

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101424\
 Data File : BP022394.D
 Acq On : 15 Oct 2024 18:19
 Operator : RC/JU
 Sample : P4349-11
 Misc :
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EOAG4

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: BP022394.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.117	20	22	29	rVB7	4024	5740	0.17%	0.020%
2	3.222	36	40	47	rVB	20419	28138	0.86%	0.100%
3	3.640	107	111	124	rBV	55797	93337	2.84%	0.330%
4	3.952	160	164	173	rVB6	5985	11756	0.36%	0.042%
5	4.117	187	192	194	rBV6	2117	3589	0.11%	0.013%
6	4.181	198	203	208	rBV8	2803	5291	0.16%	0.019%
7	4.322	223	227	231	rBV4	6305	8608	0.26%	0.030%
8	4.452	245	249	253	rBV2	3948	6304	0.19%	0.022%
9	4.505	253	258	266	rVB3	12685	20551	0.62%	0.073%
10	4.869	312	320	326	rBV8	3372	8120	0.25%	0.029%
11	5.540	430	434	443	rBV9	3060	6274	0.19%	0.022%
12	5.740	460	468	476	rVB2	7948	16078	0.49%	0.057%
13	5.987	507	510	515	rVB6	2738	4038	0.12%	0.014%
14	6.105	520	530	537	rVB2	28770	55839	1.70%	0.198%
15	6.528	597	602	605	rBV6	2673	3862	0.12%	0.014%
16	6.569	605	609	617	rVB3	6457	9858	0.30%	0.035%
17	6.840	651	655	656	rBV3	2878	3588	0.11%	0.013%
18	6.881	656	662	673	rVB	165603	271096	8.24%	0.960%
19	7.034	682	688	700	rBV	257472	408258	12.41%	1.445%
20	7.234	715	722	733	rBV	326703	549510	16.71%	1.945%
21	7.693	793	800	809	rBV	327671	548957	16.69%	1.943%
22	7.805	815	819	822	rVB6	2592	3350	0.10%	0.012%
23	8.393	911	919	930	rBV	352321	628723	19.12%	2.225%
24	8.834	986	994	1004	rBV	356775	581395	17.68%	2.058%
25	9.246	1058	1064	1069	rBV8	4006	6790	0.21%	0.024%
26	9.393	1084	1089	1096	rBV2	7446	12056	0.37%	0.043%
27	9.546	1105	1115	1122	rBV	300956	517455	15.73%	1.832%
28	10.081	1199	1206	1220	rBV	513222	840619	25.56%	2.975%
29	10.452	1260	1269	1278	rBV	443890	769746	23.40%	2.725%
30	10.587	1283	1292	1307	rBV	261269	497930	15.14%	1.762%
31	11.275	1400	1409	1417	rVB	1781	5720	0.17%	0.020%
32	11.369	1421	1425	1430	rBV7	2247	4432	0.13%	0.016%
33	11.722	1480	1485	1491	rBV8	2914	5843	0.18%	0.021%
34	12.034	1531	1538	1543	rBV2	8014	13443	0.41%	0.048%
35	12.910	1680	1687	1691	rVB9	2021	3490	0.11%	0.012%
36	13.710	1817	1823	1832	rBV	923319	1295803	39.40%	4.587%

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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-BP100724.MA.M
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37	14.004	1864	1873	1883	rBV	1013280	1438375	43.74%	5.091%
38	14.310	1917	1925	1933	rBV	798754	1191777	36.24%	4.218%
39	14.534	1956	1963	1976	rBV	79822	188355	5.73%	0.667%
40	14.687	1983	1989	1993	rBV5	8256	15189	0.46%	0.054%
41	14.728	1993	1996	2003	rVB9	5613	10736	0.33%	0.038%
42	15.322	2089	2097	2104	rBV	1322185	1970709	59.92%	6.976%
43	15.445	2112	2118	2129	rBV	561124	672616	20.45%	2.381%
44	15.557	2134	2137	2144	rVB6	6773	10037	0.31%	0.036%
45	15.704	2155	2162	2167	rBV	289844	407762	12.40%	1.443%
46	15.887	2189	2193	2198	rVB8	2229	3770	0.11%	0.013%
47	16.039	2213	2219	2229	rBV	45761	73379	2.23%	0.260%
48	16.263	2250	2257	2264	rVB2	33855	54831	1.67%	0.194%
49	16.504	2293	2298	2304	rBV10	5650	9859	0.30%	0.035%
50	16.710	2328	2333	2339	rBV10	3131	6000	0.18%	0.021%
51	16.786	2343	2346	2350	rVB6	3126	3996	0.12%	0.014%
52	16.828	2350	2353	2357	rBV5	2247	4051	0.12%	0.014%
53	16.981	2375	2379	2382	rVB5	3886	4860	0.15%	0.017%
54	17.116	2394	2402	2408	rBV	1185895	1648559	50.13%	5.835%
55	17.210	2412	2418	2429	rVB	1671108	2394523	72.81%	8.476%
56	17.516	2468	2470	2477	rVB8	3294	5133	0.16%	0.018%
57	17.904	2532	2536	2546	rVB8	6514	11332	0.34%	0.040%
58	18.051	2555	2561	2569	rBV2	94834	145466	4.42%	0.515%
59	18.345	2608	2611	2613	rBV4	4513	6679	0.20%	0.024%
60	18.551	2644	2646	2650	rVB4	7501	8352	0.25%	0.030%
61	18.922	2705	2709	2714	rBV5	12992	22660	0.69%	0.080%
62	19.133	2740	2745	2749	rBV4	20670	32583	0.99%	0.115%
63	19.216	2755	2759	2762	rBV	26582	38528	1.17%	0.136%
64	19.380	2783	2787	2792	rBV3	47900	66779	2.03%	0.236%
65	19.528	2809	2812	2815	rBV5	11904	14979	0.46%	0.053%
66	19.592	2816	2823	2830	rBV	2204668	3098008	94.20%	10.966%
67	21.539	3148	3154	3164	rBV	1367726	2049285	62.31%	7.254%
68	24.598	3665	3674	3689	rVB	1253831	3288826	100.00%	11.641%
69	24.839	3706	3715	3726	rVB2	805320	2097864	63.79%	7.426%

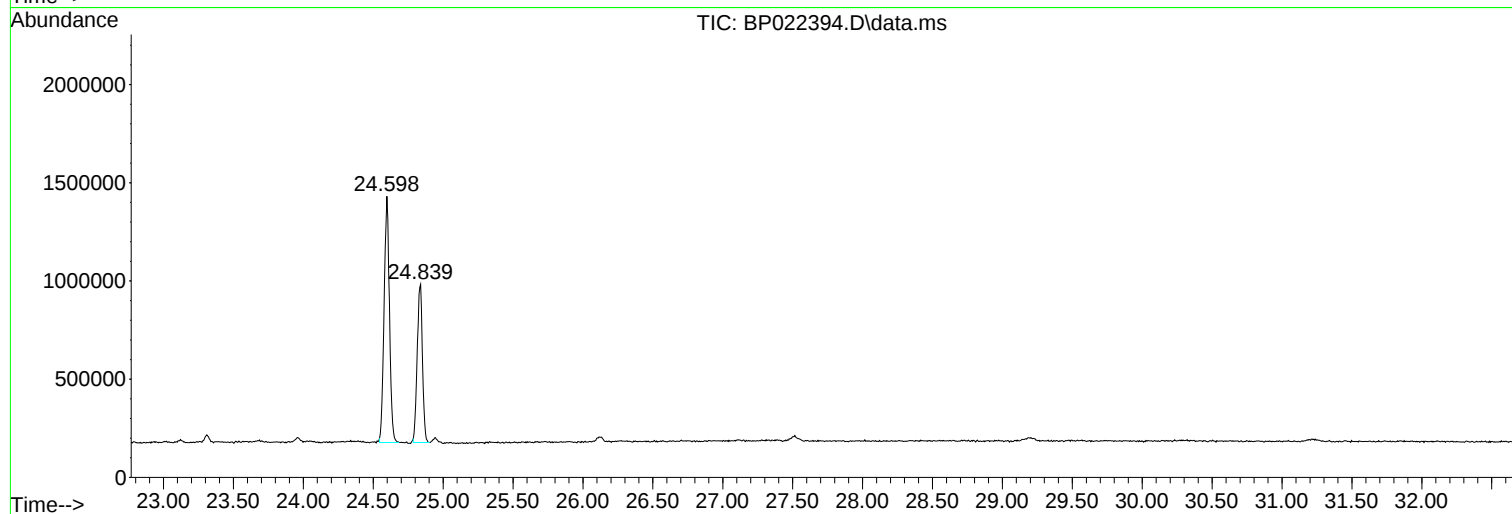
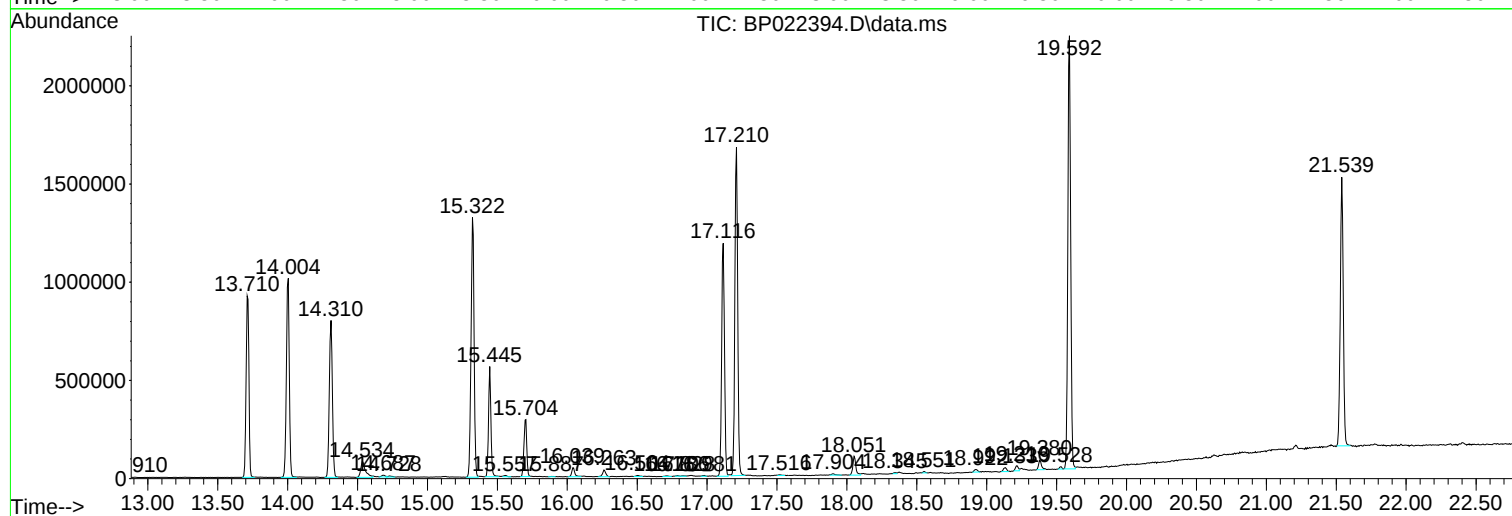
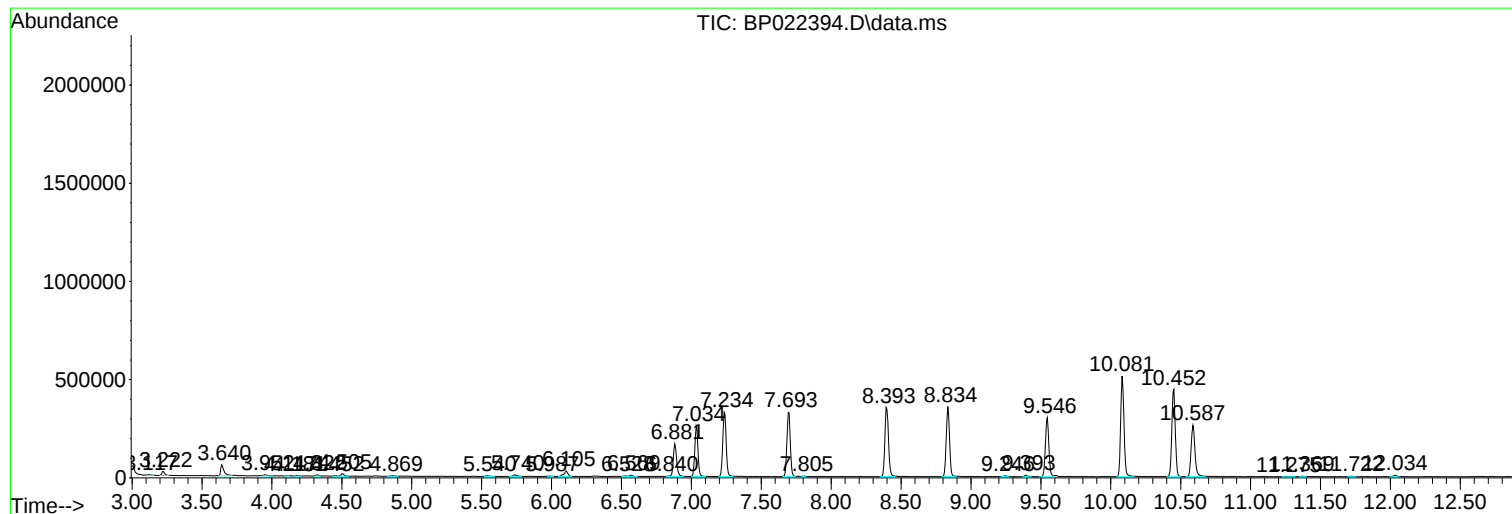
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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



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Misc :
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Instrument :
BNA_P
ClientSampleId :
EOAG4

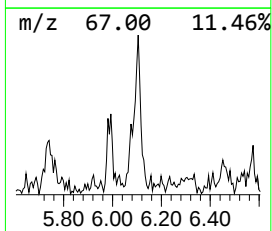
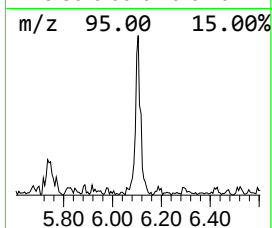
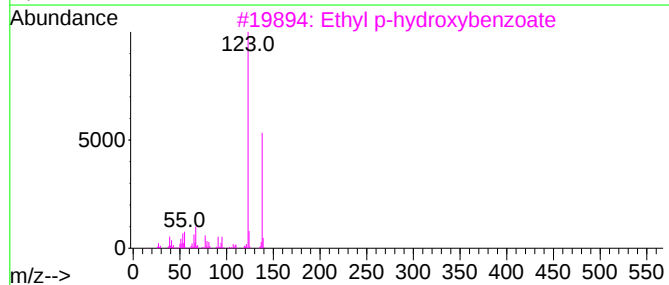
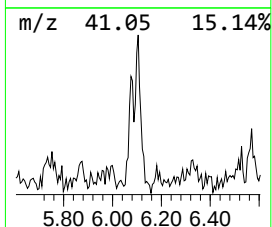
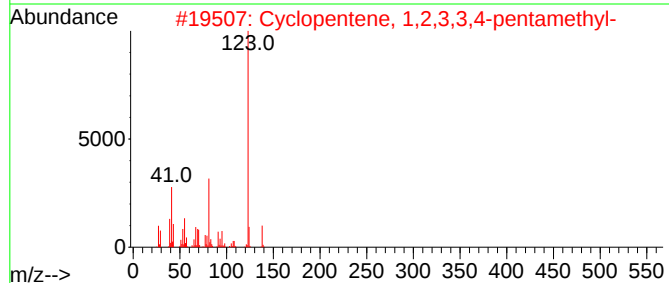
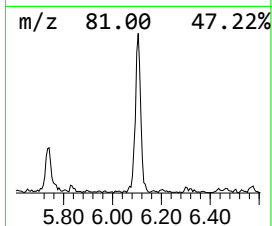
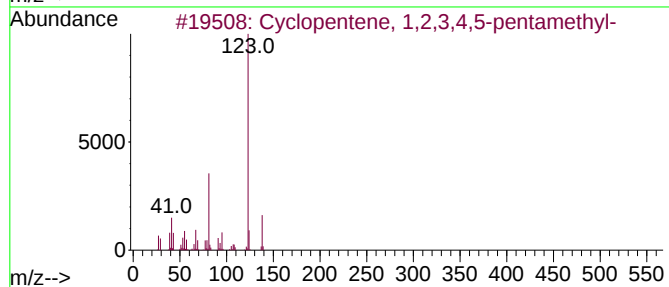
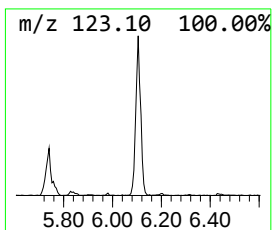
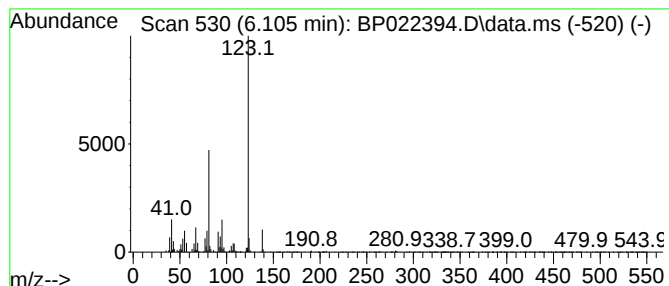
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Cyclopentene, 1,2,3,4,5-pen... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.105	2.03 ng/ul	55839	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentene, 1,2,3,4,5-pentamet...	138	C10H18	1000154-28-6	91
2			Cyclopentene, 1,2,3,3,4-pentamet...	138	C10H18	197390-29-7	81
3			Ethyl p-hydroxybenzoate	138	C8H10O2	002349-70-4	64
4			Phorone	138	C9H14O	000504-20-1	53
5			4-Ethylcatechol	138	C8H10O2	001124-39-6	50



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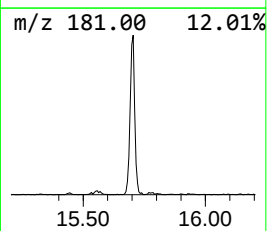
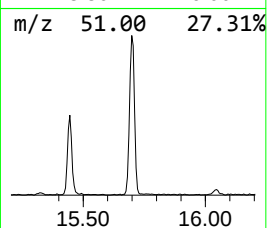
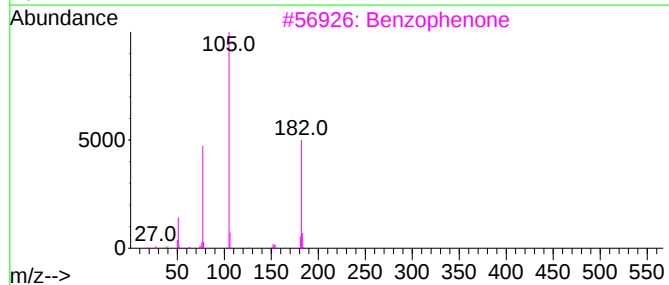
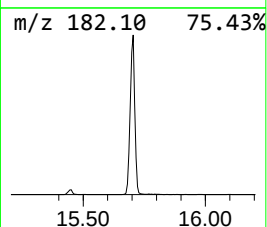
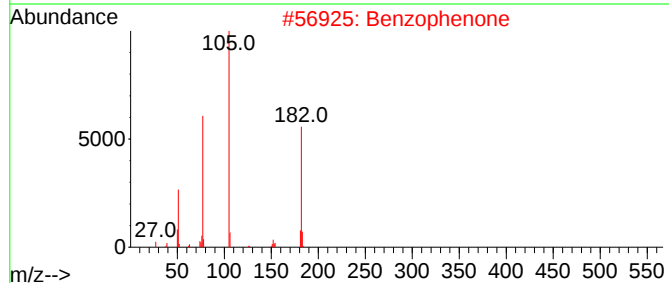
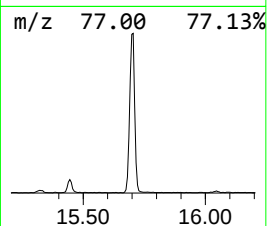
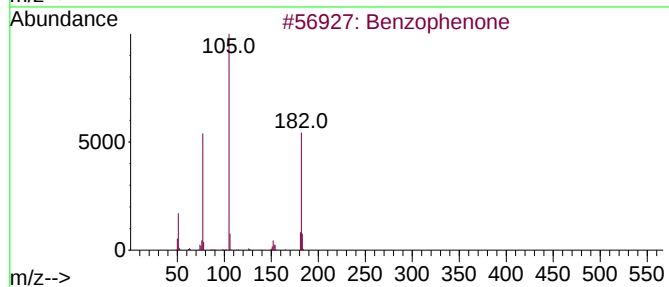
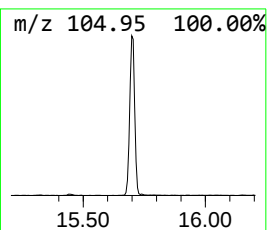
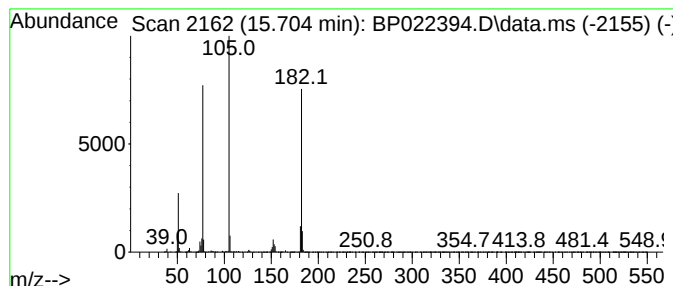
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzophenone Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.704	6.84 ng/ul	407762	Acenaphthene-d10	14.310

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	91
4			Benzophenone	182	C13H10O	000119-61-9	81
5			1,2,4,5-Tetroxane, 3,3,6,6-tetra...	396	C26H20O4	016204-36-7	74



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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Cyclopentene, 1...	6.105	2.0	ng/ul	55839	1	7.699	548957	20.0
Benzophenone	15.704	6.8	ng/ul	407762	3	14.310	1191780	20.0