

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX012722\
 Data File : VX026632.D
 Acq On : 27 Jan 2022 18:12
 Operator : JC/MD
 Sample : N1296-09
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-10R

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\82X011122W.M
 Title : SW846 8260

Signal : TIC: VX026632.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.374	43	47	53	rVB	37807	36423	2.41%	0.186%
2	1.758	104	110	126	rVB	438071	619967	41.01%	3.162%
3	1.959	136	143	153	rVB	16973	25594	1.69%	0.131%
4	2.221	181	186	195	rVB	14741	22097	1.46%	0.113%
5	2.349	198	207	220	rBV	33134	70607	4.67%	0.360%
6	2.788	270	279	282	rBV	116360	248999	16.47%	1.270%
7	2.831	282	286	292	rVB	174962	319823	21.16%	1.631%
8	3.111	322	332	348	rBV	208957	476075	31.49%	2.428%
9	3.733	427	434	443	rBV	9712	22767	1.51%	0.116%
10	3.843	443	452	458	rBV3	14144	36953	2.44%	0.188%
11	3.910	458	463	472	rVB2	6678	15334	1.01%	0.078%
12	4.105	489	495	503	rVB4	7362	18828	1.25%	0.096%
13	4.221	503	514	520	rVV	12997	37715	2.49%	0.192%
14	4.306	520	528	541	rVV	77330	224132	14.83%	1.143%
15	4.434	541	549	563	rVB4	7810	26914	1.78%	0.137%
16	5.117	649	661	662	rBV4	6612	16038	1.06%	0.082%
17	5.196	668	674	684	rVB3	7982	26286	1.74%	0.134%
18	5.391	694	706	714	rBV	78511	227147	15.03%	1.158%
19	5.483	714	721	725	rVV	34008	104578	6.92%	0.533%
20	5.556	726	733	740	rVV	134539	404749	26.77%	2.064%
21	5.629	741	745	763	rVV2	47058	128538	8.50%	0.655%
22	5.830	767	778	790	rVV4	35715	113965	7.54%	0.581%
23	5.964	790	800	807	rVV	85174	226527	14.99%	1.155%
24	6.044	807	813	822	rVV2	28737	79331	5.25%	0.405%
25	6.166	823	833	842	rVV5	33186	129965	8.60%	0.663%
26	6.263	842	849	858	rVV3	45140	134224	8.88%	0.684%
27	6.361	858	865	875	rVB3	22861	61603	4.08%	0.314%
28	6.763	921	931	942	rBV	208106	513850	33.99%	2.620%
29	6.848	942	945	955	rVB5	12114	29201	1.93%	0.149%
30	7.178	990	999	1006	rBV2	10448	22053	1.46%	0.112%
31	7.385	1021	1033	1044	rBV3	50966	135026	8.93%	0.689%
32	7.659	1072	1078	1091	rVB2	11593	23201	1.53%	0.118%
33	7.854	1104	1110	1112	rBV2	9161	15337	1.01%	0.078%
34	7.885	1112	1115	1122	rVB3	11445	21282	1.41%	0.109%
35	7.988	1122	1132	1141	rVB3	12906	32872	2.17%	0.168%
36	8.110	1143	1152	1155	rBV	27270	55741	3.69%	0.284%

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37	8.147	1155	1158	1163	rVB	20419	30681	2.03%	0.156%
38	8.507	1211	1217	1226	rVB3	24622	44755	2.96%	0.228%
39	8.604	1226	1233	1235	rBV4	9314	16341	1.08%	0.083%
40	8.653	1235	1241	1247	rVV	418510	663820	43.91%	3.385%
41	8.720	1247	1252	1258	rVB	184855	277681	18.37%	1.416%
42	8.842	1264	1272	1280	rVB4	10135	22443	1.48%	0.114%
43	9.165	1320	1325	1336	rVB	15728	24466	1.62%	0.125%
44	10.055	1461	1471	1477	rBV	429946	568342	37.60%	2.898%
45	10.195	1489	1494	1500	rBV	1055877	1337322	88.47%	6.820%
46	10.305	1506	1512	1523	rVB	1165203	1511683	100.00%	7.709%
47	10.640	1562	1567	1578	rVB	304194	396690	26.24%	2.023%
48	10.963	1613	1620	1628	rVB	191735	239944	15.87%	1.224%
49	11.079	1634	1639	1649	rVB	310749	408836	27.05%	2.085%
50	11.305	1666	1676	1683	rBV	290854	347012	22.96%	1.770%
51	11.384	1683	1689	1690	rVV	332164	449999	29.77%	2.295%
52	11.402	1690	1692	1696	rVV	381914	389840	25.79%	1.988%
53	11.451	1696	1700	1708	rVB	323503	398939	26.39%	2.034%
54	11.616	1721	1727	1736	rBV	567683	725638	48.00%	3.700%
55	11.756	1744	1750	1756	rBV	1165109	1468559	97.15%	7.489%
56	11.866	1763	1768	1770	rBV	18794	24403	1.61%	0.124%
57	11.890	1770	1772	1779	rVB	25676	29268	1.94%	0.149%
58	11.969	1779	1785	1789	rBV	25482	33169	2.19%	0.169%
59	12.024	1789	1794	1800	rVB	405017	507973	33.60%	2.590%
60	12.091	1800	1805	1810	rBV	289280	368330	24.37%	1.878%
61	12.244	1823	1830	1838	rBV3	458535	666291	44.08%	3.398%
62	12.317	1838	1842	1852	rVB3	144818	226433	14.98%	1.155%
63	12.408	1852	1857	1862	rBV2	39670	52949	3.50%	0.270%
64	12.463	1862	1866	1872	rVB	54549	65532	4.34%	0.334%
65	12.530	1872	1877	1880	rBV	131752	174487	11.54%	0.890%
66	12.616	1886	1891	1894	rBV	278725	330759	21.88%	1.687%
67	12.646	1894	1896	1900	rVB	37982	39351	2.60%	0.201%
68	12.701	1900	1905	1913	rBV	212751	283150	18.73%	1.444%
69	12.847	1924	1929	1935	rVB2	24021	32580	2.16%	0.166%
70	12.920	1935	1941	1945	rBV	172018	221912	14.68%	1.132%
71	12.963	1945	1948	1954	rVB	224616	258710	17.11%	1.319%
72	13.170	1978	1982	1987	rVB	140947	159820	10.57%	0.815%
73	13.256	1992	1996	1997	rBV	12798	15196	1.01%	0.077%
74	13.292	1997	2002	2007	rVB2	331965	428510	28.35%	2.185%
75	13.341	2007	2010	2013	rBV3	16733	18825	1.25%	0.096%
76	13.426	2019	2024	2029	rVB	47391	61819	4.09%	0.315%

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77	13.506	2032	2037	2039	rBV2	13707	18157	1.20%	0.093%
78	13.542	2039	2043	2046	rVV	35315	51292	3.39%	0.262%
79	13.585	2046	2050	2056	rVV3	33861	67781	4.48%	0.346%
80	13.664	2060	2063	2070	rVB2	34642	53857	3.56%	0.275%
81	13.774	2075	2081	2092	rVV	662855	816218	53.99%	4.162%
82	13.871	2092	2097	2101	rVV	21720	29941	1.98%	0.153%
83	13.926	2101	2106	2110	rVV2	11038	15663	1.04%	0.080%
84	14.079	2126	2131	2138	rBV	23250	28694	1.90%	0.146%
85	14.213	2149	2153	2159	rBV	18646	31540	2.09%	0.161%
86	14.371	2176	2179	2184	rVB2	13448	15930	1.05%	0.081%
87	14.640	2218	2223	2232	rVB	250426	306967	20.31%	1.565%
88	14.780	2241	2246	2257	rVB	114852	147477	9.76%	0.752%

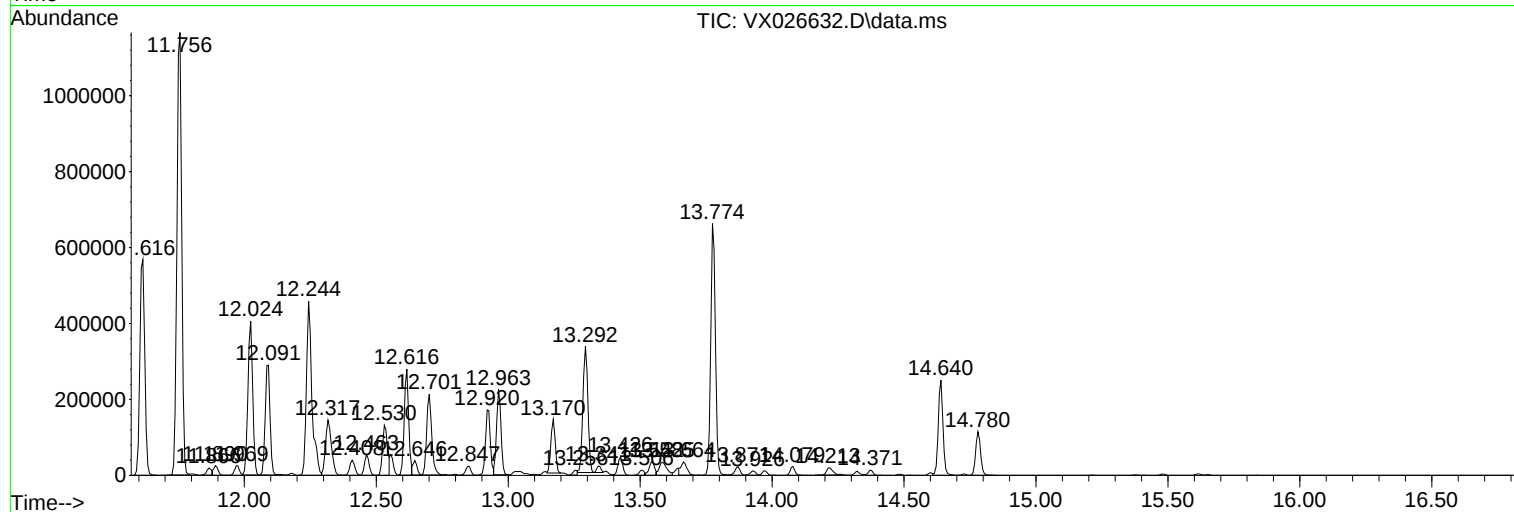
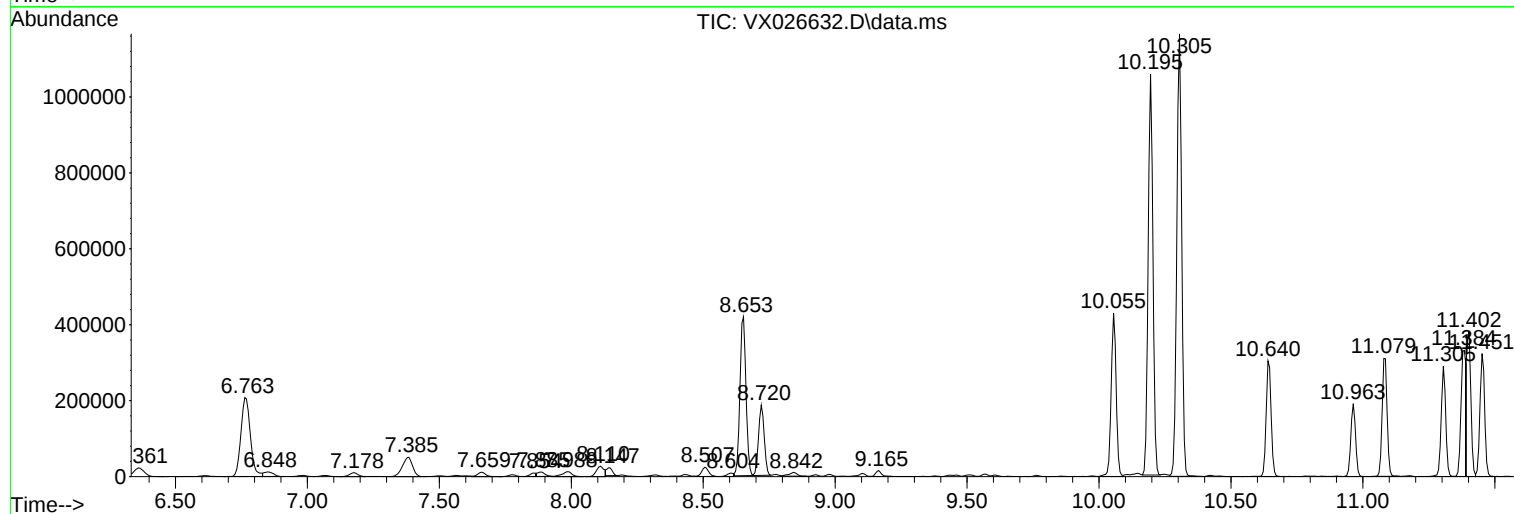
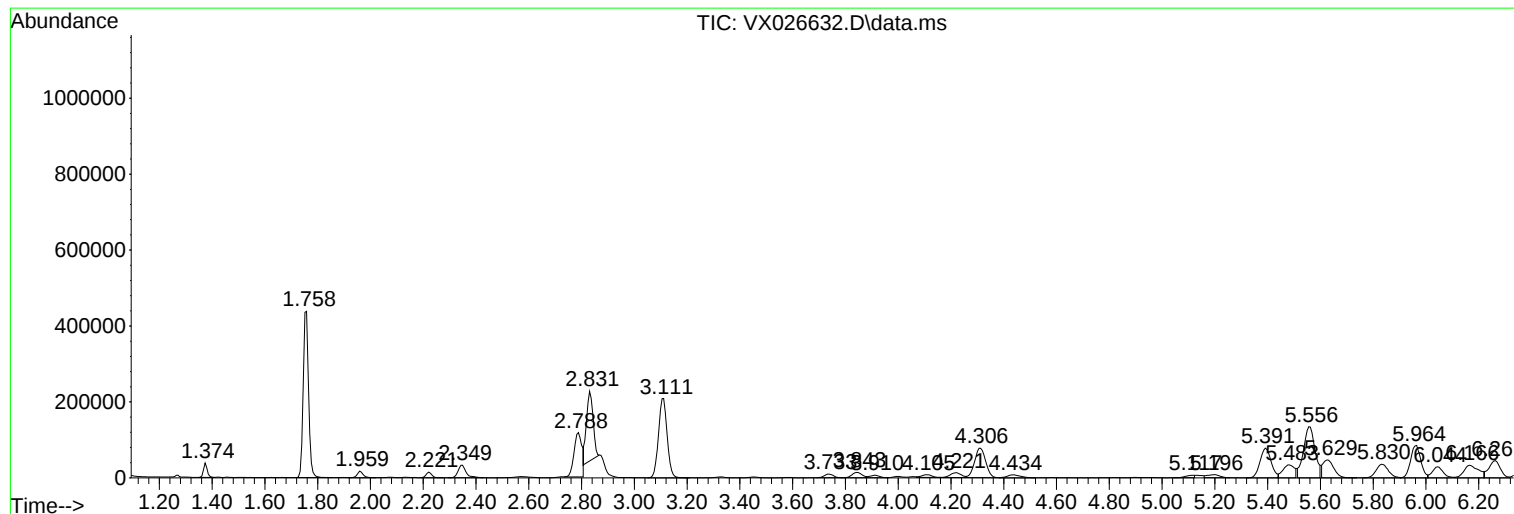
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X011122W.M
Quant Title : SW846 8260

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



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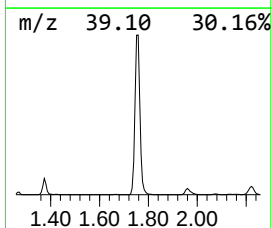
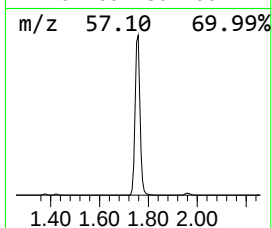
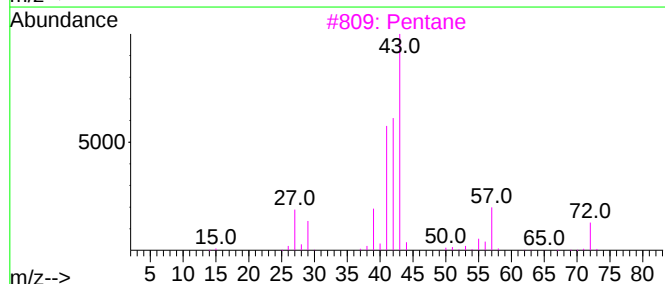
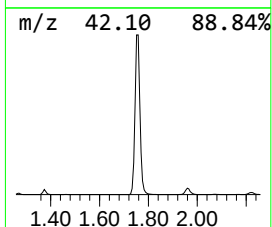
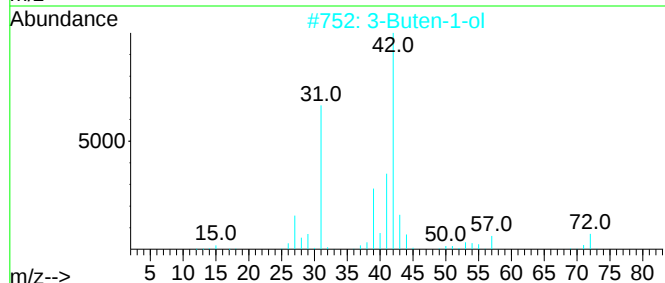
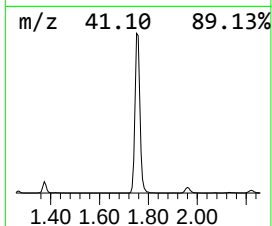
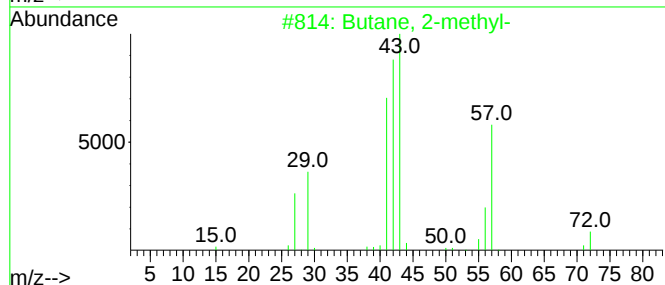
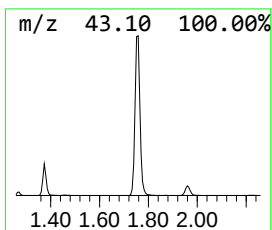
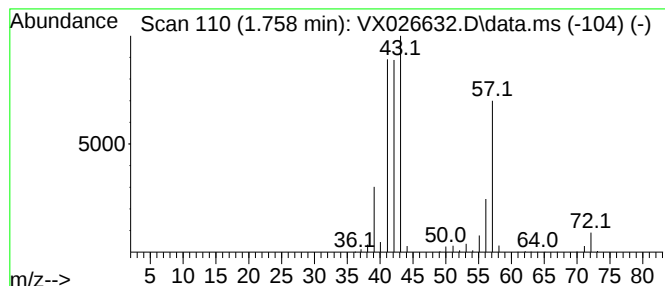
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X011122W.M
Quant Title : SW846 8260

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

Peak Number 1 Butane, 2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.758	76.59 ug/l	619967	Pentafluorobenzene	5.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methyl-	72	C5H12	000078-78-4	90
2			3-Buten-1-ol	72	C4H8O	000627-27-0	50
3			Pentane	72	C5H12	000109-66-0	36
4			1-Propene, 2-methyl-	56	C4H8	000115-11-7	10
5			1-Butene	56	C4H8	000106-98-9	9



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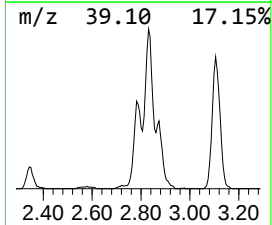
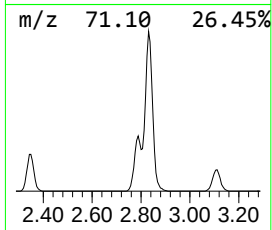
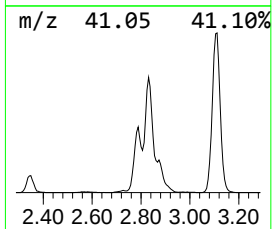
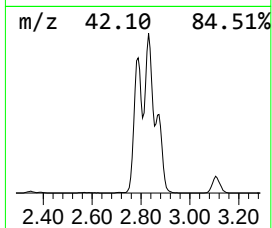
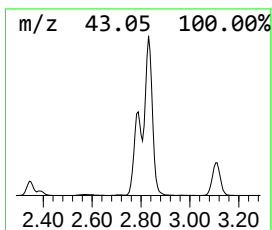
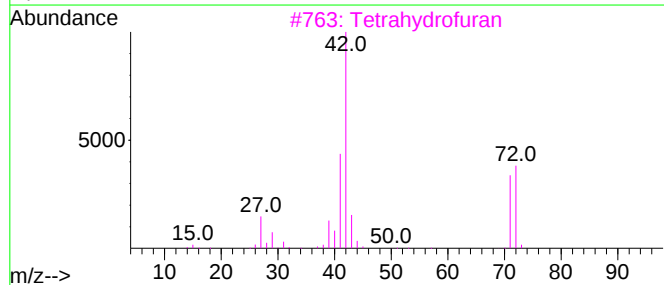
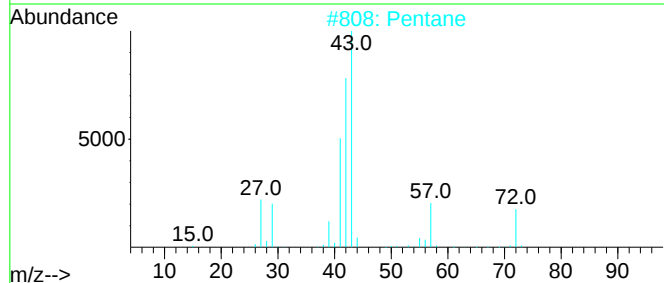
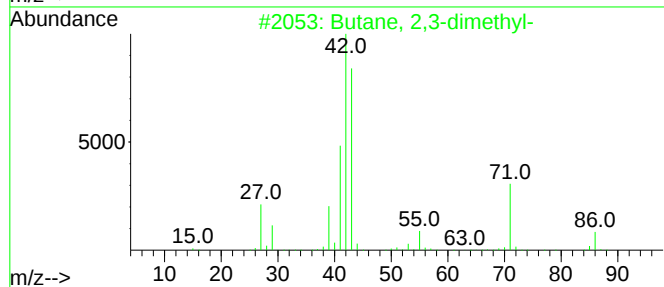
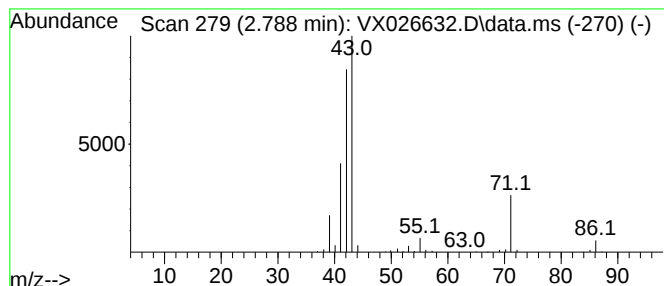
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X011122W.M
Quant Title : SW846 8260

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

Peak Number 2 Butane, 2,3-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.788	30.76 ug/l	248999	Pentafluorobenzene	5.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	90
2			Pentane	72	C5H12	000109-66-0	45
3			Tetrahydrofuran	72	C4H8O	000109-99-9	38
4			Pentane, 2-methyl-	86	C6H14	000107-83-5	37
5			Propanoyl chloride, 2-methyl-	106	C4H7ClO	000079-30-1	27



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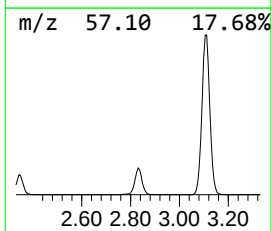
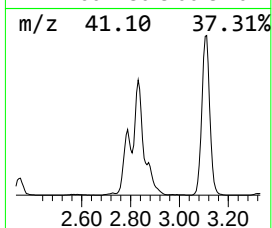
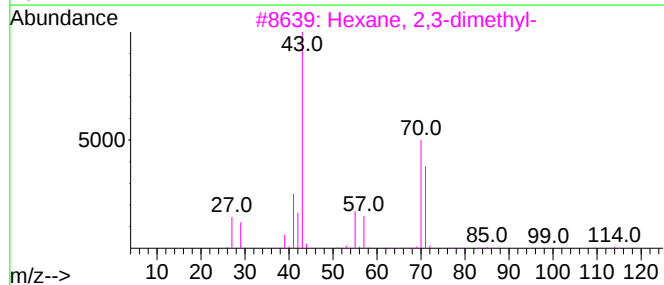
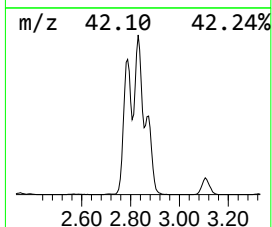
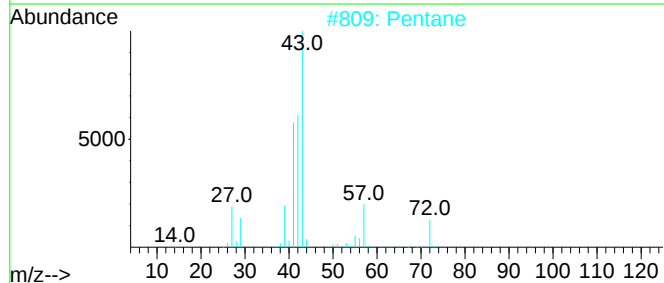
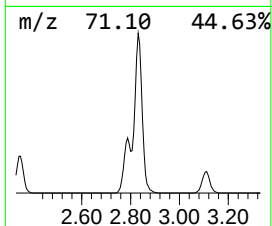
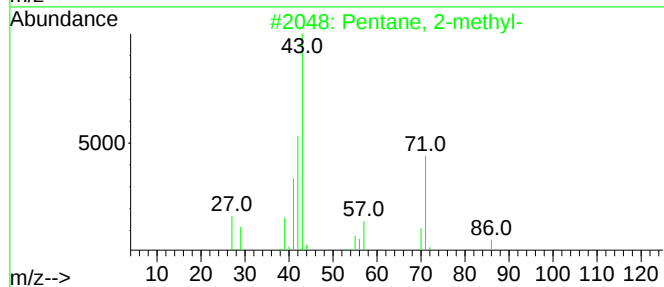
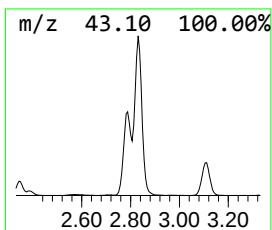
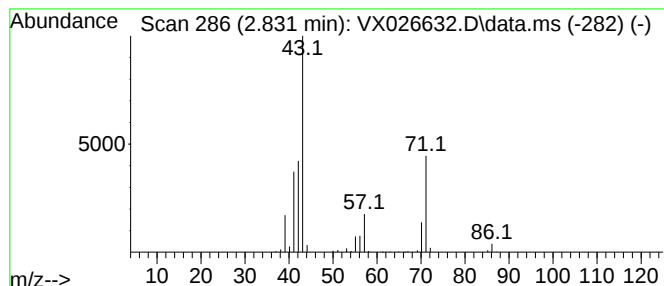
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X011122W.M
Quant Title : SW846 8260

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

Peak Number 3 Pentane, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.831	39.51 ug/l	319823	Pentafluorobenzene	5.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2-methyl-	86	C6H14	000107-83-5	91
2			Pentane	72	C5H12	000109-66-0	38
3			Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	38
4			Pentane, 3,3-dimethyl-	100	C7H16	000562-49-2	23
5			Propanoic acid, 2-methyl-, 2-pro...	128	C7H12O2	015727-77-2	23



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MW-10R

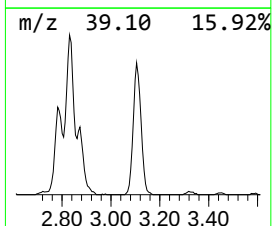
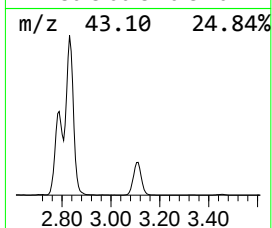
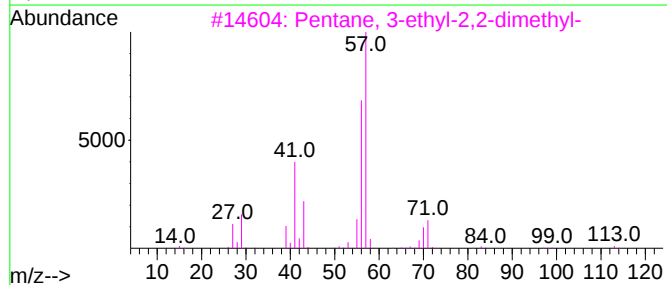
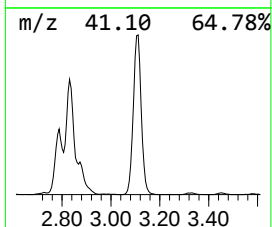
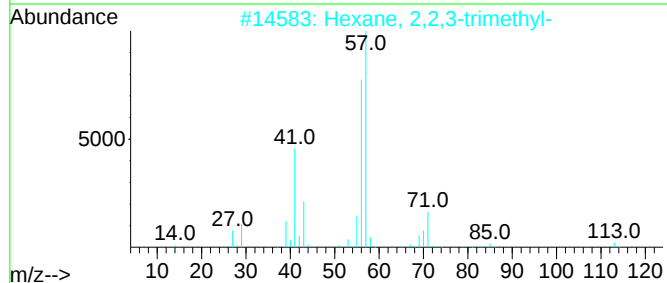
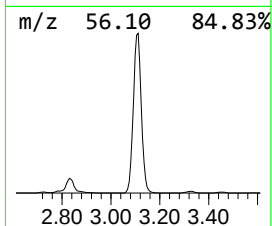
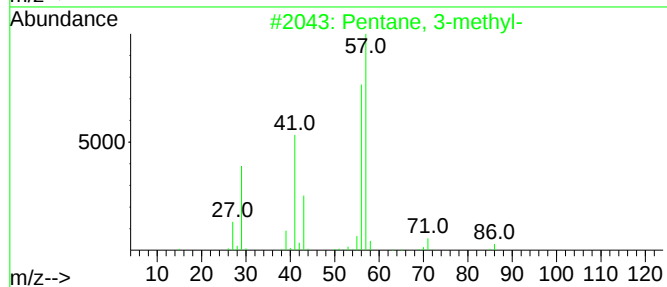
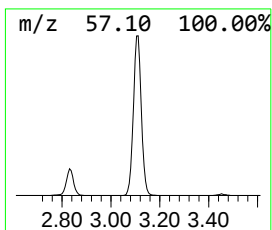
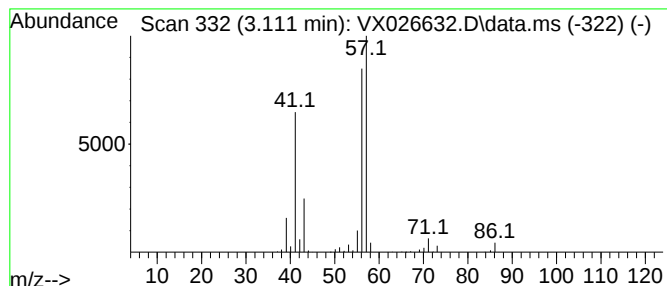
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Quant Title : SW846 8260

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

Peak Number 4 Pentane, 3-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.111	58.81 ug/l	476075	Pentafluorobenzene	5.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 3-methyl-	86	C6H14	000096-14-0	91
2			Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	78
3			Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	64
4			Pentane, 2,2,3-trimethyl-	114	C8H18	000564-02-3	50
5			2-Ethyl-oxetane	86	C5H10O	1010386-40-2	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX012722\
Data File : VX026632.D
Acq On : 27 Jan 2022 18:12
Operator : JC/MD
Sample : N1296-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-10R

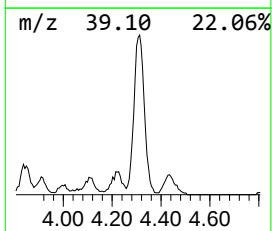
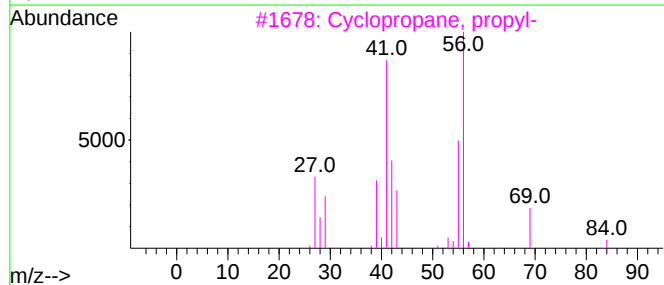
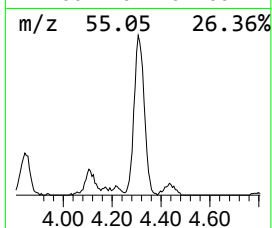
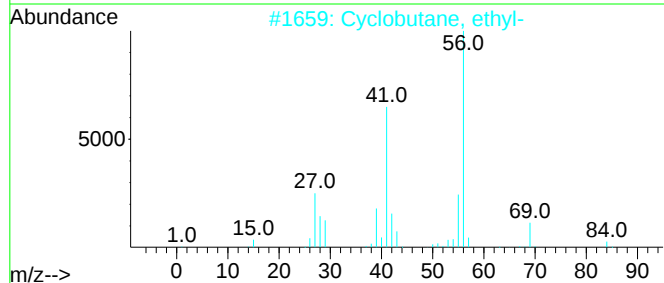
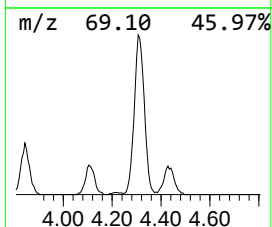
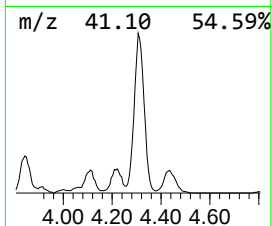
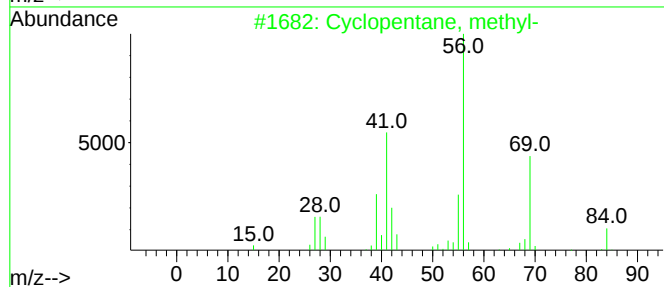
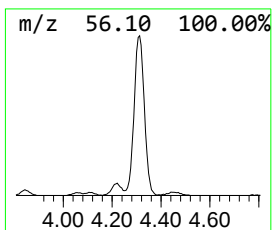
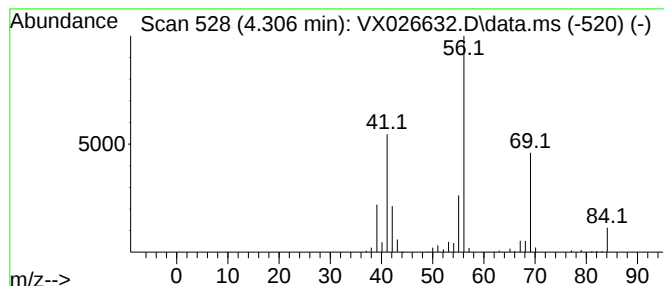
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Quant Title : SW846 8260

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

Peak Number 5 Cyclopentane, methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.306	27.69 ug/l	224132	Pentafluorobenzene	5.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, methyl-	84	C6H12	000096-37-7	91
2			Cyclobutane, ethyl-	84	C6H12	004806-61-5	72
3			Cyclopropane, propyl-	84	C6H12	002415-72-7	72
4			1-Hexene	84	C6H12	000592-41-6	52
5			1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX012722\
Data File : VX026632.D
Acq On : 27 Jan 2022 18:12
Operator : JC/MD
Sample : N1296-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-10R

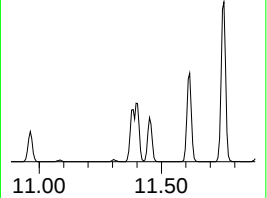
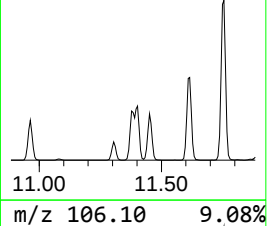
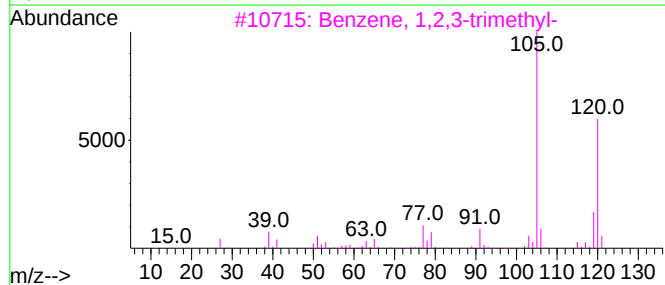
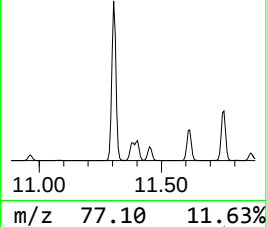
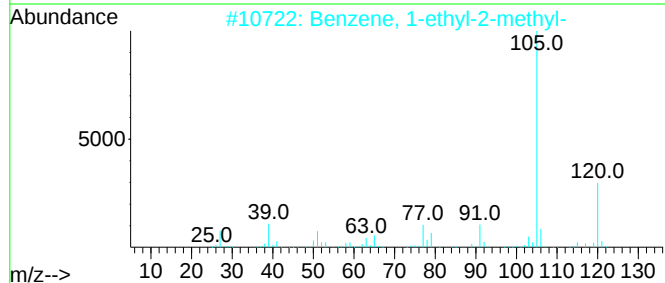
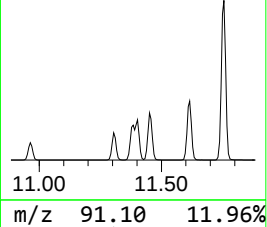
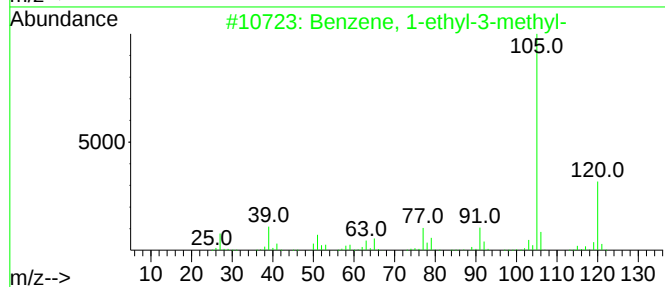
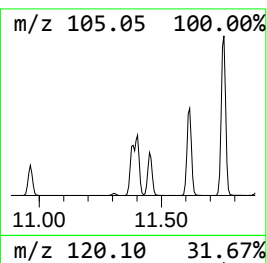
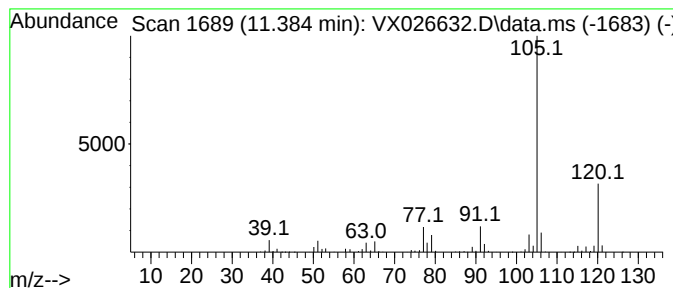
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1-ethyl-3-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.384	44.29 ug/l	449999	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	95
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
4			Mesitylene	120	C9H12	000108-67-8	91
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX012722\
Data File : VX026632.D
Acq On : 27 Jan 2022 18:12
Operator : JC/MD
Sample : N1296-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-10R

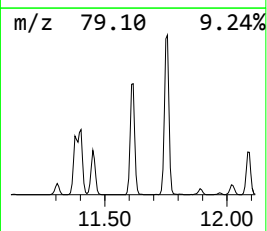
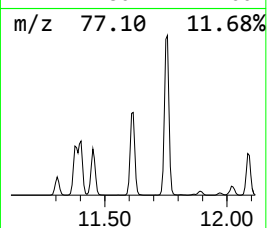
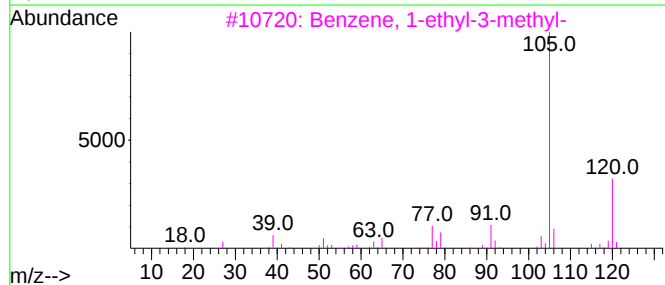
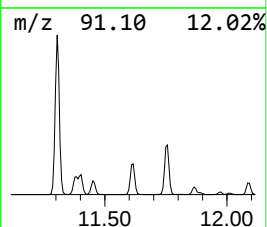
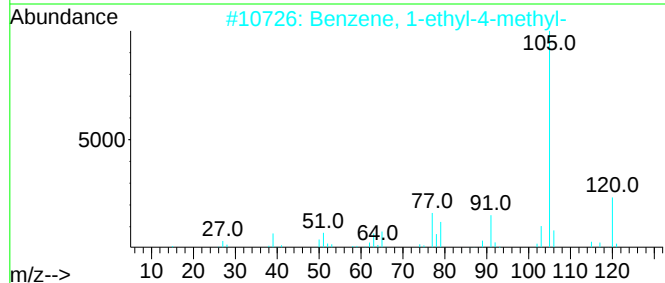
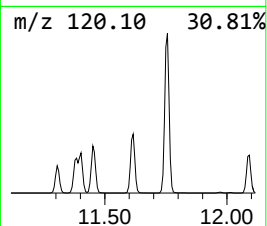
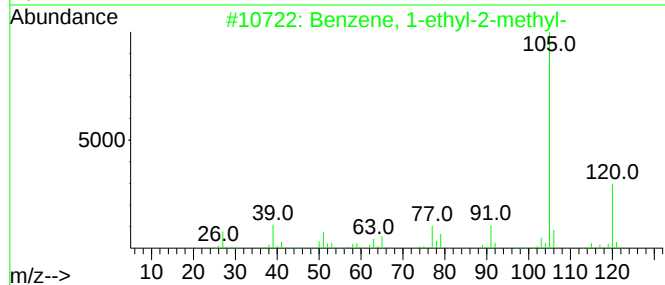
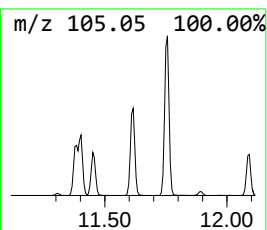
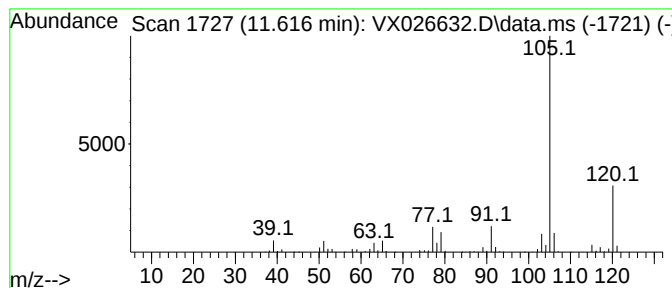
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.616	71.42 ug/l	725638	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	94
2			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	94
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX012722\
Data File : VX026632.D
Acq On : 27 Jan 2022 18:12
Operator : JC/MD
Sample : N1296-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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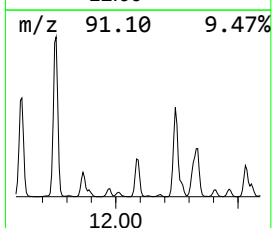
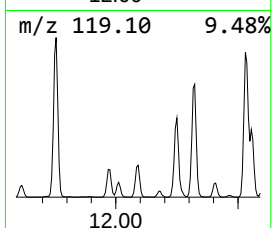
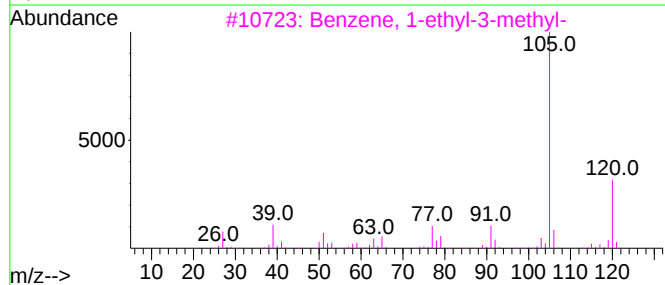
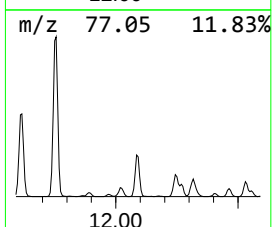
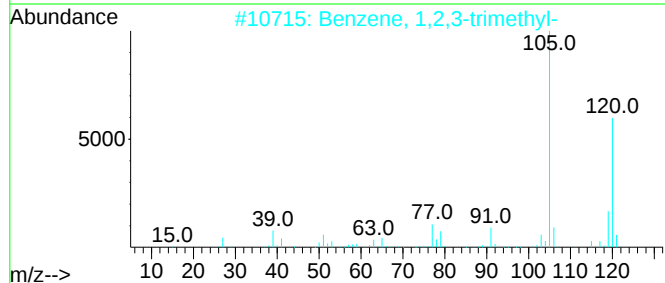
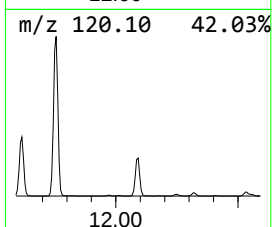
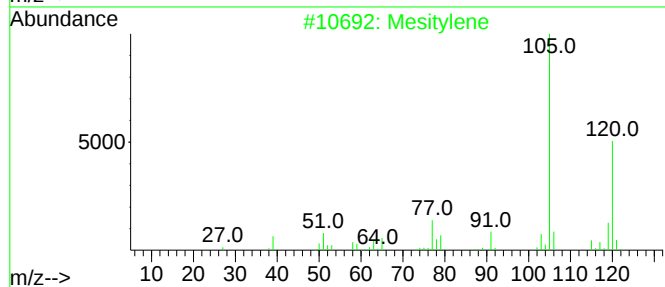
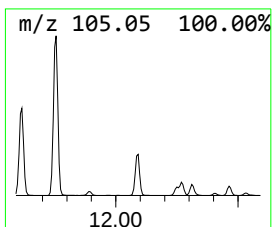
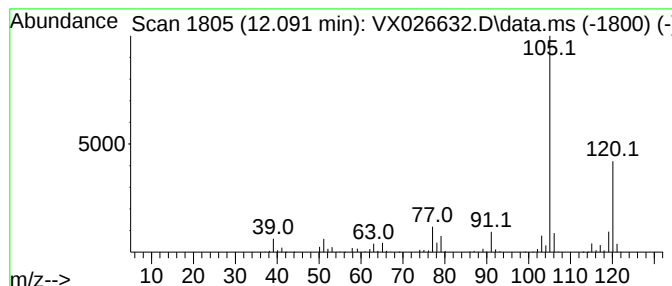
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1,2,3-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.091	36.25 ug/l	368330	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Mesitylene	120	C9H12	000108-67-8	97
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX012722\
Data File : VX026632.D
Acq On : 27 Jan 2022 18:12
Operator : JC/MD
Sample : N1296-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-10R

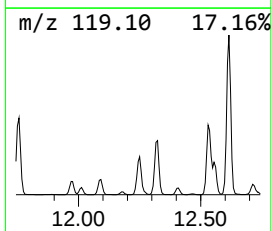
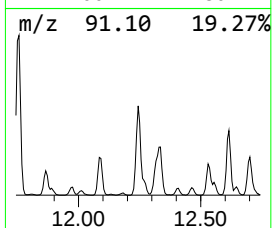
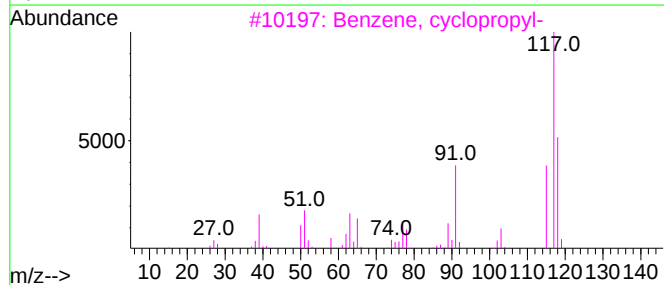
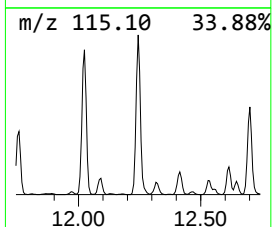
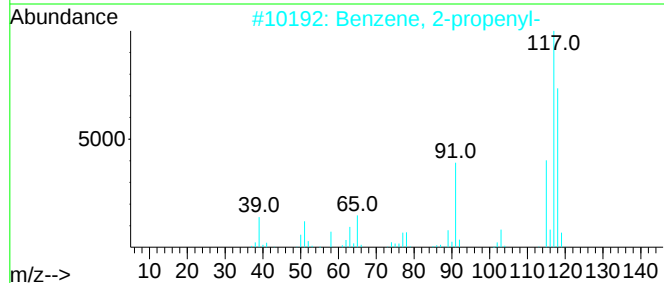
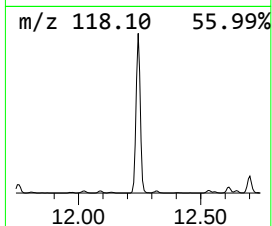
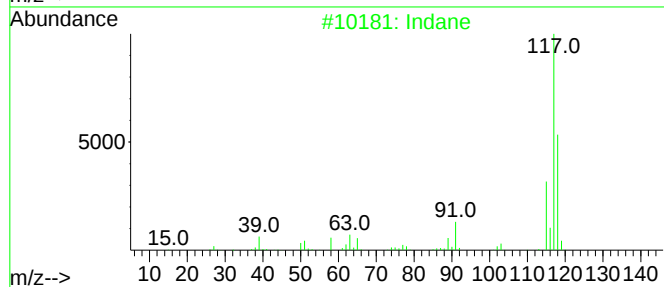
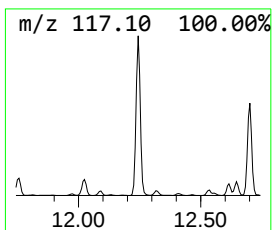
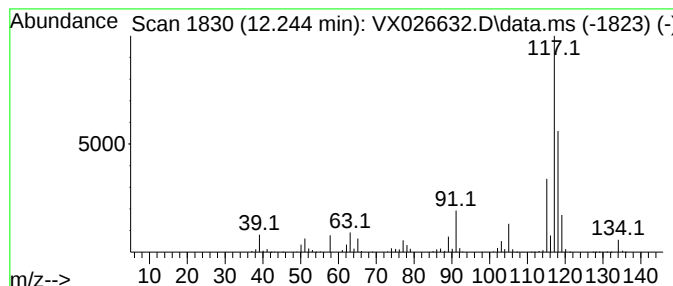
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Quant Title : SW846 8260

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

Peak Number 9 Indane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	65.58 ug/l	666291	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	76
2			Benzene, 2-propenyl-	118	C9H10	000300-57-2	76
3			Benzene, cyclopropyl-	118	C9H10	000873-49-4	76
4			Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	64
5			Deltacyclene	118	C9H10	007785-10-6	52



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Data File : VX026632.D
Acq On : 27 Jan 2022 18:12
Operator : JC/MD
Sample : N1296-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-10R

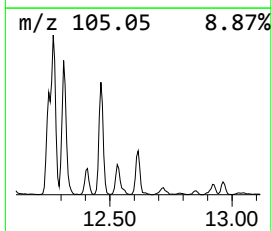
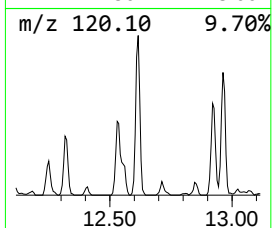
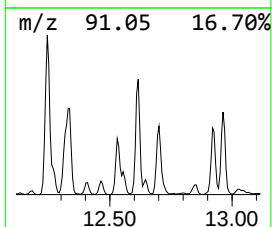
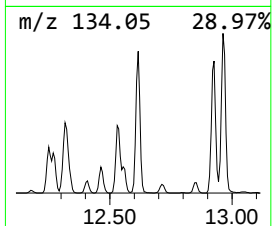
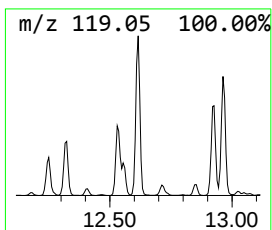
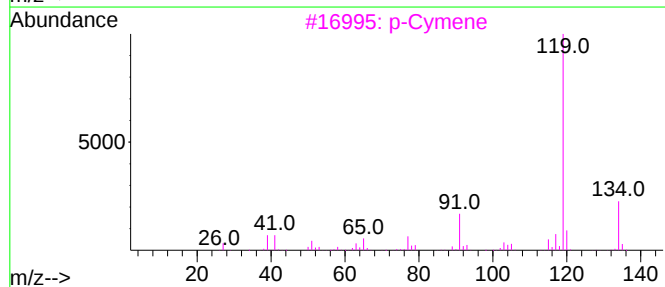
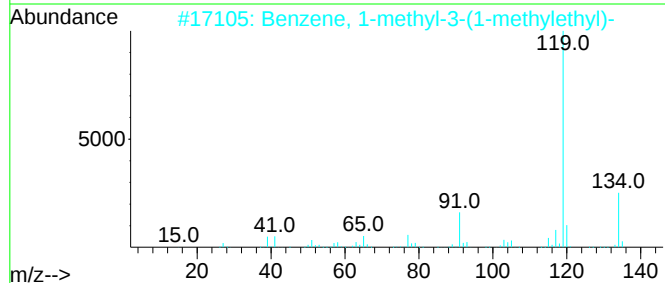
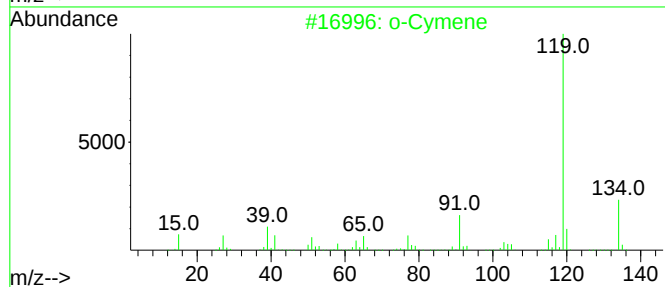
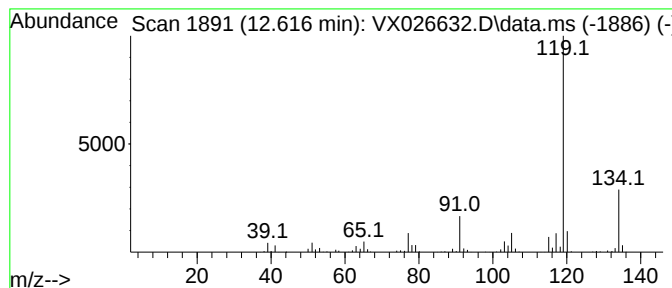
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Quant Title : SW846 8260

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

Peak Number 10 o-Cymene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.616	32.56 ug/l	330759	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Cymene	134	C10H14	000527-84-4	97
2			Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	97
3			p-Cymene	134	C10H14	000099-87-6	97
4			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX012722\
Data File : VX026632.D
Acq On : 27 Jan 2022 18:12
Operator : JC/MD
Sample : N1296-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-10R

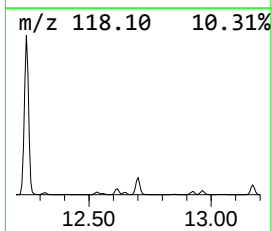
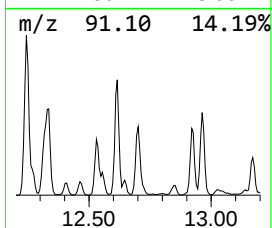
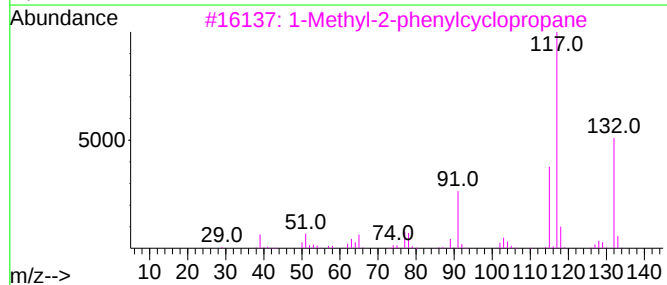
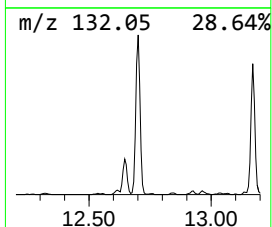
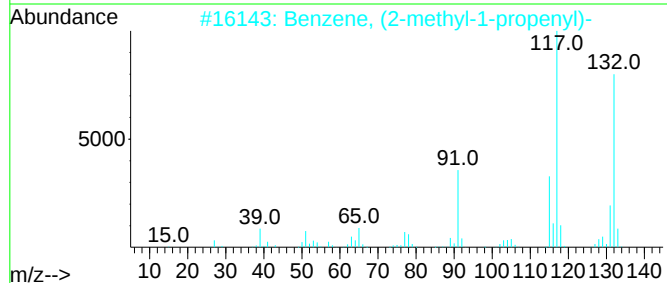
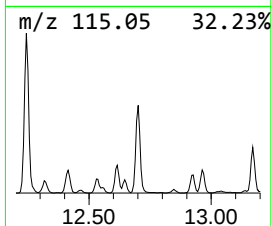
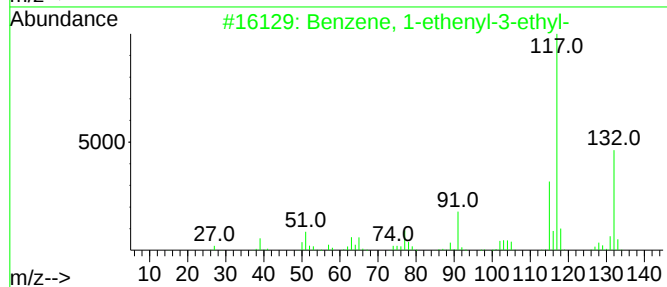
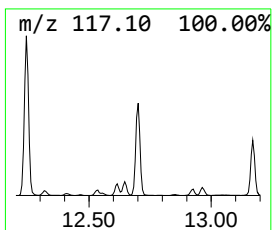
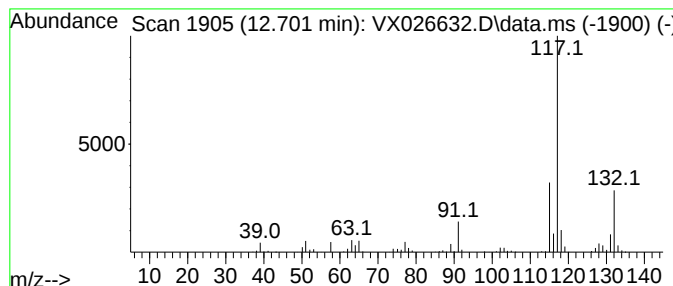
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Quant Title : SW846 8260

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

Peak Number 11 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.701	27.87 ug/l	283150	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			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	90
4			Indan, 1-methyl-	132	C10H12	000767-58-8	90
5			1-Phenyl-1-butene	132	C10H12	000824-90-8	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX012722\
Data File : VX026632.D
Acq On : 27 Jan 2022 18:12
Operator : JC/MD
Sample : N1296-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 22 Sample Multiplier: 1

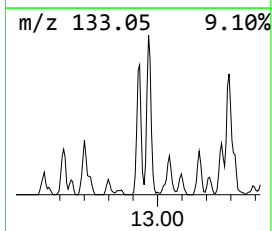
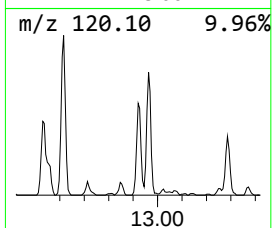
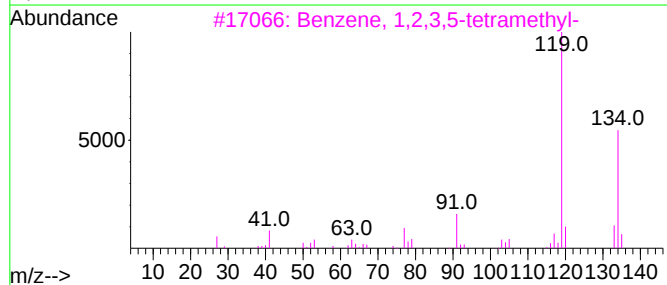
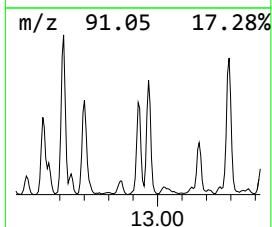
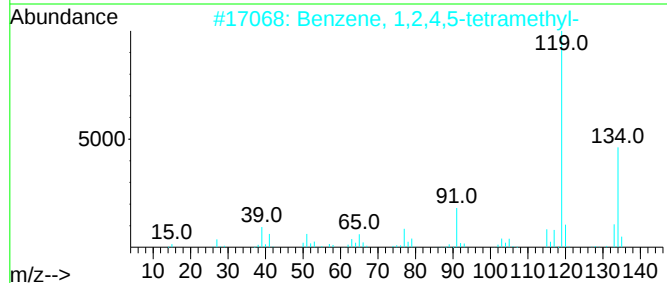
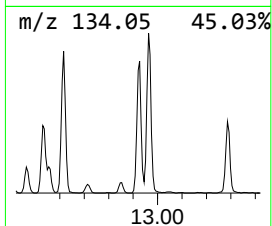
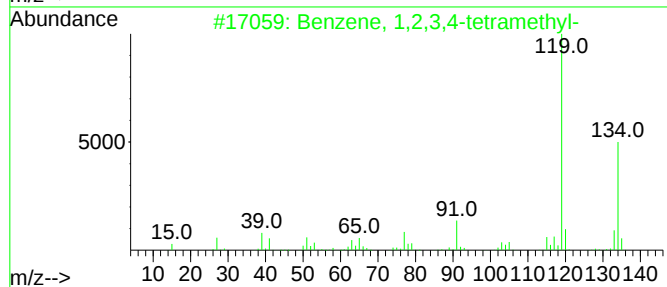
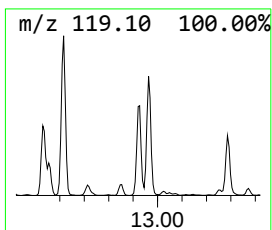
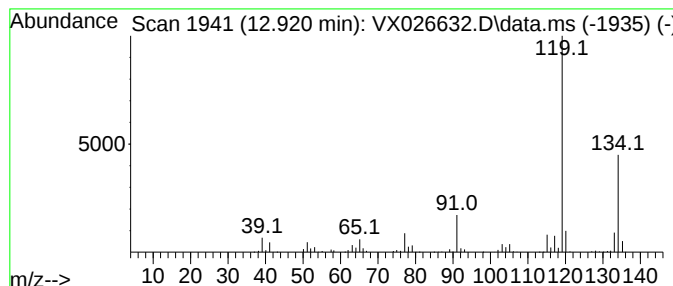
Instrument :
MSVOA_X
ClientSampleId :
MW-10R

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X011122W.M
Quant Title : SW846 8260

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

Peak Number 12 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD			R.T.
12.920	21.84 ug/l	221912	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	96
3	Benzene, 1,2,3,5-tetramethyl-		134	C10H14	000527-53-7	96
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\VX012722\
Data File : VX026632.D
Acq On : 27 Jan 2022 18:12
Operator : JC/MD
Sample : N1296-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 22 Sample Multiplier: 1

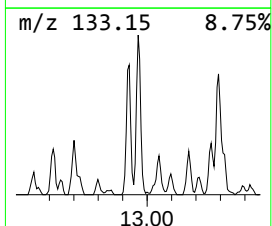
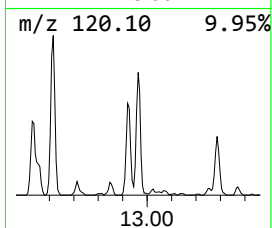
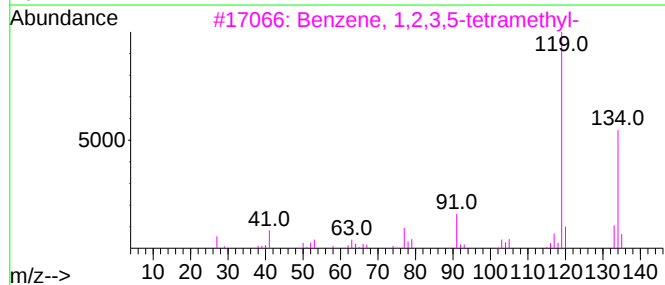
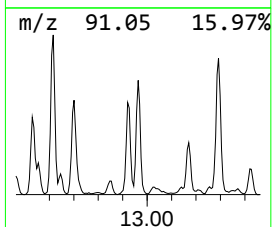
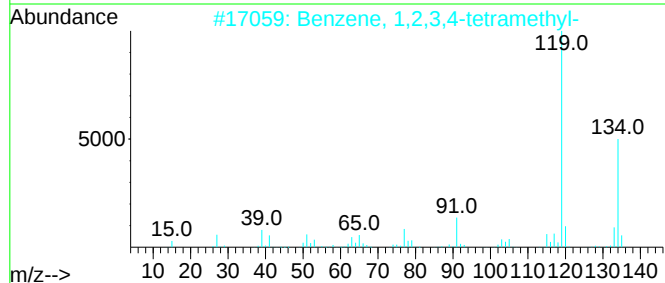
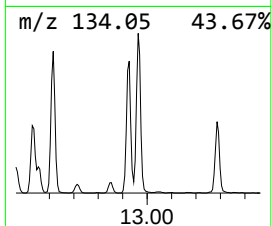
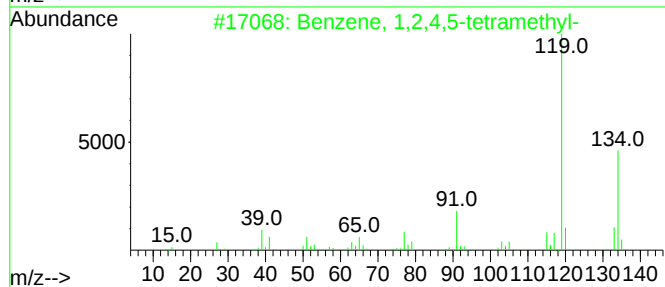
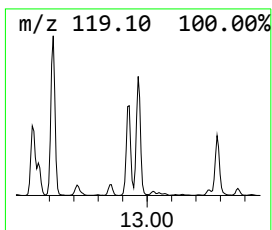
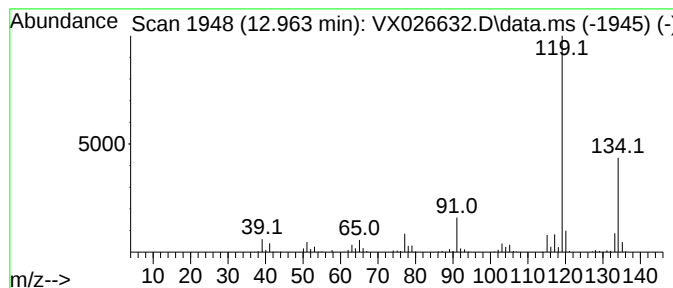
Instrument :
MSVOA_X
ClientSampleId :
MW-10R

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X011122W.M
Quant Title : SW846 8260

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

Peak Number 13 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD		R.T.	
12.963	25.46 ug/l	258710	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	96
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\VX012722\
Data File : VX026632.D
Acq On : 27 Jan 2022 18:12
Operator : JC/MD
Sample : N1296-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-10R

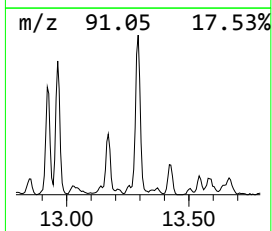
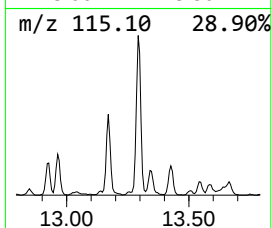
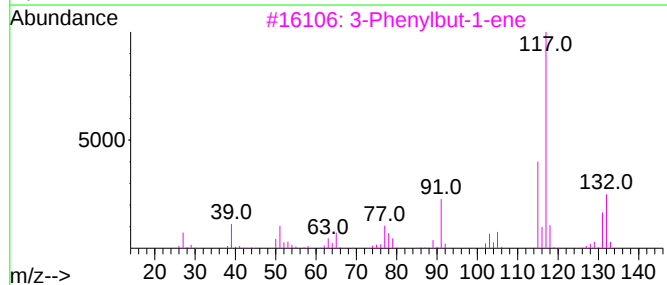
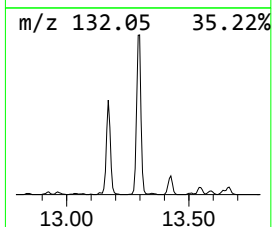
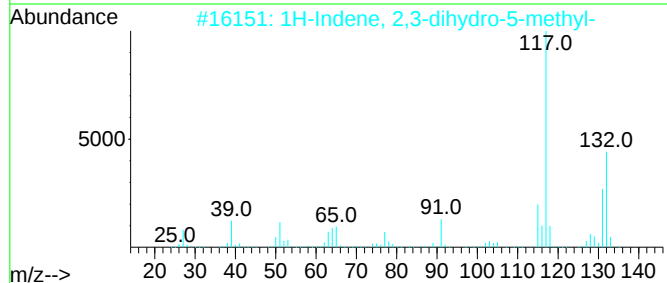
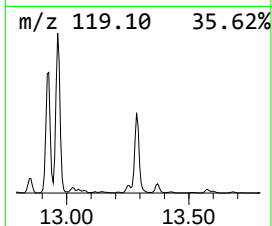
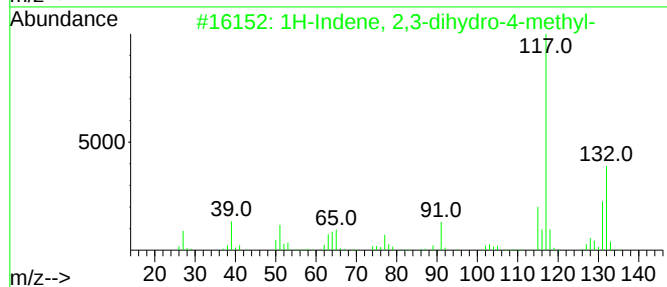
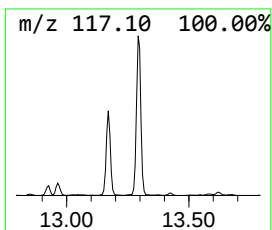
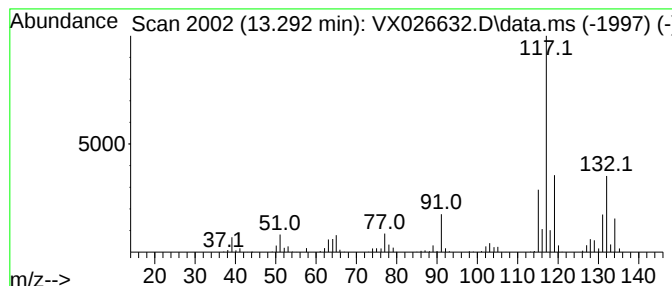
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Quant Title : SW846 8260

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

Peak Number 14 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	42.18 ug/l	428510	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	93
2		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	87
3		3-Phenylbut-1-ene	132	C10H12	000934-10-1	87
4		Benzene, 2-butenyl-	132	C10H12	001560-06-1	87
5		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX012722\
Data File : VX026632.D
Acq On : 27 Jan 2022 18:12
Operator : JC/MD
Sample : N1296-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 22 Sample Multiplier: 1

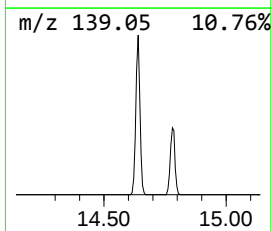
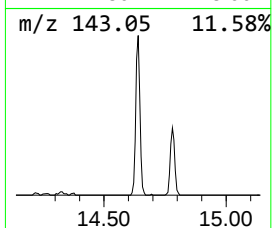
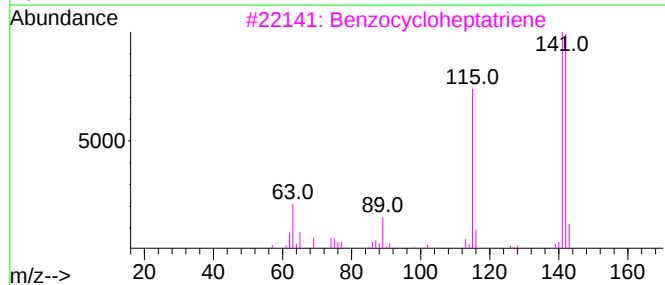
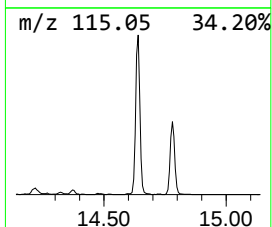
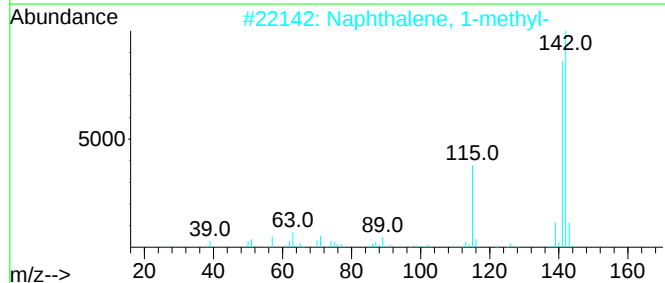
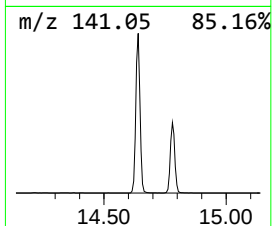
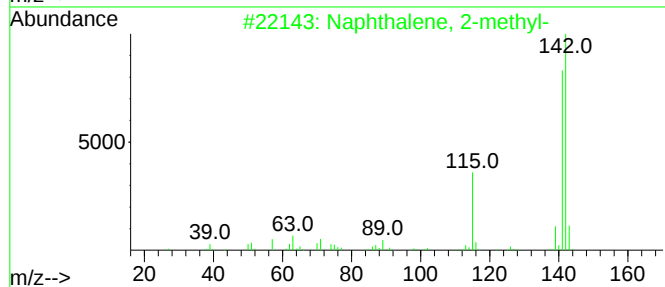
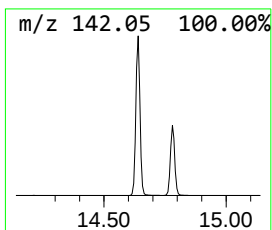
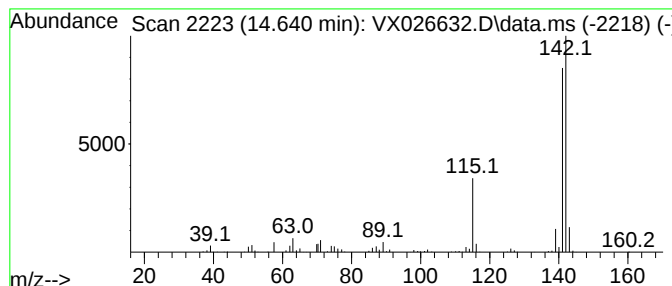
Instrument :
MSVOA_X
ClientSampleId :
MW-10R

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X011122W.M
Quant Title : SW846 8260

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

Peak Number 15 Naphthalene, 2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.640	30.21 ug/l	306967	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	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX012722\
Data File : VX026632.D
Acq On : 27 Jan 2022 18:12
Operator : JC/MD
Sample : N1296-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-10R

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X011122W.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
Butane, 2-methyl-	1.758	76.6	ug/l	619967	1	5.556	404749	50.0
Butane, 2,3-dim...	2.788	30.8	ug/l	248999	1	5.556	404749	50.0
Pentane, 2-methyl-	2.831	39.5	ug/l	319823	1	5.556	404749	50.0
Pentane, 3-methyl-	3.111	58.8	ug/l	476075	1	5.556	404749	50.0
Cyclopentane, m...	4.306	27.7	ug/l	224132	1	5.556	404749	50.0
Benzene, 1-ethy...	11.384	44.3	ug/l	449999	4	12.024	507973	50.0
Benzene, 1-ethy...	11.616	71.4	ug/l	725638	4	12.024	507973	50.0
Benzene, 1,2,3-...	12.091	36.3	ug/l	368330	4	12.024	507973	50.0
Indane	12.244	65.6	ug/l	666291	4	12.024	507973	50.0
o-Cymene	12.616	32.6	ug/l	330759	4	12.024	507973	50.0
Benzene, 1-ethe...	12.701	27.9	ug/l	283150	4	12.024	507973	50.0
Benzene, 1,2,3,...	12.920	21.8	ug/l	221912	4	12.024	507973	50.0
Benzene, 1,2,4,...	12.963	25.5	ug/l	258710	4	12.024	507973	50.0
1H-Indene, 2,3-...	13.292	42.2	ug/l	428510	4	12.024	507973	50.0
Naphthalene, 2-...	14.640	30.2	ug/l	306967	4	12.024	507973	50.0