

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
 Data File : BP013196.D
 Acq On : 21 Dec 2022 08:24
 Operator : CG/JU
 Sample : N5984-13
 Misc :
 ALS Vial : 78 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 COLE4

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-BP120722.M
 Title : SVOA CALIBRATION

Signal : TIC: BP013196.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.334	21	25	31	rVB	130160	163703	1.53%	0.139%
2	3.764	95	98	115	rBV	1115866	1674202	15.62%	1.417%
3	3.993	135	137	141	rVB3	16144	17002	0.16%	0.014%
4	4.093	150	154	158	rBV	43020	55655	0.52%	0.047%
5	5.046	312	316	322	rVB	29558	50903	0.47%	0.043%
6	6.928	632	636	639	rBV5	7129	11683	0.11%	0.010%
7	7.111	660	667	677	rBV	2106881	3332119	31.08%	2.820%
8	7.275	689	695	707	rVB	1759172	2713630	25.31%	2.296%
9	7.481	723	730	741	rBV	2135045	3376958	31.50%	2.858%
10	7.952	803	810	817	rBV	1878558	2916556	27.21%	2.468%
11	8.011	817	820	828	rVB	21342	35540	0.33%	0.030%
12	8.658	923	930	942	rBV	2571267	4166584	38.87%	3.526%
13	9.116	1001	1008	1016	rBV	2035295	3318190	30.95%	2.808%
14	9.275	1031	1035	1043	rVB	10880	19830	0.18%	0.017%
15	9.522	1069	1077	1084	rVB	60822	101512	0.95%	0.086%
16	9.840	1124	1131	1144	rBV	1691010	2828128	26.38%	2.393%
17	10.069	1164	1170	1177	rBV3	23057	46168	0.43%	0.039%
18	10.381	1213	1223	1237	rBV	2632777	4359682	40.67%	3.689%
19	10.763	1280	1288	1297	rBV	2564160	4233497	39.49%	3.583%
20	10.904	1303	1312	1329	rBV	2193310	3816457	35.60%	3.230%
21	11.116	1343	1348	1351	rBV6	8535	11186	0.10%	0.009%
22	11.263	1366	1373	1378	rBV3	11674	32528	0.30%	0.028%
23	12.175	1522	1528	1532	rBV	31384	48895	0.46%	0.041%
24	12.346	1551	1557	1562	rBV2	48375	77928	0.73%	0.066%
25	12.863	1638	1645	1649	rBV8	14331	27285	0.25%	0.023%
26	12.981	1662	1665	1670	rVB5	11408	12945	0.12%	0.011%
27	13.063	1676	1679	1685	rVB5	9324	12387	0.12%	0.010%
28	13.122	1685	1689	1694	rVB4	11534	17766	0.17%	0.015%
29	13.404	1728	1737	1745	rBV	64048	102334	0.95%	0.087%
30	13.569	1758	1765	1770	rVB7	12056	24569	0.23%	0.021%
31	13.998	1831	1838	1847	rBV	4399600	5963920	55.64%	5.047%
32	14.104	1852	1856	1865	rVB	17889	34342	0.32%	0.029%
33	14.181	1865	1869	1872	rBV6	7976	11146	0.10%	0.009%
34	14.240	1874	1879	1880	rBV4	10401	14690	0.14%	0.012%
35	14.287	1880	1887	1898	rBV	5126481	7449181	69.49%	6.304%
36	14.475	1915	1919	1926	rVB3	23442	34573	0.32%	0.029%

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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-BP120722.M
 Title : SVOA CALIBRATION

37	14.593	1932	1939	1947	rBV	3693270	5259170	49.06%	4.451%
38	14.681	1952	1954	1961	rVB8	8266	12407	0.12%	0.010%
39	14.787	1966	1972	1986	rBV	1294109	2112629	19.71%	1.788%
40	14.998	2003	2008	2016	rBV3	95596	154841	1.44%	0.131%
41	15.381	2067	2073	2075	rBV6	10164	18002	0.17%	0.015%
42	15.434	2078	2082	2085	rVB4	17095	22251	0.21%	0.019%
43	15.587	2101	2108	2120	rBV	6489114	8933316	83.34%	7.560%
44	15.704	2122	2128	2143	rVV	1375695	1931327	18.02%	1.634%
45	15.828	2143	2149	2155	rVB4	34866	63757	0.59%	0.054%
46	15.951	2165	2170	2182	rVB	395101	536831	5.01%	0.454%
47	16.292	2225	2228	2231	rBV5	14667	17086	0.16%	0.014%
48	16.357	2237	2239	2244	rBV5	6004	10828	0.10%	0.009%
49	16.522	2263	2267	2270	rVB4	9299	13538	0.13%	0.011%
50	16.734	2301	2303	2308	rVB5	18085	22054	0.21%	0.019%
51	16.998	2344	2348	2351	rBV5	13576	22949	0.21%	0.019%
52	17.098	2361	2365	2368	rBV5	14387	19122	0.18%	0.016%
53	17.351	2401	2408	2418	rBV	3836127	5385476	50.24%	4.558%
54	17.451	2418	2425	2436	rVV	5868068	8423910	78.58%	7.129%
55	17.539	2436	2440	2448	rVB3	52946	84914	0.79%	0.072%
56	17.722	2468	2471	2477	rBV7	29893	50267	0.47%	0.043%
57	17.957	2509	2511	2516	rVB6	13134	16283	0.15%	0.014%
58	18.004	2516	2519	2524	rBV7	16995	22594	0.21%	0.019%
59	18.216	2550	2555	2565	rBV	318841	444154	4.14%	0.376%
60	18.633	2623	2626	2628	rVB3	27562	29504	0.28%	0.025%
61	19.257	2728	2732	2738	rVB2	66249	86915	0.81%	0.074%
62	19.328	2739	2744	2749	rBV	75492	135090	1.26%	0.114%
63	19.451	2762	2765	2769	rBV4	55498	66936	0.62%	0.057%
64	19.680	2798	2804	2815	rBV	7354979	9496647	88.59%	8.037%
65	21.127	3046	3050	3057	rBV2	714609	1039779	9.70%	0.880%
66	21.433	3097	3102	3109	rBV	4041109	5552337	51.80%	4.699%
67	23.757	3490	3497	3509	rVB2	5152097	10719708	100.00%	9.072%
68	23.916	3517	3524	3534	rVB2	3087621	6346535	59.20%	5.371%

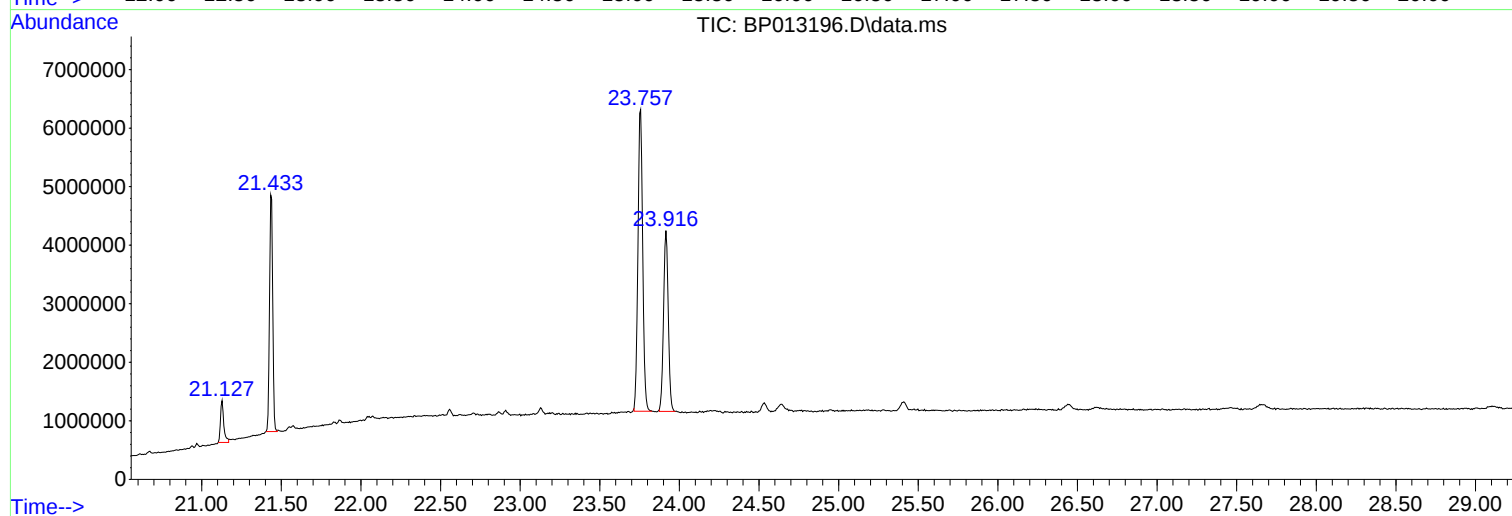
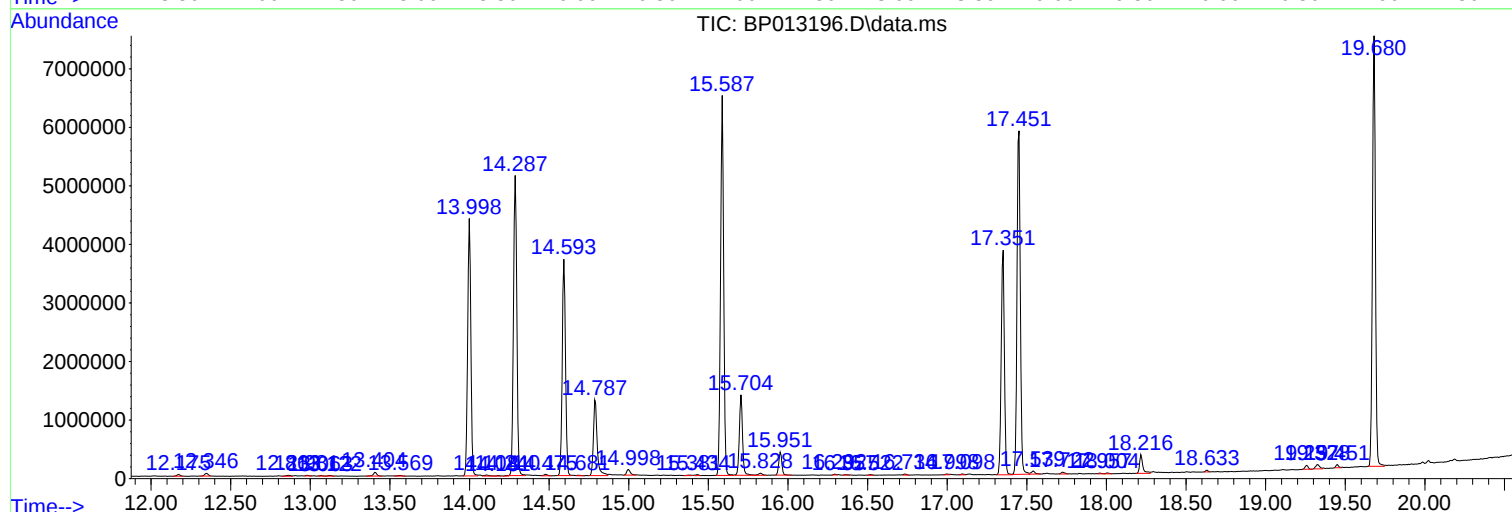
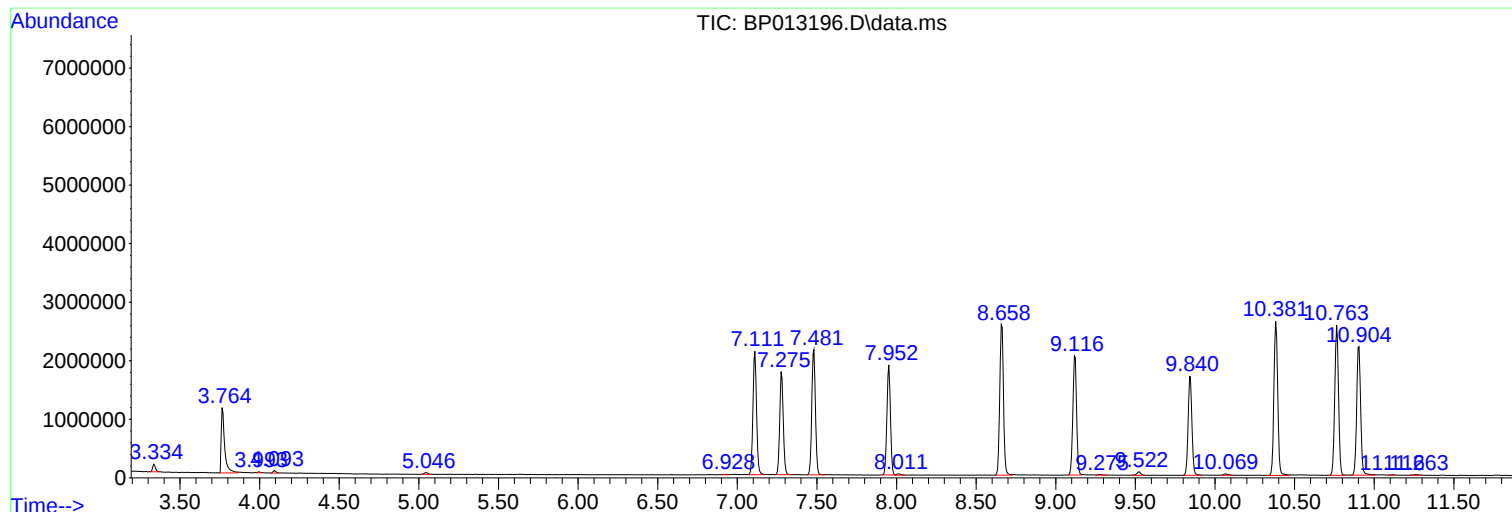
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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



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Operator : CG/JU
Sample : N5984-13
Misc :
ALS Vial : 78 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
COLE4

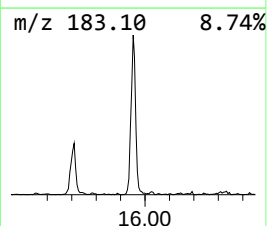
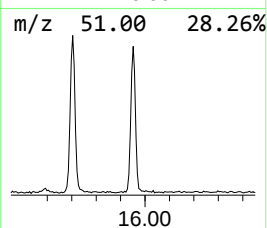
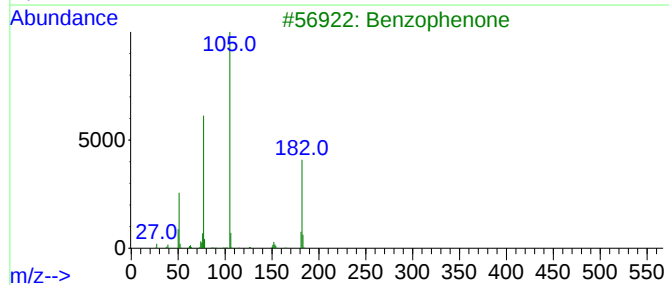
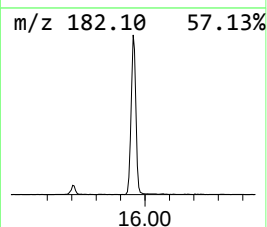
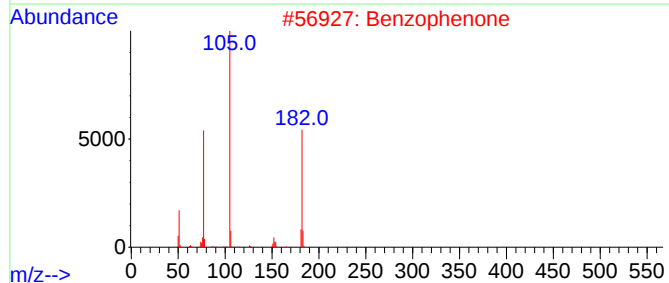
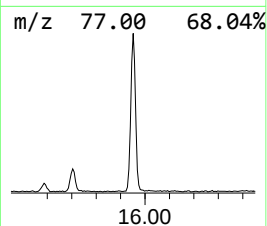
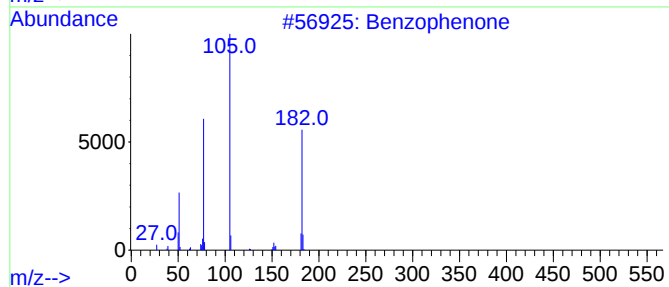
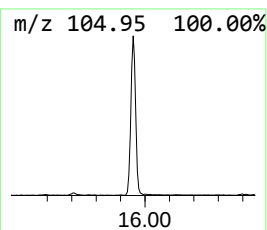
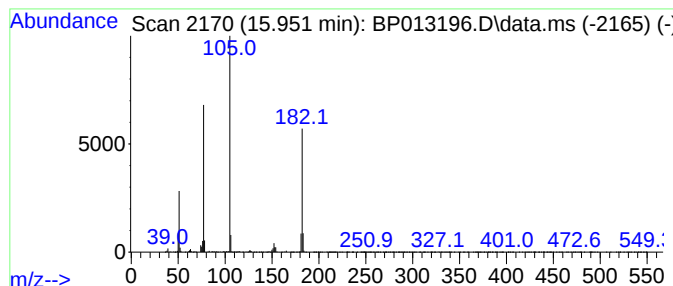
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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.
15.951	2.04 ng/ul	536831	Acenaphthene-d10	14.592

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzophenone	182	C13H10O	000119-61-9	95
3			Benzophenone	182	C13H10O	000119-61-9	94
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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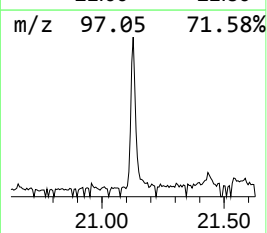
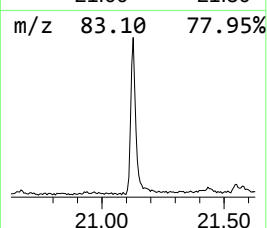
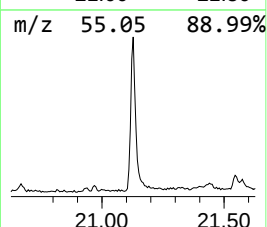
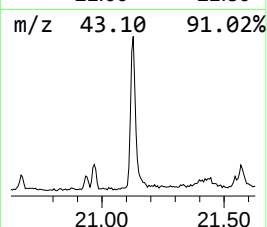
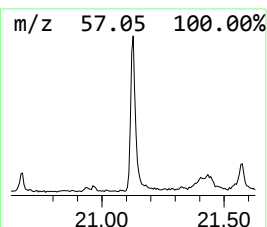
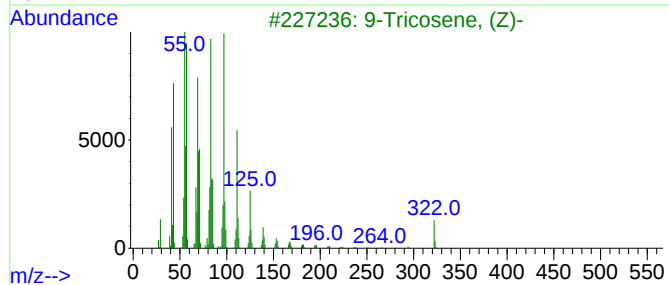
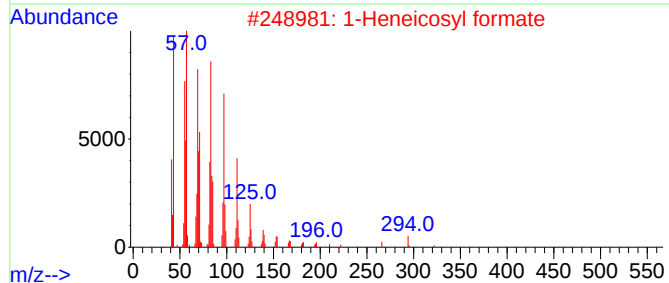
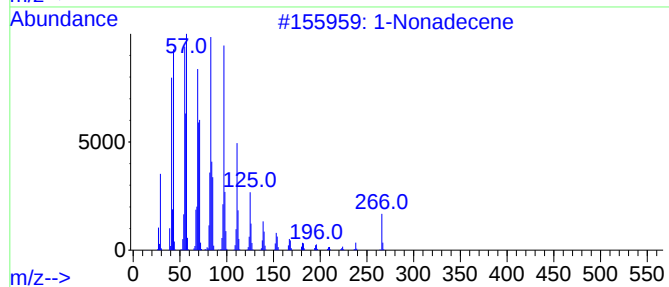
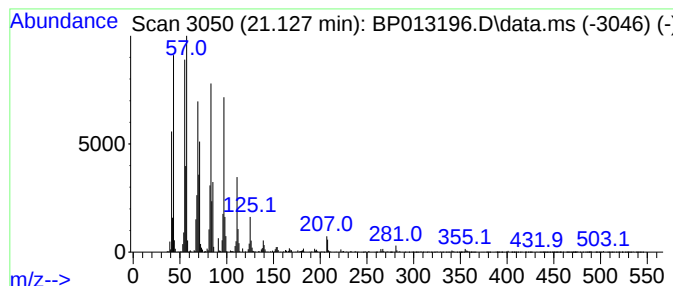
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.128	3.75 ng/ul	1039780	Chrysene-d12	21.433

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Nonadecene			266	C19H38	018435-45-5	98
2	1-Heneicosyl formate			340	C22H44O2	077899-03-7	97
3	9-Tricosene, (Z)-			322	C23H46	027519-02-4	96
4	1-Nonadecene			266	C19H38	018435-45-5	95
5	Dichloroacetic acid, heptadecyl ...			366	C19H36Cl2O2	1000282-98-2	94



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzophenone	15.951	2.0	ng/u1	536831	3	14.592	5259170	20.0
(DEL) Alkane: S...	21.128	3.8	ng/u1	1039780	5	21.433	5552340	20.0