LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX111418\
Data File : VX005996.D
Acq On : 15 Nov 2018 01:54
Operator : JC/MD
Sample : J5939-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 33 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\82X110718W.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	1.605	65	74	77	rBV4	3305	6709	1.30%	0.095%
2	2.446	206	212	217	rBV	3386	5826	1.13%	0.082%
3	5.501	702	713	728	rBV	73413	202243	39.19%	2.858%
4	5.665	728	740	754	rBV	126666	345505	66.95%	4.882%
5	6.068	795	806	820	rBV2	74683	200209	38.79%	2.829%
6	6.860	926	936	955	rBV	172259	420360	81.45%	5.940%
7	8.713	1233	1240	1249	rBV	329126	516071	100.00%	7.292%
8	10.116	1464	1470	1478	rBV	331132	451027	87.40%	6.373%
9	10.640	1550	1556	1563	rBV4	3216	6907	1.34%	0.098%
10	10.835	1578	1588	1592	rBV4	5800	9630	1.87%	0.136%
11	10.914	1597	1601	1605	rBV3	4881	7748	1.50%	0.109%
12	11.018	1613	1618	1622	rBV	10150	12748	2.47%	0.180%
13	11.140	1632	1638	1643	rBV	279184	357275	69.23%	5.048%
14	11.280	1653	1661	1668	rVB4	6019	10377	2.01%	0.147%
15	11.359	1670	1674	1677	rBV	4491	5415	1.05%	0.077%
16	11.670	1720	1725	1732	rVB4	5436	9733	1.89%	0.138%
17	11.804	1744	1747	1757	rVB3	4672	9743	1.89%	0.138%
18	11.920	1757	1766	1767	rBV2	9227	13780	2.67%	0.195%
19	11.945	1767	1770	1775	rVB	34483	43748	8.48%	0.618%
20	11.999	1775	1779	1782	rBV3	4462	5661	1.10%	0.080%
21	12.079	1787	1792	1800	rBV	376872	466281	90.35%	6.588%
22	12.231	1813	1817	1824	rVB	10076	14798	2.87%	0.209%
23	12.304	1824	1829	1835	rBV	24952	32847	6.36%	0.464%
24	12.371	1835	1840	1842	rBV3	13276	22059	4.27%	0.312%
25	12.463	1847	1855	1862	rBV	26244	39127	7.58%	0.553%
26	12.524	1862	1865	1870	rVB4	4626	7306	1.42%	0.103%
27	12.609	1870	1879	1881	rBV6	3551	8249	1.60%	0.117%
28	12.670	1881	1889	1892	rVV	62225	88719	17.19%	1.254%
29	12.700	1892	1894	1898	rVV	21849	26258	5.09%	0.371%
30	12.755	1898	1903	1911	rVV5	76864	151421	29.34%	2.140%
31	12.816	1911	1913	1918	rVV3	9163	12881	2.50%	0.182%
32	12.896	1919	1926	1930	rVV	27462	46077	8.93%	0.651%
33	12.975	1930	1939	1944	rVV2	72182	117102	22.69%	1.655%
34	13.042	1944	1950	1953	rVV4	7573	14080	2.73%	0.199%

LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX111418\
Data File : VX005996.D
Acq On : 15 Nov 2018 01:54
Operator : JC/MD
Sample : J5939-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 33 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\82X110718W.M
Title : SW846 8260

35	13.085	1953	1957	1963	rVB3	23539	37736	7.31%	0.533%
36	13.152	1963	1968	1971	rBV	25179	32928	6.38%	0.465%
37	13.194	1971	1975	1979	rVV2	51422	66620	12.91%	0.941%
38	13.255	1981	1985	1989	rVB	12435	17288	3.35%	0.244%
39	13.347	1995	2000	2006	rVB2	252540	369401	71.58%	5.219%
40	13.426	2006	2013	2016	rBV2	19889	38158	7.39%	0.539%
41	13.481	2018	2022	2030	rVB	92379	121751	23.59%	1.720%
42	13.566	2032	2036	2038	rBV2	15017	20738	4.02%	0.293%
43	13.597	2038	2041	2044	rVV	23315	30055	5.82%	0.425%
44	13.639	2044	2048	2052	rVV2	76951	107817	20.89%	1.523%
45	13.700	2052	2058	2059	rVV2	51382	89299	17.30%	1.262%
46	13.719	2059	2061	2067	rVB2	78570	113650	22.02%	1.606%
47	13.810	2070	2076	2079	rBV	29872	42413	8.22%	0.599%
48	13.853	2079	2083	2088	rVB2	41857	57077	11.06%	0.806%
49	13.914	2088	2093	2097	rVB	111176	145287	28.15%	2.053%
50	13.987	2097	2105	2109	rBV2	151037	219991	42.63%	3.108%
51	14.030	2109	2112	2116	rBV2	29730	40642	7.88%	0.574%
52	14.060	2116	2117	2121	rVB	14501	14574	2.82%	0.206%
53	14.127	2124	2128	2131	rBV2	23675	32240	6.25%	0.456%
54	14.164	2131	2134	2138	rBV2	22014	22456	4.35%	0.317%
55	14.219	2141	2143	2146	rBV2	6463	7723	1.50%	0.109%
56	14.286	2146	2154	2160	rVB2	107093	210222	40.74%	2.970%
57	14.347	2161	2164	2166	rBV2	6957	7805	1.51%	0.110%
58	14.377	2166	2169	2173	rVB	40626	48773	9.45%	0.689%
59	14.432	2173	2178	2182	rVB2	48444	60419	11.71%	0.854%
60	14.475	2182	2185	2190	rVB2	25054	32098	6.22%	0.454%
61	14.542	2190	2196	2200	rBV2	238362	313778	60.80%	4.434%
62	14.633	2209	2211	2214	rBV2	6203	7313	1.42%	0.103%
63	14.676	2214	2218	2224	rVV2	68291	127505	24.71%	1.802%
64	14.731	2224	2227	2232	rVB2	68137	83474	16.17%	1.179%
65	14.785	2232	2236	2240	rBV2	18871	26221	5.08%	0.370%
66	14.840	2240	2245	2249	rVV2	70588	116006	22.48%	1.639%
67	14.877	2249	2251	2253	rVV2	37483	45105	8.74%	0.637%
68	14.907	2253	2256	2260	rVV2	43930	59293	11.49%	0.838%
69	14.981	2260	2268	2273	rVV5	30463	76086	14.74%	1.075%
70	15.042	2275	2278	2282	rVB2	9210	13734	2.66%	0.194%
71	15.115	2283	2290	2297	rBV3	16373	30710	5.95%	0.434%

LSC Area Percent Report

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Operator : JC/MD
Sample : J5939-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 33 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\82X110718W.M
Title : SW846 8260

72	15.249	2306	2312	2320	rBV	53566	106497	20.64%	1.505%
73	15.322	2320	2324	2331	rVB6	13228	24819	4.81%	0.351%
74	15.389	2331	2335	2339	rVB5	3573	6264	1.21%	0.089%
75	15.444	2339	2344	2347	rBV2	14364	22903	4.44%	0.324%
76	15.480	2347	2350	2353	rVV	10661	13706	2.66%	0.194%
77	15.548	2353	2361	2366	rVV2	23640	43810	8.49%	0.619%
78	15.596	2366	2369	2375	rVV3	16192	33119	6.42%	0.468%
79	15.688	2378	2384	2388	rVV	58436	103534	20.06%	1.463%
80	15.724	2388	2390	2398	rVB	24665	38062	7.38%	0.538%
81	15.804	2398	2403	2412	rBV4	8877	23019	4.46%	0.325%
82	15.889	2412	2417	2419	rVV3	8628	15644	3.03%	0.221%
83	15.926	2419	2423	2437	rVB4	15177	38261	7.41%	0.541%
84	16.072	2443	2447	2450	rBV3	3727	5673	1.10%	0.080%
85	16.102	2450	2452	2459	rVB4	4236	6460	1.25%	0.091%
86	16.175	2459	2464	2470	rBV7	2948	6611	1.28%	0.093%
87	16.255	2473	2477	2486	rVB9	2496	6643	1.29%	0.094%
88	16.389	2490	2499	2505	rBV9	3162	8022	1.55%	0.113%

Sum of corrected areas: 7077353

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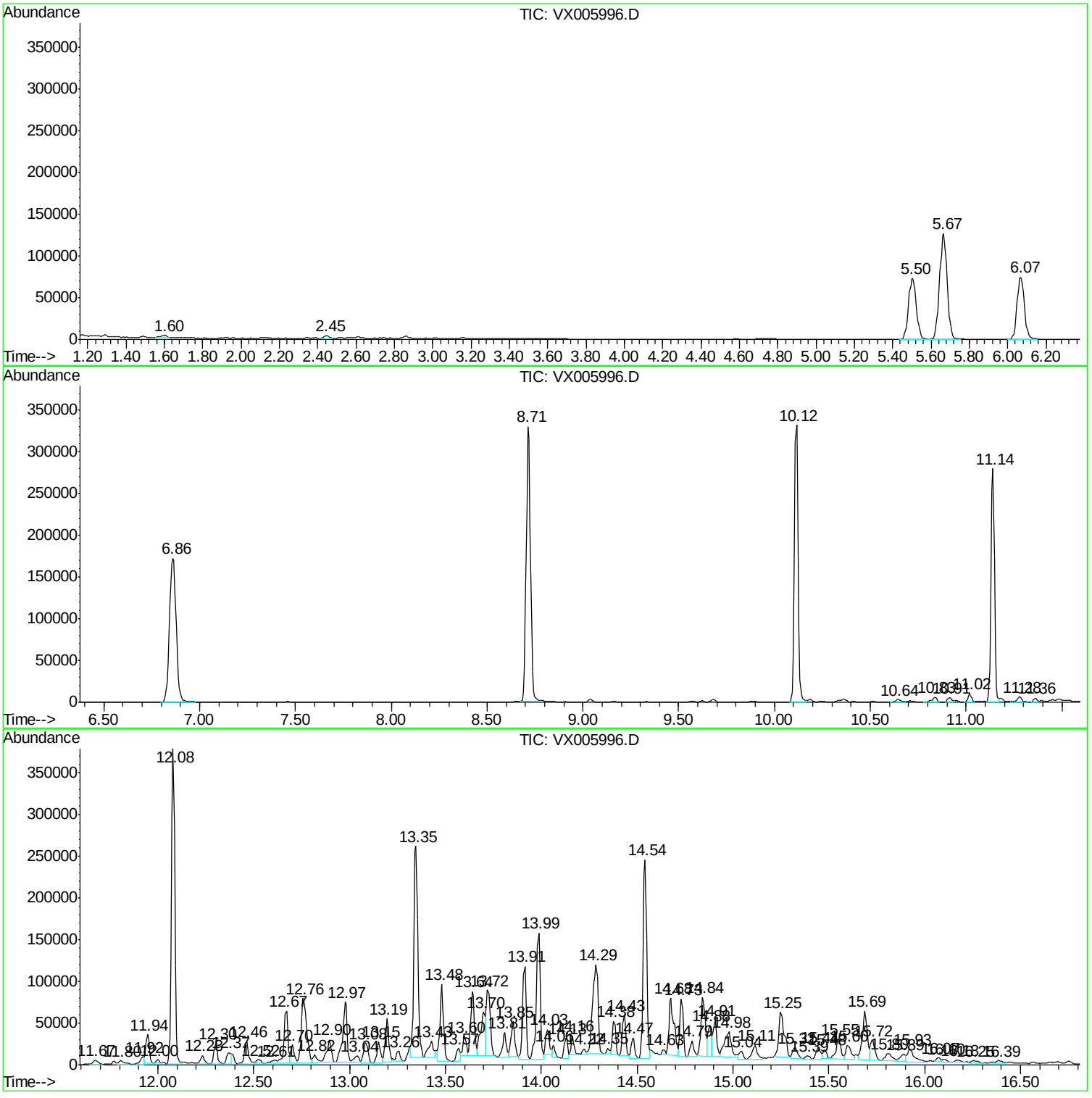
Operator : JC/MD
Sample : J5939-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 33 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
OILY-WATER

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.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\VX111418\
Data File : VX005996.D
Acq On : 15 Nov 2018 01:54
Operator : JC/MD
Sample : J5939-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 33 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : OILY-WATER

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
Quant Title : SW846 8260

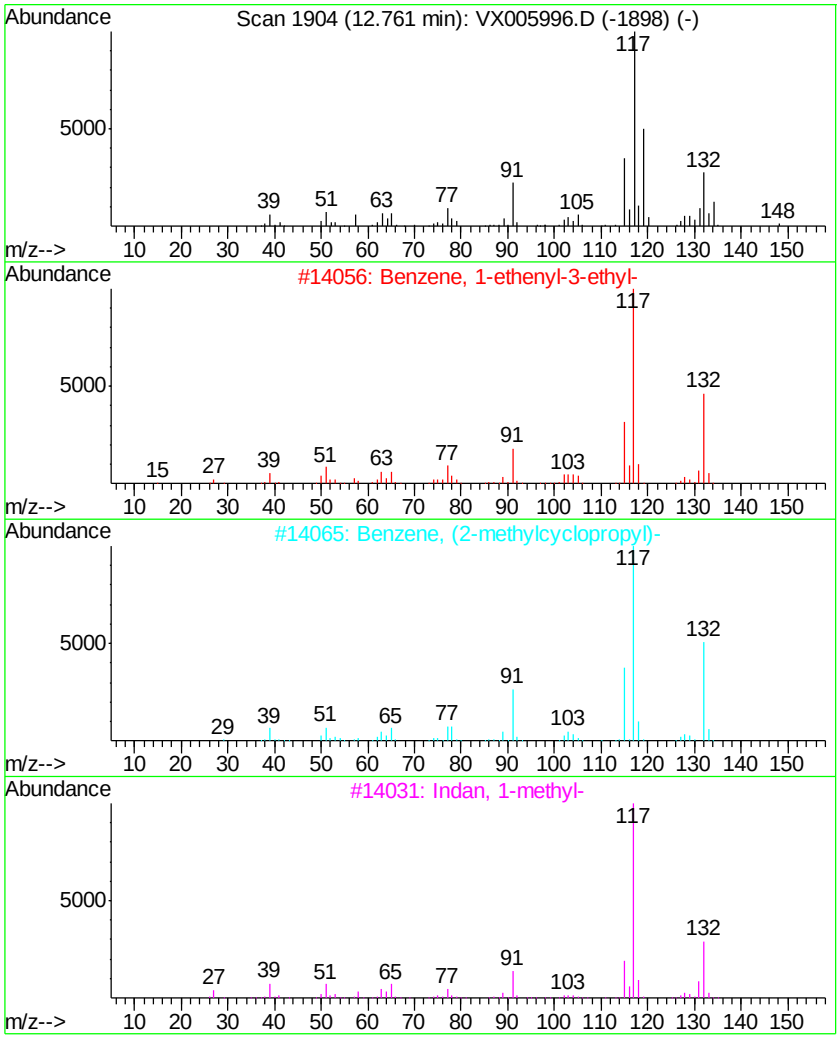
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

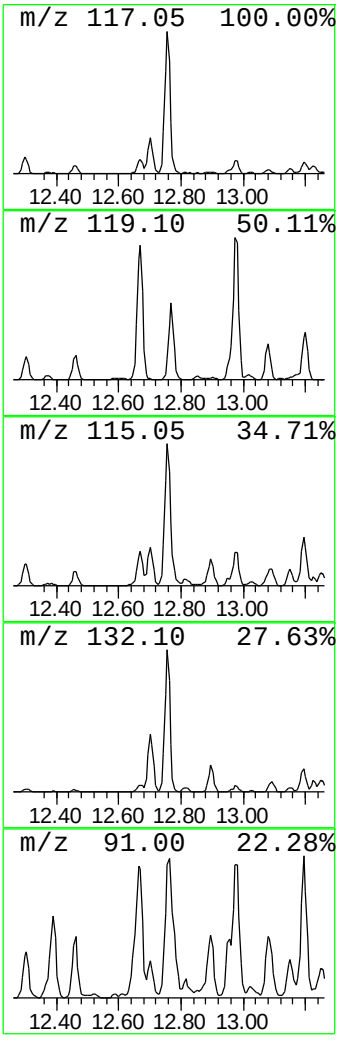
****

Peak Number 1 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	16.24 ug/l	151421	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	83
2	Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	81
3	Indan, 1-methyl-	132	C10H12	000767-58-8	81
4	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	80
5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	80





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX111418\
Data File : VX005996.D
Acq On : 15 Nov 2018 01:54
Operator : JC/MD
Sample : J5939-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 33 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : OILY-WATER

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
Quant Title : SW846 8260

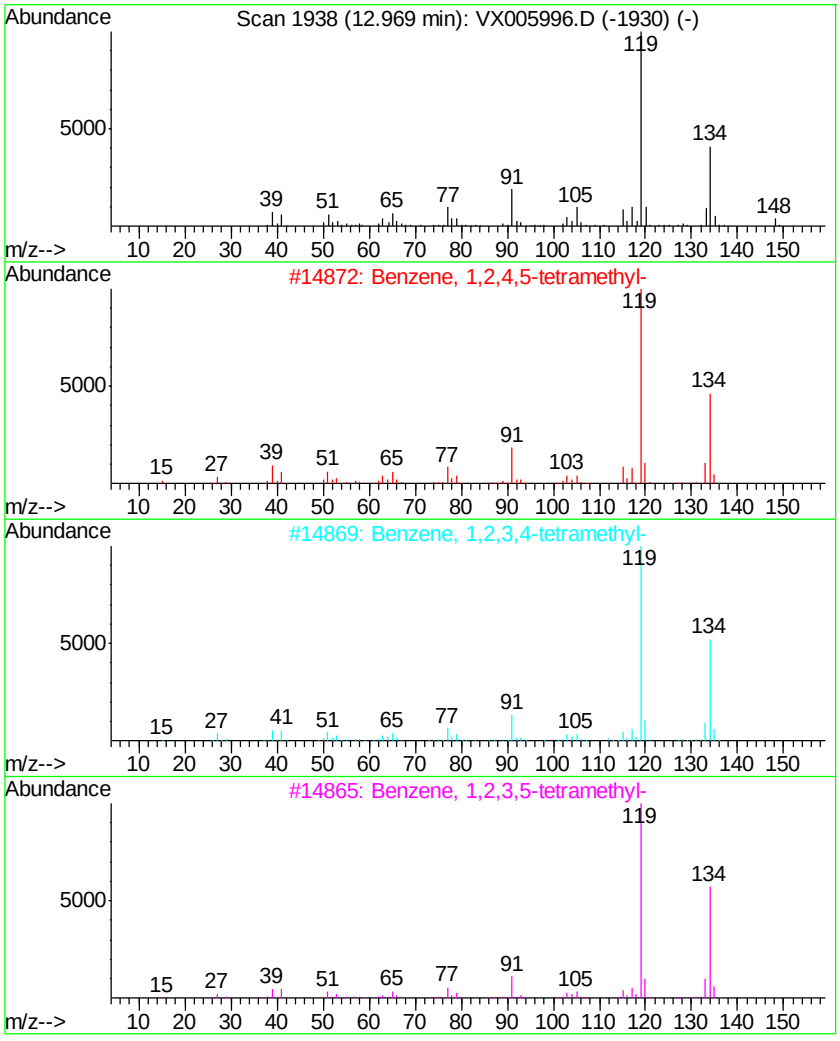
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

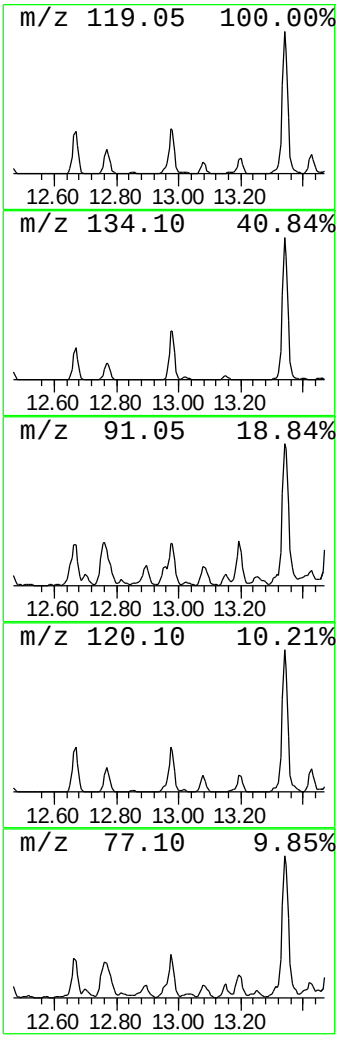
****

Peak Number 2 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.97	12.56 ug/l	117102	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
2	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
3	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
4	o-Cymene	134	C10H14	000527-84-4	93
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	91





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX111418\
Data File : VX005996.D
Acq On : 15 Nov 2018 01:54
Operator : JC/MD
Sample : J5939-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 33 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : OILY-WATER

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
Quant Title : SW846 8260

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

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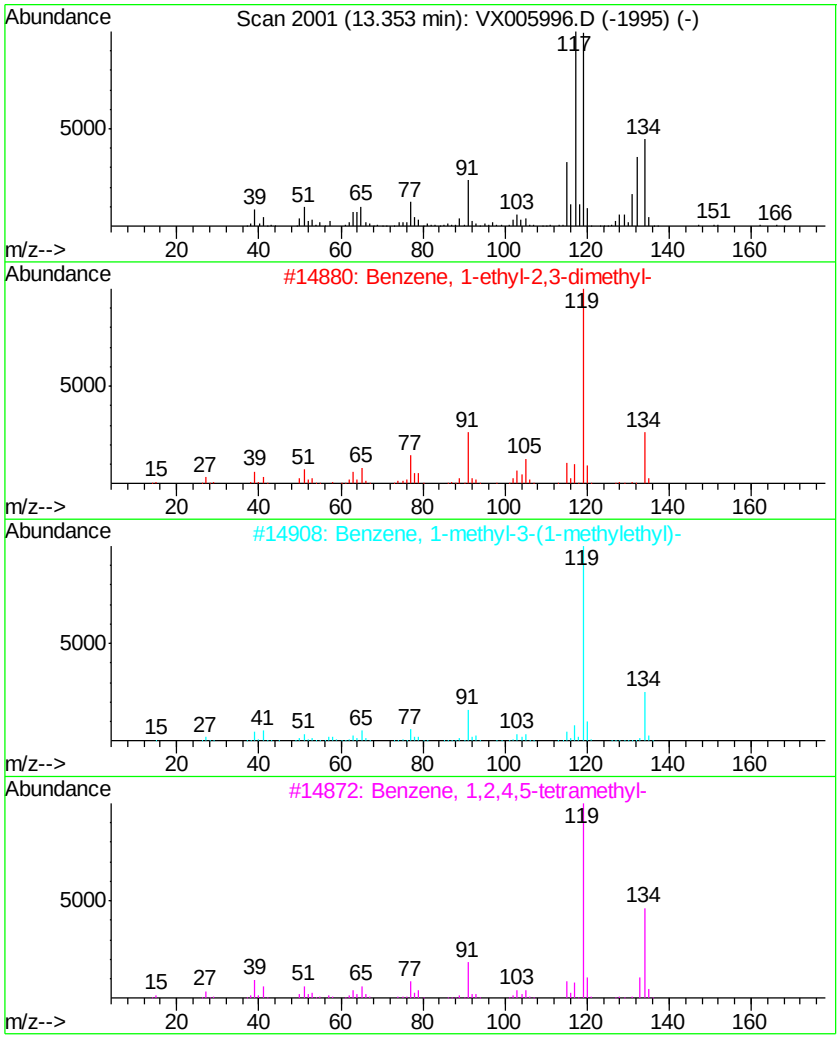
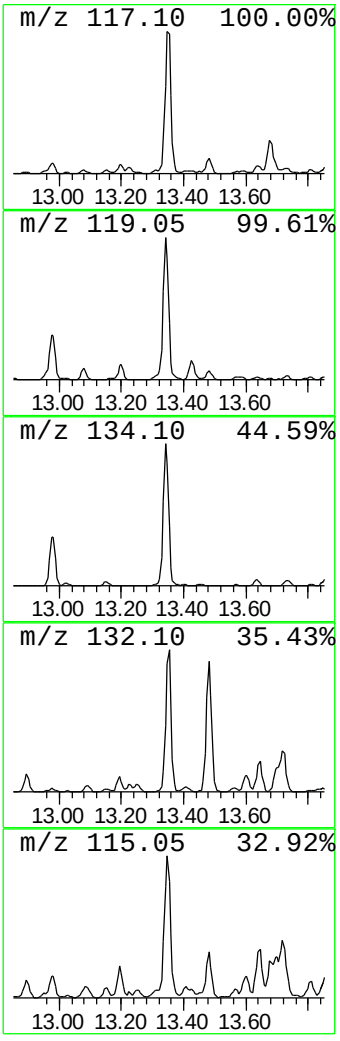
Peak Number 3 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	39.61 ug/l	369401	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	70
2	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	70
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	70
4	o-Cymene	134	C10H14	000527-84-4	68
5	p-Cymene	134	C10H14	000099-87-6	62





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX111418\
Data File : VX005996.D
Acq On : 15 Nov 2018 01:54
Operator : JC/MD
Sample : J5939-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 33 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : OILY-WATER

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
Quant Title : SW846 8260

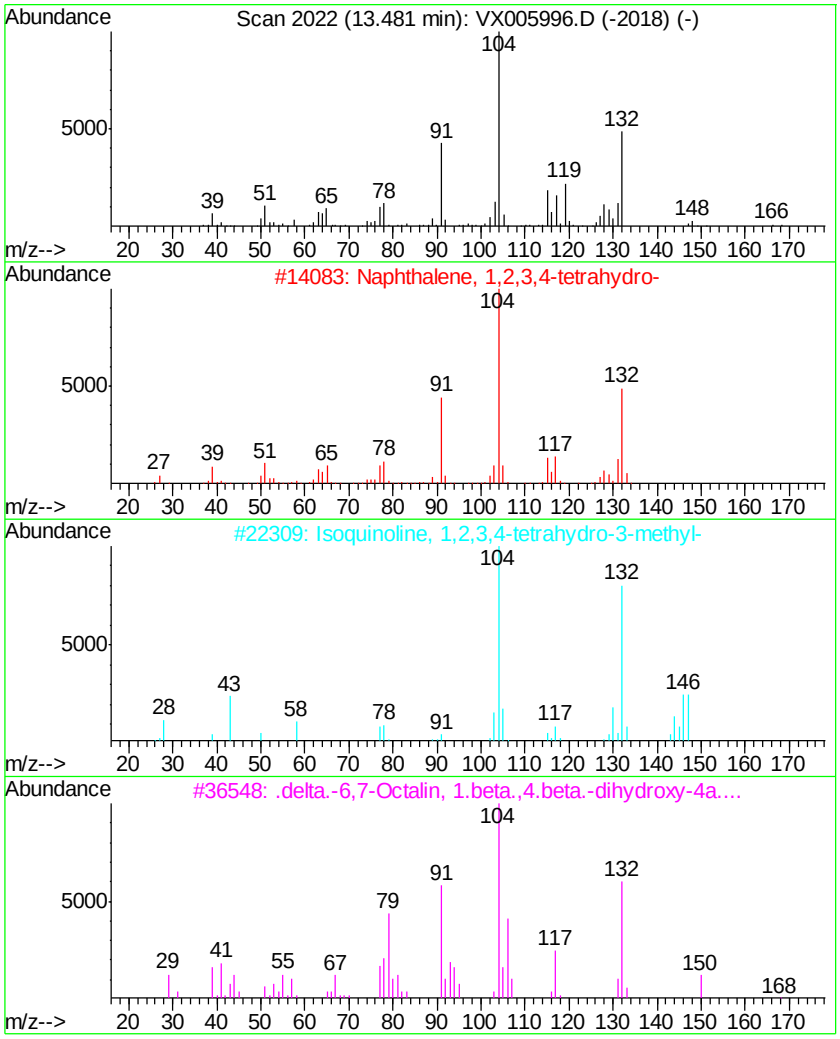
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

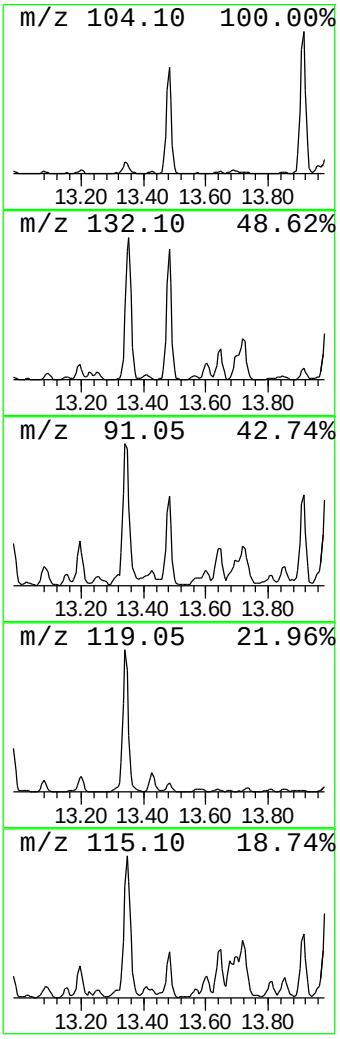
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.48	13.06 ug/l	121751	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	90
2	Isoquinoline, 1,2,3,4-tetrahydro...	147	C10H13N	029726-60-1	42
3	.delta.-6,7-Octalin, 1.beta.,4.b...	168	C10H16O2	051921-13-2	38
4	2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	38
5	2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	35





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX111418\
Data File : VX005996.D
Acq On : 15 Nov 2018 01:54
Operator : JC/MD
Sample : J5939-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 33 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : OILY-WATER

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
Quant Title : SW846 8260

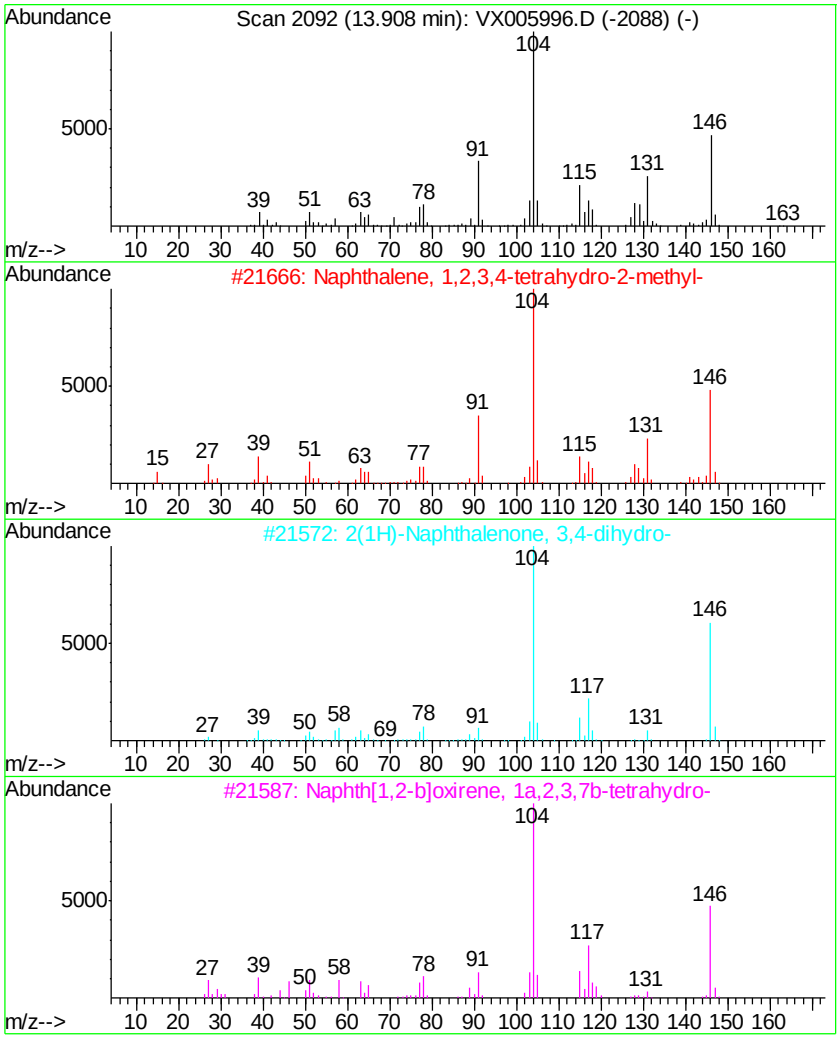
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

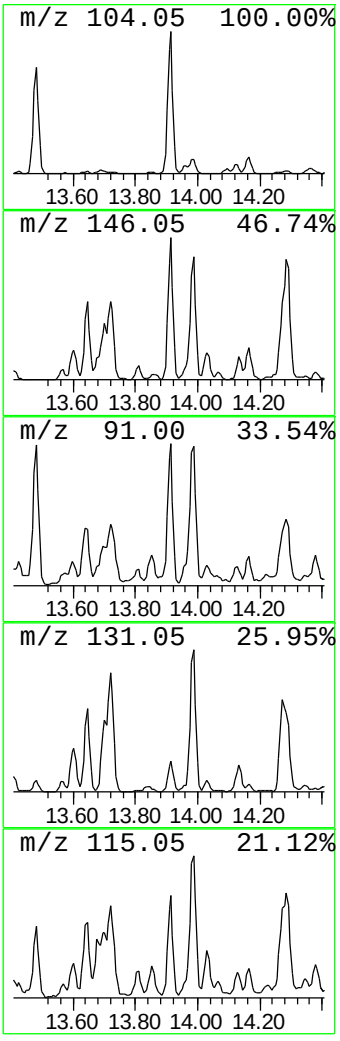
****

Peak Number 5 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.91	15.58 ug/l	145287	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	003877-19-8	94
2	2(1H)-Naphthalenone, 3,4-dihydro-	146	C10H10O	000530-93-8	64
3	Naphth[1,2-b]oxirene, 1a,2,3,7b-...	146	C10H10O	002461-34-9	62
4	2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	38
5	Naphthalene, 2-ethyl-1,2,3,4-tet...	160	C12H16	032367-54-7	38





Library Search Compound Report

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Data File : VX005996.D
Acq On : 15 Nov 2018 01:54
Operator : JC/MD
Sample : J5939-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 33 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : OILY-WATER

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
Quant Title : SW846 8260

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

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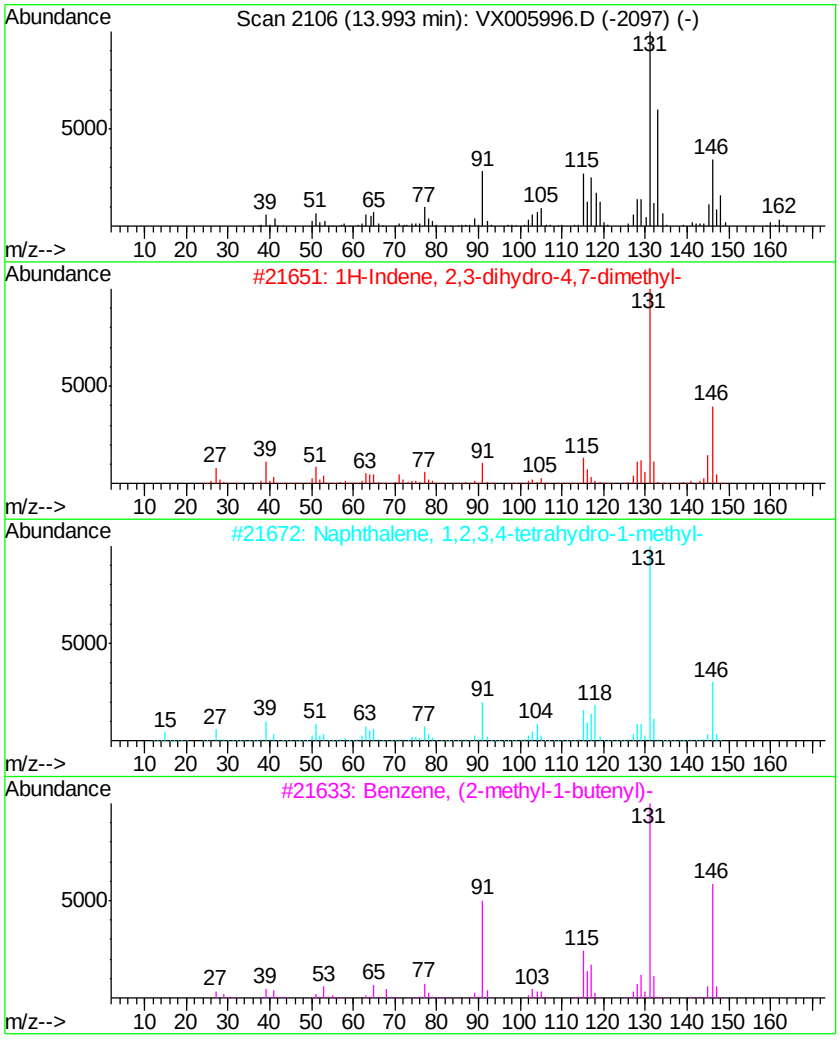
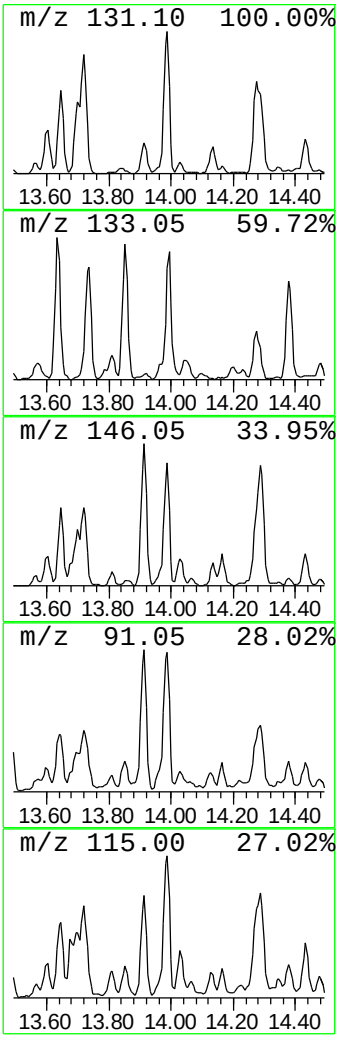
Peak Number 6 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.99	23.59 ug/l	219991	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	93
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	93
3	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	89
4	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	86
5	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	86





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX111418\
Data File : VX005996.D
Acq On : 15 Nov 2018 01:54
Operator : JC/MD
Sample : J5939-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 33 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : OILY-WATER

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
Quant Title : SW846 8260

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

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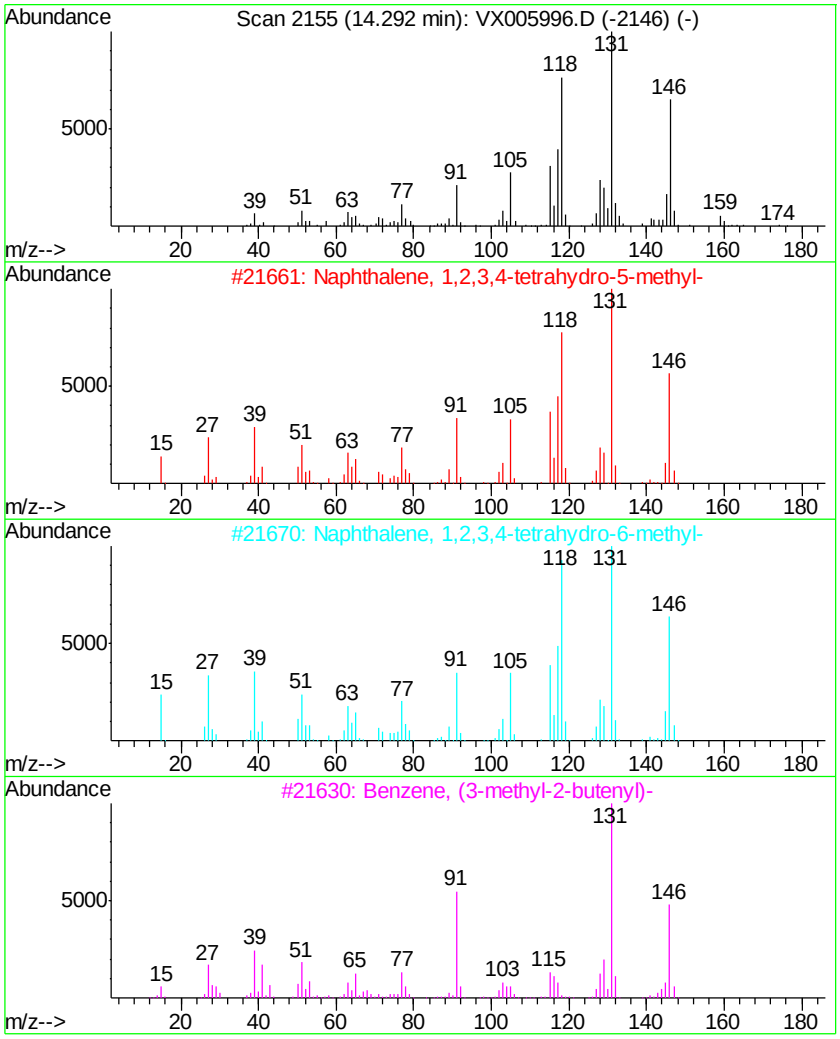
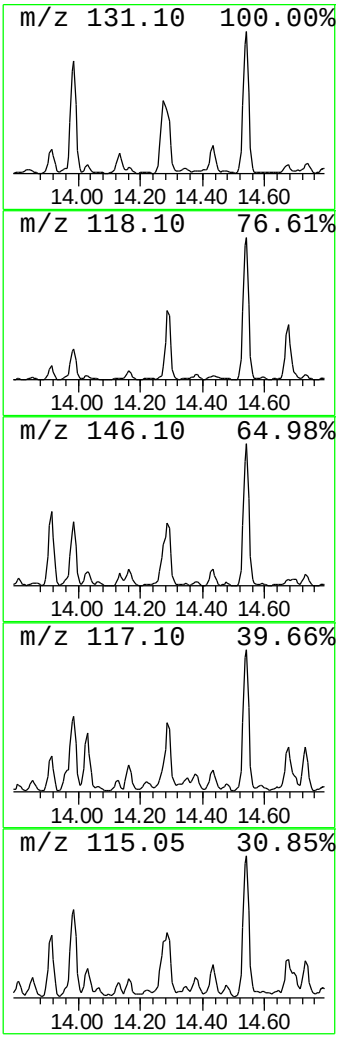
Peak Number 7 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.29	22.54 ug/l	210222	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	96
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	95
3	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	70
4	Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	70
5	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	64





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX111418\
Data File : VX005996.D
Acq On : 15 Nov 2018 01:54
Operator : JC/MD
Sample : J5939-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 33 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : OILY-WATER

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
Quant Title : SW846 8260

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

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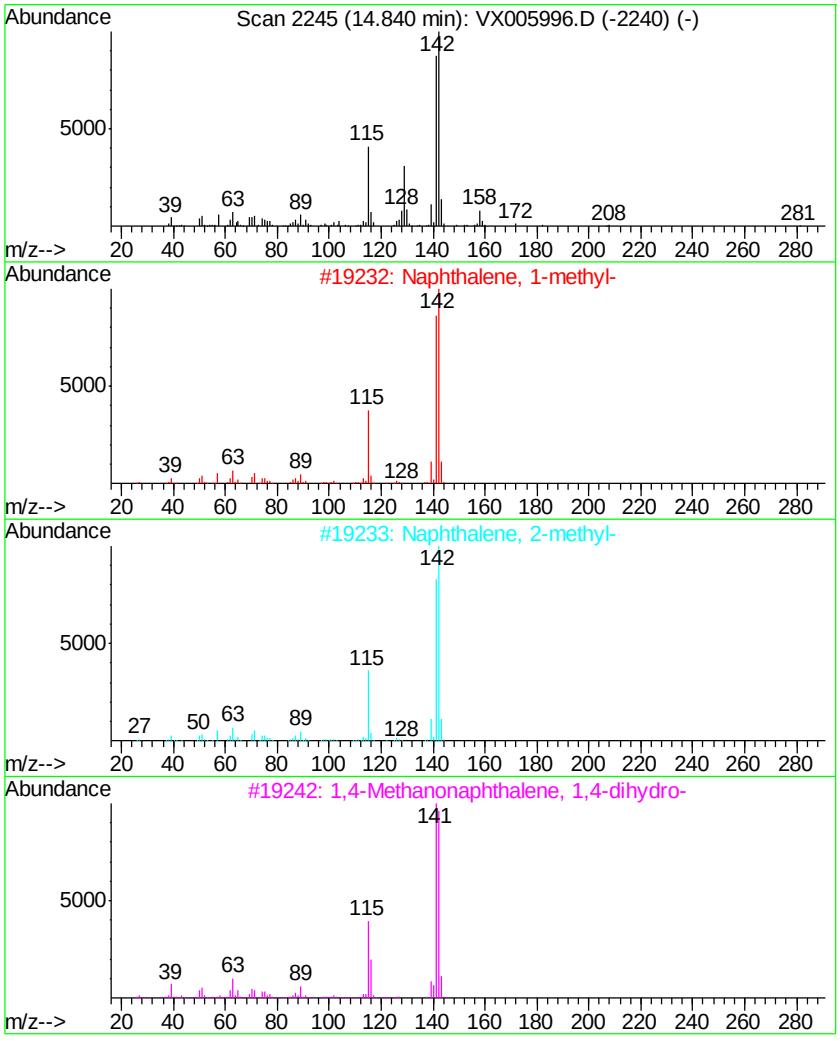
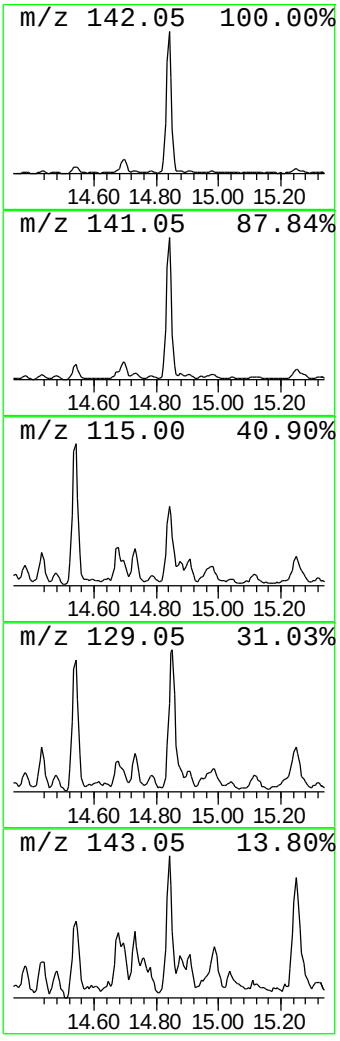
Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.84	12.44 ug/l	116006	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
4	Benzocycloheptatriene	142	C11H10	000264-09-5	87
5	1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	53





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX111418\
Data File : VX005996.D
Acq On : 15 Nov 2018 01:54
Operator : JC/MD
Sample : J5939-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 33 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : OILY-WATER

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.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
Benzene, 1-etheny...	12.76	16.2	ug/l	151421	4	12.08	466281	50.0
Benzene, 1,2,4,5-...	12.97	12.6	ug/l	117102	4	12.08	466281	50.0
Benzene, 1-ethyl-...	13.35	39.6	ug/l	369401	4	12.08	466281	50.0
Naphthalene, 1,2,...	13.48	13.1	ug/l	121751	4	12.08	466281	50.0
Naphthalene, 1,2,...	13.91	15.6	ug/l	145287	4	12.08	466281	50.0
1H-Indene, 2,3-di...	13.99	23.6	ug/l	219991	4	12.08	466281	50.0
Naphthalene, 1,2,...	14.29	22.5	ug/l	210222	4	12.08	466281	50.0
Naphthalene, 1-me...	14.84	12.4	ug/l	116006	4	12.08	466281	50.0