LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX092618\
Data File : VX004747.D
Acq On : 27 Sep 2018 02:24
Operator : JC/MD
Sample : J5029-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 30 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T11-MW-1-8.0-092018

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\82X092618W.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	1.636	67	79	83	rBV7	17076	52188	1.64%	0.191%
2	2.446	207	212	218	rBV	49674	79287	2.49%	0.290%
3	2.623	235	241	251	rVB	93562	162272	5.09%	0.593%
4	3.074	310	315	322	rVB	22864	51013	1.60%	0.186%
5	3.202	324	336	349	rVB	75049	200046	6.28%	0.731%
6	4.397	520	532	544	rVB4	22076	69471	2.18%	0.254%
7	5.507	702	714	721	rBV2	201081	578822	18.16%	2.114%
8	5.574	722	725	731	rVV2	36508	86761	2.72%	0.317%
9	5.659	731	739	763	rVB	267890	760959	23.88%	2.780%
10	6.068	796	806	826	rBV2	208260	547442	17.18%	2.000%
11	6.299	828	844	849	rBV6	10524	41735	1.31%	0.152%
12	6.860	924	936	945	rBV	384282	930740	29.20%	3.400%
13	7.458	1024	1034	1045	rBV3	43543	113089	3.55%	0.413%
14	8.214	1144	1158	1164	rBV	32091	71413	2.24%	0.261%
15	8.573	1211	1217	1227	rVB	34184	62909	1.97%	0.230%
16	8.714	1234	1240	1258	rBV	869804	1501806	47.12%	5.486%
17	9.220	1319	1323	1334	rVB	21878	38008	1.19%	0.139%
18	10.116	1460	1470	1480	rBV	791612	1170393	36.72%	4.275%
19	10.360	1500	1510	1516	rVB2	24724	49644	1.56%	0.181%
20	11.140	1632	1638	1653	rBV	828801	1090516	34.22%	3.984%
21	11.670	1717	1725	1733	rBV	31827	54802	1.72%	0.200%
22	11.768	1733	1741	1745	rBV	25845	42413	1.33%	0.155%
23	11.920	1761	1766	1769	rBV	56471	77290	2.43%	0.282%
24	12.079	1786	1792	1800	rBV	1035929	1331741	41.79%	4.865%
25	12.231	1813	1817	1821	rBV	62804	77803	2.44%	0.284%
26	12.304	1821	1829	1836	rVV	1334477	1747429	54.83%	6.383%
27	12.371	1836	1840	1848	rVB2	91587	140116	4.40%	0.512%
28	12.463	1848	1855	1861	rBV	176699	229782	7.21%	0.839%
29	12.670	1881	1889	1892	rBV	1310040	1545614	48.50%	5.646%
30	12.701	1892	1894	1899	rVV	363420	428451	13.44%	1.565%
31	12.756	1899	1903	1910	rVV2	1114961	1615091	50.68%	5.900%
32	12.816	1910	1913	1918	rVV	32015	41656	1.31%	0.152%
33	12.896	1920	1926	1932	rVB	111415	147924	4.64%	0.540%
34	12.981	1932	1940	1945	rBV	1027324	1284830	40.31%	4.693%

LSC Area Percent Report

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Acq On : 27 Sep 2018 02:24
Operator : JC/MD
Sample : J5029-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 30 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T11-MW-1-8.0-092018

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\82X092618W.M
Title : SW846 8260

35	13.085	1952	1957	1963	rBV2	86030	139967	4.39%	0.511%
36	13.152	1964	1968	1971	rVV	52783	66725	2.09%	0.244%
37	13.194	1971	1975	1977	rVV2	119293	156909	4.92%	0.573%
38	13.225	1977	1980	1990	rVB	655649	832260	26.11%	3.040%
39	13.347	1995	2000	2006	rVV2	2258314	3186991	100.00%	11.642%
40	13.408	2006	2010	2011	rVV2	62350	86492	2.71%	0.316%
41	13.426	2011	2013	2018	rVV	73748	91968	2.89%	0.336%
42	13.481	2018	2022	2027	rVB	68639	90616	2.84%	0.331%
43	13.566	2031	2036	2038	rVV	72798	96045	3.01%	0.351%
44	13.603	2038	2042	2045	rVV	206220	293208	9.20%	1.071%
45	13.640	2045	2048	2052	rVV2	258717	376941	11.83%	1.377%
46	13.700	2052	2058	2059	rVV2	179329	276571	8.68%	1.010%
47	13.725	2059	2062	2071	rVB3	291056	491879	15.43%	1.797%
48	13.810	2071	2076	2079	rBV	40751	56981	1.79%	0.208%
49	13.853	2079	2083	2089	rVV	71353	91133	2.86%	0.333%
50	13.914	2089	2093	2097	rVB	114228	136416	4.28%	0.498%
51	13.987	2097	2105	2109	rBV2	215507	303757	9.53%	1.110%
52	14.030	2109	2112	2116	rVB	101293	118216	3.71%	0.432%
53	14.133	2124	2129	2132	rBV	159081	205219	6.44%	0.750%
54	14.164	2132	2134	2138	rVB	57110	58142	1.82%	0.212%
55	14.274	2147	2152	2162	rBV3	242255	488981	15.34%	1.786%
56	14.377	2162	2169	2174	rBV	104644	145499	4.57%	0.532%
57	14.432	2174	2178	2183	rVV3	182228	241203	7.57%	0.881%
58	14.542	2190	2196	2201	rBV2	257438	342228	10.74%	1.250%
59	14.664	2210	2216	2220	rBV3	38833	84326	2.65%	0.308%
60	14.731	2223	2227	2232	rVV2	44687	66927	2.10%	0.244%
61	14.853	2241	2247	2254	rBV3	71474	146423	4.59%	0.535%
62	14.969	2259	2266	2273	rBV8	12384	32481	1.02%	0.119%
63	15.249	2306	2312	2320	rBV	42380	73116	2.29%	0.267%
64	15.444	2338	2344	2347	rBV	68060	104139	3.27%	0.380%
65	15.481	2347	2350	2355	rVB	73728	97499	3.06%	0.356%
66	15.548	2355	2361	2367	rBV	244989	401662	12.60%	1.467%
67	15.688	2375	2384	2397	rVB	529068	855775	26.85%	3.126%
68	15.810	2398	2404	2410	rBV3	23297	42541	1.33%	0.155%
69	15.920	2412	2422	2434	rVB	149455	395078	12.40%	1.443%
70	16.072	2441	2447	2459	rVB	90911	158778	4.98%	0.580%
71	16.419	2497	2504	2512	rVB2	58421	118220	3.71%	0.432%

LSC Area Percent Report

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Acq On : 27 Sep 2018 02:24
Operator : JC/MD
Sample : J5029-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 30 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T11-MW-1-8.0-092018

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\82X092618W.M
Title : SW846 8260

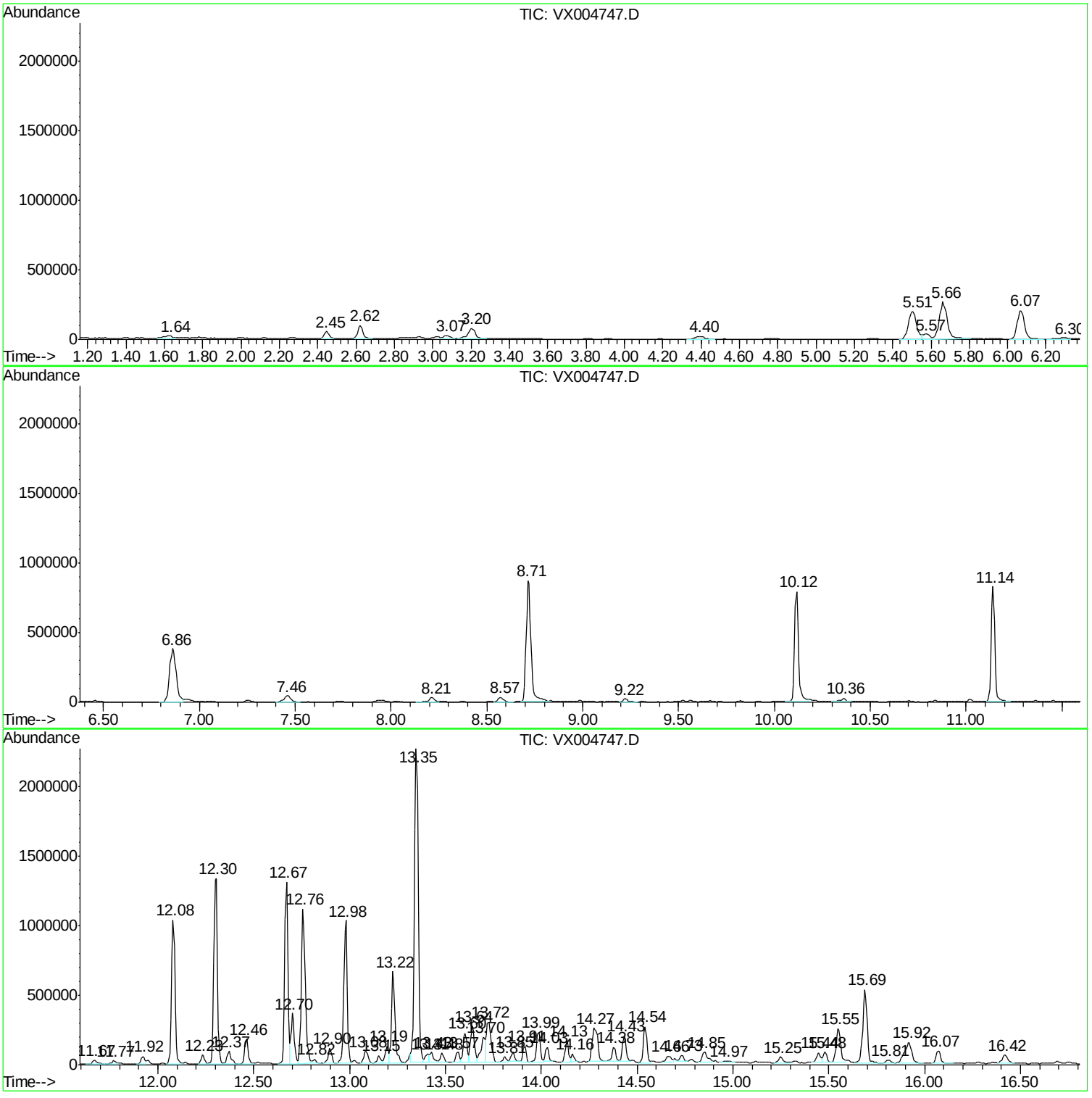
Sum of corrected areas: 27374768

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX092618\
Data File : VX004747.D
Acq On : 27 Sep 2018 02:24
Operator : JC/MD
Sample : J5029-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 30 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
T11-MW-1-8.0-092018

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

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



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX092618\
Data File : VX004747.D
Acq On : 27 Sep 2018 02:24
Operator : JC/MD
Sample : J5029-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 30 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T11-MW-1-8.0-092018

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

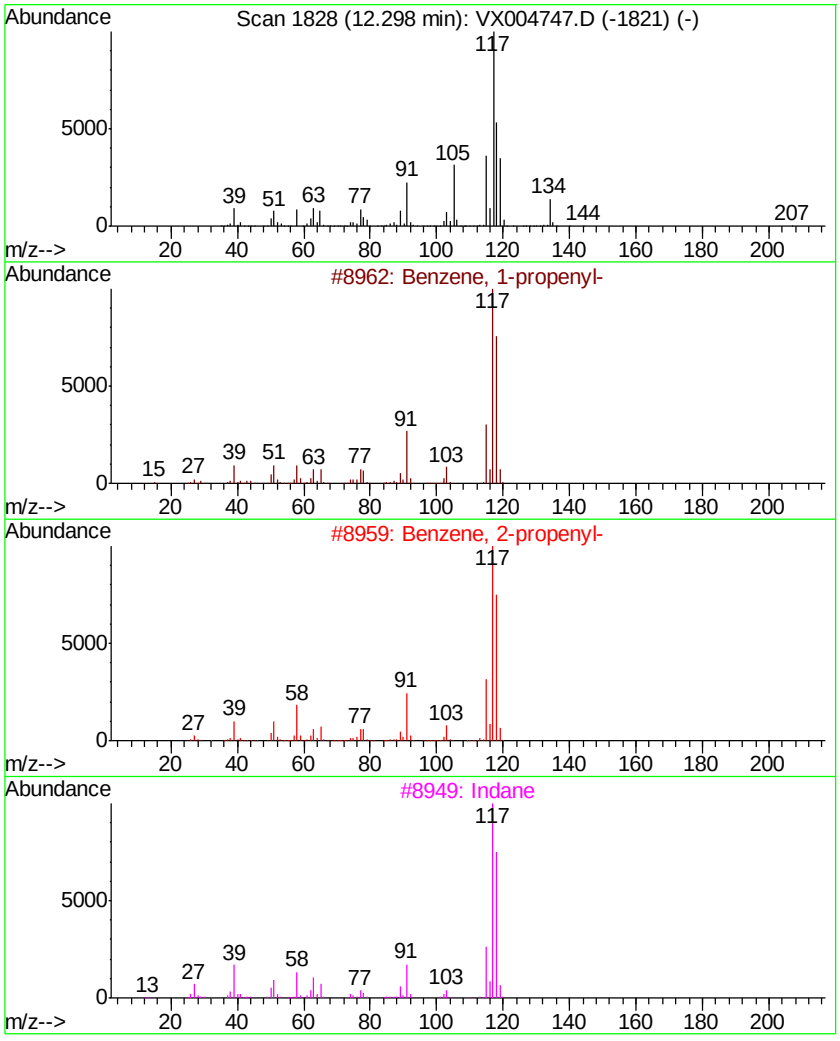
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

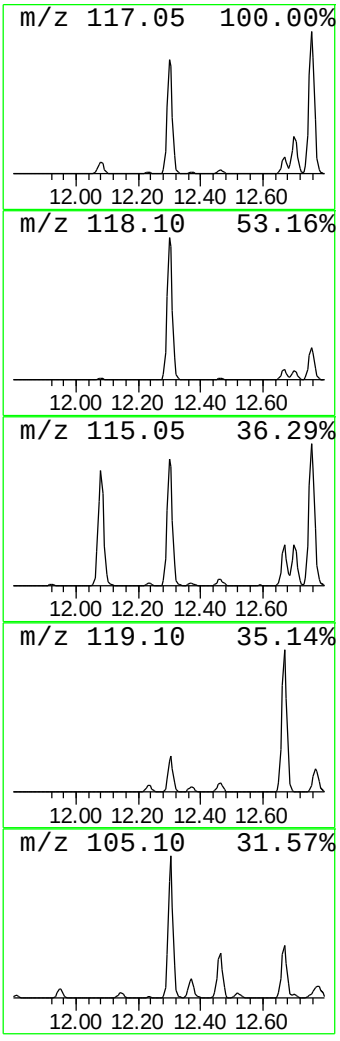
****

Peak Number 1 Benzene, 1-propenyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	65.61 ug/l	1747430	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-propenyl-	118	C9H10	000637-50-3	90
2	Benzene, 2-propenyl-	118	C9H10	000300-57-2	55
3	Indane	118	C9H10	000496-11-7	55
4	(Z)-1-Phenylpropene	118	C9H10	000766-90-5	55
5	Bicyclo[4.2.0]octa-1,3,5-triene,...	118	C9H10	055337-80-9	55





Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX092618\
Data File : VX004747.D
Acq On : 27 Sep 2018 02:24
Operator : JC/MD
Sample : J5029-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 30 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T11-MW-1-8.0-092018

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

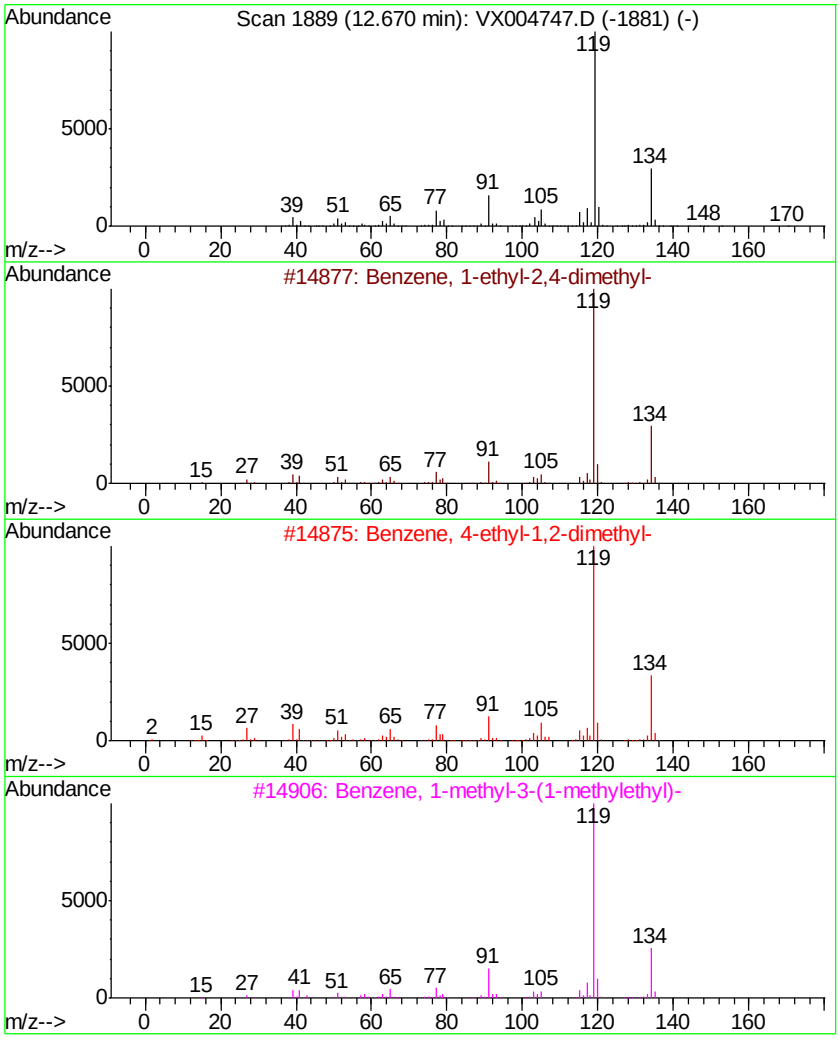
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

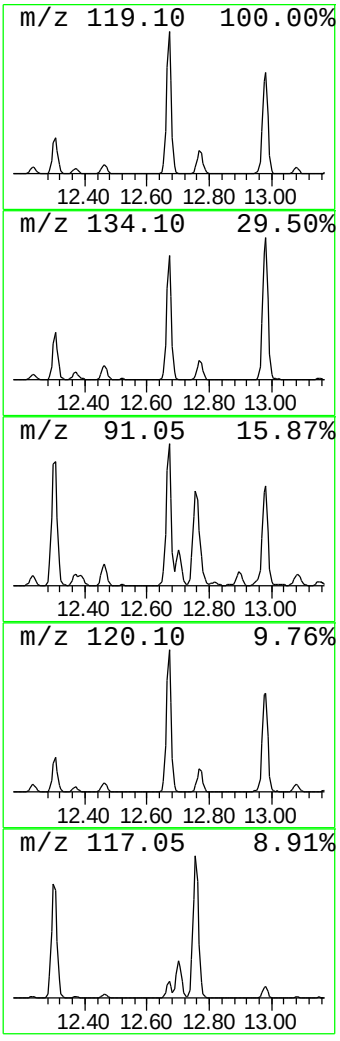
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.67	58.03 ug/l	1545610	1,4-Dichlorobenzene-d4	12.08

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, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
4	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
5	o-Cymene	134	C10H14	000527-84-4	95





Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX092618\
Data File : VX004747.D
Acq On : 27 Sep 2018 02:24
Operator : JC/MD
Sample : J5029-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 30 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T11-MW-1-8.0-092018

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

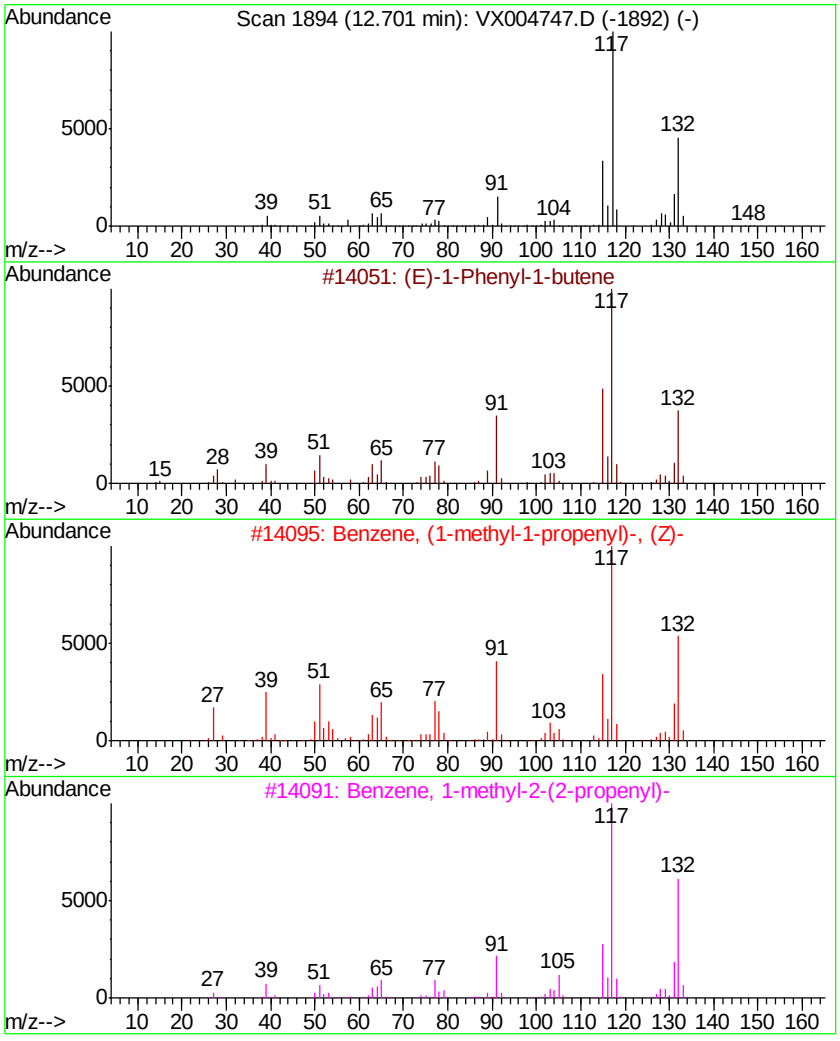
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

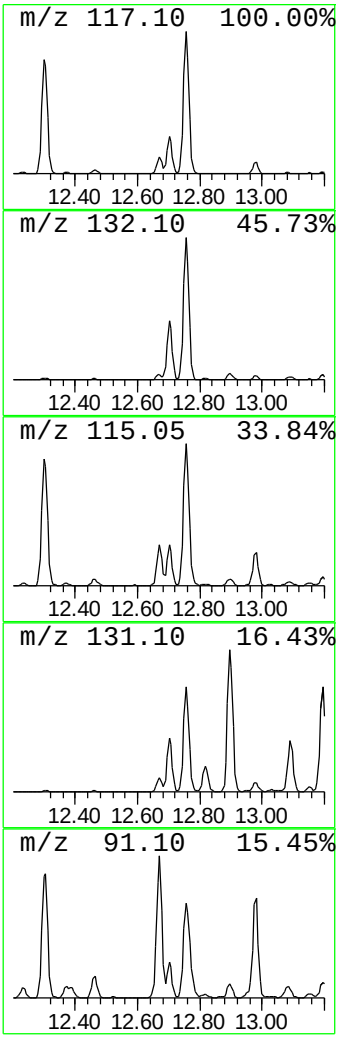
****

Peak Number 3 (E)-1-Phenyl-1-butene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.70	16.09 ug/l	428451	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(E)-1-Phenyl-1-butene	132	C10H12	001005-64-7	93
2	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	90
3	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	90
4	1-Phenyl-1-butene	132	C10H12	000824-90-8	90
5	Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	90





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX092618\
Data File : VX004747.D
Acq On : 27 Sep 2018 02:24
Operator : JC/MD
Sample : J5029-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 30 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T11-MW-1-8.0-092018

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

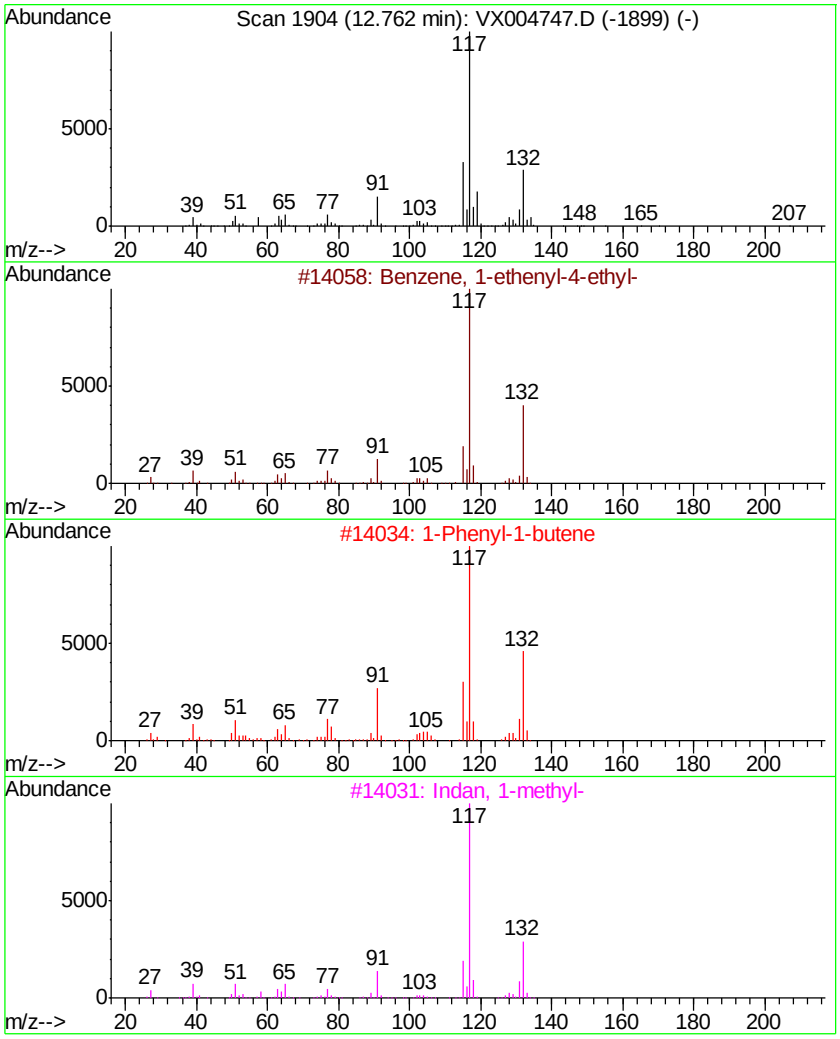
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

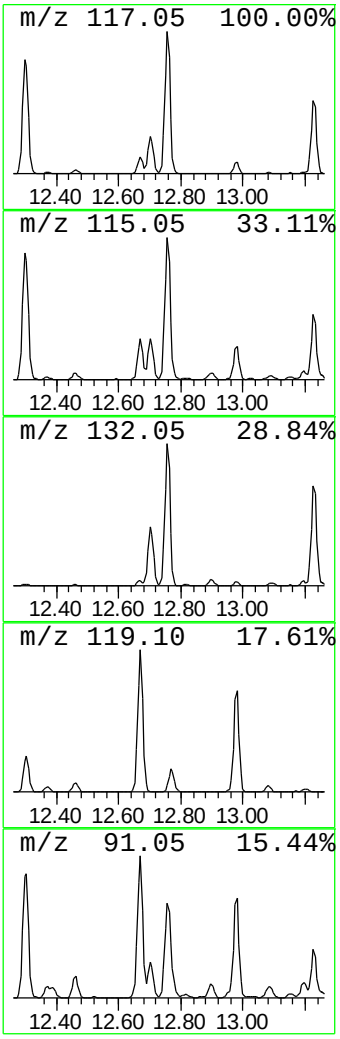
****

Peak Number 4 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	60.64 ug/l	1615090	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87
2	1-Phenyl-1-butene	132	C10H12	000824-90-8	87
3	Indan, 1-methyl-	132	C10H12	000767-58-8	87
4	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
5	Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	87





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX092618\
Data File : VX004747.D
Acq On : 27 Sep 2018 02:24
Operator : JC/MD
Sample : J5029-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 30 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T11-MW-1-8.0-092018

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

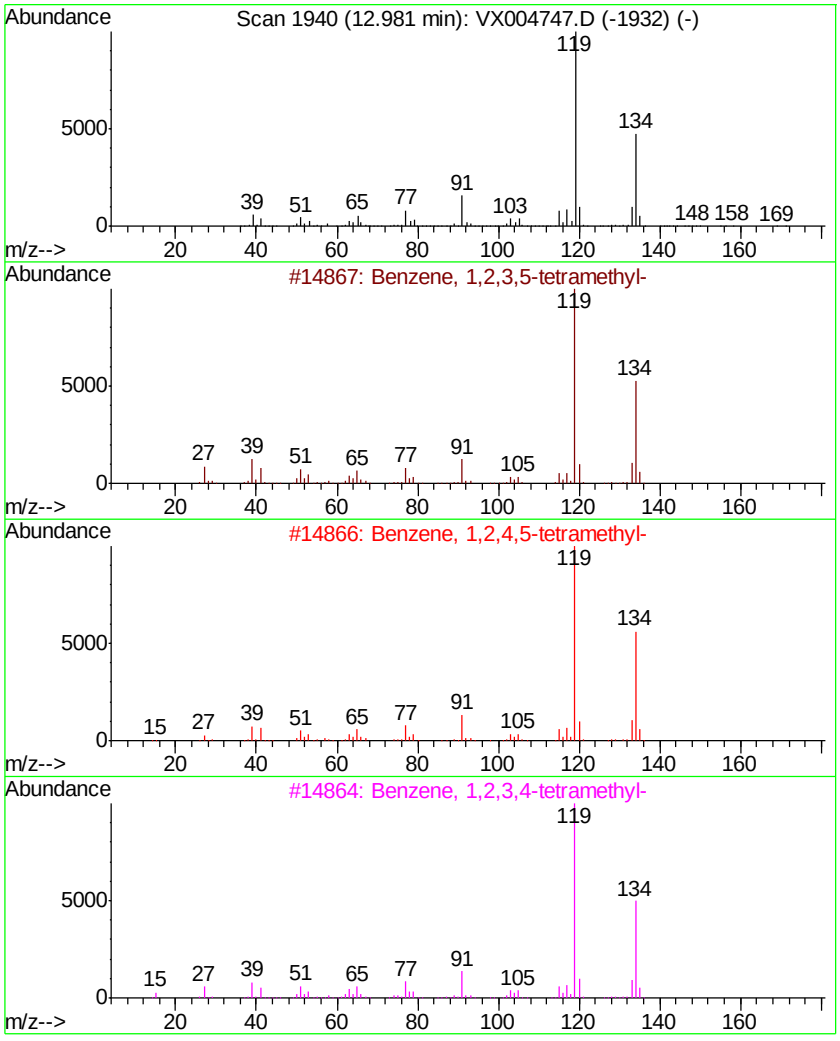
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

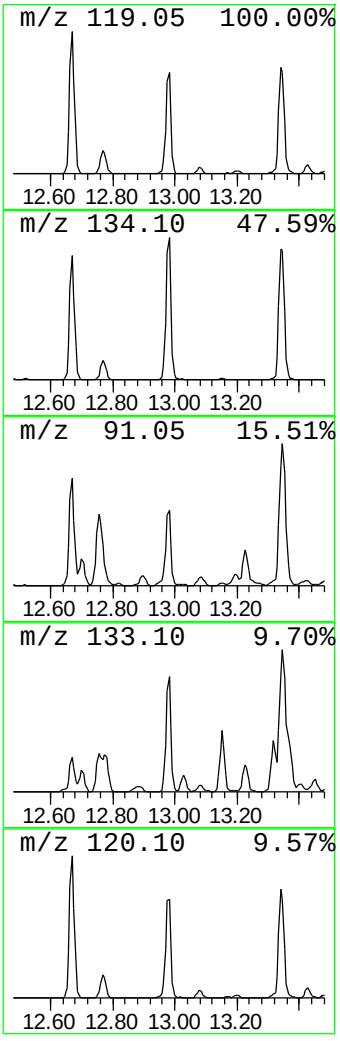
****

Peak Number 5 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.98	48.24 ug/l	1284830	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
3	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
4	o-Cymene	134	C10H14	000527-84-4	95
5	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX092618\
Data File : VX004747.D
Acq On : 27 Sep 2018 02:24
Operator : JC/MD
Sample : J5029-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 30 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T11-MW-1-8.0-092018

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

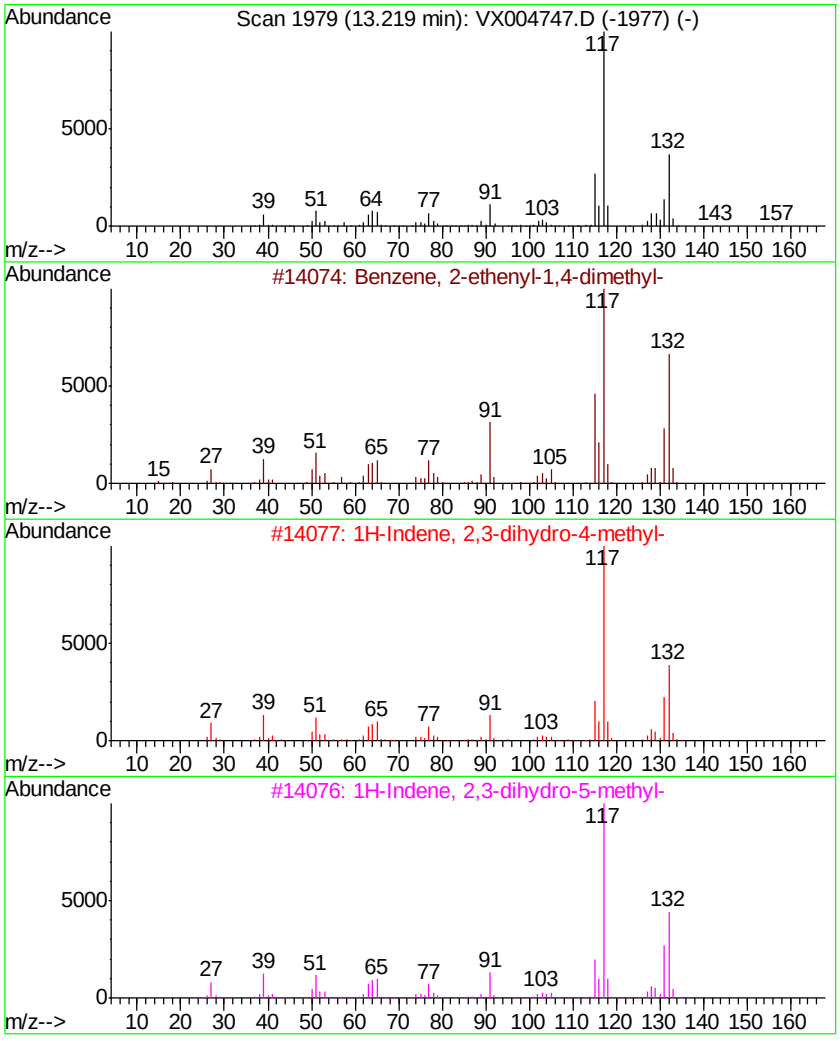
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

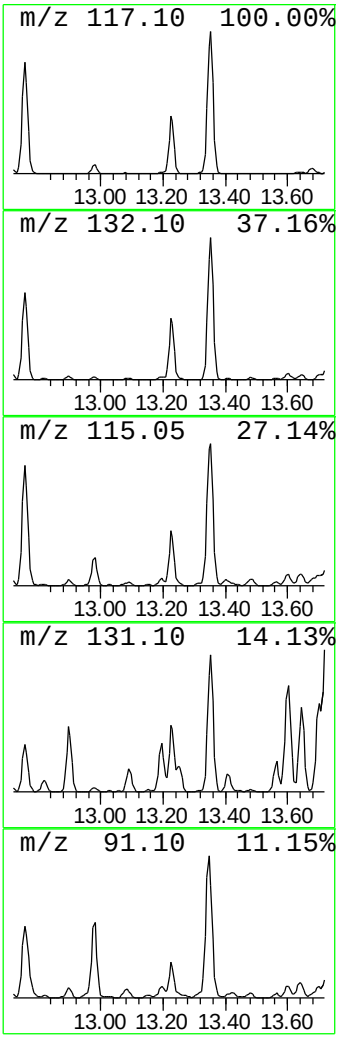
****

Peak Number 6 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.22	31.25 ug/l	832260	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	92
2	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90
3	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	90
4	Indan, 1-methyl-	132	C10H12	000767-58-8	90
5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX092618\
Data File : VX004747.D
Acq On : 27 Sep 2018 02:24
Operator : JC/MD
Sample : J5029-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 30 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T11-MW-1-8.0-092018

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

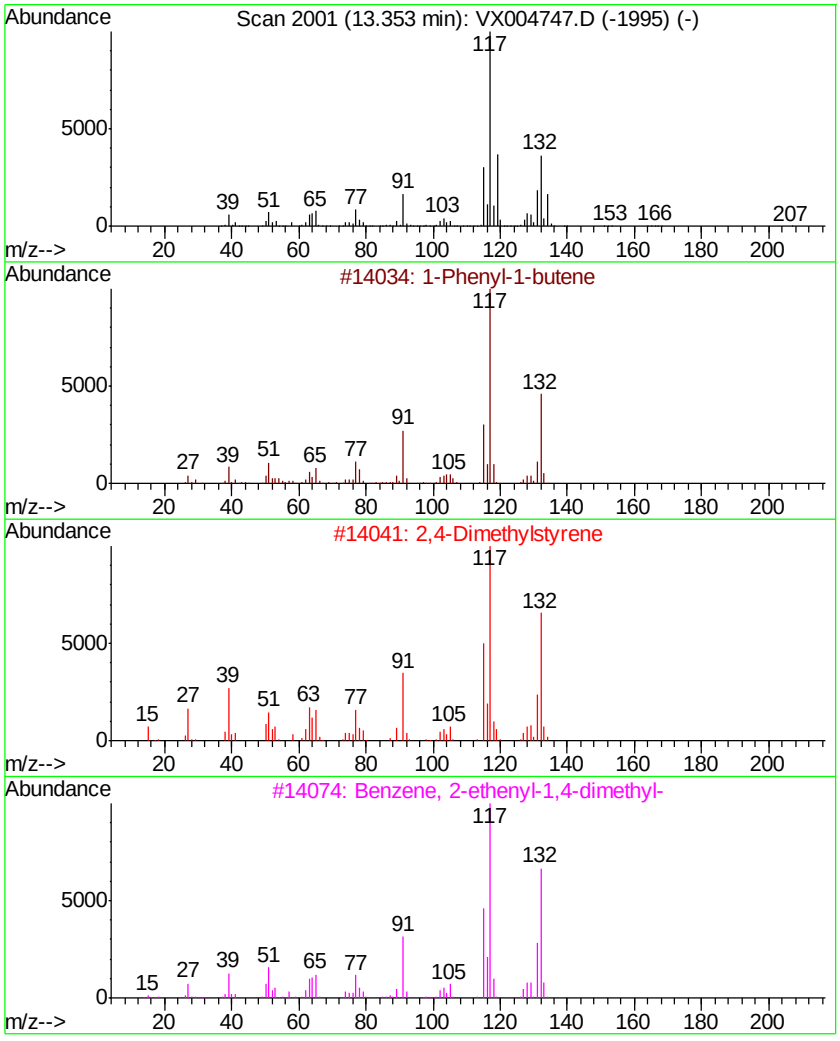
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

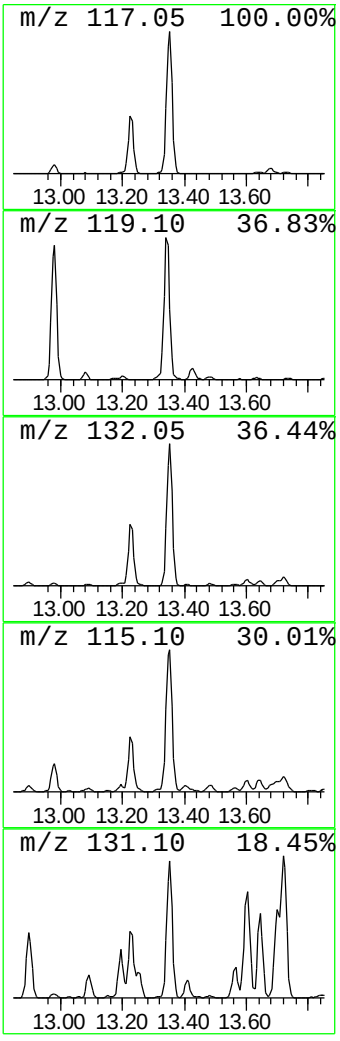
****

Peak Number 7 1-Phenyl-1-butene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	119.66 ug/l	3186990	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Phenyl-1-butene	132	C10H12	000824-90-8	95
2	2,4-Dimethylstyrene	132	C10H12	002234-20-0	94
3	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	94
4	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
5	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	87





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX092618\
Data File : VX004747.D
Acq On : 27 Sep 2018 02:24
Operator : JC/MD
Sample : J5029-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 30 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T11-MW-1-8.0-092018

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

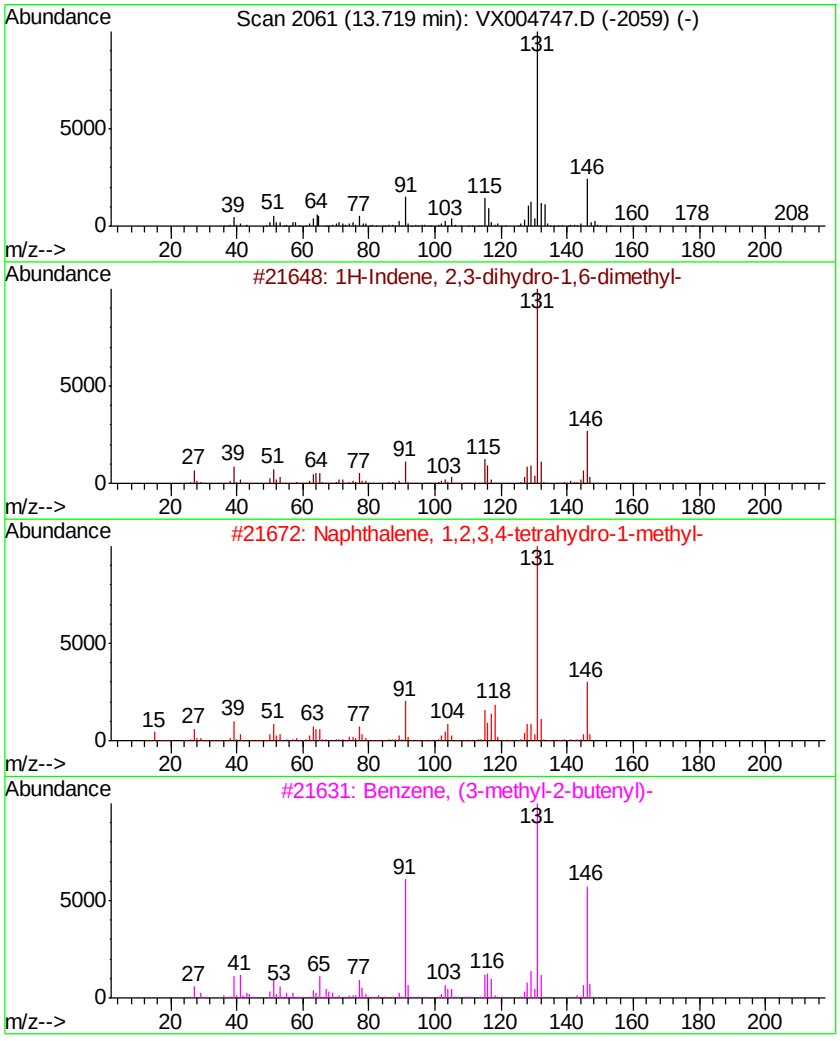
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

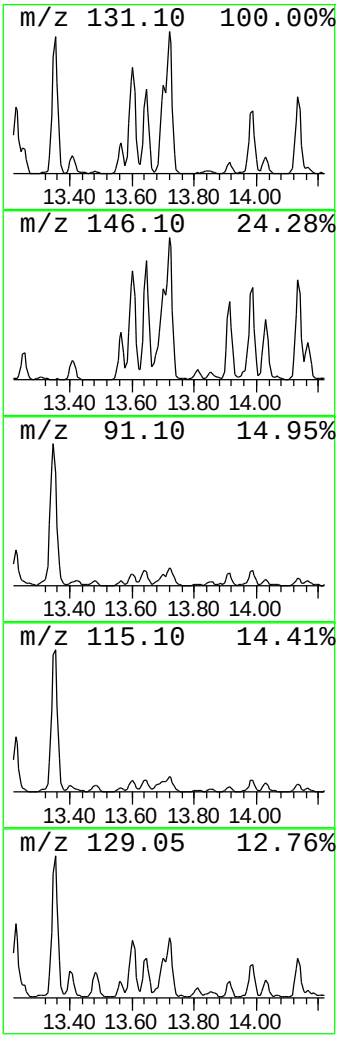
****

Peak Number 8 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.72	18.47 ug/l	491879	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	86
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	70
3	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	70
4	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	70
5	Benzene, 1-methyl-2-(1-methyl-2-...	146	C11H14	097664-19-2	70





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Acq On : 27 Sep 2018 02:24
Operator : JC/MD
Sample : J5029-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 30 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T11-MW-1-8.0-092018

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

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

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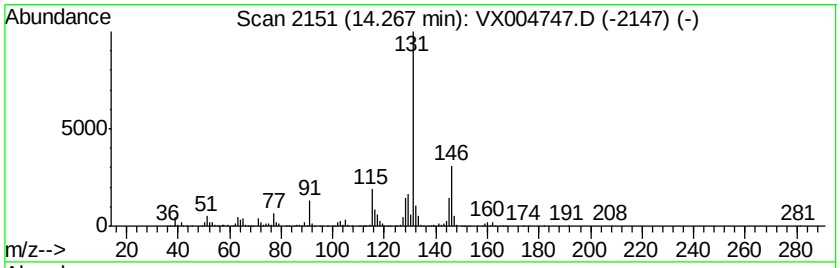
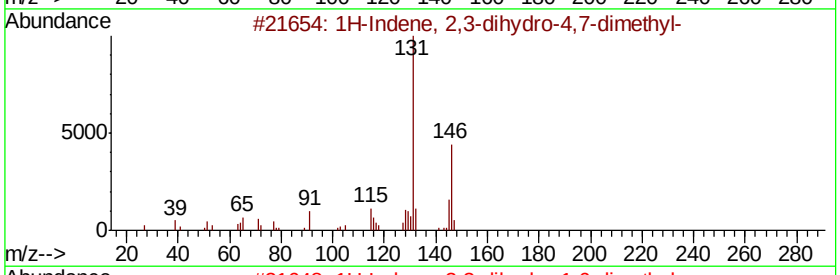
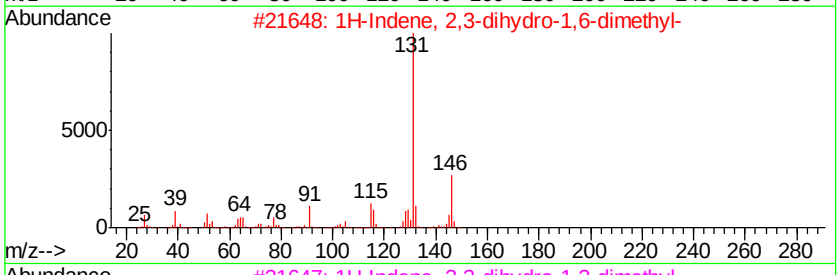
Peak Number 9 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 9

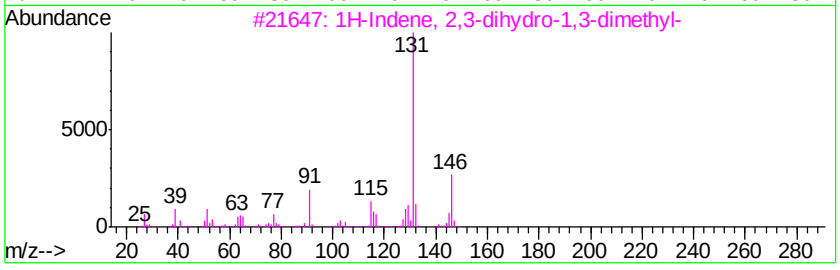
R.T.	EstConc	Area	Relative to ISTD	R.T.
14.27	18.36 ug/l	488981	1,4-Dichlorobenzene-d4	12.08

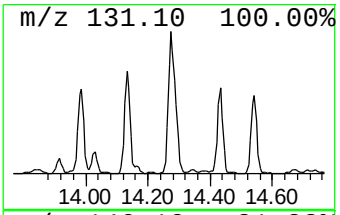
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	95
2	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	94
3	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	91
4	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	91
5	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	91

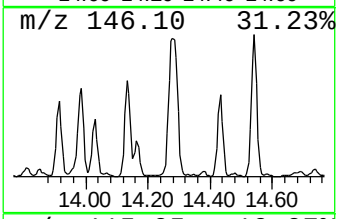


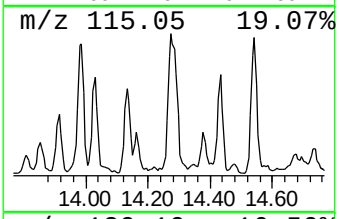


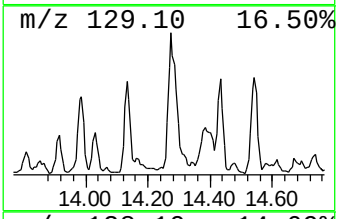


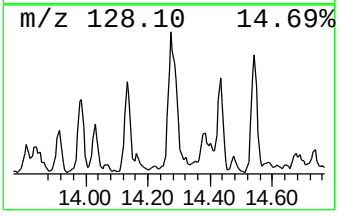












Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX092618\
Data File : VX004747.D
Acq On : 27 Sep 2018 02:24
Operator : JC/MD
Sample : J5029-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 30 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T11-MW-1-8.0-092018

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

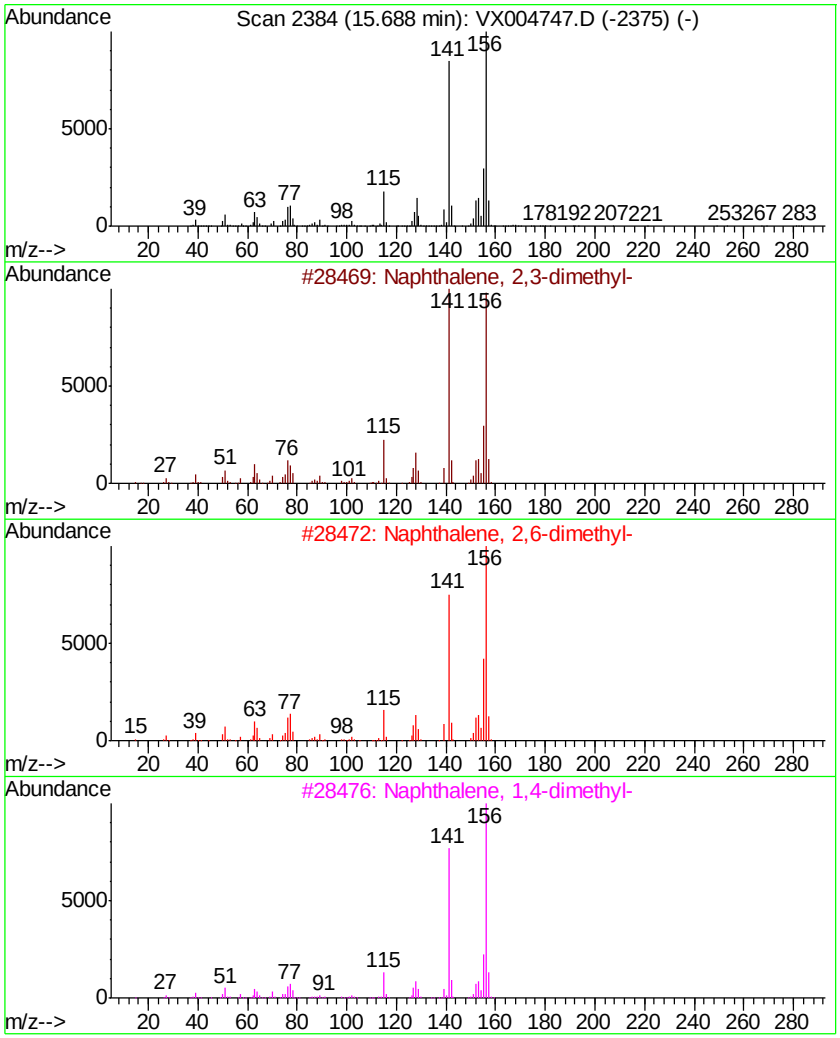
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

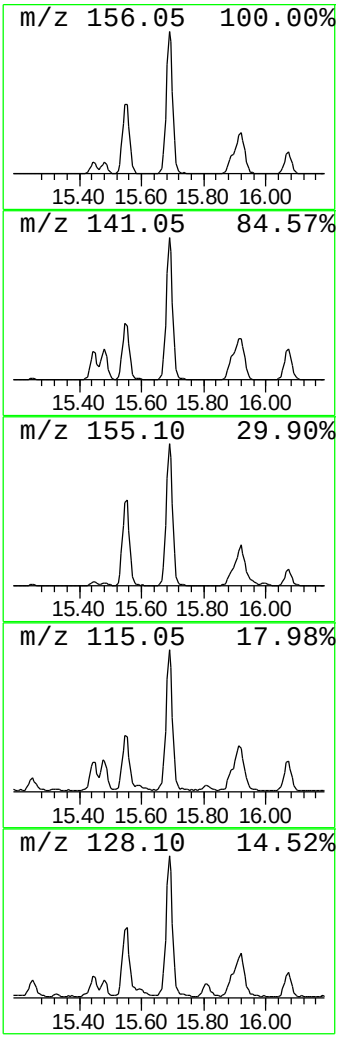
****

Peak Number 10 Naphthalene, 2,3-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.69	32.13 ug/l	855775	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
2	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
3	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97
4	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
5	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX092618\
Data File : VX004747.D
Acq On : 27 Sep 2018 02:24
Operator : JC/MD
Sample : J5029-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 30 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T11-MW-1-8.0-092018

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.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-propenyl-	12.30	65.6	ug/l	1747430	4	12.08	1331740	50.0
Benzene, 1-ethyl-...	12.67	58.0	ug/l	1545610	4	12.08	1331740	50.0
(E)-1-Phenyl-1-bu...	12.70	16.1	ug/l	428451	4	12.08	1331740	50.0
Benzene, 1-etheny...	12.76	60.6	ug/l	1615090	4	12.08	1331740	50.0
Benzene, 1,2,3,5-...	12.98	48.2	ug/l	1284830	4	12.08	1331740	50.0
Benzene, 2-etheny...	13.22	31.3	ug/l	832260	4	12.08	1331740	50.0
1-Phenyl-1-butene	13.35	119.7	ug/l	3186990	4	12.08	1331740	50.0
1H-Indene, 2,3-di...	13.72	18.5	ug/l	491879	4	12.08	1331740	50.0
1H-Indene, 2,3-di...	14.27	18.4	ug/l	488981	4	12.08	1331740	50.0
Naphthalene, 2,3-...	15.69	32.1	ug/l	855775	4	12.08	1331740	50.0