

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070623\
 Data File : VX036442.D
 Acq On : 06 Jul 2023 18:16
 Operator : JC/MD
 Sample : 03500-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-1

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\82X062323W.M
 Title : SW846 8260

Signal : TIC: VX036442.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	52	rBV	439720	477402	4.88%	0.671%
2	1.745	103	108	123	rVB	7608758	9788093	100.00%	13.757%
3	1.953	136	142	157	rVB	1447677	1930271	19.72%	2.713%
4	2.215	180	185	194	rVB	80806	119505	1.22%	0.168%
5	2.337	196	205	225	rVB	677378	1338729	13.68%	1.882%
6	2.782	270	278	281	rBV	1038118	2188740	22.36%	3.076%
7	2.824	281	285	309	rVB	2198085	5078912	51.89%	7.138%
8	3.105	319	331	344	rBV	1621379	3607847	36.86%	5.071%
9	3.446	376	387	401	rBV	400471	886768	9.06%	1.246%
10	3.733	426	434	442	rVB	65553	150847	1.54%	0.212%
11	3.836	443	451	459	rBV3	50421	128175	1.31%	0.180%
12	4.056	474	487	491	rBV3	43557	134143	1.37%	0.189%
13	4.214	502	513	519	rBV2	202068	588094	6.01%	0.827%
14	4.306	519	528	541	rVV	857636	2432083	24.85%	3.418%
15	4.434	541	549	565	rVB6	33887	132326	1.35%	0.186%
16	5.135	645	664	689	rVB6	68374	352765	3.60%	0.496%
17	5.391	689	706	711	rBV2	113757	352924	3.61%	0.496%
18	5.482	711	721	730	rBV4	302869	1294481	13.23%	1.819%
19	5.623	737	744	765	rVB2	534936	1755950	17.94%	2.468%
20	5.830	765	778	790	rBV4	354281	1101345	11.25%	1.548%
21	5.958	790	799	808	rBV2	111791	282895	2.89%	0.398%
22	6.165	817	833	840	rBV2	213079	748820	7.65%	1.052%
23	6.257	841	848	857	rVV3	342145	1032297	10.55%	1.451%
24	6.360	857	865	877	rVB2	191575	540949	5.53%	0.760%
25	6.610	891	906	918	rBV2	91806	230127	2.35%	0.323%
26	6.763	922	931	938	rBV	288639	704413	7.20%	0.990%
27	6.848	938	945	955	rVB3	97650	296272	3.03%	0.416%
28	7.171	991	998	1006	rBV3	56252	124031	1.27%	0.174%
29	7.379	1020	1032	1044	rBV3	665012	1703076	17.40%	2.394%
30	7.659	1072	1078	1088	rVV	143843	286590	2.93%	0.403%
31	7.775	1090	1097	1104	rVB2	61778	127233	1.30%	0.179%
32	7.988	1122	1132	1143	rVB	144648	354776	3.62%	0.499%
33	8.110	1143	1152	1155	rBV	238603	498912	5.10%	0.701%
34	8.141	1155	1157	1162	rVB	160113	236763	2.42%	0.333%
35	8.433	1200	1205	1210	rBV2	71133	120781	1.23%	0.170%
36	8.506	1212	1217	1227	rVB3	176928	332381	3.40%	0.467%

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37	8.604	1227	1233	1236	rBV3	115918	234772	2.40%	0.330%
38	8.647	1236	1240	1249	rVB	583619	958139	9.79%	1.347%
39	8.842	1265	1272	1280	rVV2	95190	214592	2.19%	0.302%
40	8.921	1280	1285	1290	rVV	95703	151558	1.55%	0.213%
41	9.104	1306	1315	1320	rBV	83885	163744	1.67%	0.230%
42	9.165	1320	1325	1336	rVB	208880	363272	3.71%	0.511%
43	9.451	1364	1372	1376	rBV5	55139	148047	1.51%	0.208%
44	9.506	1376	1381	1386	rVV3	57460	114963	1.17%	0.162%
45	10.024	1460	1466	1467	rBV2	86600	126508	1.29%	0.178%
46	10.055	1467	1471	1476	rVB	597878	773913	7.91%	1.088%
47	10.146	1481	1486	1490	rVV2	139089	223288	2.28%	0.314%
48	10.195	1490	1494	1498	rVV	283106	346100	3.54%	0.486%
49	10.305	1508	1512	1517	rVB	95620	140167	1.43%	0.197%
50	10.963	1614	1620	1628	rBV	764361	946628	9.67%	1.330%
51	11.079	1635	1639	1644	rVB	471006	576810	5.89%	0.811%
52	11.305	1671	1676	1685	rVB	2236601	2659923	27.18%	3.738%
53	11.402	1685	1692	1695	rVV	171061	241658	2.47%	0.340%
54	11.451	1695	1700	1708	rVB3	51868	100572	1.03%	0.141%
55	11.616	1720	1727	1735	rBV	227013	315392	3.22%	0.443%
56	11.865	1762	1768	1770	rBV	221210	287876	2.94%	0.405%
57	11.890	1770	1772	1778	rVB	237006	261919	2.68%	0.368%
58	12.024	1788	1794	1800	rVB	596475	789197	8.06%	1.109%
59	12.085	1800	1804	1810	rBV	74325	113820	1.16%	0.160%
60	12.243	1824	1830	1837	rBV2	1658157	2093632	21.39%	2.943%
61	12.317	1837	1842	1843	rBV	359115	467008	4.77%	0.656%
62	12.402	1851	1856	1861	rBV	171961	215699	2.20%	0.303%
63	12.463	1861	1866	1873	rVB	213787	245823	2.51%	0.345%
64	12.615	1885	1891	1894	rBV	1820341	2255821	23.05%	3.170%
65	12.646	1894	1896	1900	rVV	324236	325847	3.33%	0.458%
66	12.701	1900	1905	1912	rVV2	1160698	1586825	16.21%	2.230%
67	12.920	1933	1941	1946	rBV	1358220	1687146	17.24%	2.371%
68	13.024	1953	1958	1967	rBV3	137163	237154	2.42%	0.333%
69	13.140	1973	1977	1978	rBV	95034	108646	1.11%	0.153%
70	13.170	1978	1982	1992	rVB	944046	1235593	12.62%	1.737%
71	13.292	1997	2002	2008	rVV2	1645840	2198251	22.46%	3.090%
72	13.371	2012	2015	2020	rVB	115301	136981	1.40%	0.193%
73	13.420	2020	2023	2032	rVB	187436	245445	2.51%	0.345%
74	13.505	2032	2037	2040	rBV2	181295	252927	2.58%	0.355%
75	13.542	2040	2043	2046	rVV	425718	550726	5.63%	0.774%
76	13.585	2046	2050	2054	rVV3	423570	645055	6.59%	0.907%

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77	13.640	2054	2059	2060	rVV2	216662	302354	3.09%	0.425%
78	13.664	2060	2063	2070	rVB2	434037	644681	6.59%	0.906%
79	13.774	2074	2081	2090	rBV	672933	917048	9.37%	1.289%
80	13.926	2099	2106	2110	rBV2	198132	294561	3.01%	0.414%
81	13.969	2110	2113	2117	rVB	138934	161828	1.65%	0.227%
82	14.072	2126	2130	2135	rBV	329876	417658	4.27%	0.587%
83	14.213	2148	2153	2163	rBV	250552	440474	4.50%	0.619%
84	14.322	2167	2171	2176	rVV	129076	196727	2.01%	0.276%
85	14.371	2176	2179	2184	rVB	196659	236555	2.42%	0.332%
86	14.633	2218	2222	2228	rVB	658777	812646	8.30%	1.142%
87	14.780	2241	2246	2255	rBV	404353	524980	5.36%	0.738%

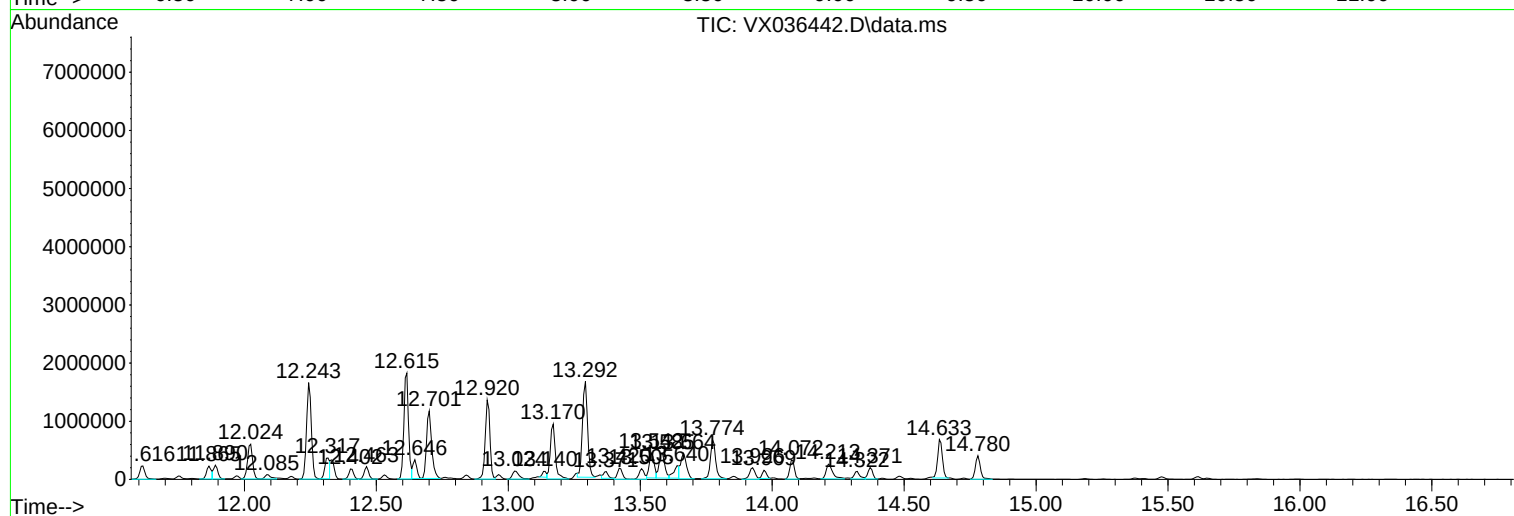
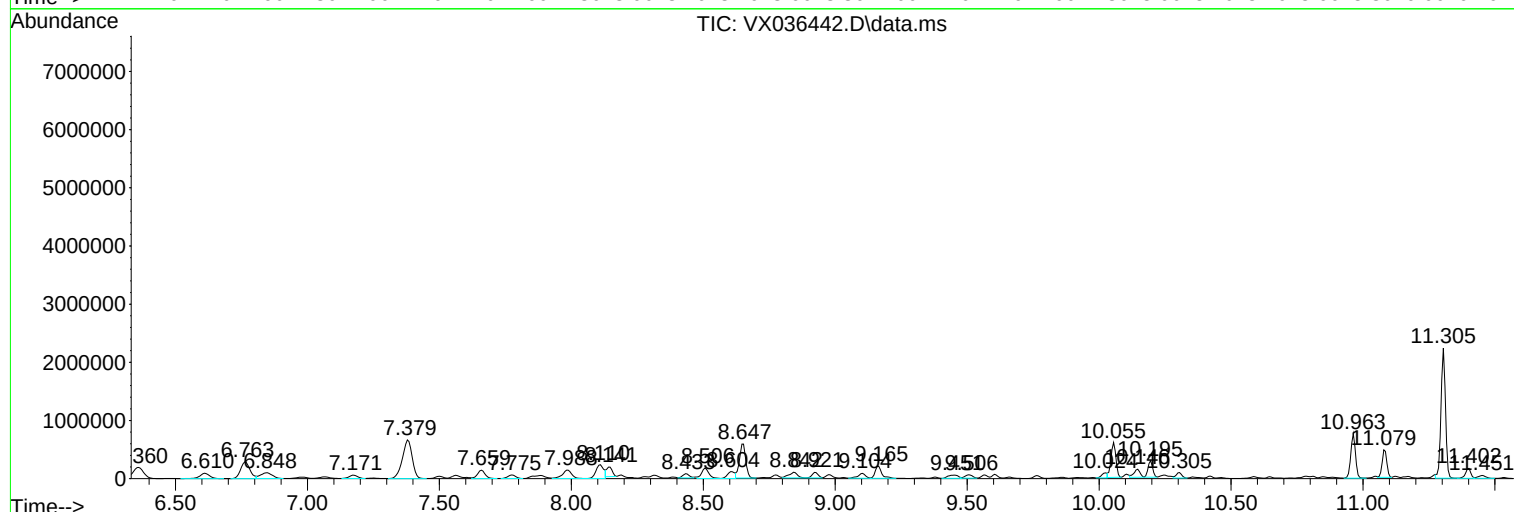
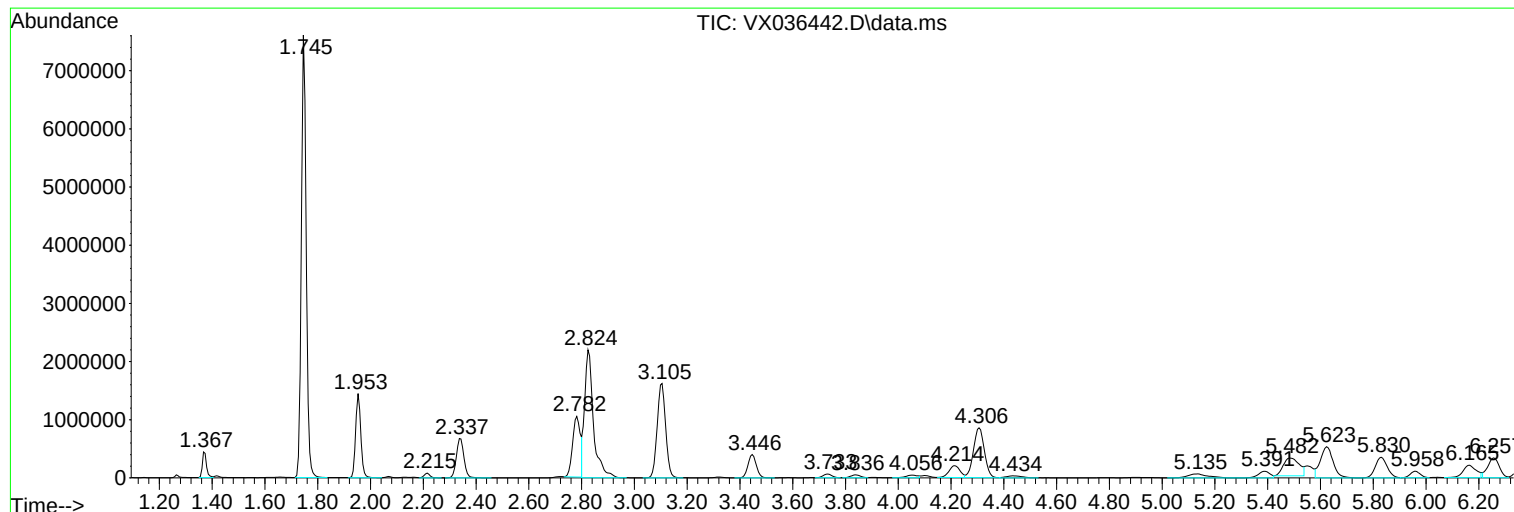
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062323W.M
Quant Title : SW846 8260

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



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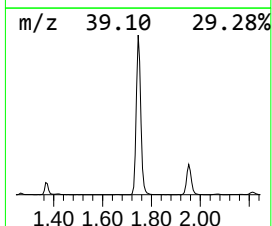
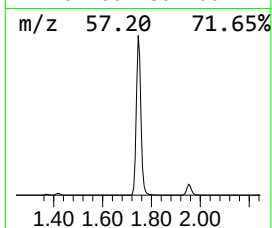
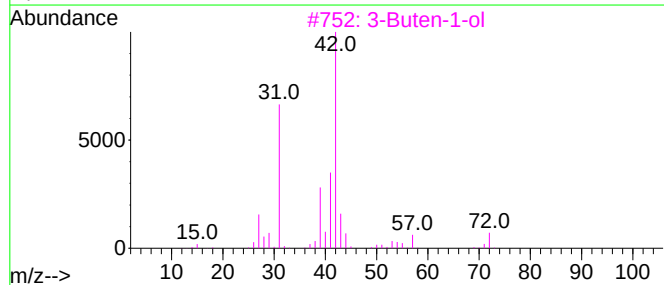
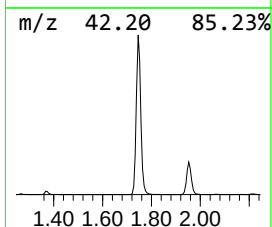
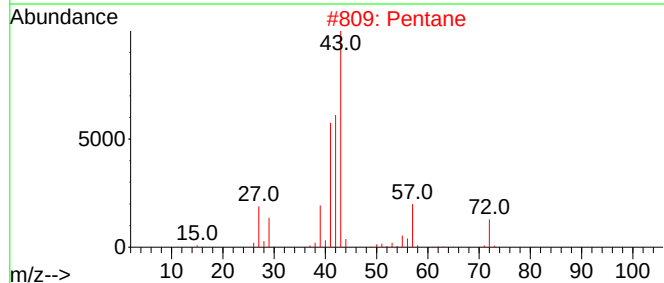
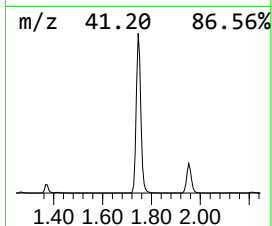
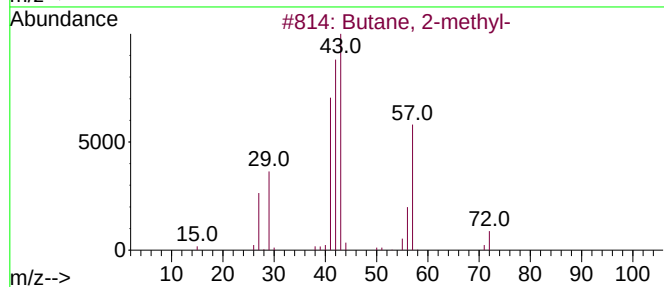
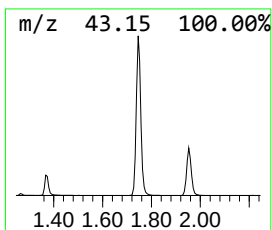
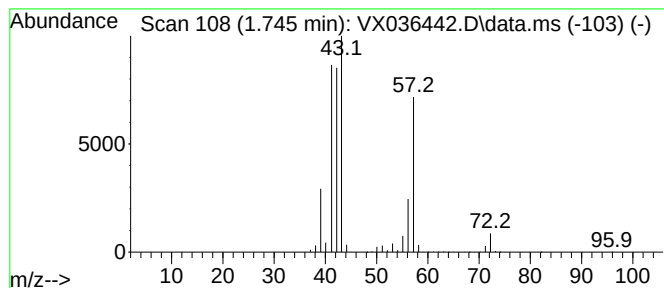
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062323W.M
Quant Title : SW846 8260

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.
1.745	278.71 ug/l	9788090	Pentafluorobenzene	5.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methyl-	72	C5H12	000078-78-4	91
2			Pentane	72	C5H12	000109-66-0	37
3			3-Buten-1-ol	72	C4H8O	000627-27-0	28
4			1-Butene	56	C4H8	000106-98-9	10
5			Methane, isocyanato-	57	C2H3NO	000624-83-9	9



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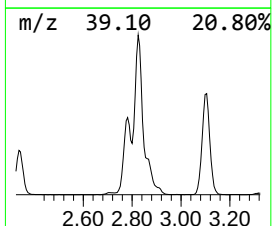
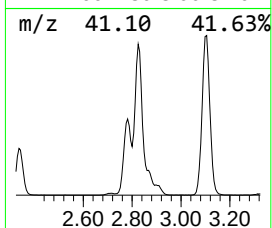
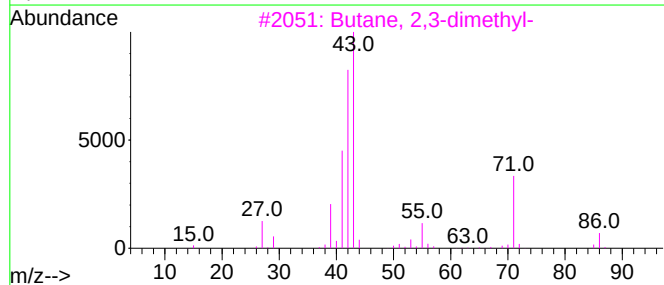
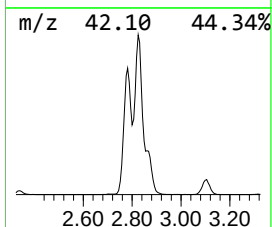
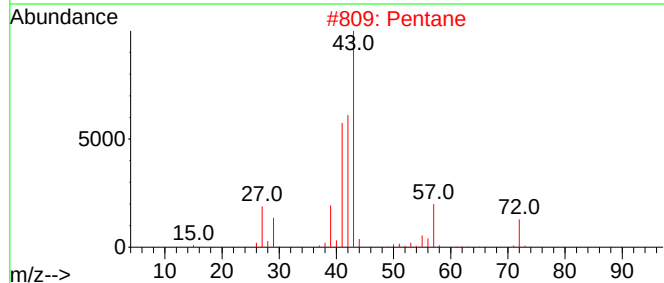
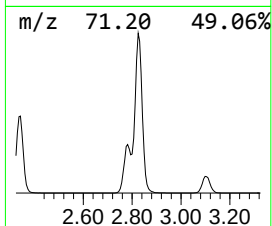
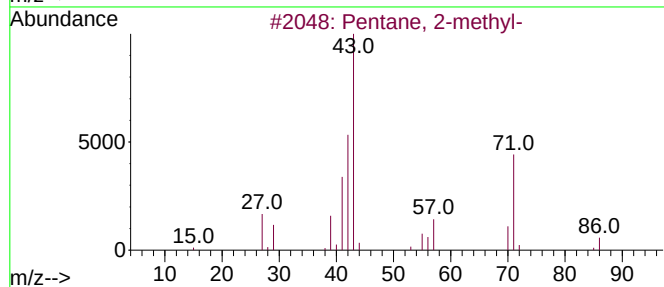
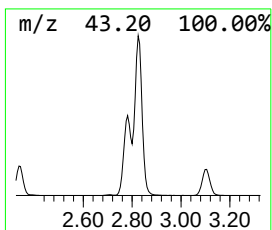
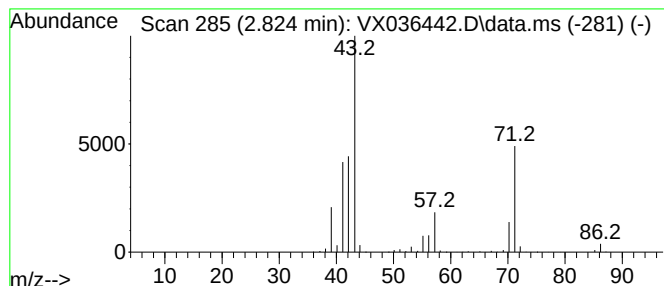
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062323W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.824	144.62 ug/l	5078910	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			Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	43
4			Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	33
5			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	25



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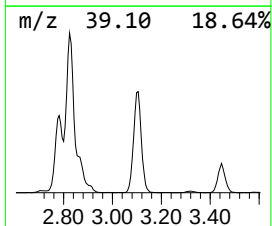
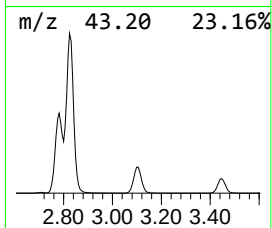
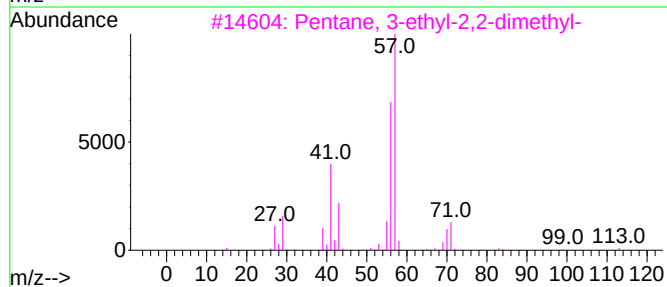
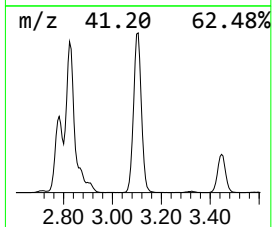
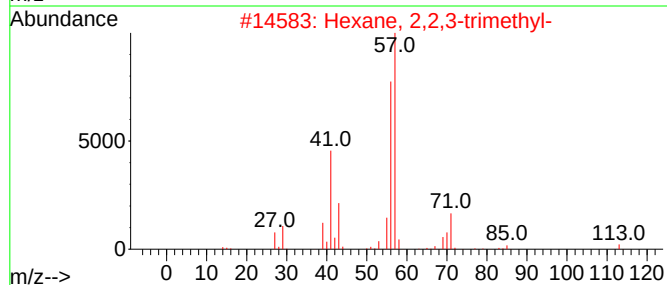
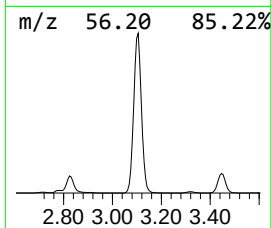
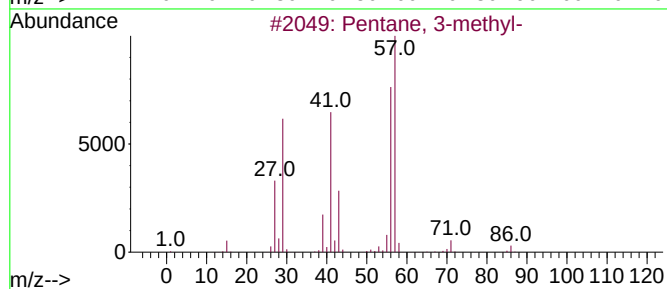
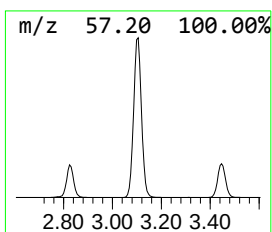
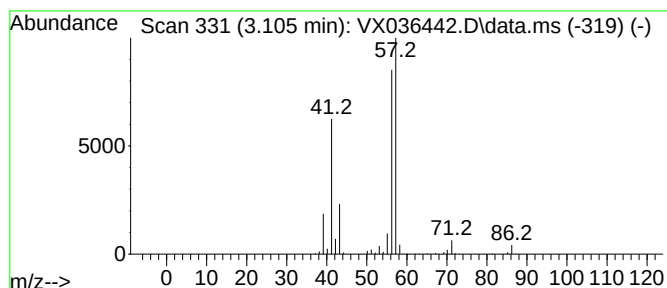
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062323W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.105	102.73 ug/l	3607850	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	83
3			Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	64
4			Sulphuric acid dibutyl ester	210	C8H18O4S	000625-22-9	40
5			Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	38



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ClientSampleId :
MW-1

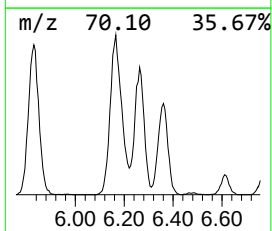
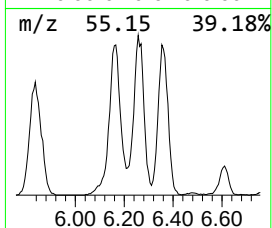
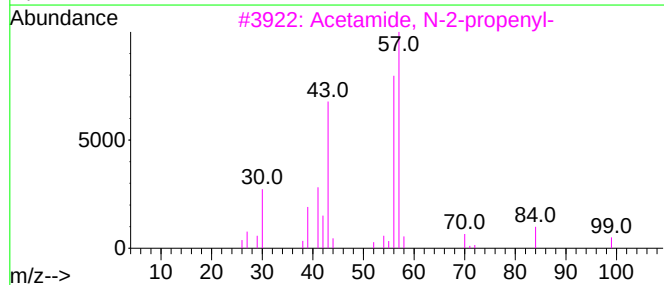
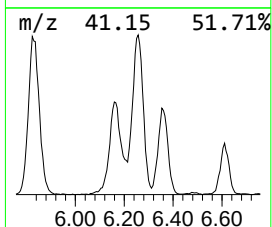
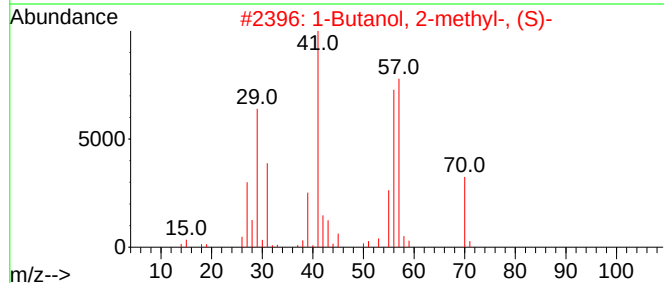
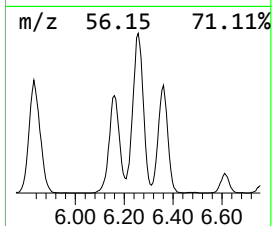
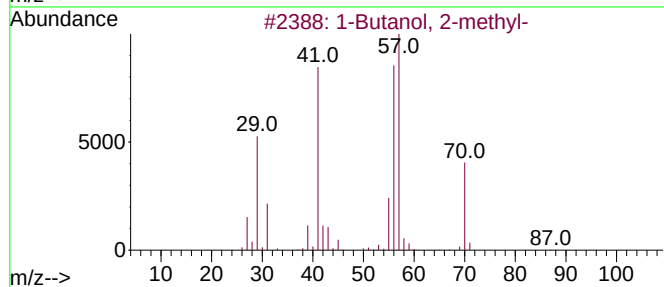
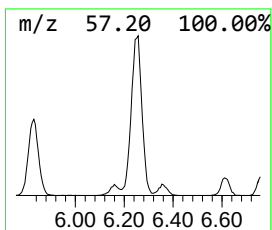
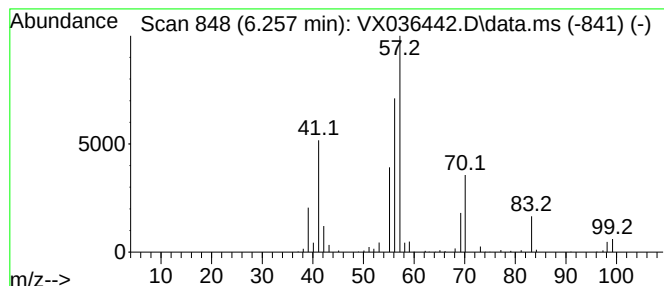
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Quant Title : SW846 8260

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

Peak Number 4 1-Butanol, 2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.257	73.27 ug/l	1032300	1,4-Difluorobenzene	6.763

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Butanol, 2-methyl-			88	C5H12O	000137-32-6	53
2	1-Butanol, 2-methyl-, (S)-			88	C5H12O	001565-80-6	50
3	Acetamide, N-2-propenyl-			99	C5H9NO	000692-33-1	47
4	Cyclopentane, 1,2-dimethyl-, cis-			98	C7H14	001192-18-3	47
5	1-Hexene, 5-methyl-			98	C7H14	003524-73-0	45



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070623\
Data File : VX036442.D
Acq On : 06 Jul 2023 18:16
Operator : JC/MD
Sample : 03500-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-1

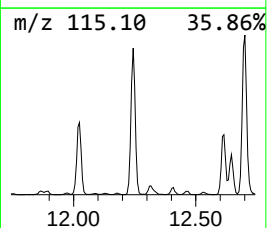
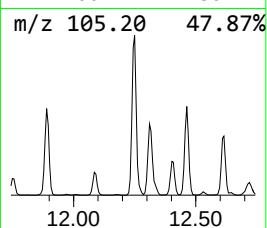
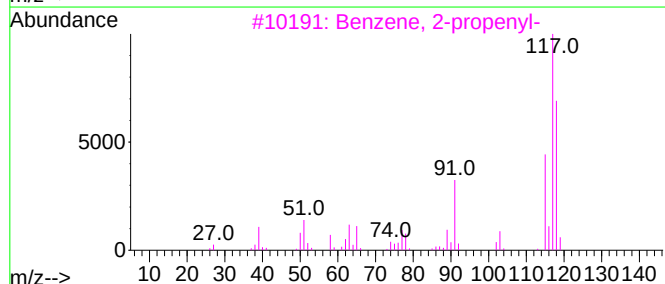
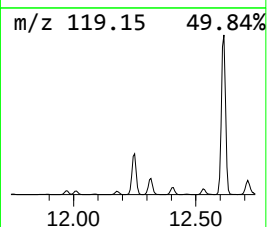
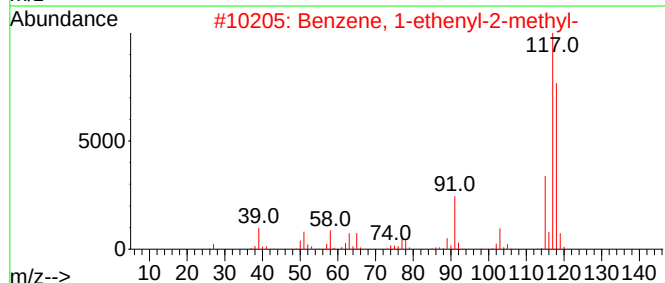
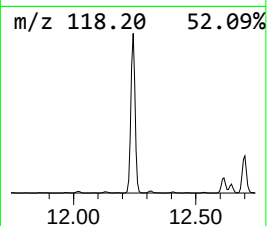
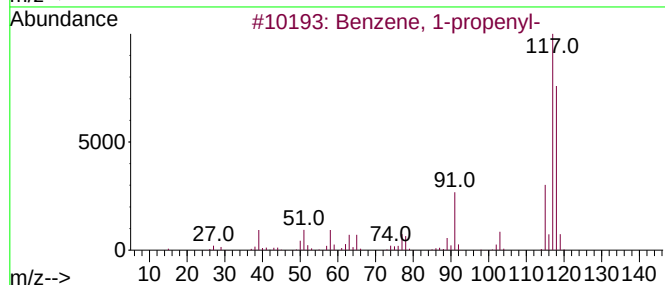
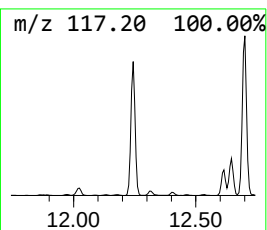
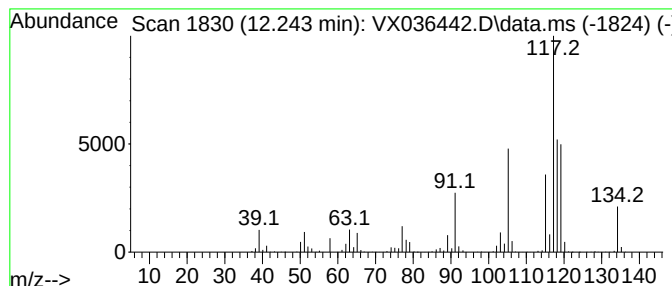
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1-propenyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.243	132.64 ug/l	2093630	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-propenyl-	118	C9H10	000637-50-3	60
2			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	60
3			Benzene, 2-propenyl-	118	C9H10	000300-57-2	60
4			Benzene, cyclopropyl-	118	C9H10	000873-49-4	60
5			Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	55



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070623\
Data File : VX036442.D
Acq On : 06 Jul 2023 18:16
Operator : JC/MD
Sample : 03500-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-1

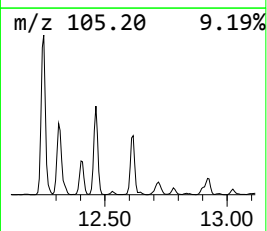
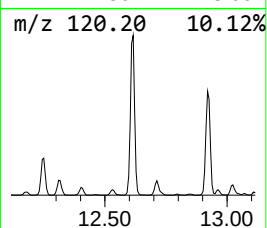
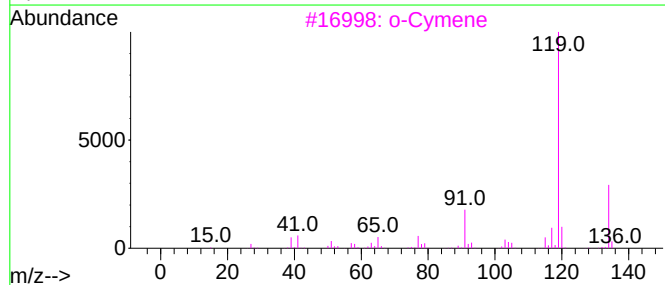
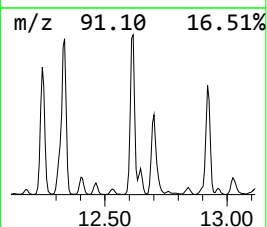
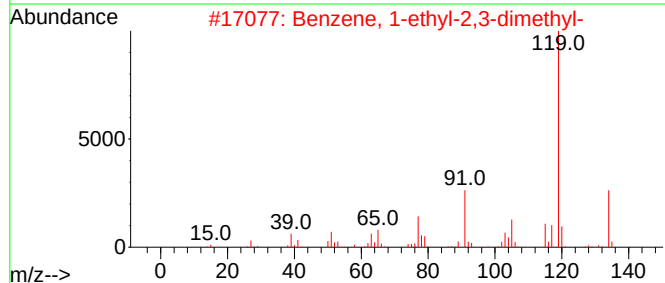
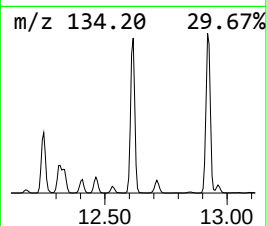
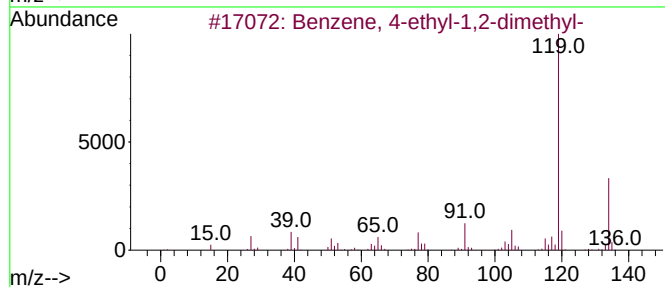
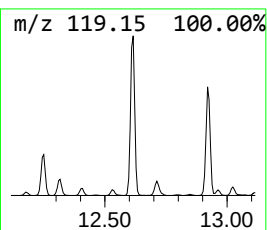
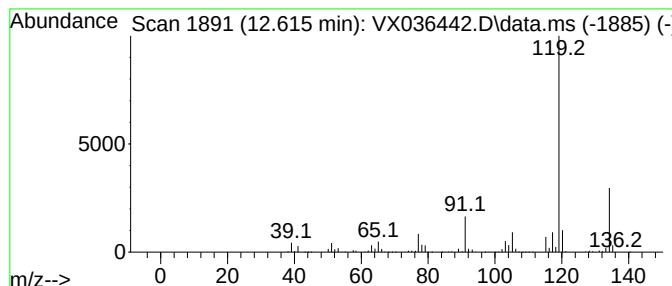
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.615	142.92 ug/l	2255820	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
2			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
3			o-Cymene	134	C10H14	000527-84-4	95
4			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
5			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070623\
Data File : VX036442.D
Acq On : 06 Jul 2023 18:16
Operator : JC/MD
Sample : 03500-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-1

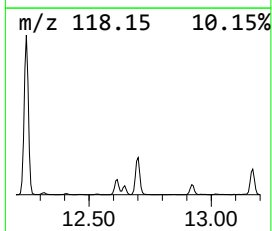
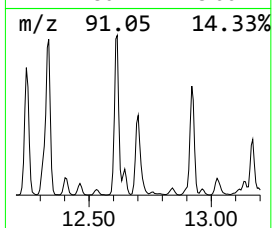
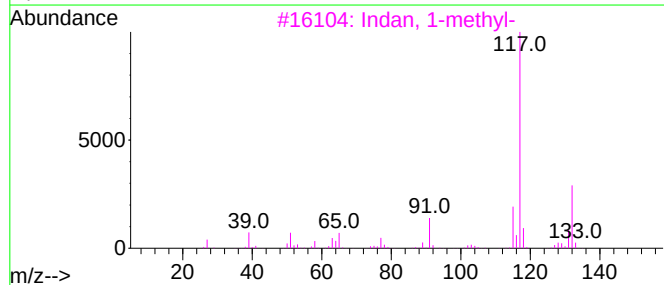
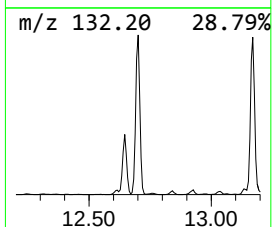
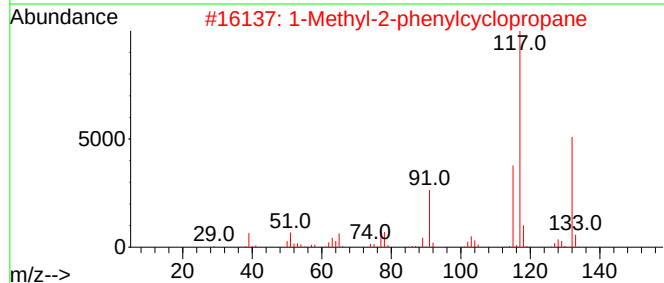
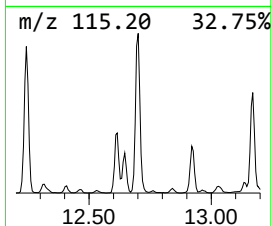
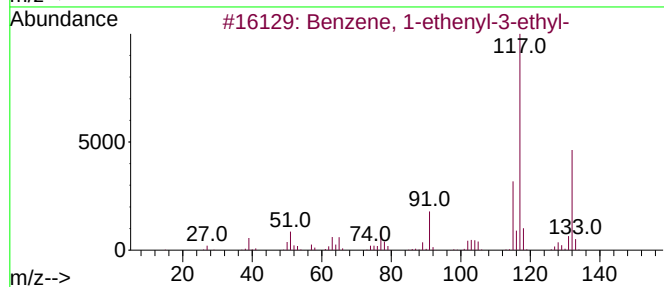
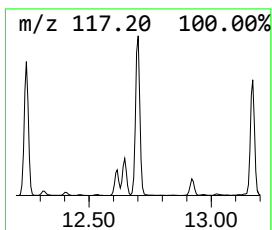
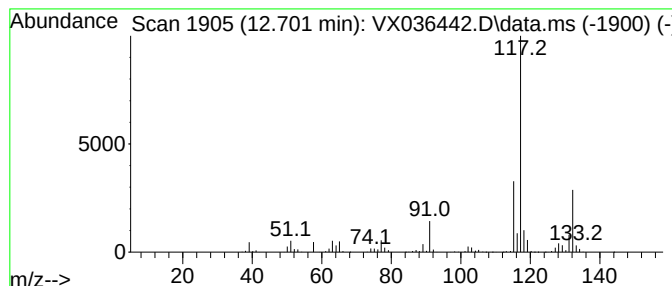
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.701	100.53 ug/l	1586830	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	90
2			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	87
3			Indan, 1-methyl-	132	C10H12	000767-58-8	87
4			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	83
5			Benzene, 2-butenyl-	132	C10H12	001560-06-1	83



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070623\
Data File : VX036442.D
Acq On : 06 Jul 2023 18:16
Operator : JC/MD
Sample : 03500-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-1

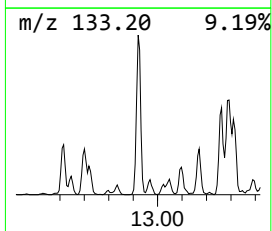
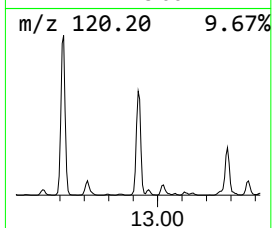
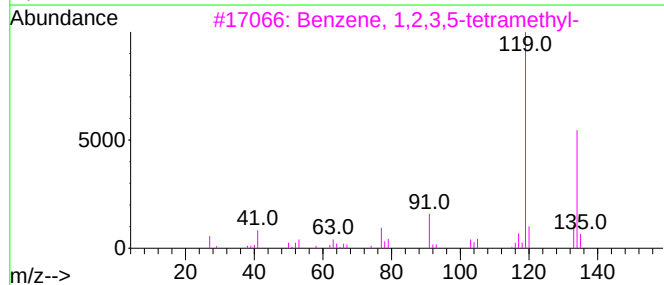
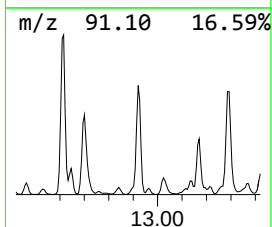
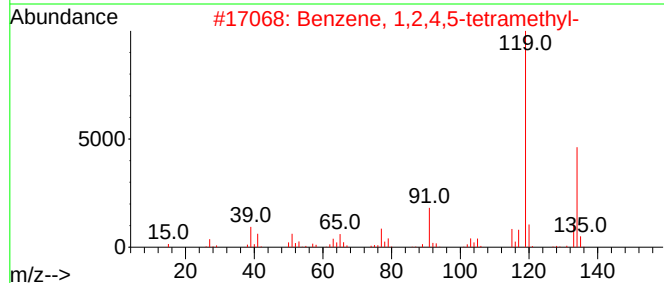
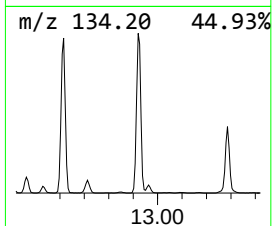
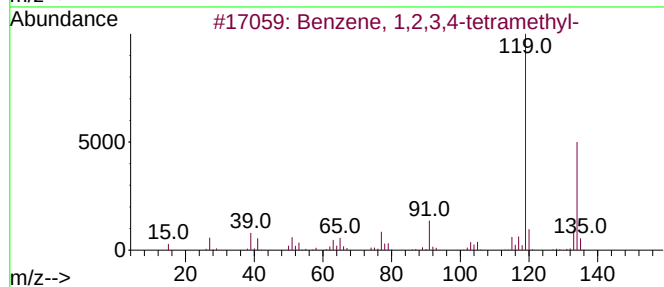
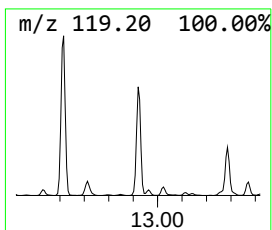
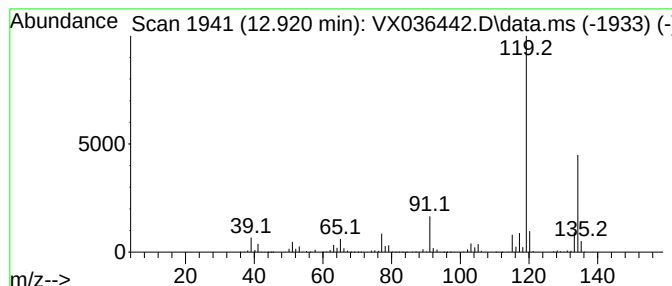
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.920	106.89 ug/l	1687150	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	93



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070623\
Data File : VX036442.D
Acq On : 06 Jul 2023 18:16
Operator : JC/MD
Sample : 03500-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-1

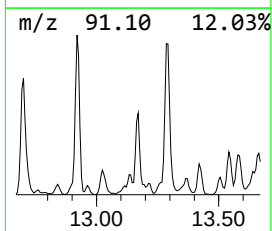
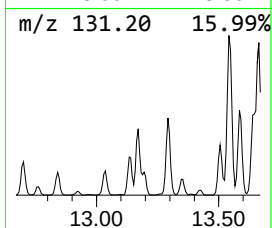
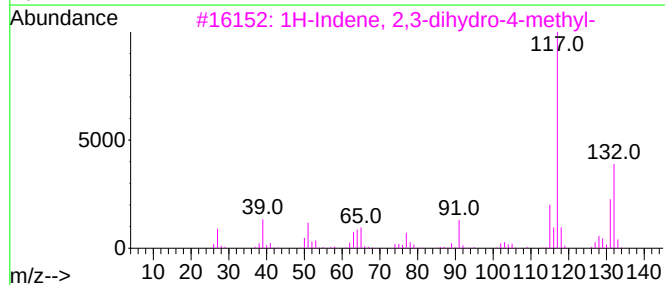
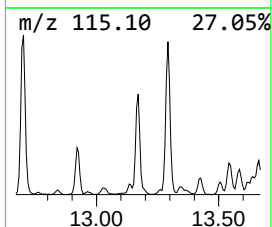
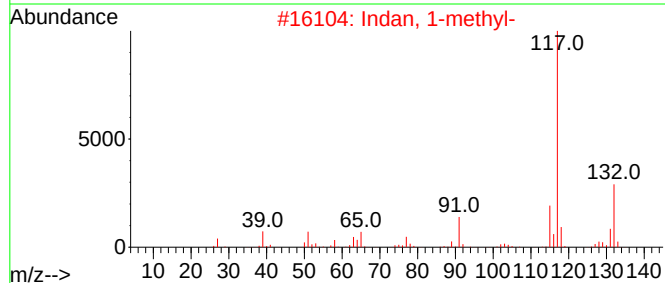
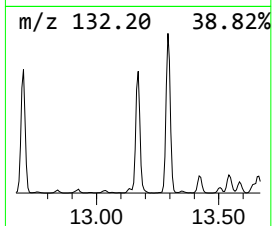
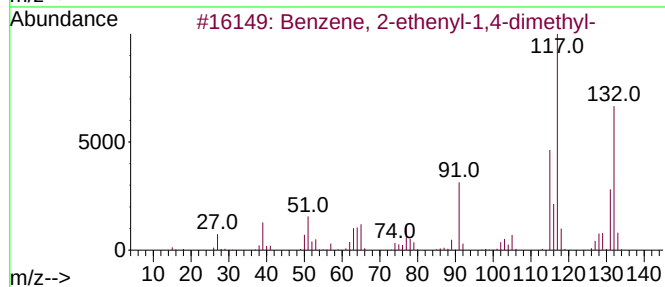
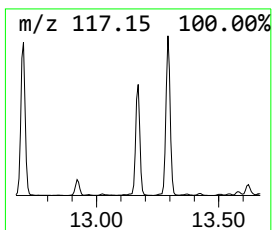
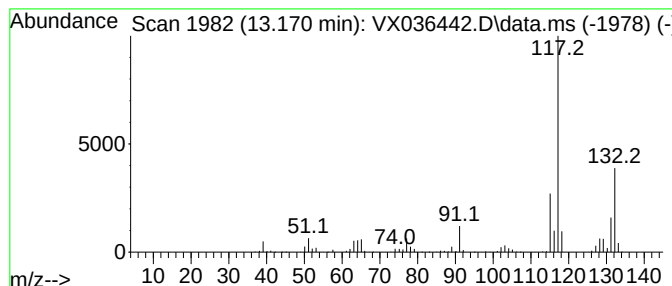
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.170	78.28 ug/l	1235590	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92
2			Indan, 1-methyl-	132	C10H12	000767-58-8	91
3			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90
4			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87
5			1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070623\
Data File : VX036442.D
Acq On : 06 Jul 2023 18:16
Operator : JC/MD
Sample : 03500-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-1

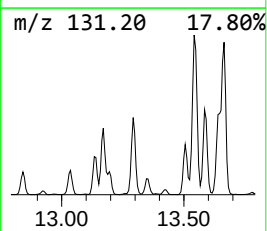
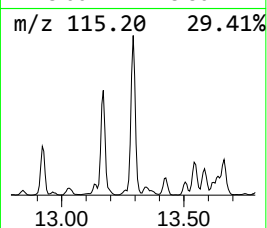
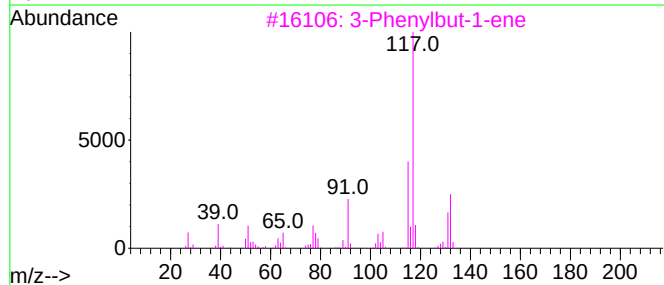
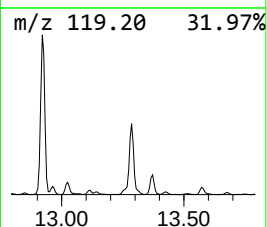
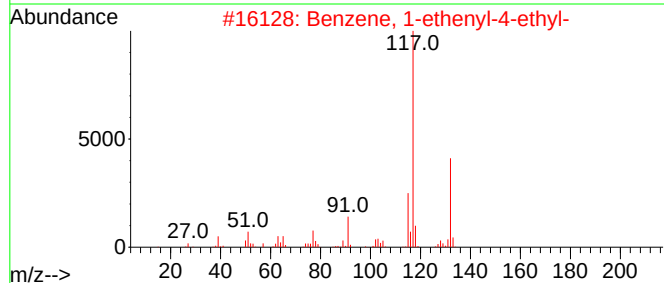
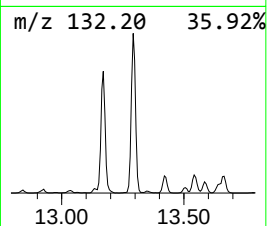
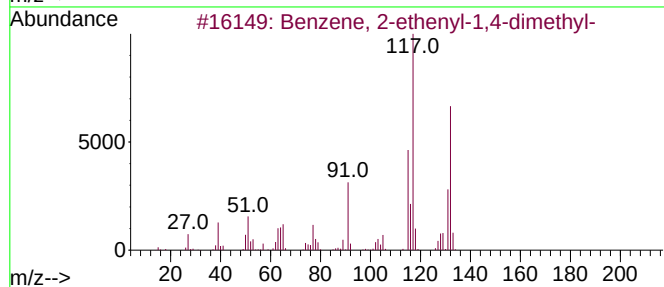
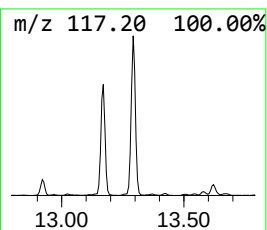
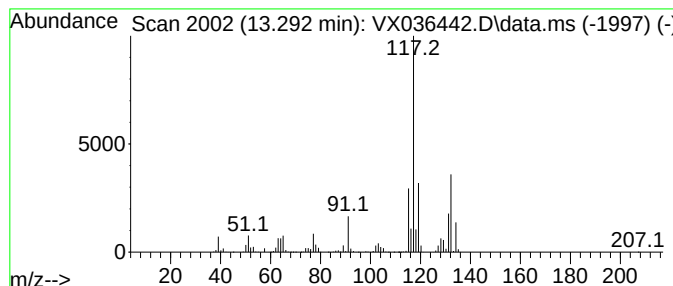
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062323W.M
Quant Title : SW846 8260

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

Peak Number 10 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	139.27 ug/l	2198250	1,4-Dichlorobenzene-d4	12.024

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92
2		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	89
3		3-Phenylbut-1-ene	132	C10H12	000934-10-1	81
4		Indan, 1-methyl-	132	C10H12	000767-58-8	81
5		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	76



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070623\
Data File : VX036442.D
Acq On : 06 Jul 2023 18:16
Operator : JC/MD
Sample : 03500-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062323W.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	278.7	ug/l	9788090	1	5.556	1755950	50.0
Pentane, 2-methyl-	2.824	144.6	ug/l	5078910	1	5.556	1755950	50.0
Pentane, 3-methyl-	3.105	102.7	ug/l	3607850	1	5.556	1755950	50.0
1-Butanol, 2-me...	6.257	73.3	ug/l	1032300	2	6.763	704413	50.0
Benzene, 1-prop...	12.243	132.6	ug/l	2093630	4	12.024	789197	50.0
Benzene, 4-ethy...	12.615	142.9	ug/l	2255820	4	12.024	789197	50.0
Benzene, 1-ethe...	12.701	100.5	ug/l	1586830	4	12.024	789197	50.0
Benzene, 1,2,3,...	12.920	106.9	ug/l	1687150	4	12.024	789197	50.0
Benzene, 2-ethe...	13.170	78.3	ug/l	1235590	4	12.024	789197	50.0
Benzene, 1-ethe...	13.292	139.3	ug/l	2198250	4	12.024	789197	50.0