

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX102423\
 Data File : VX038456.D
 Acq On : 24 Oct 2023 15:31
 Operator : JC/MD
 Sample : 04975-01
 Misc : 1.12g/10mL/100uL/5.00mL/MSVOA_X/MEOH
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 54

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\82X100523W.M
 Title : SW846 8260

Signal : TIC: VX038456.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.551	66	76	77	rBV3	25342	51313	5.38%	0.545%
2	5.385	692	705	716	rBV	94115	278827	29.22%	2.964%
3	5.544	716	731	751	rVB	159287	529491	55.48%	5.628%
4	5.806	763	774	788	rVB	30425	88338	9.26%	0.939%
5	5.958	788	799	821	rVB	102515	274452	28.76%	2.917%
6	6.598	895	904	921	rVB2	31005	77975	8.17%	0.829%
7	6.757	921	930	954	rBV	249124	611097	64.03%	6.495%
8	7.220	997	1006	1015	rBV6	3324	10413	1.09%	0.111%
9	7.373	1020	1031	1038	rBV4	7618	18700	1.96%	0.199%
10	8.647	1234	1240	1246	rBV	533194	836444	87.65%	8.890%
11	8.720	1246	1252	1272	rVB	511602	837456	87.75%	8.901%
12	10.055	1466	1471	1486	rVB	561714	794618	83.26%	8.446%
13	10.305	1501	1512	1517	rBV2	16448	30151	3.16%	0.320%
14	10.647	1564	1568	1576	rVB2	6473	14095	1.48%	0.150%
15	10.781	1586	1590	1595	rVB2	7883	10017	1.05%	0.106%
16	10.854	1595	1602	1607	rBV3	7761	15364	1.61%	0.163%
17	11.085	1634	1640	1651	rBV	516217	696255	72.96%	7.400%
18	11.189	1653	1657	1660	rBV3	7194	10120	1.06%	0.108%
19	11.305	1673	1676	1684	rVB	9296	15642	1.64%	0.166%
20	11.384	1684	1689	1694	rBV	21733	39915	4.18%	0.424%
21	11.616	1723	1727	1734	rVB3	15589	27627	2.89%	0.294%
22	11.683	1734	1738	1740	rBV2	8604	10762	1.13%	0.114%
23	11.713	1740	1743	1746	rBV2	17629	21430	2.25%	0.228%
24	11.750	1746	1749	1759	rVB2	60142	97010	10.17%	1.031%
25	11.841	1759	1764	1768	rBV	14981	24925	2.61%	0.265%
26	11.890	1768	1772	1775	rBV3	11202	13307	1.39%	0.141%
27	11.939	1775	1780	1783	rBV2	31319	47973	5.03%	0.510%
28	11.976	1783	1786	1789	rVB3	11075	13083	1.37%	0.139%
29	12.024	1789	1794	1799	rBV	765238	954340	100.00%	10.143%
30	12.073	1799	1802	1804	rBV2	9942	13047	1.37%	0.139%
31	12.146	1812	1814	1817	rVB	14562	13294	1.39%	0.141%
32	12.244	1824	1830	1832	rBV3	40733	75878	7.95%	0.806%
33	12.317	1837	1842	1849	rVV5	48917	114182	11.96%	1.214%
34	12.421	1849	1859	1863	rVV2	56124	88189	9.24%	0.937%
35	12.463	1863	1866	1873	rVB3	30403	52012	5.45%	0.553%
36	12.536	1873	1878	1879	rBV	21074	28970	3.04%	0.308%

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37	12.555	1879	1881	1884	rVV2	21942	28869	3.03%	0.307%
38	12.610	1888	1890	1893	rVB	13947	14947	1.57%	0.159%
39	12.683	1899	1902	1907	rBV3	19964	42335	4.44%	0.450%
40	12.732	1907	1910	1913	rVV3	18676	27285	2.86%	0.290%
41	12.817	1917	1924	1928	rVV5	29326	49565	5.19%	0.527%
42	12.890	1931	1936	1943	rBV5	34762	89968	9.43%	0.956%
43	12.975	1943	1950	1955	rVB5	31201	77002	8.07%	0.818%
44	13.042	1955	1961	1964	rBV4	30473	42845	4.49%	0.455%
45	13.164	1974	1981	1985	rBV3	41200	70499	7.39%	0.749%
46	13.213	1985	1989	1991	rVV2	23031	25619	2.68%	0.272%
47	13.268	1993	1998	2000	rVV	114325	159573	16.72%	1.696%
48	13.292	2000	2002	2007	rVV	105557	129765	13.60%	1.379%
49	13.341	2007	2010	2013	rVV3	13840	19434	2.04%	0.207%
50	13.384	2013	2017	2020	rVV2	34697	48532	5.09%	0.516%
51	13.420	2020	2023	2032	rVB	77604	135415	14.19%	1.439%
52	13.500	2032	2036	2040	rBV3	22501	36140	3.79%	0.384%
53	13.542	2040	2043	2046	rBV2	19164	24566	2.57%	0.261%
54	13.585	2046	2050	2055	rBV4	34227	56305	5.90%	0.598%
55	13.676	2060	2065	2067	rVV3	20663	37645	3.94%	0.400%
56	13.737	2071	2075	2078	rVB4	21199	25696	2.69%	0.273%
57	13.798	2082	2085	2088	rVB3	11939	11540	1.21%	0.123%
58	13.847	2088	2093	2099	rVB5	52923	108928	11.41%	1.158%
59	13.926	2099	2106	2111	rBV4	41640	79495	8.33%	0.845%
60	13.969	2111	2113	2116	rBV	11130	11503	1.21%	0.122%
61	14.036	2120	2124	2127	rVB	95824	98311	10.30%	1.045%
62	14.073	2127	2130	2133	rBV2	36700	42995	4.51%	0.457%
63	14.225	2147	2155	2161	rVB2	109925	192004	20.12%	2.041%
64	14.317	2167	2170	2174	rBV4	14015	19872	2.08%	0.211%
65	14.371	2176	2179	2183	rVV	30777	35534	3.72%	0.378%
66	14.439	2183	2190	2193	rVV2	50032	71990	7.54%	0.765%
67	14.475	2193	2196	2206	rVB4	60913	116286	12.18%	1.236%
68	14.567	2206	2211	2215	rVB8	10663	21028	2.20%	0.223%
69	14.615	2215	2219	2221	rBV	83443	108581	11.38%	1.154%
70	14.670	2225	2228	2233	rVB3	35722	48105	5.04%	0.511%
71	14.725	2233	2237	2239	rBV2	17729	23969	2.51%	0.255%
72	14.756	2239	2242	2245	rVB	45363	43235	4.53%	0.460%
73	14.817	2250	2252	2254	rVV2	18783	22685	2.38%	0.241%
74	14.847	2254	2257	2260	rVV3	27635	33942	3.56%	0.361%
75	14.902	2263	2266	2277	rVB2	18717	35245	3.69%	0.375%
76	15.048	2285	2290	2293	rBV5	8516	14981	1.57%	0.159%

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77	15.182	2306	2312	2320	rBV3	46242	106423	11.15%	1.131%
78	15.274	2324	2327	2331	rVV5	13563	19707	2.06%	0.209%
79	15.316	2331	2334	2339	rVB6	7997	12941	1.36%	0.138%
80	15.377	2339	2344	2348	rBV6	14267	24108	2.53%	0.256%
81	15.487	2356	2362	2366	rBV3	27722	46792	4.90%	0.497%
82	15.560	2371	2374	2378	rVB	19139	23934	2.51%	0.254%
83	15.609	2378	2382	2386	rBV5	14624	21250	2.23%	0.226%
84	15.993	2440	2445	2454	rVB5	7114	24361	2.55%	0.259%
85	16.085	2457	2460	2465	rBV6	5031	10093	1.06%	0.107%

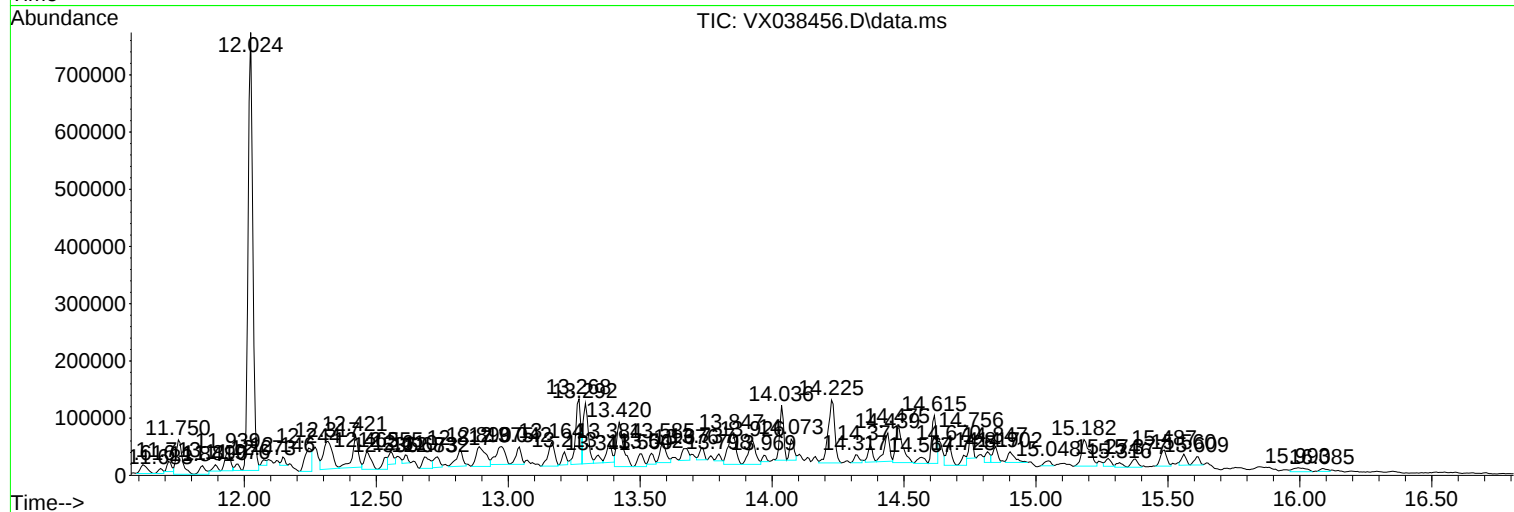
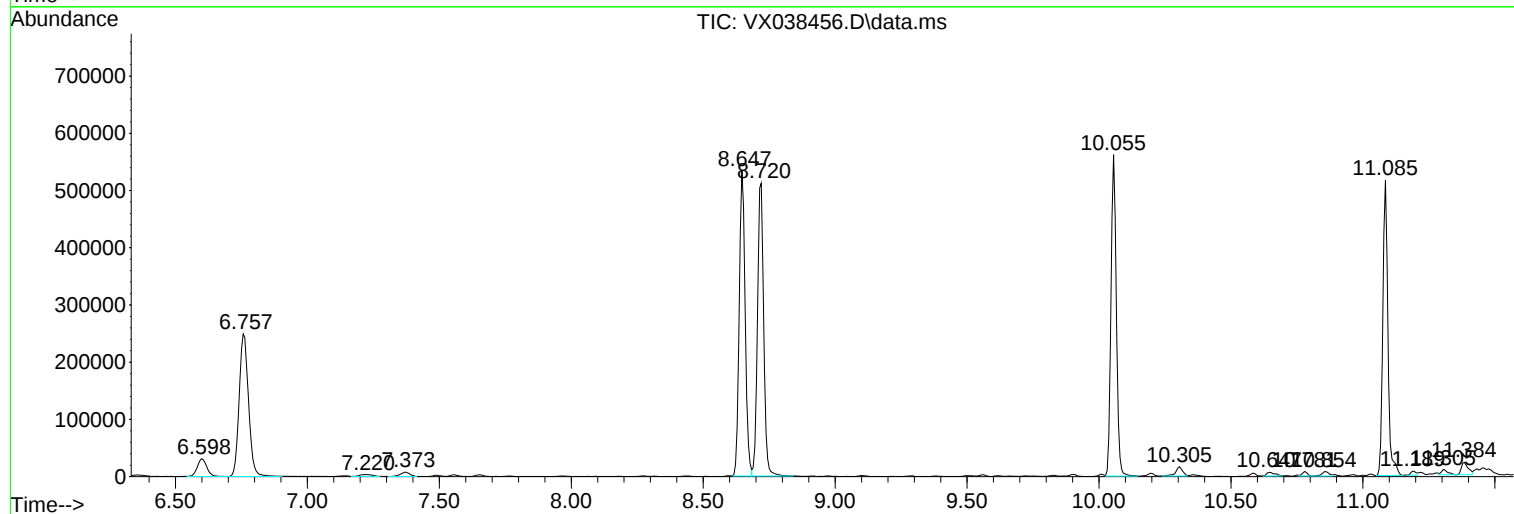
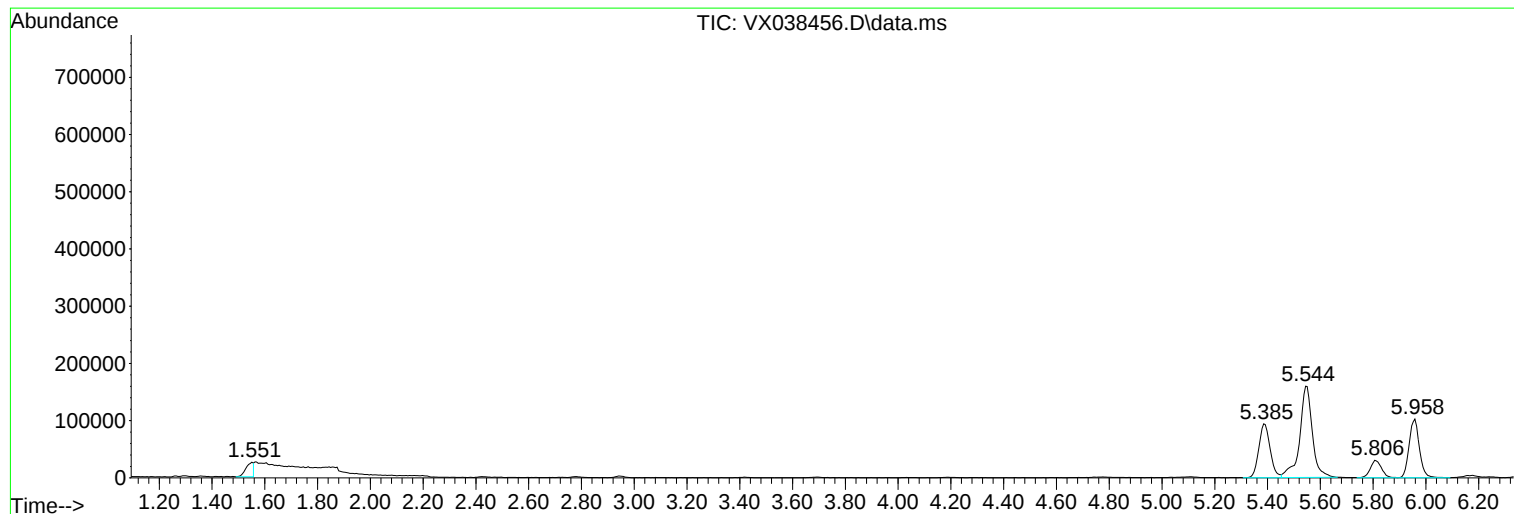
Sum of corrected areas: 9408555

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100523W.M
Quant Title : SW846 8260

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



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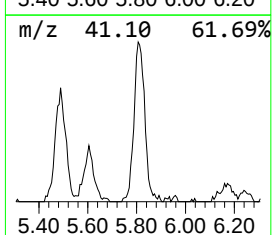
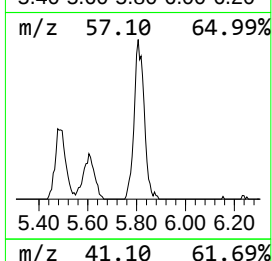
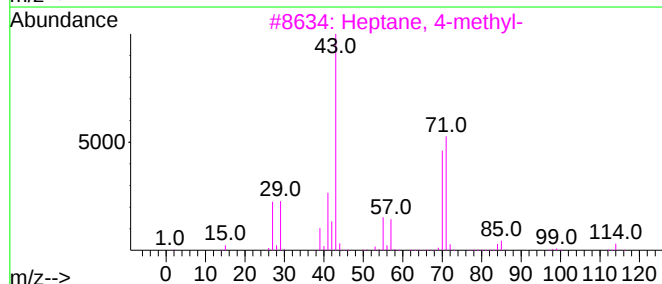
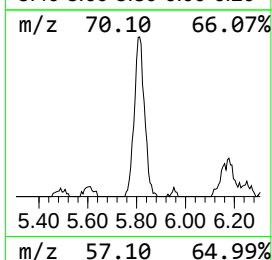
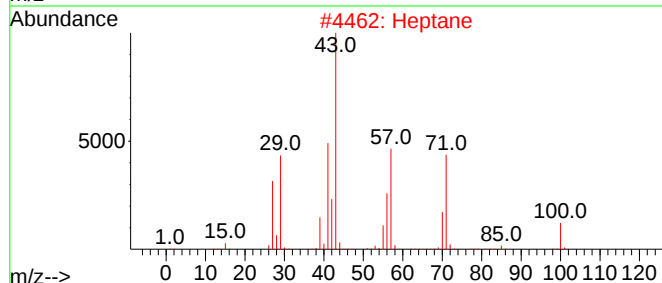
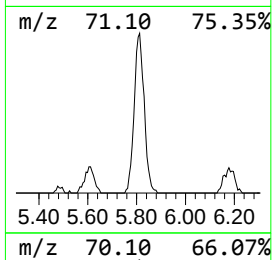
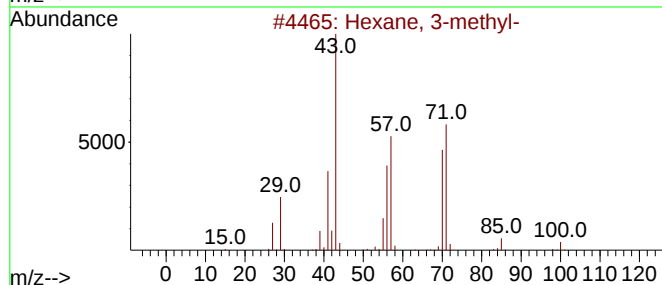
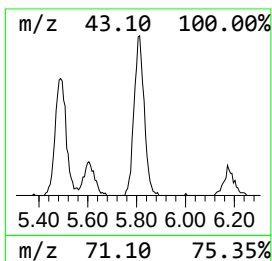
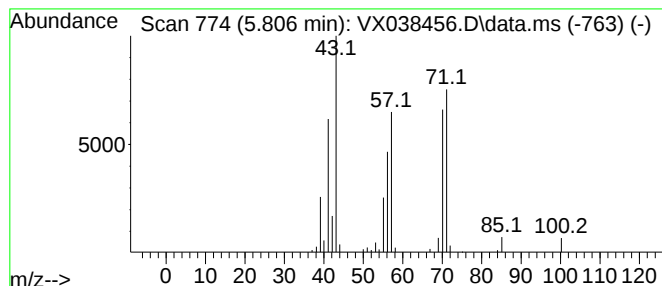
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100523W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Hexane, 3-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.806	8.34 ug/l	88338	Pentafluorobenzene	5.544

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	72
3			Heptane, 4-methyl-	114	C8H18	000589-53-7	64
4			Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	59
5			1-Butanol, 2-ethyl-	102	C6H14O	000097-95-0	59



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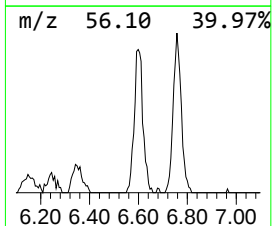
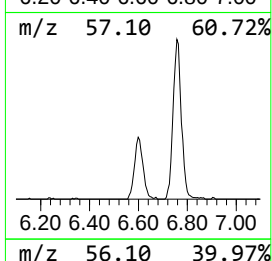
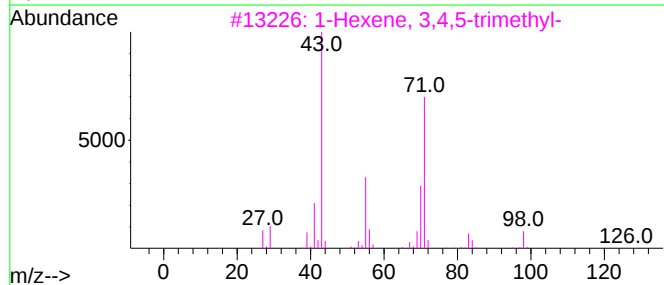
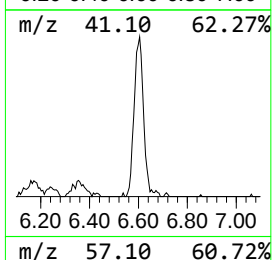
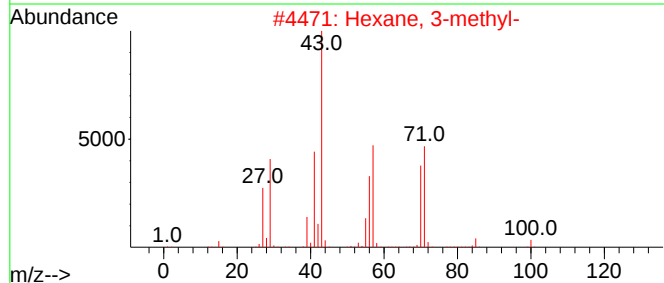
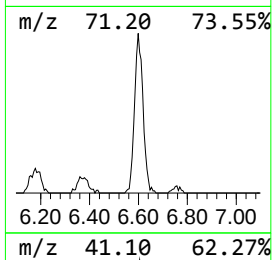
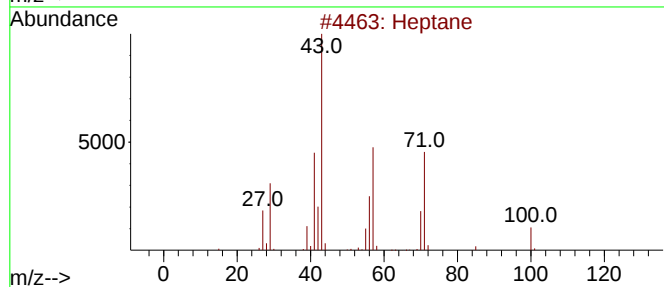
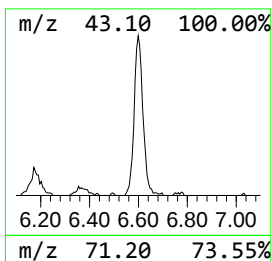
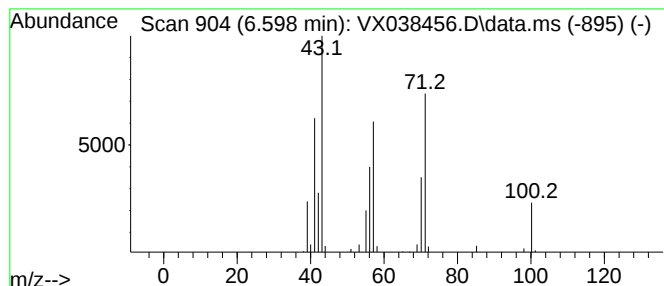
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100523W.M
Quant Title : SW846 8260

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

Peak Number 2 Heptane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.598	6.38 ug/l	77975	1,4-Difluorobenzene	6.757

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane	100	C7H16	000142-82-5	90
2			Hexane, 3-methyl-	100	C7H16	000589-34-4	59
3			1-Hexene, 3,4,5-trimethyl-	126	C9H18	056728-10-0	47
4			Pentane, 2-methyl-	86	C6H14	000107-83-5	47
5			Butane, 2,2-dimethyl-	86	C6H14	000075-83-2	43



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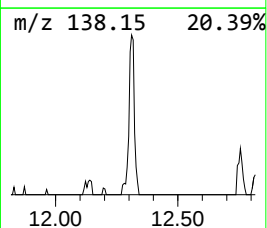
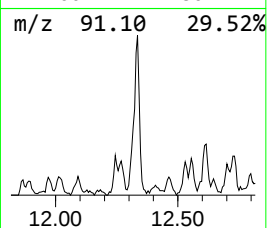
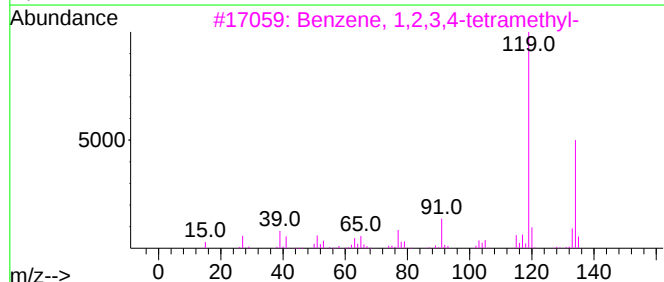
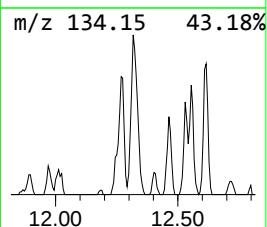
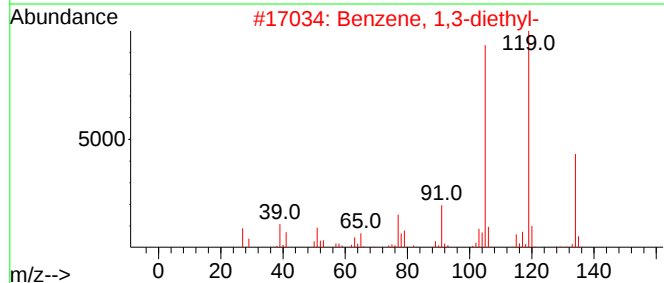
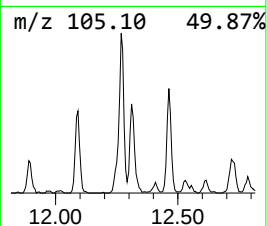
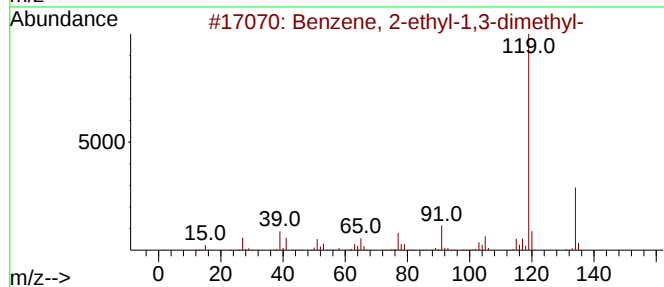
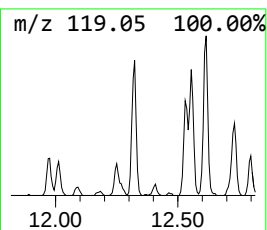
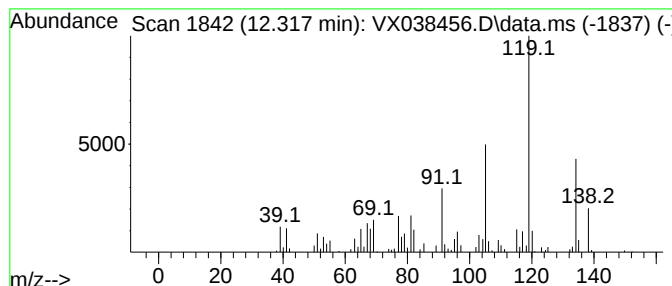
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.317	5.98 ug/l	114182	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	93
2			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	90
3			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	90
4			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	90
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	90



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ClientSampleId :
54

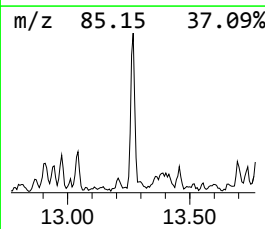
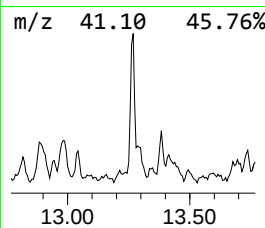
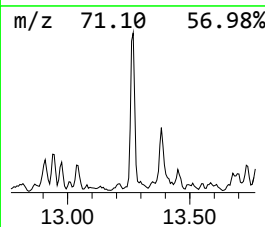
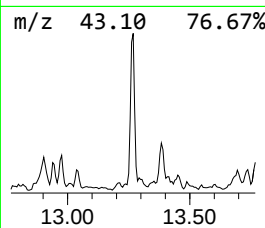
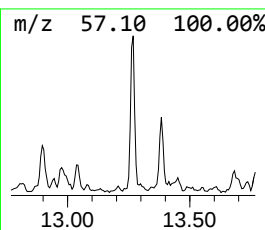
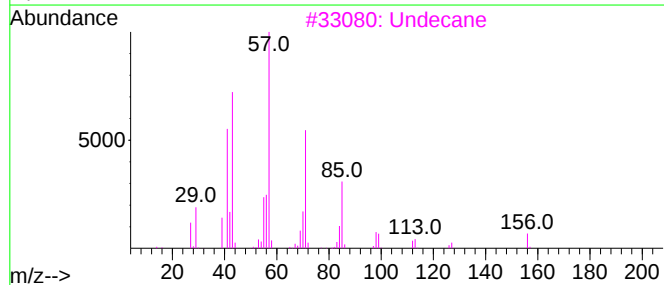
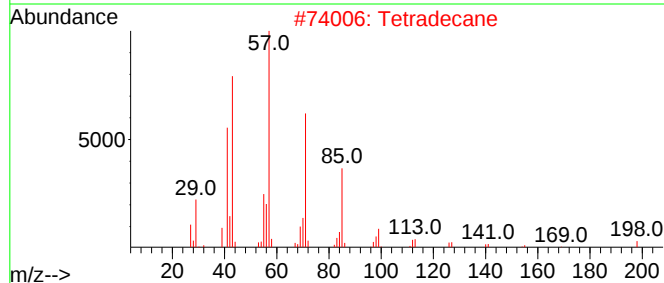
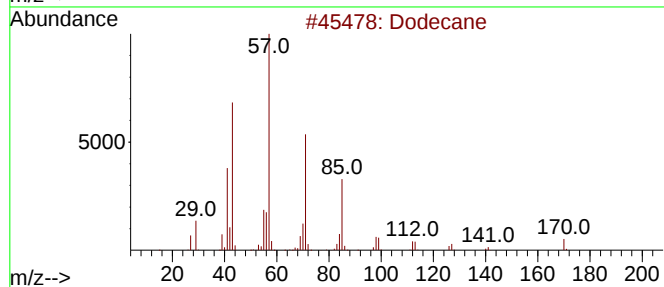
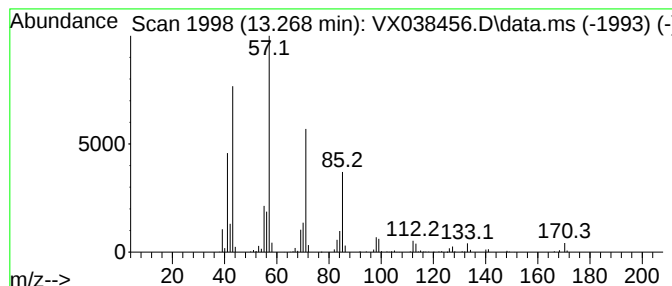
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Dodecane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.268	8.36 ug/l	159573	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane	170	C12H26	000112-40-3	96
2			Tetradecane	198	C14H30	000629-59-4	87
3			Undecane	156	C11H24	001120-21-4	86
4			Hexadecane	226	C16H34	000544-76-3	80
5			Carbonic acid, octadecyl prop-1-...	354	C22H42O3	1000383-11-5	80



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX102423\
Data File : VX038456.D
Acq On : 24 Oct 2023 15:31
Operator : JC/MD
Sample : 04975-01
Misc : 1.12g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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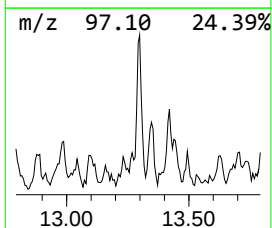
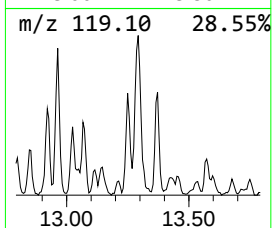
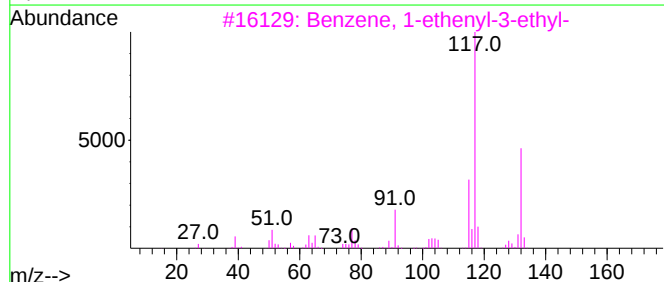
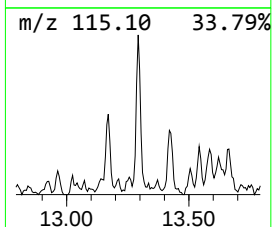
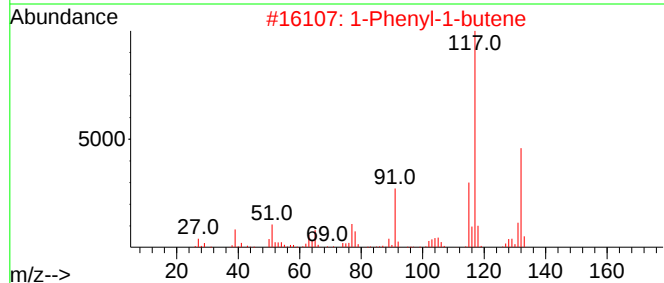
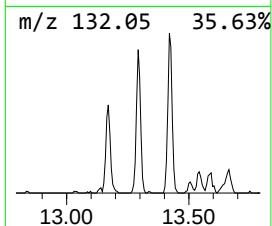
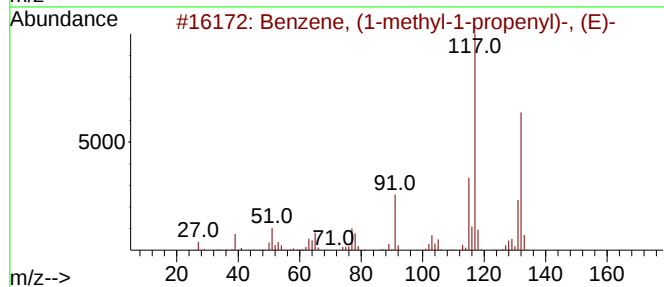
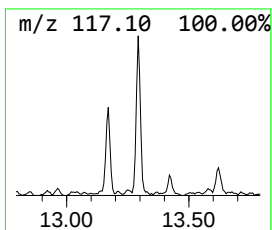
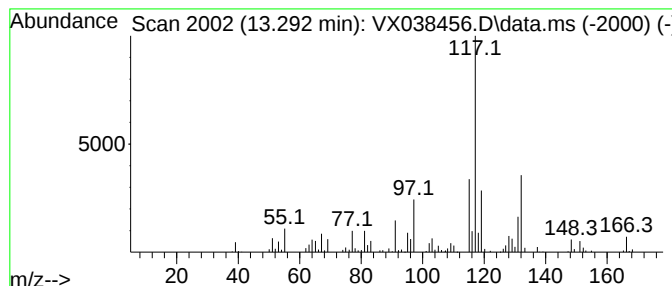
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, (1-methyl-1-propen... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	6.80 ug/l	129765	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	76
2			1-Phenyl-1-butene	132	C10H12	000824-90-8	76
3			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	70
4			Benzene, 2-butenyl-	132	C10H12	001560-06-1	68
5			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	68



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX102423\
Data File : VX038456.D
Acq On : 24 Oct 2023 15:31
Operator : JC/MD
Sample : 04975-01
Misc : 1.12g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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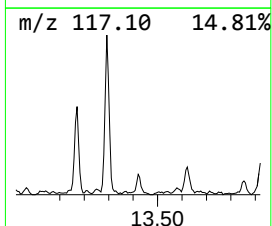
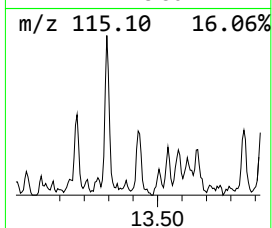
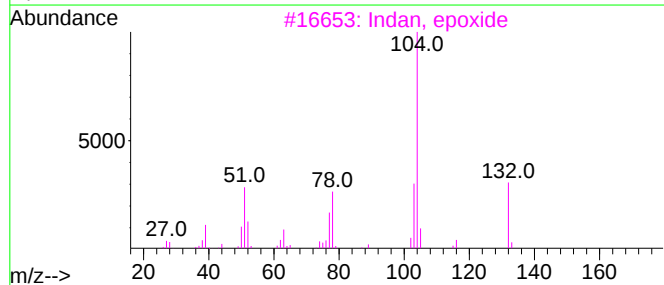
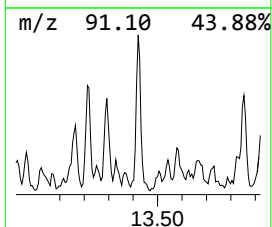
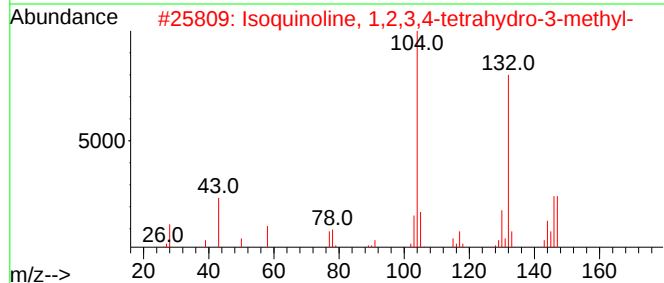
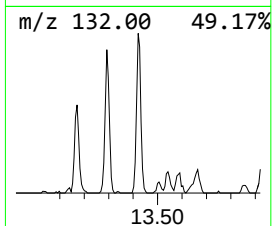
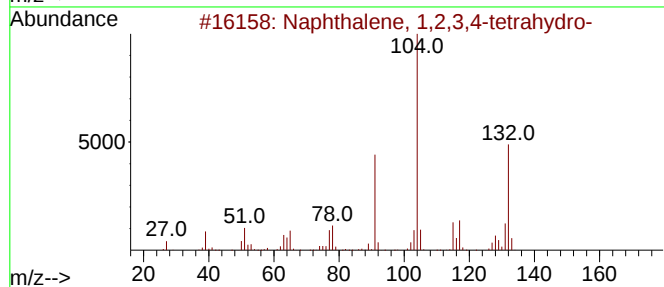
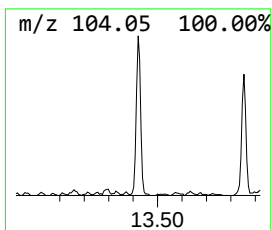
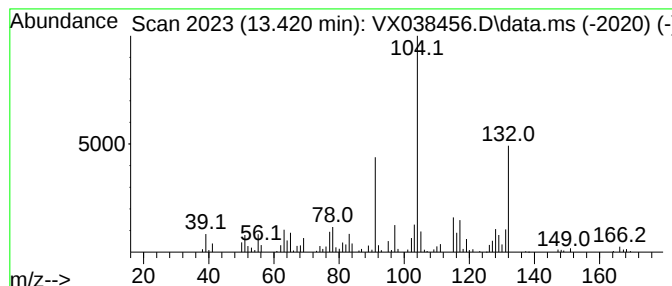
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.420	7.09 ug/l	135415	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	96
2			Isoquinoline, 1,2,3,4-tetrahydro...	147	C10H13N	029726-60-1	59
3			Indan, epoxide	132	C9H8O	000768-22-9	52
4			2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	52
5			Benzene, 3-cyclohexen-1-yl-	158	C12H14	004994-16-5	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX102423\
Data File : VX038456.D
Acq On : 24 Oct 2023 15:31
Operator : JC/MD
Sample : 04975-01
Misc : 1.12g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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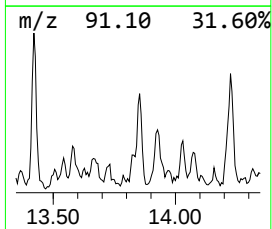
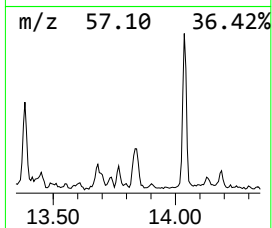
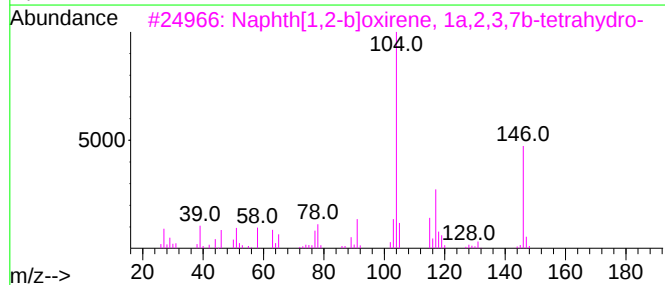
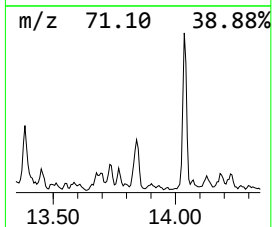
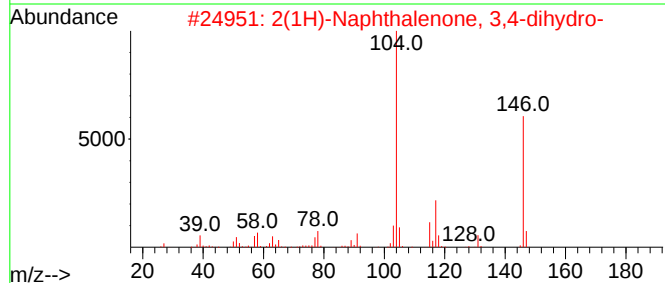
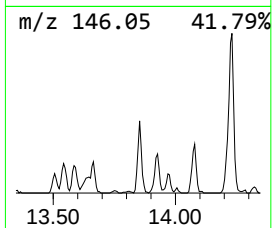
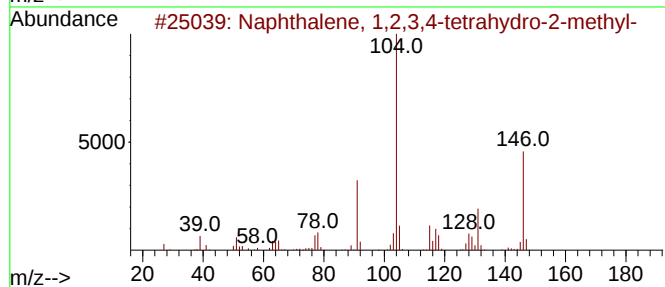
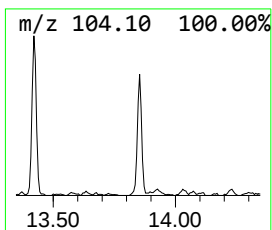
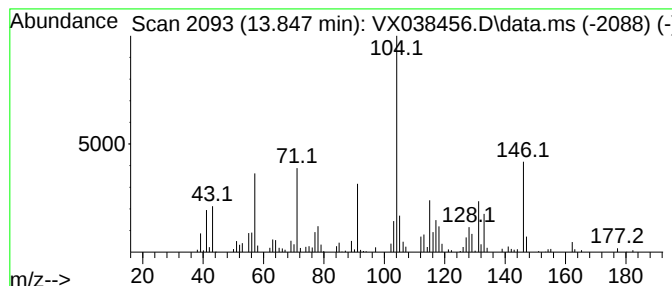
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Quant Title : SW846 8260

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

Peak Number 7 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.847	5.71 ug/l	108928	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	003877-19-8	70
2			2(1H)-Naphthalenone, 3,4-dihydro-	146	C10H10O	000530-93-8	46
3			Naphth[1,2-b]oxirene, 1a,2,3,7b-...	146	C10H10O	002461-34-9	46
4			Naphthalene, 2-ethyl-1,2,3,4-tet...	160	C12H16	032367-54-7	38
5			.beta.-Phenylethyl butyrate	192	C12H16O2	000103-52-6	35



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX102423\
Data File : VX038456.D
Acq On : 24 Oct 2023 15:31
Operator : JC/MD
Sample : 04975-01
Misc : 1.12g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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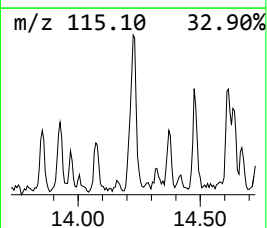
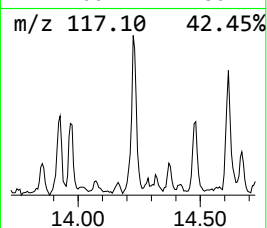
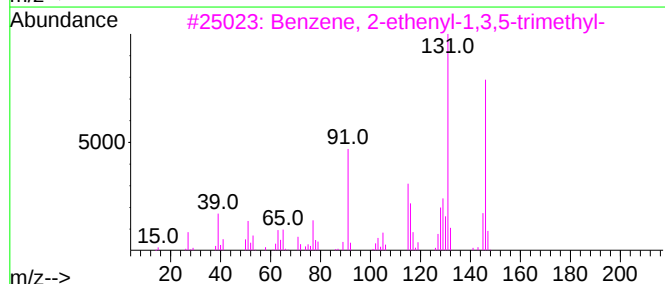
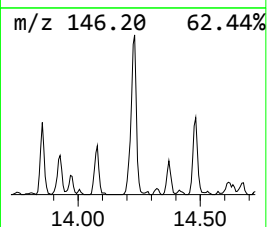
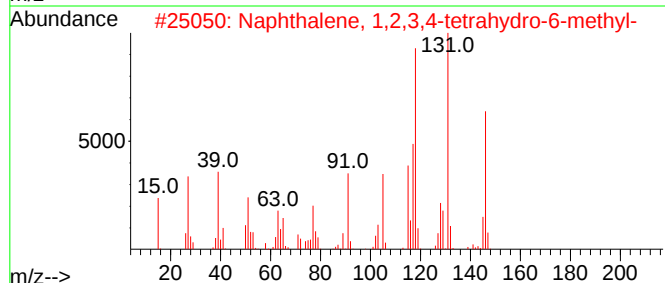
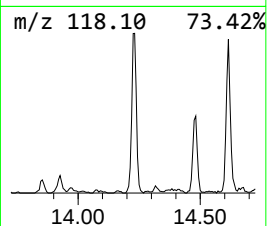
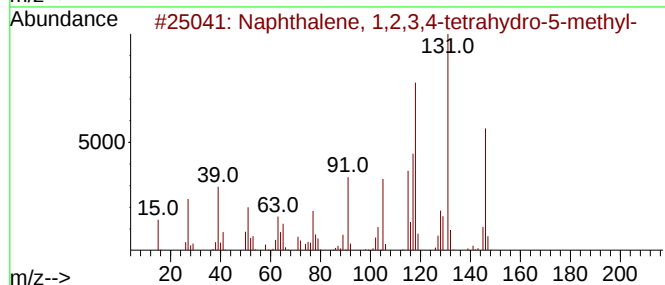
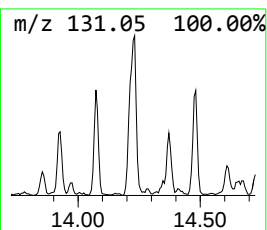
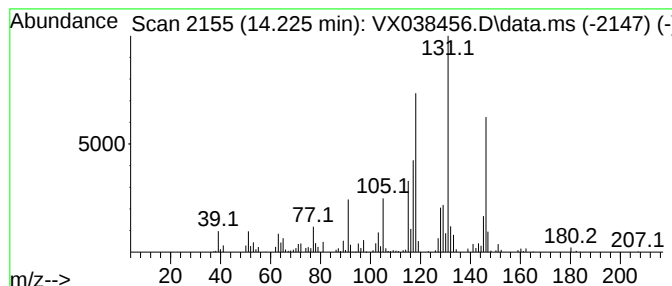
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100523W.M
Quant Title : SW846 8260

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

Peak Number 8 Naphthalene, 1,2,3,4-tetra... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.225	10.06 ug/l	192004	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	95
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	93
3			Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	78
4			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	60
5			Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	55



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX102423\
Data File : VX038456.D
Acq On : 24 Oct 2023 15:31
Operator : JC/MD
Sample : 04975-01
Misc : 1.12g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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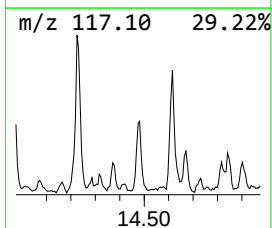
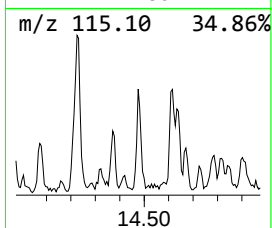
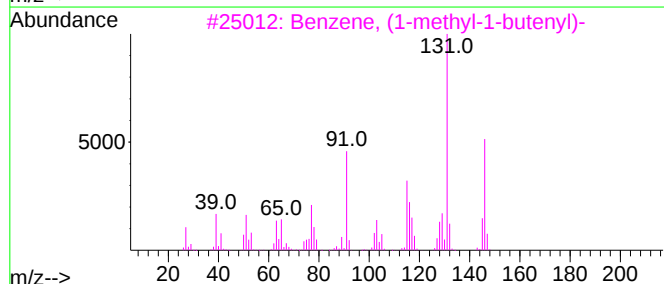
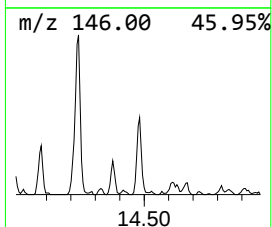
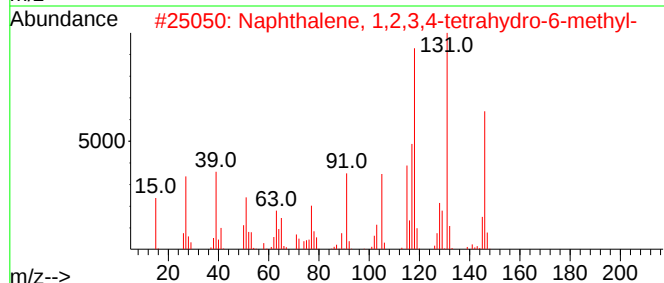
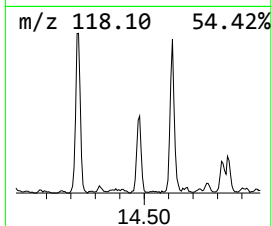
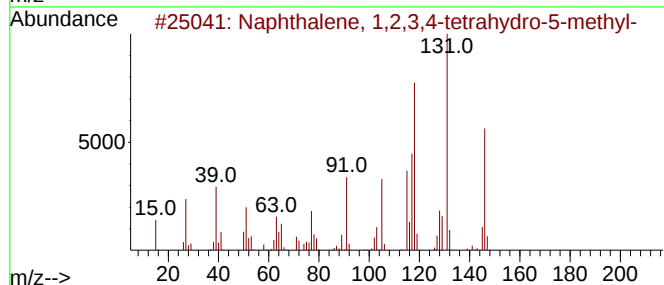
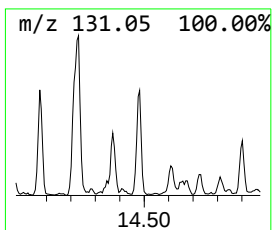
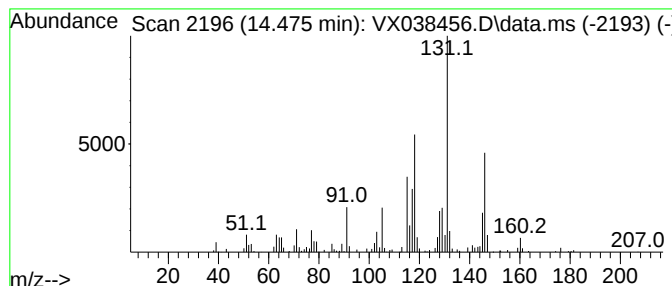
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Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1,2,3,4-tetra... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.475	6.09 ug/l	116286	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	94
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	89
3			Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	76
4			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	70
5			1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	68



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX102423\
Data File : VX038456.D
Acq On : 24 Oct 2023 15:31
Operator : JC/MD
Sample : 04975-01
Misc : 1.12g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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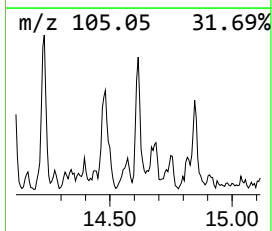
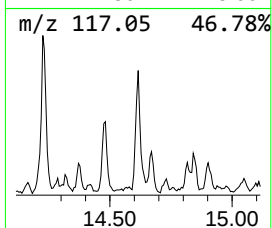
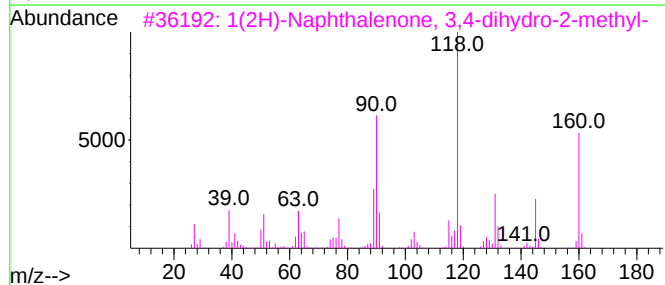
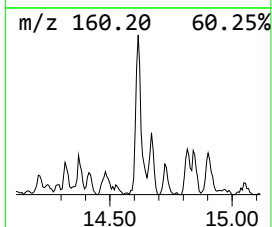
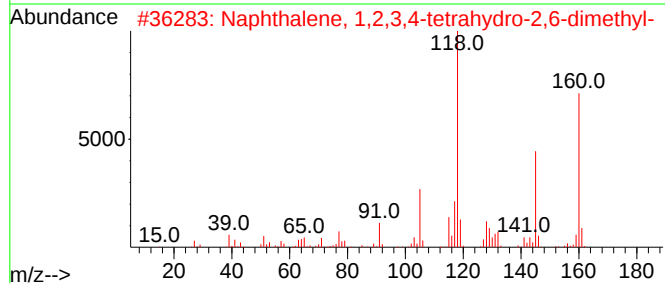
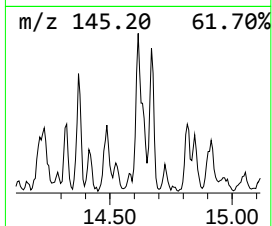
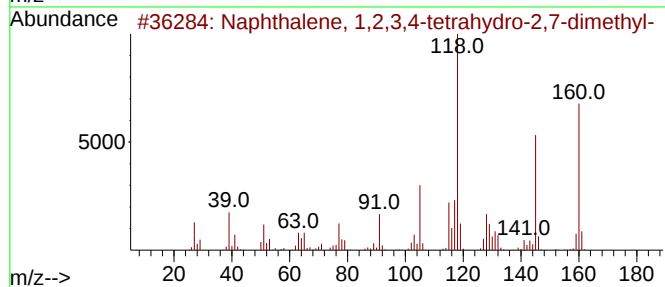
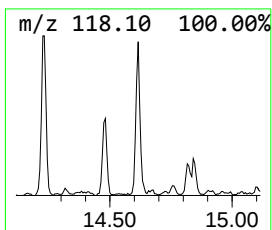
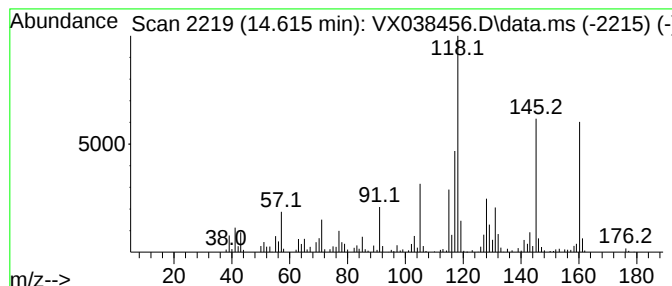
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100523W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.615	5.69 ug/l	108581	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	013065-07-1	91
2			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	007524-63-2	68
3			1(2H)-Naphthalenone, 3,4-dihydro...	160	C11H12O	001590-08-5	53
4			Benzene, (3-methyl-1-methylenebu...	160	C12H16	038212-14-5	52
5			Diboroxide, trimethylphenyl-	160	C9H14B2O	1000162-60-0	46



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX102423\
Data File : VX038456.D
Acq On : 24 Oct 2023 15:31
Operator : JC/MD
Sample : 04975-01
Misc : 1.12g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
54

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100523W.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
Hexane, 3-methyl-	5.806	8.3	ug/l	88338	1	5.544	529491	50.0
Heptane	6.598	6.4	ug/l	77975	2	6.757	611097	50.0
Benzene, 2-ethy...	12.317	6.0	ug/l	114182	4	12.024	954340	50.0
Dodecane	13.268	8.4	ug/l	159573	4	12.024	954340	50.0
Benzene, (1-met...	13.292	6.8	ug/l	129765	4	12.024	954340	50.0
Naphthalene, 1,...	13.420	7.1	ug/l	135415	4	12.024	954340	50.0
Naphthalene, 1,...	13.847	5.7	ug/l	108928	4	12.024	954340	50.0
Naphthalene, 1,...	14.225	10.1	ug/l	192004	4	12.024	954340	50.0
Naphthalene, 1,...	14.475	6.1	ug/l	116286	4	12.024	954340	50.0
Naphthalene, 1,...	14.615	5.7	ug/l	108581	4	12.024	954340	50.0