LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110419\
Data File : VX013324.D
Acq On : 04 Nov 2019 15:07
Operator : JC/SP
Sample : K5683-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : OILY-WATER

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_X\METHOD\82X103119W.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	1.064	25	29	40	rBV	2373682	2910729	24.06%	2.570%
2	1.265	58	62	67	rBV	100804	155195	1.28%	0.137%
3	1.479	93	97	102	rBV	119880	129893	1.07%	0.115%
4	2.424	246	252	268	rBV	3355738	5409568	44.71%	4.776%
5	2.588	269	279	289	rVV	607819	1121011	9.27%	0.990%
6	3.015	339	349	358	rVB2	105071	272966	2.26%	0.241%
7	4.319	555	563	571	rBV2	49203	131279	1.09%	0.116%
8	4.649	605	617	630	rBV	641466	1825367	15.09%	1.611%
9	5.478	743	753	764	rBV	410458	1171339	9.68%	1.034%
10	5.642	770	780	794	rVB	740908	2049548	16.94%	1.809%
11	6.045	836	846	860	rBV	456607	1202863	9.94%	1.062%
12	6.203	865	872	881	rBV2	49222	134337	1.11%	0.119%
13	6.715	942	956	968	rBV	355060	959282	7.93%	0.847%
14	6.843	968	977	989	rVB	1149570	2649054	21.89%	2.339%
15	7.282	1041	1049	1057	rBV2	71275	165991	1.37%	0.147%
16	7.526	1082	1089	1098	rVB	705738	1372208	11.34%	1.211%
17	7.709	1114	1119	1129	rVB	95882	181484	1.50%	0.160%
18	7.965	1152	1161	1168	rBV2	105016	216145	1.79%	0.191%
19	8.630	1264	1270	1277	rBV	3784681	5974023	49.37%	5.274%
20	8.709	1277	1283	1289	rVB	2361348	3690770	30.50%	3.258%
21	8.770	1289	1293	1298	rBV2	149869	222827	1.84%	0.197%
22	9.081	1336	1344	1350	rBV	71311	127311	1.05%	0.112%
23	9.288	1373	1378	1382	rBV2	106759	178587	1.48%	0.158%
24	9.361	1386	1390	1400	rVB	151096	249658	2.06%	0.220%
25	9.483	1405	1410	1418	rBV	617374	843544	6.97%	0.745%
26	9.575	1419	1425	1428	rBV	102672	146504	1.21%	0.129%
27	9.995	1489	1494	1500	rBV	581485	799886	6.61%	0.706%
28	10.105	1502	1512	1517	rBV	2343989	3238252	26.76%	2.859%
29	10.264	1532	1538	1540	rBV4	113070	194012	1.60%	0.171%
30	10.294	1540	1543	1546	rVB	106294	134691	1.11%	0.119%
31	10.434	1563	1566	1569	rBV	186597	225222	1.86%	0.199%
32	10.556	1581	1586	1590	rBV2	88923	122379	1.01%	0.108%
33	10.629	1594	1598	1606	rBV	111033	193443	1.60%	0.171%
34	10.733	1611	1615	1622	rVB	176154	238875	1.97%	0.211%

LSC Area Percent Report

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Operator : JC/SP
Sample : K5683-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : OILY-WATER

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_X\METHOD\82X103119W.M
Title : SW846 8260

35	10.818	1623	1629	1635	rBV	724257	933206	7.71%	0.824%
36	10.995	1654	1658	1667	rVB4	65214	123859	1.02%	0.109%
37	11.129	1675	1680	1687	rVB	1990051	2511667	20.76%	2.217%
38	11.361	1713	1718	1722	rVB2	204824	297674	2.46%	0.263%
39	11.477	1733	1737	1740	rBV3	102954	126940	1.05%	0.112%
40	11.526	1740	1745	1753	rVV	1127839	1409415	11.65%	1.244%
41	11.605	1755	1758	1762	rVV	111451	134559	1.11%	0.119%
42	11.666	1762	1768	1772	rVV2	389026	574939	4.75%	0.508%
43	11.708	1772	1775	1779	rVB3	89611	141400	1.17%	0.125%
44	11.806	1784	1791	1795	rBV	410921	517353	4.28%	0.457%
45	11.855	1795	1799	1801	rBV2	121697	157965	1.31%	0.139%
46	11.885	1801	1804	1810	rVB	1148773	1303606	10.77%	1.151%
47	11.946	1810	1814	1816	rBV	263452	336387	2.78%	0.297%
48	11.977	1816	1819	1824	rVB	1028432	1136239	9.39%	1.003%
49	12.074	1829	1835	1840	rVB	2648485	3279141	27.10%	2.895%
50	12.153	1842	1848	1859	rBV	9536839	12099405	100.00%	10.681%
51	12.263	1862	1866	1874	rBV	717458	1101317	9.10%	0.972%
52	12.397	1884	1888	1892	rVB	210864	277792	2.30%	0.245%
53	12.470	1896	1900	1902	rBV	163213	189510	1.57%	0.167%
54	12.507	1902	1906	1912	rVB4	223880	406891	3.36%	0.359%
55	12.562	1912	1915	1919	rBV	152899	185007	1.53%	0.163%
56	12.629	1919	1926	1939	rVB4	616424	1258239	10.40%	1.111%
57	12.739	1939	1944	1949	rBV	155314	216634	1.79%	0.191%
58	12.818	1952	1957	1960	rBV	492195	639189	5.28%	0.564%
59	12.903	1967	1971	1976	rVB	922251	1038537	8.58%	0.917%
60	13.019	1985	1990	2000	rBV	3073453	3722093	30.76%	3.286%
61	13.117	2002	2006	2008	rBV	3149456	3521100	29.10%	3.108%
62	13.147	2008	2011	2020	rVB	6302233	7653602	63.26%	6.757%
63	13.312	2035	2038	2049	rVB3	254894	487497	4.03%	0.430%
64	13.470	2060	2064	2074	rBV	10289312	11813025	97.63%	10.428%
65	13.562	2074	2079	2089	rVB	3484305	4634769	38.31%	4.092%
66	13.653	2091	2094	2101	rVB2	197443	301640	2.49%	0.266%
67	13.757	2104	2111	2115	rBV	1205643	2088141	17.26%	1.843%
68	13.800	2115	2118	2119	rVV	506043	616208	5.09%	0.544%
69	13.824	2119	2122	2129	rVV2	1803864	2148302	17.76%	1.897%
70	13.909	2131	2136	2138	rVV2	451388	668699	5.53%	0.590%
71	13.940	2138	2141	2152	rVB	735496	1156051	9.55%	1.021%

LSC Area Percent Report

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Operator : JC/SP
Sample : K5683-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : OILY-WATER

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_X\METHOD\82X103119W.M
Title : SW846 8260

72	14.245	2187	2191	2199	rBV3	543751	821181	6.79%	0.725%
73	14.458	2222	2226	2230	rBV3	462527	686868	5.68%	0.606%
74	14.574	2243	2245	2248	rVB2	319811	303455	2.51%	0.268%
75	14.671	2257	2261	2267	rVB4	378062	650147	5.37%	0.574%
76	14.805	2281	2283	2286	rBV	381000	376961	3.12%	0.333%
77	14.848	2287	2290	2293	rVB4	260500	289049	2.39%	0.255%
78	15.269	2357	2359	2363	rVB5	378476	430203	3.56%	0.380%
79	15.482	2391	2394	2401	rVB	638591	906017	7.49%	0.800%
80	15.811	2446	2448	2457	rVB	678209	966356	7.99%	0.853%

Sum of corrected areas: 113276478

LSC Report - Integrated Chromatogram

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110419\
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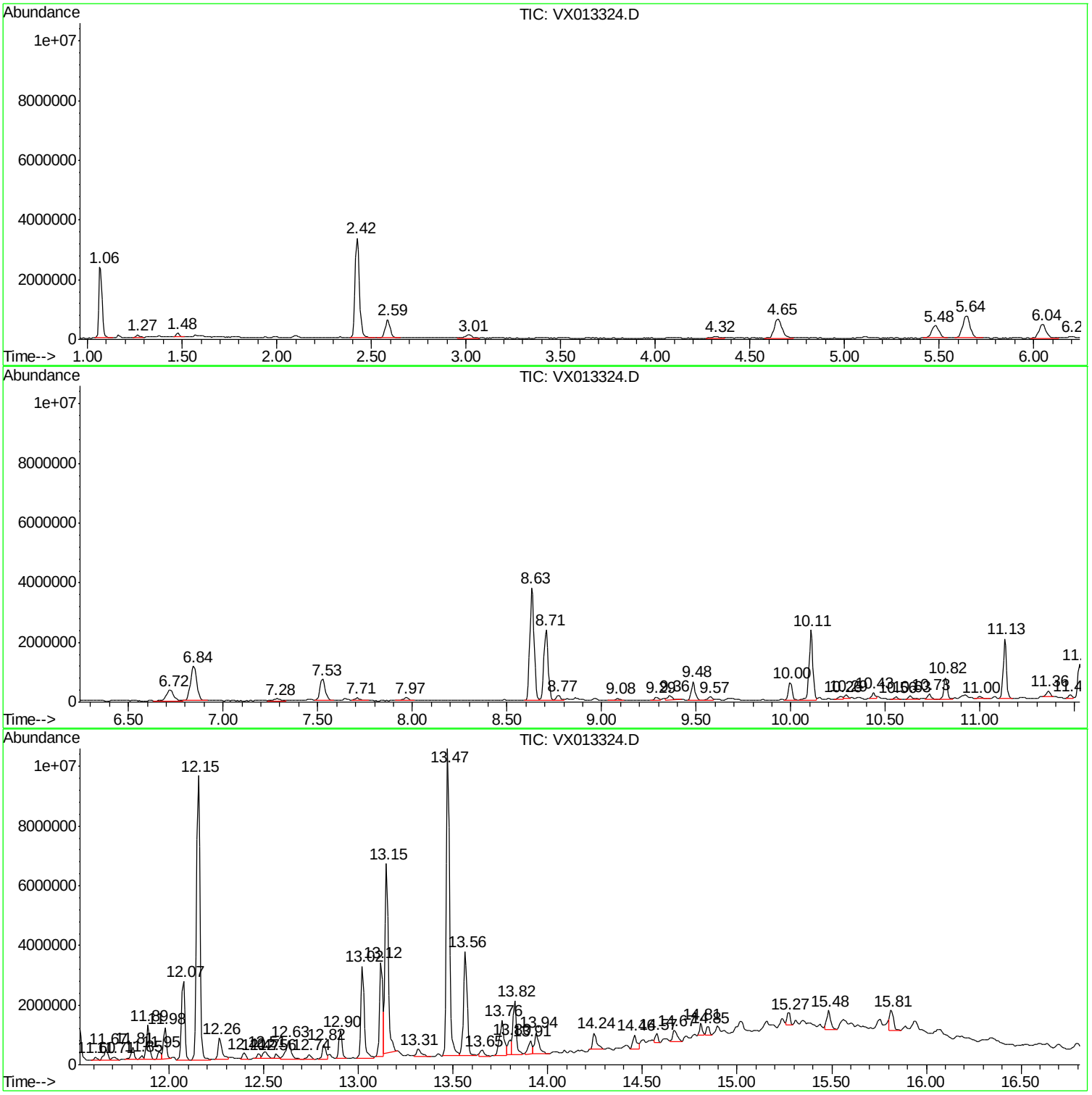
Acq On : 04 Nov 2019 15:07
Operator : JC/SP
Sample : K5683-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
OILY-WATER

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X103119W.M
Quant Title : SW846 8260

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



Library Search Compound Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110419\
Data File : VX013324.D
Acq On : 04 Nov 2019 15:07
Operator : JC/SP
Sample : K5683-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : OILY-WATER

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X103119W.M
Quant Title : SW846 8260

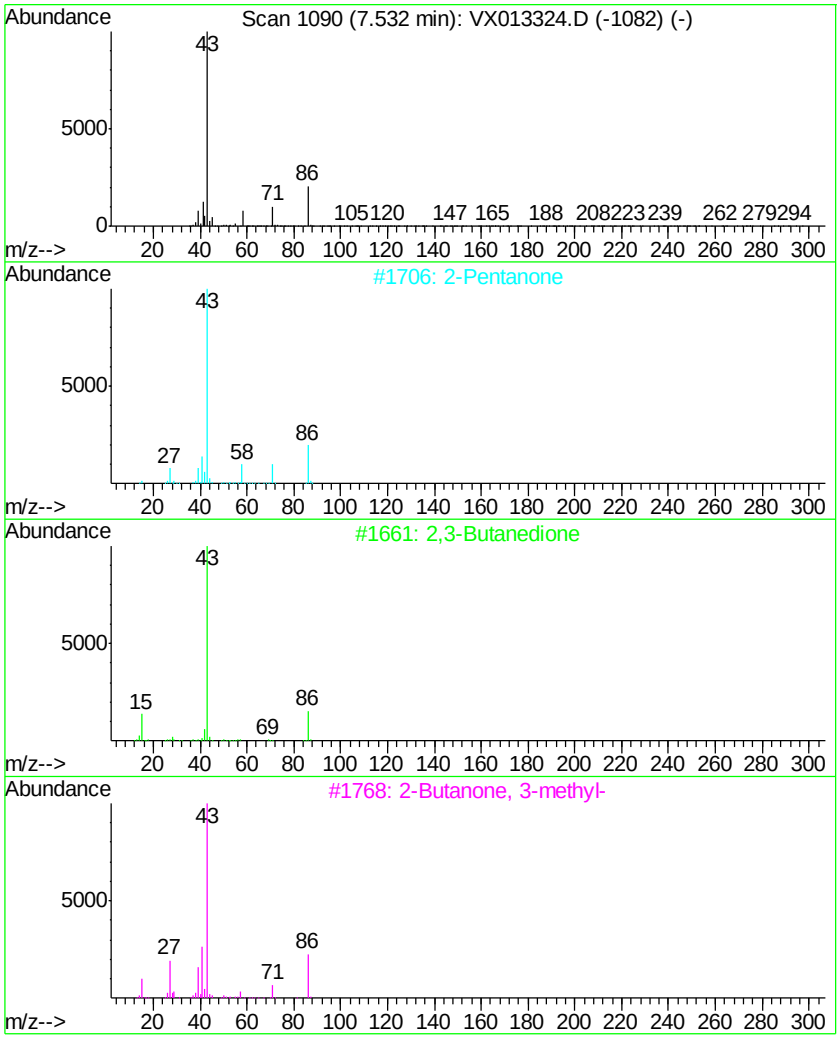
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

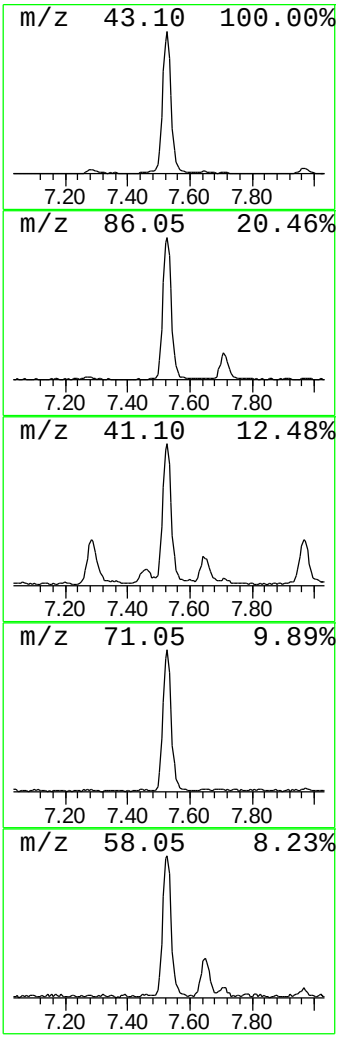
****

Peak Number 2 2-Pentanone Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.53	25.90 ug/l	1372210	1,4-Difluorobenzene	6.84

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone	86	C5H10O	000107-87-9	91
2	2,3-Butanedione	86	C4H6O2	000431-03-8	9
3	2-Butanone, 3-methyl-	86	C5H10O	000563-80-4	9
4	Cyclobutylamine	71	C4H9N	002516-34-9	9
5	2-Pentanone, 3-ethyl-	114	C7H14O	006137-03-7	9





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110419\
Data File : VX013324.D
Acq On : 04 Nov 2019 15:07
Operator : JC/SP
Sample : K5683-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : OILY-WATER

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X103119W.M
Quant Title : SW846 8260

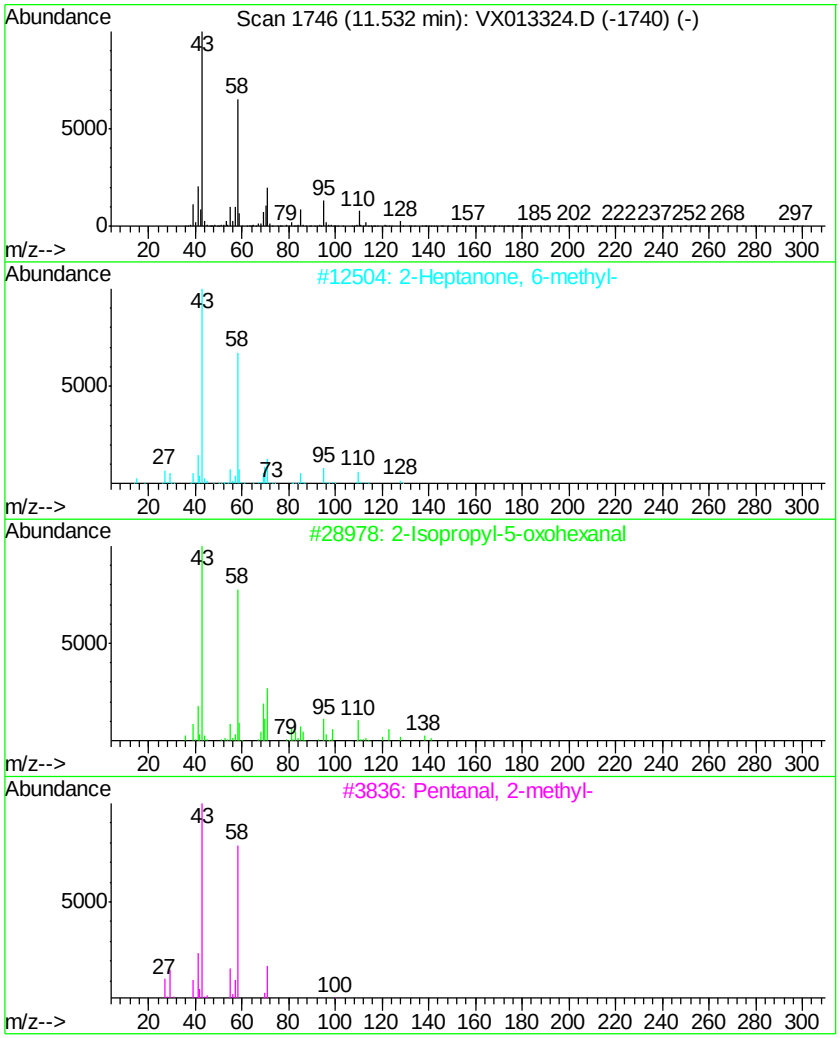
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

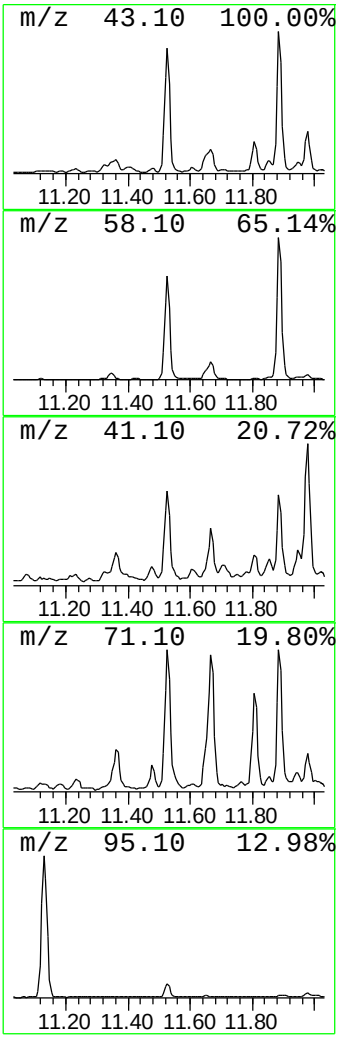
****

Peak Number 3 2-Heptanone, 6-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.53	21.49 ug/l	1409420	1,4-Dichlorobenzene-d4	12.07

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Heptanone, 6-methyl-	128	C8H16O	000928-68-7	90
2	2-Isopropyl-5-oxohexanal	156	C9H16O2	015303-46-5	64
3	Pentanal, 2-methyl-	100	C6H12O	000123-15-9	53
4	Butanimidamide	86	C4H10N2	000107-90-4	50
5	2-Octanone	128	C8H16O	000111-13-7	45





Library Search Compound Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110419\
Data File : VX013324.D
Acq On : 04 Nov 2019 15:07
Operator : JC/SP
Sample : K5683-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : OILY-WATER

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X103119W.M
Quant Title : SW846 8260

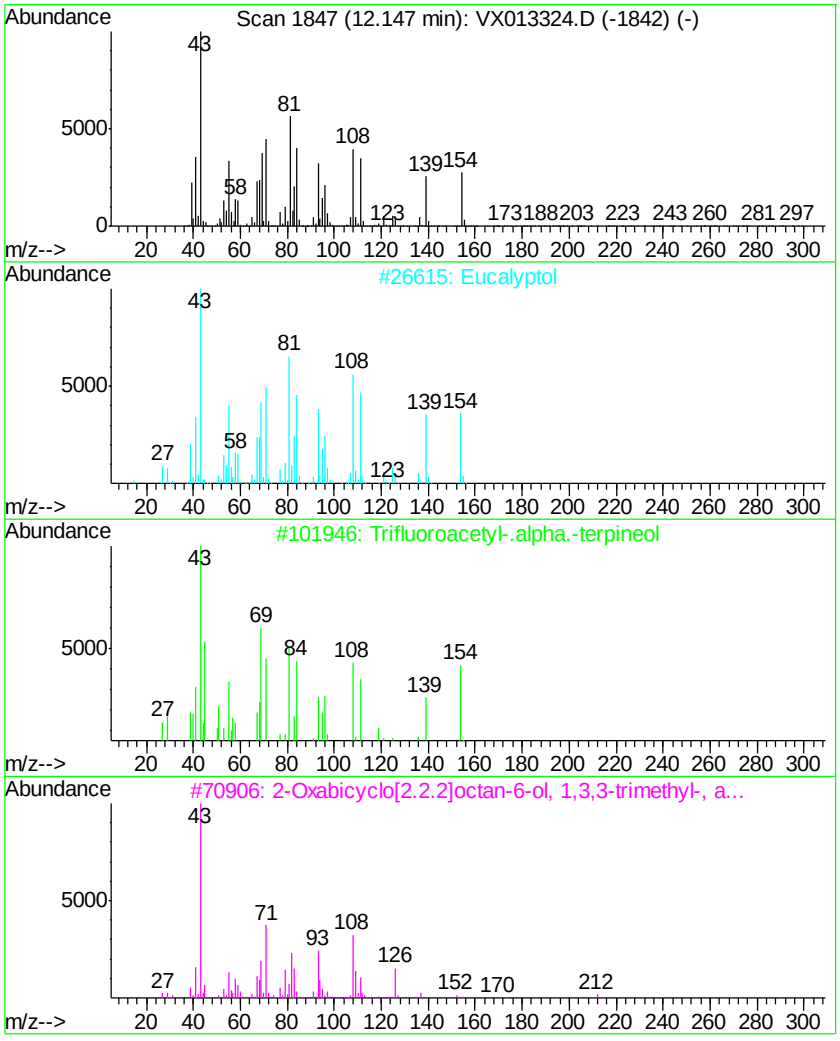
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

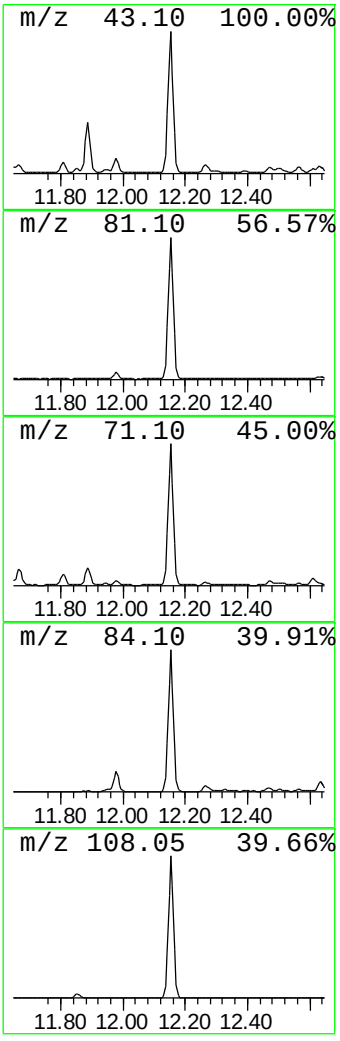
****

Peak Number 4 Eucalyptol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.15	184.49 ug/l	12099400	1,4-Dichlorobenzene-d4	12.07

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eucalyptol	154	C10H18O	000470-82-6	99
2	Trifluoroacetyl-.alpha.-terpineol	250	C12H17F3O2	1000058-17-6	68
3	2-Oxabicyclo[2.2.2]octan-6-ol, 1...	212	C12H20O3	057709-95-2	32
4	2-Isopropyl-5-methylhex-2-enal	154	C10H18O	035158-25-9	27
5	3-Hepten-2-one, 3-propyl-	154	C10H18O	032064-69-0	25





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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X103119W.M
Quant Title : SW846 8260

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

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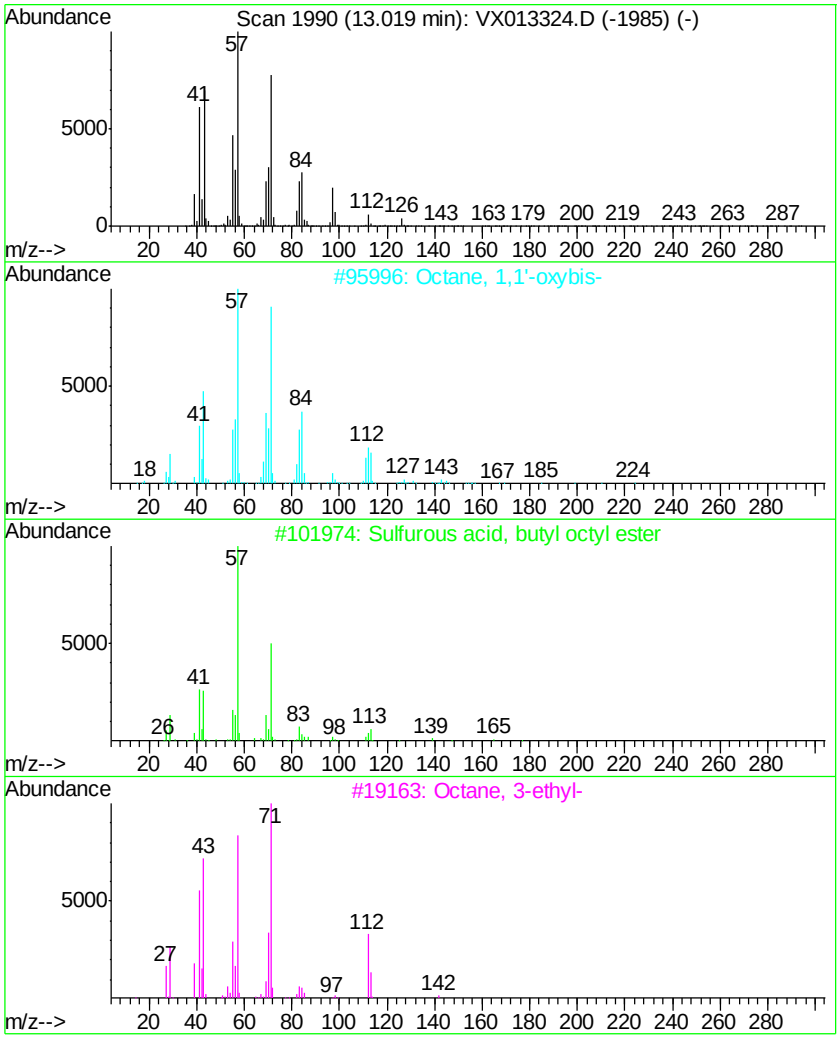
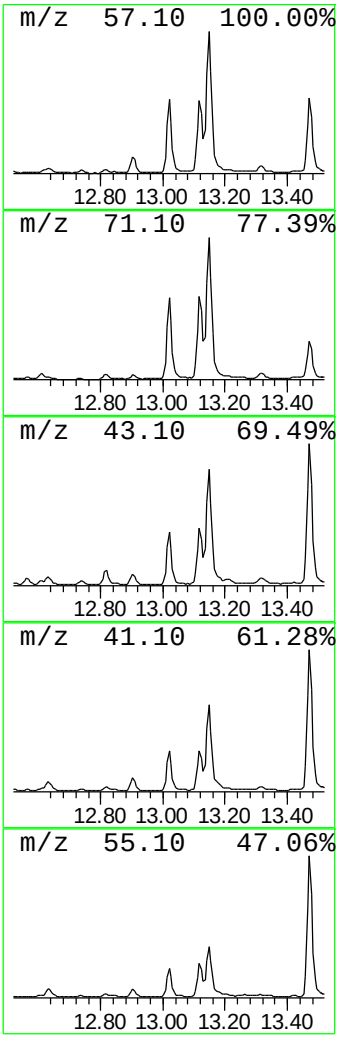
Peak Number 5 Octane, 1,1'-oxybis- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.02	56.75 ug/l	3722090	1,4-Dichlorobenzene-d4	12.07

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 1,1'-oxybis-	242	C16H34O	000629-82-3	80
2	Sulfurous acid, butyl octyl ester	250	C12H26O3S	1000309-17-5	50
3	Octane, 3-ethyl-	142	C10H22	005881-17-4	50
4	Isobutyl octyl carbonate	230	C13H26O3	1000372-74-6	49
5	1-Decene, 4-methyl-	154	C11H22	013151-29-6	47





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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : OILY-WATER

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X103119W.M
Quant Title : SW846 8260

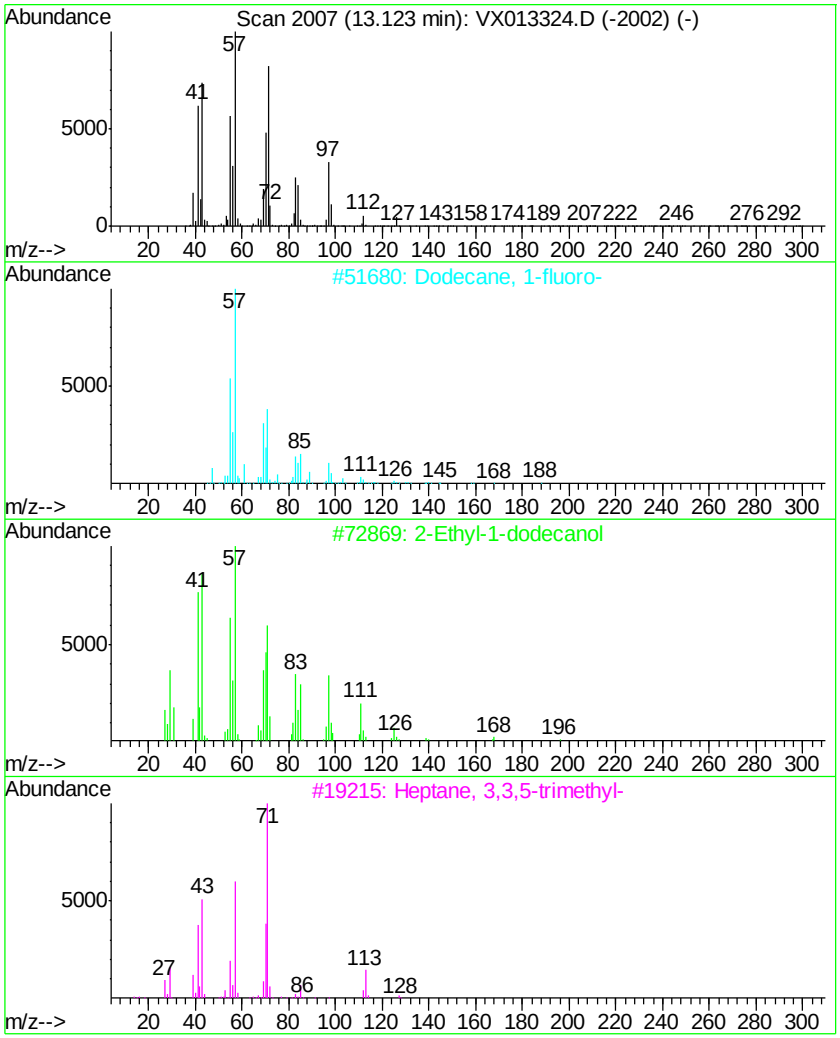
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

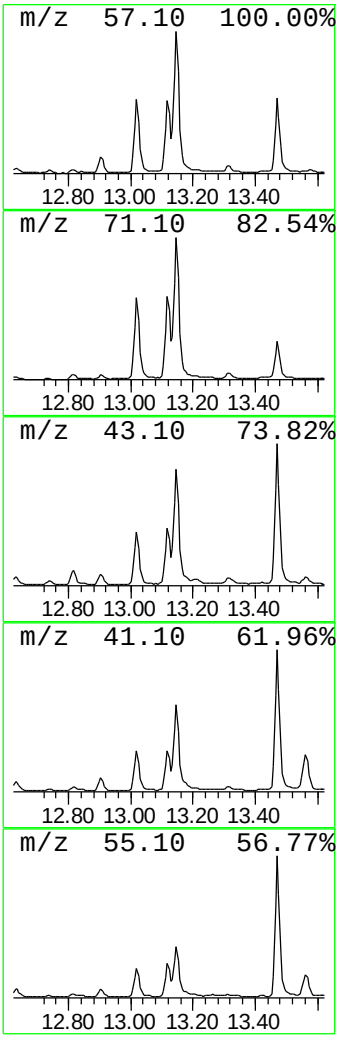
****

Peak Number 6 Dodecane, 1-fluoro- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.12	53.69 ug/l	3521100	1,4-Dichlorobenzene-d4	12.07

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane, 1-fluoro-	188	C12H25F	000334-68-9	64
2	2-Ethyl-1-dodecanol	214	C14H30O	019780-33-7	59
3	Heptane, 3,3,5-trimethyl-	142	C10H22	007154-80-5	47
4	n-Butyric acid 2-ethylhexyl ester	200	C12H24O2	025415-84-3	47
5	2-Ethyl-1-hexanol, pentafluoropr...	276	C11H17F5O2	1000365-51-1	43





Library Search Compound Report

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Sample : K5683-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : OILY-WATER

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X103119W.M
Quant Title : SW846 8260

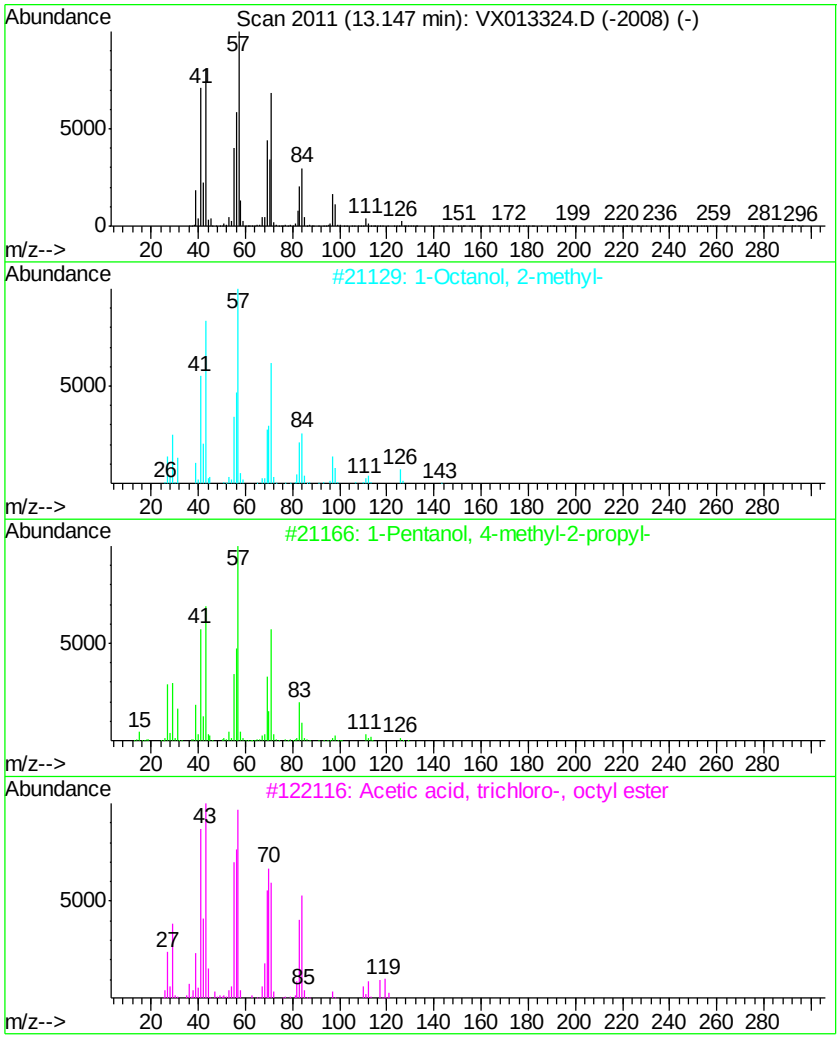
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

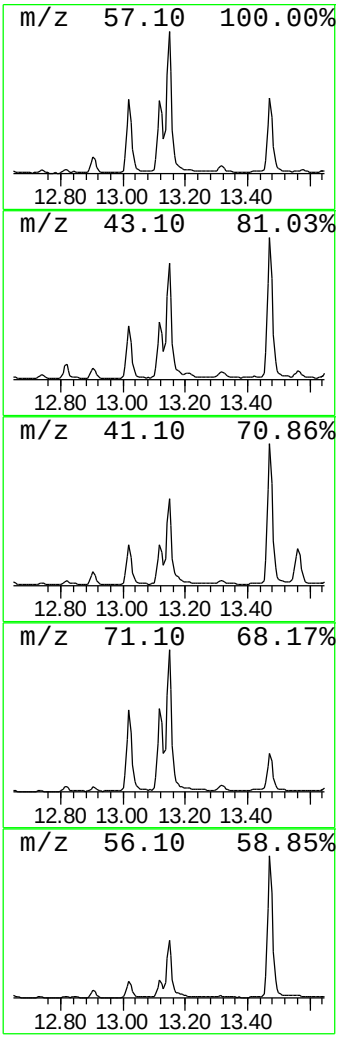
****

Peak Number 7 1-Octanol, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.15	116.70 ug/l	7653600	1,4-Dichlorobenzene-d4	12.07

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Octanol, 2-methyl-	144	C9H20O	000818-81-5	90
2	1-Pentanol, 4-methyl-2-propyl-	144	C9H20O	054004-41-0	59
3	Acetic acid, trichloro-, octyl e...	274	C10H17Cl3O2	016958-78-4	53
4	Cyclobutane, 1,2-diethyl-, cis-	112	C8H16	061141-50-2	52
5	1-Nonene	126	C9H18	000124-11-8	43





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110419\
Data File : VX013324.D
Acq On : 04 Nov 2019 15:07
Operator : JC/SP
Sample : K5683-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : OILY-WATER

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X103119W.M
Quant Title : SW846 8260

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

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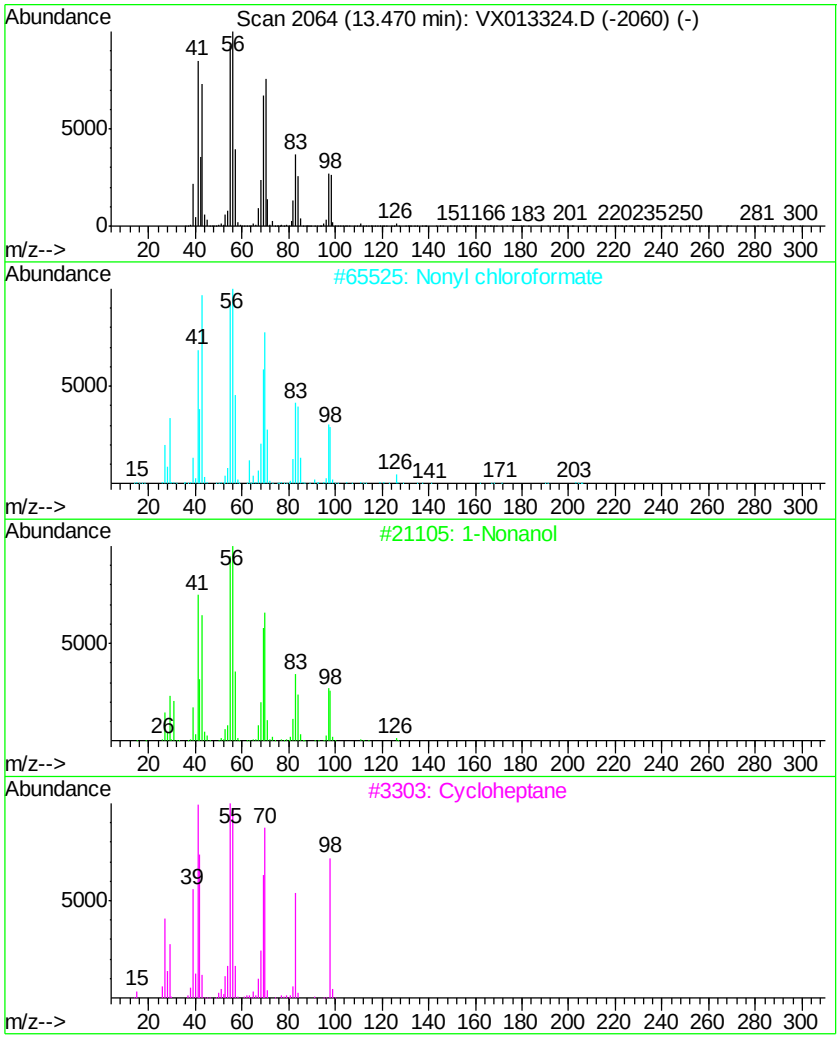
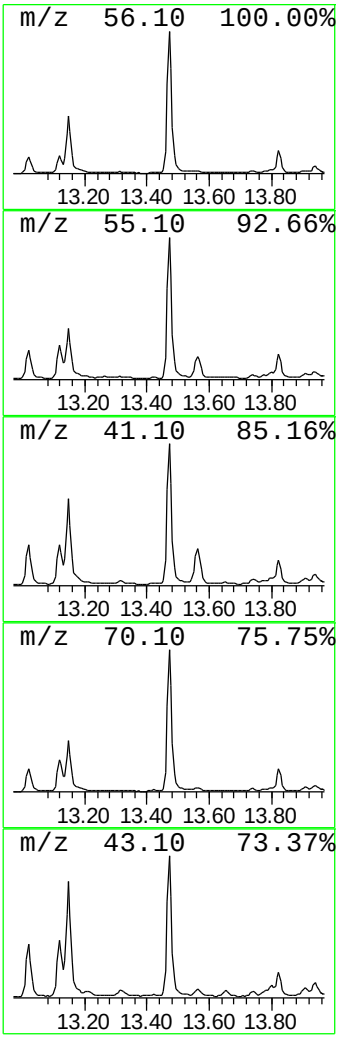
Peak Number 8 Nonyl chloroformate Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.47	180.12 ug/l	11813000	1,4-Dichlorobenzene-d4	12.07

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonyl chloroformate	206	C10H19ClO2	057045-82-6	93
2	1-Nonanol	144	C9H20O	000143-08-8	91
3	Cycloheptane	98	C7H14	000291-64-5	87
4	Cyclopropane, 1-methyl-2-octyl-	168	C12H24	037617-26-8	80
5	Isopropylcyclobutane	98	C7H14	000872-56-0	68





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110419\
Data File : VX013324.D
Acq On : 04 Nov 2019 15:07
Operator : JC/SP
Sample : K5683-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : OILY-WATER

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X103119W.M
Quant Title : SW846 8260

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

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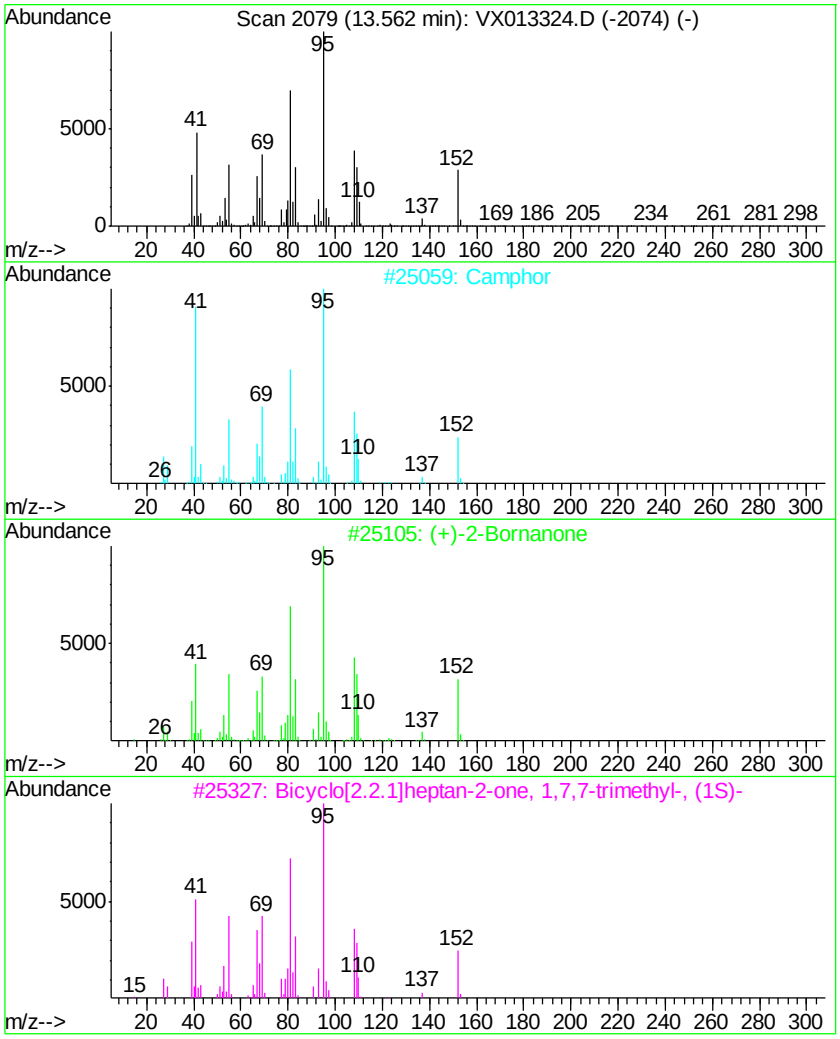
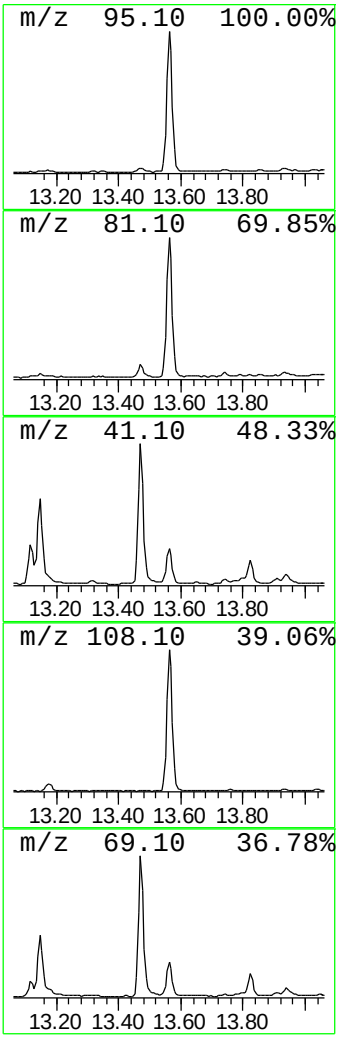
Peak Number 9 Camphor Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.56	70.67 ug/l	4634770	1,4-Dichlorobenzene-d4	12.07

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Camphor	152	C10H16O	000076-22-2	98
2	(+)-2-Bornanone	152	C10H16O	000464-49-3	98
3	Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-48-2	96
4	2-Butanone, 4-(5-methyl-2-furanyl)-	152	C9H12O2	013679-56-6	46
5	Bicyclo[3.1.0]hexane, 1,5-dimethyl-	110	C8H14	1000142-17-5	38





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110419\
Data File : VX013324.D
Acq On : 04 Nov 2019 15:07
Operator : JC/SP
Sample : K5683-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : OILY-WATER

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X103119W.M
Quant Title : SW846 8260

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

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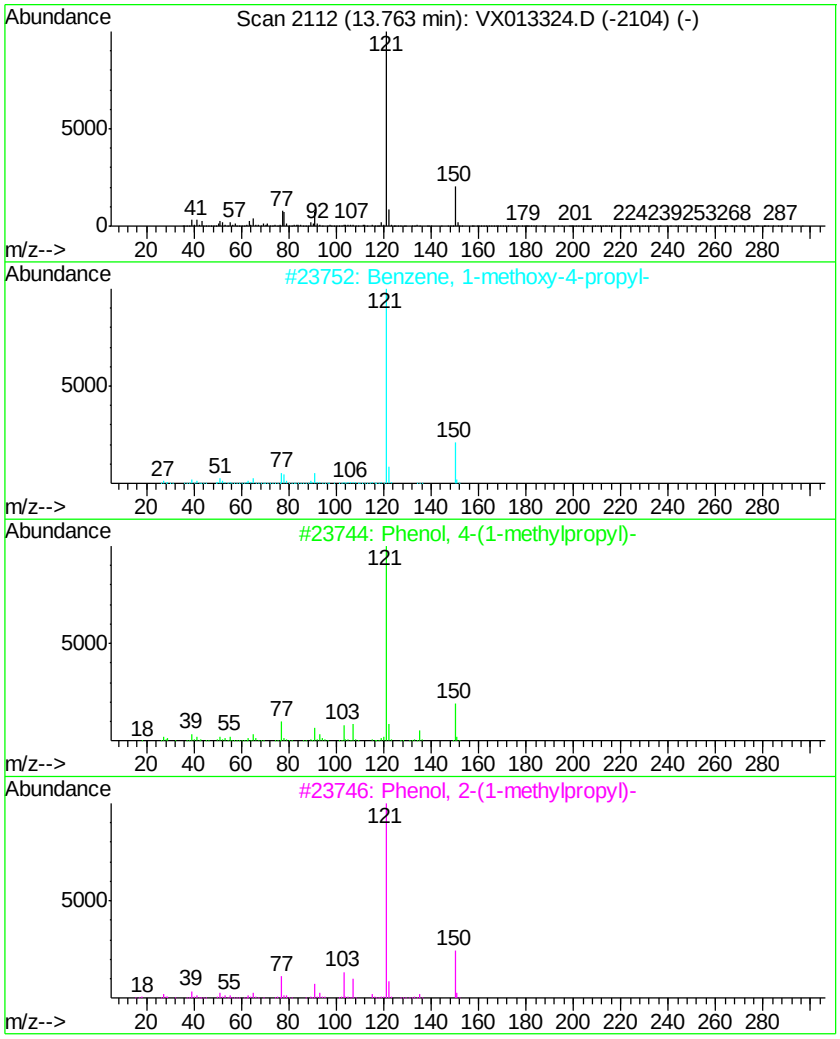
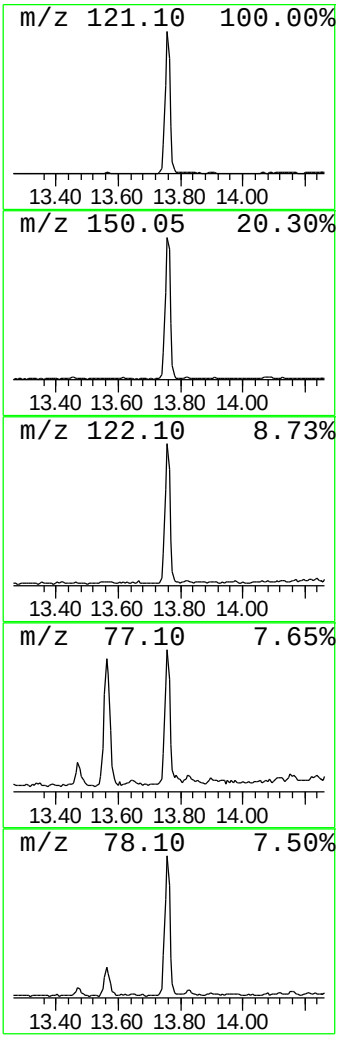
Peak Number 10 Benzene, 1-methoxy-4-propyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.76	31.84 ug/l	2088140	1,4-Dichlorobenzene-d4	12.07

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methoxy-4-propyl-	150	C10H14O	000104-45-0	90
2	Phenol, 4-(1-methylpropyl)-	150	C10H14O	000099-71-8	86
3	Phenol, 2-(1-methylpropyl)-	150	C10H14O	000089-72-5	78
4	Paroxypropione	150	C9H10O2	000070-70-2	72
5	Propanedinitrile, [3-(4-methoxyp...	226	C14H14N2O	069244-84-4	72





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110419\
Data File : VX013324.D
Acq On : 04 Nov 2019 15:07
Operator : JC/SP
Sample : K5683-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : OILY-WATER

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X103119W.M
Quant Title : SW846 8260

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

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
2-Pentanone	7.53	25.9	ug/l	1372210	2	6.84	2649050	50.0
2-Heptanone, 6-me...	11.53	21.5	ug/l	1409420	4	12.07	3279140	50.0
Eucalyptol	12.15	184.5	ug/l	12099400	4	12.07	3279140	50.0
Octane, 1,1'-oxybis-	13.02	56.8	ug/l	3722090	4	12.07	3279140	50.0
Dodecane, 1-fluoro-	13.12	53.7	ug/l	3521100	4	12.07	3279140	50.0
1-Octanol, 2-methyl-	13.15	116.7	ug/l	7653600	4	12.07	3279140	50.0
Nonyl chloroformate	13.47	180.1	ug/l	11813000	4	12.07	3279140	50.0
Camphor	13.56	70.7	ug/l	4634770	4	12.07	3279140	50.0
Benzene, 1-methox...	13.76	31.8	ug/l	2088140	4	12.07	3279140	50.0