

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX102924\
 Data File : VX043589.D
 Acq On : 29 Oct 2024 12:10
 Operator : JC/MD
 Sample : P4596-07 4X
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TOTE-WATER-COMP

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\82X102824W.M
 Title : SW846 8260

Signal : TIC: VX043589.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.258	22	28	35	rBV2	70699	79316	9.02%	0.864%
2	1.581	71	81	82	rBV6	8063	17205	1.96%	0.187%
3	2.386	207	213	221	rBV	20514	33687	3.83%	0.367%
4	2.569	235	243	250	rVB	8240	18990	2.16%	0.207%
5	2.941	296	304	319	rBV	22422	60223	6.85%	0.656%
6	3.428	377	384	392	rVB4	3560	9063	1.03%	0.099%
7	4.599	565	576	581	rBV4	3885	12442	1.42%	0.135%
8	5.379	693	704	718	rBV2	97337	291360	33.14%	3.172%
9	5.544	720	731	749	rBV	158151	465162	52.90%	5.065%
10	5.952	787	798	811	rBV	111905	299199	34.03%	3.258%
11	6.757	921	930	944	rBV	262045	647515	73.64%	7.050%
12	8.647	1233	1240	1248	rBV	547015	879260	100.00%	9.573%
13	8.720	1248	1252	1258	rVV2	107939	194509	22.12%	2.118%
14	8.775	1258	1261	1280	rVB	19090	46604	5.30%	0.507%
15	10.049	1465	1470	1488	rBV	521989	774686	88.11%	8.434%
16	10.195	1490	1494	1502	rVV	31113	44213	5.03%	0.481%
17	10.299	1505	1511	1525	rVB	118354	187570	21.33%	2.042%
18	10.640	1562	1567	1574	rBV	100195	133062	15.13%	1.449%
19	10.964	1615	1620	1627	rVB	13132	17493	1.99%	0.190%
20	11.079	1634	1639	1653	rBV	450109	593427	67.49%	6.461%
21	11.305	1670	1676	1683	rBV	21381	28842	3.28%	0.314%
22	11.378	1683	1688	1696	rVV	102770	184330	20.96%	2.007%
23	11.451	1696	1700	1707	rVB	56874	73704	8.38%	0.802%
24	11.610	1721	1726	1734	rVB	83414	104825	11.92%	1.141%
25	11.750	1744	1749	1761	rVB	244370	298952	34.00%	3.255%
26	11.890	1764	1772	1776	rVB2	12100	17776	2.02%	0.194%
27	11.969	1778	1785	1788	rBV	28187	34826	3.96%	0.379%
28	12.018	1788	1793	1800	rVV	578912	750751	85.38%	8.174%
29	12.085	1800	1804	1809	rVB	145120	176817	20.11%	1.925%
30	12.244	1823	1830	1833	rBV	131755	192583	21.90%	2.097%
31	12.317	1837	1842	1852	rVB3	66239	102496	11.66%	1.116%
32	12.402	1852	1856	1862	rBV4	12831	21208	2.41%	0.231%
33	12.463	1862	1866	1872	rVB	33017	41414	4.71%	0.451%
34	12.530	1872	1877	1879	rBV	47474	60746	6.91%	0.661%
35	12.555	1879	1881	1885	rVB	44016	44525	5.06%	0.485%
36	12.610	1885	1890	1893	rBV	56159	70160	7.98%	0.764%

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37	12.646	1893	1896	1900	rVB	21017	24775	2.82%	0.270%
38	12.695	1900	1904	1915	rBV2	61647	113180	12.87%	1.232%
39	12.792	1916	1920	1925	rBV3	7550	10491	1.19%	0.114%
40	12.847	1925	1929	1934	rVB2	30709	38542	4.38%	0.420%
41	12.921	1934	1941	1944	rBV	40982	51328	5.84%	0.559%
42	12.963	1944	1948	1954	rVB	67925	87690	9.97%	0.955%
43	13.024	1954	1958	1959	rBV	10322	11998	1.36%	0.131%
44	13.164	1973	1981	1986	rBV	82016	117902	13.41%	1.284%
45	13.213	1986	1989	1992	rVV3	12525	15708	1.79%	0.171%
46	13.250	1992	1995	1997	rVV2	15227	20171	2.29%	0.220%
47	13.292	1997	2002	2007	rVV2	209099	304150	34.59%	3.311%
48	13.329	2007	2008	2012	rVV3	8992	10631	1.21%	0.116%
49	13.372	2012	2015	2018	rVV	7964	9546	1.09%	0.104%
50	13.420	2018	2023	2032	rVB	195964	241692	27.49%	2.631%
51	13.506	2032	2037	2039	rBV3	13059	18479	2.10%	0.201%
52	13.542	2039	2043	2046	rVV	27681	39134	4.45%	0.426%
53	13.585	2046	2050	2055	rVV3	35904	67051	7.63%	0.730%
54	13.664	2055	2063	2069	rVB2	35995	83437	9.49%	0.908%
55	13.774	2074	2081	2088	rBV	73838	103233	11.74%	1.124%
56	13.853	2088	2094	2098	rVB2	42408	58110	6.61%	0.633%
57	13.926	2098	2106	2110	rBV2	49885	72933	8.29%	0.794%
58	13.969	2110	2113	2117	rVB2	16161	20027	2.28%	0.218%
59	14.073	2126	2130	2134	rBV	31890	38630	4.39%	0.421%
60	14.225	2148	2155	2161	rBV	109944	175063	19.91%	1.906%
61	14.317	2167	2170	2174	rBV2	11632	15321	1.74%	0.167%
62	14.371	2174	2179	2184	rVV	30927	39207	4.46%	0.427%
63	14.475	2191	2196	2203	rBV2	76847	103233	11.74%	1.124%
64	14.634	2215	2222	2226	rBV3	48106	96669	10.99%	1.052%
65	14.670	2226	2228	2233	rVB	18778	21046	2.39%	0.229%
66	14.780	2240	2246	2250	rVV2	28230	41843	4.76%	0.456%
67	14.817	2250	2252	2255	rVV3	11916	15791	1.80%	0.172%
68	14.902	2260	2266	2272	rBV4	10496	19300	2.20%	0.210%
69	15.054	2285	2291	2297	rBV4	4934	8987	1.02%	0.098%
70	15.182	2306	2312	2321	rBV3	20984	42564	4.84%	0.463%
71	15.475	2356	2360	2364	rBV6	9891	15996	1.82%	0.174%
72	15.609	2377	2382	2386	rBV4	7623	12949	1.47%	0.141%
73	15.652	2386	2389	2397	rVB6	4961	8999	1.02%	0.098%

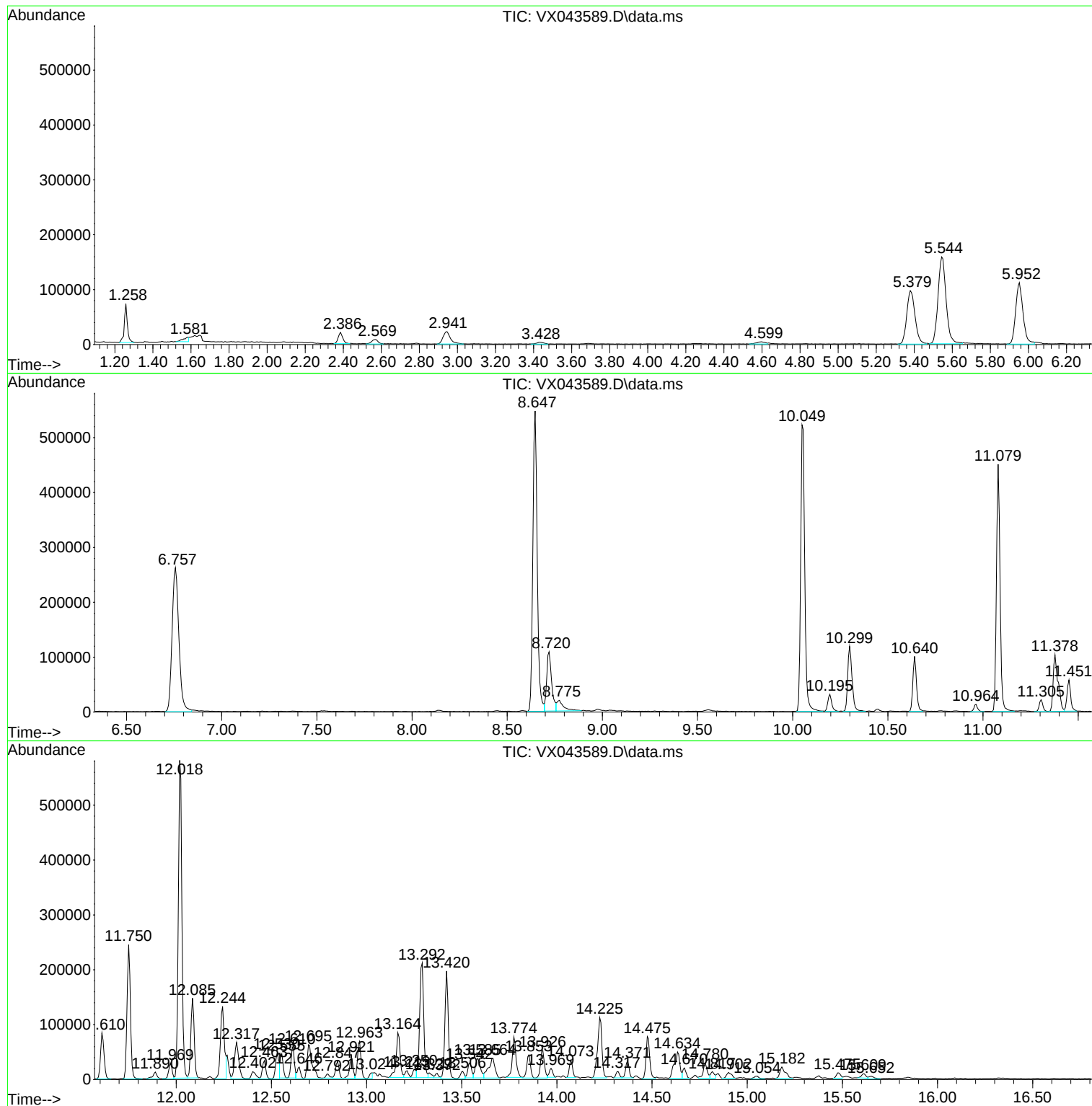
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ClientSampleId :
TOTE-WATER-COMP

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102824W.M
Quant Title : SW846 8260

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



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MSVOA_X
ClientSampleId :
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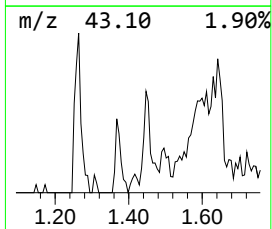
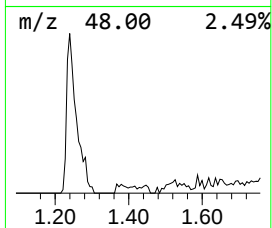
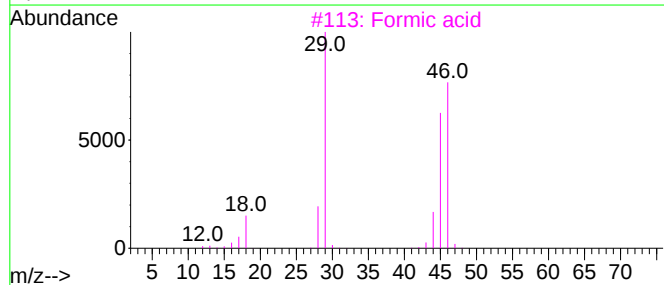
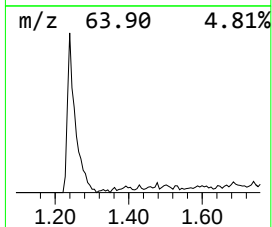
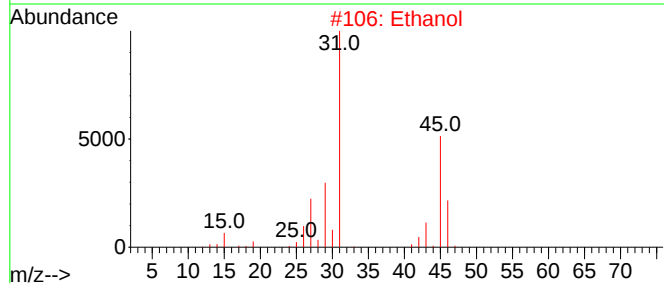
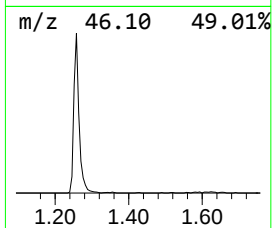
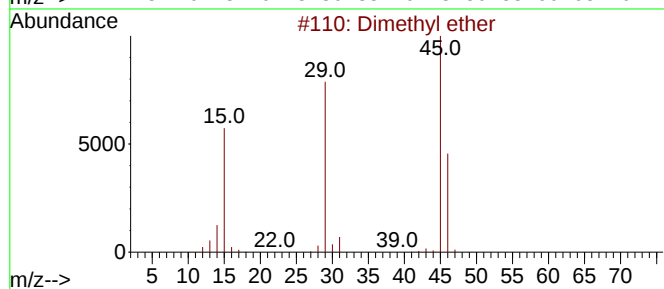
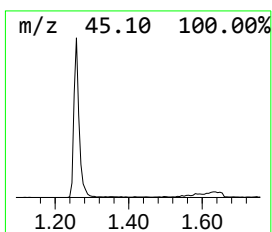
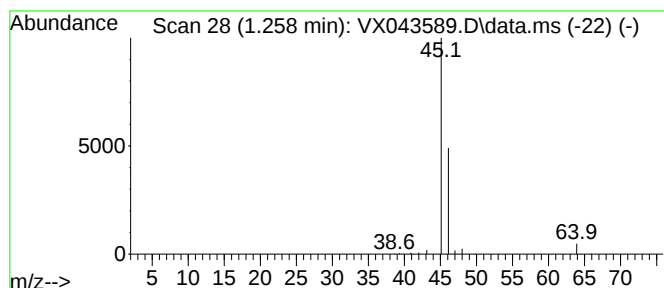
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102824W.M
Quant Title : SW846 8260

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

Peak Number 1 Dimethyl ether Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.258	8.53 ug/l	79316	Pentafluorobenzene	5.544

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dimethyl ether	46	C2H6O	000115-10-6	64
2			Ethanol	46	C2H6O	000064-17-5	7
3			Formic acid	46	CH2O2	000064-18-6	5
4			Oxalic acid	90	C2H2O4	000144-62-7	4
5			Methane, nitroso-	45	CH3NO	000865-40-7	3



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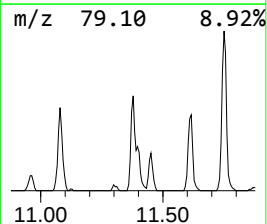
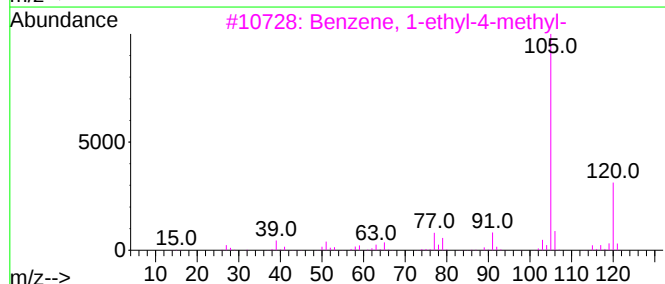
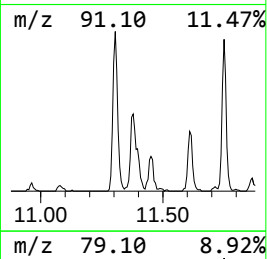
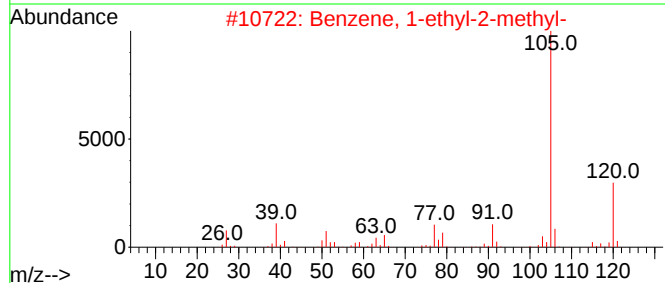
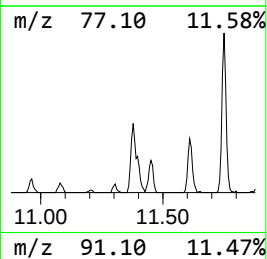
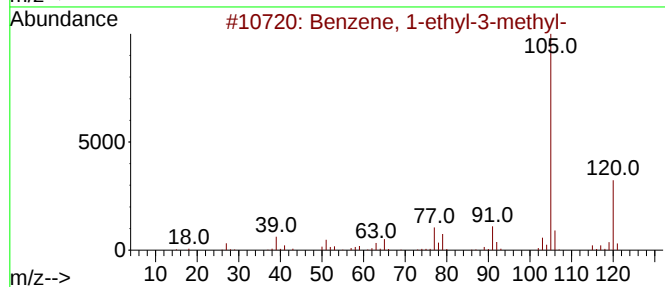
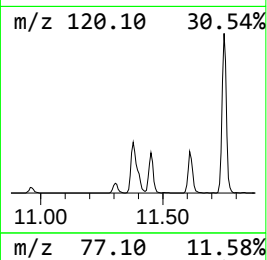
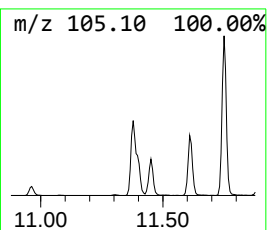
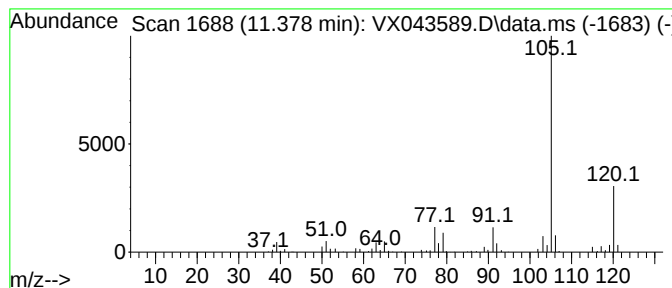
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102824W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.378	12.28 ug/l	184330	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	97
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5			Mesitylene	120	C9H12	000108-67-8	91



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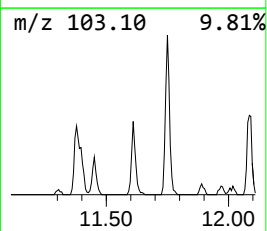
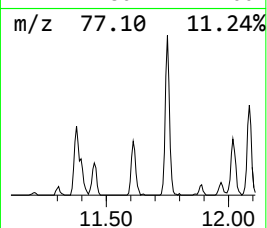
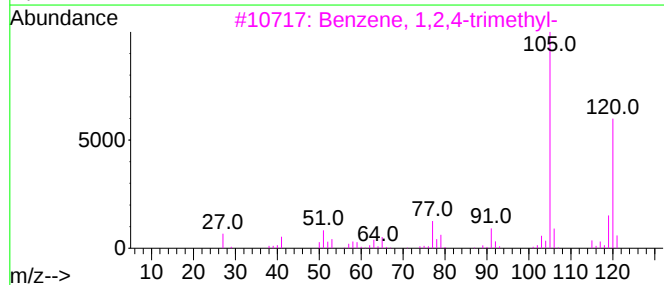
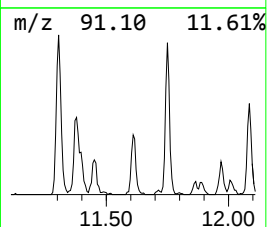
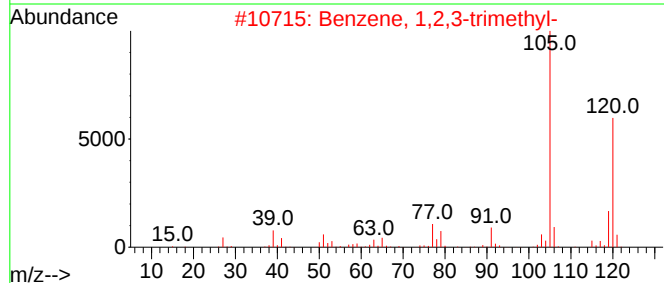
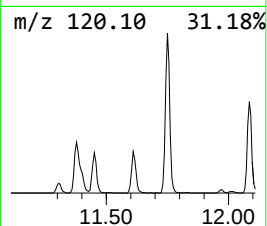
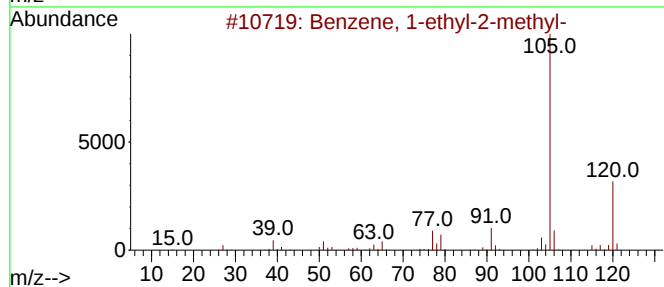
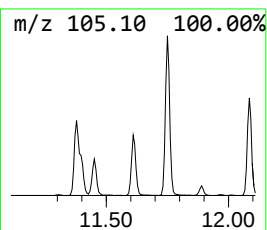
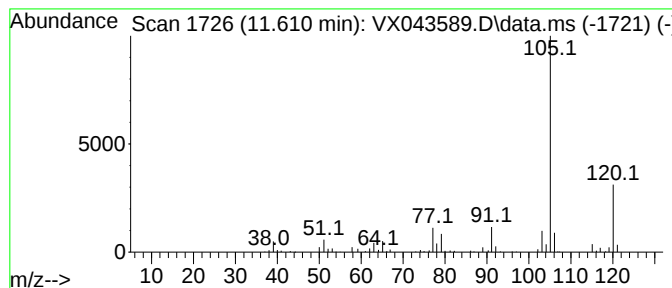
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TIC Library : C:\Database\NIST20.L
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Peak Number 3 Benzene, 1-ethyl-2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.610	6.98 ug/l	104825	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, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
3			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4			Mesitylene	120	C9H12	000108-67-8	91
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



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ClientSampleId :
TOTE-WATER-COMP

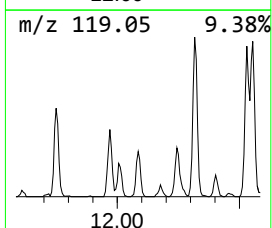
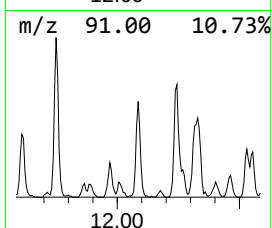
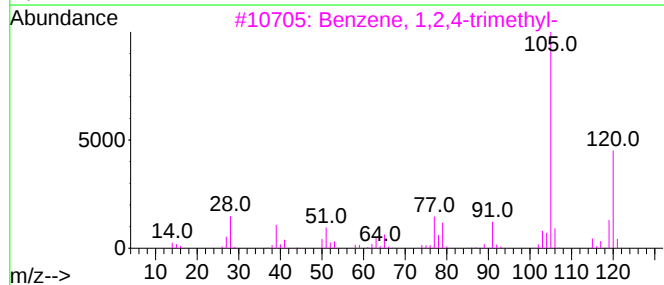
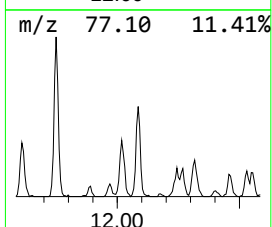
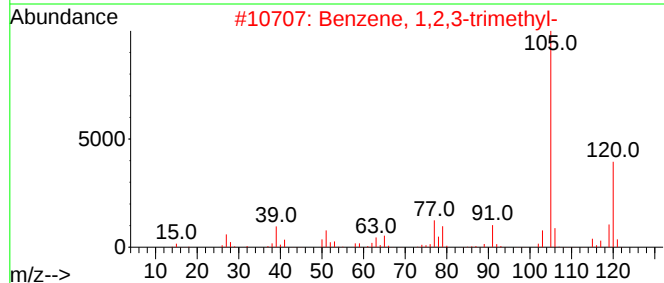
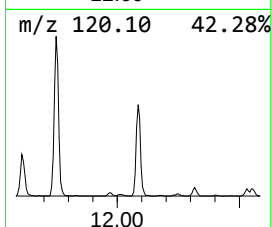
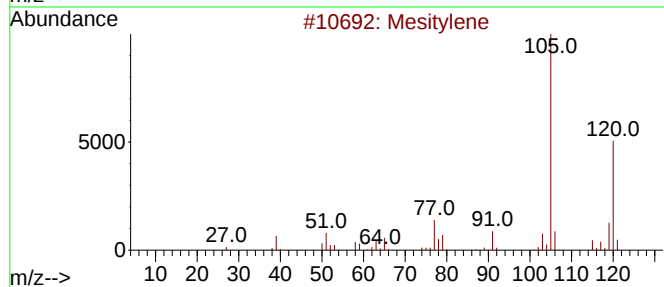
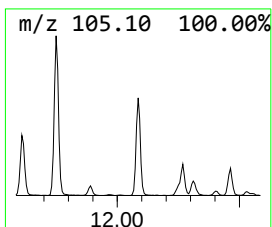
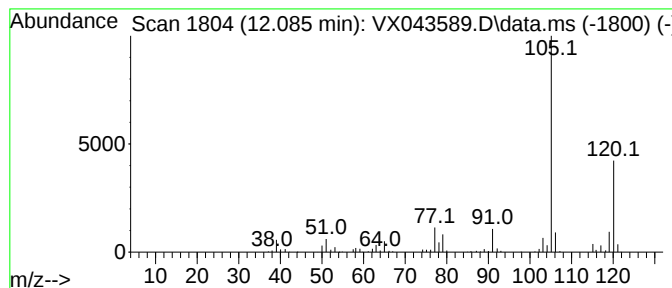
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102824W.M
Quant Title : SW846 8260

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

Peak Number 4 Mesitylene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.085	11.78 ug/l	176817	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	97
3			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX102924\
Data File : VX043589.D
Acq On : 29 Oct 2024 12:10
Operator : JC/MD
Sample : P4596-07 4X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TOTE-WATER-COMP

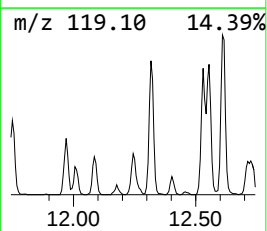
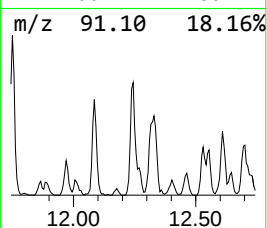
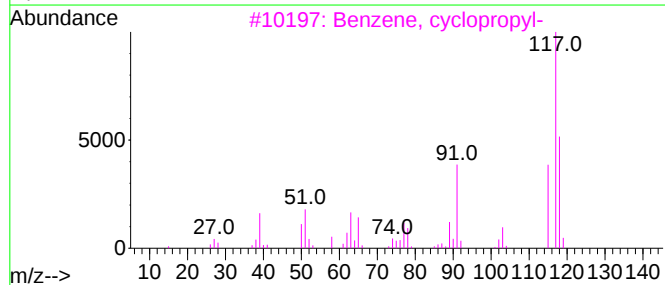
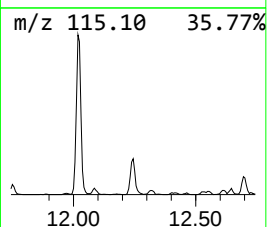
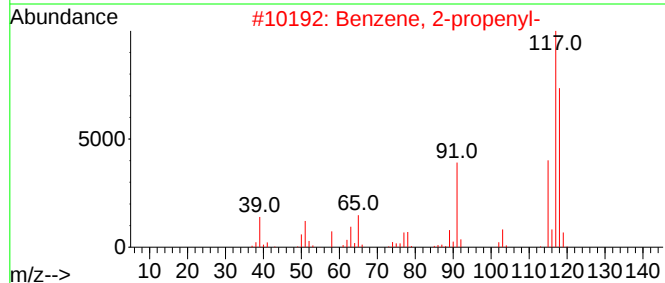
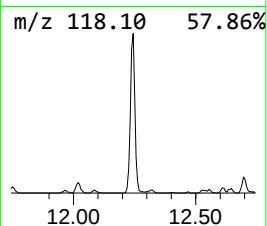
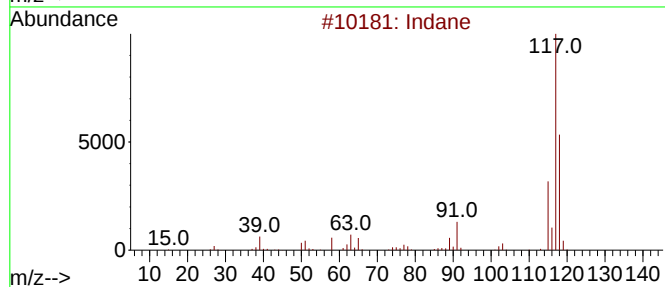
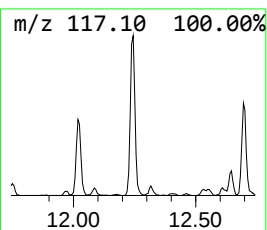
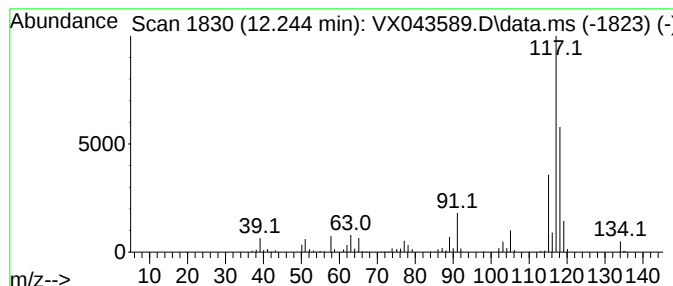
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Quant Title : SW846 8260

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

Peak Number 5 Indane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	12.83 ug/l	192583	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	89
2			Benzene, 2-propenyl-	118	C9H10	000300-57-2	76
3			Benzene, cyclopropyl-	118	C9H10	000873-49-4	76
4			Delatacyclene	118	C9H10	007785-10-6	58
5			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	58



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX102924\
Data File : VX043589.D
Acq On : 29 Oct 2024 12:10
Operator : JC/MD
Sample : P4596-07 4X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TOTE-WATER-COMP

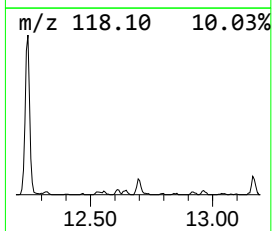
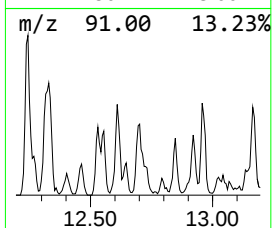
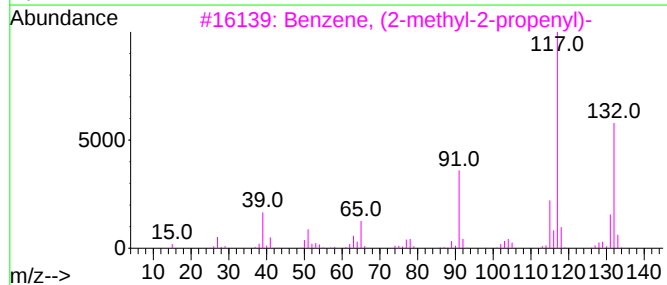
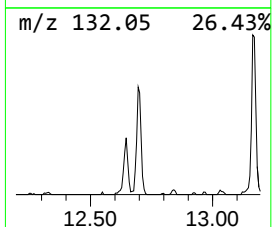
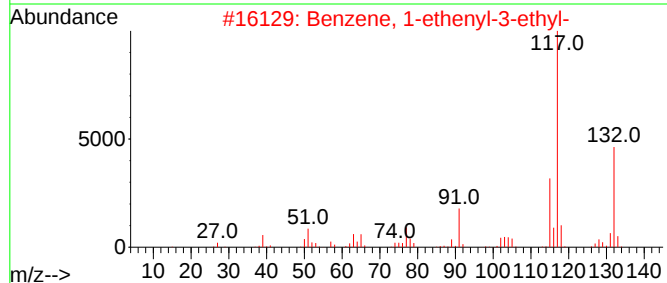
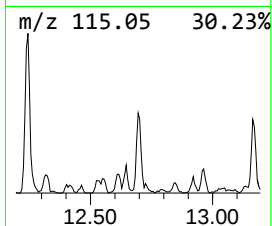
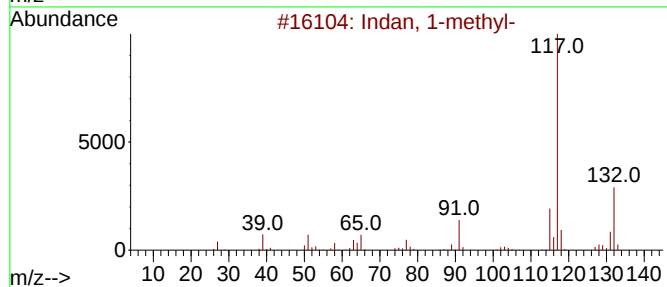
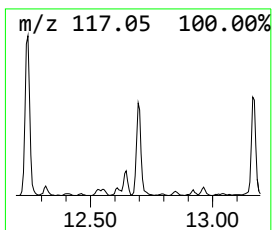
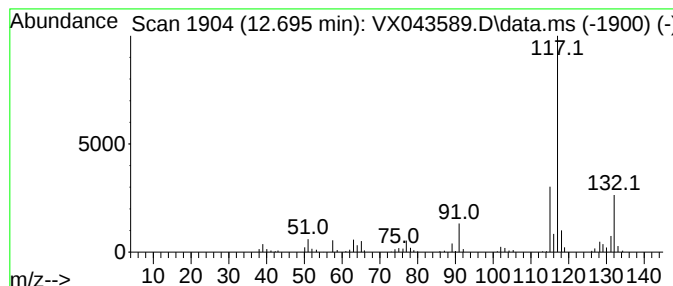
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Quant Title : SW846 8260

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

Peak Number 6 Indan, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.695	7.54 ug/l	113180	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	91
3			Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	80
4			Benzene, 2-butenyl-	132	C10H12	001560-06-1	80
5			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	80



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX102924\
Data File : VX043589.D
Acq On : 29 Oct 2024 12:10
Operator : JC/MD
Sample : P4596-07 4X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TOTE-WATER-COMP

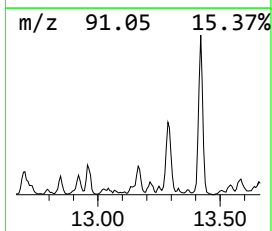
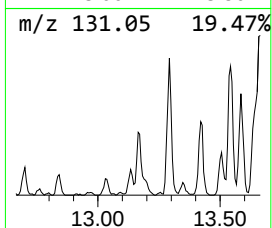
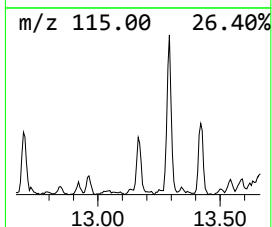
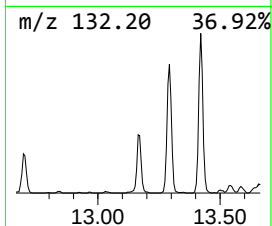
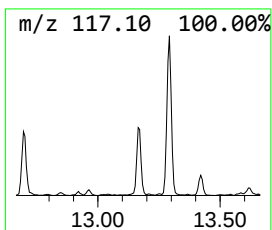
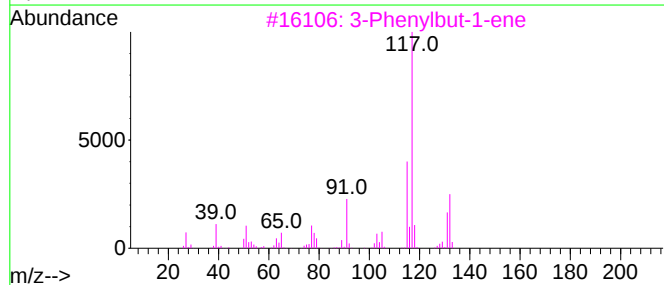
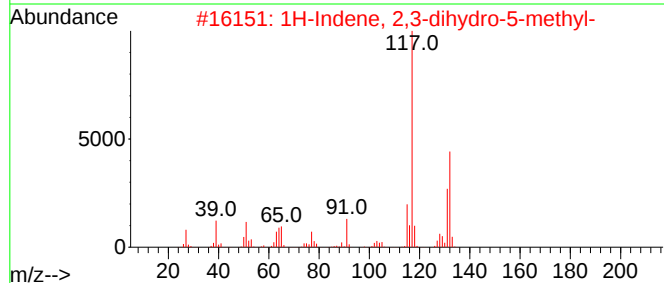
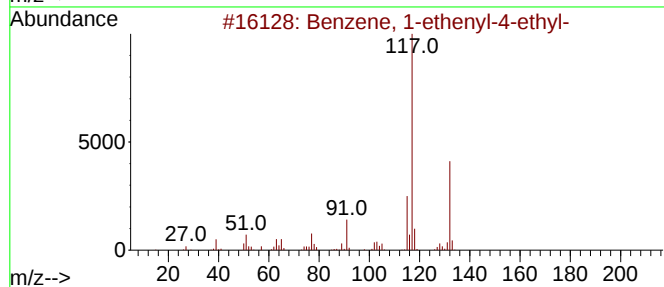
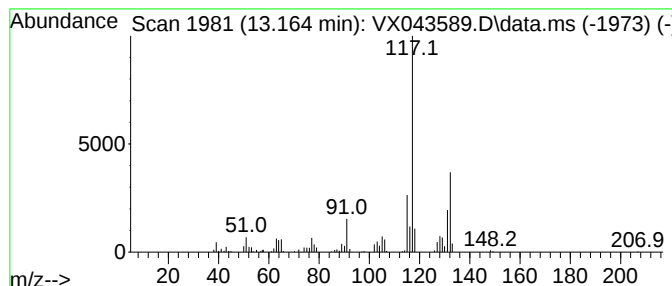
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.164	7.85 ug/l	117902	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			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	91
3			3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
4			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	87
5			Benzene, 2-butenyl-	132	C10H12	001560-06-1	83



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX102924\
Data File : VX043589.D
Acq On : 29 Oct 2024 12:10
Operator : JC/MD
Sample : P4596-07 4X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TOTE-WATER-COMP

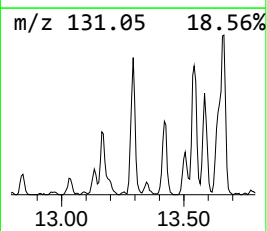
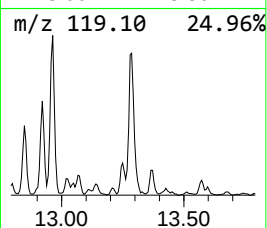
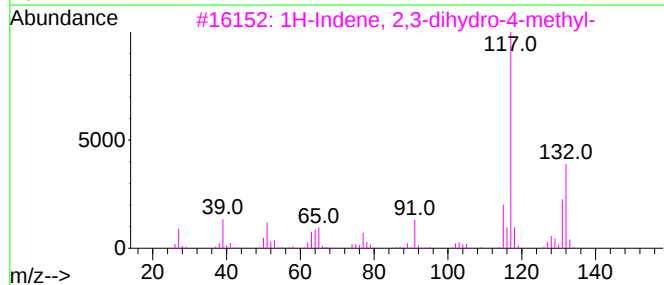
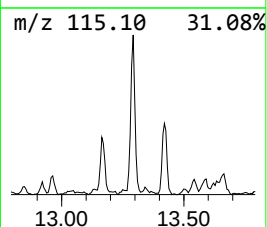
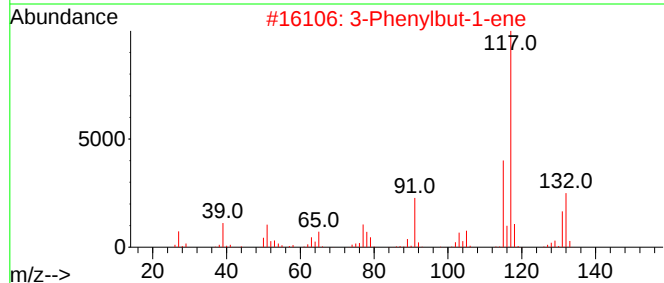
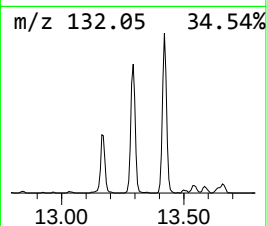
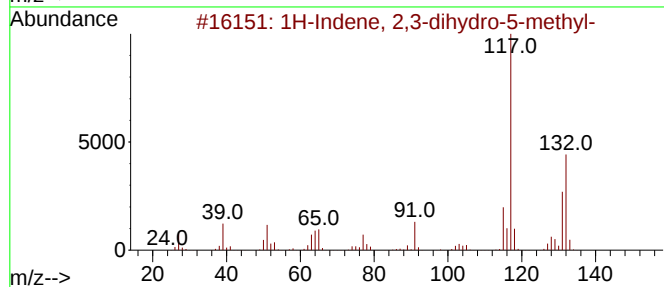
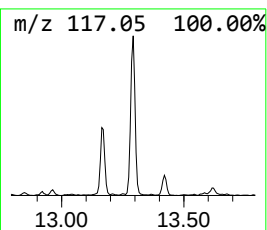
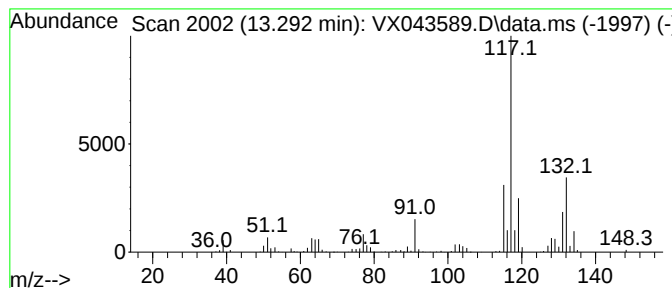
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Quant Title : SW846 8260

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

Peak Number 8 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	20.26 ug/l	304150	1,4-Dichlorobenzene-d4	12.024

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	89	
2	3-Phenylbut-1-ene	132	C10H12	000934-10-1	81	
3	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	76	
4	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	76	
5	Indan, 1-methyl-	132	C10H12	000767-58-8	76	



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX102924\
Data File : VX043589.D
Acq On : 29 Oct 2024 12:10
Operator : JC/MD
Sample : P4596-07 4X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TOTE-WATER-COMP

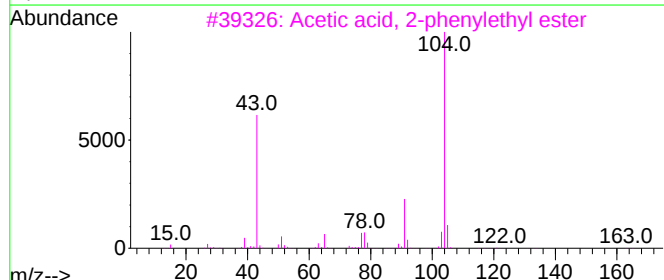
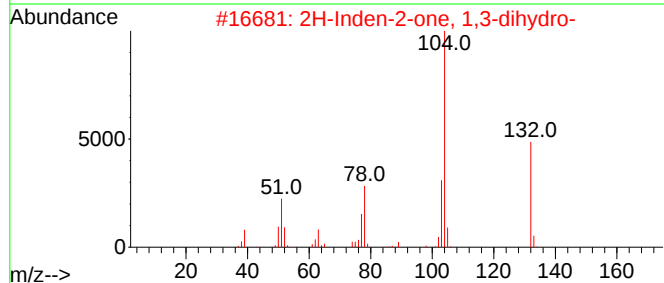
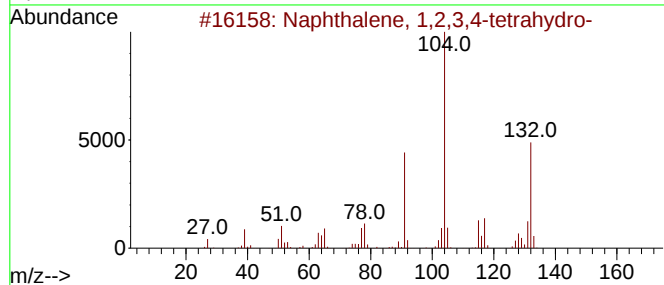
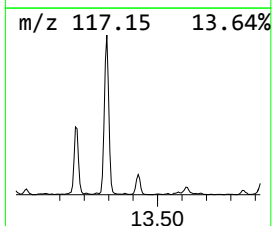
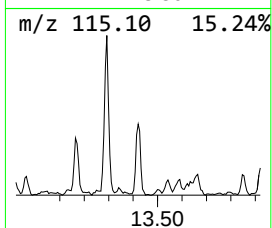
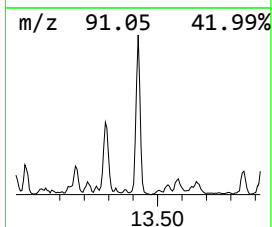
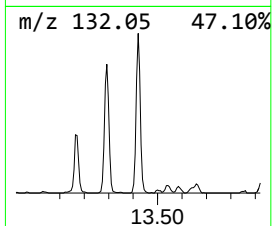
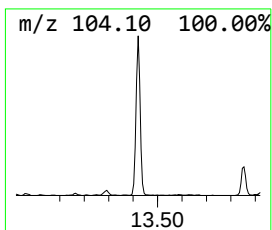
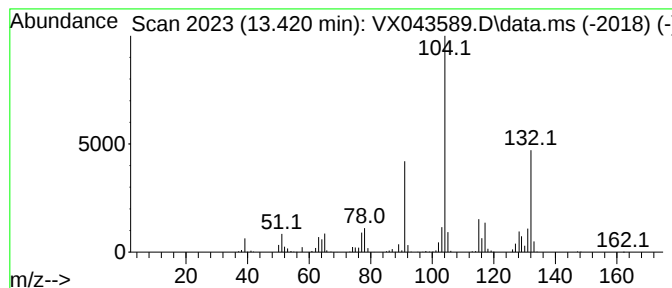
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.420	16.10 ug/l	241692	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	97
2			2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	58
3			Acetic acid, 2-phenylethyl ester	164	C10H12O2	000103-45-7	38
4			Glutaric acid, 2-methylbenzyl pr...	278	C16H22O4	1000371-32-4	38
5			2-Bromopropionic acid, 2-phenyle...	256	C11H13BrO2	043216-34-8	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX102924\
Data File : VX043589.D
Acq On : 29 Oct 2024 12:10
Operator : JC/MD
Sample : P4596-07 4X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 1 Sample Multiplier: 1

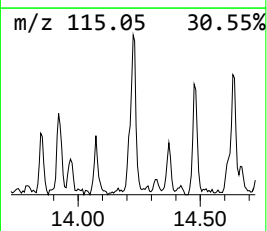
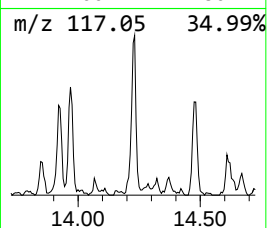
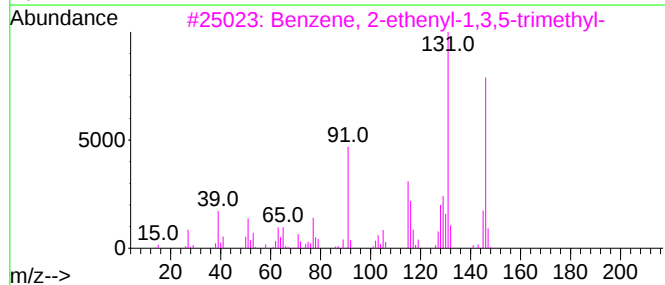
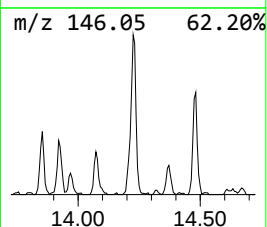
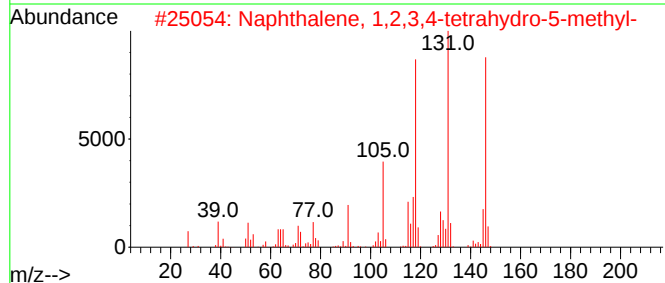
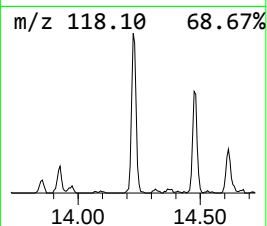
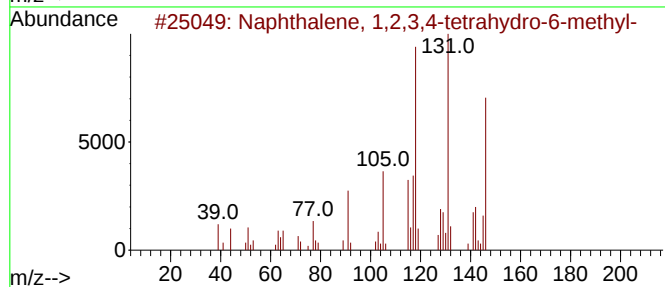
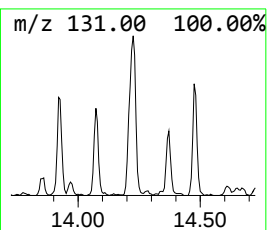
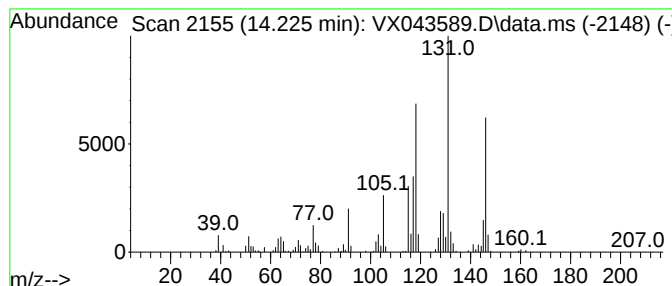
Instrument :
MSVOA_X
ClientSampleId :
TOTE-WATER-COMP

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102824W.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.	
14.225	11.66 ug/l	175063	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	98
2	Naphthalene, 1,2,3,4-tetrahydro-...		146	C11H14	002809-64-5	95
3	Benzene, 2-ethenyl-1,3,5-trimethyl-		146	C11H14	000769-25-5	86
4	Benzene, (1,2-dimethyl-1-propenyl)-		146	C11H14	000769-57-3	68
5	Benzene, (3-methyl-2-butenyl)-		146	C11H14	004489-84-3	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX102924\
Data File : VX043589.D
Acq On : 29 Oct 2024 12:10
Operator : JC/MD
Sample : P4596-07 4X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TOTE-WATER-COMP

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102824W.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
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Benzene, 1-ethy...	11.378	12.3	ug/l	184330	4	12.024	750751	50.0
Benzene, 1-ethy...	11.610	7.0	ug/l	104825	4	12.024	750751	50.0
Mesitylene	12.085	11.8	ug/l	176817	4	12.024	750751	50.0
Indane	12.244	12.8	ug/l	192583	4	12.024	750751	50.0
Indan, 1-methyl-	12.695	7.5	ug/l	113180	4	12.024	750751	50.0
Benzene, 1-ethe...	13.164	7.8	ug/l	117902	4	12.024	750751	50.0
1H-Indene, 2,3-...	13.292	20.3	ug/l	304150	4	12.024	750751	50.0
Naphthalene, 1,...	13.420	16.1	ug/l	241692	4	12.024	750751	50.0
Naphthalene, 1,...	14.225	11.7	ug/l	175063	4	12.024	750751	50.0