

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX032323\
 Data File : VX034708.D
 Acq On : 23 Mar 2023 18:58
 Operator : JC/MD
 Sample : 02055-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-4B

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\82X031023W.M
 Title : SW846 8260

Signal : TIC: VX034708.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.367	43	46	51	rBV	237781	241337	5.56%	0.328%
2	1.752	103	109	125	rVB	3178052	4263574	98.30%	5.802%
3	1.953	137	142	155	rVB	1154986	1651925	38.09%	2.248%
4	2.343	199	206	211	rBV	118344	235543	5.43%	0.321%
5	2.715	258	267	271	rBV4	46965	119107	2.75%	0.162%
6	2.782	271	278	280	rVV	353194	689069	15.89%	0.938%
7	2.825	280	285	308	rVB	1227199	3062761	70.62%	4.168%
8	3.099	320	330	346	rBV	827616	1865401	43.01%	2.539%
9	3.318	356	366	377	rBV	96653	218403	5.04%	0.297%
10	3.446	377	387	400	rBV	325536	717008	16.53%	0.976%
11	3.581	400	409	416	rBV2	42463	110765	2.55%	0.151%
12	3.733	428	434	443	rVB	129800	304082	7.01%	0.414%
13	3.837	443	451	458	rBV2	44272	114528	2.64%	0.156%
14	4.099	489	494	504	rVB	75547	193973	4.47%	0.264%
15	4.215	504	513	518	rVV2	79069	219256	5.06%	0.298%
16	4.306	518	528	542	rVB	995764	2849199	65.69%	3.877%
17	5.135	645	664	666	rBV6	33794	133622	3.08%	0.182%
18	5.385	691	705	711	rBV	199906	608567	14.03%	0.828%
19	5.477	711	720	728	rVV3	529311	2062199	47.55%	2.806%
20	5.550	728	732	739	rVV	349208	1041923	24.02%	1.418%
21	5.617	739	743	765	rVV	210605	701018	16.16%	0.954%
22	5.824	765	777	790	rVV3	363319	1129146	26.03%	1.537%
23	5.952	790	798	812	rVV	215675	575621	13.27%	0.783%
24	6.159	815	832	841	rVV2	216341	760846	17.54%	1.035%
25	6.257	841	848	856	rVV2	210869	604849	13.95%	0.823%
26	6.354	856	864	877	rVB2	256563	736515	16.98%	1.002%
27	6.610	892	906	914	rBV3	81473	216042	4.98%	0.294%
28	6.763	922	931	939	rBV	508766	1320881	30.45%	1.798%
29	7.171	990	998	1005	rBV2	47039	111500	2.57%	0.152%
30	7.379	1019	1032	1043	rBV	1178063	2895129	66.75%	3.940%
31	7.501	1043	1052	1057	rVV3	49397	128325	2.96%	0.175%
32	7.561	1057	1062	1071	rVV3	57727	138995	3.20%	0.189%
33	7.659	1071	1078	1088	rVV	172436	340039	7.84%	0.463%
34	7.775	1089	1097	1104	rVB2	106070	212516	4.90%	0.289%
35	7.958	1121	1127	1144	rVB4	79858	249104	5.74%	0.339%
36	8.104	1144	1151	1156	rBV3	63159	140520	3.24%	0.191%

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37	8.311	1182	1185	1192	rVB4	77115	142318	3.28%	0.194%
38	8.439	1201	1206	1209	rBV	71425	116207	2.68%	0.158%
39	8.506	1212	1217	1226	rVB3	135914	269172	6.21%	0.366%
40	8.598	1226	1232	1235	rBV2	231358	410330	9.46%	0.558%
41	8.647	1235	1240	1248	rVV	1068947	1771133	40.84%	2.410%
42	8.775	1256	1261	1265	rVV	99838	166582	3.84%	0.227%
43	8.848	1270	1273	1279	rVB4	71437	110856	2.56%	0.151%
44	8.976	1289	1294	1306	rVB	184783	342335	7.89%	0.466%
45	9.098	1306	1314	1320	rBV	131697	249559	5.75%	0.340%
46	9.159	1320	1324	1329	rBV2	78635	122922	2.83%	0.167%
47	9.512	1376	1382	1386	rVV3	111551	233863	5.39%	0.318%
48	9.561	1386	1390	1395	rVB	179540	250557	5.78%	0.341%
49	9.616	1395	1399	1403	rBV	89163	135309	3.12%	0.184%
50	10.055	1466	1471	1476	rBV	1023821	1362060	31.40%	1.854%
51	10.195	1488	1494	1501	rVV	709730	961362	22.17%	1.308%
52	10.299	1502	1511	1517	rVV	2298741	3206030	73.92%	4.363%
53	10.585	1551	1558	1563	rBV3	64372	136511	3.15%	0.186%
54	10.640	1563	1567	1580	rVB	137624	213089	4.91%	0.290%
55	10.780	1581	1590	1595	rBV2	75887	126777	2.92%	0.173%
56	10.860	1598	1603	1615	rVB4	89633	159818	3.68%	0.217%
57	10.963	1615	1620	1624	rBV	215417	286186	6.60%	0.389%
58	11.079	1634	1639	1644	rBV	876950	1102480	25.42%	1.500%
59	11.305	1668	1676	1681	rBV	484887	594479	13.71%	0.809%
60	11.378	1683	1688	1689	rVV	1117265	1219154	28.11%	1.659%
61	11.396	1689	1691	1696	rVV	1339645	1690887	38.99%	2.301%
62	11.451	1696	1700	1710	rVB	1922395	2278607	52.54%	3.101%
63	11.610	1721	1726	1734	rVB	1307879	1673525	38.59%	2.277%
64	11.750	1744	1749	1758	rVB	3664414	4337165	100.00%	5.902%
65	11.969	1781	1785	1789	rVV	163448	189460	4.37%	0.258%
66	12.018	1789	1793	1799	rVB	1027360	1382064	31.87%	1.881%
67	12.085	1799	1804	1810	rBV	1429441	1717776	39.61%	2.338%
68	12.244	1824	1830	1832	rBV2	933954	1200344	27.68%	1.633%
69	12.268	1832	1834	1838	rVV	613794	702299	16.19%	0.956%
70	12.317	1838	1842	1850	rVV2	1149430	1621739	37.39%	2.207%
71	12.402	1851	1856	1861	rVV2	79566	119666	2.76%	0.163%
72	12.463	1861	1866	1872	rVV	247057	306086	7.06%	0.417%
73	12.530	1872	1877	1879	rVV	461890	569405	13.13%	0.775%
74	12.554	1879	1881	1885	rVV	455975	478763	11.04%	0.652%
75	12.609	1885	1890	1894	rVV	879365	1137646	26.23%	1.548%
76	12.646	1894	1896	1900	rVV	217939	232867	5.37%	0.317%

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77	12.701	1900	1905	1913	rVV2	590259	960479	22.15%	1.307%
78	12.847	1925	1929	1934	rVV2	257716	323579	7.46%	0.440%
79	12.920	1934	1941	1944	rVV	533537	661257	15.25%	0.900%
80	12.963	1944	1948	1954	rVB	787068	915348	21.10%	1.246%
81	13.024	1954	1958	1960	rBV	120691	173291	4.00%	0.236%
82	13.042	1960	1961	1964	rVV	125200	150852	3.48%	0.205%
83	13.067	1964	1965	1969	rVV	103241	108689	2.51%	0.148%
84	13.134	1973	1976	1978	rVV	83319	113324	2.61%	0.154%
85	13.170	1978	1982	1986	rVV	531111	728009	16.79%	0.991%
86	13.256	1992	1996	1998	rVV2	116032	168429	3.88%	0.229%
87	13.292	1998	2002	2008	rVV2	1123551	1539687	35.50%	2.095%
88	13.371	2012	2015	2019	rVV	102184	116379	2.68%	0.158%
89	13.420	2019	2023	2032	rVB2	153870	216859	5.00%	0.295%
90	13.506	2032	2037	2040	rBV2	121828	168857	3.89%	0.230%
91	13.542	2040	2043	2046	rVV	223779	298701	6.89%	0.406%
92	13.585	2046	2050	2056	rVV3	262665	507340	11.70%	0.690%
93	13.664	2056	2063	2069	rVB2	227134	477576	11.01%	0.650%
94	13.774	2075	2081	2091	rBV	222906	311748	7.19%	0.424%
95	13.926	2099	2106	2110	rBV2	104980	163688	3.77%	0.223%
96	14.073	2126	2130	2135	rBV	153810	192320	4.43%	0.262%
97	14.213	2148	2153	2162	rBV	132240	236587	5.45%	0.322%
98	14.371	2175	2179	2184	rVB	90170	109546	2.53%	0.149%
99	14.639	2218	2223	2233	rVB	216576	314092	7.24%	0.427%
100	14.780	2241	2246	2255	rVB	176763	236620	5.46%	0.322%

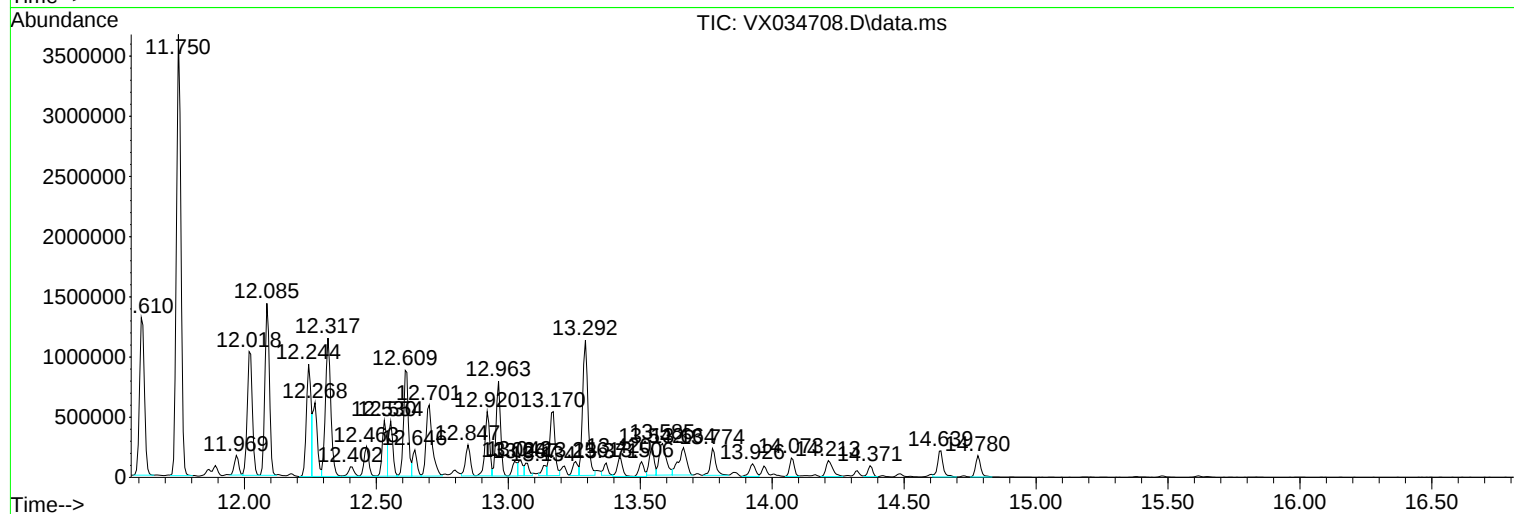
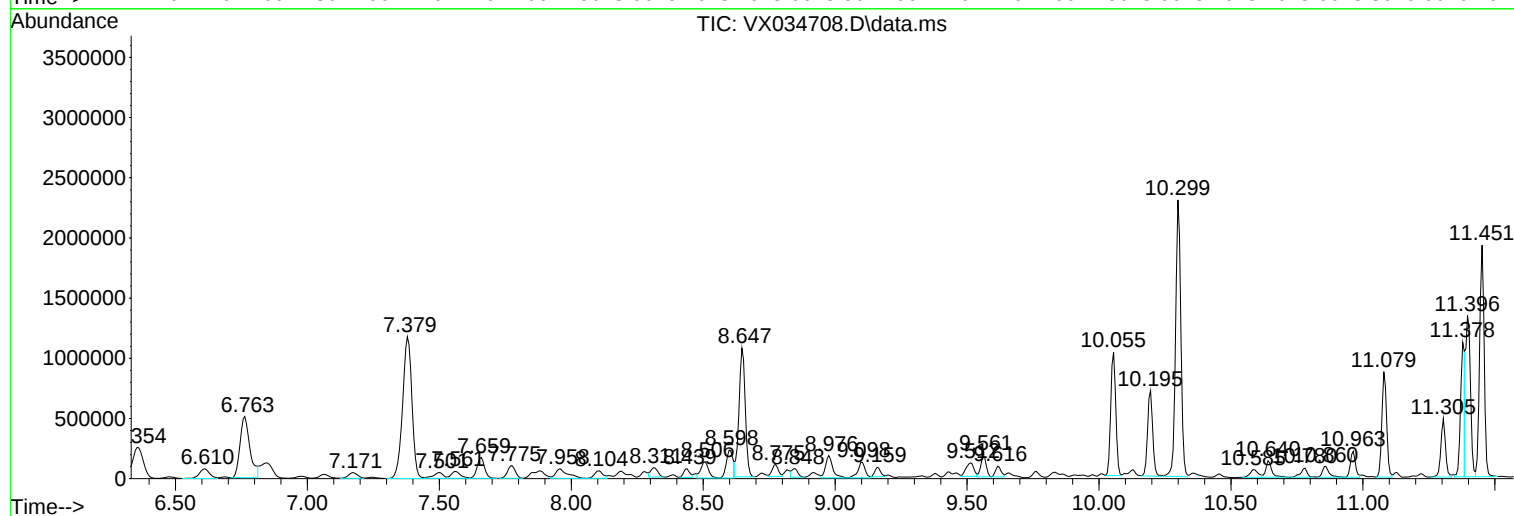
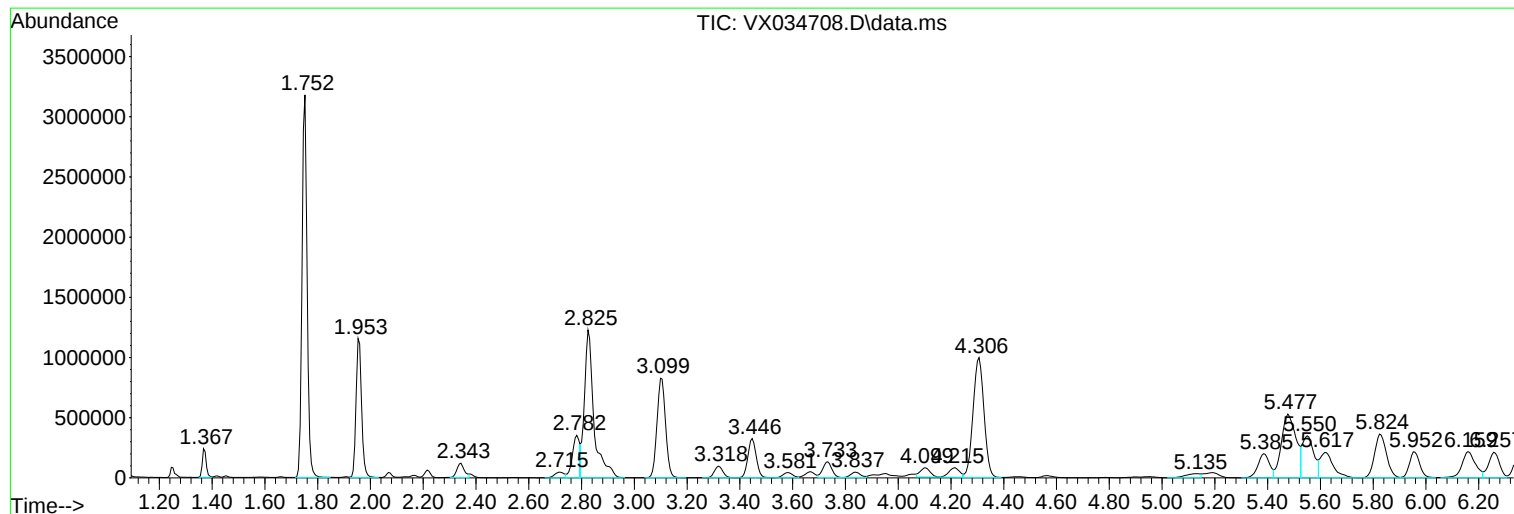
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X031023W.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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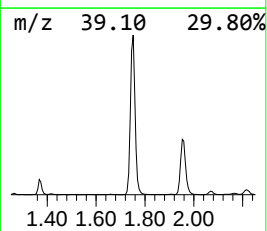
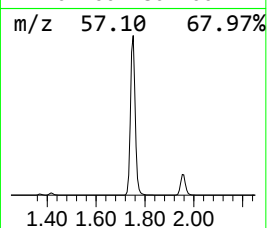
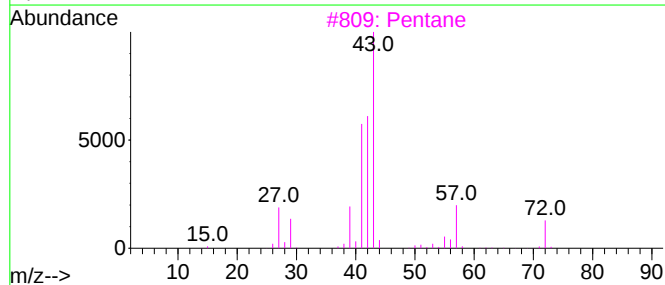
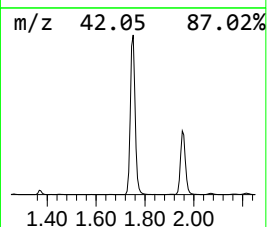
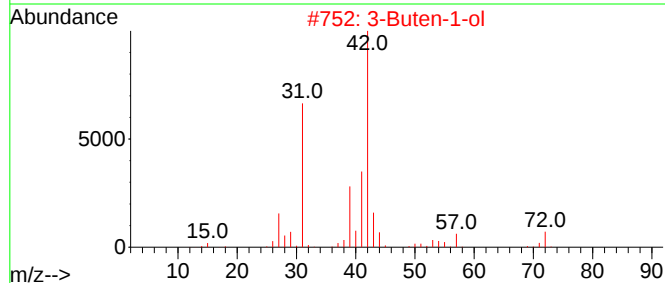
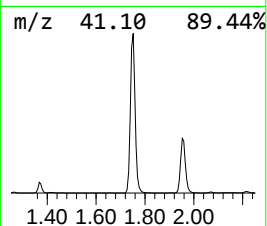
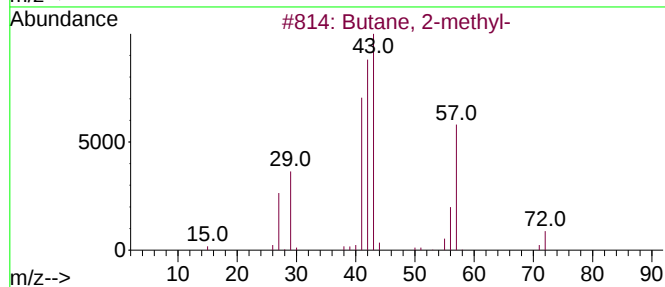
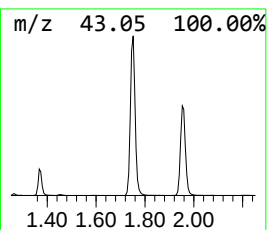
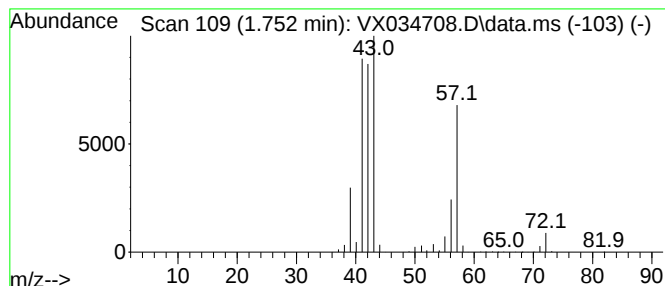
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X031023W.M
Quant Title : SW846 8260

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.
1.752	204.60 ug/l	4263570	Pentafluorobenzene	5.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methyl-	72	C5H12	000078-78-4	91
2			3-Buten-1-ol	72	C4H8O	000627-27-0	47
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	10



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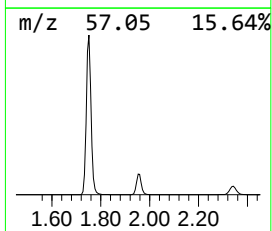
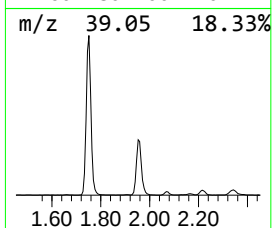
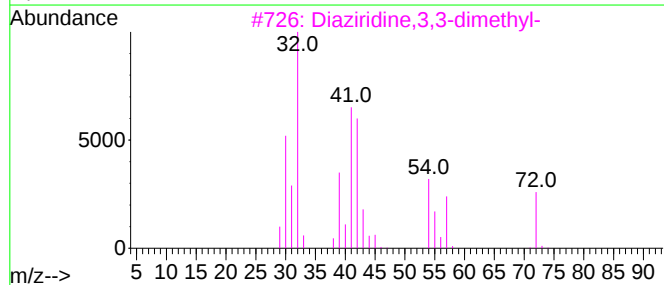
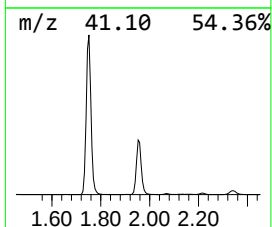
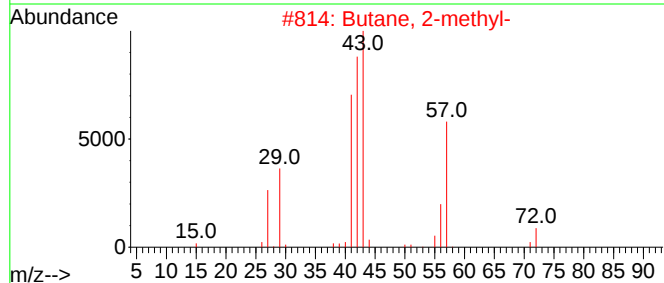
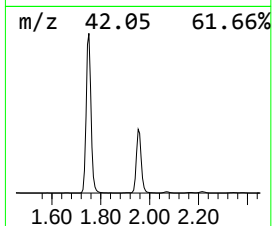
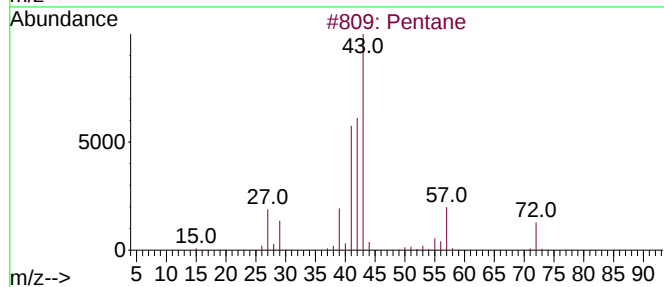
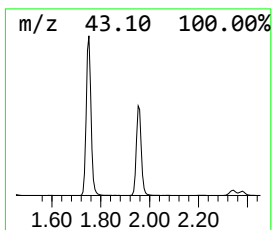
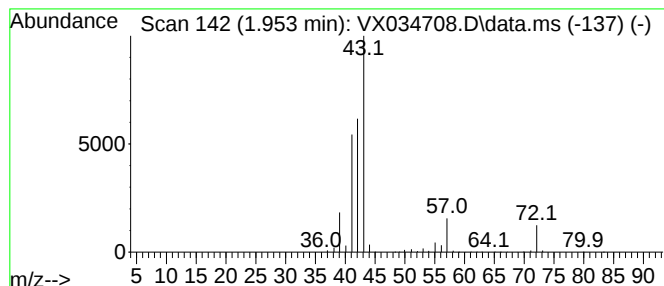
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Pentane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.953	79.27 ug/l	1651930	Pentafluorobenzene	5.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane	72	C5H12	000109-66-0	91
2			Butane, 2-methyl-	72	C5H12	000078-78-4	50
3			Diaziridine,3,3-dimethyl-	72	C3H8N2	004901-76-2	38
4			2-Propanone, hydrazone	72	C3H8N2	005281-20-9	9
5			Oxirane, ethyl-	72	C4H8O	000106-88-7	9



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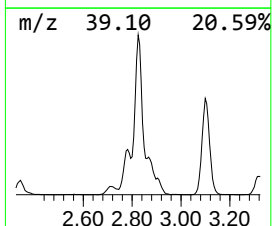
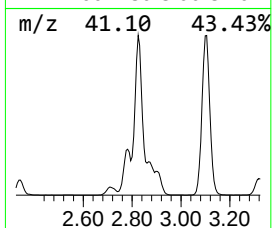
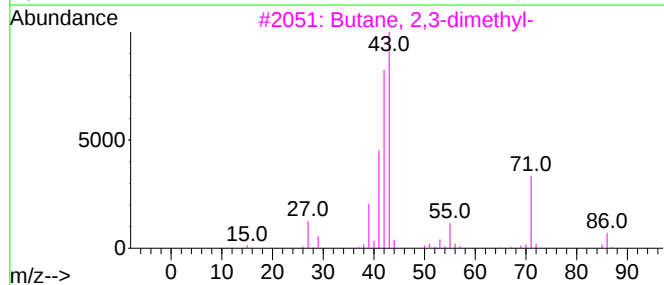
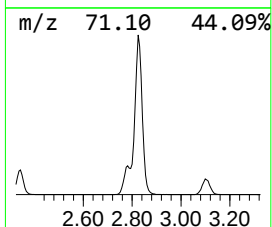
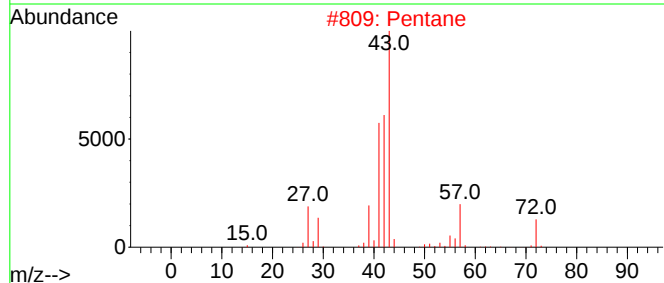
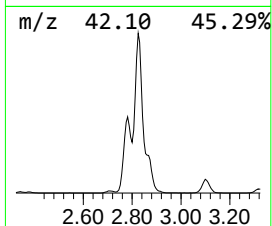
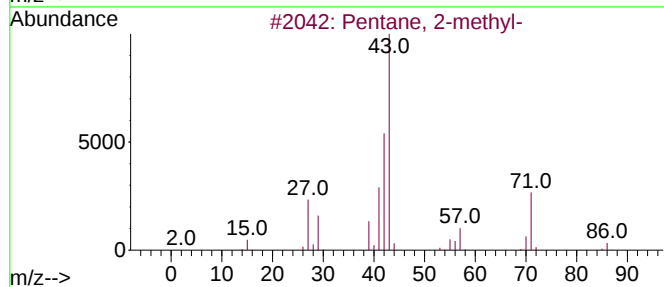
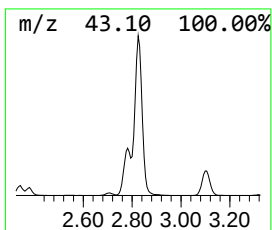
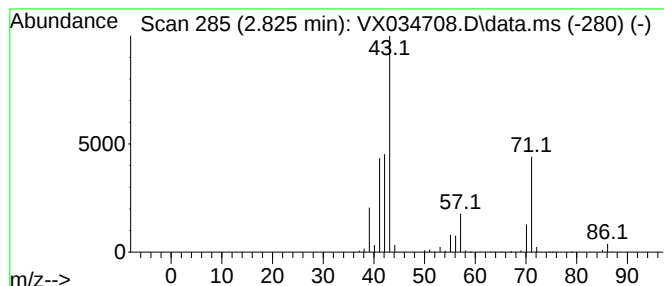
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X031023W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.825	146.98 ug/l	3062760	Pentafluorobenzene	5.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2-methyl-	86	C6H14	000107-83-5	90
2			Pentane	72	C5H12	000109-66-0	49
3			Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	46
4			Pentane, 3,3-dimethyl-	100	C7H16	000562-49-2	38
5			1-Butanol, 2,3-dimethyl-	102	C6H14O	019550-30-2	36



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX032323\
Data File : VX034708.D
Acq On : 23 Mar 2023 18:58
Operator : JC/MD
Sample : 02055-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-4B

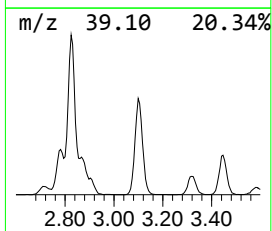
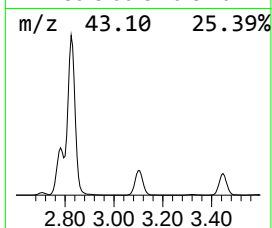
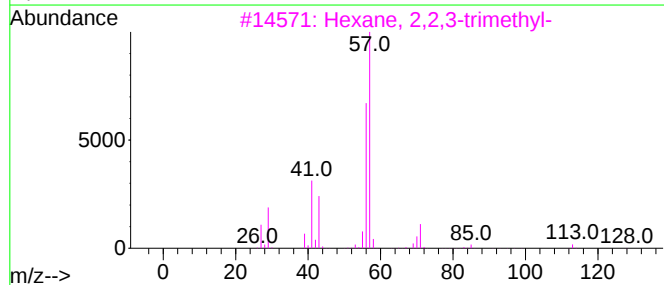
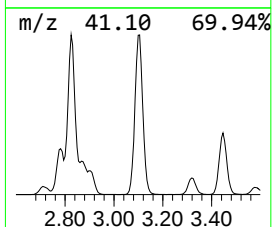
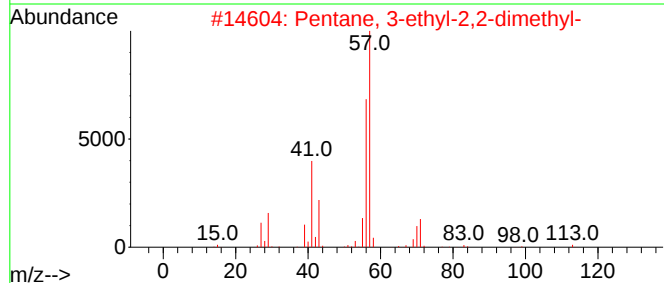
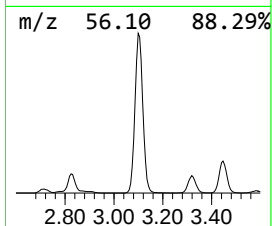
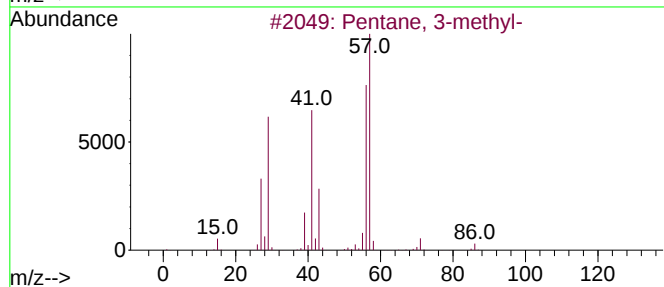
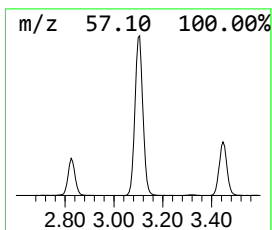
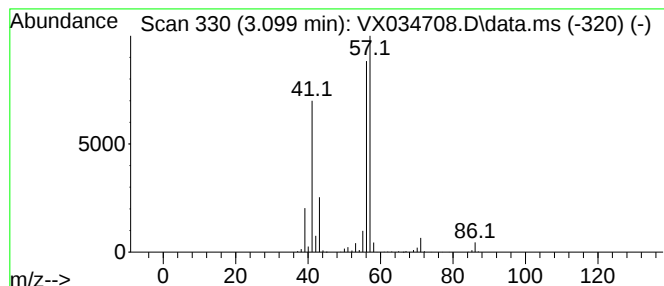
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Quant Title : SW846 8260

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.
3.099	89.52 ug/l	1865400	Pentafluorobenzene	5.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 3-methyl-	86	C6H14	000096-14-0	91
2			Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	43
3			Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	43
4			Sulphuric acid dibutyl ester	210	C8H18O4S	000625-22-9	40
5			Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX032323\
Data File : VX034708.D
Acq On : 23 Mar 2023 18:58
Operator : JC/MD
Sample : 02055-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-4B

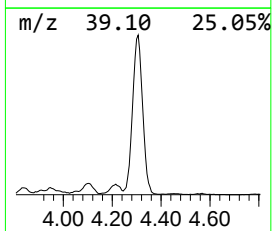
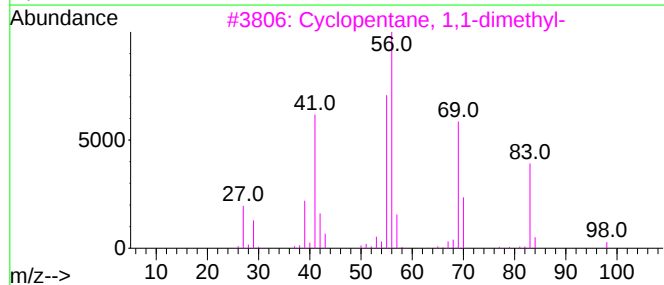
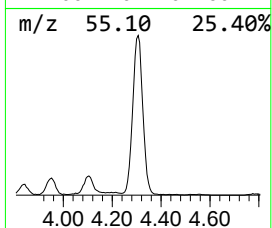
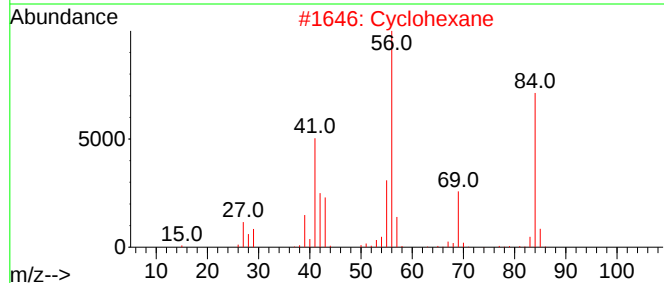
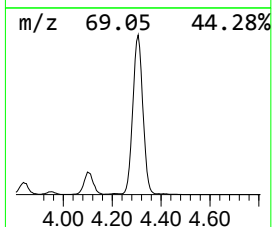
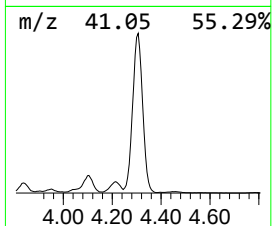
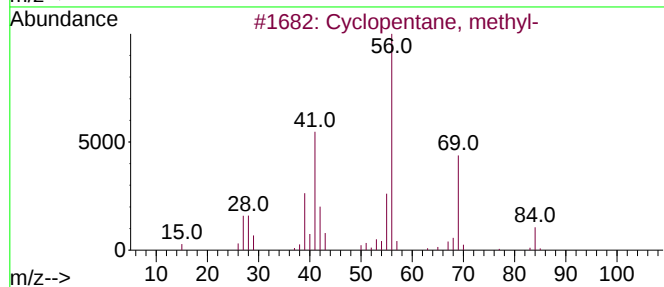
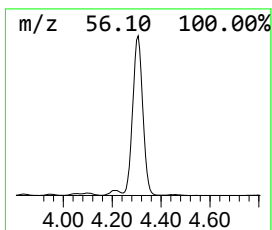
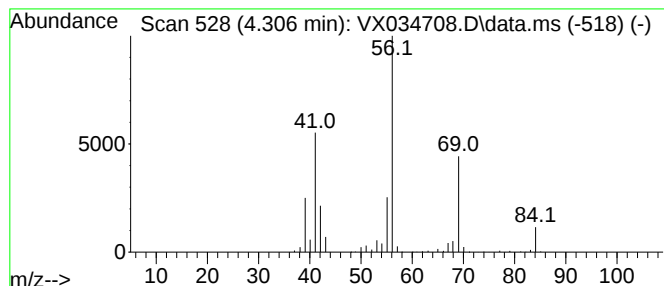
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Quant Title : SW846 8260

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

Peak Number 5 Cyclopentane, methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.306	136.73 ug/l	2849200	Pentafluorobenzene	5.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, methyl-	84	C6H12	000096-37-7	91
2			Cyclohexane	84	C6H12	000110-82-7	83
3			Cyclopentane, 1,1-dimethyl-	98	C7H14	001638-26-2	78
4			Cyclobutane, ethyl-	84	C6H12	004806-61-5	72
5			1-Pentene, 2-methyl-	84	C6H12	000763-29-1	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX032323\
Data File : VX034708.D
Acq On : 23 Mar 2023 18:58
Operator : JC/MD
Sample : 02055-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-4B

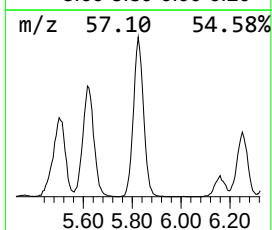
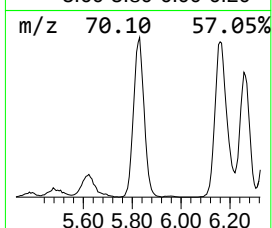
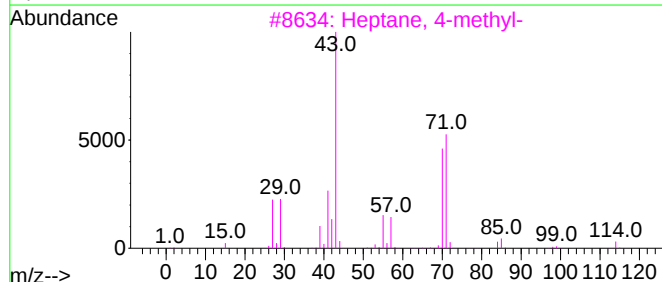
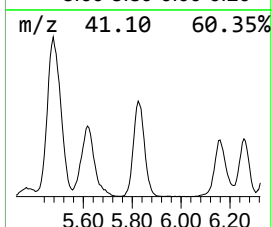
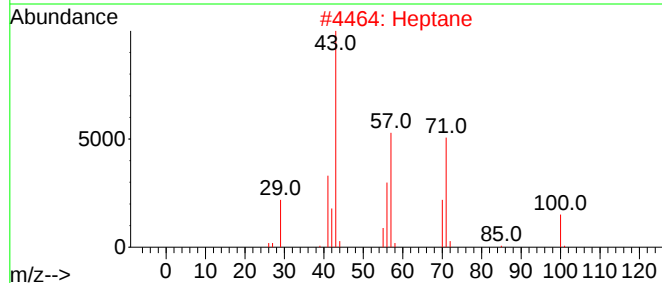
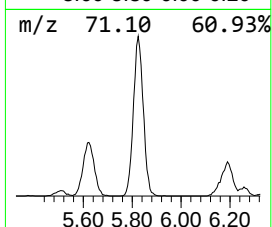
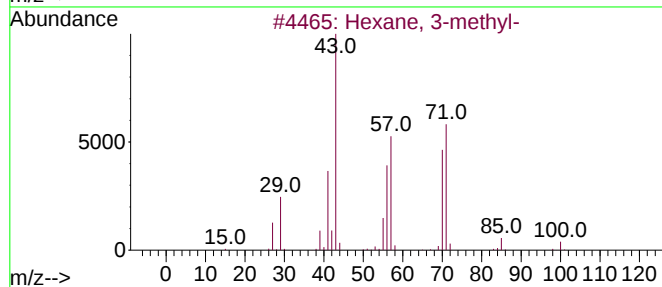
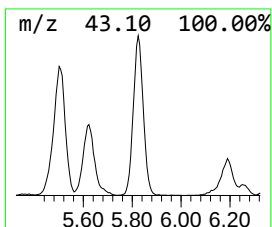
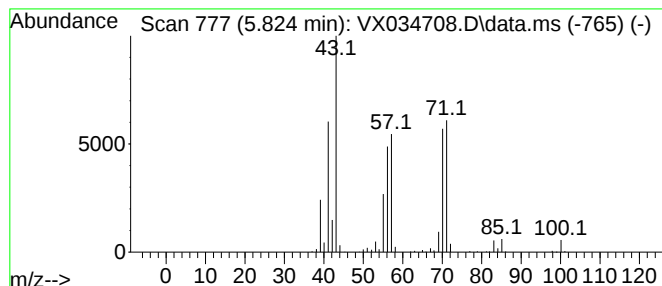
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Quant Title : SW846 8260

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

Peak Number 6 Hexane, 3-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.824	54.19 ug/l	1129150	Pentafluorobenzene	5.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 3-methyl-	100	C7H16	000589-34-4	91
2			Heptane	100	C7H16	000142-82-5	58
3			Heptane, 4-methyl-	114	C8H18	000589-53-7	53
4			Decane, 3,3,4-trimethyl-	184	C13H28	049622-18-6	53
5			Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX032323\
Data File : VX034708.D
Acq On : 23 Mar 2023 18:58
Operator : JC/MD
Sample : 02055-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-4B

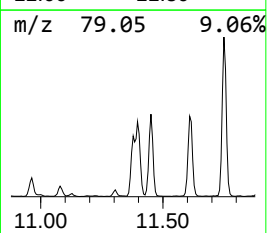
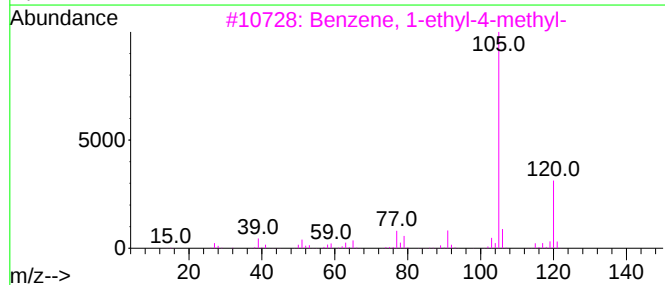
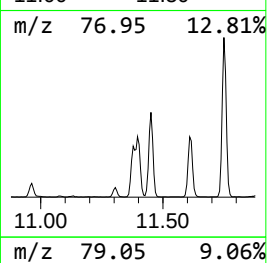
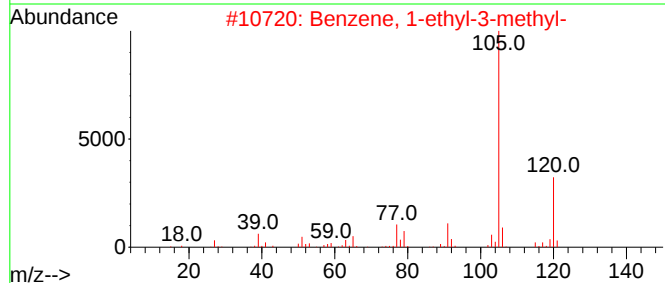
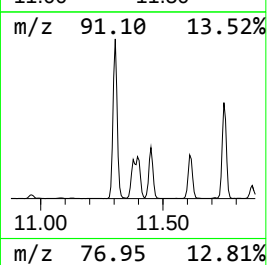
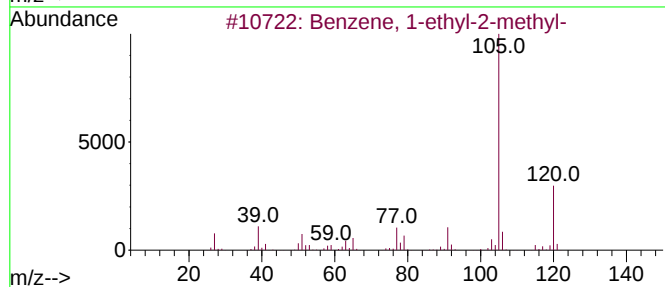
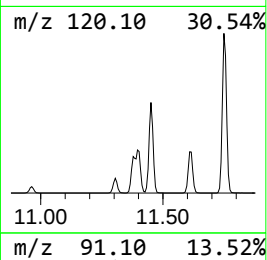
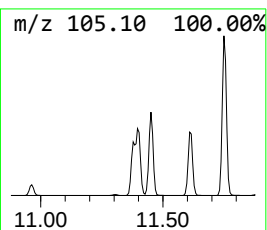
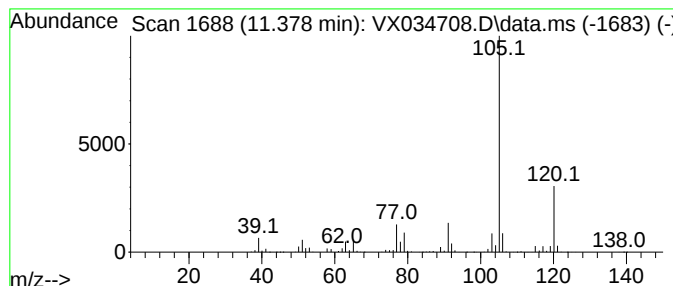
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1-ethyl-2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.378	44.11 ug/l	1219150	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX032323\
Data File : VX034708.D
Acq On : 23 Mar 2023 18:58
Operator : JC/MD
Sample : 02055-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-4B

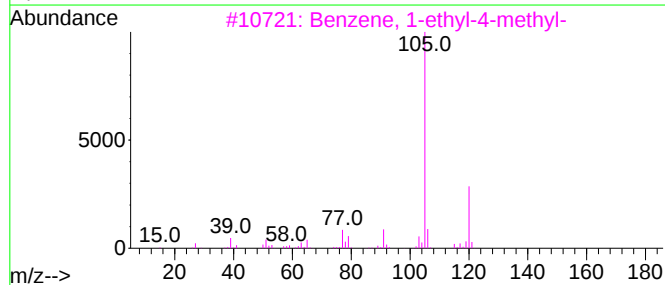
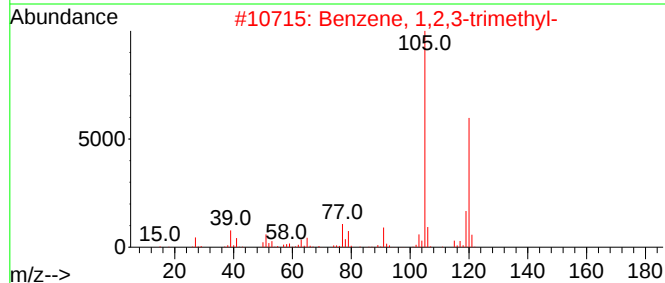
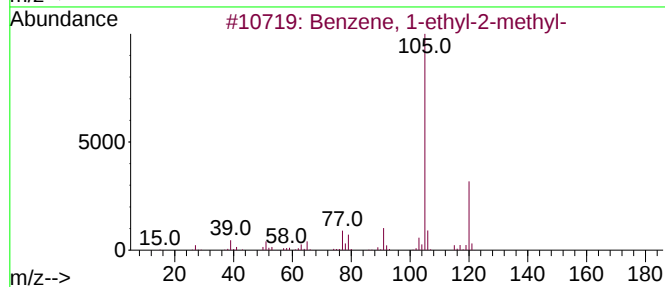
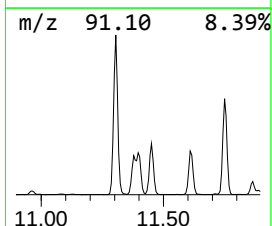
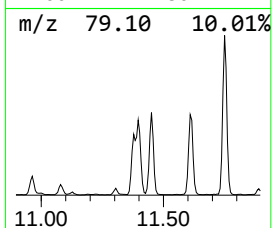
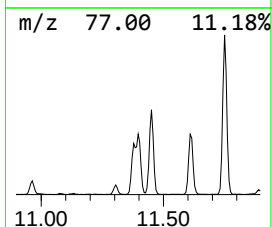
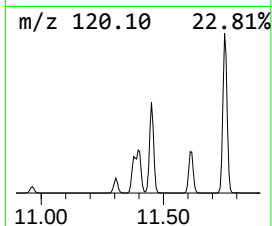
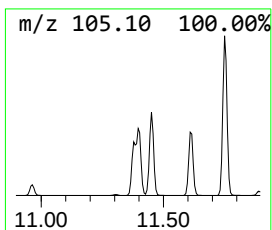
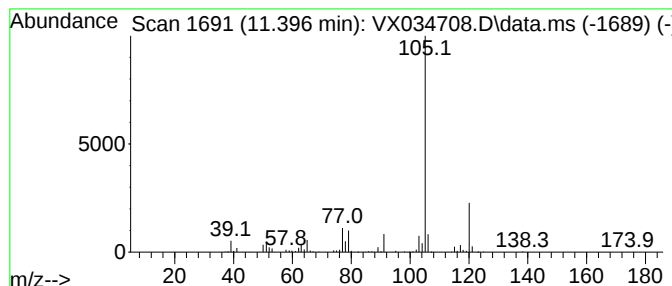
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1-ethyl-4-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.396	61.17 ug/l	1690890	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	91
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4			Mesitylene	120	C9H12	000108-67-8	91
5			Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX032323\
Data File : VX034708.D
Acq On : 23 Mar 2023 18:58
Operator : JC/MD
Sample : 02055-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-4B

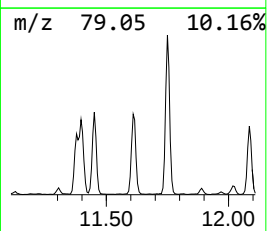
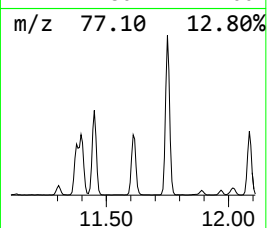
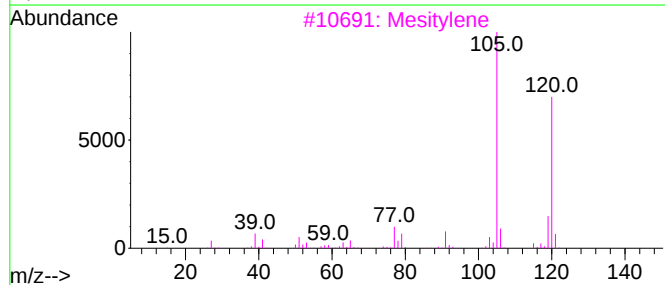
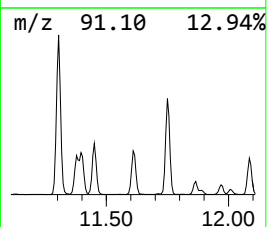
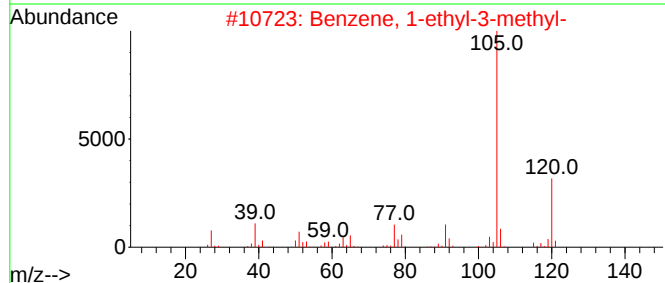
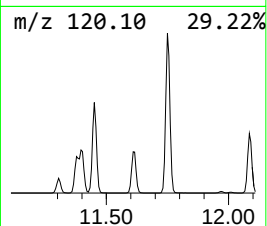
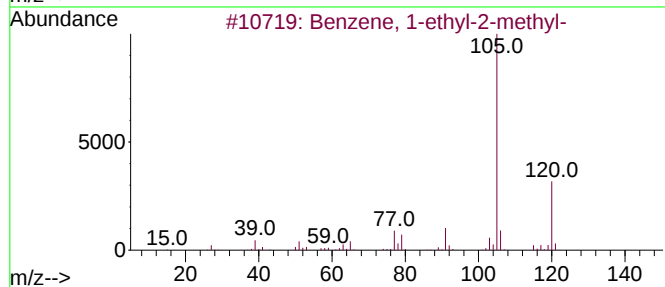
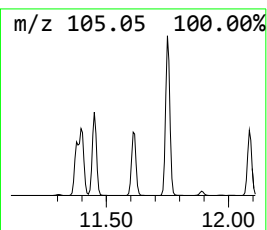
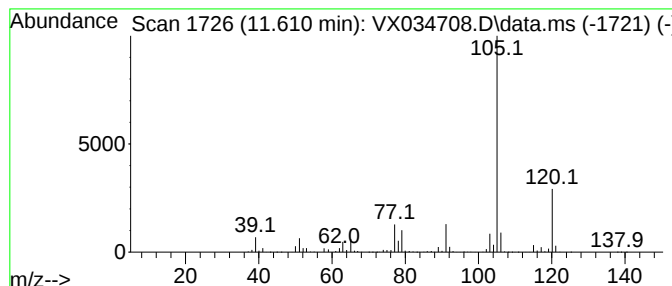
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1-ethyl-3-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.610	60.54 ug/l	1673530	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-3-methyl-	120	C9H12	000620-14-4	91
3			Mesitylene	120	C9H12	000108-67-8	91
4			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX032323\
Data File : VX034708.D
Acq On : 23 Mar 2023 18:58
Operator : JC/MD
Sample : 02055-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 26 Sample Multiplier: 1

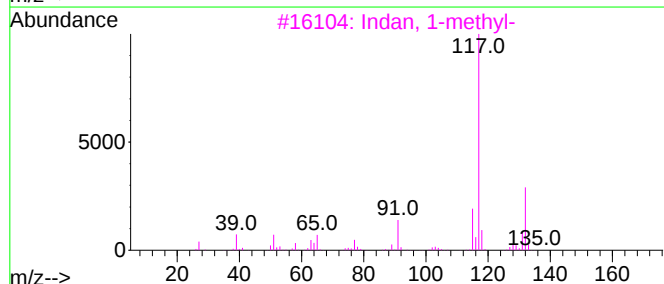
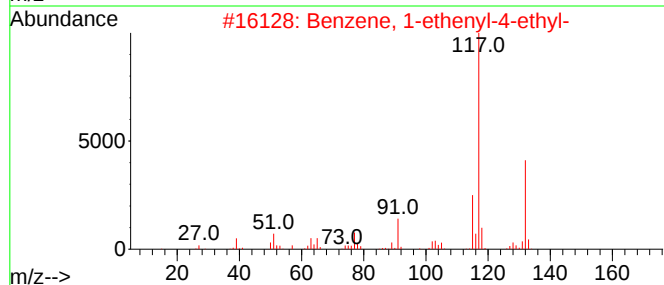
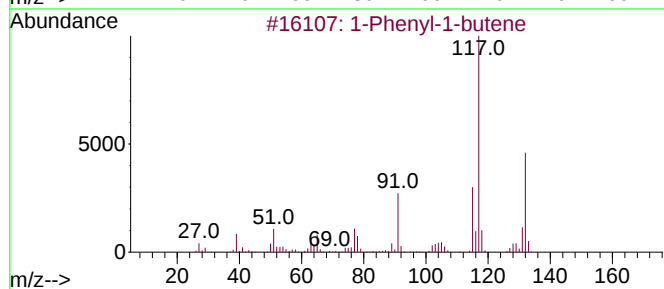
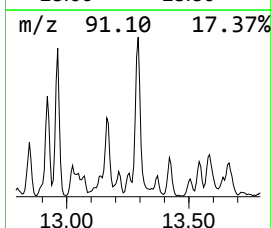
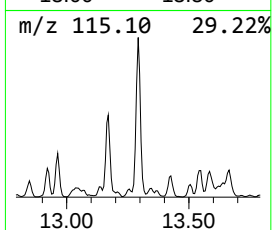
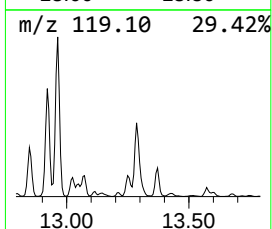
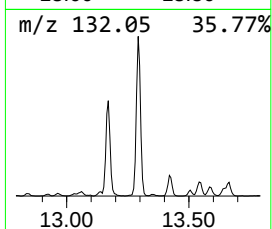
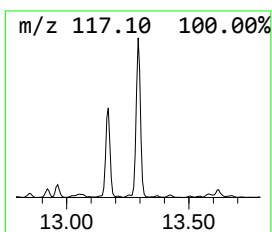
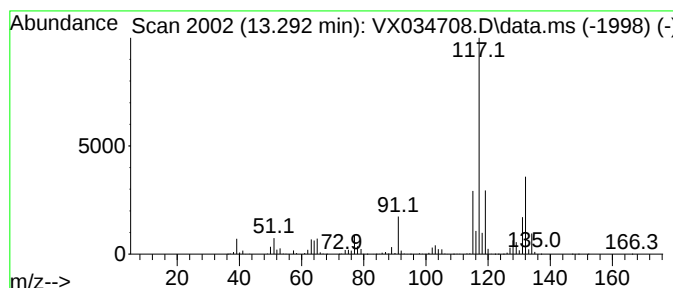
Instrument :
MSVOA_X
ClientSampleId :
MW-4B

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X031023W.M
Quant Title : SW846 8260

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

Peak Number 10 1-Phenyl-1-butene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.292	55.70 ug/l	1539690	1,4-Dichlorobenzene-d4	12.024	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Phenyl-1-butene	132	C10H12	000824-90-8	90
2	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	89
3	Indan, 1-methyl-	132	C10H12	000767-58-8	81
4	Benzene, 2-butenyl-	132	C10H12	001560-06-1	81
5	3-Phenylbut-1-ene	132	C10H12	000934-10-1	81



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX032323\
Data File : VX034708.D
Acq On : 23 Mar 2023 18:58
Operator : JC/MD
Sample : 02055-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-4B

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X031023W.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-	1.752	204.6	ug/l	4263570	1	5.556	1041920	50.0
Pentane	1.953	79.3	ug/l	1651930	1	5.556	1041920	50.0
Pentane, 2-methyl-	2.825	147.0	ug/l	3062760	1	5.556	1041920	50.0
Pentane, 3-methyl-	3.099	89.5	ug/l	1865400	1	5.556	1041920	50.0
Cyclopentane, m...	4.306	136.7	ug/l	2849200	1	5.556	1041920	50.0
Hexane, 3-methyl-	5.824	54.2	ug/l	1129150	1	5.556	1041920	50.0
Benzene, 1-ethy...	11.378	44.1	ug/l	1219150	4	12.024	1382060	50.0
Benzene, 1-ethy...	11.396	61.2	ug/l	1690890	4	12.024	1382060	50.0
Benzene, 1-ethy...	11.610	60.5	ug/l	1673530	4	12.024	1382060	50.0
1-Phenyl-1-butene	13.292	55.7	ug/l	1539690	4	12.024	1382060	50.0