

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111423\
 Data File : BP018166.D
 Acq On : 14 Nov 2023 21:14
 Operator : MA/JU
 Sample : 05337-12
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCDJ7

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-BP111023.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP018166.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.116	3	5	10	rVB	71833	70175	2.62%	0.310%
2	3.163	10	13	17	rVB	13477	17748	0.66%	0.078%
3	3.622	89	91	103	rBV	17133	50316	1.88%	0.222%
4	3.899	136	138	145	rVB7	6483	8003	0.30%	0.035%
5	4.010	153	157	158	rBV4	2197	2849	0.11%	0.013%
6	4.181	184	186	189	rBV3	2043	2765	0.10%	0.012%
7	4.416	222	226	232	rVB8	8476	13417	0.50%	0.059%
8	5.040	328	332	336	rBV7	2240	4387	0.16%	0.019%
9	5.769	454	456	462	rBV7	1919	3226	0.12%	0.014%
10	6.663	603	608	610	rBV6	2894	4357	0.16%	0.019%
11	6.857	639	641	646	rVB6	1766	2955	0.11%	0.013%
12	6.963	655	659	671	rBV	37159	81304	3.04%	0.359%
13	7.140	683	689	703	rBV	238854	391316	14.63%	1.729%
14	7.310	711	718	731	rBV	223442	397121	14.85%	1.755%
15	7.663	774	778	781	rVB6	1745	2701	0.10%	0.012%
16	7.787	793	799	816	rBV	304623	519453	19.43%	2.296%
17	8.516	916	923	941	rBV	139977	294450	11.01%	1.301%
18	8.992	998	1004	1017	rBV	297800	541427	20.25%	2.393%
19	9.163	1032	1033	1039	rVB6	3855	3570	0.13%	0.016%
20	9.710	1119	1126	1140	rBV	233891	465408	17.40%	2.057%
21	10.234	1208	1215	1240	rBV	256893	599990	22.44%	2.652%
22	10.604	1270	1278	1287	rBV	437592	719434	26.90%	3.180%
23	10.786	1303	1309	1333	rBV	234187	567748	21.23%	2.509%
24	10.934	1333	1334	1341	rVV7	6107	9522	0.36%	0.042%
25	11.651	1452	1456	1458	rVV5	2415	4641	0.17%	0.021%
26	12.239	1551	1556	1558	rBV6	3306	4694	0.18%	0.021%
27	12.651	1624	1626	1632	rVB7	2233	4000	0.15%	0.018%
28	12.710	1632	1636	1639	rBV6	1577	2699	0.10%	0.012%
29	13.304	1730	1737	1743	rBV2	18962	32068	1.20%	0.142%
30	13.680	1797	1801	1805	rBV7	1887	3109	0.12%	0.014%
31	13.886	1826	1836	1850	rBV	804289	1171948	43.83%	5.179%
32	14.163	1877	1883	1897	rBV	760546	1152585	43.10%	5.094%
33	14.469	1928	1935	1947	rBV	619322	934944	34.96%	4.132%
34	14.792	1984	1990	1995	rBV9	11432	28680	1.07%	0.127%
35	15.469	2097	2105	2114	rBV	1112278	1572380	58.80%	6.949%
36	15.627	2126	2132	2144	rBV	327817	536623	20.07%	2.372%

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37	15.886	2170	2176	2180	rBV2	12729	16509	0.62%	0.073%
38	15.927	2180	2183	2186	rVV5	7229	9583	0.36%	0.042%
39	16.186	2223	2227	2231	rBV	26744	34190	1.28%	0.151%
40	16.280	2236	2243	2247	rBV	82635	111124	4.16%	0.491%
41	16.339	2247	2253	2259	rVV2	94832	196787	7.36%	0.870%
42	16.404	2259	2264	2268	rVV3	81554	142946	5.35%	0.632%
43	16.445	2268	2271	2276	rVV	51173	68389	2.56%	0.302%
44	16.498	2276	2280	2282	rVV2	27835	39276	1.47%	0.174%
45	16.521	2282	2284	2287	rVV	23713	35319	1.32%	0.156%
46	16.551	2287	2289	2292	rVV	25447	26519	0.99%	0.117%
47	16.604	2292	2298	2305	rVV5	58252	114239	4.27%	0.505%
48	16.674	2305	2310	2315	rVV	76466	118243	4.42%	0.523%
49	16.716	2315	2317	2319	rVV2	22207	25332	0.95%	0.112%
50	16.745	2319	2322	2330	rVB2	36647	54320	2.03%	0.240%
51	16.892	2341	2347	2348	rBV6	4535	6765	0.25%	0.030%
52	16.916	2348	2351	2355	rVB5	3367	5811	0.22%	0.026%
53	17.004	2360	2366	2371	rVV7	4255	9378	0.35%	0.041%
54	17.051	2371	2374	2377	rVV4	3669	4725	0.18%	0.021%
55	17.086	2377	2380	2383	rVV5	3716	5036	0.19%	0.022%
56	17.116	2383	2385	2390	rVB6	2094	2694	0.10%	0.012%
57	17.198	2396	2399	2402	rBV4	2012	3274	0.12%	0.014%
58	17.245	2402	2407	2418	rVV	838515	1191291	44.55%	5.265%
59	17.351	2418	2425	2442	rBV	1338383	1972255	73.75%	8.716%
60	17.498	2447	2450	2452	rVV3	3281	3566	0.13%	0.016%
61	17.580	2462	2464	2468	rVB5	2610	3058	0.11%	0.014%
62	17.880	2511	2515	2522	rBV4	40805	56657	2.12%	0.250%
63	18.098	2548	2552	2557	rBV8	19443	31884	1.19%	0.141%
64	18.186	2565	2567	2570	rVB3	3101	3396	0.13%	0.015%
65	18.310	2583	2588	2594	rBV	12935	22298	0.83%	0.099%
66	19.174	2731	2735	2741	rBV3	19038	28575	1.07%	0.126%
67	19.251	2741	2748	2754	rBV3	19186	42605	1.59%	0.188%
68	19.351	2761	2765	2771	rBV7	12100	18253	0.68%	0.081%
69	19.480	2784	2787	2791	rVB6	9791	9337	0.35%	0.041%
70	19.604	2802	2808	2818	rBV	1915390	2492206	93.20%	11.014%
71	21.380	3105	3110	3122	rBV	1002638	1388982	51.94%	6.139%
72	23.686	3494	3502	3512	rBV2	1314667	2674095	100.00%	11.818%
73	23.845	3522	3529	3540	rVB	669352	1435910	53.70%	6.346%

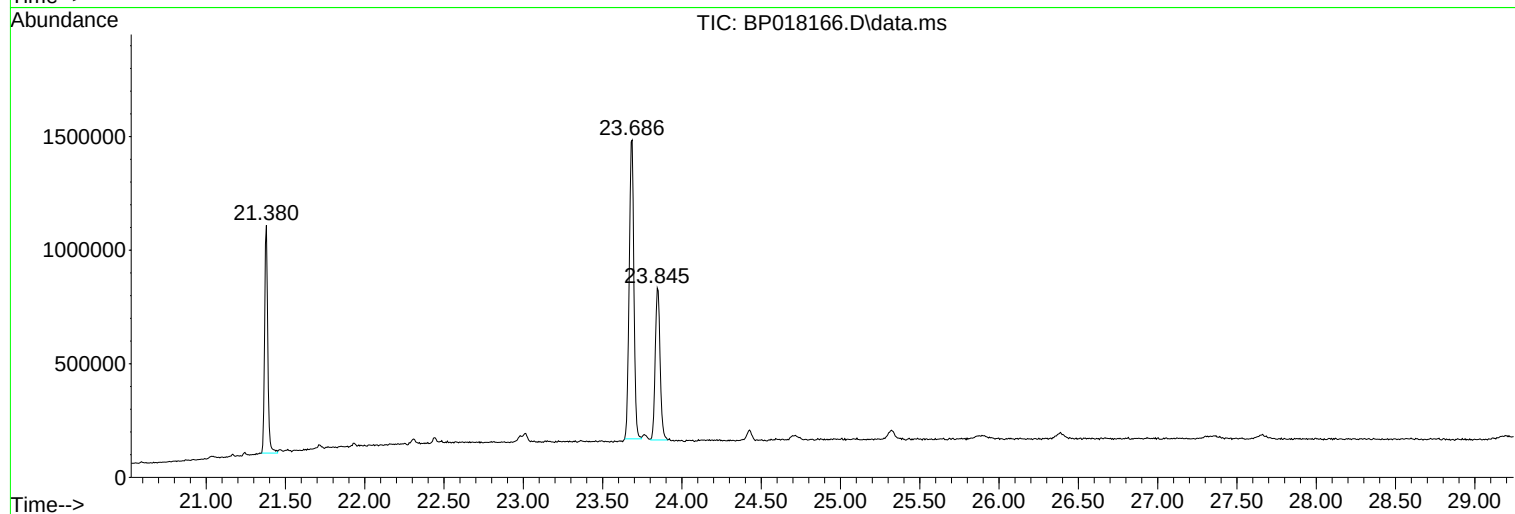
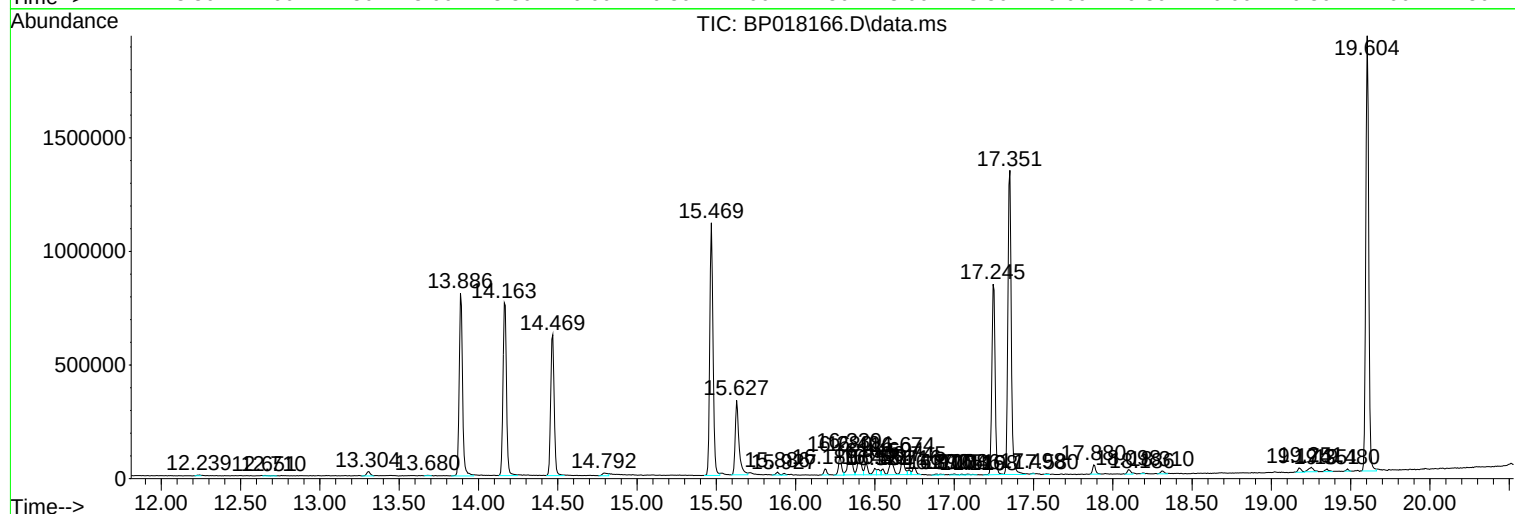
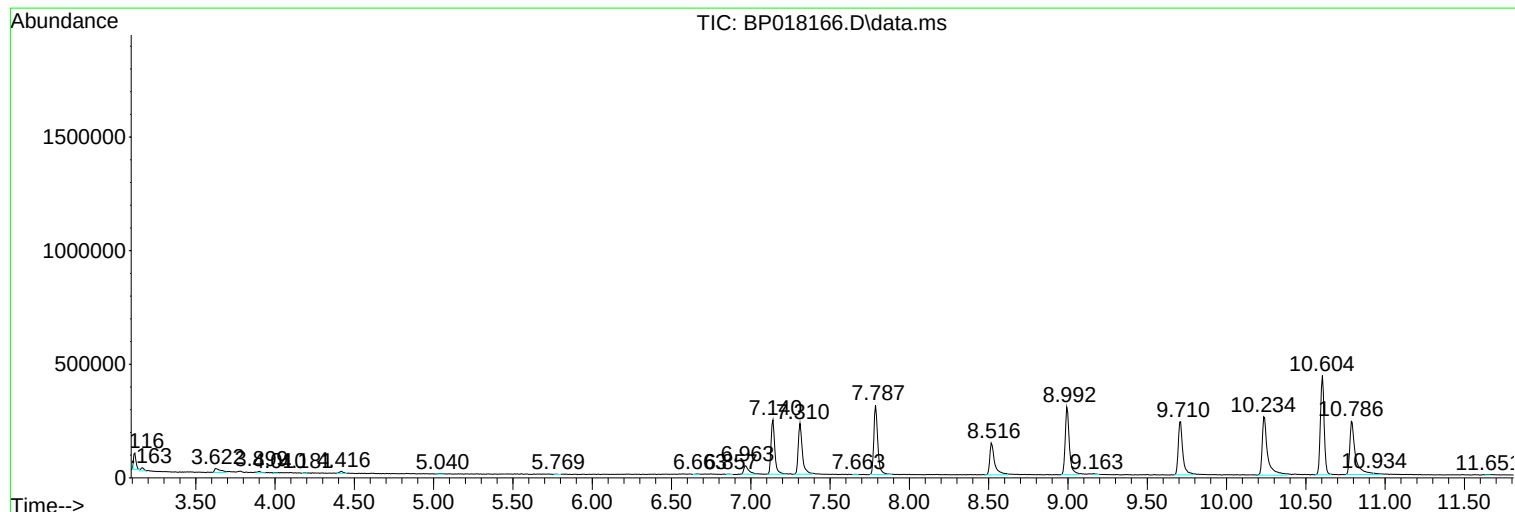
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
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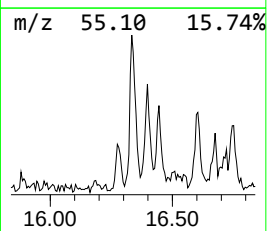
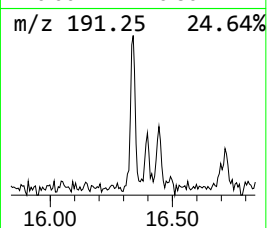
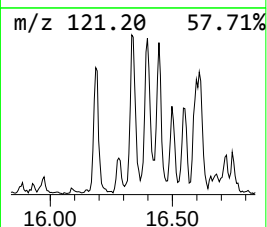
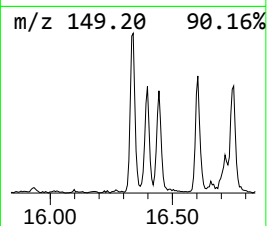
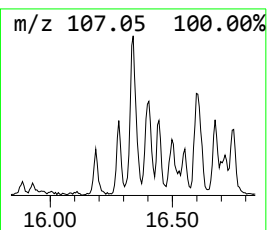
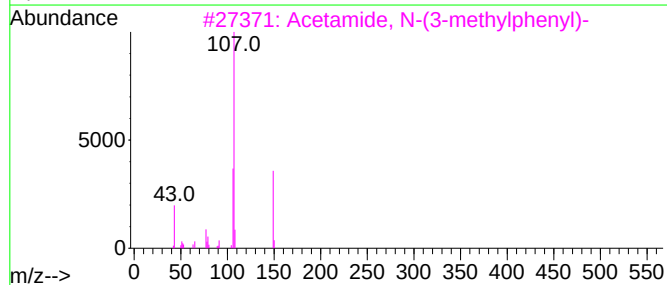
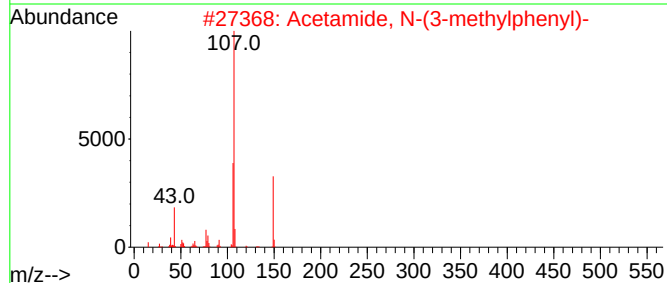
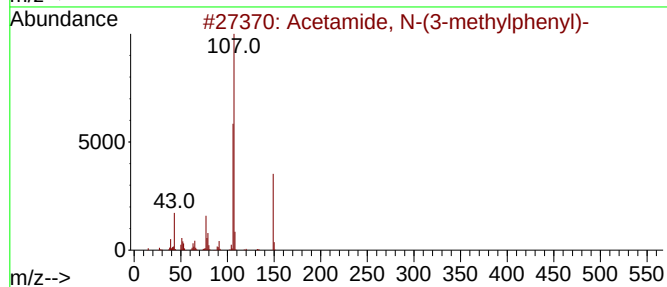
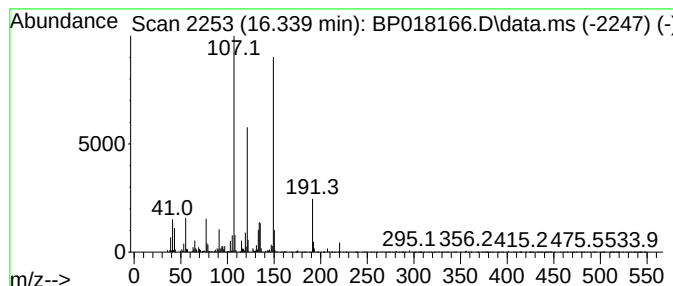
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Acetamide, N-(3-methylphenyl)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.339	3.30 ng/ul	196787	Phenanthrene-d10	17.245

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	58
2			Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	58
3			Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	46
4			Pyridine-3-carbonitrile, 1,2-dih...	191	C9H9N3O2	220098-92-0	43
5			3,4-(methylenedioxy)-phenylbutan...	192	C11H12O3	1000378-99-2	38



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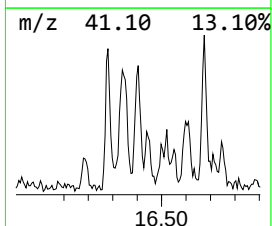
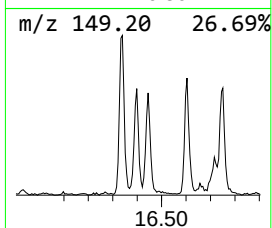
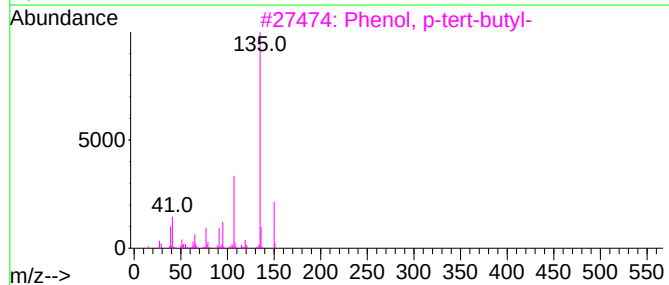
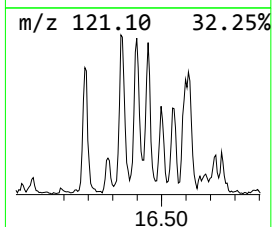
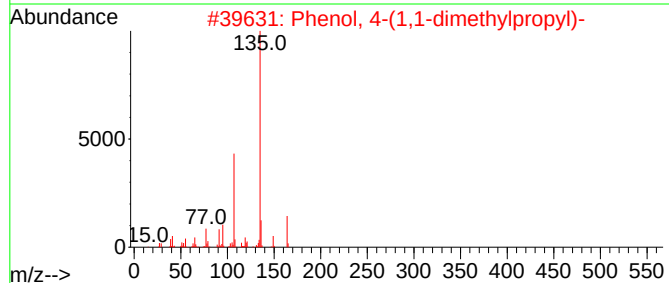
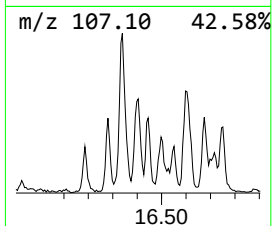
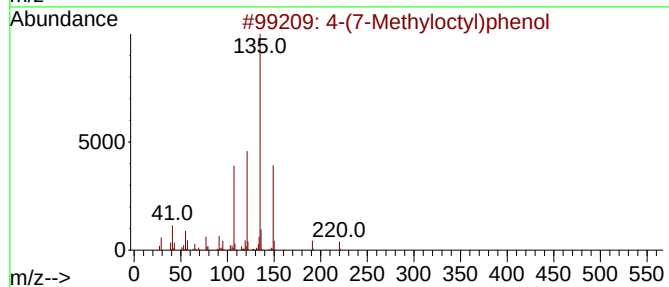
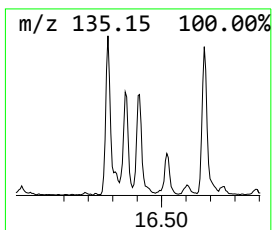
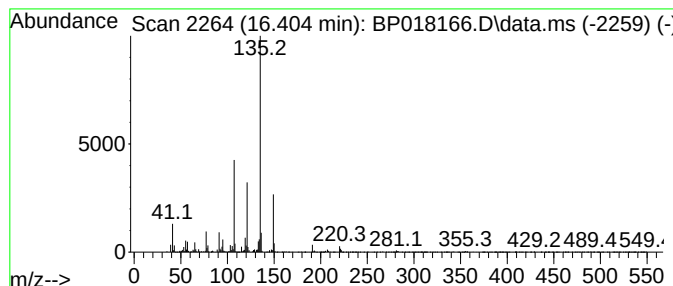
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 4-(7-Methyloctyl)phenol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.404	2.40 ng/ul	142946	Phenanthrene-d10	17.245

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-(7-Methyloctyl)phenol	220	C15H24O	024518-48-7	91
2			Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	64
3			Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	59
4			meso-Hexestrol, 0,0'-bis(n-propy...	442	C26H34O6	1000487-65-0	59
5			((1,2-Diethylethylene)bis(p-phen...	354	C22H26O4	004547-76-6	59



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Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Acetamide, N-(3...	16.339	3.3	ng/u1	196787	4	17.245	1191290	20.0
4-(7-Methylocty...	16.404	2.4	ng/u1	142946	4	17.245	1191290	20.0