

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091418\
 Data File : BM016754.D
 Acq On : 14 Sep 2018 15:35
 Operator : SJ/JU
 Sample : J4860-02
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A41T2

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0.1 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.022	3	6	23	rVB	4965991	5857276	68.77%	6.283%
2	3.269	44	48	58	rVB	84985	110801	1.30%	0.119%
3	3.540	91	94	97	rBV2	12003	14950	0.18%	0.016%
4	3.899	149	155	159	rBV	35890	53668	0.63%	0.058%
5	3.987	167	170	176	rVB	25294	32411	0.38%	0.035%
6	4.199	201	206	210	rBV3	13105	16893	0.20%	0.018%
7	4.599	267	274	281	rBV5	8750	15424	0.18%	0.017%
8	5.840	481	485	492	rBV	24411	37741	0.44%	0.040%
9	6.404	575	581	588	rVB8	6930	18120	0.21%	0.019%
10	6.899	658	665	678	rBV2	1352940	2322702	27.27%	2.491%
11	7.075	686	695	702	rBV	1017503	1574071	18.48%	1.688%
12	7.263	721	727	740	rVB	1344169	2090281	24.54%	2.242%
13	7.734	799	807	814	rVV	832393	1326570	15.58%	1.423%
14	7.798	814	818	825	rVB3	19763	33061	0.39%	0.035%
15	8.428	915	925	933	rBV	1521852	2446848	28.73%	2.625%
16	8.881	995	1002	1013	rVB	1205442	1885734	22.14%	2.023%
17	9.057	1024	1032	1033	rBV5	9517	16808	0.20%	0.018%
18	9.316	1070	1076	1081	rVB3	17062	28578	0.34%	0.031%
19	9.440	1091	1097	1104	rVB8	7997	15605	0.18%	0.017%
20	9.598	1113	1124	1131	rBV	815093	1321869	15.52%	1.418%
21	9.698	1136	1141	1145	rVV5	10496	15624	0.18%	0.017%
22	10.128	1206	1214	1223	rBV	1521190	2482138	29.14%	2.663%
23	10.510	1272	1279	1286	rBV	1141252	1896317	22.27%	2.034%
24	10.557	1286	1287	1294	rVB	16312	18946	0.22%	0.020%
25	10.645	1294	1302	1316	rBV	1455901	2462523	28.91%	2.641%
26	11.004	1360	1363	1368	rBV4	7864	11777	0.14%	0.013%
27	11.551	1451	1456	1462	rBV	28393	46145	0.54%	0.049%
28	11.810	1493	1500	1502	rBV7	7189	11700	0.14%	0.013%
29	12.104	1545	1550	1555	rVV2	28891	53546	0.63%	0.057%
30	12.175	1555	1562	1567	rVB8	17243	40535	0.48%	0.043%
31	12.257	1570	1576	1578	rBV5	7977	14428	0.17%	0.015%
32	12.398	1595	1600	1604	rBV5	12968	25337	0.30%	0.027%
33	12.651	1637	1643	1650	rBV8	10022	23094	0.27%	0.025%
34	12.728	1652	1656	1661	rVV5	10466	16238	0.19%	0.017%

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Integration Parameters: LSCINT.P

Integrator: RTE
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 Title : SVOA CALIBRATION

35	12.804	1667	1669	1675	rVB7	7760	14014	0.16%	0.015%
36	12.869	1675	1680	1684	rBV6	9027	16994	0.20%	0.018%
37	12.916	1684	1688	1692	rVV	43736	60076	0.71%	0.064%
38	12.975	1693	1698	1703	rVB8	12107	27277	0.32%	0.029%
39	13.181	1730	1733	1735	rVV4	8249	11768	0.14%	0.013%
40	13.439	1772	1777	1782	rVB8	6358	12996	0.15%	0.014%
41	13.692	1816	1820	1824	rBV5	12039	18780	0.22%	0.020%
42	13.786	1830	1836	1840	rBV	2558171	3460095	40.63%	3.712%
43	13.833	1840	1844	1848	rVV	486443	690828	8.11%	0.741%
44	13.869	1848	1850	1855	rVV3	54196	66747	0.78%	0.072%
45	13.916	1856	1858	1864	rVB5	16952	25400	0.30%	0.027%
46	14.063	1873	1883	1896	rBV	2973933	4502353	52.86%	4.830%
47	14.204	1904	1907	1915	rVB8	28901	54478	0.64%	0.058%
48	14.280	1915	1920	1924	rBV7	15543	29925	0.35%	0.032%
49	14.369	1927	1935	1942	rBV2	1543337	2418670	28.40%	2.594%
50	14.427	1942	1945	1951	rVB	45921	67660	0.79%	0.073%
51	14.498	1951	1957	1962	rVB	316597	452079	5.31%	0.485%
52	14.563	1962	1968	1976	rBV	1047610	1465886	17.21%	1.572%
53	14.698	1987	1991	1993	rBV4	10565	14129	0.17%	0.015%
54	14.769	1999	2003	2004	rBV2	21909	27784	0.33%	0.030%
55	14.792	2005	2007	2011	rVB5	34387	41985	0.49%	0.045%
56	14.910	2024	2027	2035	rVB8	27935	42221	0.50%	0.045%
57	15.310	2091	2095	2098	rBV2	51574	67373	0.79%	0.072%
58	15.369	2098	2105	2110	rVV	3798996	5531068	64.94%	5.933%
59	15.422	2110	2114	2118	rVB2	51699	70733	0.83%	0.076%
60	15.480	2118	2124	2132	rBV	771910	1097928	12.89%	1.178%
61	15.633	2140	2150	2159	rBV	906897	1380346	16.21%	1.481%
62	15.857	2185	2188	2196	rVB4	26337	51820	0.61%	0.056%
63	16.104	2227	2230	2231	rBV2	23464	27689	0.33%	0.030%
64	16.133	2231	2235	2242	rVB	99042	155325	1.82%	0.167%
65	16.239	2248	2253	2257	rBV	161083	223455	2.62%	0.240%
66	16.357	2267	2273	2278	rBV	797112	1055725	12.40%	1.132%
67	16.474	2290	2293	2297	rBV4	27819	41763	0.49%	0.045%
68	16.651	2319	2323	2328	rBV	158050	230324	2.70%	0.247%
69	16.951	2369	2374	2378	rVB3	63109	112697	1.32%	0.121%
70	17.004	2379	2383	2385	rVV2	119485	144526	1.70%	0.155%
71	17.039	2385	2389	2393	rVB	241945	322173	3.78%	0.346%

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72	17.116	2396	2402	2406	rBV	2011759	3020581	35.47%	3.240%
73	17.157	2406	2409	2413	rVV	612965	838987	9.85%	0.900%
74	17.216	2413	2419	2429	rVV	4297897	6086730	71.47%	6.529%
75	17.298	2429	2433	2442	rVB2	222340	322670	3.79%	0.346%
76	17.574	2475	2480	2483	rBV2	115661	179487	2.11%	0.193%
77	17.604	2483	2485	2489	rVB2	81632	79601	0.93%	0.085%
78	17.715	2499	2504	2512	rVB2	224811	462522	5.43%	0.496%
79	17.839	2522	2525	2532	rVB5	64678	92024	1.08%	0.099%
80	17.916	2536	2538	2541	rVB4	38272	41574	0.49%	0.045%
81	18.027	2553	2557	2563	rVB2	259307	398784	4.68%	0.428%
82	18.192	2580	2585	2588	rBV5	156108	251664	2.95%	0.270%
83	18.251	2593	2595	2599	rVB	103789	105838	1.24%	0.114%
84	18.298	2599	2603	2606	rBV3	96313	125277	1.47%	0.134%
85	18.910	2704	2707	2713	rVB3	75784	107149	1.26%	0.115%
86	19.174	2748	2752	2759	rVB	961702	1262280	14.82%	1.354%
87	19.315	2773	2776	2781	rVB4	95762	140646	1.65%	0.151%
88	19.510	2804	2809	2821	rBV3	5391555	8517023	100.00%	9.136%
89	20.039	2896	2899	2904	rVB	135120	180231	2.12%	0.193%
90	20.139	2913	2916	2920	rBV	88450	108716	1.28%	0.117%
91	20.921	3046	3049	3053	rBV4	274465	401642	4.72%	0.431%
92	21.039	3066	3069	3080	rVB6	342805	653431	7.67%	0.701%
93	21.180	3087	3093	3097	rBV3	534167	993240	11.66%	1.065%
94	21.221	3097	3100	3107	rVV2	773447	1364324	16.02%	1.463%
95	21.304	3107	3114	3118	rVV	2987178	4460180	52.37%	4.784%
96	21.339	3118	3120	3126	rVB	402422	488451	5.73%	0.524%
97	22.521	3318	3321	3327	rVB	322146	427388	5.02%	0.458%
98	22.874	3377	3381	3383	rBV2	229148	378937	4.45%	0.406%
99	23.403	3465	3471	3478	rBV	4225223	7803411	91.62%	8.370%
100	23.545	3489	3495	3502	rVB	1946652	3559574	41.79%	3.818%

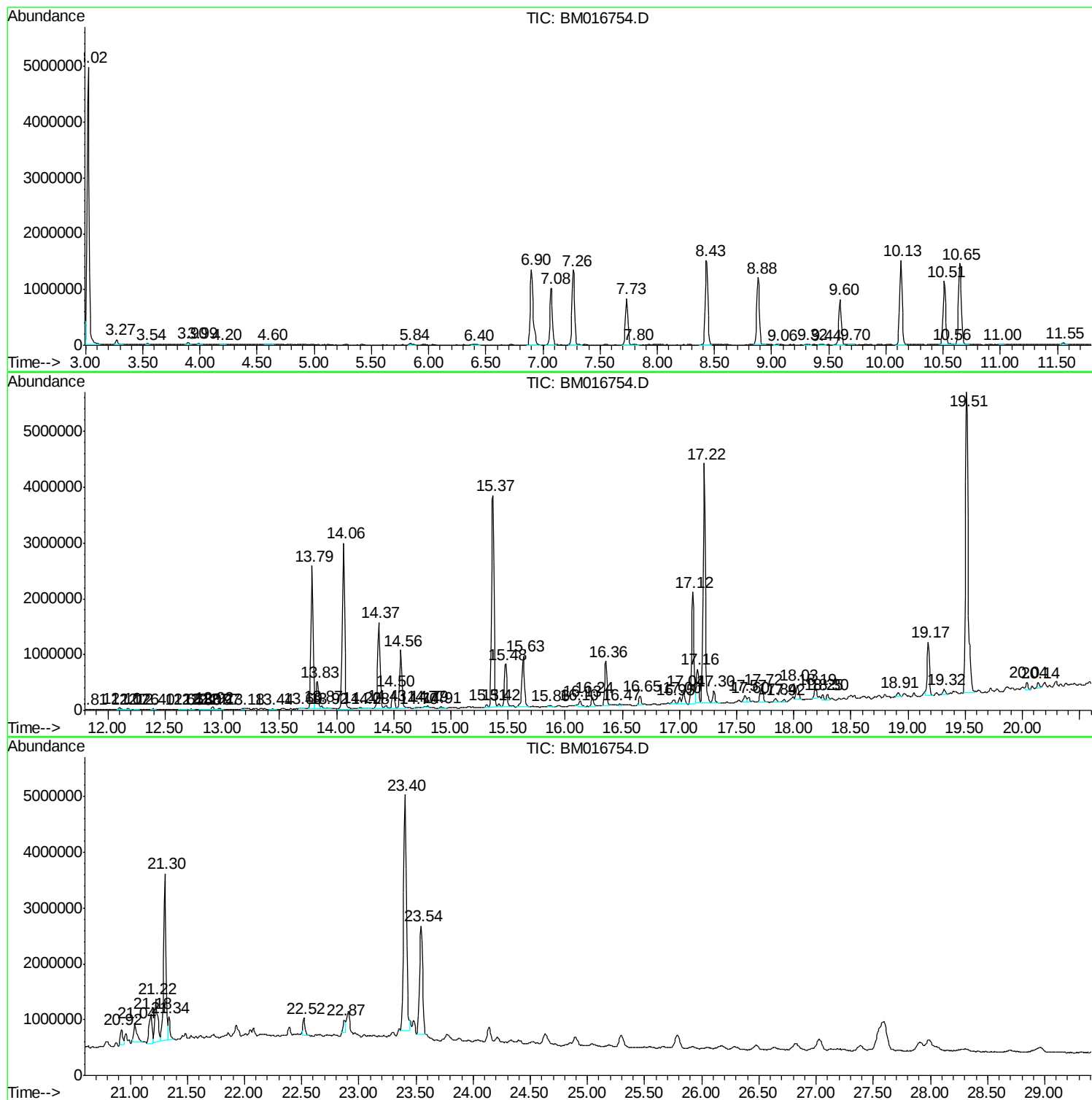
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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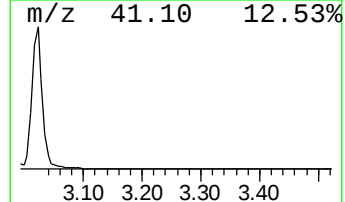
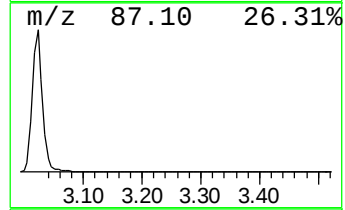
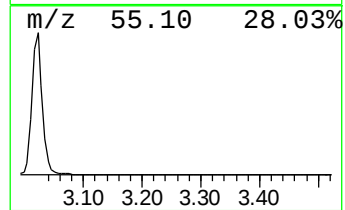
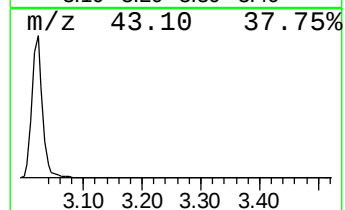
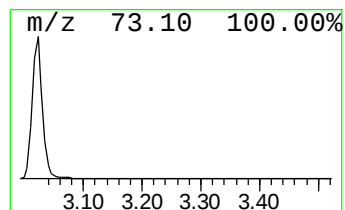
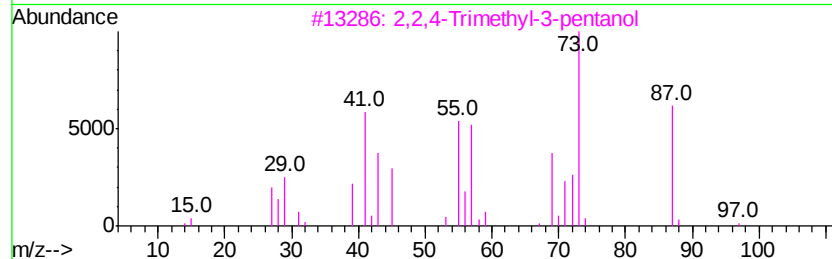
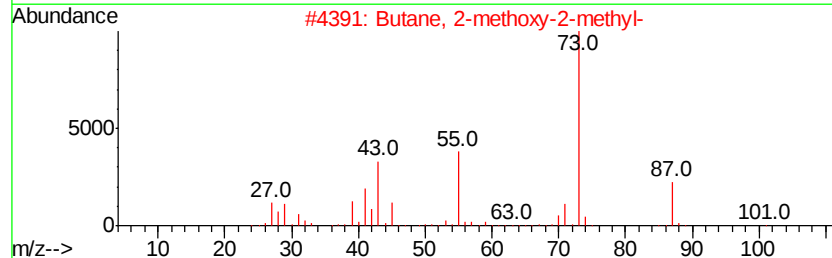
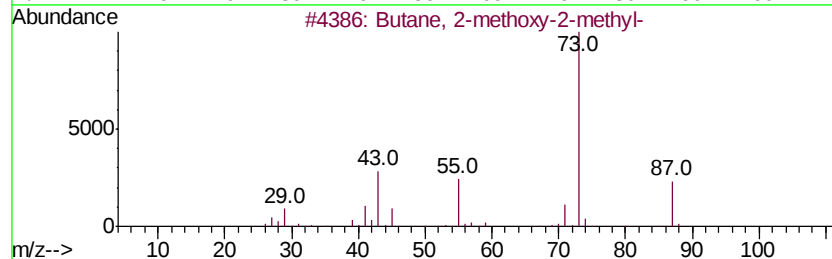
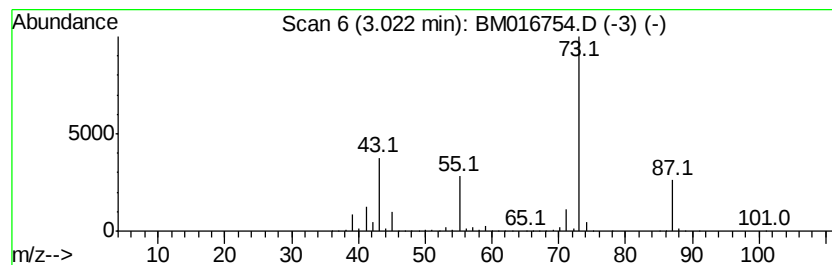
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.02	88.31 ng/ul	5857280	1,4-Dichlorobenzene-d4	7.73

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	74
3		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
4		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
5		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	22



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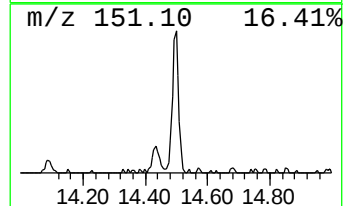
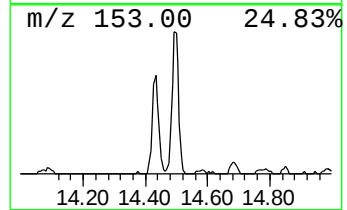
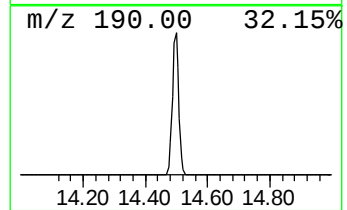
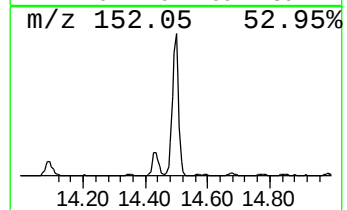
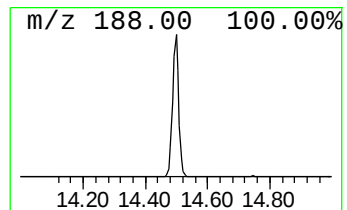
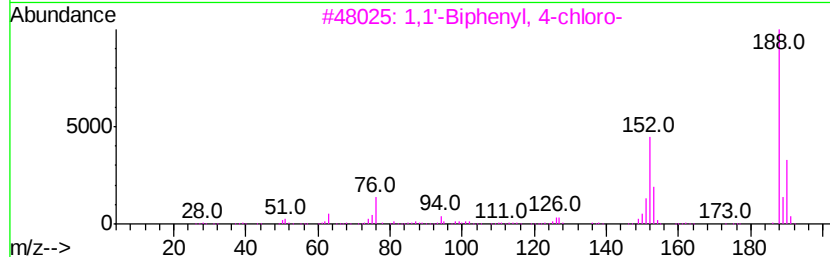
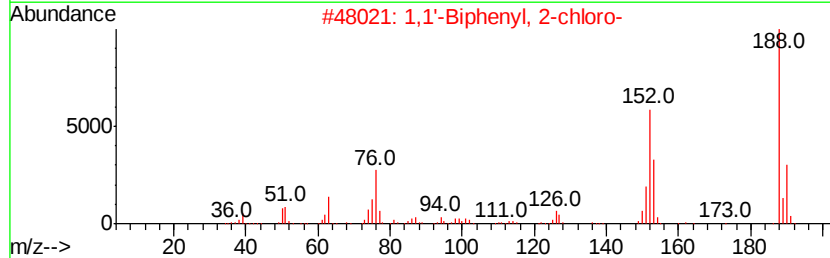
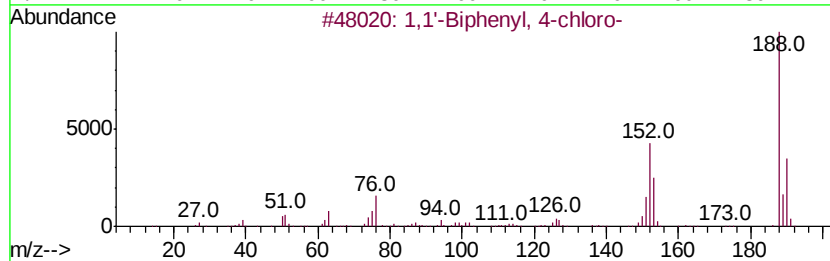
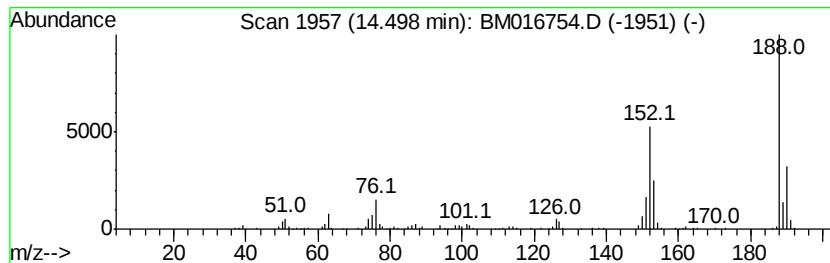
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 1,1'-Biphenyl, 4-chloro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.
14.50	3.74 ng/ul	452079	Acenaphthene-d10		14.37
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 4-chloro-	188	C12H9Cl	002051-62-9	98
2	1,1'-Biphenyl, 2-chloro-	188	C12H9Cl	002051-60-7	97
3	1,1'-Biphenyl, 4-chloro-	188	C12H9Cl	002051-62-9	95
4	1,1'-Biphenyl, 4-chloro-	188	C12H9Cl	002051-62-9	95
5	1,1'-Biphenyl, 2-chloro-	188	C12H9Cl	002051-60-7	94



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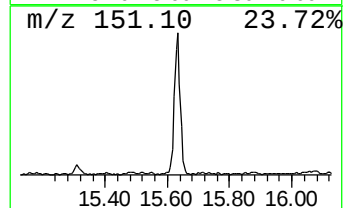
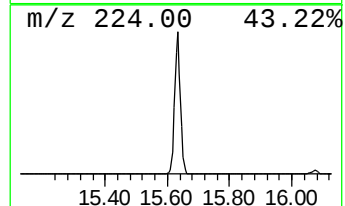
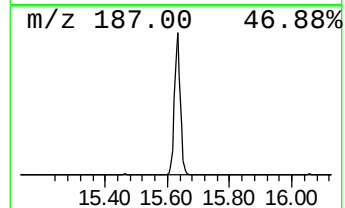
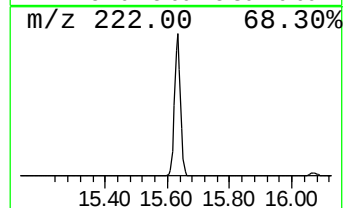
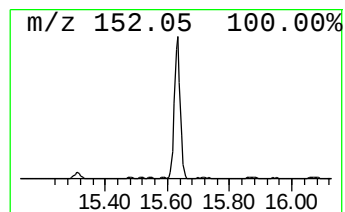
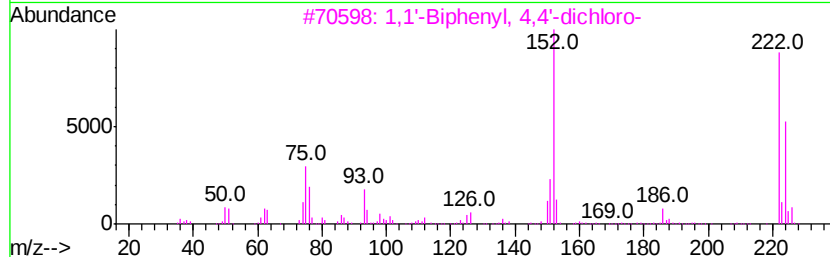
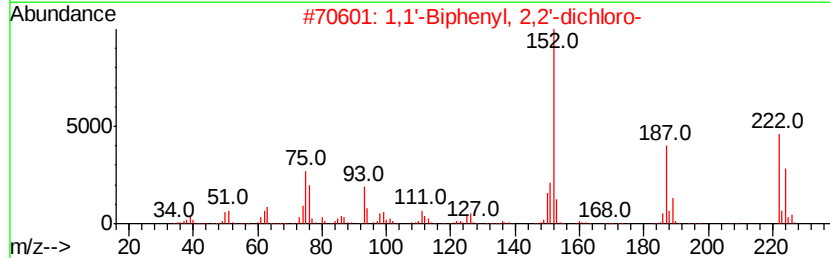
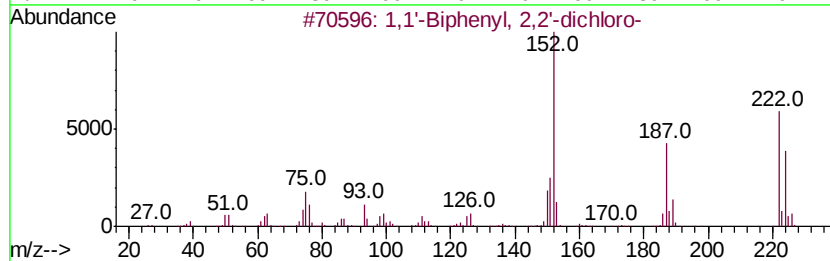
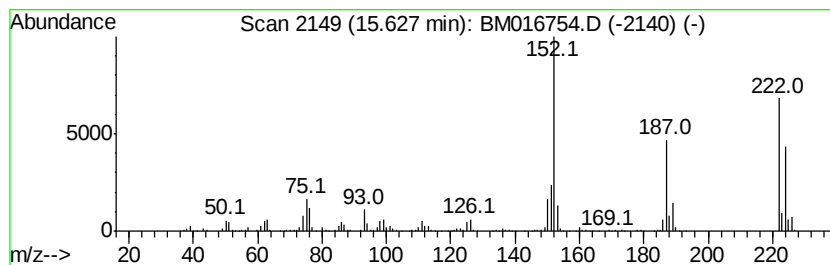
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 1,1'-Biphenyl, 2,2'-dichloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.63	11.41 ng/ul	1380350	Acenaphthene-d10	14.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,2'-dichloro-	222	C12H8Cl2	013029-08-8	99
2		1,1'-Biphenyl, 2,2'-dichloro-	222	C12H8Cl2	013029-08-8	99
3		1,1'-Biphenyl, 4,4'-dichloro-	222	C12H8Cl2	002050-68-2	98
4		1,1'-Biphenyl, 2,3-dichloro-	222	C12H8Cl2	016605-91-7	95
5		1,1'-Biphenyl, 4,4'-dichloro-	222	C12H8Cl2	002050-68-2	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091418\
Data File : BM016754.D
Acq On : 14 Sep 2018 15:35
Operator : SJ/JU
Sample : J4860-02
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A41T2

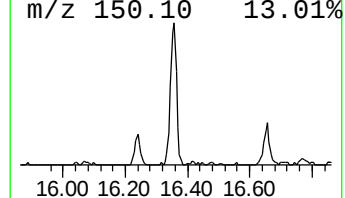
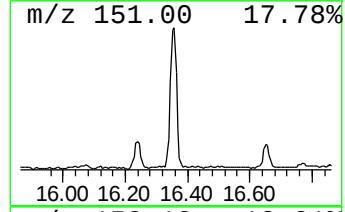
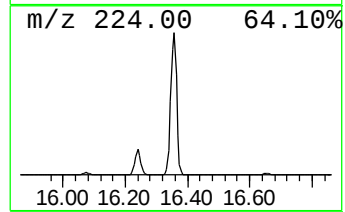
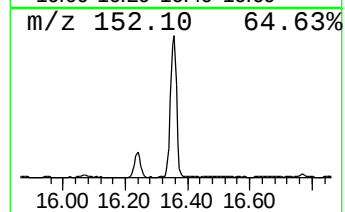
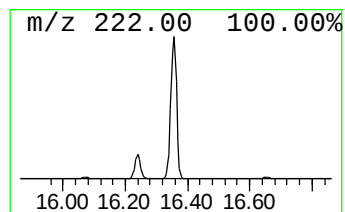
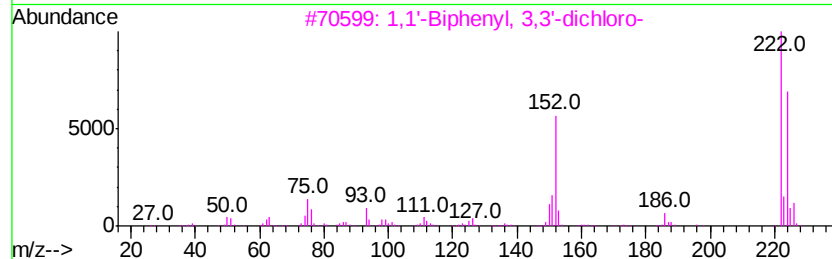
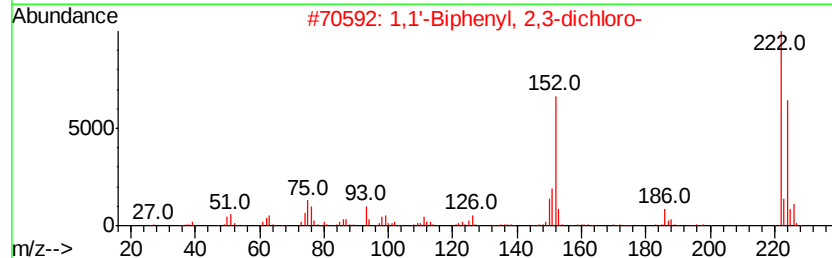
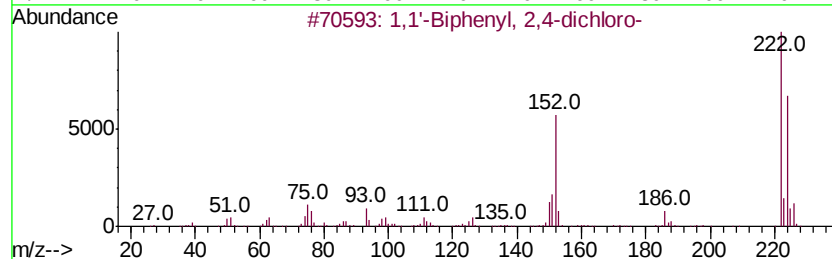
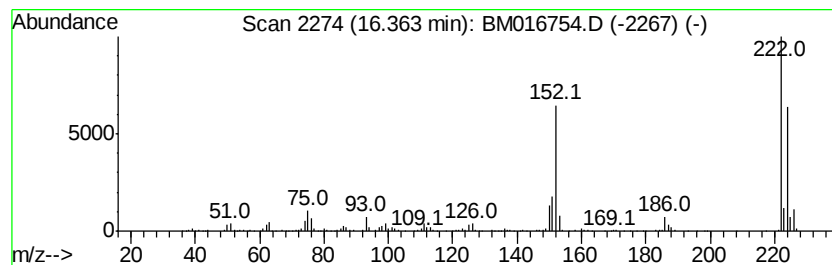
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 1,1'-Biphenyl, 2,4-dichloro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.36	6.99 ng/ul	1055730	Phenanthrene-d10	17.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,4-dichloro-	222	C12H8Cl2	033284-50-3	99
2		1,1'-Biphenyl, 2,3-dichloro-	222	C12H8Cl2	016605-91-7	99
3		1,1'-Biphenyl, 3,3'-dichloro-	222	C12H8Cl2	002050-67-1	99
4		1,1'-Biphenyl, 4,4'-dichloro-	222	C12H8Cl2	002050-68-2	98
5		1,1'-Biphenyl, 2,4-dichloro-	222	C12H8Cl2	033284-50-3	98



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091418\
Data File : BM016754.D
Acq On : 14 Sep 2018 15:35
Operator : SJ/JU
Sample : J4860-02
Misc :
ALS Vial : 11 Sample Multiplier: 1

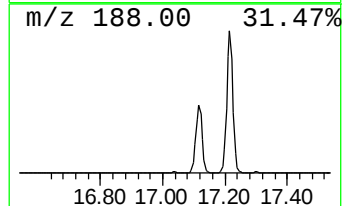
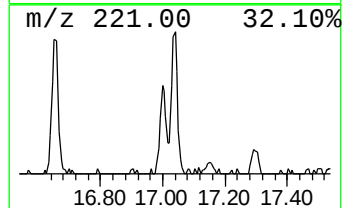
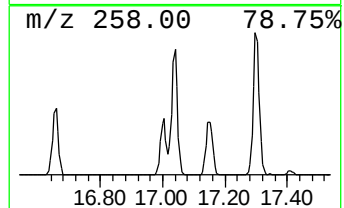
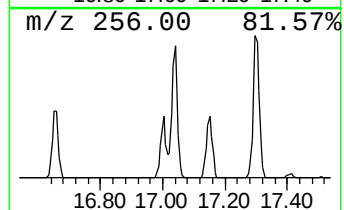
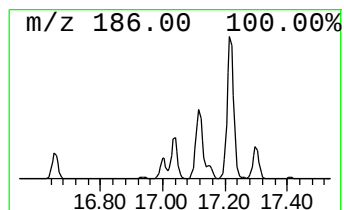
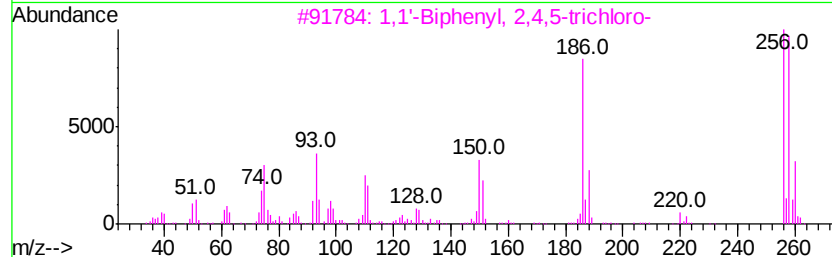
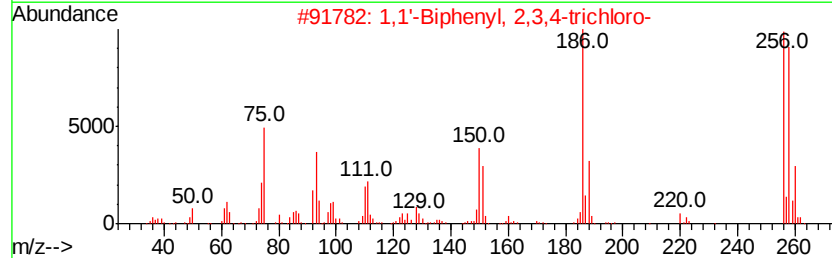
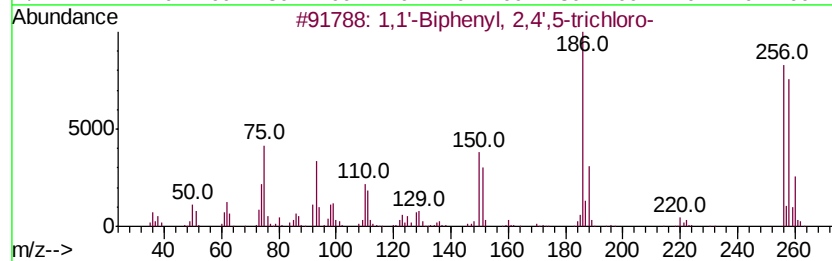
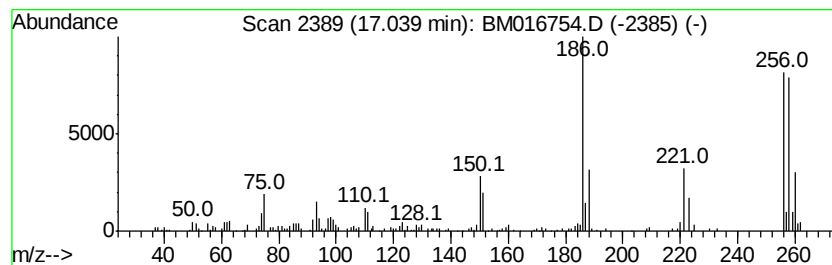
Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 1,1'-Biphenyl, 2,4',5-trich... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.		
17.04	2.13 ng/ul	322173	Phenanthrene-d10		17.12		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'	-Biphenyl,	2,4',5-trichloro-	256	C12H7Cl3	016606-02-3	98
2	1,1'	-Biphenyl,	2,3,4-trichloro-	256	C12H7Cl3	055702-46-0	98
3	1,1'	-Biphenyl,	2,4,5-trichloro-	256	C12H7Cl3	015862-07-4	98
4	1,1'	-Biphenyl,	3,4,4'-Trichloro-	256	C12H7Cl3	038444-90-5	97
5	1,1'	-Biphenyl,	2,4',5-trichloro-	256	C12H7Cl3	016606-02-3	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091418\
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Operator : SJ/JU
Sample : J4860-02
Misc :
ALS Vial : 11 Sample Multiplier: 1

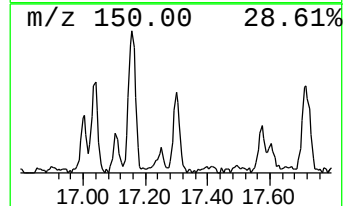
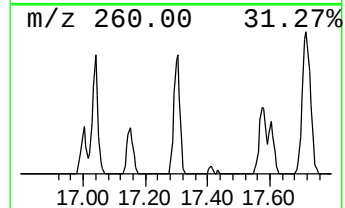
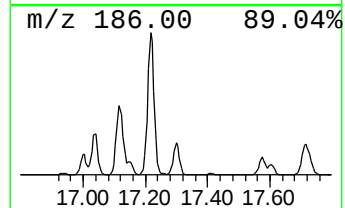
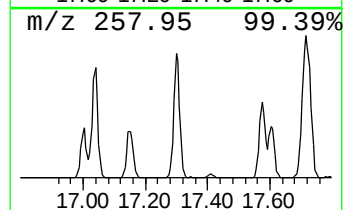
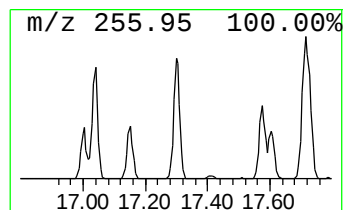
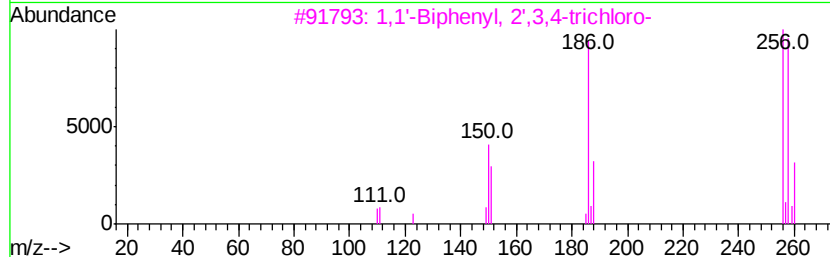
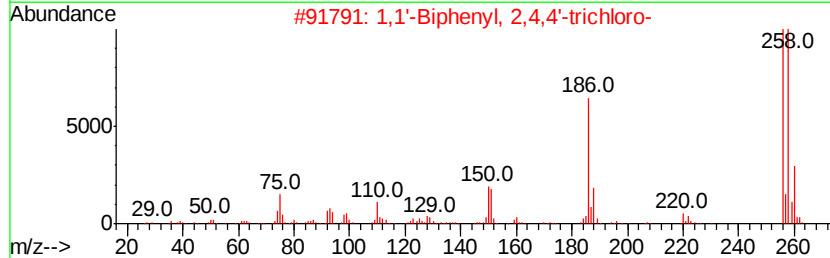
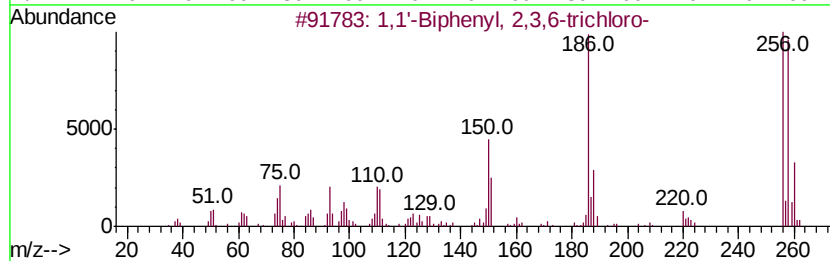
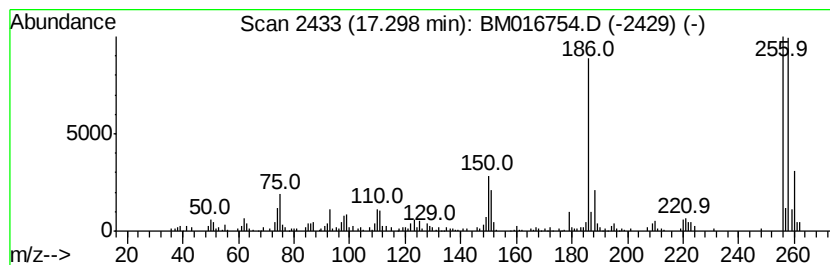
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 1,1'-Biphenyl, 2,3,6-trichl... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.30	2.14 ng/ul	322670	Phenanthrene-d10	17.12	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3,6-trichloro-	256	C12H7Cl3	055702-45-9	99
2	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	98
3	1,1'-Biphenyl, 2',3,4-trichloro-	256	C12H7Cl3	038444-86-9	97
4	1,1'-Biphenyl, 2,4,5-trichloro-	256	C12H7Cl3	015862-07-4	96
5	1,1'-Biphenyl, 2,3,4-trichloro-	256	C12H7Cl3	055702-46-0	96



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091418\
Data File : BM016754.D
Acq On : 14 Sep 2018 15:35
Operator : SJ/JU
Sample : J4860-02
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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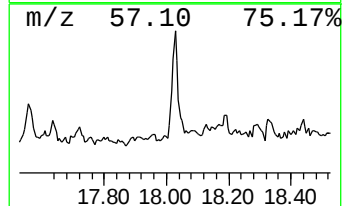
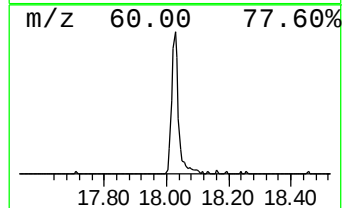
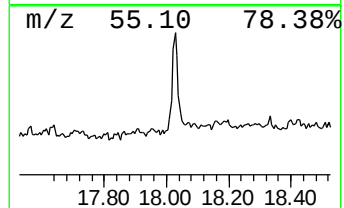
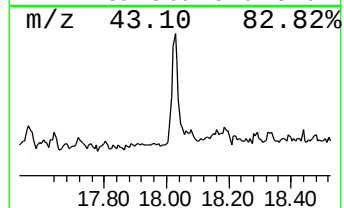
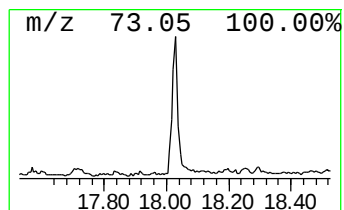
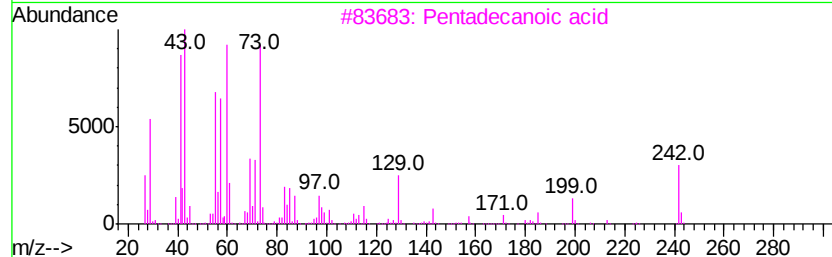
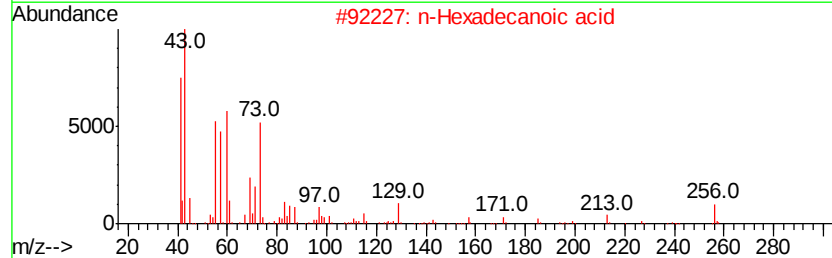
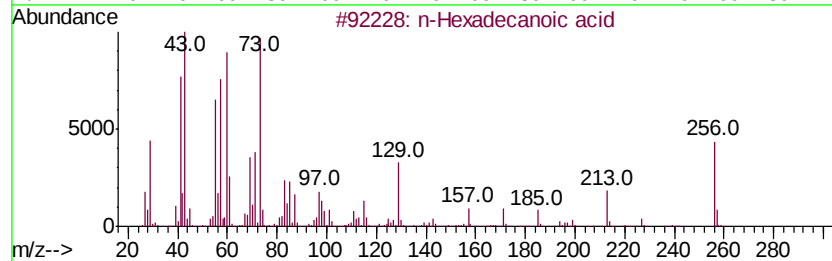
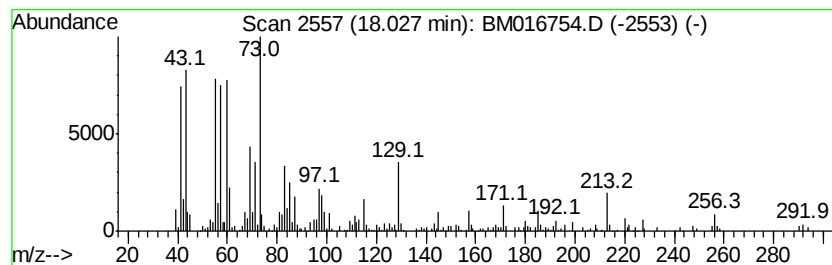
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 n-Hexadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.03	2.64 ng/ul	398784	Phenanthrene-d10	17.12

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98	
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91	
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	90	
4	Tridecanoic acid	214	C13H26O2	000638-53-9	90	
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	86	



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091418\
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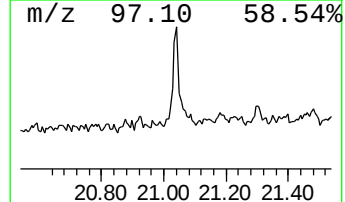
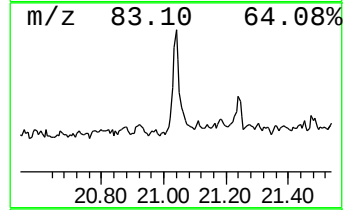
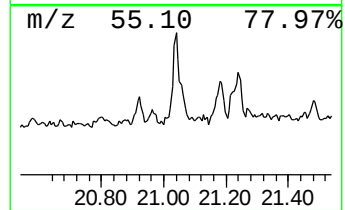
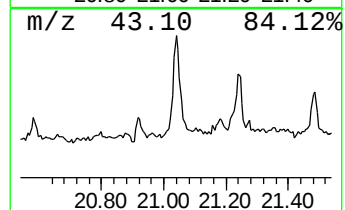
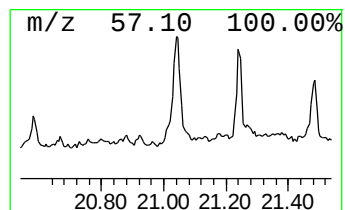
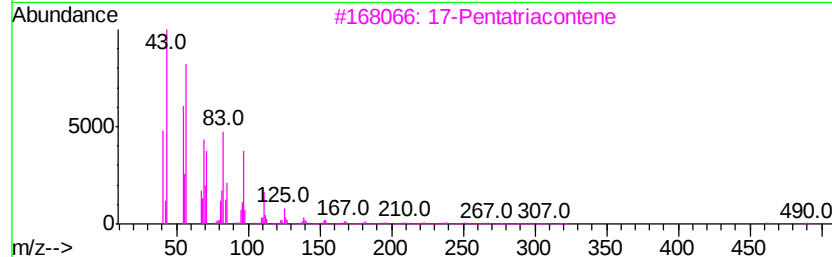
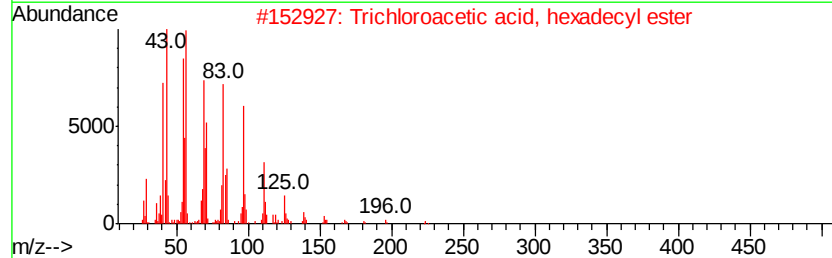
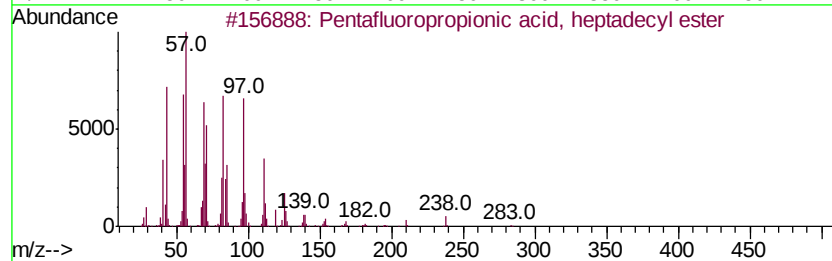
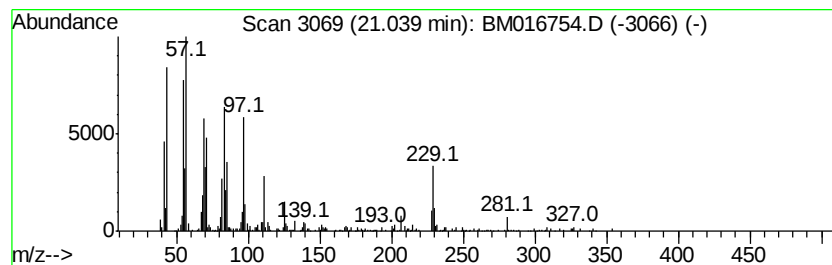
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Pentafluoropropionic acid, ... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.		
21.04	2.93 ng/ul	653431	Chrysene-d12	21.30		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentafluoropropionic acid, hepta...	402	C20H35F5O2	1000283-04-2	81
2		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	74
3		17-Pentatriacontene	491	C35H70	006971-40-0	74
4		1-Tricosene	322	C23H46	018835-32-0	74
5		Pentafluoropropionic acid, octad...	416	C21H37F5O2	1000280-07-7	74



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091418\
Data File : BM016754.D
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ClientSampleId :
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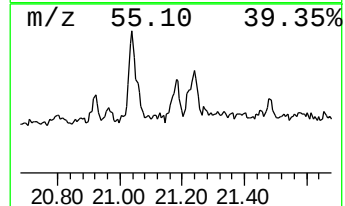
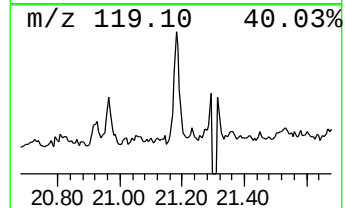
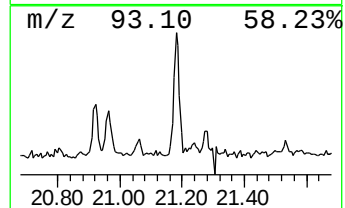
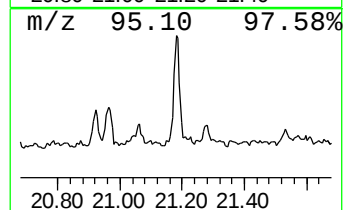
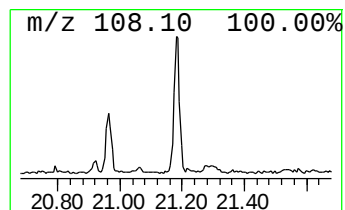
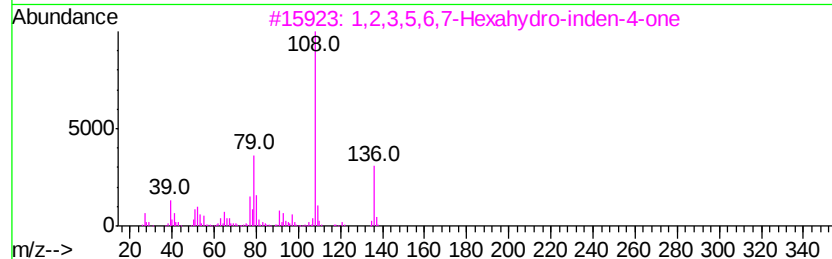
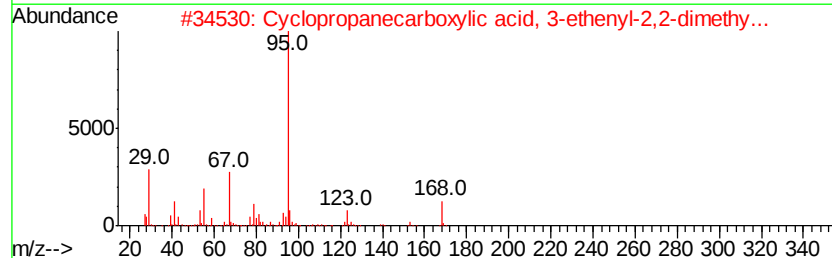
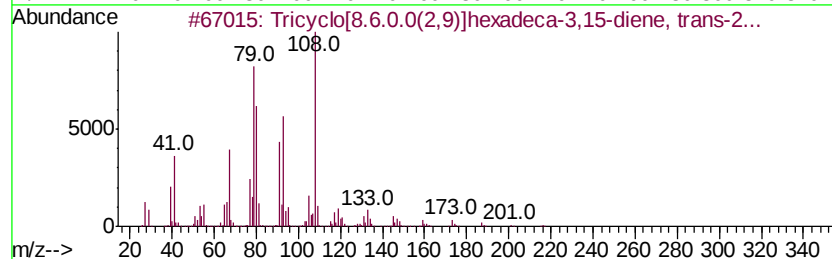
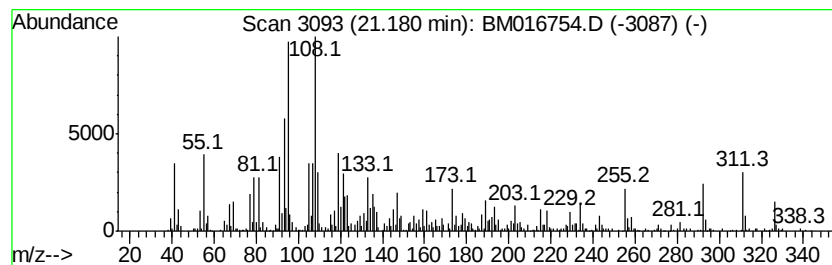
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 unknown-01 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.18	4.45 ng/ul	993240	Chrysene-d12	21.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tricyclo[8.6.0.0(2,9)]hexadeca-3...	216	C16H24	1000150-40-3	12
2		Cyclopropanecarboxylic acid, 3-e...	168	C10H16O2	060066-50-4	11
3		1,2,3,5,6,7-Hexahydro-inden-4-one	136	C9H12O	022118-01-0	10
4		Cycloheptane, 1,3-dichloro-, trans-	166	C7H12Cl2	032616-84-5	10
5		Borneol	154	C10H18O	010385-78-1	10



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091418\
Data File : BM016754.D
Acq On : 14 Sep 2018 15:35
Operator : SJ/JU
Sample : J4860-02
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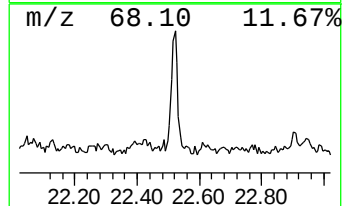
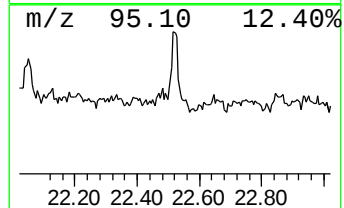
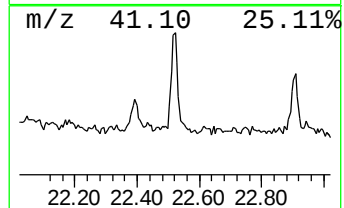
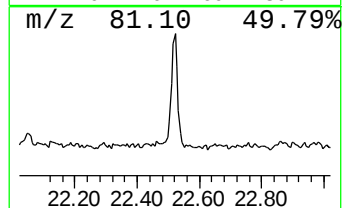
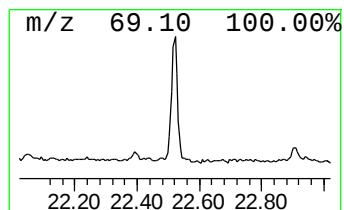
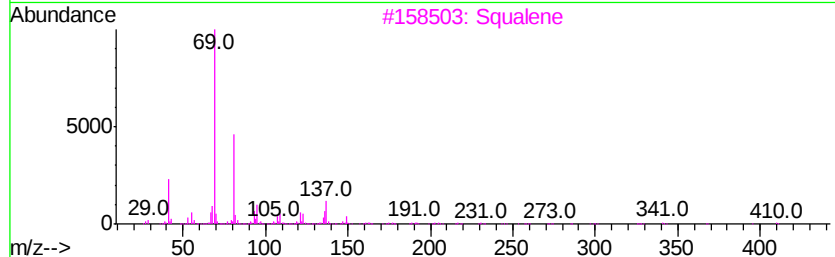
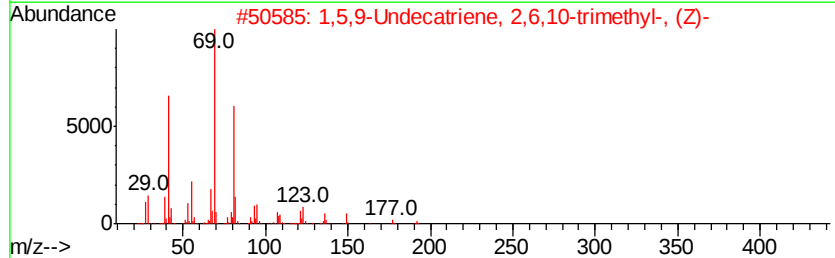
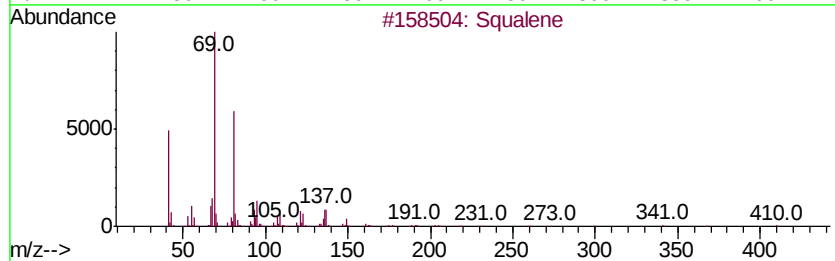
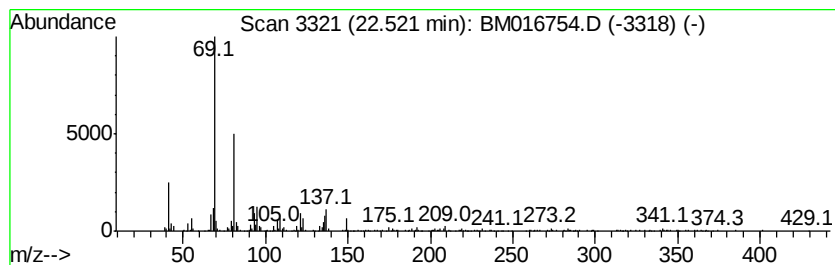
Instrument :
BNA_M
ClientSampleId :
A41T2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Squalene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.52	2.40 ng/ul	427388	Perylene-d12	23.54
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Squalene		410 C30H50	007683-64-9 87
2	1,5,9-Undecatriene, 2,6,10-trime...		192 C14H24	062951-96-6 81
3	Squalene		410 C30H50	007683-64-9 74
4	4,8,12-Tetradecatrienal, 5,9,13-...		248 C17H28O	066408-55-7 72
5	2,6,10,14-Hexadecatetraenoic aci...		332 C22H36O2	024035-35-6 72



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM091418\
Data File : BM016754.D
Acq On : 14 Sep 2018 15:35
Operator : SJ/JU
Sample : J4860-02
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A41T2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	3.02	88.3	ng/ul	5857280	1	7.73	1326570	20.0
1,1'-Biphenyl, 4-...	14.50	3.7	ng/ul	452079	3	14.37	2418670	20.0
1,1'-Biphenyl, 2,...	15.63	11.4	ng/ul	1380350	3	14.37	2418670	20.0
1,1'-Biphenyl, 2,...	16.36	7.0	ng/ul	1055730	4	17.12	3020580	20.0
1,1'-Biphenyl, 2,...	17.04	2.1	ng/ul	322173	4	17.12	3020580	20.0
1,1'-Biphenyl, 2,...	17.30	2.1	ng/ul	322670	4	17.12	3020580	20.0
n-Hexadecanoic acid	18.03	2.6	ng/ul	398784	4	17.12	3020580	20.0
Pentafluoropropio...	21.04	2.9	ng/ul	653431	5	21.30	4460180	20.0
unknown-01	21.18	4.5	ng/ul	993240	5	21.30	4460180	20.0
Squalene	22.52	2.4	ng/ul	427388	6	23.54	3559570	20.0