

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY100323\
 Data File : VY015815.D
 Acq On : 03 Oct 2023 18:53
 Operator : SY/MD
 Sample : 04704-03
 Misc : 4.93g/5.0mL/MSVOA_Y/SOIL
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 R-2(5-10)

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\82Y091923S.M
 Title : SW846 8260

Signal : TIC: VY015815.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.186	377	388	400	rVB	19096	51522	1.16%	0.210%
2	4.710	461	474	488	rBV	50432	161429	3.64%	0.659%
3	5.741	633	643	656	rBV4	21579	67835	1.53%	0.277%
4	7.716	957	967	972	rBV	130975	311351	7.02%	1.270%
5	7.783	972	978	989	rVB	234253	529916	11.95%	2.162%
6	8.143	1026	1037	1051	rBV2	179081	462894	10.44%	1.888%
7	8.685	1117	1126	1140	rBV	366613	713422	16.09%	2.910%
8	10.179	1363	1371	1377	rBV	572779	1005344	22.67%	4.101%
9	10.240	1377	1381	1388	rVB	54707	96585	2.18%	0.394%
10	11.490	1579	1586	1594	rBV	530407	859856	19.39%	3.508%
11	11.593	1596	1603	1609	rBV	85262	163407	3.69%	0.667%
12	11.697	1615	1620	1628	rVB	67169	120409	2.72%	0.491%
13	12.026	1669	1674	1687	rVB2	148709	287666	6.49%	1.173%
14	12.337	1713	1725	1727	rBV4	112039	275531	6.21%	1.124%
15	12.477	1743	1748	1769	rVB	479704	933786	21.06%	3.809%
16	12.733	1783	1790	1793	rBV	53768	105755	2.38%	0.431%
17	12.806	1799	1802	1806	rBV	50164	71463	1.61%	0.292%
18	12.934	1806	1823	1834	rVV4	252972	1469702	33.14%	5.995%
19	13.062	1835	1844	1849	rVV2	135512	502332	11.33%	2.049%
20	13.117	1849	1853	1858	rVB	189812	316698	7.14%	1.292%
21	13.276	1867	1879	1880	rBV10	14950	56719	1.28%	0.231%
22	13.422	1897	1903	1907	rBV	466982	782932	17.66%	3.194%
23	13.556	1913	1925	1930	rBV5	50329	216330	4.88%	0.882%
24	13.623	1931	1936	1951	rVB2	499401	1054976	23.79%	4.304%
25	13.806	1961	1966	1973	rVB	201391	311967	7.04%	1.273%
26	13.910	1980	1983	1987	rVB2	31202	45549	1.03%	0.186%
27	13.971	1987	1993	1996	rBV3	76757	152393	3.44%	0.622%
28	14.074	2006	2010	2013	rBV2	40240	61007	1.38%	0.249%
29	14.209	2027	2032	2035	rBV2	46613	85392	1.93%	0.348%
30	14.263	2036	2041	2043	rVV3	80113	173677	3.92%	0.708%
31	14.288	2043	2045	2049	rVV2	94448	158452	3.57%	0.646%
32	14.330	2049	2052	2057	rVV	128881	195883	4.42%	0.799%
33	14.434	2063	2069	2075	rVV5	22045	52356	1.18%	0.214%
34	14.562	2085	2090	2094	rVB	62371	90526	2.04%	0.369%
35	14.684	2101	2110	2113	rBV3	285698	697235	15.72%	2.844%
36	14.745	2116	2120	2128	rVB	271537	528898	11.93%	2.158%

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37	14.830	2128	2134	2142	rBV	311648	545527	12.30%	2.225%
38	14.940	2147	2152	2164	rVV6	89674	246085	5.55%	1.004%
39	15.050	2165	2170	2177	rVB3	49152	103262	2.33%	0.421%
40	15.233	2190	2200	2207	rBV	2684621	4434292	100.00%	18.089%
41	15.330	2212	2216	2221	rBV2	61708	101172	2.28%	0.413%
42	15.397	2222	2227	2242	rVB4	69825	261651	5.90%	1.067%
43	15.519	2242	2247	2252	rBV2	31484	59337	1.34%	0.242%
44	15.660	2265	2270	2272	rBV4	30407	50013	1.13%	0.204%
45	15.714	2274	2279	2283	rVV4	59928	122022	2.75%	0.498%
46	15.757	2283	2286	2291	rVB	52973	86143	1.94%	0.351%
47	15.830	2291	2298	2303	rBV4	78033	189031	4.26%	0.771%
48	16.031	2324	2331	2336	rBV4	29802	65000	1.47%	0.265%
49	16.098	2336	2342	2348	rBV3	29317	81171	1.83%	0.331%
50	16.226	2355	2363	2367	rBV	89212	196599	4.43%	0.802%
51	16.287	2367	2373	2387	rVB	750341	1520438	34.29%	6.202%
52	16.483	2397	2405	2417	rBV	1529963	3222321	72.67%	13.145%
53	16.586	2418	2422	2436	rVB10	17226	58656	1.32%	0.239%

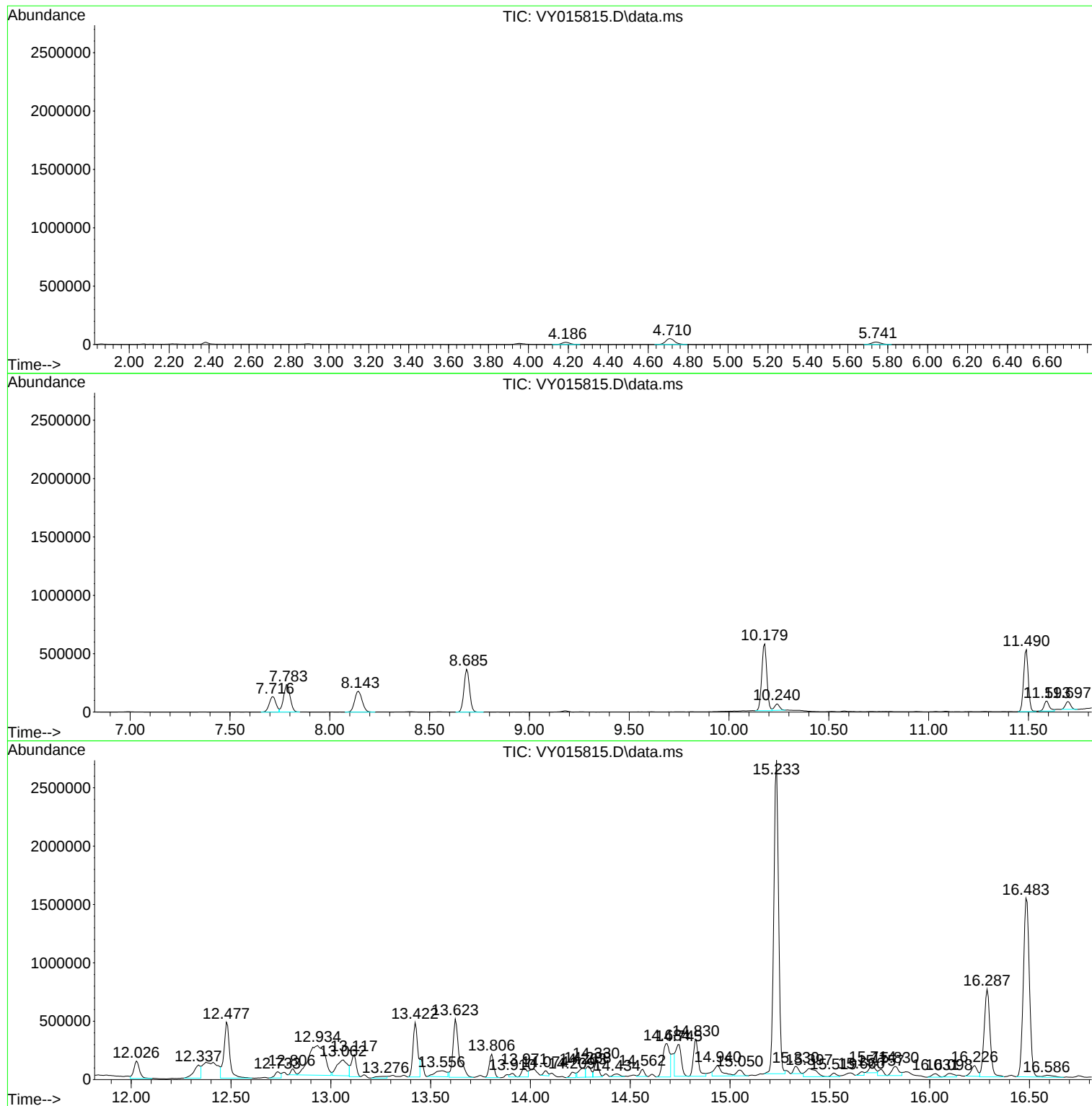
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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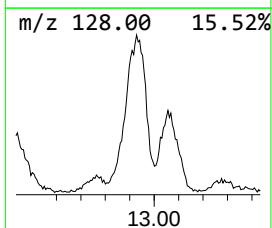
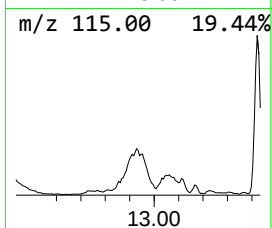
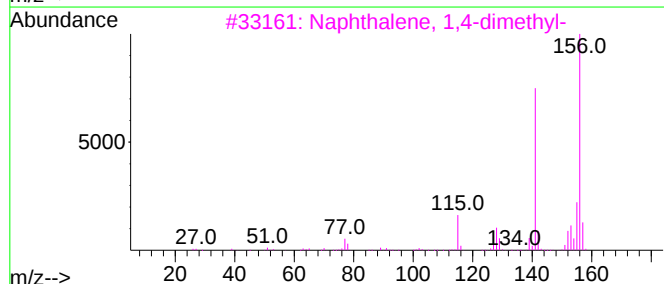
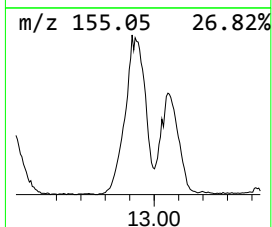
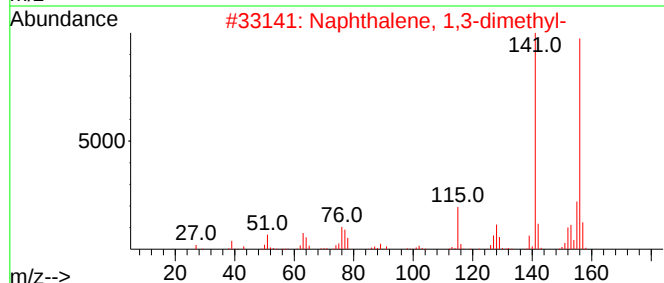
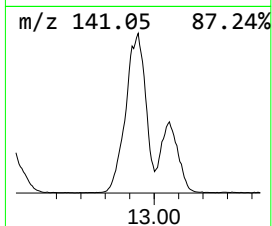
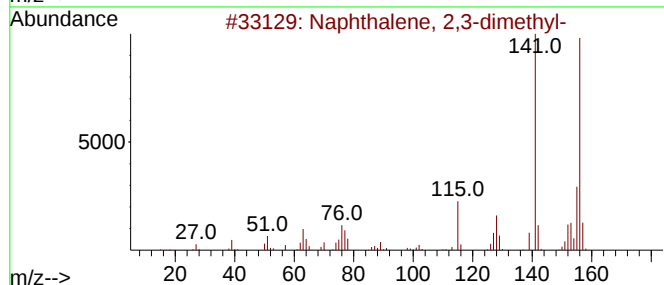
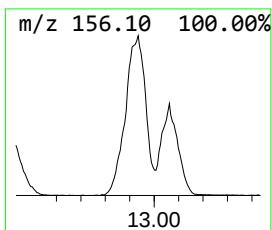
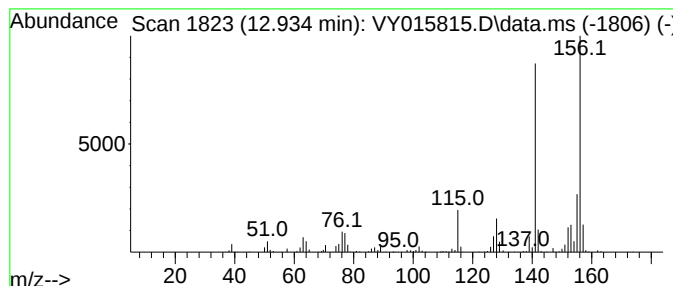
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y091923S.M
Quant Title : SW846 8260

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

Peak Number 1 Naphthalene, 2,3-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.934	93.86 ug/l	1469700	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
2			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	97
3			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97
4			Naphthalene, 1,2-dimethyl-	156	C12H12	000573-98-8	97
5			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96



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Instrument :
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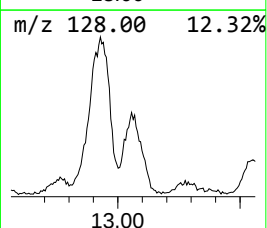
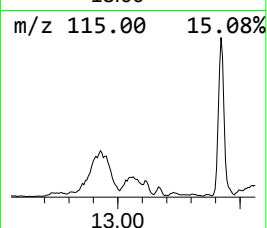
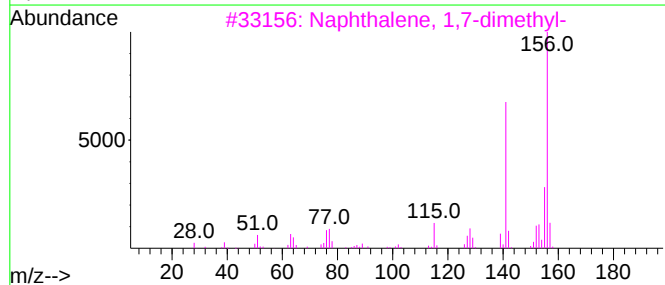
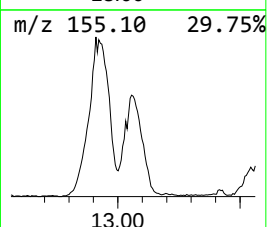
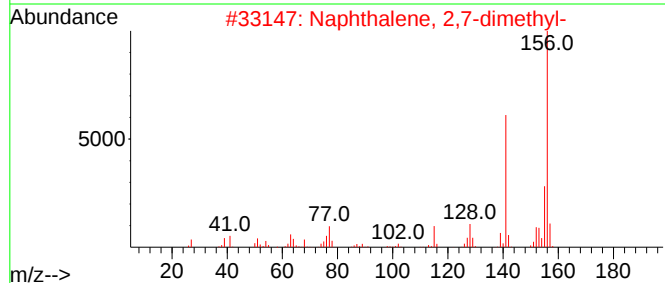
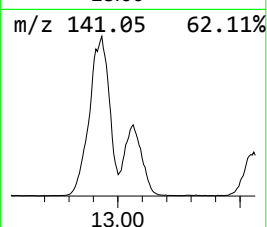
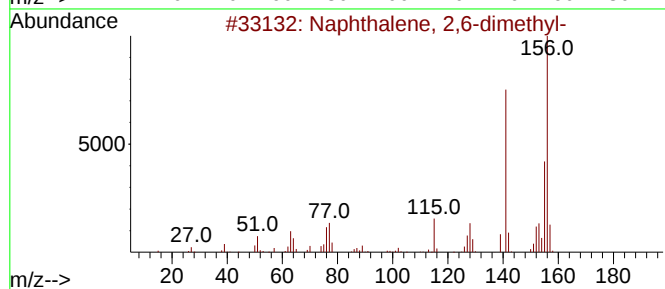
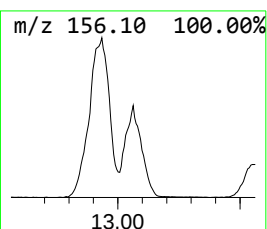
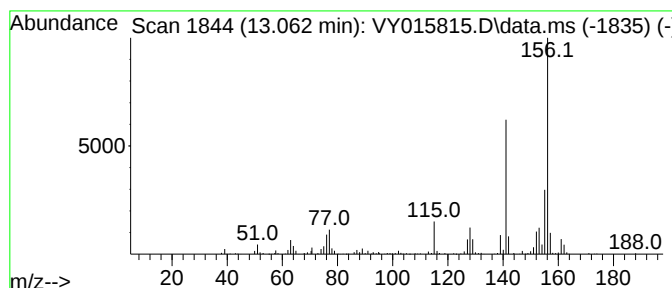
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y091923S.M
Quant Title : SW846 8260

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

Peak Number 2 Naphthalene, 2,6-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.062	32.08 ug/l	502332	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	98
2			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
3			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96
4			Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	96
5			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	96



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ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
R-2(5-10)

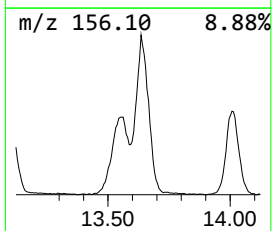
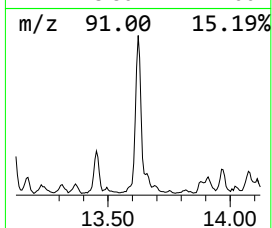
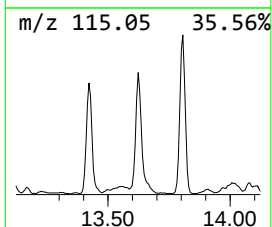
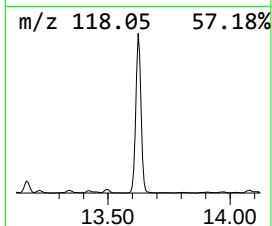
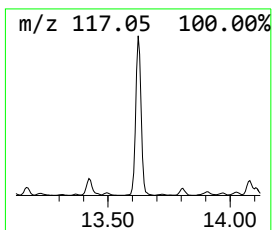
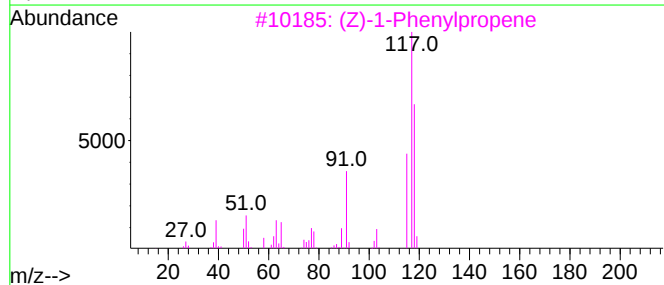
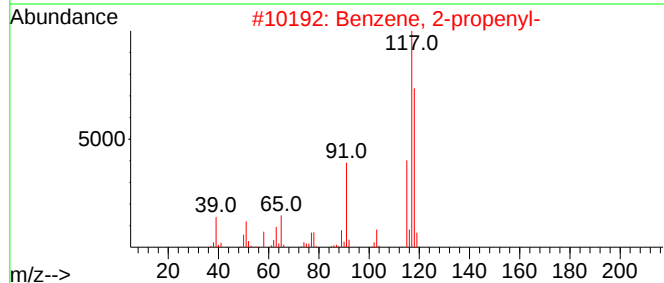
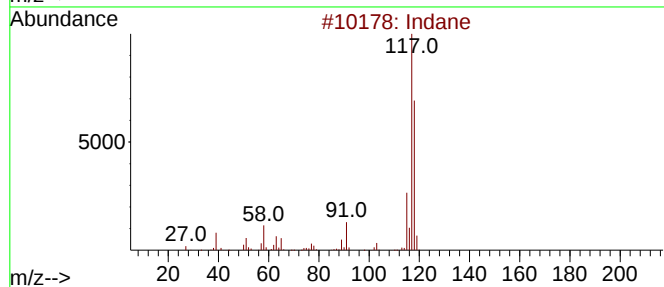
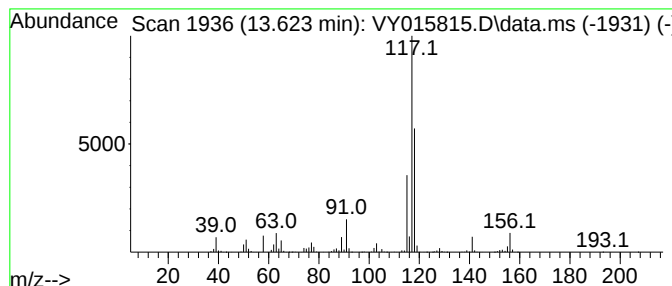
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Indane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.623	67.37 ug/l	1054980	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	90
2			Benzene, 2-propenyl-	118	C9H10	000300-57-2	72
3			(Z)-1-Phenylpropene	118	C9H10	000766-90-5	68
4			Benzene, 1-(chloromethyl)-4-ethe...	152	C9H9Cl	001592-20-7	53
5			Benzene, 1-(chloromethyl)-3-ethe...	152	C9H9Cl	039833-65-3	53



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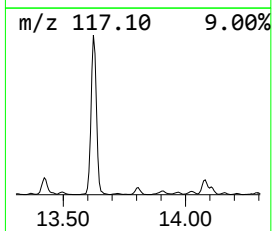
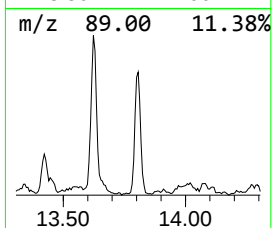
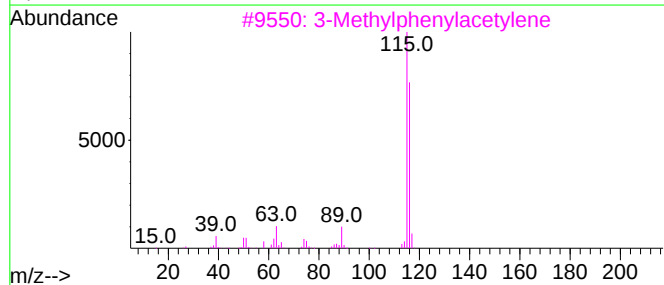
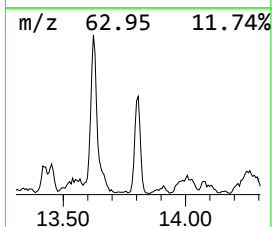
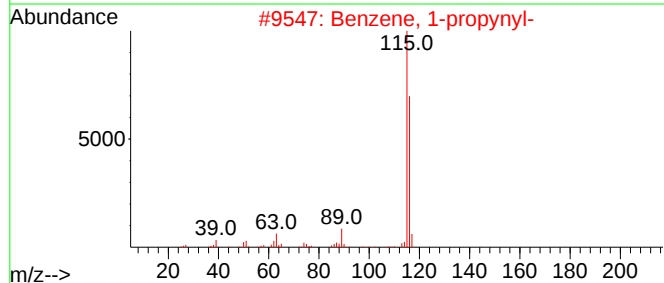
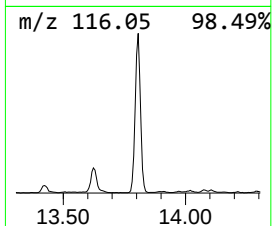
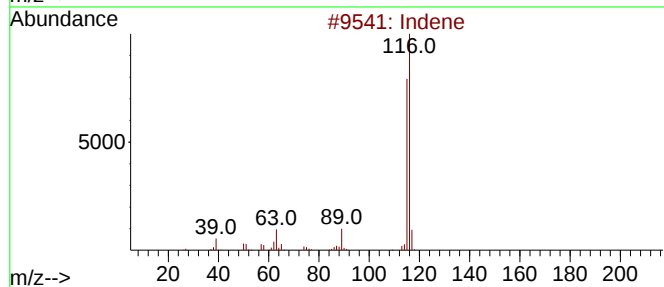
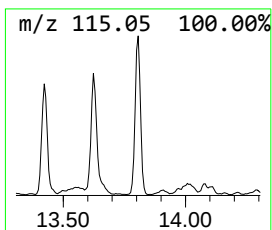
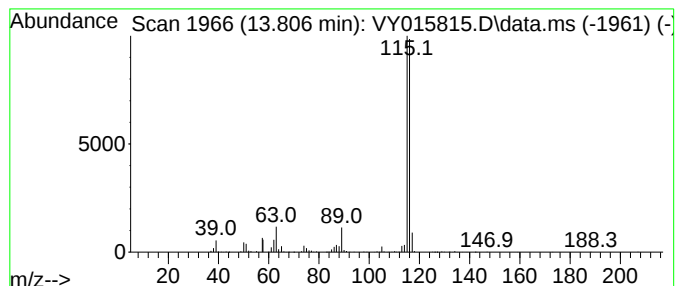
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Quant Title : SW846 8260

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

Peak Number 4 Indene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.806	19.92 ug/l	311967	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indene	116	C9H8	000095-13-6	97
2			Benzene, 1-propynyl-	116	C9H8	000673-32-5	94
3			3-Methylphenylacetylene	116	C9H8	000766-82-5	91
4			Benzene, 1-ethynyl-4-methyl-	116	C9H8	000766-97-2	91
5			2-Methylphenylacetylene	116	C9H8	000766-47-2	83



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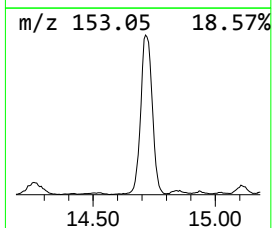
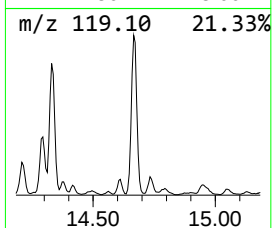
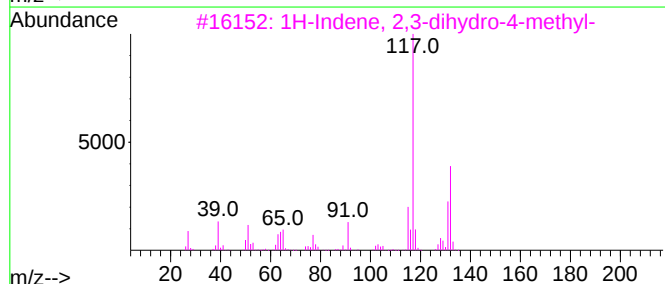
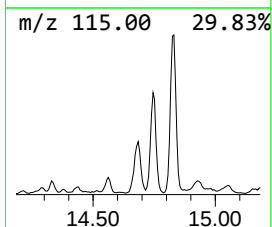
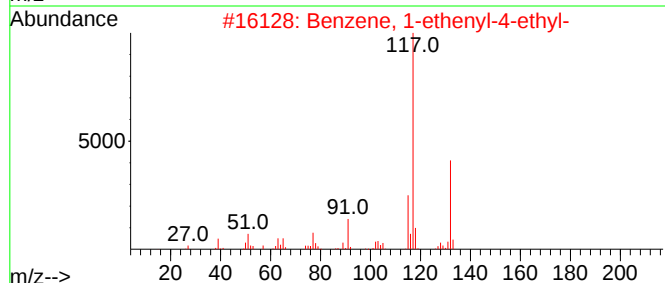
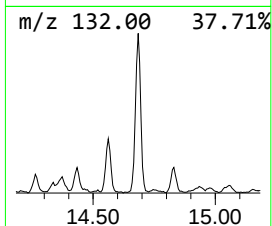
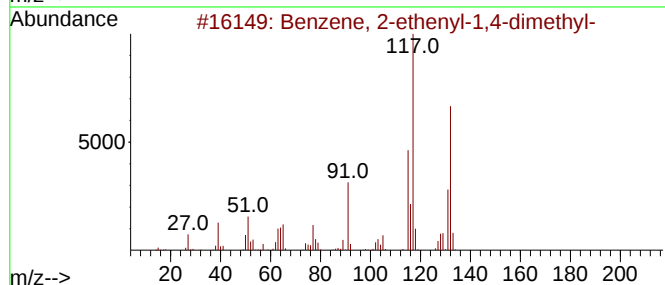
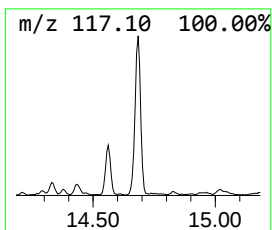
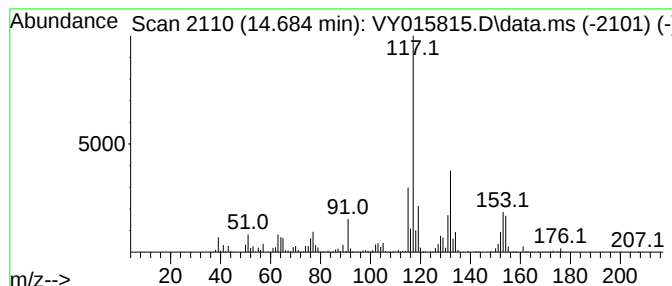
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.684	44.53 ug/l	697235	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	94
2			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	92
3			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	91
4			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	83
5			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	83



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY100323\
Data File : VY015815.D
Acq On : 03 Oct 2023 18:53
Operator : SY/MD
Sample : 04704-03
Misc : 4.93g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
R-2(5-10)

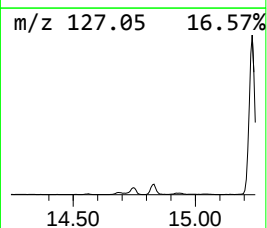
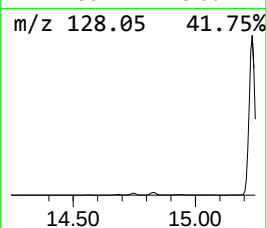
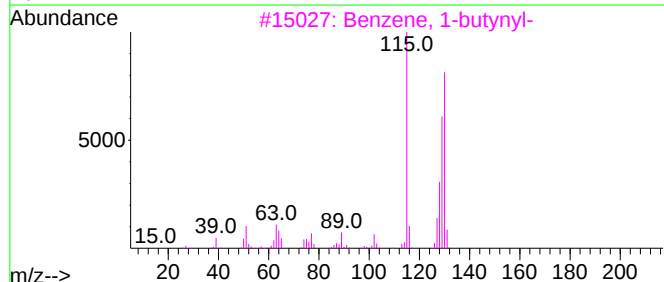
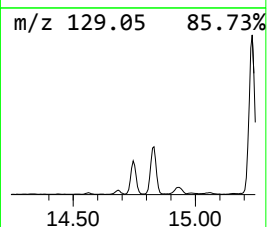
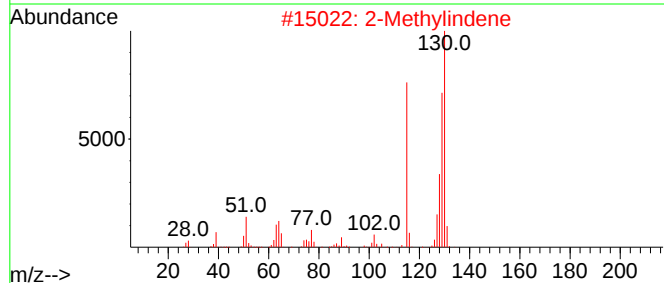
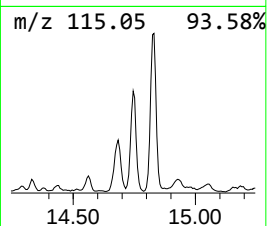
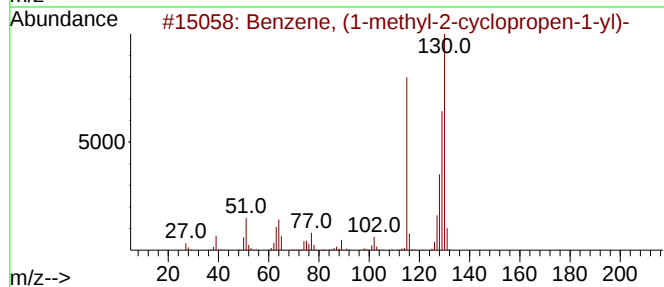
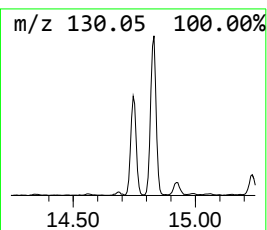
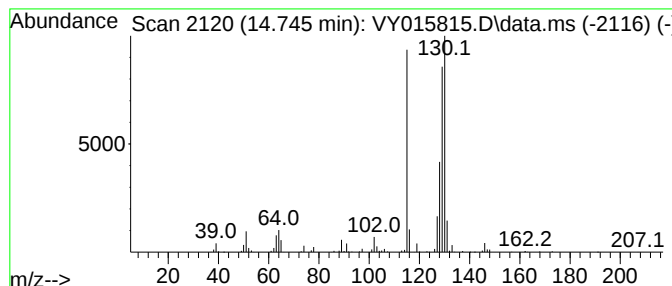
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, (1-methyl-2-cyclop... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.745	33.78 ug/l	528898	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methyl-2-cyclopropen...	130	C10H10	065051-83-4	87
2			2-Methylindene	130	C10H10	002177-47-1	87
3			Benzene, 1-butyryl-	130	C10H10	000622-76-4	87
4			Benzene,1-methyl-1,2-propadienyl-	130	C10H10	022433-39-2	81
5			1H-Indene, 3-methyl-	130	C10H10	000767-60-2	74



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY100323\
Data File : VY015815.D
Acq On : 03 Oct 2023 18:53
Operator : SY/MD
Sample : 04704-03
Misc : 4.93g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
R-2(5-10)

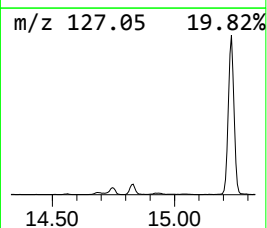
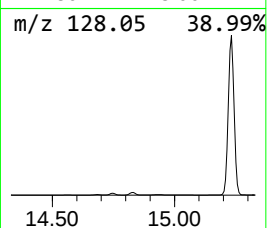
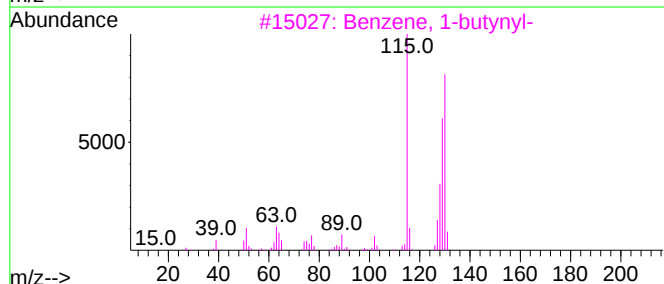
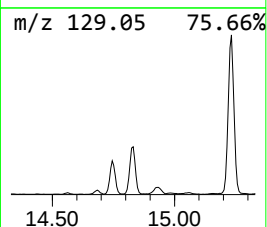
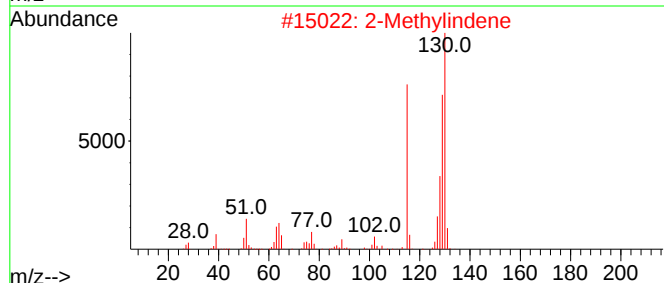
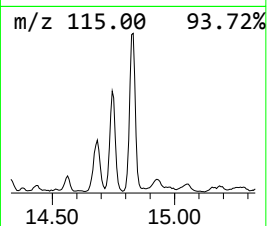
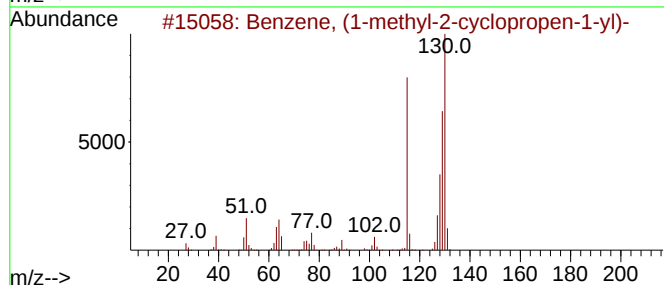
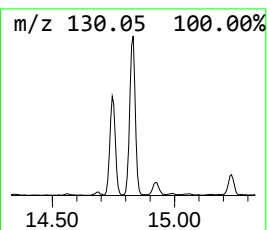
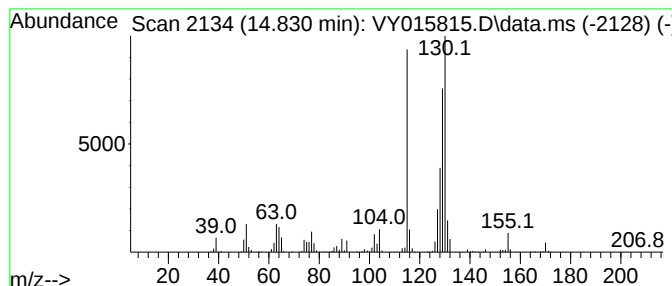
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Quant Title : SW846 8260

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

Peak Number 7 2-Methylindene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.830	34.84 ug/l	545527	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methyl-2-cyclopropen...	130	C10H10	065051-83-4	96
2			2-Methylindene	130	C10H10	002177-47-1	95
3			Benzene, 1-butyryl-	130	C10H10	000622-76-4	95
4			1H-Indene, 1-methyl-	130	C10H10	000767-59-9	94
5			1H-Indene, 3-methyl-	130	C10H10	000767-60-2	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY100323\
Data File : VY015815.D
Acq On : 03 Oct 2023 18:53
Operator : SY/MD
Sample : 04704-03
Misc : 4.93g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
R-2(5-10)

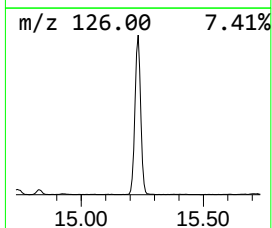
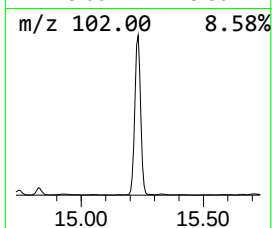
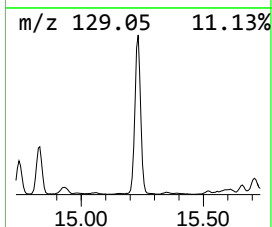
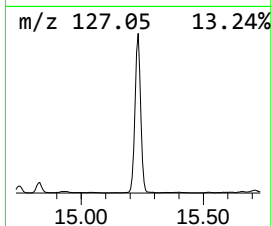
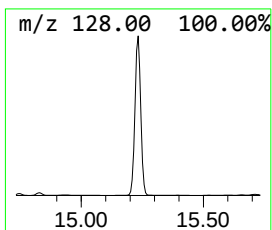
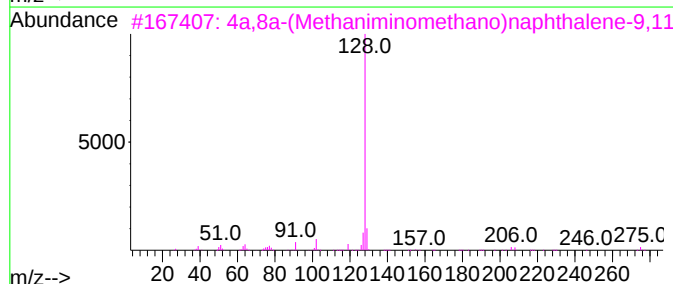
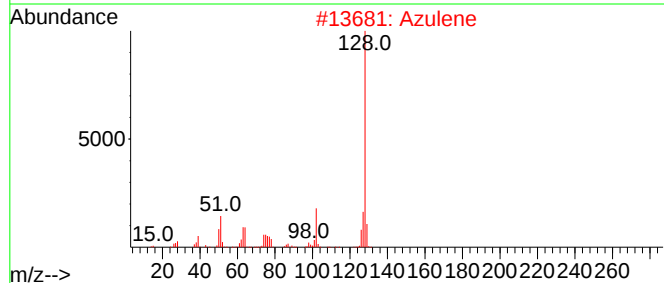
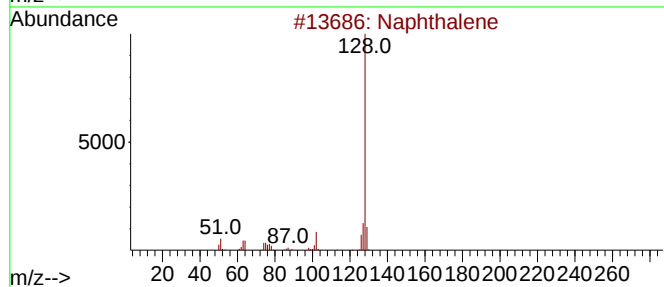
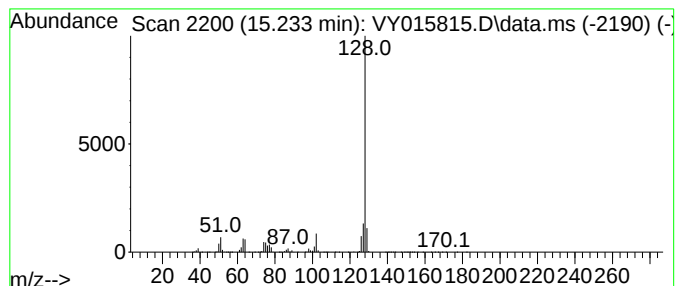
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Quant Title : SW846 8260

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

Peak Number 8 Azulene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.233	283.19 ug/l	4434290	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene	128	C10H8	000091-20-3	95
2			Azulene	128	C10H8	000275-51-4	91
3			4a,8a-(Methaniminomethano)naphth...	275	C18H13NO2	069915-10-2	78
4			1H-Indene, 1-methylene-	128	C10H8	002471-84-3	47
5			1,3-Dicyanobenzene	128	C8H4N2	000626-17-5	45



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY100323\
Data File : VY015815.D
Acq On : 03 Oct 2023 18:53
Operator : SY/MD
Sample : 04704-03
Misc : 4.93g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
R-2(5-10)

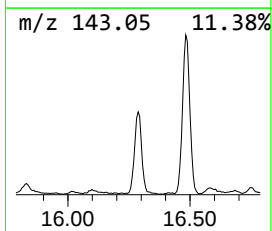
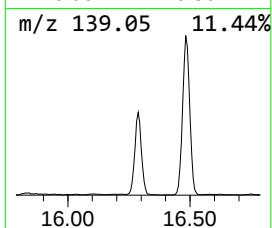
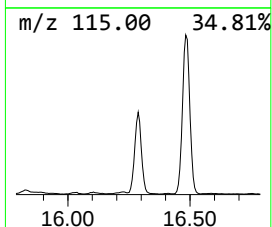
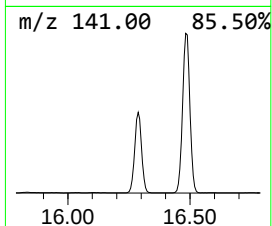
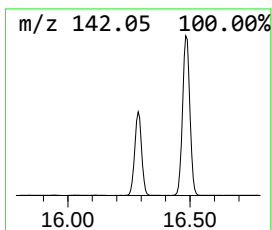
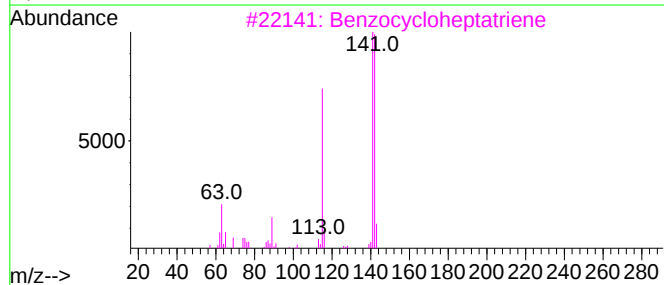
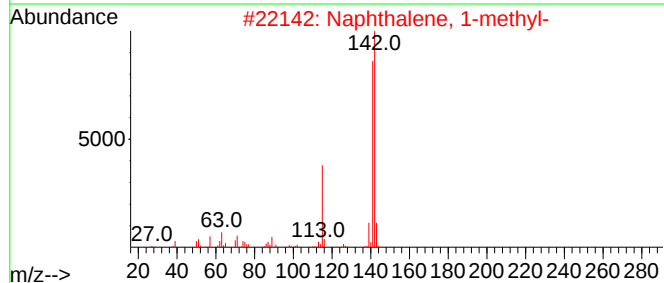
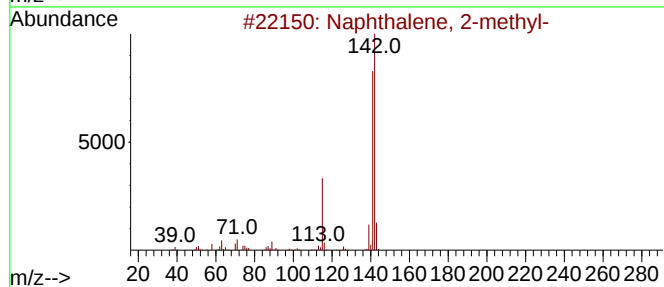
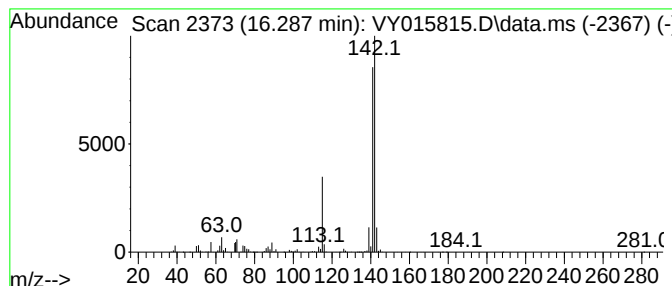
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y091923S.M
Quant Title : SW846 8260

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

Peak Number 9 Benzocycloheptatriene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.287	97.10 ug/l	1520440	1,4-Dichlorobenzene-d4	13.422

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			Benzocycloheptatriene	142	C11H10	000264-09-5	91
4			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
5			Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY100323\
Data File : VY015815.D
Acq On : 03 Oct 2023 18:53
Operator : SY/MD
Sample : 04704-03
Misc : 4.93g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
R-2(5-10)

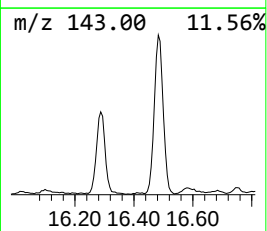
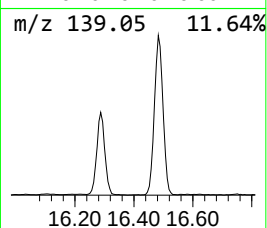
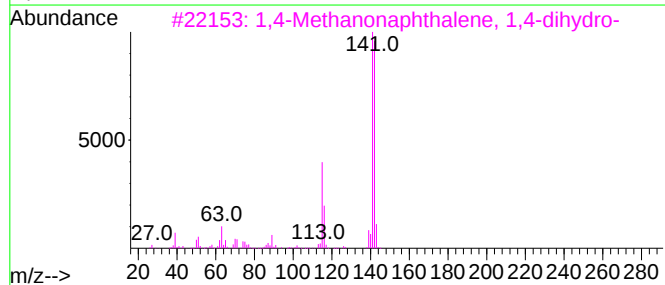
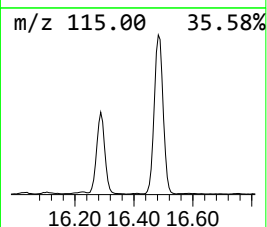
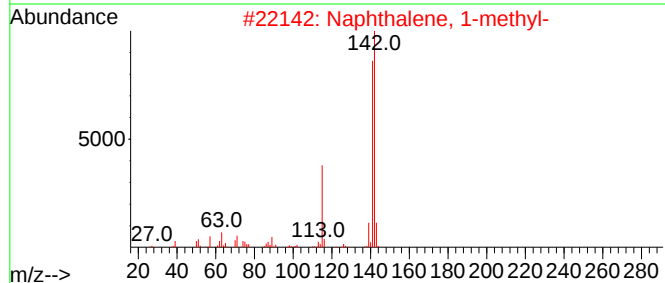
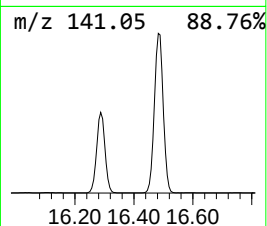
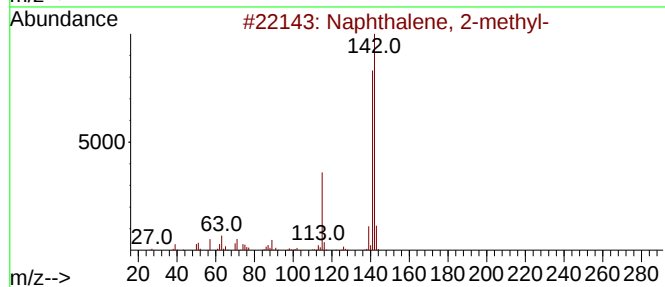
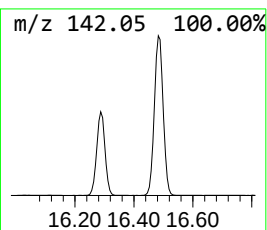
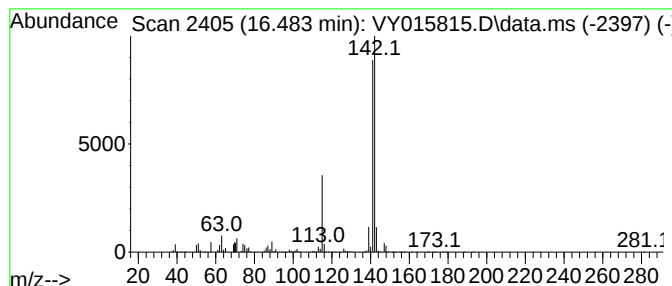
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y091923S.M
Quant Title : SW846 8260

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

Peak Number 10 1,4-Methanonaphthalene, 1,4... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.483	205.79 ug/l	3222320	1,4-Dichlorobenzene-d4	13.422

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		Benzocycloheptatriene	142	C11H10	000264-09-5	91
5		1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY100323\
Data File : VY015815.D
Acq On : 03 Oct 2023 18:53
Operator : SY/MD
Sample : 04704-03
Misc : 4.93g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
R-2(5-10)

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y091923S.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
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Naphthalene, 2,...	13.062	32.1	ug/l	502332	4	13.422	782932	50.0
Indane	13.623	67.4	ug/l	1054980	4	13.422	782932	50.0
Indene	13.806	19.9	ug/l	311967	4	13.422	782932	50.0
Benzene, 2-ethe...	14.684	44.5	ug/l	697235	4	13.422	782932	50.0
Benzene, (1-met...	14.745	33.8	ug/l	528898	4	13.422	782932	50.0
2-Methylindene	14.830	34.8	ug/l	545527	4	13.422	782932	50.0
Azulene	15.233	283.2	ug/l	4434290	4	13.422	782932	50.0
Benzocyclohepta...	16.287	97.1	ug/l	1520440	4	13.422	782932	50.0
1,4-Methanonaph...	16.483	205.8	ug/l	3222320	4	13.422	782932	50.0