

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111722\
 Data File : VX032895.D
 Acq On : 17 Nov 2022 13:21
 Operator : JC/MD
 Sample : N5650-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 34271

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\82X102622W.M
 Title : SW846 8260

Signal : TIC: VX032895.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.392	207	214	224	rVB2	5617	12364	2.03%	0.182%
2	5.385	693	705	721	rBV	76198	218565	35.92%	3.214%
3	5.550	721	732	746	rVB2	116073	333168	54.75%	4.899%
4	5.958	788	799	815	rBV	78749	210459	34.59%	3.094%
5	6.763	922	931	946	rVB	178560	429446	70.57%	6.314%
6	8.647	1234	1240	1252	rBV	390371	608501	100.00%	8.947%
7	10.055	1465	1471	1481	rBV	378769	506871	83.30%	7.453%
8	10.640	1562	1567	1573	rBV	17041	23033	3.79%	0.339%
9	11.079	1633	1639	1652	rBV	369167	470982	77.40%	6.925%
10	11.384	1682	1689	1690	rBV	8202	10914	1.79%	0.160%
11	11.402	1690	1692	1696	rVB	8998	9572	1.57%	0.141%
12	11.451	1696	1700	1704	rBV	14059	16556	2.72%	0.243%
13	11.616	1721	1727	1734	rVB	27078	36337	5.97%	0.534%
14	11.756	1746	1750	1757	rVB	8412	12053	1.98%	0.177%
15	11.890	1768	1772	1776	rVB	5063	6855	1.13%	0.101%
16	11.969	1781	1785	1788	rVV	17073	18609	3.06%	0.274%
17	12.024	1788	1794	1800	rVB	391161	499251	82.05%	7.341%
18	12.085	1800	1804	1809	rBV	62089	77134	12.68%	1.134%
19	12.250	1823	1831	1832	rBV2	31683	41995	6.90%	0.617%
20	12.268	1832	1834	1838	rVB	29887	31337	5.15%	0.461%
21	12.317	1838	1842	1850	rVB2	57202	83256	13.68%	1.224%
22	12.408	1850	1857	1862	rBV3	13557	21959	3.61%	0.323%
23	12.463	1862	1866	1871	rVB	29870	36433	5.99%	0.536%
24	12.530	1871	1877	1879	rBV	13008	17199	2.83%	0.253%
25	12.555	1879	1881	1885	rVB2	16511	16869	2.77%	0.248%
26	12.616	1885	1891	1894	rBV	62824	79414	13.05%	1.168%
27	12.646	1894	1896	1900	rVB	9702	9878	1.62%	0.145%
28	12.701	1900	1905	1907	rBV	32565	47787	7.85%	0.703%
29	12.799	1917	1921	1925	rBV	12309	15666	2.57%	0.230%
30	12.847	1925	1929	1934	rVB2	35629	44561	7.32%	0.655%
31	12.920	1934	1941	1945	rBV	45487	60941	10.01%	0.896%
32	12.963	1945	1948	1954	rVB	49714	60363	9.92%	0.888%
33	13.024	1954	1958	1960	rBV	18496	26137	4.30%	0.384%
34	13.073	1964	1966	1968	rVB	8544	7550	1.24%	0.111%
35	13.140	1973	1977	1978	rVV2	19556	24861	4.09%	0.366%
36	13.164	1978	1981	1985	rVV3	25659	40310	6.62%	0.593%

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37	13.213	1985	1989	1992	rVV3	17659	25073	4.12%	0.369%
38	13.256	1992	1996	1998	rVV2	24696	37961	6.24%	0.558%
39	13.292	1998	2002	2007	rVV2	152189	228048	37.48%	3.353%
40	13.335	2007	2009	2012	rVV2	16162	22313	3.67%	0.328%
41	13.372	2012	2015	2018	rVV	18423	22341	3.67%	0.328%
42	13.420	2019	2023	2031	rVB	81855	113670	18.68%	1.671%
43	13.506	2032	2037	2040	rBV2	16245	24202	3.98%	0.356%
44	13.542	2040	2043	2046	rVV	31728	42727	7.02%	0.628%
45	13.585	2046	2050	2055	rVV3	67242	121275	19.93%	1.783%
46	13.640	2056	2059	2060	rVV	29848	35043	5.76%	0.515%
47	13.664	2060	2063	2069	rVB2	60173	90109	14.81%	1.325%
48	13.750	2074	2077	2080	rVB2	7095	7709	1.27%	0.113%
49	13.792	2080	2084	2088	rBV3	9395	12457	2.05%	0.183%
50	13.853	2088	2094	2099	rVB	67379	96713	15.89%	1.422%
51	13.926	2099	2106	2110	rBV2	88533	127117	20.89%	1.869%
52	13.969	2110	2113	2117	rVV	26447	31564	5.19%	0.464%
53	14.036	2121	2124	2127	rVB4	8492	8569	1.41%	0.126%
54	14.073	2127	2130	2134	rBV	42647	52925	8.70%	0.778%
55	14.164	2142	2145	2148	rBV2	7302	7876	1.29%	0.116%
56	14.225	2148	2155	2163	rVV2	122852	237880	39.09%	3.498%
57	14.323	2167	2171	2175	rBV	30888	40277	6.62%	0.592%
58	14.371	2175	2179	2183	rVB2	69019	82664	13.58%	1.215%
59	14.420	2183	2187	2192	rVB	15448	21365	3.51%	0.314%
60	14.481	2192	2197	2202	rBV2	144099	190858	31.37%	2.806%
61	14.615	2215	2219	2225	rBV2	74739	128311	21.09%	1.887%
62	14.670	2225	2228	2233	rVB	49650	67309	11.06%	0.990%
63	14.725	2233	2237	2240	rBV2	17672	25015	4.11%	0.368%
64	14.755	2240	2242	2244	rVV2	9428	8211	1.35%	0.121%
65	14.780	2244	2246	2249	rVV4	5122	7831	1.29%	0.115%
66	14.816	2249	2252	2255	rVV	33649	49488	8.13%	0.728%
67	14.847	2255	2257	2262	rVB2	30930	32640	5.36%	0.480%
68	14.902	2262	2266	2272	rBV2	33658	59942	9.85%	0.881%
69	15.048	2284	2290	2295	rBV2	16833	28506	4.68%	0.419%
70	15.182	2308	2312	2320	rVB	71575	129155	21.23%	1.899%
71	15.249	2320	2323	2326	rBV3	9457	13563	2.23%	0.199%
72	15.316	2331	2334	2339	rVV6	8152	14161	2.33%	0.208%
73	15.377	2339	2344	2352	rVV6	19597	41484	6.82%	0.610%
74	15.450	2352	2356	2357	rVV3	6093	9005	1.48%	0.132%
75	15.481	2357	2361	2365	rVV3	36177	63963	10.51%	0.940%
76	15.524	2365	2368	2373	rVV3	20525	42665	7.01%	0.627%

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77	15.572	2373	2376	2378	rVV4	12602	19034	3.13%	0.280%
78	15.615	2378	2383	2386	rVV2	30460	57698	9.48%	0.848%
79	15.652	2386	2389	2398	rVB6	15090	28974	4.76%	0.426%
80	15.755	2403	2406	2411	rBV3	3798	6338	1.04%	0.093%
81	15.847	2417	2421	2425	rBV5	8606	15571	2.56%	0.229%
82	15.889	2425	2428	2434	rVB5	6475	10589	1.74%	0.156%
83	16.024	2447	2450	2456	rVB5	6929	11985	1.97%	0.176%
84	16.091	2456	2461	2467	rVV3	8417	19367	3.18%	0.285%
85	16.170	2471	2474	2483	rVB9	4917	12953	2.13%	0.190%
86	16.298	2483	2495	2499	rBV10	4874	12079	1.99%	0.178%
87	16.377	2502	2508	2512	rVB4	7760	15231	2.50%	0.224%
88	16.523	2527	2532	2535	rBV5	3208	7167	1.18%	0.105%
89	16.597	2541	2544	2550	rVB3	4365	6317	1.04%	0.093%
90	16.658	2550	2554	2559	rVV3	4630	8650	1.42%	0.127%

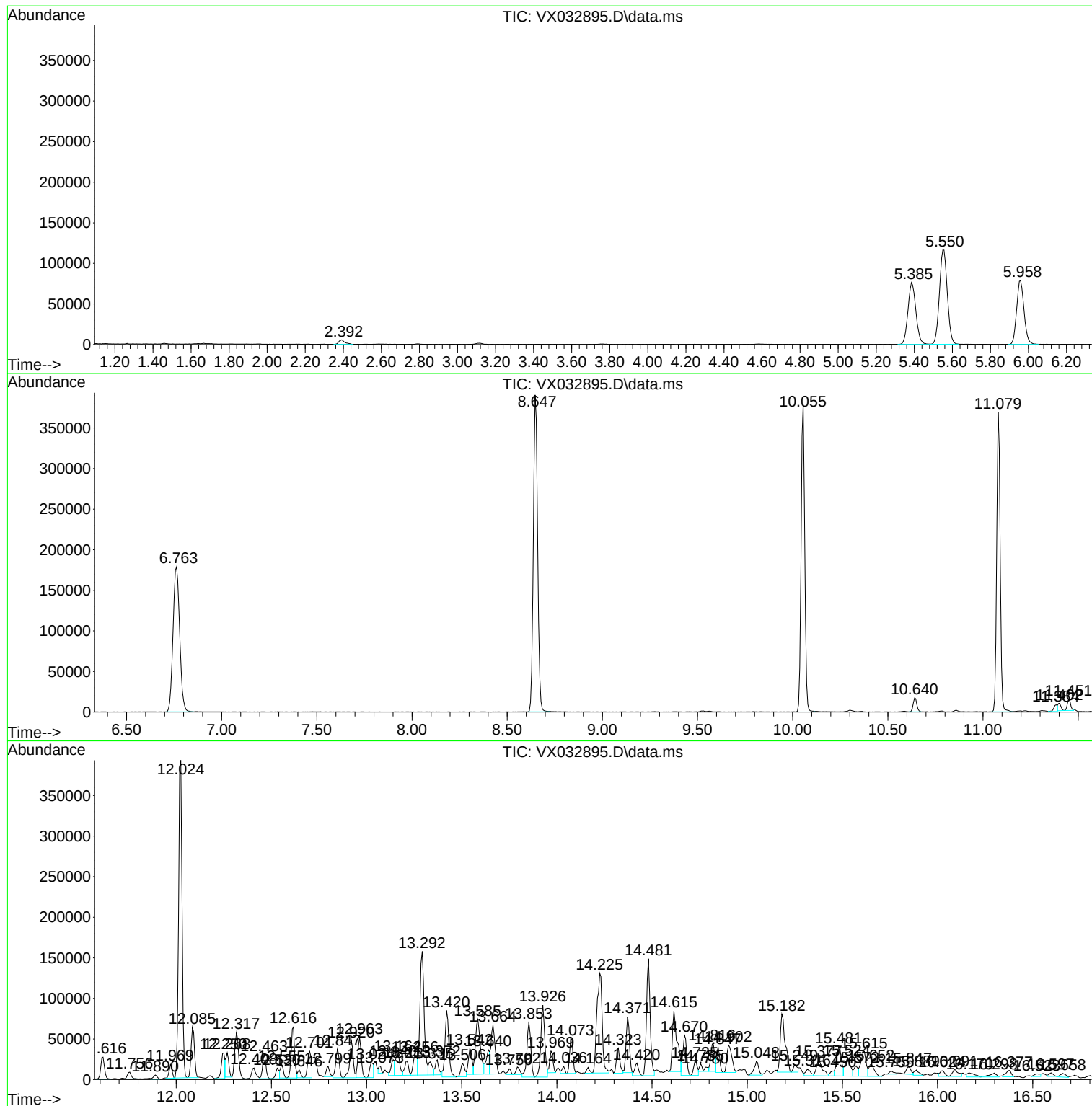
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title : SW846 8260

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



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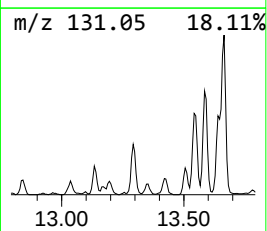
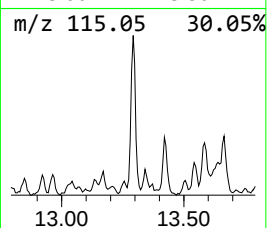
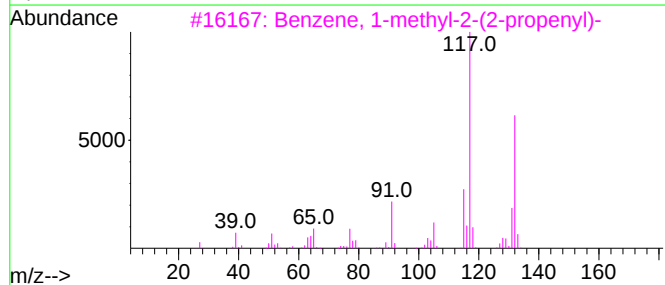
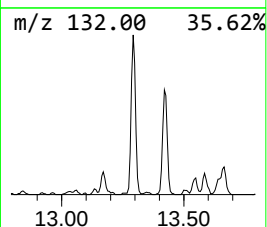
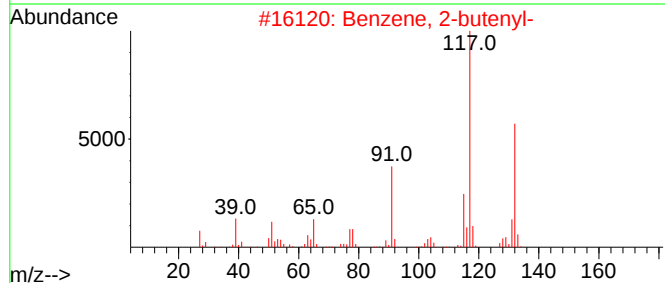
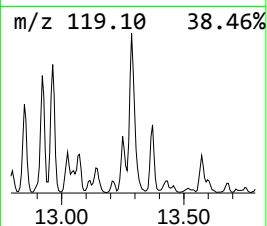
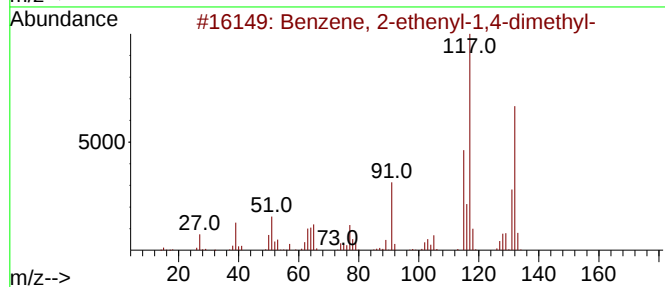
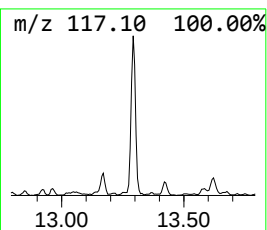
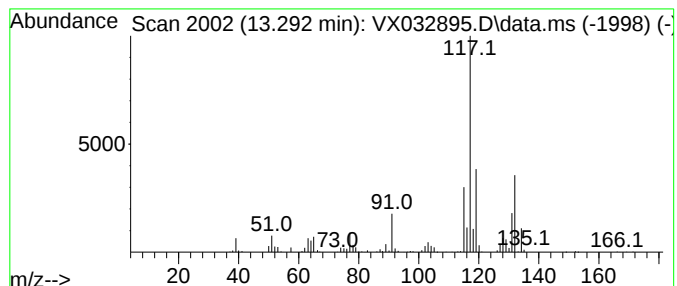
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	22.84 ug/l	228048	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	91
2			Benzene, 2-butenyl-	132	C10H12	001560-06-1	70
3			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	70
4			3-Phenylbut-1-ene	132	C10H12	000934-10-1	70
5			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	70



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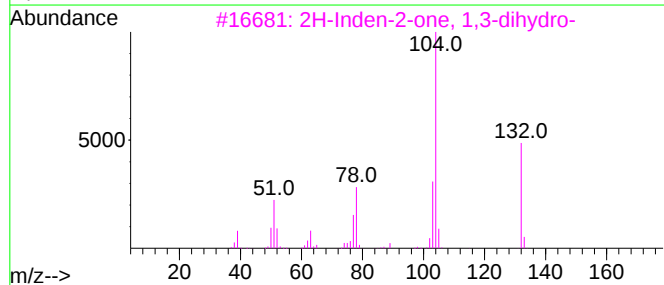
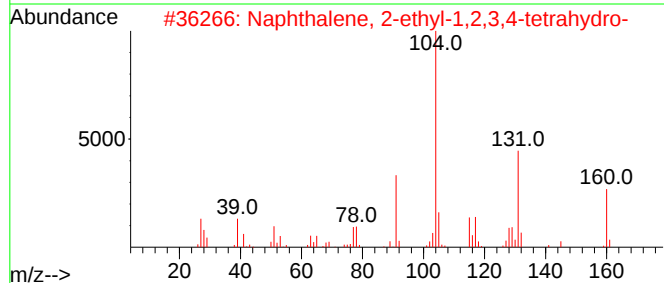
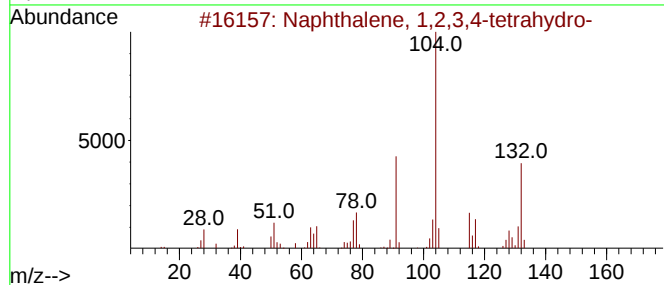
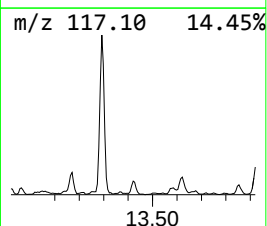
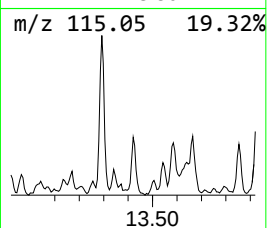
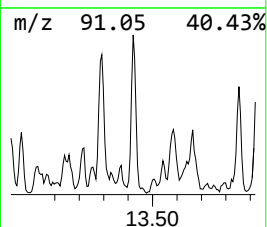
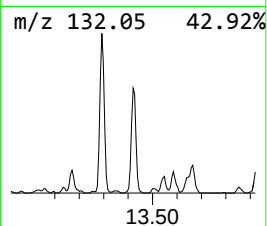
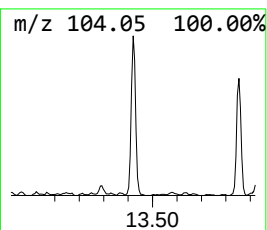
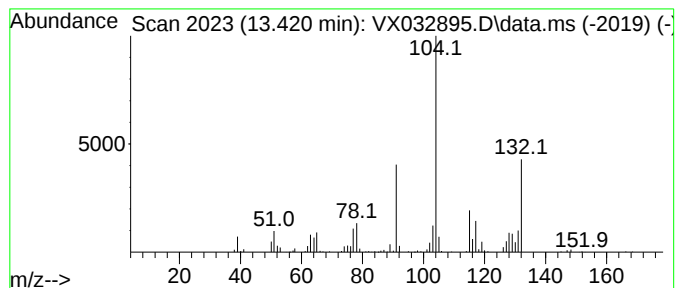
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.420	11.38 ug/l	113670	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			Naphthalene, 2-ethyl-1,2,3,4-tetrahydro-	160	C12H16	032367-54-7	50
3			2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	49
4			2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	43
5			3(2H)-Furanone, dihydro-2,2-dimethyl-	190	C12H14O2	063678-00-2	35



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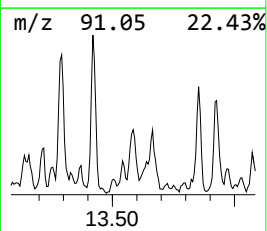
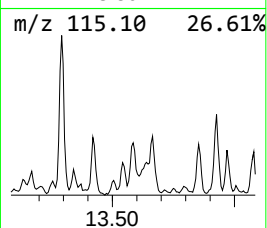
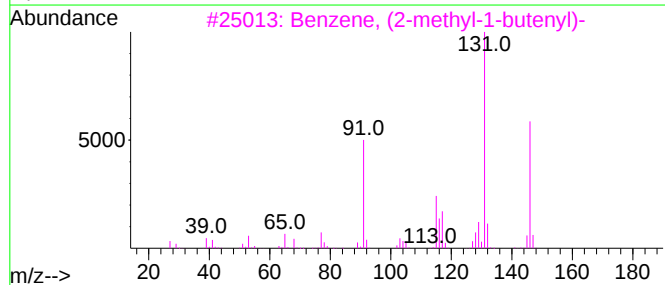
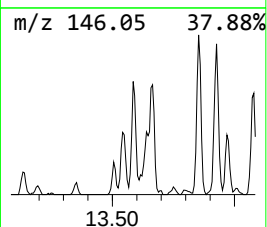
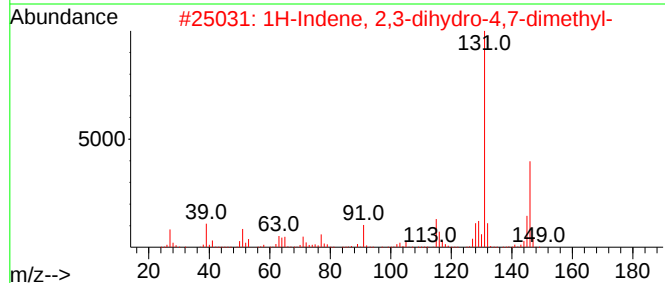
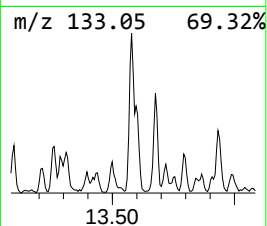
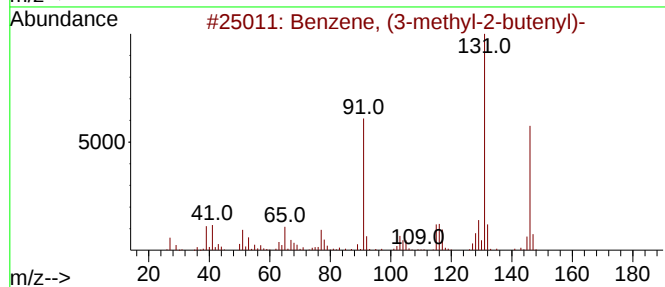
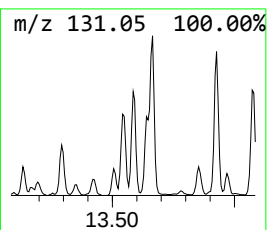
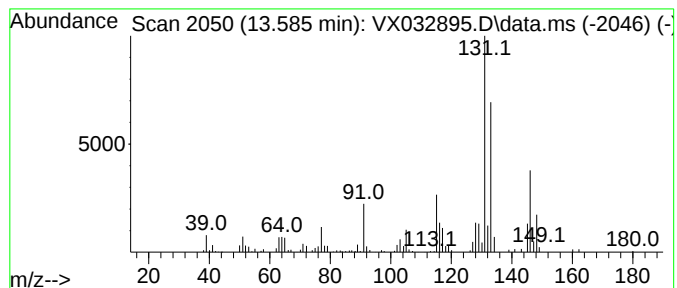
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title : SW846 8260

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

Peak Number 3 Benzene, (3-methyl-2-butenyl)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.585	12.15 ug/l	121275	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	92
2			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	86
3			Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	86
4			Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	76
5			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	76



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Operator : JC/MD
Sample : N5650-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
34271

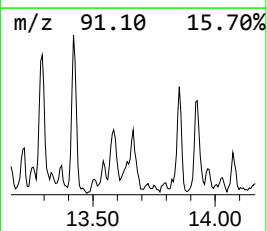
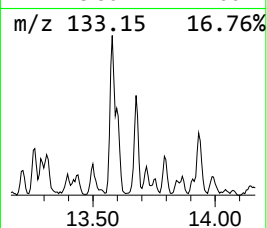
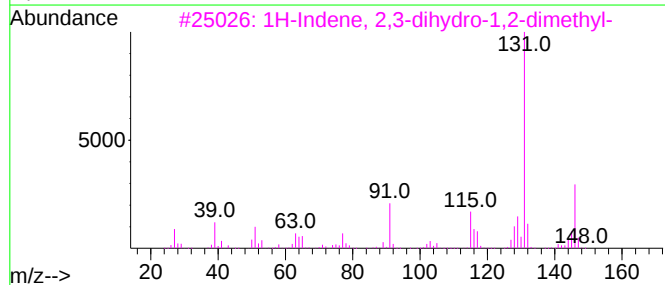
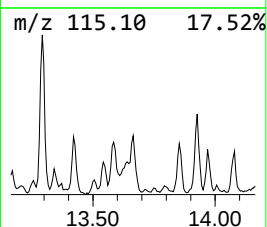
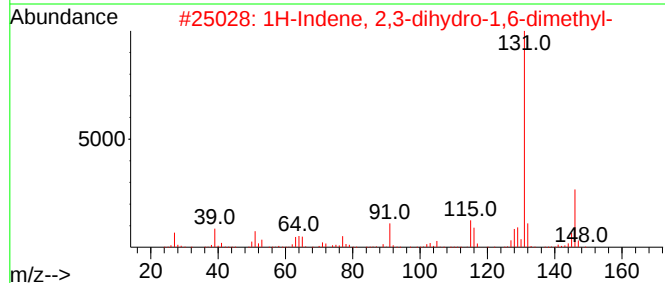
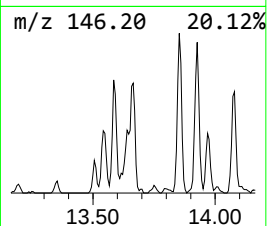
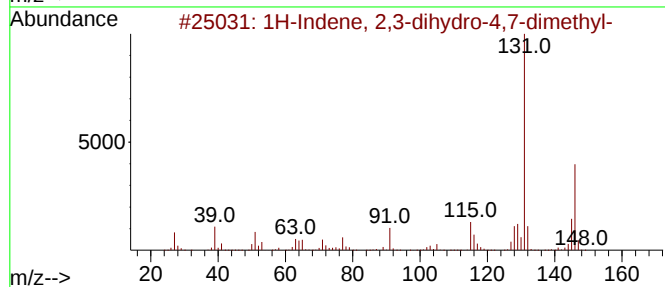
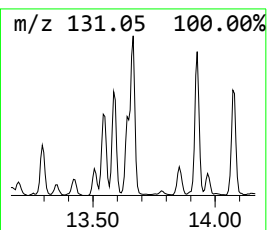
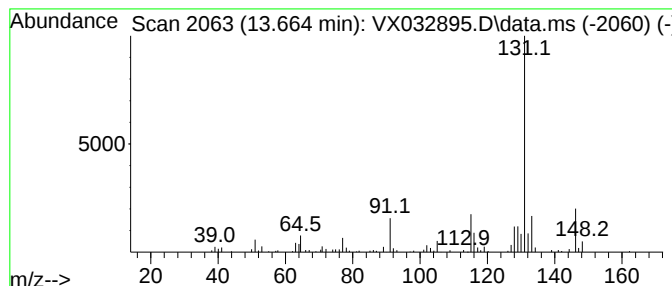
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Quant Title : SW846 8260

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

Peak Number 4 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.664	9.02 ug/l	90109	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	93
2			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	93
3			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	87
4			Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	87
5			1H-Indene, 2,3-dihydro-4,6-dimet...	146	C11H14	001685-82-1	81



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111722\
Data File : VX032895.D
Acq On : 17 Nov 2022 13:21
Operator : JC/MD
Sample : N5650-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
34271

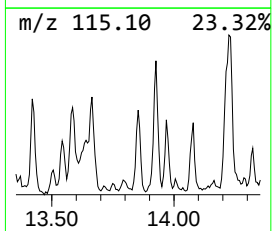
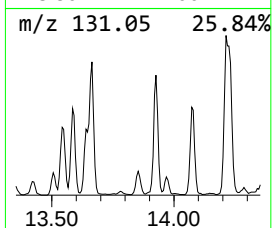
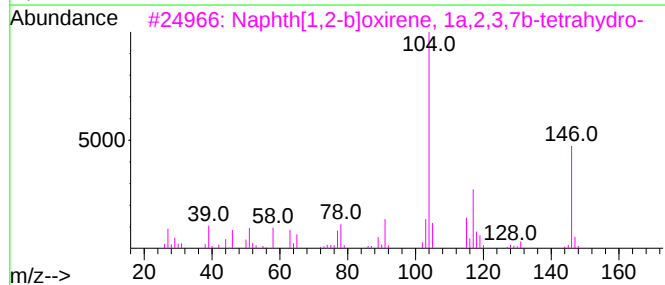
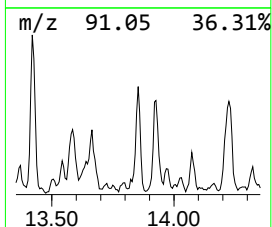
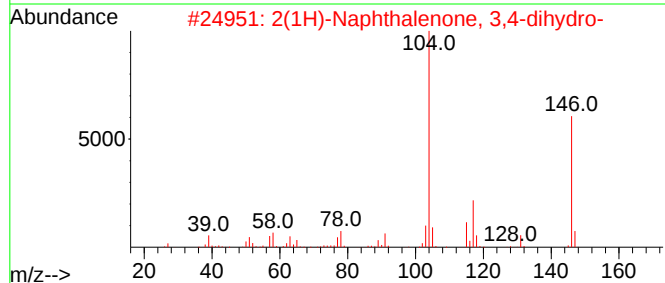
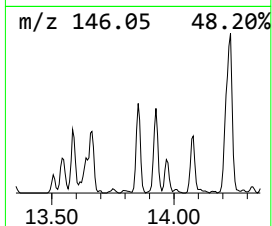
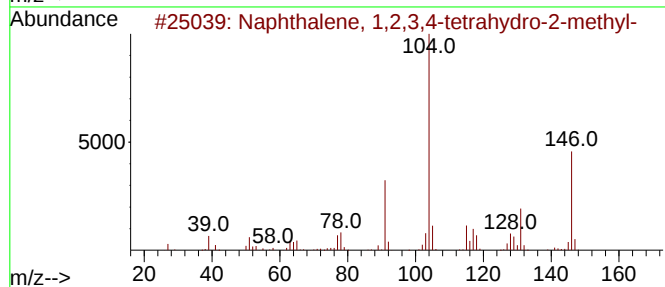
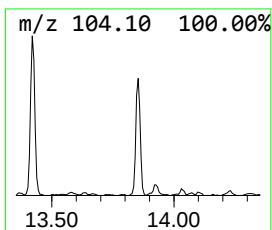
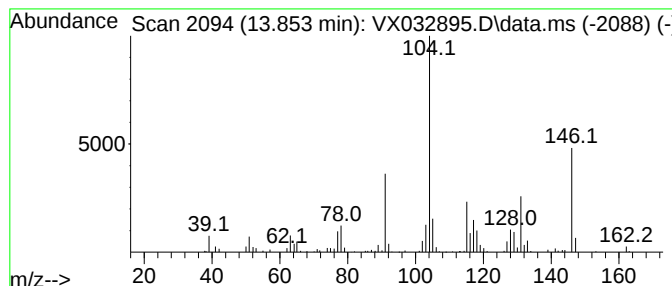
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Quant Title : SW846 8260

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

Peak Number 5 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.853	9.69 ug/l	96713	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	64
3			Naphth[1,2-b]oxirene, 1a,2,3,7b...	146	C10H10O	002461-34-9	58
4			2-Furanone, 4-phenyltetrahydro	162	C10H10O2	1000197-38-4	35
5			Acetic acid, 2-phenylethyl ester	164	C10H12O2	000103-45-7	27



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111722\
Data File : VX032895.D
Acq On : 17 Nov 2022 13:21
Operator : JC/MD
Sample : N5650-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
34271

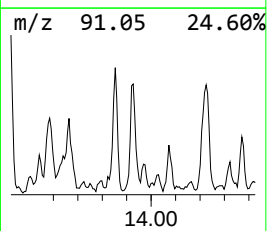
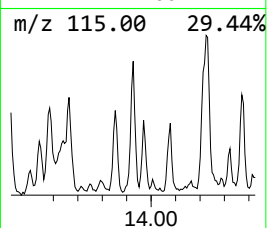
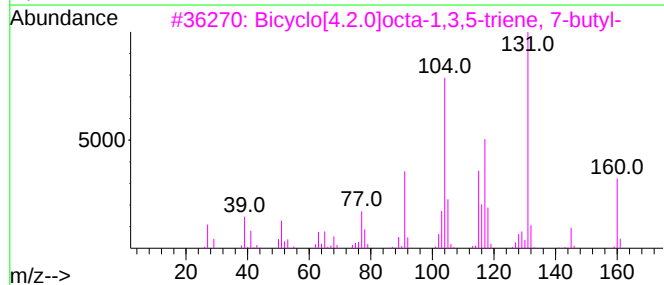
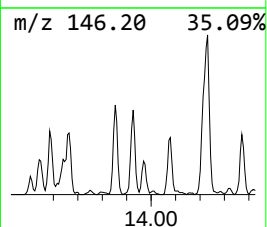
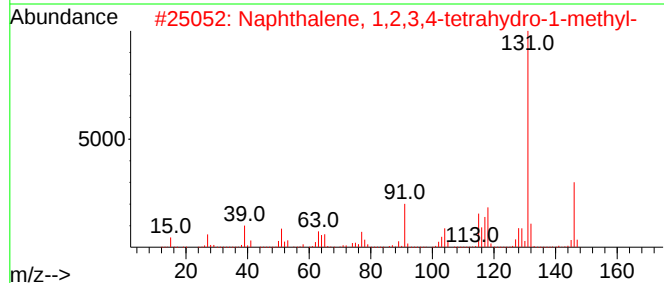
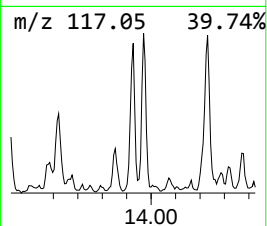
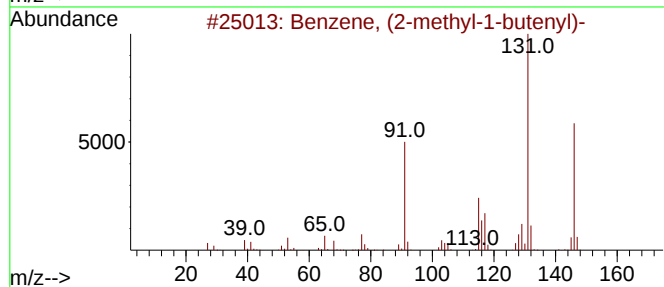
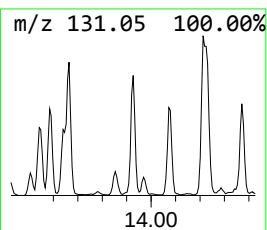
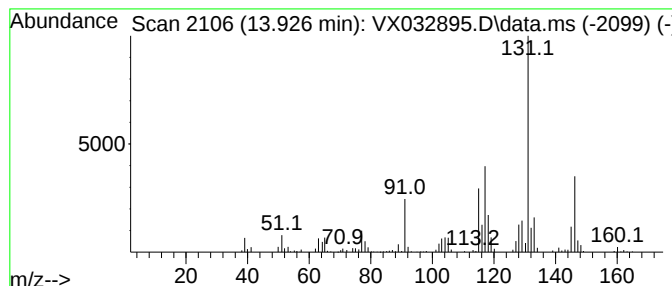
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, (2-methyl-1-butenyl)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.926	12.73 ug/l	127117	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			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	93
3			Bicyclo[4.2.0]octa-1,3,5-triene,...	160	C12H16	078920-29-3	93
4			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	92
5			1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111722\
Data File : VX032895.D
Acq On : 17 Nov 2022 13:21
Operator : JC/MD
Sample : N5650-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
34271

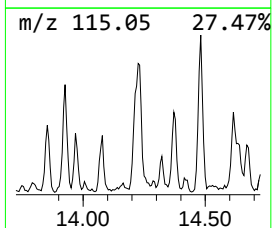
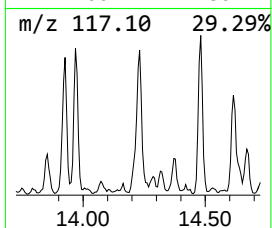
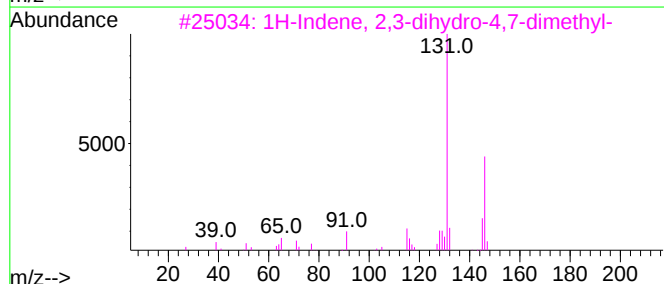
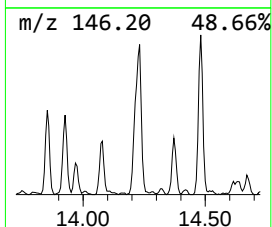
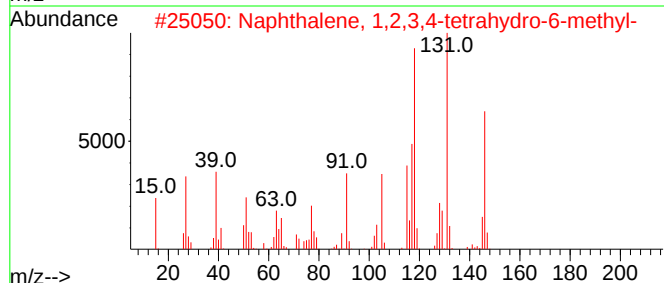
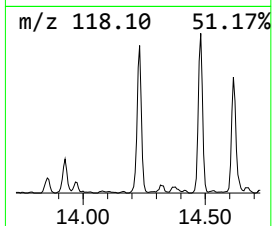
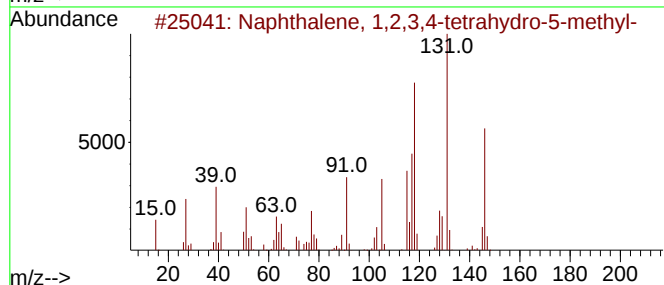
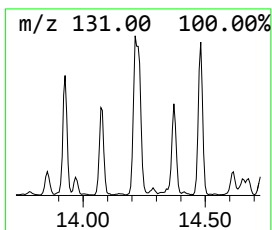
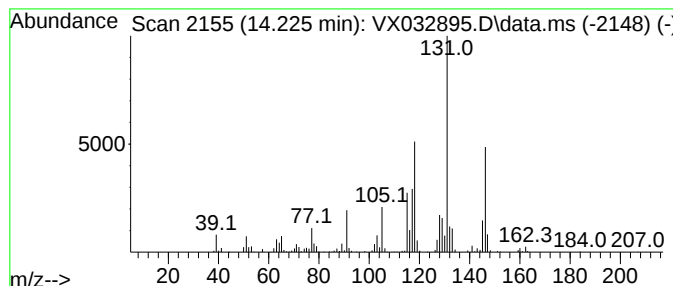
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.225	23.82 ug/l	237880	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	95
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	95
3			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	93
4			1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	76
5			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	70



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111722\
Data File : VX032895.D
Acq On : 17 Nov 2022 13:21
Operator : JC/MD
Sample : N5650-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
34271

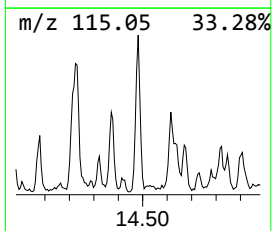
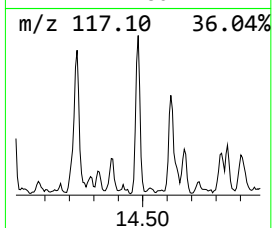
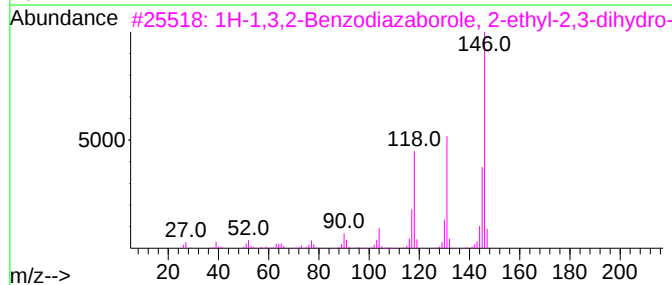
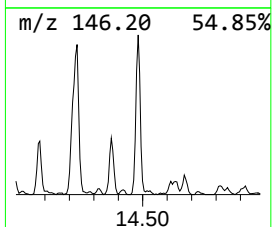
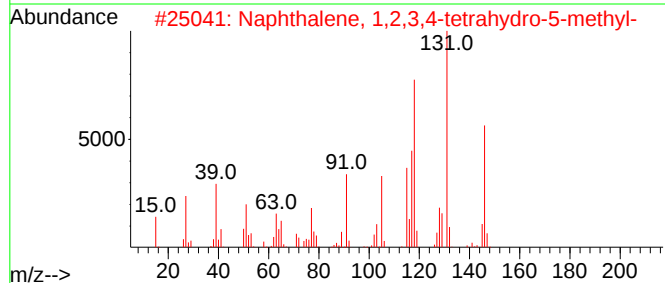
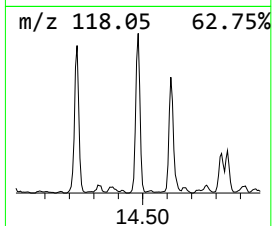
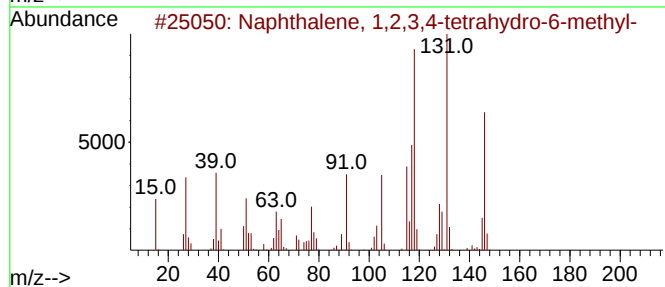
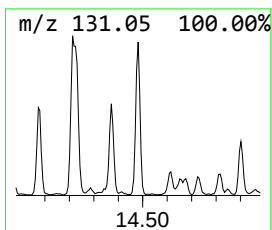
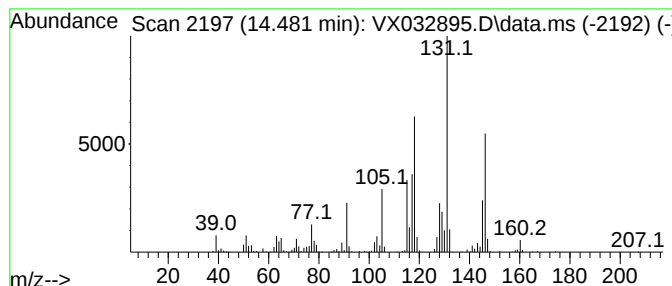
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.481	19.11 ug/l	190858	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	93
3			1H-1,3,2-Benzodiazaborole, 2-eth...	146	C8H11BN2	074663-81-3	87
4			Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	70
5			2-Propyn-1-ol, 3-(4-methylphenyl)-	146	C10H10O	016017-24-6	68



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111722\
Data File : VX032895.D
Acq On : 17 Nov 2022 13:21
Operator : JC/MD
Sample : N5650-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

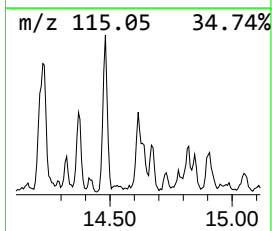
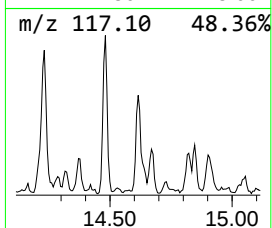
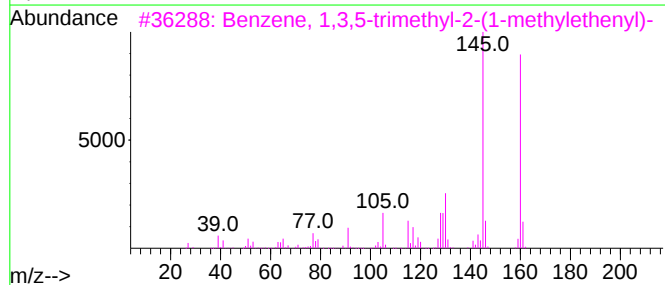
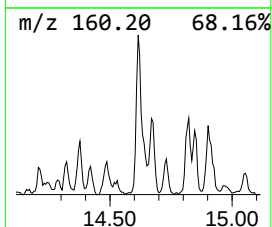
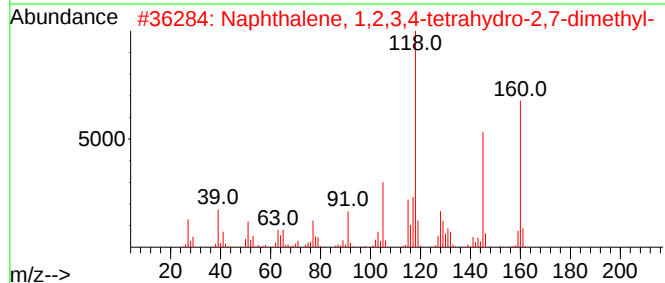
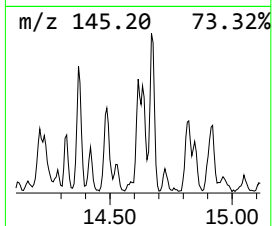
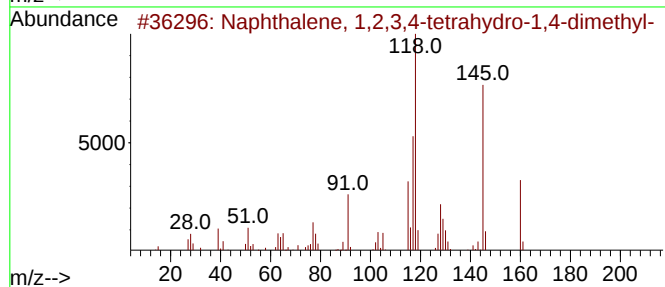
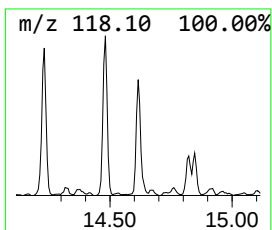
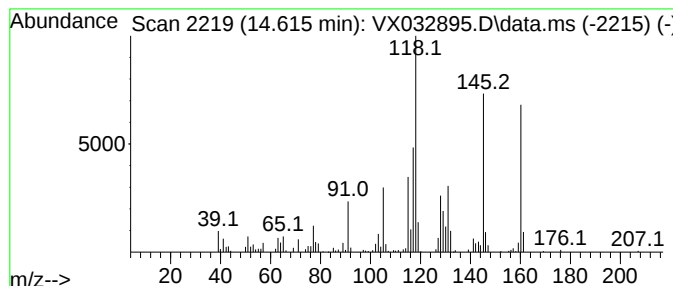
Instrument :
MSVOA_X
ClientSampleId :
34271

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD			R.T.
14.615	12.85 ug/l	128311	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	90
3	Benzene, 1,3,5-trimethyl-2-(1-me...		160	C12H16	014679-13-1	70
4	Benzene, (3-methyl-1-methylenebu...		160	C12H16	038212-14-5	68
5	Naphthalene, 1,2,3,4-tetrahydro-...		160	C12H16	007524-63-2	62



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111722\
Data File : VX032895.D
Acq On : 17 Nov 2022 13:21
Operator : JC/MD
Sample : N5650-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
34271

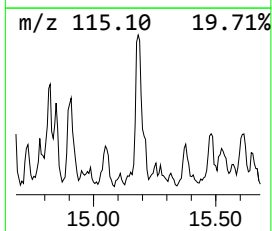
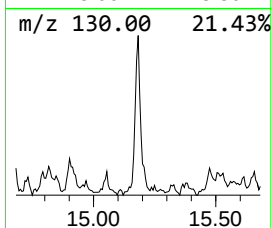
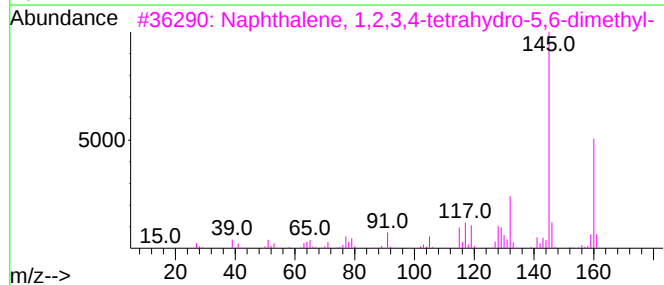
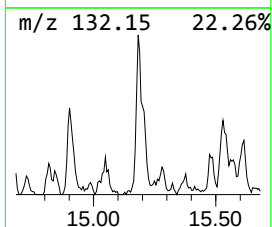
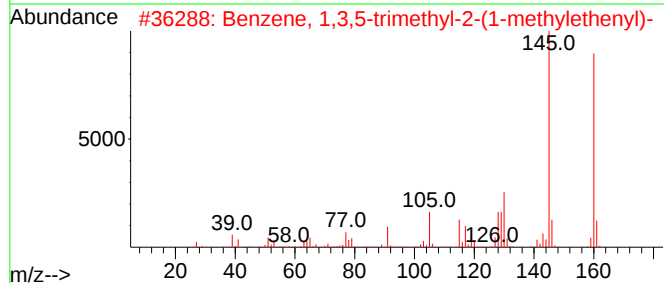
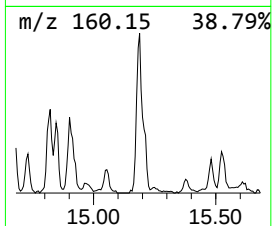
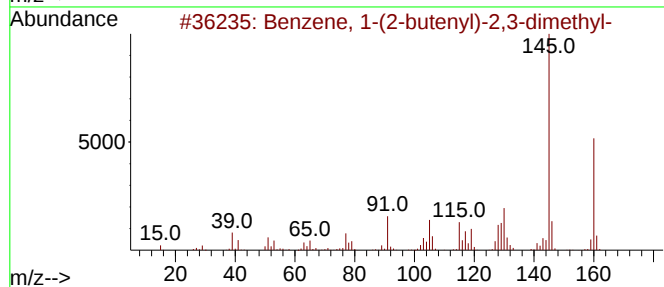
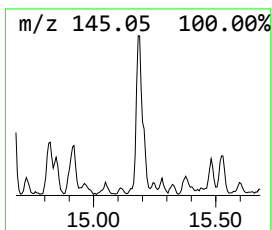
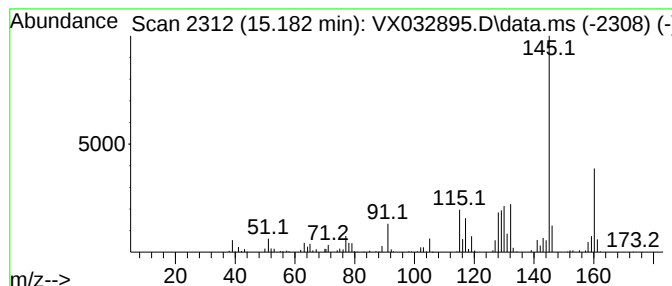
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title : SW846 8260

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

Peak Number 10 Benzene, 1-(2-butenyl)-2,3-... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.182	12.93 ug/l	129155	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-(2-butenyl)-2,3-dimet...	160	C12H16	054340-85-1	95
2			Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	95
3			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	020027-77-4	93
4			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	021564-91-0	90
5			Benzene, 4-(2-butenyl)-1,2-dimet...	160	C12H16	054340-86-2	89



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111722\
Data File : VX032895.D
Acq On : 17 Nov 2022 13:21
Operator : JC/MD
Sample : N5650-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
34271

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.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
Benzene, 2-ethe...	13.292	22.8	ug/l	228048	4	12.024	499251	50.0
Naphthalene, 1,...	13.420	11.4	ug/l	113670	4	12.024	499251	50.0
Benzene, (3-met...	13.585	12.2	ug/l	121275	4	12.024	499251	50.0
1H-Indene, 2,3-...	13.664	9.0	ug/l	90109	4	12.024	499251	50.0
Naphthalene, 1,...	13.853	9.7	ug/l	96713	4	12.024	499251	50.0
Benzene, (2-met...	13.926	12.7	ug/l	127117	4	12.024	499251	50.0
Naphthalene, 1,...	14.225	23.8	ug/l	237880	4	12.024	499251	50.0
Naphthalene, 1,...	14.481	19.1	ug/l	190858	4	12.024	499251	50.0
Naphthalene, 1,...	14.615	12.9	ug/l	128311	4	12.024	499251	50.0
Benzene, 1-(2-b...	15.182	12.9	ug/l	129155	4	12.024	499251	50.0