LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN121318\
Data File : VN052857.D
Acq On : 13 Dec 2018 12:01
Operator : MD\SY
Sample : J6319-01
Misc : 4.98g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 7 Sample Multiplier: 28

Instrument : MSVOA_N
ClientSampleId : 055_BT-1-1

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_N\METHODS\82N121218W.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	4.584	827	874	927	rBV3	264201	1126589	2.16%	0.133%
2	5.047	991	1018	1050	rBV2	223993	820091	1.58%	0.097%
3	7.651	1781	1828	1845	rBV4	1669900	8324356	15.99%	0.986%
4	7.741	1845	1856	1878	rVB3	618436	1653108	3.18%	0.196%
5	7.873	1878	1897	1911	rBV	1930432	4718858	9.07%	0.559%
6	8.024	1929	1944	1967	rVB	659312	1599091	3.07%	0.189%
7	8.156	1967	1985	1997	rBV2	1077640	2705701	5.20%	0.320%
8	8.233	1997	2009	2019	rVV3	1045237	2556018	4.91%	0.303%
9	8.301	2020	2030	2052	rVB	1090998	2648899	5.09%	0.314%
10	8.432	2052	2071	2095	rVB4	412597	1150105	2.21%	0.136%
11	8.584	2098	2118	2140	rBV3	2568448	6370477	12.24%	0.754%
12	8.818	2178	2191	2201	rBV	968265	2117140	4.07%	0.251%
13	9.079	2245	2272	2284	rBV4	6316068	17062406	32.78%	2.020%
14	9.143	2284	2292	2307	rVB	1453312	2879218	5.53%	0.341%
15	9.275	2307	2333	2344	rBV4	1305719	3370249	6.48%	0.399%
16	9.365	2345	2361	2376	rVB2	2126393	5308122	10.20%	0.629%
17	9.509	2395	2406	2418	rBV4	2593742	4993846	9.60%	0.591%
18	9.616	2430	2439	2445	rBV2	993752	1680232	3.23%	0.199%
19	9.664	2446	2454	2462	rVV2	2056486	3323259	6.39%	0.393%
20	9.728	2463	2474	2498	rVB2	11229958	29688038	57.04%	3.515%
21	9.866	2498	2517	2540	rBV	8415350	21125889	40.59%	2.501%
22	9.973	2541	2550	2553	rBV4	596068	932128	1.79%	0.110%
23	10.024	2555	2566	2574	rVB2	3668788	6373251	12.25%	0.755%
24	10.085	2574	2585	2594	rBV2	13625659	26479398	50.88%	3.135%
25	10.214	2610	2625	2630	rBV2	2384170	4752044	9.13%	0.563%
26	10.281	2631	2646	2659	rVB7	6451162	18528999	35.60%	2.194%
27	10.349	2660	2667	2670	rBV3	612769	860346	1.65%	0.102%
28	10.387	2672	2679	2685	rVV4	2379812	3957659	7.60%	0.469%
29	10.432	2685	2693	2709	rVB	6106344	10906082	20.95%	1.291%
30	10.526	2710	2722	2734	rBV4	6573427	15043899	28.90%	1.781%
31	10.625	2745	2753	2763	rVB4	3364582	5789953	11.12%	0.686%
32	10.718	2763	2782	2792	rBV	9239914	16778420	32.24%	1.987%
33	10.821	2793	2814	2826	rBV	6298407	15751320	30.26%	1.865%
34	10.886	2826	2834	2839	rBV6	2716888	4447154	8.54%	0.527%

LSC Area Percent Report

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Operator : MD\SY
Sample : J6319-01
Misc : 4.98g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 7 Sample Multiplier: 28

Instrument : MSVOA_N
ClientSampleId : 055_BT-1-1

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_N\METHODS\82N121218W.M
Title : SW846 8260

35	10.953	2848	2855	2863	rVB	8315865	12136549	23.32%	1.437%
36	11.005	2863	2871	2882	rVB2	11601857	20269986	38.95%	2.400%
37	11.063	2882	2889	2897	rVB3	2874850	4105950	7.89%	0.486%
38	11.120	2898	2907	2916	rVB6	6207258	10299033	19.79%	1.219%
39	11.201	2916	2932	2952	rVB5	21111675	52046283	100.00%	6.163%
40	11.300	2952	2963	2988	rBV	16401425	31529179	60.58%	3.733%
41	11.410	2989	2997	3006	rBV3	3970320	7133340	13.71%	0.845%
42	11.593	3045	3054	3063	rBV5	4730950	9123134	17.53%	1.080%
43	11.654	3063	3073	3081	rVV	9561054	17576424	33.77%	2.081%
44	11.699	3083	3087	3096	rVB2	6590684	9123372	17.53%	1.080%
45	11.754	3096	3104	3112	rBV7	1867938	3449963	6.63%	0.408%
46	11.812	3114	3122	3128	rBV5	1974105	3287357	6.32%	0.389%
47	11.857	3129	3136	3144	rVB5	2304216	3519980	6.76%	0.417%
48	11.924	3145	3157	3169	rBV6	8981672	21116096	40.57%	2.500%
49	12.046	3188	3195	3202	rVB	8734114	11402601	21.91%	1.350%
50	12.120	3210	3218	3224	rBV4	6525720	10308248	19.81%	1.221%
51	12.162	3224	3231	3241	rVB5	9886050	17373492	33.38%	2.057%
52	12.445	3313	3319	3327	rVB2	6566074	8556872	16.44%	1.013%
53	12.564	3349	3356	3360	rBV3	4346478	6546223	12.58%	0.775%
54	12.725	3395	3406	3416	rBV6	7714944	16741738	32.17%	1.982%
55	12.924	3461	3468	3473	rBV4	4094104	6418329	12.33%	0.760%
56	12.963	3474	3480	3493	rVB3	9028906	14337117	27.55%	1.698%
57	13.156	3533	3540	3542	rBV3	2163077	3010144	5.78%	0.356%
58	13.426	3618	3624	3635	rVB3	3211492	4991201	9.59%	0.591%
59	13.513	3644	3651	3657	rBV	6299054	9433132	18.12%	1.117%
60	13.596	3669	3677	3679	rBV2	4455777	5617293	10.79%	0.665%
61	13.667	3691	3699	3708	rVB5	6312622	10598546	20.36%	1.255%
62	13.889	3759	3768	3778	rVV	22414419	33976463	65.28%	4.023%
63	13.943	3780	3785	3792	rVV	3258638	4815238	9.25%	0.570%
64	13.995	3792	3801	3812	rVB2	16873130	27791836	53.40%	3.291%
65	14.210	3862	3868	3884	rVB	13312570	23512222	45.18%	2.784%
66	14.294	3886	3894	3901	rBV2	6141798	9848329	18.92%	1.166%
67	14.477	3943	3951	3961	rBV2	10214867	17820337	34.24%	2.110%
68	14.593	3975	3987	3997	rBV4	19557112	39778600	76.43%	4.710%
69	14.651	3998	4005	4015	rVB	5894217	9560930	18.37%	1.132%
70	14.850	4050	4067	4074	rBV4	12102997	26146968	50.24%	3.096%
71	14.959	4091	4101	4113	rVB2	10490969	21777390	41.84%	2.579%

LSC Area Percent Report

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Misc : 4.98g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 7 Sample Multiplier: 28

Instrument : MSVOA_N
ClientSampleId : 055_BT-1-1

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_N\METHODS\82N121218W.M
Title : SW846 8260

72	15.111	4139	4148	4156	rBV6	1568051	2994615	5.75%	0.355%
73	15.239	4180	4188	4197	rBV3	4139778	7137186	13.71%	0.845%
74	15.291	4198	4204	4212	rVB2	2168677	3301050	6.34%	0.391%
75	15.419	4233	4244	4257	rBV	4824063	9463222	18.18%	1.120%
76	15.574	4277	4292	4313	rBV6	2741604	9873461	18.97%	1.169%
77	15.699	4324	4331	4340	rVB5	1169485	1882613	3.62%	0.223%
78	15.763	4342	4351	4363	rVB6	1853598	4050370	7.78%	0.480%
79	15.908	4386	4396	4406	rBV5	807196	1617616	3.11%	0.192%
80	16.352	4522	4534	4558	rVB	1433642	3337224	6.41%	0.395%

Sum of corrected areas: 844559597

LSC Report - Integrated Chromatogram

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Acq On : 13 Dec 2018 12:01
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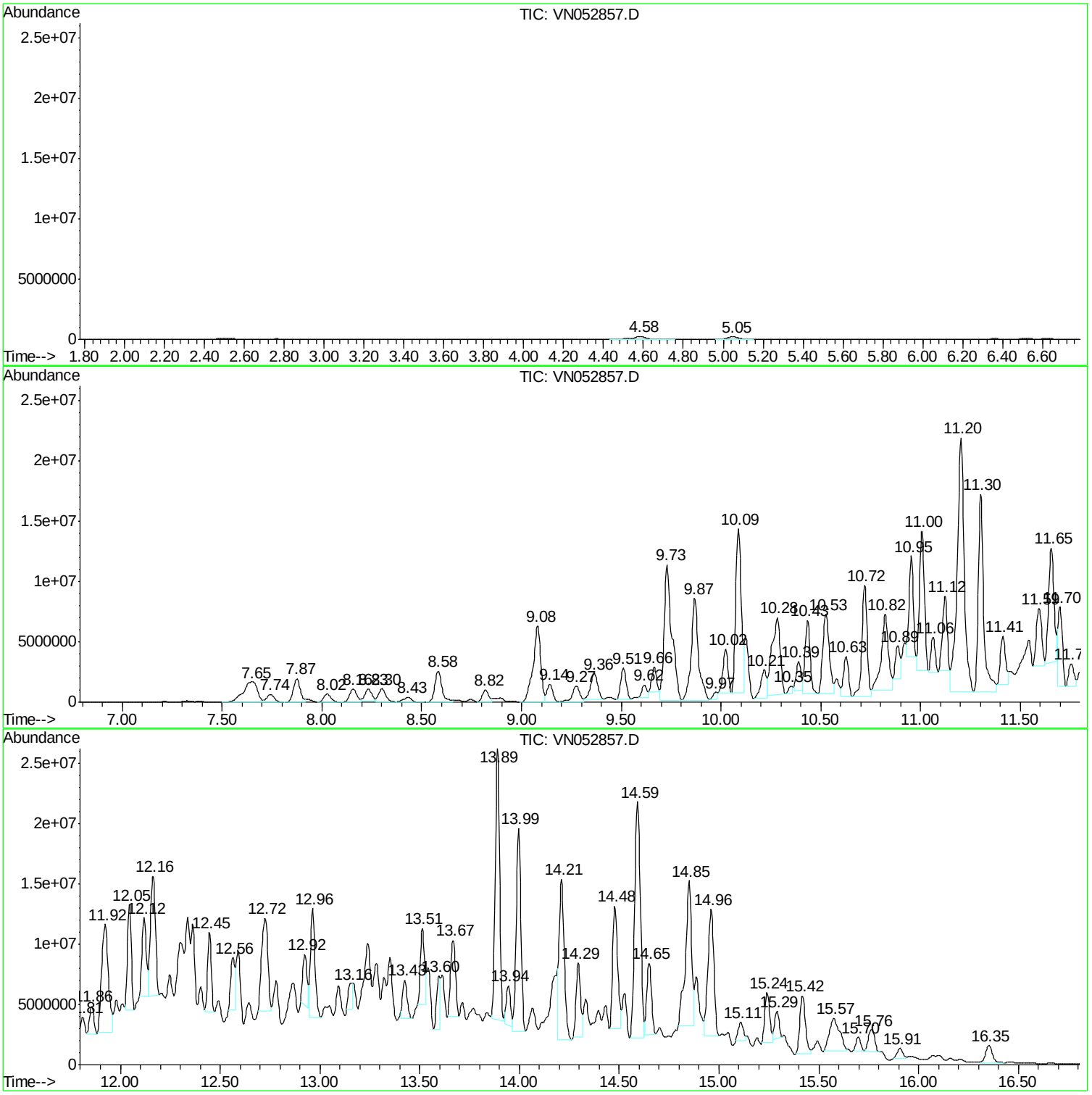
ALS Vial : 7 Sample Multiplier: 28

Instrument : MSVOA_N
ClientSampleId : 055_BT-1-1

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N121218W.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_N\DATA\VN121318\
Data File : VN052857.D
Acq On : 13 Dec 2018 12:01
Operator : MD\SY
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Misc : 4.98g/10mL/100uL/5.00mL/MSVOA_N/MEOH
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Instrument : MSVOA_N
ClientSampleId : 055_BT-1-1

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N121218W.M
Quant Title : SW846 8260

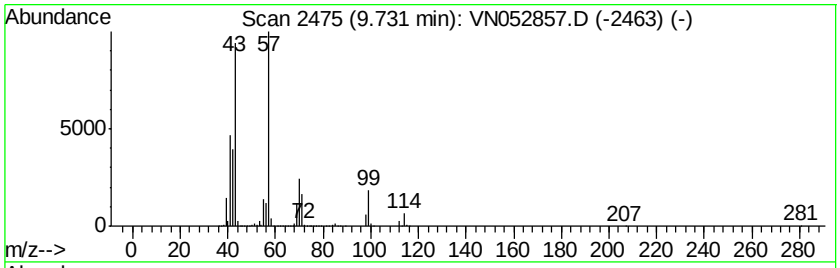
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

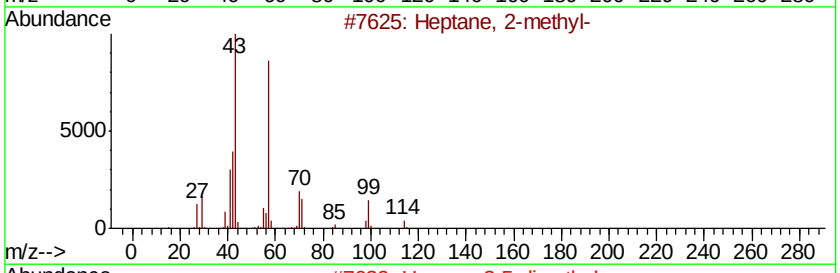
****

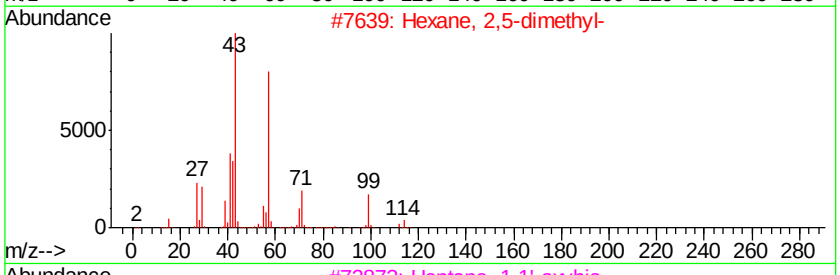
Peak Number 1 Heptane, 2-methyl- Concentration Rank 7

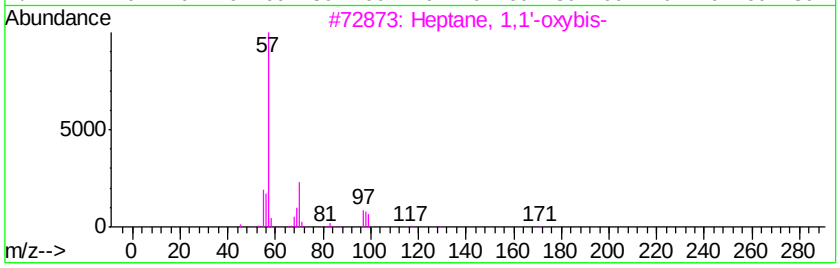
R.T.	EstConc	Area	Relative to ISTD	R.T.
9.73	233.01 ug/l	29688000	1,4-Difluorobenzene	8.58

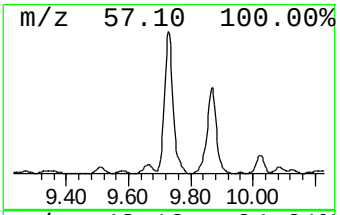
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 2-methyl-	114	C8H18	000592-27-8	94
2	Hexane, 2,5-dimethyl-	114	C8H18	000592-13-2	68
3	Heptane, 1,1'-oxybis-	214	C14H30O	000629-64-1	50
4	Oxalic acid, heptyl propyl ester	230	C12H22O4	1000309-26-0	47
5	Heptane, 3-ethyl-	128	C9H20	015869-80-4	38

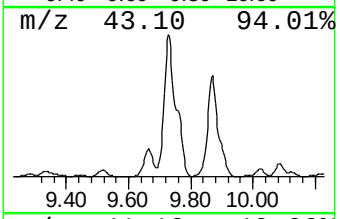


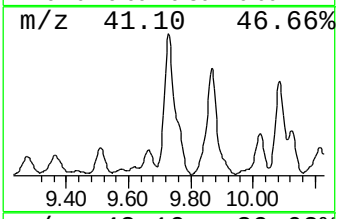


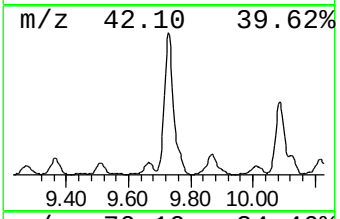


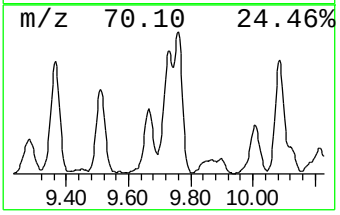












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Instrument : MSVOA_N
ClientSampleId : 055_BT-1-1

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N121218W.M
Quant Title : SW846 8260

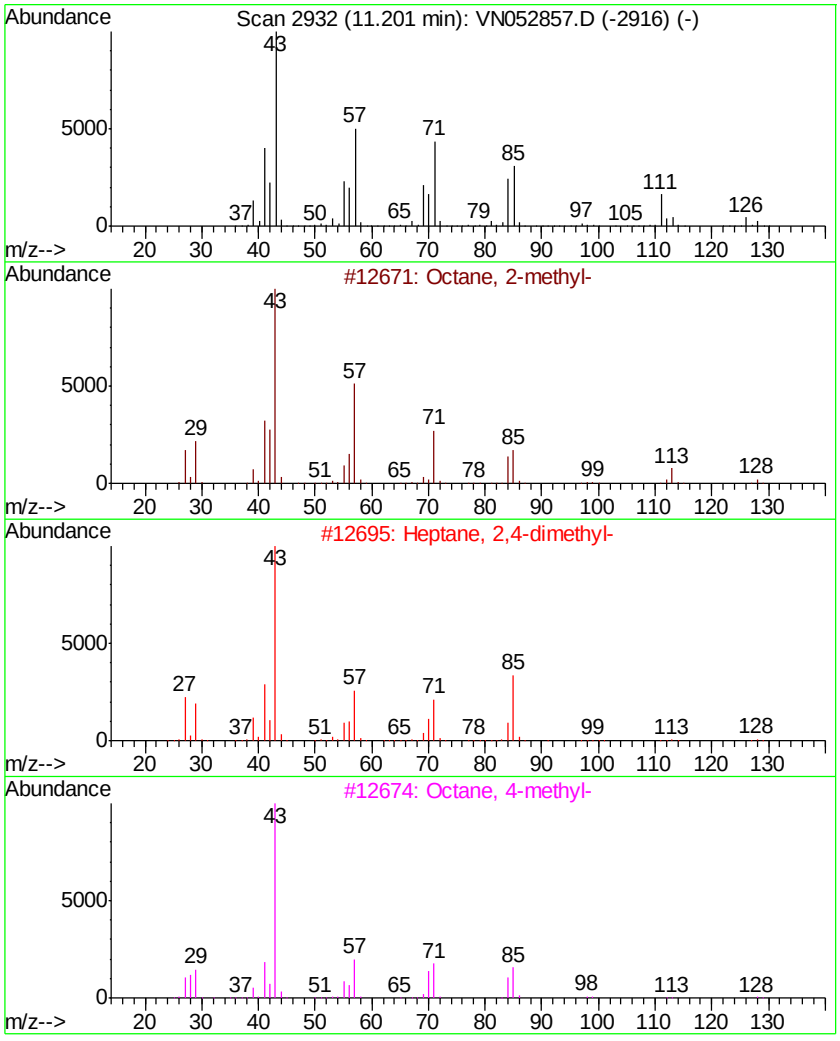
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

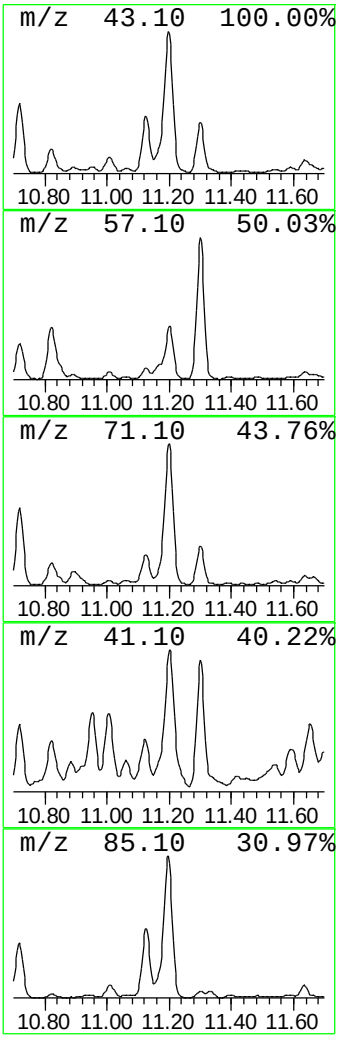
****

Peak Number 2 Octane, 2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.20	364.81 ug/l	52046300	Chlorobenzene-d5	11.41

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 2-methyl-	128	C9H20	003221-61-2	72
2	Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	52
3	Octane, 4-methyl-	128	C9H20	002216-34-4	47
4	Hexyl octyl ether	214	C14H30O	017071-54-4	47
5	1-Butanol, 2-ethyl-	102	C6H14O	000097-95-0	47





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Instrument : MSVOA_N
ClientSampleId : 055_BT-1-1

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N121218W.M
Quant Title : SW846 8260

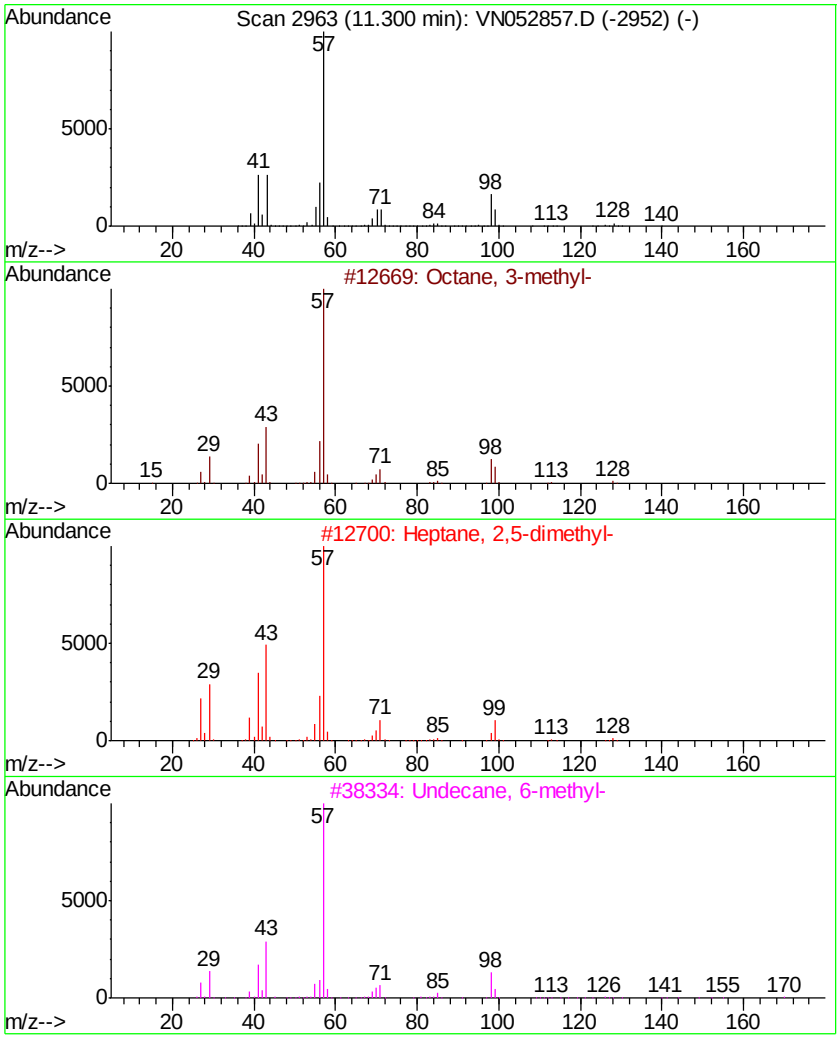
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

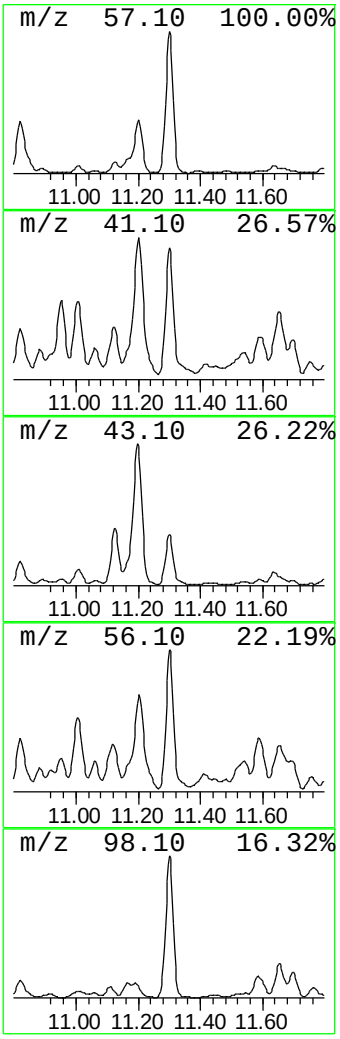
****

Peak Number 3 Octane, 3-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.30	221.00 ug/l	31529200	Chlorobenzene-d5	11.41

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 3-methyl-	128	C9H20	002216-33-3	90
2	Heptane, 2,5-dimethyl-	128	C9H20	002216-30-0	74
3	Undecane, 6-methyl-	170	C12H26	017302-33-9	53
4	Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	50
5	Heptane, 3-ethyl-	128	C9H20	015869-80-4	50





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ALS Vial : 7 Sample Multiplier: 28

Instrument : MSVOA_N
ClientSampleId : 055_BT-1-1

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N121218W.M
Quant Title : SW846 8260

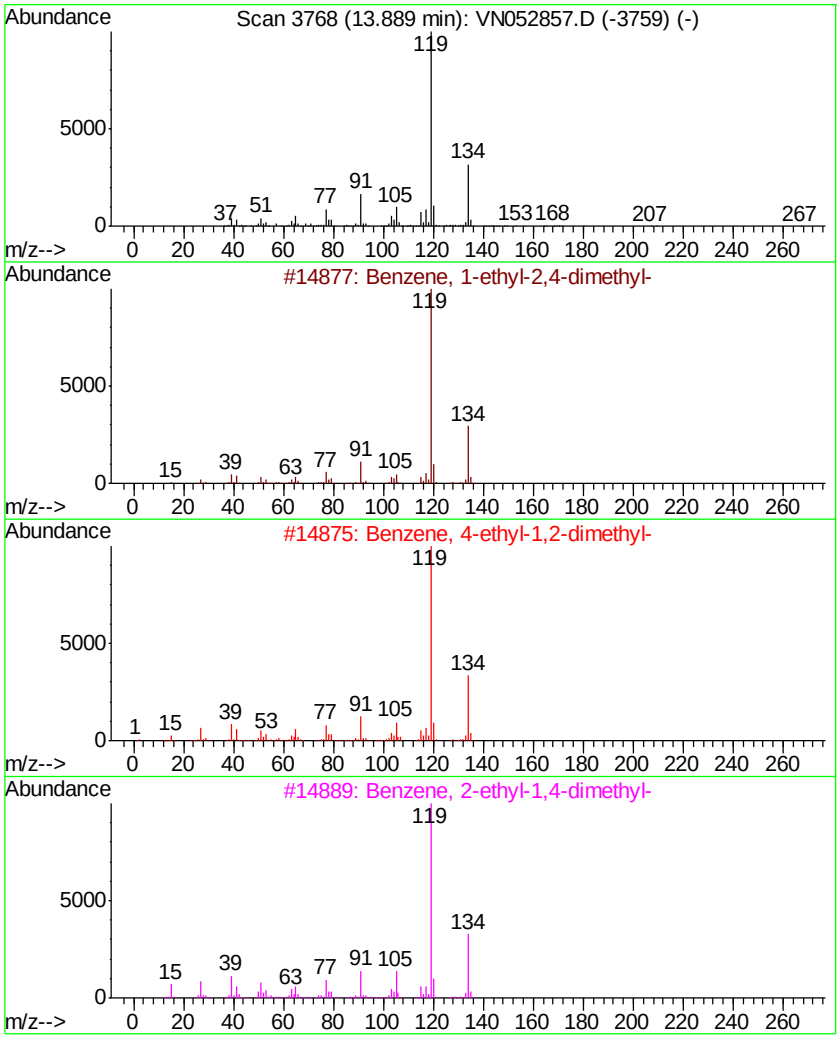
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

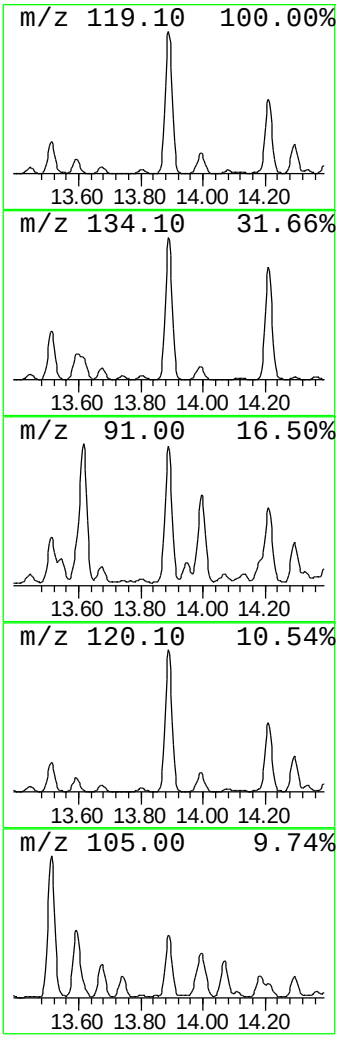
****

Peak Number 4 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.89	340.36 ug/l	33976500	1,4-Dichlorobenzene-d4	13.35

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	97
2	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
3	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
4	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
5	o-Cymene	134	C10H14	000527-84-4	95





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN121318\
Data File : VN052857.D
Acq On : 13 Dec 2018 12:01
Operator : MD\SY
Sample : J6319-01
Misc : 4.98g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 7 Sample Multiplier: 28

Instrument : MSVOA_N
ClientSampleId : 055_BT-1-1

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N121218W.M
Quant Title : SW846 8260

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

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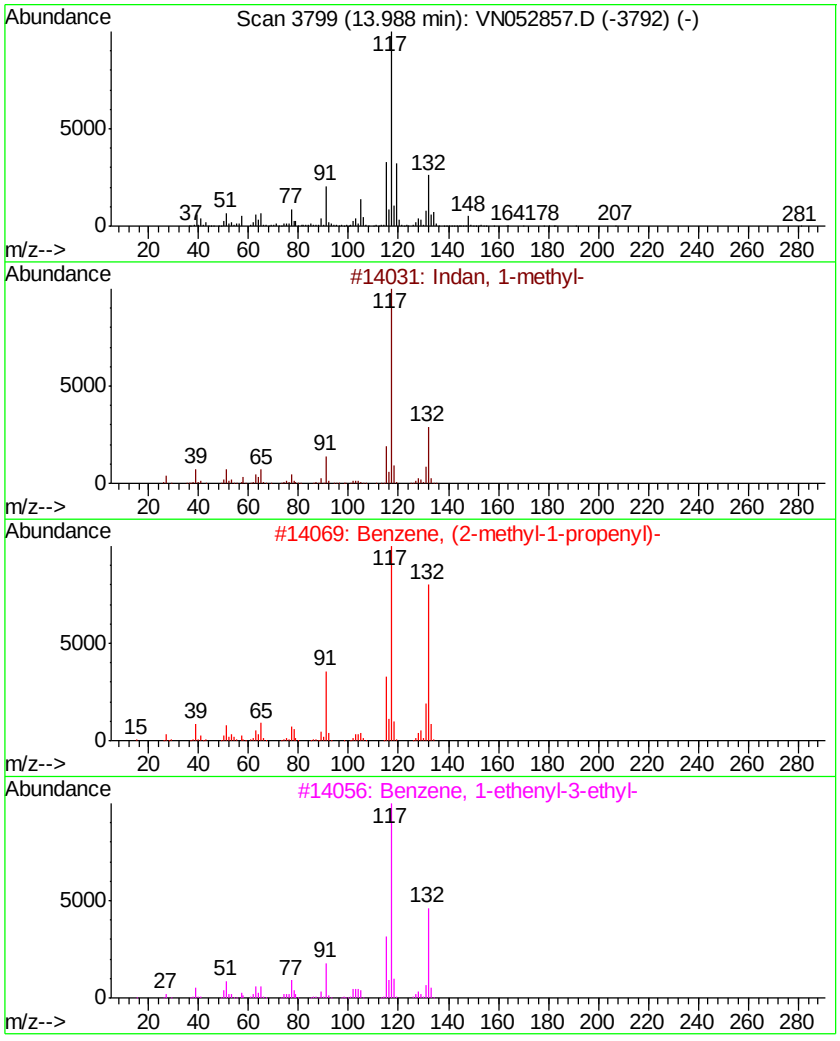
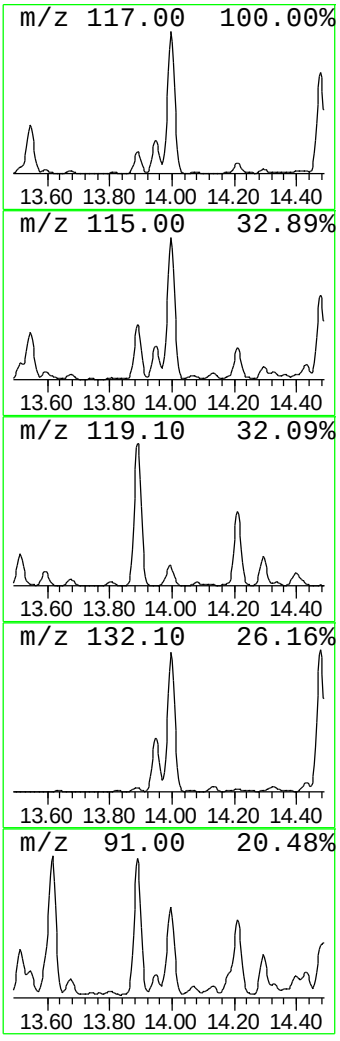
Peak Number 5 Indan, 1-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.99	278.41 ug/l	27791800	1,4-Dichlorobenzene-d4	13.35

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indan, 1-methyl-	132	C10H12	000767-58-8	93
2	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	74
3	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	74
4	Benzene, 2-butenyl-	132	C10H12	001560-06-1	72
5	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	68





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN121318\
Data File : VN052857.D
Acq On : 13 Dec 2018 12:01
Operator : MD\SY
Sample : J6319-01
Misc : 4.98g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 7 Sample Multiplier: 28

Instrument : MSVOA_N
ClientSampleId : 055_BT-1-1

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N121218W.M
Quant Title : SW846 8260

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

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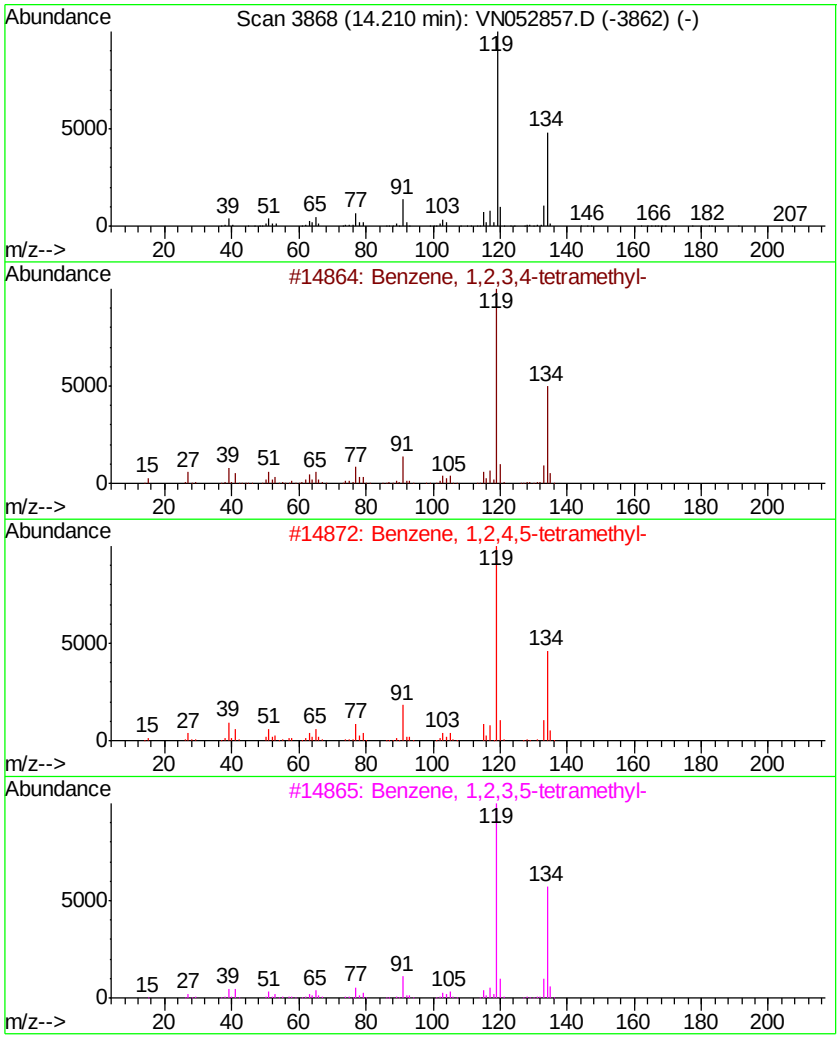
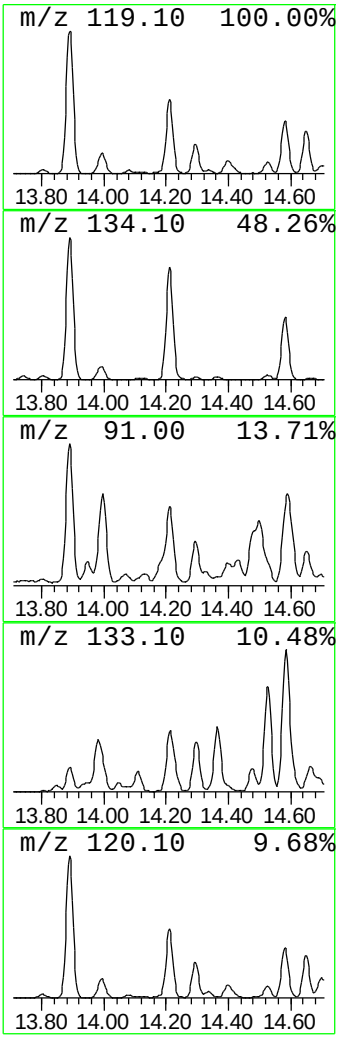
Peak Number 6 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.21	235.54 ug/l	23512200	1,4-Dichlorobenzene-d4	13.35

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	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-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91





Library Search Compound Report

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Data File : VN052857.D
Acq On : 13 Dec 2018 12:01
Operator : MD\SY
Sample : J6319-01
Misc : 4.98g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 7 Sample Multiplier: 28

Instrument : MSVOA_N
ClientSampleId : 055_BT-1-1

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N121218W.M
Quant Title : SW846 8260

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

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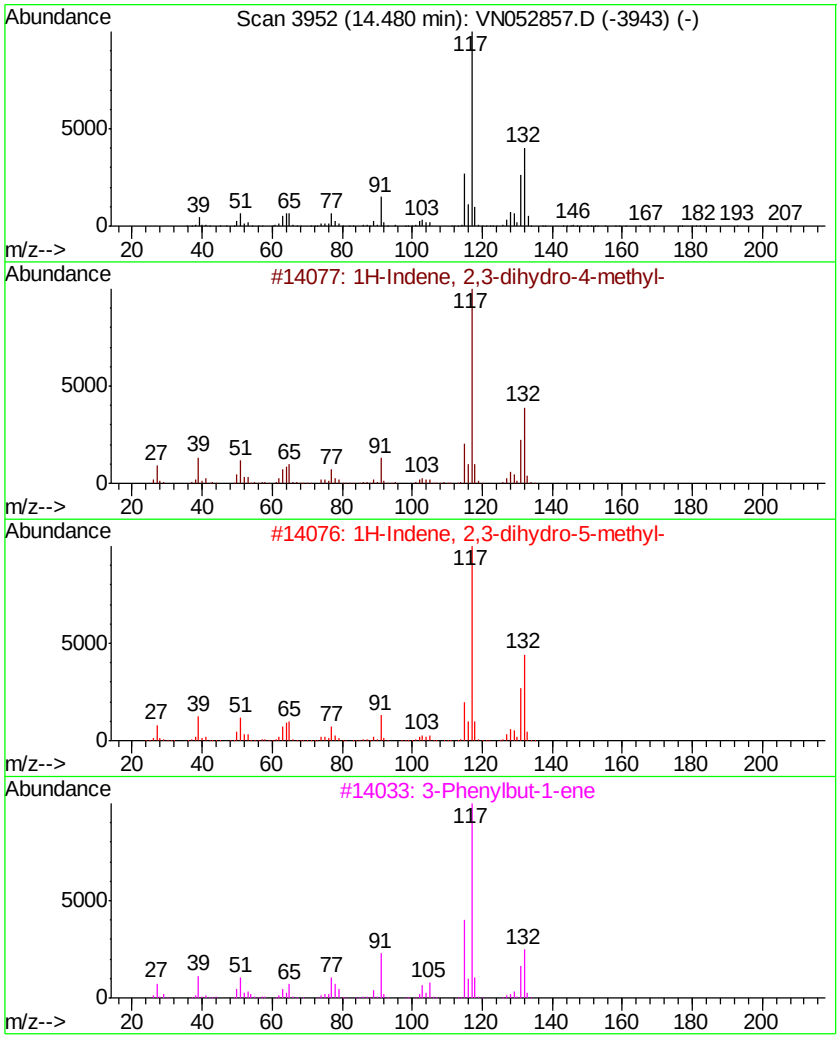
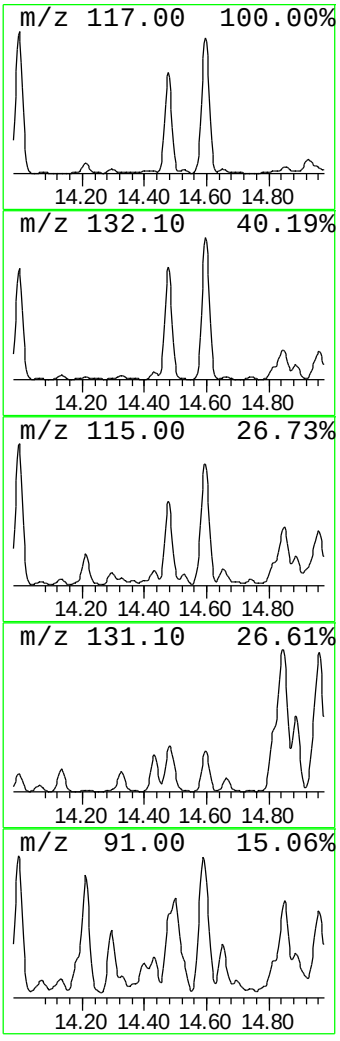
Peak Number 7 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.48	178.52 ug/l	17820300	1,4-Dichlorobenzene-d4	13.35

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	94
2	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
3	3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
4	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	83





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN121318\
Data File : VN052857.D
Acq On : 13 Dec 2018 12:01
Operator : MD\SY
Sample : J6319-01
Misc : 4.98g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 7 Sample Multiplier: 28

Instrument : MSVOA_N
ClientSampleId : 055_BT-1-1

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N121218W.M
Quant Title : SW846 8260

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

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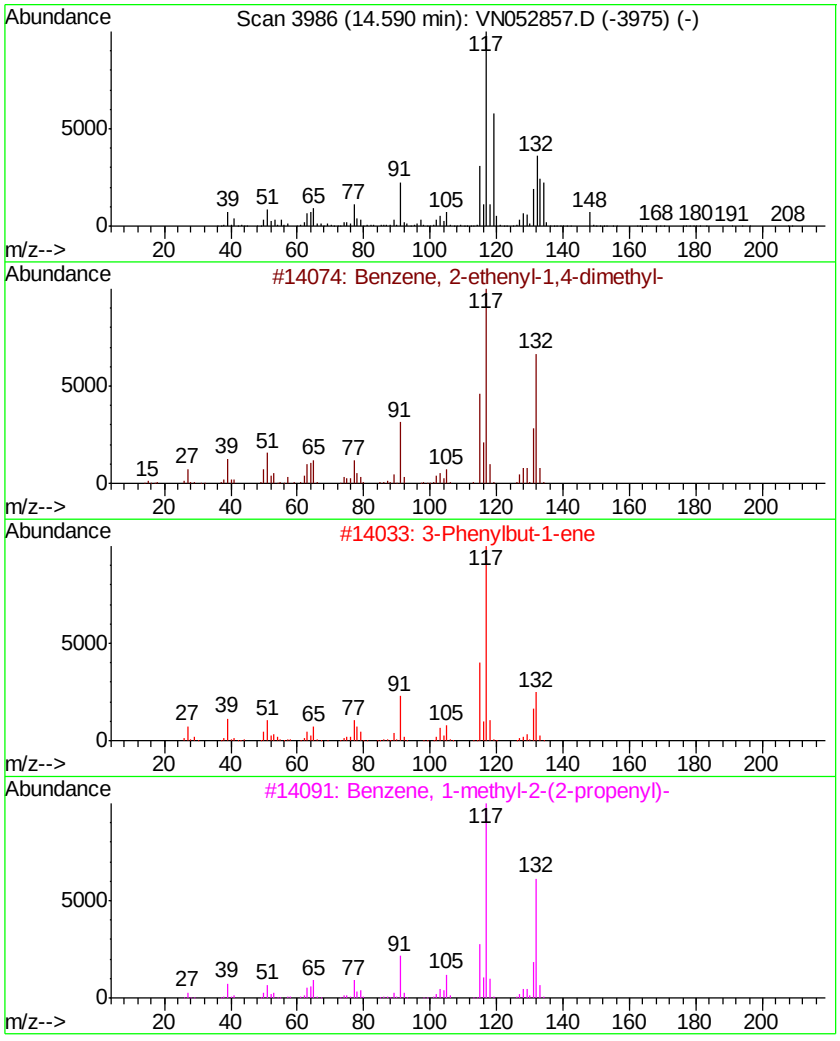
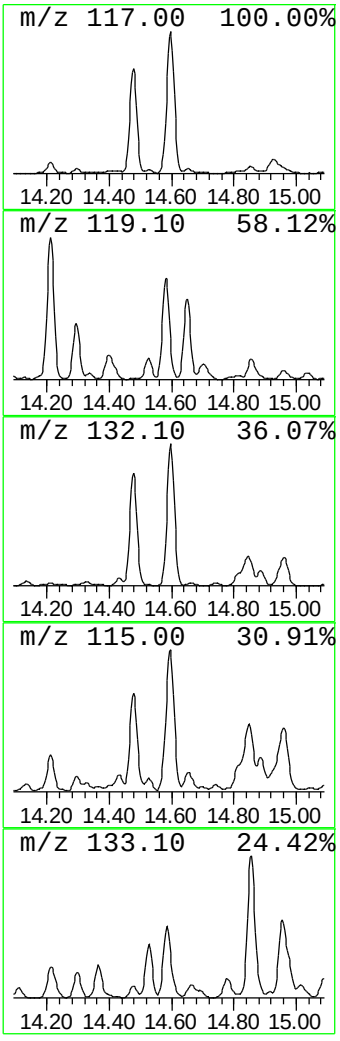
Peak Number 8 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.59	398.49 ug/l	39778600	1,4-Dichlorobenzene-d4	13.35

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	90
2	3-Phenylbut-1-ene	132	C10H12	000934-10-1	70
3	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	70
4	1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	62
5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	62





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Data File : VN052857.D
Acq On : 13 Dec 2018 12:01
Operator : MD\SY
Sample : J6319-01
Misc : 4.98g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 7 Sample Multiplier: 28

Instrument : MSVOA_N
ClientSampleId : 055_BT-1-1

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N121218W.M
Quant Title : SW846 8260

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

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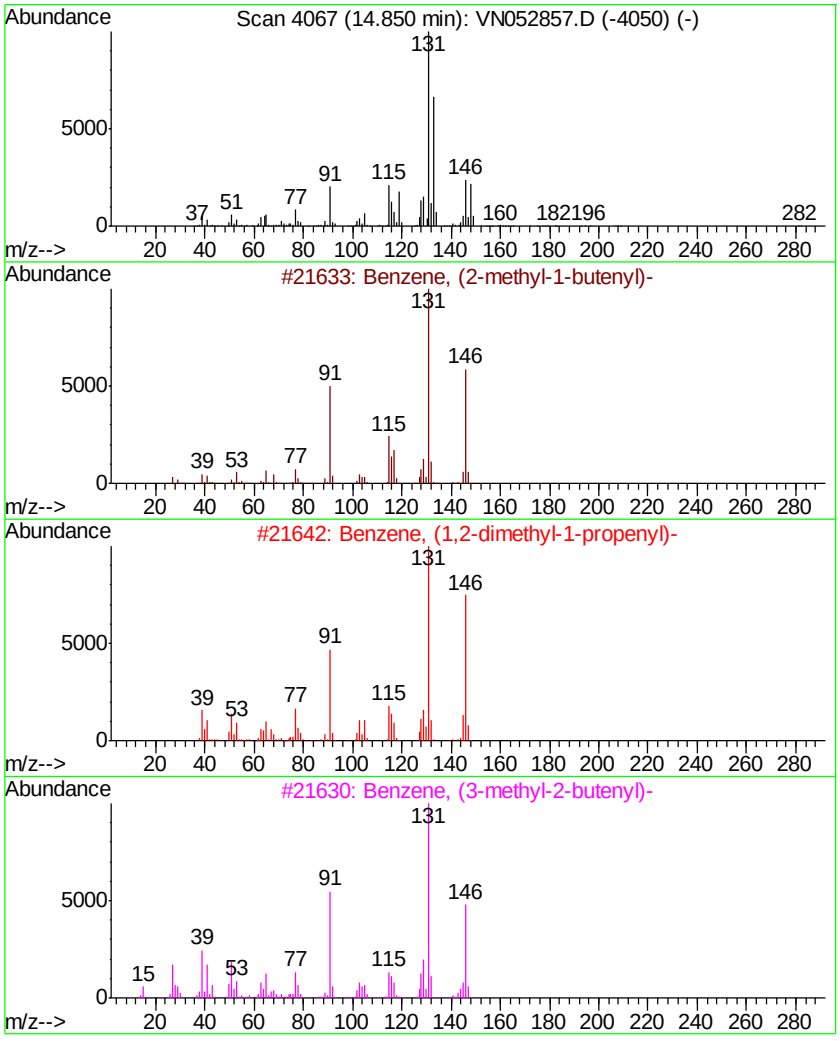
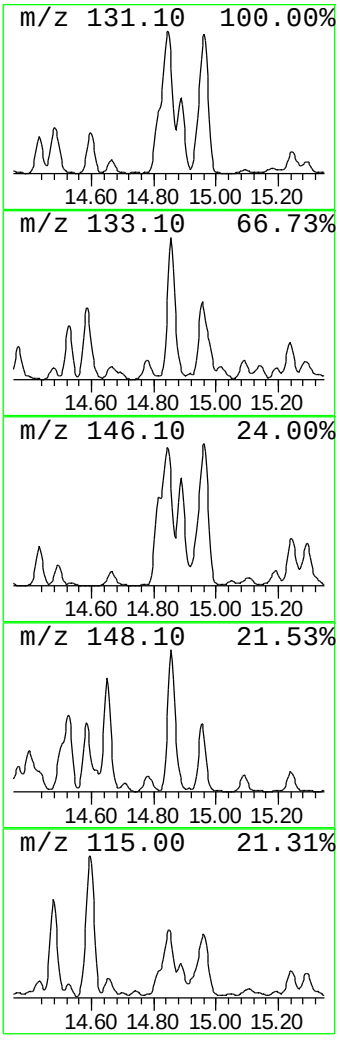
Peak Number 9 Benzene, (2-methyl-1-butenyl)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.85	261.93 ug/l	26147000	1,4-Dichlorobenzene-d4	13.35

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	86
2	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	86
3	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	78
4	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	70
5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	70





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN121318\
Data File : VN052857.D
Acq On : 13 Dec 2018 12:01
Operator : MD\SY
Sample : J6319-01
Misc : 4.98g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 7 Sample Multiplier: 28

Instrument : MSVOA_N
ClientSampleId : 055_BT-1-1

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N121218W.M
Quant Title : SW846 8260

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

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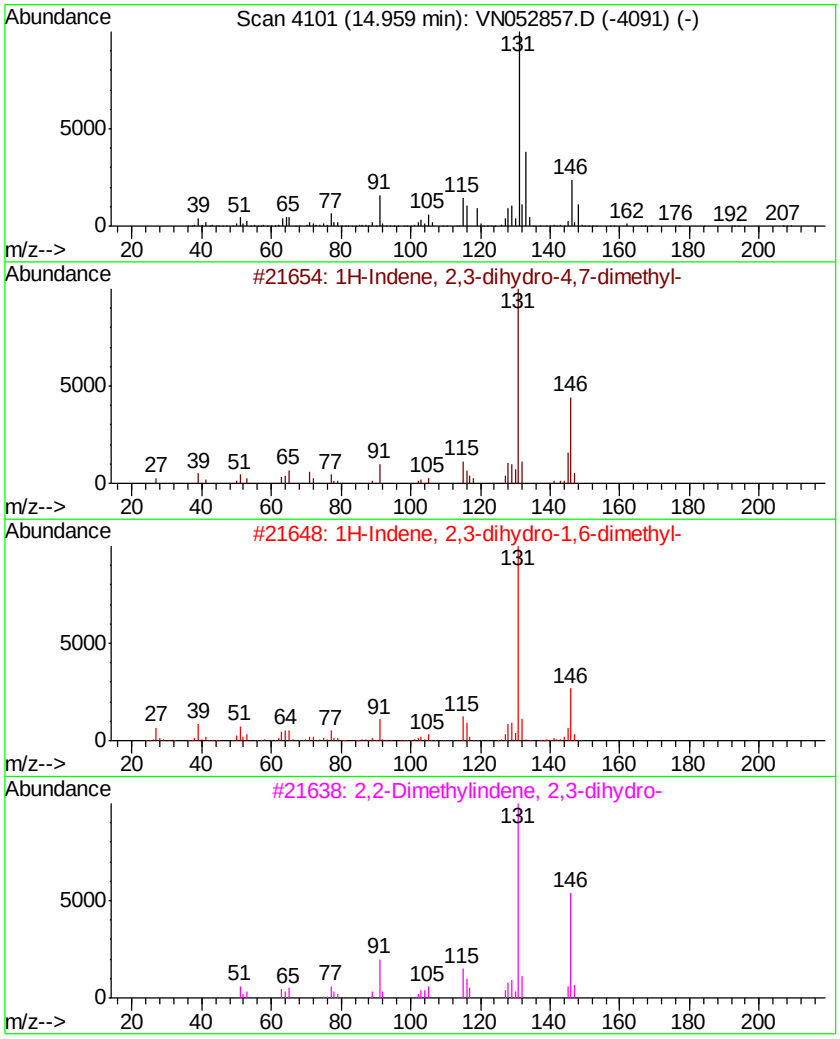
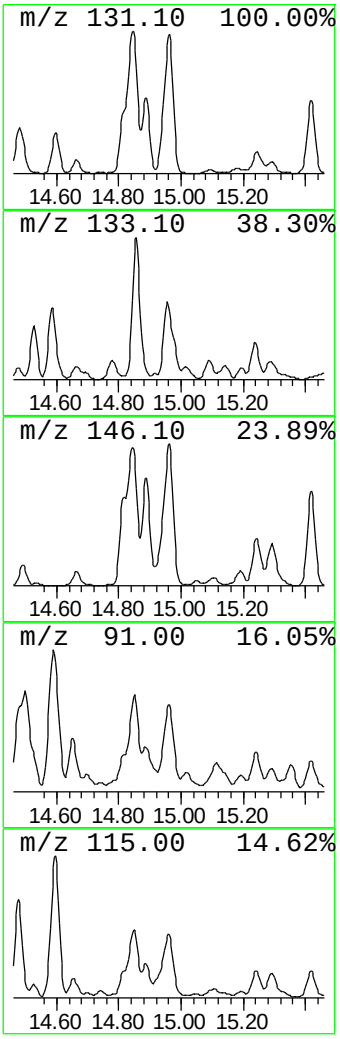
Peak Number 10 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.96	218.16 ug/l	21777400	1,4-Dichlorobenzene-d4	13.35

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	64
2	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	64
3	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	64
4	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	64
5	1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	58





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN121318\
Data File : VN052857.D
Acq On : 13 Dec 2018 12:01
Operator : MD\SY
Sample : J6319-01
Misc : 4.98g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 7 Sample Multiplier: 28

Instrument : MSVOA_N
ClientSampleId : 055_BT-1-1

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N121218W.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
Heptane, 2-methyl-	9.73	233.0	ug/l	29688000	2	8.58	6370480	50.0
Octane, 2-methyl-	11.20	364.8	ug/l	52046300	3	11.41	7133340	50.0
Octane, 3-methyl-	11.30	221.0	ug/l	31529200	3	11.41	7133340	50.0
Benzene, 1-ethyl-...	13.89	340.4	ug/l	33976500	4	13.35	4991200	50.0
Indan, 1-methyl-	13.99	278.4	ug/l	27791800	4	13.35	4991200	50.0
Benzene, 1,2,3,4-...	14.21	235.5	ug/l	23512200	4	13.35	4991200	50.0
1H-Indene, 2,3-di...	14.48	178.5	ug/l	17820300	4	13.35	4991200	50.0
Benzene, 2-etheny...	14.59	398.5	ug/l	39778600	4	13.35	4991200	50.0
Benzene, (2-methy...	14.85	261.9	ug/l	26147000	4	13.35	4991200	50.0
1H-Indene, 2,3-di...	14.96	218.2	ug/l	21777400	4	13.35	4991200	50.0