

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY030322\
 Data File : VY007715.D
 Acq On : 03 Mar 2022 19:26
 Operator : SY/MD
 Sample : N1762-02
 Misc : 7.44g/5.0mL/MSVOA_Y/SOIL
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 282

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\82Y022822S.M

Title : SW846 8260

Signal : TIC: VY007715.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	7.722	955	968	973	rBV	393522	949453	2.07%	0.577%
2	7.789	973	979	988	rVB	588477	1338743	2.92%	0.813%
3	8.143	1028	1037	1048	rBV	363318	798228	1.74%	0.485%
4	8.691	1118	1127	1139	rBV	969838	1874753	4.08%	1.138%
5	10.179	1364	1371	1385	rBV	1539908	2631940	5.73%	1.598%
6	11.490	1578	1586	1596	rBV	1508361	2426538	5.29%	1.474%
7	11.697	1612	1620	1631	rBV	540299	1172744	2.55%	0.712%
8	12.331	1713	1724	1730	rBV2	511573	1029135	2.24%	0.625%
9	12.489	1743	1750	1755	rBV2	2384682	4297804	9.36%	2.610%
10	12.672	1773	1780	1784	rVB	311404	512147	1.12%	0.311%
11	12.733	1784	1790	1793	rBV	1050723	1646459	3.59%	1.000%
12	12.764	1793	1795	1799	rVV	547325	786960	1.71%	0.478%
13	12.812	1799	1803	1810	rVB	862946	1332137	2.90%	0.809%
14	13.117	1848	1853	1859	rVB	2359672	3393993	7.39%	2.061%
15	13.251	1868	1875	1882	rBV	1083102	1905690	4.15%	1.157%
16	13.422	1897	1903	1907	rBV	1511161	2487955	5.42%	1.511%
17	13.593	1925	1931	1934	rBV	2363750	3699358	8.06%	2.246%
18	13.623	1934	1936	1939	rVV	1709740	2219304	4.83%	1.348%
19	13.666	1939	1943	1952	rVB2	1921306	3488058	7.60%	2.118%
20	13.751	1952	1957	1963	rBV2	507693	808559	1.76%	0.491%
21	13.885	1973	1979	1981	rBV	566388	855402	1.86%	0.519%
22	13.916	1981	1984	1988	rVB	520771	705021	1.54%	0.428%
23	13.971	1988	1993	1997	rBV	1364323	1957143	4.26%	1.188%
24	14.081	2006	2011	2016	rBV	826078	1245055	2.71%	0.756%
25	14.294	2037	2046	2049	rBV	834226	1377525	3.00%	0.837%
26	14.331	2049	2052	2057	rVV	926588	1400795	3.05%	0.851%
27	14.562	2086	2090	2095	rVV	1213055	1882580	4.10%	1.143%
28	14.611	2095	2098	2103	rVB	344489	469651	1.02%	0.285%
29	14.684	2103	2110	2115	rBV2	1960334	3681101	8.02%	2.235%
30	14.745	2115	2120	2129	rVB4	421958	855907	1.86%	0.520%
31	14.946	2148	2153	2156	rVV3	524210	982038	2.14%	0.596%
32	15.007	2158	2163	2167	rVV2	1671628	2710831	5.90%	1.646%
33	15.056	2168	2171	2181	rVB2	454568	810058	1.76%	0.492%
34	15.233	2192	2200	2209	rVB2	27843995	45907918	100.00%	27.878%
35	15.336	2210	2217	2223	rBV4	791548	1716470	3.74%	1.042%
36	15.525	2242	2248	2254	rBV	658011	1135002	2.47%	0.689%

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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\82Y022822S.M

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37	15.690	2268	2275	2290	rVB3	366378	1183939	2.58%	0.719%
38	15.812	2290	2295	2300	rBV3	268652	538483	1.17%	0.327%
39	15.885	2302	2307	2313	rVB2	276212	574466	1.25%	0.349%
40	16.220	2354	2362	2366	rBV4	336827	900785	1.96%	0.547%
41	16.287	2366	2373	2387	rVV	17681194	33841241	73.72%	20.550%
42	16.409	2387	2393	2398	rVV2	233033	560376	1.22%	0.340%
43	16.489	2398	2406	2419	rVB	10036822	20582331	44.83%	12.499%

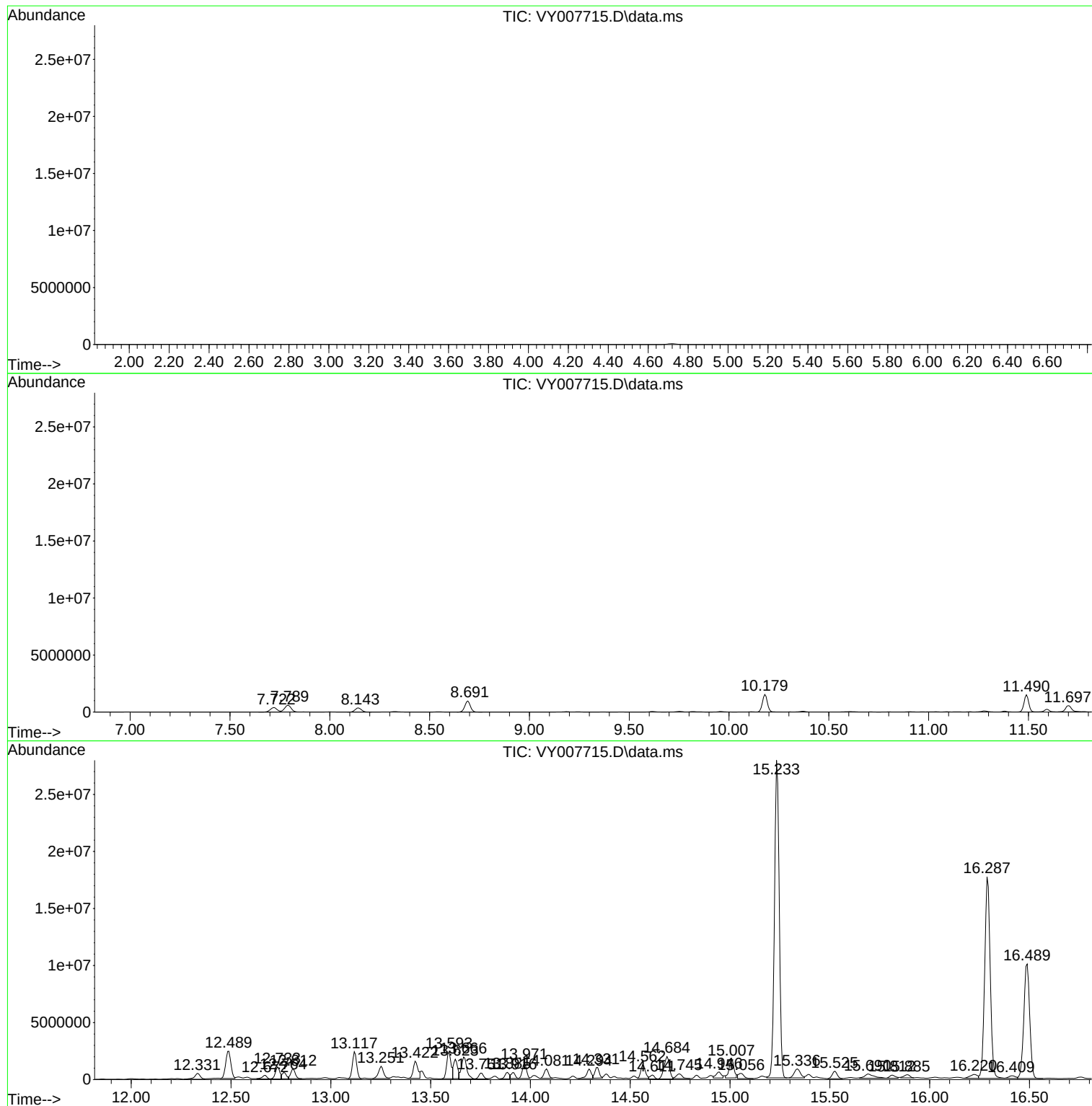
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y022822S.M
Quant Title : SW846 8260

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



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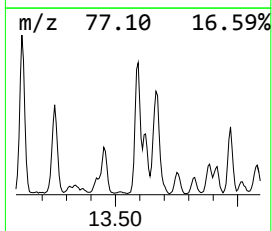
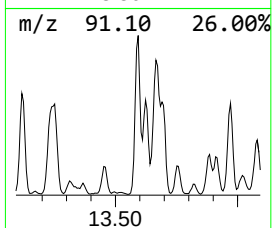
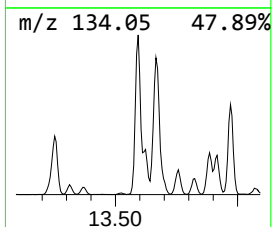
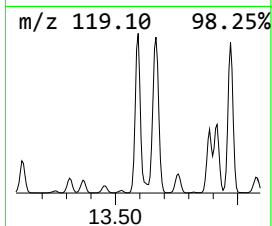
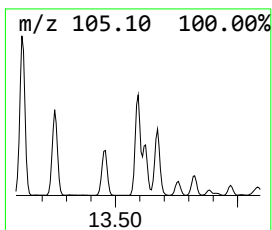
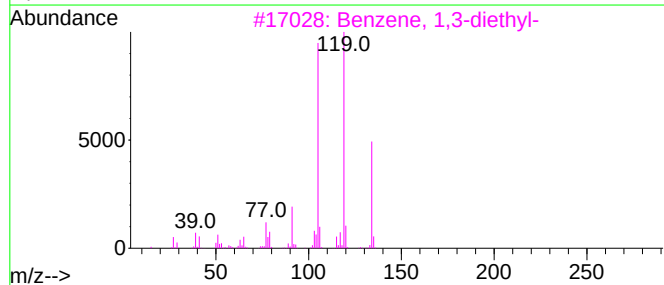
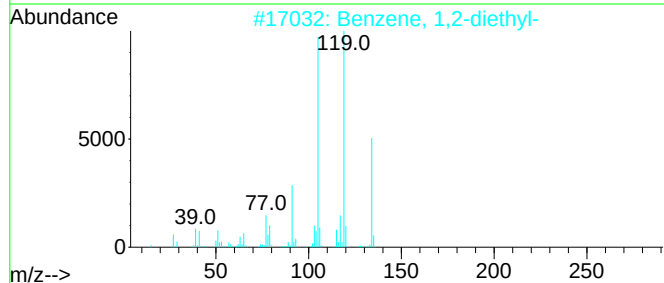
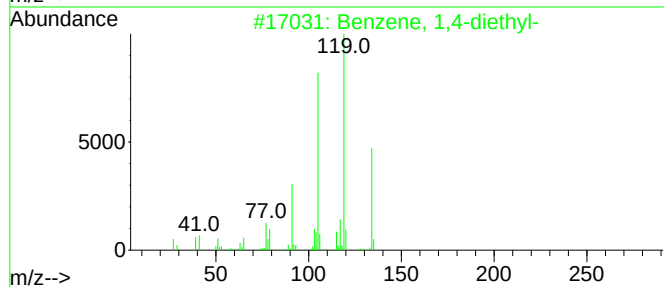
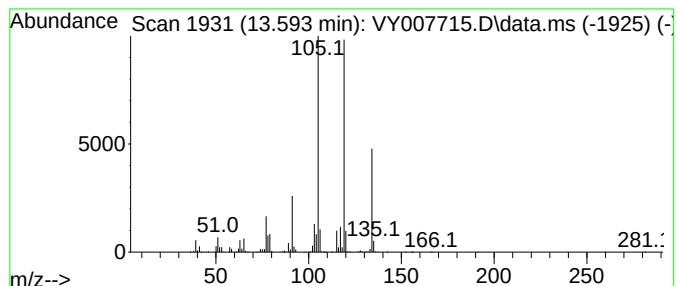
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y022822S.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1,4-diethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.593	74.35 ug/l	3699360	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	97
2			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
3			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	96
4			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91
5			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	87



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Instrument :
MSVOA_Y
ClientSampleId :
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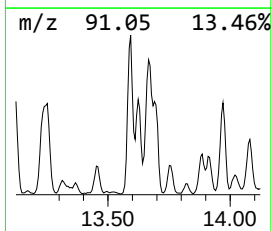
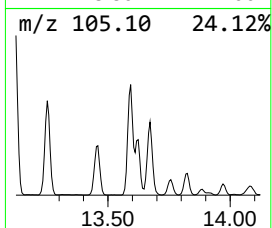
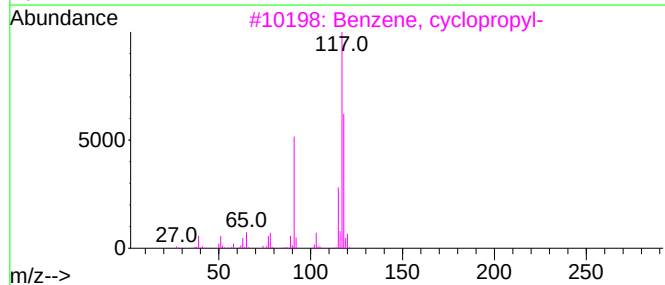
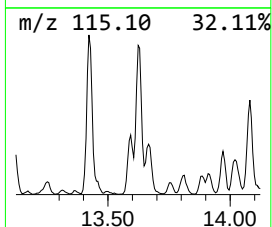
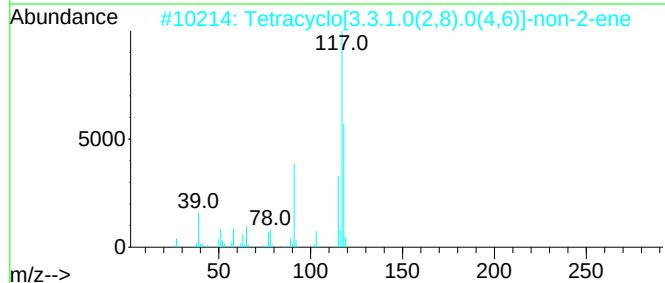
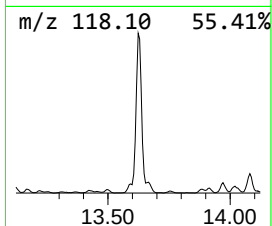
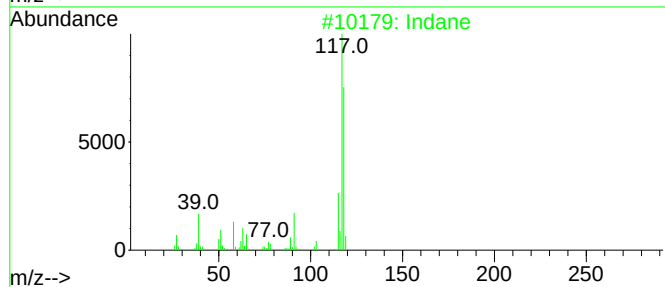
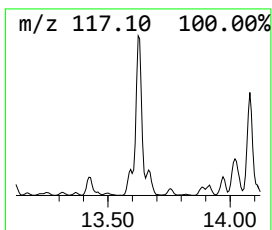
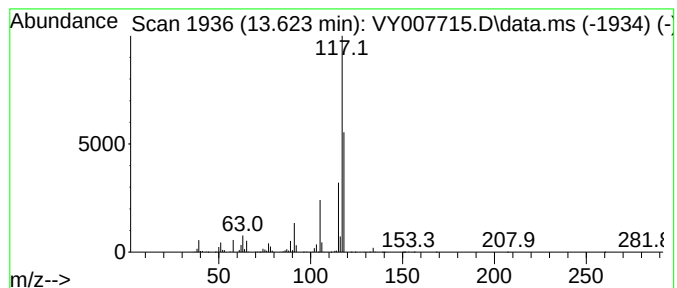
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y022822S.M
Quant Title : SW846 8260

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

Peak Number 2 Indane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.623	44.60 ug/l	2219300	1,4-Dichlorobenzene-d4	13.422

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Indane	118	C9H10	000496-11-7	68
2		Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	64
3		Benzene, cyclopropyl-	118	C9H10	000873-49-4	64
4		Benzene, 2-propenyl-	118	C9H10	000300-57-2	52
5		Benzene, 1-propenyl-	118	C9H10	000637-50-3	50



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ClientSampleId :
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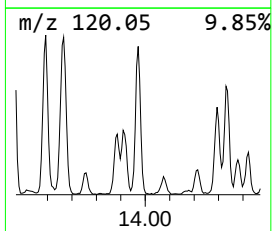
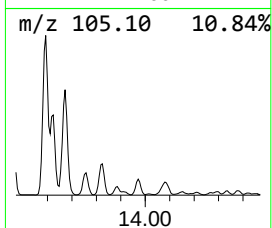
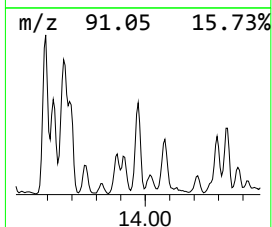
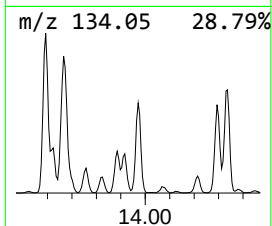
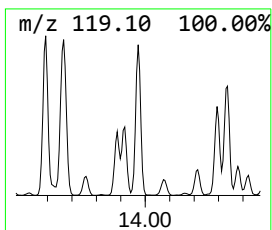
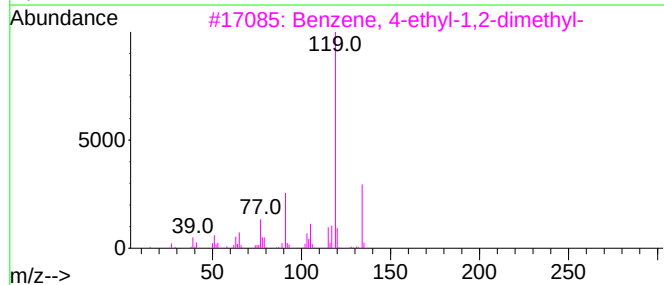
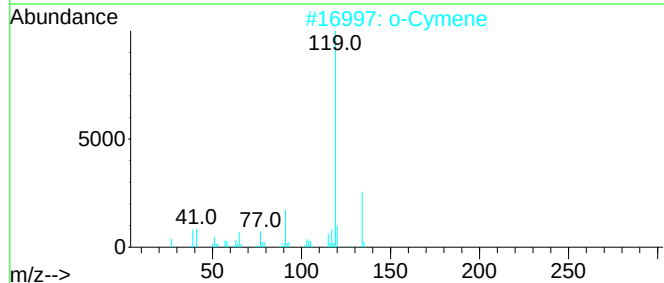
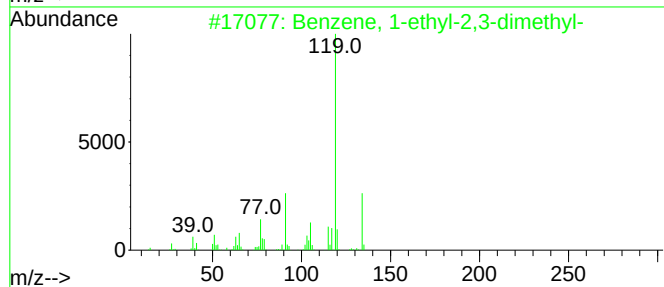
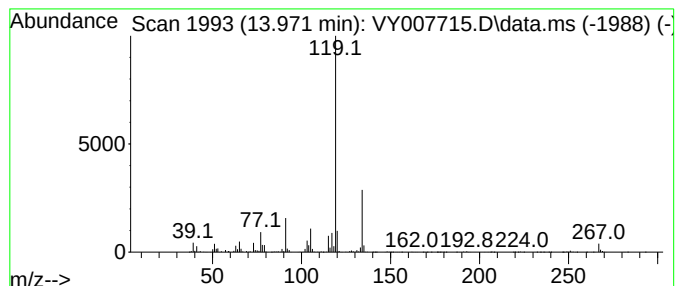
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.971	39.33 ug/l	1957140	1,4-Dichlorobenzene-d4	13.422

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



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ALS Vial : 26 Sample Multiplier: 1

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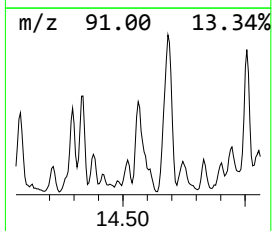
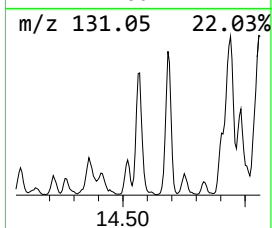
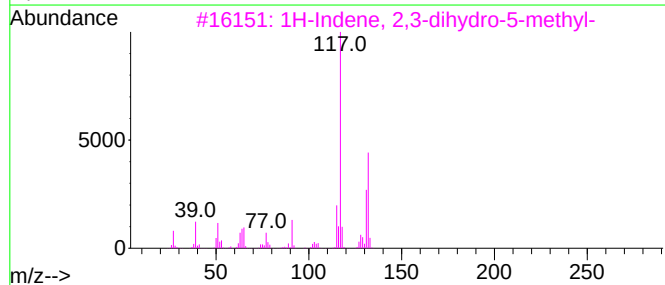
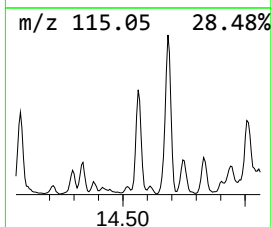
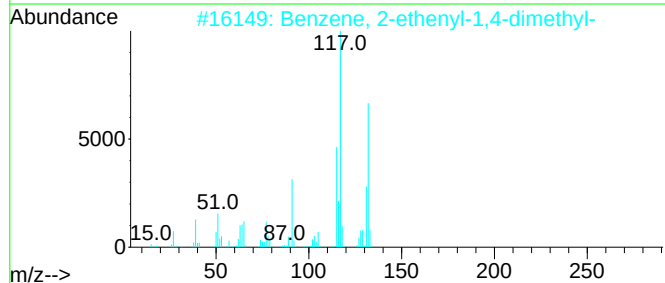
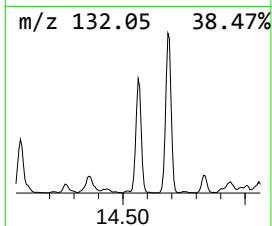
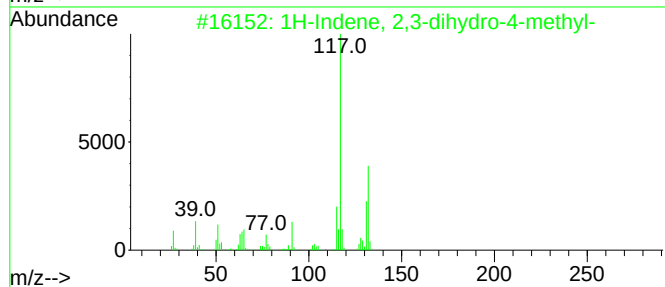
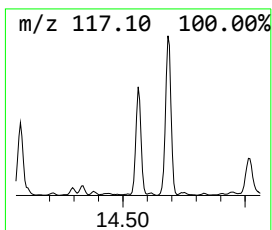
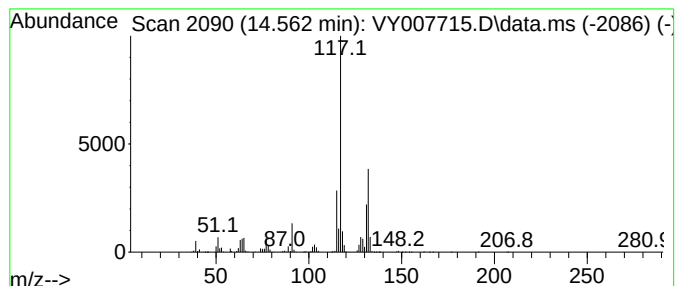
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.562	37.83 ug/l	1882580	1,4-Dichlorobenzene-d4	13.422

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	94
2		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	93
3		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
4		3-Phenylbut-1-ene	132	C10H12	000934-10-1	91
5		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87



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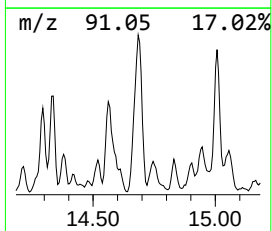
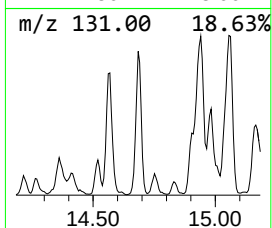
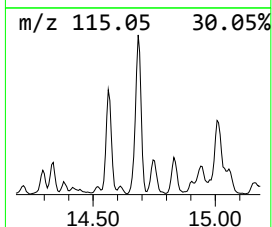
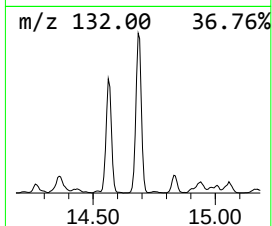
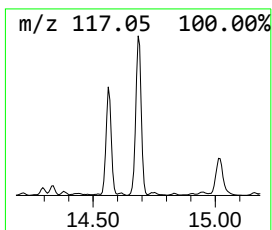
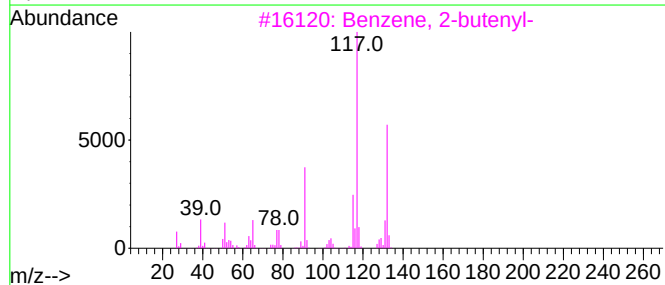
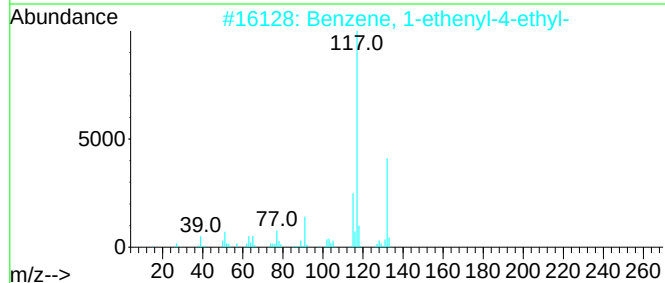
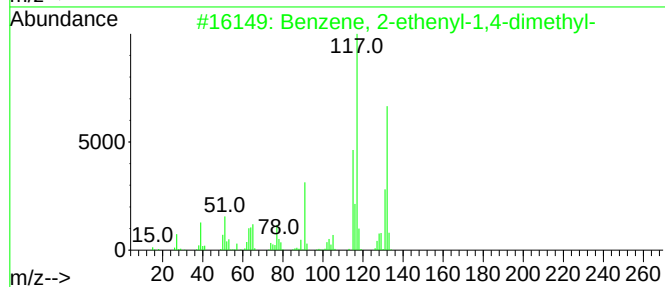
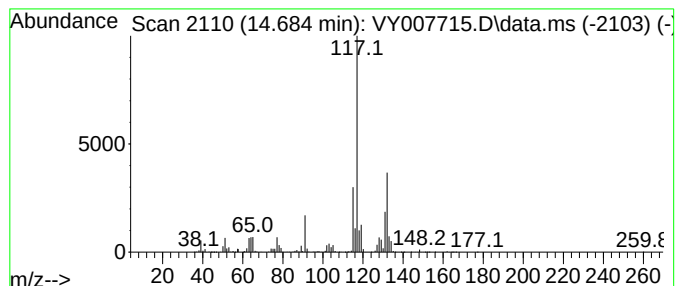
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.684	73.98 ug/l	3681100	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	96
2			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
3			Benzene, 2-butenyl-	132	C10H12	001560-06-1	87
4			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
5			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY030322\
Data File : VY007715.D
Acq On : 03 Mar 2022 19:26
Operator : SY/MD
Sample : N1762-02
Misc : 7.44g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
282

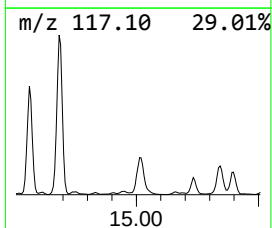
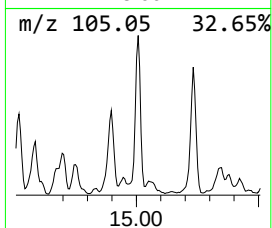
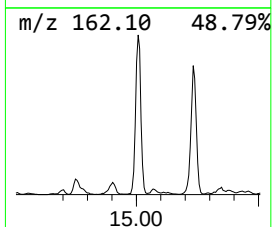
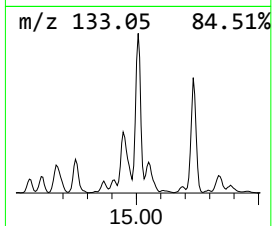
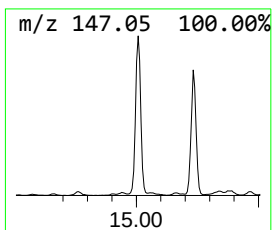
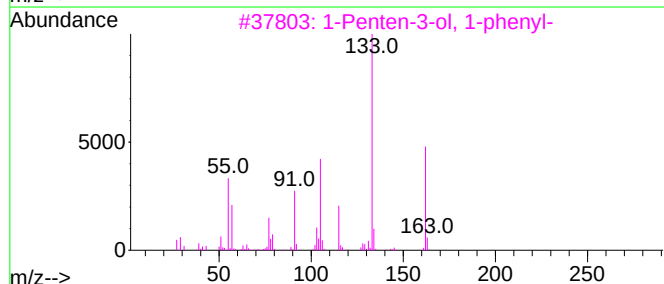
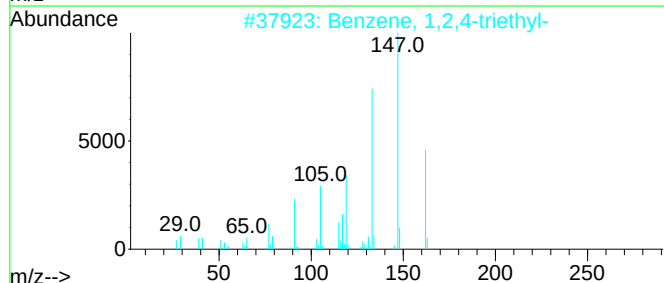
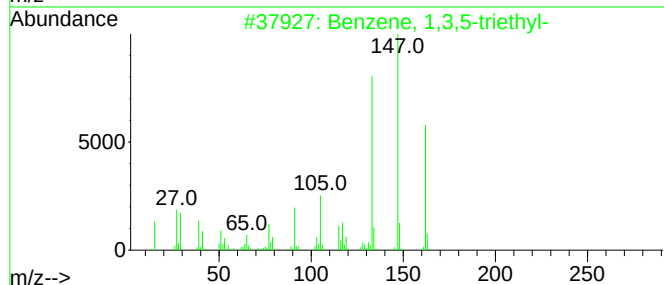
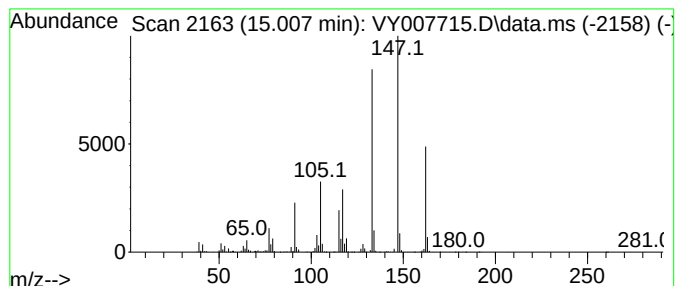
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1,3,5-triethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.007	54.48 ug/l	2710830	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3,5-triethyl-	162	C12H18	000102-25-0	93
2			Benzene, 1,2,4-triethyl-	162	C12H18	000877-44-1	93
3			1-Penten-3-ol, 1-phenyl-	162	C11H14O	034862-94-7	70
4			Benzene, 3-ethyl-1,2,4,5-tetrame...	162	C12H18	031365-98-7	53
5			4,5-Dimethyl-3H-isobenzofuran-1-one	162	C10H10O2	1000188-08-0	46



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY030322\
Data File : VY007715.D
Acq On : 03 Mar 2022 19:26
Operator : SY/MD
Sample : N1762-02
Misc : 7.44g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
282

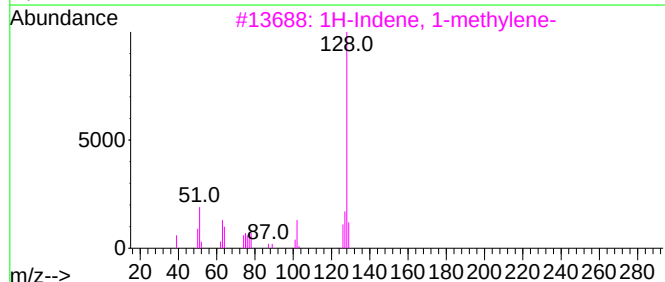
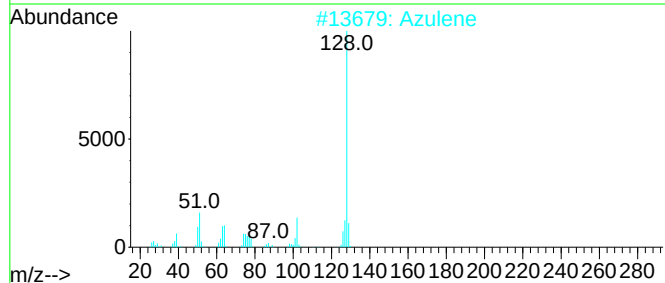
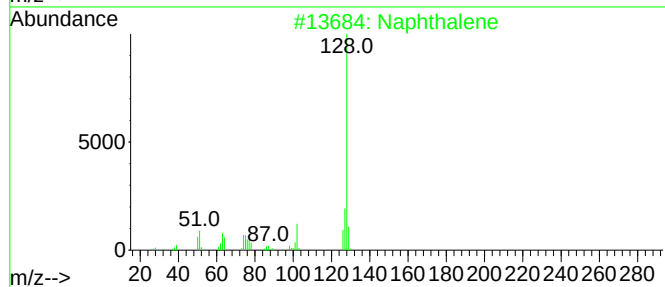
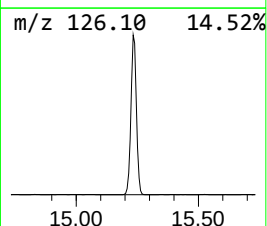
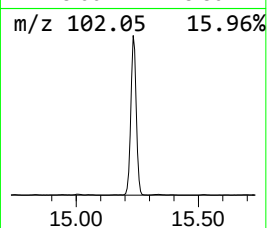
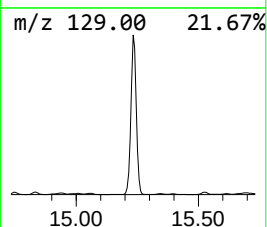
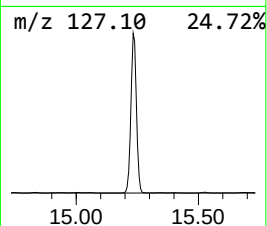
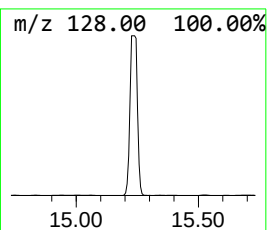
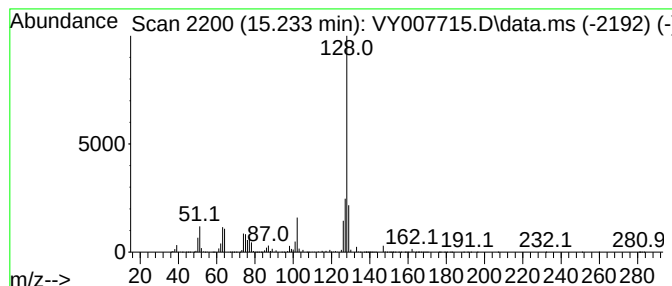
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Quant Title : SW846 8260

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

Peak Number 7 Azulene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.233	922.60 ug/l	45907900	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene	128	C10H8	000091-20-3	93
2			Azulene	128	C10H8	000275-51-4	87
3			1H-Indene, 1-methylene-	128	C10H8	002471-84-3	83
4			4-Phenylbut-3-ene-1-yne	128	C10H8	1000222-12-0	74
5			[4.2.2]Propella-2,4,7,9-tetraene	128	C10H8	088090-34-0	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY030322\
Data File : VY007715.D
Acq On : 03 Mar 2022 19:26
Operator : SY/MD
Sample : N1762-02
Misc : 7.44g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
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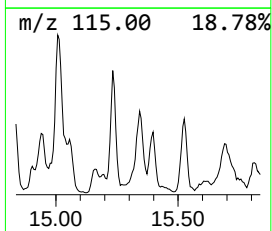
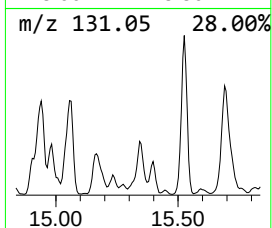
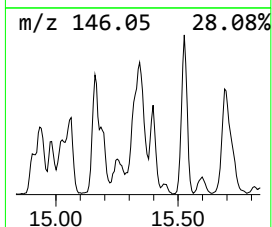
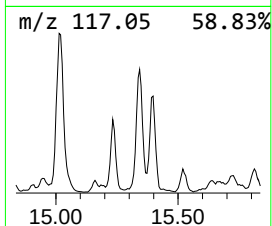
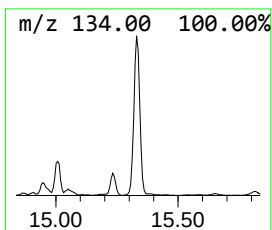
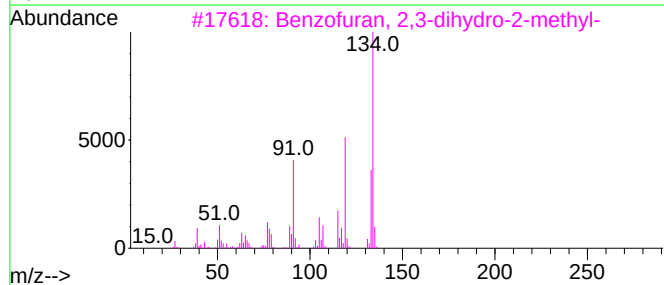
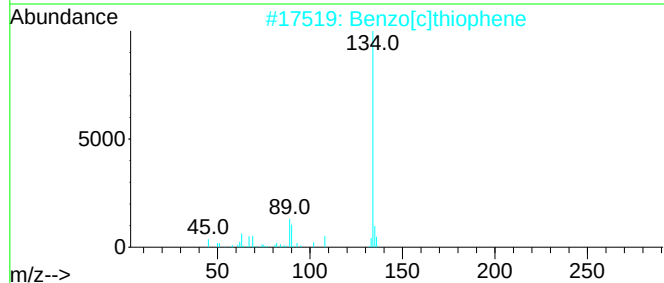
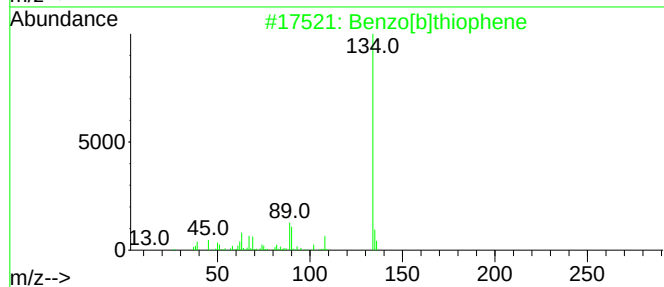
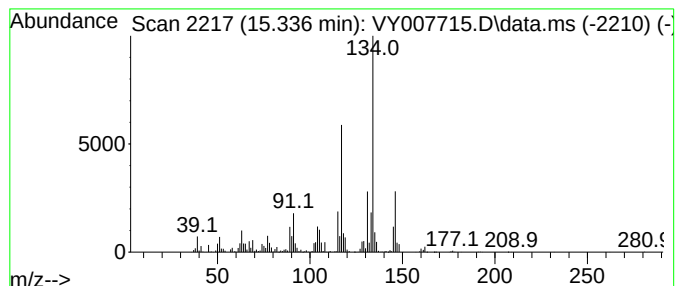
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Quant Title : SW846 8260

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

Peak Number 8 unknown15.336 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.336	34.50 ug/l	1716470	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[b]thiophene	134	C8H6S	000095-15-8	46
2			Benzo[c]thiophene	134	C8H6S	000270-82-6	46
3			Benzofuran, 2,3-dihydro-2-methyl-	134	C9H10O	001746-11-8	43
4			1H-Imidazole, 4,5-dihydro-2-phenyl-	146	C9H10N2	000936-49-2	38
5			Bicyclo[4.2.1]nona-2,4,7-triene,...	146	C11H14	1000164-42-6	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY030322\
Data File : VY007715.D
Acq On : 03 Mar 2022 19:26
Operator : SY/MD
Sample : N1762-02
Misc : 7.44g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
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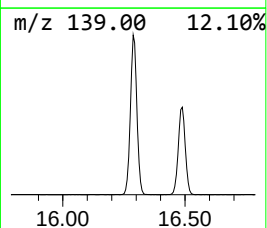
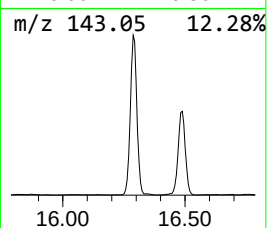
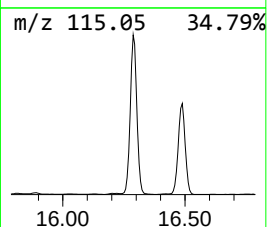
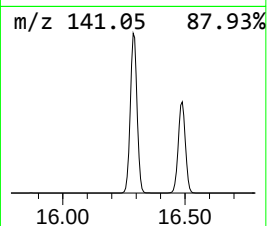
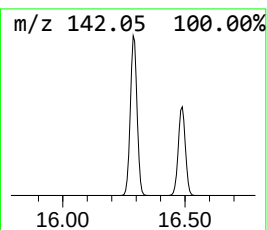
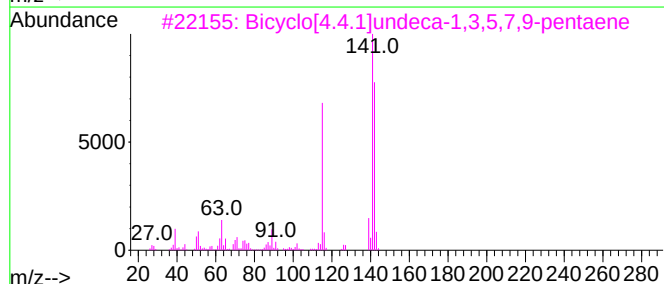
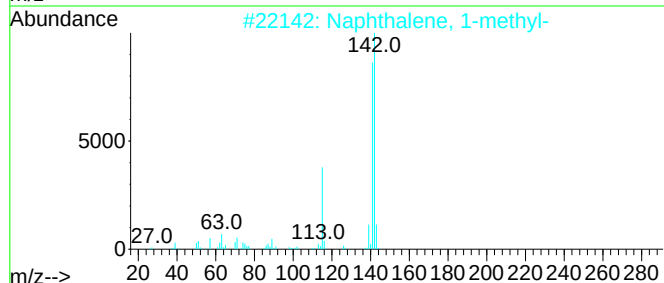
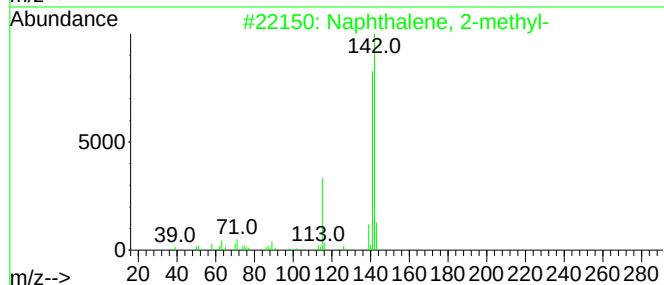
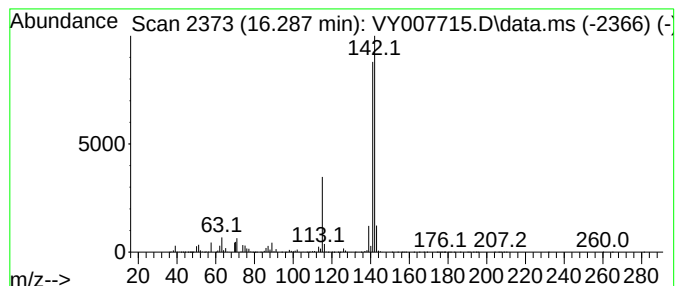
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Quant Title : SW846 8260

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

Peak Number 9 Bicyclo[4.4.1]undeca-1,3,5,... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.288	680.10 ug/l	33841200	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			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	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY030322\
Data File : VY007715.D
Acq On : 03 Mar 2022 19:26
Operator : SY/MD
Sample : N1762-02
Misc : 7.44g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
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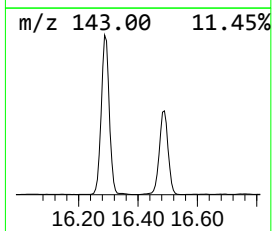
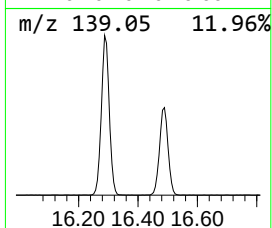
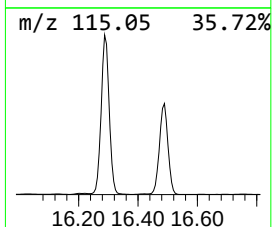
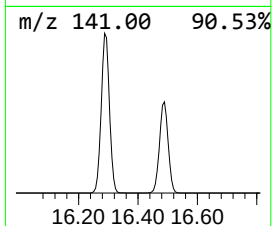
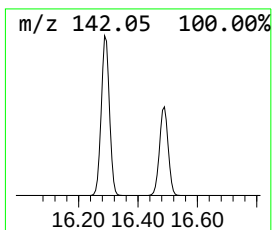
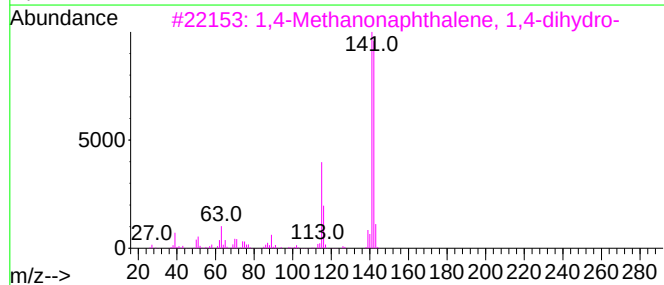
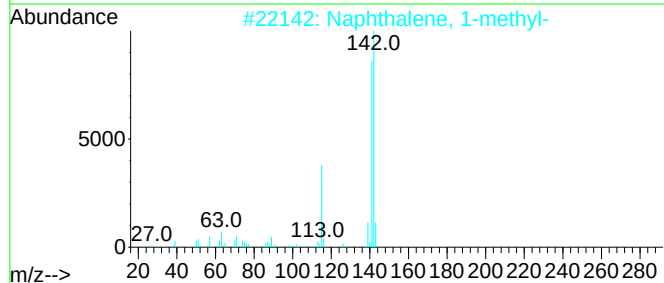
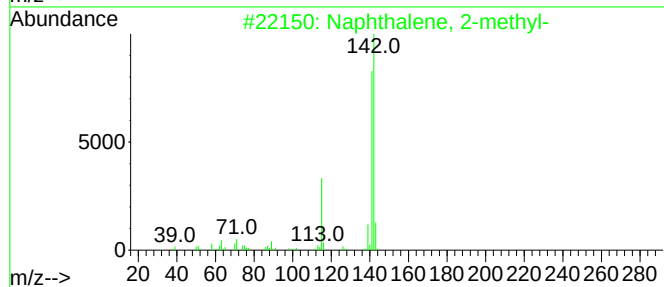
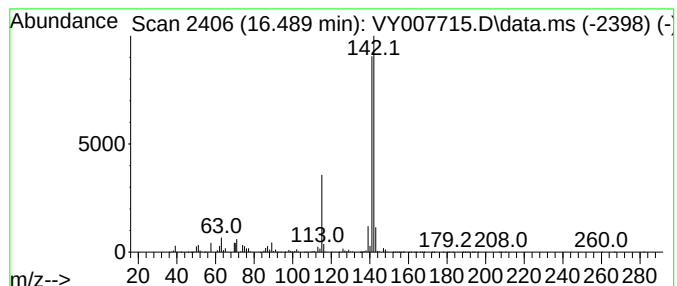
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y022822S.M
Quant Title : SW846 8260

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

Peak Number 10 1,4-Methanonaphthalene, 1,4... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.489	413.64 ug/l	20582300	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	94
4			Benzocycloheptatriene	142	C11H10	000264-09-5	91
5			Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY030322\
Data File : VY007715.D
Acq On : 03 Mar 2022 19:26
Operator : SY/MD
Sample : N1762-02
Misc : 7.44g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y022822S.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
Benzene, 1,4-di...	13.593	74.3	ug/l	3699360	4	13.422	2487960	50.0
Indane	13.623	44.6	ug/l	2219300	4	13.422	2487960	50.0
Benzene, 1-ethy...	13.971	39.3	ug/l	1957140	4	13.422	2487960	50.0
1H-Indene, 2,3-...	14.562	37.8	ug/l	1882580	4	13.422	2487960	50.0
Benzene, 2-ethe...	14.684	74.0	ug/l	3681100	4	13.422	2487960	50.0
Benzene, 1,3,5-...	15.007	54.5	ug/l	2710830	4	13.422	2487960	50.0
Azulene	15.233	922.6	ug/l	45907900	4	13.422	2487960	50.0
unknown15.336	15.336	34.5	ug/l	1716470	4	13.422	2487960	50.0
Bicyclo[4.4.1]u...	16.288	680.1	ug/l	33841200	4	13.422	2487960	50.0
1,4-Methanonaph...	16.489	413.6	ug/l	20582300	4	13.422	2487960	50.0