

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041422\
 Data File : VX028104.D
 Acq On : 14 Apr 2022 14:04
 Operator : JC/MD
 Sample : N2403-03
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-31R

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\82X041222W.M
 Title : SW846 8260

Signal : TIC: VX028104.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.752	103	109	118	rBV	269904	377310	2.27%	0.285%
2	2.782	270	278	281	rBV2	173771	356698	2.14%	0.269%
3	2.825	281	285	303	rVB	244481	547098	3.29%	0.413%
4	3.105	322	331	345	rVB2	347286	823736	4.95%	0.622%
5	4.306	519	528	540	rVB2	181227	518730	3.12%	0.392%
6	5.391	690	706	714	rBV2	159391	483692	2.90%	0.365%
7	5.495	714	723	726	rVV5	87471	305588	1.84%	0.231%
8	5.556	727	733	739	rVV	238754	700621	4.21%	0.529%
9	5.623	740	744	763	rVB	144701	429442	2.58%	0.324%
10	5.830	763	778	790	rBV4	142670	456927	2.74%	0.345%
11	5.958	790	799	807	rBV2	165790	413137	2.48%	0.312%
12	6.165	823	833	841	rBV2	119487	391697	2.35%	0.296%
13	6.257	841	848	857	rVV3	162661	479361	2.88%	0.362%
14	6.361	857	865	876	rVB2	96385	277280	1.67%	0.209%
15	6.763	922	931	939	rBV	383349	959463	5.76%	0.725%
16	6.848	939	945	958	rVB2	100197	300733	1.81%	0.227%
17	7.379	1020	1032	1043	rBV4	283364	785662	4.72%	0.593%
18	7.988	1122	1132	1143	rVB5	62134	235548	1.41%	0.178%
19	8.110	1143	1152	1155	rBV2	113387	241758	1.45%	0.183%
20	8.506	1213	1217	1227	rVB5	84863	181575	1.09%	0.137%
21	8.604	1227	1233	1236	rBV2	146562	279154	1.68%	0.211%
22	8.653	1236	1241	1250	rVB	858681	1428101	8.58%	1.078%
23	8.775	1250	1261	1265	rBV4	83219	180920	1.09%	0.137%
24	8.976	1290	1294	1300	rVB	110766	170931	1.03%	0.129%
25	9.104	1306	1315	1320	rBV	150214	275578	1.65%	0.208%
26	9.165	1320	1325	1335	rVB	212644	353745	2.12%	0.267%
27	9.506	1377	1381	1387	rVB2	94633	184959	1.11%	0.140%
28	10.055	1466	1471	1476	rVB	883649	1139222	6.84%	0.860%
29	10.195	1490	1494	1503	rVB	539491	705239	4.24%	0.533%
30	10.305	1503	1512	1518	rVB	647963	898165	5.39%	0.678%
31	10.963	1615	1620	1628	rBV	749286	945167	5.68%	0.714%
32	11.085	1635	1640	1644	rBV	724875	917815	5.51%	0.693%
33	11.305	1666	1676	1683	rBV	2082826	2509083	15.07%	1.895%
34	11.402	1683	1692	1696	rVV	2005407	3463183	20.80%	2.615%
35	11.451	1696	1700	1707	rVB	765780	949553	5.70%	0.717%
36	11.616	1721	1727	1736	rBV	2133490	2570658	15.44%	1.941%

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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\82X041222W.M
 Title : SW846 8260

37	11.756	1745	1750	1755	rVB	15209856	16652587	100.00%	12.576%
38	11.866	1762	1768	1770	rBV	437278	539208	3.24%	0.407%
39	11.890	1770	1772	1778	rVB	480143	567934	3.41%	0.429%
40	11.975	1781	1786	1789	rBV	201994	269207	1.62%	0.203%
41	12.024	1789	1794	1799	rVB2	1201425	1671297	10.04%	1.262%
42	12.091	1799	1805	1810	rBV	2311362	2829735	16.99%	2.137%
43	12.250	1825	1831	1837	rBV	4787370	6206011	37.27%	4.687%
44	12.317	1837	1842	1849	rVB3	2585957	3794761	22.79%	2.866%
45	12.408	1852	1857	1861	rVB	537801	634650	3.81%	0.479%
46	12.469	1861	1867	1873	rVB2	1537534	2124557	12.76%	1.604%
47	12.536	1873	1878	1880	rBV	3271788	4421781	26.55%	3.339%
48	12.555	1880	1881	1886	rVB	2414475	2301615	13.82%	1.738%
49	12.616	1886	1891	1894	rBV	6267811	7062402	42.41%	5.333%
50	12.646	1894	1896	1900	rVB	1390750	1526490	9.17%	1.153%
51	12.701	1900	1905	1913	rBV	4850338	6466572	38.83%	4.883%
52	12.847	1924	1929	1934	rVB2	904173	1208893	7.26%	0.913%
53	12.926	1934	1942	1945	rBV	3644637	4468658	26.83%	3.375%
54	12.963	1945	1948	1954	rVB	4890600	5854693	35.16%	4.421%
55	13.030	1954	1959	1967	rBV4	509229	934029	5.61%	0.705%
56	13.140	1971	1977	1978	rBV2	383281	412455	2.48%	0.311%
57	13.170	1978	1982	1987	rVB	3698211	4160509	24.98%	3.142%
58	13.256	1992	1996	1998	rVV2	587264	816735	4.90%	0.617%
59	13.292	1998	2002	2008	rVV2	5422441	7581283	45.53%	5.725%
60	13.347	2008	2011	2013	rVV3	386799	571605	3.43%	0.432%
61	13.371	2013	2015	2020	rVV	384460	488344	2.93%	0.369%
62	13.426	2020	2024	2030	rVB	613513	784765	4.71%	0.593%
63	13.506	2031	2037	2040	rBV2	752352	1015245	6.10%	0.767%
64	13.548	2040	2044	2047	rVV	1738187	2482925	14.91%	1.875%
65	13.585	2047	2050	2056	rVV3	1721271	3276055	19.67%	2.474%
66	13.646	2056	2060	2061	rVV	930952	1329168	7.98%	1.004%
67	13.664	2061	2063	2070	rVB2	1843129	2457583	14.76%	1.856%
68	13.780	2074	2082	2091	rVB2	521440	905176	5.44%	0.684%
69	13.859	2091	2095	2100	rVB2	160463	242490	1.46%	0.183%
70	13.926	2100	2106	2110	rBV2	671855	913594	5.49%	0.690%
71	13.975	2110	2114	2118	rVB	471596	563425	3.38%	0.425%
72	14.079	2126	2131	2137	rBV	1362683	1641258	9.86%	1.239%
73	14.219	2149	2154	2159	rBV	973004	1457248	8.75%	1.101%
74	14.323	2167	2171	2176	rBV	452537	601785	3.61%	0.454%
75	14.371	2176	2179	2184	rVB	635138	764219	4.59%	0.577%
76	14.481	2192	2197	2202	rBV2	158961	245889	1.48%	0.186%

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ClientSampleId :
MW-31R

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\82X041222W.M
Title : SW846 8260

77	14.603	2210	2217	2220	rBV	160554	261970	1.57%	0.198%
78	14.640	2220	2223	2227	rVB	250123	314712	1.89%	0.238%
79	14.780	2241	2246	2262	rVB	2202141	2837209	17.04%	2.143%
80	15.481	2355	2361	2365	rBV	116155	189480	1.14%	0.143%
81	15.615	2375	2383	2387	rBV	206050	329048	1.98%	0.248%

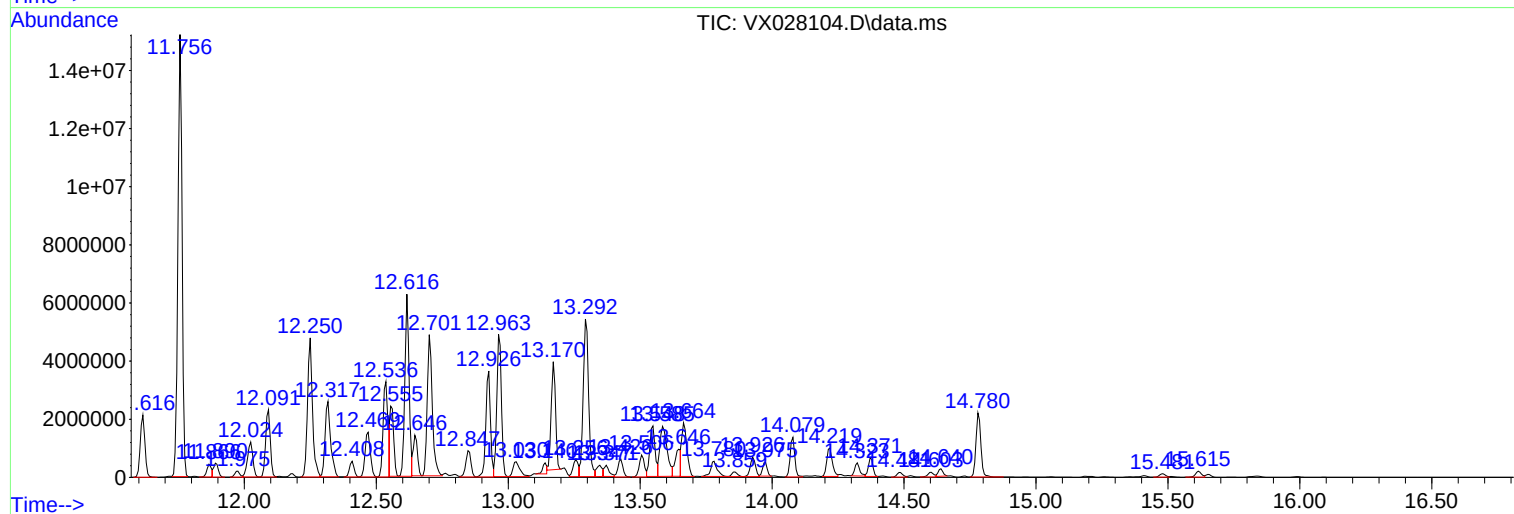
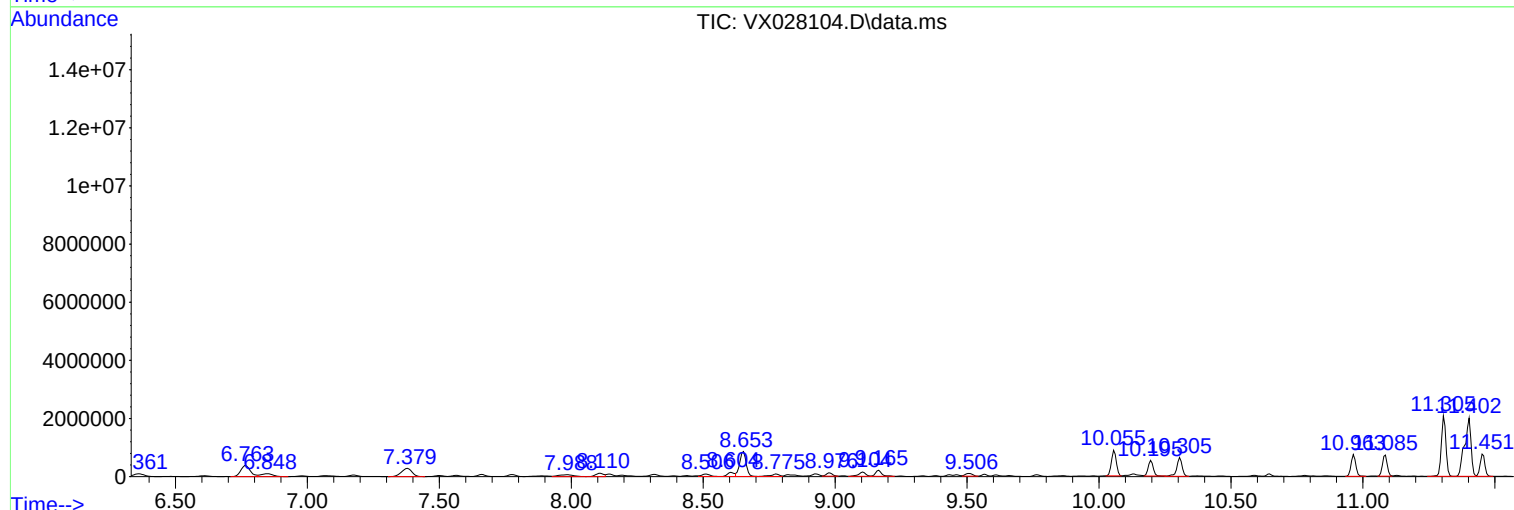
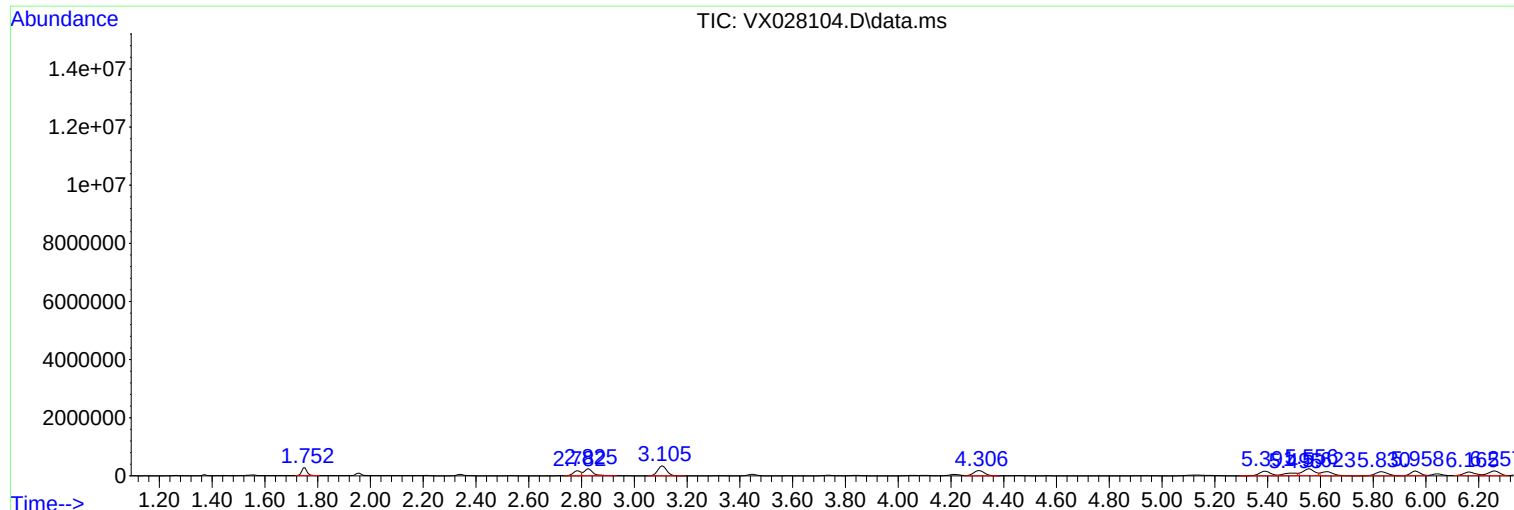
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Instrument :
MSVOA_X
ClientSampleId :
MW-31R

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041222W.M
Quant Title : SW846 8260

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



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Operator : JC/MD
Sample : N2403-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-31R

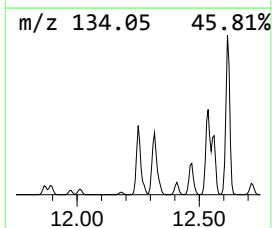
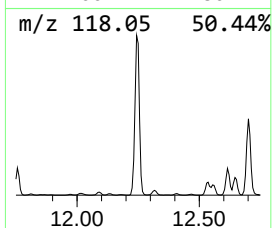
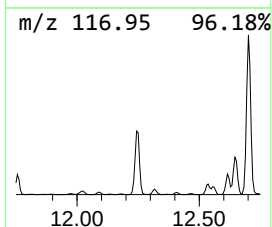
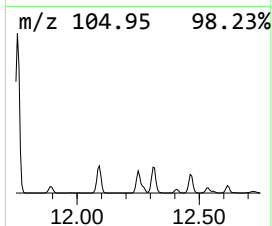
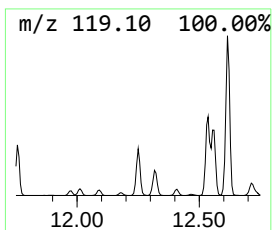
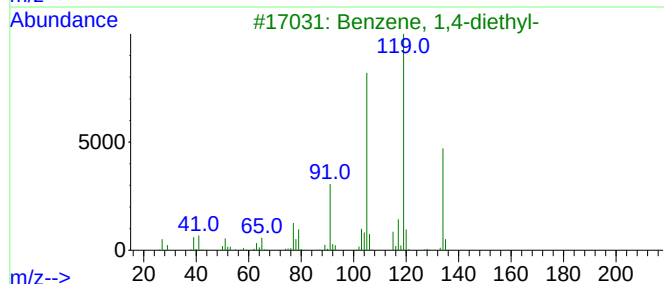
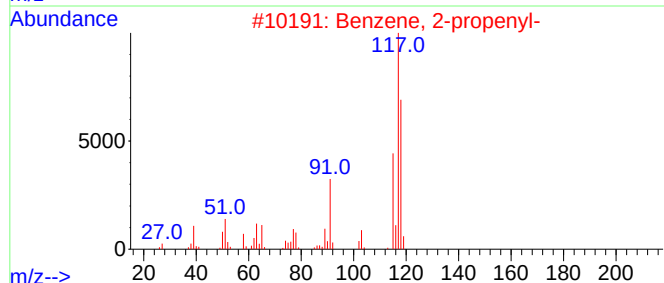
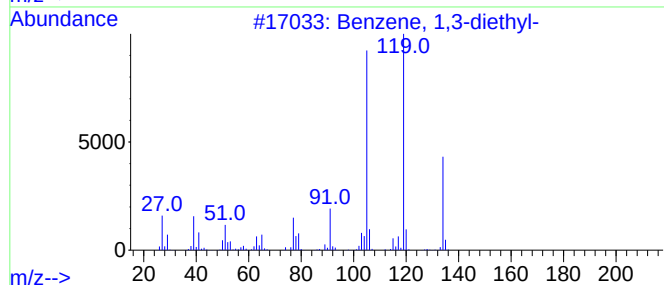
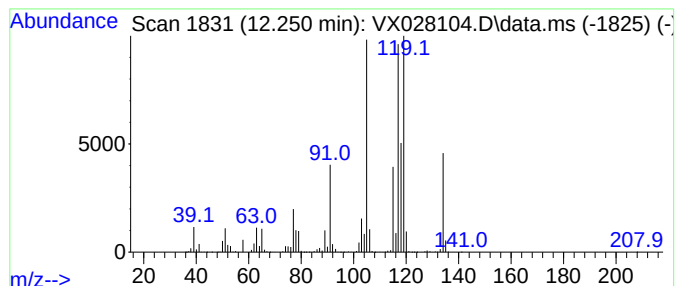
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041222W.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1,3-diethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.250	185.66 ug/l	6206010	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	86
2			Benzene, 2-propenyl-	118	C9H10	000300-57-2	80
3			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	55
4			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	55
5			Benzene, cyclopropyl-	118	C9H10	000873-49-4	42



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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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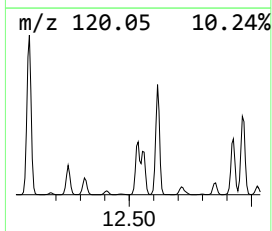
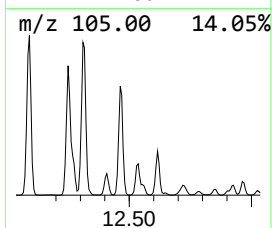
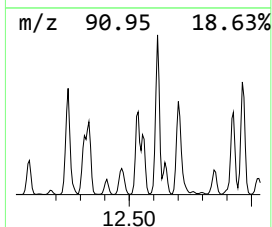
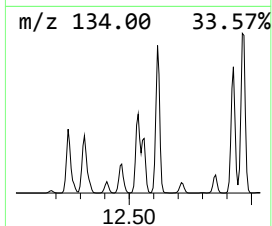
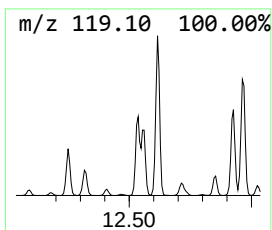
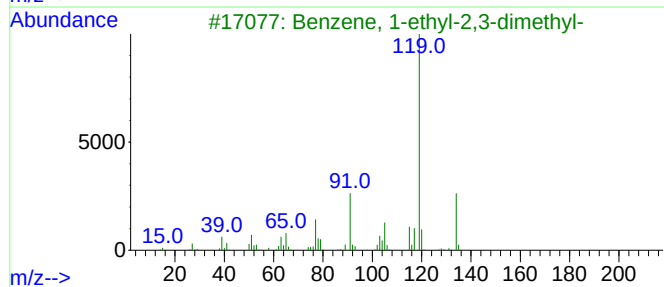
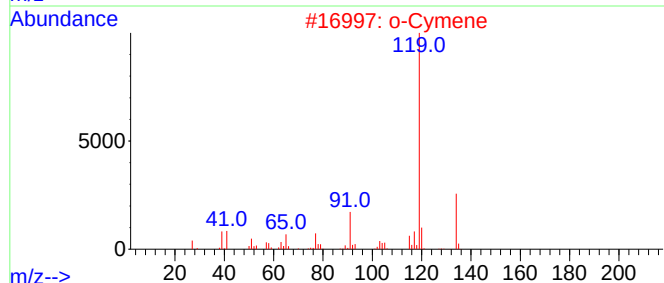
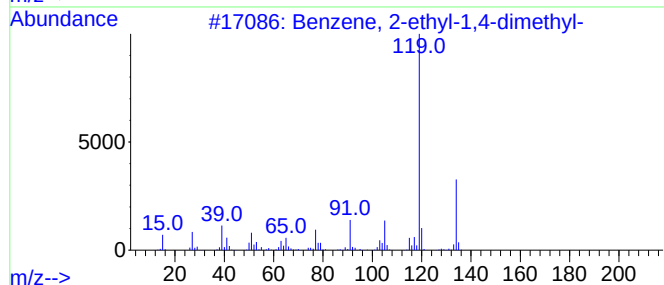
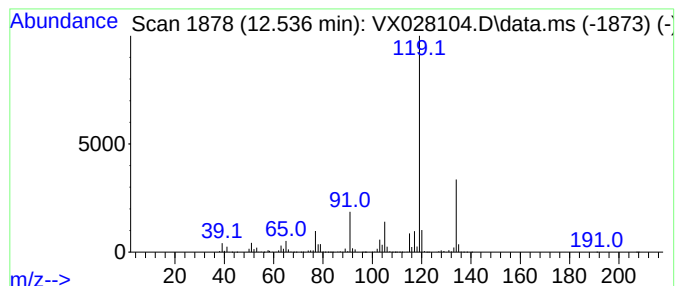
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041222W.M
Quant Title : SW846 8260

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

Peak Number 2 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.536	132.29 ug/l	4421780	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			o-Cymene	134	C10H14	000527-84-4	95
3			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
4			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94



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Instrument :
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ClientSampleId :
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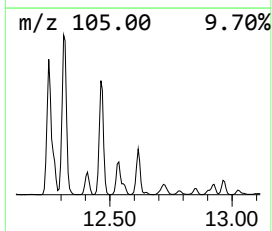
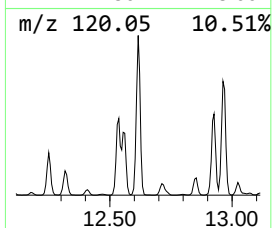
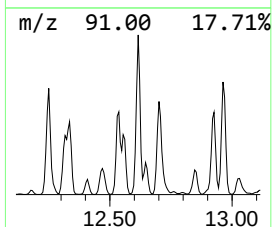
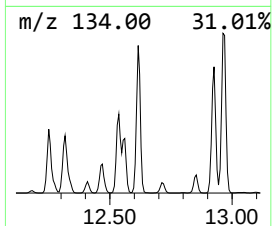
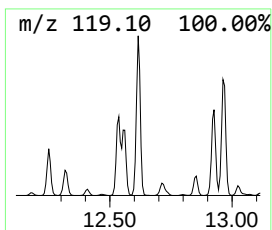
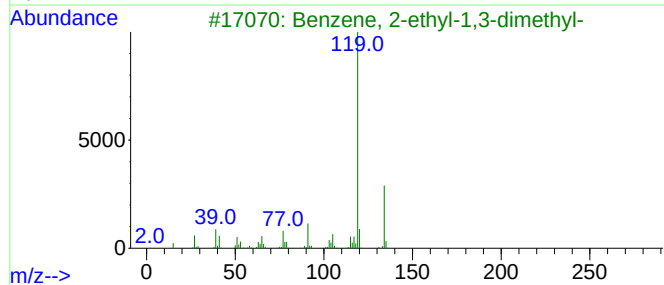
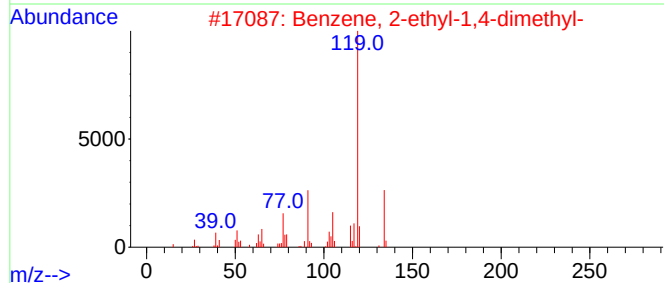
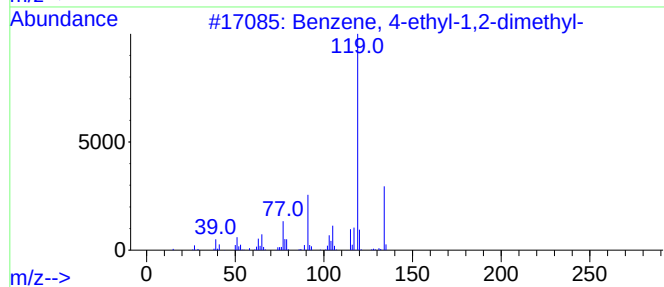
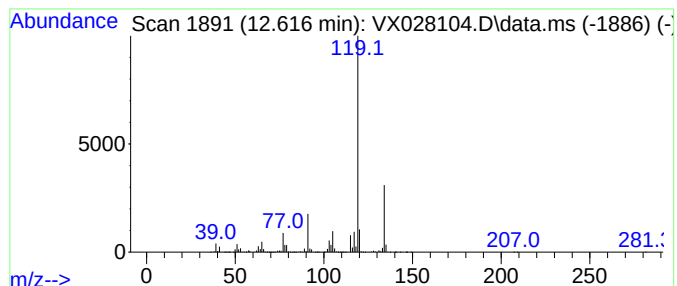
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041222W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.616	211.29 ug/l	7062400	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
3			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
4			o-Cymene	134	C10H14	000527-84-4	95
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95



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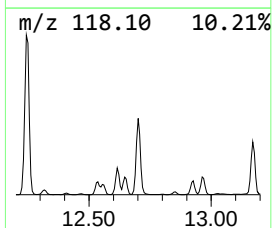
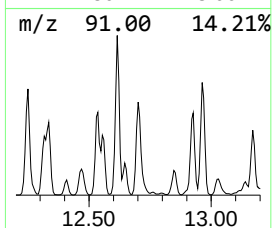
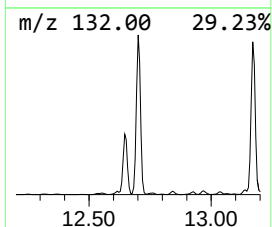
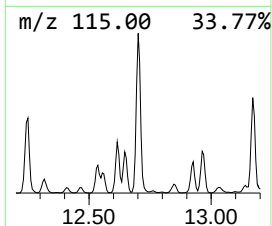
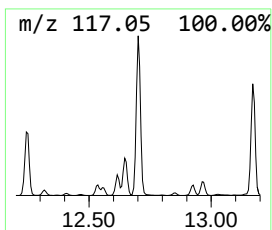
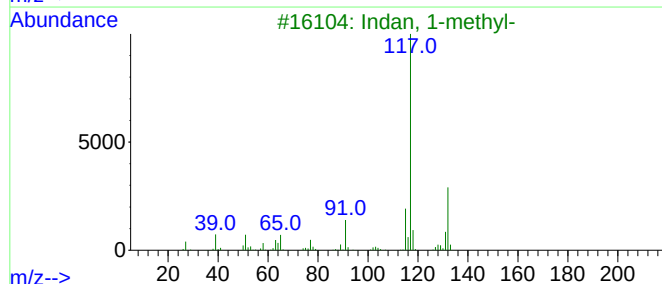
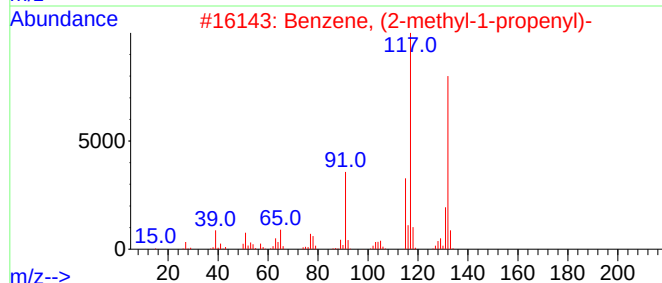
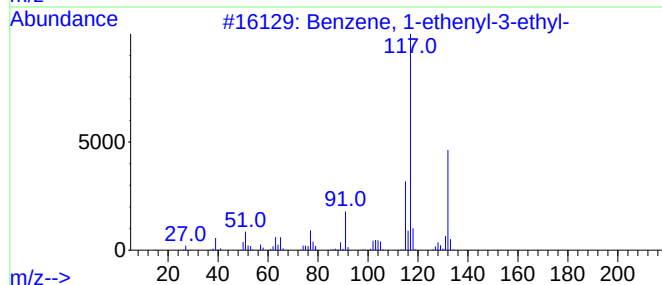
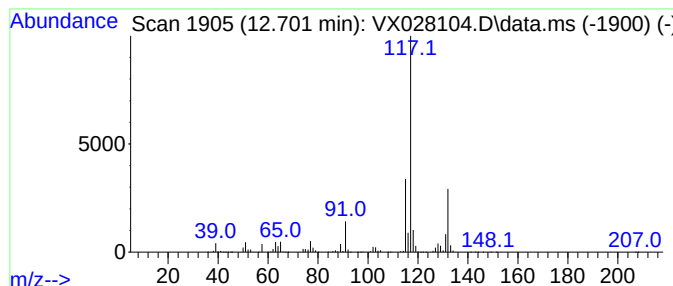
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.701	193.46 ug/l	6466570	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	91
2			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	90
3			Indan, 1-methyl-	132	C10H12	000767-58-8	90
4			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	87
5			1-Phenyl-1-butene	132	C10H12	000824-90-8	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041422\
Data File : VX028104.D
Acq On : 14 Apr 2022 14:04
Operator : JC/MD
Sample : N2403-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-31R

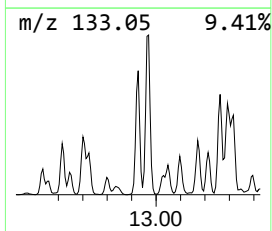
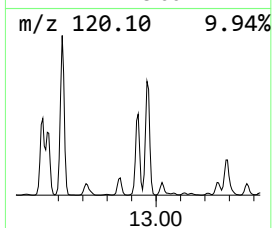
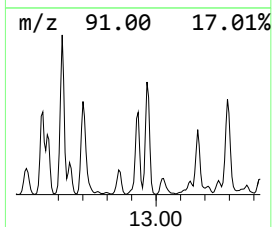
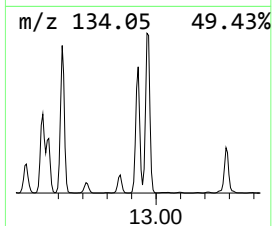
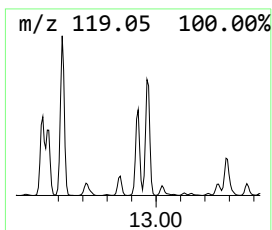
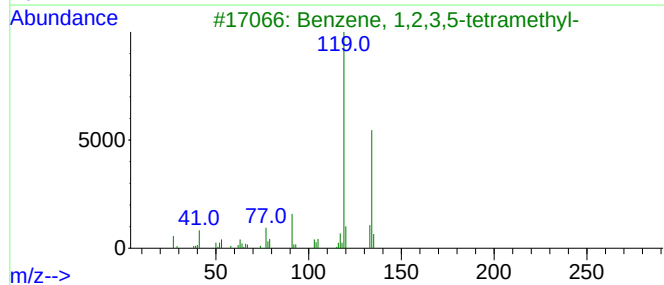
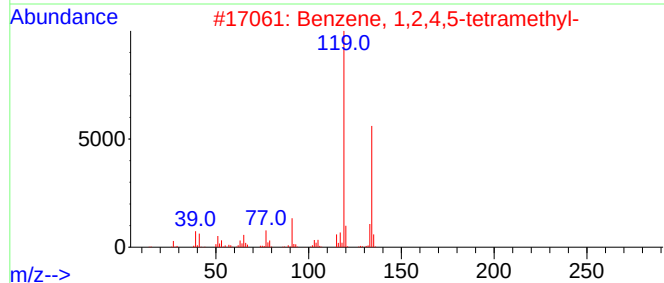
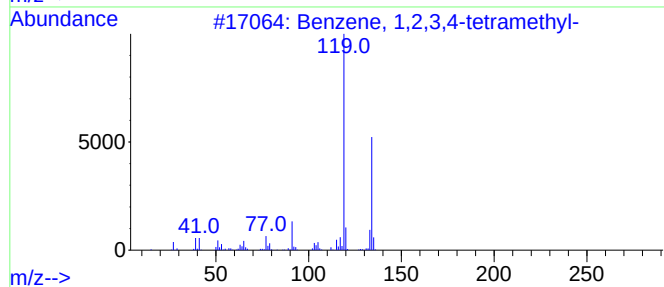
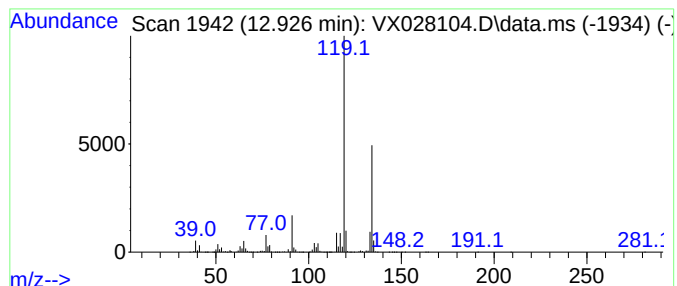
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.926	133.69 ug/l	4468660	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
3			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
4			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	94
5			1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041422\
Data File : VX028104.D
Acq On : 14 Apr 2022 14:04
Operator : JC/MD
Sample : N2403-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-31R

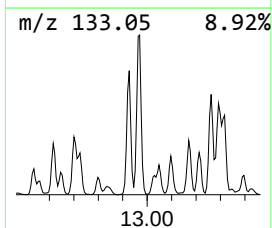
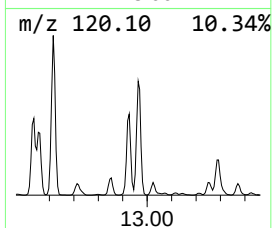
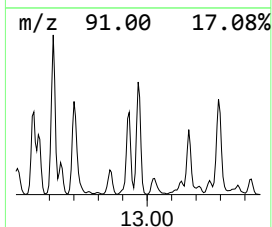
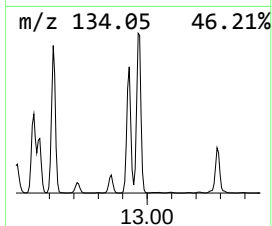
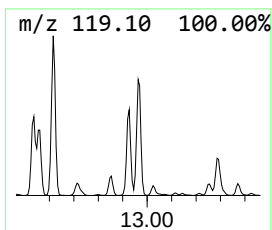
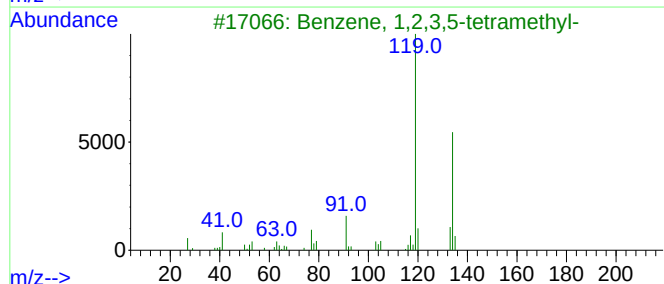
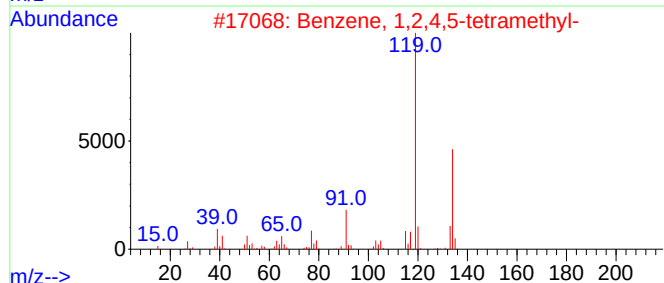
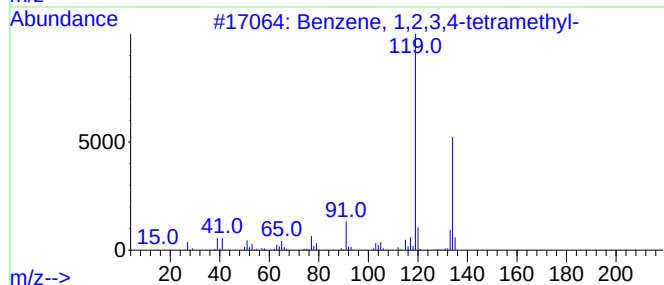
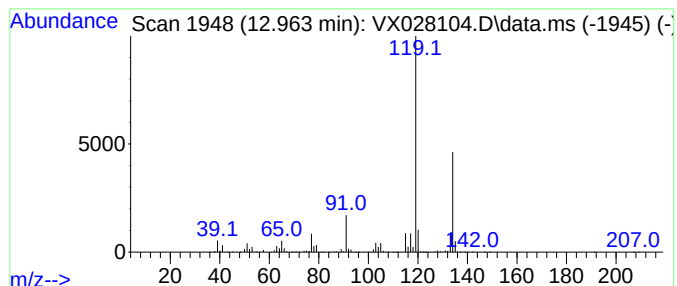
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.963	175.15 ug/l	5854690	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
3			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
4			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	95
5			o-Cymene	134	C10H14	000527-84-4	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041422\
Data File : VX028104.D
Acq On : 14 Apr 2022 14:04
Operator : JC/MD
Sample : N2403-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-31R

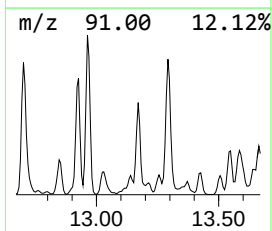
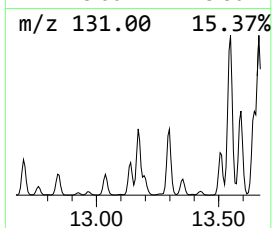
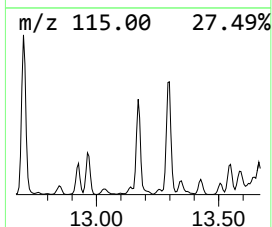
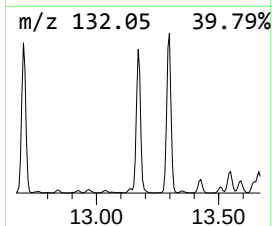
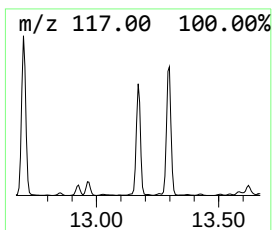
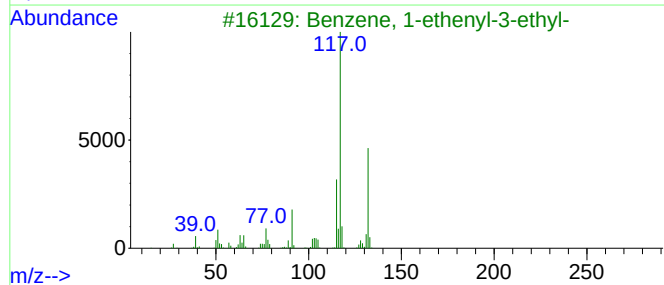
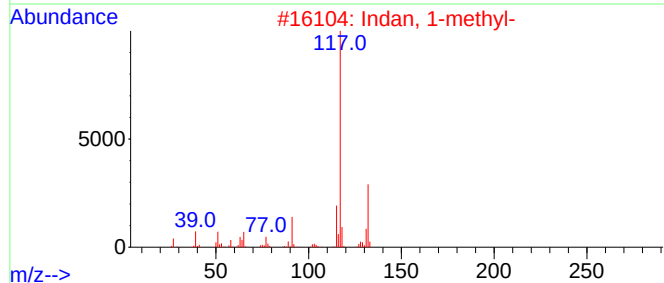
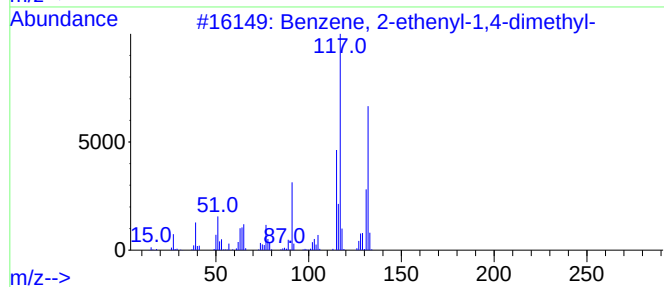
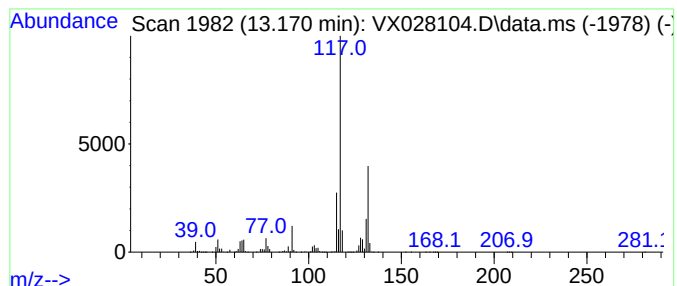
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.170	124.47 ug/l	4160510	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92
2			Indan, 1-methyl-	132	C10H12	000767-58-8	90
3			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	90
4			1-Phenyl-1-butene	132	C10H12	000824-90-8	90
5			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VO041422\
Data File : VX028104.D
Acq On : 14 Apr 2022 14:04
Operator : JC/MD
Sample : N2403-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-31R

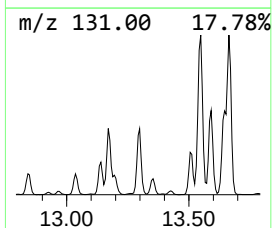
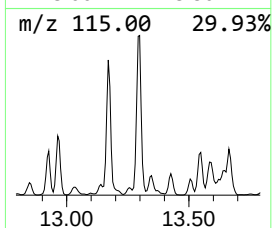
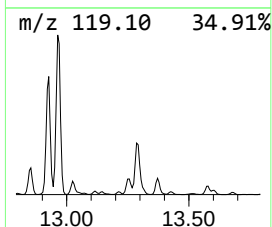
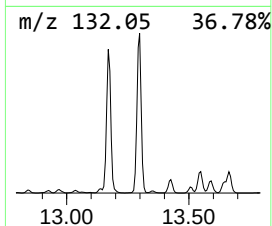
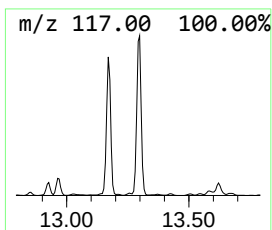
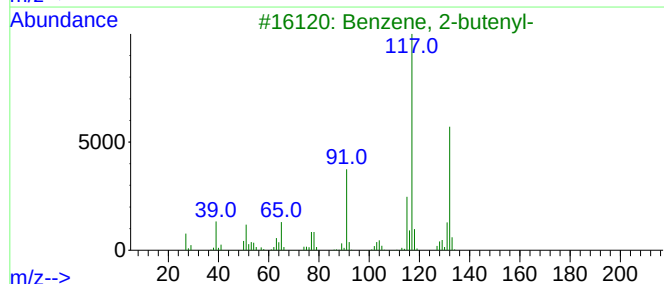
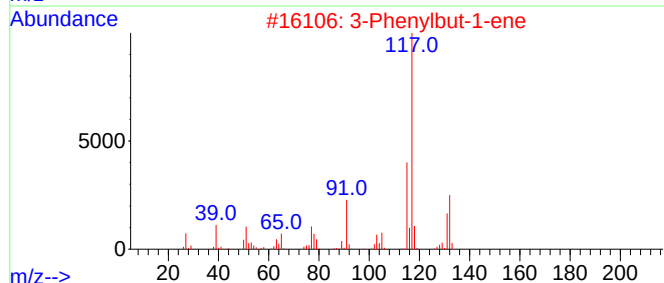
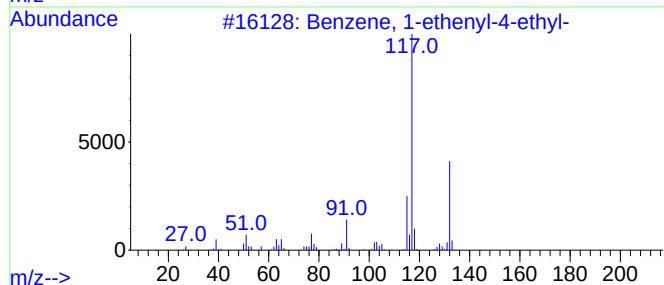
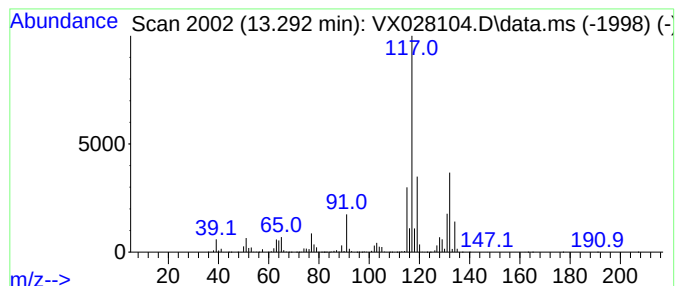
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	226.81 ug/l	7581280	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			3-Phenylbut-1-ene	132	C10H12	000934-10-1	81
3			Benzene, 2-butenyl-	132	C10H12	001560-06-1	81
4			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	81
5			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	78



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041422\
Data File : VX028104.D
Acq On : 14 Apr 2022 14:04
Operator : JC/MD
Sample : N2403-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-31R

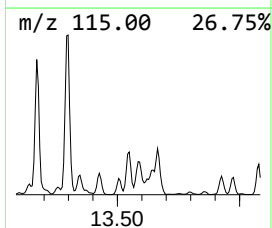
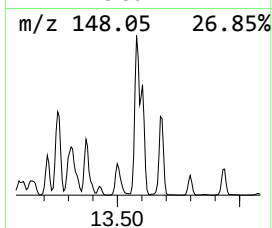
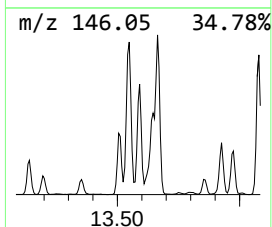
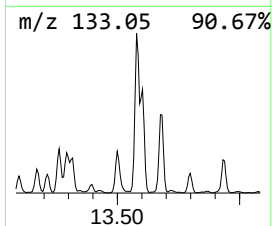
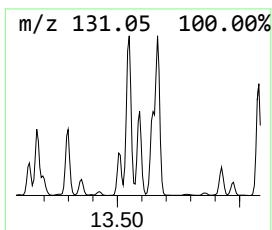
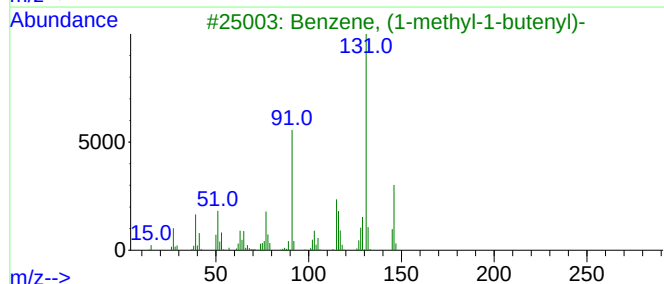
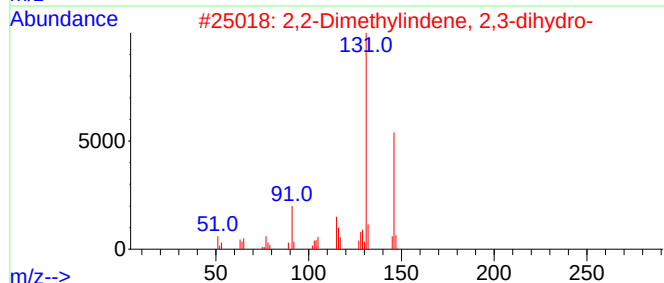
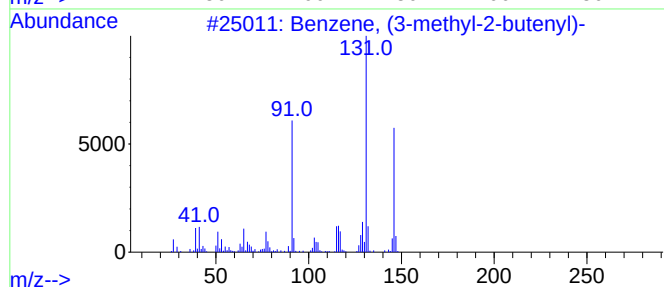
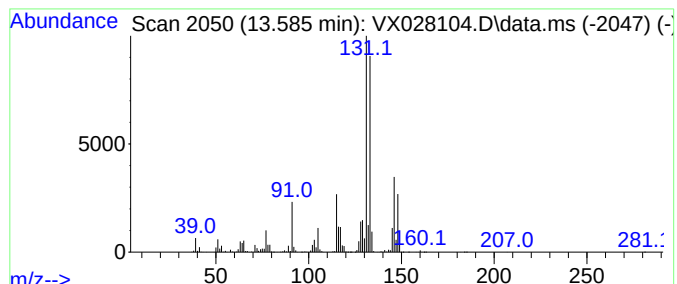
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, (3-methyl-2-butenyl)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.585	98.01 ug/l	3276060	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	70
2			2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	70
3			Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	70
4			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	55
5			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	55



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041422\
Data File : VX028104.D
Acq On : 14 Apr 2022 14:04
Operator : JC/MD
Sample : N2403-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-31R

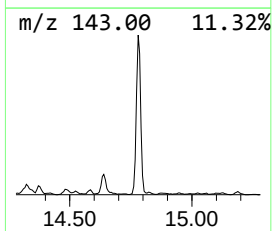
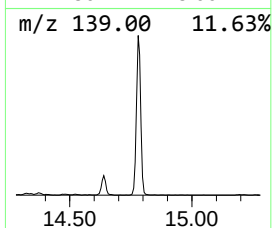
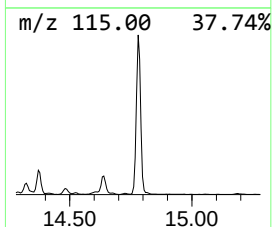
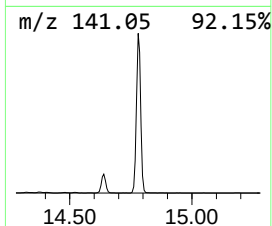
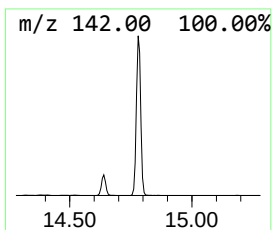
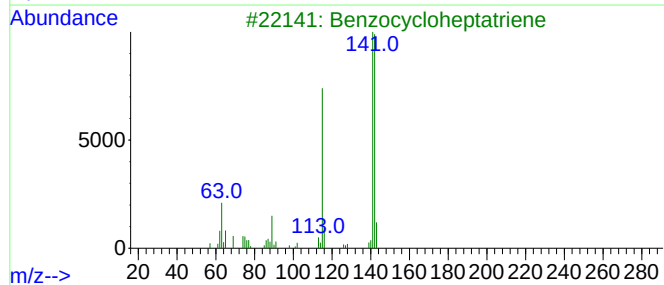
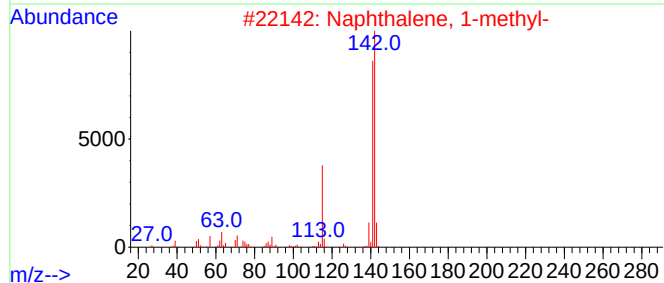
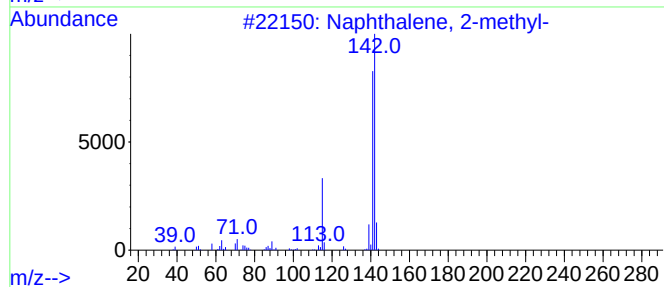
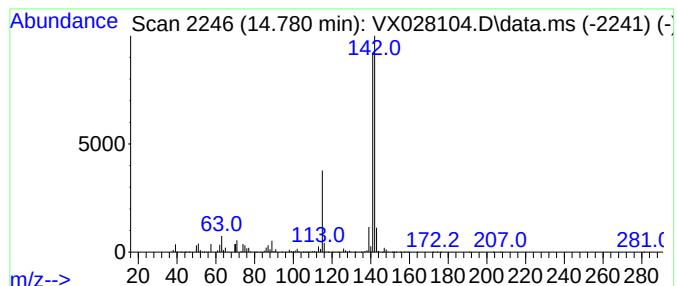
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Quant Title : SW846 8260

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

Peak Number 10 Benzocycloheptatriene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.780	84.88 ug/l	2837210	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			Benzocycloheptatriene	142	C11H10	000264-09-5	91
4			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
5			1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041422\
Data File : VX028104.D
Acq On : 14 Apr 2022 14:04
Operator : JC/MD
Sample : N2403-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-31R

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041222W.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, 1,3-di...	12.250	185.7	ug/l	6206010	4	12.024	1671300	50.0
Benzene, 2-ethy...	12.536	132.3	ug/l	4421780	4	12.024	1671300	50.0
Benzene, 4-ethy...	12.616	211.3	ug/l	7062400	4	12.024	1671300	50.0
Benzene, 1-ethe...	12.701	193.5	ug/l	6466570	4	12.024	1671300	50.0
Benzene, 1,2,3,...	12.926	133.7	ug/l	4468660	4	12.024	1671300	50.0
Benzene, 1,2,4,...	12.963	175.2	ug/l	5854690	4	12.024	1671300	50.0
Benzene, 2-ethe...	13.170	124.5	ug/l	4160510	4	12.024	1671300	50.0
Benzene, 1-ethe...	13.292	226.8	ug/l	7581280	4	12.024	1671300	50.0
Benzene, (3-met...	13.585	98.0	ug/l	3276060	4	12.024	1671300	50.0
Benzocyclohepta...	14.780	84.9	ug/l	2837210	4	12.024	1671300	50.0