

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111324\
 Data File : BP023092.D
 Acq On : 13 Nov 2024 22:15
 Operator : RC/JU
 Sample : P4724-04
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BH9B9

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF Filtering: 5
 Sampling : 1 Min Area: 0.1 % 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:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP023092.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.140	22	26	34	rVB3	7839	13404	0.72%	0.087%
2	3.564	94	98	110	rBV	38757	73517	3.97%	0.475%
3	3.652	110	113	119	rVB2	8649	11476	0.62%	0.074%
4	3.858	143	148	158	rBV	71543	100234	5.42%	0.648%
5	4.228	209	211	217	rVV7	1929	3192	0.17%	0.021%
6	4.293	220	222	226	rVB5	1736	2164	0.12%	0.014%
7	4.752	296	300	304	rBV3	4705	6606	0.36%	0.043%
8	5.616	443	447	448	rBV3	1536	1927	0.10%	0.012%
9	5.634	448	450	453	rVB4	2066	2263	0.12%	0.015%
10	6.164	537	540	545	rVB7	1190	1904	0.10%	0.012%
11	6.699	625	631	632	rBV6	1968	2797	0.15%	0.018%
12	6.787	638	646	665	rBV	110123	242952	13.13%	1.570%
13	6.928	665	670	682	rVB	99130	164947	8.91%	1.066%
14	7.128	698	704	715	rBV	127353	218983	11.84%	1.415%
15	7.405	745	751	754	rBV8	1954	3917	0.21%	0.025%
16	7.581	771	781	797	rBV	318378	519891	28.10%	3.359%
17	8.299	897	903	915	rBV	89372	185281	10.01%	1.197%
18	8.399	915	920	930	rVB	49766	92218	4.98%	0.596%
19	8.728	969	976	991	rBV	126175	237842	12.85%	1.537%
20	9.110	1037	1041	1046	rVB8	2047	3674	0.20%	0.024%
21	9.299	1067	1073	1074	rBV5	3312	4427	0.24%	0.029%
22	9.440	1089	1097	1108	rBV2	98884	196843	10.64%	1.272%
23	9.687	1134	1139	1144	rBV9	4411	11495	0.62%	0.074%
24	9.804	1154	1159	1164	rVB3	3751	5800	0.31%	0.037%
25	9.981	1183	1189	1205	rBV	133534	305244	16.50%	1.972%
26	10.328	1241	1248	1259	rBV	418528	711136	38.44%	4.594%
27	10.487	1266	1275	1292	rBV	87906	234927	12.70%	1.518%
28	10.640	1299	1301	1305	rVB5	1548	1948	0.11%	0.013%
29	10.916	1340	1348	1353	rBV5	3512	9427	0.51%	0.061%
30	11.169	1387	1391	1393	rBV5	2397	3355	0.18%	0.022%
31	11.751	1486	1490	1495	rBV7	2141	4285	0.23%	0.028%
32	11.928	1514	1520	1526	rBV4	2574	5185	0.28%	0.033%
33	12.522	1613	1621	1622	rBV8	1878	3022	0.16%	0.020%
34	12.798	1664	1668	1673	rVB8	1019	1975	0.11%	0.013%
35	13.004	1698	1703	1708	rBV7	1730	3591	0.19%	0.023%
36	13.522	1786	1791	1796	rBV9	2686	5088	0.27%	0.033%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % 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:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
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37	13.598	1796	1804	1817	rBV	362755	516304	27.90%	3.336%
38	13.893	1845	1854	1862	rBV	397515	571399	30.88%	3.692%
39	14.104	1885	1890	1892	rBV5	1638	2180	0.12%	0.014%
40	14.198	1898	1906	1925	rBV	689237	1069712	57.82%	6.911%
41	14.492	1950	1956	1969	rBV2	56079	178067	9.62%	1.150%
42	14.657	1978	1984	1993	rVB8	13333	30603	1.65%	0.198%
43	15.004	2039	2043	2045	rBV4	2017	2210	0.12%	0.014%
44	15.063	2050	2053	2060	rVB9	1990	3811	0.21%	0.025%
45	15.216	2070	2079	2089	rBV	468913	800579	43.27%	5.172%
46	15.363	2098	2104	2115	rBV	92523	158890	8.59%	1.027%
47	15.592	2134	2143	2153	rBV	102692	163328	8.83%	1.055%
48	15.945	2199	2203	2205	rBV5	2237	3100	0.17%	0.020%
49	16.128	2230	2234	2236	rBV5	2418	3959	0.21%	0.026%
50	16.539	2299	2304	2309	rBV4	9908	17682	0.96%	0.114%
51	16.739	2334	2338	2340	rBV5	2324	3787	0.20%	0.024%
52	17.010	2376	2384	2390	rBV	957320	1381056	74.64%	8.923%
53	17.104	2395	2400	2413	rVV	708362	939151	50.76%	6.068%
54	17.222	2413	2420	2429	rVB10	18201	36875	1.99%	0.238%
55	17.404	2447	2451	2452	rBV4	3266	4521	0.24%	0.029%
56	17.516	2468	2470	2474	rBV5	2908	4079	0.22%	0.026%
57	17.704	2498	2502	2507	rBV8	3665	7310	0.40%	0.047%
58	17.810	2517	2520	2523	rBV5	3539	5158	0.28%	0.033%
59	17.939	2536	2542	2550	rBV	121805	188437	10.18%	1.217%
60	18.533	2640	2643	2650	rVB3	8696	13321	0.72%	0.086%
61	18.775	2681	2684	2688	rBV2	12587	18545	1.00%	0.120%
62	19.151	2744	2748	2754	rVB	38538	60820	3.29%	0.393%
63	19.280	2766	2770	2778	rBV2	36893	62022	3.35%	0.401%
64	19.492	2800	2806	2816	rBV	744387	1140088	61.62%	7.366%
65	21.116	3079	3082	3083	rBV3	44141	48284	2.61%	0.312%
66	21.439	3131	3137	3144	rBV	1254652	1798574	97.21%	11.620%
67	24.439	3640	3647	3656	rVB	405044	997040	53.89%	6.442%
68	24.657	3674	3684	3695	rBV	674911	1850226	100.00%	11.954%

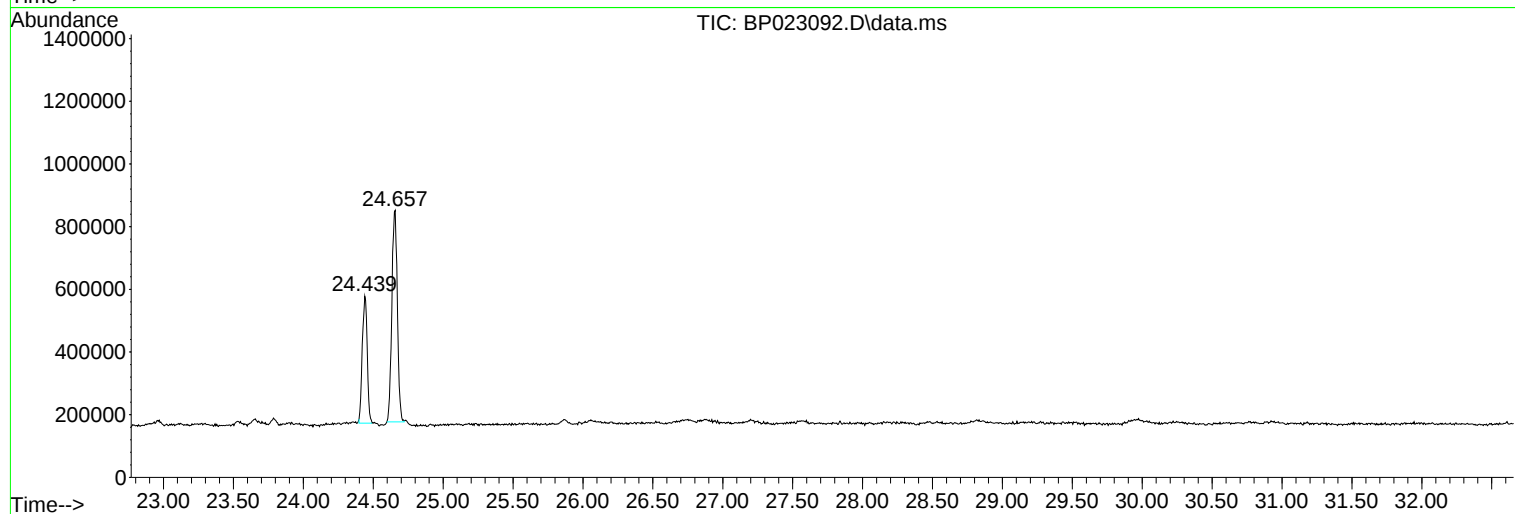
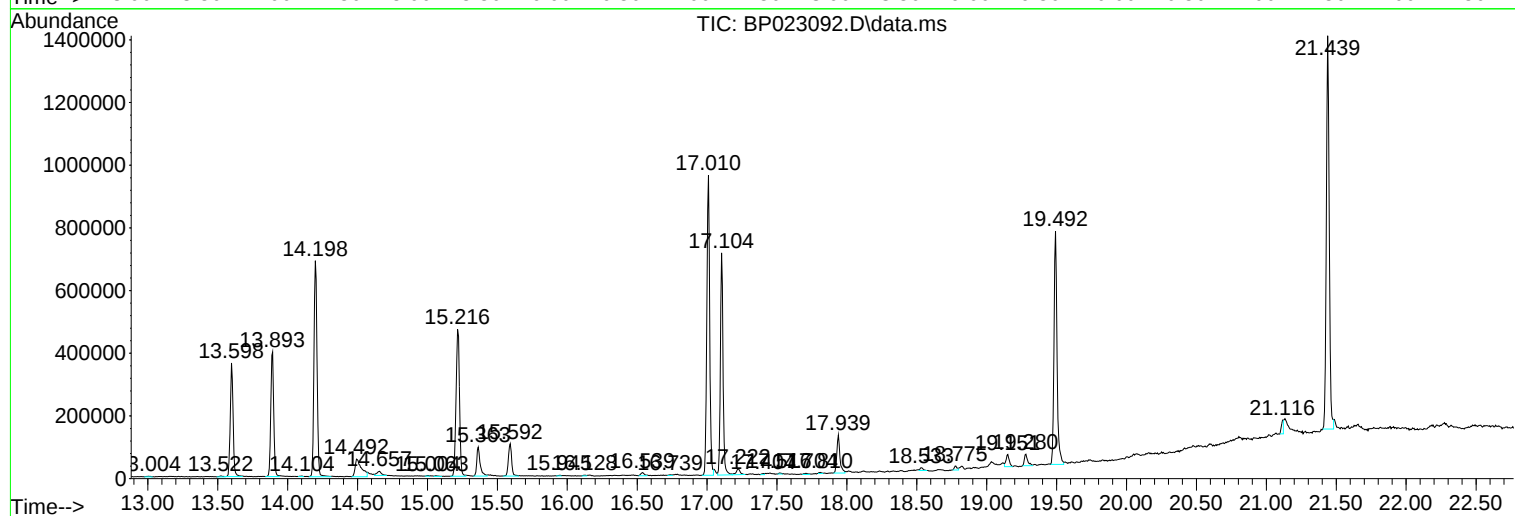
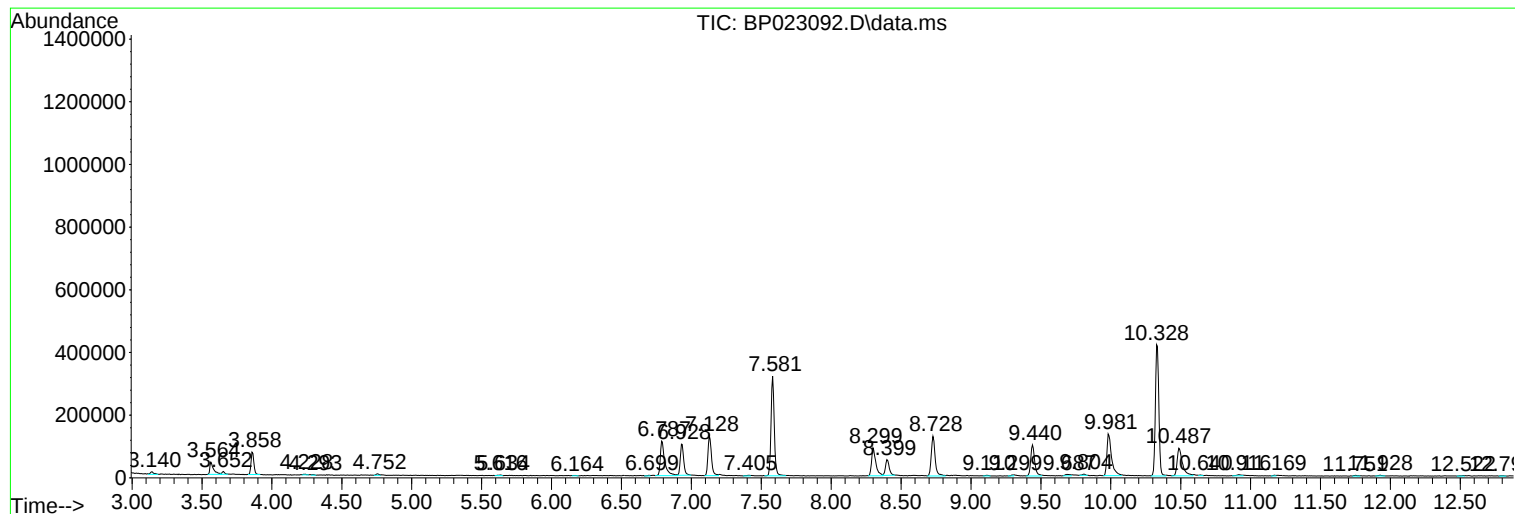
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P



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Misc :
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Instrument :
BNA_P
ClientSampleId :
BH9B9

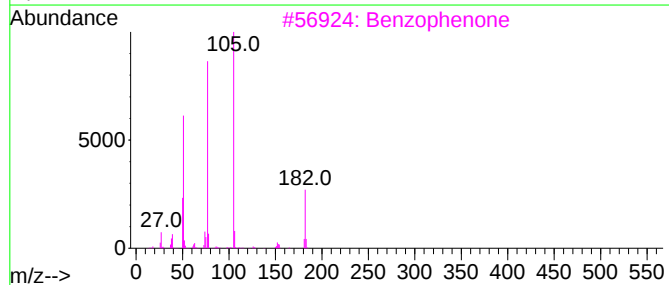
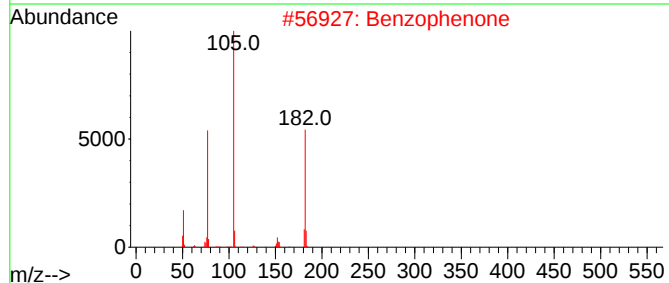
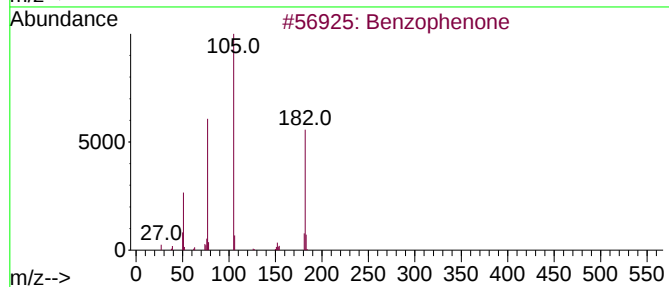
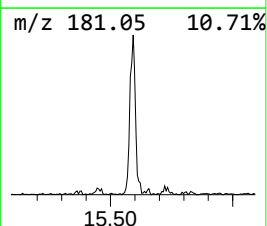
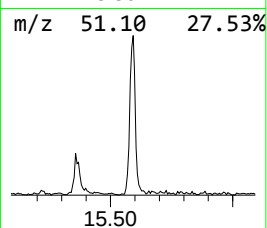
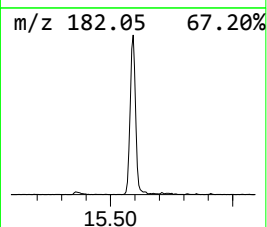
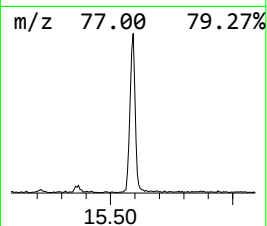
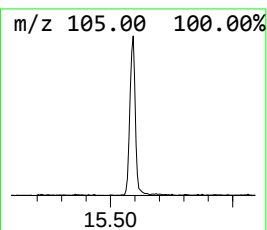
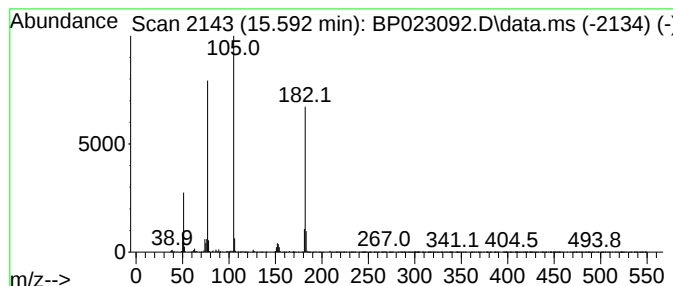
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzophenone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.592	3.05 ng/ul	163328	Acenaphthene-d10	14.198

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	97
2			Benzophenone	182	C13H10O	000119-61-9	97
3			Benzophenone	182	C13H10O	000119-61-9	93
4			Benzophenone	182	C13H10O	000119-61-9	91
5			1,2,4-Trioxolane, 3,3,5-triphenyl-	304	C20H16O3	023246-12-0	83



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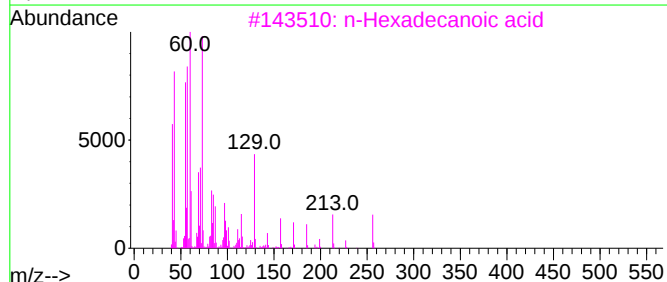
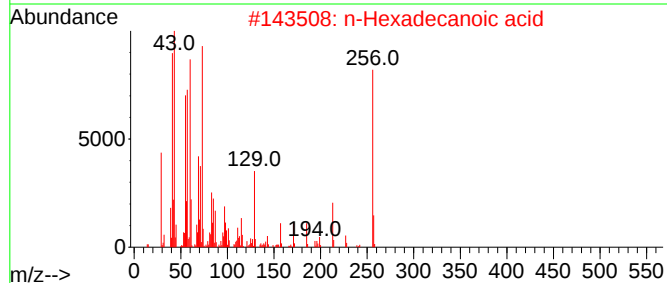
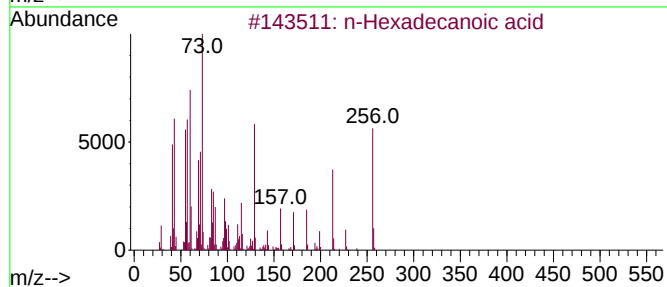
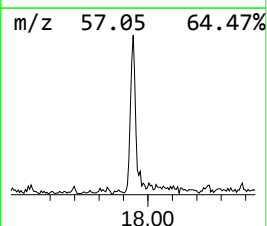
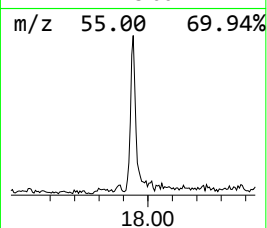
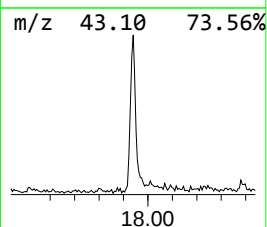
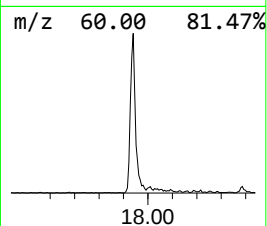
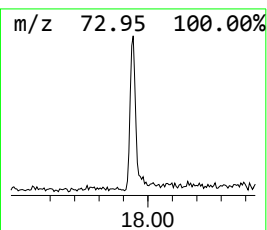
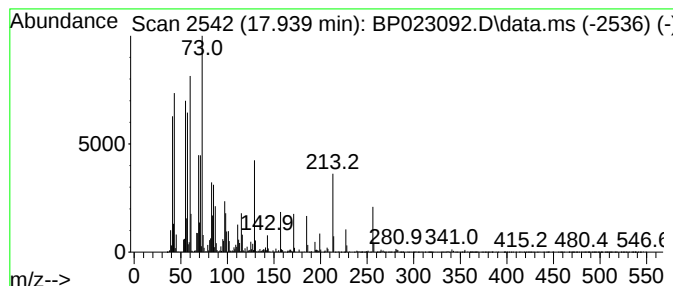
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.939	2.73 ng/ul	188437	Phenanthrene-d10	17.010

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	98
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	93



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzophenone	15.592	3.0	ng/ul	163328	3	14.198	1069710	20.0
n-Hexadecanoic ...	17.939	2.7	ng/ul	188437	4	17.010	1381060	20.0