

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110424\
 Data File : BP022900.D
 Acq On : 06 Nov 2024 19:17
 Operator : RC/JU
 Sample : P4622-01
 Misc :
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BH6M7

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-BP100724.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP022900.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.170	28	31	37	rBV	7059	9067	0.70%	0.063%
2	3.575	96	100	103	rBV5	1059	1910	0.15%	0.013%
3	3.611	103	106	112	rBV2	7845	14410	1.12%	0.100%
4	3.893	151	154	161	rVB5	4608	7371	0.57%	0.051%
5	4.069	182	184	189	rVB6	1176	1749	0.14%	0.012%
6	4.122	189	193	196	rVB5	1429	1687	0.13%	0.012%
7	4.393	236	239	244	rVB6	2071	2779	0.22%	0.019%
8	5.446	415	418	421	rBV5	1394	1529	0.12%	0.011%
9	6.681	626	628	633	rVB6	1061	1394	0.11%	0.010%
10	6.822	646	652	668	rBV2	121302	241812	18.73%	1.676%
11	6.969	670	677	687	rBV	103222	163410	12.66%	1.132%
12	7.169	703	711	722	rBV	132767	238138	18.44%	1.650%
13	7.616	779	787	797	rBV	254513	418106	32.38%	2.897%
14	8.328	902	908	926	rVV	166022	329816	25.54%	2.286%
15	8.763	973	982	994	rBV	133922	248035	19.21%	1.719%
16	8.922	1005	1009	1011	rBV5	1558	2181	0.17%	0.015%
17	9.163	1044	1050	1056	rVB9	3073	5918	0.46%	0.041%
18	9.346	1077	1081	1083	rBV5	1364	1532	0.12%	0.011%
19	9.475	1097	1103	1120	rBV	115073	217384	16.84%	1.506%
20	9.746	1143	1149	1151	rBV7	733	1467	0.11%	0.010%
21	10.022	1188	1196	1214	rBV	159738	365526	28.31%	2.533%
22	10.375	1246	1256	1269	rBV	328151	579142	44.86%	4.013%
23	10.522	1274	1281	1298	rBV	139975	306102	23.71%	2.121%
24	10.904	1342	1346	1348	rBV5	1434	1703	0.13%	0.012%
25	11.551	1453	1456	1461	rVB7	710	1364	0.11%	0.009%
26	11.787	1489	1496	1501	rBV7	1324	3314	0.26%	0.023%
27	11.840	1501	1505	1511	rVB9	761	1393	0.11%	0.010%
28	11.951	1518	1524	1525	rBV6	1178	1396	0.11%	0.010%
29	11.975	1525	1528	1535	rVB8	2365	4515	0.35%	0.031%
30	12.069	1542	1544	1548	rVB5	1042	1459	0.11%	0.010%
31	12.675	1644	1647	1650	rVB4	1177	1724	0.13%	0.012%
32	13.057	1708	1712	1716	rBV7	2600	4386	0.34%	0.030%
33	13.128	1721	1724	1730	rVB8	759	1434	0.11%	0.010%
34	13.540	1789	1794	1796	rBV5	1636	2322	0.18%	0.016%
35	13.645	1805	1812	1823	rBV	329637	538540	41.71%	3.732%
36	13.928	1852	1860	1870	rBV	427798	668590	51.78%	4.633%

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37	14.140	1893	1896	1900	rBV5	1637	2175	0.17%	0.015%
38	14.240	1905	1913	1928	rVV	614108	875840	67.84%	6.069%
39	14.351	1928	1932	1935	rVB6	1022	1617	0.13%	0.011%
40	14.457	1948	1950	1953	rBV4	1389	1400	0.11%	0.010%
41	14.522	1955	1961	1979	rBV2	56005	178317	13.81%	1.236%
42	14.704	1987	1992	1995	rVB6	3330	4735	0.37%	0.033%
43	14.822	2008	2012	2014	rVB4	991	1469	0.11%	0.010%
44	15.057	2050	2052	2057	rVB6	1331	1573	0.12%	0.011%
45	15.098	2057	2059	2061	rBV3	1250	1454	0.11%	0.010%
46	15.251	2079	2085	2099	rVB	671735	915042	70.87%	6.341%
47	15.416	2105	2113	2123	rBV3	34569	87543	6.78%	0.607%
48	15.622	2143	2148	2158	rVB	151890	194805	15.09%	1.350%
49	15.816	2177	2181	2182	rBV3	1157	1450	0.11%	0.010%
50	15.839	2184	2185	2188	rVB3	1685	1499	0.12%	0.010%
51	16.186	2240	2244	2245	rBV4	1480	1458	0.11%	0.010%
52	16.210	2245	2248	2255	rVB4	9784	13272	1.03%	0.092%
53	16.457	2288	2290	2293	rVB4	1463	1355	0.10%	0.009%
54	16.928	2366	2370	2372	rBV5	1461	2479	0.19%	0.017%
55	16.969	2374	2377	2382	rVB6	2034	2809	0.22%	0.019%
56	17.033	2383	2388	2395	rBV	787705	1121535	86.86%	7.772%
57	17.151	2401	2408	2420	rVB2	535639	952035	73.74%	6.597%
58	17.269	2425	2428	2433	rBV7	5211	7724	0.60%	0.054%
59	17.469	2460	2462	2464	rBV3	2627	2948	0.23%	0.020%
60	17.569	2477	2479	2480	rBV2	2493	2054	0.16%	0.014%
61	17.980	2544	2549	2553	rBV2	38981	57592	4.46%	0.399%
62	18.157	2577	2579	2582	rBV3	2925	3395	0.26%	0.024%
63	18.292	2599	2602	2606	rBV5	5444	9421	0.73%	0.065%
64	18.422	2622	2624	2630	rVB6	9913	12421	0.96%	0.086%
65	19.080	2732	2736	2742	rBV7	10798	18386	1.42%	0.127%
66	19.151	2745	2748	2749	rVV3	9473	12017	0.93%	0.083%
67	19.180	2749	2753	2759	rVB	32269	56085	4.34%	0.389%
68	19.316	2774	2776	2779	rBV3	8548	9568	0.74%	0.066%
69	19.527	2806	2812	2822	rBV	839869	1279129	99.07%	8.864%
70	21.492	3141	3146	3154	rBV	770804	1242992	96.27%	8.614%
71	23.015	3400	3405	3419	rVB2	214731	475044	36.79%	3.292%
72	24.510	3651	3659	3670	rVB2	429798	1194145	92.49%	8.275%
73	24.745	3691	3699	3711	rVB	438785	1291132	100.00%	8.947%

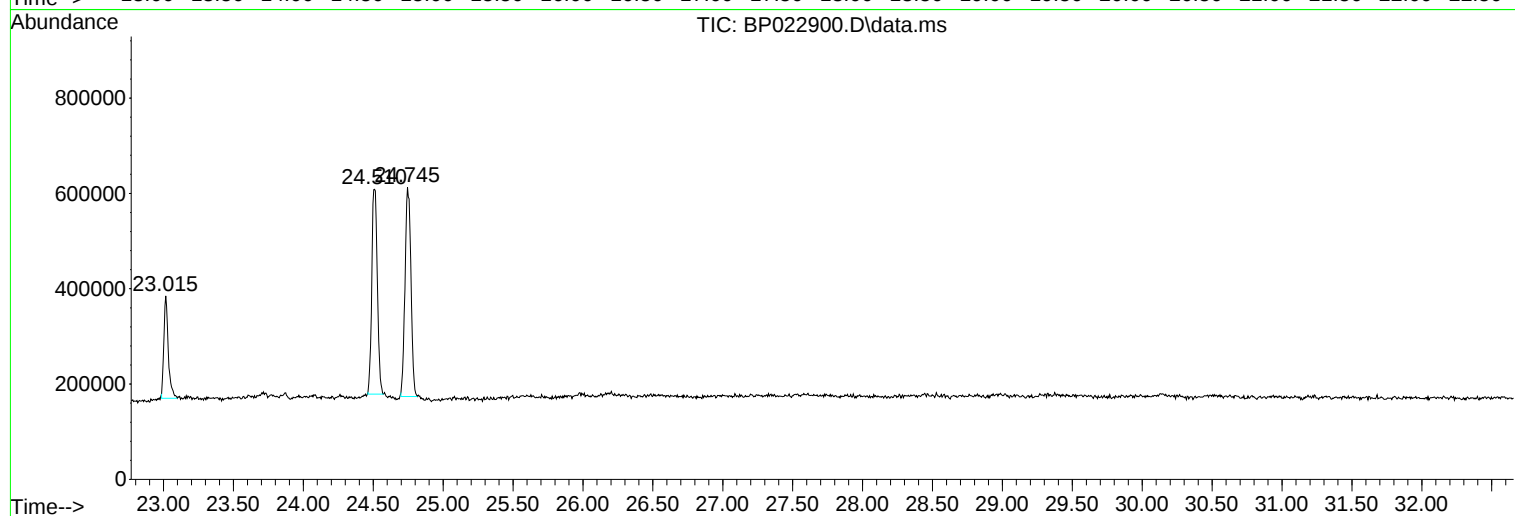
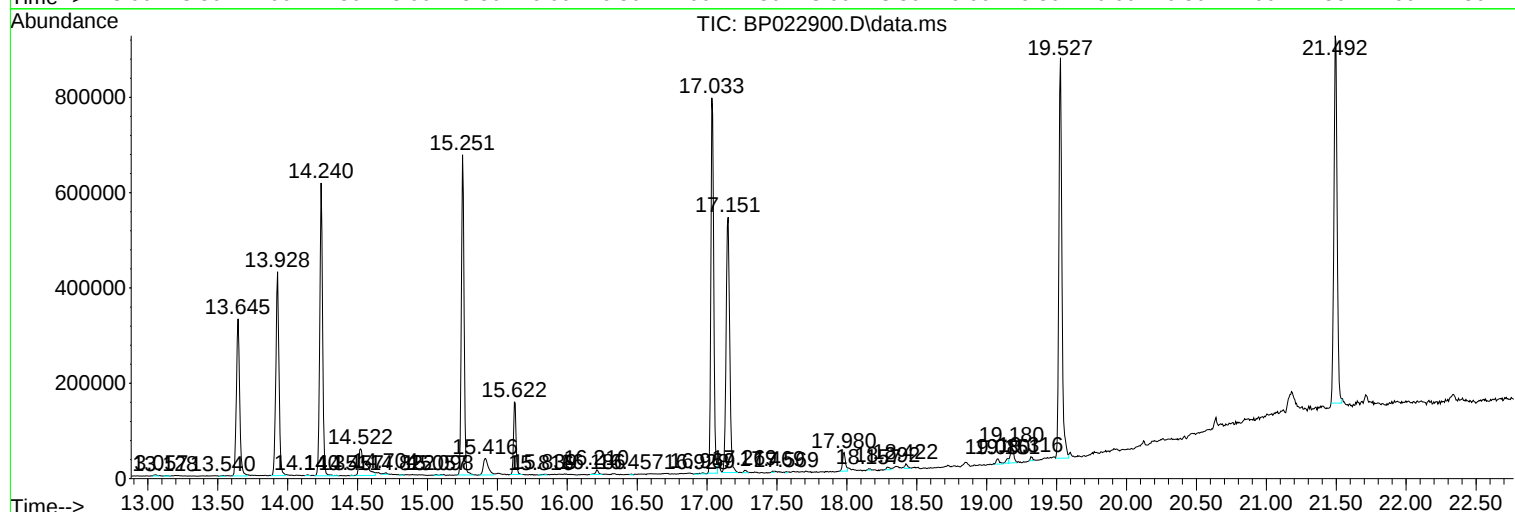
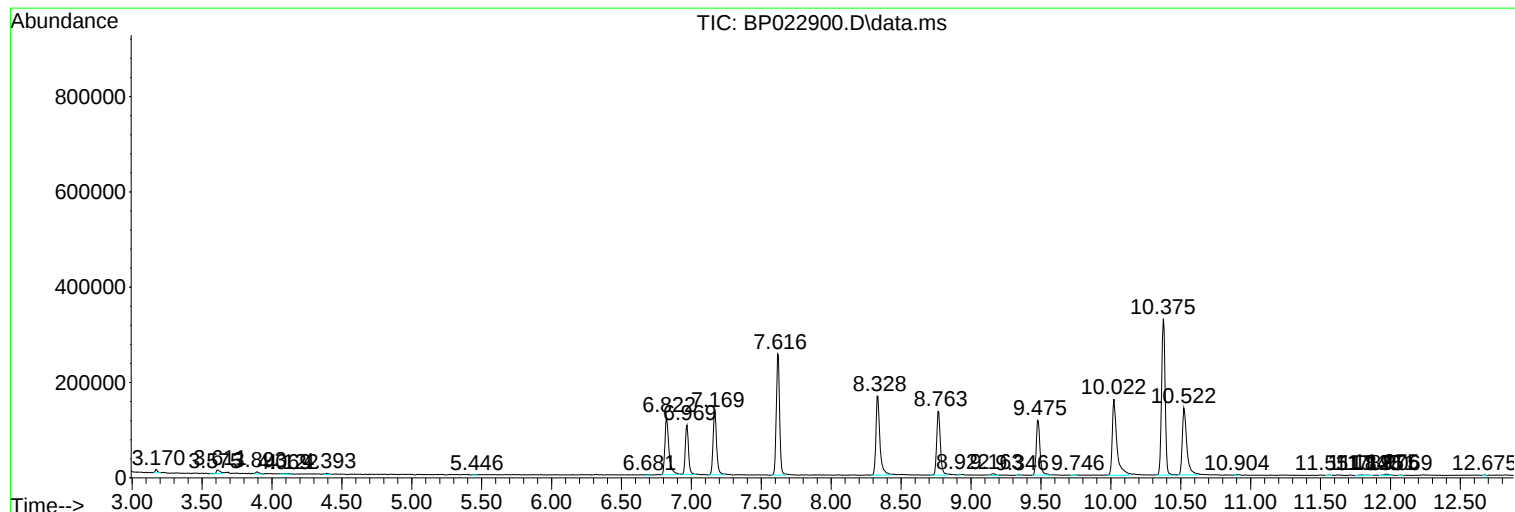
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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



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Operator : RC/JU
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Instrument :
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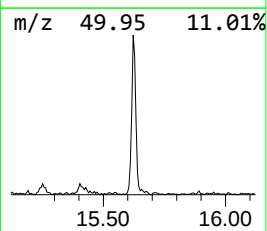
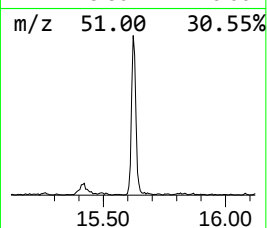
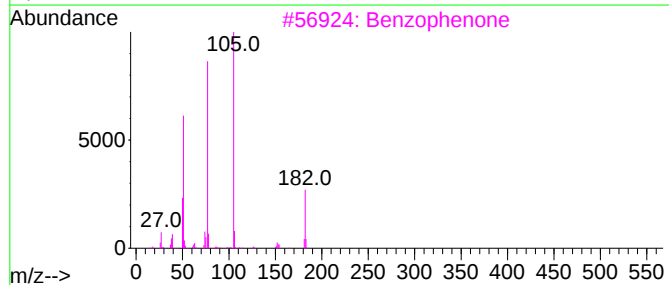
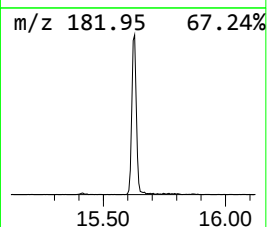
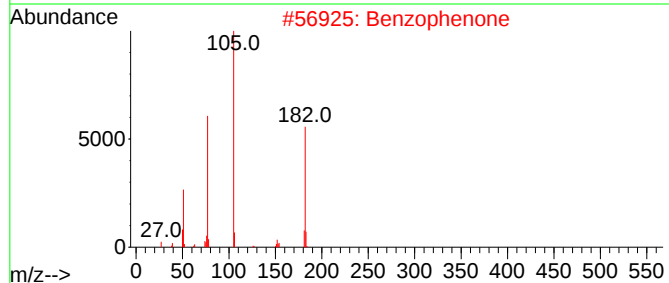
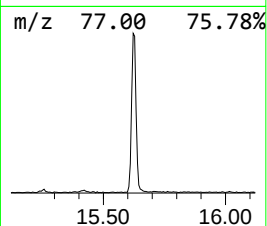
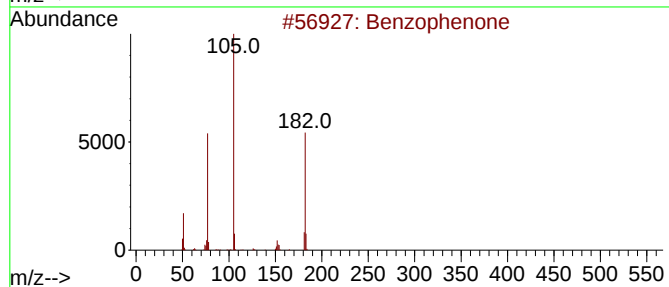
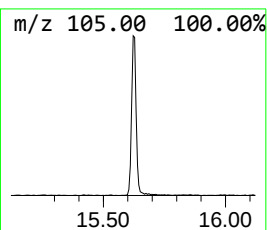
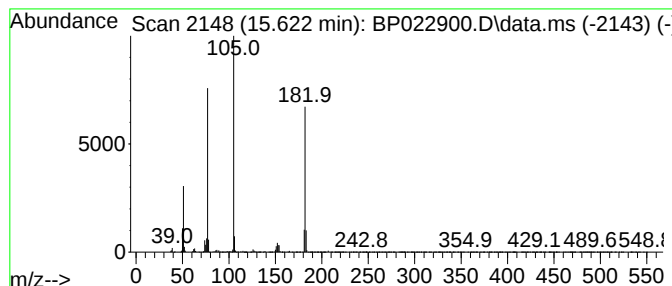
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzophenone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.622	4.45 ng/ul	194805	Acenaphthene-d10	14.240

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	97
2			Benzophenone	182	C13H10O	000119-61-9	96
3			Benzophenone	182	C13H10O	000119-61-9	95
4			Benzophenone	182	C13H10O	000119-61-9	94
5			Benzophenone	182	C13H10O	000119-61-9	91



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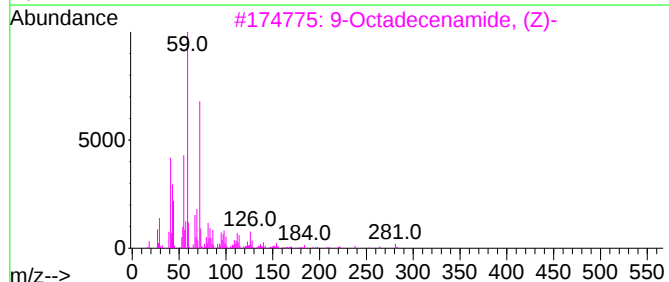
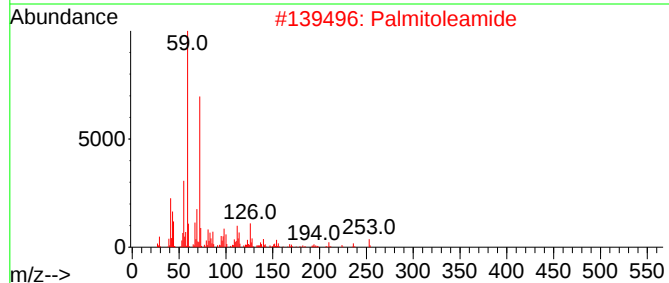
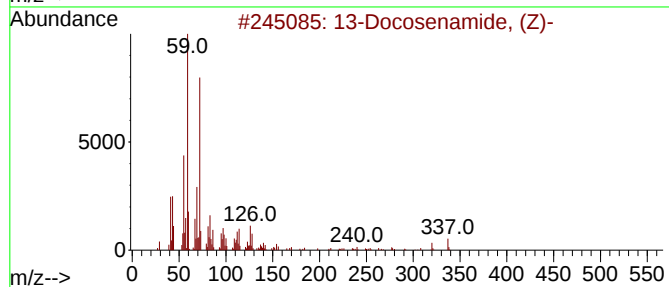
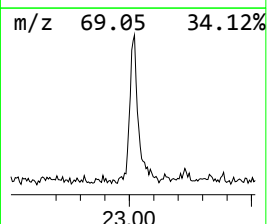
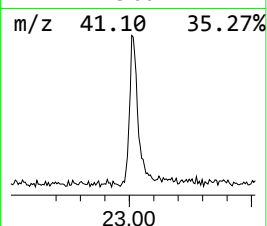
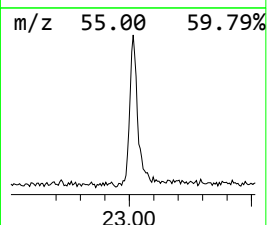
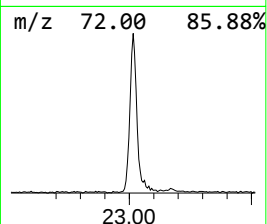
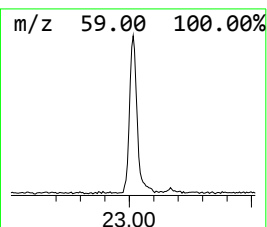
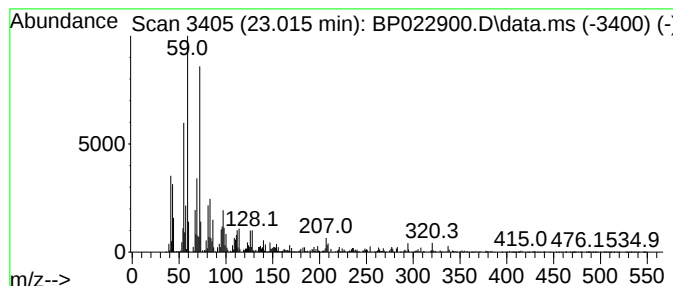
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 13-Docosenamide, (Z)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.015	7.64 ng/ul	475044	Chrysene-d12	21.492

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	96
2			Palmitoleamide	253	C16H31NO	106010-22-4	83
3			9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	83
4			13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	50
5			13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	46



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzophenone	15.622	4.5	ng/u1	194805	3	14.240	875840	20.0
13-Docosenamide...	23.015	7.6	ng/u1	475044	5	21.492	1242990	20.0