

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY102723\
 Data File : VY016118.D
 Acq On : 27 Oct 2023 20:51
 Operator : SY/MD
 Sample : 05107-10
 Misc : 5.04g/5.0mL/MSVOA_Y/SOIL
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 1080

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_Y\methods\82Y102623S.M

Title : SW846 8260

Signal : TIC: VY016118.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.077	36	42	60	rBV	890757	1234825	2.44%	0.316%
2	2.247	67	70	81	rVB	407178	708822	1.40%	0.181%
3	2.906	169	178	196	rBV	620505	1371746	2.71%	0.351%
4	3.259	227	236	255	rVB2	344718	864958	1.71%	0.221%
5	4.777	463	485	510	rVB5	249953	1336600	2.64%	0.342%
6	5.247	544	562	585	rBV	151104	537815	1.06%	0.138%
7	5.753	632	645	665	rBV	370735	1174605	2.32%	0.301%
8	6.801	807	817	837	rVB	371660	1161517	2.30%	0.297%
9	7.789	970	979	988	rVV4	1043349	3010445	5.95%	0.771%
10	8.008	1003	1015	1021	rBV	359398	870953	1.72%	0.223%
11	8.167	1030	1041	1053	rVB	14452888	32347853	63.96%	8.281%
12	8.423	1077	1083	1095	rVB2	223585	549592	1.09%	0.141%
13	8.551	1095	1104	1117	rBV	1018615	1998028	3.95%	0.511%
14	8.697	1119	1128	1149	rBV	360894	776540	1.54%	0.199%
15	9.191	1195	1209	1215	rBV	2256881	4884304	9.66%	1.250%
16	9.825	1307	1313	1325	rVB2	678553	1534441	3.03%	0.393%
17	9.965	1326	1336	1347	rBV	582344	1317926	2.61%	0.337%
18	10.185	1365	1372	1376	rBV	1195040	2249216	4.45%	0.576%
19	10.252	1376	1383	1390	rVB	22982448	43378956	85.77%	11.104%
20	10.374	1395	1403	1410	rVB2	1760020	3111628	6.15%	0.797%
21	10.526	1420	1428	1433	rVB2	373683	769269	1.52%	0.197%
22	10.910	1485	1491	1498	rBV	230890	535162	1.06%	0.137%
23	11.044	1505	1513	1517	rVV	552342	928606	1.84%	0.238%
24	11.093	1517	1521	1527	rVB	338325	553897	1.10%	0.142%
25	11.288	1544	1553	1563	rVV4	722190	1623505	3.21%	0.416%
26	11.386	1564	1569	1579	rVB	537368	905710	1.79%	0.232%
27	11.496	1581	1587	1592	rBV	495194	791218	1.56%	0.203%
28	11.599	1597	1604	1612	rBV	4530865	7171138	14.18%	1.836%
29	11.703	1612	1621	1632	rBV	18977496	35139025	69.48%	8.995%
30	12.032	1664	1675	1683	rBV	11732632	20407853	40.35%	5.224%
31	12.209	1694	1704	1707	rVV5	333827	896190	1.77%	0.229%
32	12.252	1707	1711	1715	rVB3	309846	530875	1.05%	0.136%
33	12.337	1720	1725	1728	rBV	470019	754710	1.49%	0.193%
34	12.422	1729	1739	1740	rBV5	358861	908024	1.80%	0.232%
35	12.489	1746	1750	1754	rVV3	430830	676699	1.34%	0.173%
36	12.672	1773	1780	1785	rVB	793168	1563988	3.09%	0.400%

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Integrator: RTE

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37	12.739	1785	1791	1794	rBV	3555894	5338222	10.55%	1.367%
38	12.770	1794	1796	1799	rVV	1275841	1717201	3.40%	0.440%
39	12.812	1799	1803	1809	rVB2	5520295	8691039	17.18%	2.225%
40	12.977	1817	1830	1838	rBV2	2117842	5156158	10.19%	1.320%
41	13.044	1838	1841	1844	rVV	393962	578379	1.14%	0.148%
42	13.123	1848	1854	1859	rVV	8998319	13859821	27.40%	3.548%
43	13.172	1859	1862	1867	rVB	967640	1405976	2.78%	0.360%
44	13.227	1867	1871	1880	rVB5	354896	875589	1.73%	0.224%
45	13.318	1880	1886	1890	rBV2	663200	1257624	2.49%	0.322%
46	13.373	1891	1895	1898	rVV2	420849	686461	1.36%	0.176%
47	13.459	1901	1909	1914	rVV	4303391	7143971	14.13%	1.829%
48	13.501	1914	1916	1924	rVB3	430682	655596	1.30%	0.168%
49	13.599	1924	1932	1933	rBV5	340904	596255	1.18%	0.153%
50	13.629	1933	1937	1940	rVV2	1486706	2259529	4.47%	0.578%
51	13.666	1940	1943	1952	rVB2	1556632	2521426	4.99%	0.645%
52	13.757	1953	1958	1962	rBV2	2051645	2797368	5.53%	0.716%
53	13.812	1962	1967	1973	rVB	5992761	9457436	18.70%	2.421%
54	13.891	1975	1980	1982	rBV	824524	1313433	2.60%	0.336%
55	13.916	1982	1984	1989	rVV	1024267	1421674	2.81%	0.364%
56	13.977	1989	1994	1998	rVB	1087092	1534043	3.03%	0.393%
57	14.080	2006	2011	2015	rBV2	1139386	1736230	3.43%	0.444%
58	14.129	2016	2019	2028	rVB6	256795	659023	1.30%	0.169%
59	14.221	2029	2034	2037	rBV3	534689	847024	1.67%	0.217%
60	14.300	2037	2047	2050	rVV3	1074373	2989073	5.91%	0.765%
61	14.337	2050	2053	2058	rVV	1624602	2383497	4.71%	0.610%
62	14.385	2058	2061	2065	rVV2	766106	1091732	2.16%	0.279%
63	14.428	2065	2068	2075	rVB4	561755	887689	1.76%	0.227%
64	14.568	2087	2091	2095	rBV3	460693	799506	1.58%	0.205%
65	14.617	2095	2099	2103	rVB	2234826	3008869	5.95%	0.770%
66	14.678	2103	2109	2115	rBV3	2317400	4756525	9.40%	1.218%
67	14.751	2115	2121	2126	rVB3	1132163	2375625	4.70%	0.608%
68	14.836	2130	2135	2139	rVB	1430394	2012395	3.98%	0.515%
69	14.952	2149	2154	2157	rBV4	669206	1062463	2.10%	0.272%
70	15.056	2166	2171	2177	rBV4	681960	1387061	2.74%	0.355%
71	15.141	2181	2185	2191	rVB5	590982	1147186	2.27%	0.294%
72	15.245	2191	2202	2212	rBV	25427332	50575618	100.00%	12.947%
73	15.342	2212	2218	2224	rBV2	1050824	2018431	3.99%	0.517%
74	15.440	2230	2234	2242	rVB	2079078	3101398	6.13%	0.794%
75	15.531	2242	2249	2256	rBV4	534605	1453188	2.87%	0.372%
76	15.611	2258	2262	2266	rVB3	347906	625008	1.24%	0.160%

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77	15.696	2268	2276	2285	rBV6	404304	1425721	2.82%	0.365%
78	15.891	2303	2308	2315	rVB6	314405	673362	1.33%	0.172%
79	16.031	2325	2331	2336	rBV4	368801	938106	1.85%	0.240%
80	16.092	2337	2341	2348	rVV7	240972	553732	1.09%	0.142%
81	16.159	2348	2352	2357	rVV	416906	649915	1.29%	0.166%
82	16.233	2359	2364	2368	rVV	461075	967854	1.91%	0.248%
83	16.300	2368	2375	2382	rVV	13691261	28368257	56.09%	7.262%
84	16.367	2382	2386	2391	rVV	1091459	2024420	4.00%	0.518%
85	16.415	2391	2394	2399	rVV3	267549	520930	1.03%	0.133%
86	16.495	2399	2407	2419	rVB	10394436	21760860	43.03%	5.571%

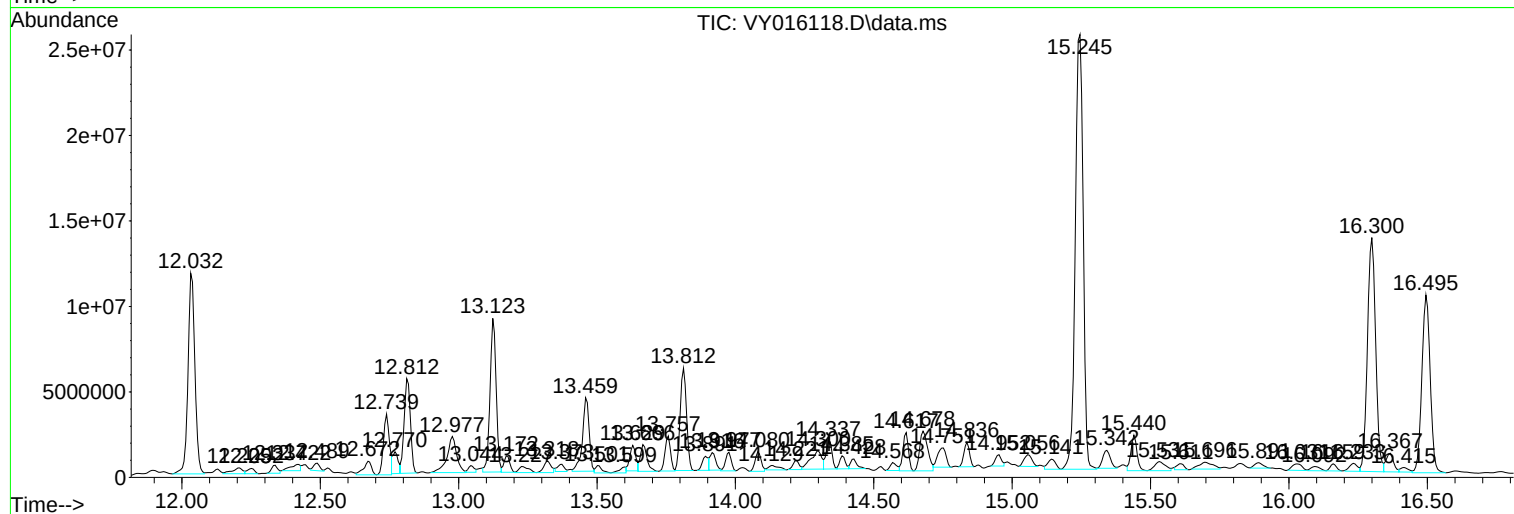
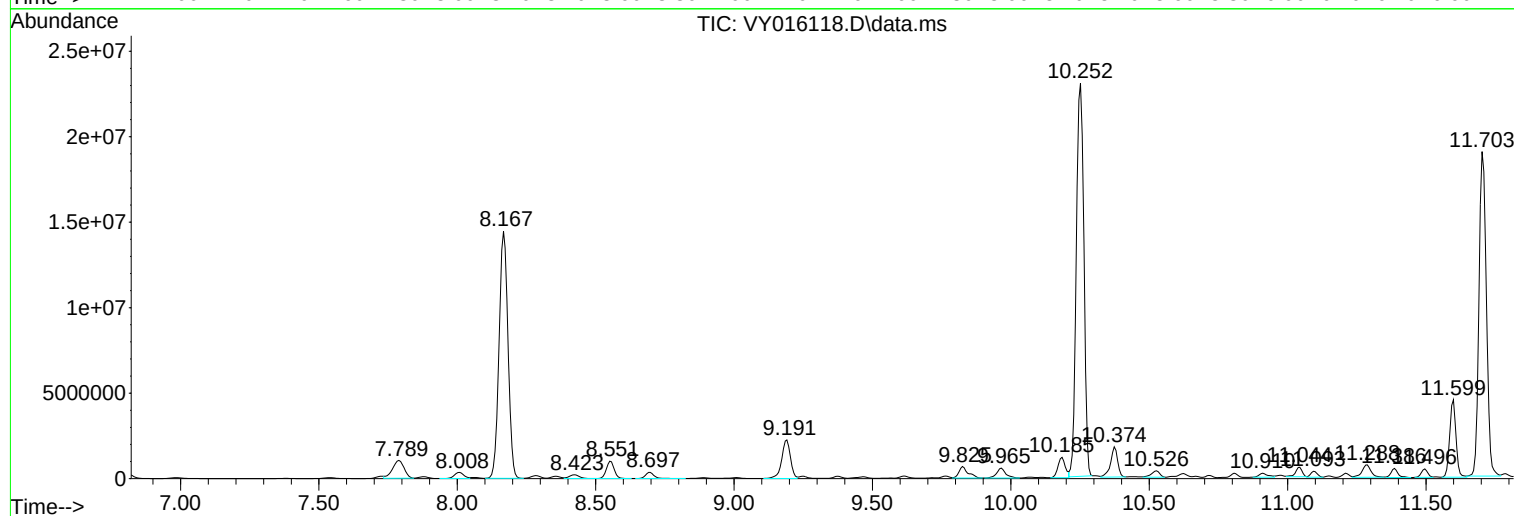
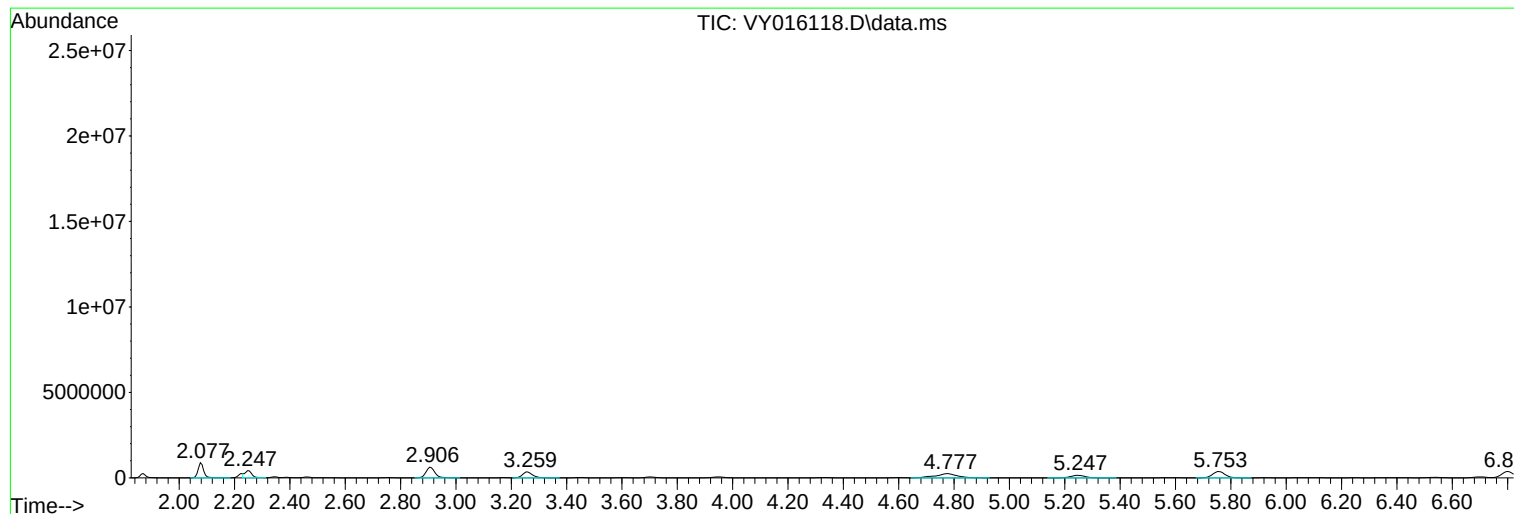
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y102623S.M
Quant Title : SW846 8260

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



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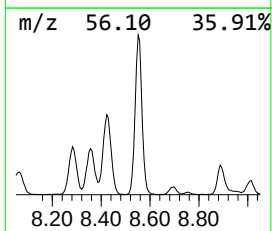
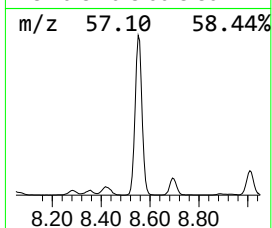
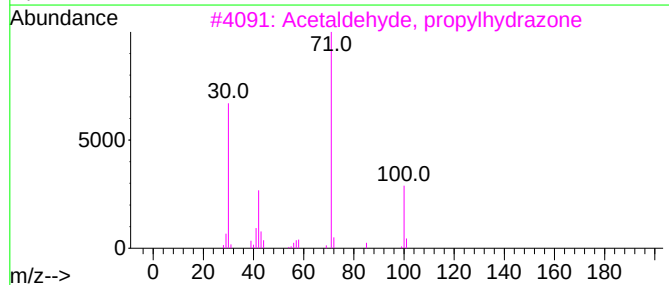
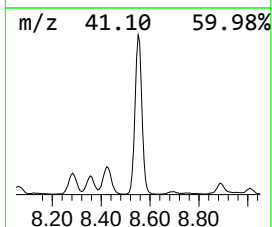
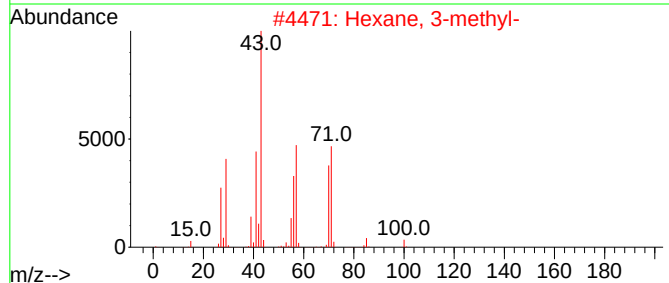
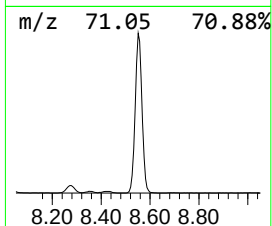
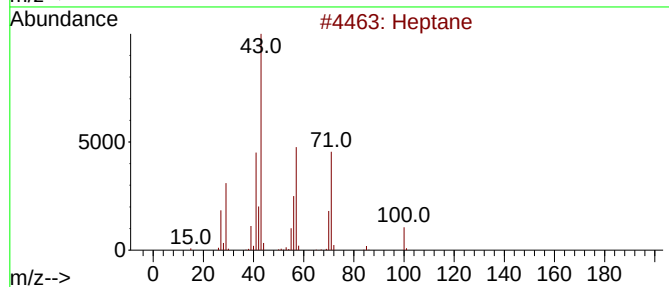
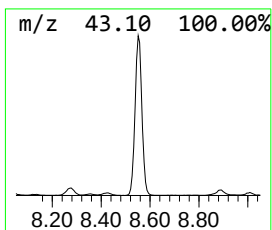
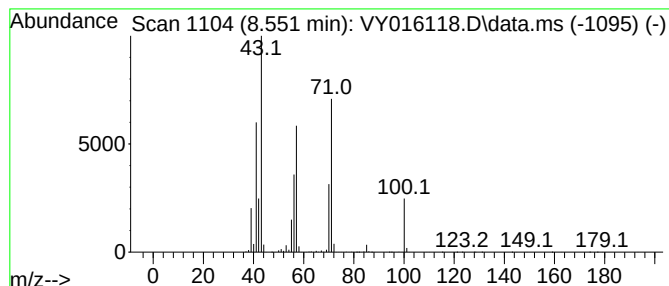
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y102623S.M
Quant Title : SW846 8260

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

Peak Number 1 Heptane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.551	128.65 ug/l	1998030	1,4-Difluorobenzene	8.697

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane	100	C7H16	000142-82-5	91
2			Hexane, 3-methyl-	100	C7H16	000589-34-4	50
3			Acetaldehyde, propylhydrazone	100	C5H12N2	007422-88-0	49
4			Pentane, 2-methyl-	86	C6H14	000107-83-5	43
5			Oxirane, butyl-	100	C6H12O	001436-34-6	38



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ALS Vial : 27 Sample Multiplier: 1

Instrument :
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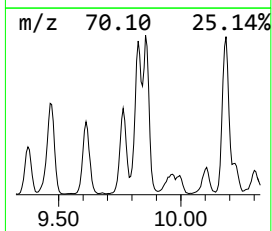
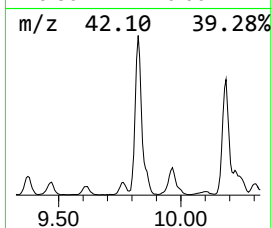
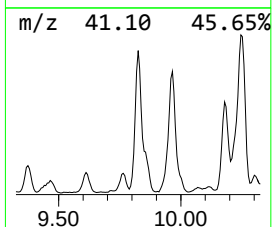
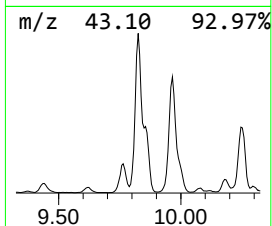
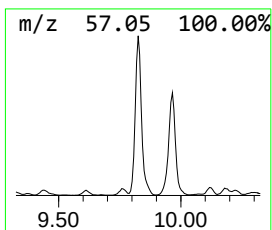
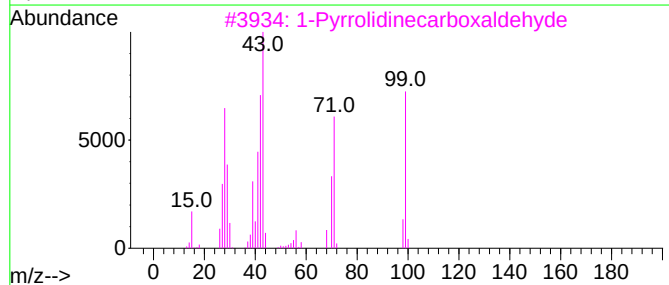
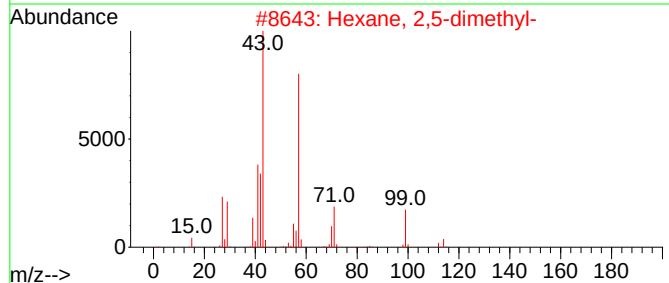
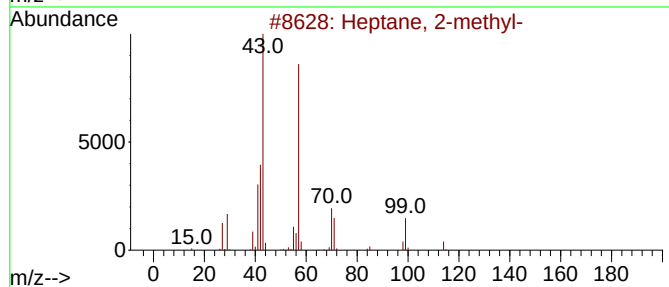
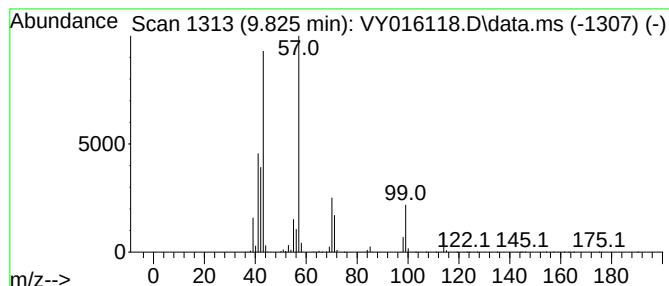
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y102623S.M
Quant Title : SW846 8260

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

Peak Number 2 Heptane, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.825	98.80 ug/l	1534440	1,4-Difluorobenzene	8.697

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane, 2-methyl-	114	C8H18	000592-27-8	94
2			Hexane, 2,5-dimethyl-	114	C8H18	000592-13-2	58
3			1-Pyrrolidinecarboxaldehyde	99	C5H9NO	003760-54-1	38
4			Hexane, 2,4,4-trimethyl-	128	C9H20	016747-30-1	37
5			Butane, 1-(ethenyloxy)-3-methyl-	114	C7H14O	039782-38-2	35



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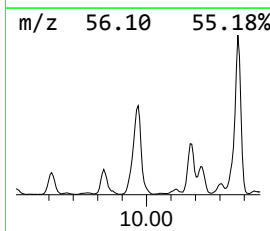
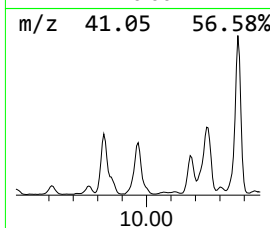
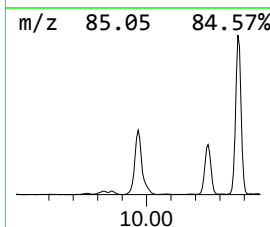
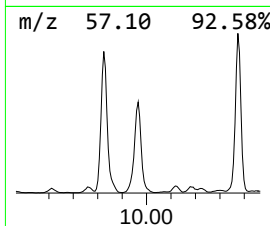
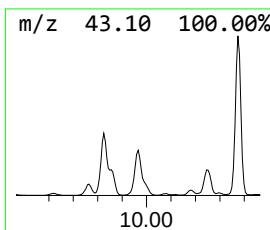
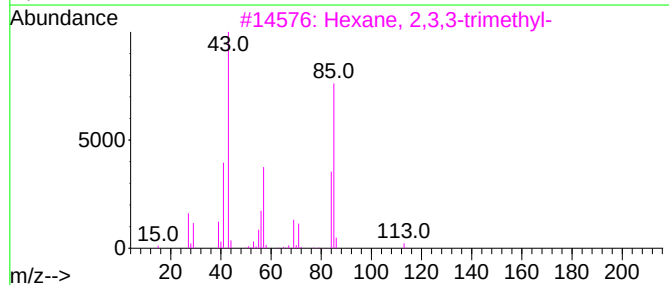
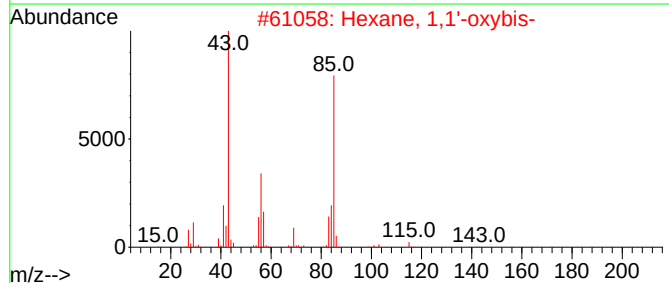
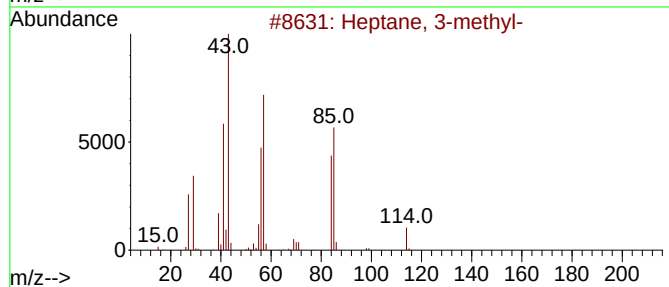
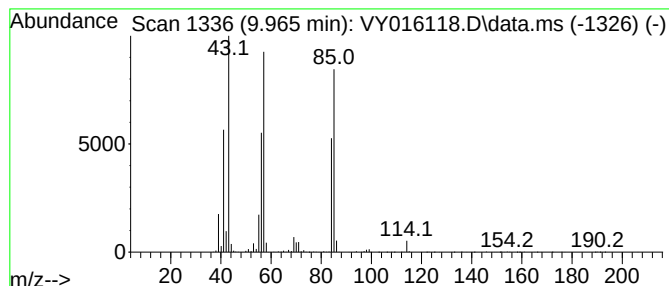
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y102623S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.965	84.86 ug/l	1317930	1,4-Difluorobenzene	8.697

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane, 3-methyl-	114	C8H18	000589-81-1	94
2			Hexane, 1,1'-oxybis-	186	C12H26O	000112-58-3	64
3			Hexane, 2,3,3-trimethyl-	128	C9H20	016747-28-7	64
4			Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	64
5			Heptane, 3,4,5-trimethyl-	142	C10H22	020278-89-1	59



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ClientSampleId :
1080

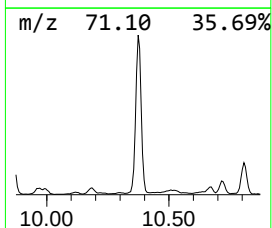
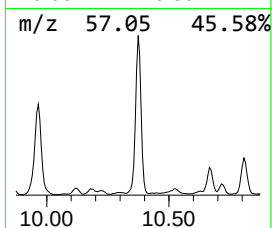
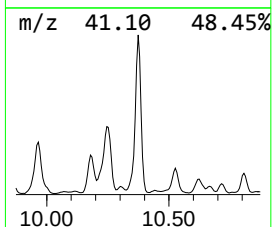
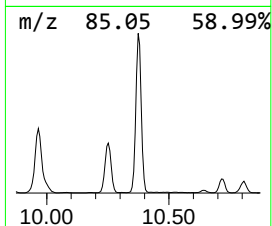
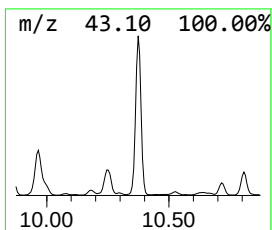
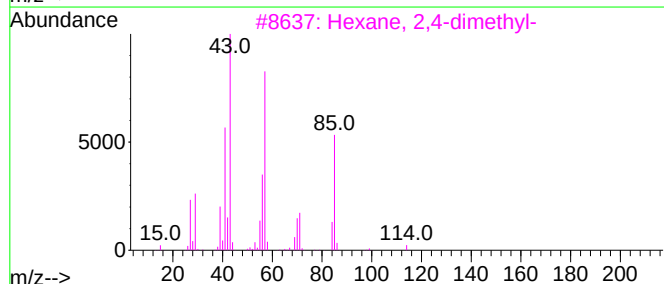
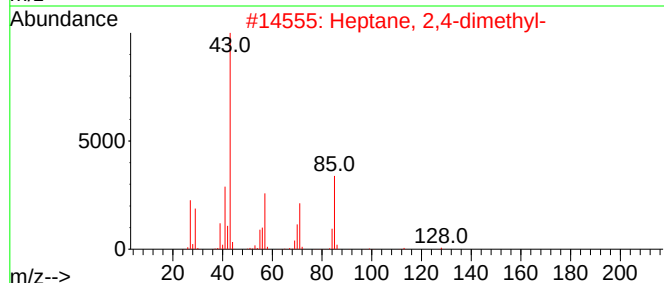
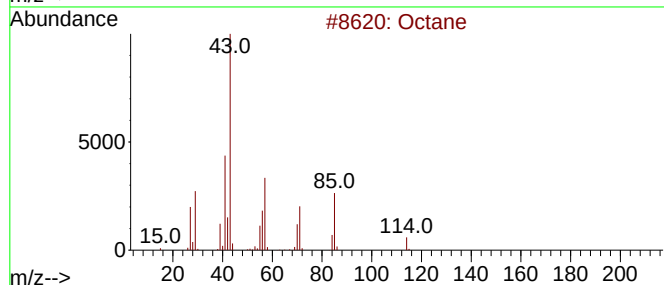
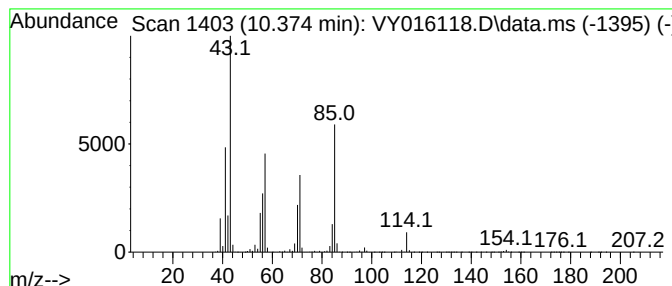
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Quant Title : SW846 8260

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

Peak Number 4 Octane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.374	196.63 ug/l	3111630	Chlorobenzene-d5	11.496

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane			114	C8H18	000111-65-9	94
2	Heptane, 2,4-dimethyl-			128	C9H20	002213-23-2	59
3	Hexane, 2,4-dimethyl-			114	C8H18	000589-43-5	59
4	Hexane, 1-(hexyloxy)-2-methyl-			200	C13H28O	074421-17-3	53
5	Undecane, 2,4-dimethyl-			184	C13H28	017312-80-0	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY102723\
Data File : VY016118.D
Acq On : 27 Oct 2023 20:51
Operator : SY/MD
Sample : 05107-10
Misc : 5.04g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
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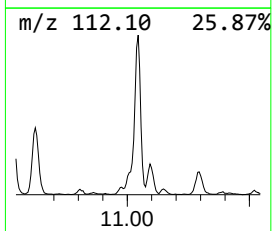
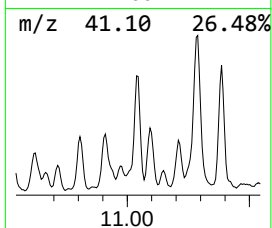
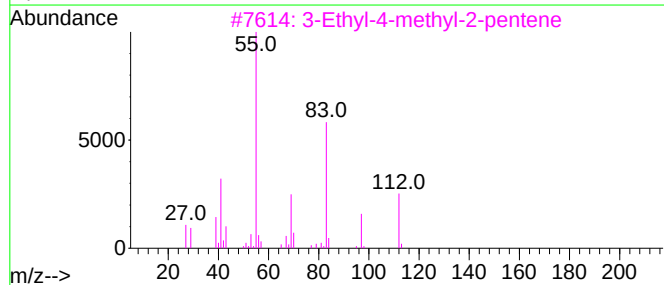
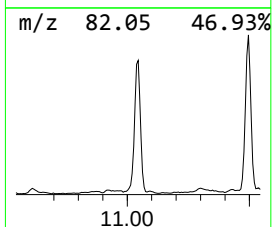
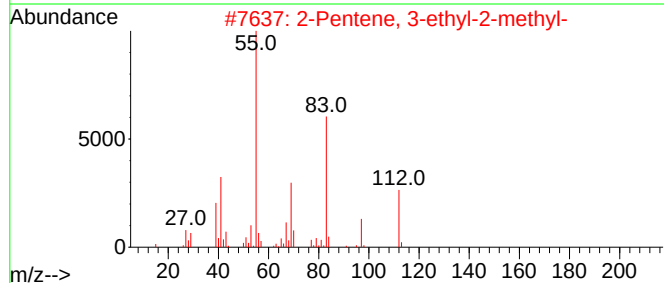
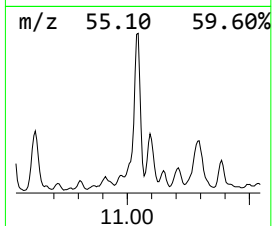
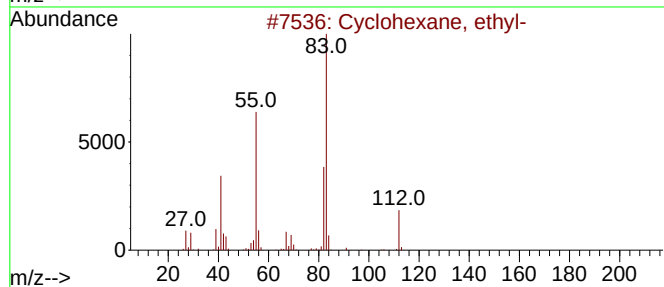
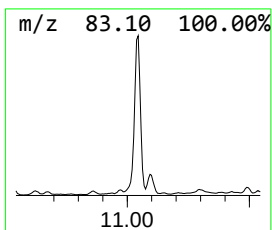
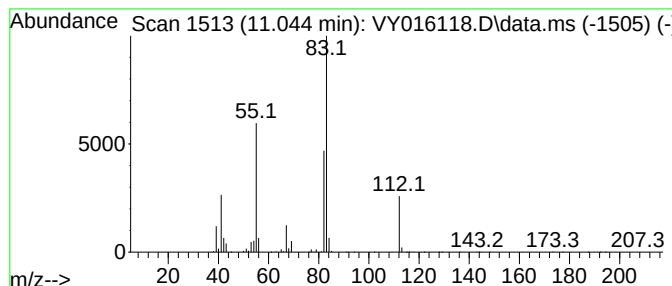
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Quant Title : SW846 8260

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

Peak Number 5 Cyclohexane, ethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.044	58.68 ug/l	928606	Chlorobenzene-d5	11.496

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, ethyl-	112	C8H16	001678-91-7	90
2			2-Pentene, 3-ethyl-2-methyl-	112	C8H16	019780-67-7	58
3			3-Ethyl-4-methyl-2-pentene	112	C8H16	019780-68-8	53
4			2-Hexene, 2,3-dimethyl-	112	C8H16	007145-20-2	53
5			4-Hexen-3-one, 5-methyl-	112	C7H12O	013905-10-7	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY102723\
Data File : VY016118.D
Acq On : 27 Oct 2023 20:51
Operator : SY/MD
Sample : 05107-10
Misc : 5.04g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
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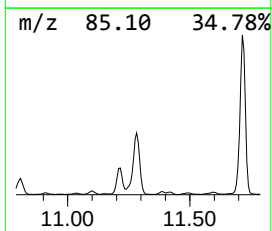
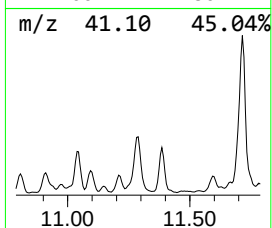
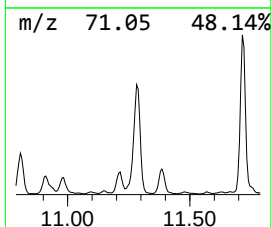
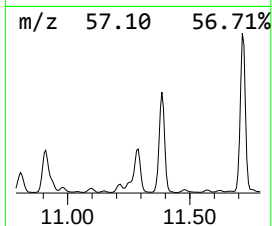
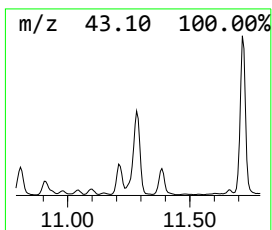
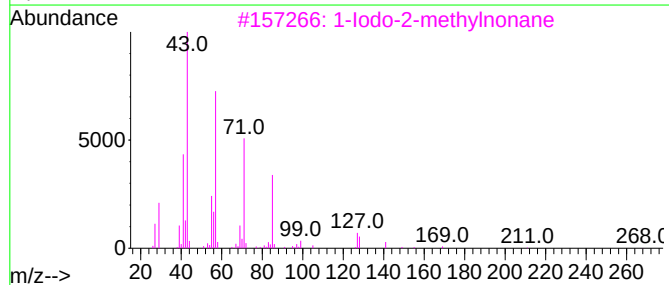
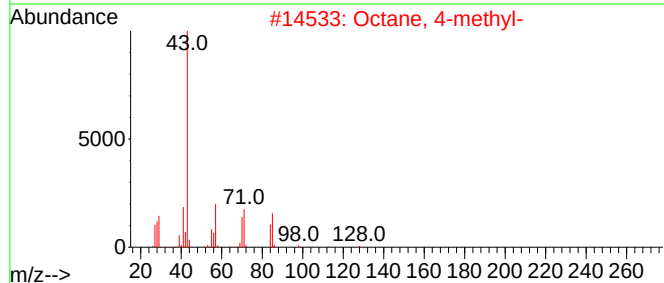
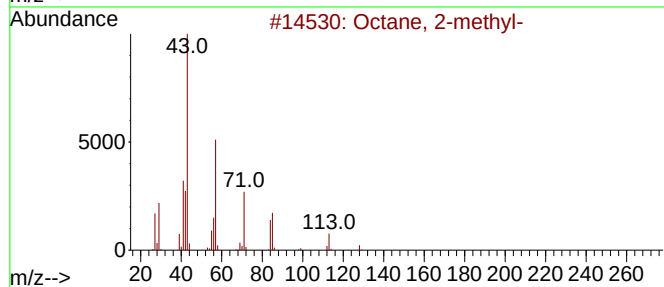
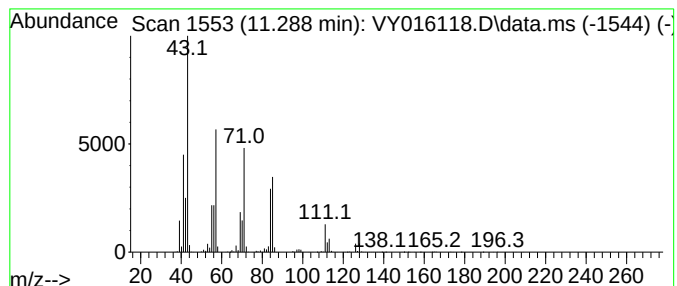
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Quant Title : SW846 8260

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

Peak Number 6 Octane, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.288	102.60 ug/l	1623510	Chlorobenzene-d5	11.496

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane, 2-methyl-	128	C9H20	003221-61-2	87
2			Octane, 4-methyl-	128	C9H20	002216-34-4	59
3			1-Iodo-2-methylnonane	268	C10H21I	1000101-47-9	50
4			Ether, hexyl pentyl	172	C11H24O	032357-83-8	47
5			Sulfurous acid, hexyl octyl ester	278	C14H30O3S	1000309-13-0	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY102723\
Data File : VY016118.D
Acq On : 27 Oct 2023 20:51
Operator : SY/MD
Sample : 05107-10
Misc : 5.04g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
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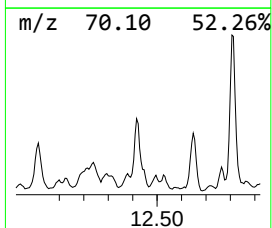
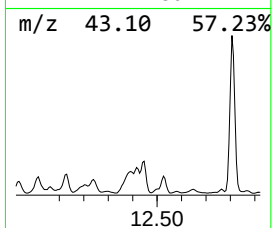
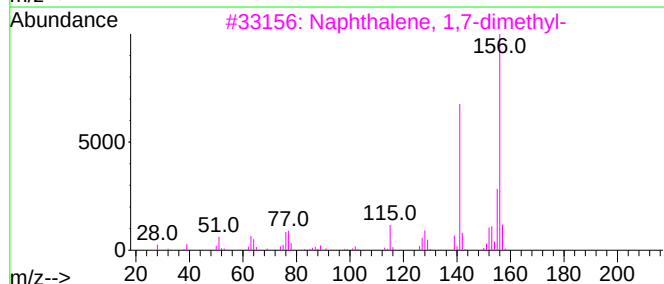
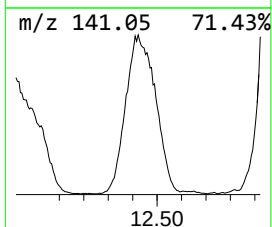
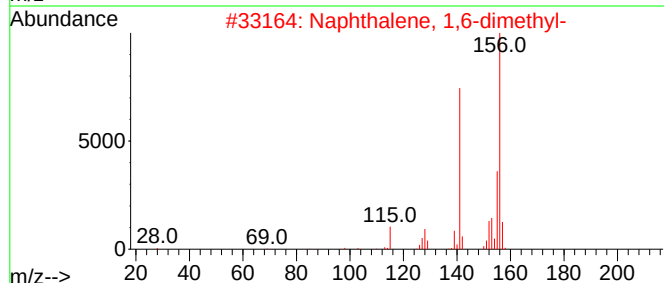
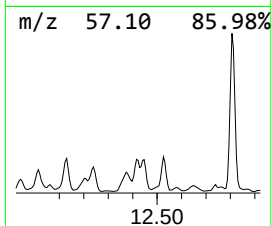
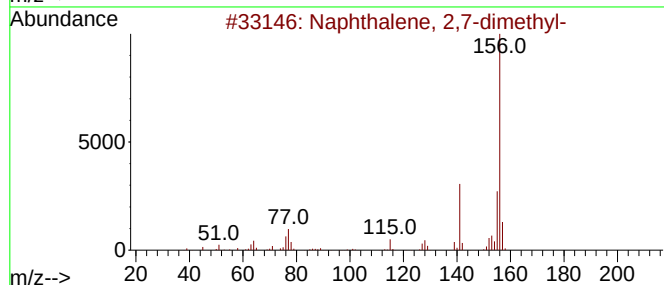
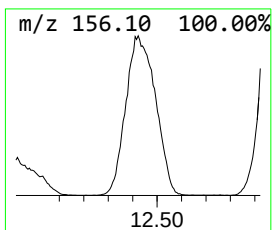
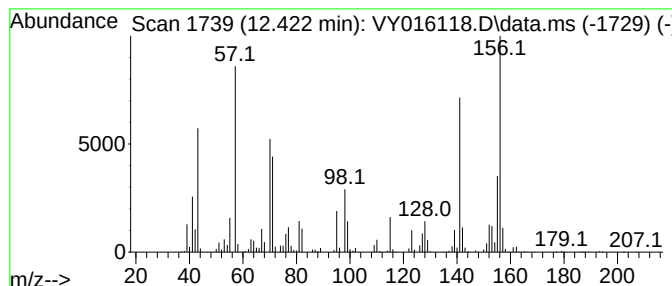
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Quant Title : SW846 8260

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

Peak Number 7 Naphthalene, 2,7-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.422	57.38 ug/l	908024	Chlorobenzene-d5	11.496

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96
2			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
3			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	95
4			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	95
5			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY102723\
Data File : VY016118.D
Acq On : 27 Oct 2023 20:51
Operator : SY/MD
Sample : 05107-10
Misc : 5.04g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
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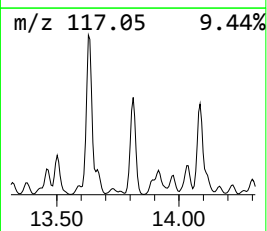
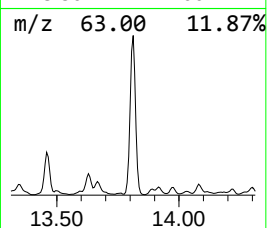
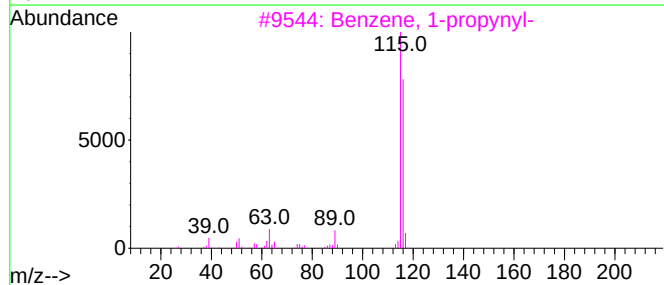
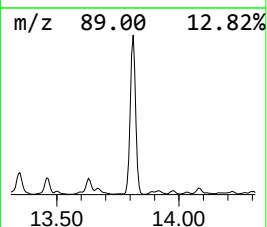
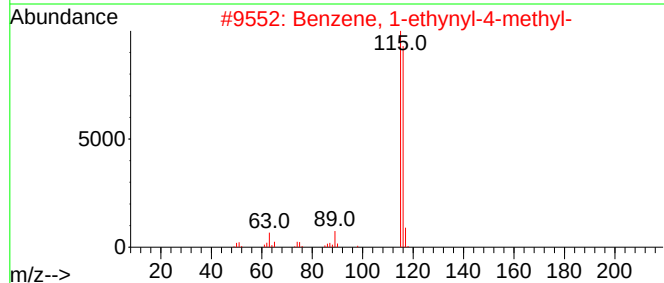
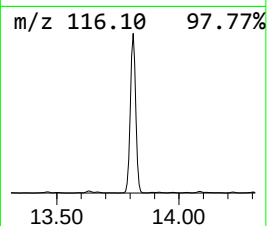
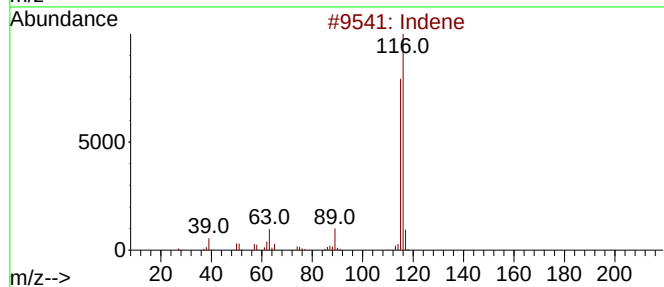
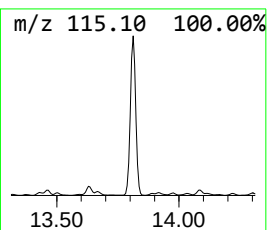
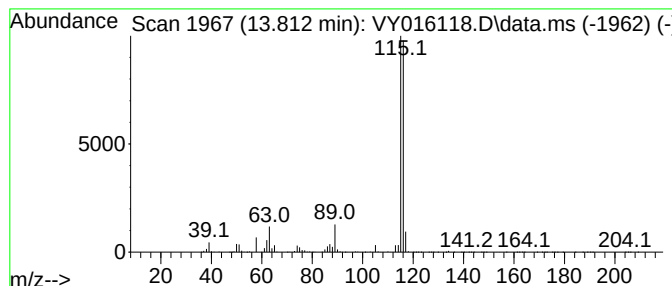
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Quant Title : SW846 8260

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

Peak Number 8 Indene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.812	66.19 ug/l	9457440	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indene	116	C9H8	000095-13-6	97
2			Benzene, 1-ethynyl-4-methyl-	116	C9H8	000766-97-2	91
3			Benzene, 1-propynyl-	116	C9H8	000673-32-5	91
4			3-Methylphenylacetylene	116	C9H8	000766-82-5	91
5			2-Methylphenylacetylene	116	C9H8	000766-47-2	83



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY102723\
Data File : VY016118.D
Acq On : 27 Oct 2023 20:51
Operator : SY/MD
Sample : 05107-10
Misc : 5.04g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
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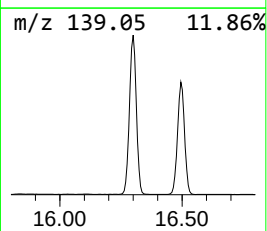
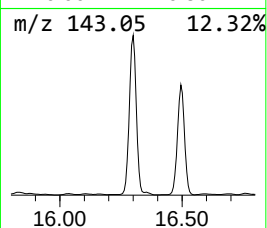
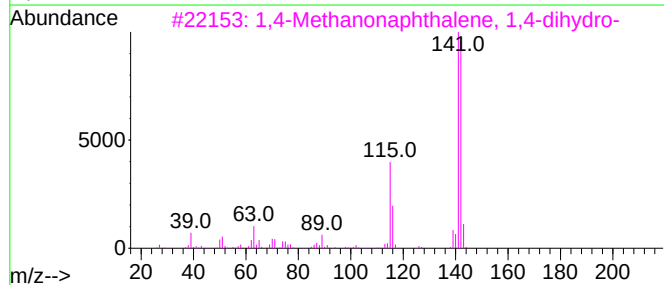
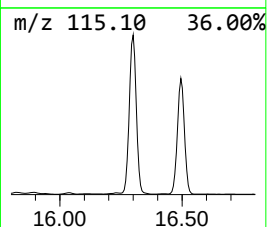
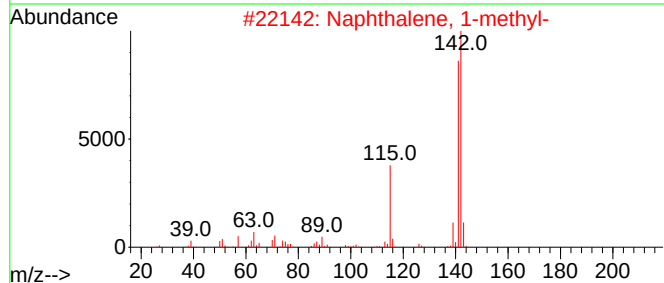
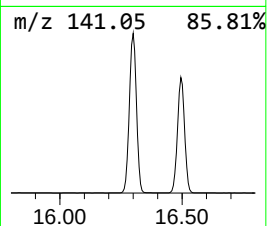
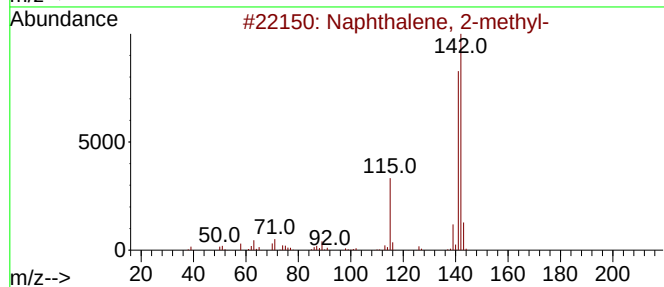
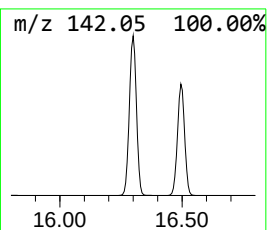
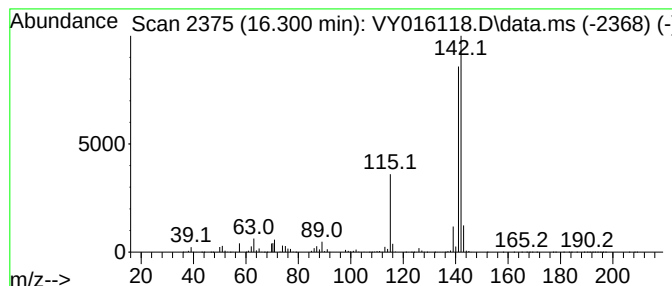
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Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.300	198.55 ug/l	28368300	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
4			Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	91
5			Benzocycloheptatriene	142	C11H10	000264-09-5	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY102723\
Data File : VY016118.D
Acq On : 27 Oct 2023 20:51
Operator : SY/MD
Sample : 05107-10
Misc : 5.04g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
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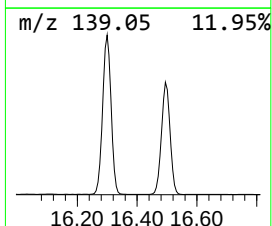
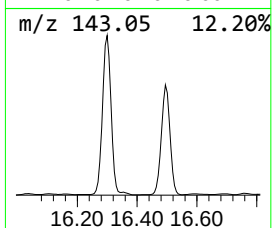
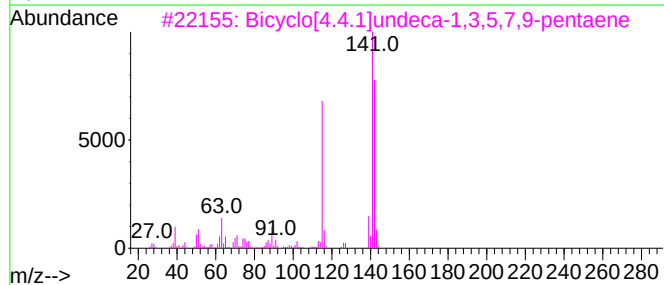
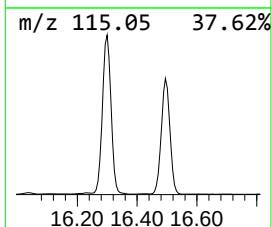
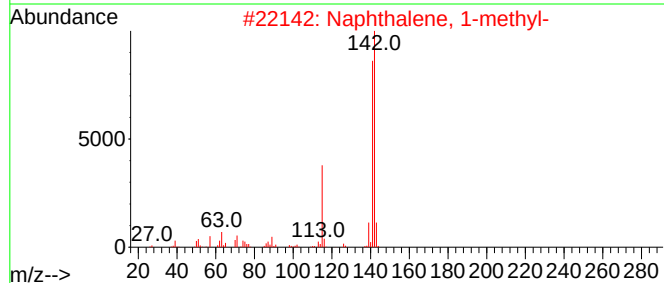
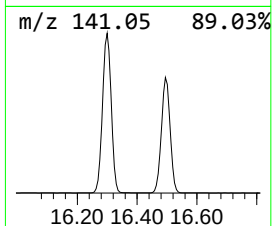
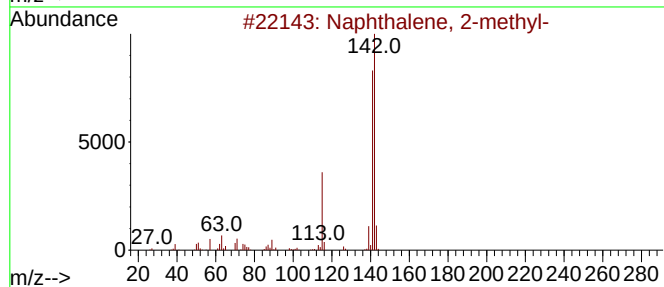
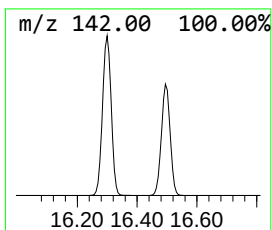
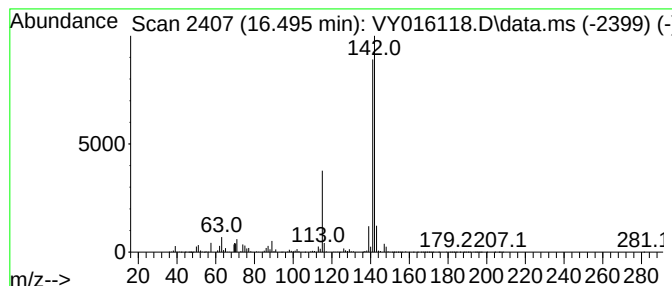
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Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.495	152.30 ug/l	21760900	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3			Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	91
4			Benzocycloheptatriene	142	C11H10	000264-09-5	91
5			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY102723\
Data File : VY016118.D
Acq On : 27 Oct 2023 20:51
Operator : SY/MD
Sample : 05107-10
Misc : 5.04g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
1080

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y102623S.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
Heptane	8.551	128.7	ug/l	1998030	2	8.697	776540	50.0
Heptane, 2-methyl-	9.825	98.8	ug/l	1534440	2	8.697	776540	50.0
Heptane, 3-methyl-	9.965	84.9	ug/l	1317930	2	8.697	776540	50.0
Octane	10.374	196.6	ug/l	3111630	3	11.496	791218	50.0
Cyclohexane, et...	11.044	58.7	ug/l	928606	3	11.496	791218	50.0
Octane, 2-methyl-	11.288	102.6	ug/l	1623510	3	11.496	791218	50.0
Naphthalene, 2,...	12.422	57.4	ug/l	908024	3	11.496	791218	50.0
Indene	13.812	66.2	ug/l	9457440	4	13.428	7143970	50.0
Naphthalene, 2-...	16.300	198.6	ug/l	28368300	4	13.428	7143970	50.0
Naphthalene, 1-...	16.495	152.3	ug/l	21760900	4	13.428	7143970	50.0