

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072423\
 Data File : VX036633.D
 Acq On : 24 Jul 2023 16:33
 Operator : JC/MD
 Sample : 03722-08
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 32416

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\82X072123W.M
 Title : SW846 8260

Signal : TIC: VX036633.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.453	57	60	66	rBV	8458	10680	1.18%	0.082%
2	1.605	73	85	92	rBV7	6524	22030	2.43%	0.168%
3	2.380	207	212	224	rVB	56348	97892	10.82%	0.747%
4	2.556	236	241	250	rVB2	7132	17964	1.99%	0.137%
5	2.965	302	308	316	rVB	4184	9871	1.09%	0.075%
6	4.568	563	571	580	rBV	9156	27466	3.04%	0.210%
7	5.385	694	705	718	rBV	103990	302175	33.39%	2.306%
8	5.556	721	733	752	rVB	172971	491890	54.36%	3.754%
9	5.958	786	799	819	rBV	114189	304827	33.69%	2.327%
10	6.763	921	931	955	rVB	278179	667044	73.71%	5.091%
11	8.580	1222	1229	1233	rBV2	10375	17801	1.97%	0.136%
12	8.647	1233	1240	1248	rVV	585073	904925	100.00%	6.907%
13	8.720	1248	1252	1257	rVB2	5935	10059	1.11%	0.077%
14	9.031	1296	1303	1310	rBV	7559	12314	1.36%	0.094%
15	9.519	1377	1383	1388	rBV	14377	21828	2.41%	0.167%
16	10.055	1465	1471	1488	rVB	582420	802034	88.63%	6.122%
17	10.195	1488	1494	1504	rBV2	10882	19600	2.17%	0.150%
18	10.299	1504	1511	1518	rVB	63966	90218	9.97%	0.689%
19	10.640	1562	1567	1572	rBV	90590	116127	12.83%	0.886%
20	10.854	1598	1602	1607	rBV2	15919	22201	2.45%	0.169%
21	10.963	1615	1620	1625	rBV	13452	16621	1.84%	0.127%
22	11.079	1634	1639	1653	rBV	477790	619719	68.48%	4.730%
23	11.305	1670	1676	1681	rBV	20879	26578	2.94%	0.203%
24	11.378	1682	1688	1696	rVV	138845	249953	27.62%	1.908%
25	11.451	1696	1700	1712	rVB	72495	93734	10.36%	0.715%
26	11.610	1722	1726	1733	rBV	134972	175233	19.36%	1.337%
27	11.750	1744	1749	1755	rVB	204848	254803	28.16%	1.945%
28	11.835	1757	1763	1766	rBV4	20593	31727	3.51%	0.242%
29	11.890	1766	1772	1776	rVB2	22523	33136	3.66%	0.253%
30	11.969	1781	1785	1788	rVV	39220	46135	5.10%	0.352%
31	12.018	1788	1793	1799	rVV	590170	758973	83.87%	5.793%
32	12.085	1799	1804	1812	rVB	257384	316958	35.03%	2.419%
33	12.244	1822	1830	1833	rBV	238341	341681	37.76%	2.608%
34	12.317	1838	1842	1851	rVB3	118548	172158	19.02%	1.314%
35	12.402	1851	1856	1861	rBV	22518	31764	3.51%	0.242%
36	12.463	1861	1866	1871	rVB	71915	86315	9.54%	0.659%

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072423\
 Data File : VX036633.D
 Acq On : 24 Jul 2023 16:33
 Operator : JC/MD
 Sample : 03722-08
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 32416

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\82X072123W.M
 Title : SW846 8260

37	12.530	1871	1877	1879	rBV	67916	90386	9.99%	0.690%
38	12.555	1879	1881	1885	rVV	66565	67964	7.51%	0.519%
39	12.610	1885	1890	1893	rVV	125162	150856	16.67%	1.151%
40	12.646	1893	1896	1900	rVV	47256	57379	6.34%	0.438%
41	12.701	1900	1905	1913	rVV2	141534	237399	26.23%	1.812%
42	12.798	1917	1921	1925	rVV	32246	46482	5.14%	0.355%
43	12.847	1925	1929	1934	rVB2	70168	90976	10.05%	0.694%
44	12.920	1934	1941	1944	rBV	97237	118226	13.06%	0.902%
45	12.963	1944	1948	1954	rVB	101333	128381	14.19%	0.980%
46	13.024	1954	1958	1960	rBV2	24922	35188	3.89%	0.269%
47	13.067	1964	1965	1968	rVV	13654	14037	1.55%	0.107%
48	13.097	1968	1970	1973	rVV2	11468	15841	1.75%	0.121%
49	13.140	1973	1977	1978	rVV2	28429	36920	4.08%	0.282%
50	13.164	1978	1981	1985	rVV	126313	161195	17.81%	1.230%
51	13.213	1985	1989	1992	rVV3	23645	31531	3.48%	0.241%
52	13.256	1992	1996	1997	rVV2	22936	30590	3.38%	0.233%
53	13.292	1997	2002	2007	rVV2	500013	673961	74.48%	5.144%
54	13.335	2007	2009	2012	rVV2	16784	24021	2.65%	0.183%
55	13.365	2012	2014	2017	rVV	17703	20051	2.22%	0.153%
56	13.420	2018	2023	2032	rVB	511171	615503	68.02%	4.698%
57	13.506	2032	2037	2039	rBV2	24879	34700	3.83%	0.265%
58	13.542	2039	2043	2046	rVV	47230	66313	7.33%	0.506%
59	13.585	2046	2050	2054	rVV3	89092	154624	17.09%	1.180%
60	13.640	2056	2059	2060	rVV	58887	70009	7.74%	0.534%
61	13.664	2060	2063	2068	rVV2	91501	137693	15.22%	1.051%
62	13.713	2068	2071	2074	rVV3	17352	24204	2.67%	0.185%
63	13.774	2074	2081	2087	rVV	146686	219739	24.28%	1.677%
64	13.853	2089	2094	2099	rVB	136674	177527	19.62%	1.355%
65	13.926	2099	2106	2110	rBV2	151248	222352	24.57%	1.697%
66	13.969	2110	2113	2117	rVV2	51221	67920	7.51%	0.518%
67	13.999	2117	2118	2122	rVV2	10718	11405	1.26%	0.087%
68	14.073	2126	2130	2134	rBV	55095	66556	7.35%	0.508%
69	14.109	2134	2136	2141	rVB4	7470	9703	1.07%	0.074%
70	14.225	2147	2155	2163	rBV	202515	356370	39.38%	2.720%
71	14.317	2167	2170	2174	rBV	33761	38446	4.25%	0.293%
72	14.371	2174	2179	2184	rBV2	82329	100347	11.09%	0.766%
73	14.420	2184	2187	2191	rVB	14972	19139	2.11%	0.146%
74	14.481	2191	2197	2202	rBV2	241098	321771	35.56%	2.456%
75	14.634	2214	2222	2225	rBV2	129909	239986	26.52%	1.832%
76	14.670	2225	2228	2233	rVB3	60750	86045	9.51%	0.657%

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072423\
 Data File : VX036633.D
 Acq On : 24 Jul 2023 16:33
 Operator : JC/MD
 Sample : 03722-08
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 32416

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\82X072123W.M
 Title : SW846 8260

77	14.725	2233	2237	2241	rBV2	17014	23388	2.58%	0.179%
78	14.780	2241	2246	2249	rBV	91273	117498	12.98%	0.897%
79	14.816	2249	2252	2255	rVV2	39570	56058	6.19%	0.428%
80	14.841	2255	2256	2261	rVB2	27456	27650	3.06%	0.211%
81	14.902	2261	2266	2273	rBV2	35298	65749	7.27%	0.502%
82	15.048	2283	2290	2296	rBV3	18146	32657	3.61%	0.249%
83	15.182	2304	2312	2314	rBV	53106	84739	9.36%	0.647%
84	15.200	2314	2315	2320	rVV	44016	48229	5.33%	0.368%
85	15.255	2320	2324	2331	rVB5	15480	29522	3.26%	0.225%
86	15.371	2338	2343	2347	rBV3	19396	30556	3.38%	0.233%
87	15.475	2353	2360	2365	rBV3	36599	61555	6.80%	0.470%
88	15.524	2365	2368	2373	rVV4	14978	28704	3.17%	0.219%
89	15.609	2377	2382	2386	rVV	47543	84935	9.39%	0.648%
90	15.646	2386	2388	2397	rVB2	28503	42870	4.74%	0.327%
91	15.841	2411	2420	2426	rBV4	13441	28982	3.20%	0.221%
92	15.987	2433	2444	2448	rBV3	5947	13862	1.53%	0.106%
93	16.084	2455	2460	2464	rBV2	7794	12922	1.43%	0.099%
94	16.475	2519	2524	2530	rBV	7929	15800	1.75%	0.121%

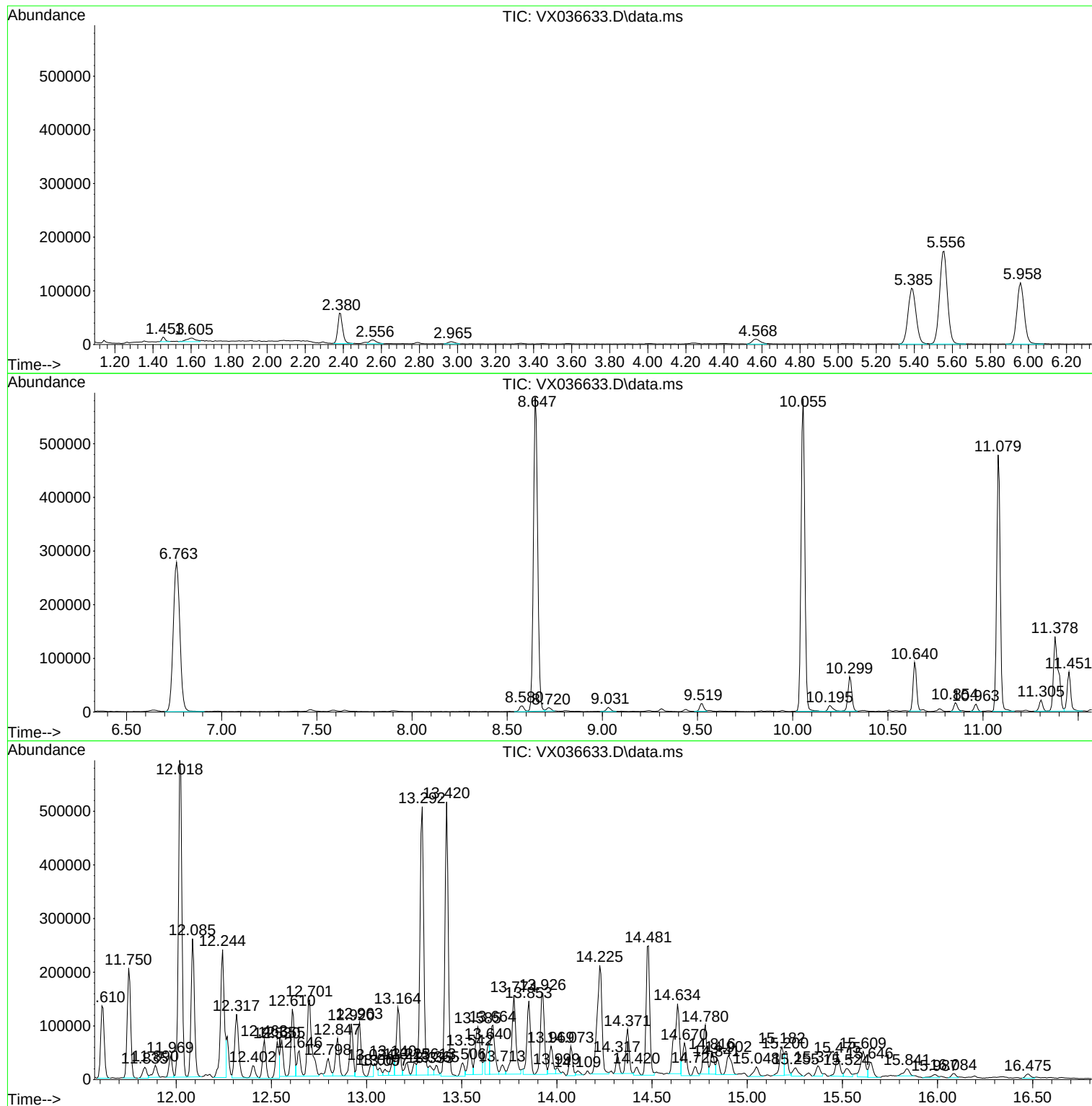
Sum of corrected areas: 13101846

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072423\
Data File : VX036633.D
Acq On : 24 Jul 2023 16:33
Operator : JC/MD
Sample : 03722-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
32416

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072423\
Data File : VX036633.D
Acq On : 24 Jul 2023 16:33
Operator : JC/MD
Sample : 03722-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
32416

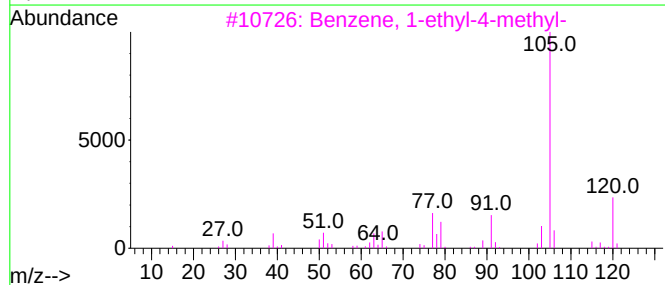
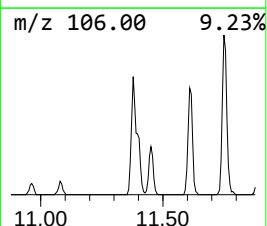
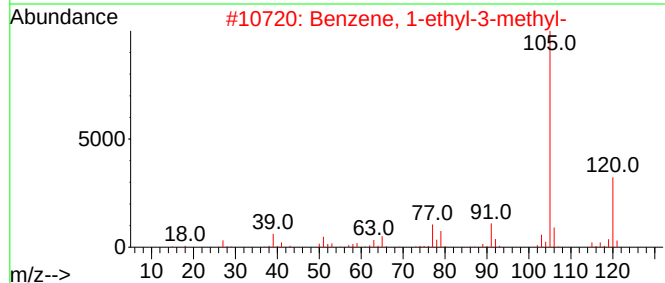
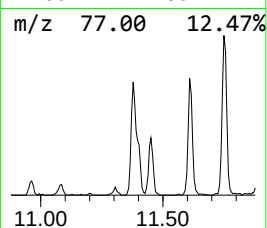
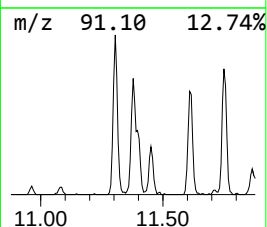
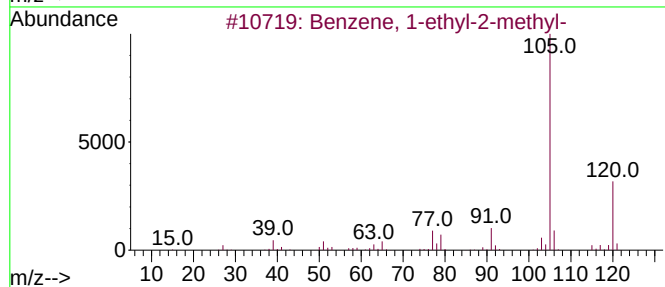
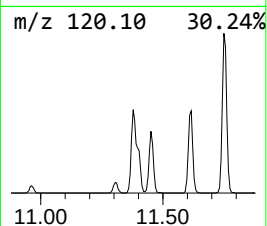
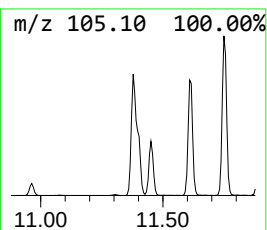
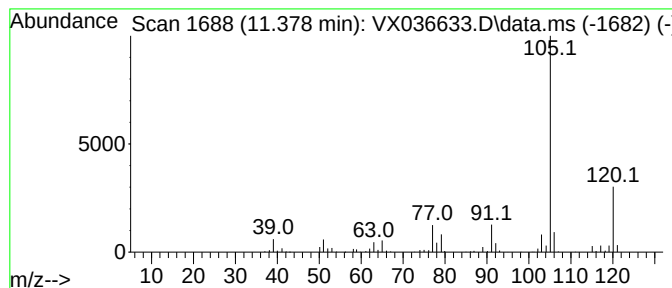
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1-ethyl-2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.378	16.47 ug/l	249953	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5			Mesitylene	120	C9H12	000108-67-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072423\
Data File : VX036633.D
Acq On : 24 Jul 2023 16:33
Operator : JC/MD
Sample : 03722-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
32416

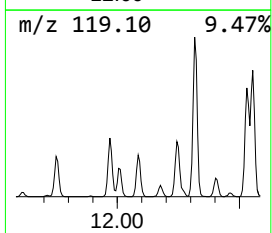
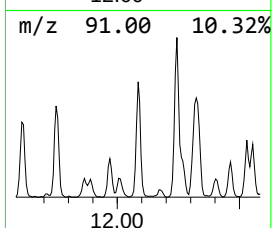
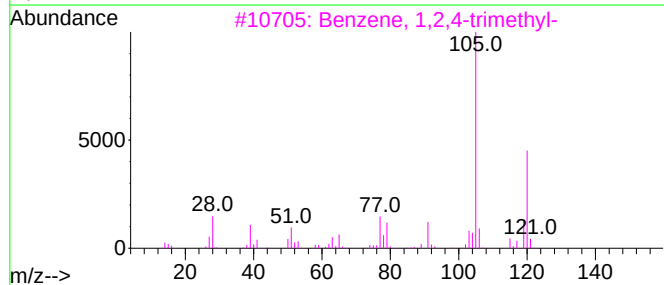
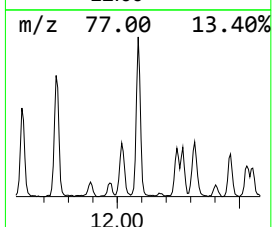
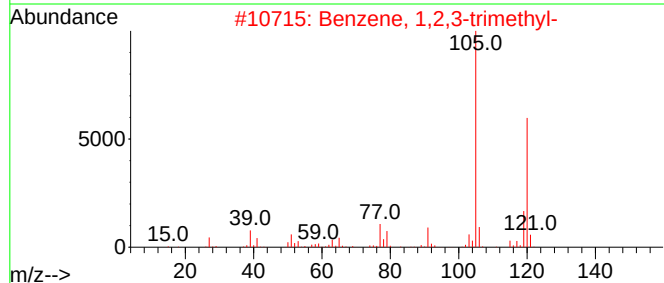
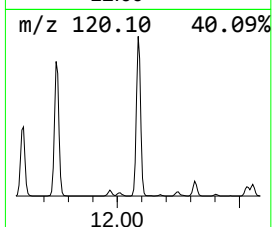
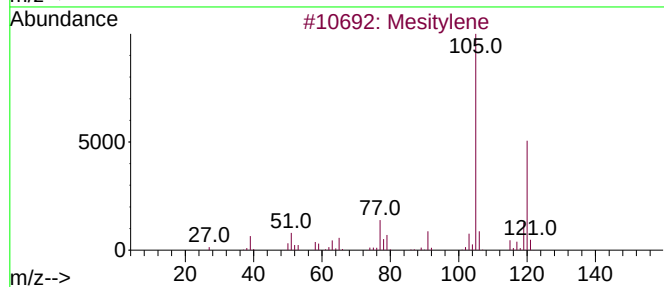
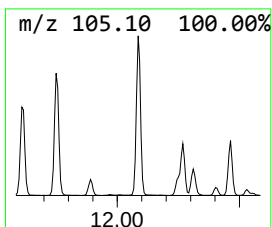
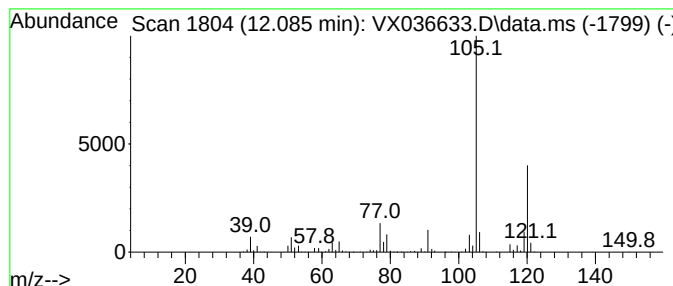
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.M
Quant Title : SW846 8260

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

Peak Number 2 Mesitylene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.085	20.88 ug/l	316958	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Mesitylene	120	C9H12	000108-67-8	97
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90
5			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072423\
Data File : VX036633.D
Acq On : 24 Jul 2023 16:33
Operator : JC/MD
Sample : 03722-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
32416

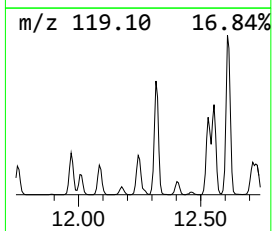
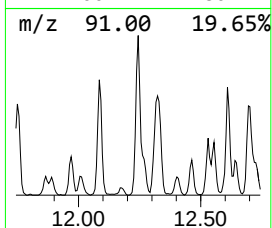
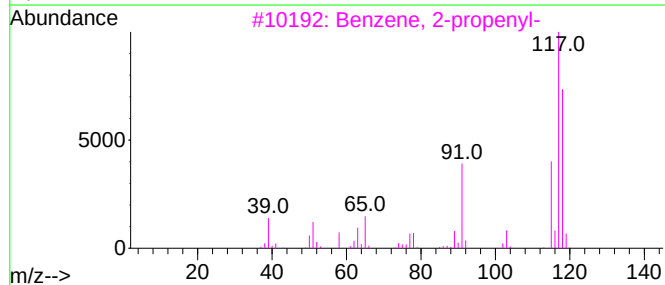
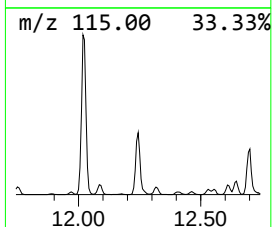
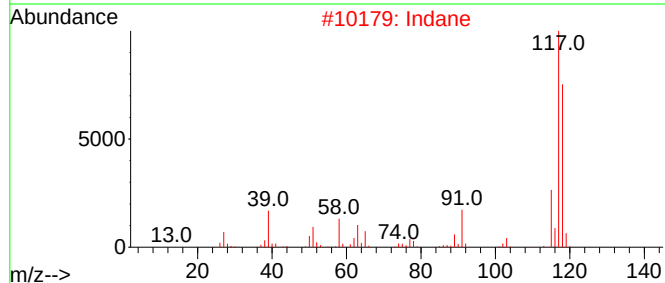
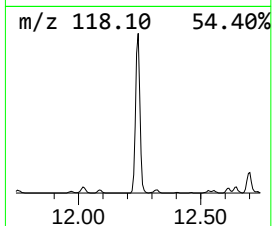
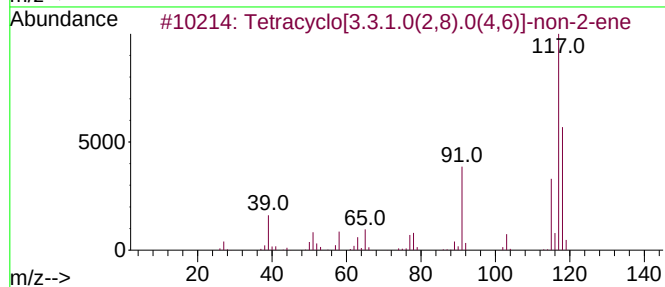
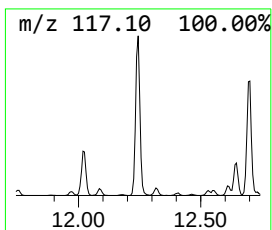
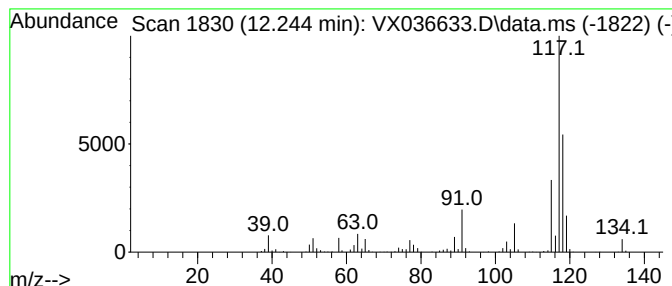
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.M
Quant Title : SW846 8260

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

Peak Number 3 Tetracyclo[3.3.1.0(2,8).0(4... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	22.51 ug/l	341681	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	76
2			Indane	118	C9H10	000496-11-7	76
3			Benzene, 2-propenyl-	118	C9H10	000300-57-2	76
4			Benzene, cyclopropyl-	118	C9H10	000873-49-4	76
5			Deltacyclene	118	C9H10	007785-10-6	62



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072423\
Data File : VX036633.D
Acq On : 24 Jul 2023 16:33
Operator : JC/MD
Sample : 03722-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
32416

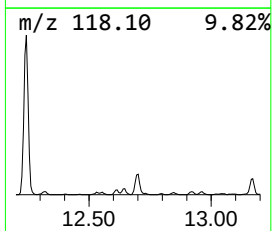
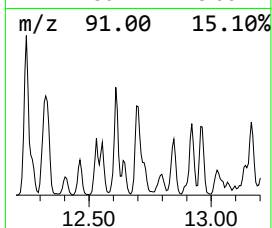
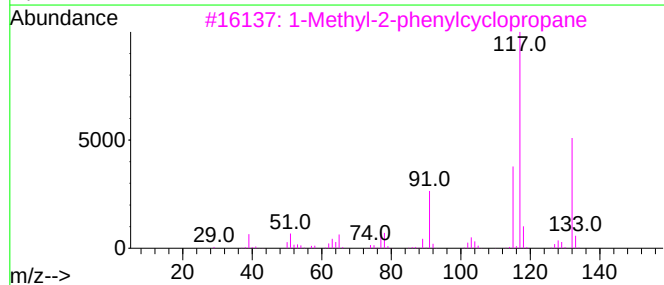
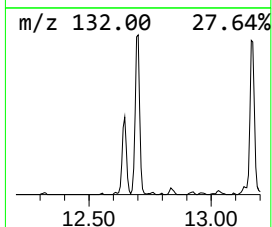
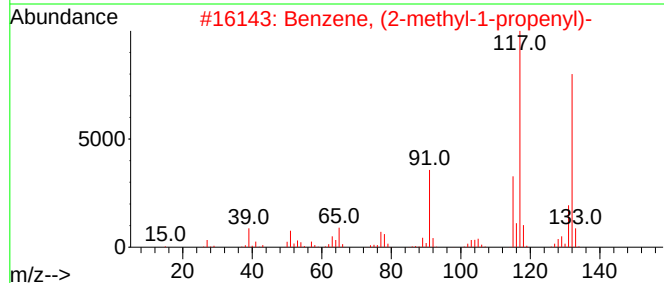
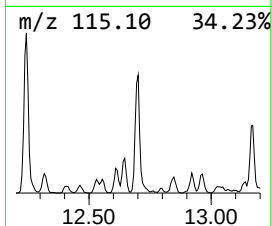
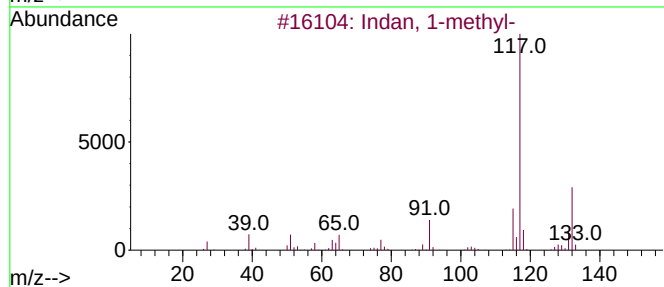
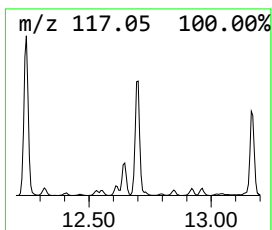
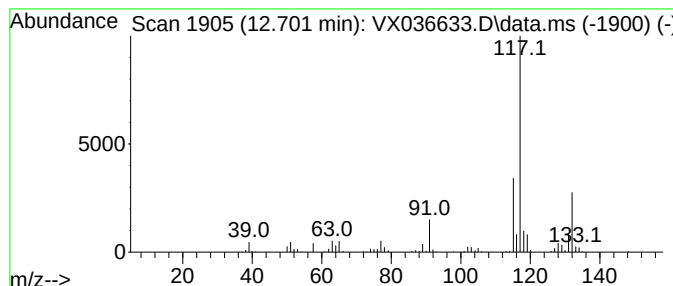
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.M
Quant Title : SW846 8260

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

Peak Number 4 Indan, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.701	15.64 ug/l	237399	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indan, 1-methyl-	132	C10H12	000767-58-8	93
2			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	87
3			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	87
4			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	86
5			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	83



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072423\
Data File : VX036633.D
Acq On : 24 Jul 2023 16:33
Operator : JC/MD
Sample : 03722-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
32416

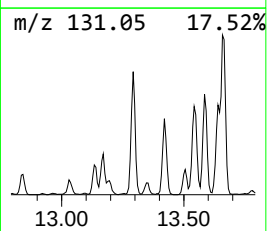
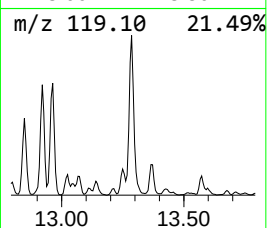
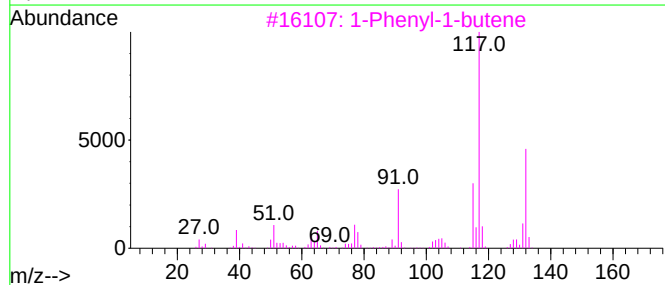
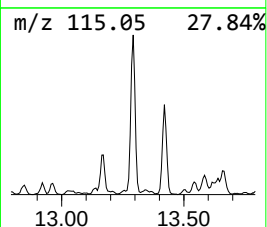
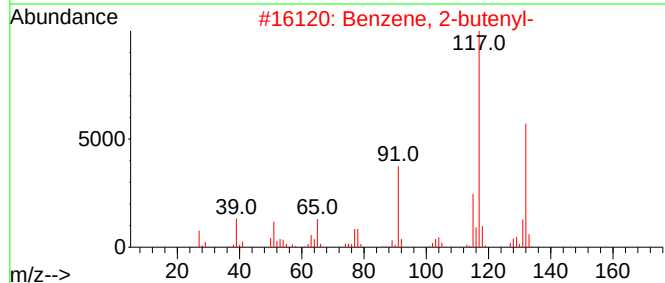
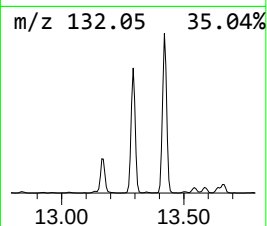
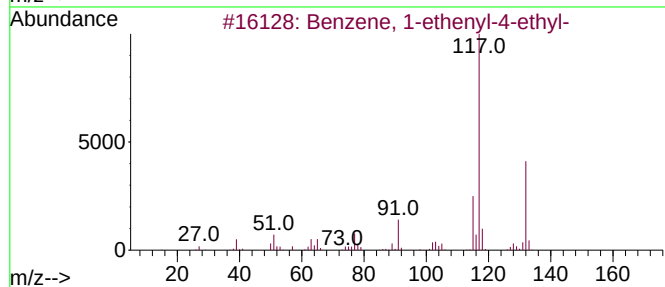
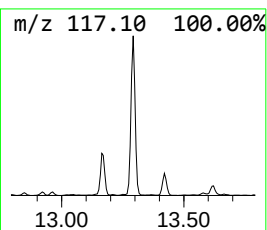
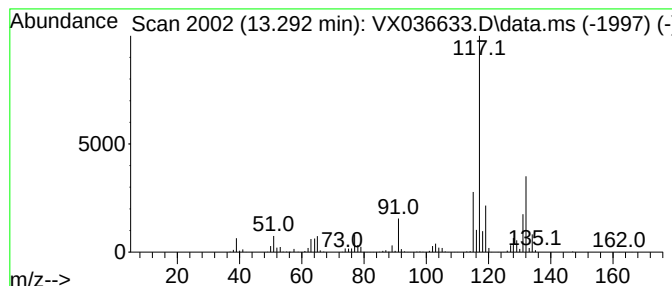
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.M
Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	44.40 ug/l	673961	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	89
2			Benzene, 2-butenyl-	132	C10H12	001560-06-1	81
3			1-Phenyl-1-butene	132	C10H12	000824-90-8	81
4			3-Phenylbut-1-ene	132	C10H12	000934-10-1	81
5			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	81



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072423\
Data File : VX036633.D
Acq On : 24 Jul 2023 16:33
Operator : JC/MD
Sample : 03722-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

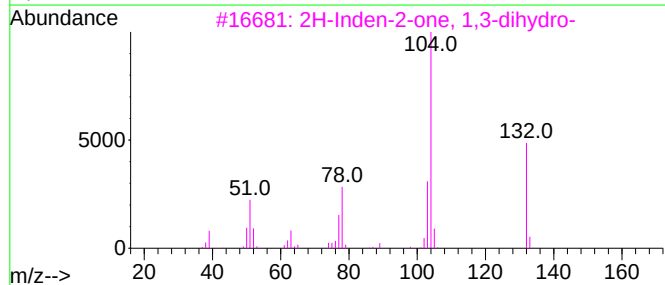
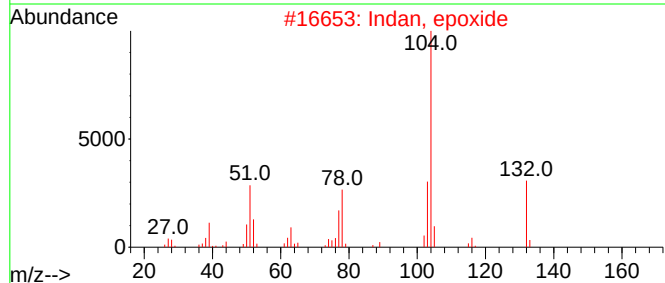
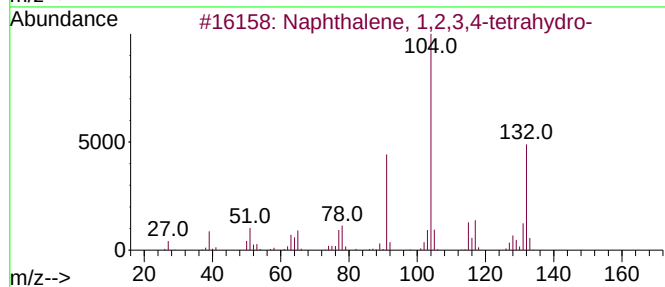
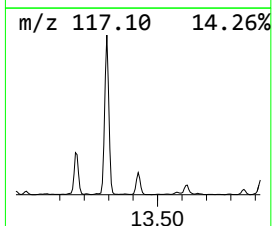
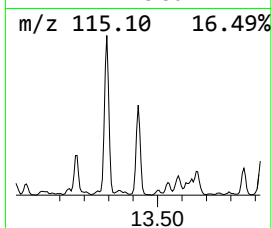
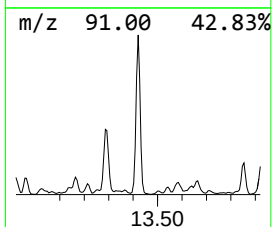
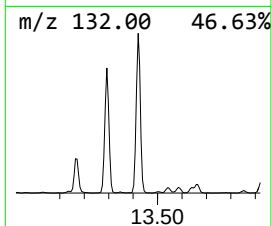
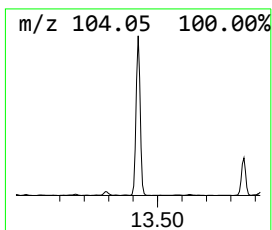
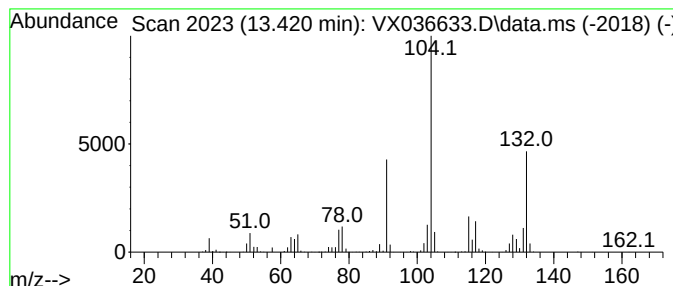
Instrument :
MSVOA_X
ClientSampleId :
32416

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.M
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 2

R.T.	EstConc	Area	Relative to ISTD			R.T.
13.420	40.55 ug/l	615503	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	96
2	Indan, epoxide		132	C9H8O	000768-22-9	62
3	2H-Inden-2-one, 1,3-dihydro-		132	C9H8O	000615-13-4	50
4	Benzene, cyclobutyl-		132	C10H12	004392-30-7	45
5	N-Ethylbenzimidoyl bromide		211	C9H10BrN	041182-87-0	40



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072423\
Data File : VX036633.D
Acq On : 24 Jul 2023 16:33
Operator : JC/MD
Sample : 03722-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
32416

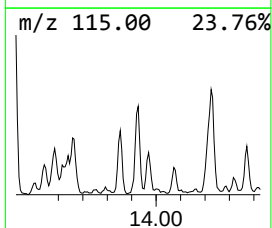
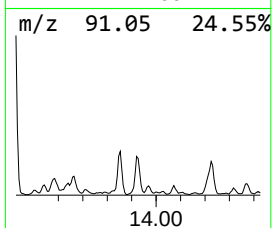
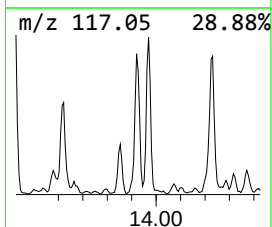
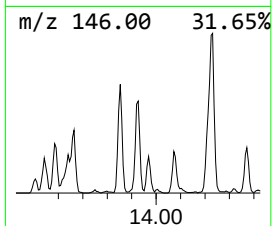
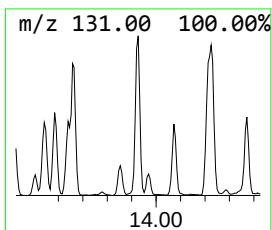
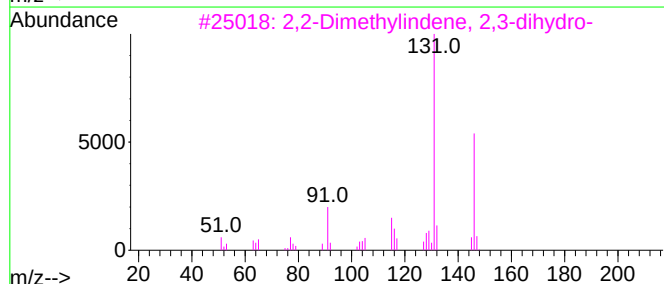
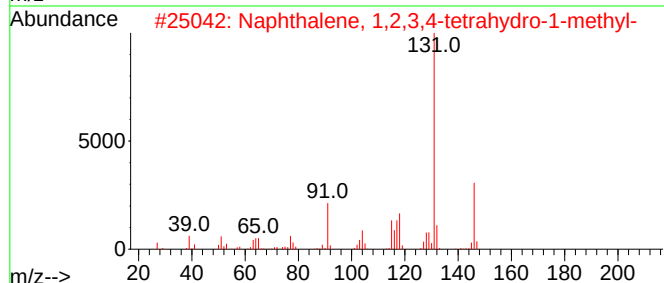
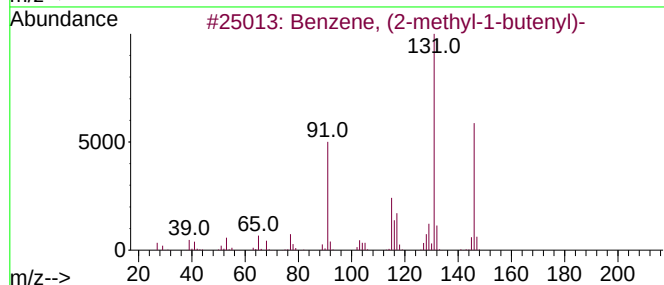
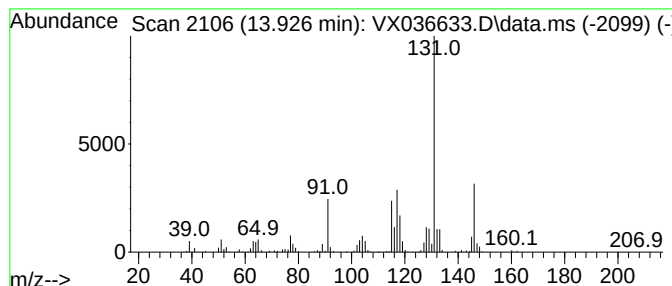
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.M
Quant Title : SW846 8260

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

Peak Number 7 Benzene, (2-methyl-1-butenyl)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.926	14.65 ug/l	222352	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			2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	90
4			1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	87
5			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	76



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072423\
Data File : VX036633.D
Acq On : 24 Jul 2023 16:33
Operator : JC/MD
Sample : 03722-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
32416

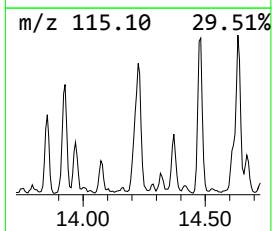
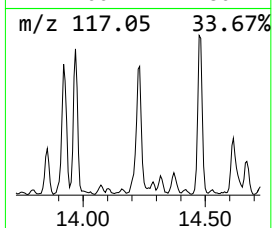
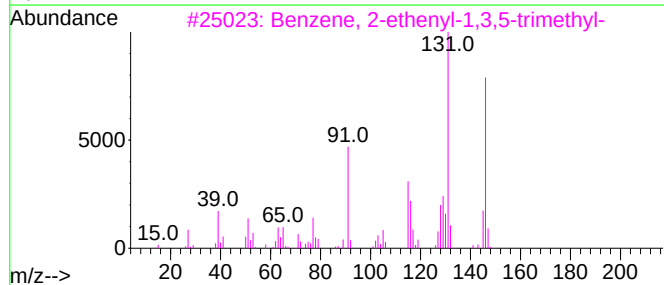
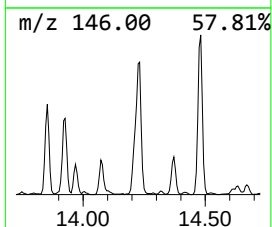
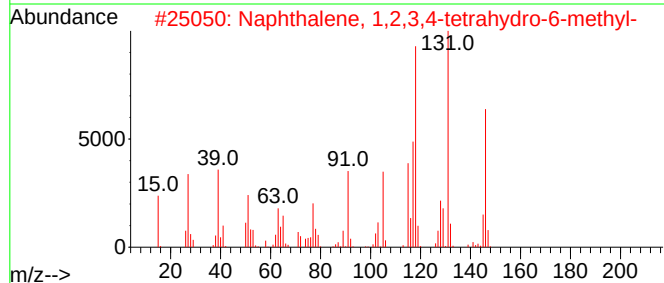
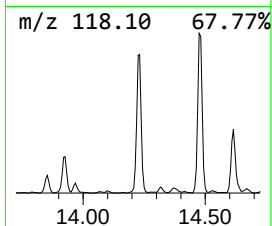
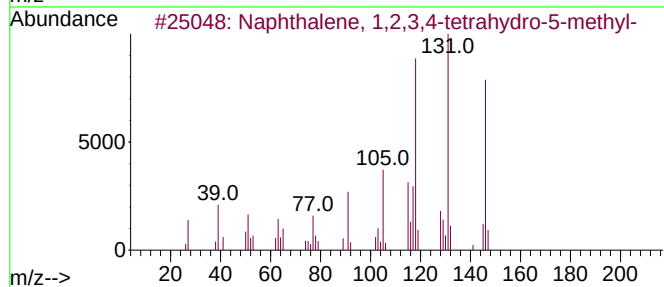
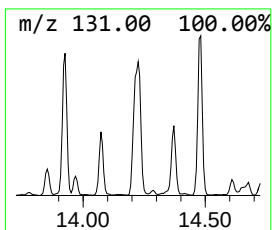
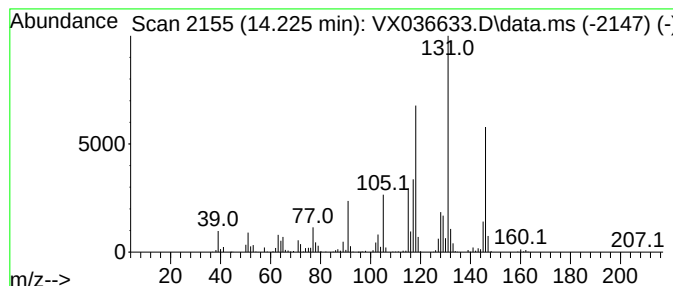
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.M
Quant Title : SW846 8260

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.
14.225	23.48 ug/l	356370	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, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	86
4			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	86
5			Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	76



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072423\
Data File : VX036633.D
Acq On : 24 Jul 2023 16:33
Operator : JC/MD
Sample : 03722-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
32416

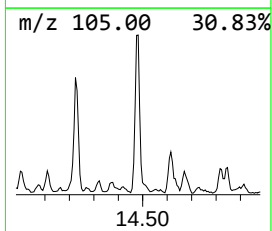
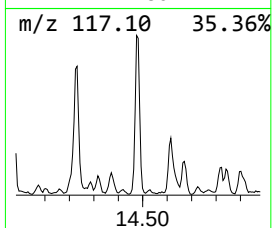
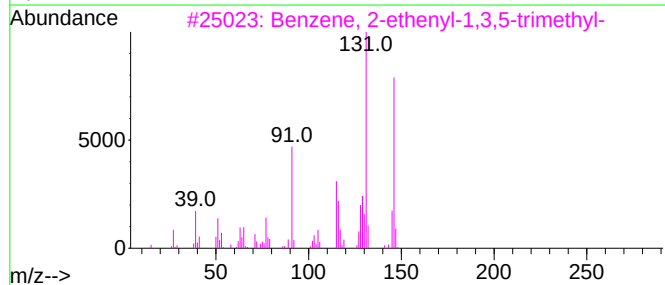
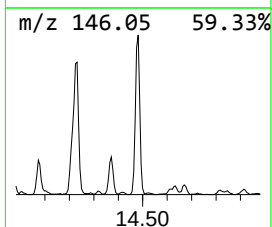
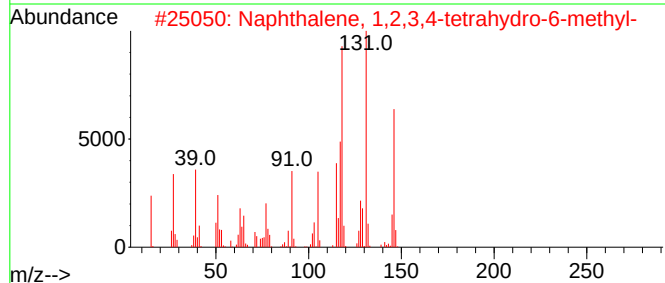
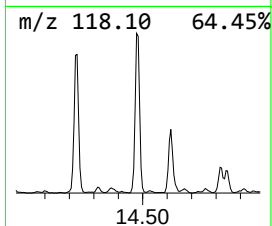
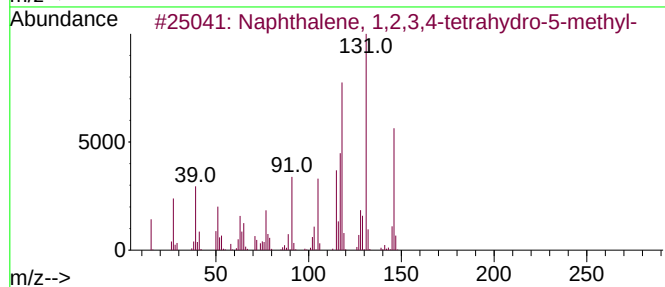
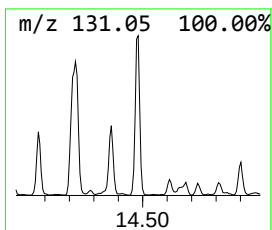
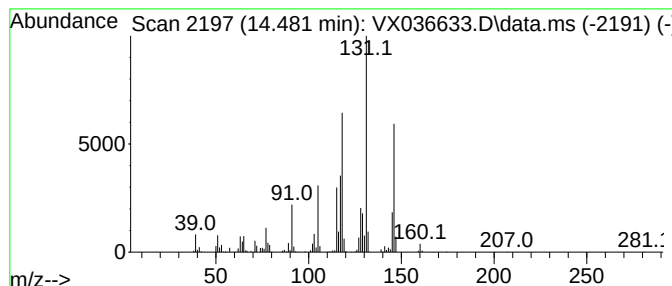
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.M
Quant Title : SW846 8260

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.
14.481	21.20 ug/l	321771	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	94
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	94
3			Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	89
4			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	70
5			Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	70



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072423\
Data File : VX036633.D
Acq On : 24 Jul 2023 16:33
Operator : JC/MD
Sample : 03722-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
32416

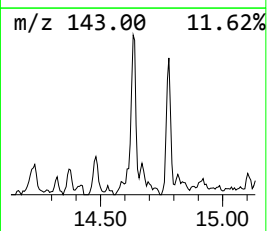
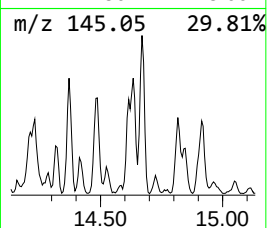
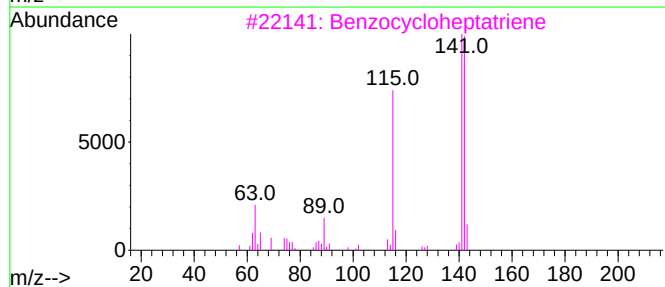
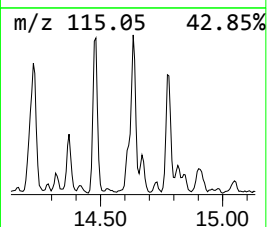
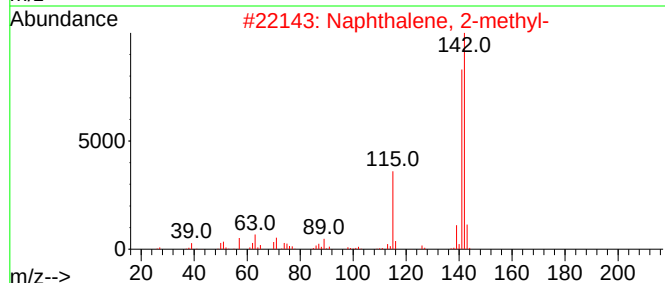
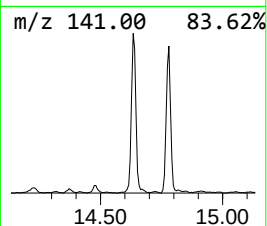
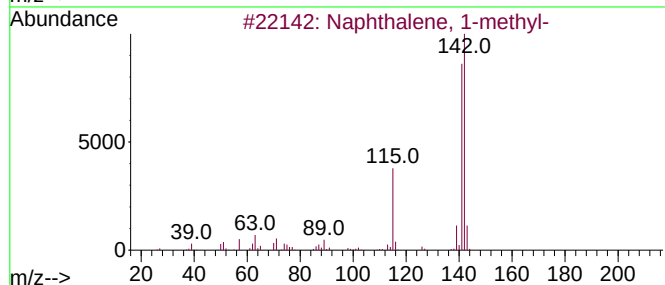
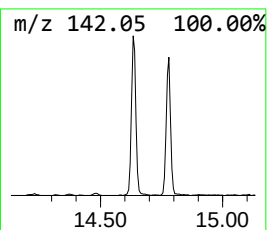
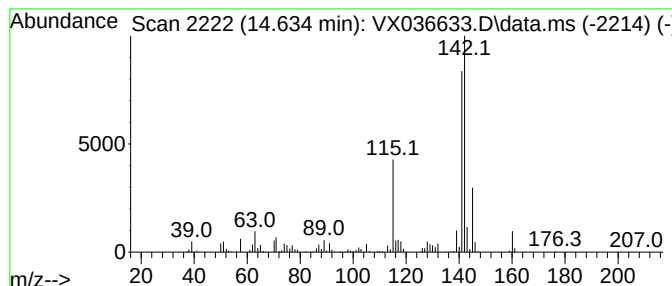
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.634	15.81 ug/l	239986	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3			Benzocycloheptatriene	142	C11H10	000264-09-5	81
4			1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	52
5			1-Naphthaleneethylamine	171	C12H13N	004735-50-6	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072423\
Data File : VX036633.D
Acq On : 24 Jul 2023 16:33
Operator : JC/MD
Sample : 03722-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
32416

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.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, 1-ethy...	11.378	16.5	ug/l	249953	4	12.024	758973	50.0
Mesitylene	12.085	20.9	ug/l	316958	4	12.024	758973	50.0
Tetracyclo[3.3....	12.244	22.5	ug/l	341681	4	12.024	758973	50.0
Indan, 1-methyl-	12.701	15.6	ug/l	237399	4	12.024	758973	50.0
Benzene, 1-ethe...	13.292	44.4	ug/l	673961	4	12.024	758973	50.0
Naphthalene, 1,...	13.420	40.5	ug/l	615503	4	12.024	758973	50.0
Benzene, (2-met...	13.926	14.7	ug/l	222352	4	12.024	758973	50.0
Naphthalene, 1,...	14.225	23.5	ug/l	356370	4	12.024	758973	50.0
Naphthalene, 1,...	14.481	21.2	ug/l	321771	4	12.024	758973	50.0
Naphthalene, 1-...	14.634	15.8	ug/l	239986	4	12.024	758973	50.0