

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
 Data File : BP018254.D
 Acq On : 18 Nov 2023 10:59
 Operator : MA/JU
 Sample : 05370-06
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCDQ2

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: BP018254.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.164	9	13	19	rVB	14756	19832	0.76%	0.089%
2	3.899	135	138	142	rBV5	3863	6114	0.24%	0.028%
3	4.263	196	200	206	rVB2	4994	8446	0.33%	0.038%
4	4.411	221	225	230	rVB4	9020	13642	0.53%	0.062%
5	5.587	421	425	428	rVB6	2208	3176	0.12%	0.014%
6	6.152	519	521	525	rBV4	1946	2899	0.11%	0.013%
7	6.575	588	593	595	rBV6	1591	2603	0.10%	0.012%
8	6.734	616	620	625	rVB2	9406	13402	0.52%	0.060%
9	6.805	629	632	635	rBV4	2446	2818	0.11%	0.013%
10	6.952	650	657	670	rBV	95190	198908	7.66%	0.897%
11	7.128	682	687	699	rBV	253682	399592	15.40%	1.803%
12	7.299	710	716	735	rBV	333728	574349	22.13%	2.592%
13	7.775	791	797	811	rBV	219596	404066	15.57%	1.823%
14	8.169	860	864	867	rBV6	1777	2822	0.11%	0.013%
15	8.504	915	921	943	rBV	244386	504256	19.43%	2.275%
16	8.987	996	1003	1017	rBV	325240	601018	23.16%	2.712%
17	9.151	1027	1031	1037	rVB2	16605	25582	0.99%	0.115%
18	9.699	1117	1124	1143	rBV	293214	573440	22.10%	2.587%
19	10.169	1199	1204	1208	rBV6	12836	20316	0.78%	0.092%
20	10.228	1208	1214	1244	rVV2	278077	718972	27.71%	3.244%
21	10.469	1250	1255	1261	rVB10	6750	11634	0.45%	0.052%
22	10.593	1269	1276	1285	rBV	327192	540707	20.84%	2.440%
23	11.393	1407	1412	1417	rBV9	4088	7864	0.30%	0.035%
24	11.534	1431	1436	1438	rBV6	2715	5147	0.20%	0.023%
25	11.645	1450	1455	1460	rBV2	6878	10637	0.41%	0.048%
26	12.240	1549	1556	1562	rBV2	3276	9925	0.38%	0.045%
27	13.028	1686	1690	1694	rBV6	2615	3640	0.14%	0.016%
28	13.298	1730	1736	1743	rBV7	8416	17401	0.67%	0.079%
29	13.669	1794	1799	1800	rBV5	2589	3328	0.13%	0.015%
30	13.787	1817	1819	1824	rVB6	1865	2777	0.11%	0.013%
31	13.881	1829	1835	1847	rVV	962624	1360213	52.42%	6.137%
32	14.157	1876	1882	1896	rBV	938787	1306790	50.36%	5.896%
33	14.392	1919	1922	1924	rVB4	2259	2680	0.10%	0.012%
34	14.416	1924	1926	1927	rBV2	2918	2827	0.11%	0.013%
35	14.457	1927	1933	1946	rVB	527409	761131	29.33%	3.434%
36	14.769	1979	1986	2001	rBV4	35443	145796	5.62%	0.658%

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37	15.104	2041	2043	2046	rVV3	3373	2756	0.11%	0.012%
38	15.145	2048	2050	2053	rVB4	3161	3161	0.12%	0.014%
39	15.245	2064	2067	2071	rVB6	2386	3590	0.14%	0.016%
40	15.463	2097	2104	2118	rBV	1268678	1806981	69.63%	8.153%
41	15.575	2118	2123	2126	rBV6	7544	11010	0.42%	0.050%
42	15.622	2126	2131	2143	rBV	405437	668084	25.74%	3.014%
43	15.881	2170	2175	2178	rBV2	6738	11080	0.43%	0.050%
44	15.922	2178	2182	2186	rVB3	5821	7362	0.28%	0.033%
45	16.181	2221	2226	2231	rBV	19606	26737	1.03%	0.121%
46	16.275	2236	2242	2246	rBV	62892	86398	3.33%	0.390%
47	16.333	2247	2252	2258	rVV2	68931	135288	5.21%	0.610%
48	16.398	2258	2263	2267	rVV3	55596	92075	3.55%	0.415%
49	16.439	2267	2270	2274	rVB2	34720	42653	1.64%	0.192%
50	16.492	2274	2279	2281	rBV3	15462	22533	0.87%	0.102%
51	16.539	2285	2287	2291	rVB	16067	19118	0.74%	0.086%
52	16.592	2291	2296	2305	rBV5	39143	75920	2.93%	0.343%
53	16.669	2305	2309	2313	rBV	61499	82214	3.17%	0.371%
54	16.745	2318	2322	2330	rVB	27693	42804	1.65%	0.193%
55	16.886	2341	2346	2347	rBV5	2200	3847	0.15%	0.017%
56	16.992	2362	2364	2368	rVB4	2669	3395	0.13%	0.015%
57	17.239	2400	2406	2417	rBV	723346	984516	37.94%	4.442%
58	17.339	2417	2423	2444	rVV	1352667	2019073	77.81%	9.110%
59	17.510	2451	2452	2458	rVB6	4418	4196	0.16%	0.019%
60	17.575	2458	2463	2467	rBV8	3527	6718	0.26%	0.030%
61	17.875	2510	2514	2520	rBV3	33120	39967	1.54%	0.180%
62	18.057	2542	2545	2548	rBV5	2652	3638	0.14%	0.016%
63	18.098	2548	2552	2561	rVV2	23814	40925	1.58%	0.185%
64	18.192	2564	2568	2571	rVB3	4300	6388	0.25%	0.029%
65	18.304	2581	2587	2600	rBV	133130	219024	8.44%	0.988%
66	18.680	2648	2651	2653	rBV4	2968	3902	0.15%	0.018%
67	18.810	2670	2673	2677	rVB6	7544	9207	0.35%	0.042%
68	19.174	2731	2735	2739	rBV3	18372	21980	0.85%	0.099%
69	19.251	2743	2748	2757	rVB	16551	27874	1.07%	0.126%
70	19.345	2760	2764	2771	rBV5	23946	38527	1.48%	0.174%
71	19.604	2802	2808	2819	rBV	1823148	2476103	95.42%	11.172%
72	21.380	3105	3110	3121	rBV	783985	1119970	43.16%	5.053%
73	23.686	3495	3502	3515	rBV	1319177	2595030	100.00%	11.709%
74	23.851	3524	3530	3540	rVB	519331	1109948	42.77%	5.008%

Sum of corrected areas: 22162742

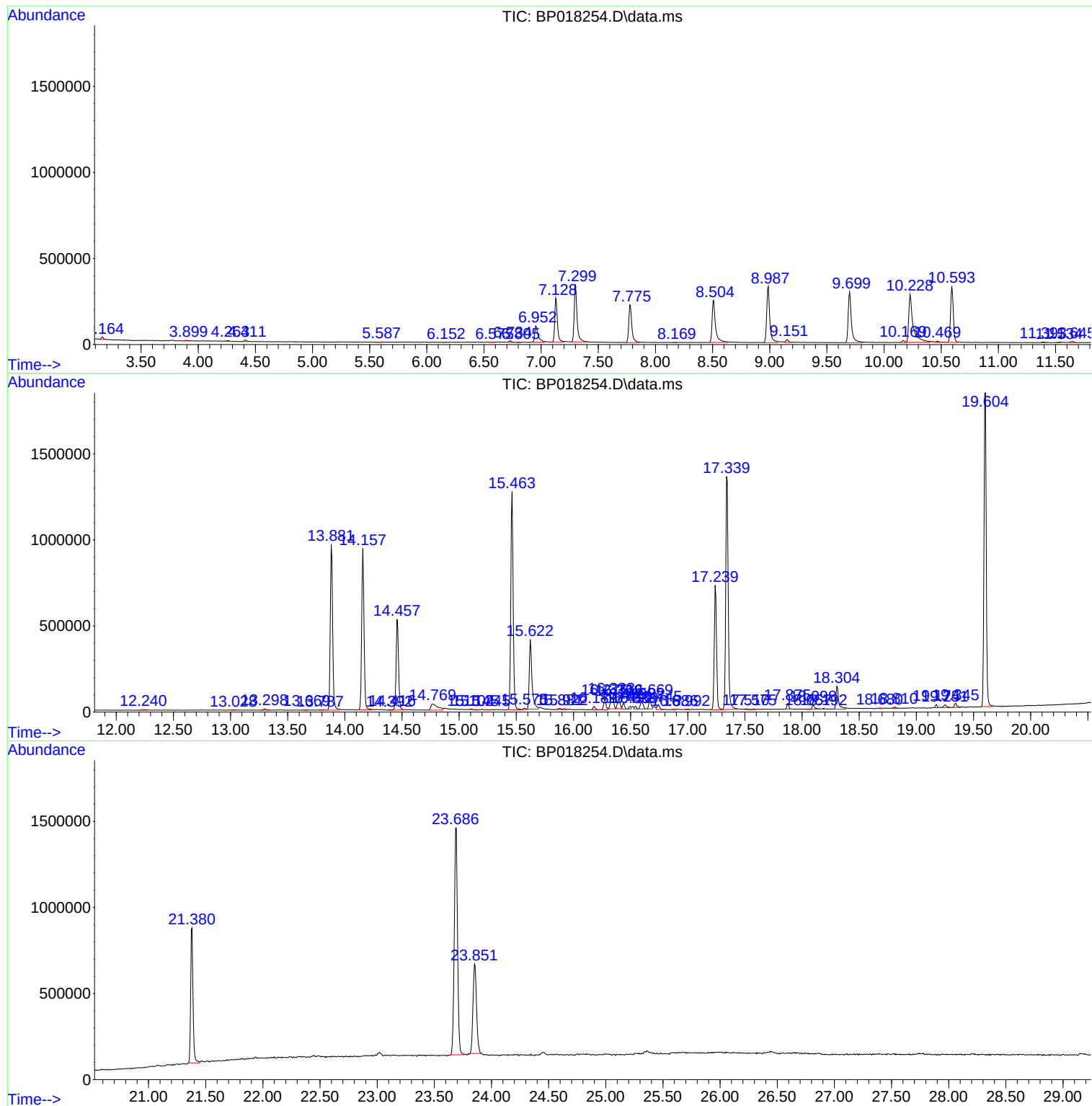
Instrument :
BNA_P
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GCDQ2

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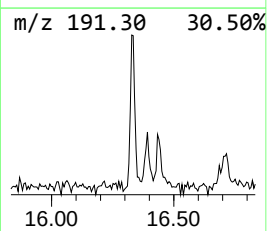
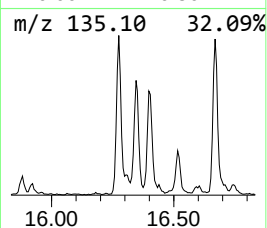
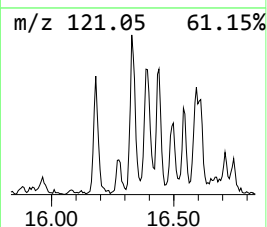
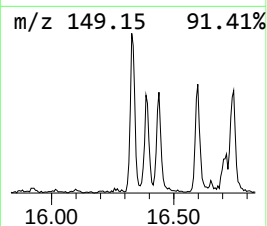
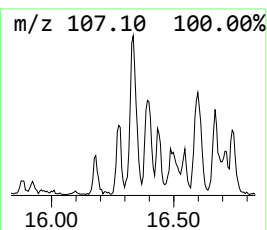
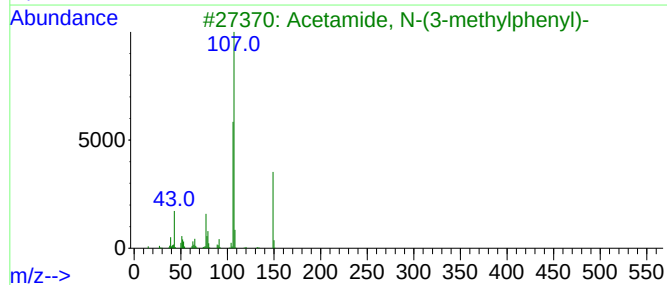
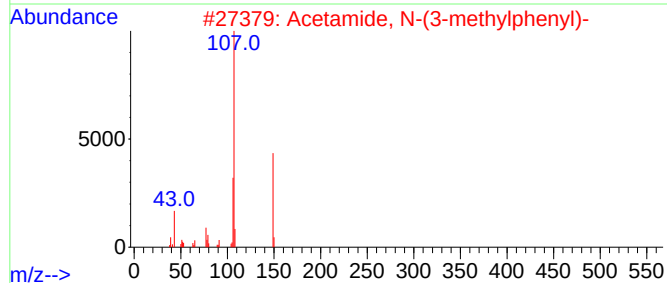
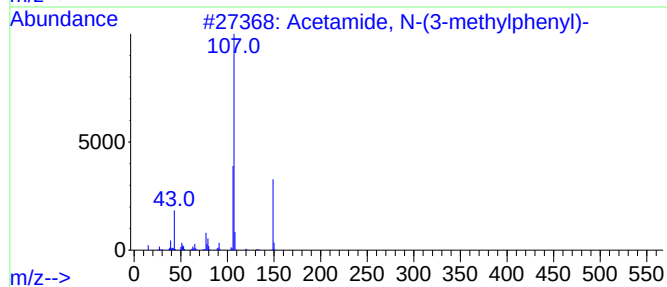
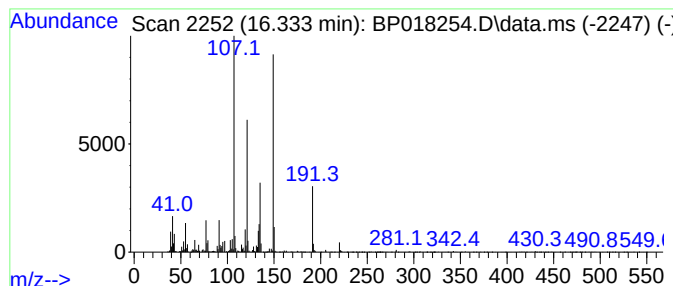
Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Acetamide, N-(3-methylphenyl)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.333	2.75 ng/ul	135288	Phenanthrene-d10	17.239	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	64
2	Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	58
3	Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	58
4	Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	52
5	Acetamide, N-(2-methylphenyl)-	149	C9H11NO	000120-66-1	52



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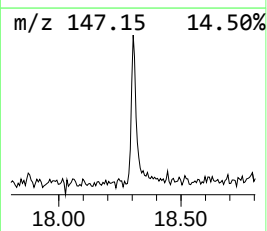
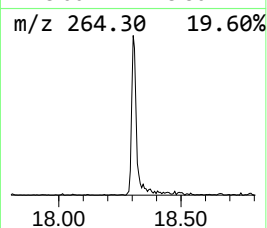
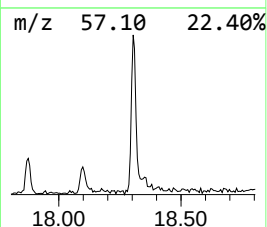
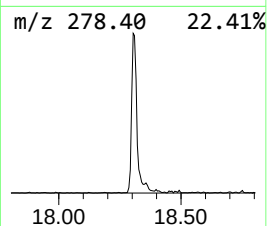
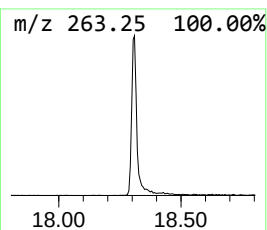
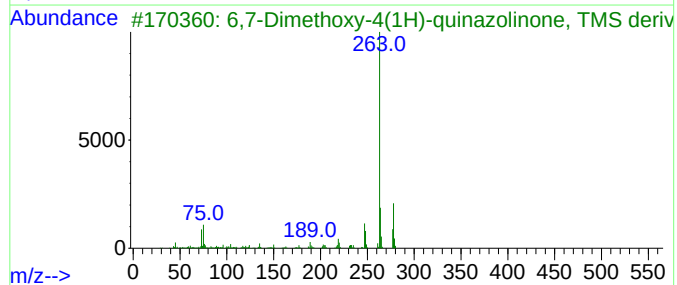
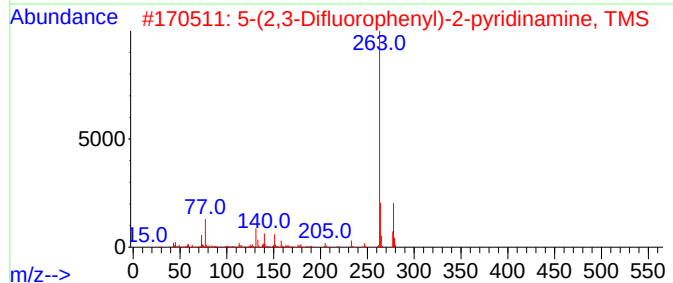
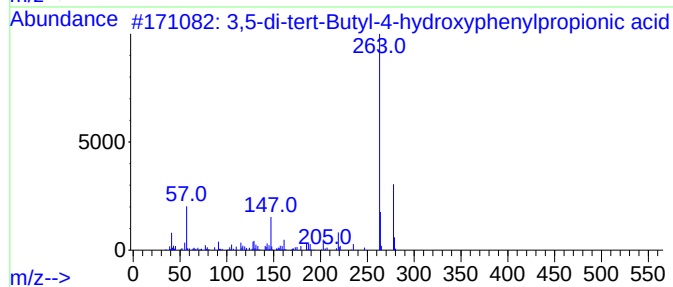
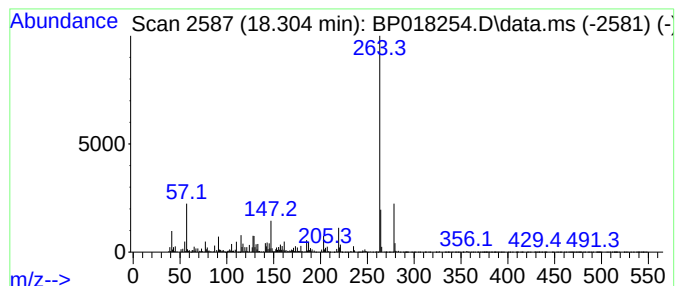
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 3,5-di-tert-Butyl-4-hydroxy... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.304	4.45 ng/ul	219024	Phenanthrene-d10	17.239

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3,5-di-tert-Butyl-4-hydroxypheny...	278	C17H26O3	020170-32-5	94
2			5-(2,3-Difluorophenyl)-2-pyridin...	278	C14H16F2N2Si	1000504-49-4	81
3			6,7-Dimethoxy-4(1H)-quinazolinon...	278	C13H18N2O3Si	1000479-65-5	72
4			2,6-Bis(tert-butyl)phenol, TMS d...	278	C17H30O5Si	010416-73-6	64
5			3,5-di-tert-Butyl-4-hydroxypheny...	278	C17H26O3	020170-32-5	58



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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...	16.333	2.8	ng/ul	135288	4	17.239	984516	20.0
3,5-di-tert-But...	18.304	4.5	ng/ul	219024	4	17.239	984516	20.0