

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072423\
 Data File : VX036634.D
 Acq On : 24 Jul 2023 16:57
 Operator : JC/MD
 Sample : 03722-09
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 33712

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: VX036634.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.593	73	83	84	rBV8	5458	12918	1.44%	0.080%
2	2.380	207	212	225	rVB	22777	40632	4.54%	0.252%
3	2.550	237	240	254	rVB	3873	10141	1.13%	0.063%
4	5.385	693	705	721	rBV	101674	298024	33.30%	1.849%
5	5.550	721	732	748	rVB	169933	483147	53.98%	2.998%
6	5.958	785	799	815	rBV	113973	300441	33.57%	1.864%
7	6.763	921	931	946	rBV	275901	657076	73.41%	4.078%
8	8.580	1222	1229	1234	rBV	7912	14393	1.61%	0.089%
9	8.647	1234	1240	1249	rVV	575385	895079	100.00%	5.554%
10	8.720	1249	1252	1258	rVB	6612	10889	1.22%	0.068%
11	9.525	1377	1384	1389	rBV2	8353	13886	1.55%	0.086%
12	10.055	1465	1471	1480	rBV	597181	799903	89.37%	4.964%
13	10.195	1488	1494	1500	rBV	20326	28353	3.17%	0.176%
14	10.299	1504	1511	1518	rVB	116149	159514	17.82%	0.990%
15	10.640	1562	1567	1582	rVB	121438	160382	17.92%	0.995%
16	10.854	1598	1602	1613	rVB2	9480	15174	1.70%	0.094%
17	10.963	1613	1620	1626	rBV	18120	24141	2.70%	0.150%
18	11.079	1634	1639	1651	rBV	491133	623538	69.66%	3.869%
19	11.305	1671	1676	1682	rBV	35583	44677	4.99%	0.277%
20	11.378	1682	1688	1696	rVV	233206	389848	43.55%	2.419%
21	11.451	1696	1700	1710	rVB	121583	147885	16.52%	0.918%
22	11.610	1722	1726	1733	rVB	169395	222353	24.84%	1.380%
23	11.750	1744	1749	1756	rVB	479621	574642	64.20%	3.566%
24	11.829	1757	1762	1766	rBV2	16450	28156	3.15%	0.175%
25	11.890	1766	1772	1777	rVV	27494	39869	4.45%	0.247%
26	11.969	1780	1785	1788	rVV	50731	58553	6.54%	0.363%
27	12.018	1788	1793	1799	rVV	598462	788413	88.08%	4.893%
28	12.085	1799	1804	1812	rVB	334127	426973	47.70%	2.650%
29	12.244	1822	1830	1833	rBV	324409	482802	53.94%	2.996%
30	12.317	1838	1842	1851	rVB3	161746	237541	26.54%	1.474%
31	12.402	1851	1856	1861	rBV3	27339	37823	4.23%	0.235%
32	12.463	1861	1866	1871	rVB	82466	102075	11.40%	0.633%
33	12.530	1871	1877	1879	rBV	114256	147916	16.53%	0.918%
34	12.555	1879	1881	1885	rVB	101057	105267	11.76%	0.653%
35	12.609	1885	1890	1893	rBV	159098	193808	21.65%	1.203%
36	12.646	1893	1896	1900	rVV	59781	71341	7.97%	0.443%

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37	12.701	1900	1905	1913	rVB2	172447	276434	30.88%	1.715%
38	12.798	1917	1921	1925	rVV2	29792	40494	4.52%	0.251%
39	12.847	1925	1929	1934	rVB2	86108	113141	12.64%	0.702%
40	12.920	1934	1941	1944	rBV	129388	153133	17.11%	0.950%
41	12.963	1944	1948	1954	rVB	177612	211894	23.67%	1.315%
42	13.024	1954	1958	1960	rBV	25277	35119	3.92%	0.218%
43	13.067	1964	1965	1968	rVV	15487	13369	1.49%	0.083%
44	13.097	1968	1970	1973	rVV3	11442	14971	1.67%	0.093%
45	13.140	1973	1977	1978	rVV2	28633	37198	4.16%	0.231%
46	13.164	1978	1981	1986	rVV	220453	281778	31.48%	1.749%
47	13.213	1986	1989	1992	rVB	27603	29415	3.29%	0.183%
48	13.250	1992	1995	1997	rBV2	30665	38525	4.30%	0.239%
49	13.292	1997	2002	2007	rVV2	677654	890745	99.52%	5.528%
50	13.329	2007	2008	2012	rVV3	14738	19194	2.14%	0.119%
51	13.365	2012	2014	2017	rVB	18048	17933	2.00%	0.111%
52	13.420	2018	2023	2032	rVB	690143	813402	90.87%	5.048%
53	13.506	2032	2037	2040	rBV2	39839	57526	6.43%	0.357%
54	13.542	2040	2043	2046	rVV	68040	85603	9.56%	0.531%
55	13.585	2046	2050	2054	rVV2	103602	174052	19.45%	1.080%
56	13.658	2055	2062	2068	rVB2	106732	234562	26.21%	1.456%
57	13.713	2068	2071	2074	rBV4	8848	11051	1.23%	0.069%
58	13.774	2074	2081	2087	rVV	271614	357184	39.91%	2.217%
59	13.853	2089	2094	2098	rVB	165956	210743	23.54%	1.308%
60	13.926	2098	2106	2110	rBV2	179384	251191	28.06%	1.559%
61	13.969	2110	2113	2117	rVB	57286	63146	7.05%	0.392%
62	14.073	2126	2130	2134	rBV	94721	112606	12.58%	0.699%
63	14.109	2134	2136	2141	rVV4	9672	13737	1.53%	0.085%
64	14.164	2141	2145	2148	rVV2	8355	12576	1.41%	0.078%
65	14.225	2148	2155	2163	rVV	363257	585177	65.38%	3.631%
66	14.316	2167	2170	2175	rVV	39965	50896	5.69%	0.316%
67	14.371	2175	2179	2183	rVV	110469	133718	14.94%	0.830%
68	14.414	2183	2186	2191	rVB	14955	19201	2.15%	0.119%
69	14.475	2191	2196	2202	rBV2	304445	400735	44.77%	2.487%
70	14.633	2214	2222	2226	rBV2	309386	484137	54.09%	3.004%
71	14.670	2226	2228	2233	rVB2	66898	78799	8.80%	0.489%
72	14.725	2233	2237	2240	rBV2	20124	24677	2.76%	0.153%
73	14.780	2240	2246	2249	rVV	199845	253945	28.37%	1.576%
74	14.816	2249	2252	2255	rVV2	45288	65630	7.33%	0.407%
75	14.847	2255	2257	2261	rVB3	31971	32839	3.67%	0.204%
76	14.902	2261	2266	2273	rBV2	40196	74822	8.36%	0.464%

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77	15.048	2283	2290	2297	rBV	17928	29829	3.33%	0.185%
78	15.182	2303	2312	2314	rBV	67457	111520	12.46%	0.692%
79	15.200	2314	2315	2320	rVV	60419	64653	7.22%	0.401%
80	15.255	2320	2324	2331	rVB4	15957	29657	3.31%	0.184%
81	15.371	2339	2343	2347	rBV2	28897	45389	5.07%	0.282%
82	15.475	2352	2360	2365	rVV	61356	104438	11.67%	0.648%
83	15.524	2365	2368	2373	rVV3	19062	35441	3.96%	0.220%
84	15.609	2377	2382	2386	rVV	72326	125643	14.04%	0.780%
85	15.652	2386	2389	2396	rVB	40255	59734	6.67%	0.371%
86	15.841	2411	2420	2425	rBV2	17852	38076	4.25%	0.236%
87	15.987	2439	2444	2448	rBV2	7230	12506	1.40%	0.078%
88	16.084	2455	2460	2466	rBV2	10337	17929	2.00%	0.111%
89	16.475	2519	2524	2529	rBV	9159	16870	1.88%	0.105%

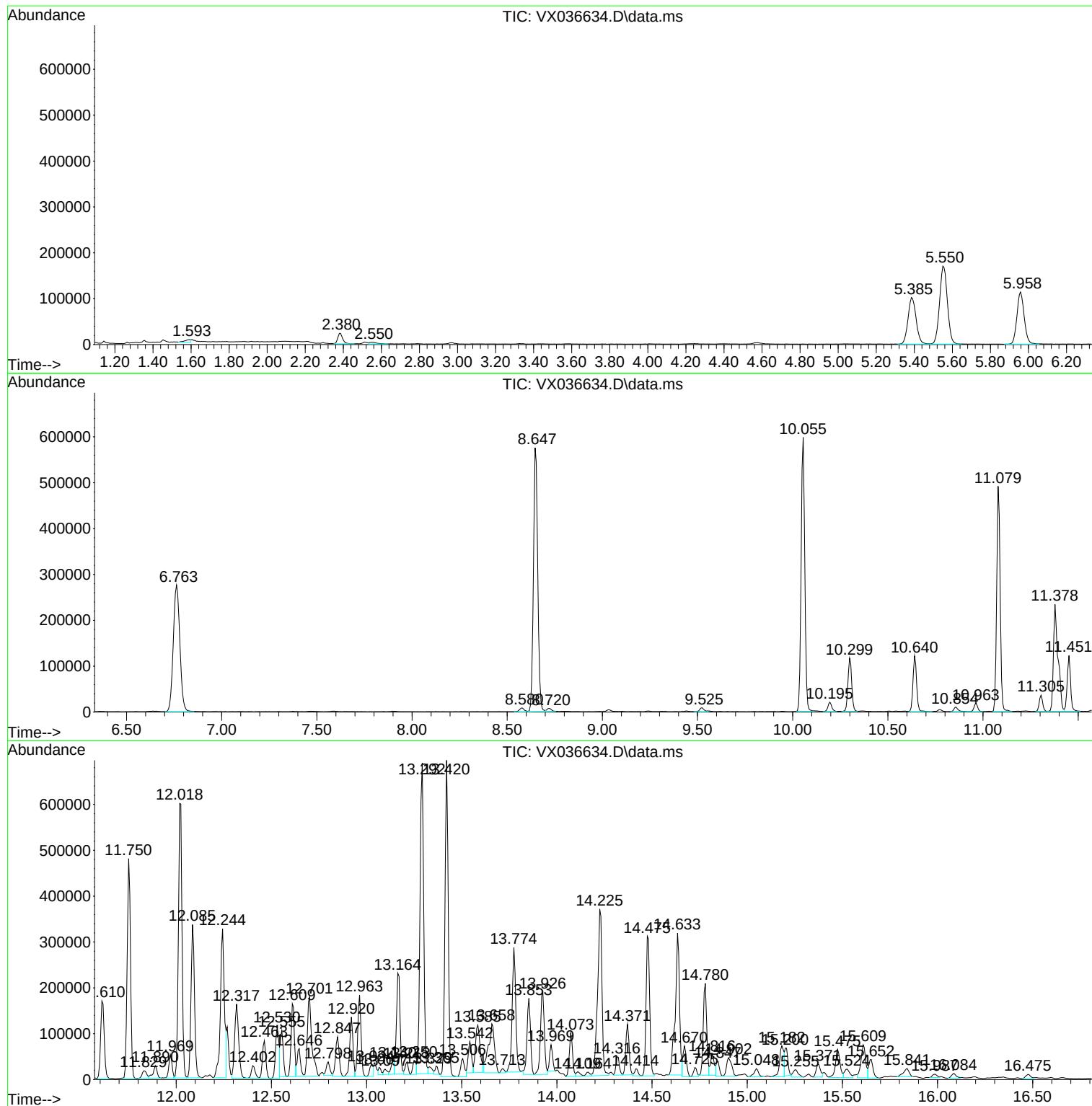
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Instrument :
MSVOA_X
ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.M
Quant Title : SW846 8260

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



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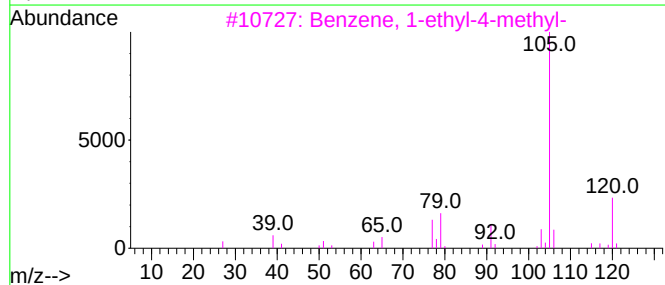
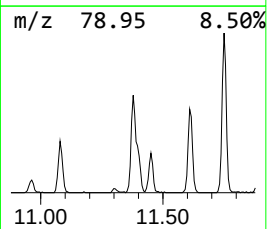
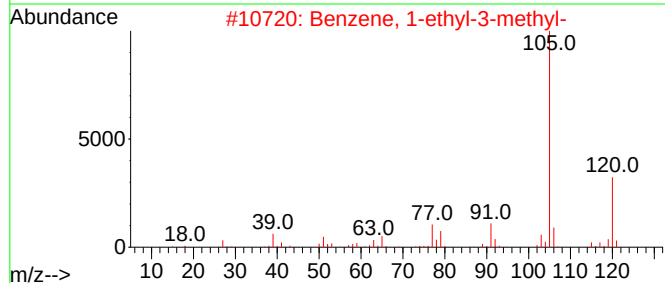
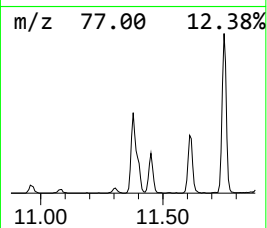
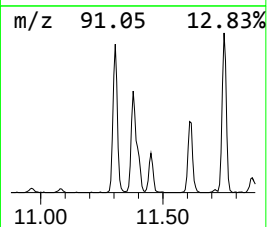
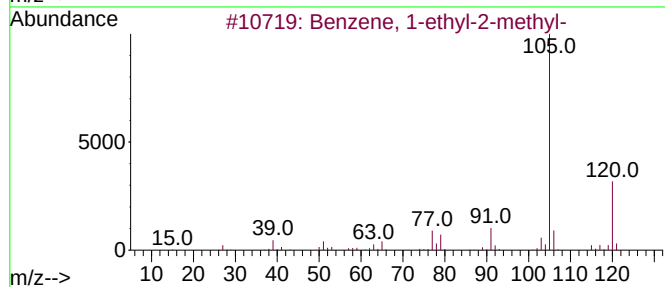
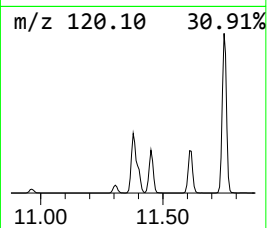
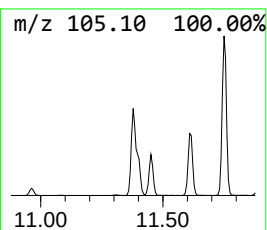
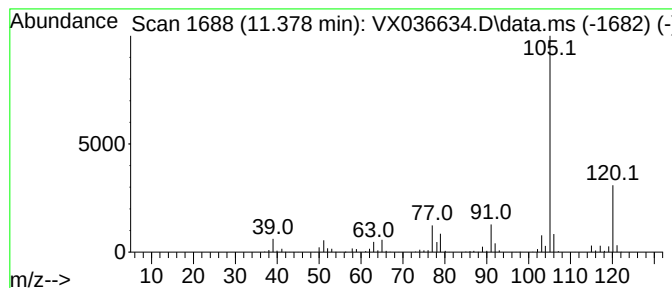
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1-ethyl-2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.378	24.72 ug/l	389848	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	94
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4			Mesitylene	120	C9H12	000108-67-8	91
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



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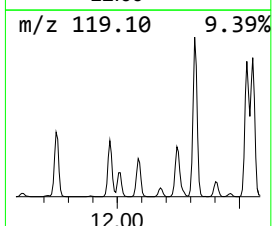
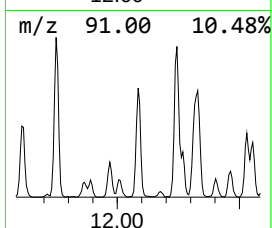
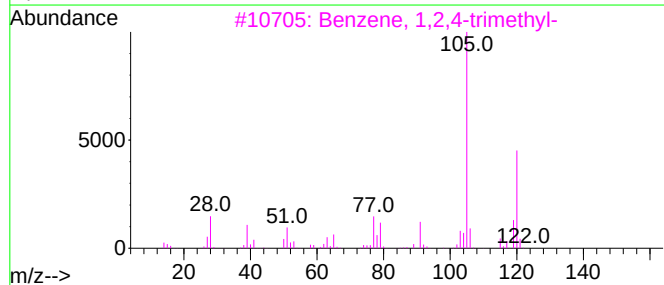
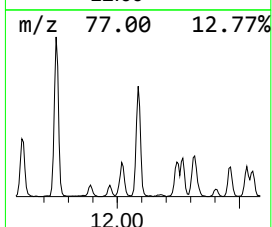
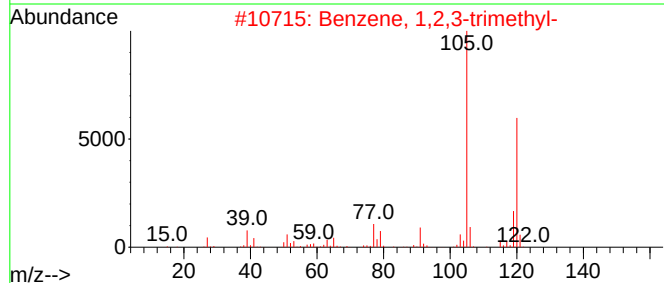
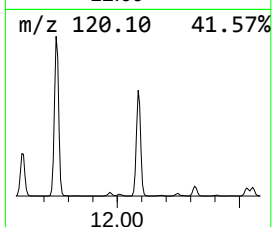
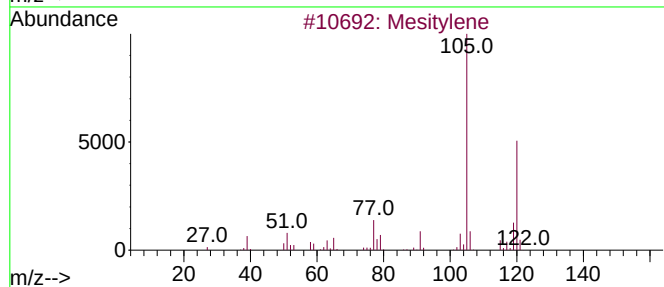
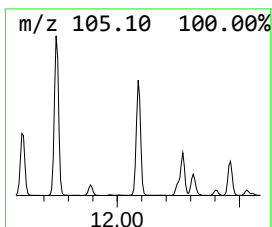
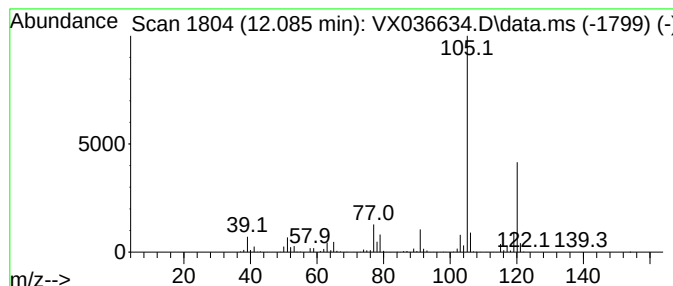
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Mesitylene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.085	27.08 ug/l	426973	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-2-methyl-	120	C9H12	000611-14-3	93
5			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91



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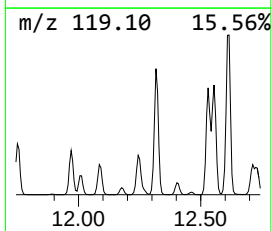
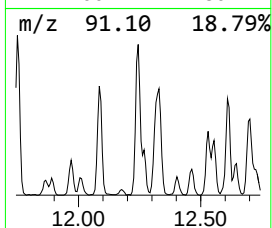
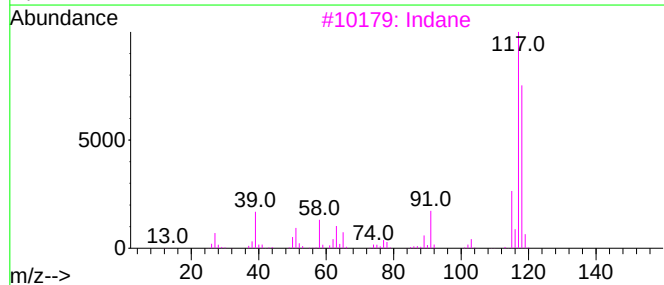
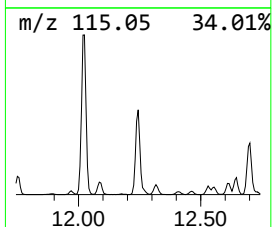
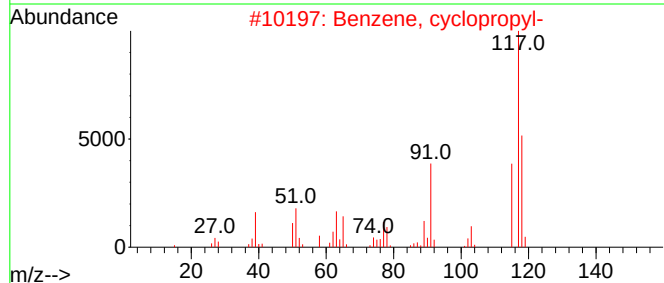
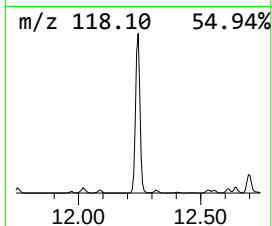
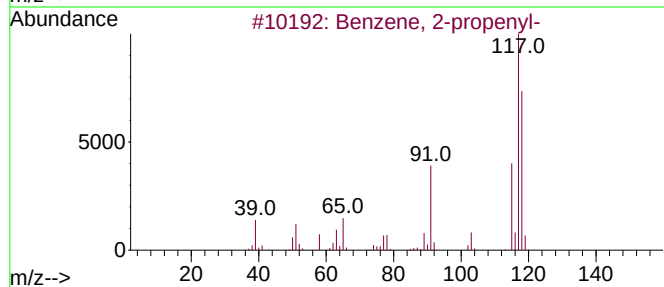
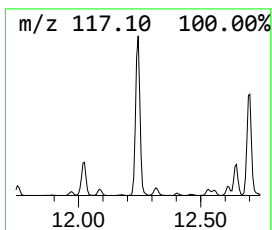
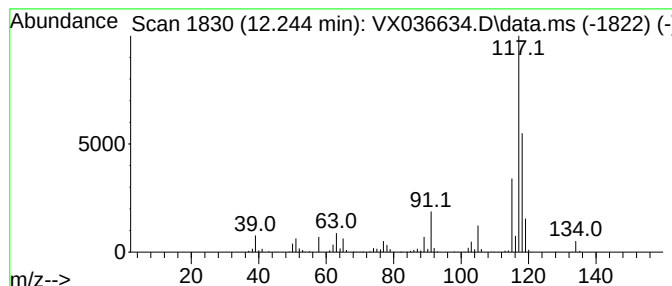
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.M
Quant Title : SW846 8260

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

Peak Number 3 Benzene, 2-propenyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	30.62 ug/l	482802	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-propenyl-	118	C9H10	000300-57-2	76
2			Benzene, cyclopropyl-	118	C9H10	000873-49-4	76
3			Indane	118	C9H10	000496-11-7	76
4			Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	64
5			Deltacyclene	118	C9H10	007785-10-6	58



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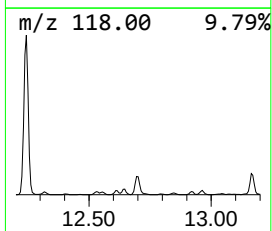
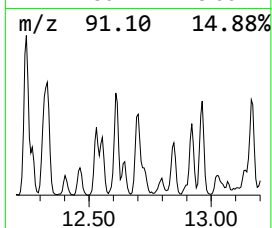
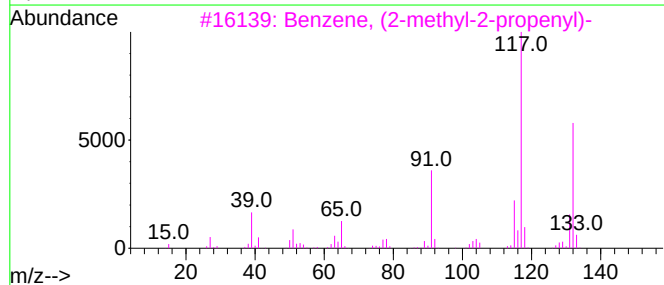
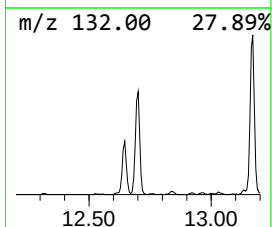
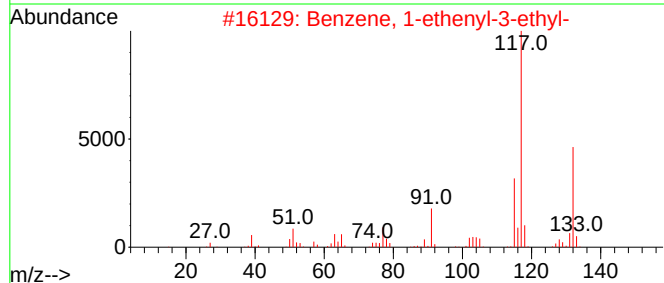
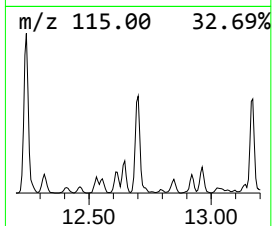
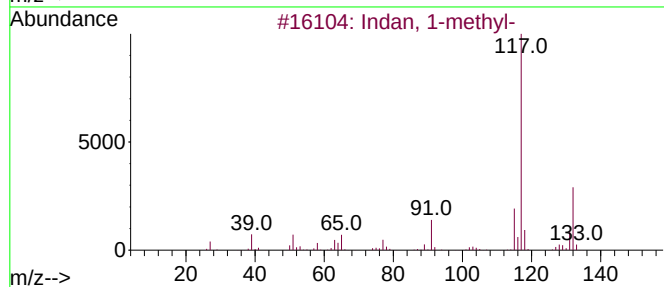
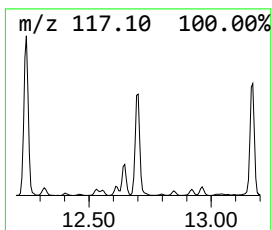
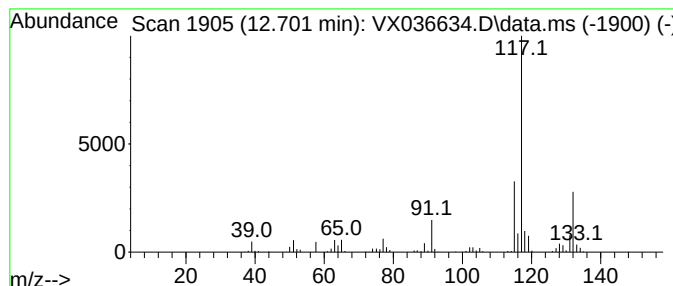
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Quant Title : SW846 8260

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

Peak Number 4 Indan, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.701	17.53 ug/l	276434	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, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	90
3		Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	87
4		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	87
5		1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072423\
Data File : VX036634.D
Acq On : 24 Jul 2023 16:57
Operator : JC/MD
Sample : 03722-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
33712

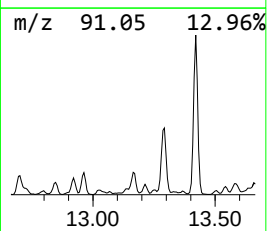
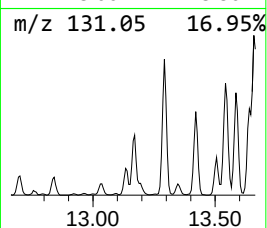
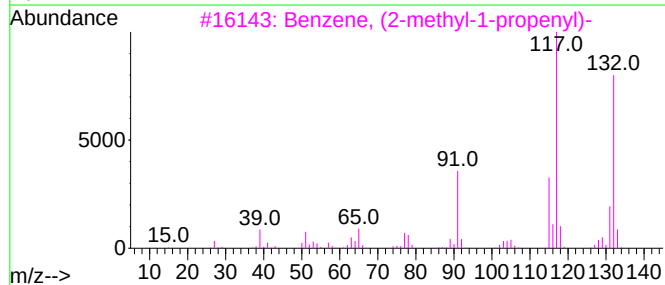
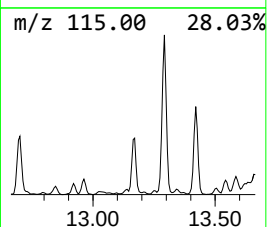
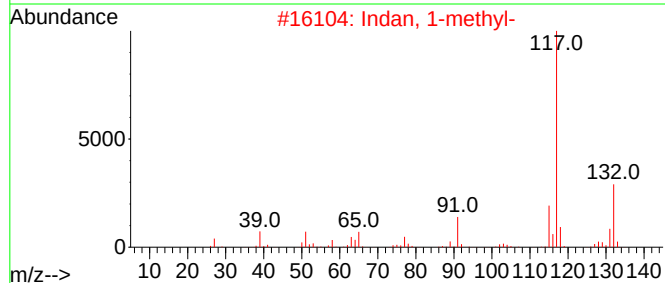
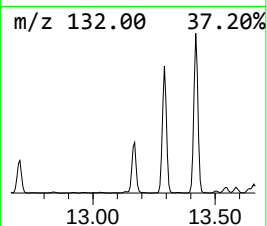
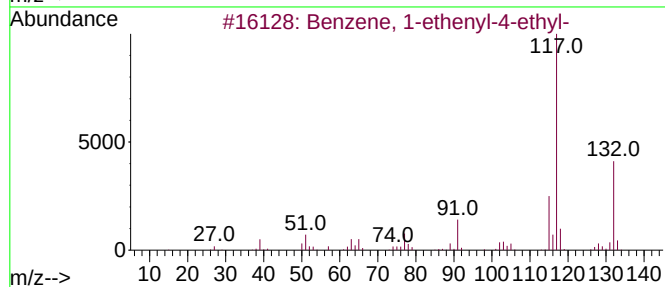
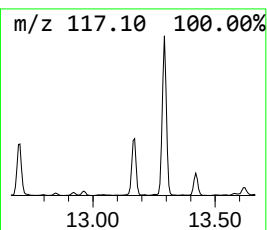
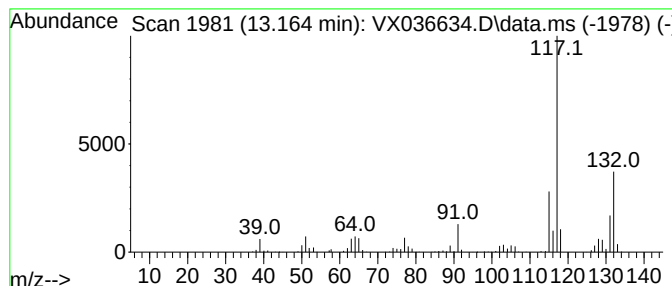
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.164	17.87 ug/l	281778	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			Indan, 1-methyl-	132	C10H12	000767-58-8	91
3			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	91
4			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	91
5			Benzene, 2-butenyl-	132	C10H12	001560-06-1	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072423\
Data File : VX036634.D
Acq On : 24 Jul 2023 16:57
Operator : JC/MD
Sample : 03722-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
33712

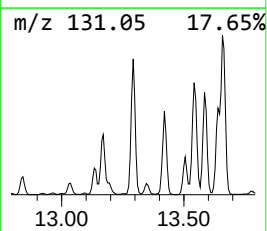
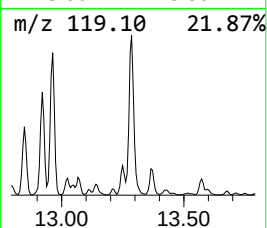
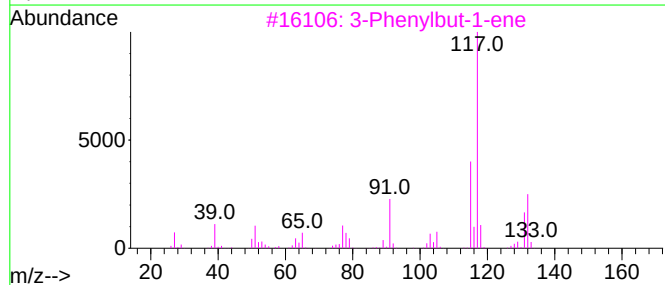
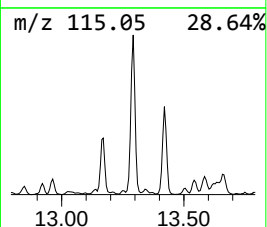
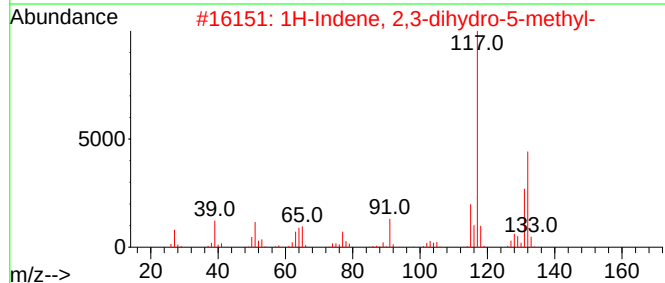
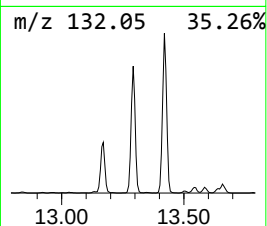
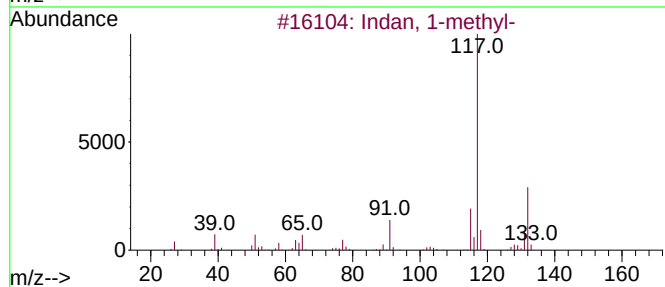
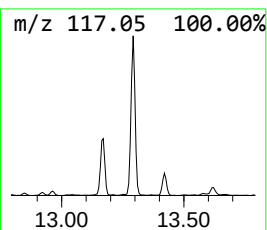
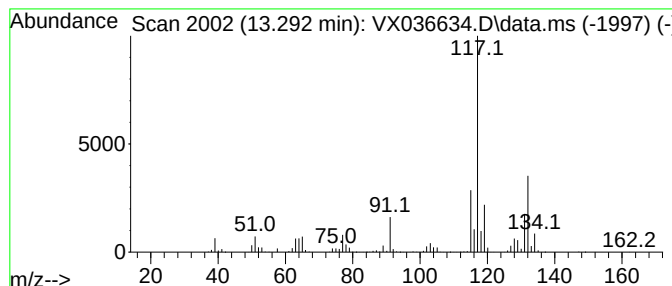
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Quant Title : SW846 8260

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

Peak Number 6 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	56.49 ug/l	890745	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		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	87
3		3-Phenylbut-1-ene	132	C10H12	000934-10-1	81
4		Benzene, 2-butenyl-	132	C10H12	001560-06-1	74
5		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	74



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072423\
Data File : VX036634.D
Acq On : 24 Jul 2023 16:57
Operator : JC/MD
Sample : 03722-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

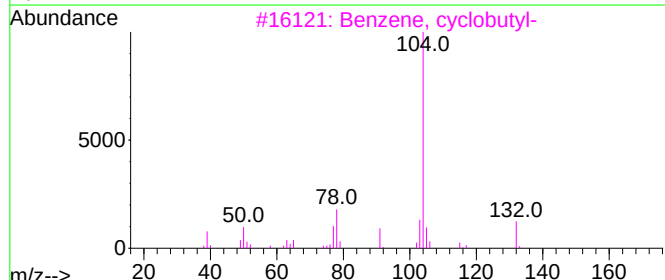
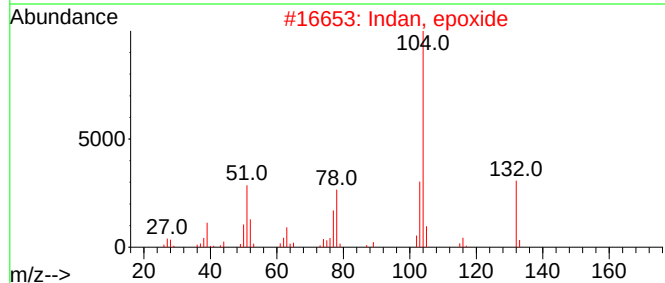
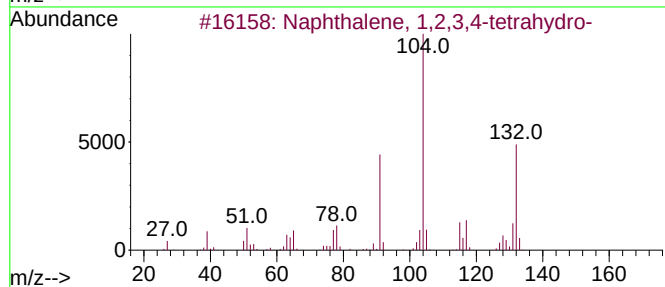
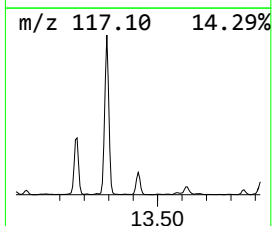
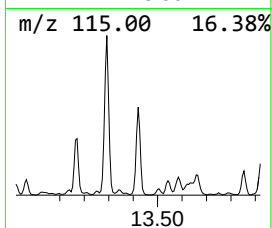
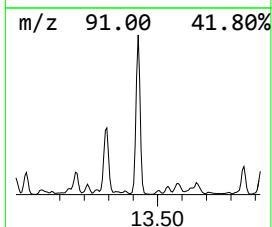
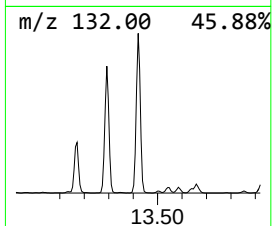
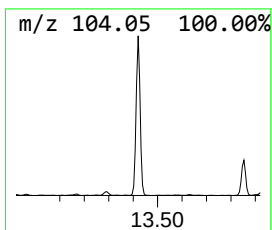
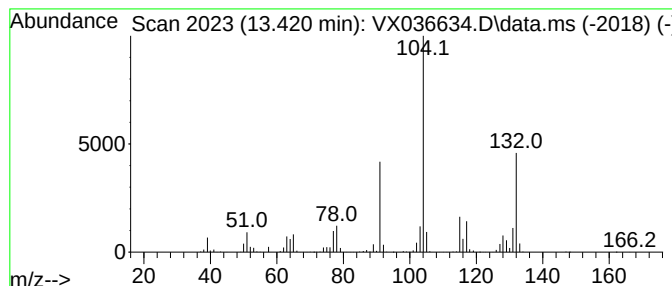
Instrument :
MSVOA_X
ClientSampleId :
33712

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 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD			R.T.
13.420	51.58 ug/l	813402	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	58
3	Benzene, cyclobutyl-		132	C10H12	004392-30-7	45
4	N-Ethylbenzimidoyl bromide		211	C9H10BrN	041182-87-0	40
5	Benzene, 1,1'-(1,2-cyclobutanedi...		208	C16H16	020071-09-4	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072423\
Data File : VX036634.D
Acq On : 24 Jul 2023 16:57
Operator : JC/MD
Sample : 03722-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
33712

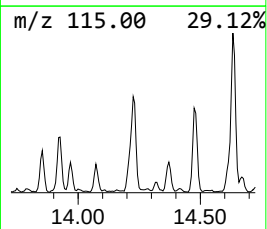
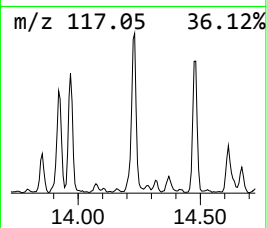
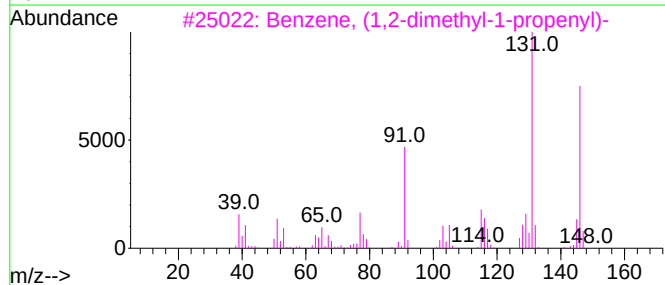
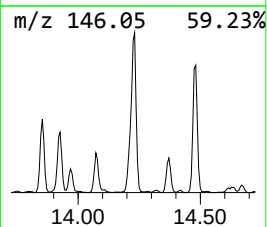
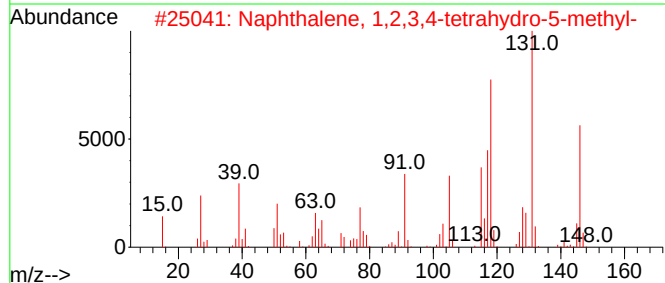
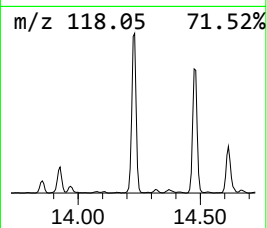
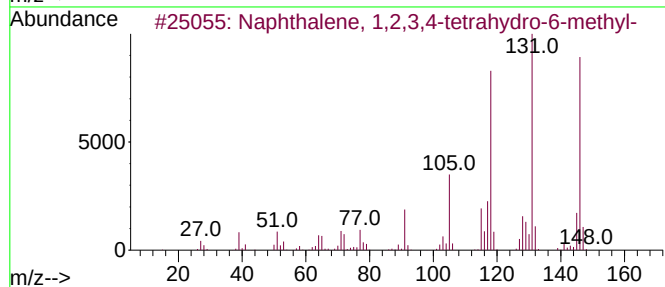
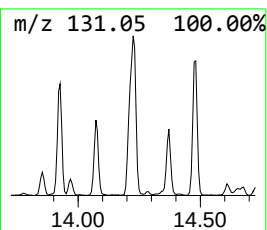
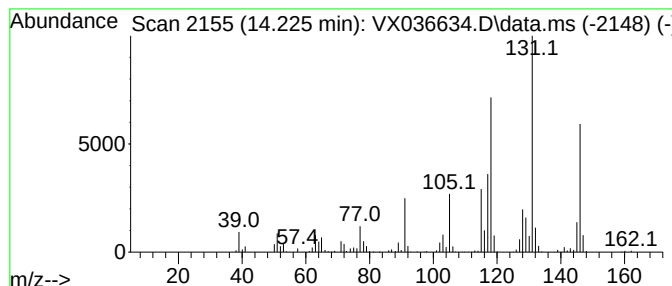
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.M
Quant Title : SW846 8260

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	37.11 ug/l	585177	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	97
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	96
3			Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	92
4			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	70
5			Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	70



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072423\
Data File : VX036634.D
Acq On : 24 Jul 2023 16:57
Operator : JC/MD
Sample : 03722-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

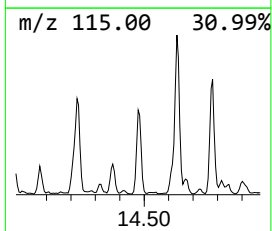
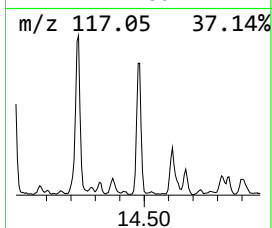
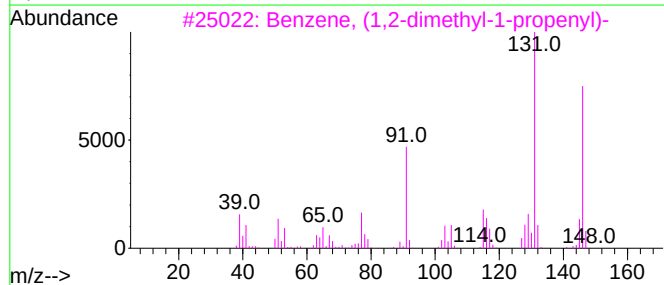
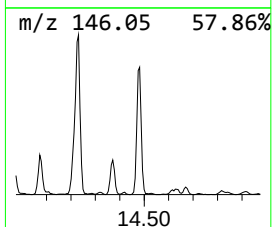
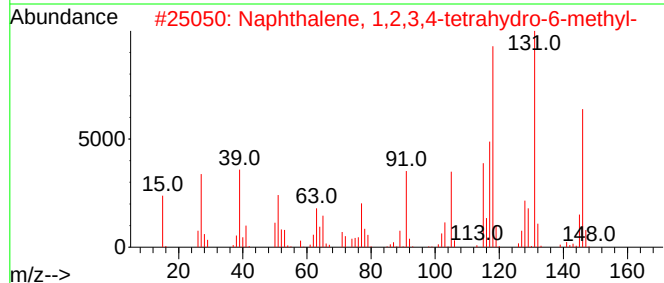
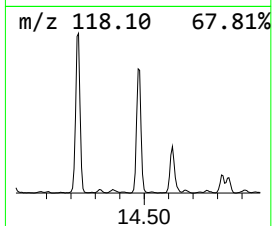
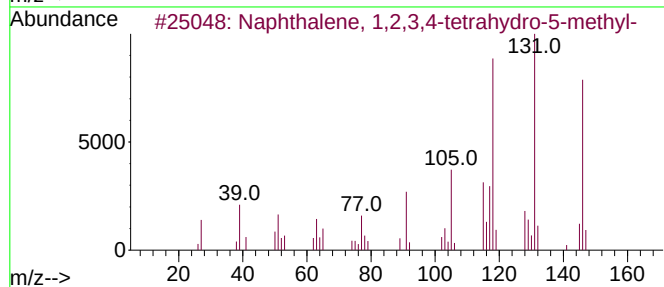
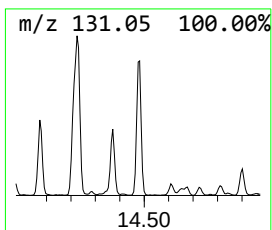
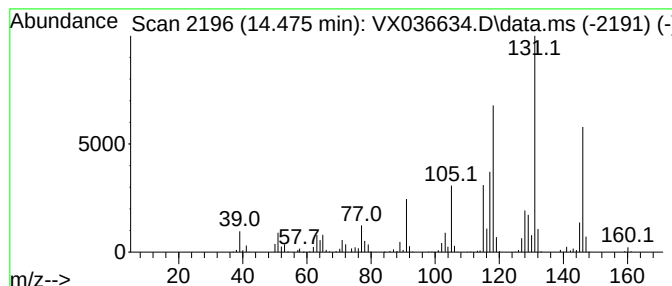
Instrument :
MSVOA_X
ClientSampleId :
33712

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 7

R.T.	EstConc	Area	Relative to ISTD			R.T.
14.475	25.41 ug/l	400735	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	96
3	Benzene, (1,2-dimethyl-1-propenyl)-		146	C11H14	000769-57-3	92
4	Benzene, (3-methyl-2-butenyl)-		146	C11H14	004489-84-3	70
5	Benzene, 2-ethenyl-1,3,5-trimethyl-		146	C11H14	000769-25-5	68



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072423\
Data File : VX036634.D
Acq On : 24 Jul 2023 16:57
Operator : JC/MD
Sample : 03722-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
33712

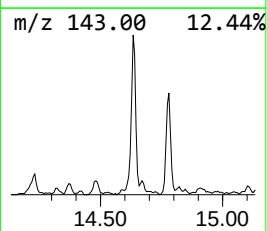
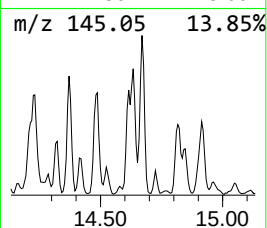
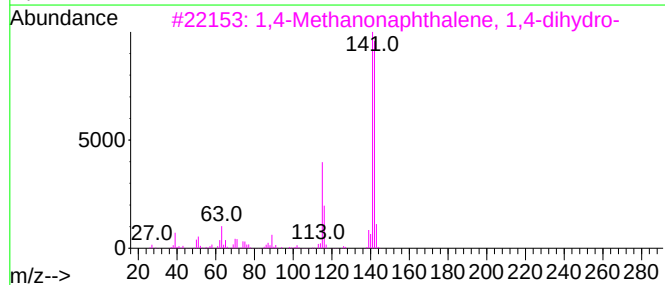
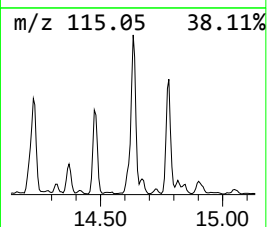
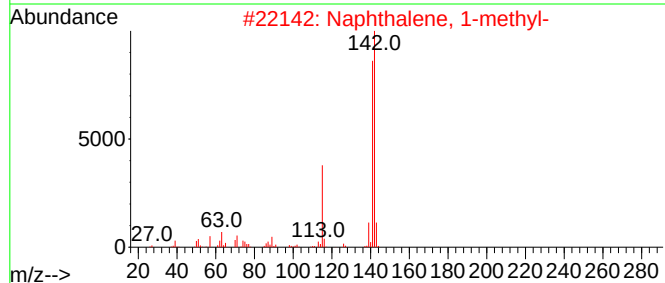
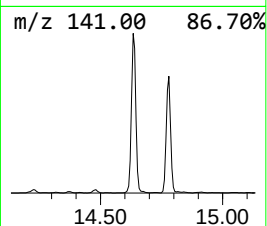
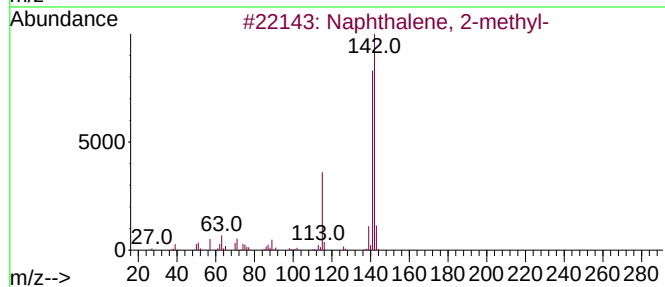
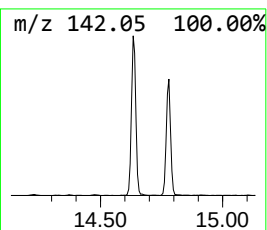
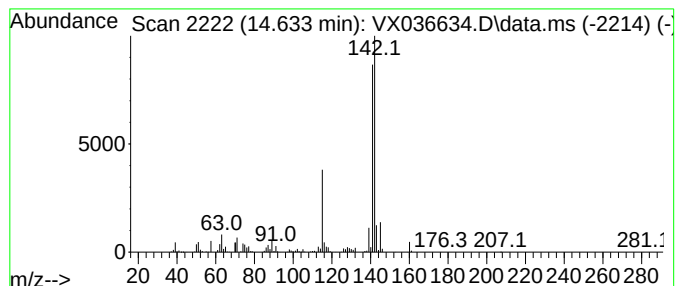
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Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.634	30.70 ug/l	484137	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	93
4			Benzocycloheptatriene	142	C11H10	000264-09-5	90
5			1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072423\
Data File : VX036634.D
Acq On : 24 Jul 2023 16:57
Operator : JC/MD
Sample : 03722-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
33712

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	24.7	ug/l	389848	4	12.024	788413	50.0
Mesitylene	12.085	27.1	ug/l	426973	4	12.024	788413	50.0
Benzene, 2-prop...	12.244	30.6	ug/l	482802	4	12.024	788413	50.0
Indan, 1-methyl-	12.701	17.5	ug/l	276434	4	12.024	788413	50.0
Benzene, 1-ethe...	13.164	17.9	ug/l	281778	4	12.024	788413	50.0
1H-Indene, 2,3-...	13.292	56.5	ug/l	890745	4	12.024	788413	50.0
Naphthalene, 1,...	13.420	51.6	ug/l	813402	4	12.024	788413	50.0
Naphthalene, 1,...	14.225	37.1	ug/l	585177	4	12.024	788413	50.0
Naphthalene, 1,...	14.475	25.4	ug/l	400735	4	12.024	788413	50.0
Naphthalene, 2-...	14.634	30.7	ug/l	484137	4	12.024	788413	50.0