LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX073019\
Data File : VX011208.D
Acq On : 30 Jul 2019 20:24
Operator : JC/SP
Sample : K4048-02
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : OWL-C7-GW

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\82X071619W.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	1.276	15	20	32	rBV	34736	49797	5.45%	0.562%
2	1.599	65	73	78	rBV8	4660	12360	1.35%	0.139%
3	2.440	208	211	220	rVB	5624	9798	1.07%	0.110%
4	2.611	234	239	248	rVB	12551	23625	2.59%	0.266%
5	5.501	702	713	725	rBV	107446	310396	33.99%	3.500%
6	5.659	728	739	759	rVB	208442	594268	65.08%	6.702%
7	6.061	796	805	818	rBV	115071	297525	32.58%	3.355%
8	6.860	926	936	947	rBV	292769	681225	74.60%	7.682%
9	8.713	1229	1240	1252	rBV	586431	913144	100.00%	10.298%
10	9.037	1286	1293	1299	rBV4	8356	17769	1.95%	0.200%
11	10.110	1463	1469	1478	rBV	569869	774008	84.76%	8.729%
12	10.189	1478	1482	1487	rVB3	6909	11884	1.30%	0.134%
13	10.335	1501	1506	1516	rVB3	5320	11555	1.27%	0.130%
14	10.652	1551	1558	1564	rBV8	4689	11432	1.25%	0.129%
15	10.835	1579	1588	1593	rBV7	7094	13956	1.53%	0.157%
16	11.018	1613	1618	1624	rBV	34692	47440	5.20%	0.535%
17	11.134	1631	1637	1643	rBV	474140	602554	65.99%	6.795%
18	11.670	1717	1725	1732	rBV7	6140	11925	1.31%	0.134%
19	11.768	1735	1741	1745	rBV	14140	19715	2.16%	0.222%
20	11.920	1758	1766	1767	rBV3	13609	21282	2.33%	0.240%
21	11.945	1767	1770	1775	rVB	30064	37248	4.08%	0.420%
22	12.073	1785	1791	1802	rBV	580049	754795	82.66%	8.512%
23	12.231	1813	1817	1821	rVB	14188	19286	2.11%	0.217%
24	12.298	1823	1828	1834	rVV	67272	85230	9.33%	0.961%
25	12.384	1834	1842	1848	rVB5	8601	23370	2.56%	0.264%
26	12.457	1849	1854	1863	rBV	41248	53656	5.88%	0.605%
27	12.664	1878	1888	1891	rBV	55796	80659	8.83%	0.910%
28	12.701	1891	1894	1898	rVV	38073	51788	5.67%	0.584%
29	12.749	1898	1902	1910	rVV	126273	191272	20.95%	2.157%
30	12.816	1910	1913	1917	rVB	12130	15241	1.67%	0.172%
31	12.896	1917	1926	1931	rBV	32848	56625	6.20%	0.639%
32	12.975	1931	1939	1945	rBV	125803	177117	19.40%	1.997%
33	13.079	1951	1956	1962	rVB2	18620	31559	3.46%	0.356%
34	13.146	1963	1967	1970	rBV2	10098	13529	1.48%	0.153%

LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX073019\
Data File : VX011208.D
Acq On : 30 Jul 2019 20:24
Operator : JC/SP
Sample : K4048-02
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : OWL-C7-GW

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\82X071619W.M
Title : SW846 8260

35	13.194	1970	1975	1978	rBV2	39652	54757	6.00%	0.618%
36	13.225	1978	1980	1982	rVV	13093	13651	1.49%	0.154%
37	13.249	1982	1984	1989	rVB2	10096	10969	1.20%	0.124%
38	13.310	1989	1994	1995	rBV3	9094	14730	1.61%	0.166%
39	13.347	1995	2000	2005	rVB2	230643	328189	35.94%	3.701%
40	13.402	2005	2009	2011	rBV2	10405	13621	1.49%	0.154%
41	13.597	2037	2041	2044	rVV2	21229	32152	3.52%	0.363%
42	13.639	2044	2048	2051	rVV3	63911	97776	10.71%	1.103%
43	13.694	2054	2057	2058	rVV	42159	51471	5.64%	0.580%
44	13.719	2058	2061	2070	rVV2	76012	119864	13.13%	1.352%
45	13.804	2072	2075	2078	rVV	11680	13937	1.53%	0.157%
46	13.847	2078	2082	2087	rVB3	26396	34001	3.72%	0.383%
47	13.908	2087	2092	2097	rBV	146734	180006	19.71%	2.030%
48	13.981	2097	2104	2108	rBV2	112137	155142	16.99%	1.750%
49	14.023	2108	2111	2115	rVV2	24434	33135	3.63%	0.374%
50	14.127	2124	2128	2132	rBV	33583	44186	4.84%	0.498%
51	14.267	2146	2151	2161	rBV2	54834	103930	11.38%	1.172%
52	14.377	2165	2169	2173	rVV	77401	100318	10.99%	1.131%
53	14.426	2173	2177	2182	rVV3	60016	93482	10.24%	1.054%
54	14.475	2182	2185	2190	rVB5	11085	17819	1.95%	0.201%
55	14.536	2190	2195	2200	rBV2	137941	184012	20.15%	2.075%
56	14.578	2200	2202	2208	rVV5	6548	11737	1.29%	0.132%
57	14.670	2212	2217	2219	rVV2	34459	59384	6.50%	0.670%
58	14.688	2219	2220	2223	rVV	35715	35380	3.87%	0.399%
59	14.731	2223	2227	2232	rVV2	31959	46281	5.07%	0.522%
60	14.786	2232	2236	2241	rVV7	12109	19141	2.10%	0.216%
61	14.847	2241	2246	2248	rVV	33664	48909	5.36%	0.552%
62	14.877	2248	2251	2253	rVV2	24445	36329	3.98%	0.410%
63	14.901	2253	2255	2260	rVB2	29647	38026	4.16%	0.429%
64	14.975	2260	2267	2272	rBV3	16051	31826	3.49%	0.359%
65	15.109	2287	2289	2296	rVB4	9380	13241	1.45%	0.149%
66	15.243	2304	2311	2318	rBV2	47172	83845	9.18%	0.946%
67	15.322	2318	2324	2331	rVV3	25252	44502	4.87%	0.502%
68	15.401	2331	2337	2338	rVV5	4317	9288	1.02%	0.105%
69	15.438	2340	2343	2345	rVV4	15147	21206	2.32%	0.239%
70	15.474	2345	2349	2354	rVV	33325	53169	5.82%	0.600%
71	15.541	2356	2360	2364	rVV2	41239	69742	7.64%	0.786%

LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX073019\
Data File : VX011208.D
Acq On : 30 Jul 2019 20:24
Operator : JC/SP
Sample : K4048-02
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : OWL-C7-GW

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\82X071619W.M
Title : SW846 8260

72	15.584	2364	2367	2373	rVV6	15563	31550	3.46%	0.356%
73	15.682	2373	2383	2388	rVV2	45341	94403	10.34%	1.065%
74	15.724	2388	2390	2397	rVB7	8790	14184	1.55%	0.160%
75	15.804	2397	2403	2409	rBV3	15408	33416	3.66%	0.377%
76	15.883	2409	2416	2418	rVV3	30663	53998	5.91%	0.609%
77	15.913	2418	2421	2429	rVB	53265	95084	10.41%	1.072%
78	16.066	2437	2446	2457	rBV	58082	108783	11.91%	1.227%
79	16.407	2493	2502	2509	rVB2	17200	40838	4.47%	0.461%
80	16.614	2531	2536	2542	rVV6	7805	21196	2.32%	0.239%
81	16.688	2544	2548	2553	rVV6	7762	16149	1.77%	0.182%
82	16.749	2553	2558	2564	rVB6	5735	11712	1.28%	0.132%

Sum of corrected areas: 8867500

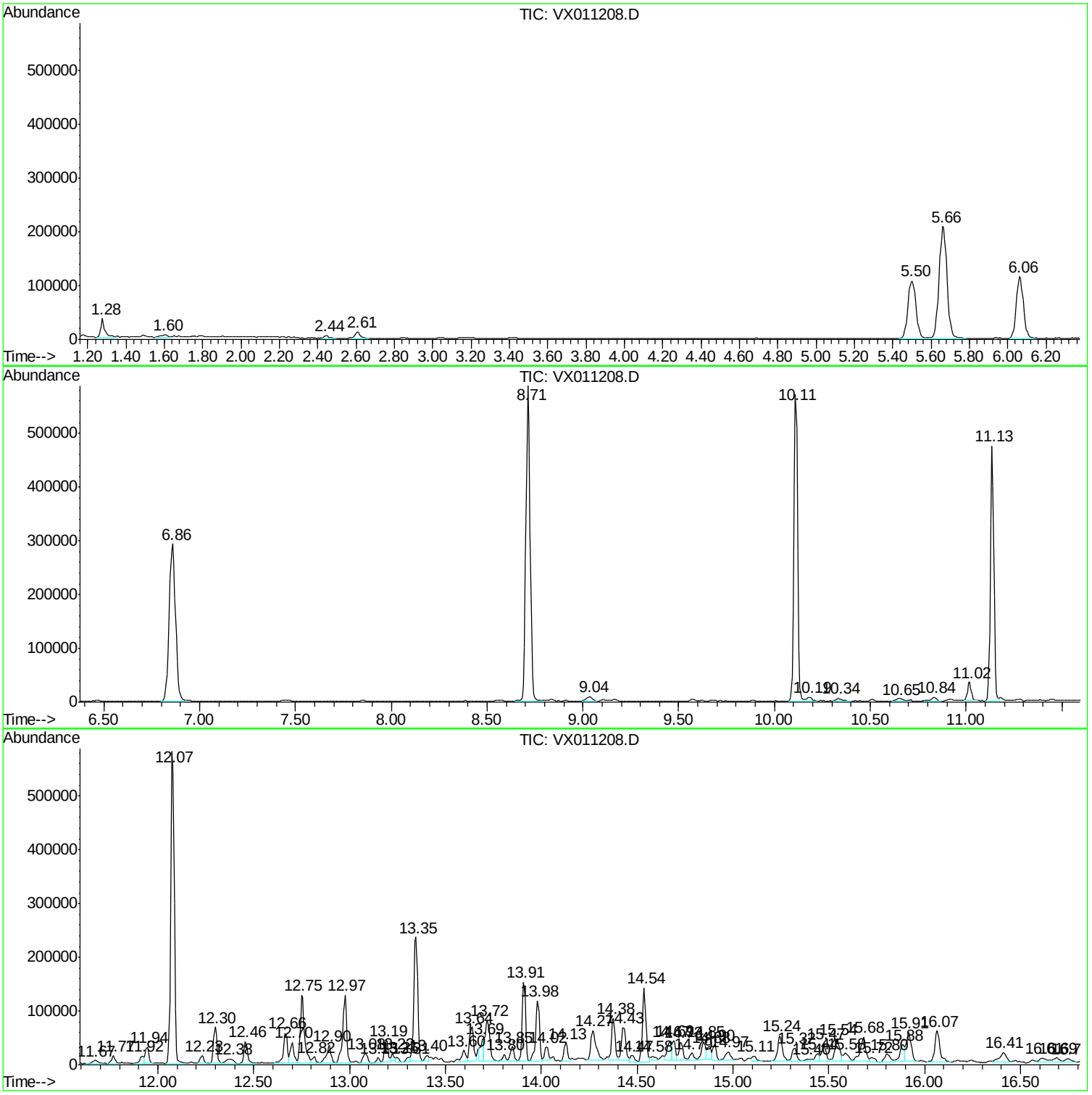
LSC Report - Integrated Chromatogram

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX073019\
Data File : VX011208.D
Acq On : 30 Jul 2019 20:24
Operator : JC/SP
Sample : K4048-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : OWL-C7-GW

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.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\VX073019\
Data File : VX011208.D
Acq On : 30 Jul 2019 20:24
Operator : JC/SP
Sample : K4048-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : OWL-C7-GW

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

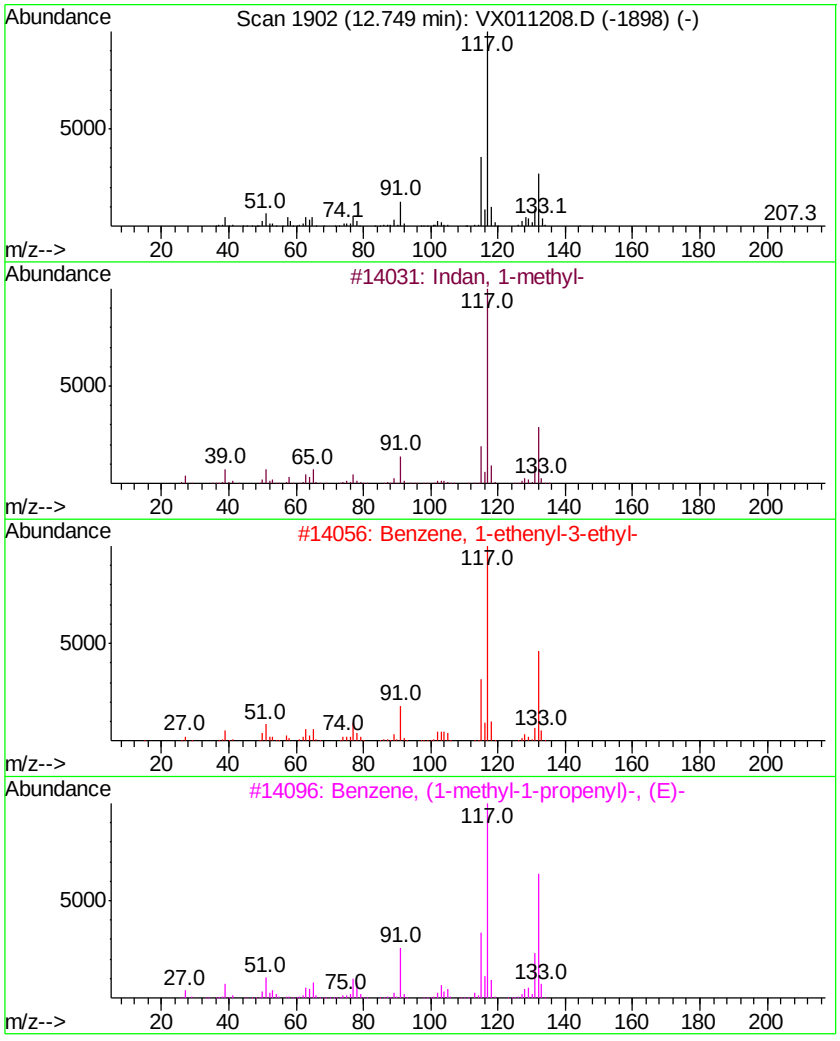
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

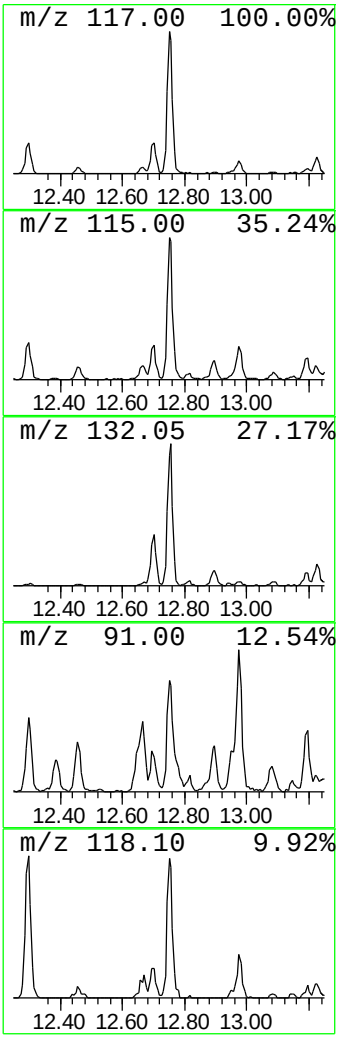
****

Peak Number 1 Indan, 1-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.75	12.67 ug/l	191272	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indan, 1-methyl-	132	C10H12	000767-58-8	90
2	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
3	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	87
4	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	86
5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	86





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX073019\
Data File : VX011208.D
Acq On : 30 Jul 2019 20:24
Operator : JC/SP
Sample : K4048-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : OWL-C7-GW

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

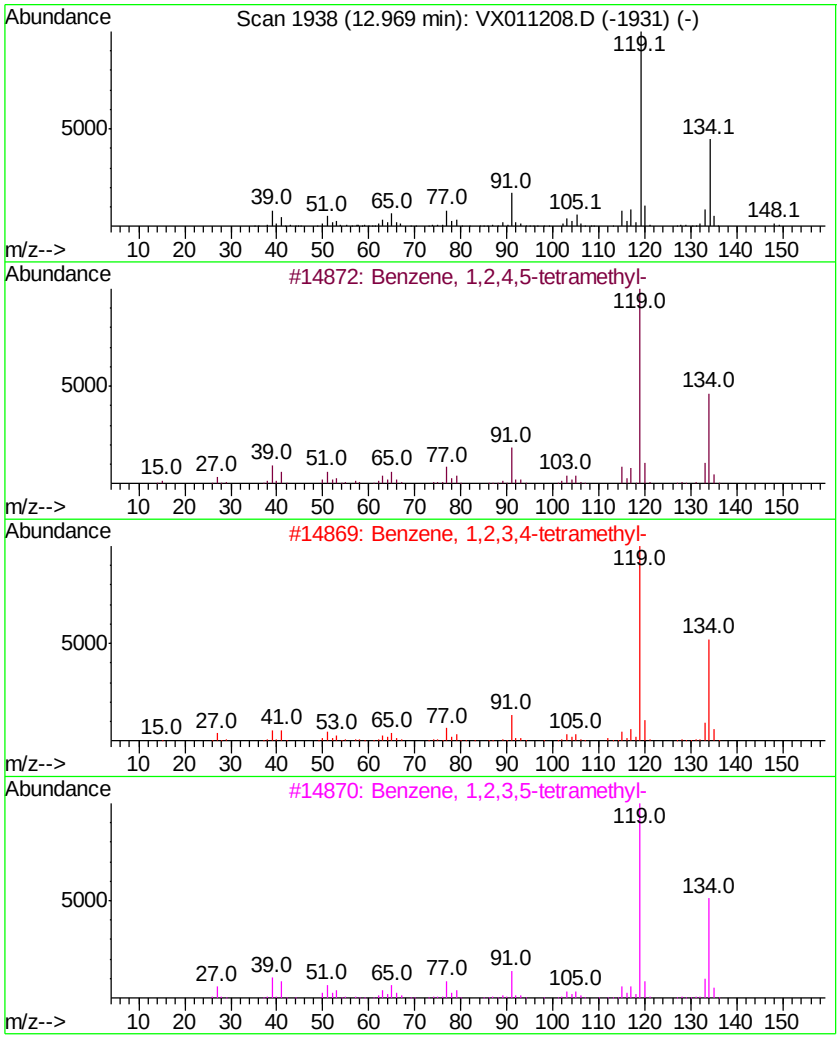
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

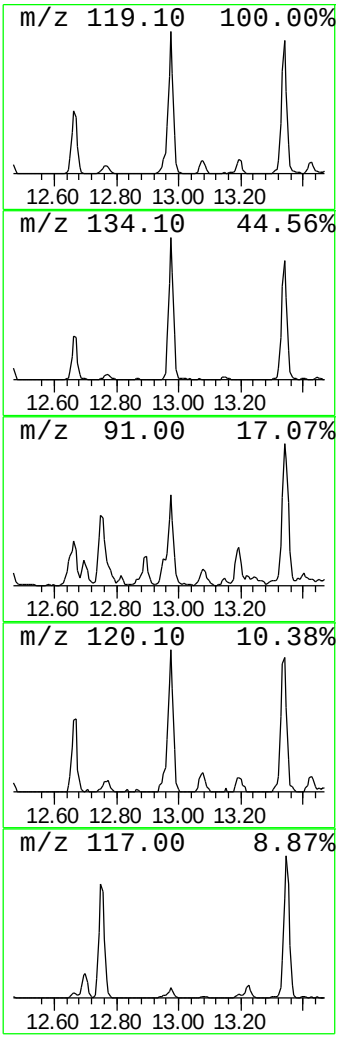
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.97	11.73 ug/l	177117	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	97
2	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
4	o-Cymene	134	C10H14	000527-84-4	95
5	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX073019\
Data File : VX011208.D
Acq On : 30 Jul 2019 20:24
Operator : JC/SP
Sample : K4048-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : OWL-C7-GW

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

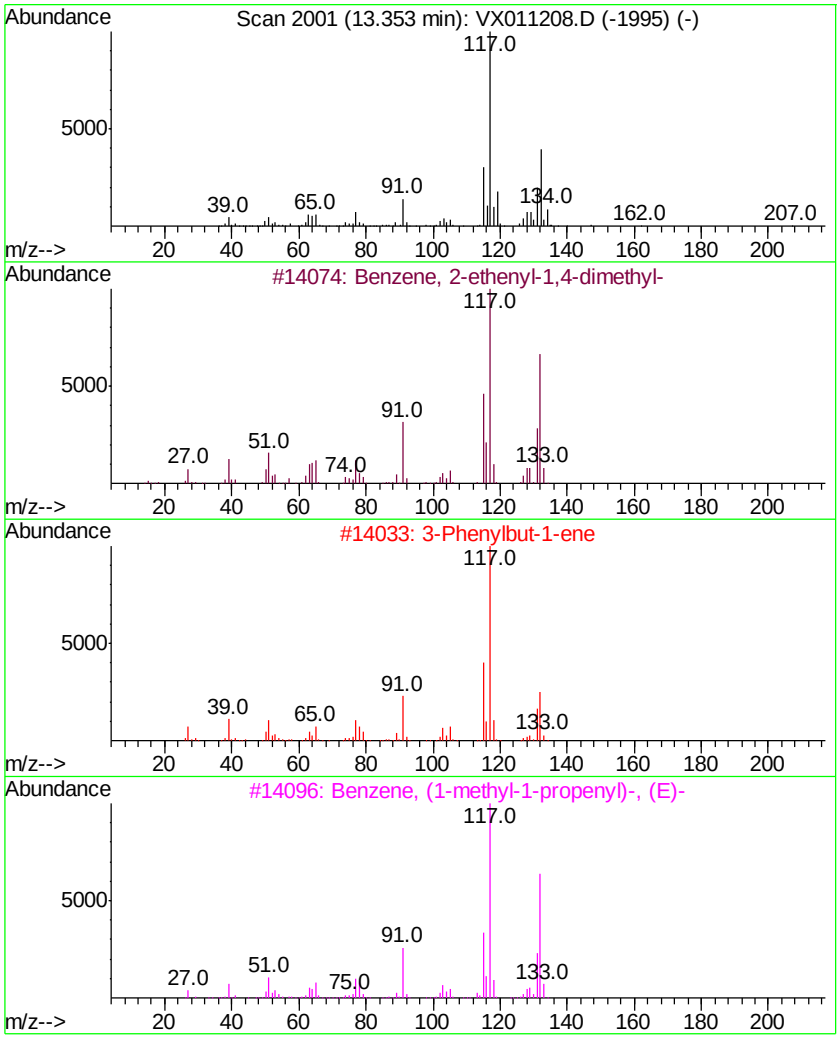
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

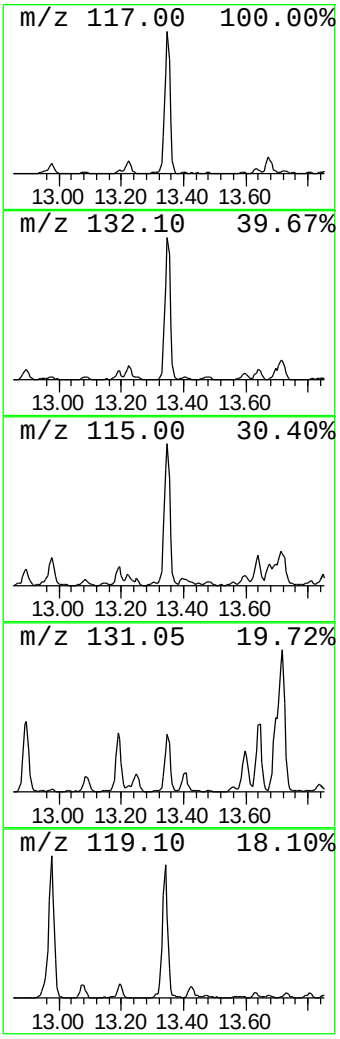
****

Peak Number 3 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	21.74 ug/l	328189	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
2	3-Phenylbut-1-ene	132	C10H12	000934-10-1	70
3	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	70
4	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	70
5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	64





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX073019\
Data File : VX011208.D
Acq On : 30 Jul 2019 20:24
Operator : JC/SP
Sample : K4048-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : OWL-C7-GW

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

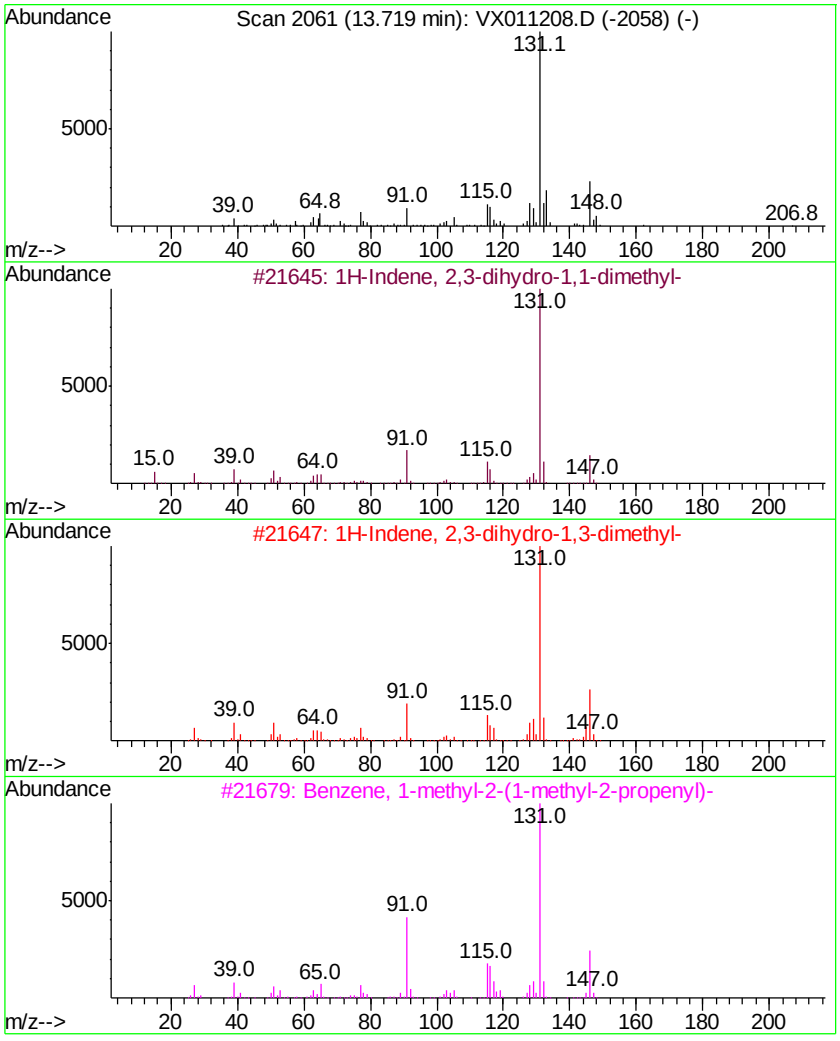
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

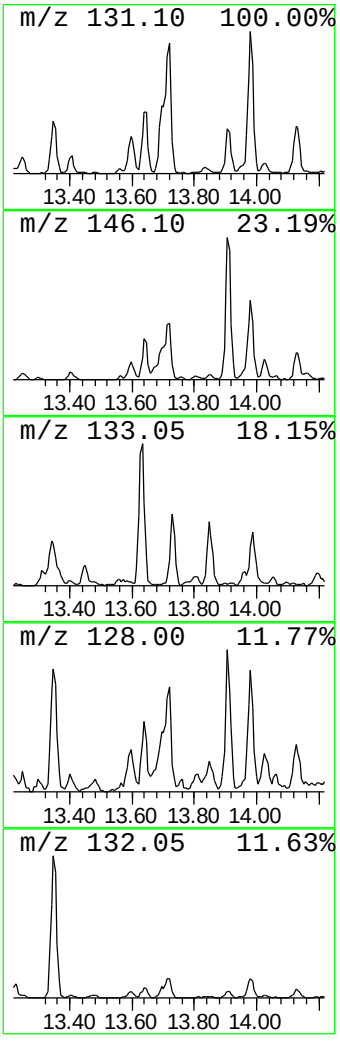
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.72	7.94 ug/l	119864	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	87
2	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	87
3	Benzene, 1-methyl-2-(1-methyl-2-...	146	C11H14	097664-19-2	83
4	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	81
5	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	81





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX073019\
Data File : VX011208.D
Acq On : 30 Jul 2019 20:24
Operator : JC/SP
Sample : K4048-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : OWL-C7-GW

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

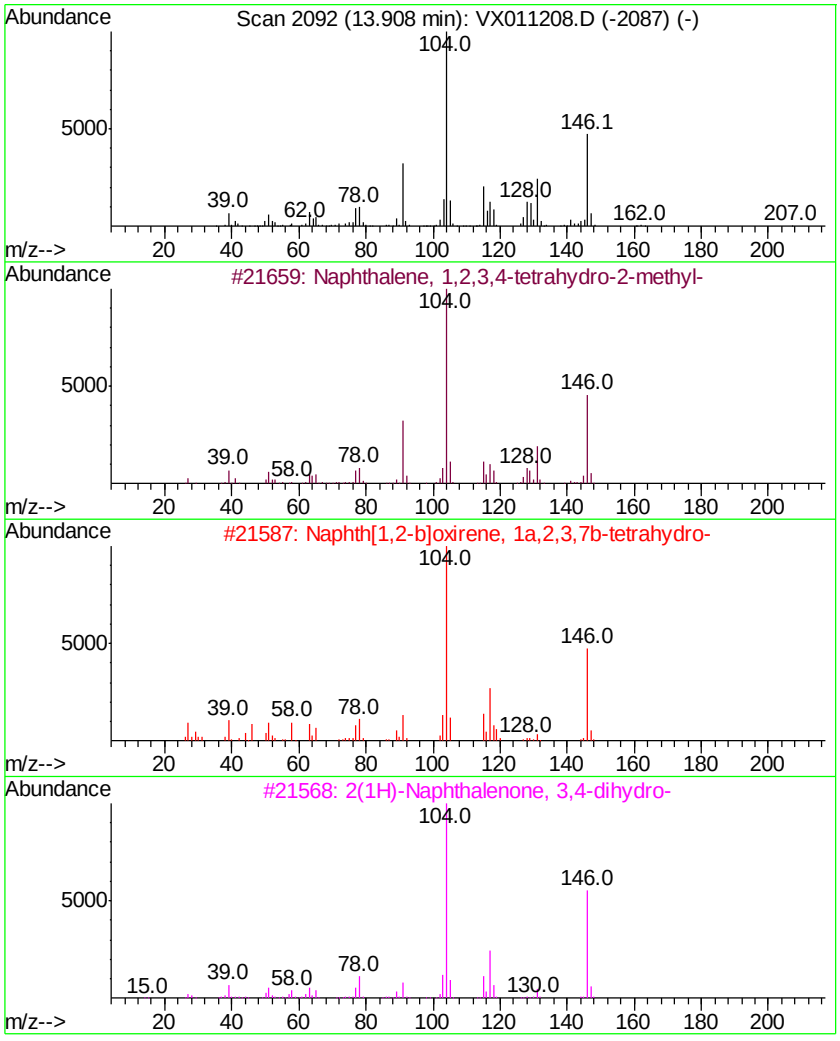
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

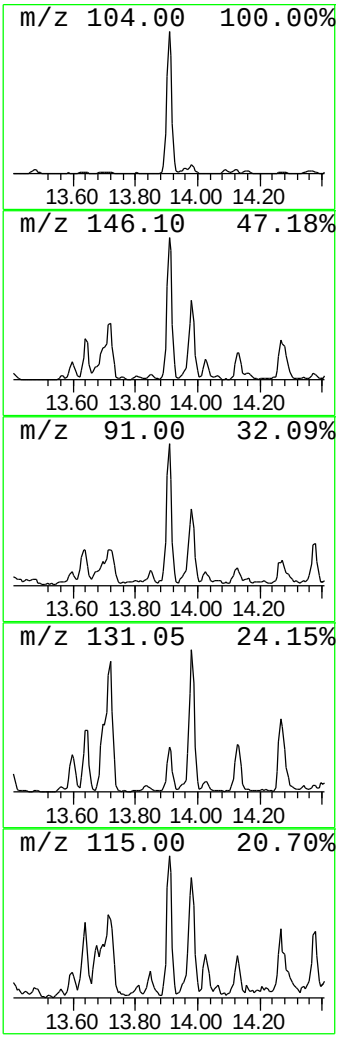
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.91	11.92 ug/l	180006	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	93
2	Naphth[1,2-b]oxirene, 1a,2,3,7b-...	146	C10H10O	002461-34-9	70
3	2(1H)-Naphthalenone, 3,4-dihydro-	146	C10H10O	000530-93-8	62
4	2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	43
5	7-Azabicyclo[4.2.2]deca-2,4,9-tr...	147	C9H9NO	017198-06-0	42





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX073019\
Data File : VX011208.D
Acq On : 30 Jul 2019 20:24
Operator : JC/SP
Sample : K4048-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : OWL-C7-GW

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

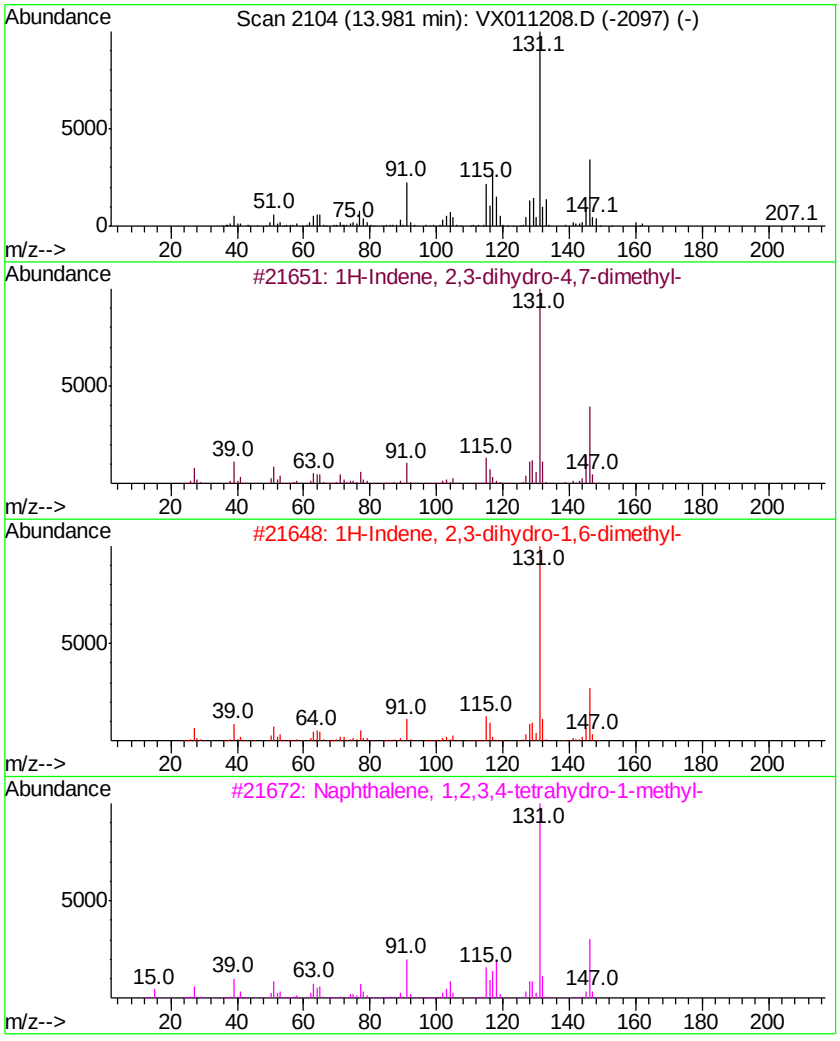
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

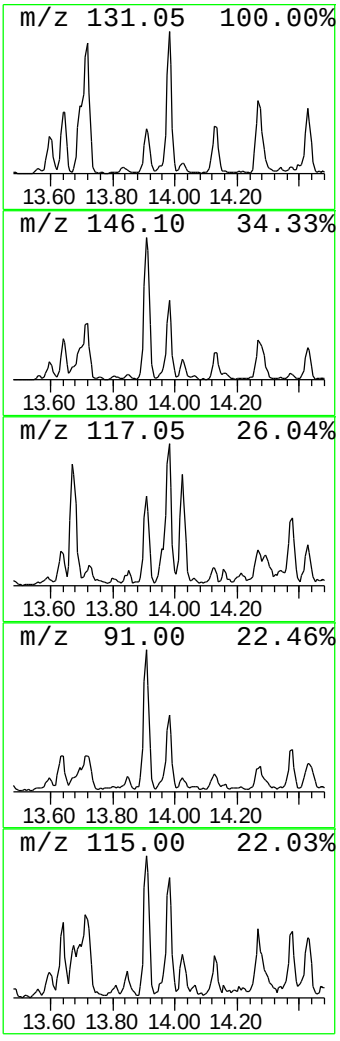
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.98	10.28 ug/l	155142	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	94
2	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	93
3	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	93
4	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	93
5	1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	91





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX073019\
Data File : VX011208.D
Acq On : 30 Jul 2019 20:24
Operator : JC/SP
Sample : K4048-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : OWL-C7-GW

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

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

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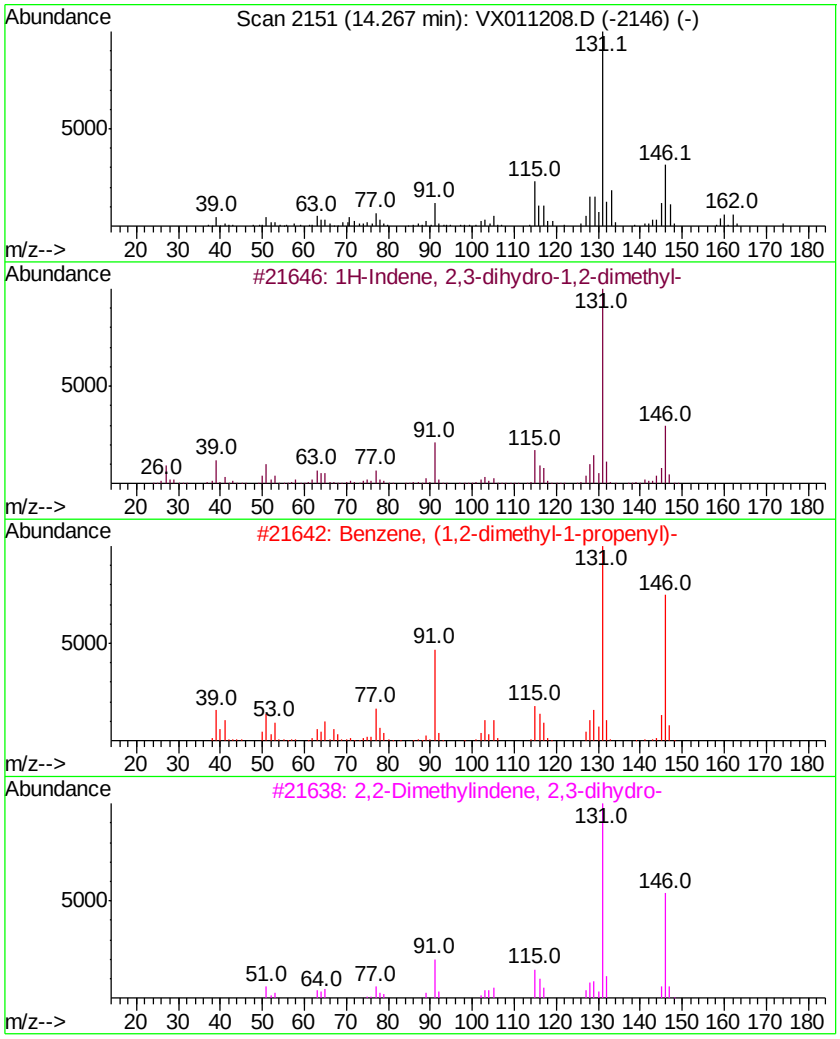
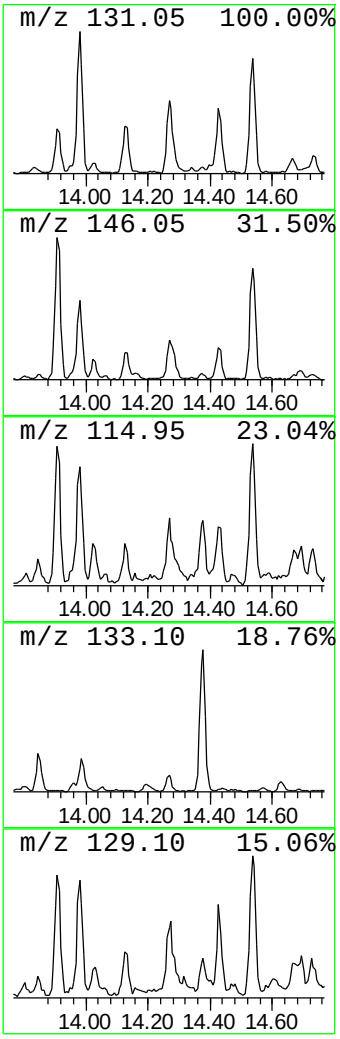
Peak Number 7 1H-Indene, 2,3-dihydro-1,2-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.27	6.88 ug/l	103930	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	93
2	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	90
3	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	87
4	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	87
5	Bicyclo[4.2.0]octa-1,3,5-triene,...	146	C11H14	027087-54-3	87





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX073019\
Data File : VX011208.D
Acq On : 30 Jul 2019 20:24
Operator : JC/SP
Sample : K4048-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : OWL-C7-GW

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

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

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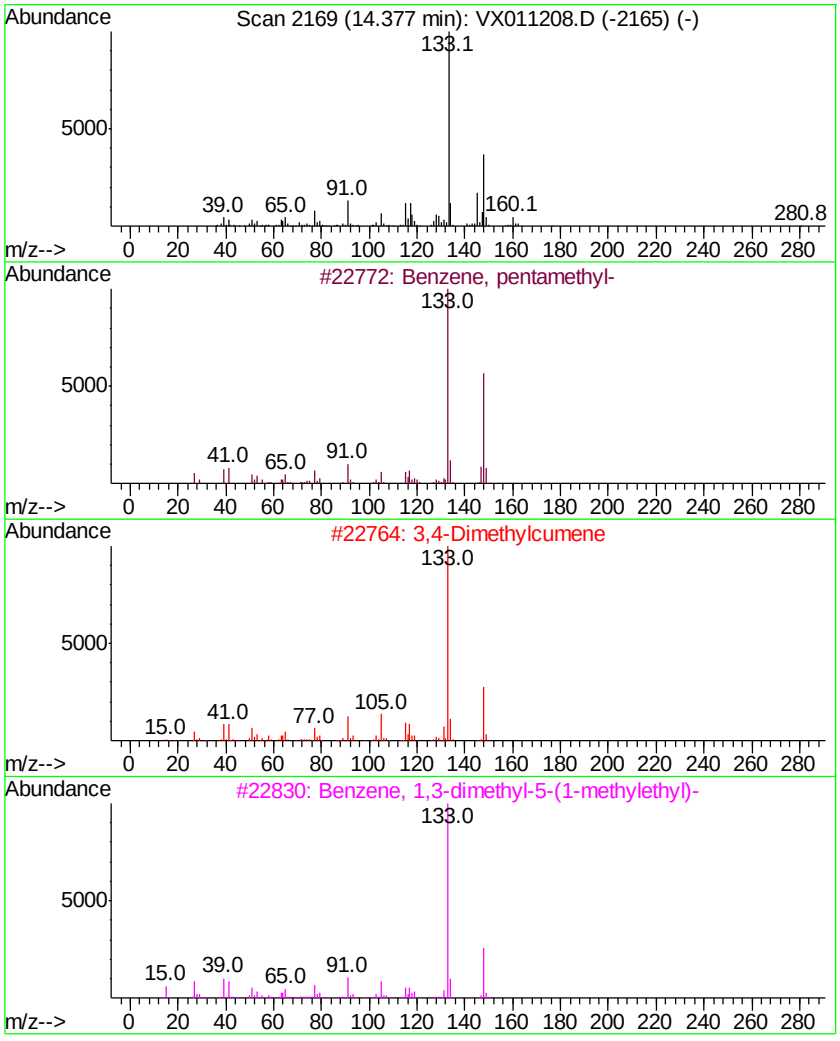
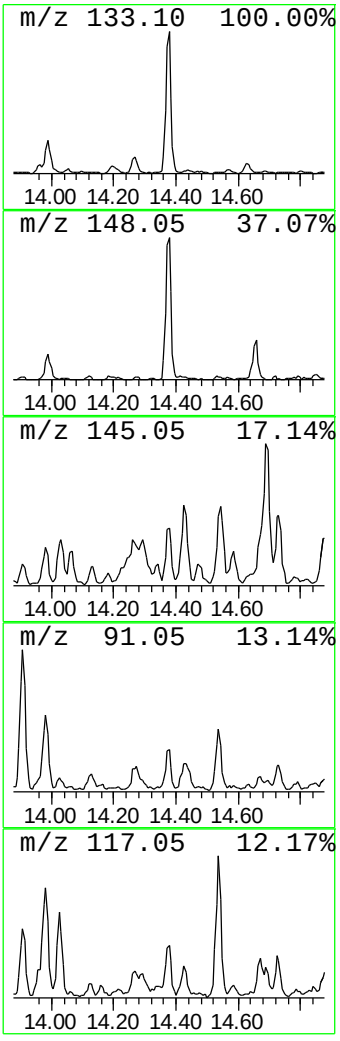
Peak Number 8 Benzene, pentamethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.38	6.65 ug/l	100318	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, pentamethyl-	148	C11H16	000700-12-9	87
2	3,4-Dimethylcumene	148	C11H16	1000370-34-1	81
3	Benzene, 1,3-dimethyl-5-(1-methy...	148	C11H16	004706-90-5	72
4	2-tert-Butyltoluene	148	C11H16	001074-92-6	72
5	Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	72





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX073019\
Data File : VX011208.D
Acq On : 30 Jul 2019 20:24
Operator : JC/SP
Sample : K4048-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : OWL-C7-GW

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

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

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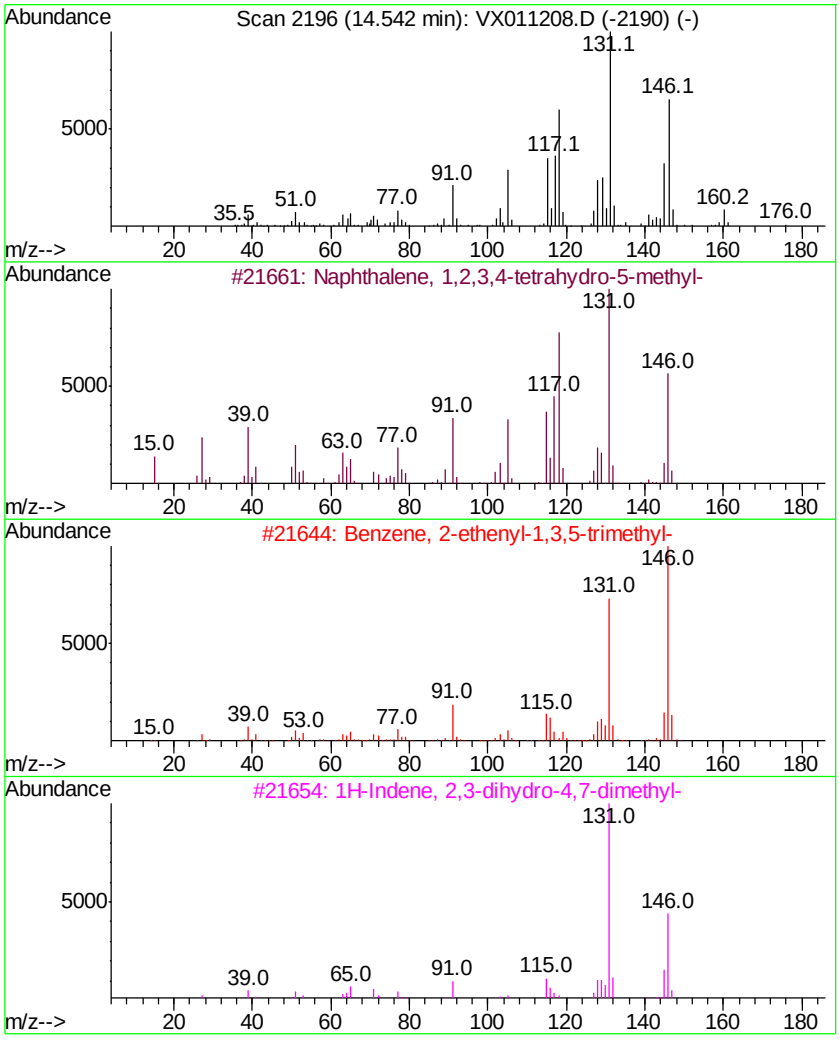
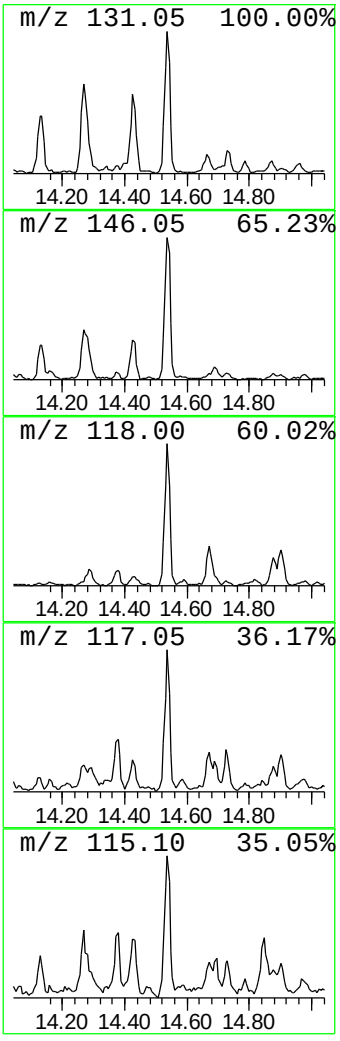
Peak Number 9 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.54	12.19 ug/l	184012	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	91
2	Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	64
3	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	64
4	1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	62
5	2-Propyn-1-ol, 3-(4-methylphenyl)-	146	C10H10O	016017-24-6	62





Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX073019\
Data File : VX011208.D
Acq On : 30 Jul 2019 20:24
Operator : JC/SP
Sample : K4048-02
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : OWL-C7-GW

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

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

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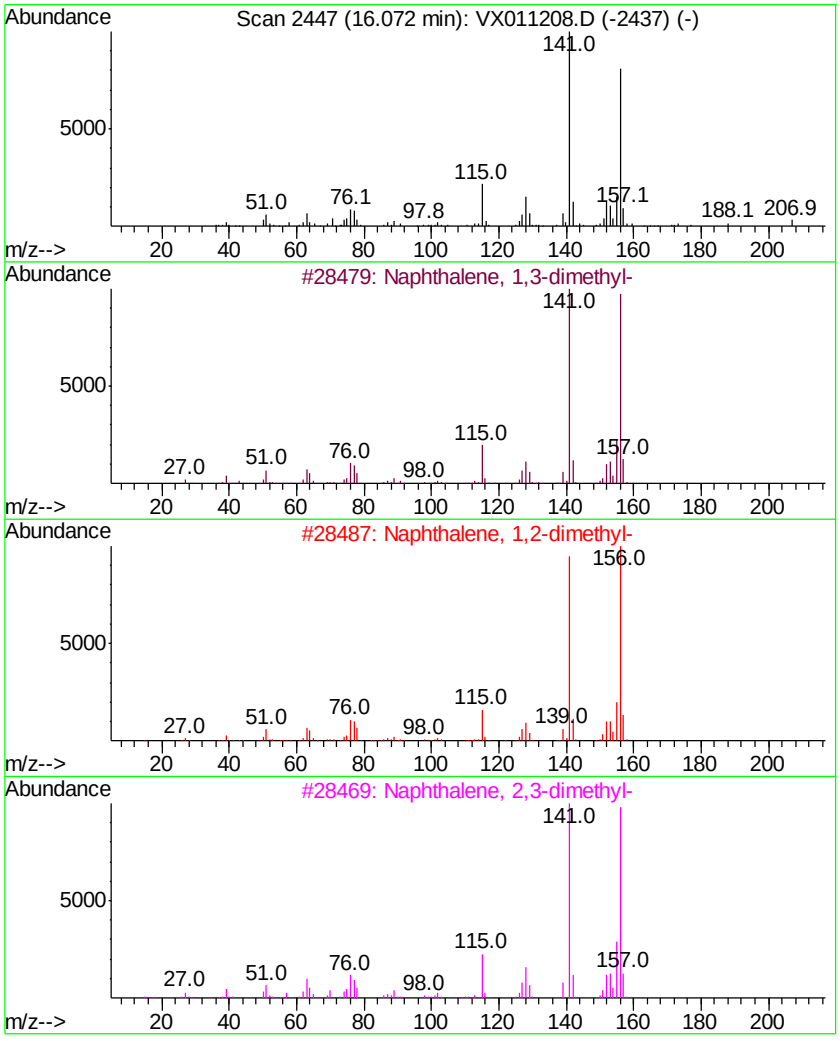
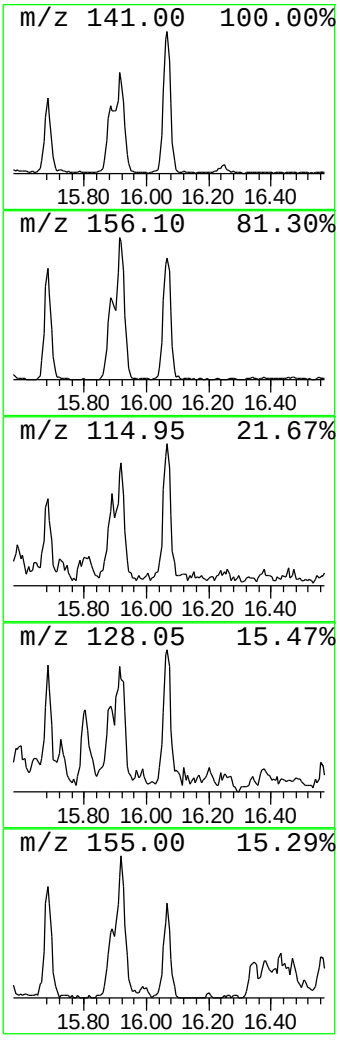
Peak Number 10 Naphthalene, 1,3-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.07	7.21 ug/l	108783	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	98
2	Naphthalene, 1,2-dimethyl-	156	C12H12	000573-98-8	96
3	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	95
4	Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	94
5	Naphthalene, 1,8-dimethyl-	156	C12H12	000569-41-5	94





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX073019\
Data File : VX011208.D
Acq On : 30 Jul 2019 20:24
Operator : JC/SP
Sample : K4048-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : OWL-C7-GW

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.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
Indan, 1-methyl-	12.75	12.7	ug/l	191272	4	12.08	754795	50.0
Benzene, 1,2,4,5-...	12.97	11.7	ug/l	177117	4	12.08	754795	50.0
Benzene, 2-etheny...	13.35	21.7	ug/l	328189	4	12.08	754795	50.0
1H-Indene, 2,3-di...	13.72	7.9	ug/l	119864	4	12.08	754795	50.0
Naphthalene, 1,2,...	13.91	11.9	ug/l	180006	4	12.08	754795	50.0
1H-Indene, 2,3-di...	13.98	10.3	ug/l	155142	4	12.08	754795	50.0
1H-Indene, 2,3-di...	14.27	6.9	ug/l	103930	4	12.08	754795	50.0
Benzene, pentamet...	14.38	6.7	ug/l	100318	4	12.08	754795	50.0
Naphthalene, 1,2,...	14.54	12.2	ug/l	184012	4	12.08	754795	50.0
Naphthalene, 1,3-...	16.07	7.2	ug/l	108783	4	12.08	754795	50.0