

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041521\
 Data File : VY004439.D
 Acq On : 15 Apr 2021 14:34
 Operator : SY/MD
 Sample : M2040-09
 Misc : 5.93G/5ML/MSVOA_Y/SOIL
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 FILL-C-2

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\82Y041221S.M
 Title : SW846 8260

Signal : TIC: VY004439.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.235	47	68	69	rBV3	17927	83358	3.26%	0.670%
2	2.607	116	129	147	rBV7	24596	177951	6.95%	1.431%
3	4.728	466	477	483	rBV2	12643	46289	1.81%	0.372%
4	7.728	959	969	974	rBV2	159556	380259	14.86%	3.058%
5	7.801	974	981	990	rVV	286477	678613	26.51%	5.457%
6	8.008	1006	1015	1023	rBV2	16151	40863	1.60%	0.329%
7	8.148	1029	1038	1050	rBV	167175	373116	14.58%	3.001%
8	8.343	1064	1070	1078	rVV4	10971	33871	1.32%	0.272%
9	8.551	1096	1104	1113	rVB3	12201	28212	1.10%	0.227%
10	8.697	1119	1128	1139	rVB	442211	883580	34.52%	7.106%
11	9.191	1196	1209	1214	rBV2	30561	74712	2.92%	0.601%
12	9.959	1326	1335	1345	rBV	17818	42165	1.65%	0.339%
13	10.185	1364	1372	1385	rBV	718362	1262701	49.33%	10.155%
14	10.374	1395	1403	1409	rVB4	22387	45480	1.78%	0.366%
15	10.617	1433	1443	1449	rVB5	9791	26104	1.02%	0.210%
16	11.282	1544	1552	1563	rVB4	26023	62935	2.46%	0.506%
17	11.380	1563	1568	1579	rBV	22044	41184	1.61%	0.331%
18	11.495	1579	1587	1595	rBV	628462	1044602	40.81%	8.401%
19	11.593	1598	1603	1612	rBV	54715	92313	3.61%	0.742%
20	11.715	1612	1623	1631	rBV2	49382	104130	4.07%	0.837%
21	12.014	1665	1672	1678	rBV5	12121	35025	1.37%	0.282%
22	12.416	1729	1738	1740	rVV6	18948	48504	1.90%	0.390%
23	12.489	1744	1750	1760	rVB2	1461738	2559497	100.00%	20.583%
24	12.672	1772	1780	1785	rBV	49331	80648	3.15%	0.649%
25	12.770	1792	1796	1799	rBV	33405	50933	1.99%	0.410%
26	12.806	1799	1802	1808	rVB2	66052	100357	3.92%	0.807%
27	12.971	1822	1829	1837	rBV	63725	127898	5.00%	1.029%
28	13.123	1848	1854	1859	rVV	137424	213230	8.33%	1.715%
29	13.428	1897	1904	1914	rVV2	592244	1009813	39.45%	8.121%
30	13.501	1914	1916	1923	rVB2	18621	26200	1.02%	0.211%
31	13.593	1926	1931	1933	rBV	29278	42889	1.68%	0.345%
32	13.629	1933	1937	1940	rVV2	70002	103209	4.03%	0.830%
33	13.672	1940	1944	1953	rVB3	50596	106300	4.15%	0.855%
34	13.751	1953	1957	1962	rBV	58877	87132	3.40%	0.701%
35	13.824	1965	1969	1974	rVB	26153	37559	1.47%	0.302%
36	13.885	1975	1979	1982	rBV	28884	46310	1.81%	0.372%

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Integrator: RTE

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Filtering: 5

Sampling : 1

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Max Peaks: 100

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Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y041221S.M
 Title : SW846 8260

37	13.916	1982	1984	1988	rVB	28583	33012	1.29%	0.265%
38	13.971	1988	1993	1999	rVB2	156399	243676	9.52%	1.960%
39	14.080	2006	2011	2017	rVB	69568	112308	4.39%	0.903%
40	14.220	2028	2034	2037	rBV3	24215	50297	1.97%	0.404%
41	14.300	2042	2047	2050	rVV	61477	111435	4.35%	0.896%
42	14.336	2050	2053	2057	rVB	52104	71183	2.78%	0.572%
43	14.379	2057	2060	2064	rBV	36416	51221	2.00%	0.412%
44	14.525	2080	2084	2087	rBV	19470	26939	1.05%	0.217%
45	14.568	2087	2091	2095	rVV	71654	120593	4.71%	0.970%
46	14.611	2095	2098	2103	rVB2	69734	98122	3.83%	0.789%
47	14.684	2103	2110	2115	rBV3	118504	265935	10.39%	2.139%
48	14.739	2115	2119	2125	rVB2	40414	71194	2.78%	0.573%
49	14.946	2148	2153	2157	rBV3	51921	84492	3.30%	0.679%
50	15.056	2165	2171	2178	rBV2	49181	98665	3.85%	0.793%
51	15.239	2193	2201	2207	rVB	103975	201446	7.87%	1.620%
52	15.348	2214	2219	2223	rBV5	31639	55621	2.17%	0.447%
53	15.397	2223	2227	2230	rVV4	18736	31434	1.23%	0.253%
54	15.434	2230	2233	2241	rVB4	36741	63861	2.50%	0.514%
55	15.525	2243	2248	2255	rBV3	35201	74873	2.93%	0.602%
56	15.885	2303	2307	2314	rVB5	17775	36573	1.43%	0.294%
57	16.293	2368	2374	2381	rBV	115287	219756	8.59%	1.767%
58	16.494	2401	2407	2416	rVB	52348	114277	4.46%	0.919%

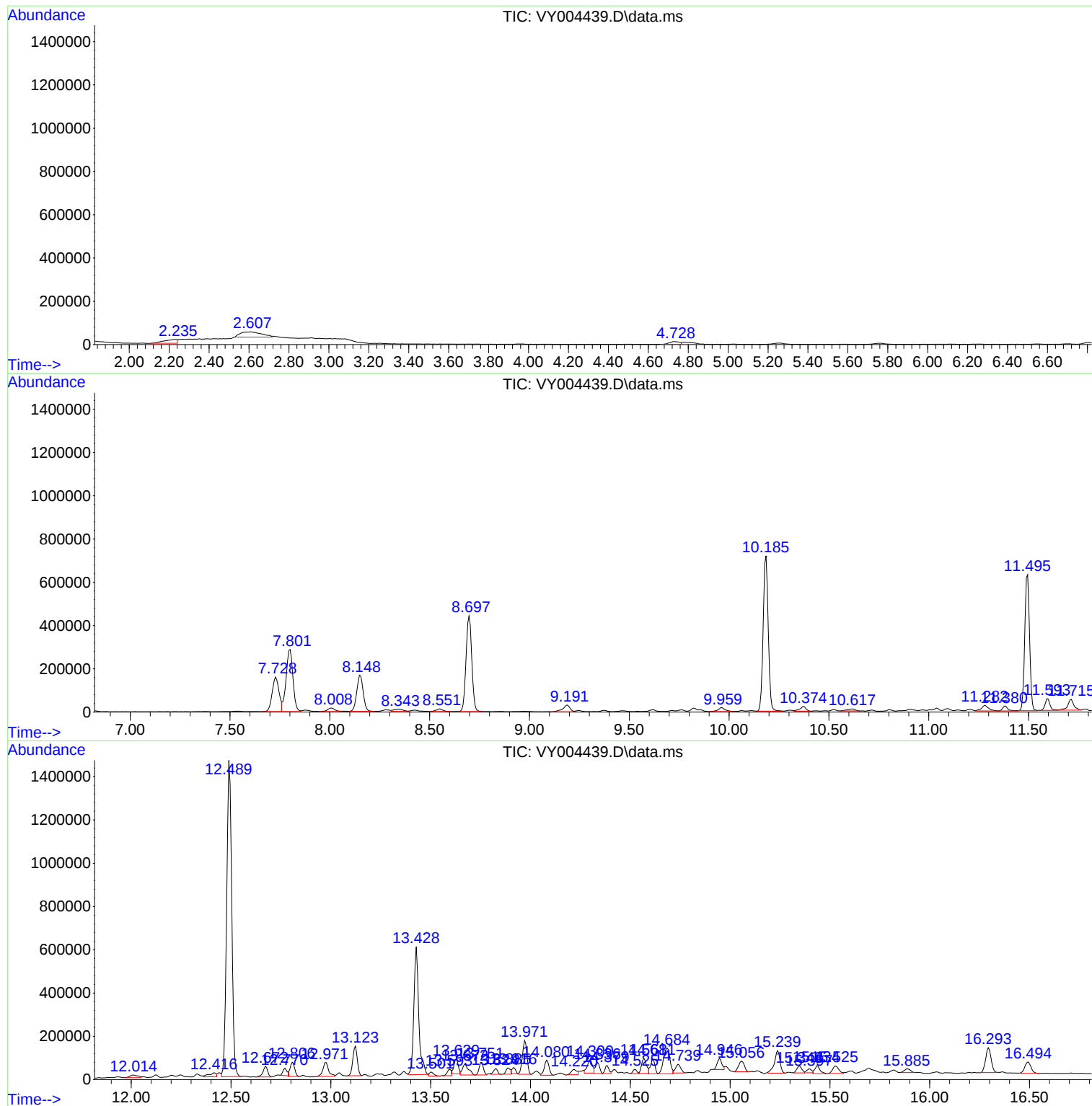
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y041221S.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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Instrument :
MSVOA_Y
ClientSampleId :
FILL-C-2

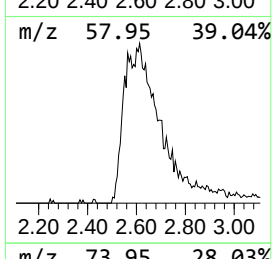
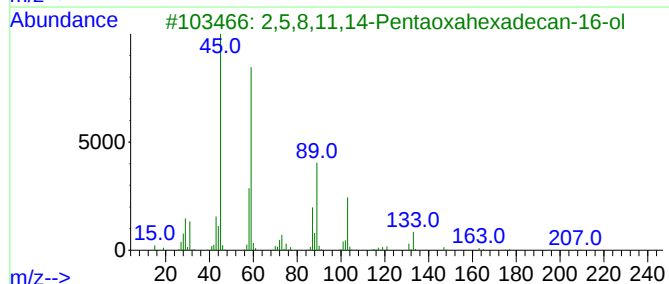
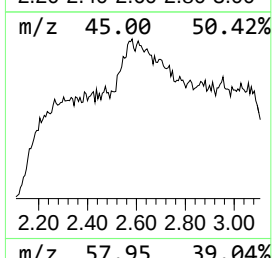
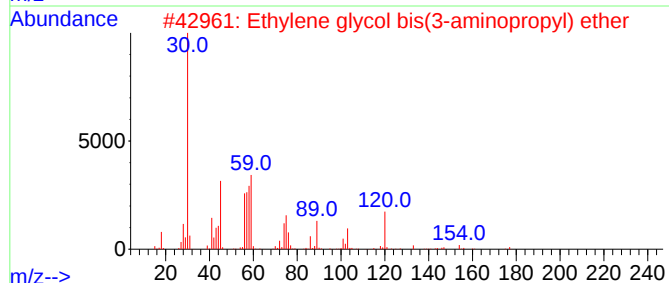
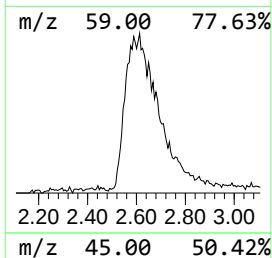
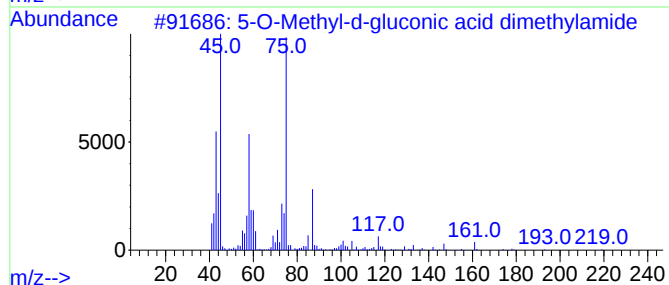
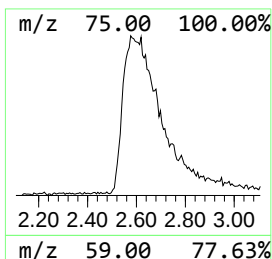
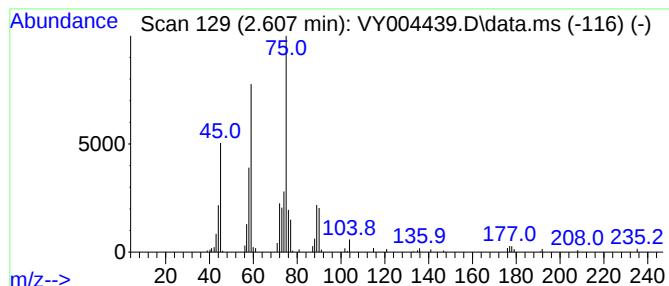
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown2.607 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.607	13.11 ug/l	177951	Pentafluorobenzene	7.801

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-O-Methyl-d-gluconic acid dimet...	237	C9H19NO6	013096-67-8	37
2			Ethylene glycol bis(3-aminopropy...	176	C8H20N2O2	002997-01-5	25
3			2,5,8,11,14-Pentaoxahexadecan-16-ol	252	C11H24O6	023778-52-1	25
4			Propane, 1,2,3-trimethoxy-	134	C6H14O3	020637-49-4	23
5			Silane, ethyldimethyl-	88	C4H12Si	000758-21-4	17



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ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
FILL-C-2

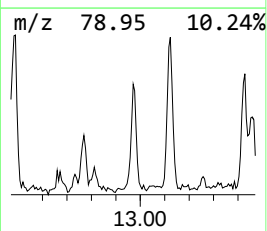
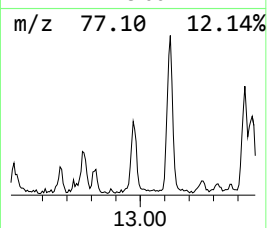
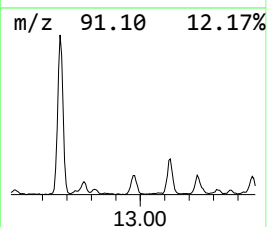
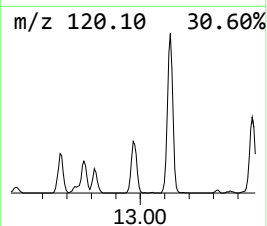
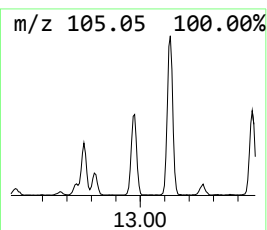
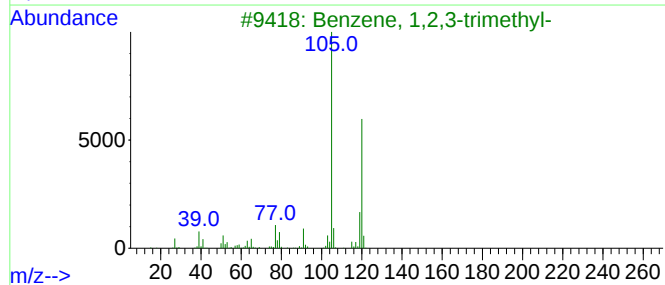
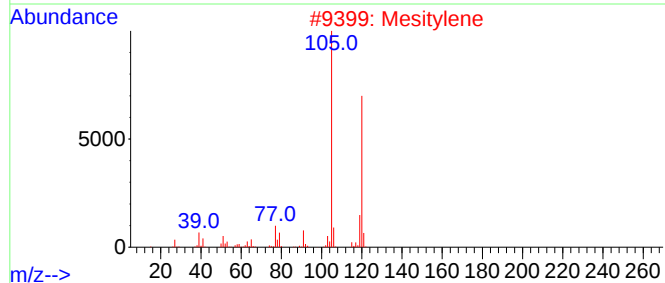
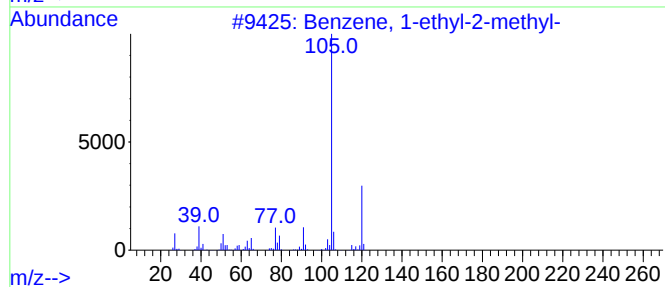
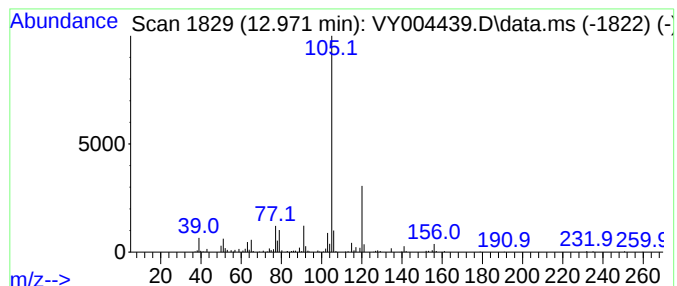
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y041221S.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 1-ethyl-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.971	6.33 ug/l	127898	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2			Mesitylene	120	C9H12	000108-67-8	91
3			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
4			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91



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Misc : 5.93G/5ML/MSVOA_Y/SOIL
ALS Vial : 12 Sample Multiplier: 1

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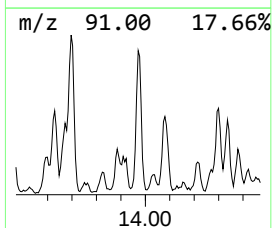
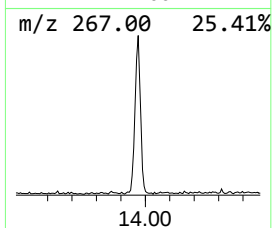
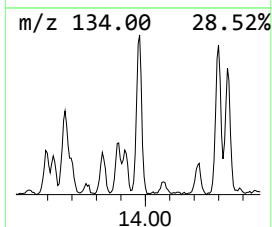
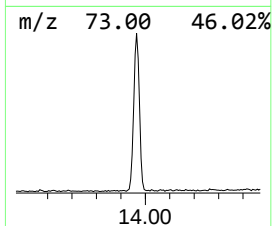
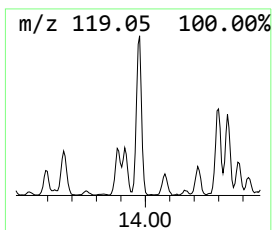
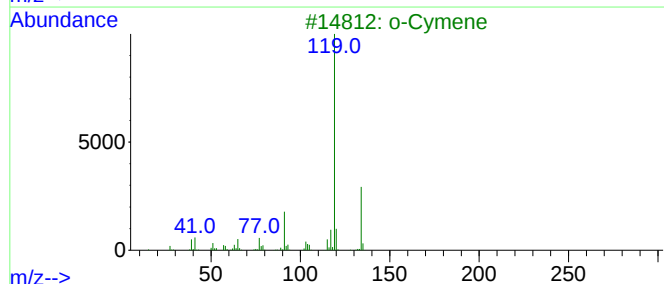
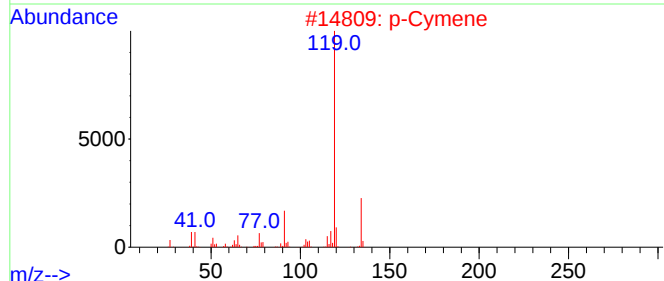
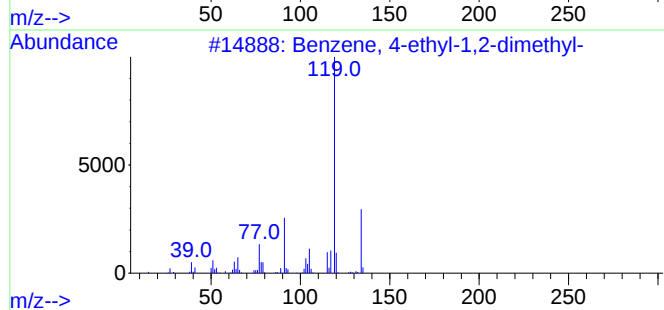
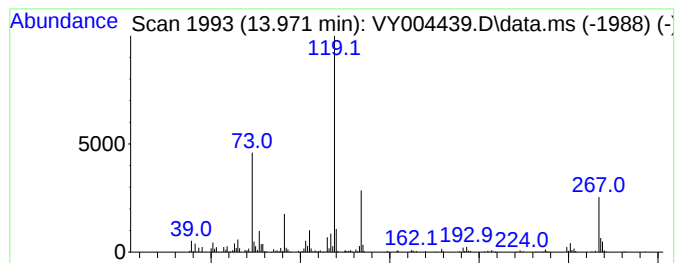
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y041221S.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.970	12.07 ug/l	243676	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2			p-Cymene	134	C10H14	000099-87-6	93
3			o-Cymene	134	C10H14	000527-84-4	93
4			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	93
5			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	92



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Instrument :
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ClientSampleId :
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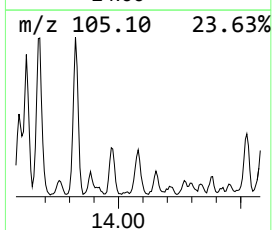
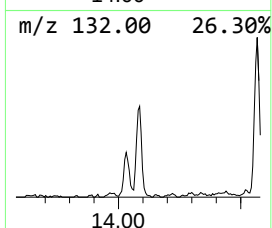
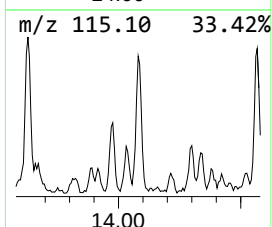
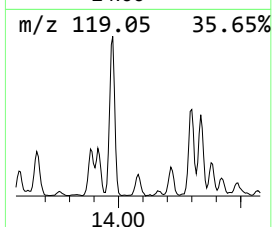
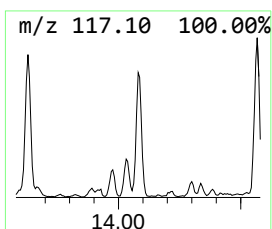
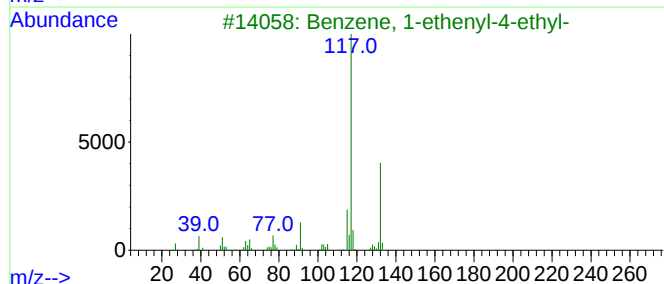
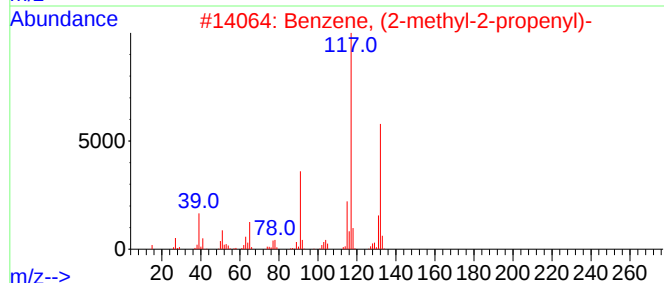
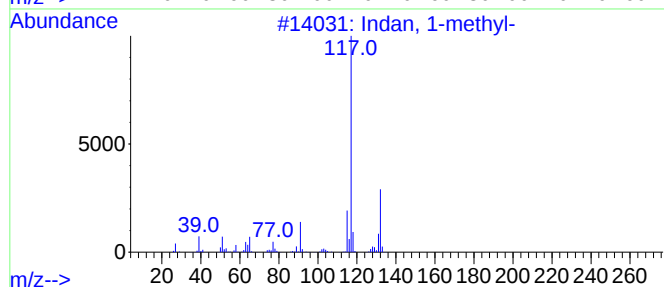
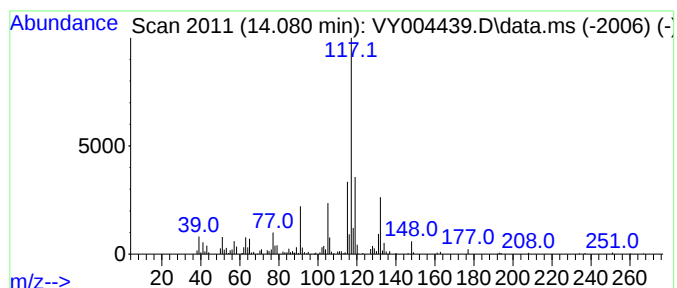
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y041221S.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Indan, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.080	5.56 ug/l	112308	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indan, 1-methyl-	132	C10H12	000767-58-8	76
2			Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	59
3			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	52
4			(E)-1-Phenyl-1-butene	132	C10H12	001005-64-7	49
5			Benzene, (1-ethyl-2-propenyl)-	146	C11H14	019947-22-9	47



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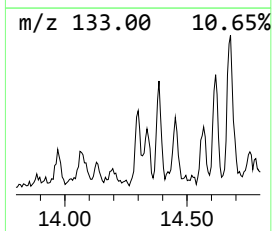
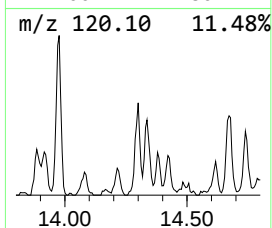
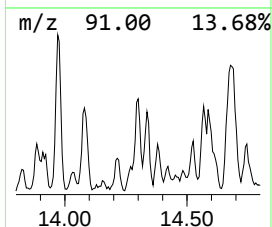
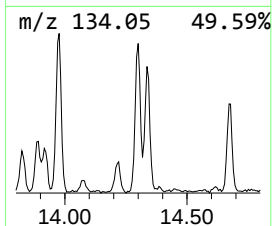
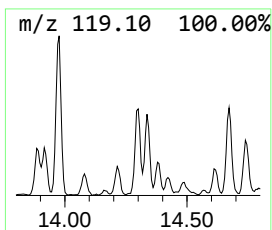
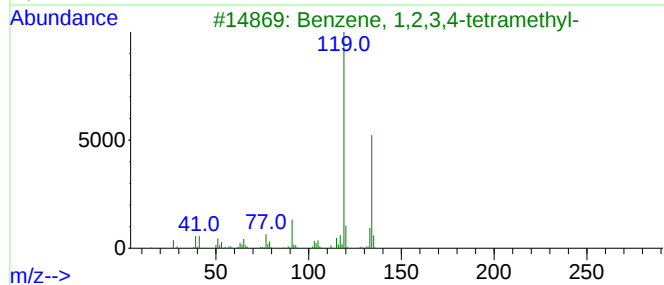
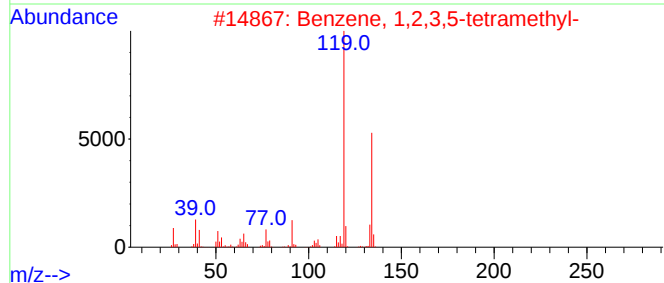
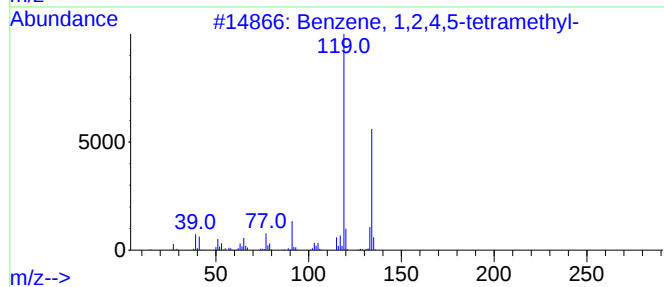
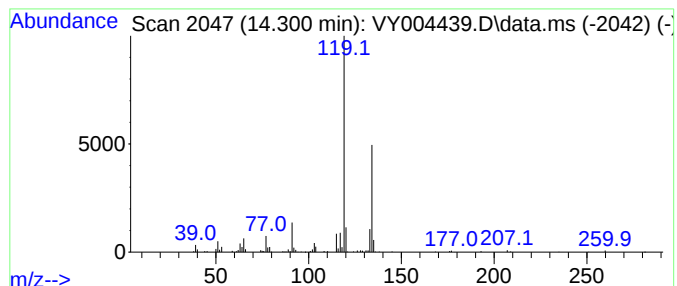
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.300	5.52 ug/l	111435	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
3			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
4			o-Cymene	134	C10H14	000527-84-4	94
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041521\
Data File : VY004439.D
Acq On : 15 Apr 2021 14:34
Operator : SY/MD
Sample : M2040-09
Misc : 5.93G/5ML/MSVOA_Y/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
FILL-C-2

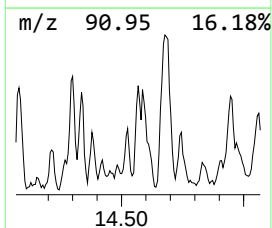
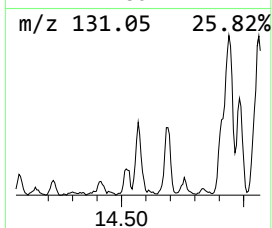
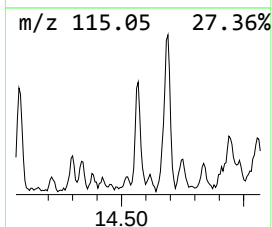
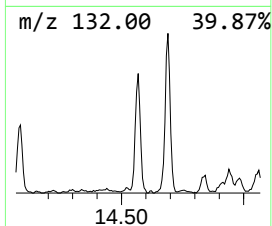
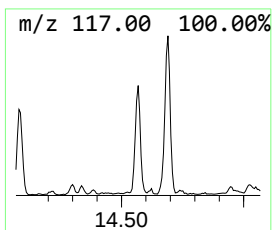
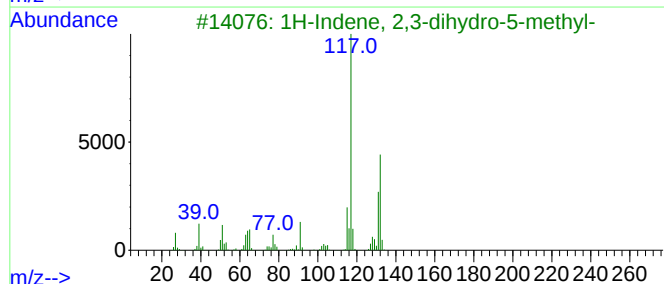
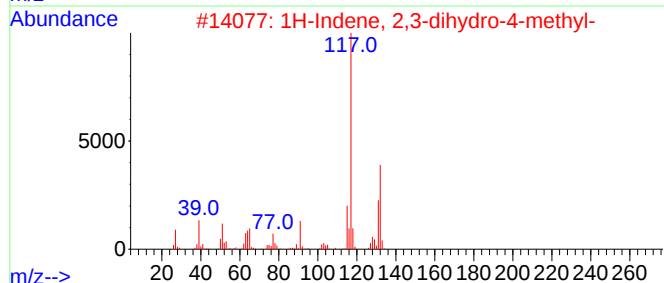
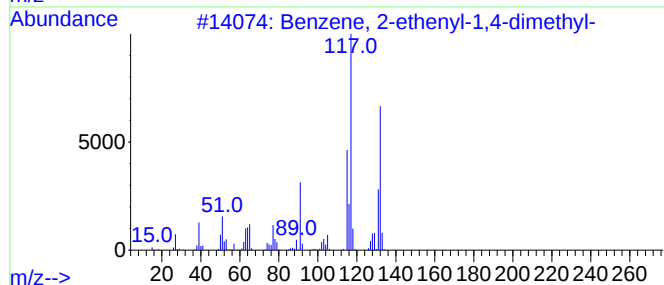
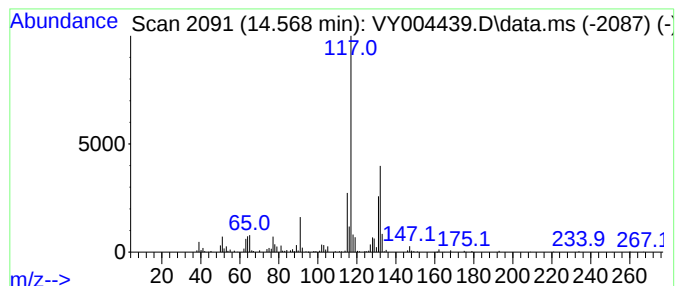
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.568	5.97 ug/l	120593	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92
2			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90
3			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	90
4			1-Phenyl-1-butene	132	C10H12	000824-90-8	87
5			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	83



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041521\
Data File : VY004439.D
Acq On : 15 Apr 2021 14:34
Operator : SY/MD
Sample : M2040-09
Misc : 5.93G/5ML/MSVOA_Y/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
FILL-C-2

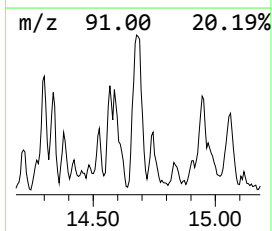
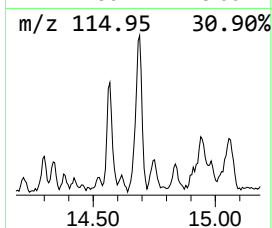
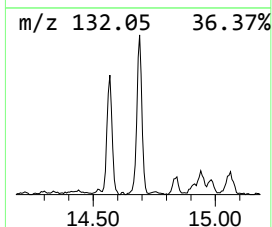
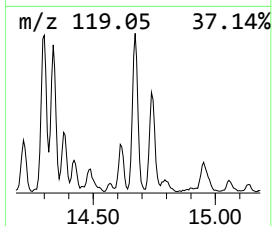
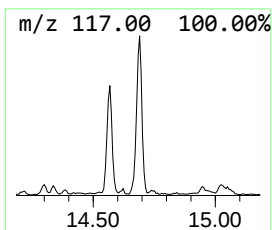
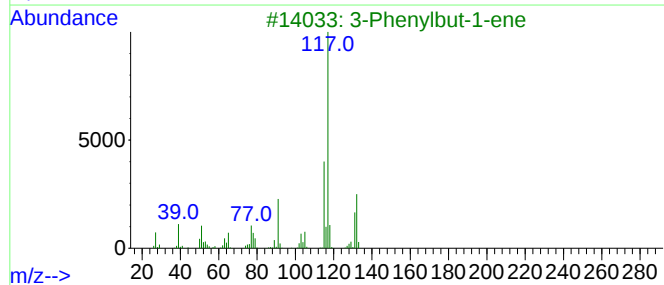
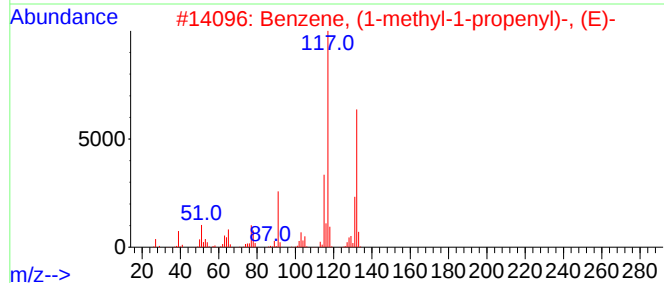
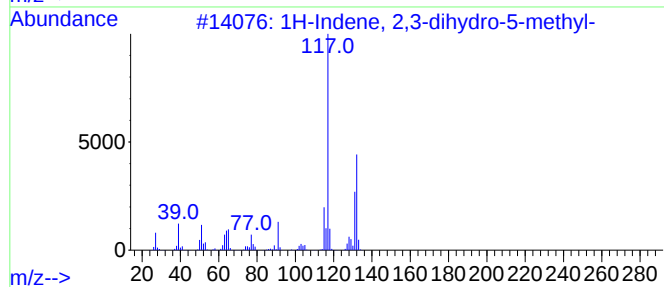
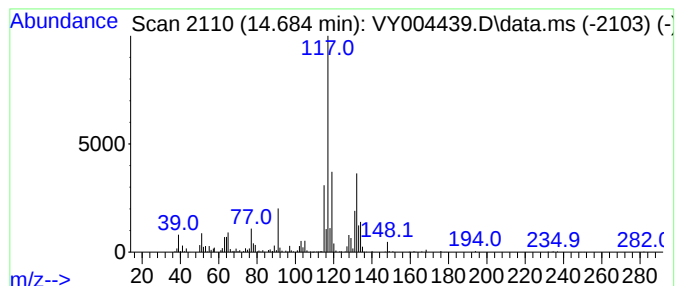
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.684	13.17 ug/l	265935	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
2			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	90
3			3-Phenylbut-1-ene	132	C10H12	000934-10-1	81
4			1-Phenyl-1-butene	132	C10H12	000824-90-8	76
5			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	76



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041521\
Data File : VY004439.D
Acq On : 15 Apr 2021 14:34
Operator : SY/MD
Sample : M2040-09
Misc : 5.93G/5ML/MSVOA_Y/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
FILL-C-2

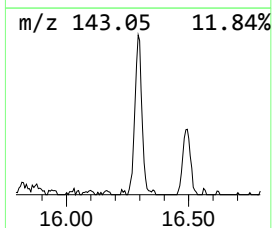
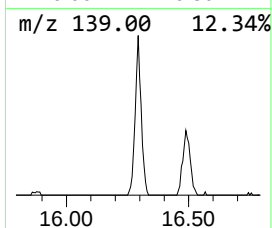
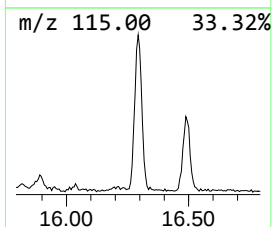
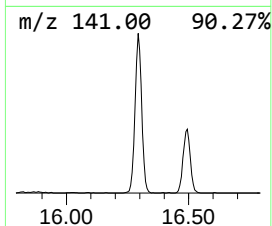
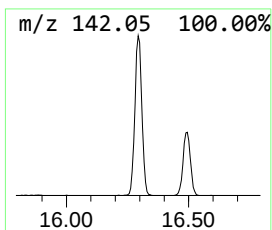
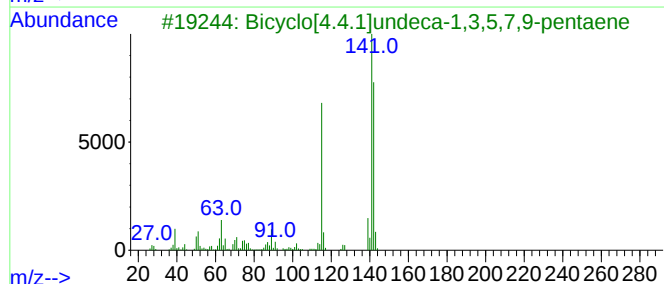
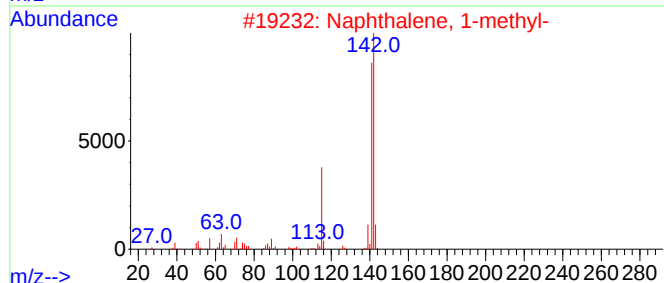
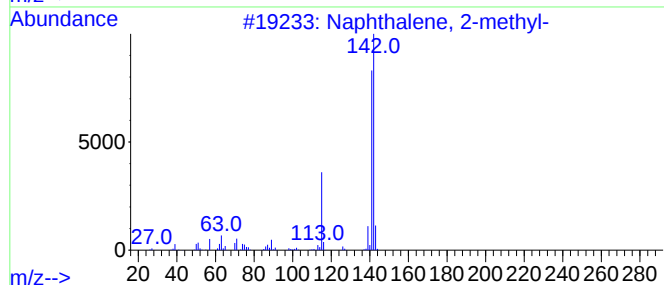
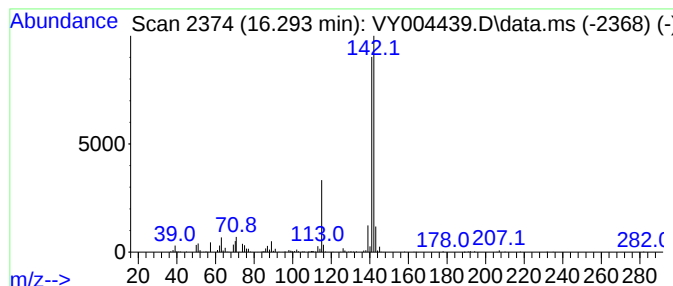
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y041221S.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Naphthalene, 2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.293	10.88 ug/l	219756	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			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
5			Benzocycloheptatriene	142	C11H10	000264-09-5	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041521\
Data File : VY004439.D
Acq On : 15 Apr 2021 14:34
Operator : SY/MD
Sample : M2040-09
Misc : 5.93G/5ML/MSVOA_Y/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
FILL-C-2

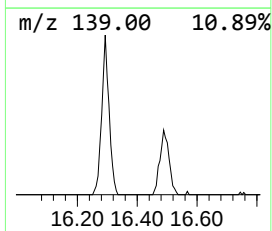
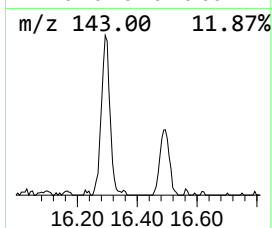
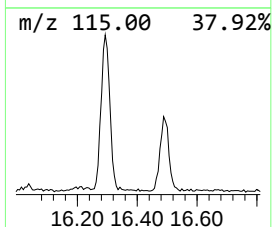
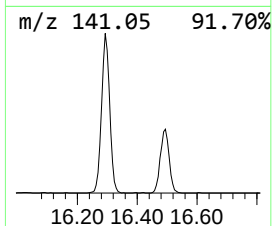
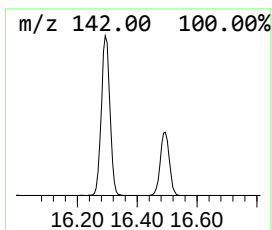
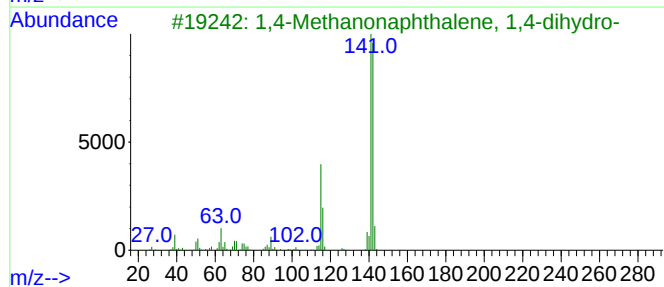
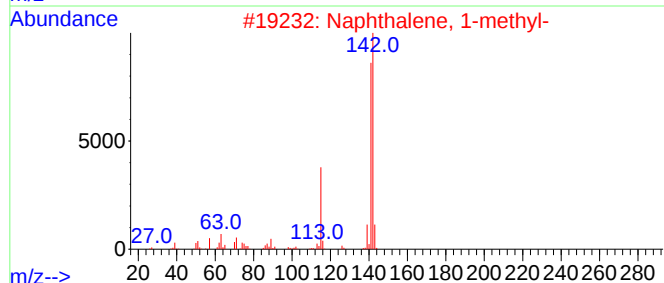
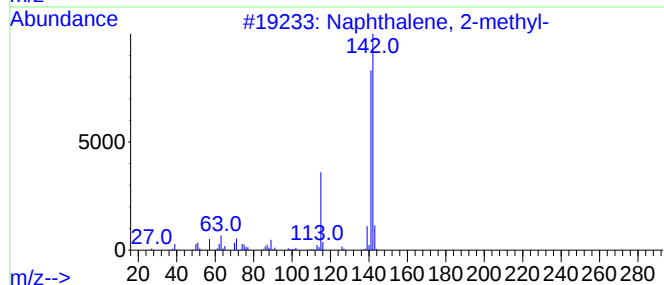
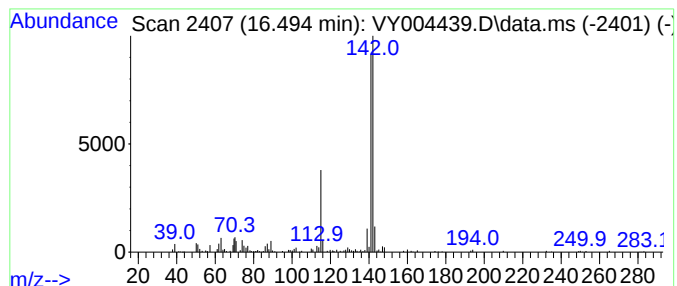
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y041221S.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.494	5.66 ug/l	114277	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	97
2			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	97
3			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
4			Benzocycloheptatriene	142	C11H10	000264-09-5	91
5			3,4-Difluorobenzaldehyde	142	C7H4F2O	034036-07-2	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041521\
Data File : VY004439.D
Acq On : 15 Apr 2021 14:34
Operator : SY/MD
Sample : M2040-09
Misc : 5.93G/5ML/MSVOA_Y/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
FILL-C-2

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y041221S.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Benzene, 1-ethy...	12.971	6.3	ug/l	127898	4	13.428	1009810	50.0
Benzene, 4-ethy...	13.970	12.1	ug/l	243676	4	13.428	1009810	50.0
Indan, 1-methyl-	14.080	5.6	ug/l	112308	4	13.428	1009810	50.0
Benzene, 1,2,4,...	14.300	5.5	ug/l	111435	4	13.428	1009810	50.0
Benzene, 2-ethe...	14.568	6.0	ug/l	120593	4	13.428	1009810	50.0
1H-Indene, 2,3-...	14.684	13.2	ug/l	265935	4	13.428	1009810	50.0
Naphthalene, 2-...	16.293	10.9	ug/l	219756	4	13.428	1009810	50.0
Naphthalene, 1-...	16.494	5.7	ug/l	114277	4	13.428	1009810	50.0