

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111324\
 Data File : BG063398.D
 Acq On : 14 Nov 2024 19:19
 Operator : RC/JU
 Sample : P4715-13
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0AK6

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_G\Methods\SFAM-EPA-BG110624.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG063398.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.304	34	37	44	rVB4	26677	39092	0.86%	0.100%
2	3.792	118	120	123	rBV3	7083	6754	0.15%	0.017%
3	4.403	221	224	228	rVB5	11997	12412	0.27%	0.032%
4	4.544	246	248	255	rVB6	6514	10045	0.22%	0.026%
5	5.073	336	338	342	rBV4	5424	6733	0.15%	0.017%
6	5.690	439	443	447	rVB2	10330	16009	0.35%	0.041%
7	7.006	658	667	677	rBV2	205501	370545	8.18%	0.947%
8	7.165	687	694	704	rBV	395892	653255	14.43%	1.669%
9	7.364	721	728	737	rBV	489658	842775	18.61%	2.153%
10	7.828	800	807	814	rBV	396527	690725	15.25%	1.765%
11	7.893	814	818	821	rVV2	20684	29295	0.65%	0.075%
12	8.539	919	928	940	rBV	540192	915183	20.21%	2.338%
13	8.657	945	948	952	rVB4	6558	9701	0.21%	0.025%
14	8.927	988	994	998	rBV	62085	110196	2.43%	0.282%
15	8.986	998	1004	1015	rVB	510833	882466	19.49%	2.254%
16	9.145	1029	1031	1037	rVB7	6880	11893	0.26%	0.030%
17	9.709	1119	1127	1134	rBV	464338	821460	18.14%	2.099%
18	10.161	1200	1204	1207	rVB4	4620	7251	0.16%	0.019%
19	10.249	1211	1219	1229	rBV	666967	1224326	27.04%	3.128%
20	10.478	1257	1258	1263	rVB5	10010	11889	0.26%	0.030%
21	10.537	1263	1268	1270	rBV4	11435	13851	0.31%	0.035%
22	10.619	1274	1282	1289	rVV	571964	999993	22.08%	2.555%
23	10.678	1289	1292	1297	rVB5	9198	15637	0.35%	0.040%
24	10.754	1298	1305	1317	rBV	304859	578400	12.77%	1.478%
25	11.166	1371	1375	1379	rVB2	5951	6261	0.14%	0.016%
26	11.495	1428	1431	1433	rBV3	7194	8914	0.20%	0.023%
27	11.806	1479	1484	1486	rBV2	4875	7031	0.16%	0.018%
28	12.212	1547	1553	1560	rBV4	12225	25815	0.57%	0.066%
29	12.523	1602	1606	1609	rBV3	6647	7931	0.18%	0.020%
30	12.805	1650	1654	1657	rBV4	5400	7761	0.17%	0.020%
31	13.069	1696	1699	1705	rBV4	13473	23065	0.51%	0.059%
32	13.387	1745	1753	1762	rBV2	71850	142817	3.15%	0.365%
33	13.604	1787	1790	1793	rBV3	4863	8378	0.19%	0.021%
34	13.874	1830	1836	1849	rVB	1372236	1984250	43.82%	5.069%
35	13.992	1853	1856	1859	rVB3	6336	7321	0.16%	0.019%
36	14.162	1876	1885	1893	rBV2	1439602	2232693	49.30%	5.704%

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37	14.333	1912	1914	1917	rBV4	5780	7165	0.16%	0.018%
38	14.468	1930	1937	1947	rBV	1046902	1618167	35.73%	4.134%
39	14.544	1947	1950	1956	rVB3	14275	24342	0.54%	0.062%
40	14.685	1965	1974	1980	rBV2	134645	264195	5.83%	0.675%
41	14.844	1997	2001	2005	rVB6	7562	10614	0.23%	0.027%
42	14.891	2005	2009	2016	rBV9	21464	41086	0.91%	0.105%
43	15.044	2029	2035	2037	rBV3	4105	6799	0.15%	0.017%
44	15.279	2071	2075	2081	rVB9	8565	16686	0.37%	0.043%
45	15.467	2100	2107	2116	rVB	2046749	2978906	65.78%	7.610%
46	15.590	2122	2128	2140	rBV	649198	1012155	22.35%	2.586%
47	15.831	2166	2169	2174	rBV5	6761	8859	0.20%	0.023%
48	15.907	2181	2182	2185	rBV	5003	6169	0.14%	0.016%
49	15.995	2195	2197	2201	rBV2	5519	7902	0.17%	0.020%
50	16.506	2282	2284	2290	rVB6	10576	14001	0.31%	0.036%
51	16.812	2334	2336	2340	rBV4	10189	11306	0.25%	0.029%
52	16.965	2360	2362	2364	rBV3	6207	6940	0.15%	0.018%
53	17.000	2364	2368	2370	rBV4	5102	7778	0.17%	0.020%
54	17.106	2380	2386	2388	rBV4	8533	8948	0.20%	0.023%
55	17.223	2395	2406	2413	rBV	1405606	2142200	47.30%	5.473%
56	17.317	2416	2422	2434	rVV	2426901	3586360	79.20%	9.162%
57	17.423	2439	2440	2446	rVB3	13851	11151	0.25%	0.028%
58	17.576	2463	2466	2467	rBV2	5836	6206	0.14%	0.016%
59	17.887	2512	2519	2525	rBV2	280828	394919	8.72%	1.009%
60	18.193	2566	2571	2572	rBV4	10748	10999	0.24%	0.028%
61	18.369	2598	2601	2605	rVB5	7192	10892	0.24%	0.028%
62	18.410	2605	2608	2614	rVB6	5503	9621	0.21%	0.025%
63	18.610	2639	2642	2645	rBV4	9497	13201	0.29%	0.034%
64	19.062	2716	2719	2722	rVB3	6113	7471	0.16%	0.019%
65	19.180	2735	2739	2745	rVB2	37591	54823	1.21%	0.140%
66	19.250	2747	2751	2759	rVB	32360	50023	1.10%	0.128%
67	19.620	2807	2814	2822	rBV	3215154	4528510	100.00%	11.569%
68	19.855	2851	2854	2857	rVB4	21062	24280	0.54%	0.062%
69	20.208	2911	2914	2917	rBV5	9152	11837	0.26%	0.030%
70	21.477	3123	3130	3139	rBV	1591629	2517264	55.59%	6.431%
71	22.646	3327	3329	3330	rBV2	12364	8456	0.19%	0.022%
72	24.303	3601	3611	3626	rVB	1689792	4321760	95.43%	11.041%
73	24.509	3635	3646	3657	rBV	980182	2640474	58.31%	6.745%

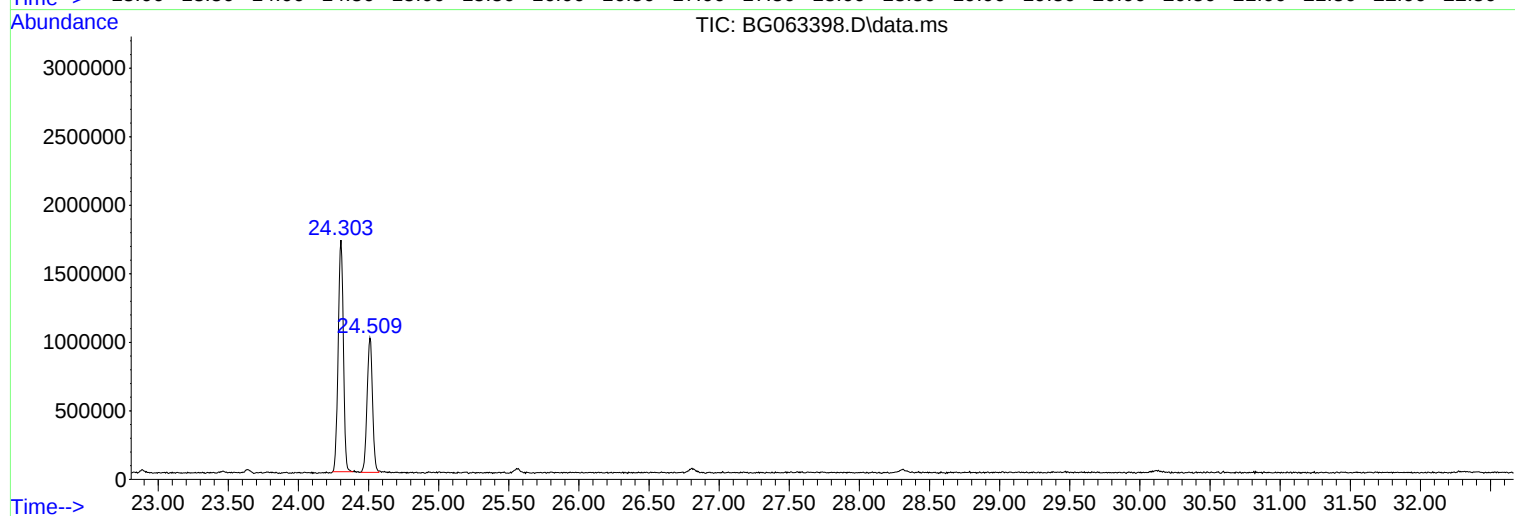
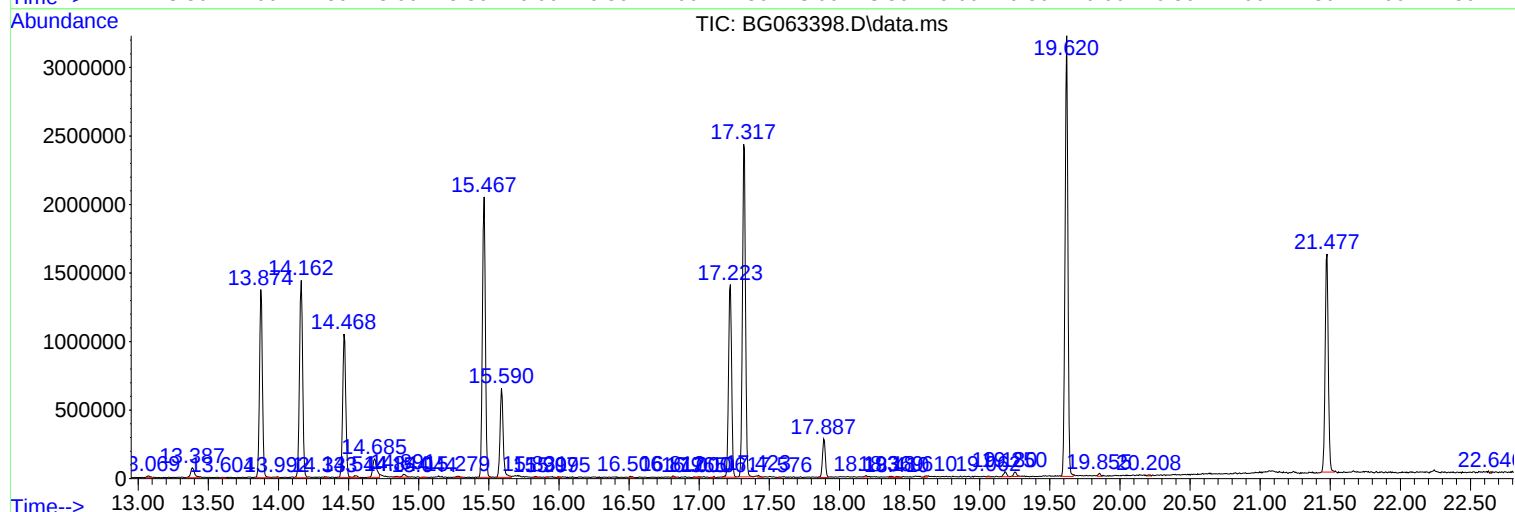
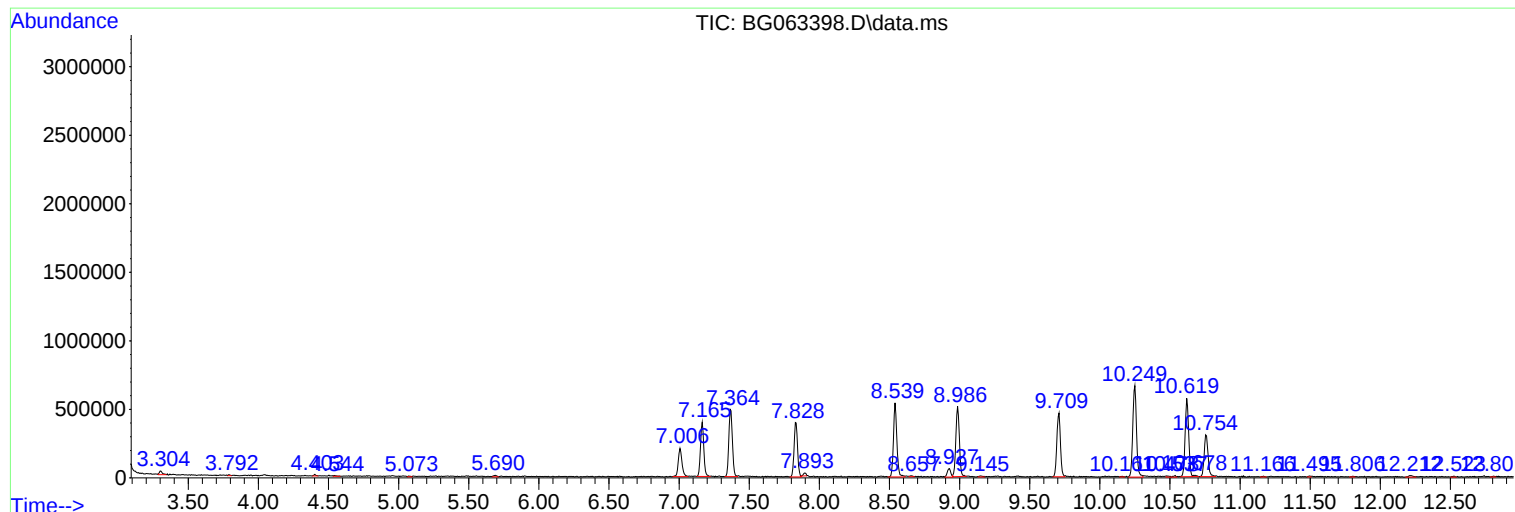
Sum of corrected areas: 39144588

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110624.MA.M
Quant Title : SVOA CALIBRATION

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



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Operator : RC/JU
Sample : P4715-13
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AK6

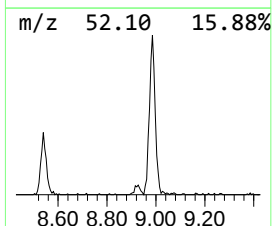
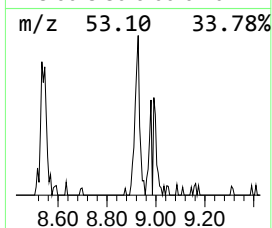
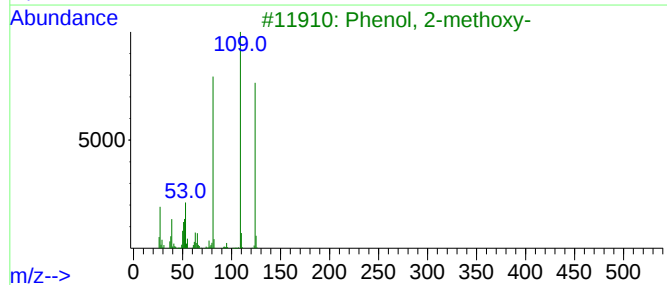
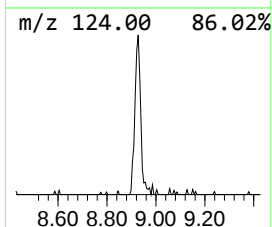
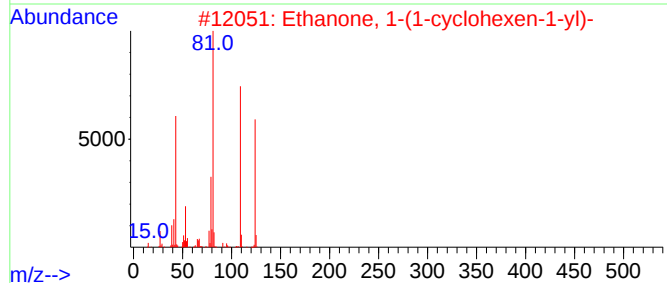
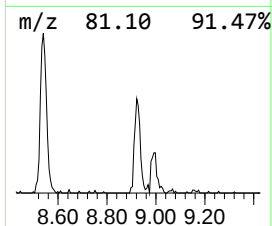
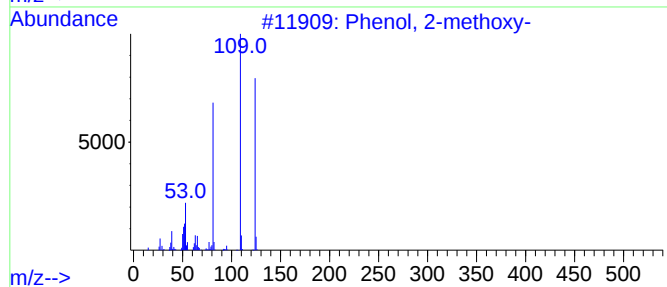
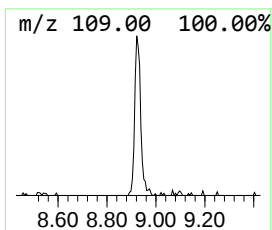
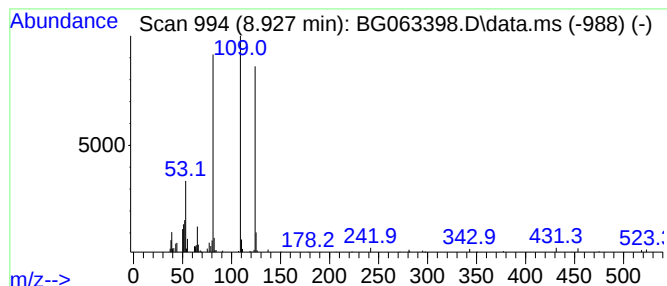
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110624.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Phenol, 2-methoxy- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.927	3.19 ng/ul	110196	1,4-Dichlorobenzene-d4	7.834

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	90
2			Ethanone, 1-(1-cyclohexen-1-yl)-	124	C8H12O	000932-66-1	90
3			Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	90
4			Mequinol	124	C7H8O2	000150-76-5	87
5			Mequinol	124	C7H8O2	000150-76-5	87



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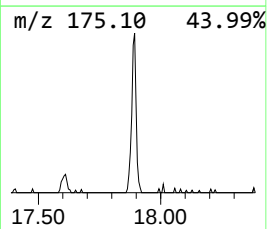
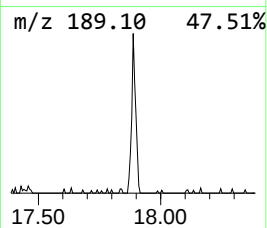
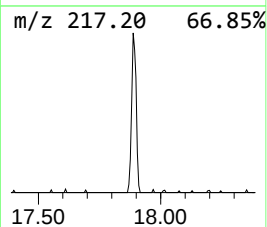
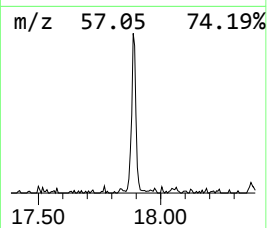
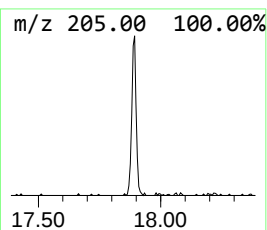
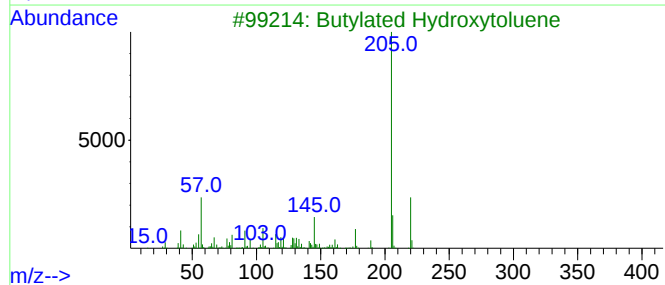
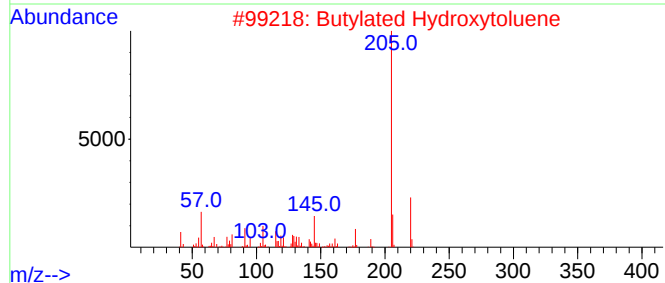
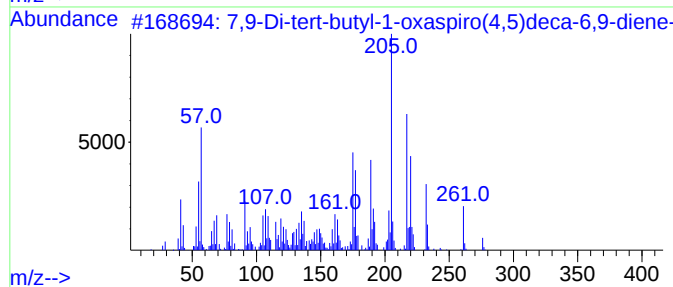
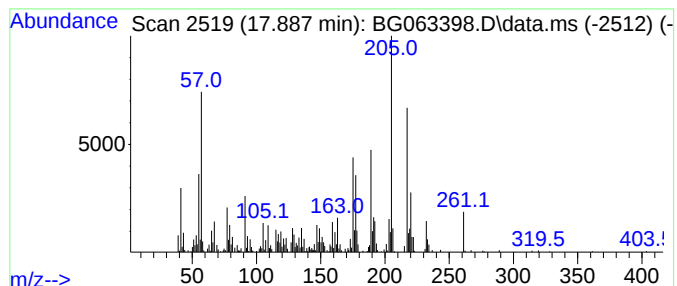
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 7,9-Di-tert-butyl-1-oxaspiro... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.887	3.69 ng/ul	394919	Phenanthrene-d10	17.223

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	95		
2	Butylated Hydroxytoluene	220	C15H24O	000128-37-0	55		
3	Butylated Hydroxytoluene	220	C15H24O	000128-37-0	42		
4	Butylated Hydroxytoluene	220	C15H24O	000128-37-0	42		
5	2,4-Diamino-6-nitroquinazoline	205	C8H7N5O2	007154-34-9	41		



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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
Phenol, 2-methoxy-	8.927	3.2	ng/ul	110196	1	7.834	690725	20.0
7,9-Di-tert-but...	17.887	3.7	ng/ul	394919	4	17.223	2142200	20.0