

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY011322\
 Data File : VY007149.D
 Acq On : 13 Jan 2022 19:00
 Operator : SY/MD
 Sample : N1119-04
 Misc : 4.85g/5.0mL/MSVOA_Y/SOIL
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 NORTH-EAST-CORNER-EXCAVATION-EAST-SIDEWALL

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\82Y011322S.M
 Title : SW846 8260

Signal : TIC: VY007149.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.058	34	39	46	rBV	2204	3585	5.08%	1.263%
2	7.722	958	968	973	rBV2	4180	9139	12.94%	3.219%
3	7.789	973	979	991	rVB2	6538	16340	23.14%	5.755%
4	8.143	1030	1037	1044	rBV3	3966	8444	11.96%	2.974%
5	8.319	1062	1066	1070	rBV4	739	1396	1.98%	0.492%
6	8.691	1119	1127	1137	rBV	10930	22070	31.26%	7.773%
7	9.612	1273	1278	1281	rVB2	680	1136	1.61%	0.400%
8	9.746	1297	1300	1305	rBV2	627	914	1.29%	0.322%
9	9.971	1333	1337	1343	rVB5	647	1425	2.02%	0.502%
10	10.179	1365	1371	1379	rBV	15015	28190	39.93%	9.929%
11	10.612	1439	1442	1446	rBV3	533	883	1.25%	0.311%
12	10.721	1455	1460	1466	rBV6	2579	4434	6.28%	1.562%
13	11.093	1517	1521	1525	rVB3	949	1468	2.08%	0.517%
14	11.209	1534	1540	1544	rBV3	495	998	1.41%	0.351%
15	11.276	1544	1551	1557	rVV7	966	2322	3.29%	0.818%
16	11.386	1564	1569	1575	rVB4	788	1602	2.27%	0.564%
17	11.489	1580	1586	1593	rVB	14407	23070	32.68%	8.125%
18	11.709	1614	1622	1628	rBV7	815	2444	3.46%	0.861%
19	12.008	1663	1671	1676	rBV6	1515	3399	4.81%	1.197%
20	12.123	1684	1690	1695	rVB3	1235	2164	3.07%	0.762%
21	12.203	1697	1703	1706	rVV5	1220	2440	3.46%	0.859%
22	12.245	1706	1710	1717	rVB7	1578	3286	4.65%	1.157%
23	12.331	1720	1724	1729	rVV2	2816	4875	6.90%	1.717%
24	12.386	1729	1733	1735	rVV4	837	1391	1.97%	0.490%
25	12.416	1735	1738	1740	rVV4	857	1301	1.84%	0.458%
26	12.489	1744	1750	1758	rVB2	41007	70602	100.00%	24.866%
27	12.648	1768	1776	1783	rVB6	1849	4376	6.20%	1.541%
28	12.721	1783	1788	1791	rBV4	782	1560	2.21%	0.549%
29	12.812	1796	1803	1807	rBV7	2201	4727	6.70%	1.665%
30	12.953	1822	1826	1829	rBV5	888	1272	1.80%	0.448%
31	13.001	1831	1834	1837	rBV2	708	833	1.18%	0.293%
32	13.044	1837	1841	1844	rBV5	1347	2009	2.85%	0.708%
33	13.166	1857	1861	1862	rBV4	699	731	1.04%	0.257%
34	13.251	1872	1875	1879	rVV3	857	1303	1.85%	0.459%
35	13.288	1879	1881	1884	rVV4	924	1368	1.94%	0.482%
36	13.325	1885	1887	1890	rVV3	1124	1473	2.09%	0.519%

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37	13.367	1890	1894	1898	rVV6	942	2037	2.89%	0.717%
38	13.428	1899	1904	1912	rVB	12229	18954	26.85%	6.676%
39	13.672	1940	1944	1946	rBV4	900	1410	2.00%	0.497%
40	13.696	1946	1948	1952	rVV4	1250	1522	2.16%	0.536%
41	13.751	1952	1957	1961	rVV6	2127	4325	6.13%	1.523%
42	13.794	1961	1964	1966	rVV3	663	909	1.29%	0.320%
43	13.831	1966	1970	1972	rVV4	582	980	1.39%	0.345%
44	13.898	1979	1981	1986	rVV5	775	847	1.20%	0.298%
45	13.971	1986	1993	1998	rVV3	1729	2895	4.10%	1.020%
46	14.087	2009	2012	2016	rVB5	570	813	1.15%	0.286%
47	14.276	2035	2043	2045	rVV6	1108	2179	3.09%	0.767%
48	14.379	2055	2060	2063	rBV5	486	937	1.33%	0.330%
49	14.422	2063	2067	2071	rVV4	1175	1764	2.50%	0.621%
50	14.489	2074	2078	2082	rVB3	1083	1547	2.19%	0.545%
51	14.678	2103	2109	2114	rBV7	1160	2062	2.92%	0.726%
52	16.056	2331	2335	2339	rBV5	572	995	1.41%	0.350%
53	16.635	2427	2430	2432	rBV3	645	781	1.11%	0.275%

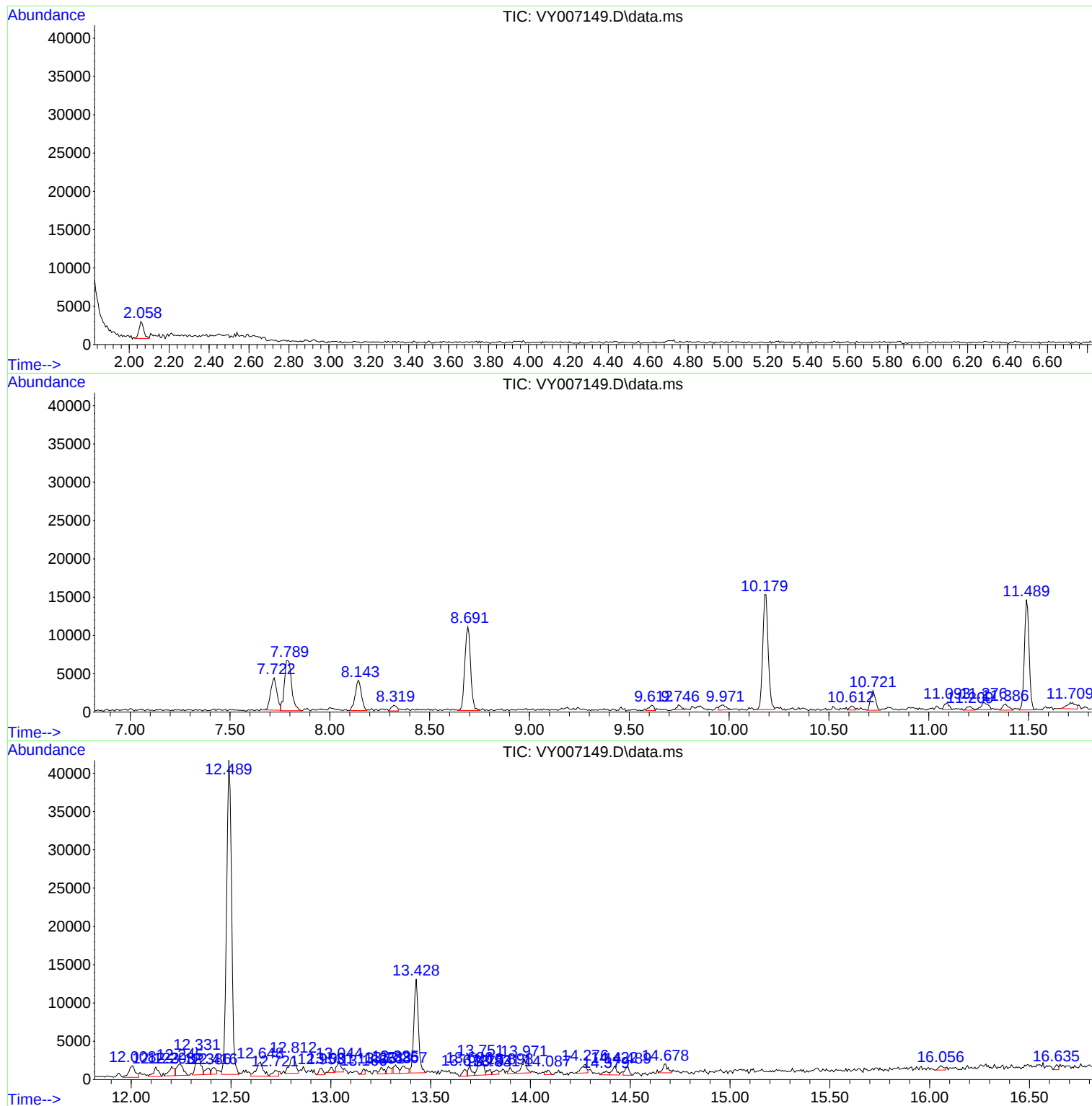
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y011322S.M
Quant Title : SW846 8260

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



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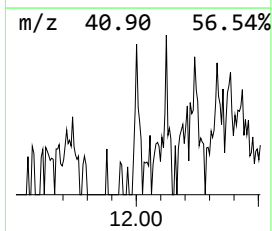
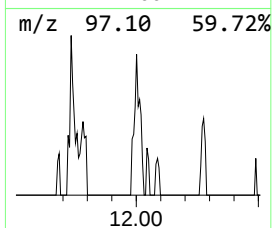
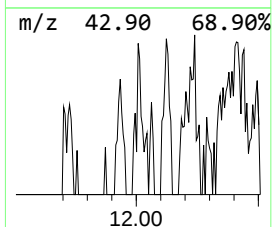
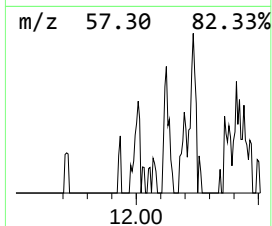
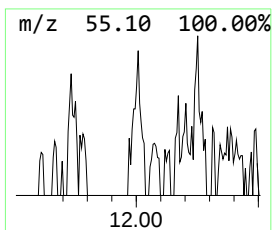
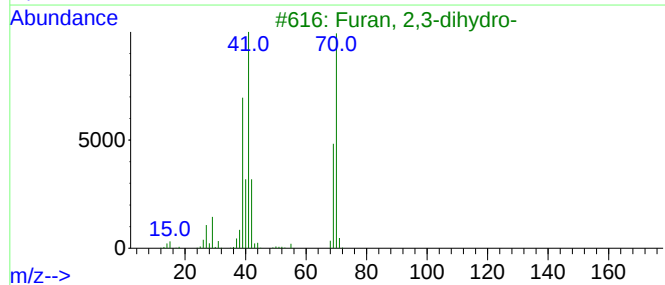
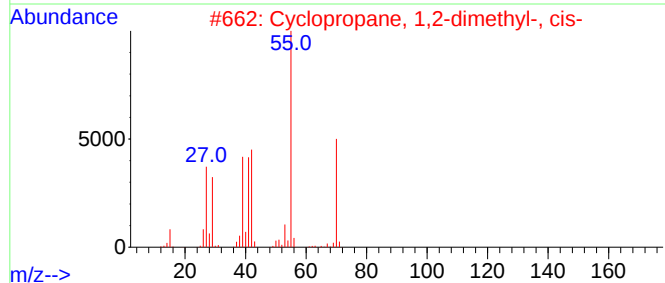
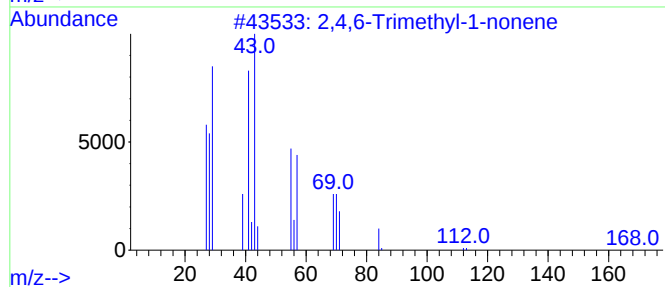
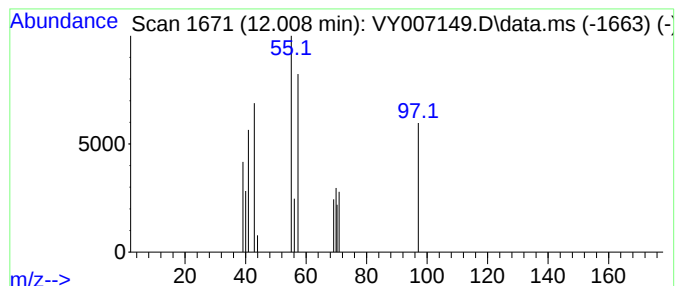
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y011322S.M
Quant Title : SW846 8260

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

Peak Number 2 unknown12.008 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.008	7.37 ug/l	3399	Chlorobenzene-d5	11.496

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,4,6-Trimethyl-1-nonene	168	C12H24	055771-40-9	45
2			Cyclopropane, 1,2-dimethyl-, cis-	70	C5H10	000930-18-7	32
3			Furan, 2,3-dihydro-	70	C4H6O	001191-99-7	27
4			(E)-1,3-Butadien-1-ol	70	C4H6O	070411-98-2	25
5			1,6-Hexanediol	118	C6H14O2	000629-11-8	25



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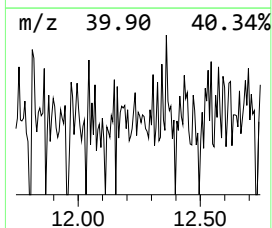
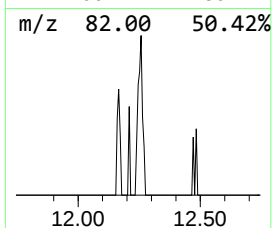
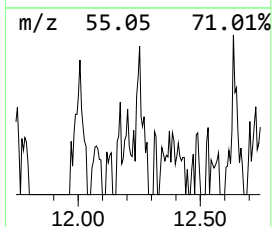
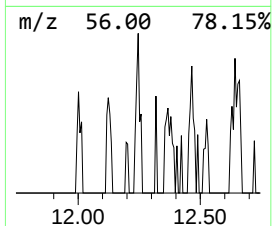
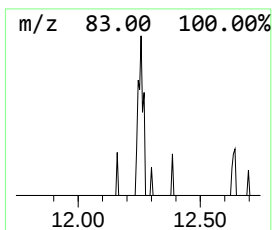
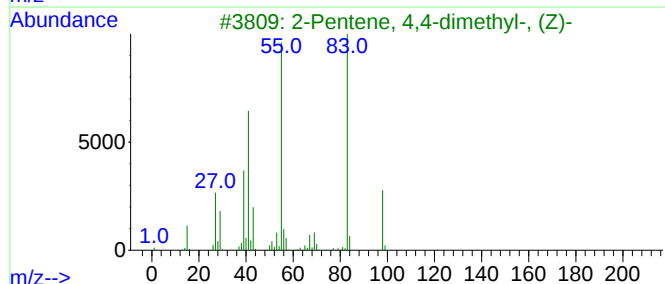
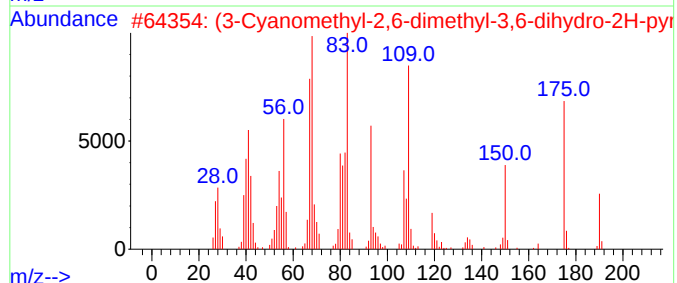
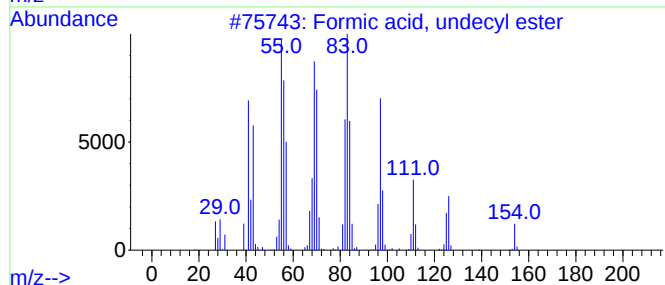
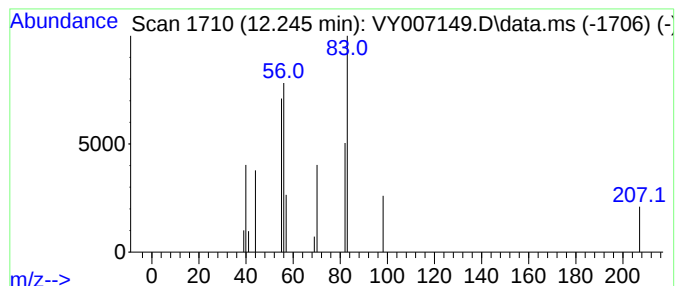
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y011322S.M
Quant Title : SW846 8260

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

Peak Number 3 unknown12.245 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.245	7.12 ug/l	3286	Chlorobenzene-d5	11.496

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Formic acid, undecyl ester	200	C12H24O2	1000368-25-0	36
2			(3-Cyanomethyl-2,6-dimethyl-3,6-...	190	C10H14N4	1000187-56-2	33
3			2-Pentene, 4,4-dimethyl-, (Z)-	98	C7H14	000762-63-0	12
4			2-Pentene, 4,4-dimethyl-, (E)-	98	C7H14	000690-08-4	12
5			1-Butene, 2,3,3-trimethyl-	98	C7H14	000594-56-9	9



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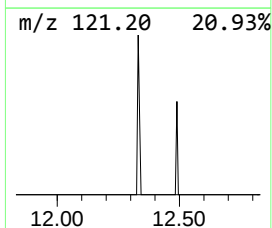
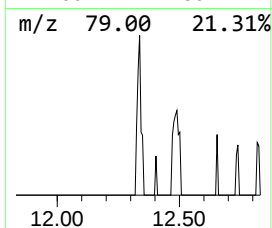
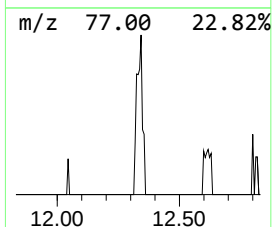
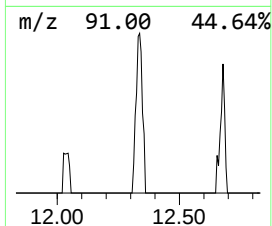
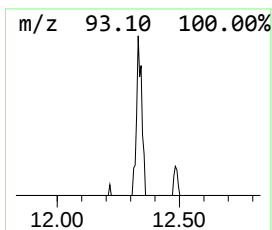
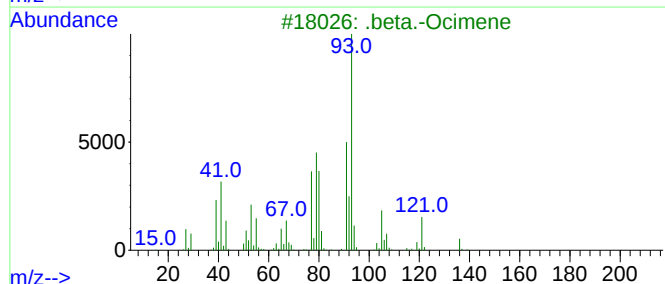
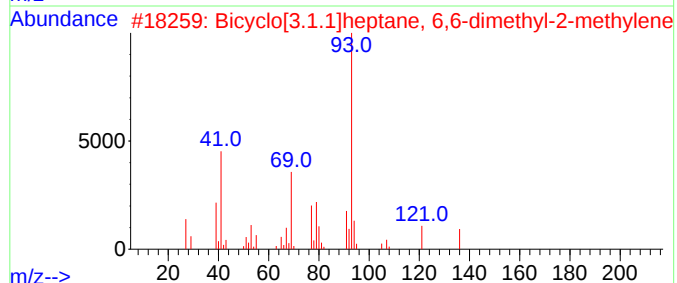
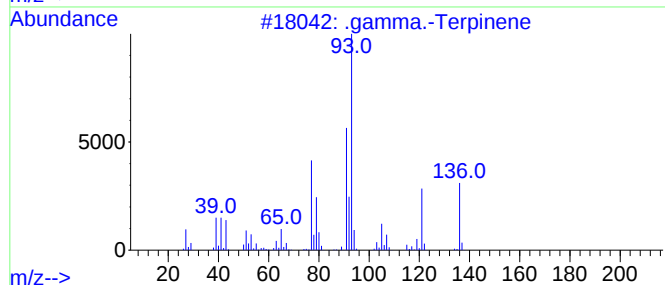
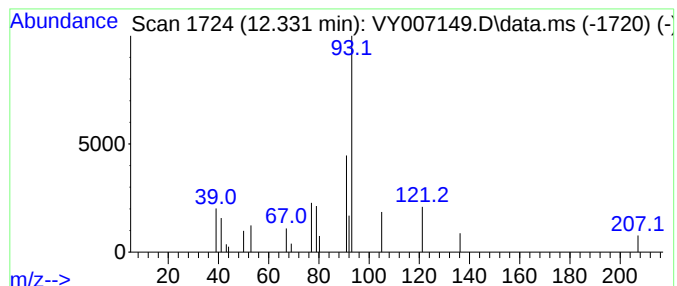
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 .gamma.-Terpinene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.331	10.57 ug/l	4875	Chlorobenzene-d5	11.496

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			.gamma.-Terpinene	136	C10H16	000099-85-4	86
2			Bicyclo[3.1.1]heptane, 6,6-dimet...	136	C10H16	018172-67-3	86
3			.beta.-Ocimene	136	C10H16	013877-91-3	83
4			Tricyclo[2.2.1.0(2,6)]heptane, 1...	136	C10H16	000508-32-7	80
5			.beta.-Pinene	136	C10H16	000127-91-3	78



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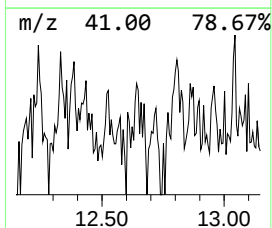
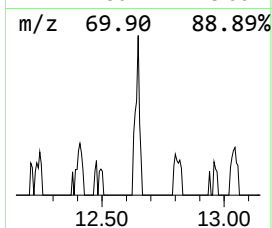
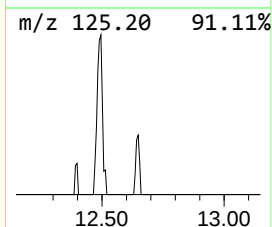
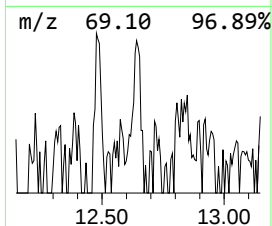
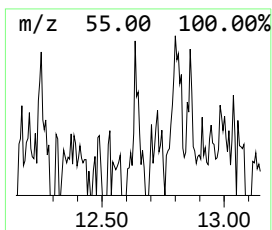
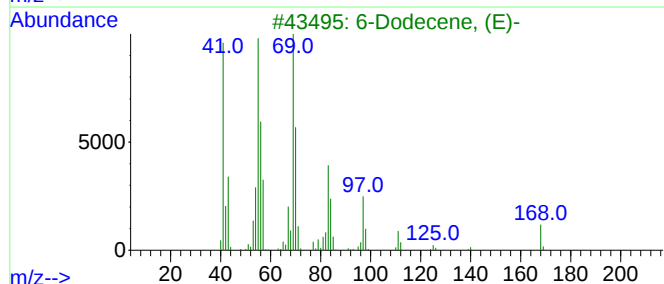
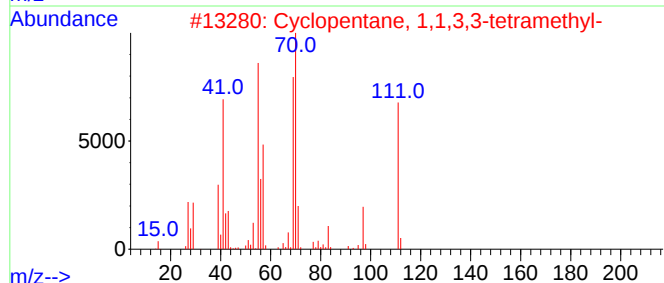
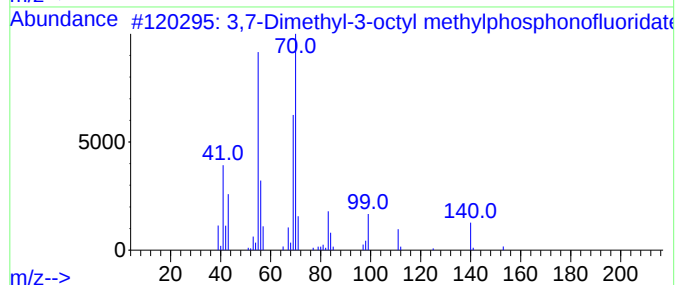
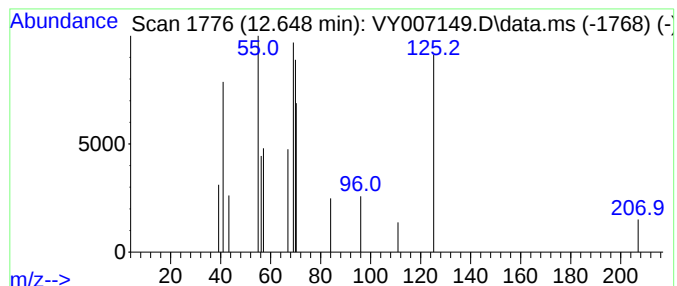
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TIC Library : C:\Database\NIST20.L
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Peak Number 5 unknown12.648 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.648	11.54 ug/l	4376	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3,7-Dimethyl-3-octyl methylphosp...	238	C11H24FO2P	159395-79-6	38
2			Cyclopentane, 1,1,3,3-tetramethyl-	126	C9H18	050876-33-0	37
3			6-Dodecene, (E)-	168	C12H24	007206-17-9	37
4			4-Heptafluorobutyryloxyhexadecane	438	C20H33F7O2	1010282-97-2	37
5			3-Cyclopentylpropionic acid, 2-e...	254	C16H30O2	1000293-47-0	33



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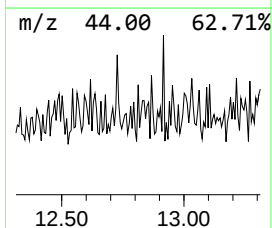
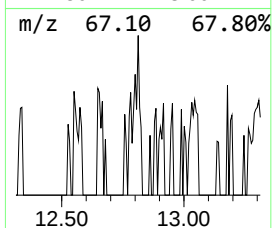
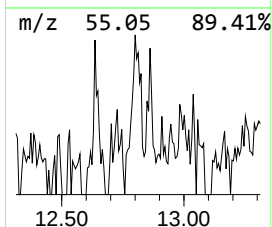
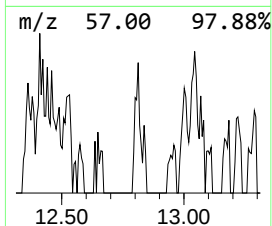
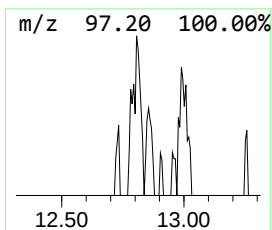
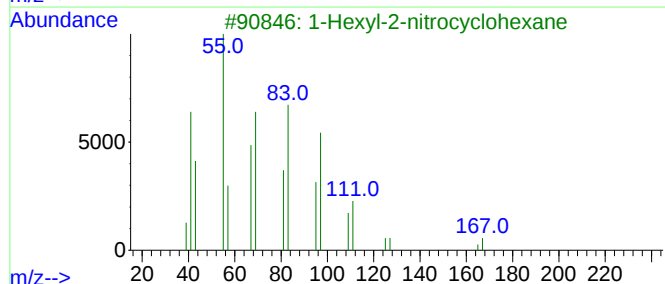
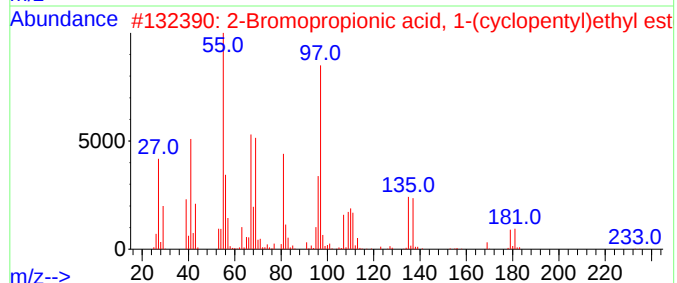
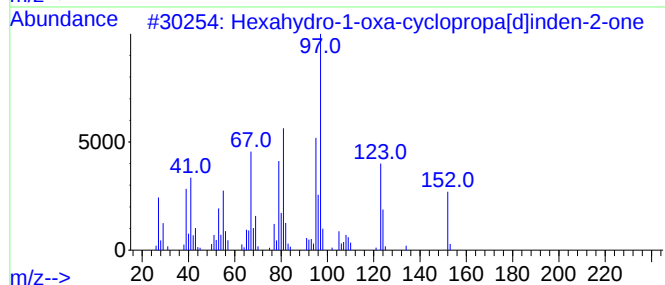
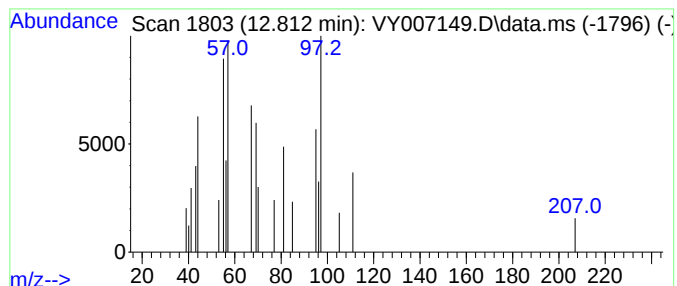
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Quant Title : SW846 8260

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

Peak Number 6 unknown12.812 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.812	12.47 ug/l	4727	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexahydro-1-oxa-cyclopropa[d]ind...	152	C9H12O2	037643-23-5	43
2			2-Bromopropionic acid, 1-(cyclop...	248	C10H17BrO2	1000293-39-1	38
3			1-Hexyl-2-nitrocyclohexane	213	C12H23NO2	118252-04-3	37
4			n-Tetracosanol-1	354	C24H50O	000506-51-4	35
5			6,10,13-Trimethyltetradecanol	256	C17H36O	1000131-71-0	32



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY011322\
Data File : VY007149.D
Acq On : 13 Jan 2022 19:00
Operator : SY/MD
Sample : N1119-04
Misc : 4.85g/5.0mL/MSVOA_Y/SoIL
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
NORTH-EAST-CORNER-EXCAVATION-EAST-SIDEWALL

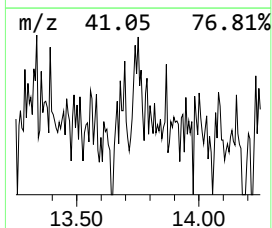
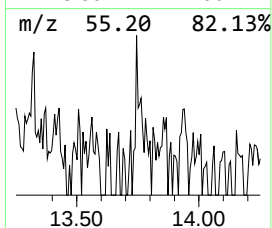
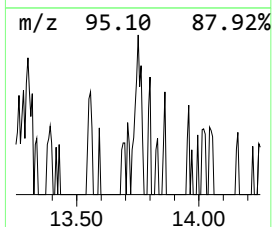
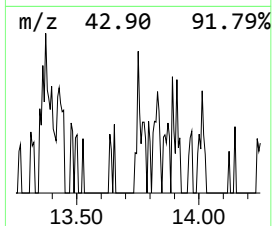
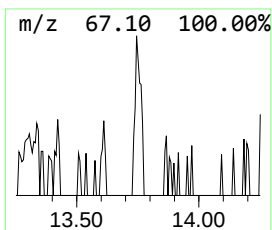
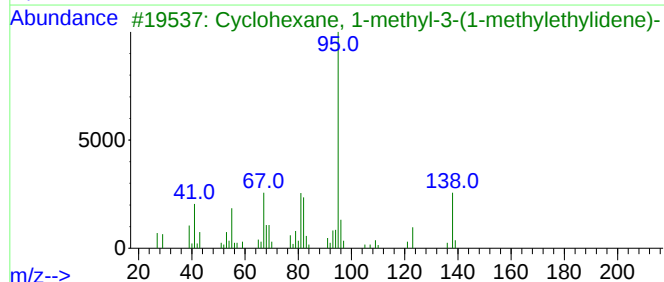
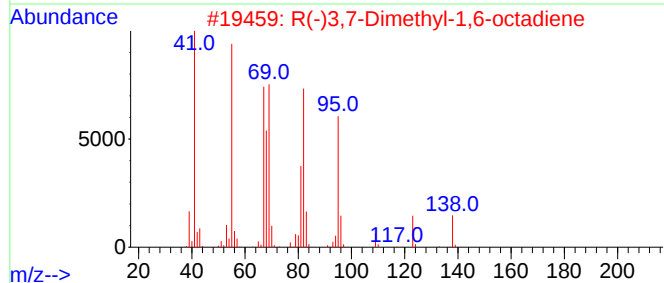
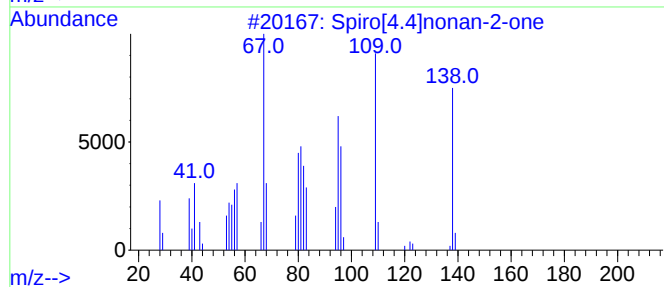
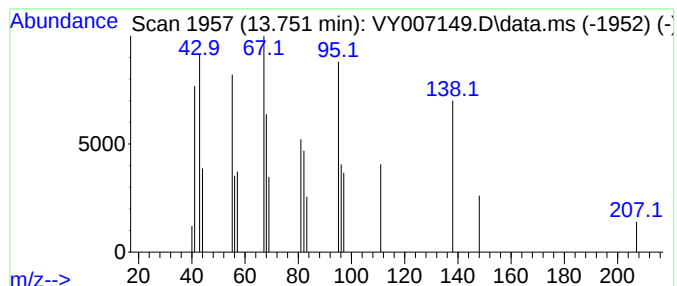
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Quant Title : SW846 8260

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

Peak Number 7 unknown13.751 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.751	11.41 ug/l	4325	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Spiro[4.4]nonan-2-one	138	C9H14O	034177-18-9	49
2			R(-)3,7-Dimethyl-1,6-octadiene	138	C10H18	010281-56-8	47
3			Cyclohexane, 1-methyl-3-(1-methy...	138	C10H18	013828-34-7	47
4			trans-2-Dodecen-1-ol, acetate	226	C14H26O2	084801-16-1	47
5			Naphthalene, decahydro-	138	C10H18	000091-17-8	46



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY011322\
Data File : VY007149.D
Acq On : 13 Jan 2022 19:00
Operator : SY/MD
Sample : N1119-04
Misc : 4.85g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
NORTH-EAST-CORNER-EXCAVATION-EAST-SIDEWALL

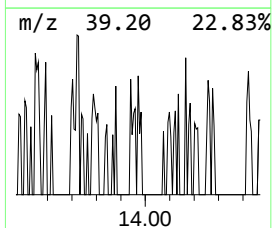
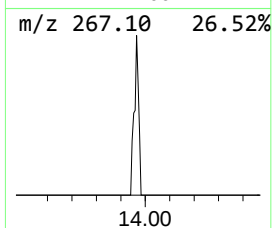
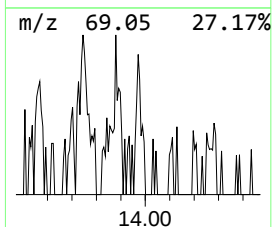
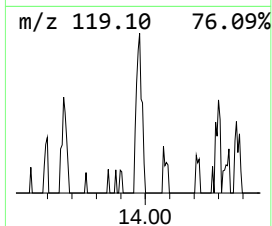
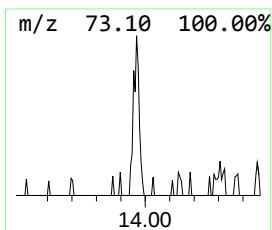
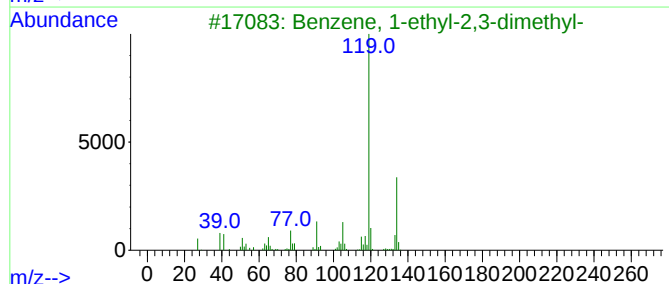
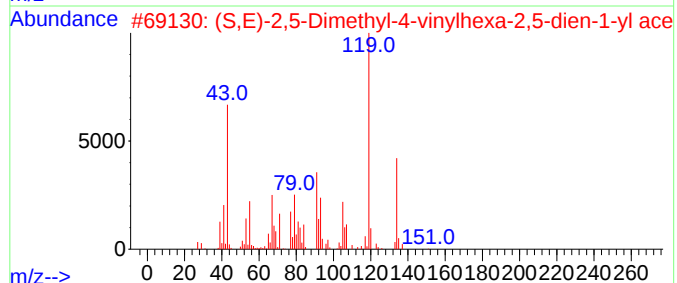
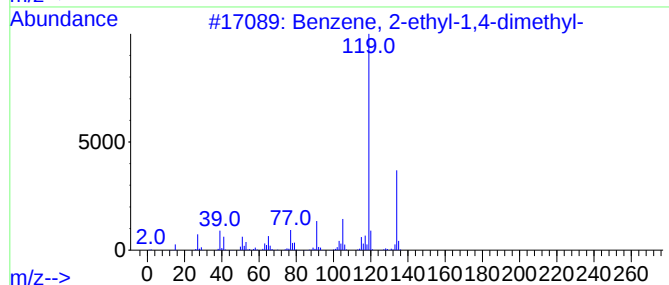
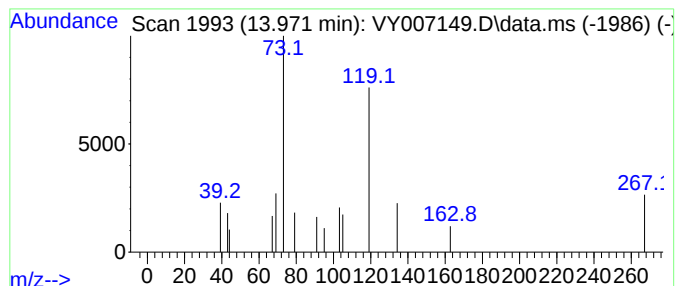
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Quant Title : SW846 8260

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

Peak Number 8 unknown13.971 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.971	7.64 ug/l	2895	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	38
2			(S,E)-2,5-Dimethyl-4-vinylhexa-2...	194	C12H18O2	020384-05-8	37
3			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	37
4			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	9
5			p-Cymene	134	C10H14	000099-87-6	9



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY011322\
Data File : VY007149.D
Acq On : 13 Jan 2022 19:00
Operator : SY/MD
Sample : N1119-04
Misc : 4.85g/5.0mL/MSVOA_Y/SoIL
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
NORTH-EAST-CORNER-EXCAVATION-EAST-SIDEWALL

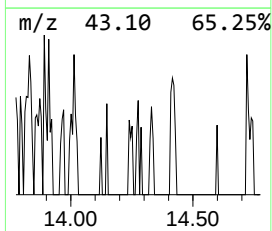
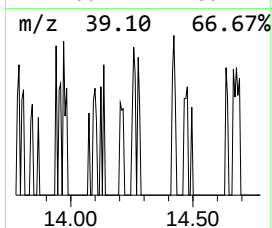
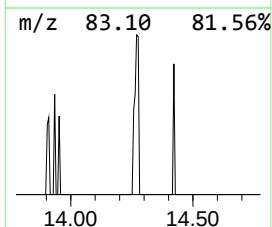
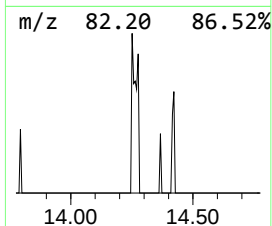
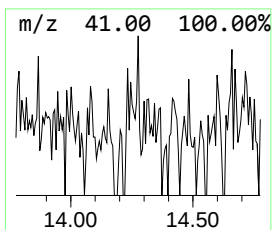
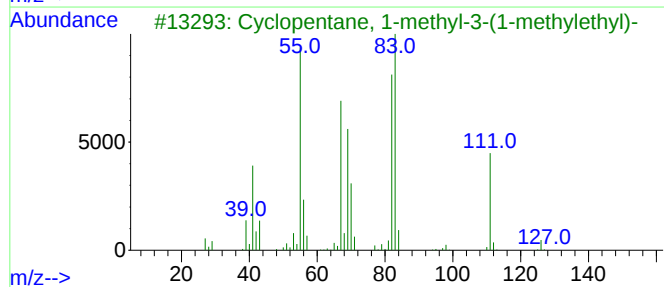
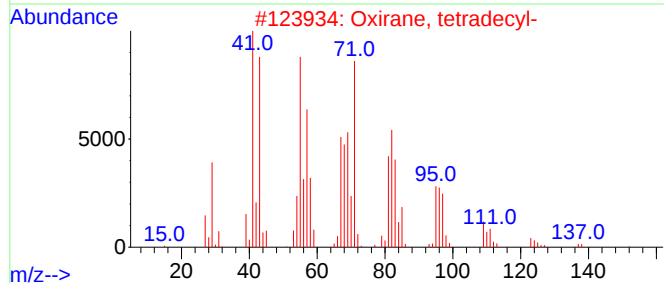
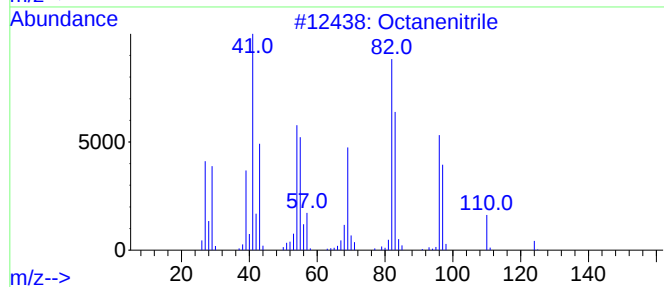
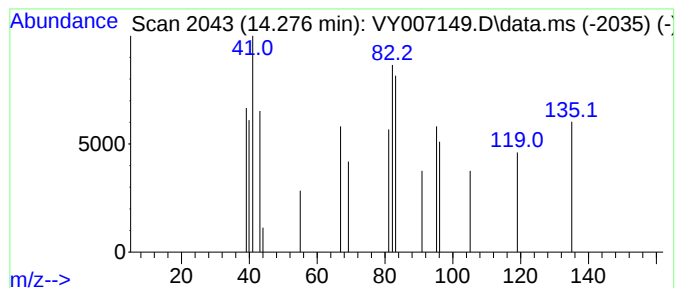
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Quant Title : SW846 8260

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

Peak Number 9 unknown14.276 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.276	5.75 ug/l	2179	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octanenitrile	125	C8H15N	000124-12-9	25
2			Oxirane, tetradecyl-	240	C16H32O	007320-37-8	25
3			Cyclopentane, 1-methyl-3-(1-meth...	126	C9H18	053771-88-3	17
4			2-Propylcyclohexanol	142	C9H18O	090676-25-8	17
5			13-Heptadecyn-1-ol	252	C17H32O	056554-77-9	17



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY011322\
Data File : VY007149.D
Acq On : 13 Jan 2022 19:00
Operator : SY/MD
Sample : N1119-04
Misc : 4.85g/5.0mL/MSVOA_Y/soil
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
NORTH-EAST-CORNER-EXCAVATION-EAST-SIDEWALL

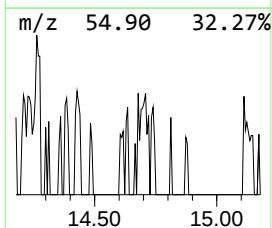
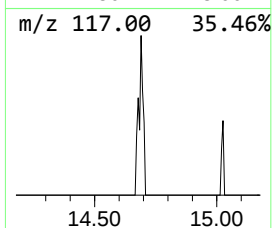
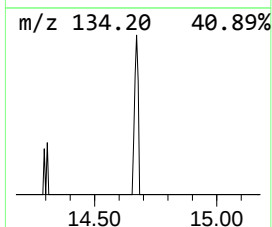
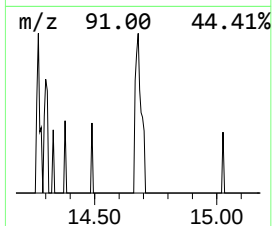
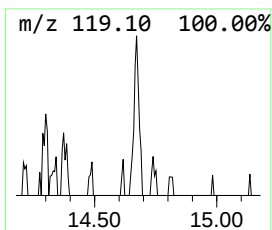
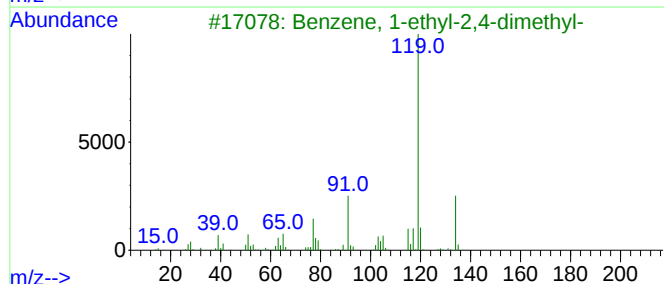
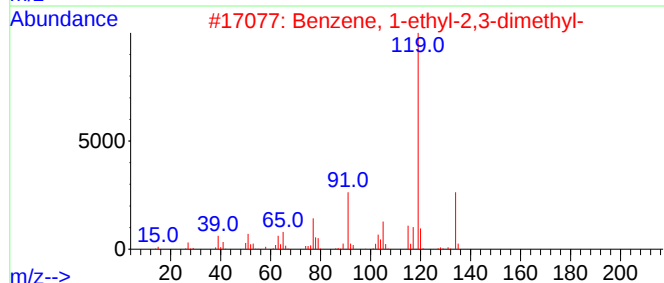
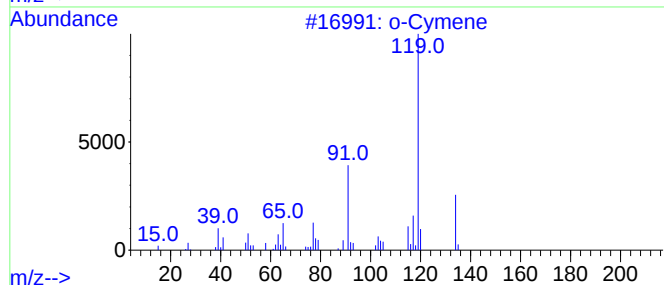
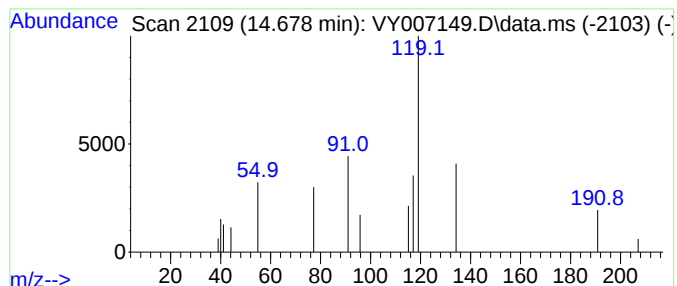
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Quant Title : SW846 8260

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

Peak Number 10 unknown14.678 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.678	5.44 ug/l	2062	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Cymene	134	C10H14	000527-84-4	17
2			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	9
3			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	9
4			1,3,5-Cycloheptatriene, 3,7,7-tr...	134	C10H14	003479-89-8	9
5			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	9



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY011322\
Data File : VY007149.D
Acq On : 13 Jan 2022 19:00
Operator : SY/MD
Sample : N1119-04
Misc : 4.85g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
NORTH-EAST-CORNER-EXCAVATION-EAST-SIDEWALL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y011322S.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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unknown12.245	12.245	7.1	ug/l	3286	3	11.496	23070	50.0
.gamma.-Terpinene	12.331	10.6	ug/l	4875	3	11.496	23070	50.0
unknown12.648	12.648	11.5	ug/l	4376	4	13.428	18954	50.0
unknown12.812	12.812	12.5	ug/l	4727	4	13.428	18954	50.0
unknown13.751	13.751	11.4	ug/l	4325	4	13.428	18954	50.0
unknown13.971	13.971	7.6	ug/l	2895	4	13.428	18954	50.0
unknown14.276	14.276	5.8	ug/l	2179	4	13.428	18954	50.0
unknown14.678	14.678	5.4	ug/l	2062	4	13.428	18954	50.0