LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW011419\
Data File : VW008248.D
Acq On : 14 Jan 2019 15:44
Operator : SY/VA
Sample : K1135-01
Misc : 5.00G/5ML/MSVOA W/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_W
ClientSampleId : CPSB-19(10-15)

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_W\METHOD\82W010819S.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	7.884	971	981	986	rBV3	67441	162718	7.40%	0.461%
2	7.945	986	991	1001	rVV	119071	263812	11.99%	0.747%
3	8.159	1018	1026	1031	rBV3	10543	23976	1.09%	0.068%
4	8.305	1043	1050	1062	rVB2	76335	190337	8.65%	0.539%
5	8.427	1064	1070	1077	rVV5	13369	36458	1.66%	0.103%
6	8.500	1077	1082	1088	rVV5	15379	35374	1.61%	0.100%
7	8.573	1088	1094	1103	rVB3	18017	44911	2.04%	0.127%
8	8.842	1129	1138	1149	rBV	183858	360998	16.41%	1.022%
9	9.335	1204	1219	1225	rBV	151112	357874	16.27%	1.014%
10	9.604	1254	1263	1271	rVB3	16546	40273	1.83%	0.114%
11	9.750	1279	1287	1296	rBV3	21953	46066	2.09%	0.130%
12	9.896	1305	1311	1317	rBV2	13566	26837	1.22%	0.076%
13	10.079	1334	1341	1346	rBV5	10948	26902	1.22%	0.076%
14	10.323	1373	1381	1395	rBV	378406	714792	32.50%	2.025%
15	10.494	1404	1409	1420	rVB3	27051	72402	3.29%	0.205%
16	10.664	1431	1437	1447	rVB2	60235	110753	5.04%	0.314%
17	10.762	1447	1453	1458	rBV3	32442	60584	2.75%	0.172%
18	10.847	1463	1467	1473	rVB2	15807	24713	1.12%	0.070%
19	10.939	1473	1482	1488	rVB	43629	76837	3.49%	0.218%
20	11.036	1492	1498	1506	rBV2	23365	61468	2.79%	0.174%
21	11.109	1506	1510	1513	rBV2	27040	39980	1.82%	0.113%
22	11.176	1513	1521	1525	rBV	136975	235401	10.70%	0.667%
23	11.231	1525	1530	1536	rVB	182956	304560	13.85%	0.863%
24	11.286	1536	1539	1542	rVB2	22967	28692	1.30%	0.081%
25	11.341	1542	1548	1552	rBV4	68165	112372	5.11%	0.318%
26	11.439	1559	1564	1571	rVB	103687	185582	8.44%	0.526%
27	11.512	1571	1576	1580	rBV2	31967	54088	2.46%	0.153%
28	11.658	1589	1600	1606	rBV2	1013945	2113849	96.11%	5.987%
29	11.725	1606	1611	1614	rBV2	38193	55762	2.54%	0.158%
30	11.768	1614	1618	1621	rBV	51654	73554	3.34%	0.208%
31	11.810	1621	1625	1630	rVB4	68001	134159	6.10%	0.380%
32	11.878	1630	1636	1639	rBV	172011	313365	14.25%	0.888%
33	11.920	1640	1643	1648	rVB2	129140	200162	9.10%	0.567%
34	11.975	1648	1652	1658	rBV4	29399	59940	2.73%	0.170%

LSC Area Percent Report

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Sample : K1135-01
Misc : 5.00G/5ML/MSVOA W/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_W
ClientSampleId : CPSB-19(10-15)

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_W\METHOD\82W010819S.M
Title : SW846 8260

35	12.054	1658	1665	1671	rVB3	46670	120463	5.48%	0.341%
36	12.140	1671	1679	1685	rBV3	250948	577674	26.27%	1.636%
37	12.256	1691	1698	1702	rVB2	226490	389539	17.71%	1.103%
38	12.304	1702	1706	1708	rBV2	85963	143104	6.51%	0.405%
39	12.341	1708	1712	1714	rBV2	218958	342271	15.56%	0.969%
40	12.463	1727	1732	1737	rBV	192045	323468	14.71%	0.916%
41	12.621	1752	1758	1763	rVB2	287481	470014	21.37%	1.331%
42	12.701	1763	1771	1776	rBV5	127166	331742	15.08%	0.940%
43	12.804	1780	1788	1793	rVB2	402725	1064113	48.38%	3.014%
44	12.859	1793	1797	1802	rBV4	100545	211090	9.60%	0.598%
45	12.944	1802	1811	1816	rVV3	308297	822044	37.38%	2.328%
46	12.993	1816	1819	1824	rVB3	169624	270181	12.28%	0.765%
47	13.048	1824	1828	1829	rBV2	24205	29424	1.34%	0.083%
48	13.085	1829	1834	1837	rBV4	146097	213670	9.72%	0.605%
49	13.127	1837	1841	1844	rBV3	92728	150293	6.83%	0.426%
50	13.170	1844	1848	1860	rVB5	271407	618593	28.13%	1.752%
51	13.268	1860	1864	1866	rBV2	65141	92204	4.19%	0.261%
52	13.298	1866	1869	1872	rBV2	64337	74127	3.37%	0.210%
53	13.353	1873	1878	1879	rBV4	97505	152527	6.94%	0.432%
54	13.383	1879	1883	1888	rVB2	321490	496766	22.59%	1.407%
55	13.450	1888	1894	1898	rBV4	302100	598294	27.20%	1.695%
56	13.560	1907	1912	1921	rVB	275792	526915	23.96%	1.492%
57	13.658	1921	1928	1930	rBV4	73704	146026	6.64%	0.414%
58	13.725	1931	1939	1942	rBV2	276017	573894	26.09%	1.625%
59	13.761	1942	1945	1949	rVB	347644	491554	22.35%	1.392%
60	13.828	1949	1956	1960	rBV2	180136	386005	17.55%	1.093%
61	13.883	1960	1965	1970	rBV2	681522	1158910	52.69%	3.282%
62	13.981	1977	1981	1985	rBV4	124469	219963	10.00%	0.623%
63	14.103	1996	2001	2005	rVB	693543	1034591	47.04%	2.930%
64	14.213	2013	2019	2024	rVB4	442280	842015	38.28%	2.385%
65	14.267	2024	2028	2034	rBV5	153792	284316	12.93%	0.805%
66	14.322	2034	2037	2040	rBV3	100748	172031	7.82%	0.487%
67	14.389	2044	2048	2051	rVV3	526255	914003	41.56%	2.589%
68	14.426	2051	2054	2062	rVB2	493372	788204	35.84%	2.232%
69	14.505	2063	2067	2070	rBV	153191	227289	10.33%	0.644%
70	14.554	2070	2075	2081	rVV4	261567	542639	24.67%	1.537%
71	14.627	2081	2087	2095	rVV3	157125	461412	20.98%	1.307%

LSC Area Percent Report

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ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_W
ClientSampleId : CPSB-19(10-15)

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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Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W010819S.M
Title : SW846 8260

72	14.700	2096	2099	2102	rVV3	73623	111383	5.06%	0.315%
73	14.743	2103	2106	2109	rVV3	140218	194108	8.83%	0.550%
74	14.798	2110	2115	2121	rVV2	1119532	2199375	100.00%	6.229%
75	14.853	2121	2124	2131	rVB4	408768	848120	38.56%	2.402%
76	15.017	2148	2151	2152	rBV2	55464	69373	3.15%	0.196%
77	15.072	2156	2160	2163	rVB	276309	411128	18.69%	1.164%
78	15.115	2163	2167	2172	rBV2	131732	260977	11.87%	0.739%
79	15.182	2173	2178	2184	rVB3	319170	686298	31.20%	1.944%
80	15.334	2198	2203	2210	rVB2	608445	1059376	48.17%	3.001%
81	15.426	2211	2218	2223	rBV	433330	1022446	46.49%	2.896%
82	15.511	2228	2232	2239	rVB2	358994	819763	37.27%	2.322%
83	15.670	2250	2258	2264	rBV8	129322	417691	18.99%	1.183%
84	15.743	2265	2270	2276	rVV4	207781	396703	18.04%	1.124%
85	15.871	2287	2291	2299	rVB	537599	943734	42.91%	2.673%
86	16.188	2336	2343	2354	rVB2	481982	1152571	52.40%	3.265%
87	16.297	2356	2361	2368	rVV	283911	523832	23.82%	1.484%
88	16.383	2369	2375	2382	rVV2	299333	653008	29.69%	1.850%
89	16.450	2382	2386	2399	rVB4	121021	384426	17.48%	1.089%
90	16.651	2412	2419	2420	rBV6	110416	240956	10.96%	0.682%

Sum of corrected areas: 35306187

LSC Report - Integrated Chromatogram

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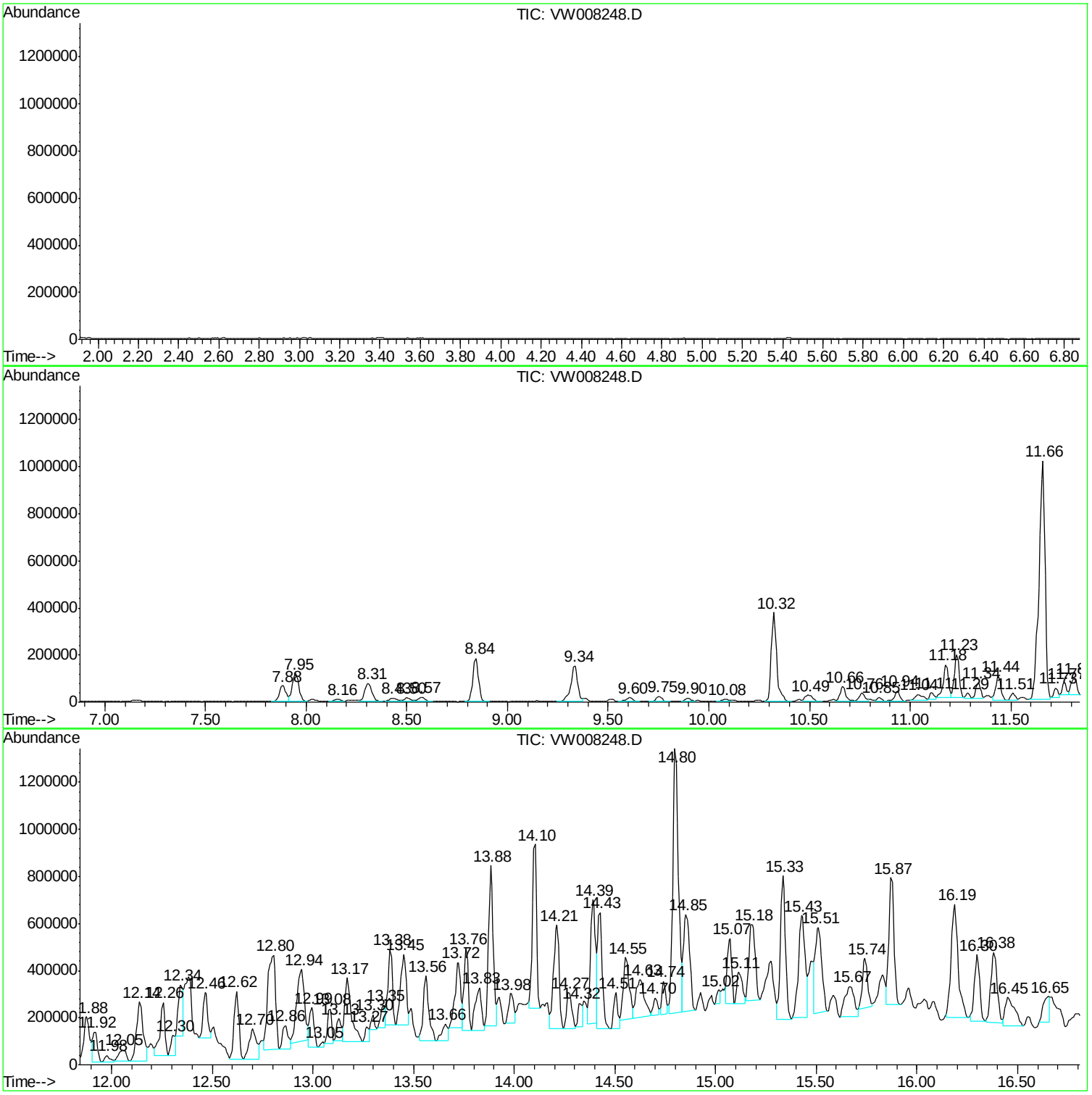
Sample : K1135-01
Misc : 5.00G/5ML/MSVOA W/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_W
ClientSampleId : CPSB-19(10-15)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W010819S.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 W\DATA\VW011419\
Data File : VW008248.D
Acq On : 14 Jan 2019 15:44
Operator : SY/VA
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Instrument : MSVOA_W
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Quant Title : SW846 8260

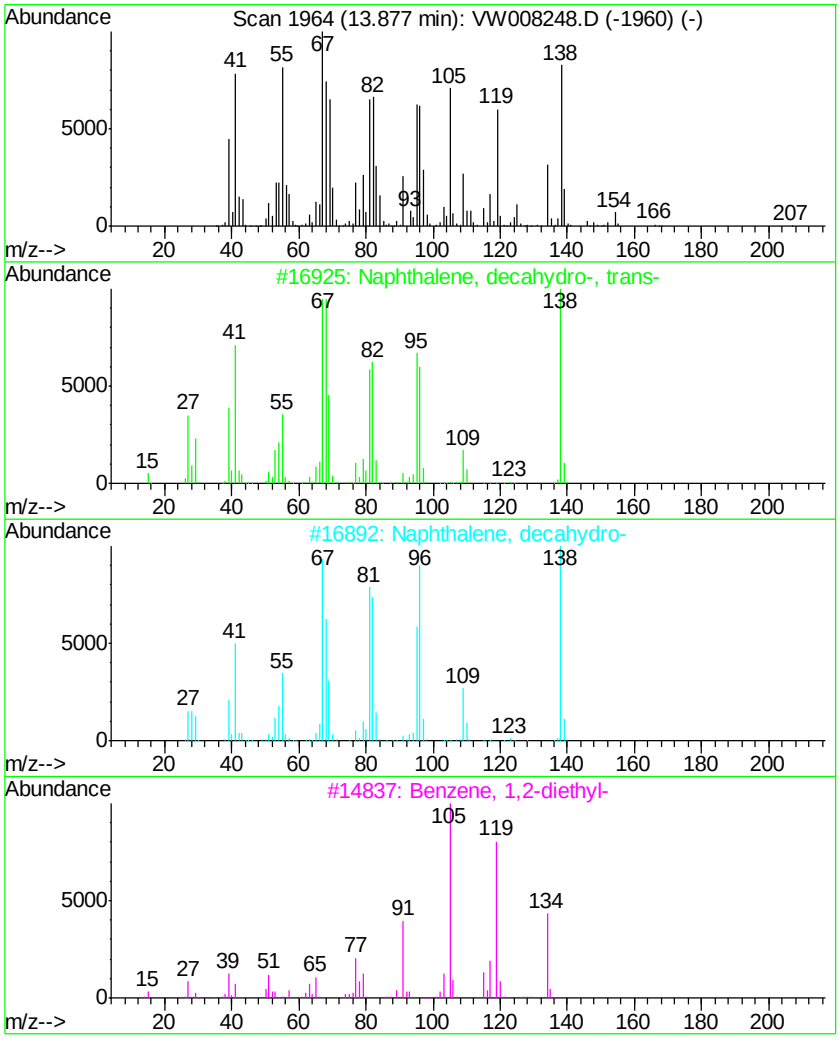
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

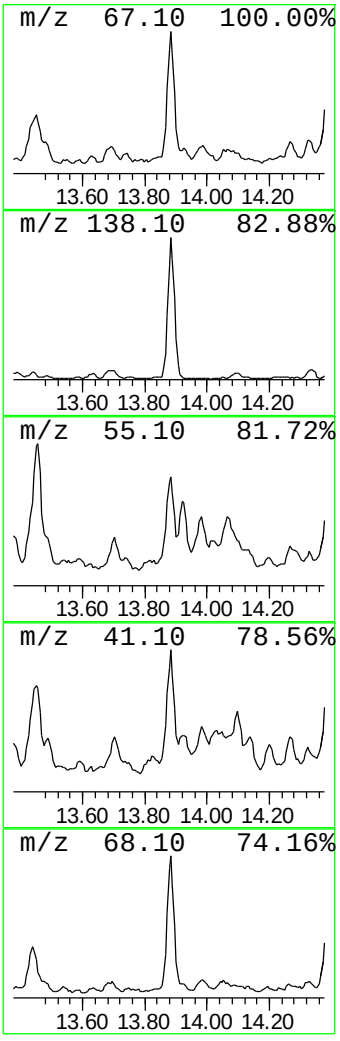
****

Peak Number 1 Naphthalene, decahydro-, tr... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.88	109.97 ug/l	1158910	1,4-Dichlorobenzene-d4	13.56

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, decahydro-, trans-	138	C10H18	000493-02-7	96
2	Naphthalene, decahydro-	138	C10H18	000091-17-8	96
3	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	66
4	Naphthalene, decahydro-, cis-	138	C10H18	000493-01-6	60
5	Spiro[4.5]decane	138	C10H18	000176-63-6	60





Library Search Compound Report

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Data File : VW008248.D
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Sample : K1135-01
Misc : 5.00G/5ML/MSVOA W/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_W
ClientSampleId : CPSB-19(10-15)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W010819S.M
Quant Title : SW846 8260

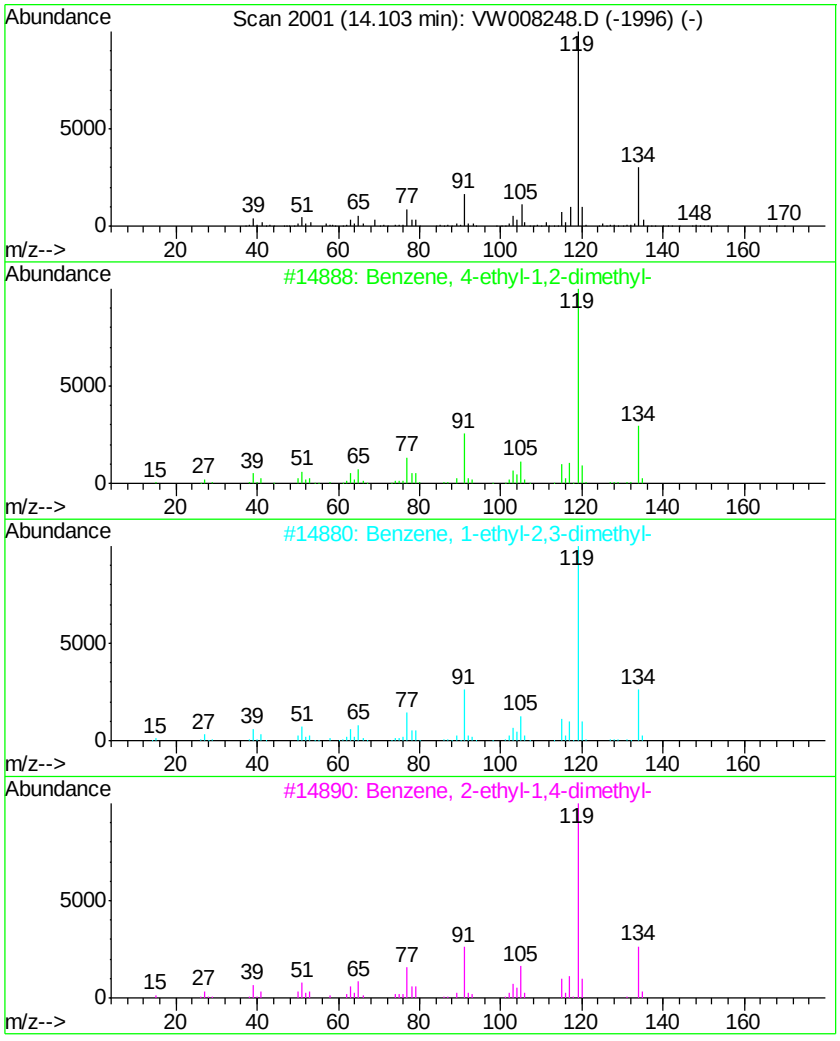
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

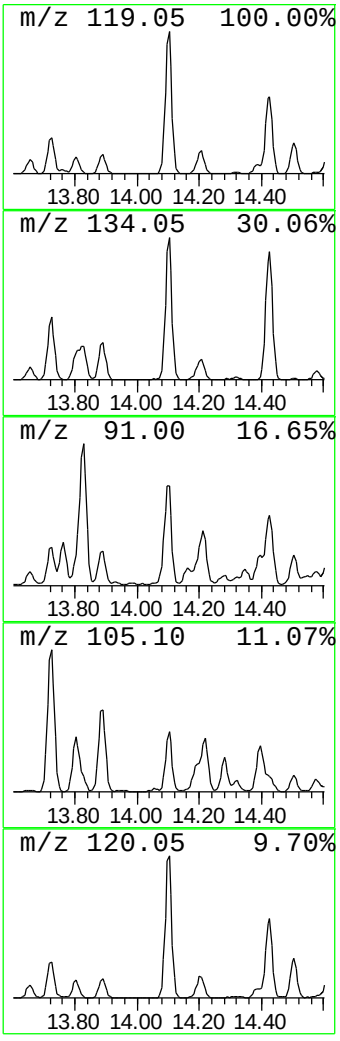
****

Peak Number 2 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.10	98.17 ug/l	1034590	1,4-Dichlorobenzene-d4	13.56

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-2,3-dimethyl-	134	C10H14	000933-98-2	96
3	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
4	o-Cymene	134	C10H14	000527-84-4	94
5	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94





Library Search Compound Report

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ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_W
ClientSampleId : CPSB-19(10-15)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W010819S.M
Quant Title : SW846 8260

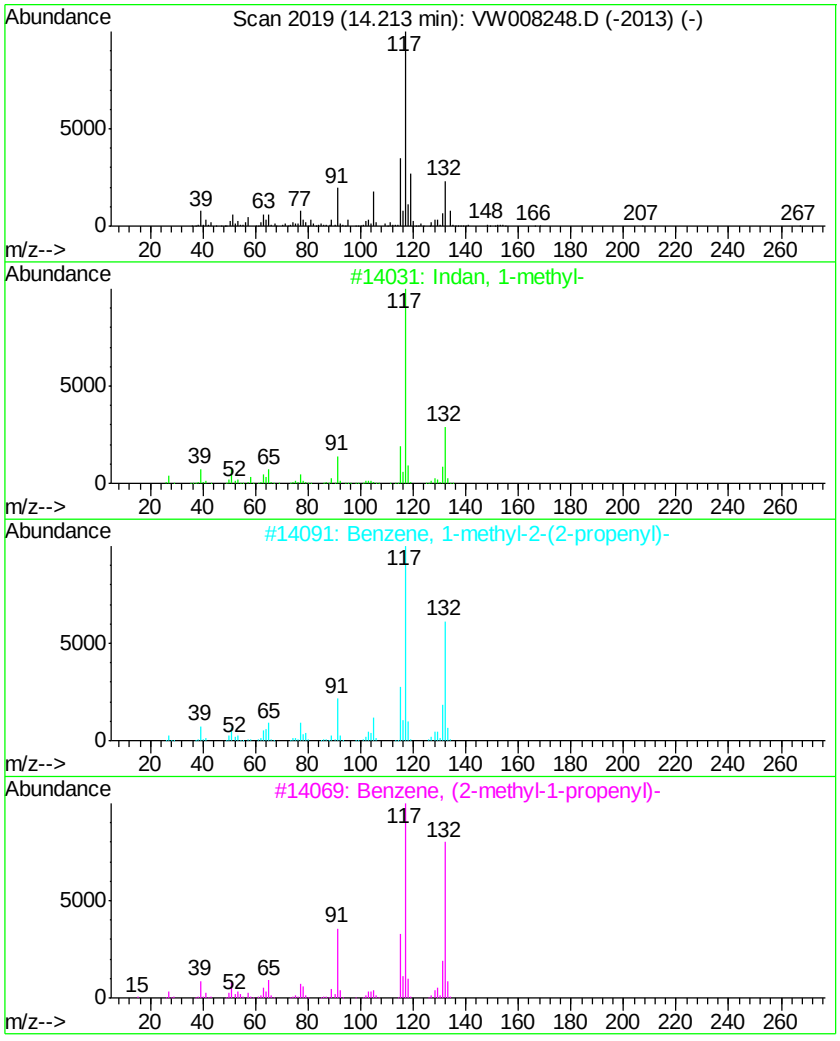
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

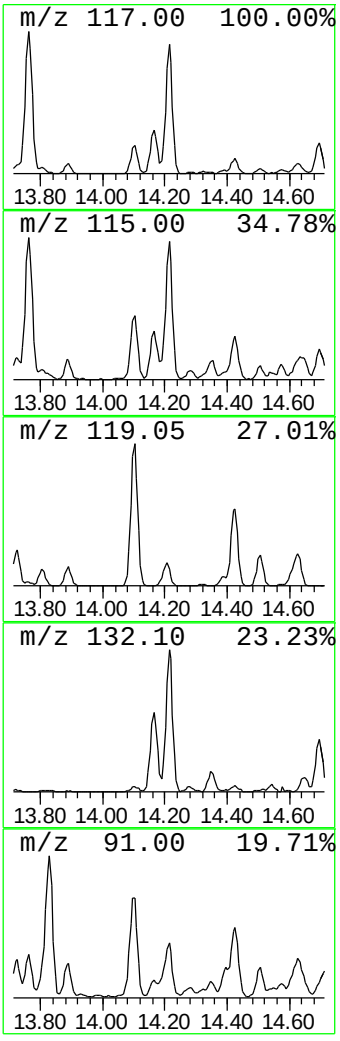
****

Peak Number 3 Indan, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.21	79.90 ug/l	842015	1,4-Dichlorobenzene-d4	13.56

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indan, 1-methyl-	132	C10H12	000767-58-8	76
2	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	74
3	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	74
4	Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	68
5	3-Phenylbut-1-ene	132	C10H12	000934-10-1	64





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Sample : K1135-01
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ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_W
ClientSampleId : CPSB-19(10-15)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W010819S.M
Quant Title : SW846 8260

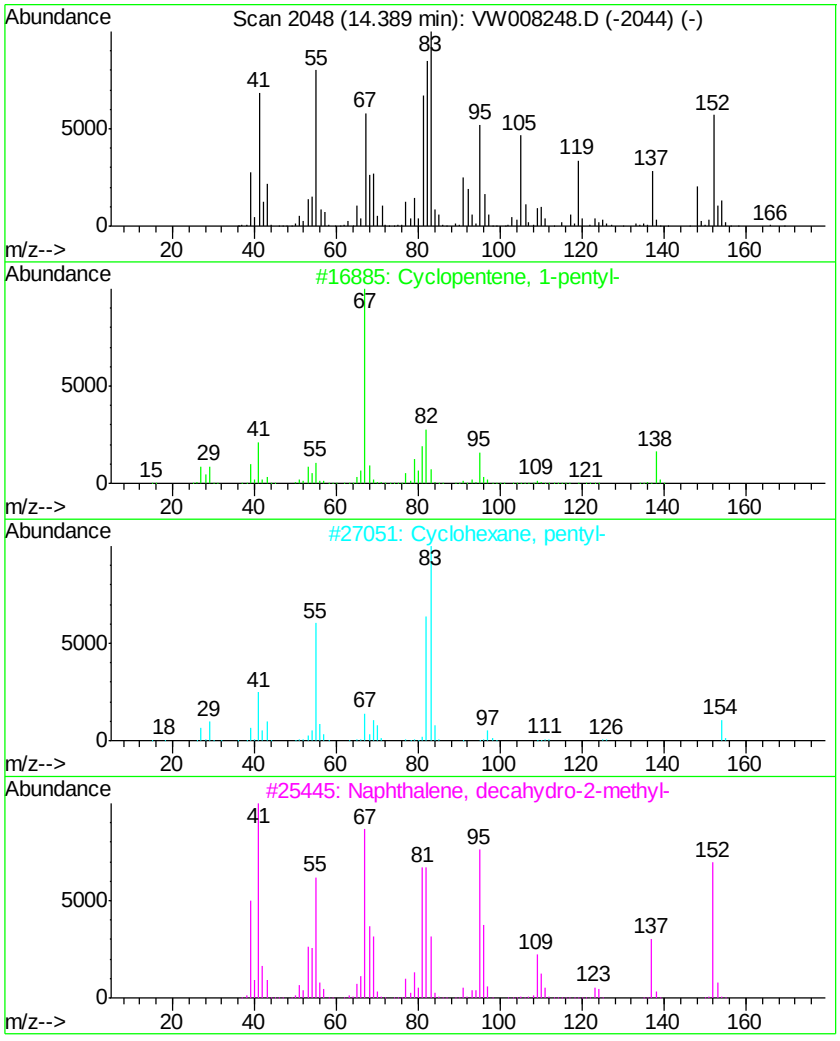
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

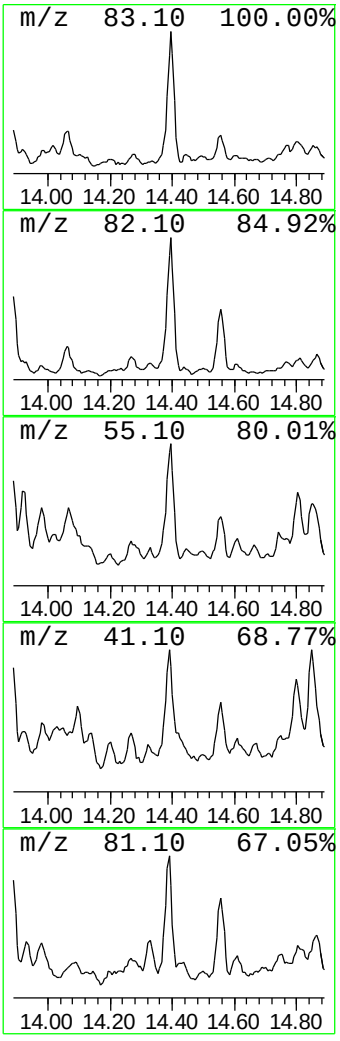
****

Peak Number 4 Cyclopentene, 1-pentyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.39	86.73 ug/l	914003	1,4-Dichlorobenzene-d4	13.56

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentene, 1-pentyl-	138	C10H18	004291-98-9	50
2	Cyclohexane, pentyl-	154	C11H22	004292-92-6	45
3	Naphthalene, decahydro-2-methyl-	152	C11H20	002958-76-1	43
4	Cyclohexane, (1-methylbutyl)-	154	C11H22	061208-94-4	43
5	1-Methyldecahydronaphthalene	152	C11H20	002958-75-0	43





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Data File : VW008248.D
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Operator : SY/VA
Sample : K1135-01
Misc : 5.00G/5ML/MSVOA W/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_W
ClientSampleId : CPSB-19(10-15)

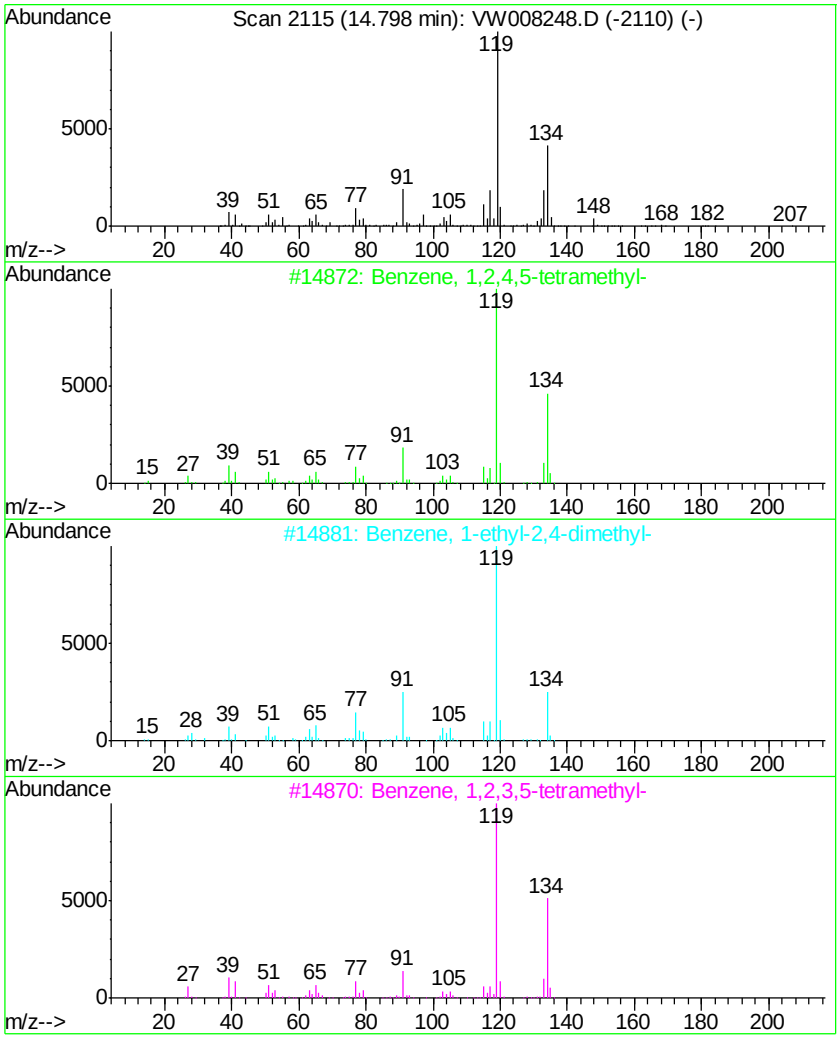
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W010819S.M
Quant Title : SW846 8260

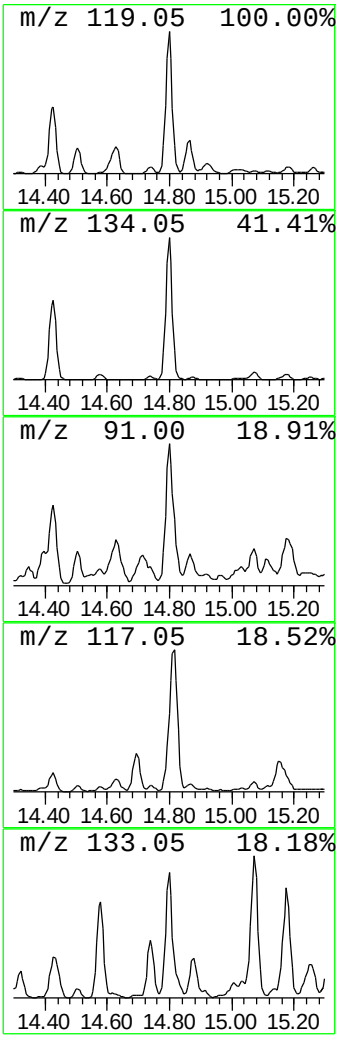
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.80	208.70 ug/l	2199380	1,4-Dichlorobenzene-d4	13.56

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	93
2	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93
3	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	87
4	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	87
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	87





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Operator : SY/VA
Sample : K1135-01
Misc : 5.00G/5ML/MSVOA W/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_W
ClientSampleId : CPSB-19(10-15)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W010819S.M
Quant Title : SW846 8260

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

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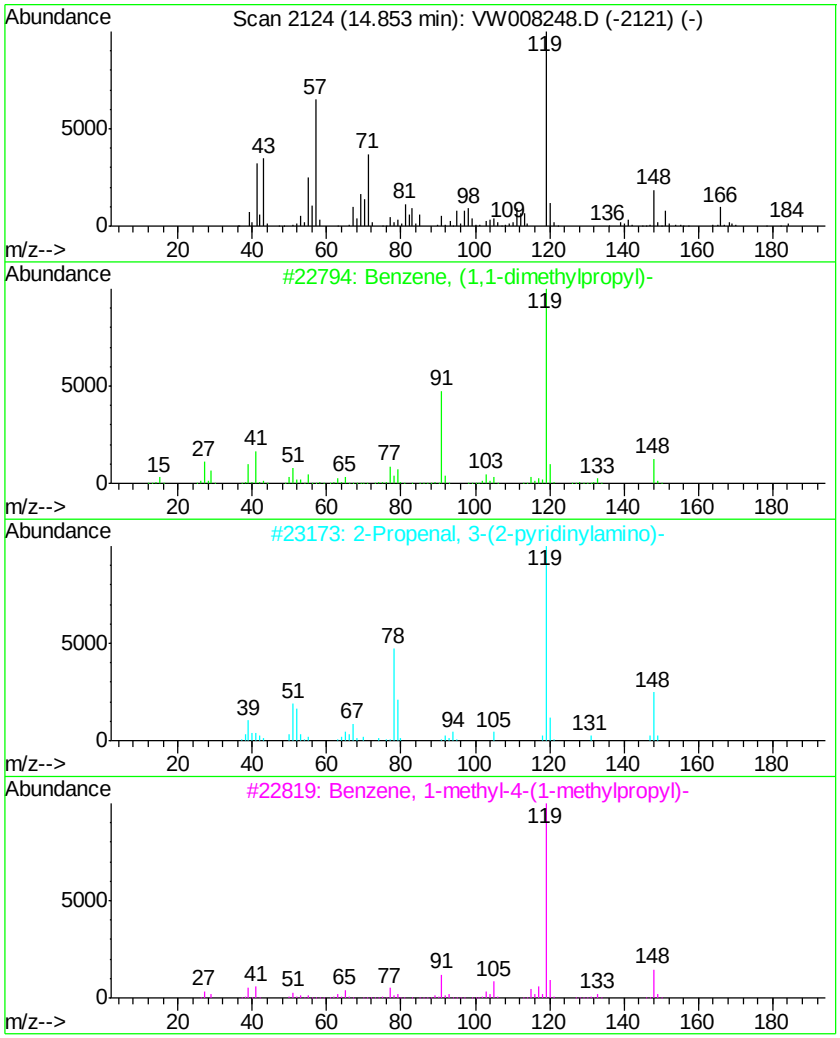
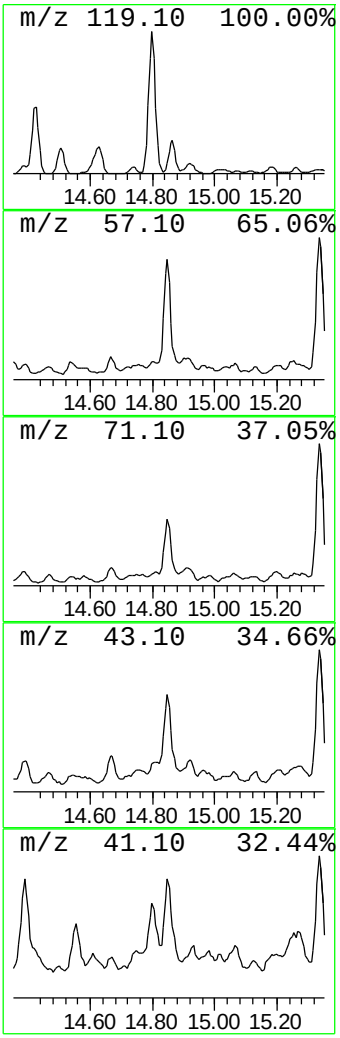
Peak Number 6 Benzene, (1,1-dimethylpropyl)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.85	80.48 ug/l	848120	1,4-Dichlorobenzene-d4	13.56

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	50
2	2-Propenal, 3-(2-pyridinylamino)-	148	C8H8N2O	068970-82-1	50
3	Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	50
4	Menthane, 3-chloro-8-phenyl-	250	C16H23Cl	184007-21-4	43
5	p-Toluic acid, 5-tridecyl ester	318	C21H34O2	1000299-80-5	38





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Data File : VW008248.D
Acq On : 14 Jan 2019 15:44
Operator : SY/VA
Sample : K1135-01
Misc : 5.00G/5ML/MSVOA_W/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_W
ClientSampleId : CPSB-19(10-15)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W010819S.M
Quant Title : SW846 8260

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

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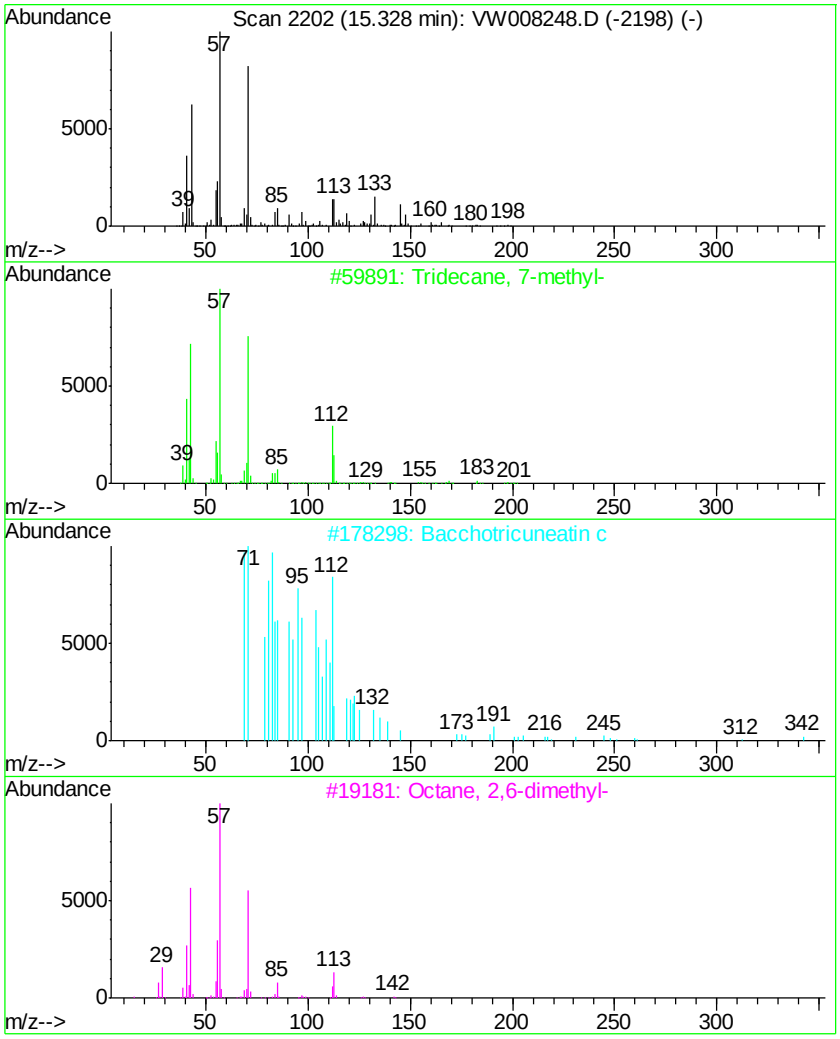
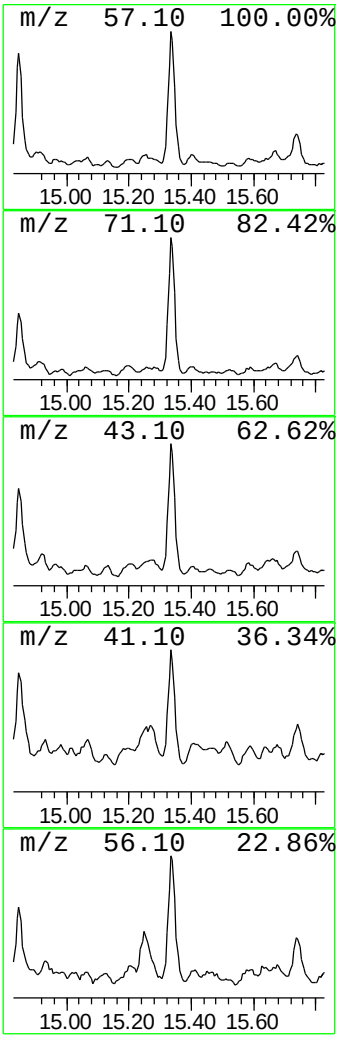
Peak Number 7 Tridecane, 7-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.33	100.53 ug/l	1059380	1,4-Dichlorobenzene-d4	13.56

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane, 7-methyl-	198	C14H30	026730-14-3	68
2	Bacchotricuneatin c	342	C20H22O5	066563-30-2	68
3	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	64
4	Decane, 2-methyl-	156	C11H24	006975-98-0	64
5	Nonane, 2,6-dimethyl-	156	C11H24	017302-28-2	64





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Data File : VW008248.D
Acq On : 14 Jan 2019 15:44
Operator : SY/VA
Sample : K1135-01
Misc : 5.00G/5ML/MSVOA W/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_W
ClientSampleId : CPSB-19(10-15)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W010819S.M
Quant Title : SW846 8260

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

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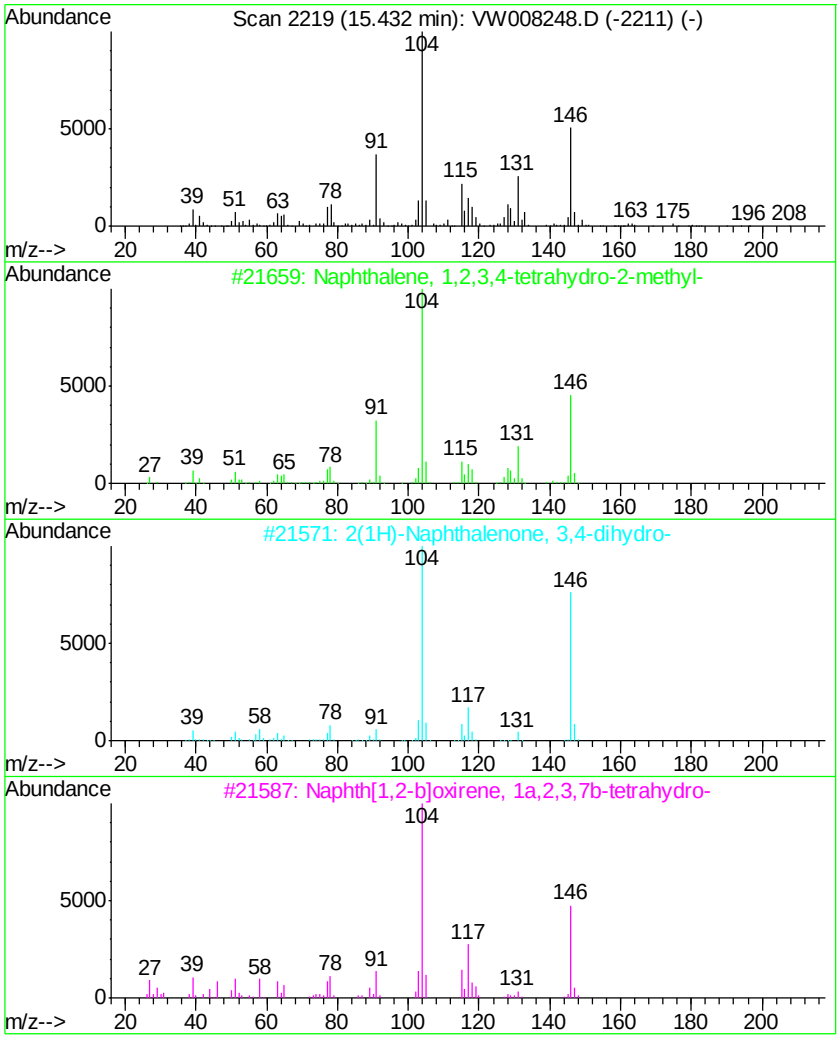
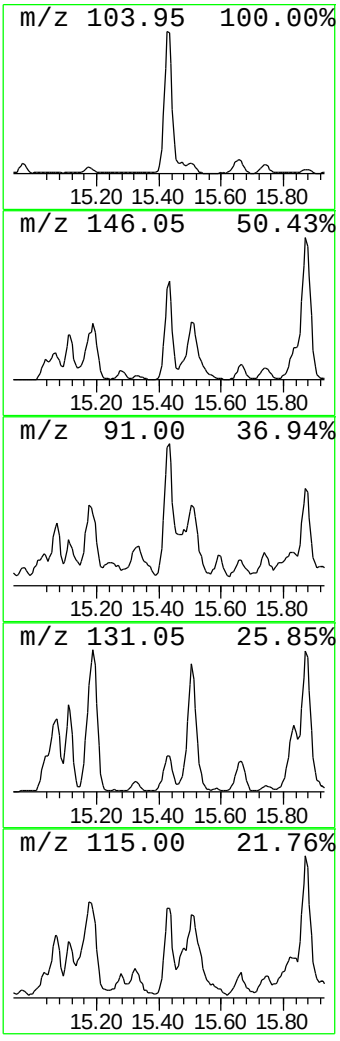
Peak Number 8 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.43	97.02 ug/l	1022450	1,4-Dichlorobenzene-d4	13.56

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	003877-19-8	94
2	2(1H)-Naphthalenone, 3,4-dihydro-	146	C10H10O	000530-93-8	62
3	Naphth[1,2-b]oxirene, 1a,2,3,7b-...	146	C10H10O	002461-34-9	62
4	Glutaric acid, heptyl 2-methylbe...	334	C20H30O4	1000371-33-0	38
5	2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	38





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_W\DATA\VW011419\
Data File : VW008248.D
Acq On : 14 Jan 2019 15:44
Operator : SY/VA
Sample : K1135-01
Misc : 5.00G/5ML/MSVOA_W/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_W
ClientSampleId : CPSB-19(10-15)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W010819S.M
Quant Title : SW846 8260

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

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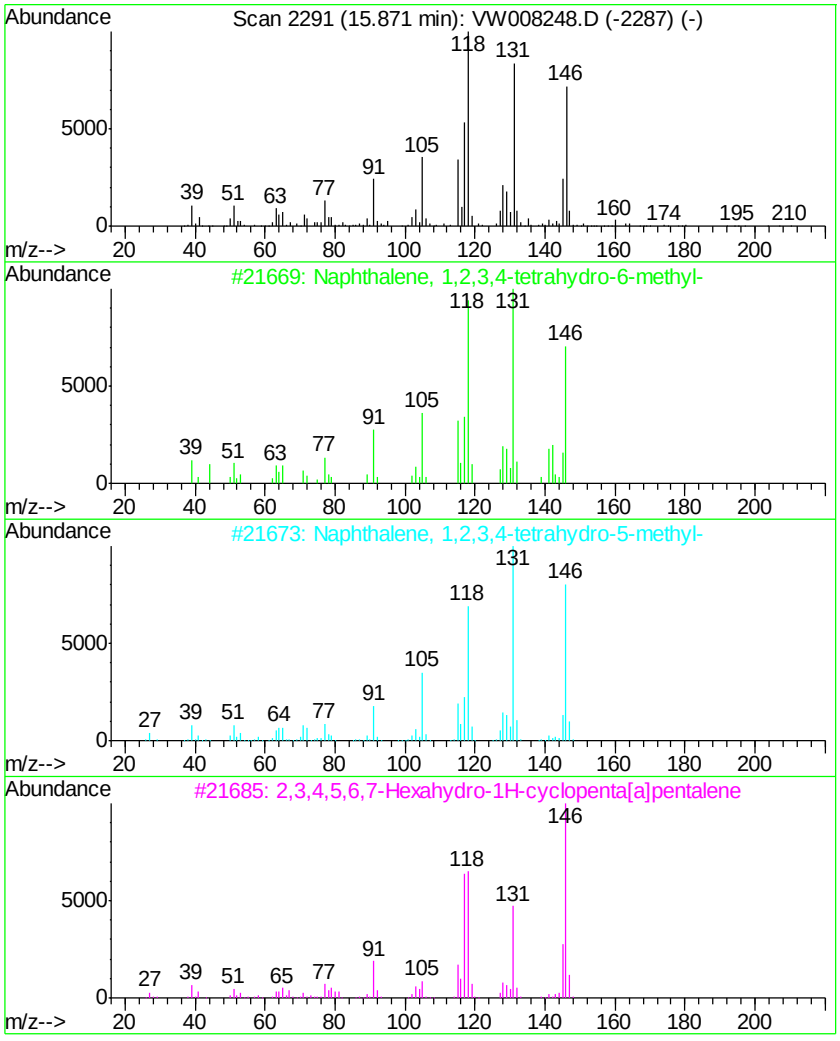
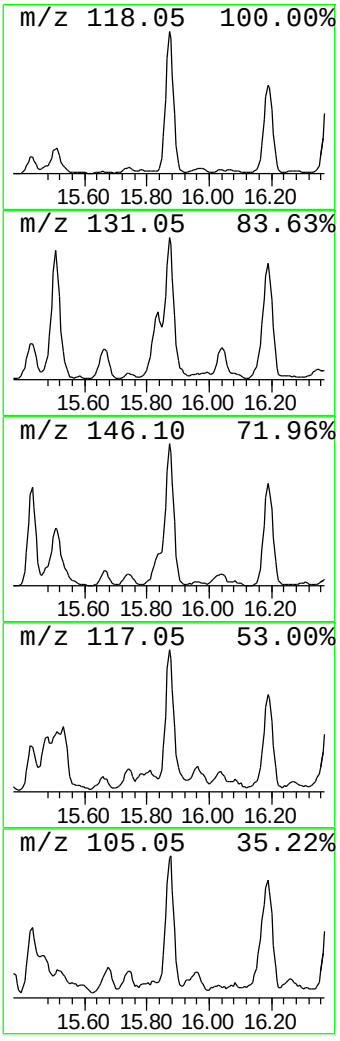
Peak Number 9 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.87	89.55 ug/l	943734	1,4-Dichlorobenzene-d4	13.56

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	93
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	70
3	2,3,4,5,6,7-Hexahydro-1H-cyclope...	146	C11H14	1000189-31-0	64
4	Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	49
5	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	45





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW011419\
Data File : VW008248.D
Acq On : 14 Jan 2019 15:44
Operator : SY/VA
Sample : K1135-01
Misc : 5.00G/5ML/MSVOA W/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_W
ClientSampleId : CPSB-19(10-15)

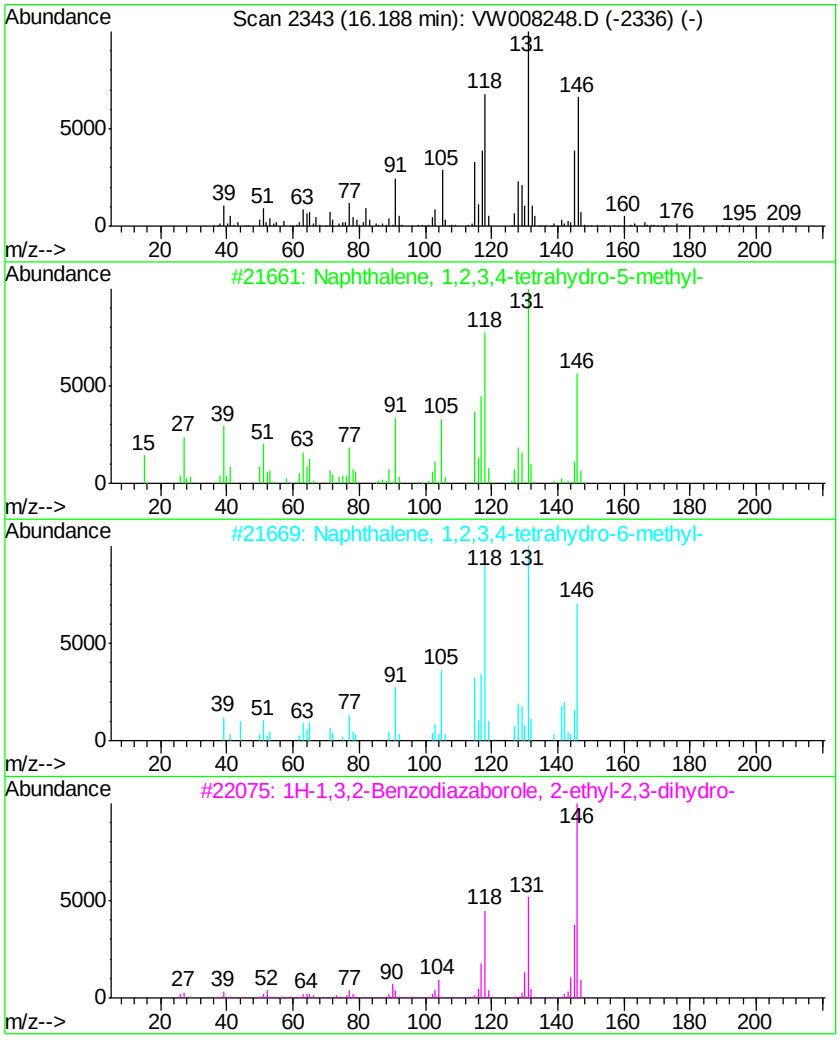
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W010819S.M
Quant Title : SW846 8260

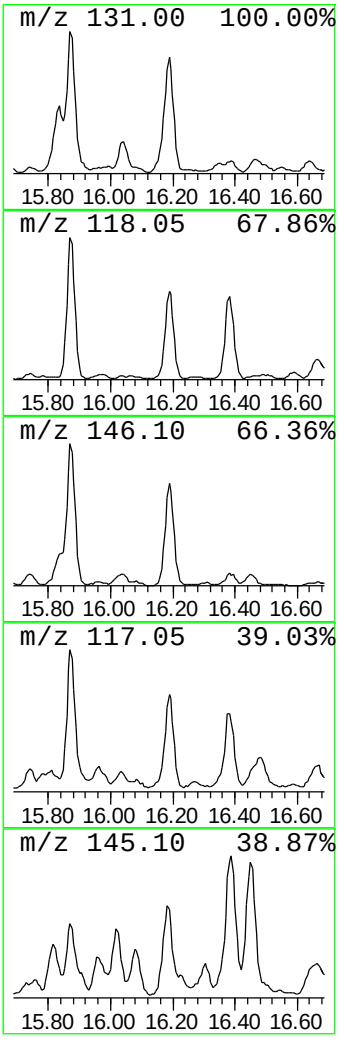
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.19	109.37 ug/l	1152570	1,4-Dichlorobenzene-d4	13.56

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	93
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	68
3	1H-1,3,2-Benzodiazaborole, 2-eth...	146	C8H11BN2	074663-81-3	62
4	1H-Benzimidazole, 5,6-dimethyl-	146	C9H10N2	000582-60-5	60
5	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	55





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_W\DATA\VW011419\
Data File : VW008248.D
Acq On : 14 Jan 2019 15:44
Operator : SY/VA
Sample : K1135-01
Misc : 5.00G/5ML/MSVOA_W/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_W
ClientSampleId : CPSB-19(10-15)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W010819S.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
Naphthalene, deca...	13.88	110.0	ug/l	1158910	4	13.56	526915	50.0
Benzene, 4-ethyl-...	14.10	98.2	ug/l	1034590	4	13.56	526915	50.0
Indan, 1-methyl-	14.21	79.9	ug/l	842015	4	13.56	526915	50.0
Cyclopentene, 1-p...	14.39	86.7	ug/l	914003	4	13.56	526915	50.0
Benzene, 1,2,4,5-...	14.80	208.7	ug/l	2199380	4	13.56	526915	50.0
Benzene, (1,1-dim...	14.85	80.5	ug/l	848120	4	13.56	526915	50.0
Tridecane, 7-methyl-	15.33	100.5	ug/l	1059380	4	13.56	526915	50.0
Naphthalene, 1,2,...	15.43	97.0	ug/l	1022450	4	13.56	526915	50.0
Naphthalene, 1,2,...	15.87	89.5	ug/l	943734	4	13.56	526915	50.0
Naphthalene, 1,2,...	16.19	109.4	ug/l	1152570	4	13.56	526915	50.0