

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD032522\
 Data File : VD072428.D
 Acq On : 25 Mar 2022 14:47
 Operator : VA/SY
 Sample : N1977-04
 Misc : 4.31G/5.0ml/MSVOA_D/SOIL
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 SB-12-20220315-16.0-17.0-A

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_D\Method\82D031122S.M
 Title : SW846 8260

Signal : TIC: VD072428.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	8.344	1080	1092	1102	rBV3	232764	604159	1.41%	0.452%
2	11.638	1645	1652	1662	rBV2	253243	525425	1.23%	0.393%
3	11.738	1662	1669	1680	rVV	506435	837492	1.96%	0.627%
4	12.473	1787	1794	1800	rBV	234598	436367	1.02%	0.327%
5	12.626	1814	1820	1830	rVB	239667	438652	1.03%	0.328%
6	12.949	1870	1875	1888	rVB	261086	533909	1.25%	0.399%
7	13.114	1895	1903	1910	rBV	384739	703414	1.65%	0.526%
8	13.255	1919	1927	1933	rBV	964108	1614521	3.78%	1.208%
9	13.567	1975	1980	1981	rVV	325707	479215	1.12%	0.359%
10	13.590	1981	1984	1990	rVB	624188	974887	2.28%	0.729%
11	13.726	2000	2007	2009	rBV	927462	1432786	3.35%	1.072%
12	13.767	2009	2014	2028	rVV2	5117283	9895083	23.17%	7.404%
13	14.020	2052	2057	2059	rVV	419599	680100	1.59%	0.509%
14	14.043	2059	2061	2066	rVB	399008	532464	1.25%	0.398%
15	14.102	2066	2071	2077	rBV	1623093	2472142	5.79%	1.850%
16	14.214	2082	2090	2095	rBV2	1158820	2214371	5.18%	1.657%
17	14.349	2106	2113	2117	rBV3	354522	730945	1.71%	0.547%
18	14.426	2123	2126	2130	rBV	418556	645236	1.51%	0.483%
19	14.467	2130	2133	2137	rVB	818976	1046296	2.45%	0.783%
20	14.508	2137	2140	2145	rVB	779048	1065227	2.49%	0.797%
21	14.579	2145	2152	2156	rBV3	321613	687037	1.61%	0.514%
22	14.696	2167	2172	2178	rBV	2001594	3291941	7.71%	2.463%
23	14.814	2184	2192	2197	rBV3	1972586	4700158	11.01%	3.517%
24	14.885	2197	2204	2212	rVB4	762350	1852501	4.34%	1.386%
25	14.967	2212	2218	2229	rBV3	1824106	3405700	7.97%	2.548%
26	15.079	2230	2237	2243	rBV3	1020980	2138190	5.01%	1.600%
27	15.190	2251	2256	2265	rBV2	632436	1431349	3.35%	1.071%
28	15.385	2279	2289	2301	rVB	21630518	42708816	100.00%	31.956%
29	15.490	2301	2307	2320	rVB3	968296	2216169	5.19%	1.658%
30	15.685	2332	2340	2343	rBV4	463352	1162363	2.72%	0.870%
31	15.879	2366	2373	2378	rBV2	702664	1492226	3.49%	1.117%
32	16.002	2386	2394	2399	rBV2	650838	1728483	4.05%	1.293%
33	16.196	2421	2427	2432	rBV9	196201	504233	1.18%	0.377%
34	16.273	2434	2440	2445	rVV5	341878	950828	2.23%	0.711%
35	16.326	2446	2449	2456	rVB2	384388	659211	1.54%	0.493%
36	16.414	2457	2464	2469	rVV	540317	1247182	2.92%	0.933%

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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_D\Method\82D031122S.M
Title : SW846 8260

37	16.479	2469	2475	2488	rVB2	1672075	4163939	9.75%	3.116%
38	16.602	2491	2496	2501	rVV3	241905	606420	1.42%	0.454%
39	16.684	2501	2510	2521	rVB	13247218	30839302	72.21%	23.075%

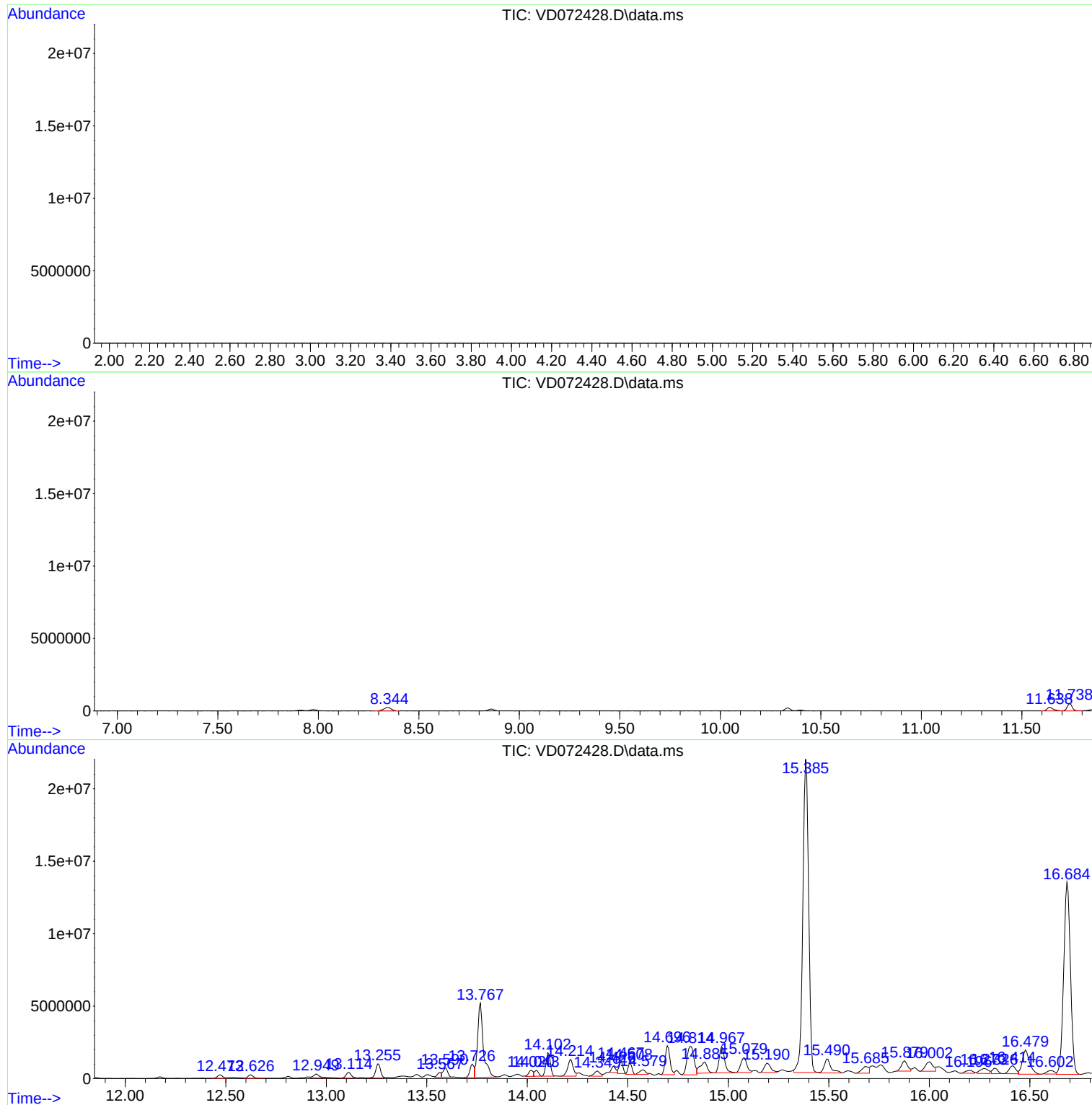
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D031122S.M
Quant Title : SW846 8260

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



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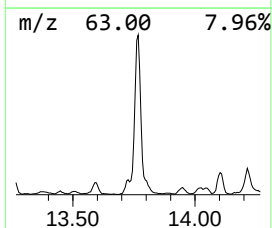
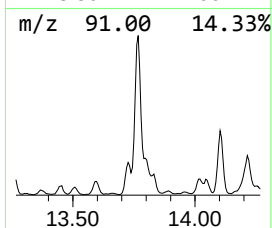
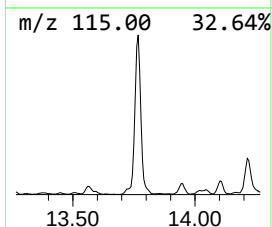
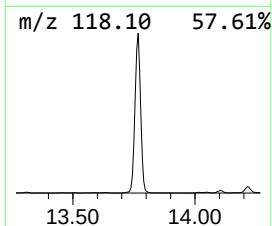
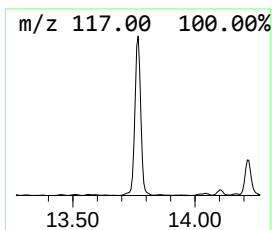
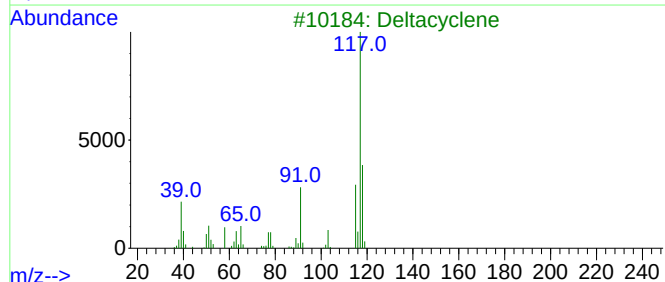
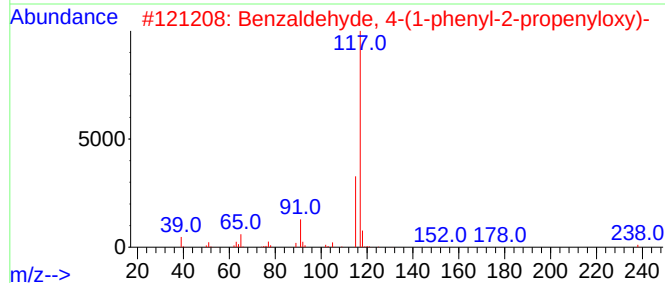
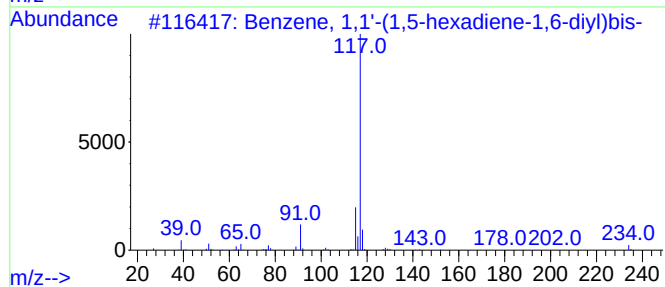
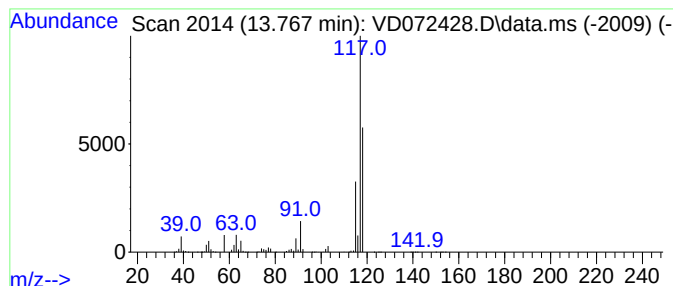
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D031122S.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1,1'-(1,5-hexadien... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.767	1032.43 ug/l	9895080	1,4-Dichlorobenzene-d4	13.567

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,1'-(1,5-hexadiene-1,6...	234	C18H18	004439-45-6	59
2			Benzaldehyde, 4-(1-phenyl-2-prop...	238	C16H14O2	1000277-56-1	53
3			Deltacyclene	118	C9H10	007785-10-6	50
4			Indole	117	C8H7N	000120-72-9	46
5			Benzene, (2-bromocyclopropyl)-	196	C9H9Br	036617-02-4	43



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ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
SB-12-20220315-16.0-17.0-A

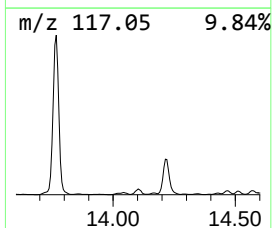
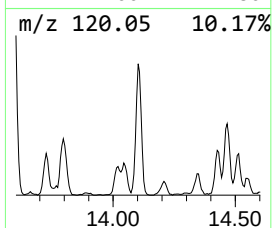
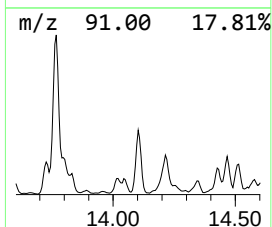
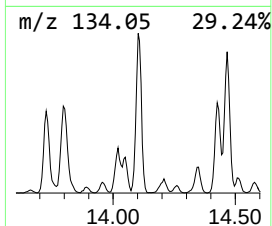
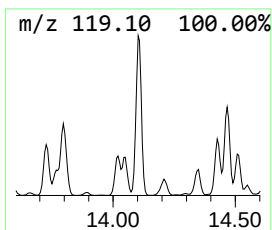
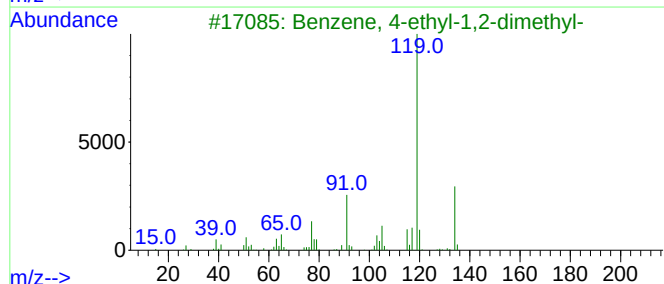
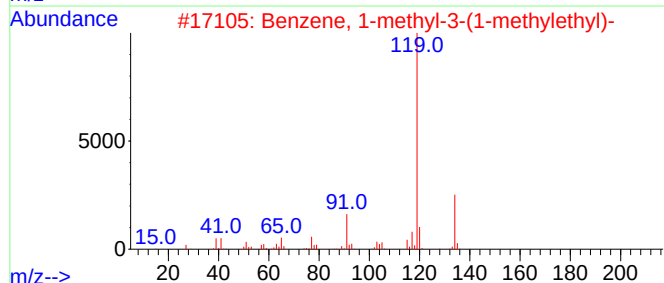
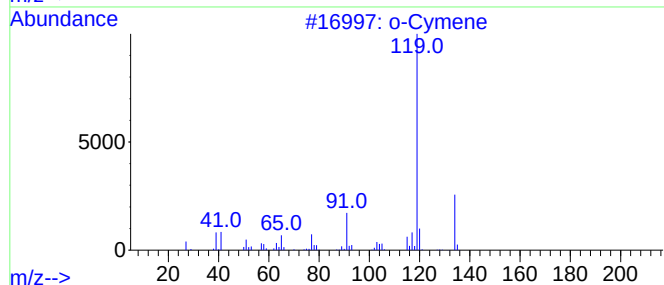
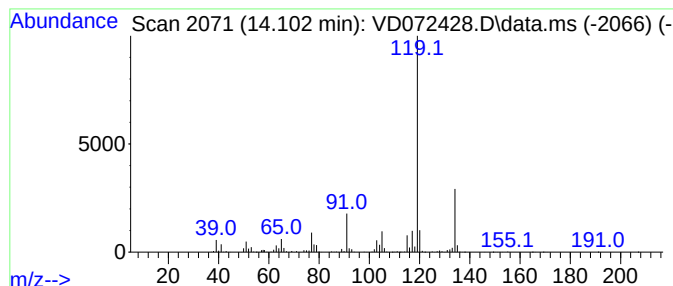
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D031122S.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 o-Cymene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.102	257.94 ug/l	2472140	1,4-Dichlorobenzene-d4	13.567

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Cymene	134	C10H14	000527-84-4	95
2			Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	95
3			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
4			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94
5			p-Cymene	134	C10H14	000099-87-6	94



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ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
SB-12-20220315-16.0-17.0-A

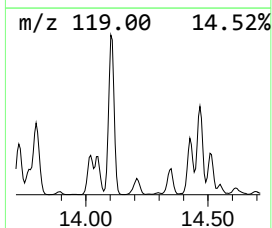
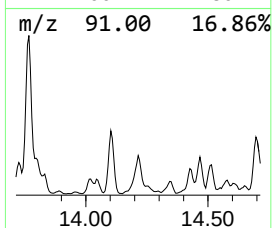
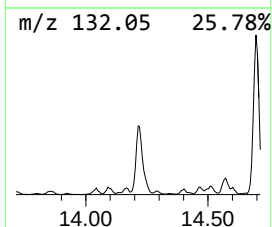
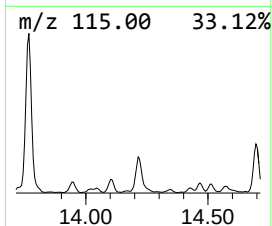
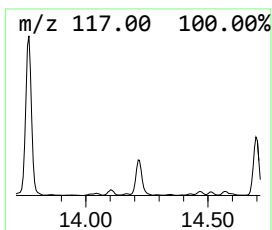
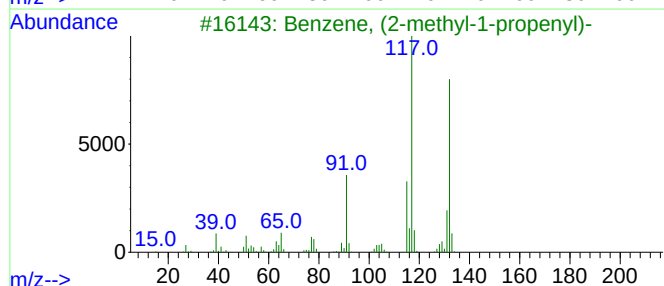
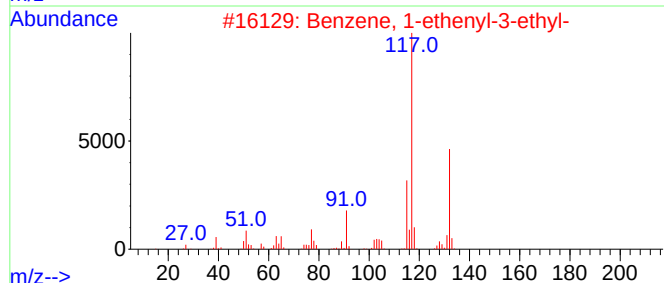
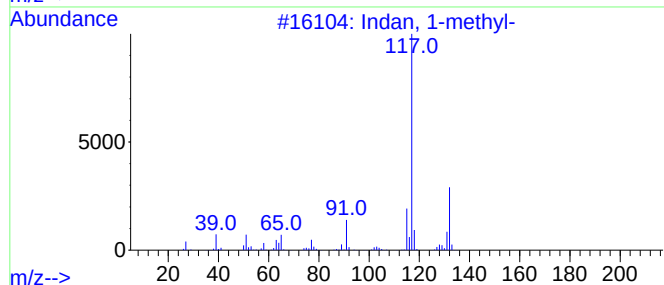
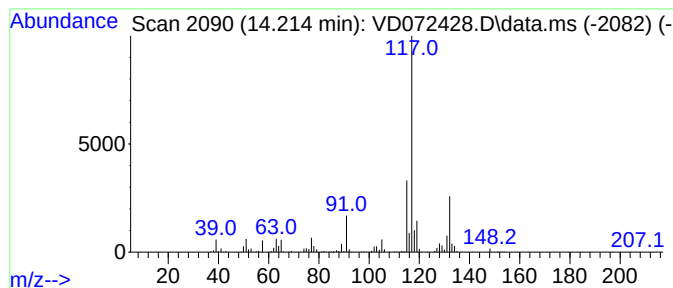
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D031122S.M
Quant Title : SW846 8260

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

Peak Number 3 Indan, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.214	231.04 ug/l	2214370	1,4-Dichlorobenzene-d4	13.567

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indan, 1-methyl-	132	C10H12	000767-58-8	93
2			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	90
3			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	87
4			Benzene, 2-butenyl-	132	C10H12	001560-06-1	83
5			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	83



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ClientSampleId :
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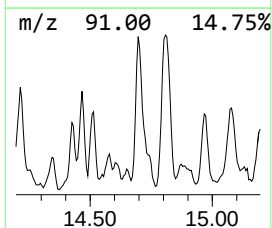
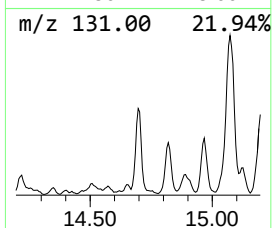
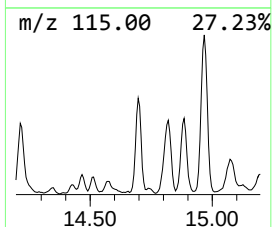
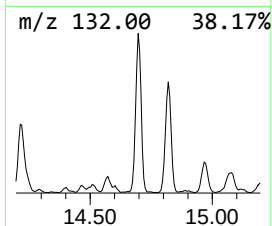
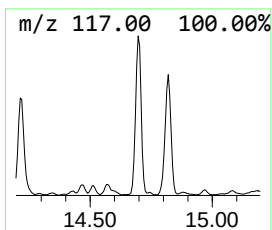
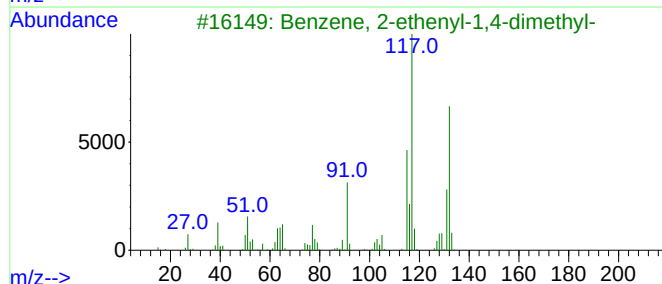
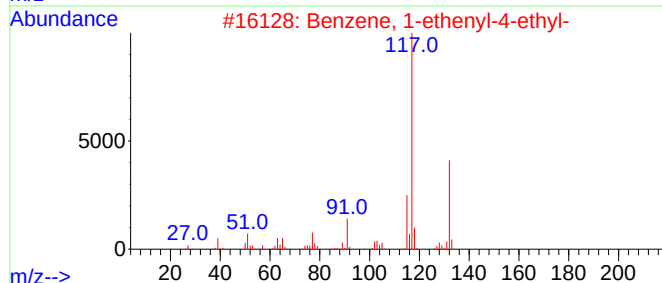
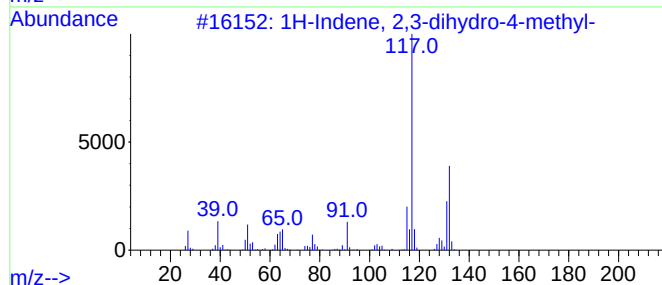
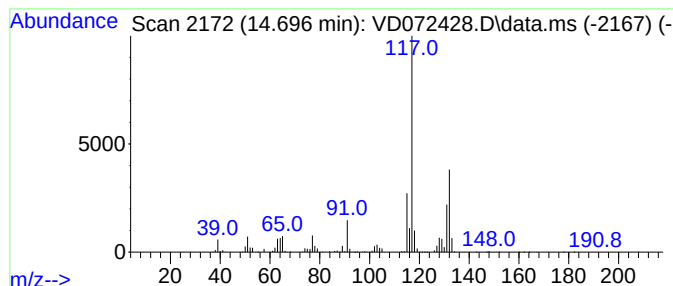
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D031122S.M
Quant Title : SW846 8260

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

Peak Number 4 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.696	343.47 ug/l	3291940	1,4-Dichlorobenzene-d4	13.567

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	95
2			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
3			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	93
4			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
5			3-Phenylbut-1-ene	132	C10H12	000934-10-1	91



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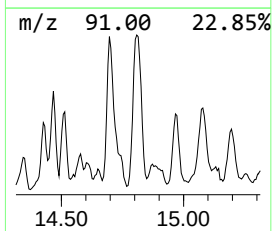
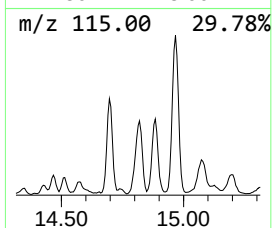
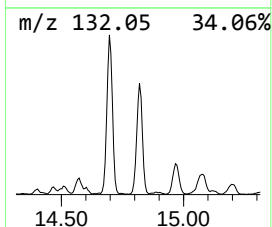
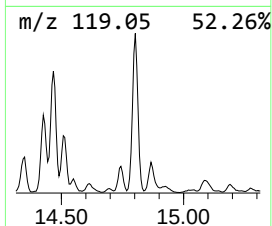
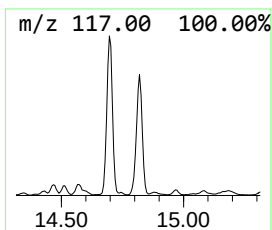
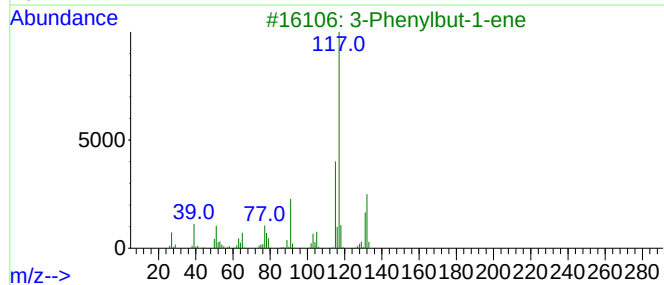
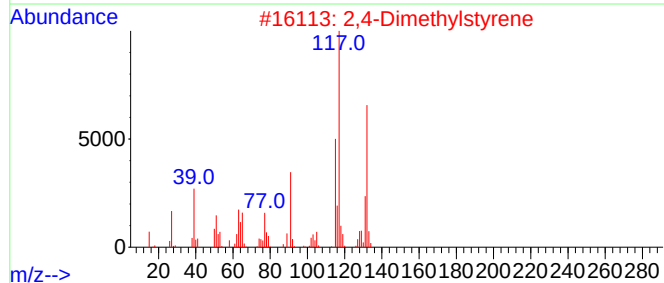
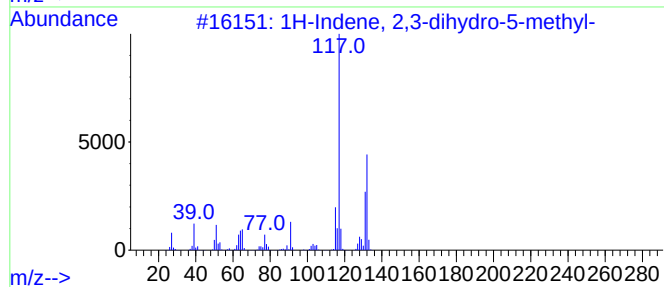
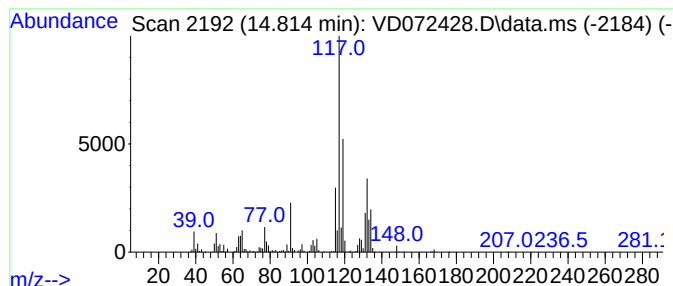
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D031122S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.814	490.40 ug/l	4700160	1,4-Dichlorobenzene-d4	13.567

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	91
2			2,4-Dimethylstyrene	132	C10H12	002234-20-0	89
3			3-Phenylbut-1-ene	132	C10H12	000934-10-1	70
4			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	70
5			1-Phenyl-1-butene	132	C10H12	000824-90-8	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD032522\
Data File : VD072428.D
Acq On : 25 Mar 2022 14:47
Operator : VA/SY
Sample : N1977-04
Misc : 4.31G/5.0ml/MSVOA_D/SOIL
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
SB-12-20220315-16.0-17.0-A

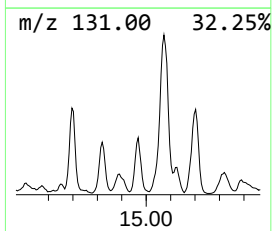
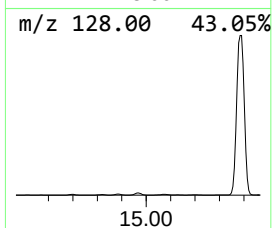
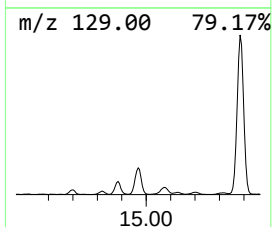
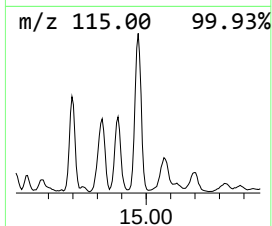
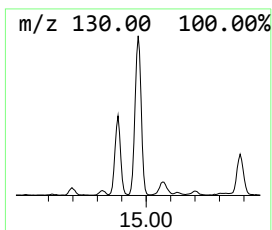
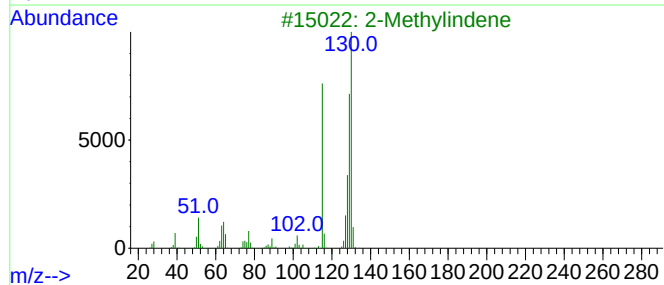
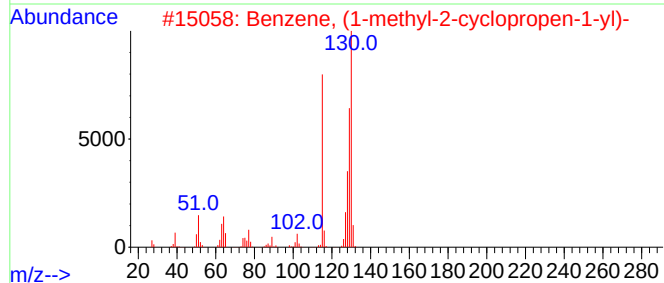
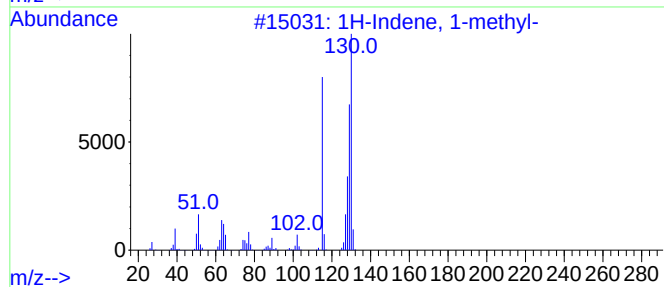
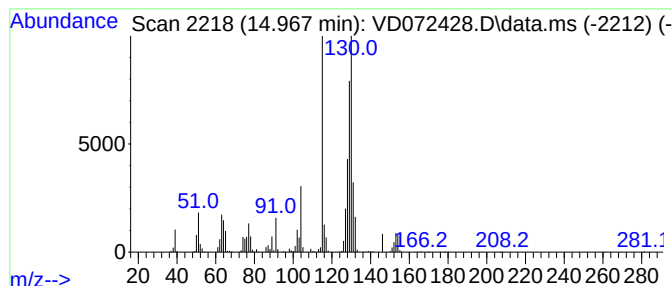
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Quant Title : SW846 8260

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

Peak Number 6 1H-Indene, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.967	355.34 ug/l	3405700	1,4-Dichlorobenzene-d4	13.567

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Indene, 1-methyl-	130	C10H10	000767-59-9	96
2		Benzene, (1-methyl-2-cyclopropen-1-yl)-	130	C10H10	065051-83-4	94
3		2-Methylindene	130	C10H10	002177-47-1	94
4		Naphthalene, 1,2-dihydro-	130	C10H10	000447-53-0	76
5		1H-Indene, 3-methyl-	130	C10H10	000767-60-2	76



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD032522\
Data File : VD072428.D
Acq On : 25 Mar 2022 14:47
Operator : VA/SY
Sample : N1977-04
Misc : 4.31G/5.0ml/MSVOA_D/SOIL
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
SB-12-20220315-16.0-17.0-A

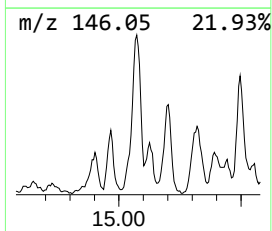
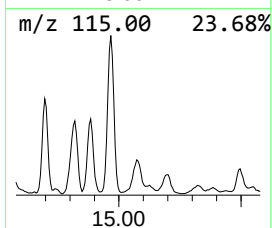
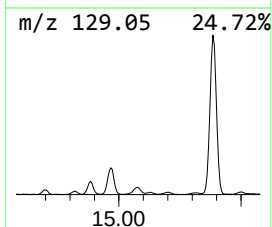
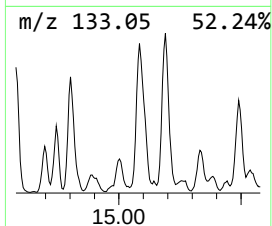
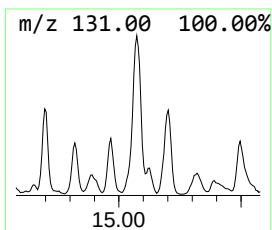
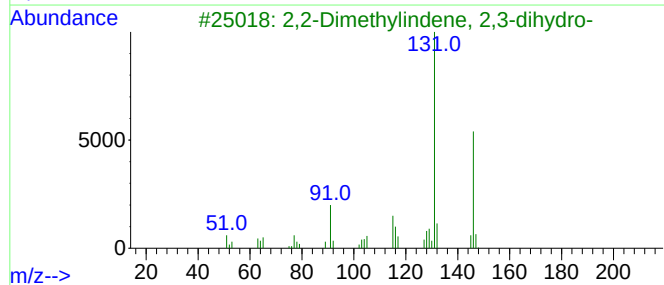
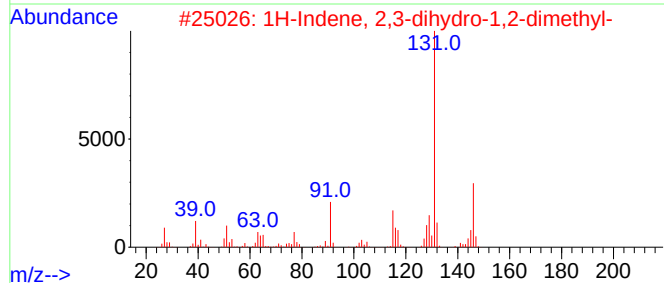
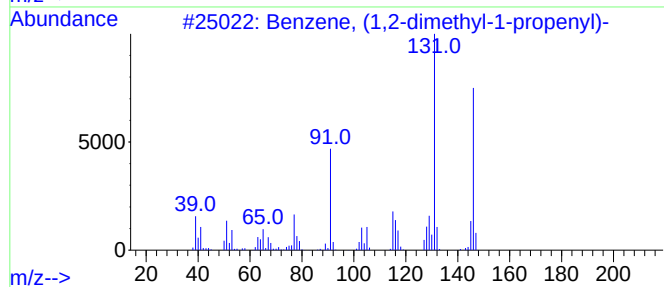
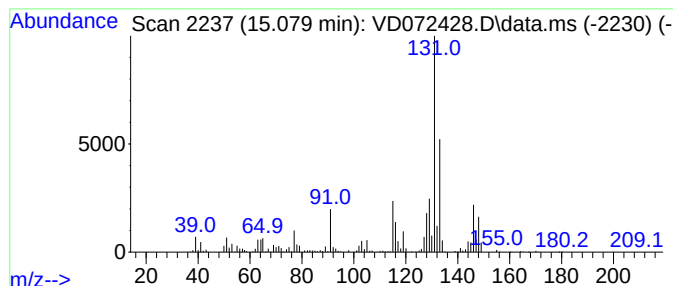
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, (1,2-dimethyl-1-propenyl)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.079	223.09 ug/l	2138190	1,4-Dichlorobenzene-d4	13.567

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	70
2			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	70
3			2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	70
4			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	70
5			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	70



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD032522\
Data File : VD072428.D
Acq On : 25 Mar 2022 14:47
Operator : VA/SY
Sample : N1977-04
Misc : 4.31G/5.0ml/MSVOA_D/SOIL
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
SB-12-20220315-16.0-17.0-A

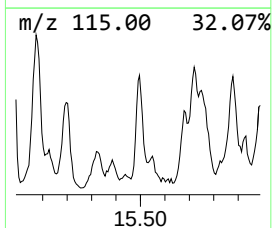
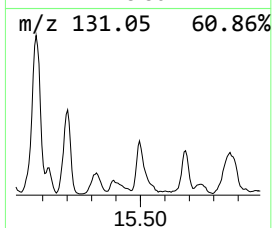
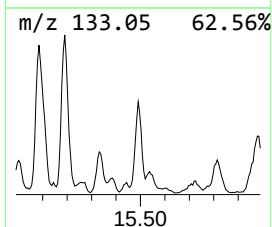
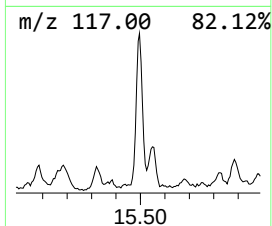
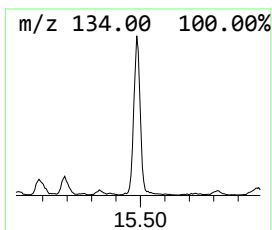
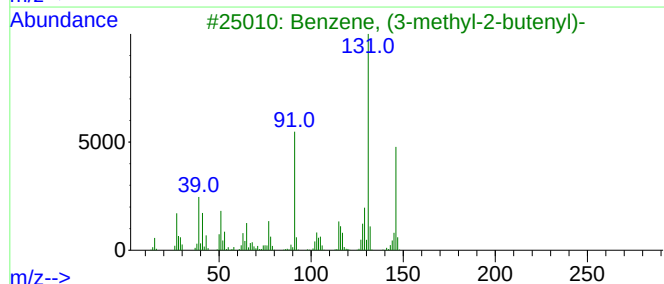
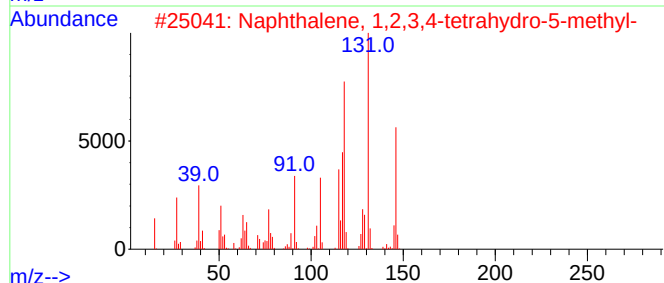
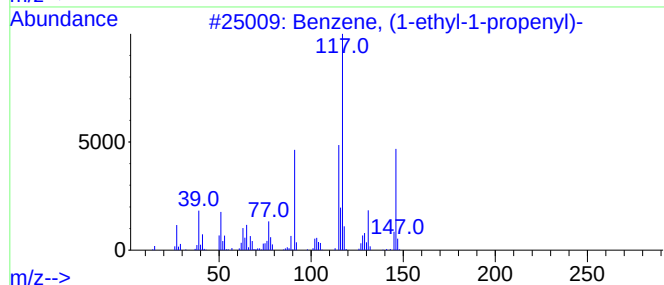
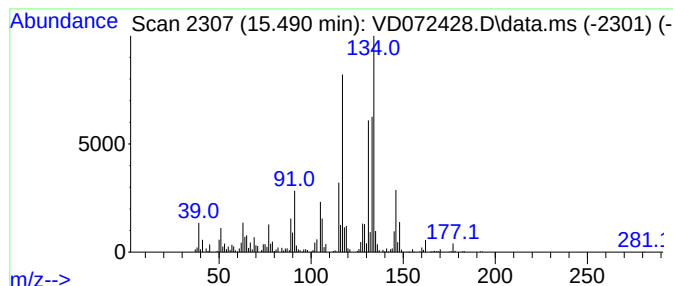
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, (1-ethyl-1-propenyl)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.490	231.23 ug/l	2216170	1,4-Dichlorobenzene-d4	13.567

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-ethyl-1-propenyl)-	146	C11H14	004701-36-4	83
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	52
3			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	47
4			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	47
5			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD032522\
Data File : VD072428.D
Acq On : 25 Mar 2022 14:47
Operator : VA/SY
Sample : N1977-04
Misc : 4.31G/5.0ml/MSVOA_D/SOIL
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
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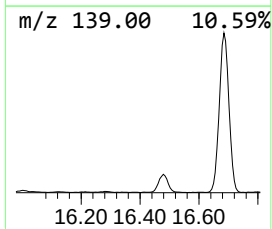
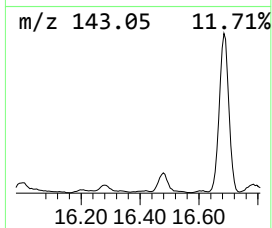
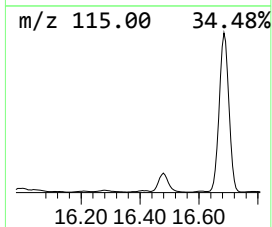
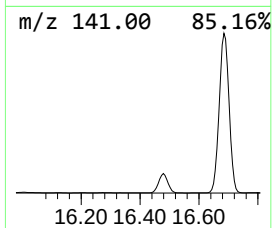
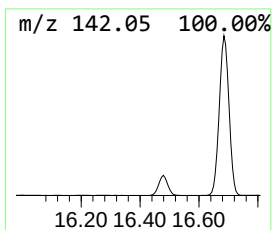
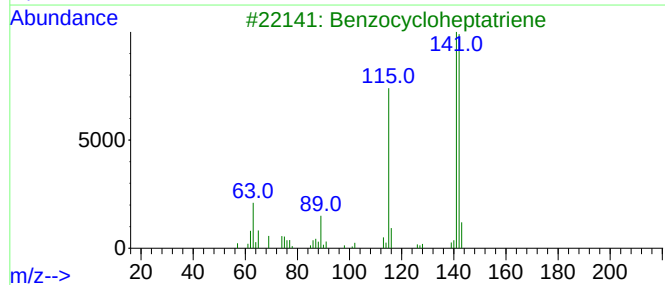
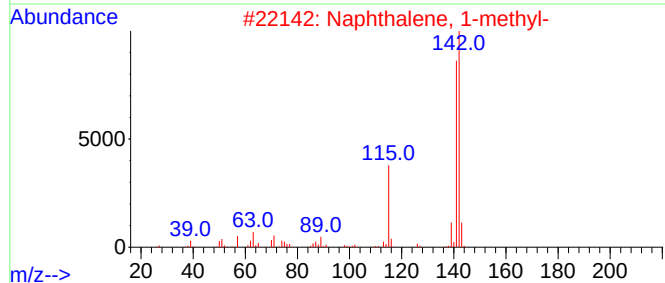
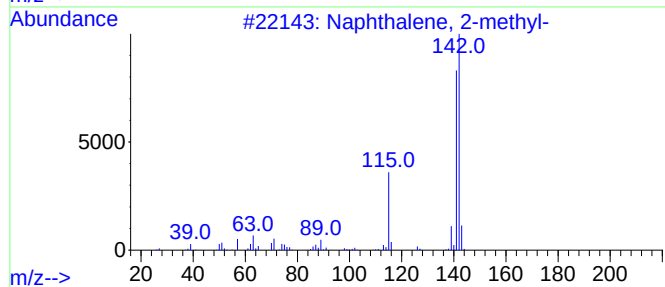
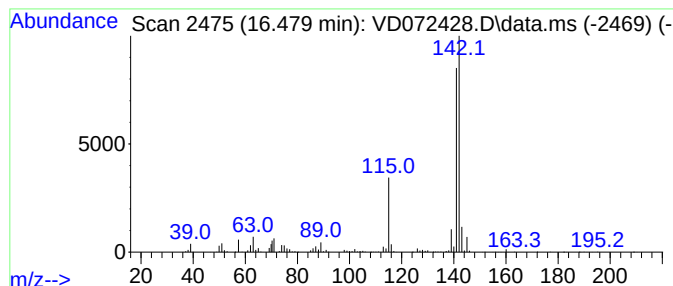
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Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.479	434.45 ug/l	4163940	1,4-Dichlorobenzene-d4	13.567

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	90
4			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	90
5			1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	46



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD032522\
Data File : VD072428.D
Acq On : 25 Mar 2022 14:47
Operator : VA/SY
Sample : N1977-04
Misc : 4.31G/5.0ml/MSVOA_D/SOIL
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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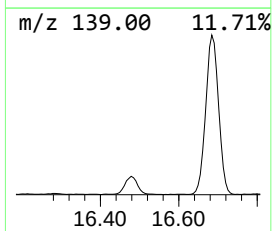
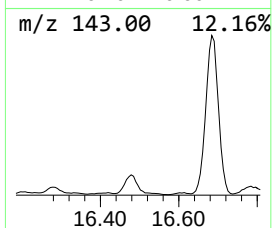
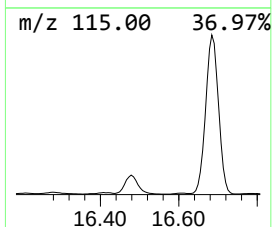
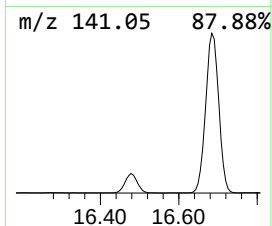
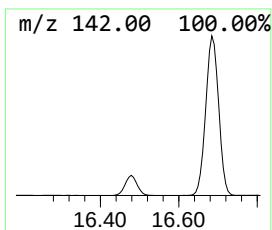
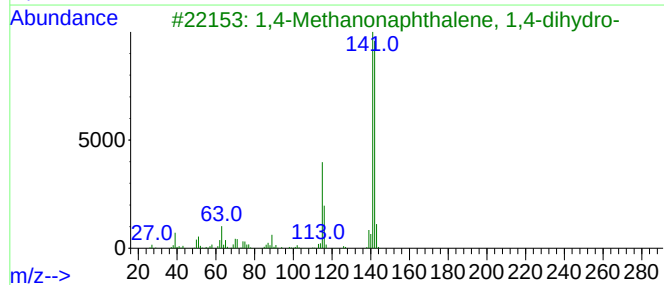
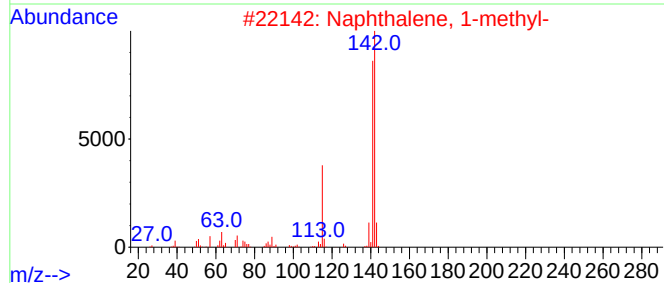
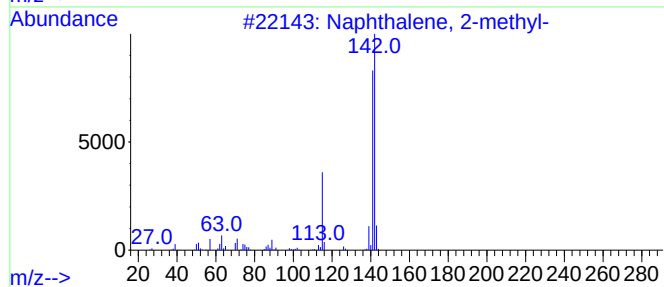
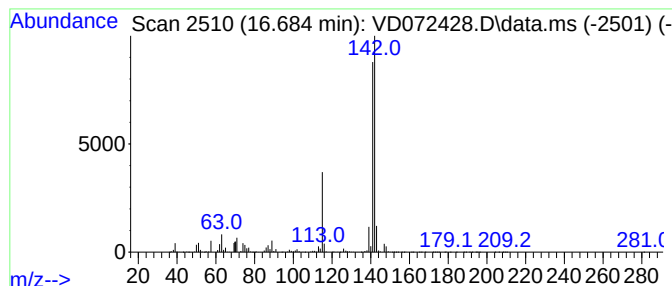
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.684	3217.69 ug/l	30839300	1,4-Dichlorobenzene-d4	13.567

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	93
4			Benzocycloheptatriene	142	C11H10	000264-09-5	90
5			1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD032522\
Data File : VD072428.D
Acq On : 25 Mar 2022 14:47
Operator : VA/SY
Sample : N1977-04
Misc : 4.31G/5.0ml/MSVOA_D/SOIL
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
SB-12-20220315-16.0-17.0-A

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D031122S.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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o-Cymene	14.102	257.9	ug/l	2472140	4	13.567	479215	50.0
Indan, 1-methyl-	14.214	231.0	ug/l	2214370	4	13.567	479215	50.0
1H-Indene, 2,3-...	14.696	343.5	ug/l	3291940	4	13.567	479215	50.0
1H-Indene, 2,3-...	14.814	490.4	ug/l	4700160	4	13.567	479215	50.0
1H-Indene, 1-me...	14.967	355.3	ug/l	3405700	4	13.567	479215	50.0
Benzene, (1,2-d...	15.079	223.1	ug/l	2138190	4	13.567	479215	50.0
Benzene, (1-eth...	15.490	231.2	ug/l	2216170	4	13.567	479215	50.0
Naphthalene, 1-...	16.479	434.4	ug/l	4163940	4	13.567	479215	50.0
1,4-Methanonaph...	16.684	3217.7	ug/l	30839300	4	13.567	479215	50.0