

Data Path : Z:\VOASRV\HPCHEM1\MSVOA Y\DATA\VY041520\
 Data File : VY002334.D
 Acq On : 15 Apr 2020 14:13
 Operator : SY/MD
 Sample : L2245-08
 Misc : 7.15G/5ML/MSVOA Y/SOIL
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 TP-26

Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_Y\METHODS\82Y040820S.M
 Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.255	77	85	92	rBV2	8000	15302	1.54%	0.142%
2	3.968	356	366	376	rBV	6122	17916	1.80%	0.166%
3	4.736	483	492	498	rBV2	8277	26903	2.71%	0.250%
4	7.736	972	984	989	rBV	122093	287176	28.88%	2.668%
5	7.803	989	995	1005	rVV	199941	454722	45.73%	4.224%
6	7.882	1005	1008	1018	rVB4	4986	11376	1.14%	0.106%
7	8.156	1043	1053	1064	rBV	114813	258269	25.97%	2.399%
8	8.339	1066	1083	1092	rBV	36691	99835	10.04%	0.927%
9	8.699	1133	1142	1151	rBV	347968	689322	69.32%	6.404%
10	9.193	1209	1223	1228	rBV6	12202	32647	3.28%	0.303%
11	9.248	1228	1232	1240	rVB3	6506	12394	1.25%	0.115%
12	9.333	1240	1246	1251	rBV	6213	12602	1.27%	0.117%
13	9.626	1286	1294	1303	rVB	41228	81015	8.15%	0.753%
14	9.754	1305	1315	1323	rBV	74496	156753	15.76%	1.456%
15	9.943	1340	1346	1352	rBV4	4813	11155	1.12%	0.104%
16	10.119	1363	1375	1379	rBV3	13243	26907	2.71%	0.250%
17	10.186	1379	1386	1399	rVV	580198	994427	100.00%	9.238%
18	10.351	1409	1413	1423	rVB10	5548	16750	1.68%	0.156%
19	10.613	1449	1456	1464	rBV8	11871	31198	3.14%	0.290%
20	11.040	1523	1526	1530	rVV	8892	14993	1.51%	0.139%
21	11.095	1531	1535	1541	rVV3	18404	34452	3.46%	0.320%
22	11.156	1541	1545	1549	rVV6	7395	13361	1.34%	0.124%
23	11.302	1564	1569	1577	rVB6	10501	26293	2.64%	0.244%
24	11.497	1594	1601	1609	rBV	507588	832751	83.74%	7.736%
25	11.686	1627	1632	1635	rBV7	7110	14787	1.49%	0.137%
26	11.741	1637	1641	1645	rBV3	14813	27837	2.80%	0.259%
27	12.009	1678	1685	1692	rVB8	19251	46251	4.65%	0.430%
28	12.125	1701	1704	1708	rVB	14808	20331	2.04%	0.189%
29	12.204	1713	1717	1720	rBV3	21265	31710	3.19%	0.295%
30	12.485	1756	1763	1771	rBV	397546	703535	70.75%	6.536%
31	12.649	1785	1790	1795	rVB2	36374	58494	5.88%	0.543%
32	12.723	1798	1802	1810	rVB7	16313	40212	4.04%	0.374%
33	12.808	1811	1816	1822	rBV7	22807	50887	5.12%	0.473%
34	12.954	1836	1840	1844	rBV4	17368	24498	2.46%	0.228%

Data Path : Z:\VOASRV\HPCHEM1\MSVOA Y\DATA\VY041520\
 Data File : VY002334.D
 Acq On : 15 Apr 2020 14:13
 Operator : SY/MD
 Sample : L2245-08
 Misc : 7.15G/5ML/MSVOA Y/SOIL
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 TP-26

Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_Y\METHODS\82Y040820S.M
 Title : SW846 8260

35	13.015	1844	1850	1851	rBV4	8206	15184	1.53%	0.141%
36	13.046	1851	1855	1862	rVV5	25871	48817	4.91%	0.453%
37	13.174	1872	1876	1881	rBV4	14559	24152	2.43%	0.224%
38	13.253	1882	1889	1892	rBV4	19774	43336	4.36%	0.403%
39	13.296	1892	1896	1898	rVV4	15491	22894	2.30%	0.213%
40	13.326	1898	1901	1904	rVB	17293	21092	2.12%	0.196%
41	13.430	1912	1918	1924	rVB	478490	715701	71.97%	6.649%
42	13.533	1931	1935	1939	rBV3	25002	46830	4.71%	0.435%
43	13.594	1942	1945	1949	rVB	33105	43820	4.41%	0.407%
44	13.686	1956	1960	1965	rBV7	17717	37829	3.80%	0.351%
45	13.753	1965	1971	1976	rBV4	89892	179343	18.03%	1.666%
46	13.857	1986	1988	1992	rBV4	11266	17558	1.77%	0.163%
47	13.972	2002	2007	2012	rVB	98830	170653	17.16%	1.585%
48	14.027	2012	2016	2020	rBV3	20003	37863	3.81%	0.352%
49	14.082	2020	2025	2030	rVB4	119796	205800	20.70%	1.912%
50	14.149	2030	2036	2041	rVB5	50838	89772	9.03%	0.834%
51	14.204	2041	2045	2047	rBV4	25824	38647	3.89%	0.359%
52	14.277	2050	2057	2059	rBV4	96420	213146	21.43%	1.980%
53	14.369	2067	2072	2077	rBV4	59812	136991	13.78%	1.273%
54	14.424	2077	2081	2085	rVV2	54574	83225	8.37%	0.773%
55	14.460	2085	2087	2089	rVV2	14351	14958	1.50%	0.139%
56	14.625	2110	2114	2117	rBV2	31116	42332	4.26%	0.393%
57	14.692	2117	2125	2129	rVV2	178509	366288	36.83%	3.403%
58	14.741	2129	2133	2139	rVB4	119316	218075	21.93%	2.026%
59	14.850	2147	2151	2154	rBV4	37524	52278	5.26%	0.486%
60	14.917	2159	2162	2163	rBV3	18545	21721	2.18%	0.202%
61	14.948	2164	2167	2170	rVB3	71266	90167	9.07%	0.838%
62	14.991	2170	2174	2176	rBV	59893	83286	8.38%	0.774%
63	15.033	2176	2181	2182	rBV2	86161	115584	11.62%	1.074%
64	15.216	2206	2211	2218	rBV2	208186	340607	34.25%	3.164%
65	15.295	2219	2224	2229	rBV	116981	204683	20.58%	1.901%
66	15.375	2229	2237	2240	rBV2	68759	180559	18.16%	1.677%
67	15.533	2256	2263	2270	rBV4	91637	236385	23.77%	2.196%
68	15.612	2271	2276	2281	rVB5	77677	135184	13.59%	1.256%
69	15.698	2283	2290	2299	rVV5	71584	220326	22.16%	2.047%
70	15.887	2317	2321	2327	rVB3	48888	103318	10.39%	0.960%
71	16.039	2340	2346	2351	rBV	133254	263835	26.53%	2.451%

Data Path : Z:\VOASRV\HPCHEM1\MSVOA Y\DATA\VY041520\
Data File : VY002334.D
Acq On : 15 Apr 2020 14:13
Operator : SY/MD
Sample : L2245-08
Misc : 7.15G/5ML/MSVOA Y/SOIL
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TP-26

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\82Y040820S.M
Title : SW846 8260

72	16.167	2361	2367	2372	rBV	167794	294921	29.66%	2.740%
73	16.240	2374	2379	2385	rVB2	82811	180174	18.12%	1.674%
74	16.490	2414	2420	2426	rBV8	78858	230354	23.16%	2.140%

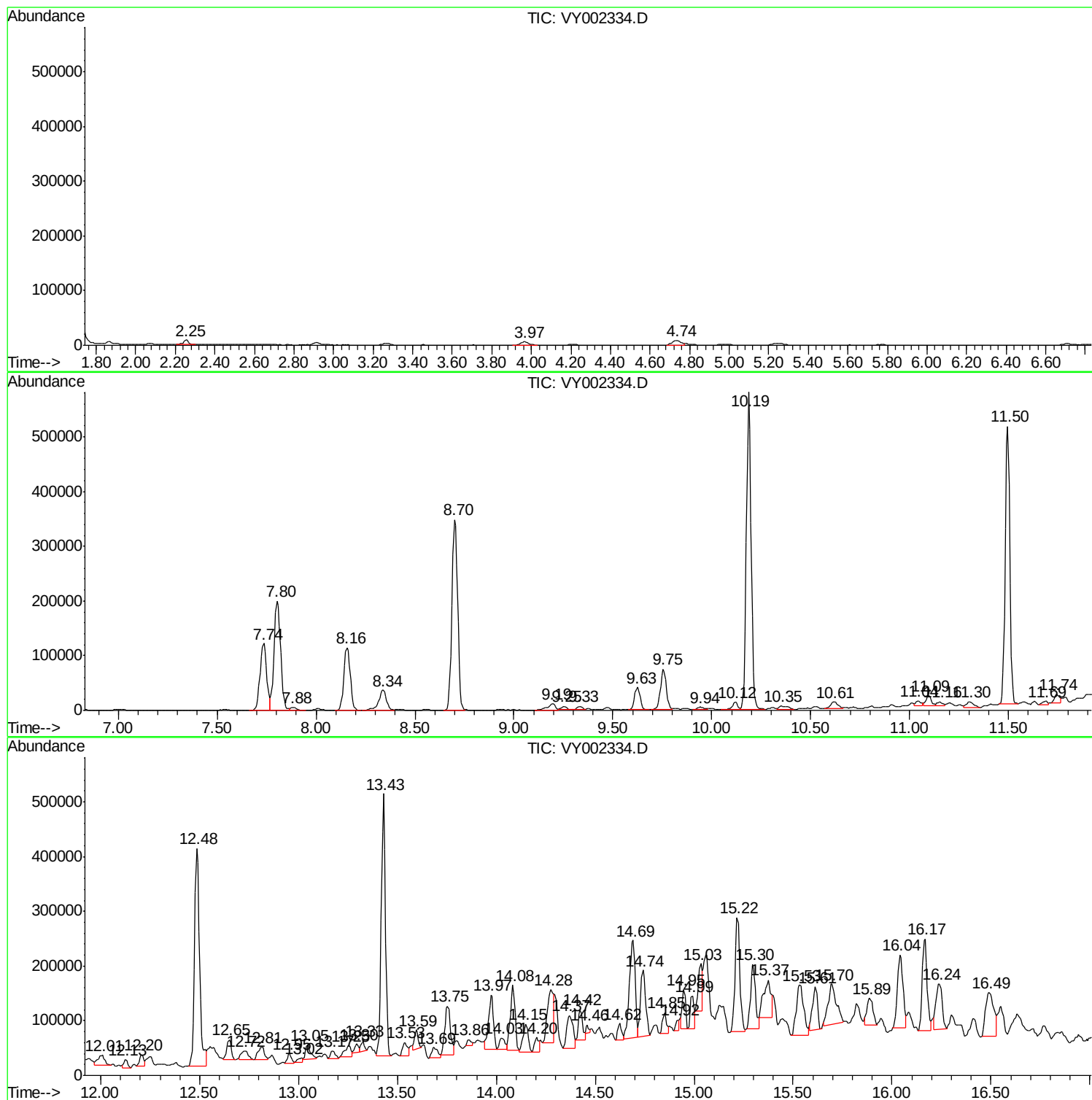
Sum of corrected areas: 10764749

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA Y\DATA\VY041520\
Data File : VY002334.D
Acq On : 15 Apr 2020 14:13
Operator : SY/MD
Sample : L2245-08
Misc : 7.15G/5ML/MSVOA Y/SOIL
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TP-26

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y040820S.M
Quant Title : SW846 8260

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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA Y\DATA\VY041520\
Data File : VY002334.D
Acq On : 15 Apr 2020 14:13
Operator : SY/MD
Sample : L2245-08
Misc : 7.15G/5ML/MSVOA Y/SOIL
ALS Vial : 11 Sample Multiplier: 1

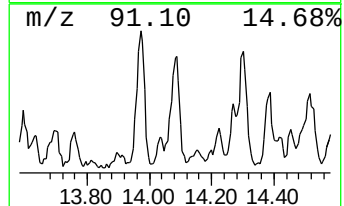
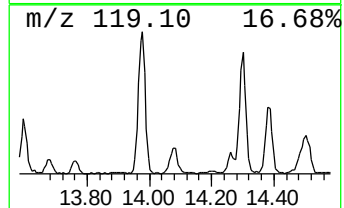
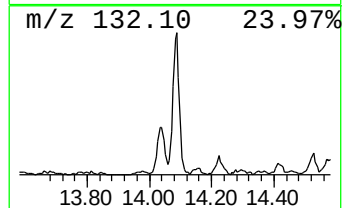
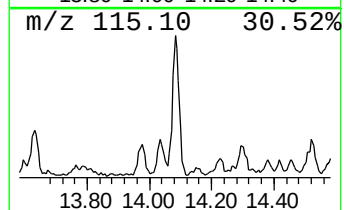
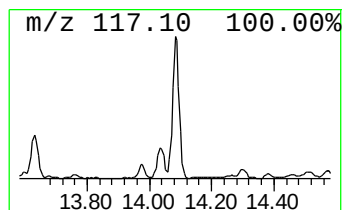
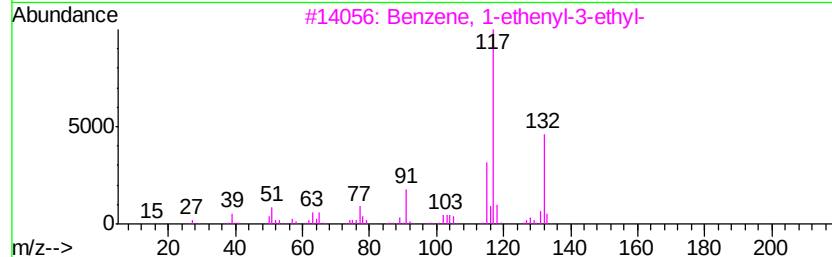
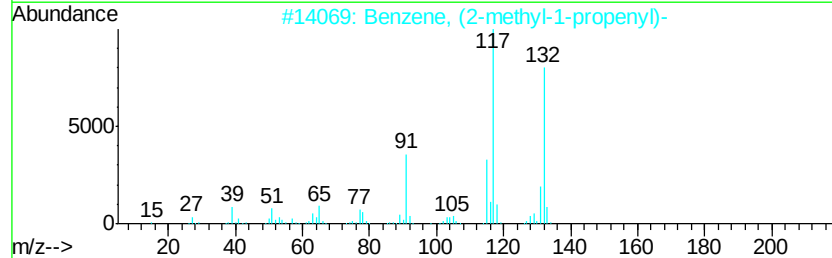
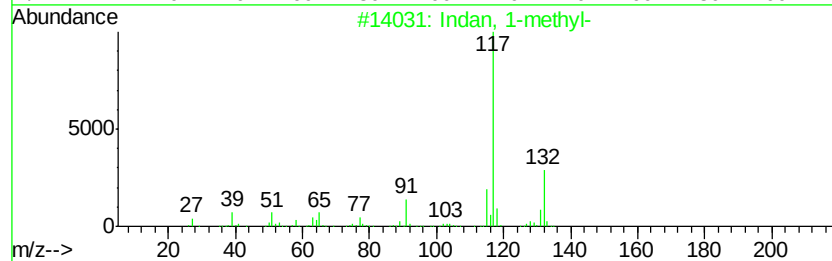
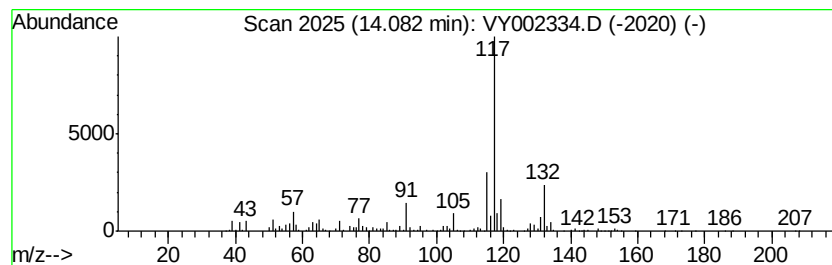
Instrument :
MSVOA_Y
ClientSampleId :
TP-26

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y040820S.M
Quant Title : SW846 8260

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

Peak Number 1 Indan, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.08	14.38 ug/l	205800	1,4-Dichlorobenzene-d4	13.43
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Indan, 1-methyl-	132 C10H12	000767-58-8	93
2	Benzene, (2-methyl-1-propenyl)-	132 C10H12	000768-49-0	87
3	Benzene, 1-ethenyl-3-ethyl-	132 C10H12	007525-62-4	83
4	3-Phenylbut-1-ene	132 C10H12	000934-10-1	83
5	Benzene, (2-methyl-2-propenyl)-	132 C10H12	003290-53-7	74



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA Y\DATA\VY041520\
Data File : VY002334.D
Acq On : 15 Apr 2020 14:13
Operator : SY/MD
Sample : L2245-08
Misc : 7.15G/5ML/MSVOA Y/SOIL
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TP-26

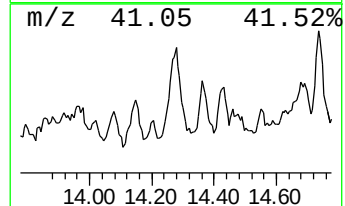
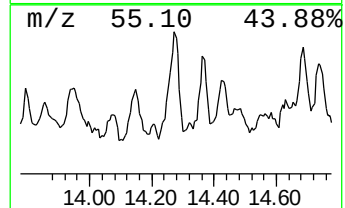
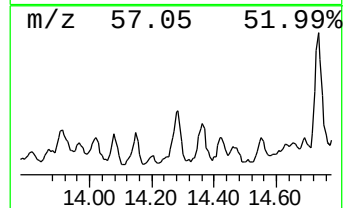
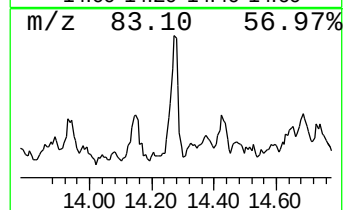
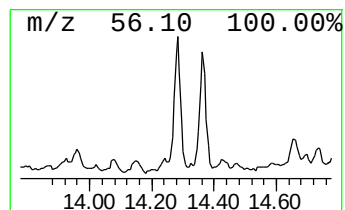
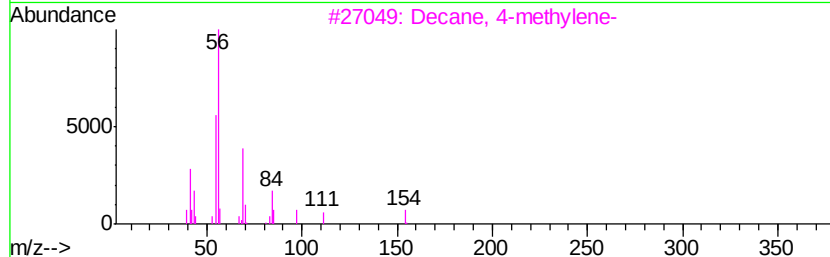
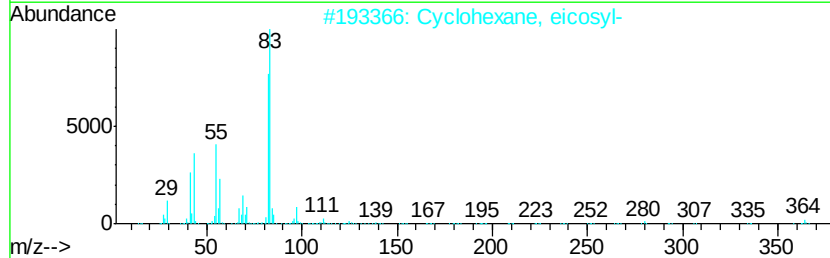
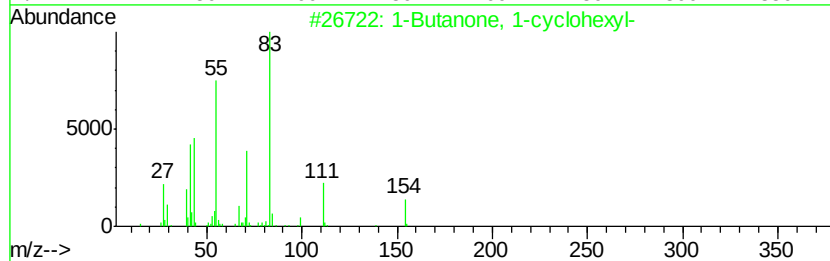
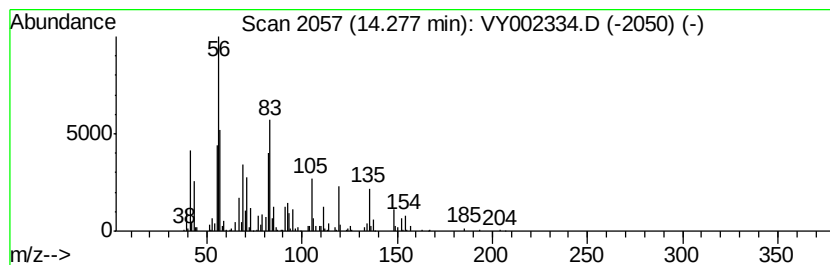
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y040820S.M
Quant Title : SW846 8260

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

Peak Number 2 unknown14.28 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.28	14.89 ug/l	213146	1,4-Dichlorobenzene-d4	13.43

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Butanone, 1-cvclohexyl-	154	C10H18O	001462-27-7	25
2		Cyclohexane, eicosyl-	364	C26H52	004443-55-4	22
3		Decane, 4-methylene-	154	C11H22	024949-41-5	22
4		Decane, 2-cvclohexyl-	224	C16H32	013151-73-0	22
5		Cyclohexane, pentyl-	154	C11H22	004292-92-6	18



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA Y\DATA\VY041520\
Data File : VY002334.D
Acq On : 15 Apr 2020 14:13
Operator : SY/MD
Sample : L2245-08
Misc : 7.15G/5ML/MSVOA Y/SOIL
ALS Vial : 11 Sample Multiplier: 1

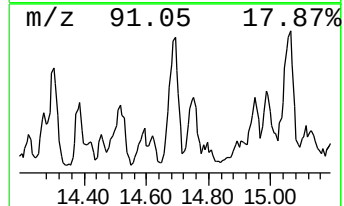
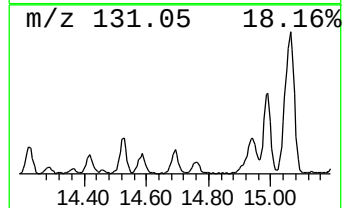
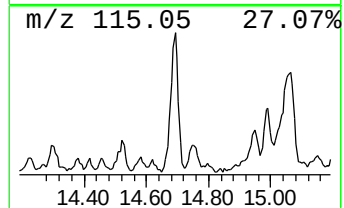
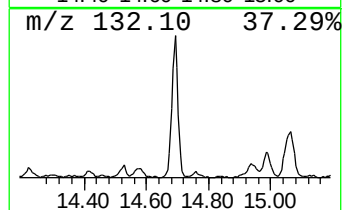
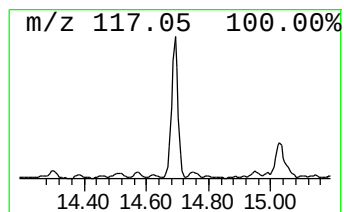
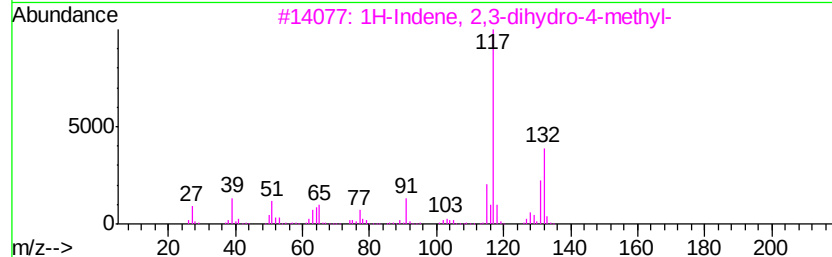
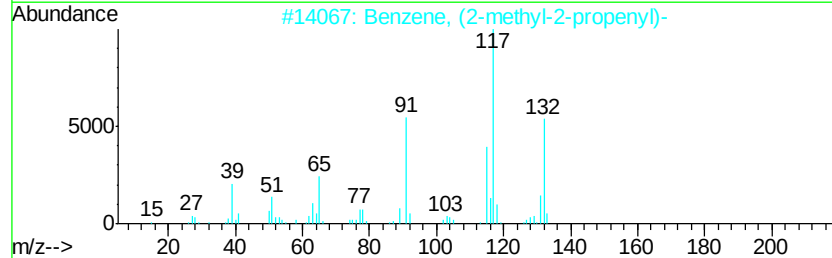
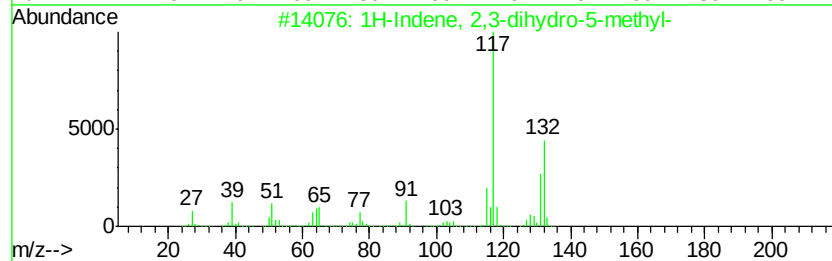
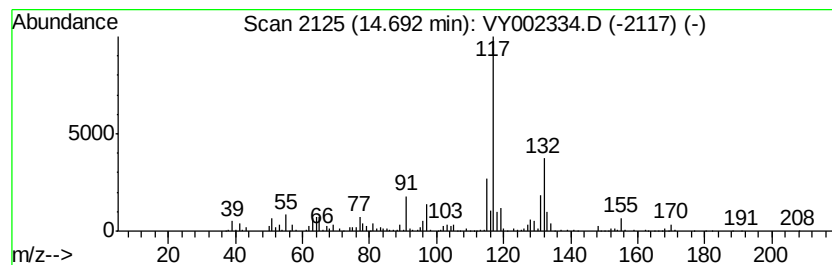
Instrument :
MSVOA_Y
ClientSampleId :
TP-26

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y040820S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.69	25.59 ug/l	366288	1,4-Dichlorobenzene-d4	13.43
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1H-Indene, 2,3-dihydro-5-methyl-	132 C10H12	000874-35-1	93
2	Benzene, (2-methyl-2-propenyl)-	132 C10H12	003290-53-7	87
3	1H-Indene, 2,3-dihydro-4-methyl-	132 C10H12	000824-22-6	87
4	Benzene, 2-butenyl-	132 C10H12	001560-06-1	83
5	Benzene, 1-ethenyl-4-ethyl-	132 C10H12	003454-07-7	81



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA Y\DATA\VY041520\
Data File : VY002334.D
Acq On : 15 Apr 2020 14:13
Operator : SY/MD
Sample : L2245-08
Misc : 7.15G/5ML/MSVOA Y/SOIL
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TP-26

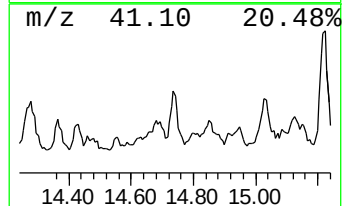
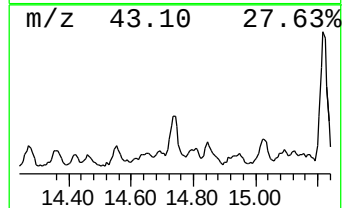
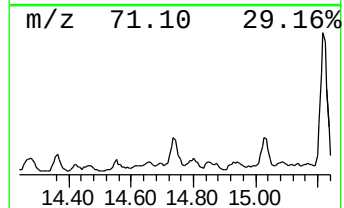
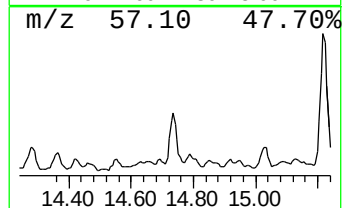
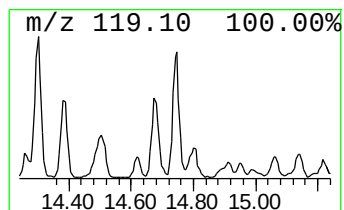
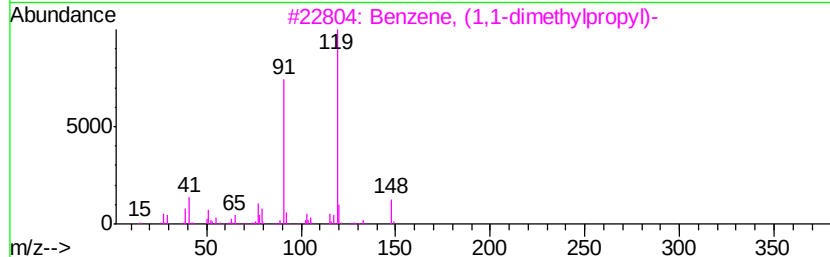
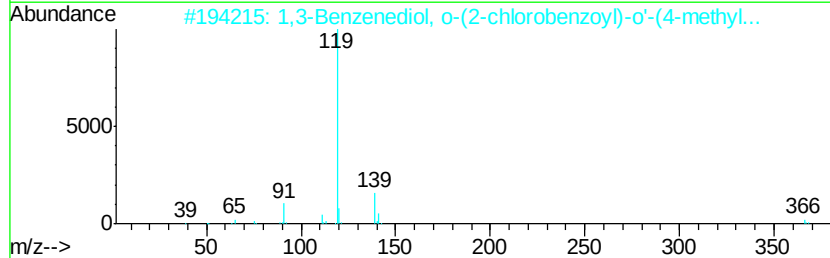
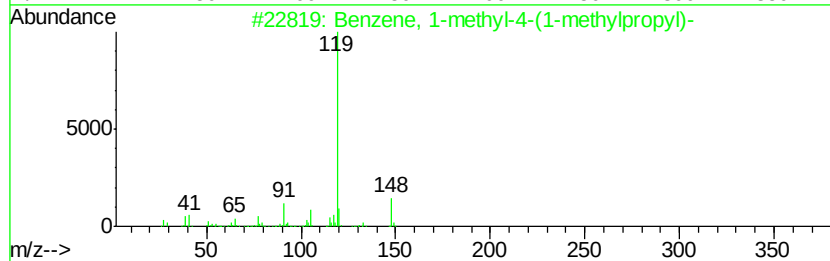
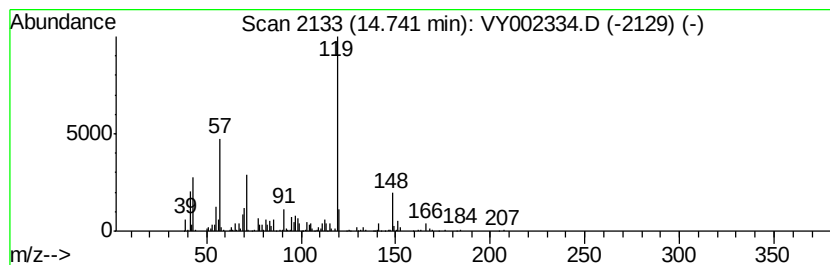
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y040820S.M
Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1-methyl-4-(1-meth... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.74	15.24 ug/l	218075	1,4-Dichlorobenzene-d4	13.43

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	52
2		1,3-Benzenediol, o-(2-chlorobenz...	366	C21H15ClO4	1000330-74-9	50
3		Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	50
4		1,2-Benzenediol, o-(2-chlorobenz...	366	C21H15ClO4	1000325-98-8	47
5		Aziridine, 2-methyl-2-(2,2,4-tri...	245	C17H27N	1000293-28-5	47



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA Y\DATA\VY041520\
Data File : VY002334.D
Acq On : 15 Apr 2020 14:13
Operator : SY/MD
Sample : L2245-08
Misc : 7.15G/5ML/MSVOA Y/SOIL
ALS Vial : 11 Sample Multiplier: 1

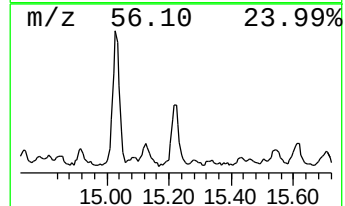
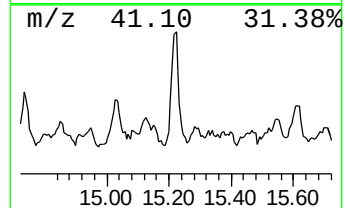
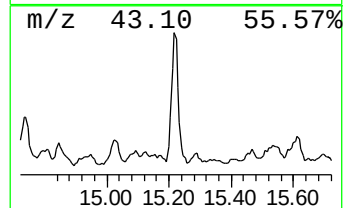
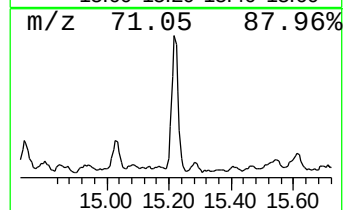
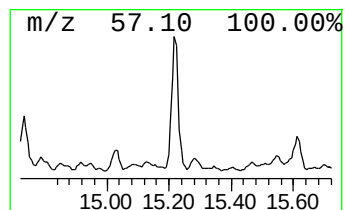
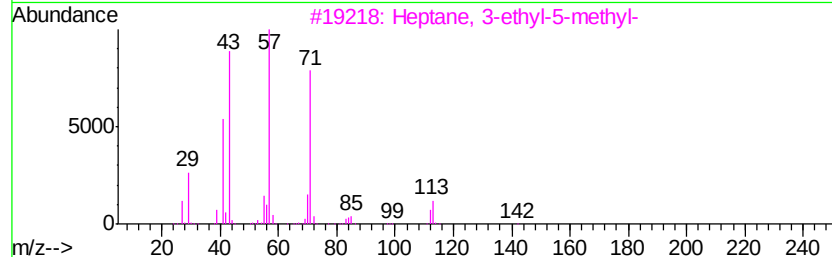
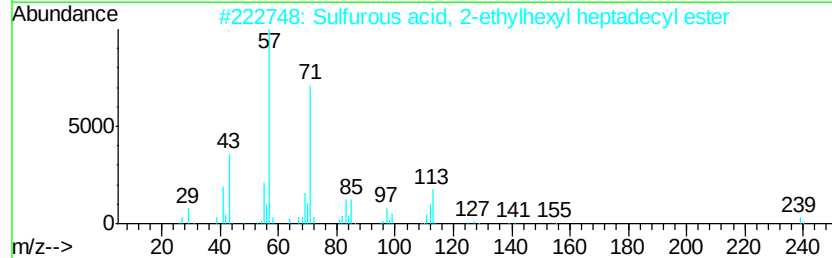
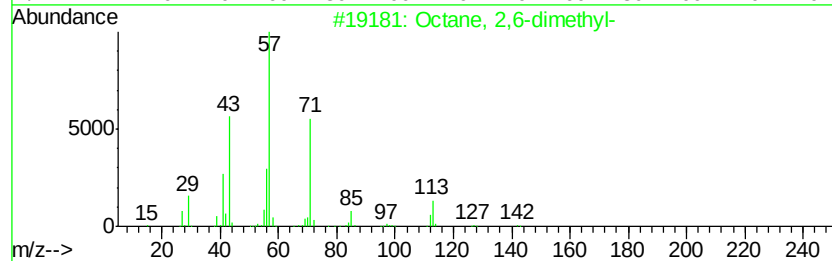
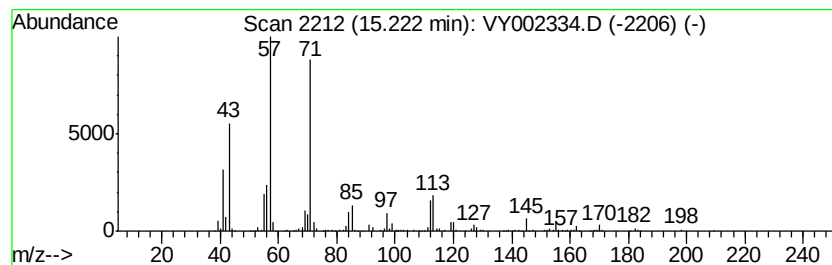
Instrument :
MSVOA_Y
ClientSampleId :
TP-26

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y040820S.M
Quant Title : SW846 8260

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

Peak Number 5 Octane, 2,6-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.22	23.80 ug/l	340607	1,4-Dichlorobenzene-d4	13.43
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Octane, 2,6-dimethyl-	142 C10H22	002051-30-1	80
2	Sulfurous acid, 2-ethylhexyl hep...	432 C25H52O3S	1000309-20-0	72
3	Heptane, 3-ethyl-5-methyl-	142 C10H22	052896-90-9	64
4	Sulfurous acid, dodecyl 2-ethylh...	362 C20H42O3S	1000309-19-5	59
5	Undecane, 5-methyl-	170 C12H26	001632-70-8	58



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA Y\DATA\VY041520\
Data File : VY002334.D
Acq On : 15 Apr 2020 14:13
Operator : SY/MD
Sample : L2245-08
Misc : 7.15G/5ML/MSVOA Y/SOIL
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TP-26

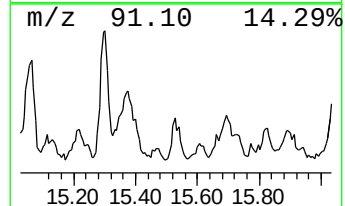
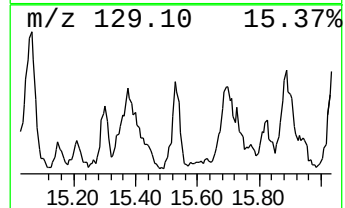
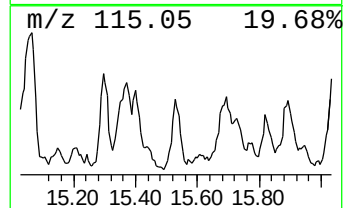
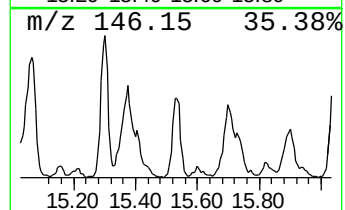
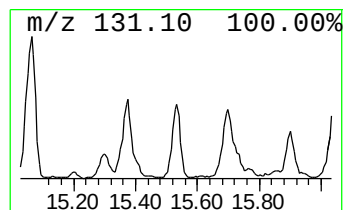
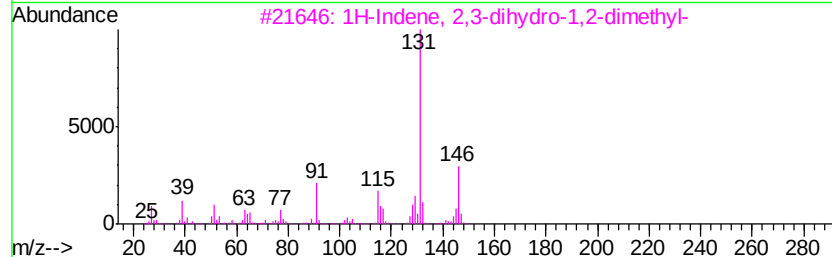
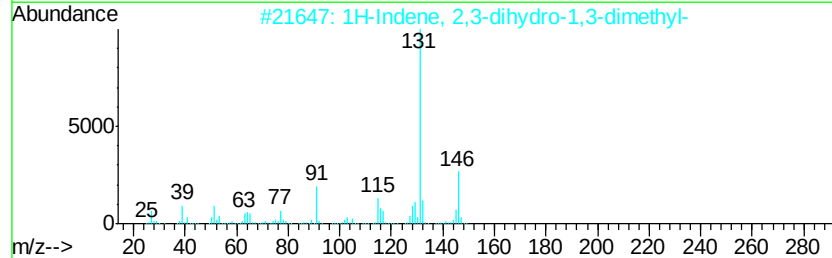
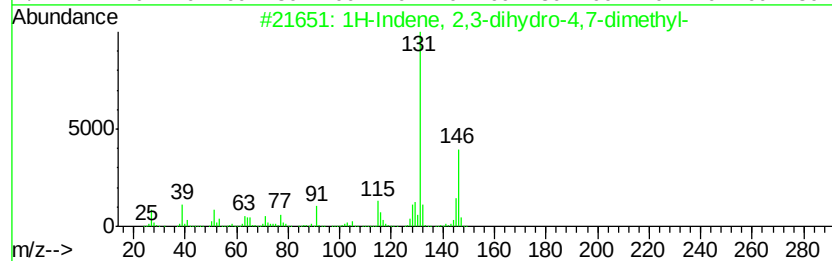
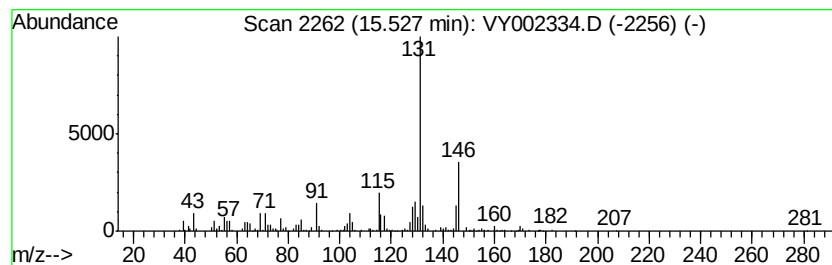
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y040820S.M
Quant Title : SW846 8260

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

Peak Number 6 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.53	16.51 ug/l	236385	1,4-Dichlorobenzene-d4	13.43

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-1,3-dimet...	146	C11H14		004175-53-5	91
3	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14		017057-82-8	91
4	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14		017059-48-2	91
5	Bicyclo[4.2.0]octa-1,3,5-triene,...	146	C11H14		027087-54-3	90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA Y\DATA\VY041520\
Data File : VY002334.D
Acq On : 15 Apr 2020 14:13
Operator : SY/MD
Sample : L2245-08
Misc : 7.15G/5ML/MSVOA Y/SOIL
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TP-26

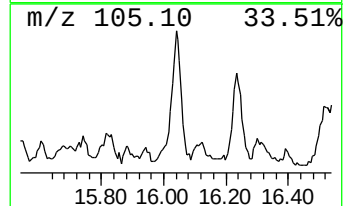
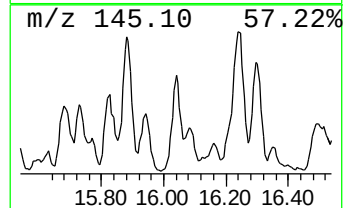
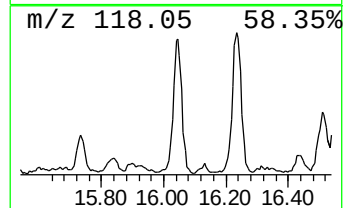
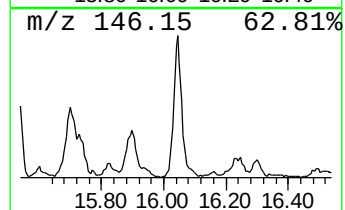
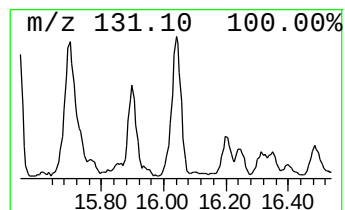
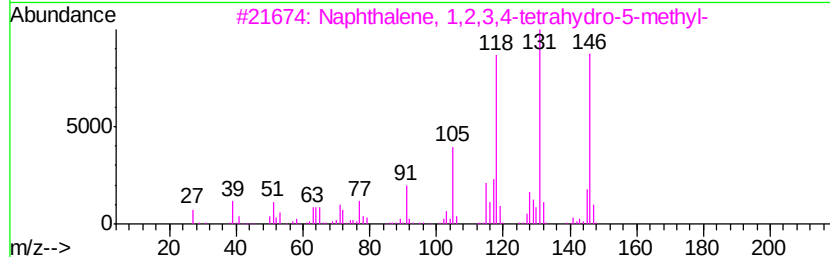
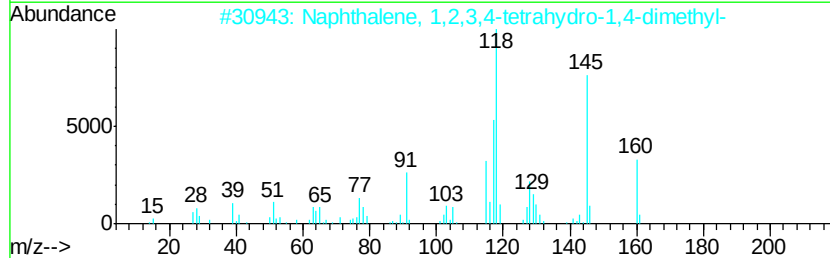
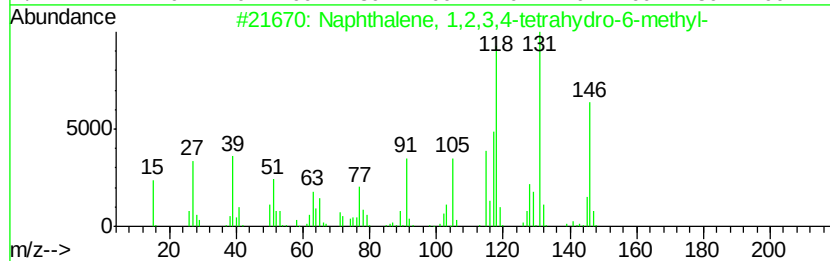
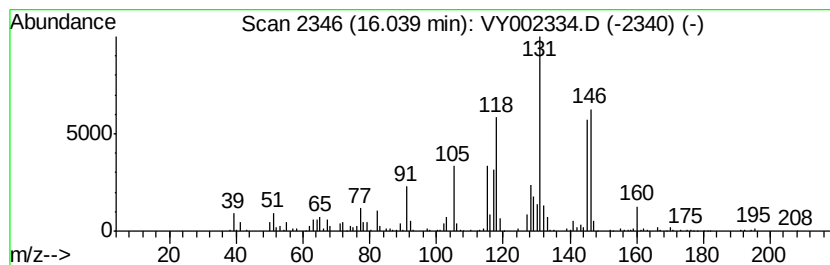
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y040820S.M
Quant Title : SW846 8260

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

Peak Number 8 Naphthalene, 1,2,3,4-tetra... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.04	18.43 ug/l	263835	1,4-Dichlorobenzene-d4	13.43

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	94
2		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	87
3		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	76
4		1H-Benzimidazole, 5,6-dimethyl-	146	C9H10N2	000582-60-5	64
5		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	53



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA Y\DATA\VY041520\
Data File : VY002334.D
Acq On : 15 Apr 2020 14:13
Operator : SY/MD
Sample : L2245-08
Misc : 7.15G/5ML/MSVOA Y/SOIL
ALS Vial : 11 Sample Multiplier: 1

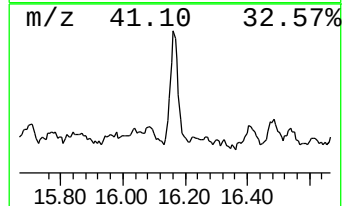
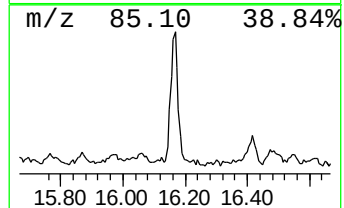
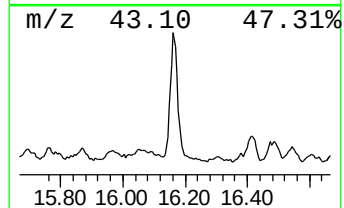
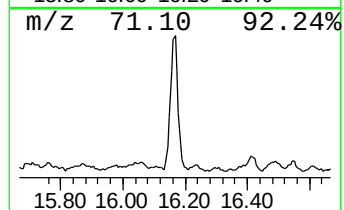
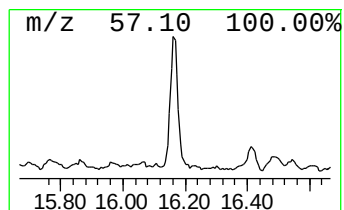
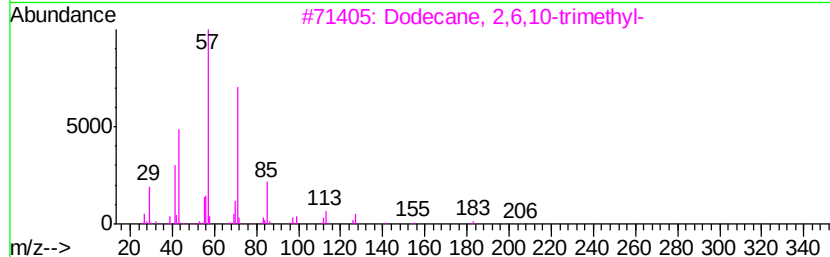
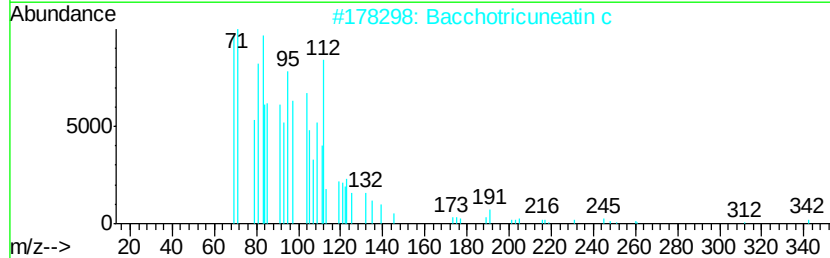
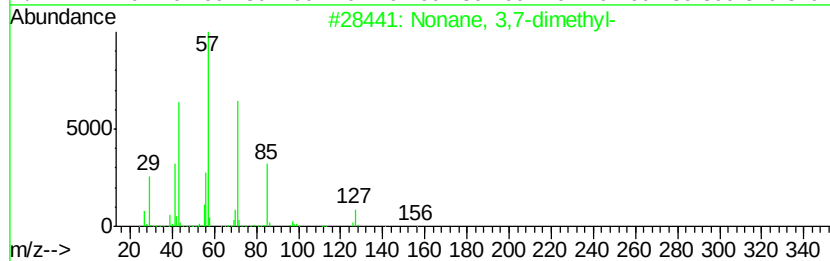
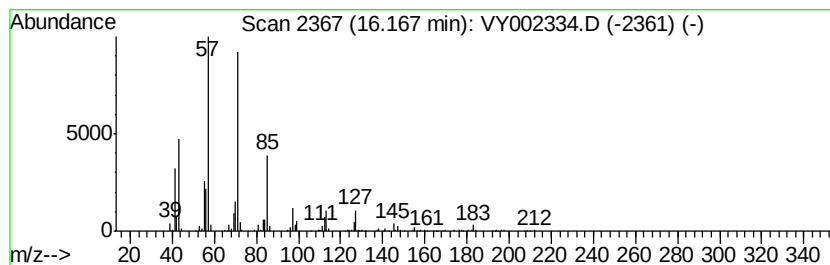
Instrument :
MSVOA_Y
ClientSampleId :
TP-26

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y040820S.M
Quant Title : SW846 8260

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

Peak Number 9 Nonane, 3,7-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.17	20.60 ug/l	294921	1,4-Dichlorobenzene-d4	13.43
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Nonane, 3,7-dimethyl-	156 C11H24	017302-32-8	83
2	Bacchotricuneatin c	342 C20H22O5	066563-30-2	81
3	Dodecane, 2,6,10-trimethyl-	212 C15H32	003891-98-3	72
4	Pentadecane, 2,6,10,14-tetramethyl-	268 C19H40	001921-70-6	72
5	Dodecane, 2,7,10-trimethyl-	212 C15H32	074645-98-0	72



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA Y\DATA\VY041520\
Data File : VY002334.D
Acq On : 15 Apr 2020 14:13
Operator : SY/MD
Sample : L2245-08
Misc : 7.15G/5ML/MSVOA Y/SOIL
ALS Vial : 11 Sample Multiplier: 1

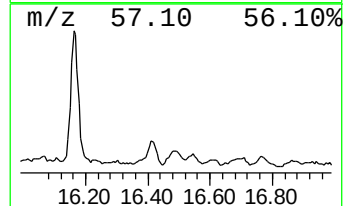
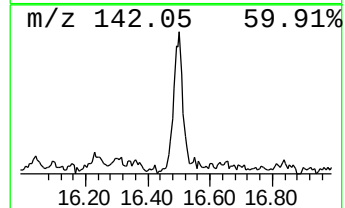
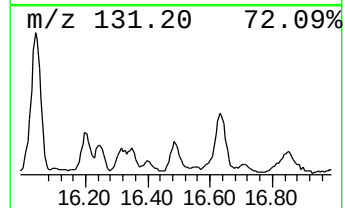
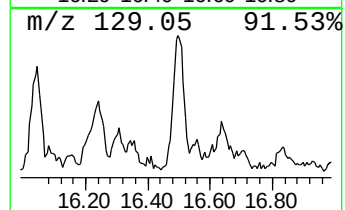
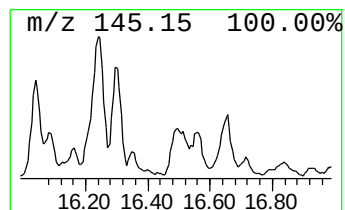
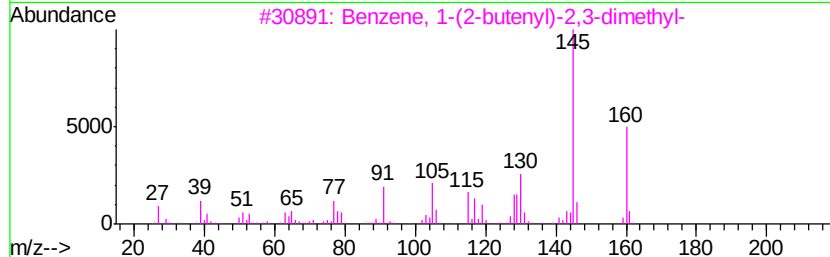
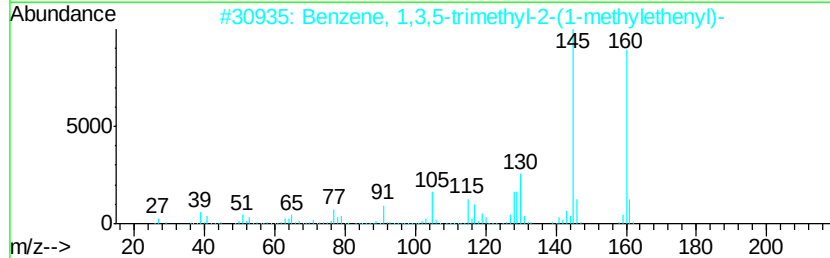
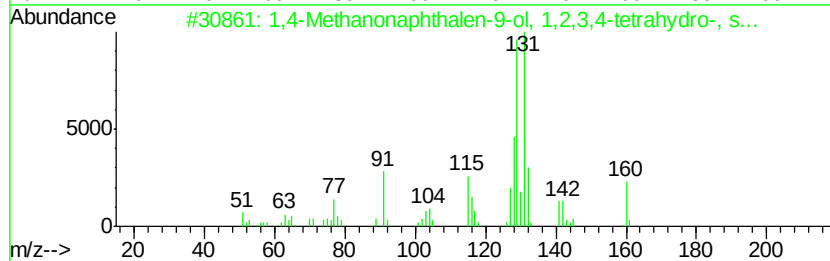
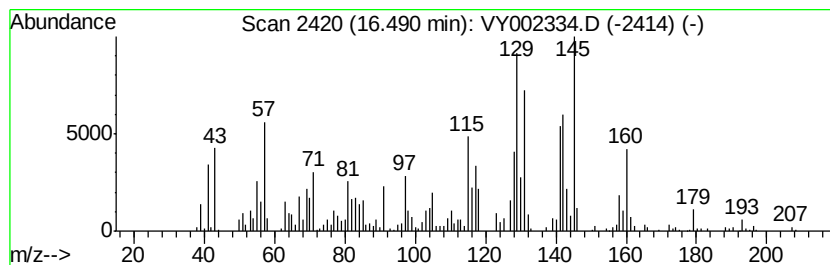
Instrument :
MSVOA_Y
ClientSampled :
TP-26

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y040820S.M
Quant Title : SW846 8260

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

Peak Number 10 unknown16.49 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.49	16.09 ug/l	230354	1,4-Dichlorobenzene-d4	13.43	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,4-Methanonaphthalen-9-ol, 1,2,...	160	C11H12O	001198-20-5	46
2	Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	38
3	Benzene, 1-(2-butenyl)-2,3-dimet...	160	C12H16	054340-85-1	30
4	2,7-Methanonaphth[2,3-b]oxirene,...	158	C11H10O	013137-34-3	30
5	1H-Pyrrolo[2,3-b]pyridine, 2-(1-...	160	C10H12N2	027257-18-7	30



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_Y\DATA\VY041520\
Data File : VY002334.D
Acq On : 15 Apr 2020 14:13
Operator : SY/MD
Sample : L2245-08
Misc : 7.15G/5ML/MSVOA_Y/SOIL
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TP-26

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y040820S.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
Indan, 1-methyl-	14.08	14.4	ug/l	205800	4	13.43	715701	50.0
unknown14.28	14.28	14.9	ug/l	213146	4	13.43	715701	50.0
1H-Indene, 2,3-di...	14.69	25.6	ug/l	366288	4	13.43	715701	50.0
Benzene, 1-methyl...	14.74	15.2	ug/l	218075	4	13.43	715701	50.0
Octane, 2,6-dimet...	15.22	23.8	ug/l	340607	4	13.43	715701	50.0
1H-Indene, 2,3-di...	15.53	16.5	ug/l	236385	4	13.43	715701	50.0
Naphthalene, 1,2,...	16.04	18.4	ug/l	263835	4	13.43	715701	50.0
Nonane, 3,7-dimet...	16.17	20.6	ug/l	294921	4	13.43	715701	50.0
unknown16.49	16.49	16.1	ug/l	230354	4	13.43	715701	50.0