

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110722\
 Data File : VX032631.D
 Acq On : 07 Nov 2022 16:48
 Operator : JC/MD
 Sample : N5484-25
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 DUP110422

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\82X102622W.M
 Title : SW846 8260

Signal : TIC: VX032631.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.745	103	108	122	rVB	220624	304873	22.18%	1.730%
2	1.953	137	142	153	rVB	165139	222351	16.18%	1.262%
3	2.337	198	205	219	rVB	31140	66209	4.82%	0.376%
4	2.715	258	267	271	rBV4	7832	18735	1.36%	0.106%
5	2.782	271	278	281	rVV	116685	248286	18.06%	1.409%
6	2.825	281	285	295	rVB	275194	542599	39.48%	3.079%
7	3.099	321	330	344	rBV	214780	484635	35.26%	2.750%
8	3.331	359	368	378	rVB2	17797	45555	3.31%	0.259%
9	3.446	378	387	401	rBV	118279	262251	19.08%	1.488%
10	3.593	401	411	415	rBV2	7720	18065	1.31%	0.103%
11	3.666	415	423	430	rVV2	21668	54227	3.95%	0.308%
12	3.837	442	451	458	rBV3	8967	23767	1.73%	0.135%
13	3.946	458	469	479	rBV4	10336	29493	2.15%	0.167%
14	4.105	488	495	502	rVB5	6965	17873	1.30%	0.101%
15	4.215	502	513	518	rVV3	17548	49723	3.62%	0.282%
16	4.306	518	528	540	rVB	175407	504138	36.68%	2.861%
17	5.117	648	661	662	rBV6	6183	18147	1.32%	0.103%
18	5.385	692	705	713	rBV	73736	227828	16.58%	1.293%
19	5.477	713	720	726	rBV4	45728	145848	10.61%	0.828%
20	5.556	726	733	739	rVB	78665	204779	14.90%	1.162%
21	5.830	766	778	790	rBV3	44041	141096	10.27%	0.801%
22	5.958	790	799	814	rBV	76199	199516	14.52%	1.132%
23	6.159	814	832	839	rBV3	38410	126561	9.21%	0.718%
24	6.257	839	848	857	rVB3	68098	203074	14.78%	1.152%
25	6.354	857	864	874	rVB2	39870	110746	8.06%	0.628%
26	6.458	874	881	894	rVB6	6293	19719	1.43%	0.112%
27	6.610	898	906	913	rBV2	15509	38728	2.82%	0.220%
28	6.763	923	931	939	rBV	177793	429608	31.26%	2.438%
29	7.068	974	981	991	rVB3	8027	18419	1.34%	0.105%
30	7.184	991	1000	1007	rBV3	13129	31914	2.32%	0.181%
31	7.379	1020	1032	1044	rBV3	163737	424262	30.87%	2.408%
32	7.501	1044	1052	1057	rBV3	8034	17058	1.24%	0.097%
33	7.562	1057	1062	1067	rBV3	8380	16083	1.17%	0.091%
34	7.659	1072	1078	1087	rVB	31481	61419	4.47%	0.349%
35	7.775	1090	1097	1104	rBV4	8335	17420	1.27%	0.099%
36	7.988	1123	1132	1142	rVB	35944	81094	5.90%	0.460%

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Integrator: RTE

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37	8.110	1142	1152	1160	rBV2	54054	118740	8.64%	0.674%
38	8.433	1201	1205	1209	rBV3	11031	18077	1.32%	0.103%
39	8.506	1212	1217	1226	rVB4	16251	32413	2.36%	0.184%
40	8.604	1226	1233	1235	rBV4	17625	33312	2.42%	0.189%
41	8.647	1235	1240	1249	rVB	372413	602020	43.80%	3.416%
42	9.104	1306	1315	1320	rBV3	11719	23432	1.70%	0.133%
43	9.165	1320	1325	1329	rBV	19389	28899	2.10%	0.164%
44	10.055	1465	1471	1476	rVV	382535	489502	35.62%	2.778%
45	10.140	1476	1485	1489	rVB6	6922	20831	1.52%	0.118%
46	10.195	1489	1494	1499	rBV	14261	19596	1.43%	0.111%
47	10.299	1505	1511	1518	rVB	438560	609886	44.37%	3.461%
48	10.640	1562	1567	1577	rVB	370680	481936	35.06%	2.735%
49	10.963	1614	1620	1626	rBV	47565	62690	4.56%	0.356%
50	11.079	1634	1639	1646	rBV	347276	444704	32.36%	2.524%
51	11.305	1673	1676	1680	rVB	17506	19303	1.40%	0.110%
52	11.378	1684	1688	1689	rBV	229510	237688	17.29%	1.349%
53	11.396	1689	1691	1696	rVV	311071	396643	28.86%	2.251%
54	11.451	1696	1700	1711	rVB	477306	566231	41.20%	3.213%
55	11.616	1720	1727	1736	rBV	311036	397966	28.96%	2.258%
56	11.750	1744	1749	1755	rBV	1108787	1374424	100.00%	7.799%
57	11.866	1762	1768	1770	rBV	25236	35349	2.57%	0.201%
58	11.890	1770	1772	1778	rVB	34419	38249	2.78%	0.217%
59	11.969	1780	1785	1789	rBV	90657	110559	8.04%	0.627%
60	12.024	1789	1794	1799	rVV	399329	531677	38.68%	3.017%
61	12.085	1799	1804	1810	rVB	317329	386328	28.11%	2.192%
62	12.250	1825	1831	1832	rBV	134242	164907	12.00%	0.936%
63	12.268	1832	1834	1838	rVV	179381	195104	14.20%	1.107%
64	12.317	1838	1842	1852	rVB2	463706	671314	48.84%	3.810%
65	12.402	1852	1856	1861	rBV2	31295	40571	2.95%	0.230%
66	12.463	1861	1866	1872	rVB	145773	173191	12.60%	0.983%
67	12.530	1872	1877	1879	rBV	247708	304745	22.17%	1.729%
68	12.555	1879	1881	1886	rVB	254157	270197	19.66%	1.533%
69	12.615	1886	1891	1895	rBV	387442	477510	34.74%	2.710%
70	12.701	1900	1905	1913	rBV3	134277	235914	17.16%	1.339%
71	12.798	1917	1921	1924	rVB2	14204	18161	1.32%	0.103%
72	12.847	1924	1929	1934	rVB	90096	111800	8.13%	0.634%
73	12.920	1934	1941	1945	rBV	242340	311176	22.64%	1.766%
74	12.963	1945	1948	1954	rVB	326081	370501	26.96%	2.102%
75	13.048	1954	1962	1964	rBV3	43932	91926	6.69%	0.522%
76	13.073	1964	1966	1968	rVB	20339	17758	1.29%	0.101%

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77	13.164	1978	1981	1985	rVB2	33890	45529	3.31%	0.258%
78	13.213	1985	1989	1992	rVB3	25477	29932	2.18%	0.170%
79	13.256	1992	1996	1998	rBV2	40165	55957	4.07%	0.318%
80	13.286	1998	2001	2008	rVV2	113058	177252	12.90%	1.006%
81	13.371	2012	2015	2020	rVB	33539	39828	2.90%	0.226%
82	13.426	2020	2024	2031	rVB2	16753	24917	1.81%	0.141%
83	13.506	2031	2037	2040	rBV3	30028	45901	3.34%	0.260%
84	13.542	2040	2043	2046	rVV	66381	88310	6.43%	0.501%
85	13.579	2046	2049	2056	rVV3	86112	157782	11.48%	0.895%
86	13.640	2056	2059	2060	rVV	32452	36744	2.67%	0.209%
87	13.664	2060	2063	2069	rVB2	68089	114573	8.34%	0.650%
88	13.774	2075	2081	2090	rVB	111306	149040	10.84%	0.846%
89	13.859	2090	2095	2099	rVB5	10300	16744	1.22%	0.095%
90	13.926	2099	2106	2110	rBV2	35226	53664	3.90%	0.305%
91	13.969	2110	2113	2117	rVB	18305	21896	1.59%	0.124%
92	14.079	2126	2131	2136	rBV	42637	56112	4.08%	0.318%
93	14.213	2148	2153	2163	rBV	50236	87730	6.38%	0.498%
94	14.323	2167	2171	2176	rVV	26823	34177	2.49%	0.194%
95	14.371	2176	2179	2183	rVB	22890	27452	2.00%	0.156%
96	14.511	2198	2202	2211	rVB	31647	47348	3.44%	0.269%
97	14.640	2211	2223	2233	rVB	157002	218054	15.87%	1.237%
98	14.780	2240	2246	2251	rBV	69760	88234	6.42%	0.501%
99	15.475	2355	2360	2371	rVB2	13490	22646	1.65%	0.129%
100	15.615	2378	2383	2386	rBV2	14756	20934	1.52%	0.119%

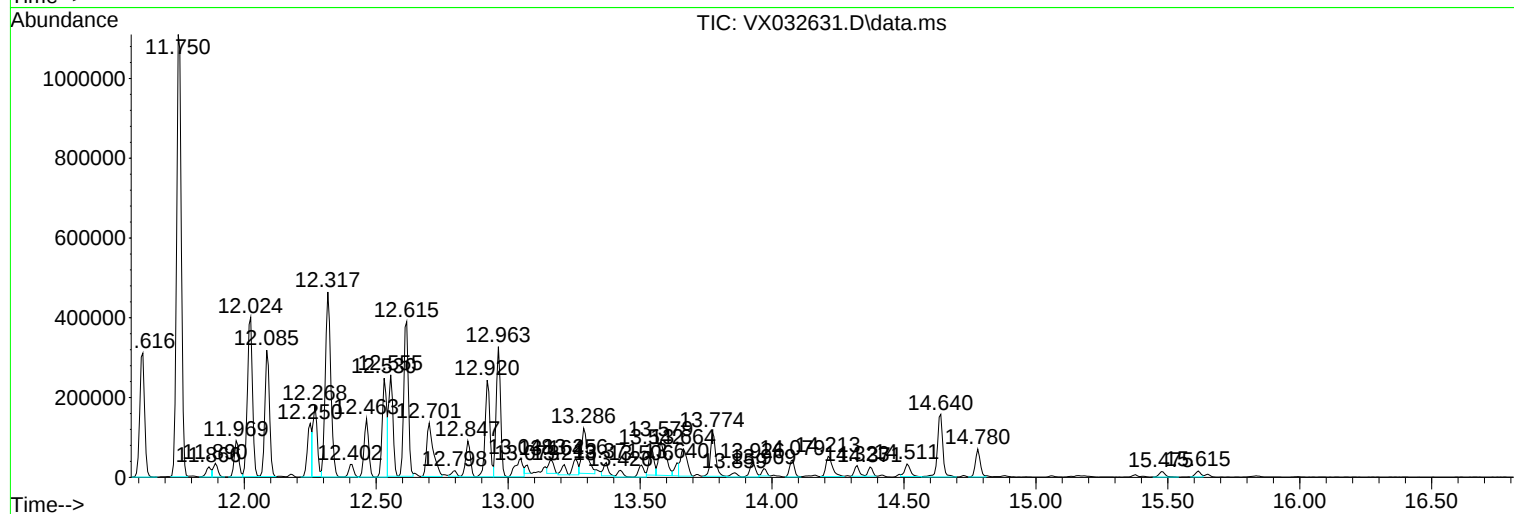
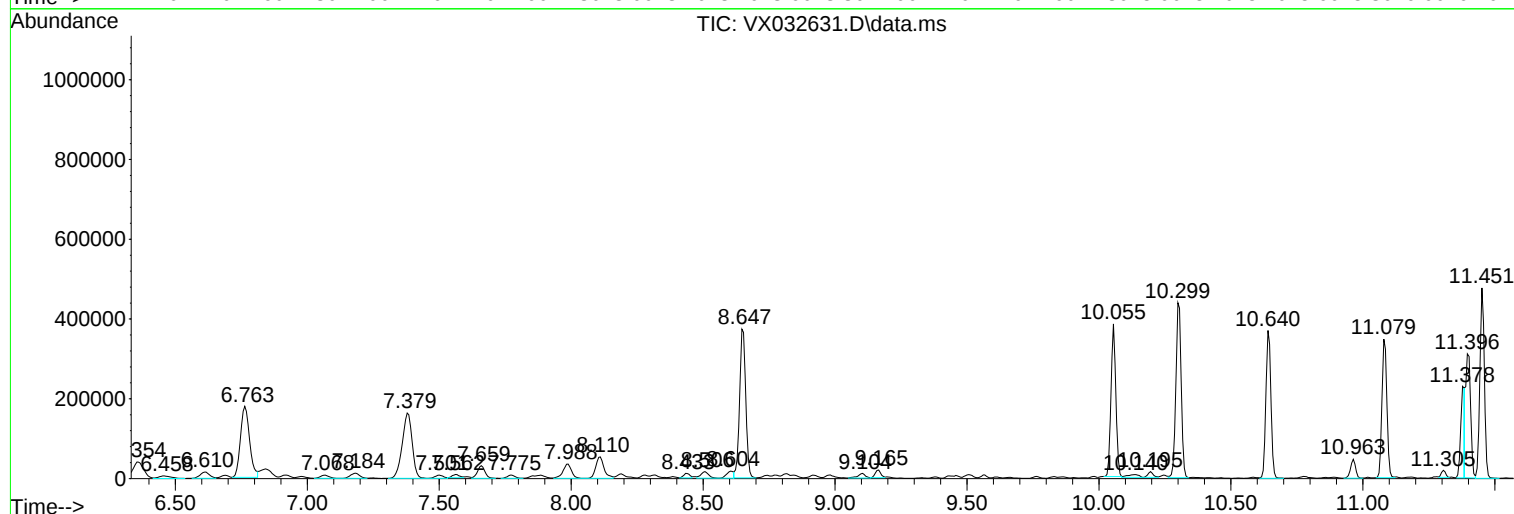
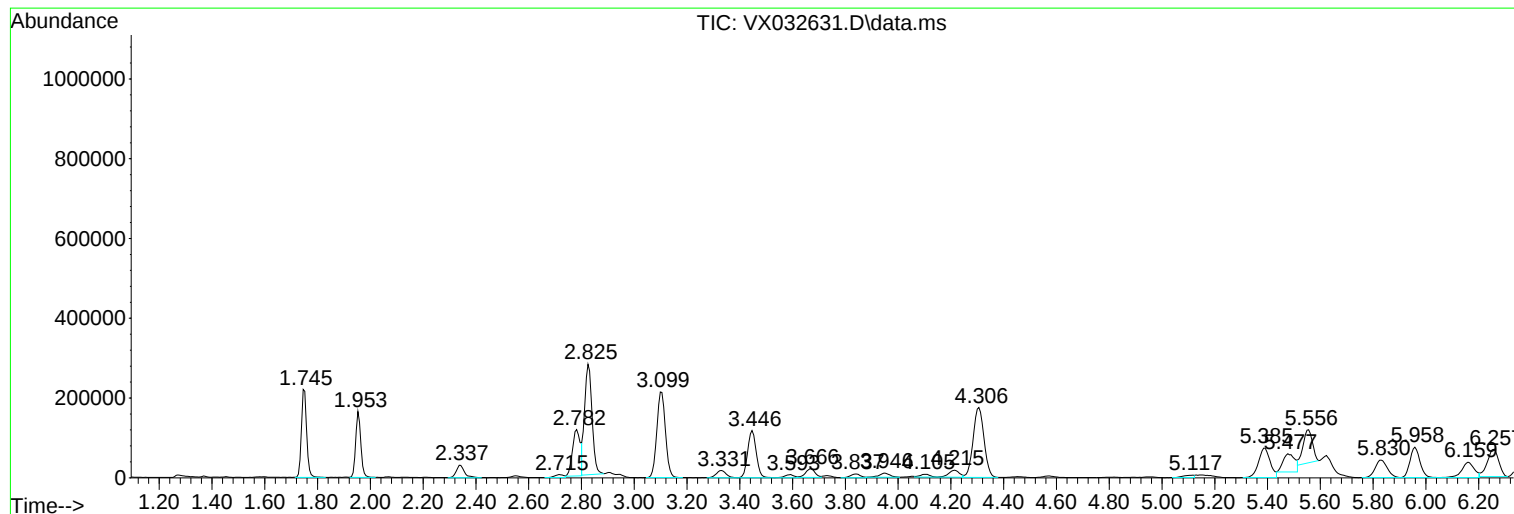
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Instrument :
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ClientSampleId :
DUP110422

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title : SW846 8260

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



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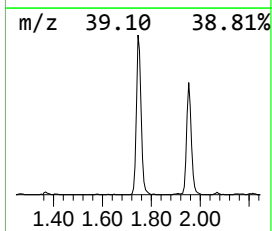
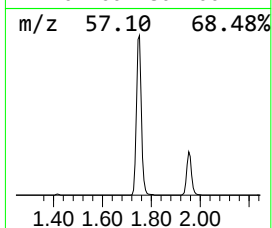
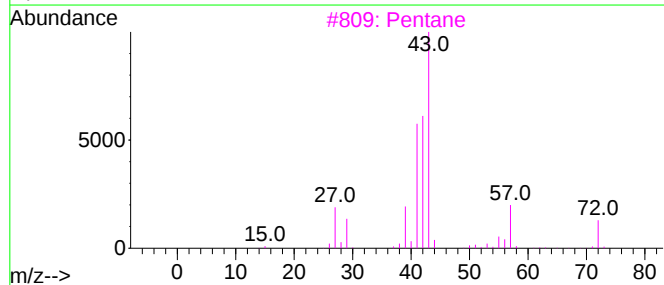
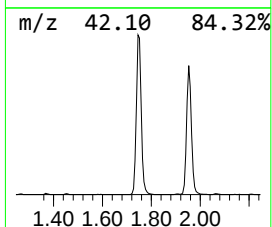
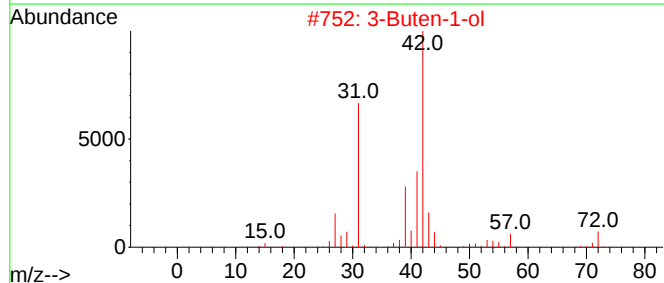
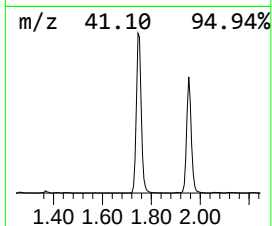
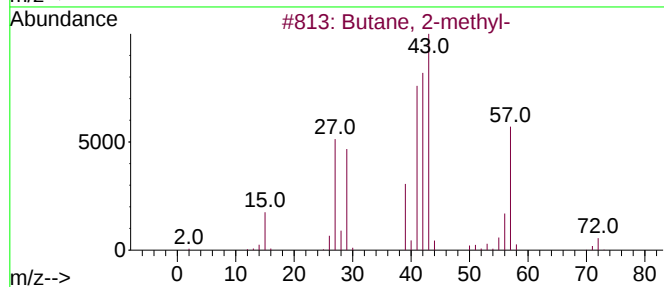
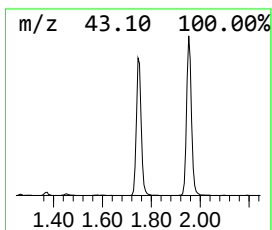
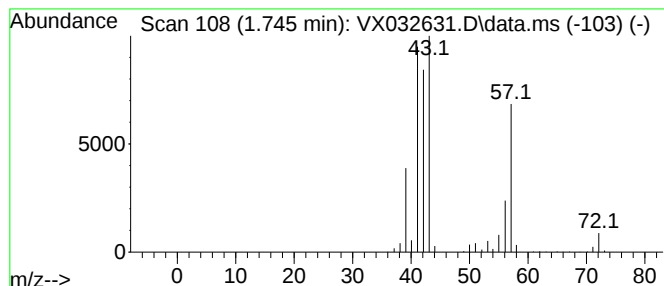
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title : SW846 8260

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

Peak Number 1 Butane, 2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.745	74.44 ug/l	304873	Pentafluorobenzene	5.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methyl-	72	C5H12	000078-78-4	64
2			3-Buten-1-ol	72	C4H8O	000627-27-0	38
3			Pentane	72	C5H12	000109-66-0	38
4			1-Butene	56	C4H8	000106-98-9	16
5			Butane, 1-chloro-2-methyl-	106	C5H11Cl	000616-13-7	10



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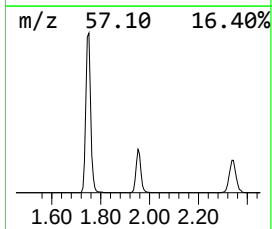
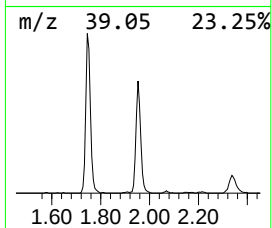
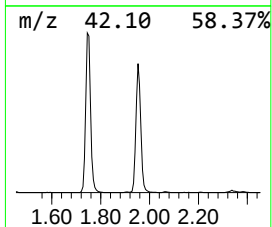
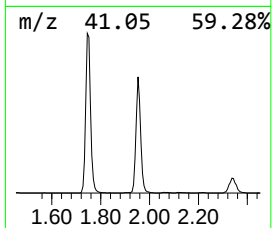
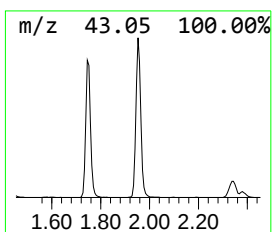
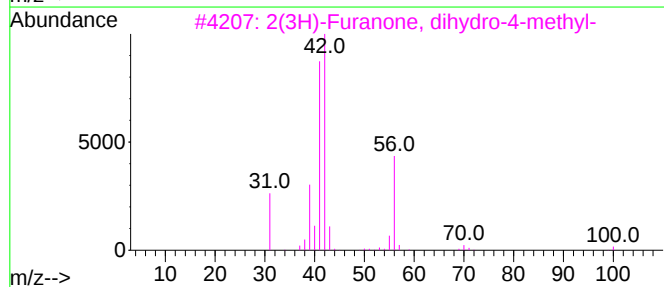
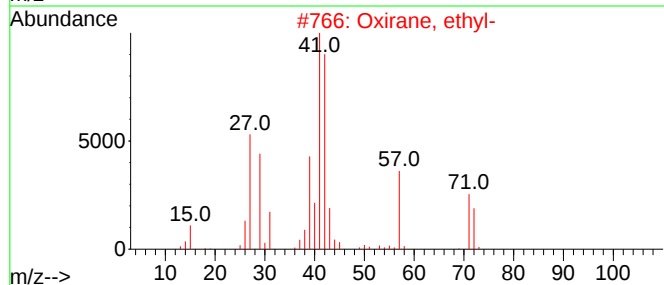
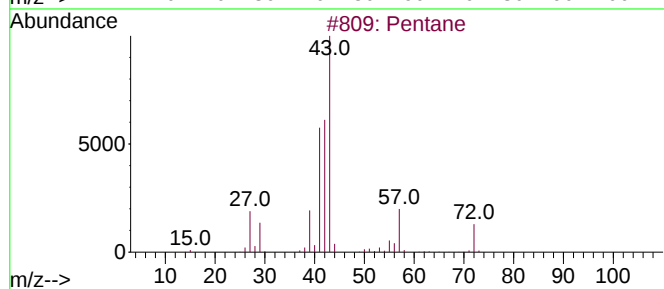
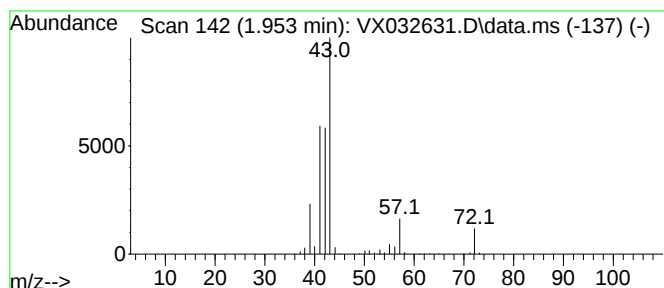
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title : SW846 8260

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

Peak Number 2 Pentane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.953	54.29 ug/l	222351	Pentafluorobenzene	5.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane	72	C5H12	000109-66-0	91
2			Oxirane, ethyl-	72	C4H8O	000106-88-7	50
3			2(3H)-Furanone, dihydro-4-methyl-	100	C5H8O2	001679-49-8	38
4			Butane, 2-methyl-	72	C5H12	000078-78-4	36
5			Diaziridine,3,3-dimethyl-	72	C3H8N2	004901-76-2	25



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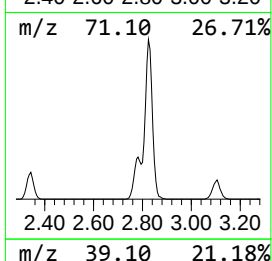
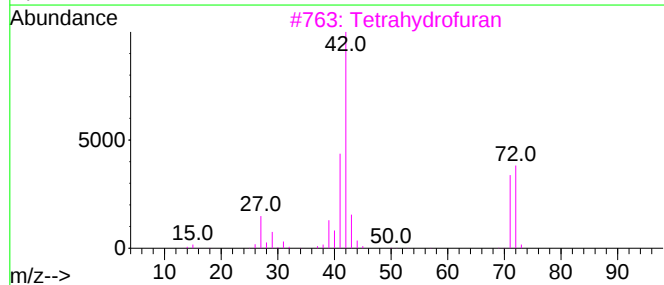
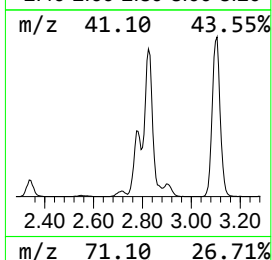
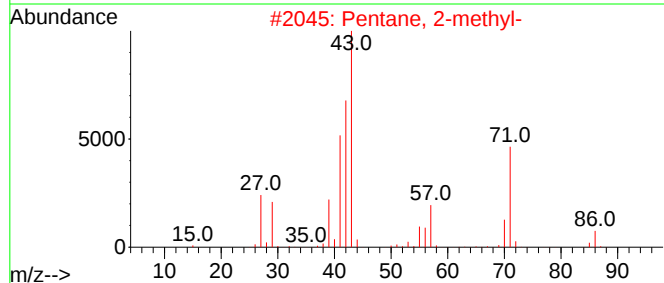
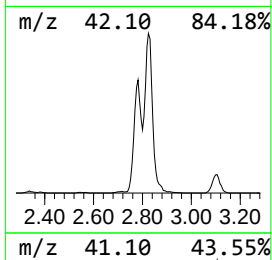
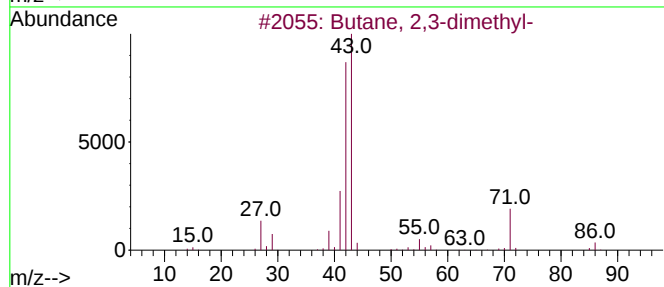
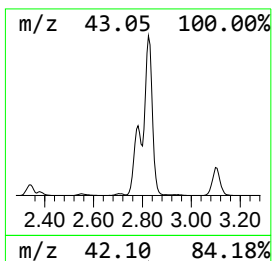
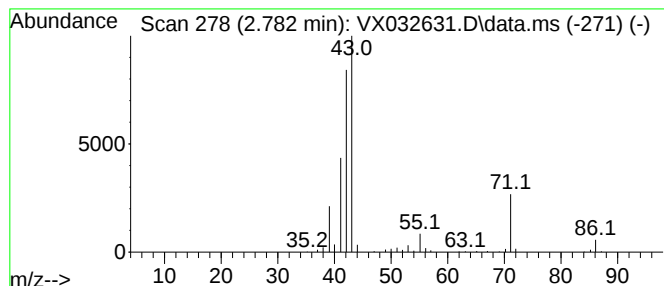
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title : SW846 8260

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

Peak Number 3 Butane, 2,3-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.782	60.62 ug/l	248286	Pentafluorobenzene	5.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	86
2			Pentane, 2-methyl-	86	C6H14	000107-83-5	78
3			Tetrahydrofuran	72	C4H8O	000109-99-9	38
4			Pentane	72	C5H12	000109-66-0	23
5			Propanoyl chloride, 2-methyl-	106	C4H7ClO	000079-30-1	12



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110722\
Data File : VX032631.D
Acq On : 07 Nov 2022 16:48
Operator : JC/MD
Sample : N5484-25
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
DUP110422

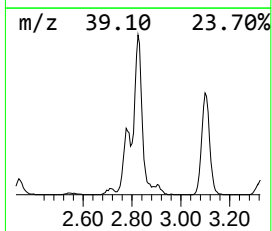
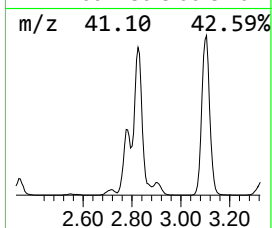
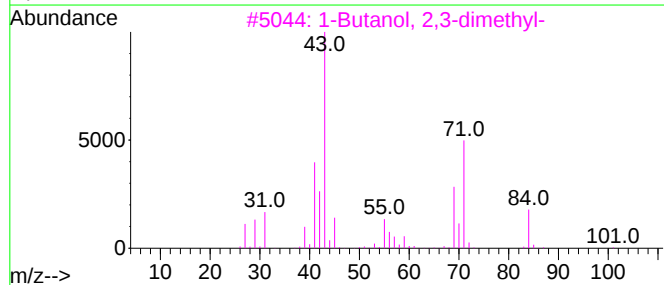
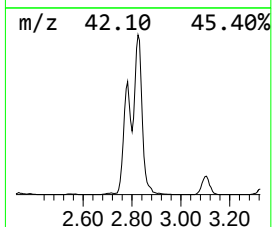
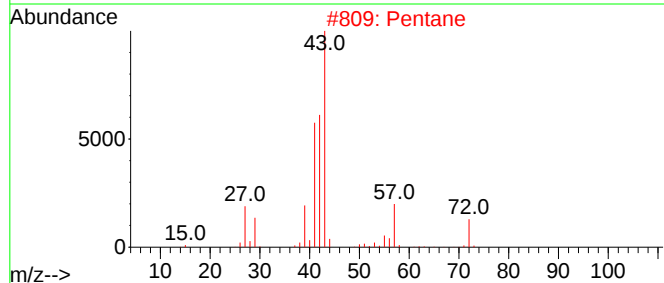
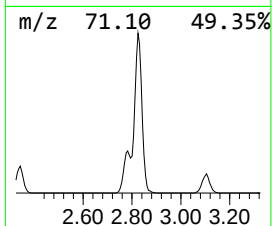
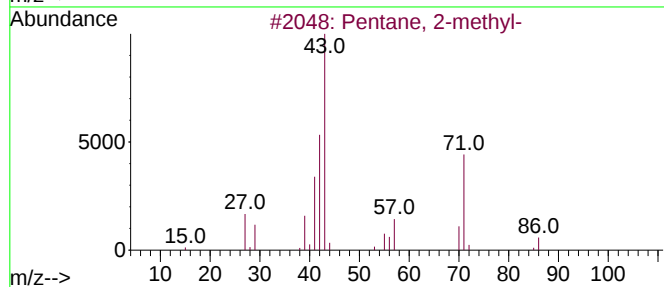
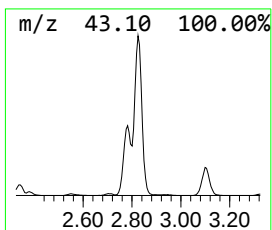
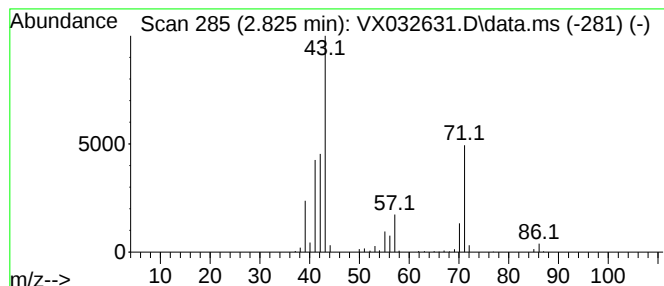
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Quant Title : SW846 8260

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

Peak Number 4 Pentane, 2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.825	132.48 ug/l	542599	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	49
3			1-Butanol, 2,3-dimethyl-	102	C6H14O	019550-30-2	42
4			Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	38
5			Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110722\
Data File : VX032631.D
Acq On : 07 Nov 2022 16:48
Operator : JC/MD
Sample : N5484-25
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
DUP110422

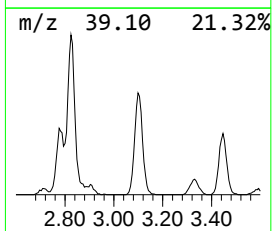
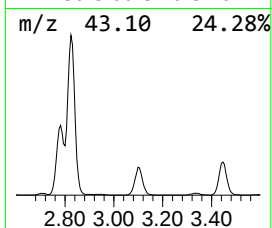
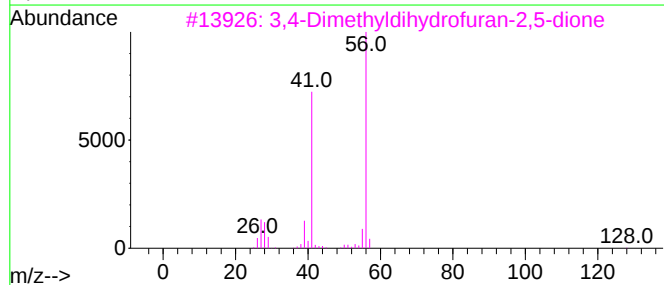
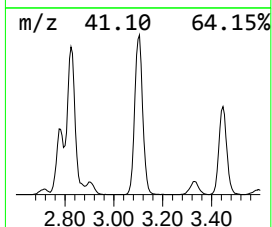
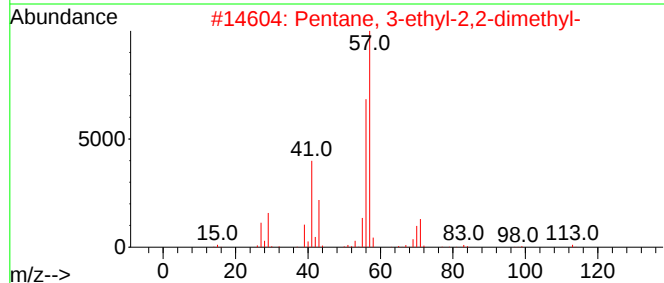
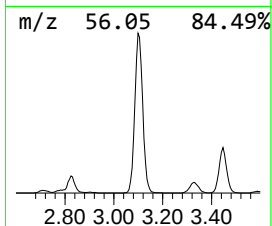
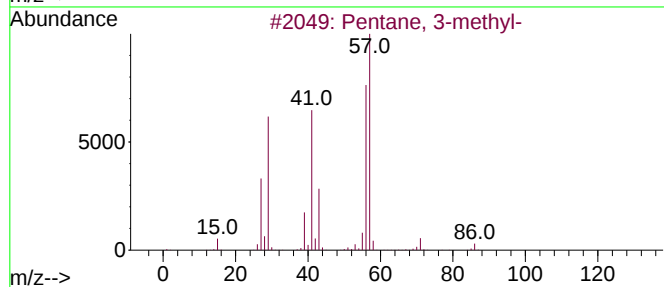
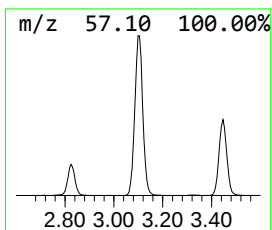
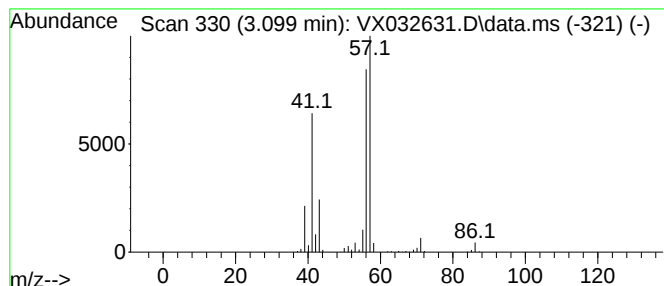
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Quant Title : SW846 8260

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

Peak Number 5 Pentane, 3-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.099	118.33 ug/l	484635	Pentafluorobenzene	5.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 3-methyl-	86	C6H14	000096-14-0	90
2			Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	64
3			3,4-Dimethyldihydrofuran-2,5-dione	128	C6H8O3	007475-92-5	53
4			Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	43
5			1-Butanol, 2-methyl-	88	C5H12O	000137-32-6	40



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110722\
Data File : VX032631.D
Acq On : 07 Nov 2022 16:48
Operator : JC/MD
Sample : N5484-25
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
DUP110422

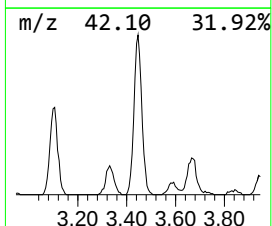
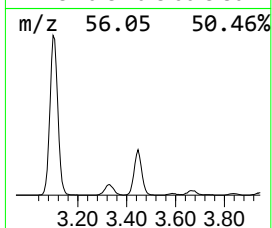
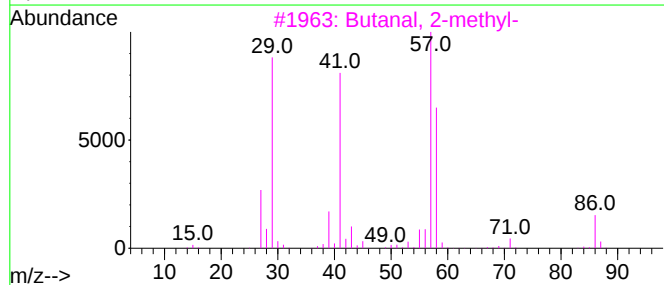
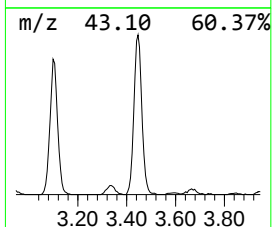
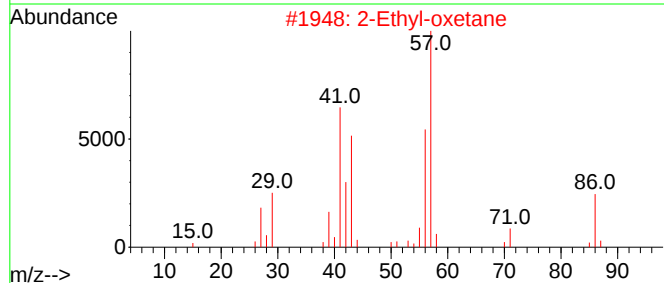
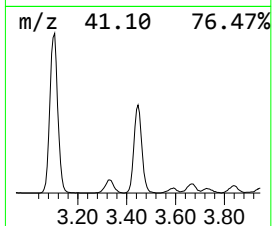
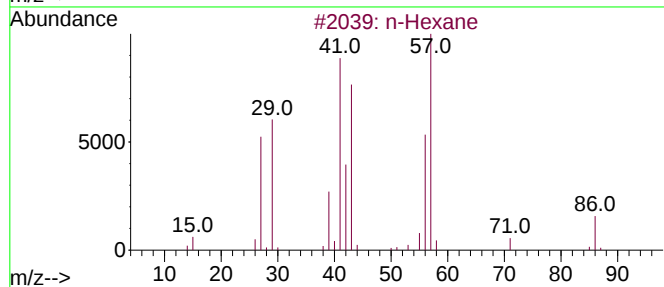
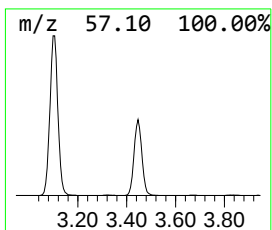
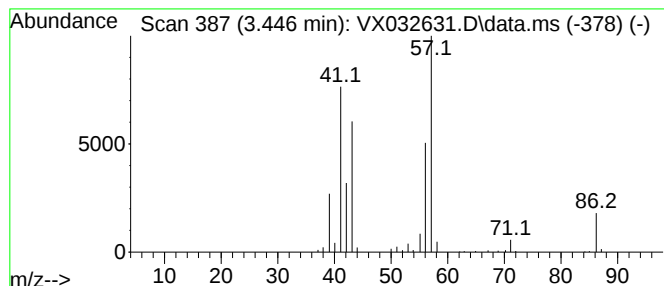
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Quant Title : SW846 8260

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

Peak Number 6 n-Hexane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.446	64.03 ug/l	262251	Pentafluorobenzene	5.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexane	86	C6H14	000110-54-3	91
2			2-Ethyl-oxetane	86	C5H10O	1010386-40-2	64
3			Butanal, 2-methyl-	86	C5H10O	000096-17-3	43
4			Propanal, 2,2-dimethyl-	86	C5H10O	000630-19-3	43
5			Pentane, 2,2,3,4-tetramethyl-	128	C9H20	001186-53-4	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110722\
Data File : VX032631.D
Acq On : 07 Nov 2022 16:48
Operator : JC/MD
Sample : N5484-25
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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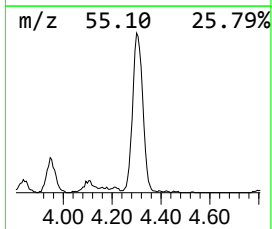
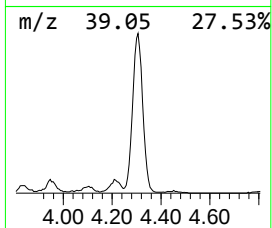
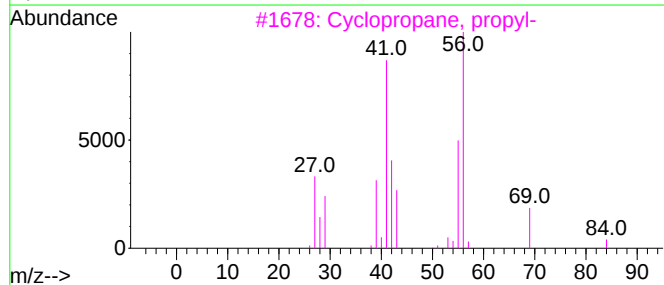
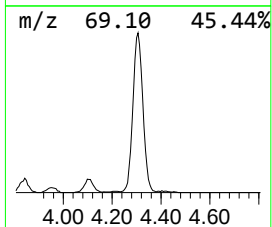
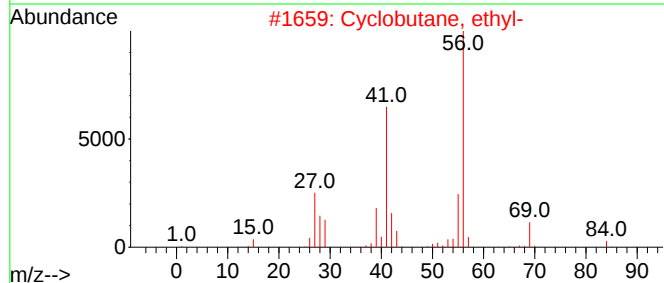
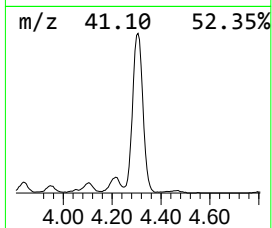
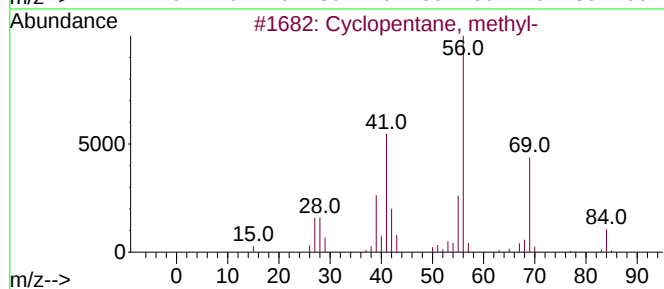
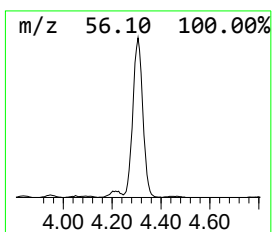
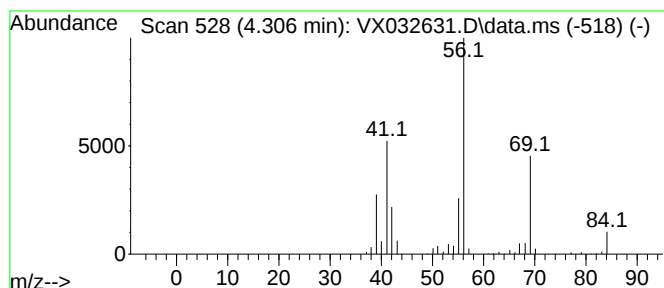
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Quant Title : SW846 8260

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

Peak Number 7 Cyclopentane, methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.306	123.09 ug/l	504138	Pentafluorobenzene	5.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, methyl-	84	C6H12	000096-37-7	90
2			Cyclobutane, ethyl-	84	C6H12	004806-61-5	72
3			Cyclopropane, propyl-	84	C6H12	002415-72-7	64
4			1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	50
5			1-Pentene, 2-methyl-	84	C6H12	000763-29-1	45



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110722\
Data File : VX032631.D
Acq On : 07 Nov 2022 16:48
Operator : JC/MD
Sample : N5484-25
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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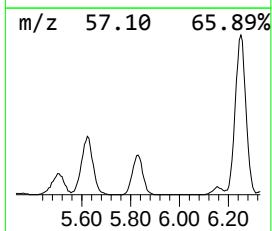
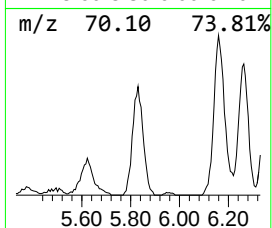
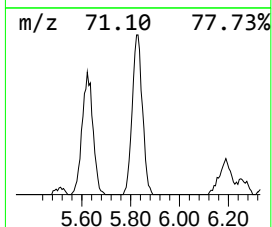
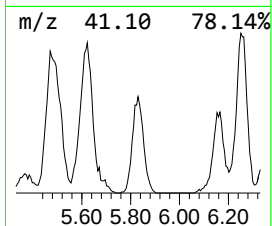
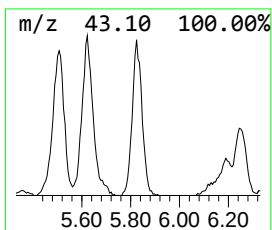
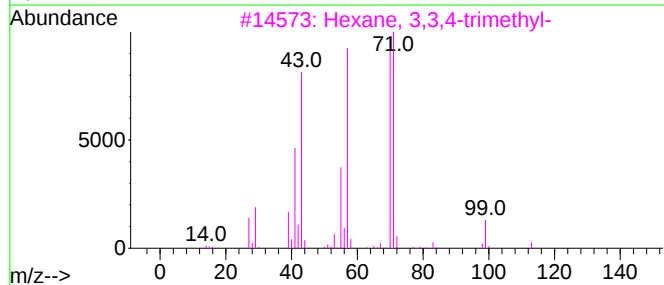
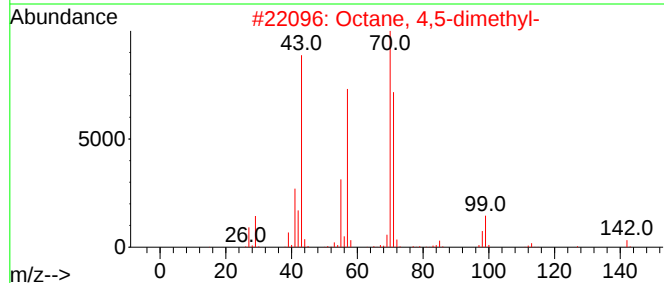
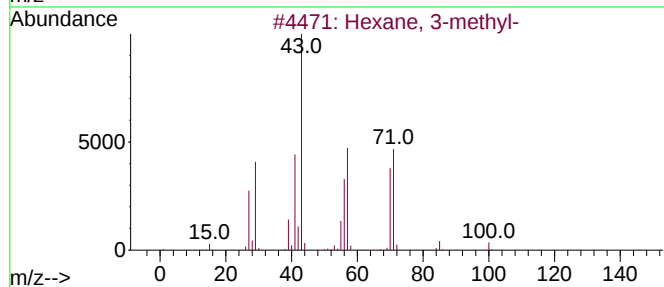
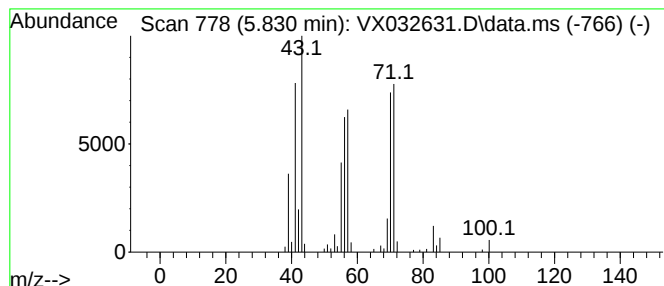
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Quant Title : SW846 8260

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

Peak Number 8 Hexane, 3-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.830	34.45 ug/l	141096	Pentafluorobenzene	5.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 3-methyl-			100	C7H16	000589-34-4	90
2	Octane, 4,5-dimethyl-			142	C10H22	015869-96-2	59
3	Hexane, 3,3,4-trimethyl-			128	C9H20	016747-31-2	53
4	Oxalic acid, isobutyl nonyl ester			272	C15H28O4	1010309-37-4	50
5	Undecane, 3-methylene-			168	C12H24	071138-64-2	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110722\
Data File : VX032631.D
Acq On : 07 Nov 2022 16:48
Operator : JC/MD
Sample : N5484-25
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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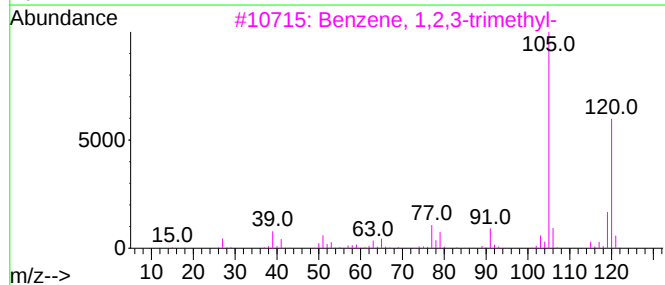
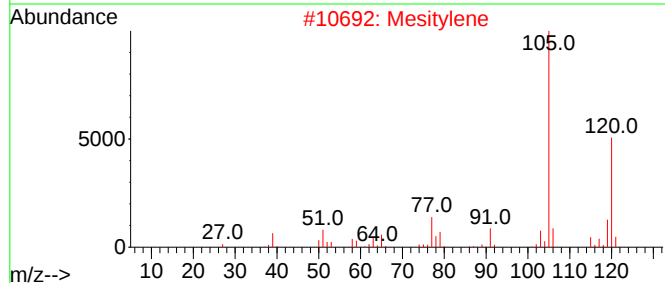
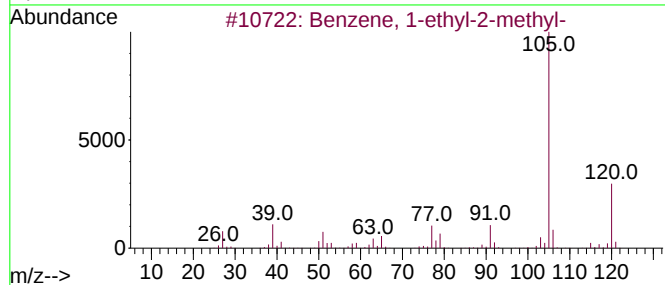
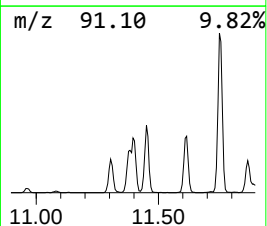
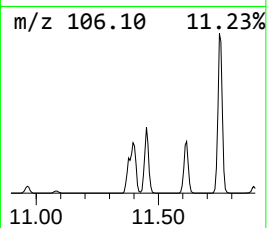
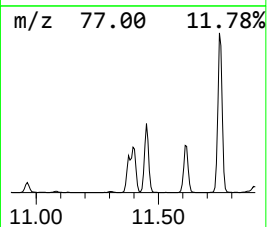
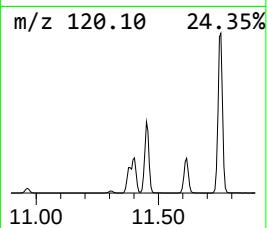
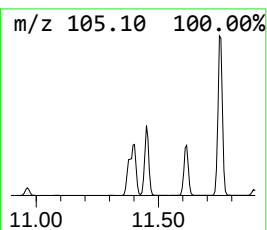
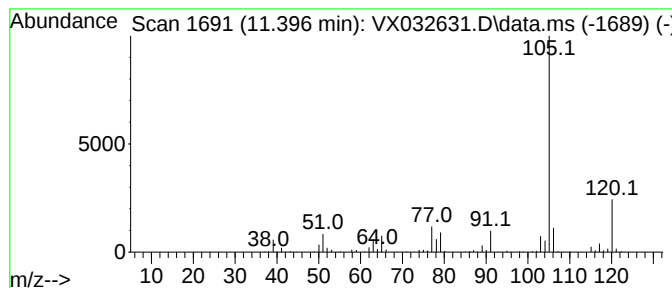
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1-ethyl-2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.396	37.30 ug/l	396643	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			Mesitylene	120	C9H12	000108-67-8	90
3			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	90
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	90
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110722\
Data File : VX032631.D
Acq On : 07 Nov 2022 16:48
Operator : JC/MD
Sample : N5484-25
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
DUP110422

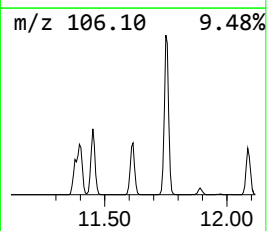
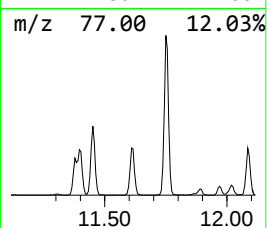
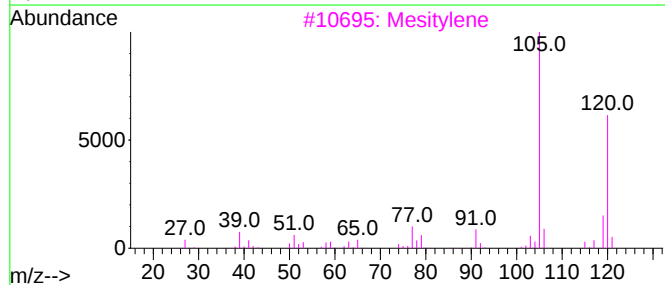
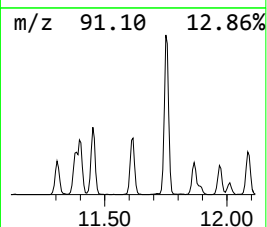
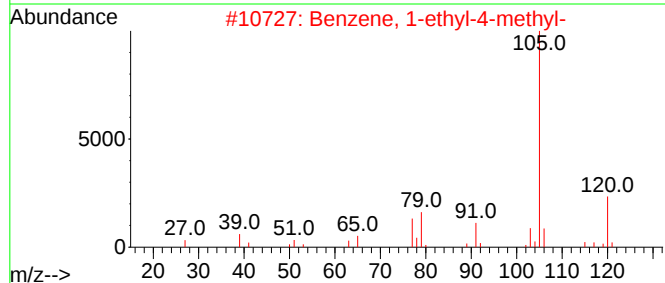
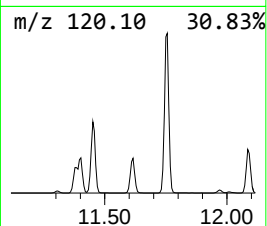
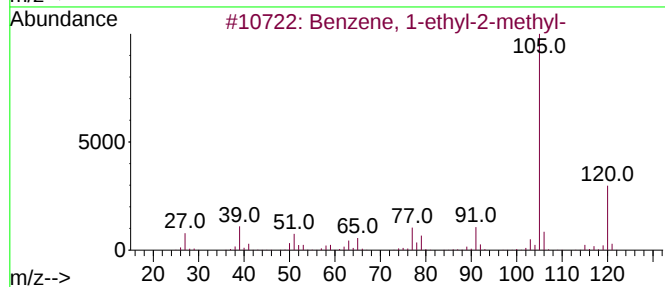
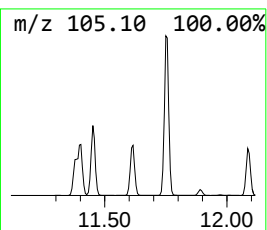
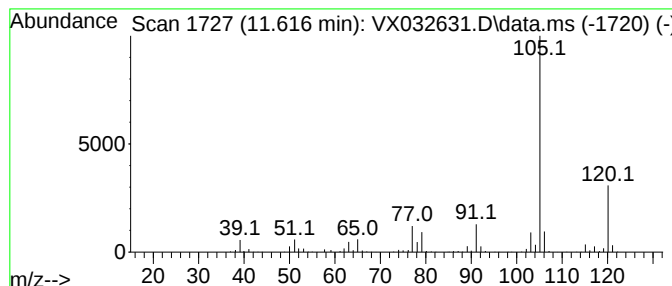
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Quant Title : SW846 8260

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

Peak Number 10 Benzene, 1-ethyl-4-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.616	37.43 ug/l	397966	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	93
3			Mesitylene	120	C9H12	000108-67-8	91
4			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110722\
Data File : VX032631.D
Acq On : 07 Nov 2022 16:48
Operator : JC/MD
Sample : N5484-25
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
DUP110422

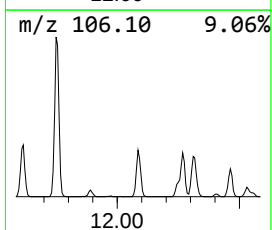
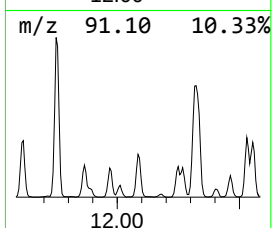
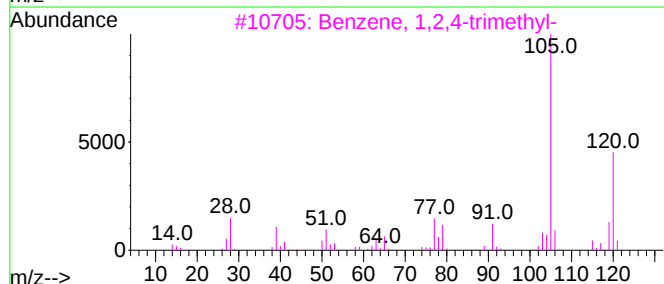
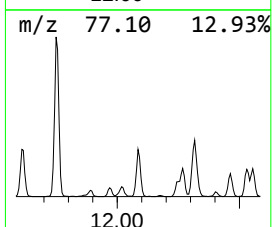
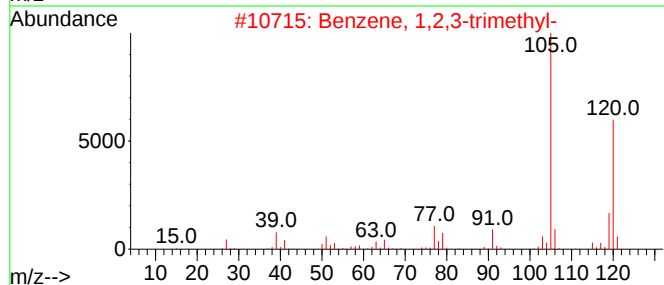
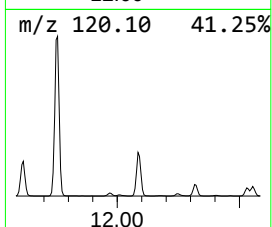
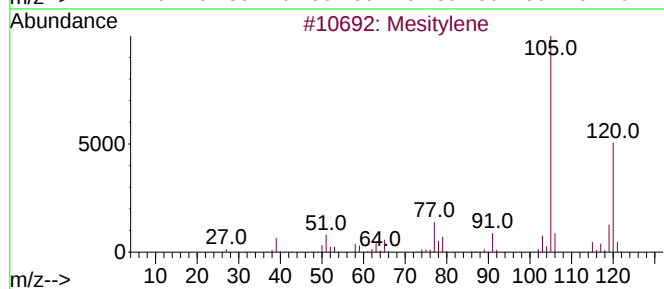
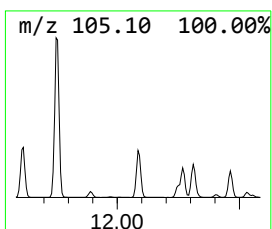
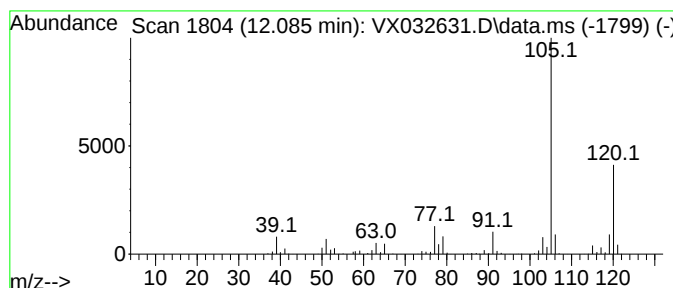
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Quant Title : SW846 8260

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

Peak Number 11 Benzene, 1,2,3-trimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.085	36.33 ug/l	386328	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, 1,2,4-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\VX110722\
Data File : VX032631.D
Acq On : 07 Nov 2022 16:48
Operator : JC/MD
Sample : N5484-25
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
DUP110422

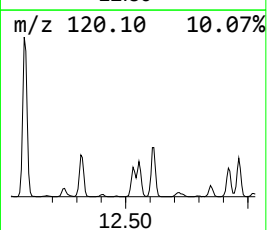
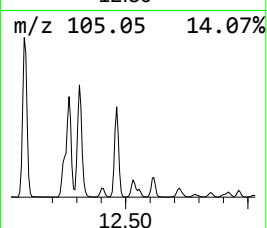
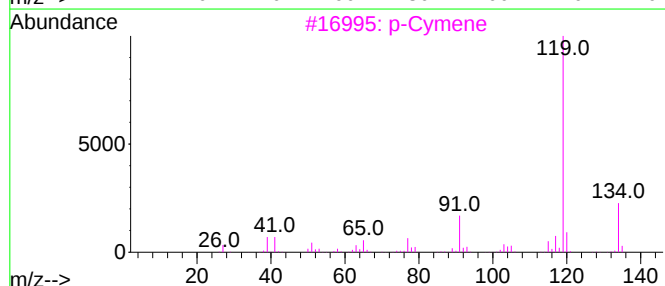
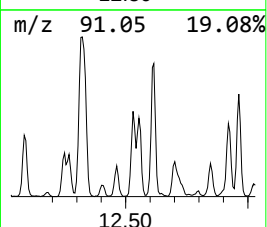
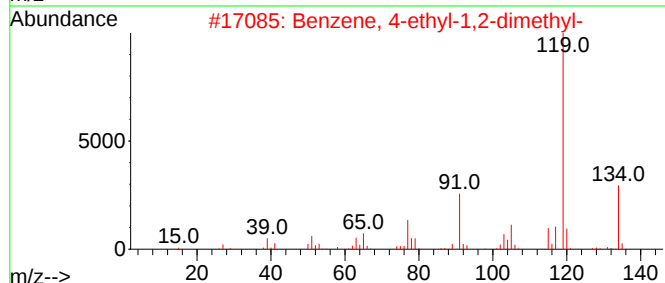
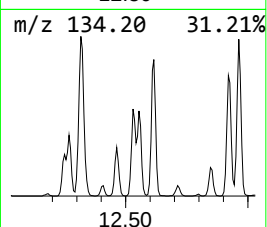
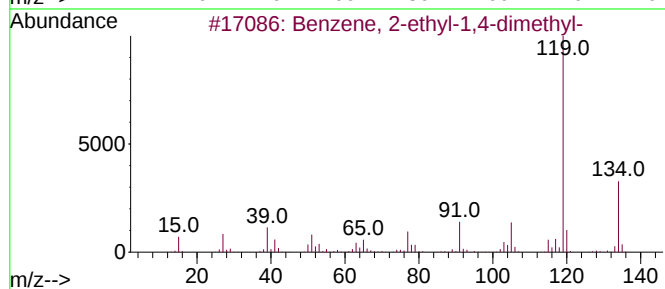
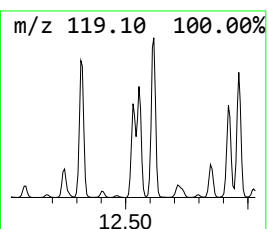
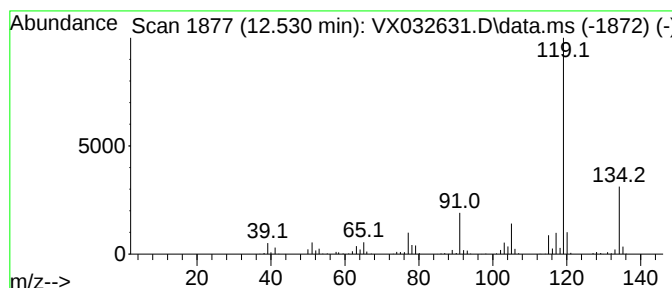
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Quant Title : SW846 8260

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

Peak Number 12 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.530	28.66 ug/l	304745	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	97
2			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
3			p-Cymene	134	C10H14	000099-87-6	95
4			o-Cymene	134	C10H14	000527-84-4	95
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110722\
Data File : VX032631.D
Acq On : 07 Nov 2022 16:48
Operator : JC/MD
Sample : N5484-25
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
DUP110422

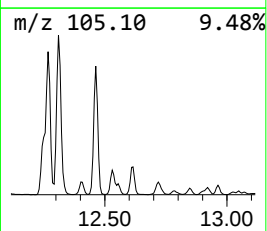
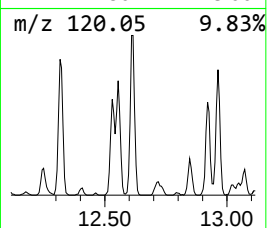
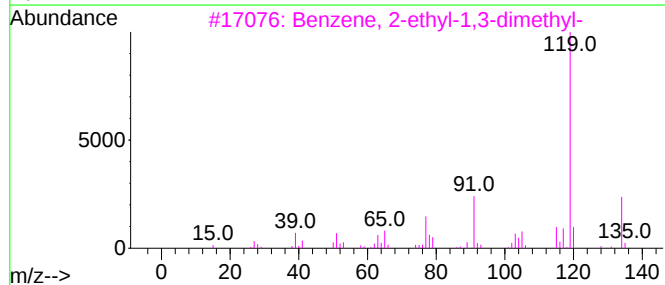
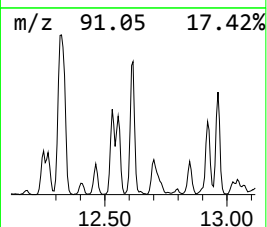
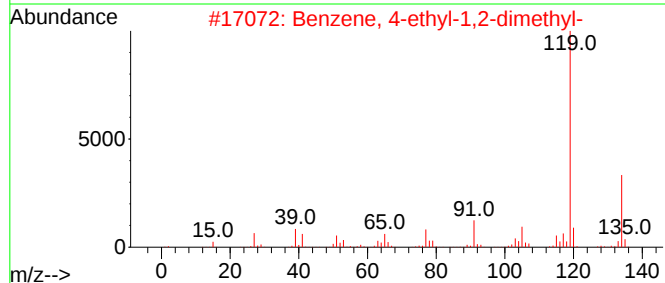
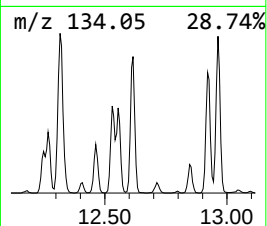
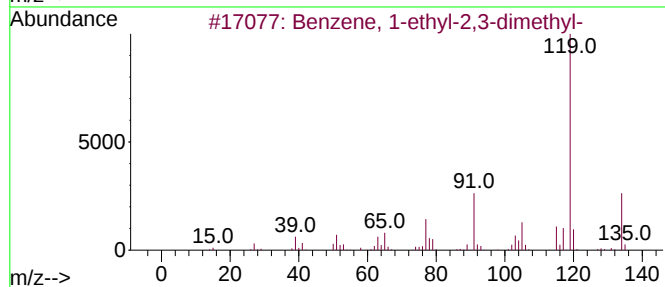
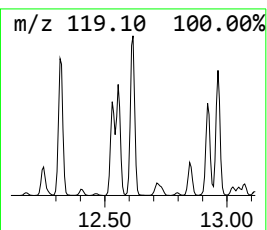
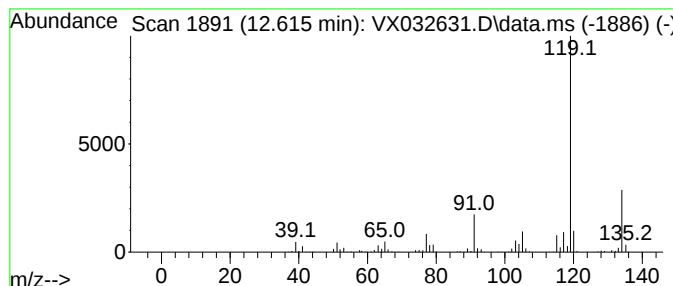
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Quant Title : SW846 8260

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

Peak Number 13 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.616	44.91 ug/l	477510	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
2			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
3			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
4			o-Cymene	134	C10H14	000527-84-4	95
5			p-Cymene	134	C10H14	000099-87-6	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110722\
Data File : VX032631.D
Acq On : 07 Nov 2022 16:48
Operator : JC/MD
Sample : N5484-25
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

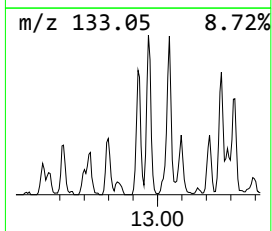
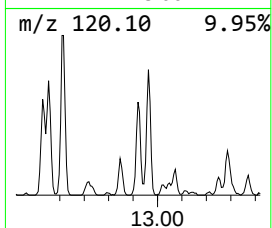
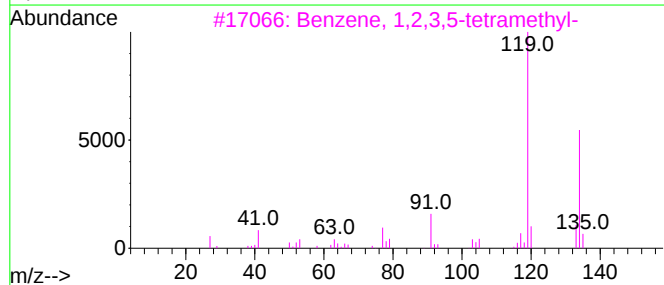
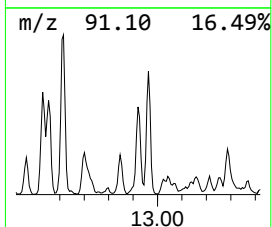
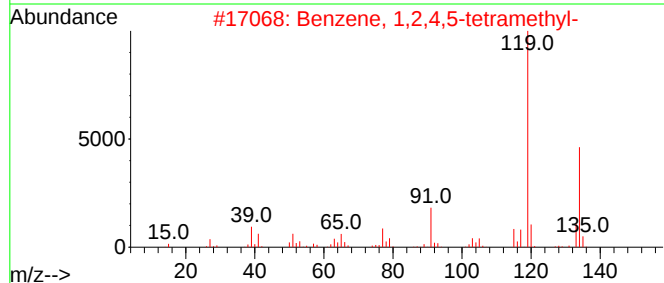
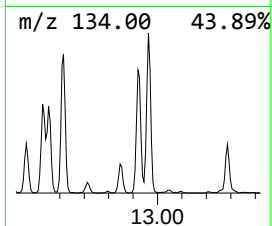
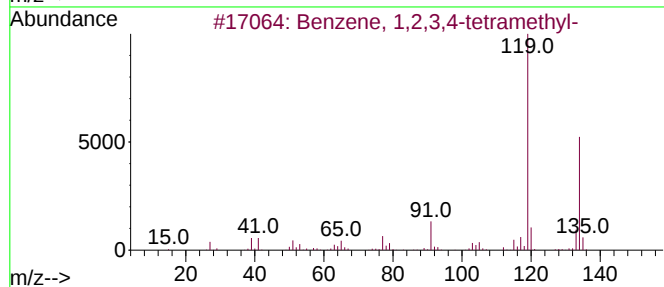
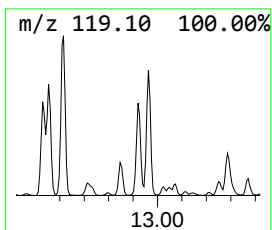
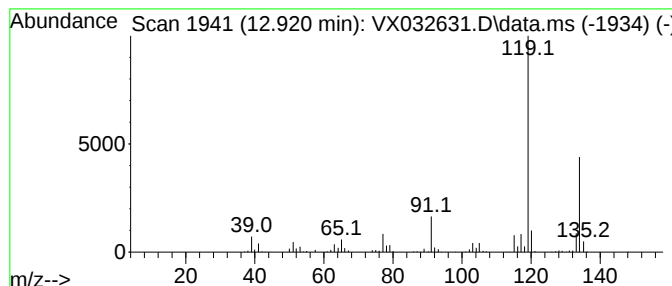
Instrument :
MSVOA_X
ClientSampleId :
DUP110422

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title : SW846 8260

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

Peak Number 14 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD			R.T.
12.920	29.26 ug/l	311176	1,4-Dichlorobenzene-d4			12.024
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,4-tetramethyl-		134	C10H14	000488-23-3	97
2	Benzene, 1,2,4,5-tetramethyl-		134	C10H14	000095-93-2	97
3	Benzene, 1,2,3,5-tetramethyl-		134	C10H14	000527-53-7	96
4	o-Cymene		134	C10H14	000527-84-4	95
5	Benzene, 1-ethyl-3,5-dimethyl-		134	C10H14	000934-74-7	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110722\
Data File : VX032631.D
Acq On : 07 Nov 2022 16:48
Operator : JC/MD
Sample : N5484-25
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
DUP110422

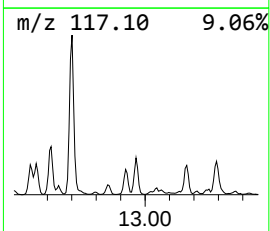
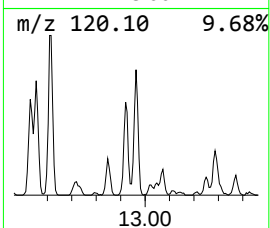
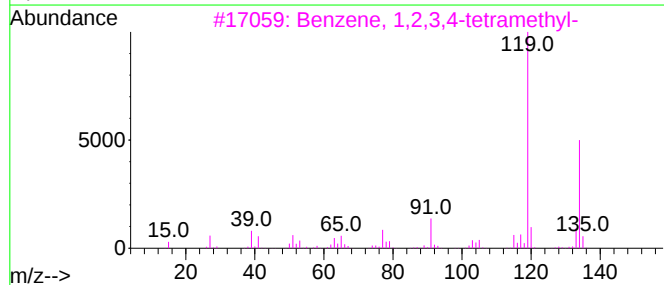
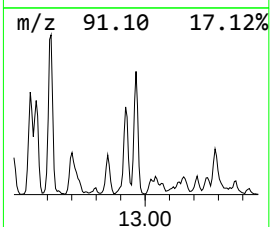
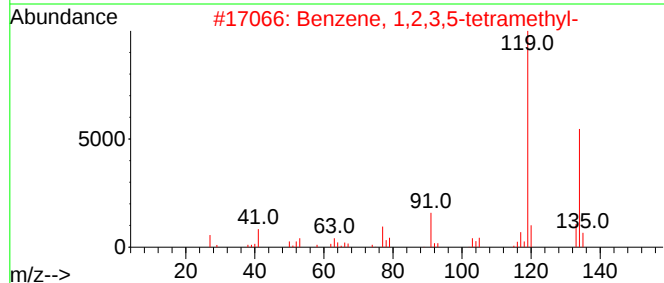
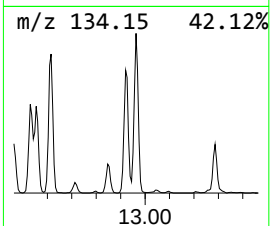
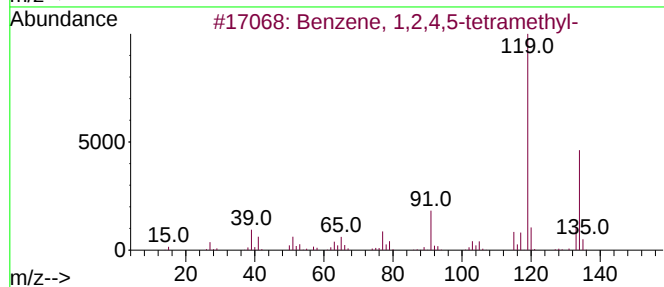
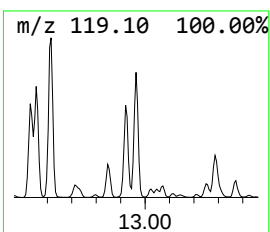
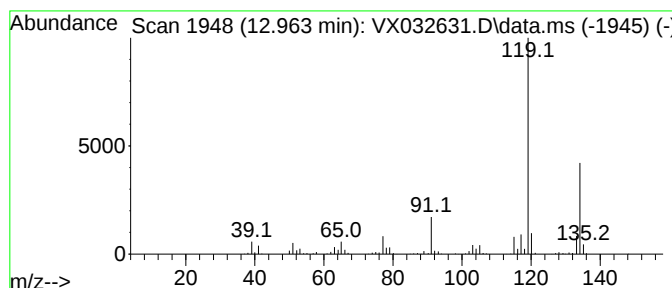
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Quant Title : SW846 8260

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

Peak Number 15 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.963	34.84 ug/l	370501	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, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
3			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
4			o-Cymene	134	C10H14	000527-84-4	95
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110722\
Data File : VX032631.D
Acq On : 07 Nov 2022 16:48
Operator : JC/MD
Sample : N5484-25
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
DUP110422

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.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-	1.745	74.4	ug/l	304873	1	5.556	204779	50.0
Pentane	1.953	54.3	ug/l	222351	1	5.556	204779	50.0
Butane, 2,3-dim...	2.782	60.6	ug/l	248286	1	5.556	204779	50.0
Pentane, 2-methyl-	2.825	132.5	ug/l	542599	1	5.556	204779	50.0
Pentane, 3-methyl-	3.099	118.3	ug/l	484635	1	5.556	204779	50.0
n-Hexane	3.446	64.0	ug/l	262251	1	5.556	204779	50.0
Cyclopentane, m...	4.306	123.1	ug/l	504138	1	5.556	204779	50.0
Hexane, 3-methyl-	5.830	34.5	ug/l	141096	1	5.556	204779	50.0
Benzene, 1-ethy...	11.396	37.3	ug/l	396643	4	12.024	531677	50.0
Benzene, 1-ethy...	11.616	37.4	ug/l	397966	4	12.024	531677	50.0
Benzene, 1,2,3-...	12.085	36.3	ug/l	386328	4	12.024	531677	50.0
Benzene, 2-ethy...	12.530	28.7	ug/l	304745	4	12.024	531677	50.0
Benzene, 1-ethy...	12.616	44.9	ug/l	477510	4	12.024	531677	50.0
Benzene, 1,2,3,...	12.920	29.3	ug/l	311176	4	12.024	531677	50.0
Benzene, 1,2,4,...	12.963	34.8	ug/l	370501	4	12.024	531677	50.0