LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY102921\
Data File : VY006533.D
Acq On : 29 Oct 2021 12:33
Operator : SY/MD
Sample : M4407-02
Misc : 2.42G/5ML/MSVOA_Y/SOIL
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
FDR-BH-7-20211028-7.5-8

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\82Y102221S.M
Title : SW846 8260

Signal : TIC: VY006533.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.869	4	8	16	rVB2	13188	24577	1.39%	0.105%
2	3.954	336	350	367	rBV	105223	291799	16.47%	1.242%
3	4.204	379	391	403	rBV2	32119	95594	5.40%	0.407%
4	4.814	479	491	502	rBV	37955	128772	7.27%	0.548%
5	4.948	502	513	530	rVB4	39487	158572	8.95%	0.675%
6	6.990	836	848	864	rBV2	23459	90203	5.09%	0.384%
7	7.722	958	968	974	rBV	84372	203929	11.51%	0.868%
8	7.795	974	980	989	rVB2	155790	349832	19.75%	1.489%
9	8.149	1028	1038	1049	rBV2	113192	265928	15.01%	1.132%
10	8.691	1119	1127	1141	rVB	259470	514940	29.07%	2.192%
11	8.892	1148	1160	1169	rBV2	15472	40114	2.26%	0.171%
12	10.179	1363	1371	1378	rBV	375855	661566	37.35%	2.816%
13	10.246	1378	1382	1388	rVB	38395	65274	3.69%	0.278%
14	10.368	1395	1402	1408	rVB6	9091	24857	1.40%	0.106%
15	10.605	1431	1441	1450	rVB4	23464	68892	3.89%	0.293%
16	10.880	1478	1486	1490	rBV	39194	73830	4.17%	0.314%
17	10.965	1494	1500	1503	rBV6	12188	20516	1.16%	0.087%
18	11.038	1509	1512	1516	rVV2	17532	23248	1.31%	0.099%
19	11.093	1516	1521	1525	rVV2	34222	56871	3.21%	0.242%
20	11.148	1525	1530	1535	rVV3	24256	45053	2.54%	0.192%
21	11.203	1535	1539	1544	rVB3	15460	28475	1.61%	0.121%
22	11.300	1549	1555	1565	rVB9	21594	73343	4.14%	0.312%
23	11.410	1565	1573	1579	rBV8	16873	47076	2.66%	0.200%
24	11.489	1580	1586	1598	rVB	295255	487297	27.51%	2.075%
25	11.599	1598	1604	1612	rBV7	39334	105824	5.97%	0.451%
26	11.685	1612	1618	1625	rBV5	74575	197661	11.16%	0.841%
27	11.837	1638	1643	1648	rBV5	19108	45873	2.59%	0.195%
28	11.892	1648	1652	1657	rBV3	21308	34543	1.95%	0.147%
29	11.953	1657	1662	1665	rBV3	19454	36960	2.09%	0.157%
30	12.002	1665	1670	1673	rBV4	35984	68743	3.88%	0.293%
31	12.123	1683	1690	1693	rBV2	74810	131801	7.44%	0.561%
32	12.166	1693	1697	1699	rBV3	36564	54059	3.05%	0.230%
33	12.203	1699	1703	1705	rVV3	71086	122260	6.90%	0.520%
34	12.239	1705	1709	1714	rVV3	180428	311455	17.58%	1.326%
35	12.325	1717	1723	1728	rVB5	55506	126192	7.12%	0.537%
36	12.392	1728	1734	1739	rVB3	162071	299454	16.91%	1.275%

LSC Area Percent Report

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Sample : M4407-02
Misc : 2.42G/5ML/MSVOA_Y/SOIL
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
FDR-BH-7-20211028-7.5-8

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

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Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y102221S.M
Title : SW846 8260

37	12.440	1739	1742	1744	rBV	26671	30243	1.71%	0.129%
38	12.489	1744	1750	1758	rBV2	756640	1350976	76.27%	5.751%
39	12.556	1758	1761	1767	rVB3	76079	127100	7.18%	0.541%
40	12.611	1767	1770	1771	rBV2	22371	24921	1.41%	0.106%
41	12.648	1772	1776	1785	rVV7	86492	187058	10.56%	0.796%
42	12.794	1797	1800	1807	rBV5	49630	89522	5.05%	0.381%
43	12.910	1815	1819	1822	rBV4	43115	66237	3.74%	0.282%
44	13.007	1830	1835	1837	rBV4	76578	116331	6.57%	0.495%
45	13.038	1837	1840	1849	rVB2	187260	332922	18.80%	1.417%
46	13.129	1850	1855	1859	rBV3	85207	145731	8.23%	0.620%
47	13.215	1866	1869	1873	rBV4	55749	96502	5.45%	0.411%
48	13.270	1875	1878	1879	rVV2	55524	71157	4.02%	0.303%
49	13.318	1882	1886	1890	rVB2	117988	188213	10.63%	0.801%
50	13.398	1895	1899	1900	rVV	94940	138420	7.82%	0.589%
51	13.428	1901	1904	1911	rVB2	228650	354494	20.01%	1.509%
52	13.501	1912	1916	1917	rBV2	38518	49155	2.78%	0.209%
53	13.617	1931	1935	1941	rVB7	69252	125859	7.11%	0.536%
54	13.696	1941	1948	1951	rBV2	134383	232723	13.14%	0.991%
55	13.751	1951	1957	1961	rVB5	156479	281744	15.91%	1.199%
56	13.806	1962	1966	1972	rBV6	85665	160351	9.05%	0.683%
57	13.971	1989	1993	1998	rVB3	109133	164973	9.31%	0.702%
58	14.062	2004	2008	2015	rBV3	311115	625670	35.32%	2.664%
59	14.135	2016	2020	2025	rVV6	59276	138170	7.80%	0.588%
60	14.190	2026	2029	2034	rVB3	66724	104062	5.88%	0.443%
61	14.306	2044	2048	2056	rVB3	152046	376277	21.24%	1.602%
62	14.379	2056	2060	2064	rBV	99180	142027	8.02%	0.605%
63	14.452	2064	2072	2077	rBV	1193358	1771197	100.00%	7.540%
64	14.586	2090	2094	2097	rBV2	207770	348588	19.68%	1.484%
65	14.666	2102	2107	2111	rVV2	388256	619979	35.00%	2.639%
66	14.733	2114	2118	2124	rVV3	438642	1000664	56.50%	4.260%
67	14.788	2124	2127	2137	rVB2	488011	763275	43.09%	3.249%
68	14.867	2137	2140	2142	rBV3	34080	45626	2.58%	0.194%
69	14.946	2148	2153	2158	rVB2	368725	567740	32.05%	2.417%
70	15.007	2158	2163	2167	rBV	551315	779140	43.99%	3.317%
71	15.056	2168	2171	2174	rVV5	65350	89645	5.06%	0.382%
72	15.099	2174	2178	2187	rVB4	146073	281217	15.88%	1.197%
73	15.214	2193	2197	2201	rVB2	212990	284046	16.04%	1.209%
74	15.342	2212	2218	2223	rBV3	130255	310451	17.53%	1.322%
75	15.446	2230	2235	2242	rVB6	145367	316259	17.86%	1.346%
76	15.525	2243	2248	2249	rBV4	94659	145161	8.20%	0.618%

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Operator : SY/MD
Sample : M4407-02
Misc : 2.42G/5ML/MSVOA_Y/SOIL
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
FDR-BH-7-20211028-7.5-8

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
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Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y102221S.M
Title : SW846 8260

77	15.617	2257	2263	2269	rBV5	163425	407239	22.99%	1.734%
78	15.720	2275	2280	2284	rBV	181626	292647	16.52%	1.246%
79	15.854	2292	2302	2317	rVB2	523494	1310574	73.99%	5.579%
80	16.025	2325	2330	2335	rBV5	85819	194504	10.98%	0.828%
81	16.111	2335	2344	2348	rVV5	193564	625086	35.29%	2.661%
82	16.153	2349	2351	2355	rVV	167820	273012	15.41%	1.162%
83	16.220	2355	2362	2370	rVV2	555906	1254294	70.82%	5.340%
84	16.293	2371	2374	2383	rVB4	98659	217908	12.30%	0.928%
85	16.428	2392	2396	2400	rBV5	65475	112193	6.33%	0.478%
86	16.482	2400	2405	2410	rBV4	77470	160817	9.08%	0.685%
87	16.745	2444	2448	2455	rVB3	107072	221759	12.52%	0.944%

Sum of corrected areas: 23489242

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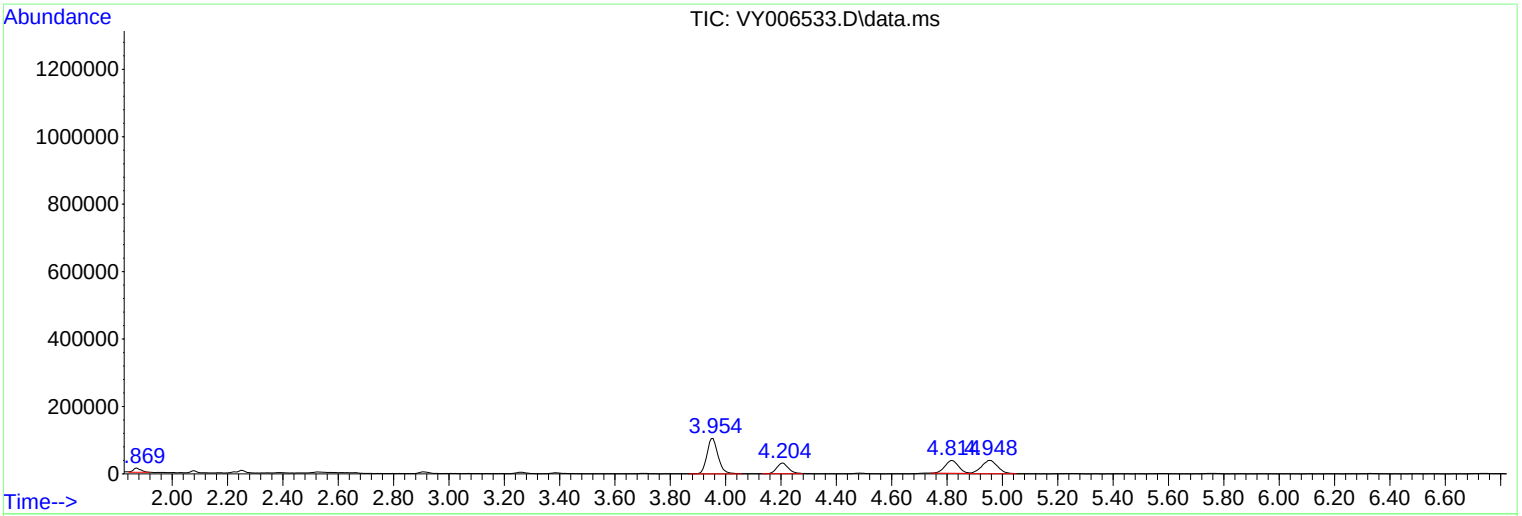
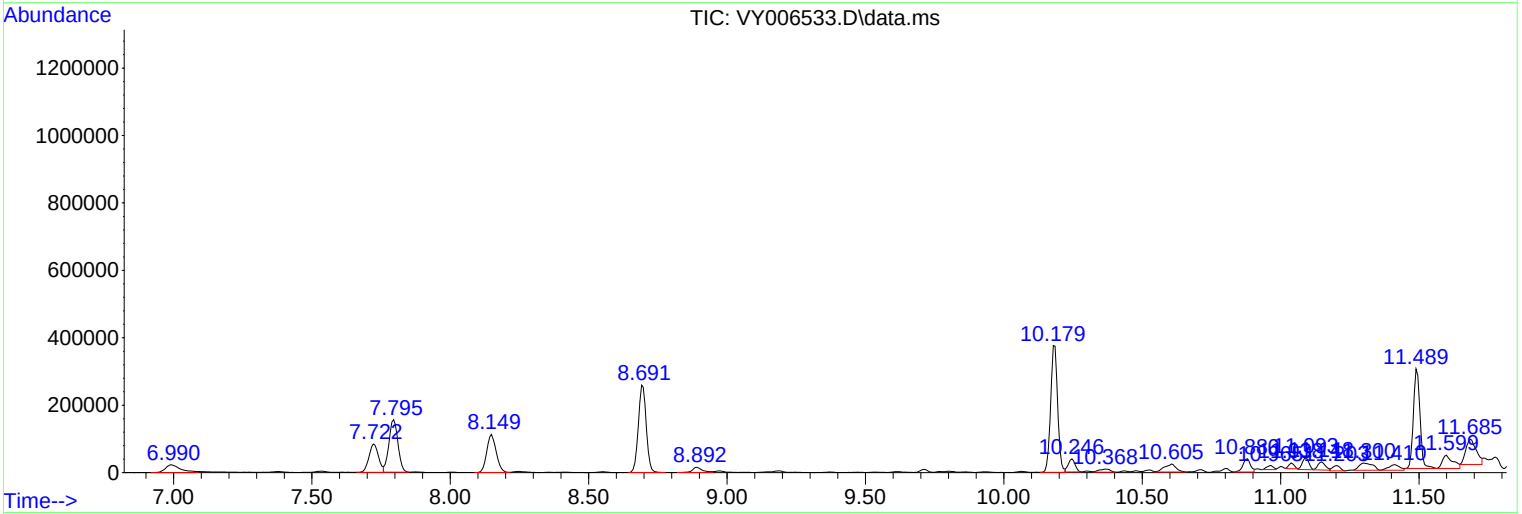
Misc : 2.42G/5ML/MSVOA_Y/SOIL
ALS Vial : 8 Sample Multiplier: 1

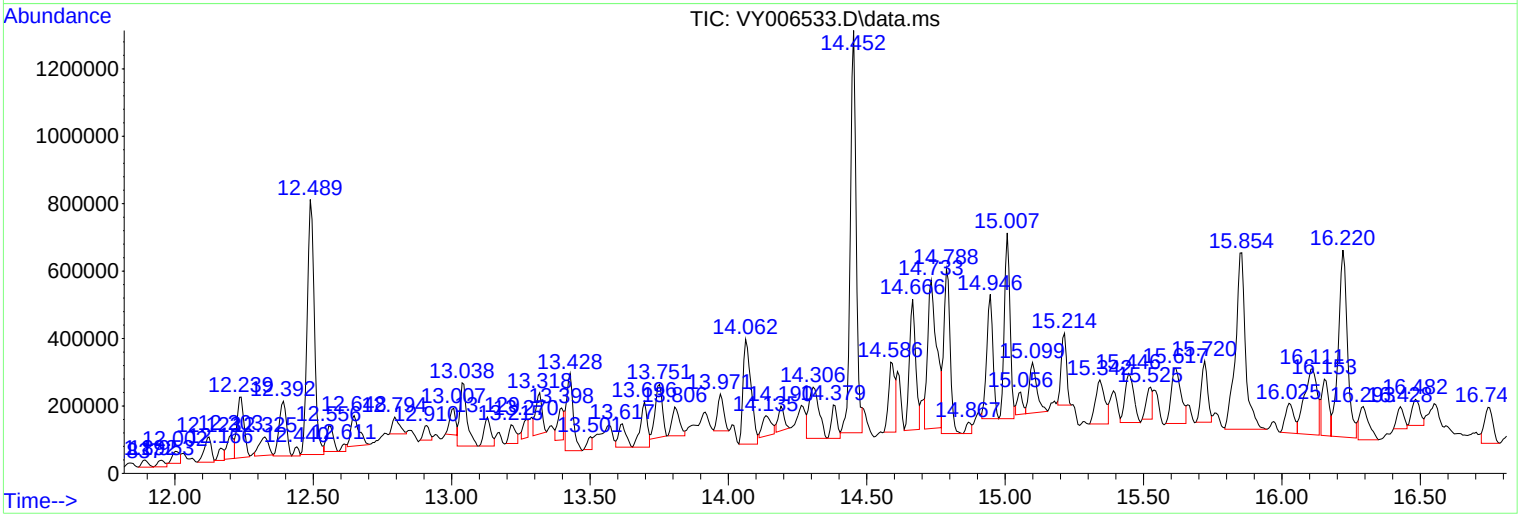
Instrument :
MSVOA_Y
ClientSampleId :
FDR-BH-7-20211028-7.5-8

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y102221S.M
Quant Title : SW846 8260

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







Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY102921\
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Instrument :
MSVOA_Y
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FDR-BH-7-20211028-7.5-8

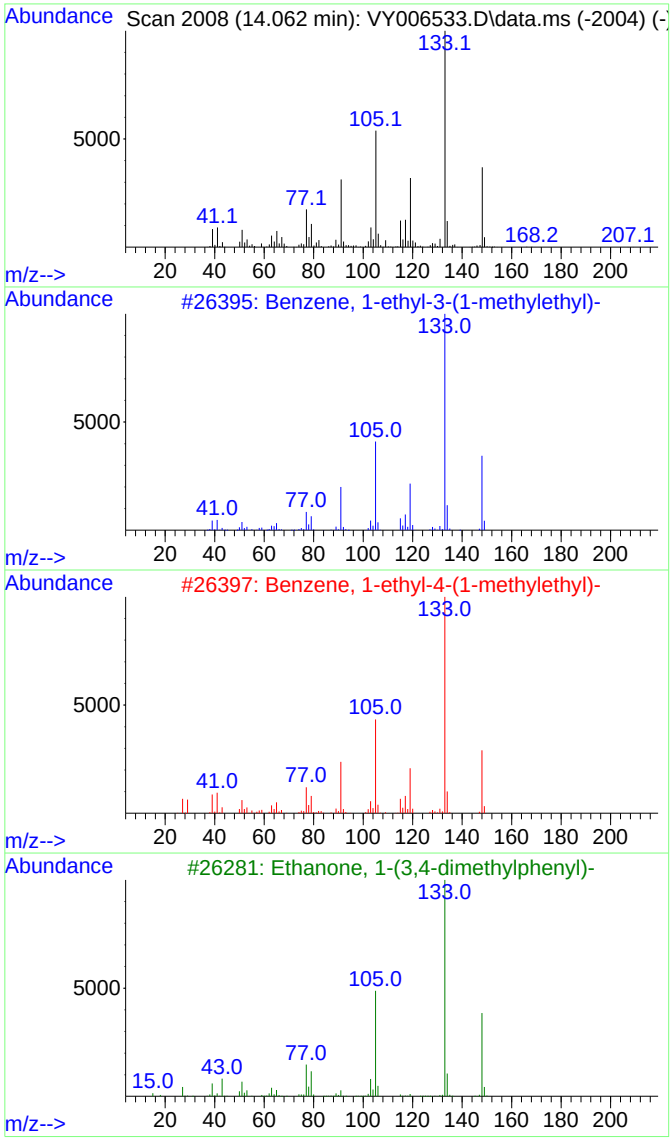
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y102221S.M
Quant Title : SW846 8260

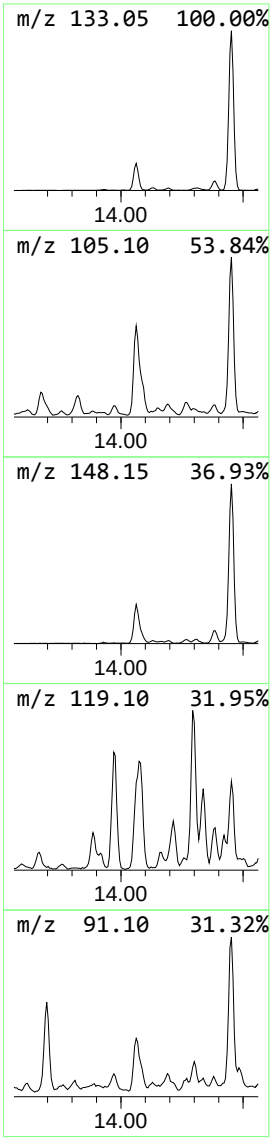
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Benzene, 1-ethyl-3-(1-methy... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.062	88.25 ug/l	625670	1,4-Dichlorobenzene-d4	13.428

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3-(1-methylethyl)-	148	C11H16	004920-99-4	97
2	Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	95
3	Ethanone, 1-(3,4-dimethylphenyl)-	148	C10H12O	003637-01-2	92
4	Benzene, 2,4-diethyl-1-methyl-	148	C11H16	001758-85-6	87
5	Benzene, pentamethyl-	148	C11H16	000700-12-9	87





Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY102921\
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Acq On : 29 Oct 2021 12:33
Operator : SY/MD
Sample : M4407-02
Misc : 2.42G/5ML/MSVOA_Y/SOIL
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
FDR-BH-7-20211028-7.5-8

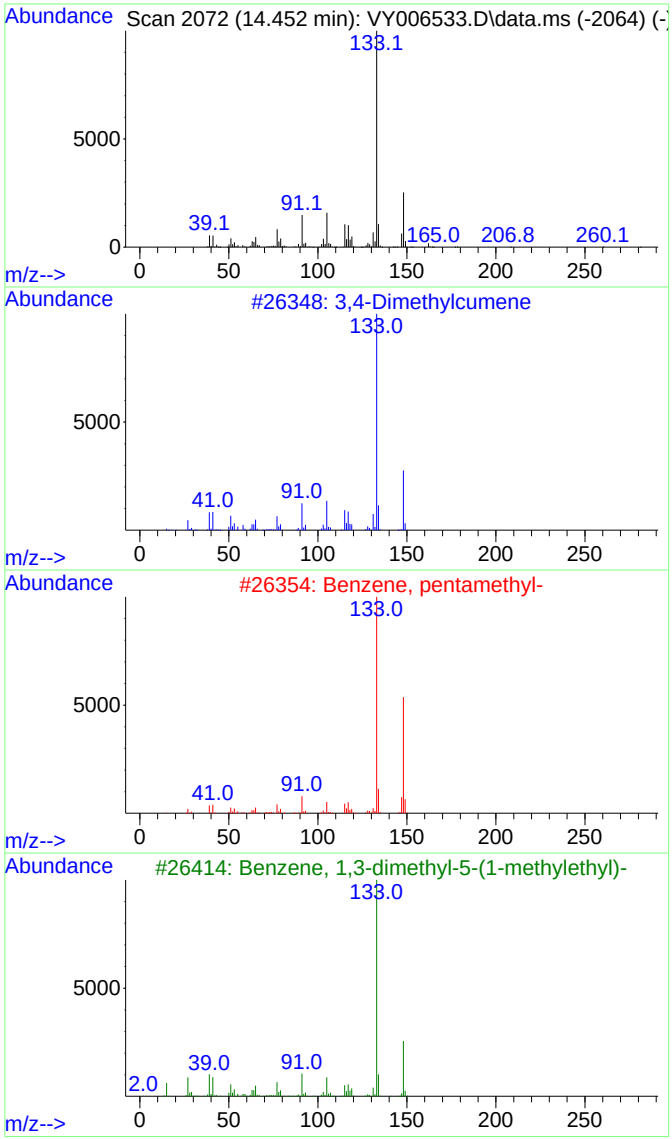
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y102221S.M
Quant Title : SW846 8260

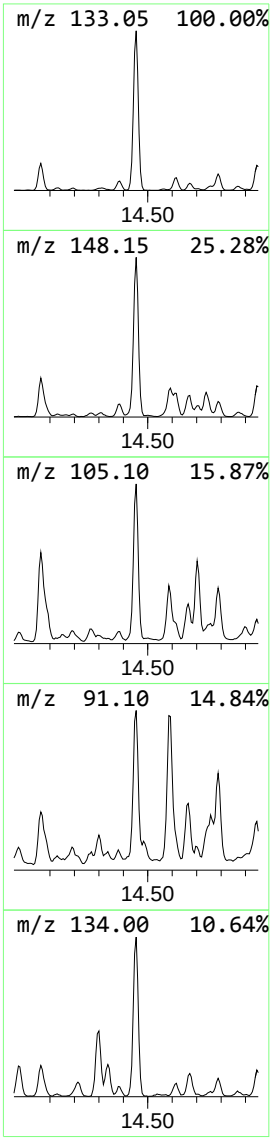
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 3,4-Dimethylcumene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.452	249.82 ug/l	1771200	1,4-Dichlorobenzene-d4	13.428

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,4-Dimethylcumene	148	C11H16	1000370-34-1	97
2	Benzene, pentamethyl-	148	C11H16	000700-12-9	94
3	Benzene, 1,3-dimethyl-5-(1-methy...	148	C11H16	004706-90-5	91
4	Benzene, 1,4-dimethyl-2-(1-methy...	148	C11H16	004132-72-3	91
5	Benzene, 2,4-dimethyl-1-(1-methy...	148	C11H16	004706-89-2	90





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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y102221S.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
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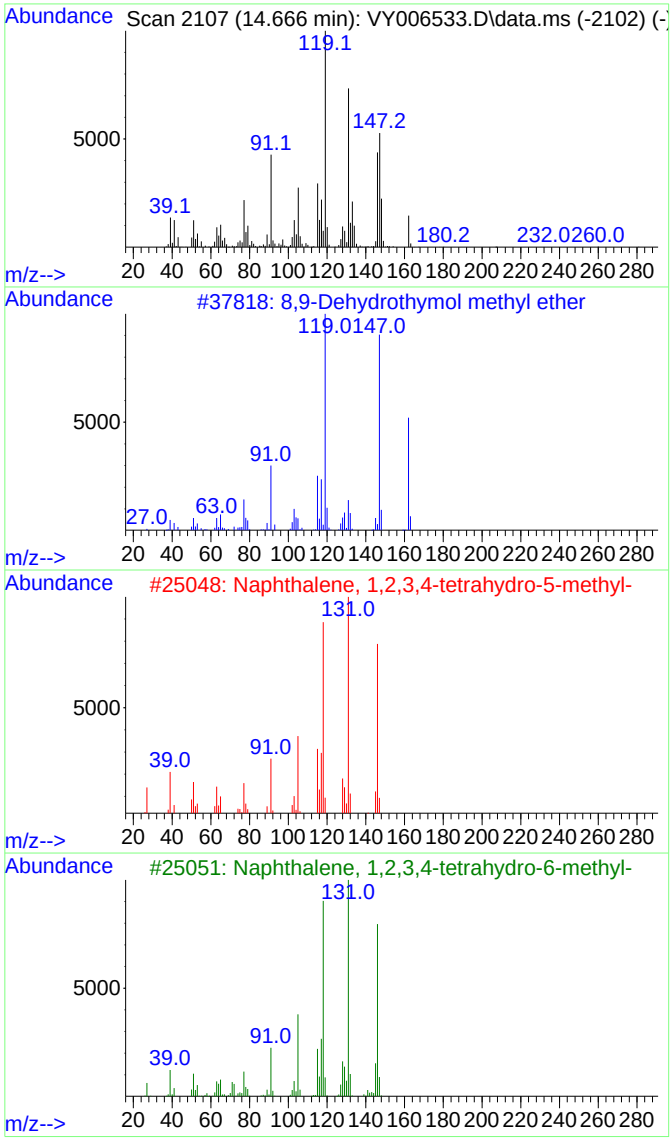
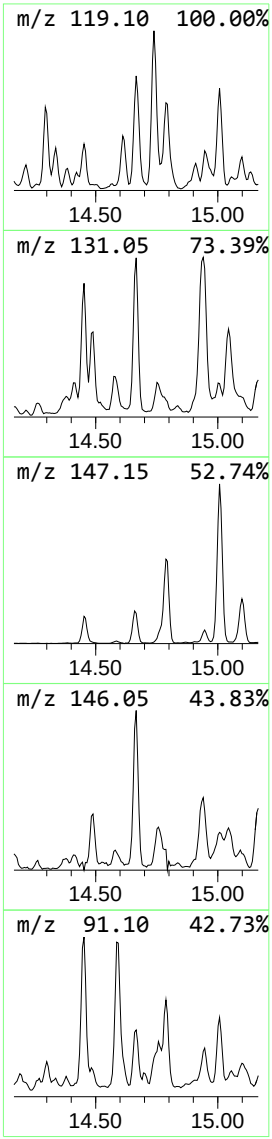
****
Peak Number 3 8,9-Dehydrothymol methyl ether Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.666	87.45 ug/l	619979	1,4-Dichlorobenzene-d4	13.428

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	8,9-Dehydrothymol methyl ether	162	C11H14O	039701-08-1	50
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	46
3	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	46
4	Benzene, 1-(1,1-dimethylethyl)-4...	162	C12H18	007364-19-4	45
5	Bicyclo[4.2.0]octa-1,3,5-triene,...	146	C11H14	027087-54-3	38





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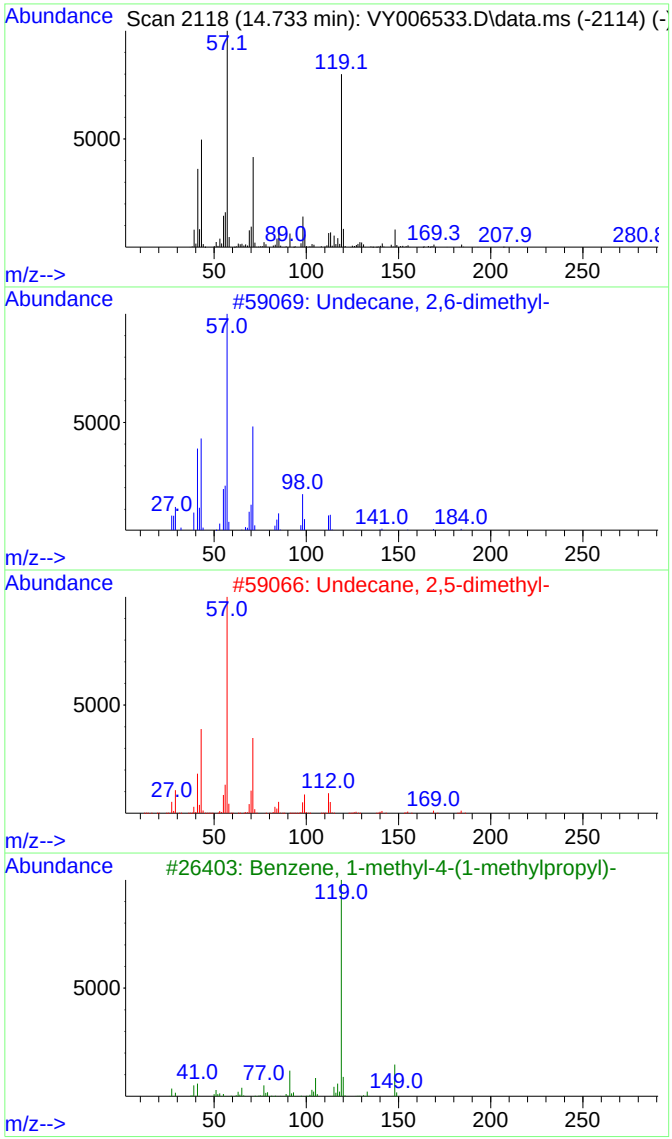
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y102221S.M
Quant Title : SW846 8260

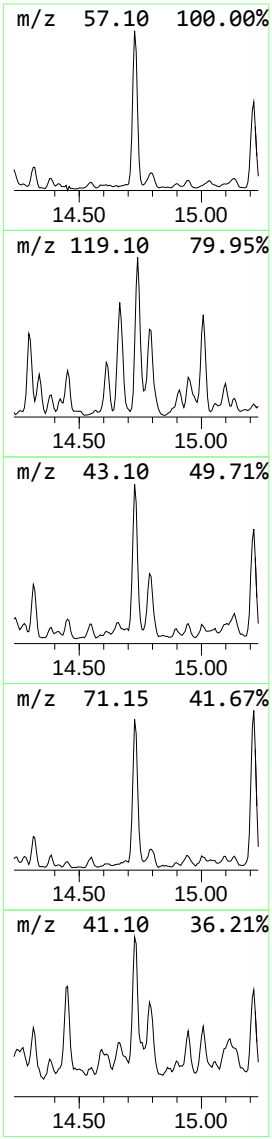
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Undecane, 2,6-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.733	141.14 ug/l	1000660	1,4-Dichlorobenzene-d4	13.428

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane, 2,6-dimethyl-	184	C13H28	017301-23-4	86
2	Undecane, 2,5-dimethyl-	184	C13H28	017301-22-3	60
3	Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	45
4	Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	43
5	4'-Methylpropiophenone	148	C10H12O	005337-93-9	38





Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY102921\
Data File : VY006533.D
Acq On : 29 Oct 2021 12:33
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Sample : M4407-02
Misc : 2.42G/5ML/MSVOA_Y/SOIL
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
FDR-BH-7-20211028-7.5-8

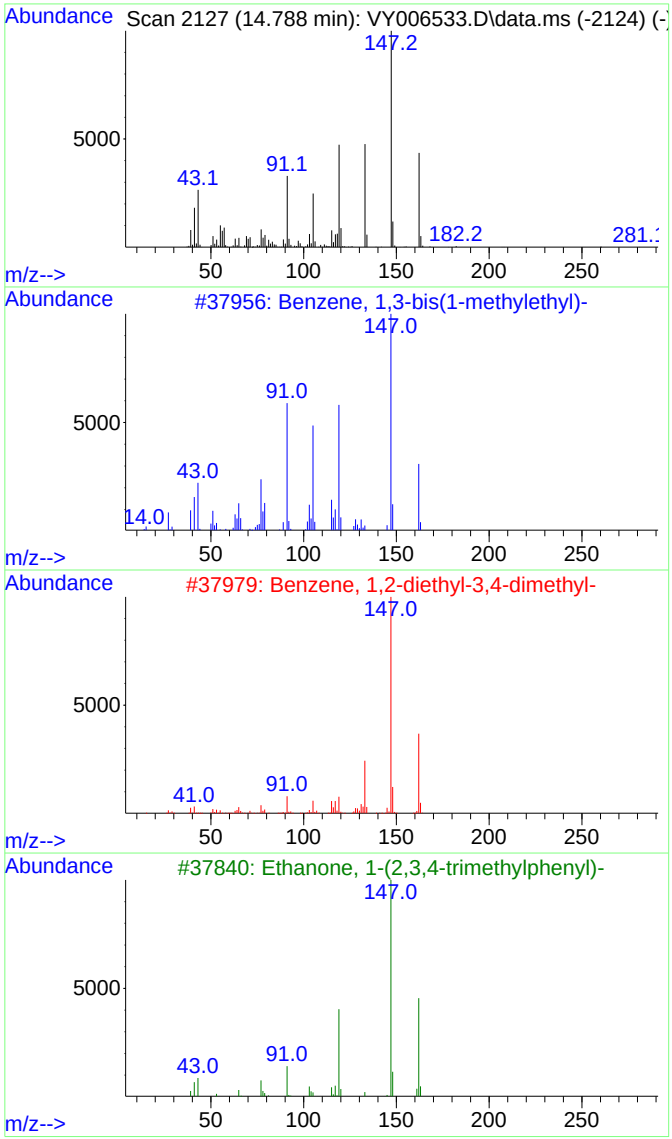
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y102221S.M
Quant Title : SW846 8260

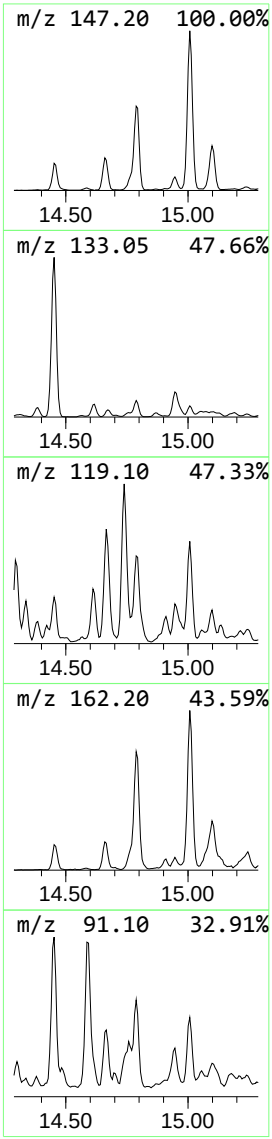
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzene, 1,3-bis(1-methylet... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.788	107.66 ug/l	763275	1,4-Dichlorobenzene-d4	13.428

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-bis(1-methylethyl)-	162	C12H18	000099-62-7	74
2	Benzene, 1,2-diethyl-3,4-dimethyl-	162	C12H18	054410-75-2	72
3	Ethanone, 1-(2,3,4-trimethylphen...	162	C11H14O	001467-36-3	70
4	Ethanone, 1-(2,4,5-trimethylphen...	162	C11H14O	002040-07-5	70
5	2,3,6-Trimethylacetophenone	162	C11H14O	1000342-30-1	70





Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY102921\
Data File : VY006533.D
Acq On : 29 Oct 2021 12:33
Operator : SY/MD
Sample : M4407-02
Misc : 2.42G/5ML/MSVOA_Y/SOIL
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
FDR-BH-7-20211028-7.5-8

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y102221S.M
Quant Title : SW846 8260

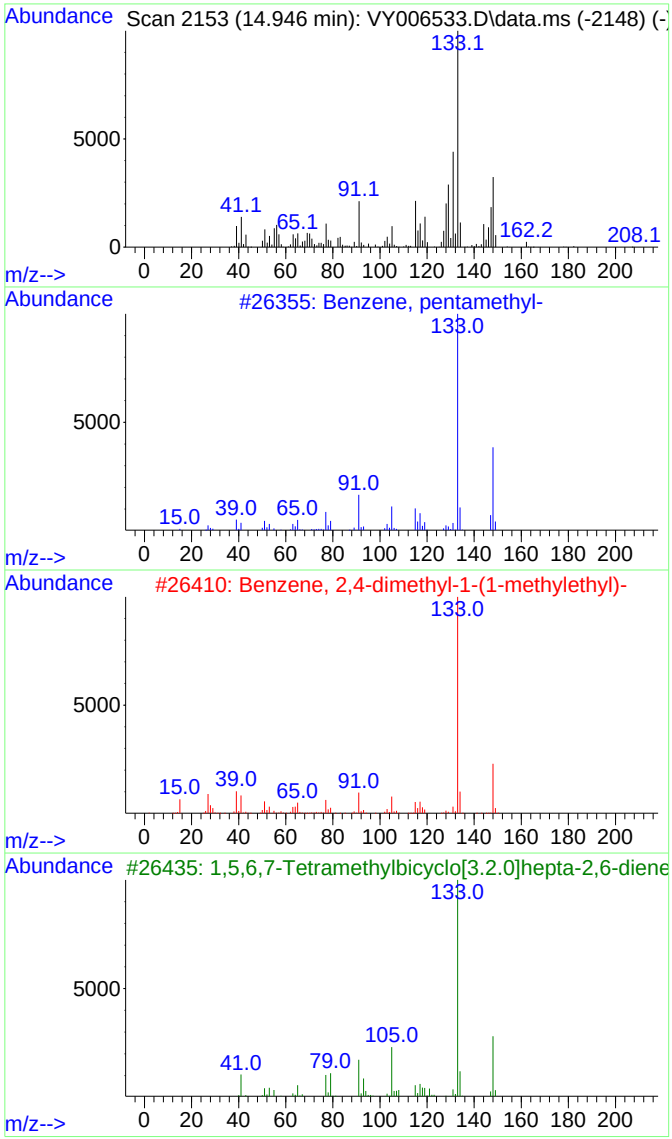
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

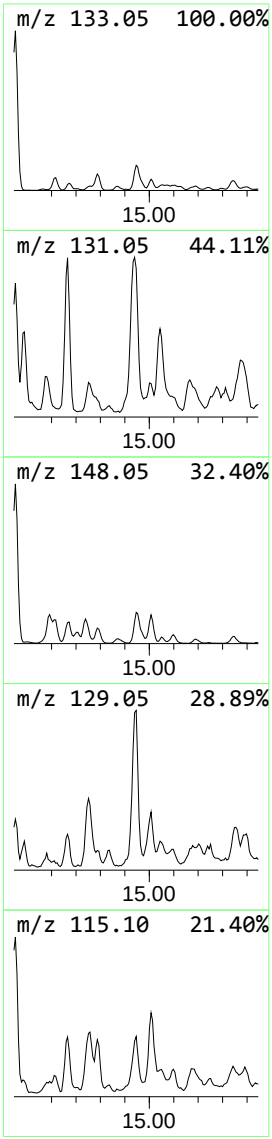
****

Peak Number 6 unknown14.946 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.946	80.08 ug/l	567740	1,4-Dichlorobenzene-d4	13.428

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, pentamethyl-	148	C11H16	000700-12-9	45
2	Benzene, 2,4-dimethyl-1-(1-methy...	148	C11H16	004706-89-2	43
3	1,5,6,7-Tetramethylbicyclo[3.2.0...	148	C11H16	134329-46-7	43
4	Benzene, 1,4-diethyl-2-methyl-	148	C11H16	013632-94-5	43
5	Benzene, 1,4-dimethyl-2-(1-methy...	148	C11H16	004132-72-3	43





Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY102921\
Data File : VY006533.D
Acq On : 29 Oct 2021 12:33
Operator : SY/MD
Sample : M4407-02
Misc : 2.42G/5ML/MSVOA_Y/SOIL
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
FDR-BH-7-20211028-7.5-8

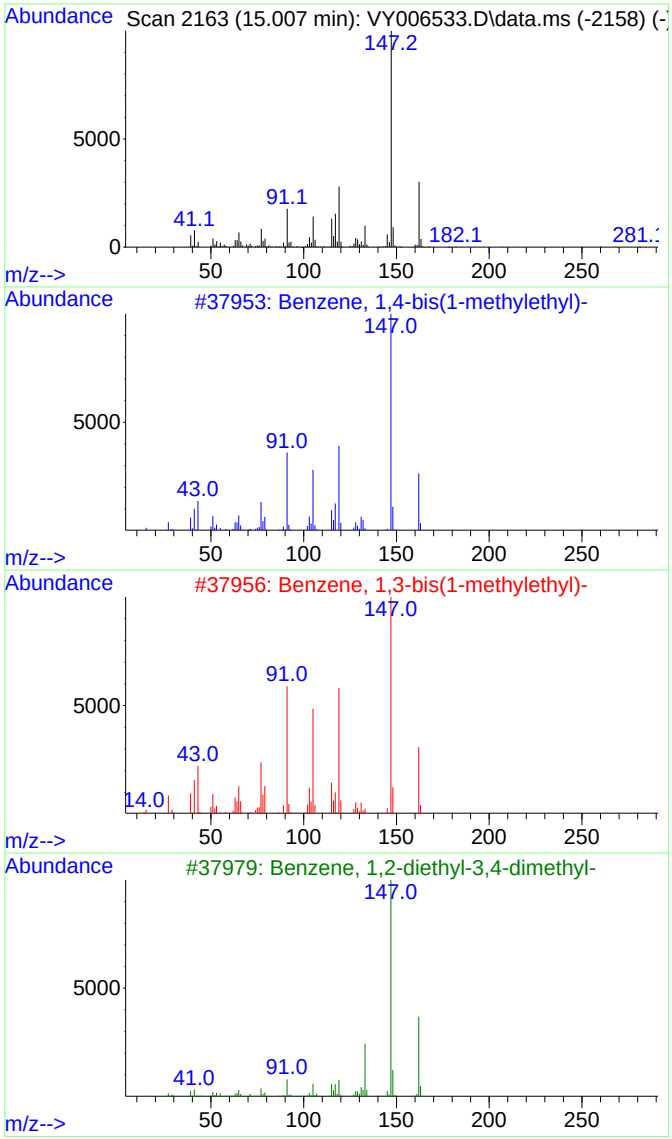
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y102221S.M
Quant Title : SW846 8260

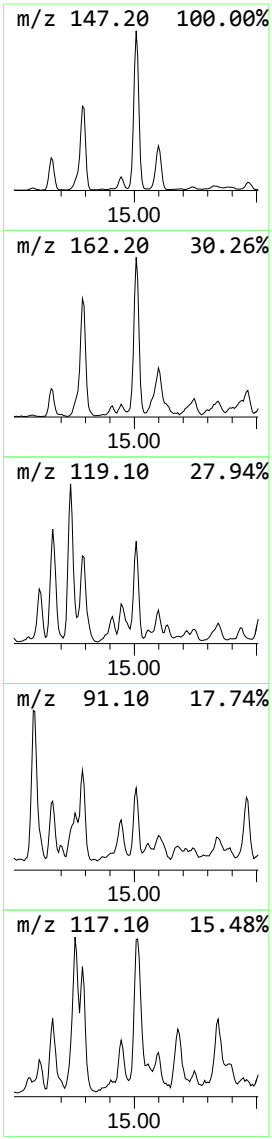
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzene, 1,4-bis(1-methylet... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.007	109.90 ug/l	779140	1,4-Dichlorobenzene-d4	13.428

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,4-bis(1-methylethyl)-	162	C12H18	000100-18-5	87
2	Benzene, 1,3-bis(1-methylethyl)-	162	C12H18	000099-62-7	87
3	Benzene, 1,2-diethyl-3,4-dimethyl-	162	C12H18	054410-75-2	80
4	Tricyclo[3.2.1.02,7]oct-3-ene, 2...	162	C12H18	062338-44-7	80
5	Benzene, 1-(1,1-dimethylethyl)-4...	162	C12H18	007364-19-4	74





Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY102921\
Data File : VY006533.D
Acq On : 29 Oct 2021 12:33
Operator : SY/MD
Sample : M4407-02
Misc : 2.42G/5ML/MSVOA_Y/SOIL
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
FDR-BH-7-20211028-7.5-8

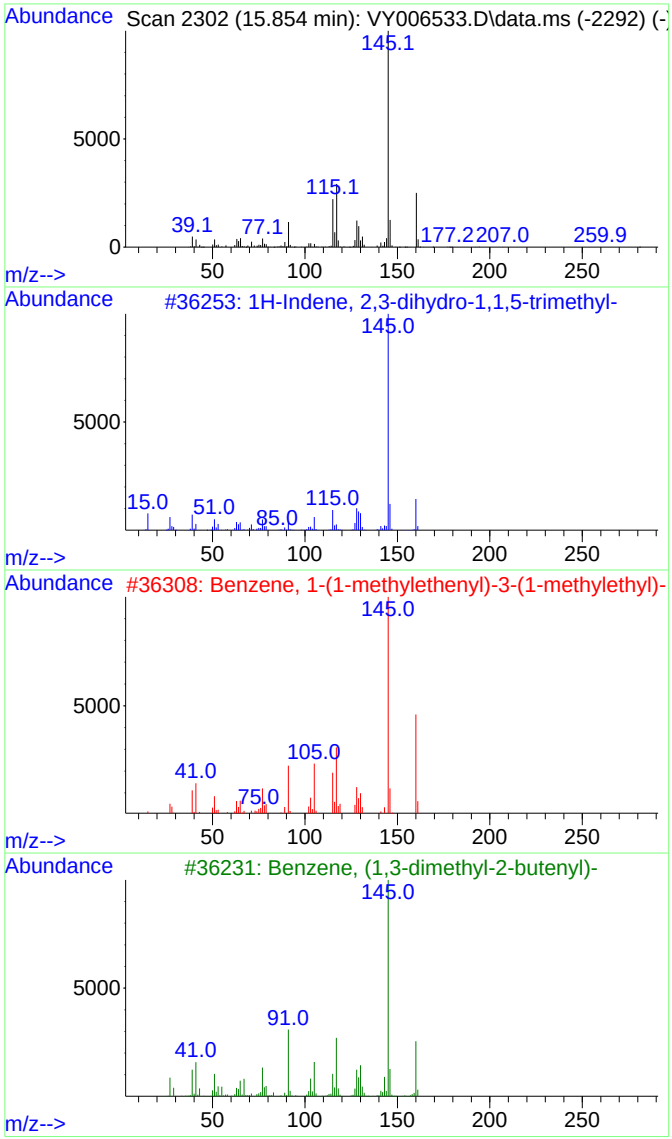
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y102221S.M
Quant Title : SW846 8260

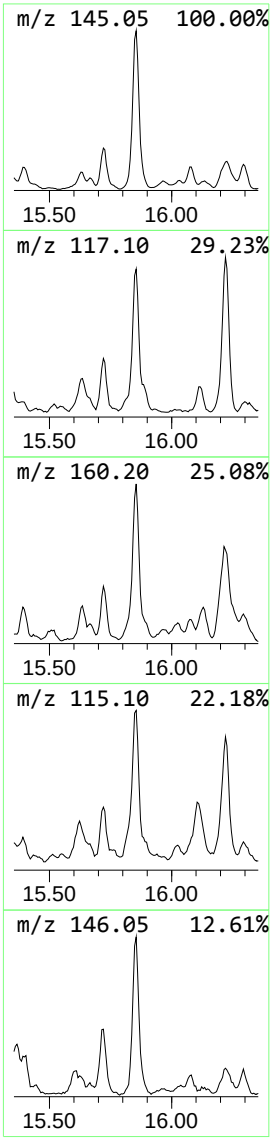
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.855	184.85 ug/l	1310570	1,4-Dichlorobenzene-d4	13.428

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,1,5-tri...	160	C12H16	040650-41-7	90
2	Benzene, 1-(1-methylethenyl)-3-(...	160	C12H16	001129-29-9	87
3	Benzene, (1,3-dimethyl-2-butenyl)-	160	C12H16	050704-01-3	83
4	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	025419-33-4	83
5	Benzene, 1-(1,1-dimethylethyl)-4...	160	C12H16	001746-23-2	83





Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY102921\
Data File : VY006533.D
Acq On : 29 Oct 2021 12:33
Operator : SY/MD
Sample : M4407-02
Misc : 2.42G/5ML/MSVOA_Y/SOIL
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
FDR-BH-7-20211028-7.5-8

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y102221S.M
Quant Title : SW846 8260

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

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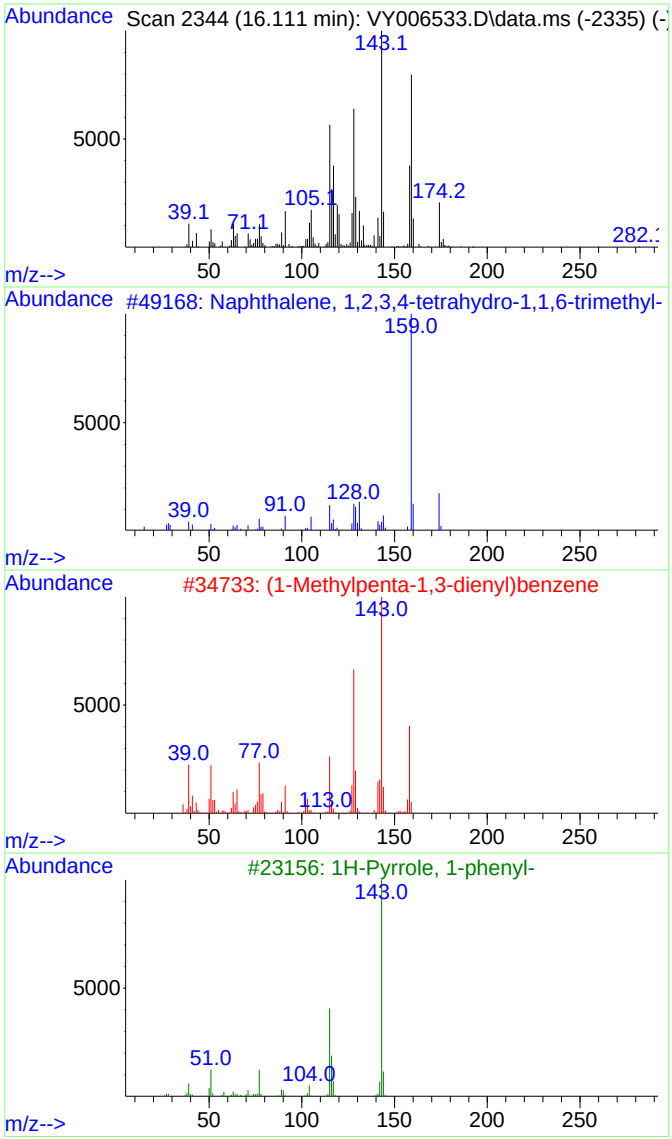
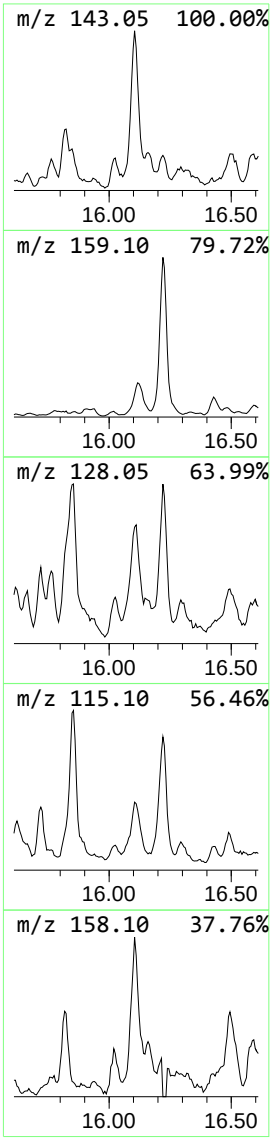
Peak Number 9 unknown16.110 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.110	88.17 ug/l	625086	1,4-Dichlorobenzene-d4	13.428

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	174	C13H18	000475-03-6	46
2	(1-Methylpenta-1,3-dienyl)benzene	158	C12H14	116669-49-9	45
3	1H-Pyrrole, 1-phenyl-	143	C10H9N	000635-90-5	38
4	(1-Methylenepent-2-enyl)benzene	158	C12H14	084313-24-6	38
5	(1-Methylpenta-2,4-dienyl)benzene	158	C12H14	1000210-01-1	38





Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY102921\
Data File : VY006533.D
Acq On : 29 Oct 2021 12:33
Operator : SY/MD
Sample : M4407-02
Misc : 2.42G/5ML/MSVOA_Y/SOIL
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
FDR-BH-7-20211028-7.5-8

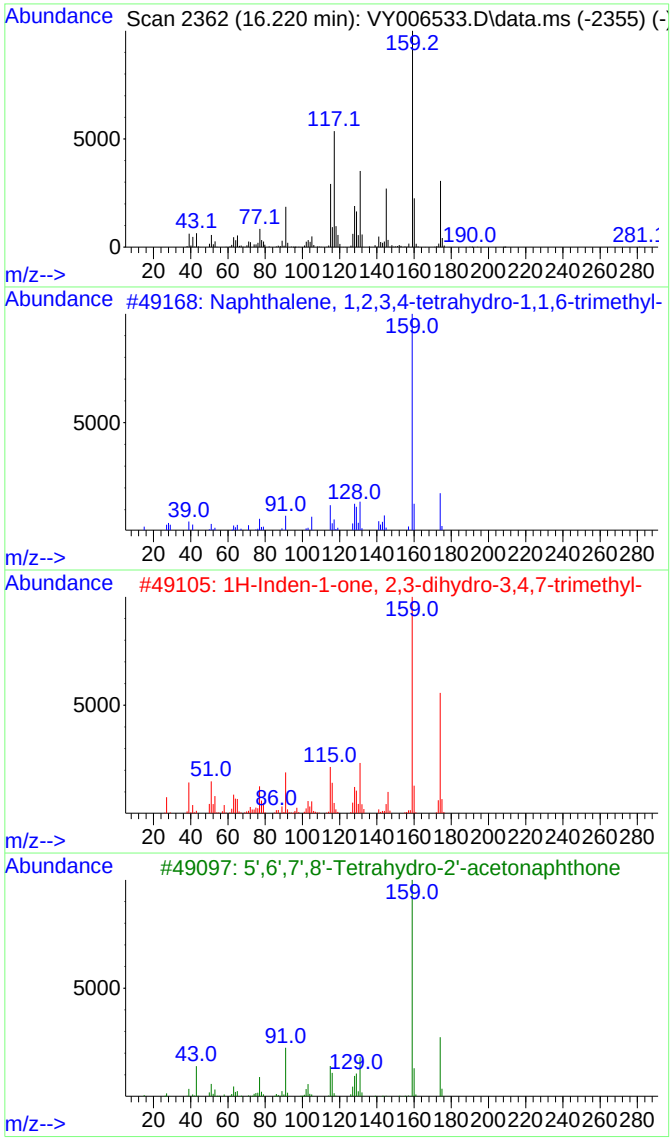
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y102221S.M
Quant Title : SW846 8260

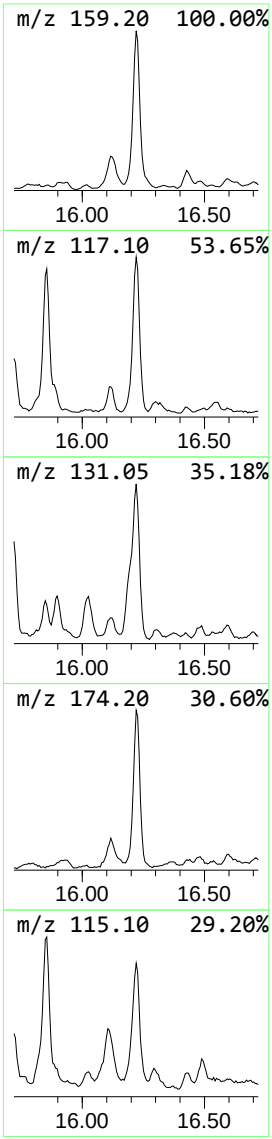
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.220	176.91 ug/l	1254290	1,4-Dichlorobenzene-d4	13.428

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	174	C13H18	000475-03-6	64
2	1H-Inden-1-one, 2,3-dihydro-3,4,...	174	C12H14O	035322-84-0	58
3	5',6',7',8'-Tetrahydro-2'-aceton...	174	C12H14O	000774-55-0	53
4	Ethanone, 1-[4-(1-methyl-2-prope...	174	C12H14O	109586-49-4	52
5	1H-Indene, 2,3-dihydro-1,1,4,6-t...	174	C13H18	000941-60-6	49





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY102921\
Data File : VY006533.D
Acq On : 29 Oct 2021 12:33
Operator : SY/MD
Sample : M4407-02
Misc : 2.42G/5ML/MSVOA_Y/SOIL
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
FDR-BH-7-20211028-7.5-8

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y102221S.M
Quant Title : SW846 8260

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Benzene, 1-ethy...	14.062	88.3	ug/l	625670	4	13.428	354494	50.0
3,4-Dimethylcumene	14.452	249.8	ug/l	1771200	4	13.428	354494	50.0
8,9-Dehydrothym...	14.666	87.5	ug/l	619979	4	13.428	354494	50.0
Undecane, 2,6-d...	14.733	141.1	ug/l	1000660	4	13.428	354494	50.0
Benzene, 1,3-bi...	14.788	107.7	ug/l	763275	4	13.428	354494	50.0
unknown14.946	14.946	80.1	ug/l	567740	4	13.428	354494	50.0
Benzene, 1,4-bi...	15.007	109.9	ug/l	779140	4	13.428	354494	50.0
1H-Indene, 2,3-...	15.855	184.8	ug/l	1310570	4	13.428	354494	50.0
unknown16.110	16.110	88.2	ug/l	625086	4	13.428	354494	50.0
Naphthalene, 1,...	16.220	176.9	ug/l	1254290	4	13.428	354494	50.0