LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY110221\
Data File : VY006574.D
Acq On : 02 Nov 2021 13:53
Operator : SY/MD
Sample : M4427-02
Misc : 4.63G/5ML/MSVOA_Y/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
FDR-BH-5-20211029-6.5-7

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: VY006574.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.954	337	350	370	rBV2	83753	259039	1.25%	0.241%
2	6.990	834	848	868	rVB	105178	362901	1.76%	0.337%
3	7.728	958	969	974	rBV	92607	219749	1.06%	0.204%
4	7.795	974	980	988	rVB	174670	388638	1.88%	0.361%
5	8.148	1029	1038	1050	rVB	110493	238031	1.15%	0.221%
6	8.331	1050	1068	1078	rBV2	80966	220413	1.07%	0.205%
7	8.697	1119	1128	1135	rBV	274007	544878	2.64%	0.506%
8	8.776	1135	1141	1149	rVB	141565	272197	1.32%	0.253%
9	8.929	1159	1166	1174	rBV	151020	325123	1.57%	0.302%
10	9.325	1224	1231	1236	rBV	103157	223842	1.08%	0.208%
11	9.392	1236	1242	1247	rVV	192689	387693	1.88%	0.360%
12	9.538	1261	1266	1273	rVB2	120964	245223	1.19%	0.228%
13	9.618	1273	1279	1288	rVB	175386	345539	1.67%	0.321%
14	9.758	1290	1302	1313	rBV2	398217	879198	4.25%	0.816%
15	10.117	1349	1361	1367	rBV	915636	1586899	7.68%	1.473%
16	10.184	1367	1372	1383	rVB	428126	763969	3.70%	0.709%
17	10.794	1467	1472	1474	rBV	190373	293333	1.42%	0.272%
18	10.831	1474	1478	1485	rVB	562605	988788	4.78%	0.918%
19	10.965	1495	1500	1504	rBV	261548	468096	2.26%	0.435%
20	11.087	1515	1520	1526	rVB2	301481	532505	2.58%	0.494%
21	11.148	1526	1530	1536	rVV2	193503	395350	1.91%	0.367%
22	11.227	1536	1543	1545	rVV3	296278	691945	3.35%	0.642%
23	11.264	1545	1549	1564	rVB2	730928	1465571	7.09%	1.361%
24	11.495	1577	1587	1591	rBV2	410819	981272	4.75%	0.911%
25	11.562	1591	1598	1604	rVV3	471936	1236598	5.98%	1.148%
26	11.635	1604	1610	1616	rVB5	111589	266970	1.29%	0.248%
27	11.751	1622	1629	1633	rBV2	336590	663329	3.21%	0.616%
28	11.855	1640	1646	1648	rBV3	179060	328385	1.59%	0.305%
29	11.892	1648	1652	1657	rVB4	225037	454181	2.20%	0.422%
30	12.135	1689	1692	1695	rBV2	149397	221843	1.07%	0.206%
31	12.209	1700	1704	1705	rBV	204030	257661	1.25%	0.239%
32	12.245	1706	1710	1713	rVV2	299997	543685	2.63%	0.505%
33	12.282	1713	1716	1725	rVB4	448996	941457	4.55%	0.874%
34	12.385	1725	1733	1739	rBV4	711955	1998243	9.66%	1.855%
35	12.483	1744	1749	1753	rBV4	453319	776100	3.75%	0.721%
36	12.532	1753	1757	1768	rVB3	847777	2529093	12.23%	2.348%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY110221\
Data File : VY006574.D
Acq On : 02 Nov 2021 13:53
Operator : SY/MD
Sample : M4427-02
Misc : 4.63G/5ML/MSVOA_Y/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
FDR-BH-5-20211029-6.5-7

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

37	12.629	1768	1773	1777	rBV4	396960	878328	4.25%	0.815%
38	12.708	1782	1786	1788	rBV	313795	436358	2.11%	0.405%
39	12.788	1794	1799	1803	rBV	677385	1182722	5.72%	1.098%
40	12.830	1803	1806	1812	rVV4	333241	741352	3.59%	0.688%
41	12.891	1812	1816	1822	rVV6	334033	688463	3.33%	0.639%
42	12.952	1823	1826	1830	rVV2	180670	361516	1.75%	0.336%
43	13.074	1834	1846	1853	rVV	11765553	20675085	100.00%	19.196%
44	13.141	1853	1857	1860	rVV2	551079	1033764	5.00%	0.960%
45	13.178	1860	1863	1870	rVB3	975771	1694536	8.20%	1.573%
46	13.257	1872	1876	1877	rBV2	250674	353779	1.71%	0.328%
47	13.288	1877	1881	1885	rVV2	589589	995091	4.81%	0.924%
48	13.355	1887	1892	1897	rVV	2752642	4705682	22.76%	4.369%
49	13.416	1897	1902	1909	rVB	1826650	3695585	17.87%	3.431%
50	13.513	1909	1918	1922	rBV5	983169	2409242	11.65%	2.237%
51	13.617	1929	1935	1941	rBV4	1342810	3253678	15.74%	3.021%
52	13.678	1941	1945	1948	rVV	813479	1176077	5.69%	1.092%
53	13.720	1949	1952	1957	rVB2	838982	1427431	6.90%	1.325%
54	13.788	1957	1963	1970	rBV2	4587144	8060597	38.99%	7.484%
55	13.855	1970	1974	1980	rVV2	1621544	2384388	11.53%	2.214%
56	13.952	1986	1990	1996	rBV3	350483	909799	4.40%	0.845%
57	14.007	1996	1999	2002	rVV3	452324	675584	3.27%	0.627%
58	14.044	2002	2005	2006	rVV	586547	721606	3.49%	0.670%
59	14.080	2006	2011	2016	rVV	7541710	11596909	56.09%	10.767%
60	14.135	2017	2020	2025	rVB2	858949	1335435	6.46%	1.240%
61	14.269	2037	2042	2043	rBV2	338458	526488	2.55%	0.489%
62	14.379	2056	2060	2064	rVB2	457876	642396	3.11%	0.596%
63	14.531	2082	2085	2089	rBV	412030	588086	2.84%	0.546%
64	14.617	2095	2099	2104	rVB3	353590	641424	3.10%	0.596%
65	14.757	2111	2122	2128	rBV	1811232	3413756	16.51%	3.170%
66	14.854	2133	2138	2141	rBV4	296893	582736	2.82%	0.541%
67	14.952	2148	2154	2163	rVB7	477911	1289021	6.23%	1.197%
68	15.220	2194	2198	2202	rVB	367320	533895	2.58%	0.496%
69	15.281	2206	2208	2214	rVB2	310596	466649	2.26%	0.433%
70	15.348	2215	2219	2223	rBV3	199159	289876	1.40%	0.269%
71	15.501	2239	2244	2248	rBV	1050975	1903502	9.21%	1.767%
72	15.568	2252	2255	2258	rVV3	267337	423927	2.05%	0.394%
73	15.879	2302	2306	2314	rVB3	279975	658220	3.18%	0.611%
74	16.007	2324	2327	2338	rVB5	341832	1055142	5.10%	0.980%

Sum of corrected areas: 107705862

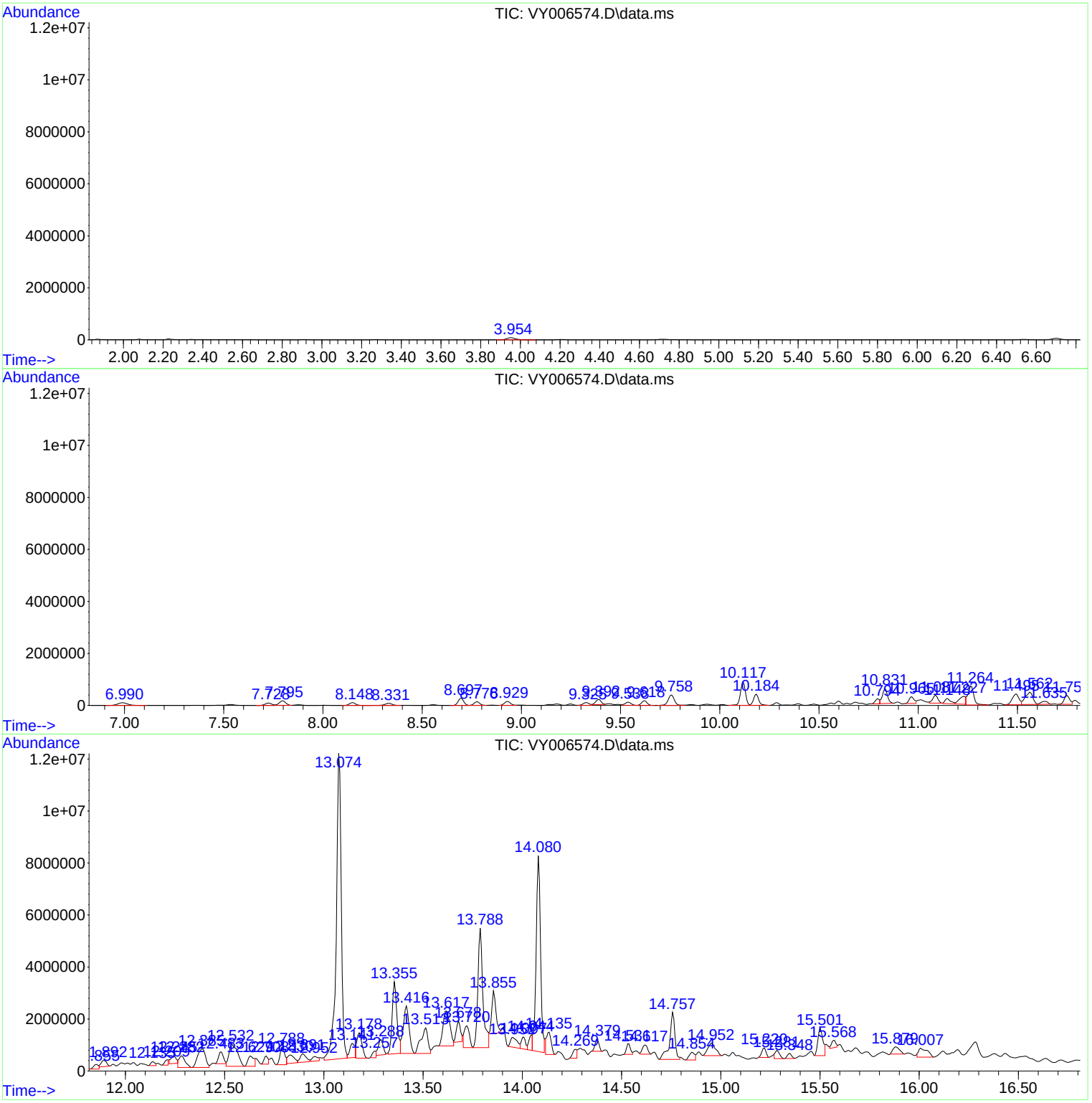
Instrument :
MSVOA_Y
ClientSampleId :
FDR-BH-5-20211029-6.5-7

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY110221\
Data File : VY006574.D
Acq On : 02 Nov 2021 13:53
Operator : SY/MD
Sample : M4427-02
Misc : 4.63G/5ML/MSVOA_Y/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
FDR-BH-5-20211029-6.5-7

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\VY110221\
Data File : VY006574.D
Acq On : 02 Nov 2021 13:53
Operator : SY/MD
Sample : M4427-02
Misc : 4.63G/5ML/MSVOA_Y/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
FDR-BH-5-20211029-6.5-7

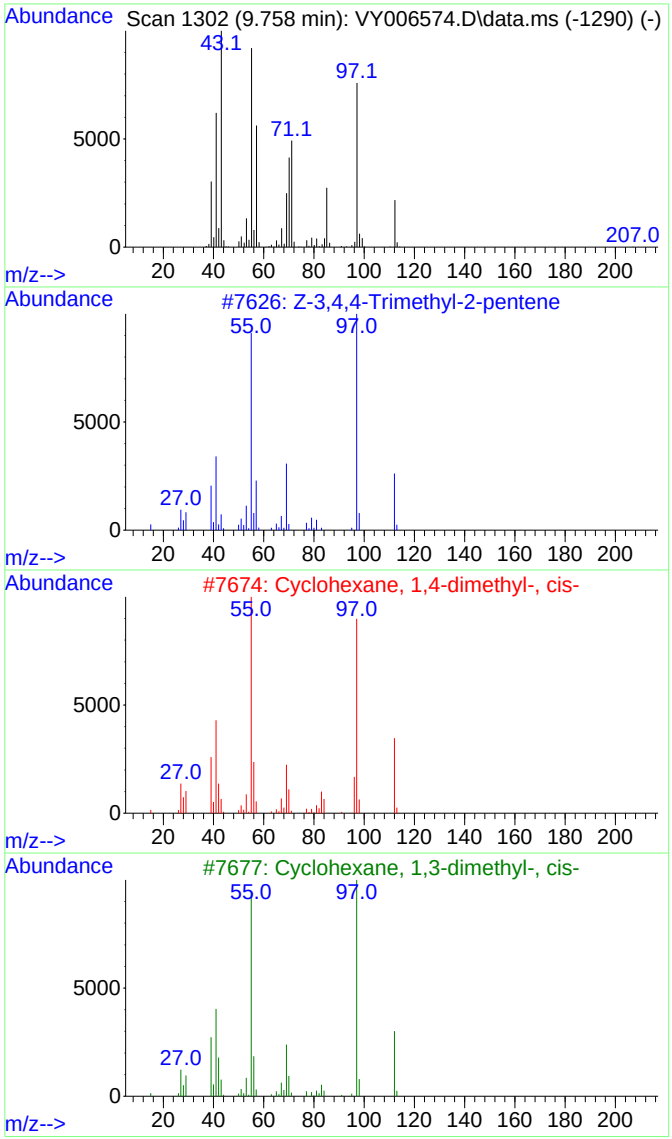
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y102221S.M
Quant Title : SW846 8260

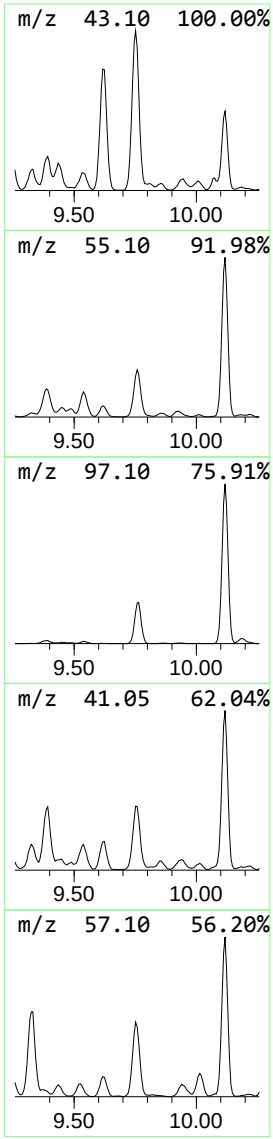
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Z-3,4,4-Trimethyl-2-pentene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.758	80.68 ug/l	879198	1,4-Difluorobenzene	8.697

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Z-3,4,4-Trimethyl-2-pentene	112	C8H16	039761-64-3	87
2	Cyclohexane, 1,4-dimethyl-, cis-	112	C8H16	000624-29-3	49
3	Cyclohexane, 1,3-dimethyl-, cis-	112	C8H16	000638-04-0	49
4	2-Pentene, 2,4,4-trimethyl-	112	C8H16	000107-40-4	46
5	2-Pentene, 2,3,4-trimethyl-	112	C8H16	000565-77-5	46





Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY110221\
Data File : VY006574.D
Acq On : 02 Nov 2021 13:53
Operator : SY/MD
Sample : M4427-02
Misc : 4.63G/5ML/MSVOA_Y/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
FDR-BH-5-20211029-6.5-7

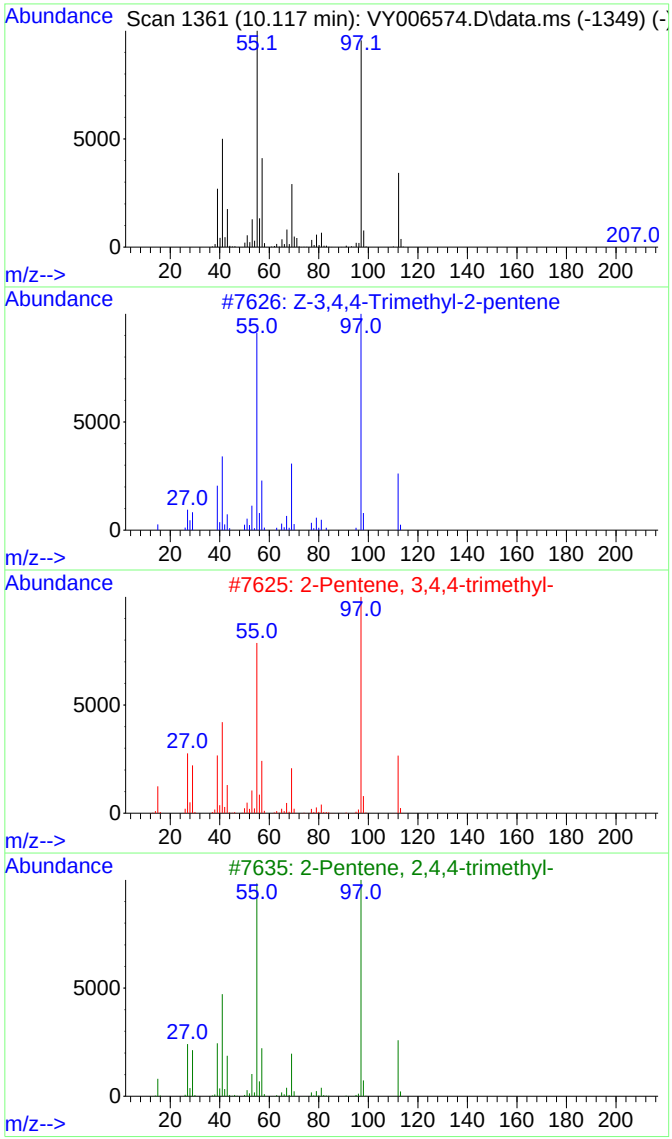
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y102221S.M
Quant Title : SW846 8260

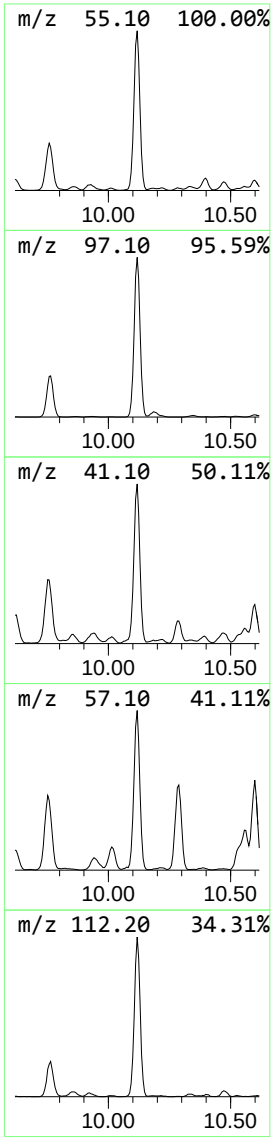
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 2-Pentene, 3,4,4-trimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.117	80.86 ug/l	1586900	Chlorobenzene-d5	11.495

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Z-3,4,4-Trimethyl-2-pentene	112	C8H16	039761-64-3	93
2	2-Pentene, 3,4,4-trimethyl-	112	C8H16	000598-96-9	91
3	2-Pentene, 2,4,4-trimethyl-	112	C8H16	000107-40-4	91
4	2-Pentene, 2,3,4-trimethyl-	112	C8H16	000565-77-5	90
5	Cyclohexane, 1,3-dimethyl-, cis-	112	C8H16	000638-04-0	80





Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY110221\
Data File : VY006574.D
Acq On : 02 Nov 2021 13:53
Operator : SY/MD
Sample : M4427-02
Misc : 4.63G/5ML/MSVOA_Y/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
FDR-BH-5-20211029-6.5-7

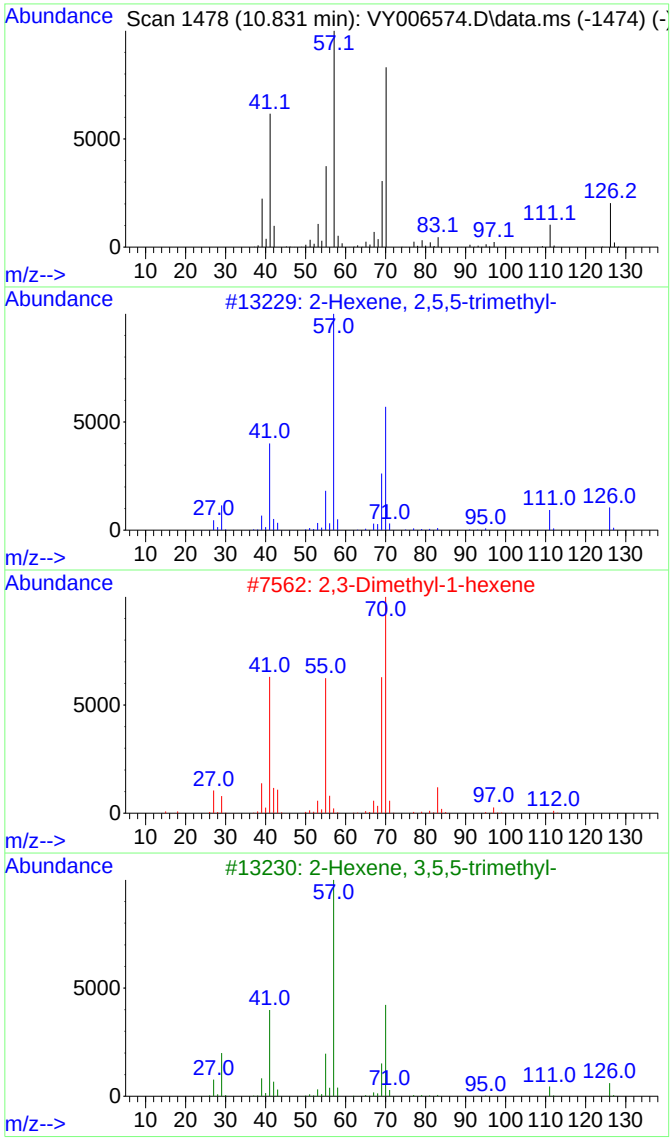
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y102221S.M
Quant Title : SW846 8260

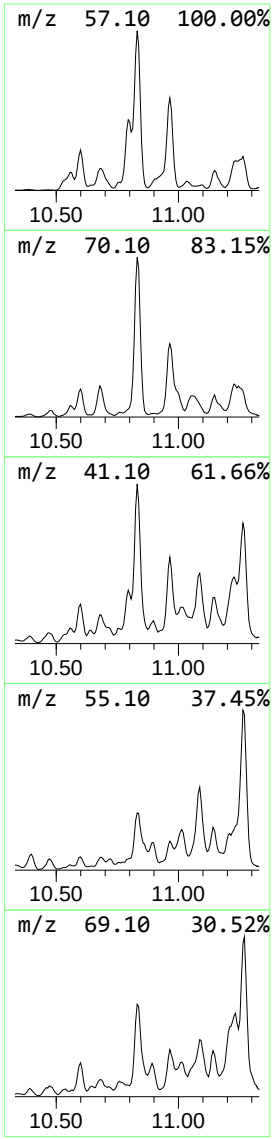
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 2-Hexene, 2,5,5-trimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.831	50.38 ug/l	988788	Chlorobenzene-d5	11.495

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexene, 2,5,5-trimethyl-	126	C9H18	040467-04-7	64
2	2,3-Dimethyl-1-hexene	112	C8H16	016746-86-4	38
3	2-Hexene, 3,5,5-trimethyl-	126	C9H18	026456-76-8	37
4	Carbonic acid, isobutyl 2-ethylh...	230	C13H26O3	1000383-13-3	36
5	Butane, 2-cyclopropyl-	98	C7H14	005750-02-7	33





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Data File : VY006574.D
Acq On : 02 Nov 2021 13:53
Operator : SY/MD
Sample : M4427-02
Misc : 4.63G/5ML/MSVOA_Y/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
FDR-BH-5-20211029-6.5-7

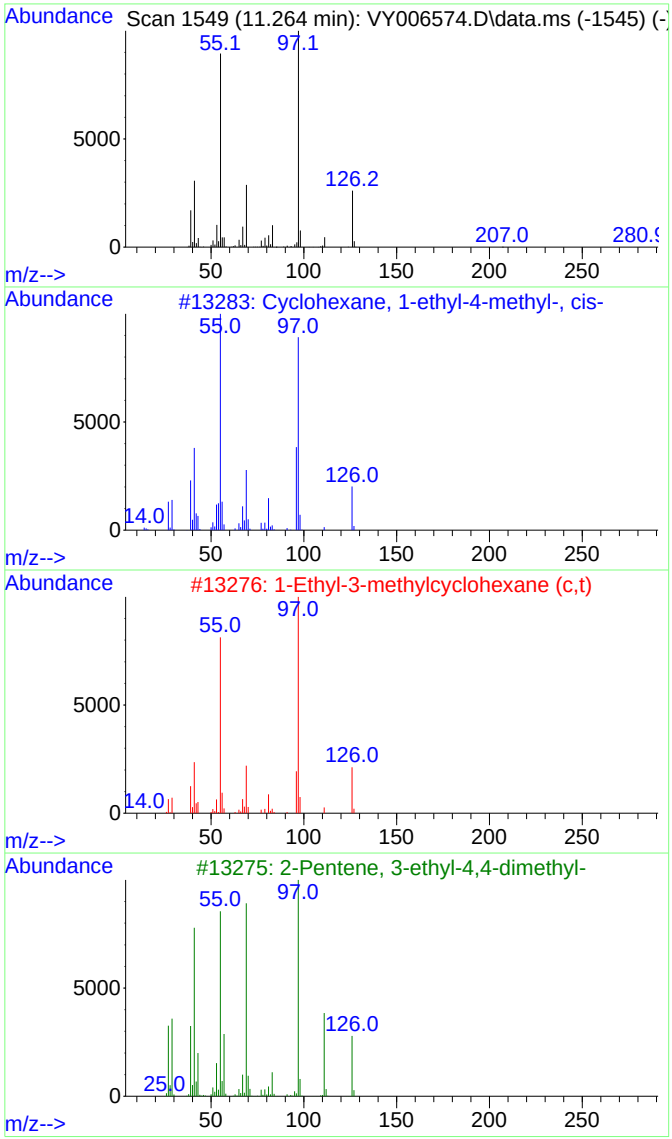
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y102221S.M
Quant Title : SW846 8260

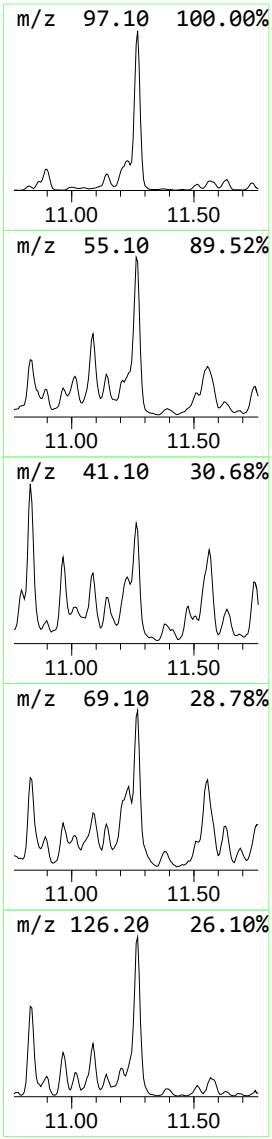
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Cyclohexane, 1-ethyl-4-meth... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.264	74.68 ug/l	1465570	Chlorobenzene-d5	11.495

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	004926-78-7	86
2	1-Ethyl-3-methylcyclohexane (c,t)	126	C9H18	003728-55-0	86
3	2-Pentene, 3-ethyl-4,4-dimethyl-	126	C9H18	053907-59-8	86
4	cis-1-Ethyl-3-methyl-cyclohexane	126	C9H18	019489-10-2	86
5	3,5-Dimethyl-3-heptene	126	C9H18	059643-68-4	80





Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY110221\
Data File : VY006574.D
Acq On : 02 Nov 2021 13:53
Operator : SY/MD
Sample : M4427-02
Misc : 4.63G/5ML/MSVOA_Y/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
FDR-BH-5-20211029-6.5-7

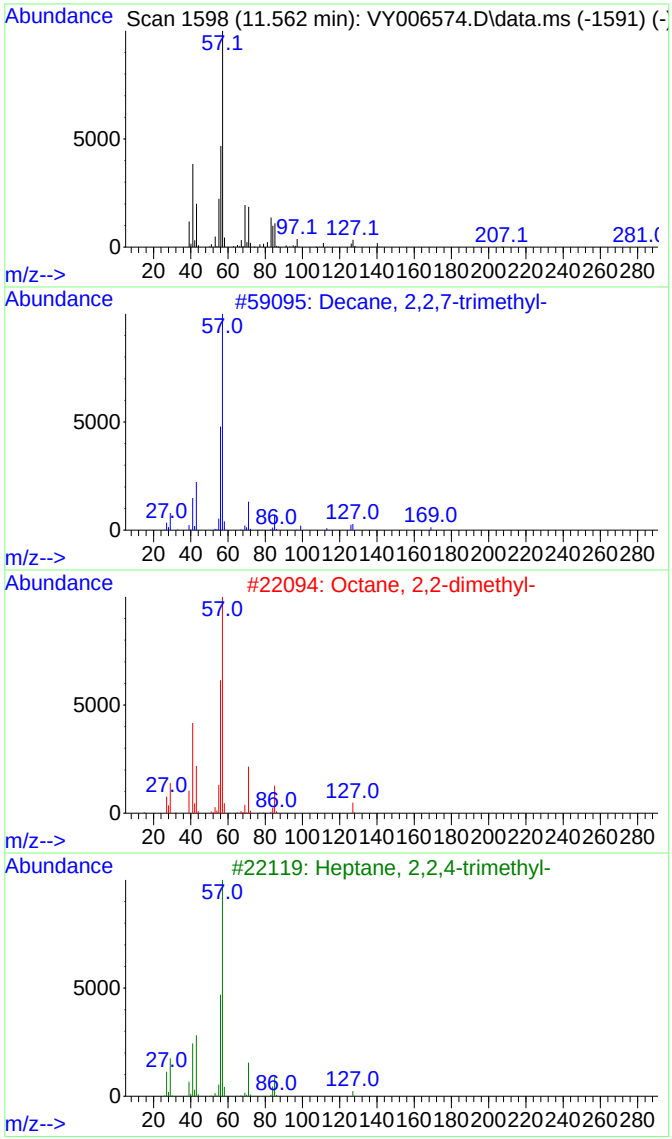
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y102221S.M
Quant Title : SW846 8260

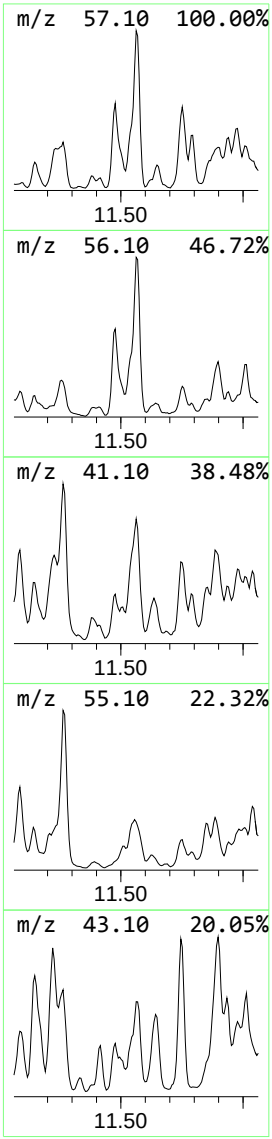
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Decane, 2,2,7-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.562	63.01 ug/l	1236600	Chlorobenzene-d5	11.495

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane, 2,2,7-trimethyl-	184	C13H28	062237-99-4	59
2	Octane, 2,2-dimethyl-	142	C10H22	015869-87-1	59
3	Heptane, 2,2,4-trimethyl-	142	C10H22	014720-74-2	59
4	Decane, 2,2,8-trimethyl-	184	C13H28	062238-01-1	56
5	Decane, 2,2,9-trimethyl-	184	C13H28	062238-00-0	53





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Data File : VY006574.D
Acq On : 02 Nov 2021 13:53
Operator : SY/MD
Sample : M4427-02
Misc : 4.63G/5ML/MSVOA_Y/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
FDR-BH-5-20211029-6.5-7

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

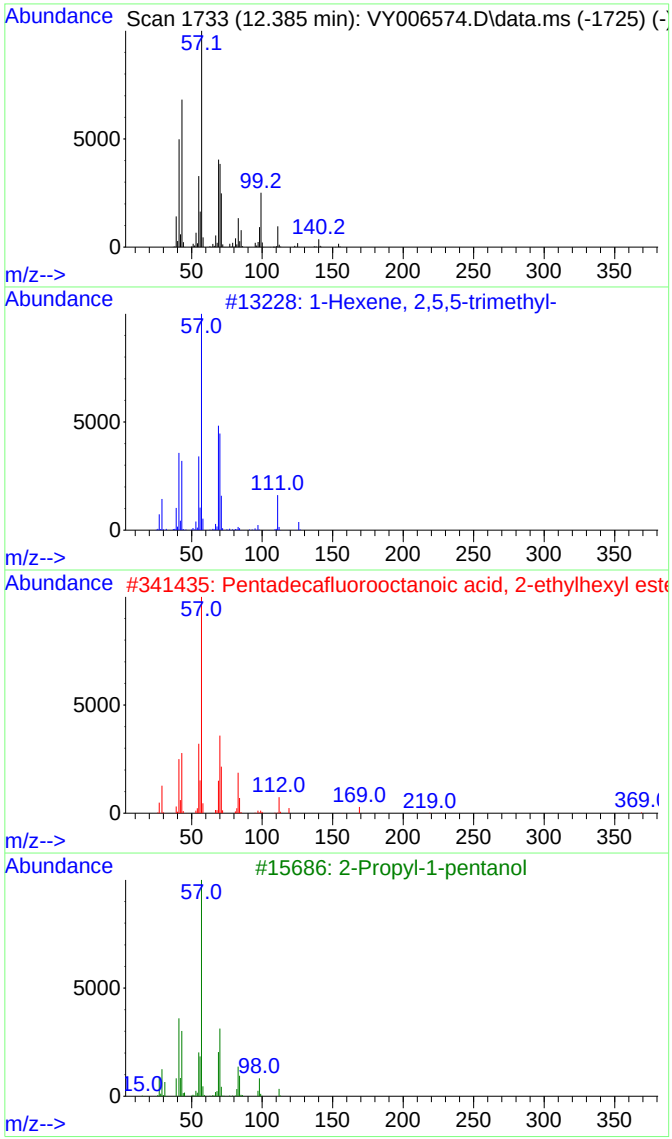
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

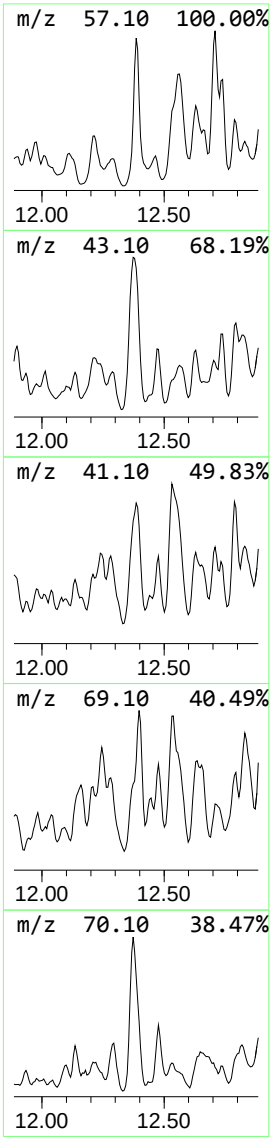
****

Peak Number 6 unknown12.385 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.385	101.82 ug/l	1998240	Chlorobenzene-d5	11.495

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexene, 2,5,5-trimethyl-	126	C9H18	062185-56-2	47
2	Pentadecafluorooctanoic acid, 2-...	526	C16H17F15O2	1000406-80-9	47
3	2-Propyl-1-pentanol	130	C8H18O	058175-57-8	47
4	Heptane, 2-methyl-	114	C8H18	000592-27-8	45
5	Heptane, 3,5-dimethyl-	128	C9H20	000926-82-9	43





Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY110221\
Data File : VY006574.D
Acq On : 02 Nov 2021 13:53
Operator : SY/MD
Sample : M4427-02
Misc : 4.63G/5ML/MSVOA_Y/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
FDR-BH-5-20211029-6.5-7

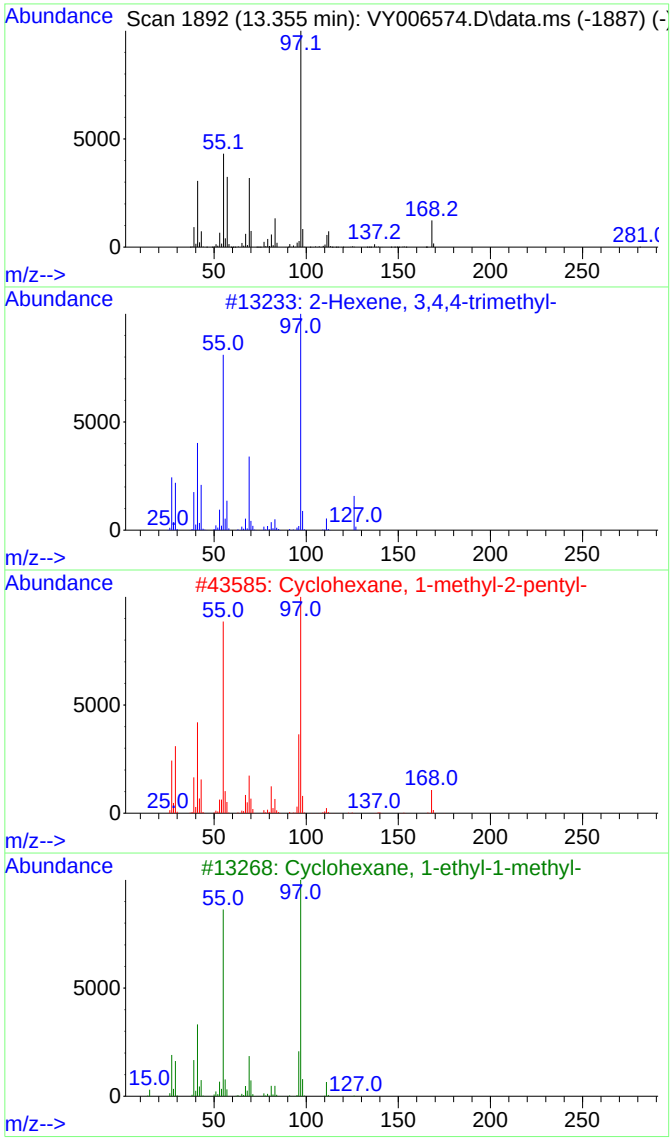
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y102221S.M
Quant Title : SW846 8260

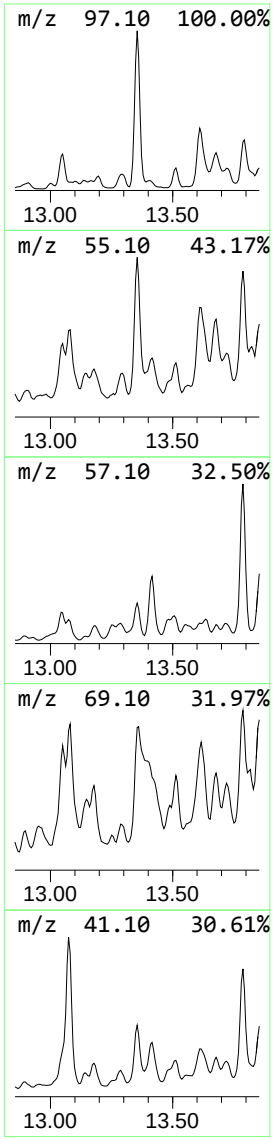
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 2-Hexene, 3,4,4-trimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.355	63.67 ug/l	4705680	1,4-Dichlorobenzene-d4	13.428

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexene, 3,4,4-trimethyl-	126	C9H18	053941-19-8	64
2	Cyclohexane, 1-methyl-2-pentyl-	168	C12H24	054411-01-7	64
3	Cyclohexane, 1-ethyl-1-methyl-	126	C9H18	004926-90-3	64
4	Thiophene, 2-hexyl-	168	C10H16S	018794-77-9	59
5	Cyclohexane, 1-methyl-3-pentyl-	168	C12H24	054411-02-8	56





Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY110221\
Data File : VY006574.D
Acq On : 02 Nov 2021 13:53
Operator : SY/MD
Sample : M4427-02
Misc : 4.63G/5ML/MSVOA_Y/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
FDR-BH-5-20211029-6.5-7

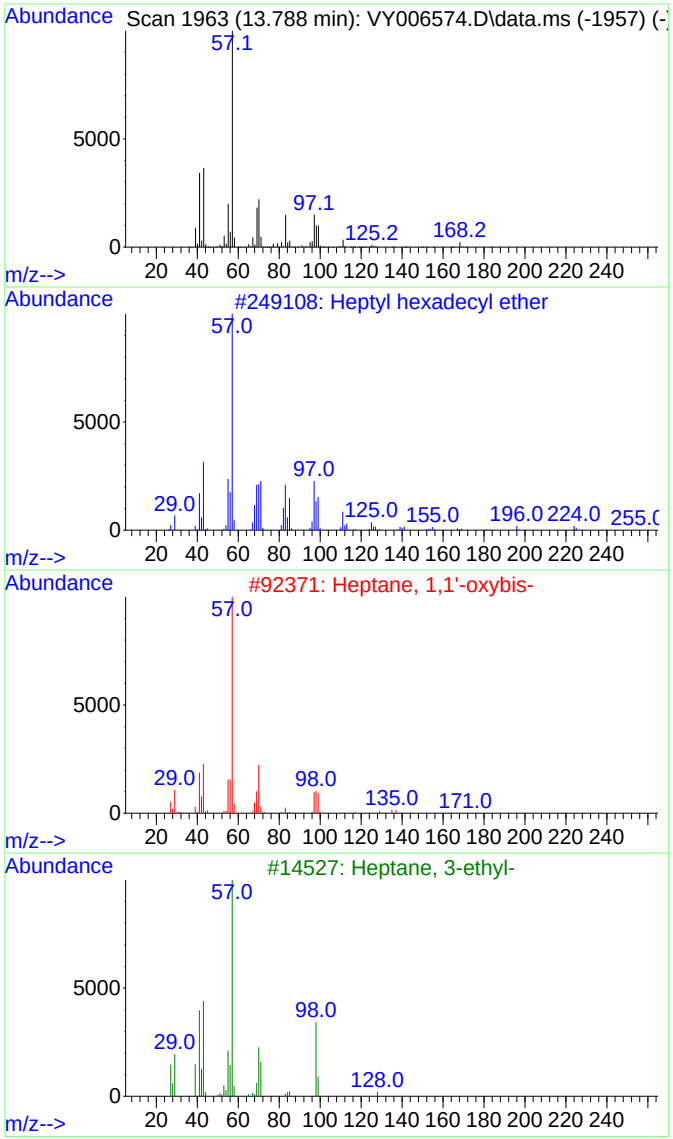
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y102221S.M
Quant Title : SW846 8260

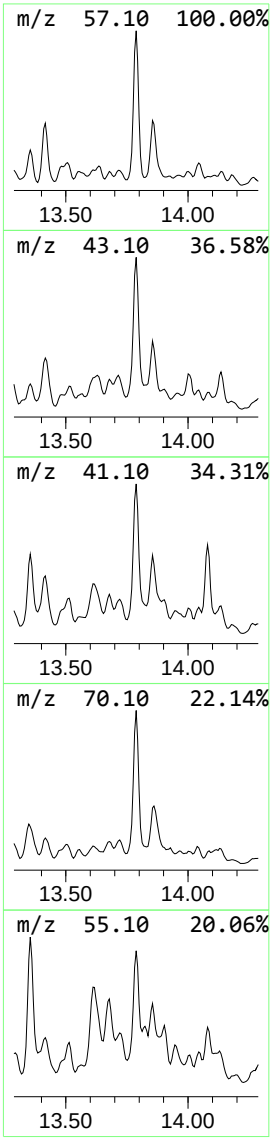
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Heptyl hexadecyl ether Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.788	109.06 ug/l	8060600	1,4-Dichlorobenzene-d4	13.428

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptyl hexadecyl ether	340	C23H48O	1000406-39-4	56
2	Heptane, 1,1'-oxybis-	214	C14H30O	000629-64-1	50
3	Heptane, 3-ethyl-	128	C9H20	015869-80-4	43
4	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	40
5	Octane, 2,3-dimethyl-	142	C10H22	007146-60-3	38





Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY110221\
Data File : VY006574.D
Acq On : 02 Nov 2021 13:53
Operator : SY/MD
Sample : M4427-02
Misc : 4.63G/5ML/MSVOA_Y/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
FDR-BH-5-20211029-6.5-7

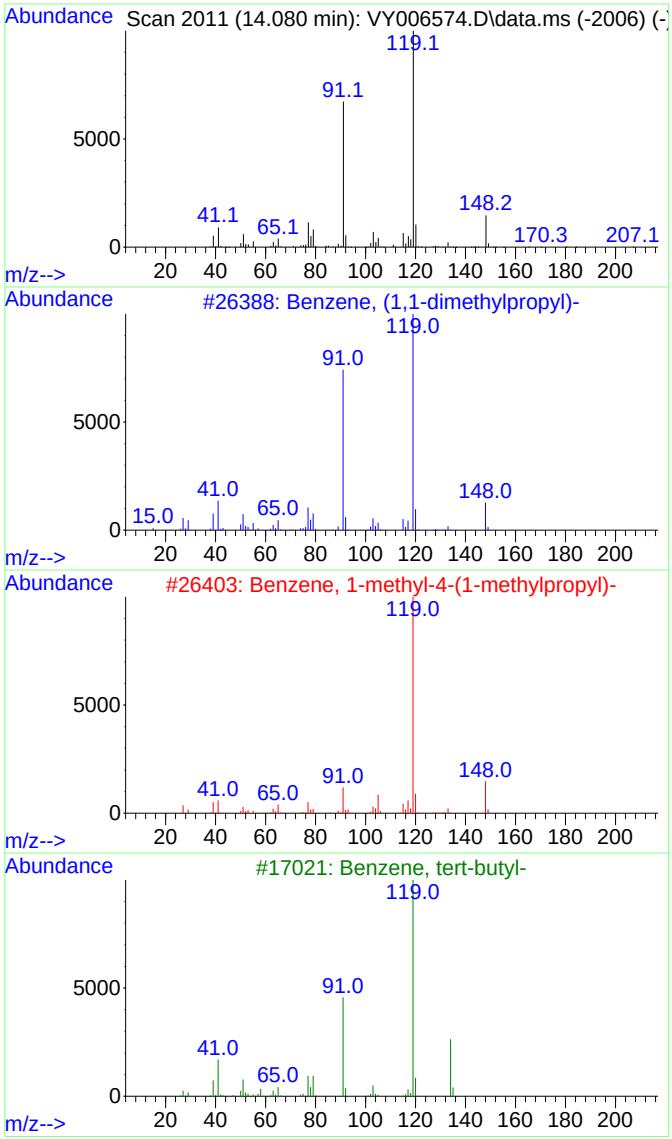
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y102221S.M
Quant Title : SW846 8260

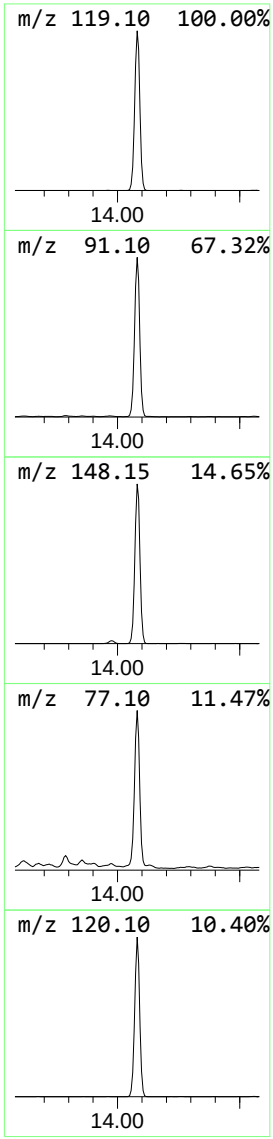
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzene, (1,1-dimethylpropyl)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.080	156.90 ug/l	11596900	1,4-Dichlorobenzene-d4	13.428

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	97
2	Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	90
3	Benzene, tert-butyl-	134	C10H14	000098-06-6	86
4	Benzene, (1-ethylpropyl)-	148	C11H16	001196-58-3	86
5	Propiophenone, 2'-methyl-	148	C10H12O	002040-14-4	72





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Data File : VY006574.D
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Operator : SY/MD
Sample : M4427-02
Misc : 4.63G/5ML/MSVOA_Y/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
FDR-BH-5-20211029-6.5-7

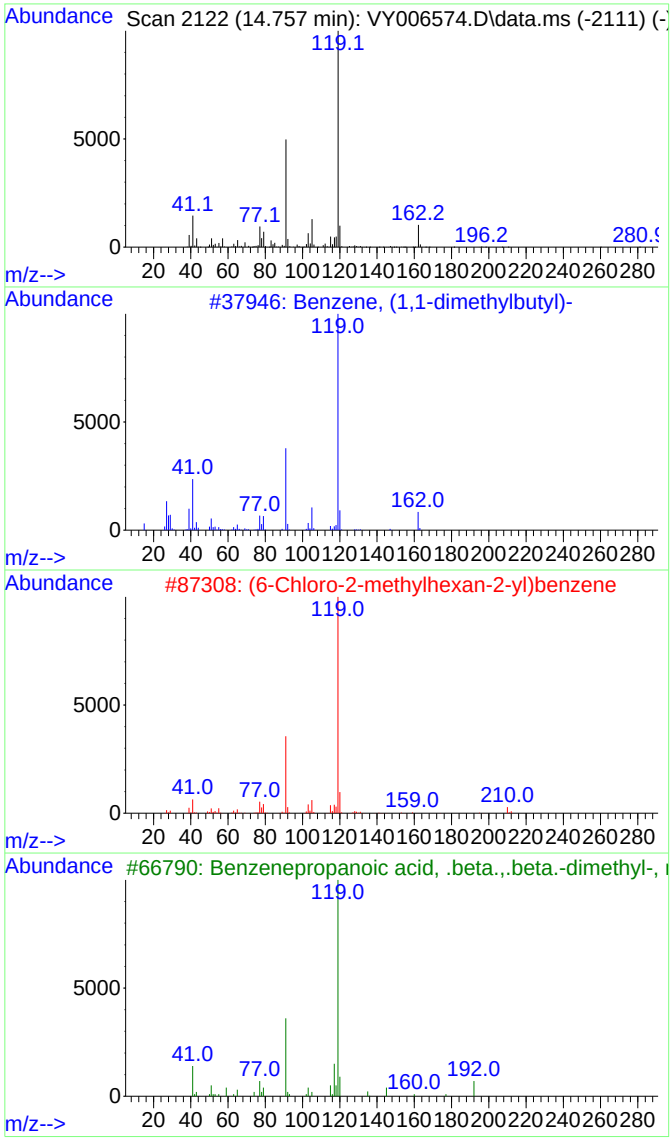
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y102221S.M
Quant Title : SW846 8260

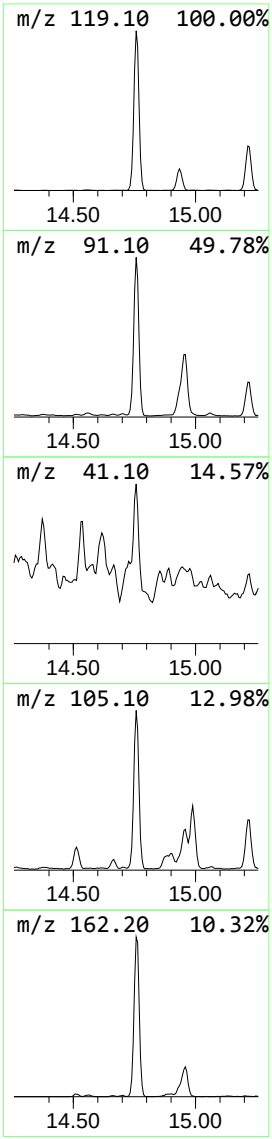
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Benzene, (1,1-dimethylbutyl)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.757	46.19 ug/l	3413760	1,4-Dichlorobenzene-d4	13.428

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1,1-dimethylbutyl)-	162	C12H18	001985-57-5	91
2	(6-Chloro-2-methylhexan-2-yl)ben...	210	C13H19Cl	1417621-45-4	86
3	Benzenepropanoic acid, .beta.,.b...	192	C12H16O2	025080-84-6	72
4	Benzene, (1,1-dimethyldecyl)-	246	C18H30	027854-40-6	72
5	Benzene, tert-butyl-	134	C10H14	000098-06-6	72





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY110221\
Data File : VY006574.D
Acq On : 02 Nov 2021 13:53
Operator : SY/MD
Sample : M4427-02
Misc : 4.63G/5ML/MSVOA_Y/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
FDR-BH-5-20211029-6.5-7

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
Z-3,4,4-Trimeth...	9.758	80.7	ug/l	879198	2	8.697	544878	50.0
2-Pentene, 3,4,...	10.117	80.9	ug/l	1586900	3	11.495	981272	50.0
2-Hexene, 2,5,5...	10.831	50.4	ug/l	988788	3	11.495	981272	50.0
Cyclohexane, 1-...	11.264	74.7	ug/l	1465570	3	11.495	981272	50.0
Decane, 2,2,7-t...	11.562	63.0	ug/l	1236600	3	11.495	981272	50.0
unknown12.385	12.385	101.8	ug/l	1998240	3	11.495	981272	50.0
2-Hexene, 3,4,4...	13.355	63.7	ug/l	4705680	4	13.428	3695590	50.0
Heptyl hexadecy...	13.788	109.1	ug/l	8060600	4	13.428	3695590	50.0
Benzene, (1,1-d...	14.080	156.9	ug/l	11596900	4	13.428	3695590	50.0
Benzene, (1,1-d...	14.757	46.2	ug/l	3413760	4	13.428	3695590	50.0