	1749	1760	rVB	537520	644174	33.47%	1.971%
22	11.890	1765	1772	1776	rVB2	39180	59549	3.09%	0.182%
23	11.945	1776	1781	1782	rBV2	17910	22129	1.15%	0.068%
24	11.969	1782	1785	1788	rVB	52703	55700	2.89%	0.170%
25	12.018	1788	1793	1799	rBV	835211	1084339	56.34%	3.317%
26	12.085	1799	1804	1811	rVB	250340	338205	17.57%	1.035%
27	12.171	1815	1818	1825	rVB3	30326	47579	2.47%	0.146%
28	12.244	1825	1830	1832	rBV	250965	329458	17.12%	1.008%
29	12.268	1832	1834	1838	rVB	171395	172784	8.98%	0.529%
30	12.317	1838	1842	1849	rVB3	281187	432587	22.48%	1.323%
31	12.420	1849	1859	1863	rBV2	234623	333436	17.33%	1.020%
32	12.463	1863	1866	1872	rVB	132786	164888	8.57%	0.504%
33	12.530	1872	1877	1879	rBV	191223	247074	12.84%	0.756%
34	12.555	1879	1881	1885	rVB	169969	170360	8.85%	0.521%
35	12.609	1886	1890	1894	rBV	234126	278151	14.45%	0.851%
36	12.646	1894	1896	1899	rVB	77918	76638	3.98%	0.234%

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37	12.701	1899	1905	1908	rBV	199213	326251	16.95%	0.998%
38	12.798	1917	1921	1922	rBV2	37812	45756	2.38%	0.140%
39	12.847	1926	1929	1933	rVB	103101	116872	6.07%	0.358%
40	12.920	1933	1941	1944	rBV2	221587	350797	18.23%	1.073%
41	12.963	1944	1948	1954	rVB	364522	514226	26.72%	1.573%
42	13.042	1954	1961	1964	rBV4	124555	235269	12.22%	0.720%
43	13.073	1964	1966	1968	rVB	42324	37529	1.95%	0.115%
44	13.164	1974	1981	1986	rBV2	427645	632726	32.88%	1.936%
45	13.213	1986	1989	1992	rVB	94220	98387	5.11%	0.301%
46	13.268	1992	1998	1999	rBV3	401975	545668	28.35%	1.669%
47	13.292	1999	2002	2007	rVV2	1049110	1343600	69.81%	4.110%
48	13.335	2007	2009	2012	rVV2	47613	62155	3.23%	0.190%
49	13.384	2012	2017	2019	rVV2	81313	147008	7.64%	0.450%
50	13.420	2019	2023	2032	rVB	992931	1312873	68.22%	4.016%
51	13.506	2032	2037	2040	rBV2	144999	215085	11.18%	0.658%
52	13.542	2040	2043	2046	rVV	222539	264221	13.73%	0.808%
53	13.585	2046	2050	2054	rVV3	279641	460824	23.94%	1.410%
54	13.640	2056	2059	2060	rVV	100986	129290	6.72%	0.396%
55	13.664	2060	2063	2071	rVV2	279201	443352	23.04%	1.356%
56	13.737	2072	2075	2077	rVV2	48509	56452	2.93%	0.173%
57	13.774	2077	2081	2088	rVB	259607	355259	18.46%	1.087%
58	13.853	2088	2094	2099	rVB2	466555	730670	37.97%	2.235%
59	13.926	2099	2106	2110	rBV2	456606	708131	36.79%	2.166%
60	13.969	2110	2113	2117	rVB	142859	150263	7.81%	0.460%
61	14.006	2117	2119	2121	rBV2	29931	29066	1.51%	0.089%
62	14.036	2121	2124	2127	rBV	393935	424083	22.04%	1.297%
63	14.073	2127	2130	2134	rVV	358914	390521	20.29%	1.195%
64	14.164	2142	2145	2147	rBV	35086	28671	1.49%	0.088%
65	14.231	2149	2156	2162	rVB	1191684	1924554	100.00%	5.887%
66	14.316	2167	2170	2175	rBV	128319	176009	9.15%	0.538%
67	14.371	2175	2179	2183	rVV2	343571	406003	21.10%	1.242%
68	14.414	2183	2186	2192	rVB5	83537	152561	7.93%	0.467%
69	14.481	2192	2197	2207	rBV2	813117	1213800	63.07%	3.713%
70	14.566	2207	2211	2215	rVB3	59515	97589	5.07%	0.299%
71	14.615	2215	2219	2220	rBV	513403	566309	29.43%	1.732%
72	14.634	2220	2222	2226	rVV	584196	758462	39.41%	2.320%
73	14.670	2226	2228	2233	rVB	239455	283302	14.72%	0.867%
74	14.725	2233	2237	2239	rBV2	96871	131025	6.81%	0.401%
75	14.755	2239	2242	2244	rVV	505797	551126	28.64%	1.686%
76	14.780	2244	2246	2250	rVV	287224	380825	19.79%	1.165%

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77	14.816	2250	2252	2255	rVV2	143076	206745	10.74%	0.632%
78	14.847	2255	2257	2260	rVV3	151547	162953	8.47%	0.498%
79	14.902	2263	2266	2273	rVB3	127488	217822	11.32%	0.666%
80	15.048	2284	2290	2294	rBV3	70471	127180	6.61%	0.389%
81	15.109	2294	2300	2302	rBV5	48215	92518	4.81%	0.283%
82	15.182	2307	2312	2314	rVV	359581	503143	26.14%	1.539%
83	15.207	2314	2316	2320	rVV2	282773	398026	20.68%	1.218%
84	15.243	2320	2322	2324	rVV3	59959	74638	3.88%	0.228%
85	15.274	2324	2327	2331	rVB4	90866	135025	7.02%	0.413%
86	15.377	2339	2344	2348	rBV2	109065	176942	9.19%	0.541%
87	15.487	2356	2362	2367	rBV2	491356	820497	42.63%	2.510%
88	15.572	2373	2376	2378	rBV3	26594	37549	1.95%	0.115%
89	15.615	2378	2383	2386	rVV2	191928	316055	16.42%	0.967%
90	15.652	2386	2389	2397	rVB2	114866	216723	11.26%	0.663%
91	15.719	2397	2400	2402	rBV4	17485	25884	1.34%	0.079%
92	15.841	2417	2420	2424	rBV2	36825	52689	2.74%	0.161%
93	15.993	2441	2445	2448	rBV5	32431	51965	2.70%	0.159%
94	16.091	2456	2461	2469	rBV4	67607	183919	9.56%	0.563%
95	16.377	2502	2508	2517	rVB	248151	434698	22.59%	1.330%

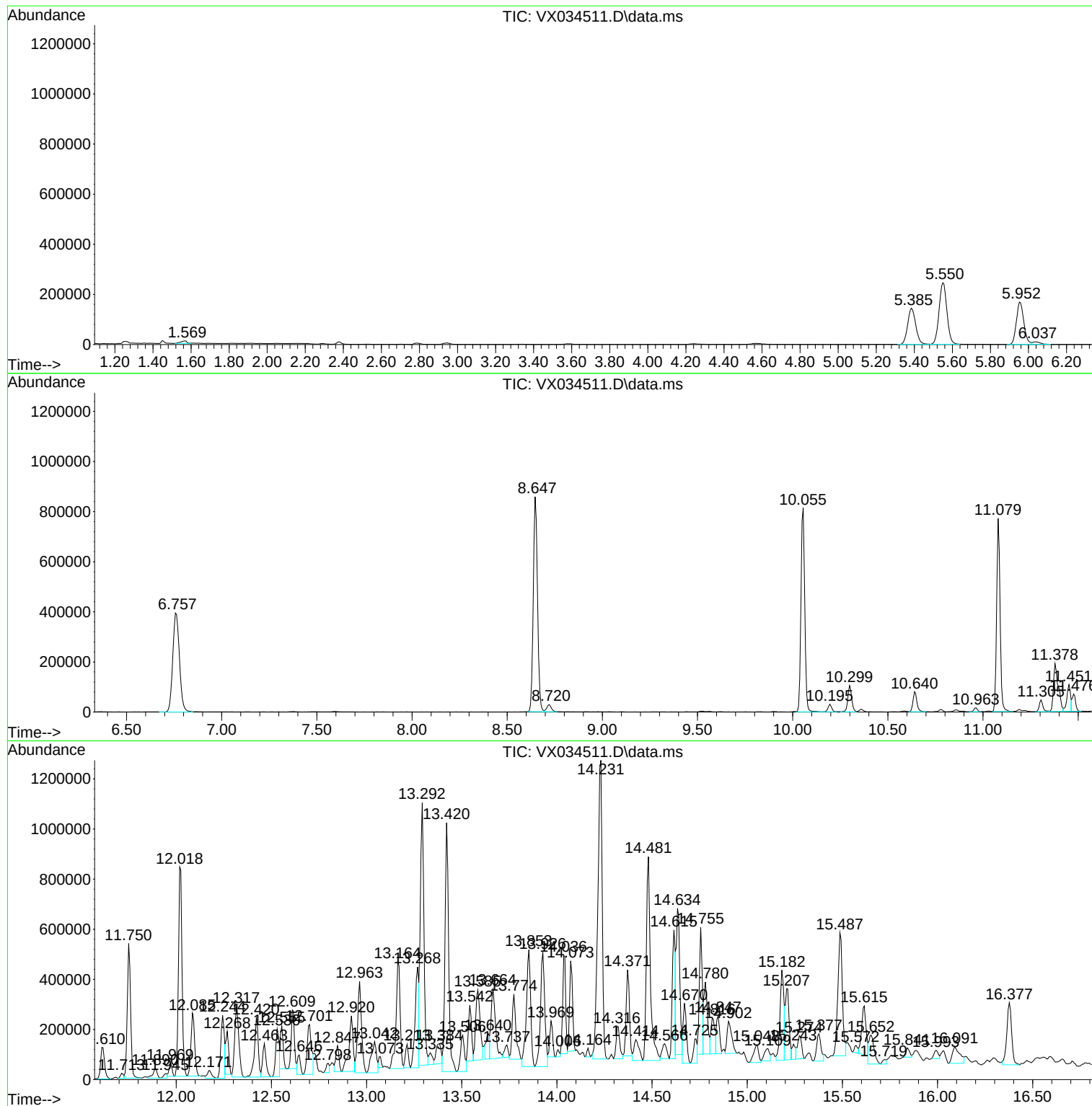
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X031023W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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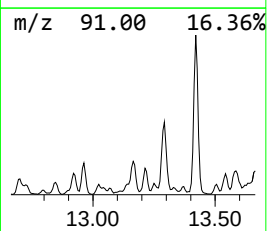
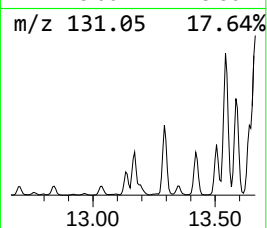
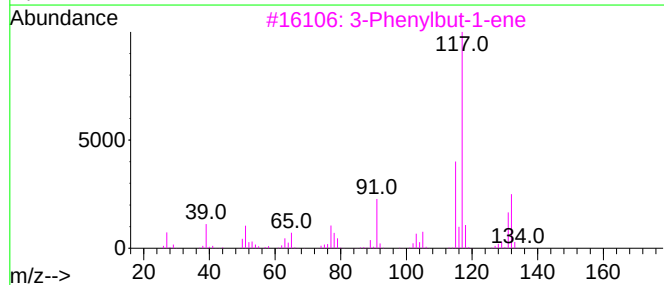
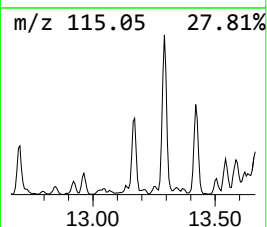
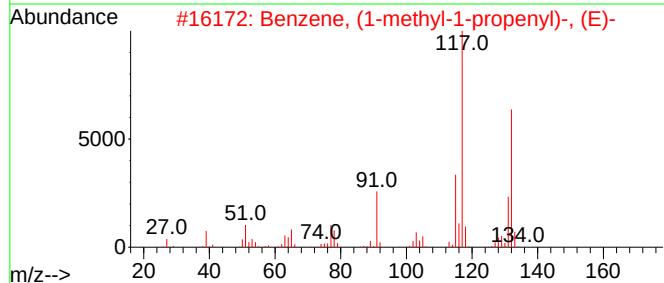
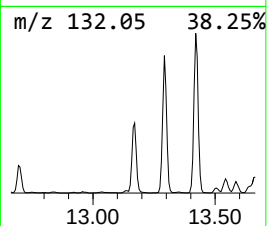
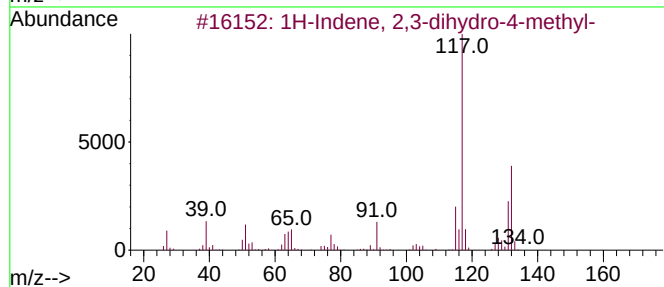
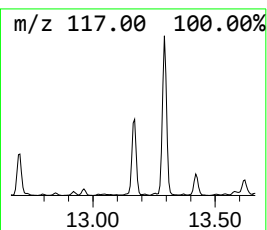
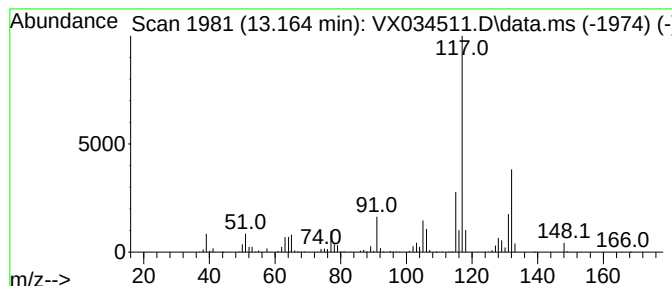
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X031023W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.164	29.18 ug/l	632726	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	93		
2	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	93		
3	3-Phenylbut-1-ene	132	C10H12	000934-10-1	93		
4	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92		
5	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	91		



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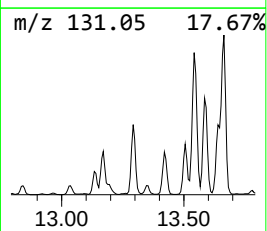
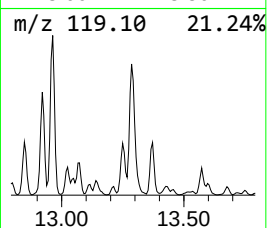
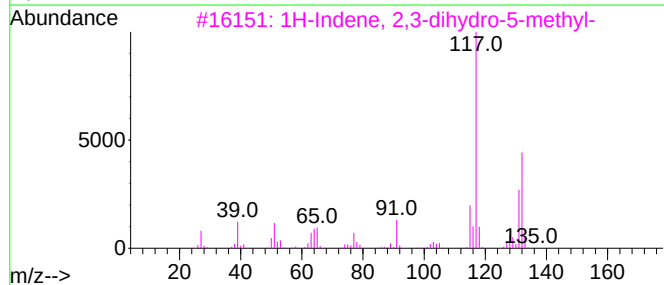
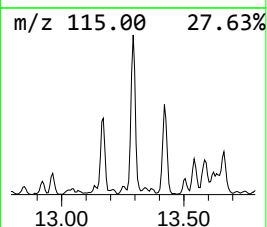
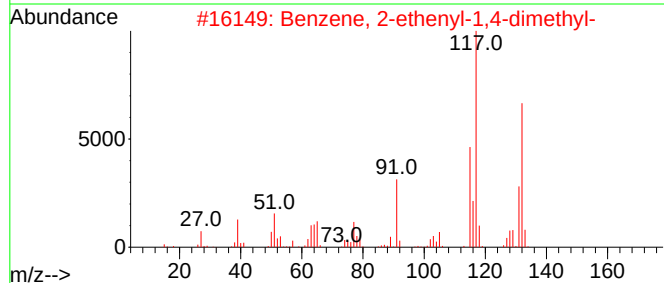
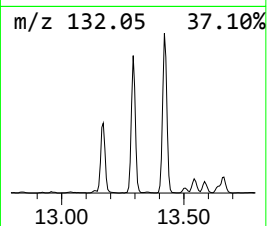
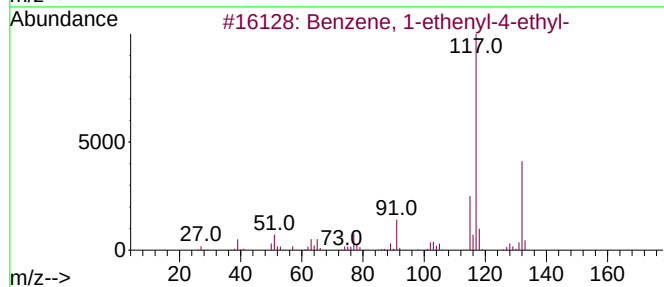
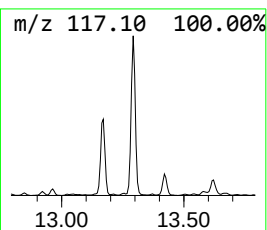
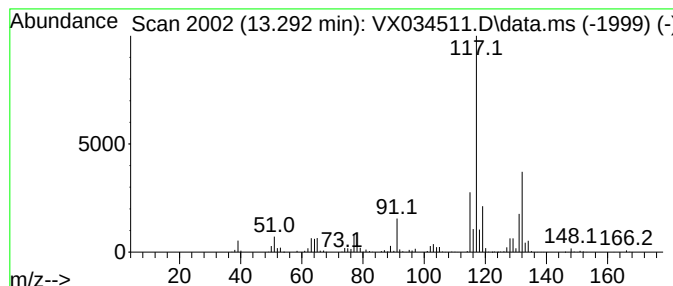
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X031023W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	61.95 ug/l	1343600	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
2			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	91
3			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	90
4			3-Phenylbut-1-ene	132	C10H12	000934-10-1	87
5			2,4-Dimethylstyrene	132	C10H12	002234-20-0	83



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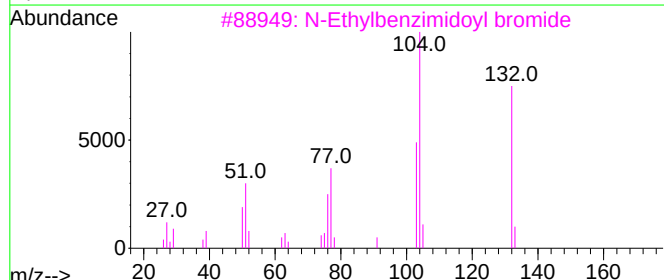
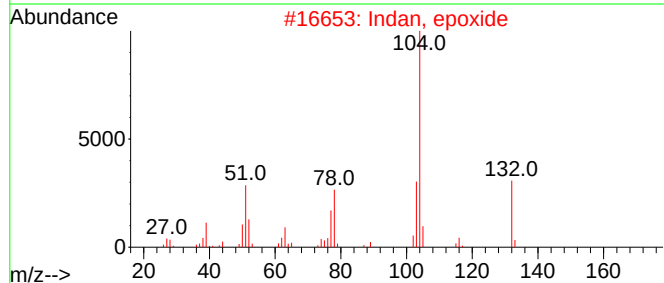
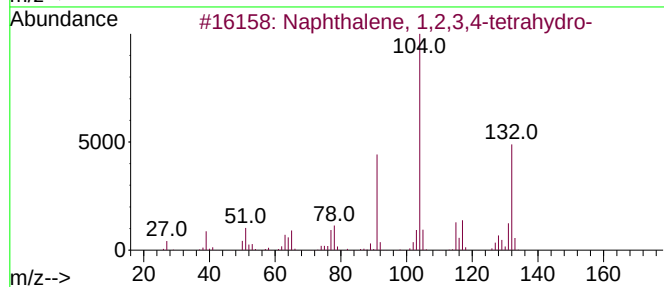
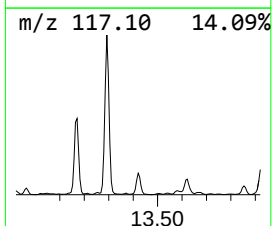
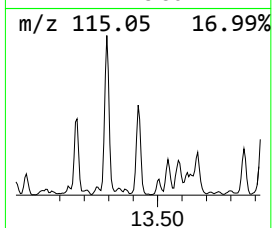
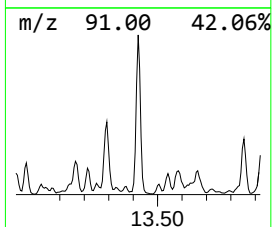
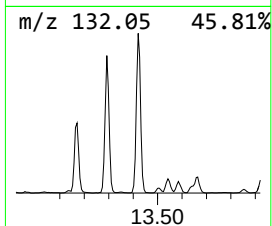
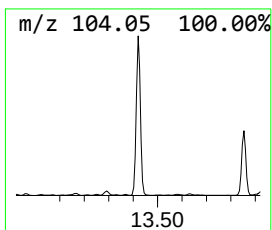
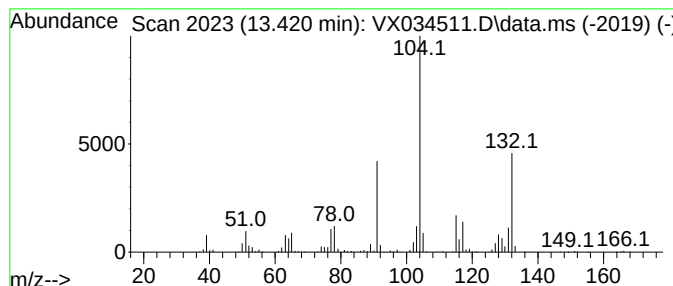
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X031023W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Naphthalene, 1,2,3,4-tetrahydro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.420	60.54 ug/l	1312870	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	95
2			Indan, epoxide	132	C9H8O	000768-22-9	58
3			N-Ethylbenzimidoyl bromide	211	C9H10BrN	041182-87-0	50
4			Isoquinoline, 1,2,3,4-tetrahydro-	133	C9H11N	000091-21-4	43
5			2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX031523\
Data File : VX034511.D
Acq On : 15 Mar 2023 16:22
Operator : JC/MD
Sample : 01892-01 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BOOM-030923

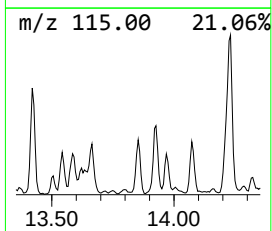
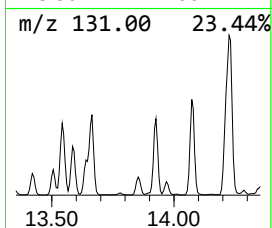
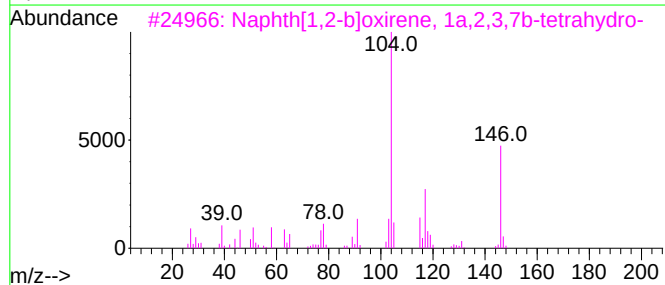
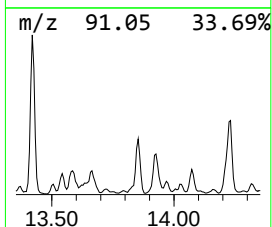
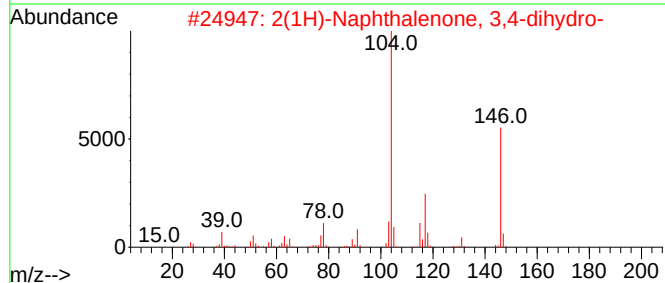
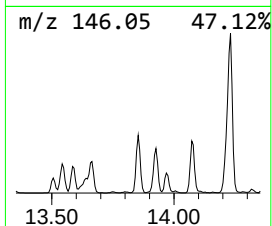
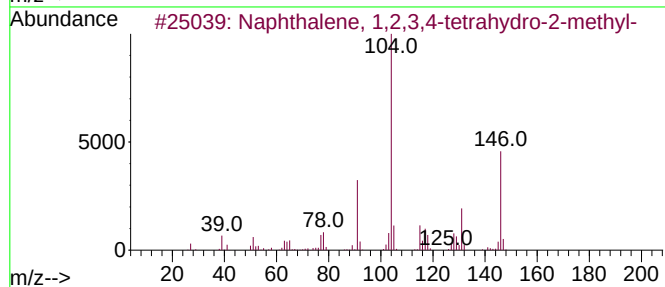
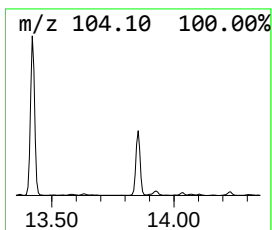
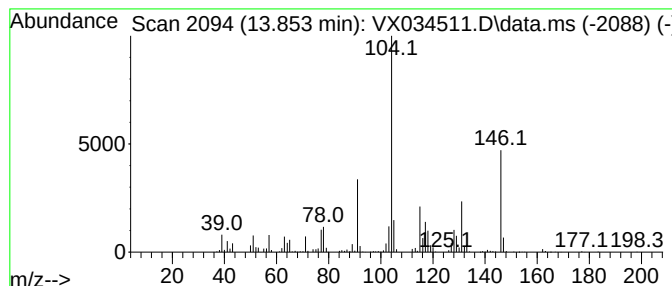
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.853	33.69 ug/l	730670	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	003877-19-8	93
2			2(1H)-Naphthalenone, 3,4-dihydro-	146	C10H10O	000530-93-8	68
3			Naphth[1,2-b]oxirene, 1a,2,3,7b-...	146	C10H10O	002461-34-9	58
4			Benzene, 3-cyclohexen-1-yl-	158	C12H14	004994-16-5	38
5			Phenethylpropanamide	177	C11H15NO	1000380-03-1	35



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX031523\
Data File : VX034511.D
Acq On : 15 Mar 2023 16:22
Operator : JC/MD
Sample : 01892-01 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BOOM-030923

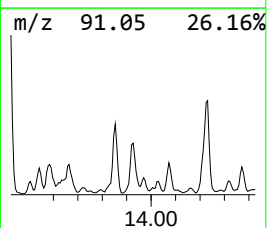
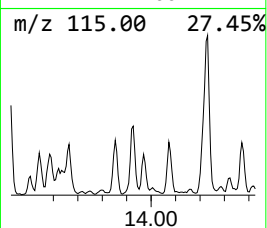
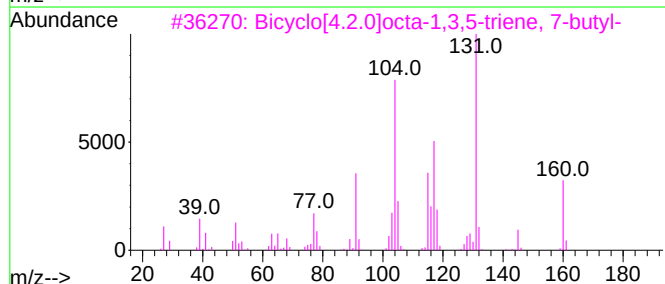
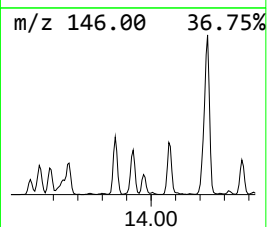
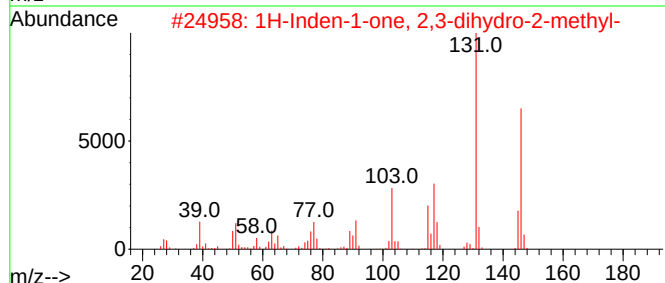
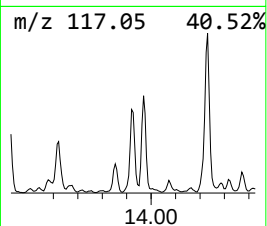
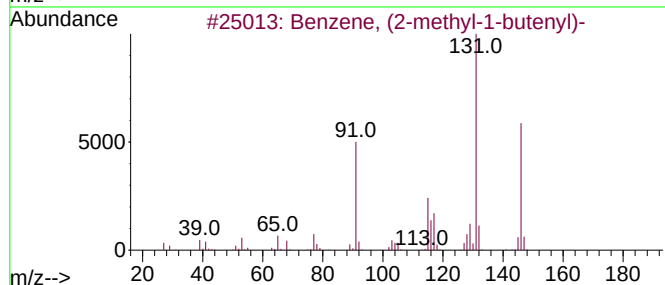
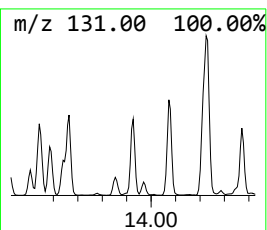
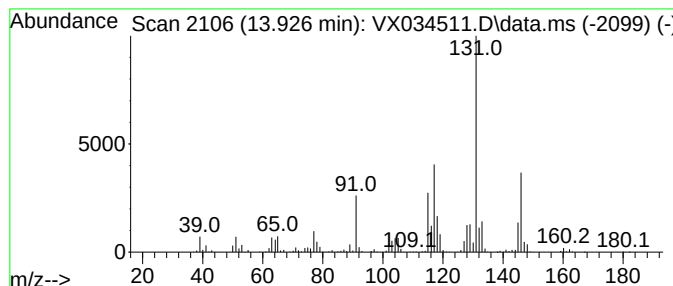
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, (2-methyl-1-butenyl)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.926	32.65 ug/l	708131	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	94
2			1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	91
3			Bicyclo[4.2.0]octa-1,3,5-triene,...	160	C12H16	078920-29-3	81
4			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	76
5			1-Methylindan-2-one	146	C10H10O	035587-60-1	74



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX031523\
Data File : VX034511.D
Acq On : 15 Mar 2023 16:22
Operator : JC/MD
Sample : 01892-01 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BOOM-030923

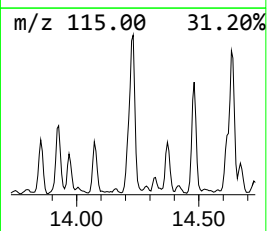
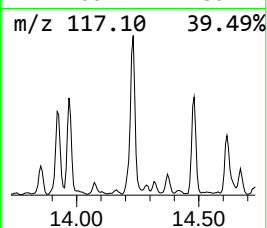
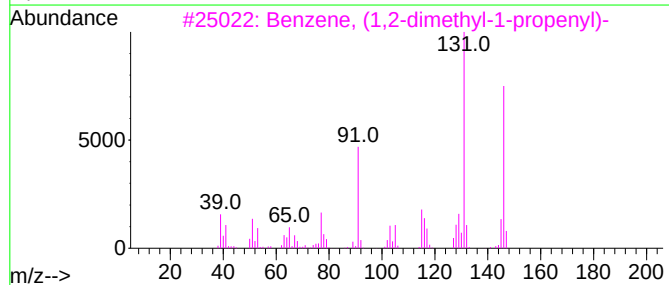
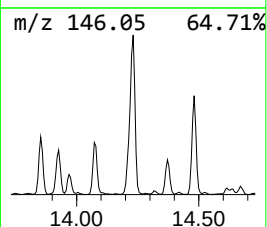
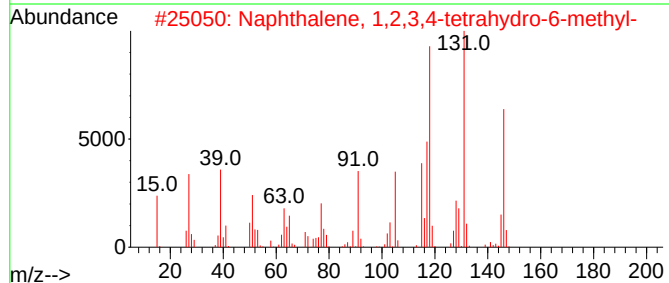
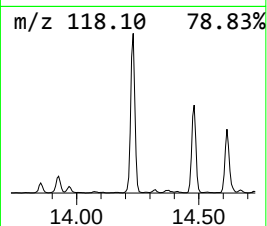
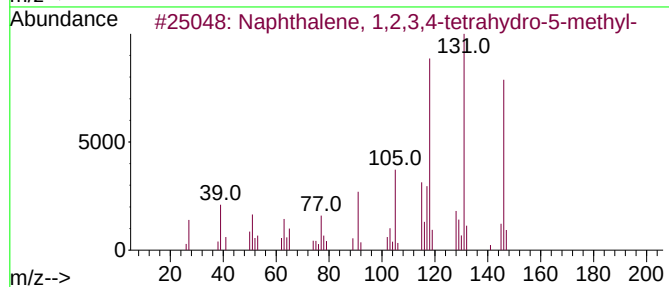
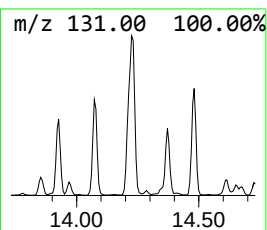
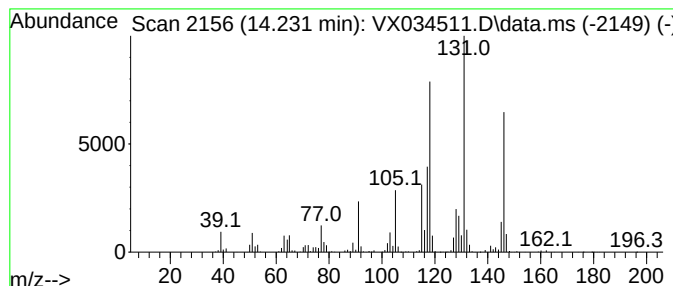
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.231	88.74 ug/l	1924550	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	97
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	97
3			Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	87
4			Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	55
5			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	55



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX031523\
Data File : VX034511.D
Acq On : 15 Mar 2023 16:22
Operator : JC/MD
Sample : 01892-01 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BOOM-030923

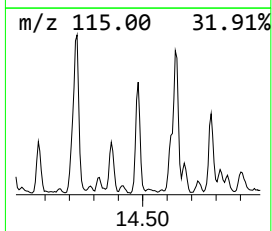
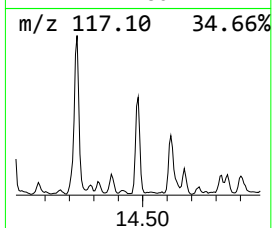
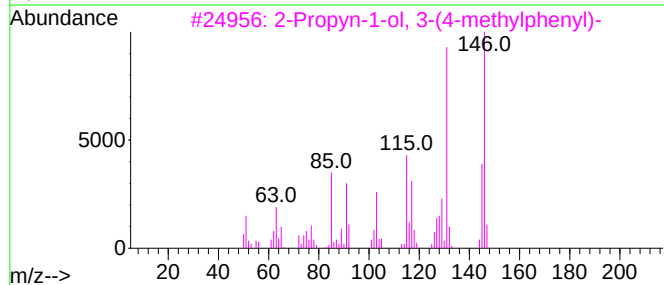
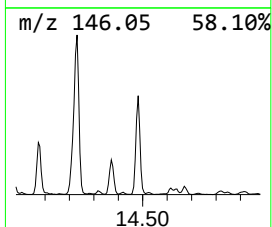
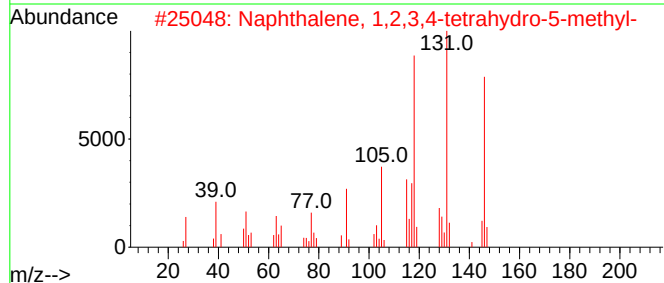
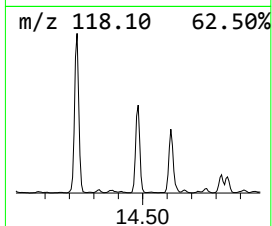
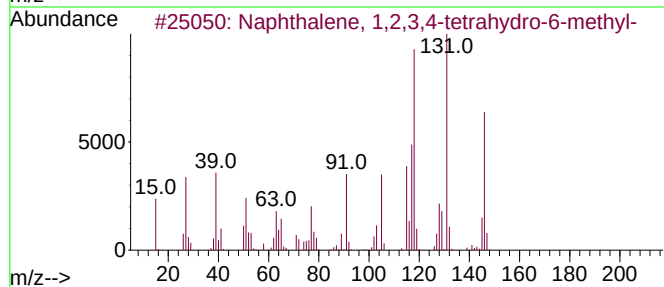
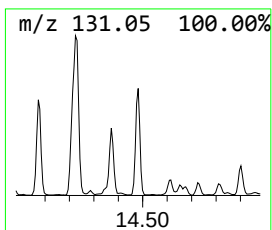
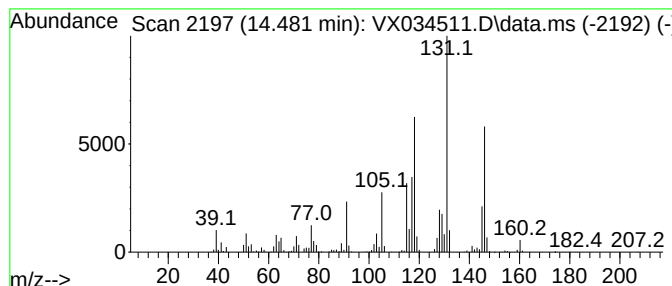
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.481	55.97 ug/l	1213800	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	95
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	95
3			2-Propyn-1-ol, 3-(4-methylphenyl)-	146	C10H10O	016017-24-6	74
4			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	70
5			Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	68



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX031523\
Data File : VX034511.D
Acq On : 15 Mar 2023 16:22
Operator : JC/MD
Sample : 01892-01 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BOOM-030923

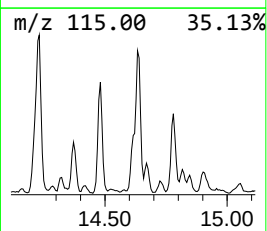
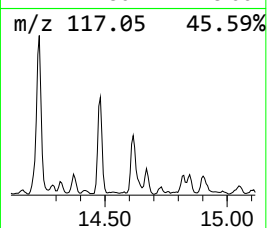
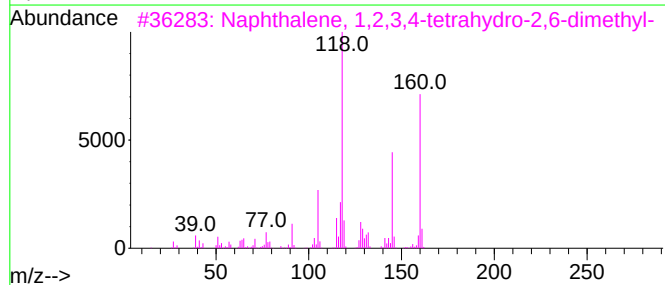
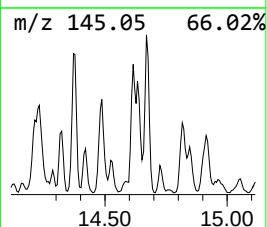
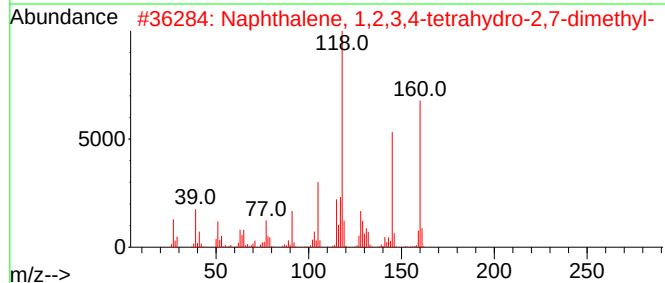
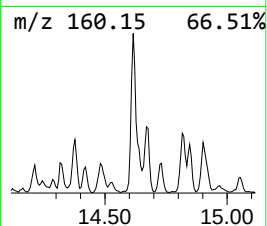
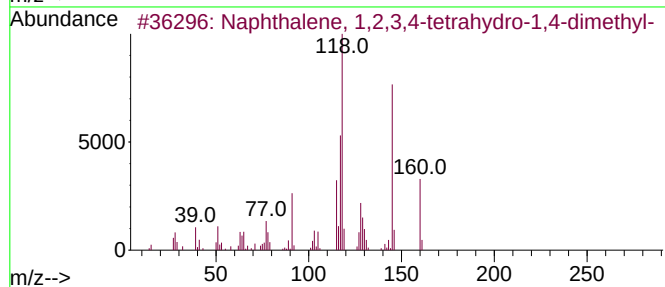
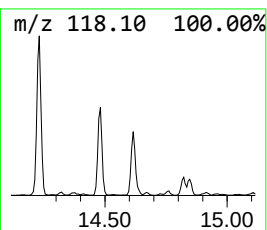
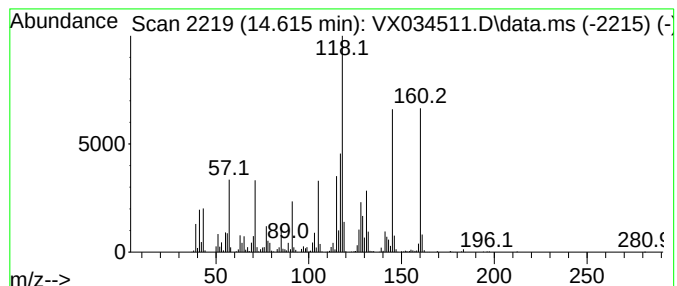
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.615	26.11 ug/l	566309	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	95
2			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	013065-07-1	89
3			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	007524-63-2	68
4			Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	64
5			Benzene, 4-(2-butenyl)-1,2-dimet...	160	C12H16	054340-86-2	60



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX031523\
Data File : VX034511.D
Acq On : 15 Mar 2023 16:22
Operator : JC/MD
Sample : 01892-01 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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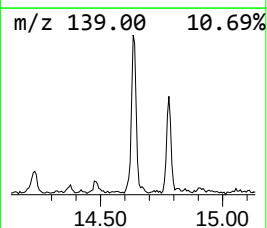
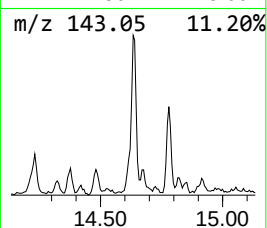
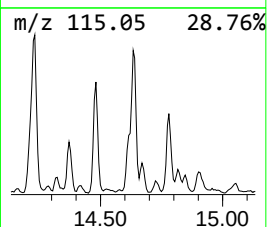
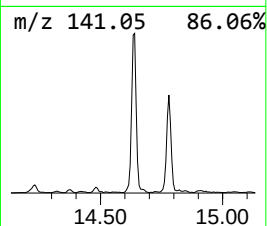
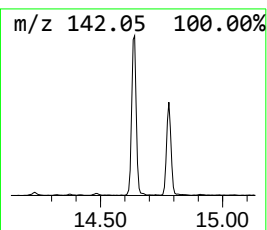
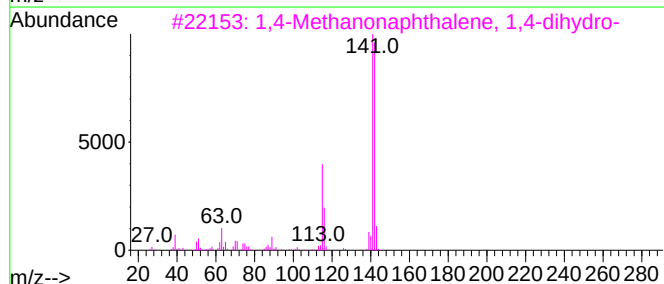
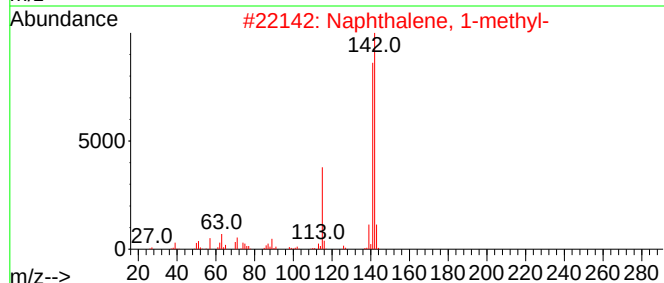
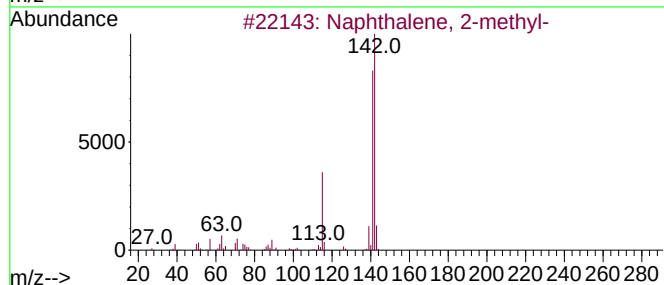
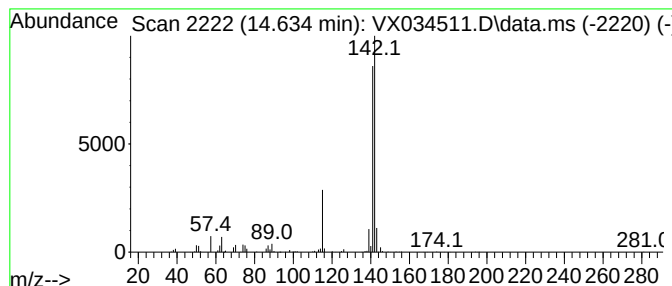
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 1,4-Methanonaphthalene, 1,4... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.634	34.97 ug/l	758462	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	94
2			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
3			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
4			Benzocycloheptatriene	142	C11H10	000264-09-5	91
5			1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	45



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX031523\
Data File : VX034511.D
Acq On : 15 Mar 2023 16:22
Operator : JC/MD
Sample : 01892-01 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
BOOM-030923

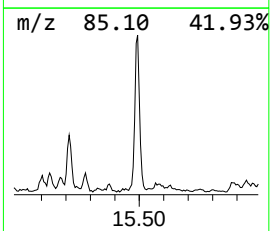
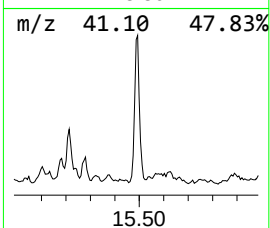
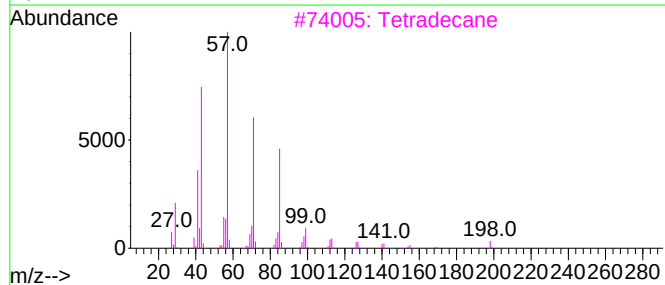
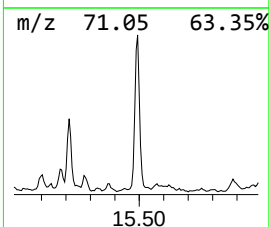
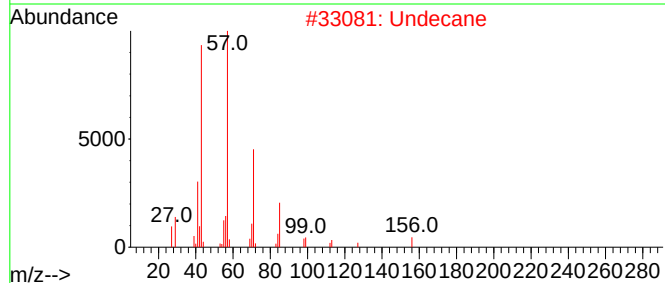
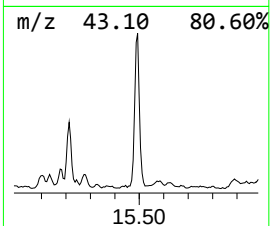
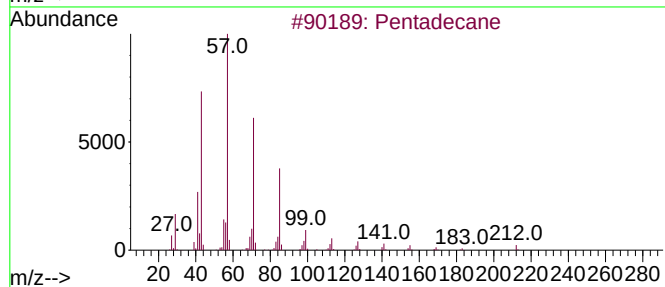
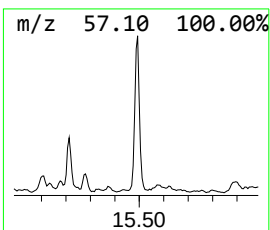
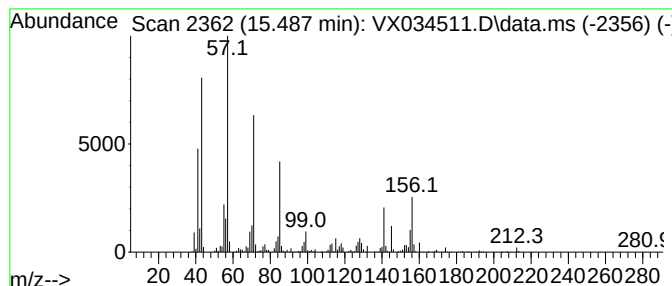
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X031023W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Pentadecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.487	37.83 ug/l	820497	1,4-Dichlorobenzene-d4	12.024

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentadecane	212	C15H32	000629-62-9	83
2		Undecane	156	C11H24	001120-21-4	58
3		Tetradecane	198	C14H30	000629-59-4	58
4		Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	52
5		Decane, 3-bromo-	220	C10H21Br	030571-71-2	52



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX031523\
Data File : VX034511.D
Acq On : 15 Mar 2023 16:22
Operator : JC/MD
Sample : 01892-01 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BOOM-030923

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X031023W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1H-Indene, 2,3-...	13.164	29.2	ug/l	632726	4	12.024	1084340	50.0
Benzene, 1-ethe...	13.292	62.0	ug/l	1343600	4	12.024	1084340	50.0
Naphthalene, 1,...	13.420	60.5	ug/l	1312870	4	12.024	1084340	50.0
Naphthalene, 1,...	13.853	33.7	ug/l	730670	4	12.024	1084340	50.0
Benzene, (2-met...	13.926	32.6	ug/l	708131	4	12.024	1084340	50.0
Naphthalene, 1,...	14.231	88.7	ug/l	1924550	4	12.024	1084340	50.0
Naphthalene, 1,...	14.481	56.0	ug/l	1213800	4	12.024	1084340	50.0
Naphthalene, 1,...	14.615	26.1	ug/l	566309	4	12.024	1084340	50.0
1,4-Methanonaph...	14.634	35.0	ug/l	758462	4	12.024	1084340	50.0
Pentadecane	15.487	37.8	ug/l	820497	4	12.024	1084340	50.0