

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD030321\
 Data File : VD068507.D
 Acq On : 03 Mar 2021 12:32
 Operator : VA/SY
 Sample : M1542-01
 Misc : 6.43G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 SU-03-030221

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\82D022621S.M
 Title : SW846 8260

Signal : TIC: VD068507.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	7.914	1009	1019	1025	rBV	280638	699645	33.40%	2.688%
2	7.979	1025	1030	1041	rVB	493332	1073201	51.24%	4.123%
3	8.332	1081	1090	1101	rBV	255511	556874	26.59%	2.139%
4	8.867	1172	1181	1197	rBV	709421	1461831	69.79%	5.616%
5	10.338	1423	1431	1439	rBV	1150165	2094525	100.00%	8.047%
6	10.402	1439	1442	1451	rVB3	13914	26544	1.27%	0.102%
7	11.643	1645	1653	1664	rBV	1043653	1794635	85.68%	6.895%
8	11.855	1682	1689	1694	rBV4	26289	49290	2.35%	0.189%
9	12.185	1735	1745	1751	rVB5	19840	46257	2.21%	0.178%
10	12.485	1790	1796	1800	rBV4	11768	24373	1.16%	0.094%
11	12.632	1814	1821	1832	rVB	1143263	1924814	91.90%	7.395%
12	12.820	1845	1853	1857	rVB2	22359	41105	1.96%	0.158%
13	12.879	1857	1863	1867	rBV	65808	119949	5.73%	0.461%
14	12.949	1870	1875	1881	rVV2	203010	353884	16.90%	1.360%
15	13.008	1881	1885	1889	rVB5	24403	38621	1.84%	0.148%
16	13.120	1896	1904	1909	rBV	42488	105846	5.05%	0.407%
17	13.184	1911	1915	1920	rVV2	56980	84415	4.03%	0.324%
18	13.267	1923	1929	1934	rVV	249698	401588	19.17%	1.543%
19	13.361	1942	1945	1948	rBV5	17466	26192	1.25%	0.101%
20	13.396	1948	1951	1955	rVB3	29723	38117	1.82%	0.146%
21	13.461	1955	1962	1966	rBV6	95275	190787	9.11%	0.733%
22	13.508	1966	1970	1973	rVV3	85658	137946	6.59%	0.530%
23	13.573	1973	1981	1990	rVV2	1137482	2080464	99.33%	7.993%
24	13.643	1990	1993	1998	rVB2	68849	91631	4.37%	0.352%
25	13.767	2010	2014	2018	rBV2	172456	285817	13.65%	1.098%
26	13.808	2018	2021	2025	rVB2	82320	119023	5.68%	0.457%
27	13.890	2030	2035	2040	rBV	651025	966424	46.14%	3.713%
28	13.943	2040	2044	2046	rBV4	28503	39548	1.89%	0.152%
29	13.967	2046	2048	2055	rVB3	37422	61573	2.94%	0.237%
30	14.032	2055	2059	2060	rBV	76310	91418	4.36%	0.351%
31	14.055	2060	2063	2068	rVB2	137941	181983	8.69%	0.699%
32	14.114	2068	2073	2078	rVB	186277	259004	12.37%	0.995%
33	14.149	2078	2079	2086	rVB6	60365	97335	4.65%	0.374%
34	14.226	2086	2092	2097	rVB	194077	322520	15.40%	1.239%
35	14.284	2097	2102	2103	rBV3	81745	99315	4.74%	0.382%
36	14.373	2110	2117	2118	rBV4	93949	183003	8.74%	0.703%

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Peak Location: TOP

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37	14.402	2118	2122	2127	rVV4	277965	555482	26.52%	2.134%
38	14.443	2127	2129	2132	rVV2	142308	208032	9.93%	0.799%
39	14.479	2132	2135	2138	rVV	142761	182621	8.72%	0.702%
40	14.520	2138	2142	2146	rVV3	147868	184020	8.79%	0.707%
41	14.567	2146	2150	2155	rVB5	140897	222227	10.61%	0.854%
42	14.620	2156	2159	2163	rBV5	68883	104306	4.98%	0.401%
43	14.661	2163	2166	2169	rVV	83155	102885	4.91%	0.395%
44	14.708	2169	2174	2177	rVV3	202811	350554	16.74%	1.347%
45	14.743	2177	2180	2186	rVV3	541235	831247	39.69%	3.194%
46	14.826	2187	2194	2198	rVV4	450330	1014095	48.42%	3.896%
47	14.867	2198	2201	2209	rVB3	240135	470152	22.45%	1.806%
48	14.984	2215	2221	2227	rVB	288634	531205	25.36%	2.041%
49	15.090	2234	2239	2243	rBV3	125436	194208	9.27%	0.746%
50	15.202	2252	2258	2267	rBV4	124208	294631	14.07%	1.132%
51	15.284	2268	2272	2279	rVB7	107331	248472	11.86%	0.955%
52	15.361	2279	2285	2288	rBV2	156338	295595	14.11%	1.136%
53	15.396	2288	2291	2296	rVB	111282	172331	8.23%	0.662%
54	15.455	2297	2301	2305	rVB	100127	146732	7.01%	0.564%
55	15.508	2305	2310	2314	rBV3	105749	223253	10.66%	0.858%
56	15.590	2320	2324	2334	rVB	368695	683096	32.61%	2.624%
57	15.696	2334	2342	2349	rBV4	138425	366762	17.51%	1.409%
58	15.767	2351	2354	2361	rVB7	66349	121612	5.81%	0.467%
59	15.873	2363	2372	2373	rVV4	76342	176302	8.42%	0.677%
60	15.902	2373	2377	2384	rVB	258055	484708	23.14%	1.862%
61	16.067	2400	2405	2413	rVB6	66126	158264	7.56%	0.608%
62	16.226	2423	2432	2437	rBV4	140449	363905	17.37%	1.398%
63	16.337	2447	2451	2458	rBV3	47774	81947	3.91%	0.315%
64	16.426	2459	2466	2472	rVV3	132153	284950	13.60%	1.095%
65	16.490	2472	2477	2482	rVV4	93501	195080	9.31%	0.749%
66	16.549	2482	2487	2494	rVB	163141	299683	14.31%	1.151%
67	16.702	2505	2513	2527	rVB	68443	311372	14.87%	1.196%

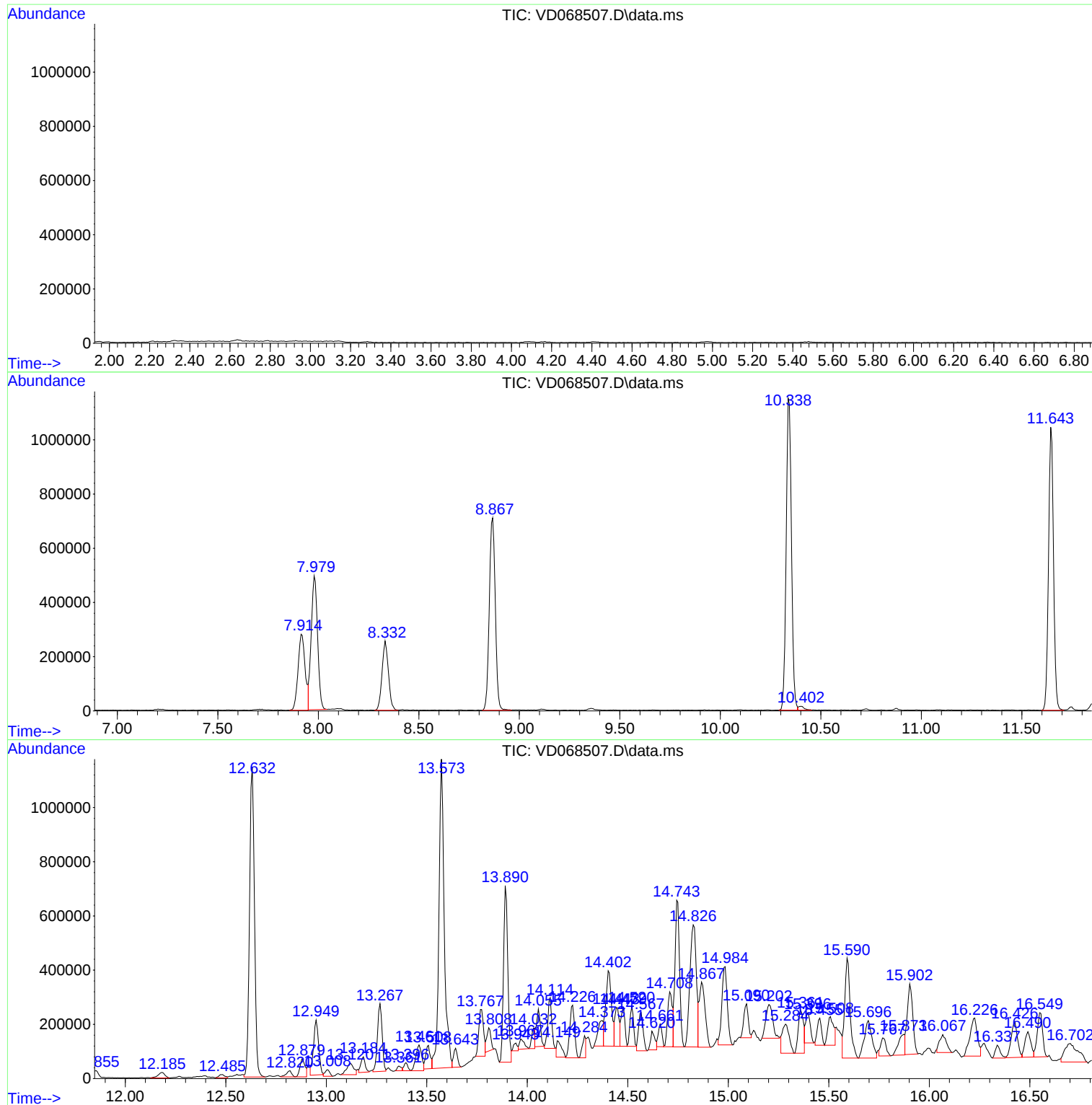
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D022621S.M
Quant Title : SW846 8260

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



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Instrument :
MSVOA_D
ClientSampleId :
SU-03-030221

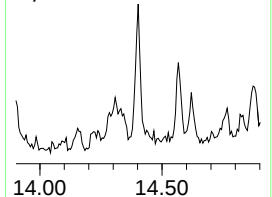
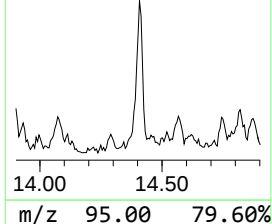
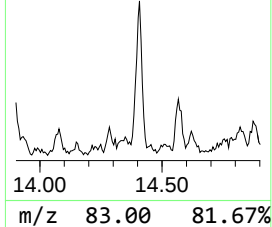
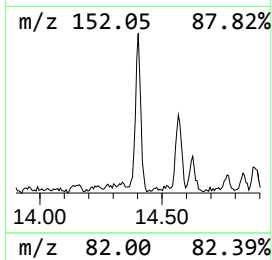
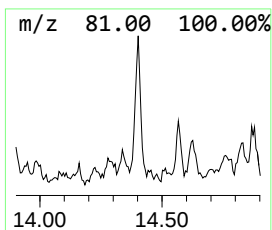
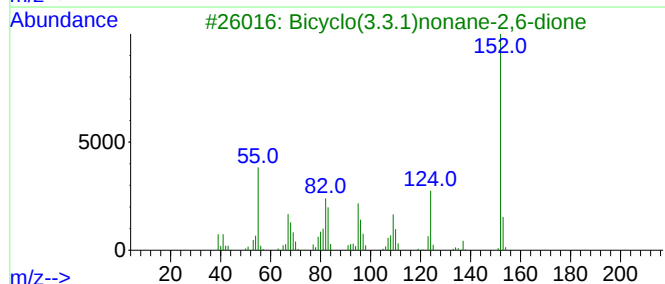
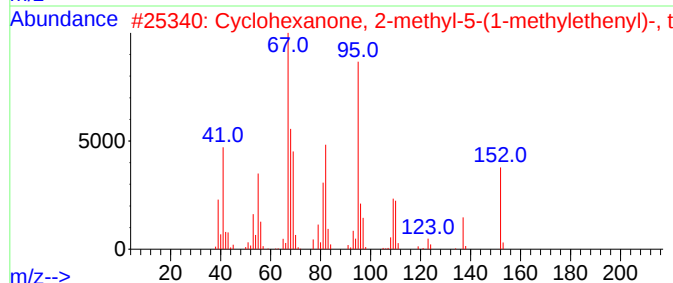
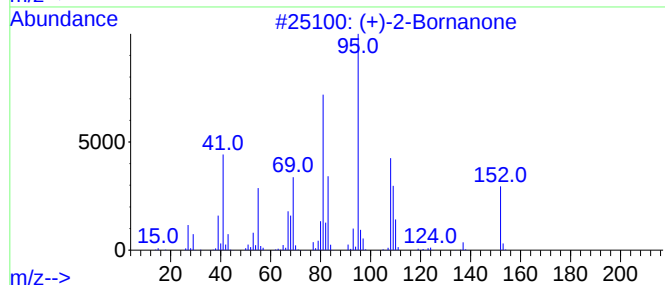
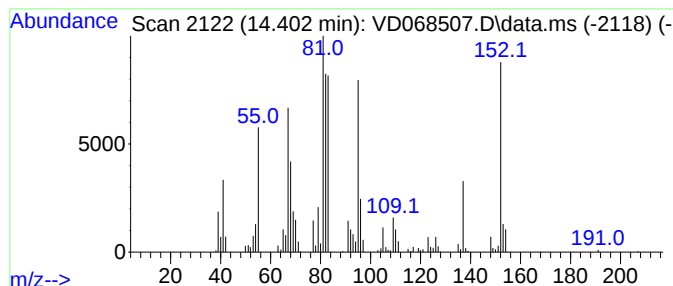
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D022621S.M
Quant Title : SW846 8260

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

Peak Number 1 (+)-2-Bornanone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.402	13.35 ug/l	555482	1,4-Dichlorobenzene-d4	13.573

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		(+)-2-Bornanone	152	C10H16O	000464-49-3	70
2		Cyclohexanone, 2-methyl-5-(1-met...	152	C10H16O	005948-04-9	46
3		Bicyclo(3.3.1)nonane-2,6-dione	152	C9H12O2	016473-11-3	46
4		Pulegone	152	C10H16O	000089-82-7	43
5		Cyclohexanone, 5-methyl-2-(1-met...	152	C10H16O	015932-80-6	43



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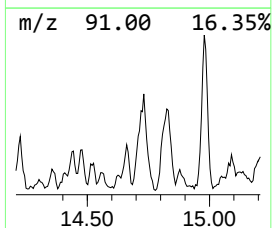
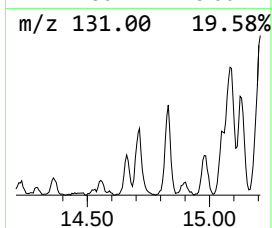
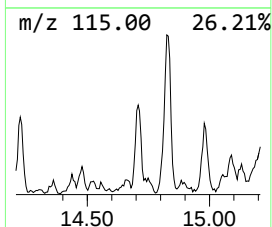
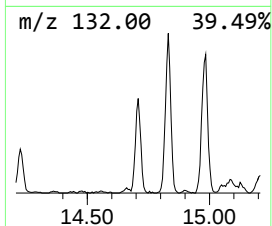
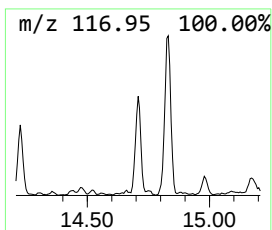
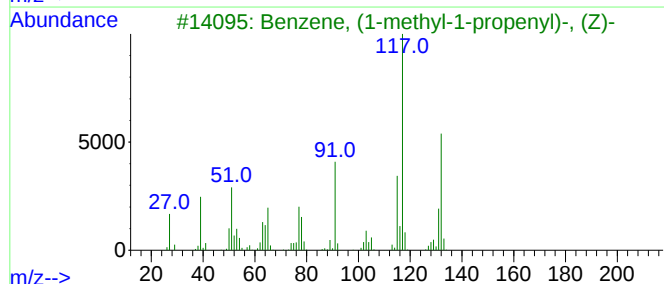
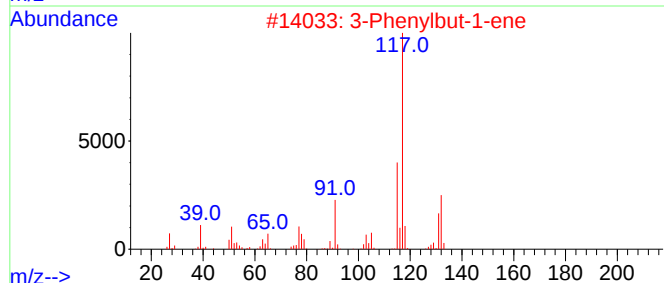
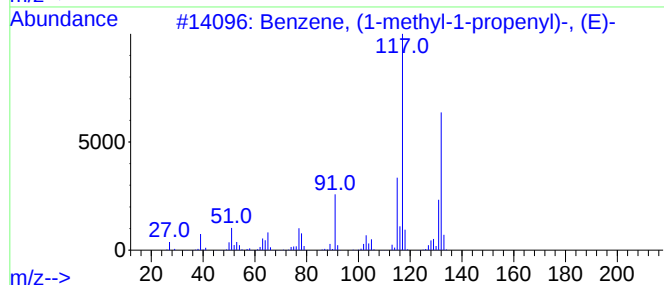
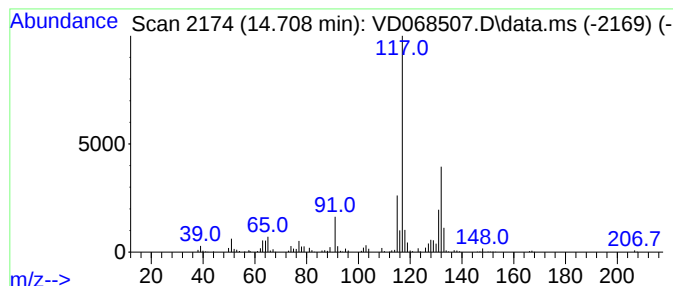
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D022621S.M
Quant Title : SW846 8260

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

Peak Number 2 Benzene, (1-methyl-1-propen... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.708	8.42 ug/l	350554	1,4-Dichlorobenzene-d4	13.573

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	87
2			3-Phenylbut-1-ene	132	C10H12	000934-10-1	87
3			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	86
4			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	80
5			Indan, 1-methyl-	132	C10H12	000767-58-8	80



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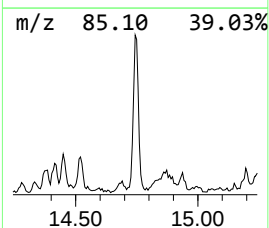
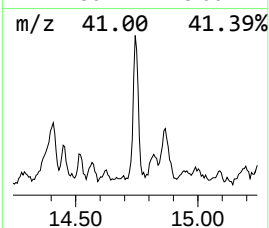
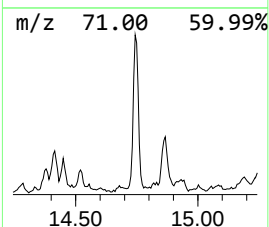
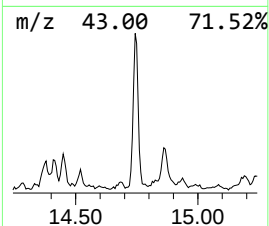
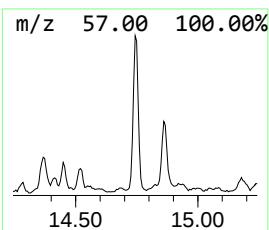
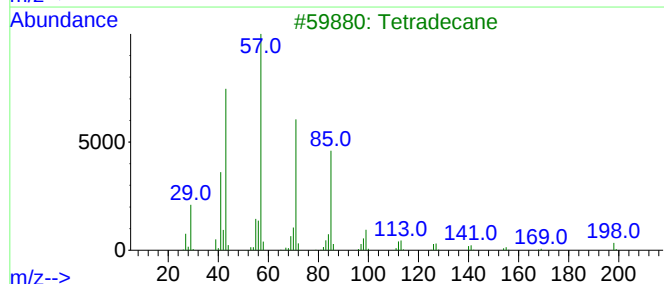
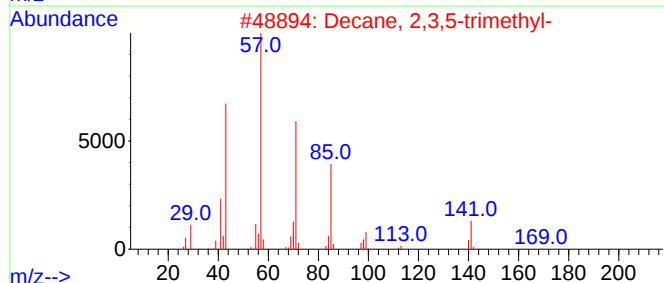
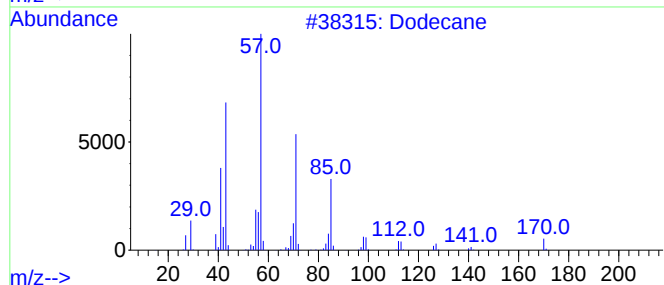
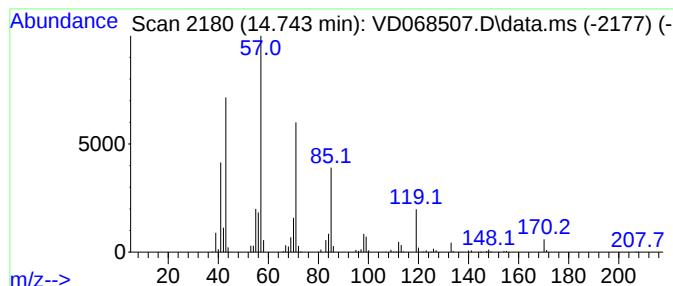
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D022621S.M
Quant Title : SW846 8260

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

Peak Number 3 Dodecane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.743	19.98 ug/l	831247	1,4-Dichlorobenzene-d4	13.573

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane	170	C12H26	000112-40-3	81
2			Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	59
3			Tetradecane	198	C14H30	000629-59-4	59
4			Sulfurous acid, 2-ethylhexyl hex...	278	C14H30O3S	1000309-20-2	53
5			Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	46



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ClientSampleId :
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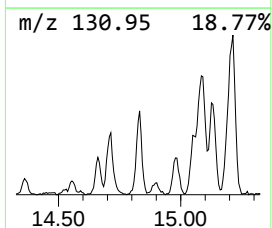
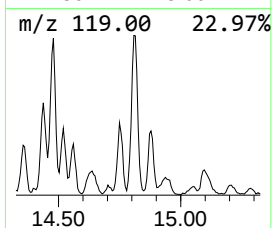
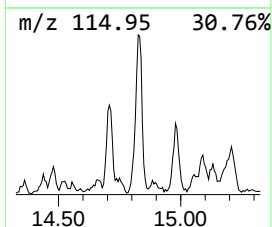
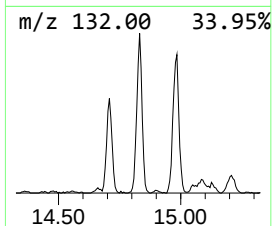
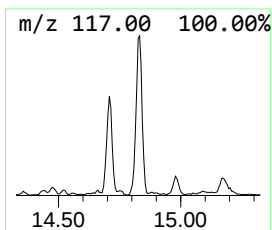
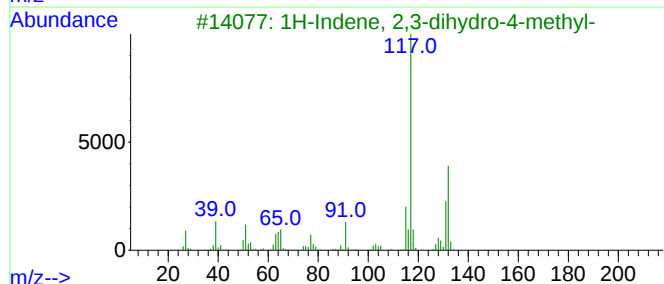
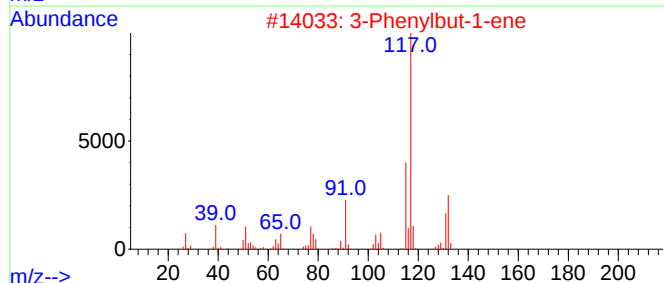
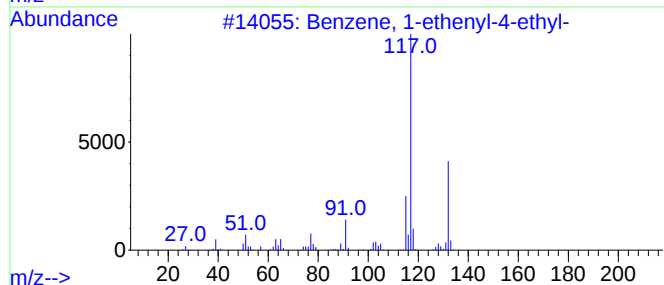
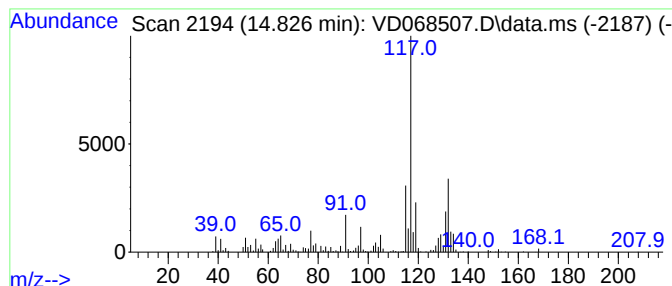
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D022621S.M
Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.826	24.37 ug/l	1014100	1,4-Dichlorobenzene-d4	13.573

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	93
2			3-Phenylbut-1-ene	132	C10H12	000934-10-1	76
3			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	76
4			1-Phenyl-1-butene	132	C10H12	000824-90-8	74
5			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	74



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD030321\
Data File : VD068507.D
Acq On : 03 Mar 2021 12:32
Operator : VA/SY
Sample : M1542-01
Misc : 6.43G/5.00ml/MSVOA_D/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
SU-03-030221

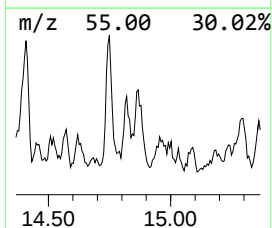
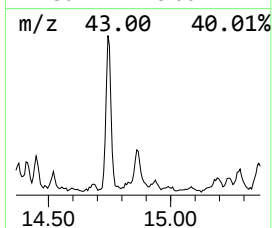
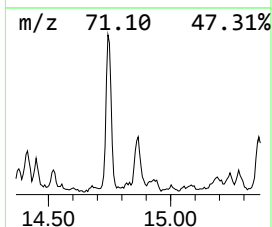
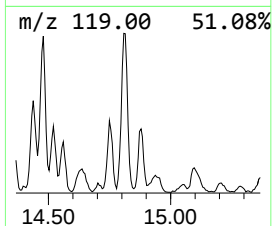
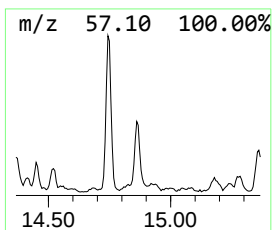
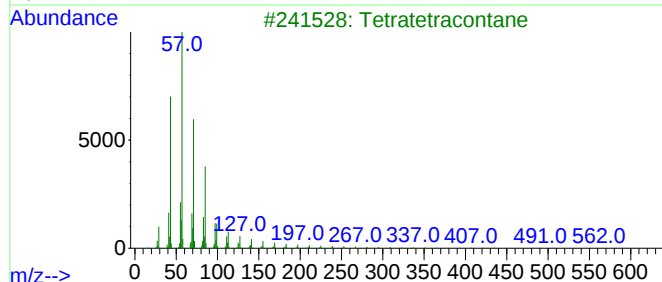
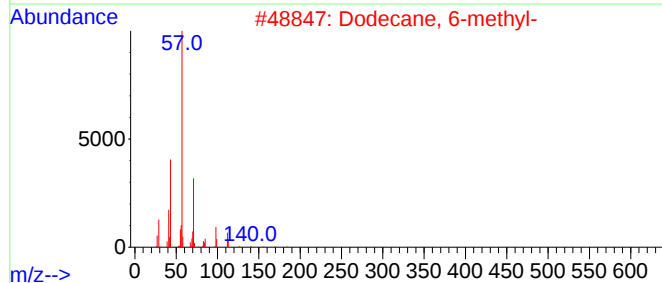
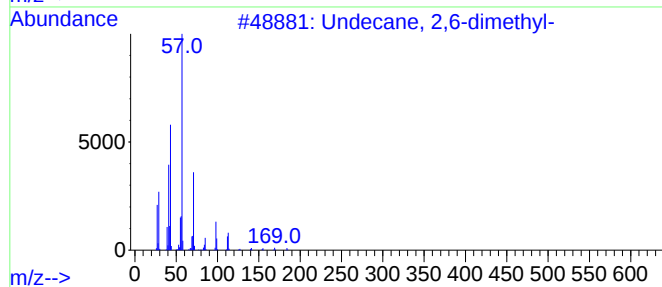
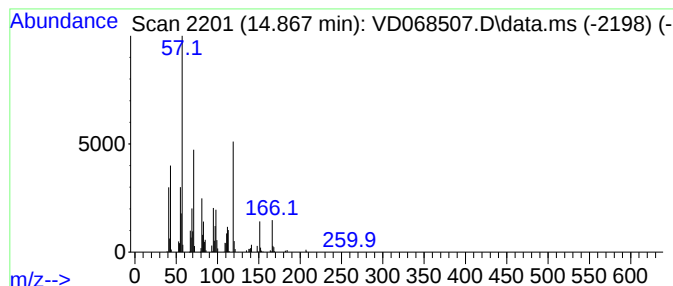
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Quant Title : SW846 8260

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

Peak Number 5 unknown14.867 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.867	11.30 ug/l	470152	1,4-Dichlorobenzene-d4	13.573

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane, 2,6-dimethyl-	184	C13H28	017301-23-4	46
2			Dodecane, 6-methyl-	184	C13H28	006044-71-9	30
3			Tetratetracontane	619	C44H90	007098-22-8	27
4			Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	27
5			Nonane, 3-methyl-	142	C10H22	005911-04-6	22



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD030321\
Data File : VD068507.D
Acq On : 03 Mar 2021 12:32
Operator : VA/SY
Sample : M1542-01
Misc : 6.43G/5.00ml/MSVOA_D/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
SU-03-030221

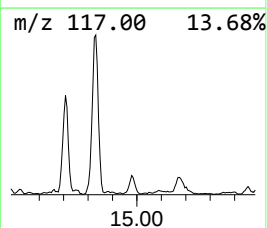
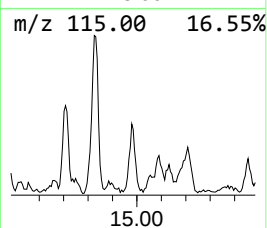
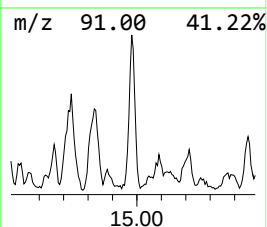
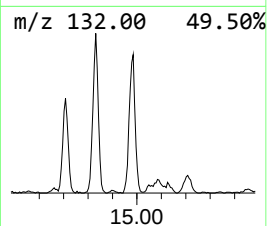
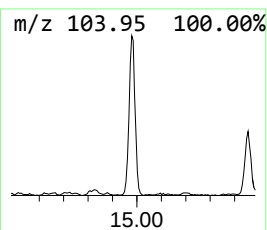
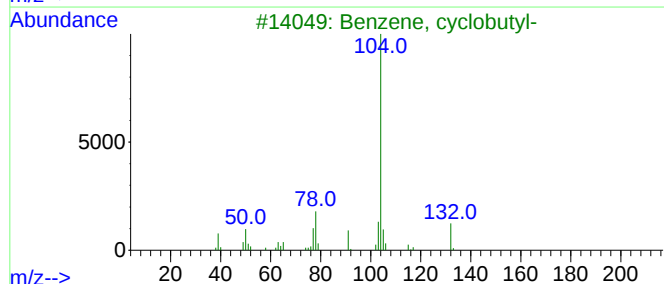
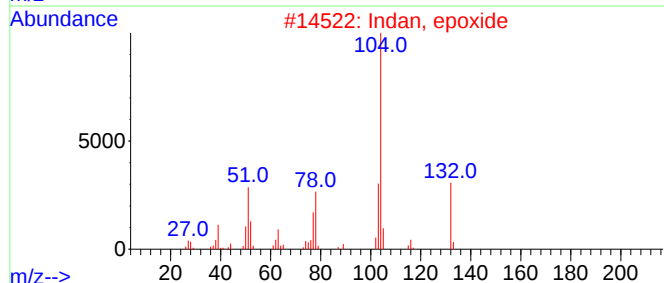
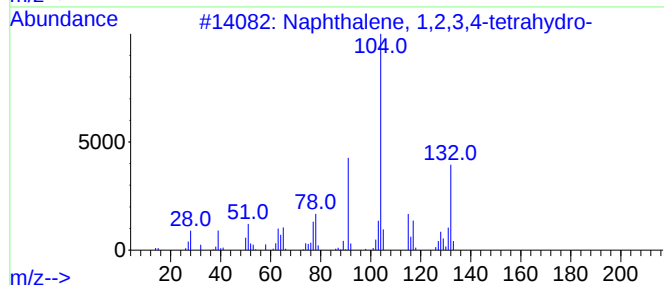
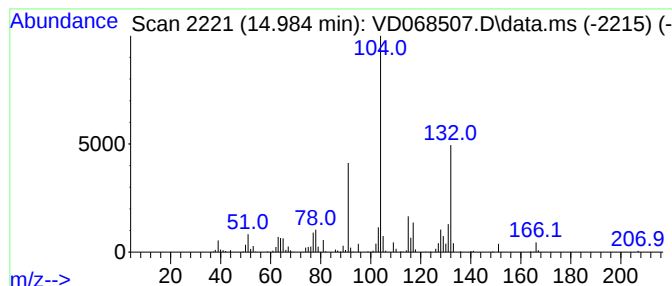
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Quant Title : SW846 8260

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

Peak Number 6 Naphthalene, 1,2,3,4-tetrahydro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.984	12.77 ug/l	531205	1,4-Dichlorobenzene-d4	13.573

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	95
2			Indan, epoxide	132	C9H8O	000768-22-9	58
3			Benzene, cyclobutyl-	132	C10H12	004392-30-7	53
4			Naphthalene, 2-ethyl-1,2,3,4-tet...	160	C12H16	032367-54-7	50
5			Phensuximide	189	C11H11NO2	000086-34-0	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD030321\
Data File : VD068507.D
Acq On : 03 Mar 2021 12:32
Operator : VA/SY
Sample : M1542-01
Misc : 6.43G/5.00ml/MSVOA_D/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
SU-03-030221

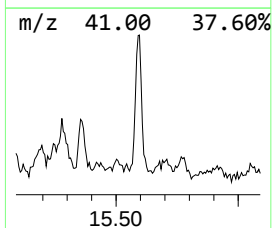
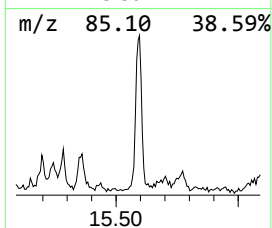
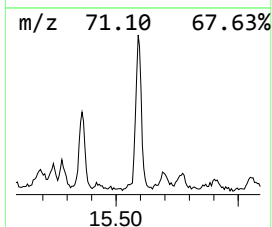
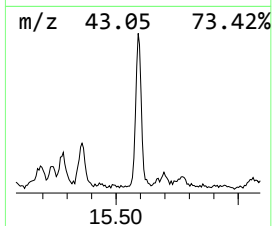
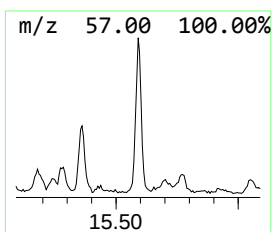
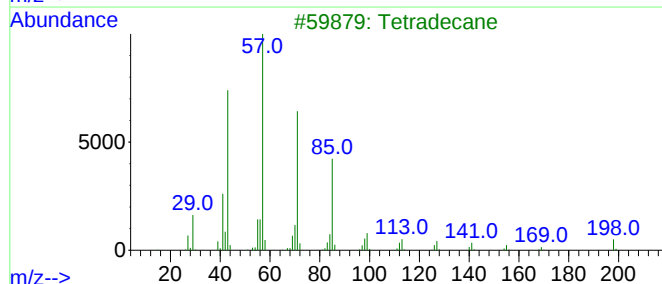
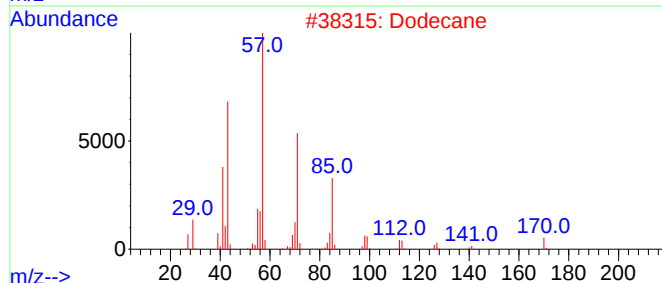
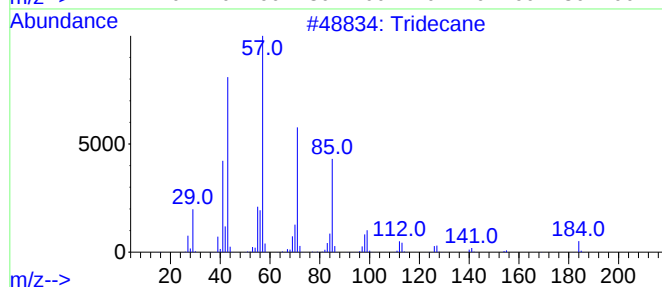
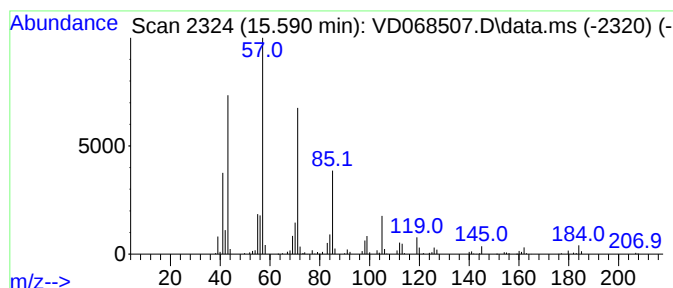
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Quant Title : SW846 8260

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

Peak Number 7 Tridecane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.590	16.42 ug/l	683096	1,4-Dichlorobenzene-d4	13.573

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane	184	C13H28	000629-50-5	96
2			Dodecane	170	C12H26	000112-40-3	72
3			Tetradecane	198	C14H30	000629-59-4	72
4			Nonadecane	268	C19H40	000629-92-5	64
5			Decane, 2,3,6-trimethyl-	184	C13H28	062238-12-4	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD030321\
Data File : VD068507.D
Acq On : 03 Mar 2021 12:32
Operator : VA/SY
Sample : M1542-01
Misc : 6.43G/5.00ml/MSVOA_D/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
SU-03-030221

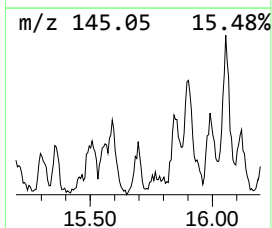
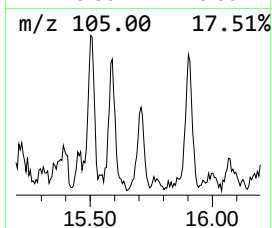
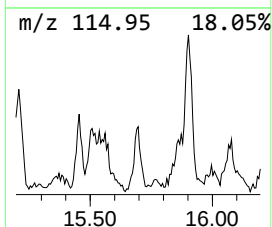
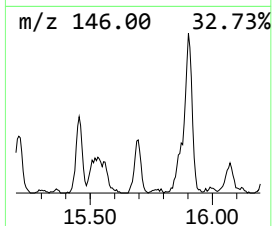
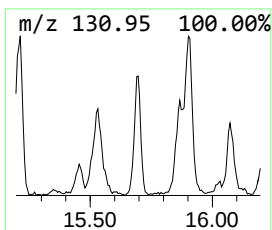
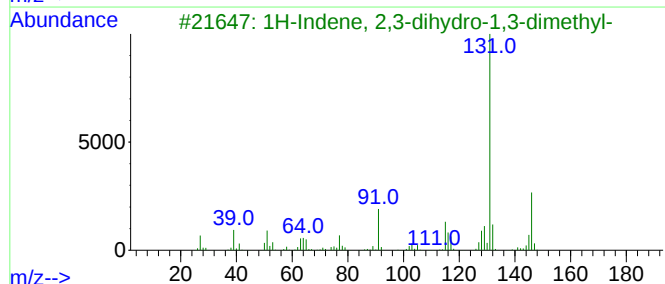
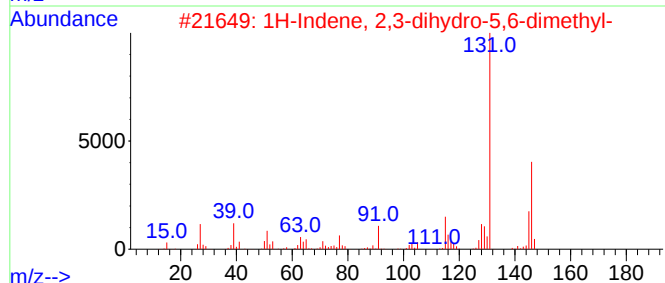
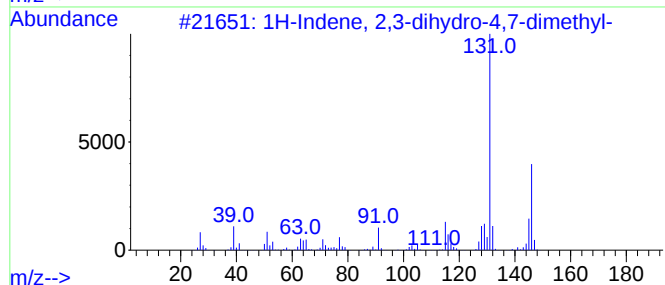
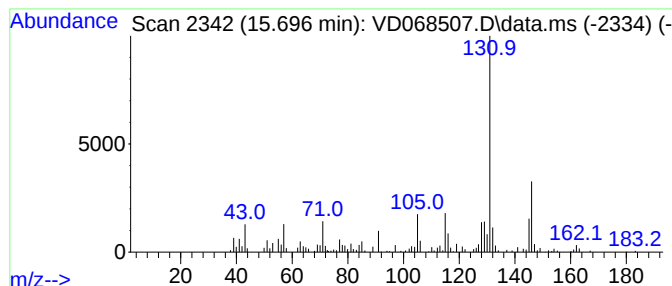
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Quant Title : SW846 8260

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

Peak Number 8 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.696	8.81 ug/l	366762	1,4-Dichlorobenzene-d4	13.573

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	94
2			1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	93
3			1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	90
4			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	90
5			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD030321\
Data File : VD068507.D
Acq On : 03 Mar 2021 12:32
Operator : VA/SY
Sample : M1542-01
Misc : 6.43G/5.00ml/MSVOA_D/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
SU-03-030221

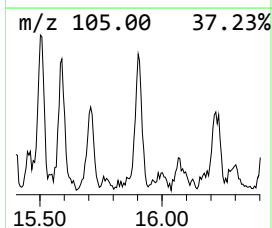
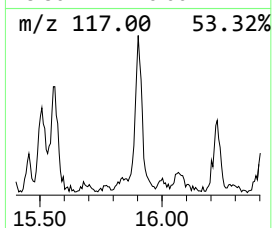
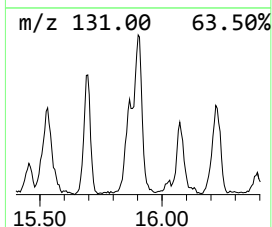
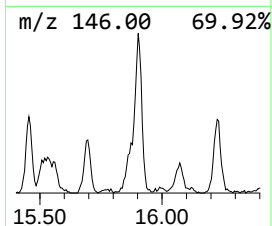
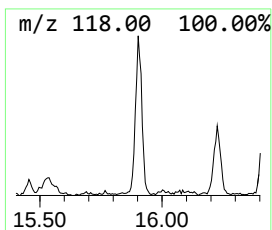
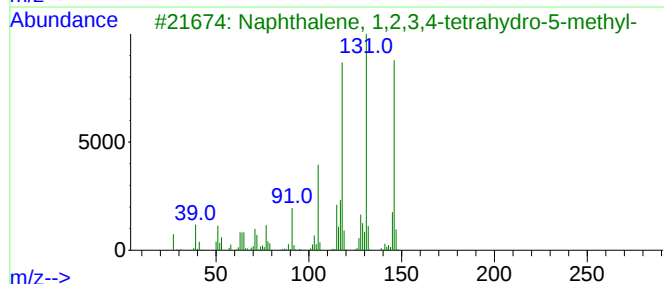
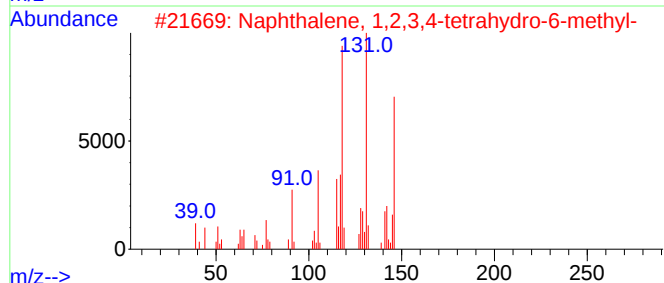
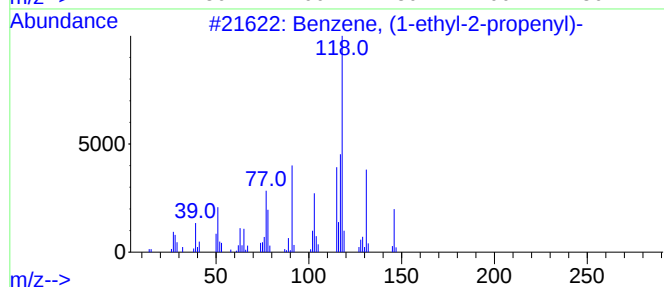
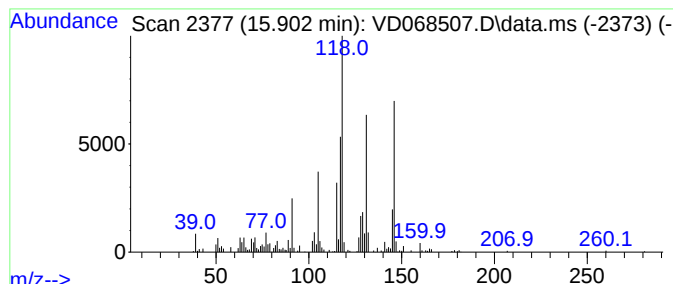
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D022621S.M
Quant Title : SW846 8260

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

Peak Number 9 Benzene, (1-ethyl-2-propenyl)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.902	11.65 ug/l	484708	1,4-Dichlorobenzene-d4	13.573

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-ethyl-2-propenyl)-	146	C11H14	019947-22-9	83
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	74
3			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	72
4			Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	64
5			Naphth[1,2-b]oxirene, 1a,2,3,7b-...	146	C10H10O	002461-34-9	49



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD030321\
Data File : VD068507.D
Acq On : 03 Mar 2021 12:32
Operator : VA/SY
Sample : M1542-01
Misc : 6.43G/5.00ml/MSVOA_D/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
SU-03-030221

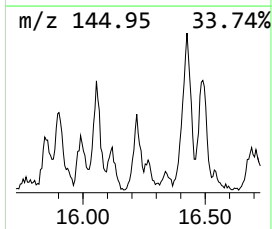
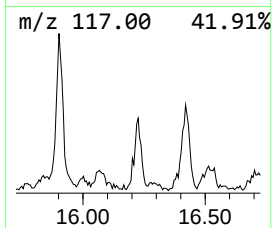
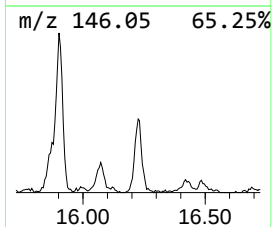
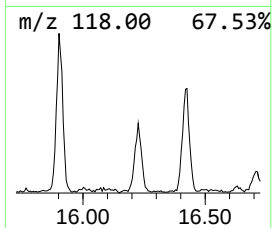
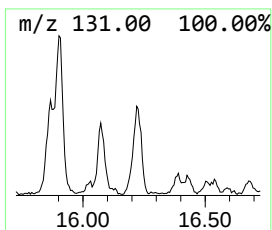
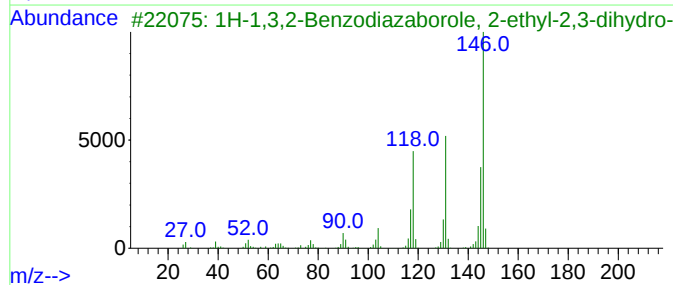
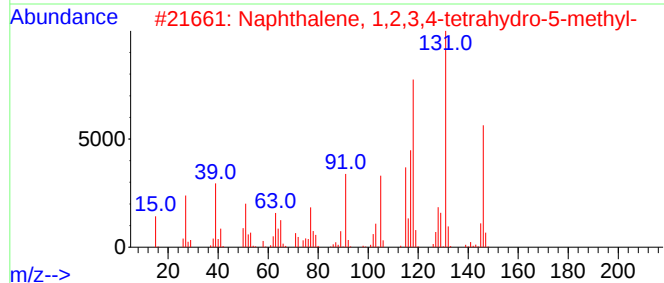
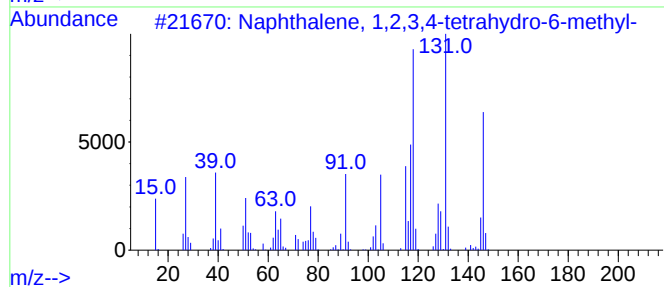
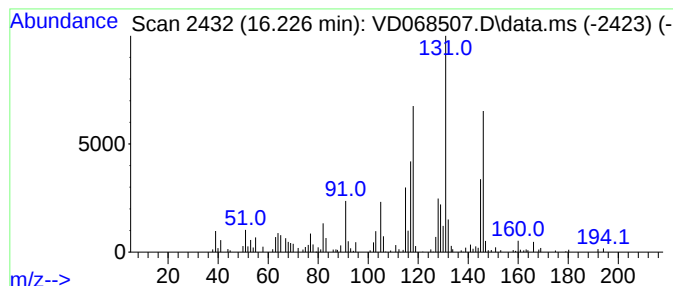
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D022621S.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.226	8.75 ug/l	363905	1,4-Dichlorobenzene-d4	13.573

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	90
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	87
3			1H-1,3,2-Benzodiazaborole, 2-eth...	146	C8H11BN2	074663-81-3	68
4			Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	53
5			Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	52



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD030321\
 Data File : VD068507.D
 Acq On : 03 Mar 2021 12:32
 Operator : VA/SY
 Sample : M1542-01
 Misc : 6.43G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 SU-03-030221

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D022621S.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-met...	14.708	8.4	ug/l	350554	4	13.573	2080460	50.0
Dodecane	14.743	20.0	ug/l	831247	4	13.573	2080460	50.0
Benzene, 1-ethe...	14.826	24.4	ug/l	1014100	4	13.573	2080460	50.0
unknown14.867	14.867	11.3	ug/l	470152	4	13.573	2080460	50.0
Naphthalene, 1,...	14.984	12.8	ug/l	531205	4	13.573	2080460	50.0
Tridecane	15.590	16.4	ug/l	683096	4	13.573	2080460	50.0
1H-Indene, 2,3-...	15.696	8.8	ug/l	366762	4	13.573	2080460	50.0
Benzene, (1-eth...	15.902	11.7	ug/l	484708	4	13.573	2080460	50.0
Naphthalene, 1,...	16.226	8.8	ug/l	363905	4	13.573	2080460	50.0