

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX090923\
 Data File : VX037630.D
 Acq On : 09 Sep 2023 16:13
 Operator : JC/MD
 Sample : 04296-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 NGTW2

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\82X090823W.M
 Title : SW846 8260

Signal : TIC: VX037630.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.385	692	705	716	rBV	77863	228023	3.67%	0.375%
2	5.550	716	732	739	rVV	146404	497796	8.02%	0.819%
3	5.611	740	742	761	rVB2	47814	122851	1.98%	0.202%
4	5.824	765	777	789	rBV2	46097	137935	2.22%	0.227%
5	5.958	789	799	808	rBV	79111	212359	3.42%	0.349%
6	6.160	821	832	837	rBV3	26204	80145	1.29%	0.132%
7	6.245	837	846	857	rVV	149638	469996	7.57%	0.773%
8	6.763	921	931	940	rBV	212236	514877	8.29%	0.847%
9	7.367	1020	1030	1040	rVB4	48828	141985	2.29%	0.234%
10	7.562	1056	1062	1067	rBV	30084	63365	1.02%	0.104%
11	7.982	1122	1131	1143	rVB	135721	290967	4.69%	0.479%
12	8.104	1143	1151	1159	rBV	160922	335215	5.40%	0.551%
13	8.184	1159	1164	1170	rVB2	36454	64288	1.04%	0.106%
14	8.604	1226	1233	1236	rBV4	64977	136922	2.20%	0.225%
15	8.647	1236	1240	1250	rVB	451573	726383	11.70%	1.195%
16	9.104	1306	1315	1320	rBV	60922	113931	1.83%	0.187%
17	9.159	1320	1324	1329	rBV	56920	82186	1.32%	0.135%
18	9.506	1376	1381	1386	rBV4	38560	73480	1.18%	0.121%
19	10.055	1466	1471	1476	rVB	443074	578889	9.32%	0.952%
20	10.195	1490	1494	1499	rVB	147365	187882	3.03%	0.309%
21	10.964	1614	1620	1628	rBV	409294	512145	8.25%	0.842%
22	11.079	1635	1639	1644	rBV	368129	453435	7.30%	0.746%
23	11.305	1671	1676	1685	rVB	1313414	1638729	26.39%	2.695%
24	11.396	1685	1691	1696	rVV	862909	1089791	17.55%	1.792%
25	11.451	1696	1700	1710	rVB	1154326	1370666	22.07%	2.254%
26	11.610	1721	1726	1735	rBV	734223	921578	14.84%	1.516%
27	11.713	1735	1743	1745	rBV2	55610	81226	1.31%	0.134%
28	11.750	1745	1749	1755	rVB	5036118	6209741	100.00%	10.214%
29	11.866	1763	1768	1770	rBV	138730	185764	2.99%	0.306%
30	11.890	1770	1772	1777	rVV	191907	200122	3.22%	0.329%
31	11.969	1781	1785	1788	rVV	151438	173284	2.79%	0.285%
32	12.018	1788	1793	1798	rVV2	477185	677794	10.92%	1.115%
33	12.085	1799	1804	1812	rVB	1072242	1317024	21.21%	2.166%
34	12.177	1816	1819	1823	rVB	67151	77804	1.25%	0.128%
35	12.244	1824	1830	1837	rBV	2119552	3058429	49.25%	5.031%
36	12.317	1837	1842	1849	rVB2	1856653	2682680	43.20%	4.412%

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Integrator: RTE

Smoothing : ON

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Sampling : 1

Min Area: 3 % of largest Peak

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Stop Thrs : 0

Peak Location: TOP

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

Peak separation: 5

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37	12.402	1852	1856	1861	rVB	221861	274283	4.42%	0.451%
38	12.463	1861	1866	1872	rBV	618640	755800	12.17%	1.243%
39	12.530	1872	1877	1879	rBV	1730588	2067111	33.29%	3.400%
40	12.555	1879	1881	1886	rVB	1178021	1246894	20.08%	2.051%
41	12.616	1886	1891	1894	rBV	2881910	3629435	58.45%	5.970%
42	12.646	1894	1896	1900	rVB	536945	546338	8.80%	0.899%
43	12.701	1900	1905	1912	rBV2	1966231	2728207	43.93%	4.487%
44	12.847	1924	1929	1934	rVB	701339	860098	13.85%	1.415%
45	12.921	1934	1941	1944	rBV	2191679	2569479	41.38%	4.226%
46	12.963	1944	1948	1954	rVB	2581976	2984834	48.07%	4.909%
47	13.024	1954	1958	1960	rBV	176024	254125	4.09%	0.418%
48	13.042	1960	1961	1964	rVV2	182485	190456	3.07%	0.313%
49	13.067	1964	1965	1968	rVV	98002	95041	1.53%	0.156%
50	13.116	1968	1973	1974	rVV2	88502	133087	2.14%	0.219%
51	13.134	1974	1976	1978	rVV	140155	180393	2.91%	0.297%
52	13.170	1978	1982	1986	rVV	1913555	2388449	38.46%	3.929%
53	13.207	1986	1988	1991	rVV2	107936	146034	2.35%	0.240%
54	13.256	1991	1996	1997	rVV3	220938	309776	4.99%	0.510%
55	13.292	1997	2002	2008	rVV2	3550189	4843559	78.00%	7.967%
56	13.372	2012	2015	2020	rVV	173735	237733	3.83%	0.391%
57	13.420	2020	2023	2031	rVB	269383	340982	5.49%	0.561%
58	13.506	2031	2037	2039	rBV2	241757	323439	5.21%	0.532%
59	13.542	2039	2043	2046	rVV	586051	824123	13.27%	1.356%
60	13.579	2046	2049	2056	rVV3	594517	1155738	18.61%	1.901%
61	13.664	2056	2063	2069	rVB2	541213	1137471	18.32%	1.871%
62	13.774	2078	2081	2091	rVB2	99631	180764	2.91%	0.297%
63	13.853	2091	2094	2099	rVB4	45807	68867	1.11%	0.113%
64	13.926	2099	2106	2110	rBV2	224737	340154	5.48%	0.559%
65	13.969	2110	2113	2117	rVB	200611	224094	3.61%	0.369%
66	14.073	2126	2130	2136	rBV	439183	527351	8.49%	0.867%
67	14.213	2148	2153	2159	rBV	361043	569490	9.17%	0.937%
68	14.317	2166	2170	2175	rBV	171933	235458	3.79%	0.387%
69	14.371	2175	2179	2184	rVB	251031	314100	5.06%	0.517%
70	14.481	2192	2197	2201	rBV2	50520	78621	1.27%	0.129%
71	14.634	2218	2222	2227	rVB	541880	643824	10.37%	1.059%
72	14.780	2240	2246	2256	rVB	1129958	1441350	23.21%	2.371%
73	15.475	2354	2360	2372	rVB	68742	112432	1.81%	0.185%
74	15.609	2374	2382	2386	rVV	76746	128358	2.07%	0.211%

Sum of corrected areas: 60797435

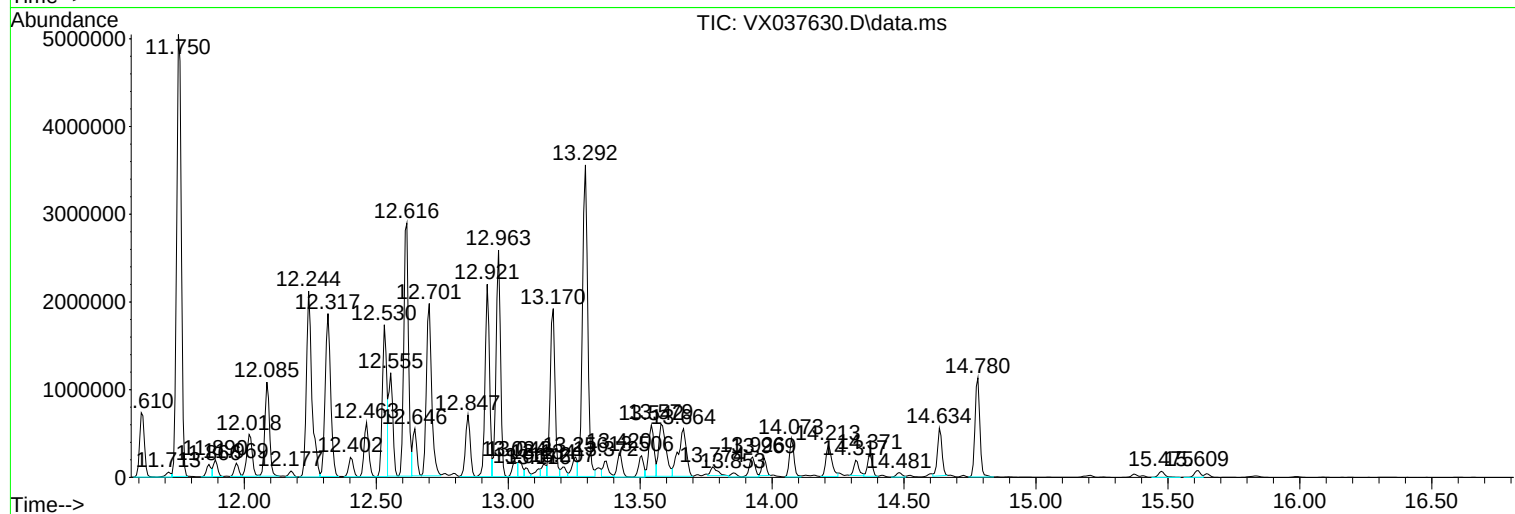
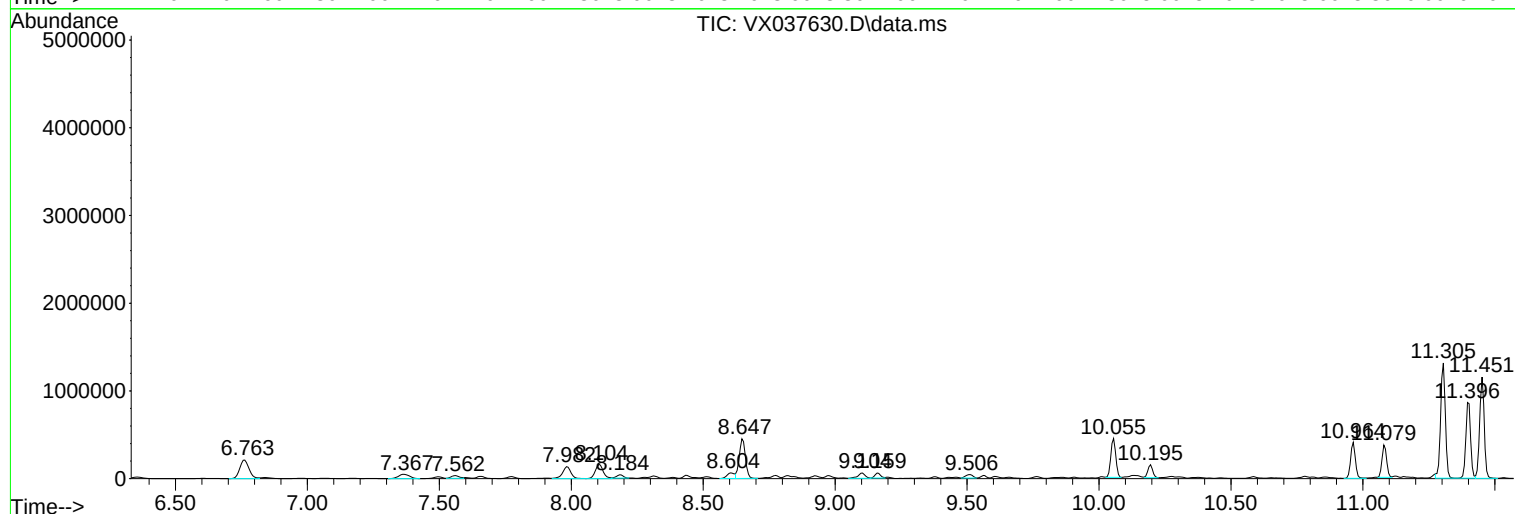
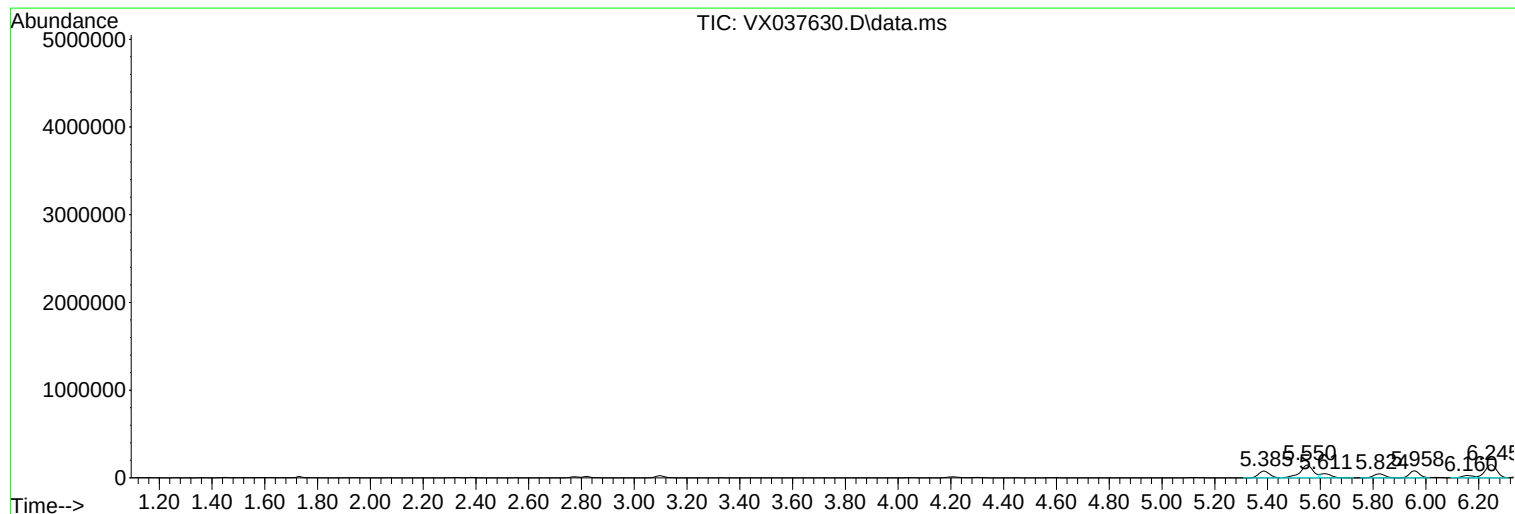
Instrument :
MSVOA_X
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Instrument :
MSVOA_X
ClientSampleId :
NGTW2

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X090823W.M
Quant Title : SW846 8260

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



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Operator : JC/MD
Sample : 04296-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
NGTW2

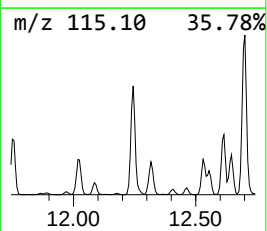
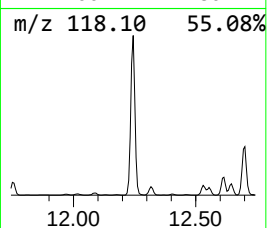
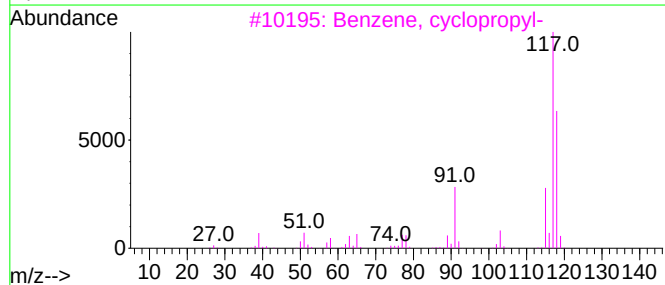
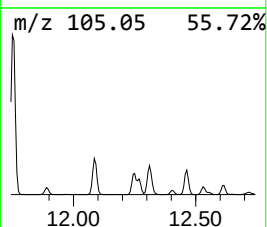
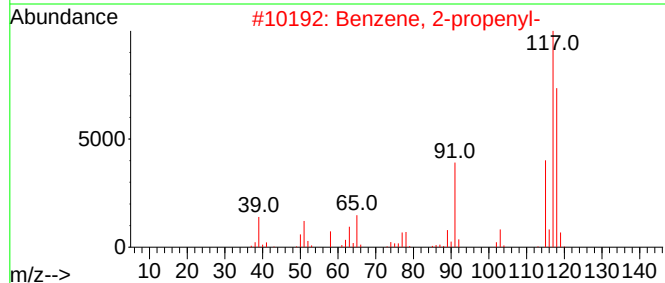
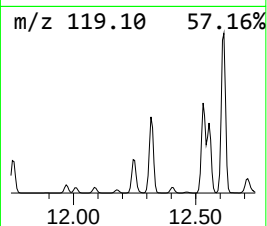
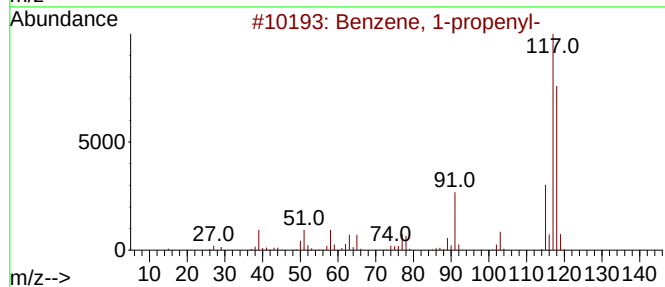
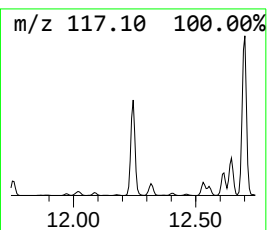
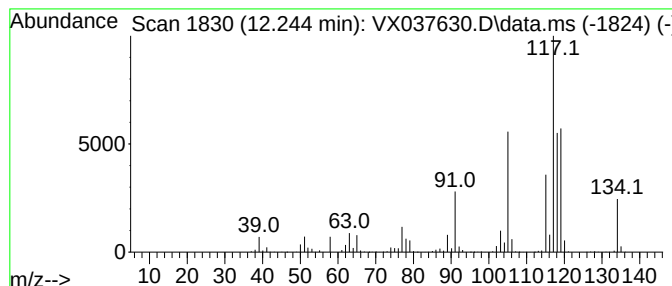
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X090823W.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1-propenyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	225.62 ug/l	3058430	1,4-Dichlorobenzene-d4	12.024

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 1-propenyl-	118	C9H10	000637-50-3	90
2		Benzene, 2-propenyl-	118	C9H10	000300-57-2	60
3		Benzene, cyclopropyl-	118	C9H10	000873-49-4	60
4		Indane	118	C9H10	000496-11-7	55
5		Deltacyclene	118	C9H10	007785-10-6	55



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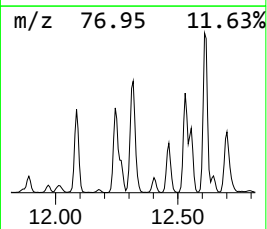
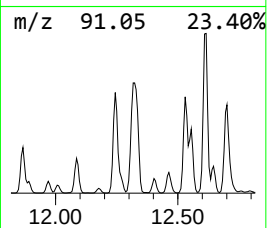
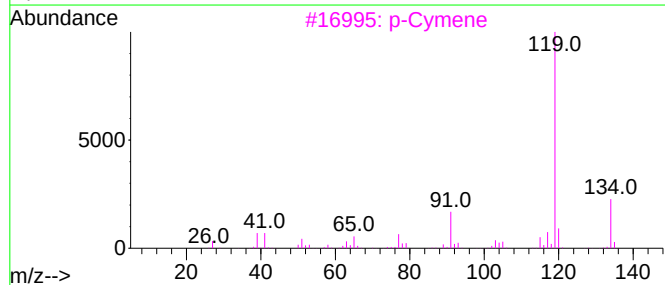
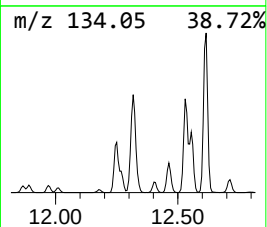
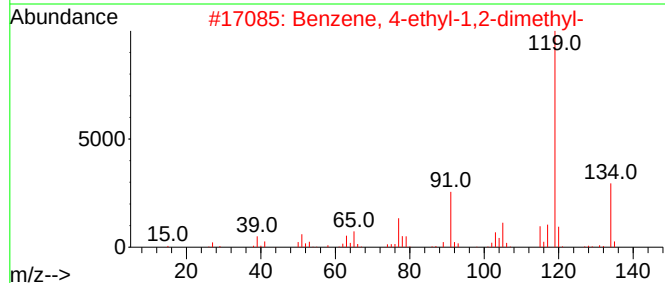
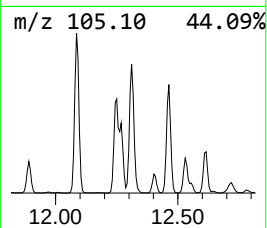
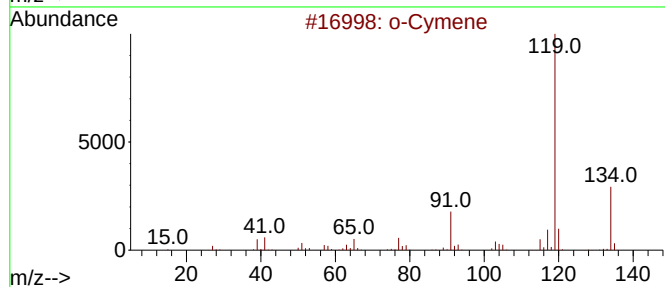
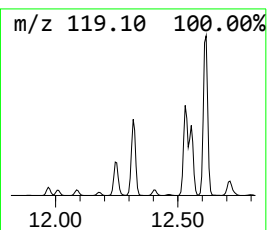
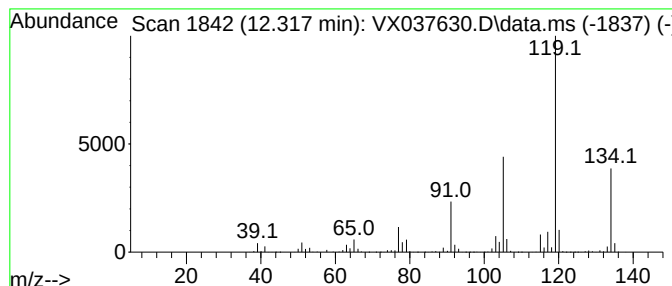
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X090823W.M
Quant Title : SW846 8260

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

Peak Number 2 o-Cymene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.317	197.90 ug/l	2682680	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Cymene	134	C10H14	000527-84-4	95
2			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94
3			p-Cymene	134	C10H14	000099-87-6	93
4			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91
5			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91



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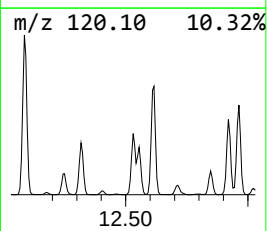
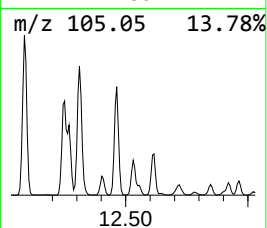
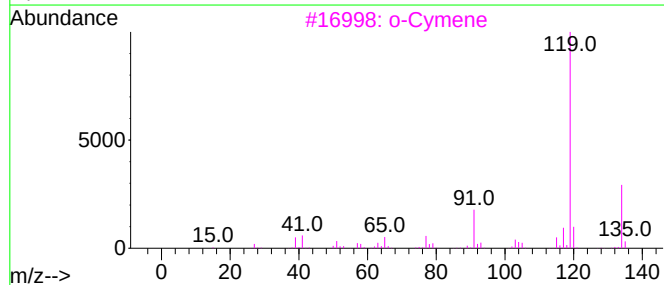
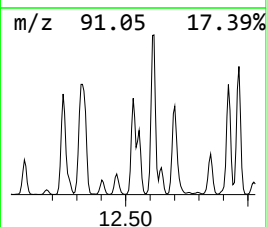
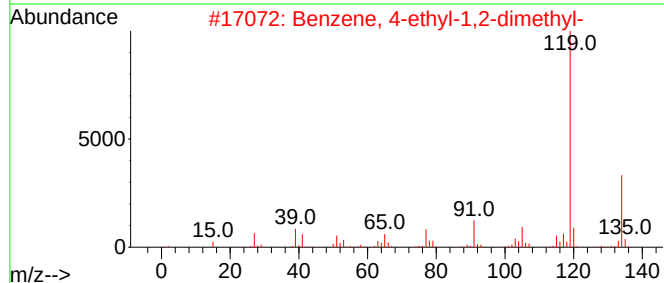
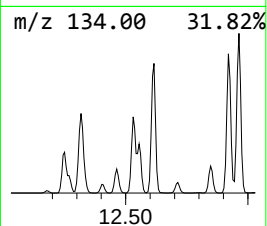
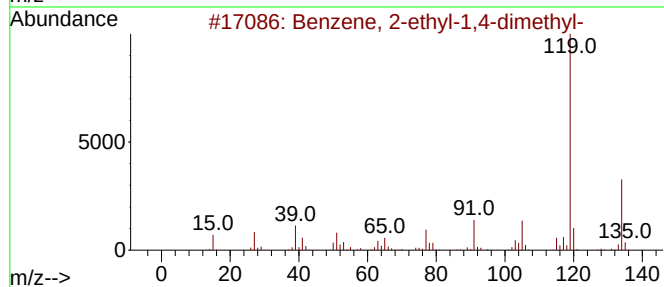
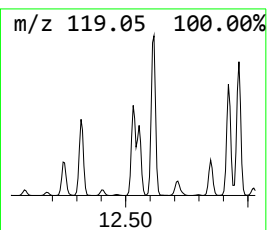
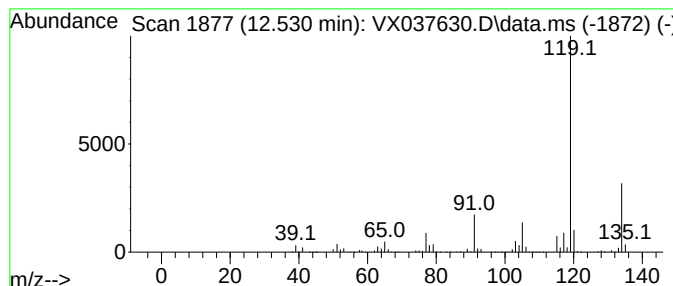
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X090823W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.530	152.49 ug/l	2067110	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	97
2			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
3			o-Cymene	134	C10H14	000527-84-4	95
4			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94



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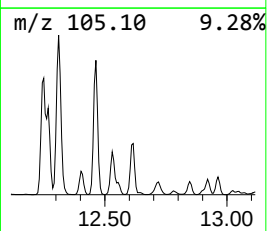
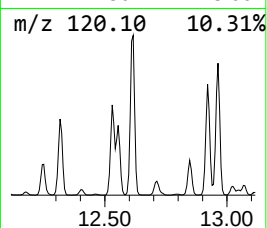
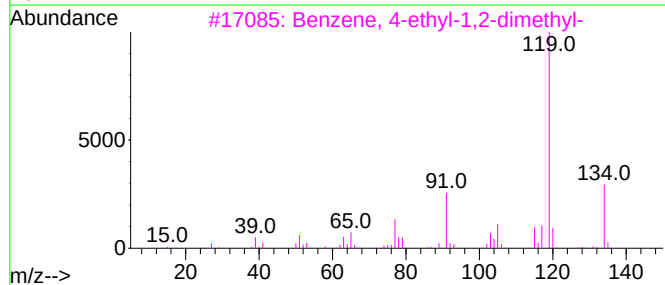
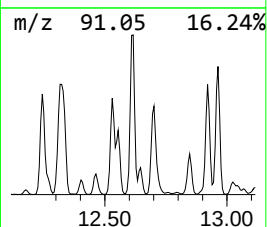
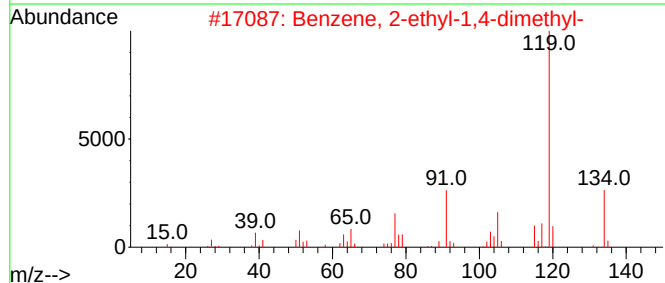
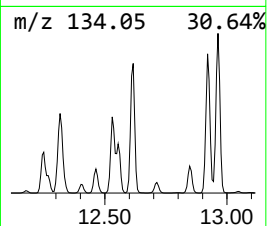
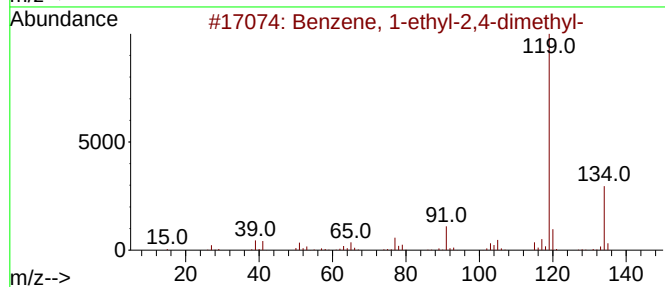
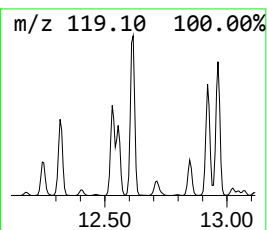
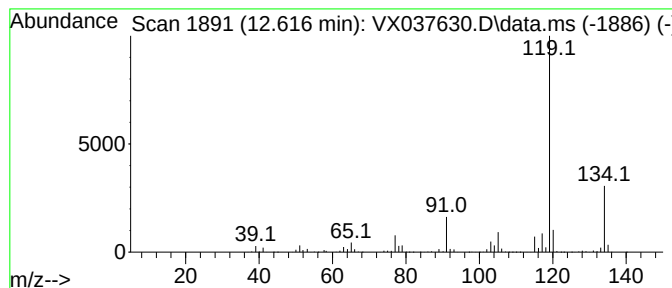
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.616	267.74 ug/l	3629440	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	97
2			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
3			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
4			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX090923\
Data File : VX037630.D
Acq On : 09 Sep 2023 16:13
Operator : JC/MD
Sample : 04296-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
NGTW2

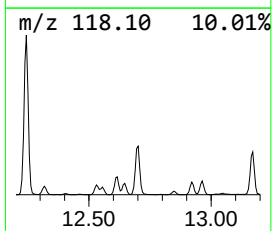
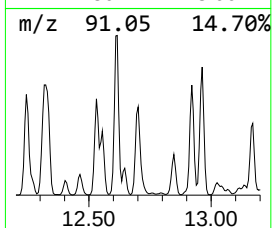
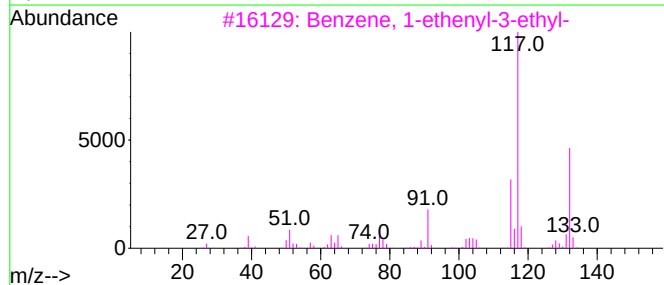
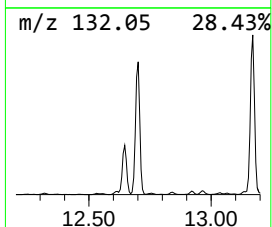
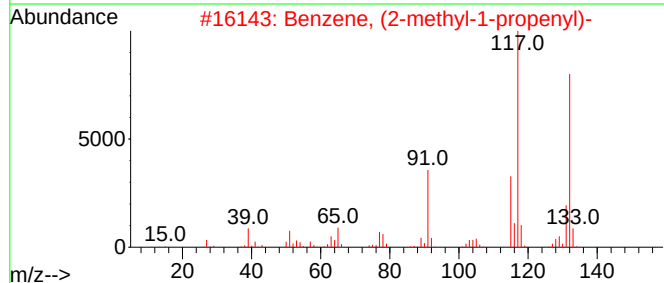
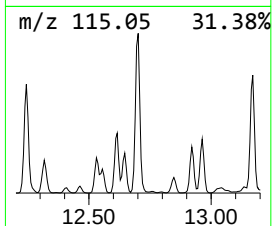
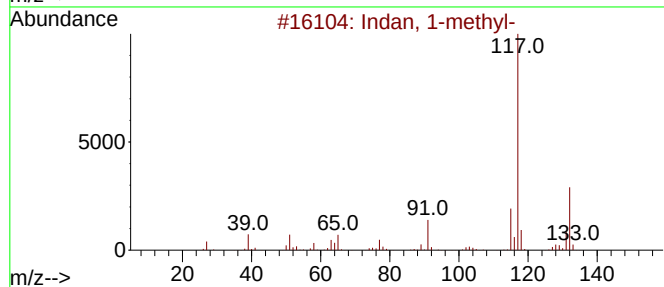
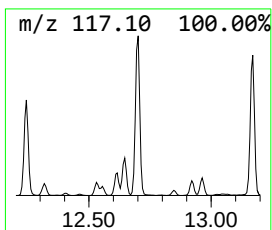
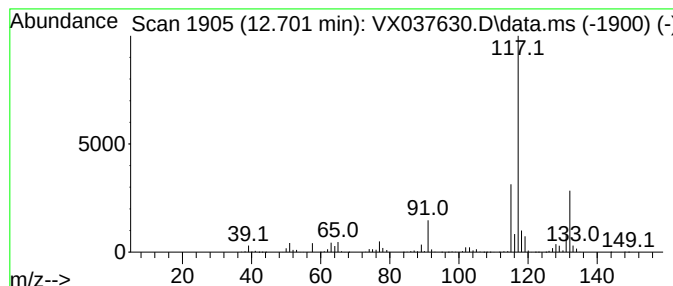
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Quant Title : SW846 8260

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

Peak Number 5 Indan, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.701	201.26 ug/l	2728210	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	90
3			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	90
4			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	87
5			Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX090923\
Data File : VX037630.D
Acq On : 09 Sep 2023 16:13
Operator : JC/MD
Sample : 04296-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

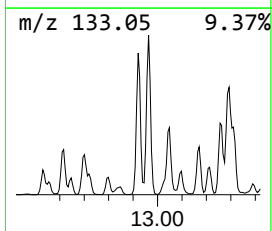
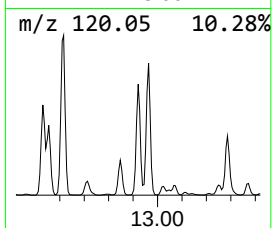
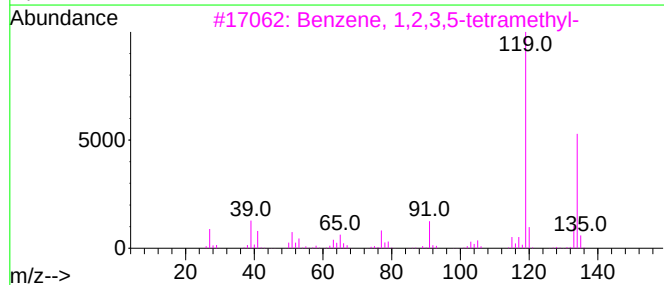
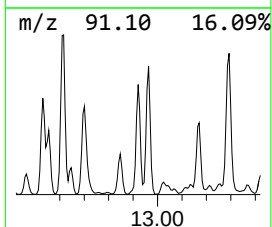
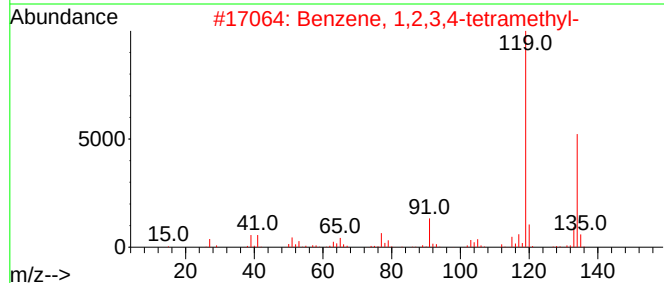
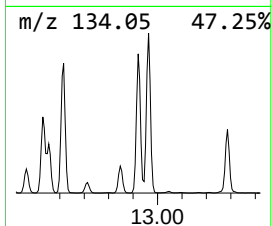
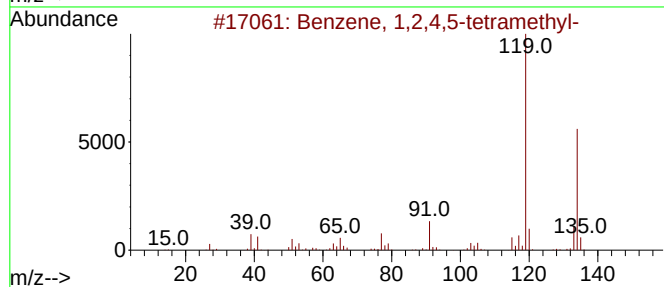
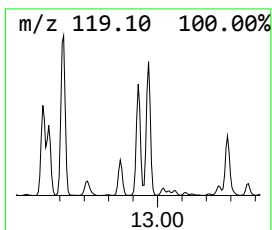
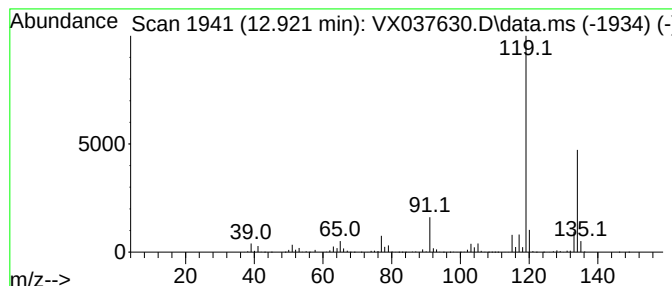
Instrument :
MSVOA_X
ClientSampleId :
NGTW2

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X090823W.M
Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD			R.T.
12.921	189.55 ug/l	2569480	1,4-Dichlorobenzene-d4			12.024
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-		134	C10H14	000095-93-2	97
2	Benzene, 1,2,3,4-tetramethyl-		134	C10H14	000488-23-3	97
3	Benzene, 1,2,3,5-tetramethyl-		134	C10H14	000527-53-7	97
4	Benzene, 1-ethyl-2,4-dimethyl-		134	C10H14	000874-41-9	94
5	Benzene, 1-ethyl-3,5-dimethyl-		134	C10H14	000934-74-7	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX090923\
Data File : VX037630.D
Acq On : 09 Sep 2023 16:13
Operator : JC/MD
Sample : 04296-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
NGTW2

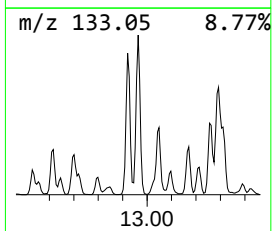
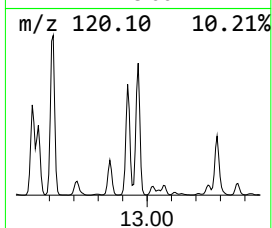
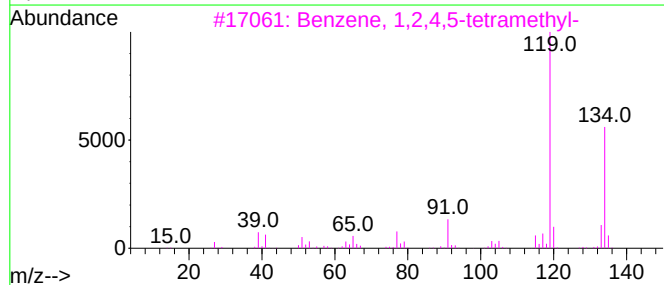
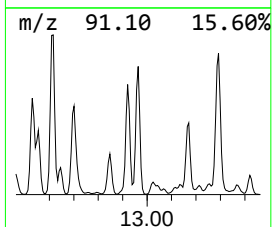
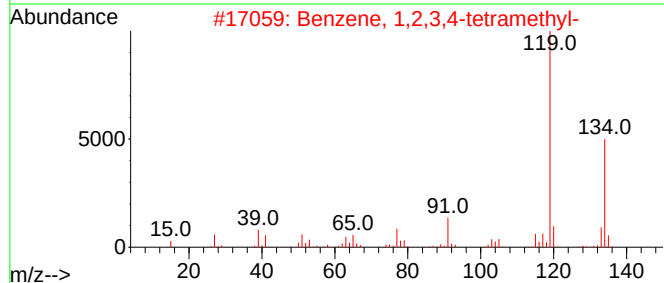
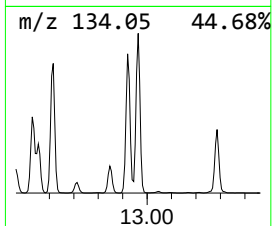
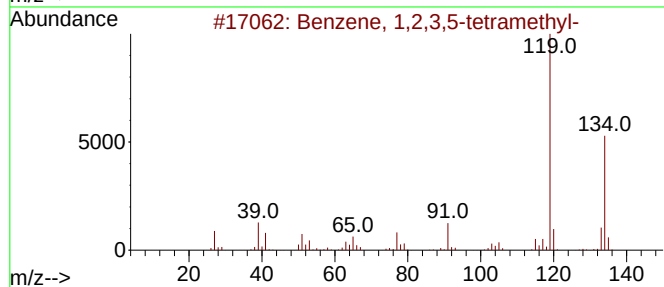
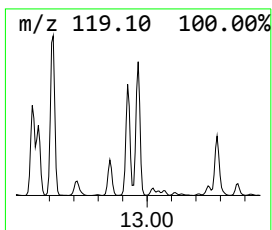
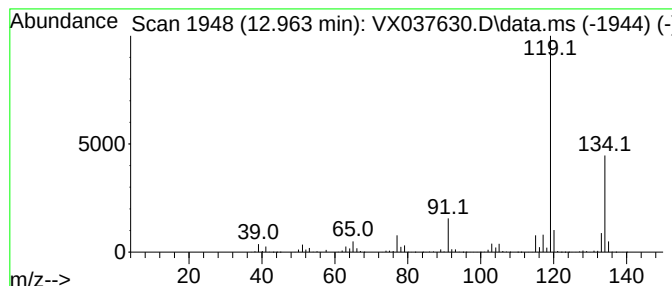
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.963	220.19 ug/l	2984830	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
2			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
4			o-Cymene	134	C10H14	000527-84-4	95
5			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX090923\
Data File : VX037630.D
Acq On : 09 Sep 2023 16:13
Operator : JC/MD
Sample : 04296-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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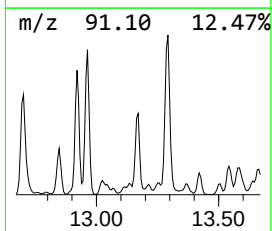
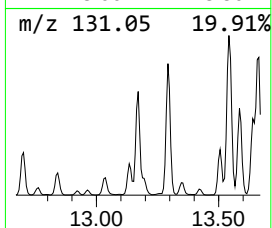
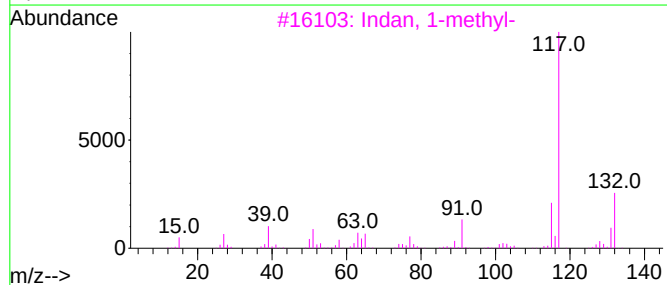
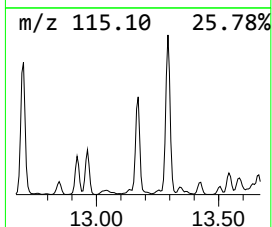
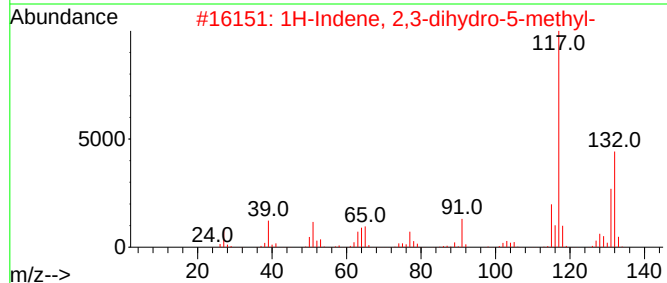
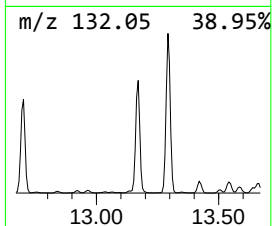
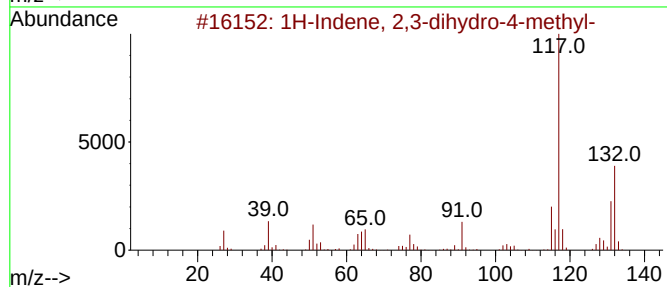
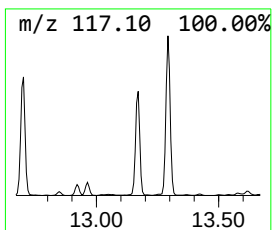
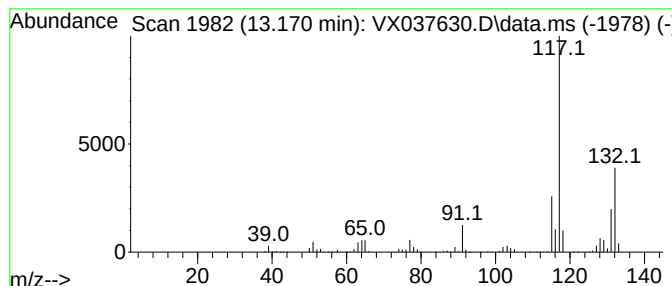
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Quant Title : SW846 8260

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

Peak Number 8 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.171	176.19 ug/l	2388450	1,4-Dichlorobenzene-d4	12.024

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	95	
2	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93	
3	Indan, 1-methyl-	132	C10H12	000767-58-8	90	
4	3-Phenylbut-1-ene	132	C10H12	000934-10-1	90	
5	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	87	



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX090923\
Data File : VX037630.D
Acq On : 09 Sep 2023 16:13
Operator : JC/MD
Sample : 04296-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
NGTW2

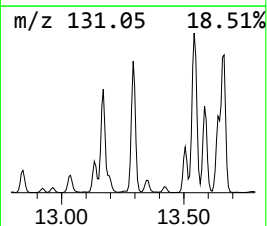
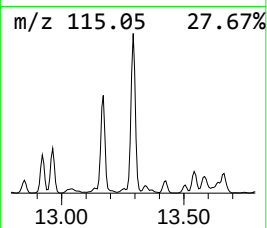
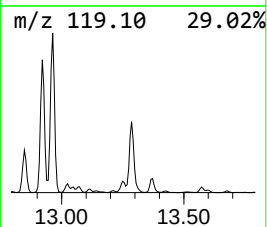
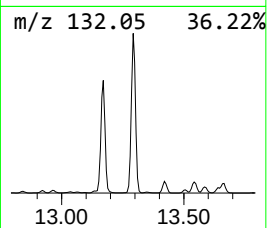
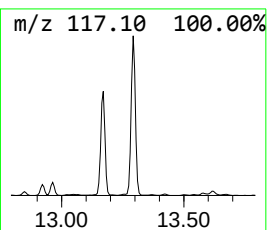
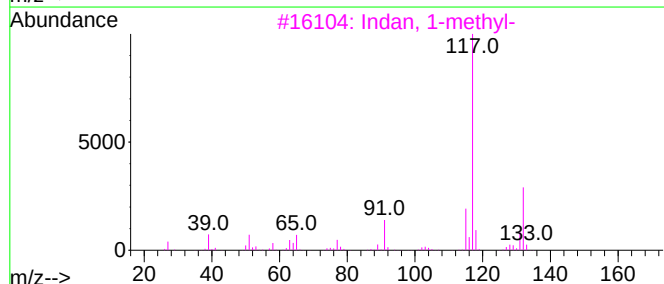
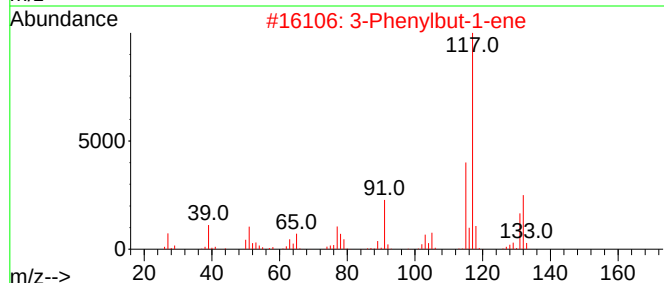
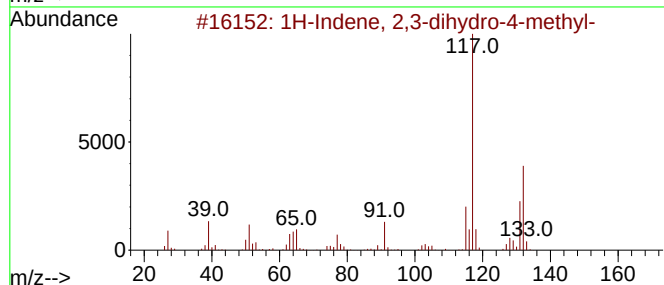
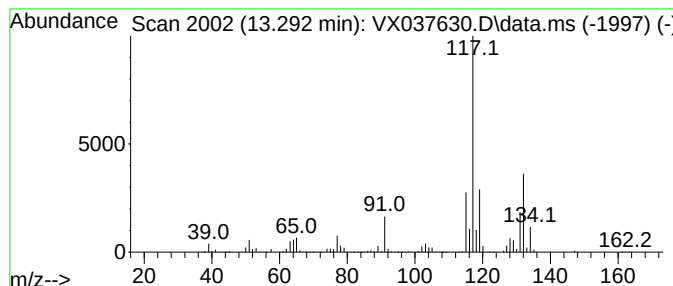
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Quant Title : SW846 8260

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

Peak Number 9 3-Phenylbut-1-ene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	357.30 ug/l	4843560	1,4-Dichlorobenzene-d4	12.024

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	81
2		3-Phenylbut-1-ene	132	C10H12	000934-10-1	81
3		Indan, 1-methyl-	132	C10H12	000767-58-8	76
4		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	76
5		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	74



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX090923\
Data File : VX037630.D
Acq On : 09 Sep 2023 16:13
Operator : JC/MD
Sample : 04296-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
NGTW2

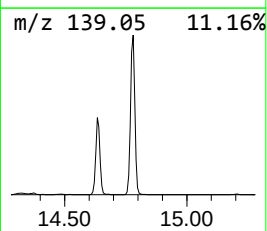
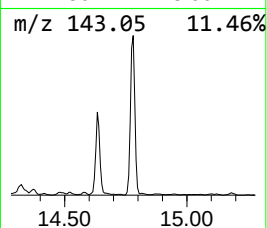
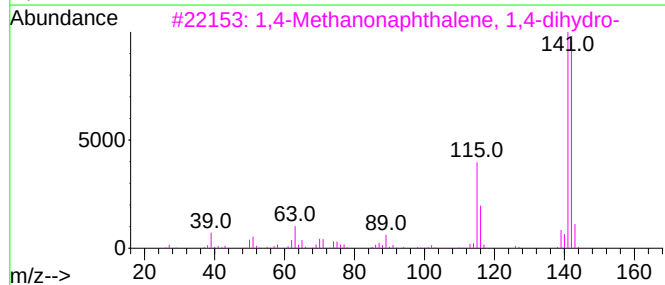
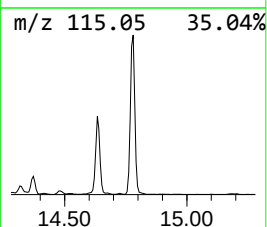
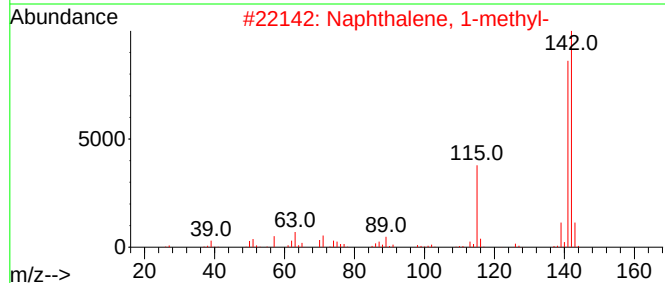
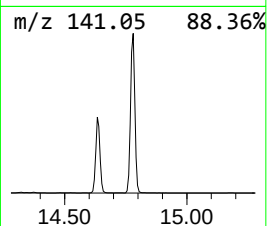
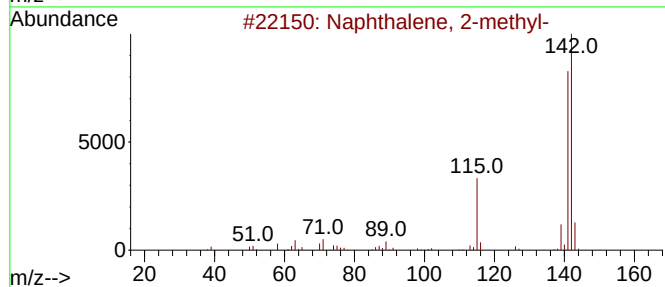
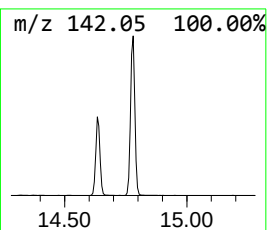
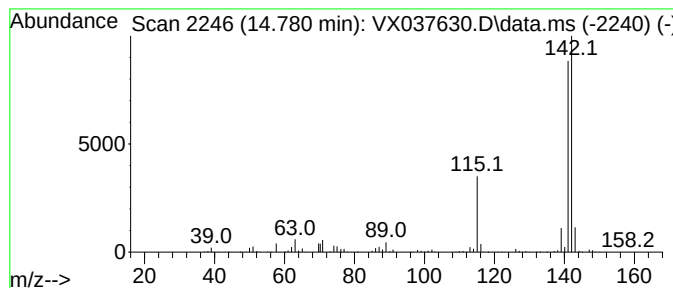
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.780	106.33 ug/l	1441350	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	91
4			Benzocycloheptatriene	142	C11H10	000264-09-5	91
5			1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX090923\
Data File : VX037630.D
Acq On : 09 Sep 2023 16:13
Operator : JC/MD
Sample : 04296-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
NGTW2

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X090823W.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-prop...	12.244	225.6	ug/l	3058430	4	12.024	677794	50.0
o-Cymene	12.317	197.9	ug/l	2682680	4	12.024	677794	50.0
Benzene, 2-ethy...	12.530	152.5	ug/l	2067110	4	12.024	677794	50.0
Benzene, 1-ethy...	12.616	267.7	ug/l	3629440	4	12.024	677794	50.0
Indan, 1-methyl-	12.701	201.3	ug/l	2728210	4	12.024	677794	50.0
Benzene, 1,2,4,...	12.921	189.6	ug/l	2569480	4	12.024	677794	50.0
Benzene, 1,2,3,...	12.963	220.2	ug/l	2984830	4	12.024	677794	50.0
1H-Indene, 2,3-...	13.171	176.2	ug/l	2388450	4	12.024	677794	50.0
3-Phenylbut-1-ene	13.292	357.3	ug/l	4843560	4	12.024	677794	50.0
Naphthalene, 2-...	14.780	106.3	ug/l	1441350	4	12.024	677794	50.0