LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX091418\
Data File : VX004536.D
Acq On : 14 Sep 2018 13:34
Operator : JC/MD
Sample : J4724-18 50X
Misc : 3.19g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : EP1-SUT-4USTS-SB

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\82X091118W.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	5.501	700	713	728	rBV2	185876	508621	5.93%	0.822%
2	5.665	728	740	757	rVB	257351	691584	8.07%	1.118%
3	6.068	795	806	825	rBV	186328	474151	5.53%	0.767%
4	6.860	926	936	964	rBV	344885	807133	9.41%	1.305%
5	8.713	1234	1240	1248	rVB	804066	1268050	14.79%	2.050%
6	9.445	1351	1360	1369	rBV2	63524	105307	1.23%	0.170%
7	9.561	1373	1379	1385	rBV5	49578	116192	1.35%	0.188%
8	9.622	1385	1389	1394	rVB	101406	141656	1.65%	0.229%
9	9.677	1394	1398	1403	rBV	154650	243403	2.84%	0.394%
10	9.890	1426	1433	1437	rBV3	111273	259362	3.02%	0.419%
11	9.957	1440	1444	1449	rVB	103539	138901	1.62%	0.225%
12	10.067	1456	1462	1465	rBV	201610	285547	3.33%	0.462%
13	10.116	1465	1470	1475	rVV	753107	1024789	11.95%	1.657%
14	10.183	1475	1481	1486	rVV3	39180	94591	1.10%	0.153%
15	10.250	1486	1492	1497	rVV2	153214	280890	3.28%	0.454%
16	10.366	1497	1511	1516	rVV3	352715	1056540	12.32%	1.708%
17	10.408	1516	1518	1530	rVB	155147	273264	3.19%	0.442%
18	10.512	1530	1535	1541	rBV	93363	140096	1.63%	0.227%
19	10.640	1549	1556	1564	rBV4	327273	644088	7.51%	1.041%
20	10.725	1564	1570	1573	rVV4	148672	247132	2.88%	0.400%
21	10.768	1574	1577	1580	rVB2	69571	89180	1.04%	0.144%
22	10.835	1580	1588	1592	rBV2	811757	1226466	14.30%	1.983%
23	10.914	1592	1601	1604	rVV2	711120	1204813	14.05%	1.948%
24	10.951	1604	1607	1612	rVV	316082	439148	5.12%	0.710%
25	11.018	1612	1618	1621	rVV	344801	506785	5.91%	0.819%
26	11.061	1621	1625	1626	rVV2	154874	212021	2.47%	0.343%
27	11.085	1626	1629	1633	rVV2	224669	339331	3.96%	0.549%
28	11.140	1633	1638	1642	rVB	1163429	1478139	17.24%	2.390%
29	11.183	1642	1645	1648	rBV	122102	137923	1.61%	0.223%
30	11.274	1653	1660	1669	rVB3	604115	1027996	11.99%	1.662%
31	11.359	1669	1674	1677	rBV	802877	947062	11.04%	1.531%
32	11.396	1677	1680	1682	rVV2	176605	262455	3.06%	0.424%
33	11.439	1682	1687	1693	rVV	1957507	3918089	45.69%	6.335%
34	11.506	1693	1698	1707	rVV	2147336	3540160	41.28%	5.724%

LSC Area Percent Report

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Sample : J4724-18 50X
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ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : EP1-SUT-4USTS-SB

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\82X091118W.M
Title : SW846 8260

35	11.579	1707	1710	1715	rVB4	98015	132228	1.54%	0.214%
36	11.676	1720	1726	1732	rVB5	249210	615328	7.18%	0.995%
37	11.737	1732	1736	1738	rBV2	153798	196881	2.30%	0.318%
38	11.768	1738	1741	1744	rVV	663834	803598	9.37%	1.299%
39	11.811	1744	1748	1757	rVB	6916424	8575126	100.00%	13.864%
40	11.890	1757	1761	1763	rBV2	109972	155717	1.82%	0.252%
41	11.920	1763	1766	1767	rVV	302569	328139	3.83%	0.531%
42	11.945	1767	1770	1775	rVV	708375	966828	11.27%	1.563%
43	11.999	1775	1779	1781	rVV	536787	691638	8.07%	1.118%
44	12.024	1781	1783	1787	rVV	909746	1092061	12.74%	1.766%
45	12.079	1787	1792	1798	rVB3	1055190	1771572	20.66%	2.864%
46	12.146	1798	1803	1809	rBV	2673143	3345612	39.02%	5.409%
47	12.231	1812	1817	1824	rVB2	167789	406171	4.74%	0.657%
48	12.304	1824	1829	1830	rBV2	855952	1076008	12.55%	1.740%
49	12.323	1830	1832	1836	rVB	1554551	1649568	19.24%	2.667%
50	12.371	1836	1840	1847	rVB3	2079913	3454788	40.29%	5.586%
51	12.463	1850	1855	1862	rBV3	260369	458845	5.35%	0.742%
52	12.524	1862	1865	1870	rVB4	198654	317380	3.70%	0.513%
53	12.585	1870	1875	1877	rBV	810437	1040346	12.13%	1.682%
54	12.609	1877	1879	1884	rVV	925766	1040051	12.13%	1.682%
55	12.670	1885	1889	1892	rVV	1024537	1112223	12.97%	1.798%
56	12.701	1892	1894	1898	rVB	228351	233245	2.72%	0.377%
57	12.774	1898	1906	1911	rBV4	624549	1545921	18.03%	2.499%
58	12.853	1914	1919	1921	rBV2	225577	269021	3.14%	0.435%
59	12.902	1921	1927	1931	rVB2	469689	746307	8.70%	1.207%
60	12.951	1931	1935	1937	rBV3	172096	281628	3.28%	0.455%
61	12.975	1937	1939	1943	rVB	430261	493685	5.76%	0.798%
62	13.018	1943	1946	1953	rVB	791977	1034068	12.06%	1.672%
63	13.085	1953	1957	1962	rBV3	207056	393254	4.59%	0.636%
64	13.152	1966	1968	1972	rVB	98551	99021	1.15%	0.160%
65	13.201	1972	1976	1977	rBV2	157875	200615	2.34%	0.324%
66	13.268	1983	1987	1990	rVB3	111588	131302	1.53%	0.212%
67	13.310	1990	1994	1996	rVV2	164861	218760	2.55%	0.354%
68	13.347	1996	2000	2006	rVV2	799698	1160083	13.53%	1.876%
69	13.426	2010	2013	2016	rVV	83382	85937	1.00%	0.139%
70	13.481	2017	2022	2031	rVB	652024	864127	10.08%	1.397%
71	13.639	2044	2048	2053	rBV2	108728	187238	2.18%	0.303%

LSC Area Percent Report

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Misc : 3.19g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : EP1-SUT-4USTS-SB

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\82X091118W.M
Title : SW846 8260

72	13.725	2059	2062	2067	rVB2	75876	115549	1.35%	0.187%
73	13.835	2072	2080	2087	rVB	306462	429844	5.01%	0.695%

Sum of corrected areas: 61850578

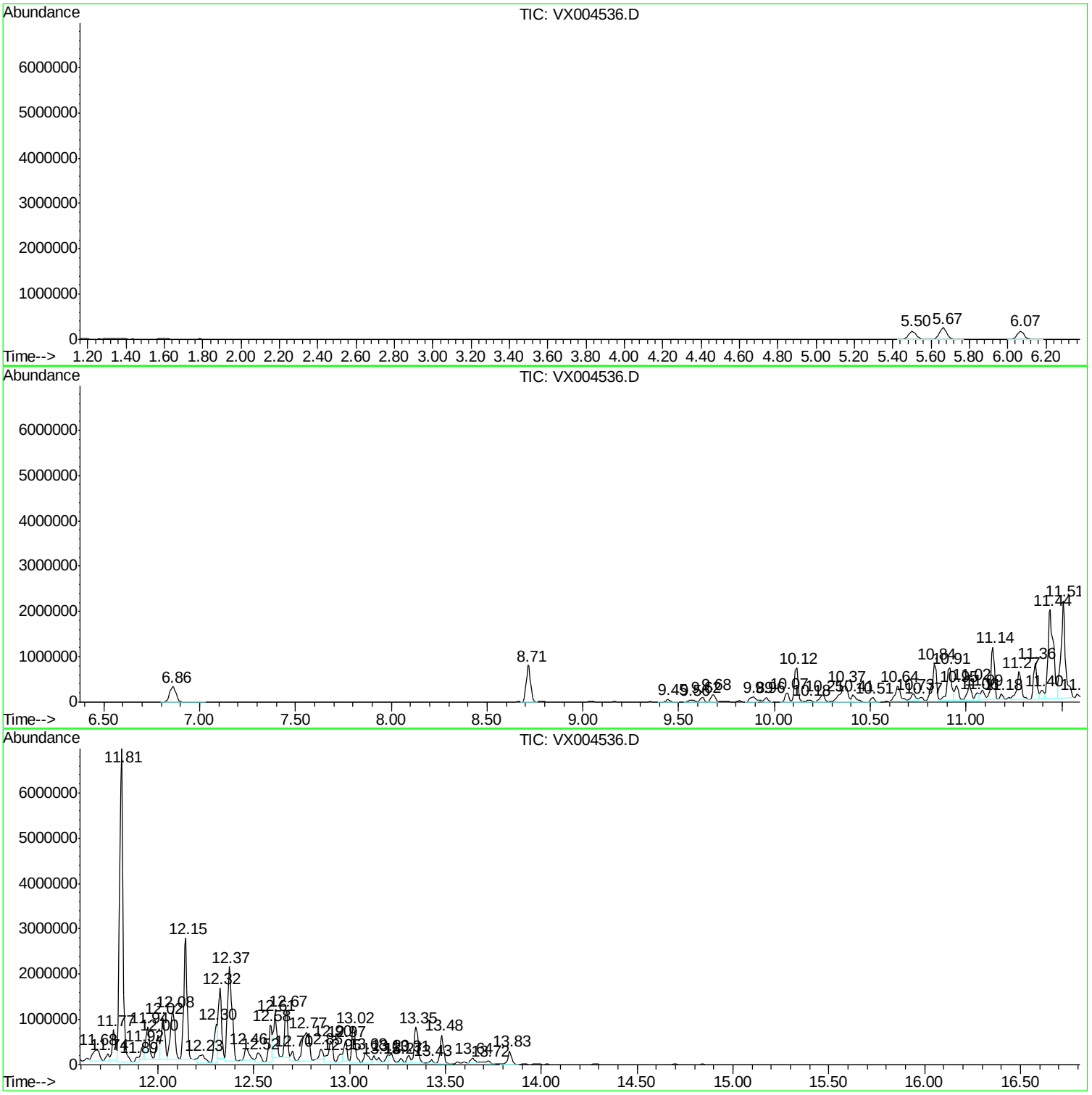
LSC Report - Integrated Chromatogram

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX091418\
Data File : VX004536.D
Acq On : 14 Sep 2018 13:34
Operator : JC/MD
Sample : J4724-18 50X
Misc : 3.19g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : EP1-SUT-4USTS-SB

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

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



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX091418\
Data File : VX004536.D
Acq On : 14 Sep 2018 13:34
Operator : JC/MD
Sample : J4724-18 50X
Misc : 3.19g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : EP1-SUT-4USTS-SB

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

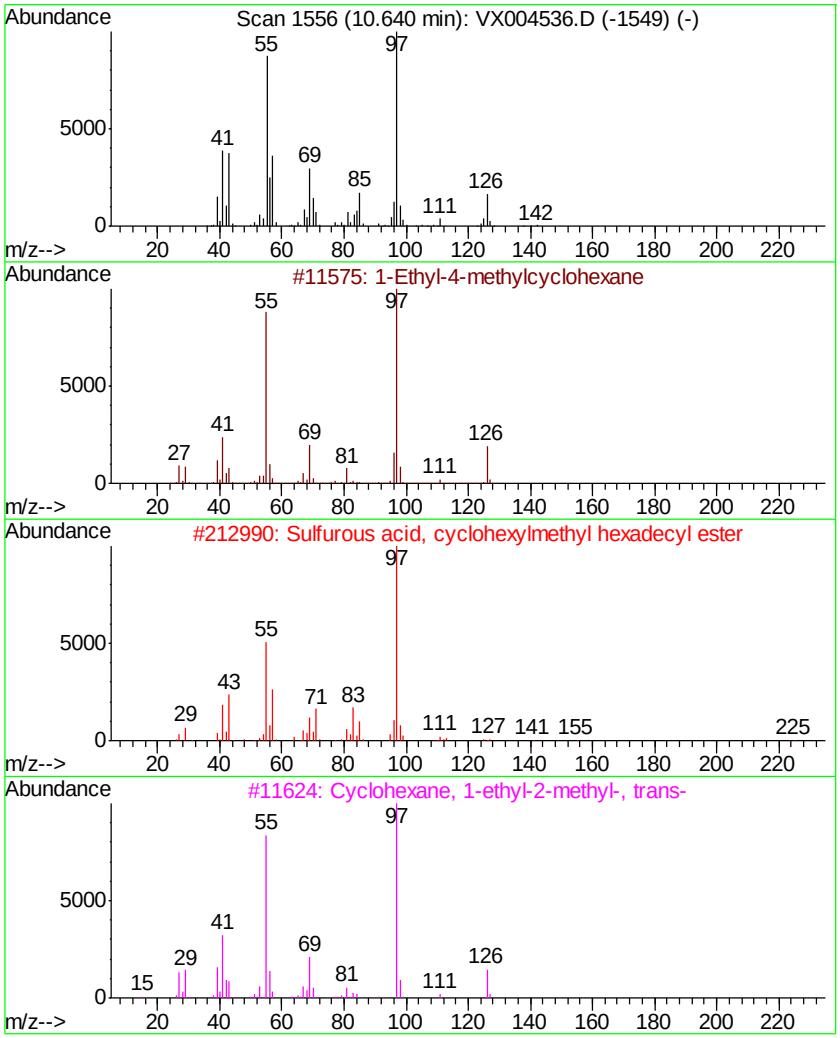
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

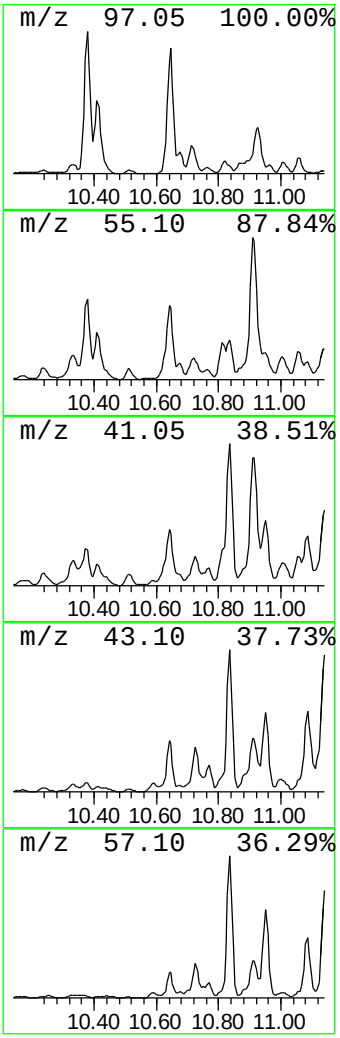
****

Peak Number 1 1-Ethyl-4-methylcyclohexane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.64	31.43 ug/l	644088	Chlorobenzene-d5	10.12

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Ethyl-4-methylcyclohexane	126	C9H18	003728-56-1	81
2	Sulfurous acid, cyclohexylmethyl...	402	C23H46O3S	1000309-22-4	80
3	Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-78-8	76
4	Cyclohexane, 1-ethyl-2-methyl-	126	C9H18	003728-54-9	74
5	3,5-Dimethyl-3-heptene	126	C9H18	059643-68-4	64





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX091418\
Data File : VX004536.D
Acq On : 14 Sep 2018 13:34
Operator : JC/MD
Sample : J4724-18 50X
Misc : 3.19g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : EP1-SUT-4USTS-SB

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

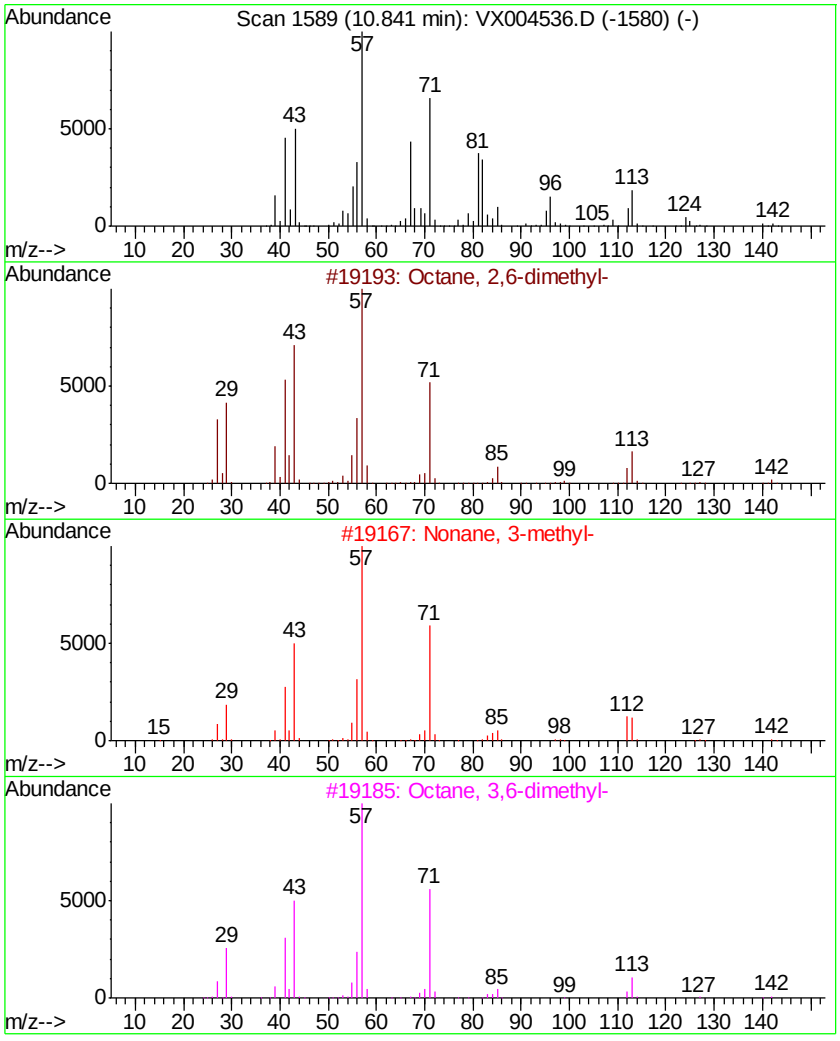
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

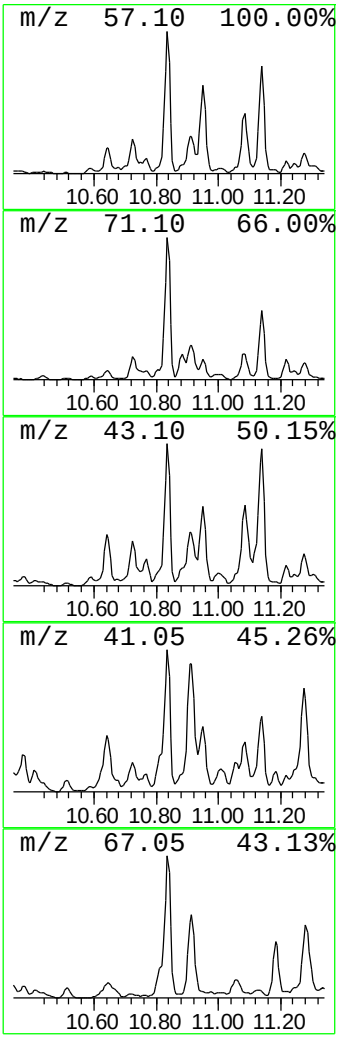
****

Peak Number 2 Octane, 2,6-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.84	59.84 ug/l	1226470	Chlorobenzene-d5	10.12

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	55
2	Nonane, 3-methyl-	142	C10H22	005911-04-6	55
3	Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	49
4	Sulfurous acid, di(2-ethylhexyl)...	306	C16H34O3S	1000309-19-1	35
5	Butane, 2,2-dimethyl-	86	C6H14	000075-83-2	35





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX091418\
Data File : VX004536.D
Acq On : 14 Sep 2018 13:34
Operator : JC/MD
Sample : J4724-18 50X
Misc : 3.19g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : EP1-SUT-4USTS-SB

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

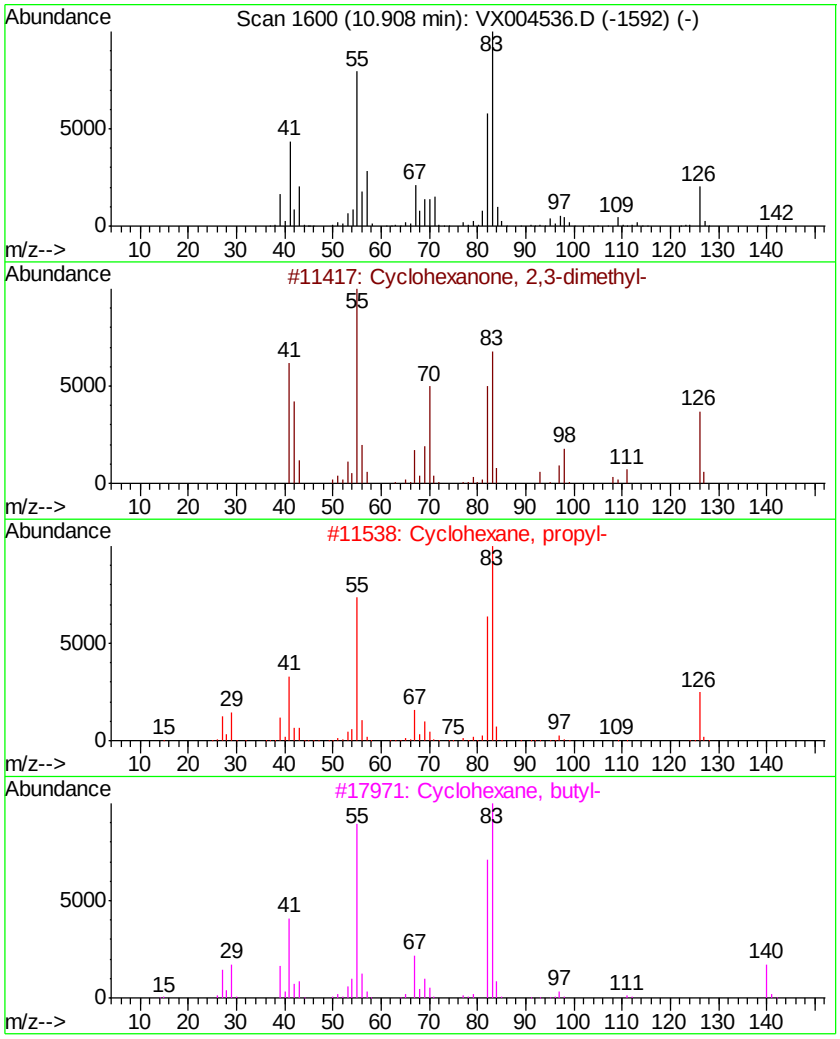
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

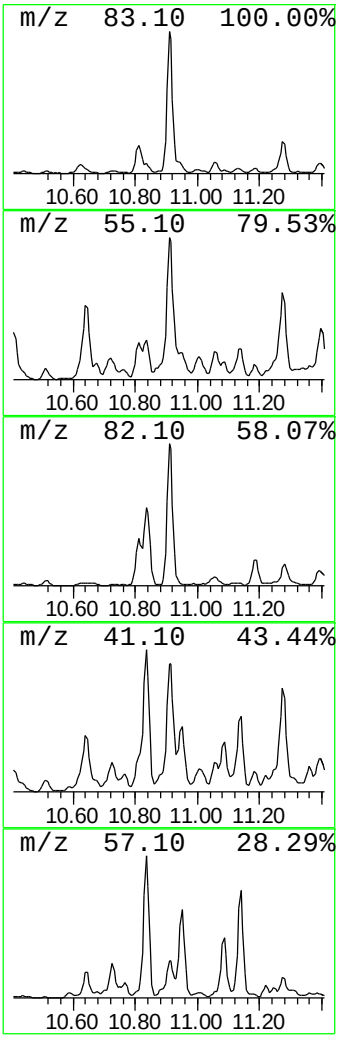
****

Peak Number 3 Cyclohexanone, 2,3-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.91	58.78 ug/l	1204810	Chlorobenzene-d5	10.12

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanone, 2,3-dimethyl-	126	C8H14O	013395-76-1	91
2	Cyclohexane, propyl-	126	C9H18	001678-92-8	74
3	Cyclohexane, butyl-	140	C10H20	001678-93-9	64
4	Cyclopentane, 1-ethyl-1-methyl-	112	C8H16	016747-50-5	62
5	Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	62





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX091418\
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Instrument : MSVOA_X
ClientSampleId : EP1-SUT-4USTS-SB

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
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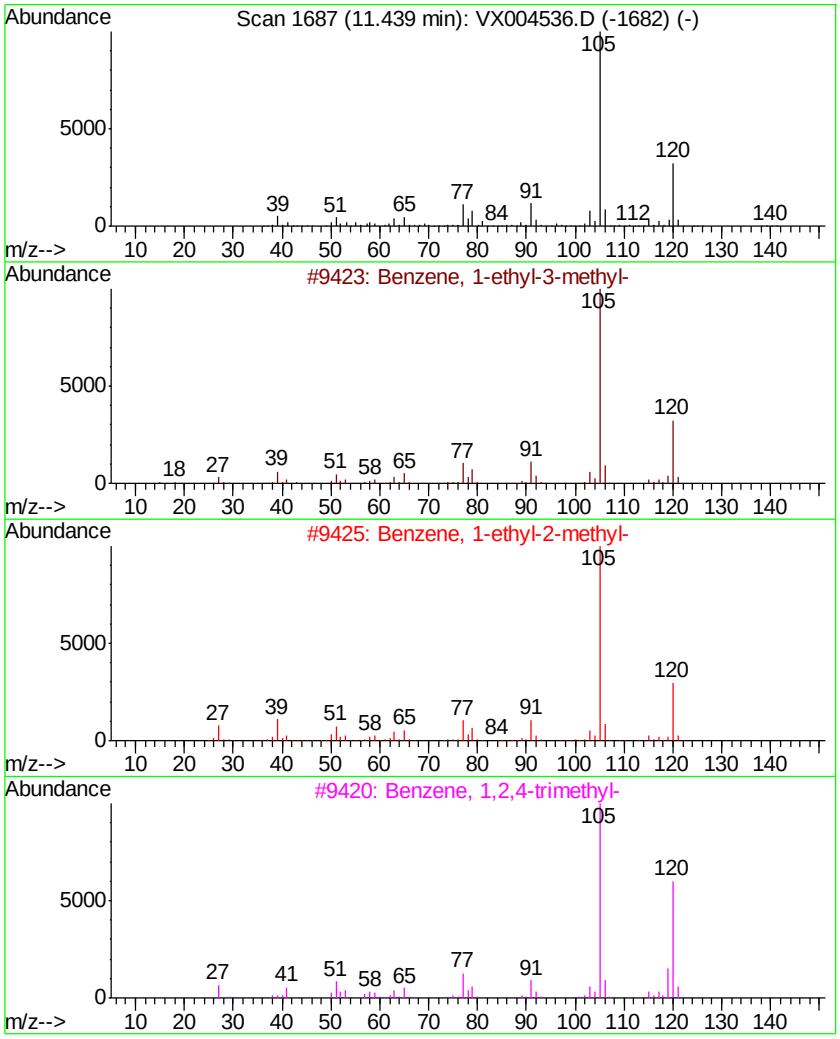
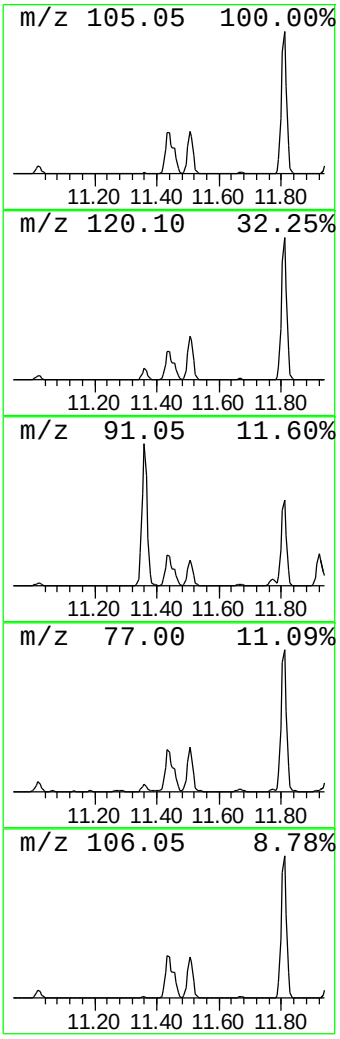
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.44	110.58 ug/l	3918090	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
2	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5	Mesitylene	120	C9H12	000108-67-8	91





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Instrument : MSVOA_X
ClientSampleId : EP1-SUT-4USTS-SB

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

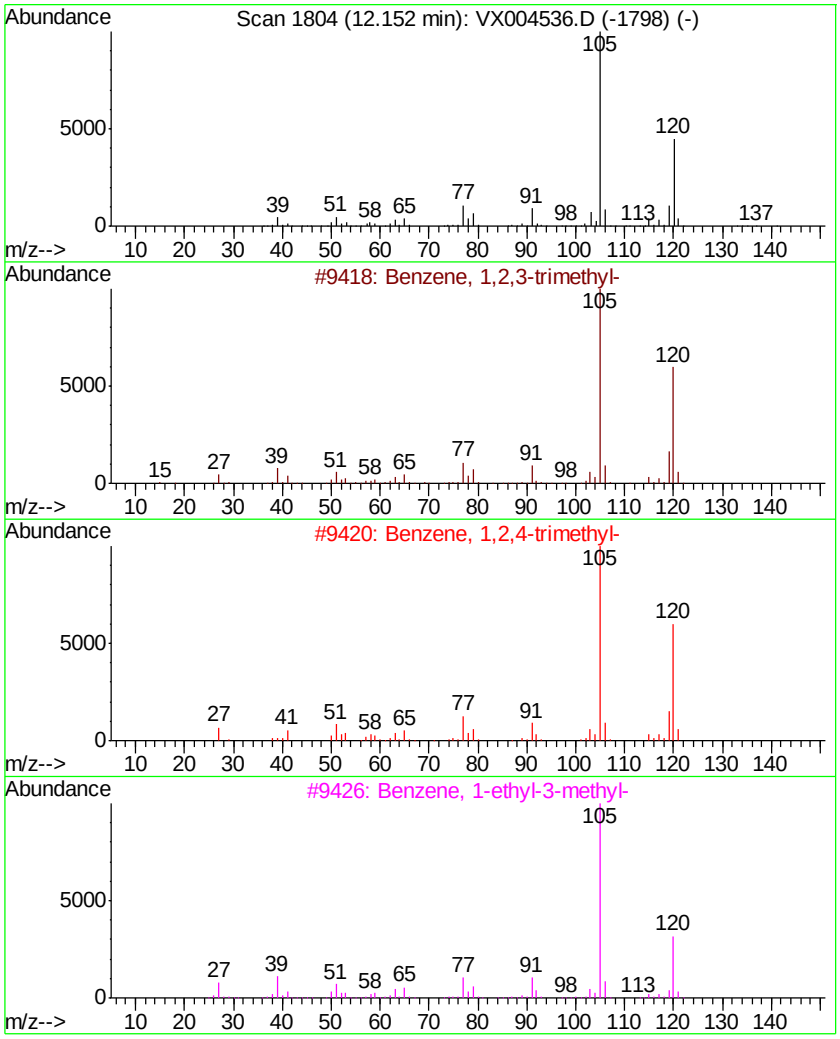
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

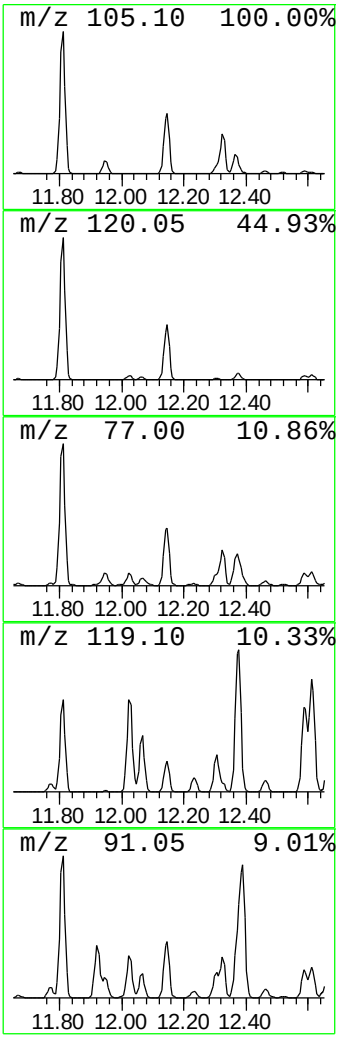
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.15	94.43 ug/l	3345610	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
3	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4	Mesitylene	120	C9H12	000108-67-8	91
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX091418\
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ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : EP1-SUT-4USTS-SB

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

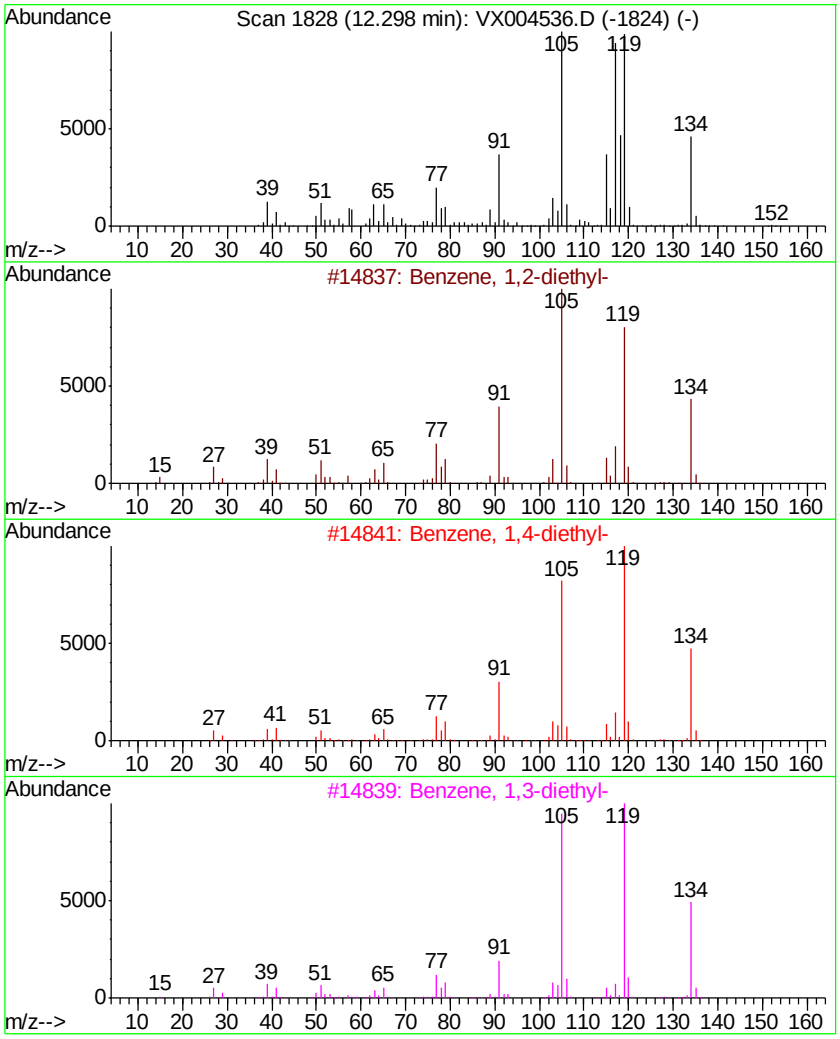
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

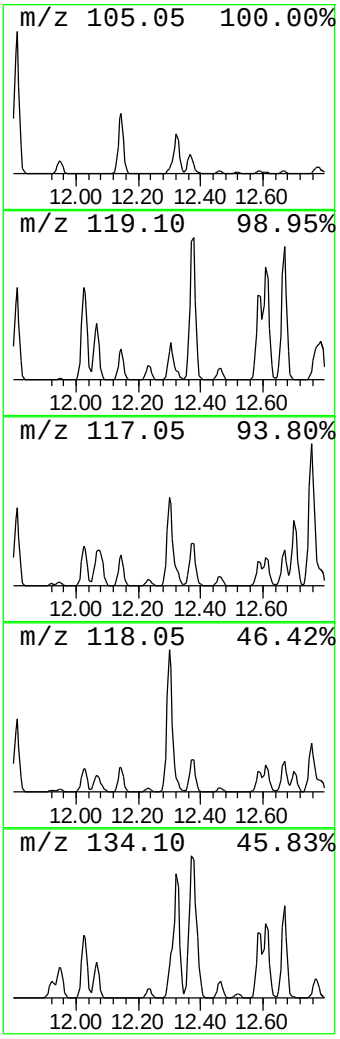
****

Peak Number 6 Benzene, 1,2-diethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	30.37 ug/l	1076010	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	94
2	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	76
3	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	70
4	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	52
5	Ethanone, 1-(3-methylphenyl)-	134	C9H10O	000585-74-0	50





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX091418\
Data File : VX004536.D
Acq On : 14 Sep 2018 13:34
Operator : JC/MD
Sample : J4724-18 50X
Misc : 3.19g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : EP1-SUT-4USTS-SB

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

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

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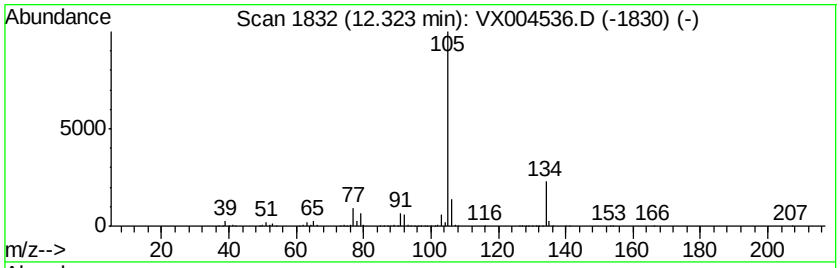
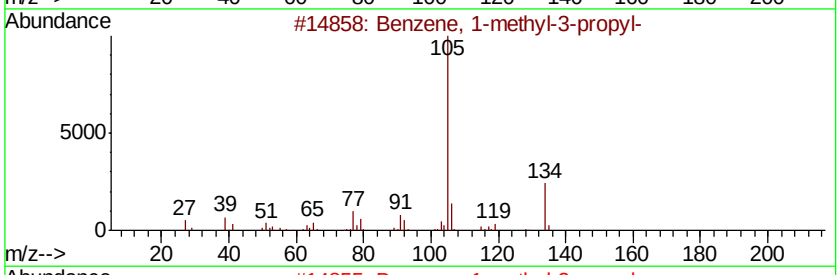
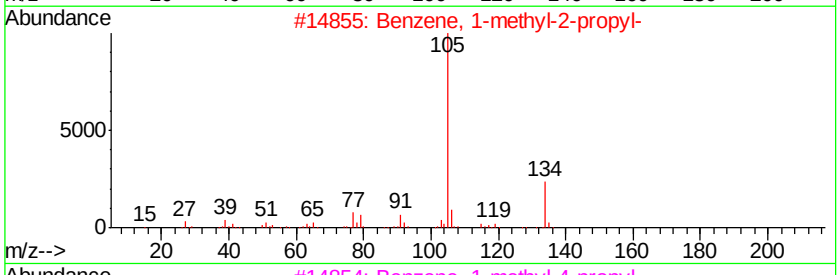
Peak Number 7 Benzene, 1-methyl-3-propyl- Concentration Rank 5

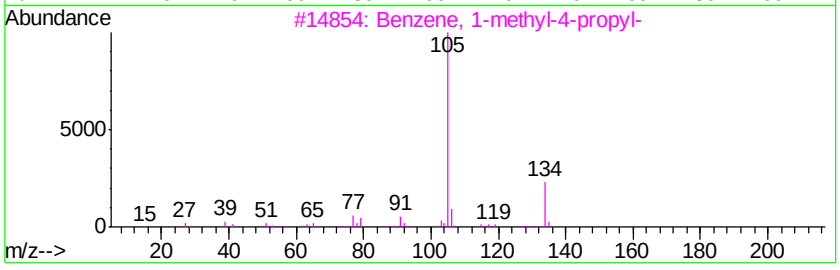
R.T.	EstConc	Area	Relative to ISTD	R.T.
12.32	46.56 ug/l	1649570	1,4-Dichlorobenzene-d4	12.08

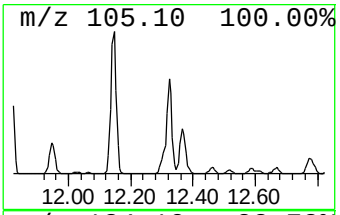
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	91
2	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	91
3	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
4	Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	78
5	2-Indanol	134	C9H10O	004254-29-9	72

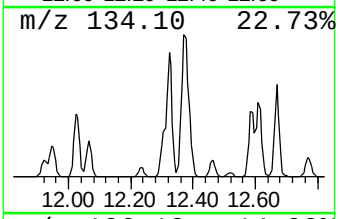


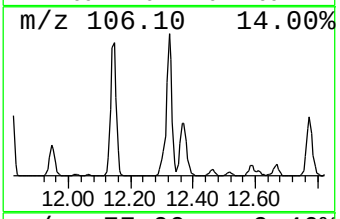


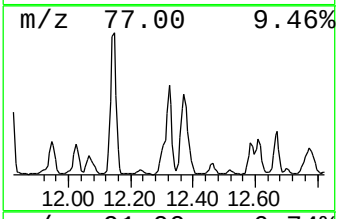


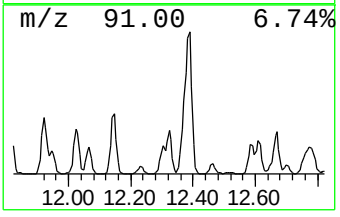












Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX091418\
Data File : VX004536.D
Acq On : 14 Sep 2018 13:34
Operator : JC/MD
Sample : J4724-18 50X
Misc : 3.19g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : EP1-SUT-4USTS-SB

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

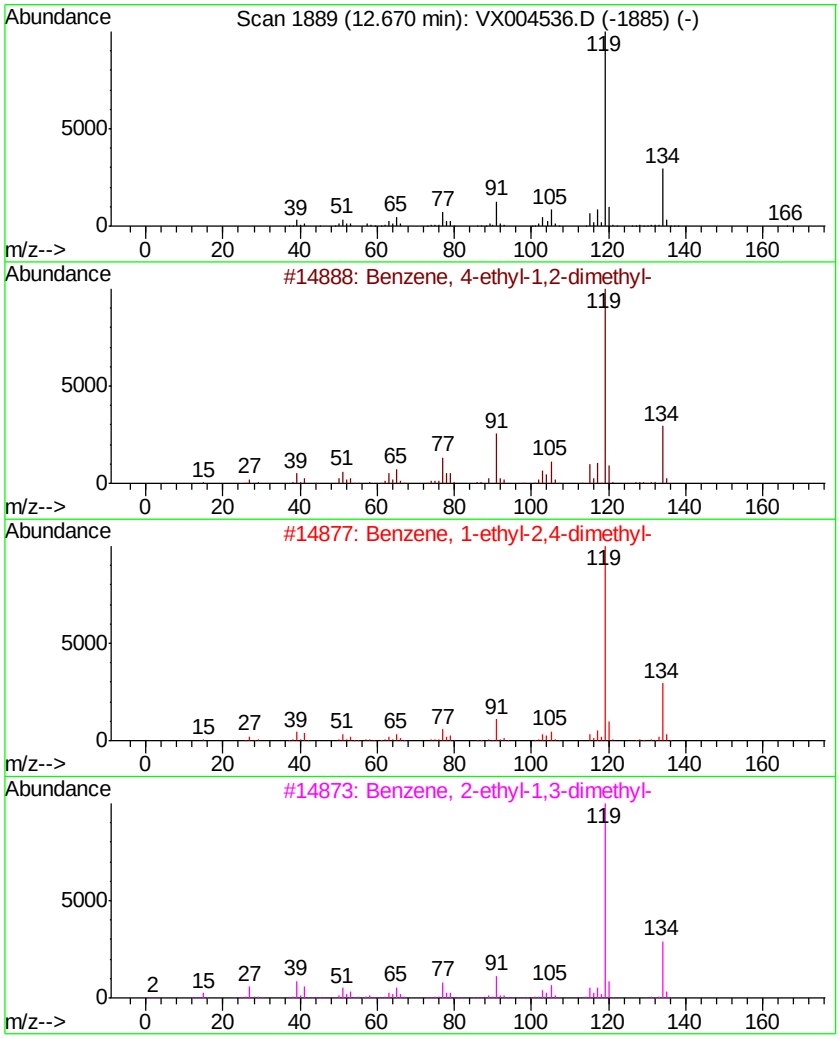
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

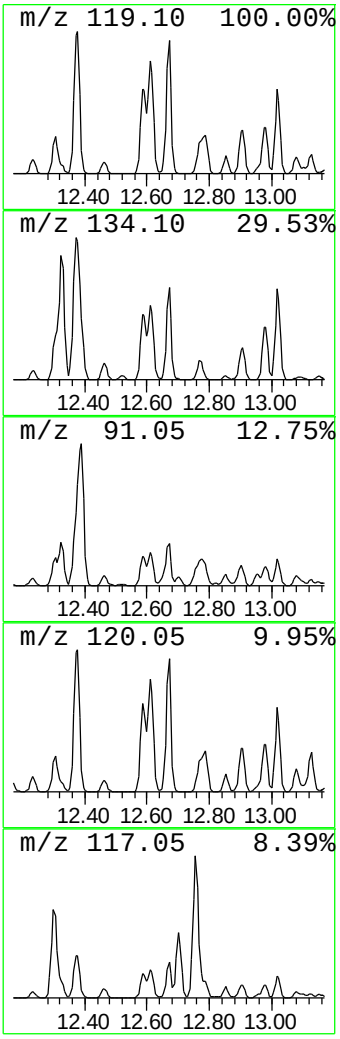
****

Peak Number 8 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.67	31.39 ug/l	1112220	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	95
3	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94
4	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX091418\
Data File : VX004536.D
Acq On : 14 Sep 2018 13:34
Operator : JC/MD
Sample : J4724-18 50X
Misc : 3.19g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : EP1-SUT-4USTS-SB

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

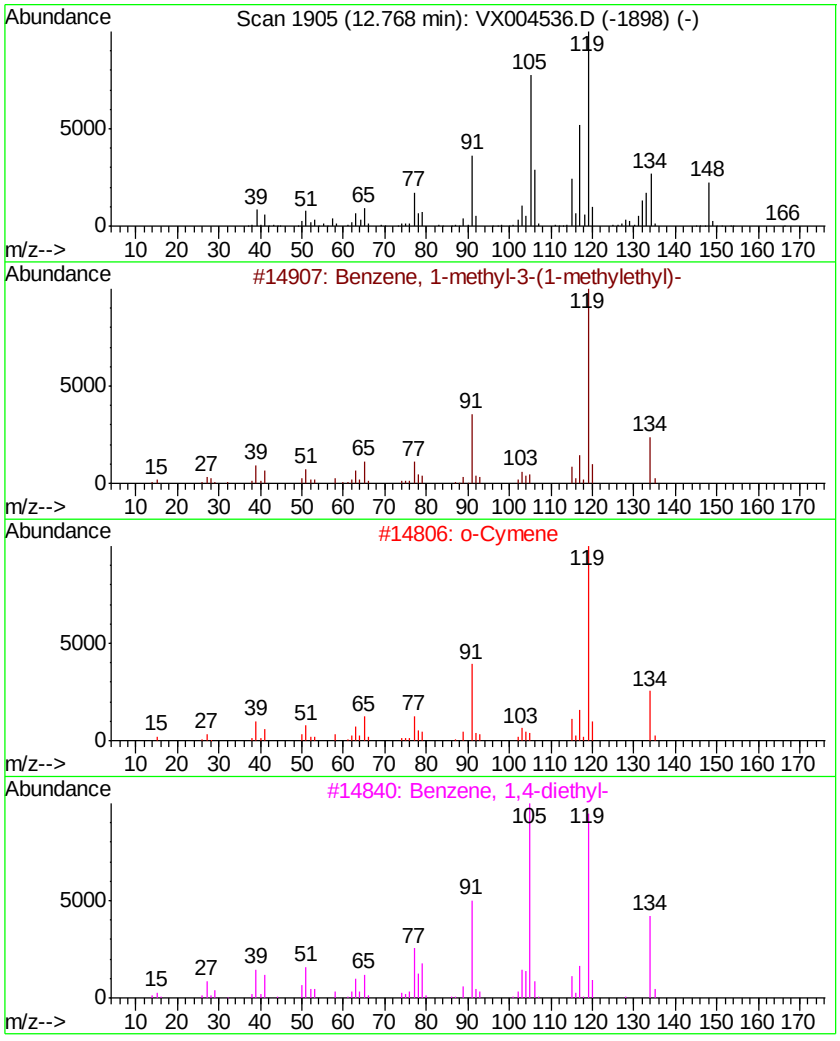
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

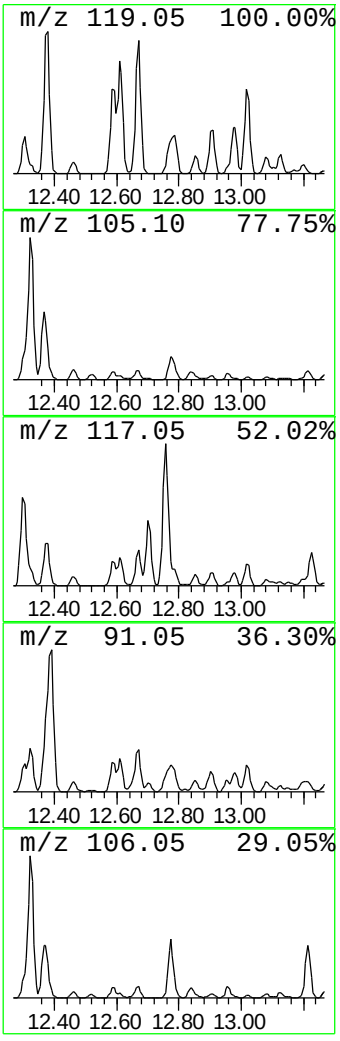
****

Peak Number 9 Benzene, 1-methyl-3-(1-meth... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.77	43.63 ug/l	1545920	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	64
2	o-Cymene	134	C10H14	000527-84-4	60
3	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	58
4	p-Cymene	134	C10H14	000099-87-6	58
5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	53





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX091418\
Data File : VX004536.D
Acq On : 14 Sep 2018 13:34
Operator : JC/MD
Sample : J4724-18 50X
Misc : 3.19g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : EP1-SUT-4USTS-SB

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

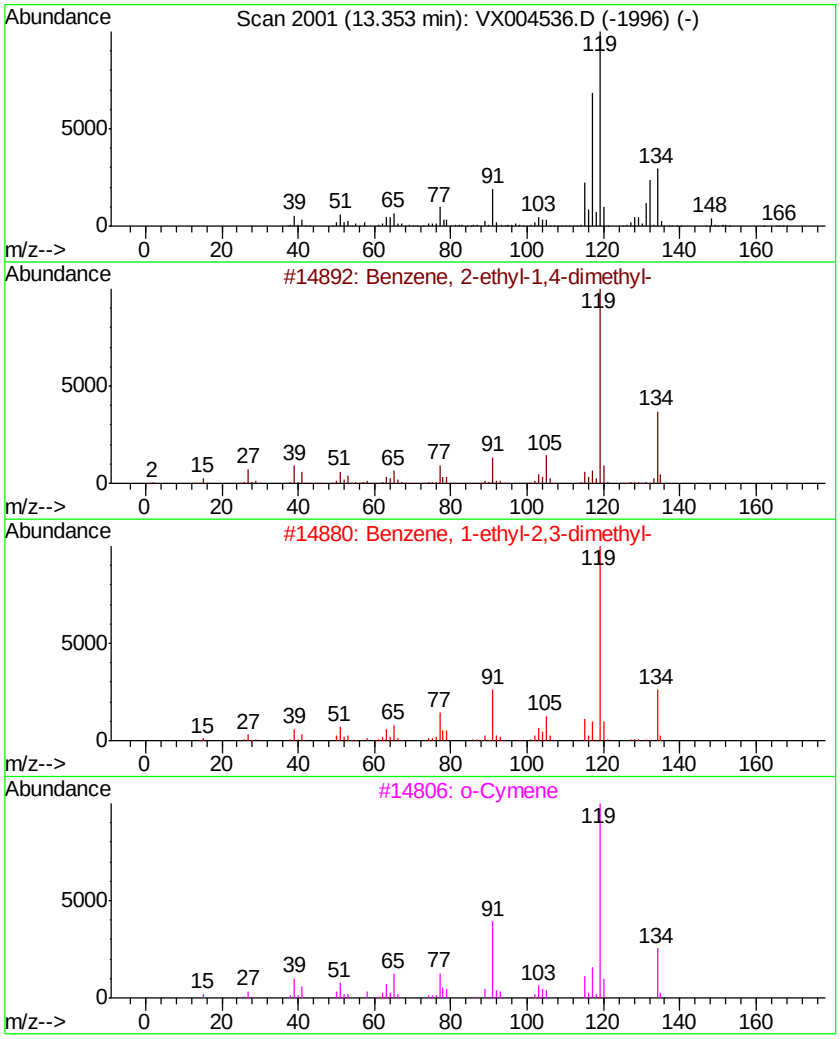
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

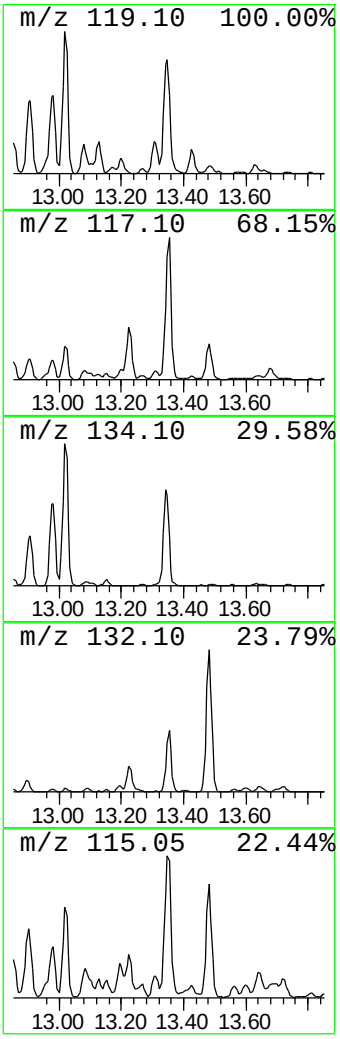
****

Peak Number 10 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	32.74 ug/l	1160080	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	70
2	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	70
3	o-Cymene	134	C10H14	000527-84-4	70
4	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	70
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	64





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX091418\
Data File : VX004536.D
Acq On : 14 Sep 2018 13:34
Operator : JC/MD
Sample : J4724-18 50X
Misc : 3.19g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : EP1-SUT-4USTS-SB

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.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
1-Ethyl-4-methylc...	10.64	31.4	ug/l	644088	3	10.12	1024790	50.0
Octane, 2,6-dimet...	10.84	59.8	ug/l	1226470	3	10.12	1024790	50.0
Cyclohexanone, 2,...	10.91	58.8	ug/l	1204810	3	10.12	1024790	50.0
Benzene, 1-ethyl-...	11.44	110.6	ug/l	3918090	4	12.08	1771570	50.0
Benzene, 1,2,3-tr...	12.15	94.4	ug/l	3345610	4	12.08	1771570	50.0
Benzene, 1,2-diet...	12.30	30.4	ug/l	1076010	4	12.08	1771570	50.0
Benzene, 1-methyl...	12.32	46.6	ug/l	1649570	4	12.08	1771570	50.0
Benzene, 4-ethyl-...	12.67	31.4	ug/l	1112220	4	12.08	1771570	50.0
Benzene, 1-methyl...	12.77	43.6	ug/l	1545920	4	12.08	1771570	50.0
Benzene, 2-ethyl-...	13.35	32.7	ug/l	1160080	4	12.08	1771570	50.0