

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111324\
 Data File : BP023139.D
 Acq On : 15 Nov 2024 14:44
 Operator : RC/JU
 Sample : P4784-08
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E27N9

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-BP110924.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP023139.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.105	16	20	23	rBV2	9622	12899	0.43%	0.053%
2	3.134	23	25	34	rVB	14341	20989	0.70%	0.086%
3	3.558	93	97	112	rBV	52133	96395	3.20%	0.397%
4	3.858	144	148	153	rBV4	3112	4557	0.15%	0.019%
5	4.281	216	220	224	rBV4	2617	3925	0.13%	0.016%
6	6.716	629	634	639	rBV6	2550	4538	0.15%	0.019%
7	6.787	639	646	664	rBV	60859	147723	4.90%	0.608%
8	6.928	664	670	679	rBV	184468	290102	9.63%	1.194%
9	7.128	697	704	725	rBV	225996	405503	13.46%	1.669%
10	7.575	773	780	791	rBV	229558	385156	12.79%	1.585%
11	8.293	895	902	922	rBV	211013	410990	13.65%	1.692%
12	8.722	966	975	995	rBV	240513	453125	15.04%	1.865%
13	9.105	1034	1040	1047	rBV8	2188	4180	0.14%	0.017%
14	9.434	1088	1096	1113	rBV	236549	410968	13.65%	1.692%
15	9.975	1181	1188	1210	rBV	255789	634793	21.08%	2.613%
16	10.328	1241	1248	1260	rBV	287935	530604	17.62%	2.184%
17	10.481	1268	1274	1294	rBV	215324	450995	14.97%	1.856%
18	10.781	1321	1325	1331	rBV5	3191	5070	0.17%	0.021%
19	10.875	1337	1341	1346	rVB7	2158	3780	0.13%	0.016%
20	11.175	1388	1392	1399	rBV9	2774	5003	0.17%	0.021%
21	11.940	1515	1522	1528	rBV8	2996	6690	0.22%	0.028%
22	12.475	1607	1613	1619	rVB9	8168	13310	0.44%	0.055%
23	12.899	1680	1685	1690	rVB3	7124	10746	0.36%	0.044%
24	13.375	1757	1766	1769	rBV	5755	11614	0.39%	0.048%
25	13.610	1800	1806	1817	rBV	781844	1046750	34.75%	4.309%
26	13.716	1817	1824	1839	rVB	498476	719950	23.90%	2.963%
27	13.851	1839	1847	1849	rBV2	43748	75643	2.51%	0.311%
28	13.893	1849	1854	1867	rVB2	649349	1097571	36.44%	4.518%
29	14.075	1879	1885	1889	rBV8	5963	8910	0.30%	0.037%
30	14.134	1891	1895	1897	rBV	9244	11931	0.40%	0.049%
31	14.198	1897	1906	1918	rVB2	562137	1021738	33.92%	4.206%
32	14.328	1925	1928	1931	rVB4	4812	5770	0.19%	0.024%
33	14.381	1931	1937	1942	rVB	31232	38210	1.27%	0.157%
34	14.434	1942	1946	1953	rVB10	4484	8289	0.28%	0.034%
35	14.504	1953	1958	1971	rBV2	38920	129343	4.29%	0.532%
36	14.734	1992	1997	2001	rBV	30124	39765	1.32%	0.164%

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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-BP110924.MA.M
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37	14.869	2013	2020	2035	rVB	431871	708040	23.51%	2.914%
38	15.051	2047	2051	2055	rBV6	3087	4558	0.15%	0.019%
39	15.204	2070	2077	2088	rBV	868036	1560781	51.82%	6.425%
40	15.369	2099	2105	2117	rBV	352679	597222	19.83%	2.458%
41	15.451	2117	2119	2125	rVB6	7230	11351	0.38%	0.047%
42	15.745	2162	2169	2174	rBV6	12709	22428	0.74%	0.092%
43	16.128	2232	2234	2242	rVB9	2296	4173	0.14%	0.017%
44	16.416	2278	2283	2291	rBV9	4905	10465	0.35%	0.043%
45	16.528	2297	2302	2306	rBV8	9784	15649	0.52%	0.064%
46	16.592	2306	2313	2323	rVB	65075	95731	3.18%	0.394%
47	16.763	2337	2342	2346	rBV8	3694	6647	0.22%	0.027%
48	16.934	2367	2371	2374	rBV5	9645	15765	0.52%	0.065%
49	16.998	2374	2382	2394	rVB	747964	1181486	39.23%	4.863%
50	17.116	2394	2402	2415	rBV	1293589	2070150	68.73%	8.521%
51	17.310	2429	2435	2446	rBV	171306	233662	7.76%	0.962%
52	17.481	2461	2464	2465	rBV3	3237	3594	0.12%	0.015%
53	17.510	2465	2469	2478	rVB	32585	46546	1.55%	0.192%
54	17.663	2492	2495	2496	rBV3	7017	7807	0.26%	0.032%
55	17.692	2496	2500	2505	rVB6	20284	29186	0.97%	0.120%
56	17.892	2530	2534	2538	rBV6	6348	9694	0.32%	0.040%
57	17.945	2538	2543	2553	rBV2	76761	125814	4.18%	0.518%
58	18.157	2576	2579	2589	rVB2	12524	17855	0.59%	0.073%
59	18.245	2591	2594	2601	rVB9	3426	6094	0.20%	0.025%
60	18.386	2616	2618	2621	rBV4	4470	5497	0.18%	0.023%
61	18.422	2621	2624	2630	rVV7	9187	13984	0.46%	0.058%
62	19.045	2726	2730	2734	rBV2	24682	32939	1.09%	0.136%
63	19.122	2737	2743	2746	rVV2	19180	38131	1.27%	0.157%
64	19.157	2746	2749	2754	rVB5	23822	31077	1.03%	0.128%
65	19.281	2762	2770	2778	rBV3	42170	96568	3.21%	0.397%
66	19.492	2800	2806	2818	rBV2	1769164	2594565	86.15%	10.680%
67	21.445	3133	3138	3150	rBV2	890317	1501834	49.86%	6.182%
68	24.445	3639	3648	3664	rVB	1265464	3011837	100.00%	12.397%
69	24.680	3678	3688	3699	rBV	606306	1655292	54.96%	6.814%

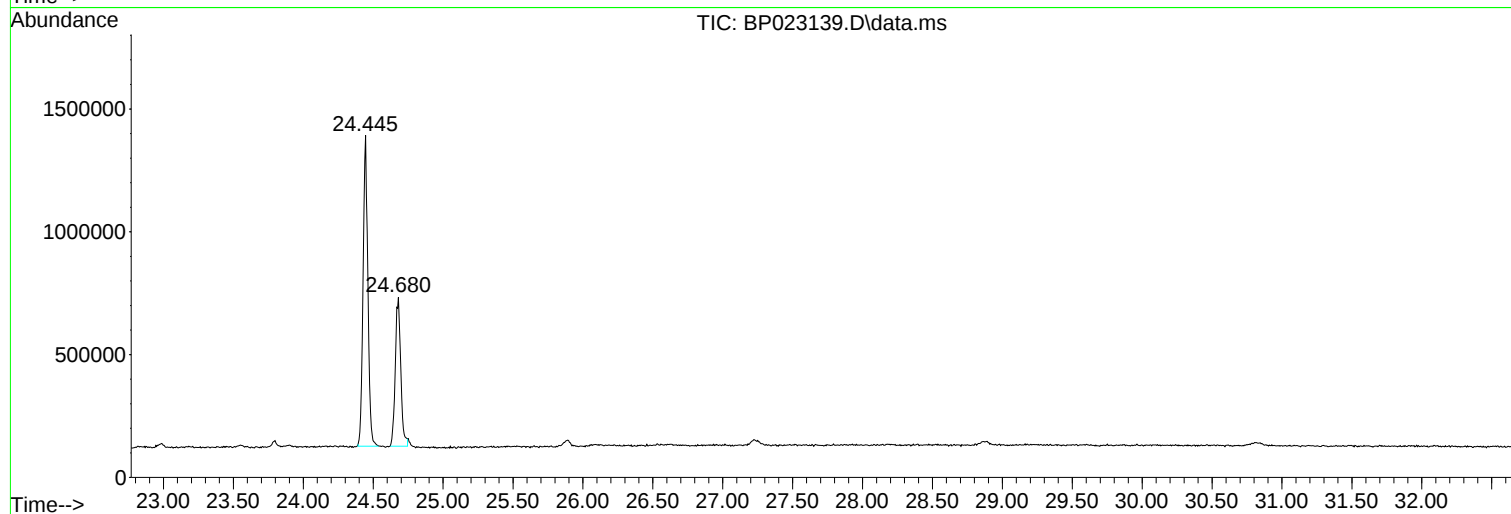
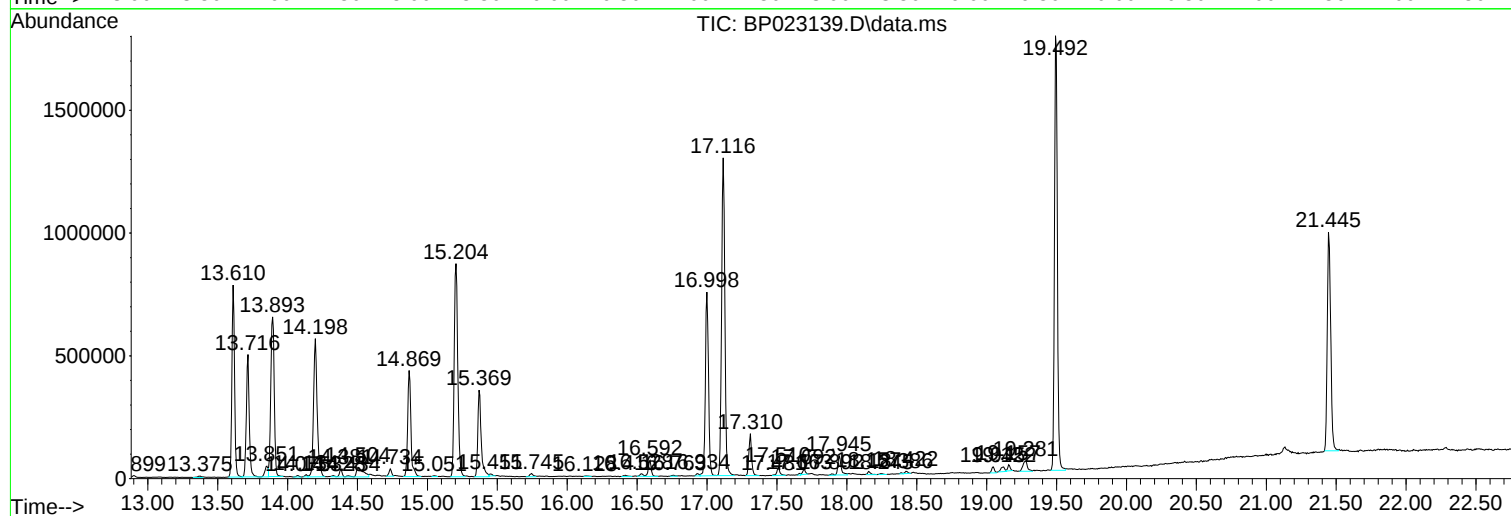
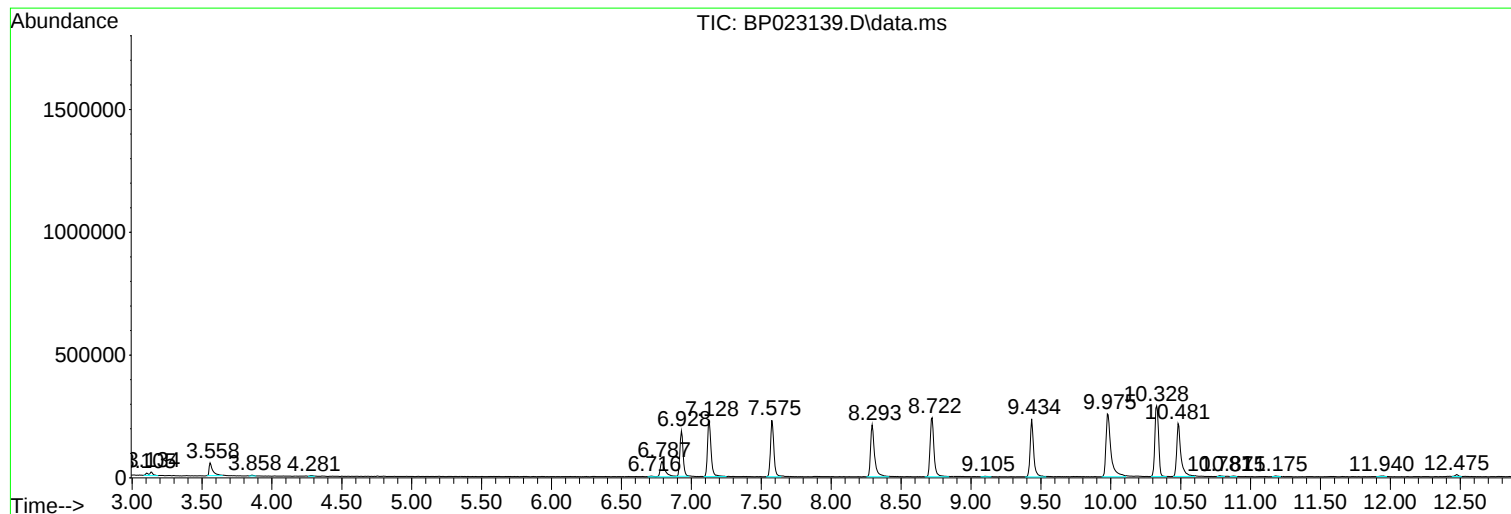
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
Quant Title : SVOA CALIBRATION

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



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ALS Vial : 7 Sample Multiplier: 1

Instrument :
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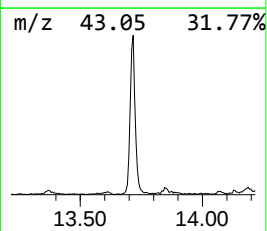
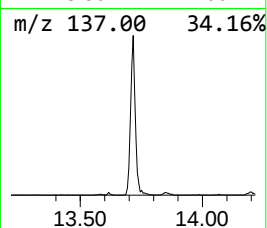
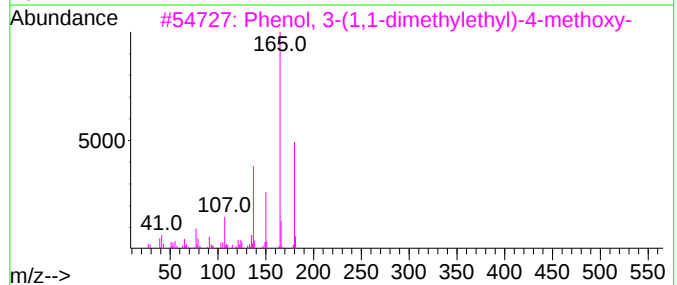
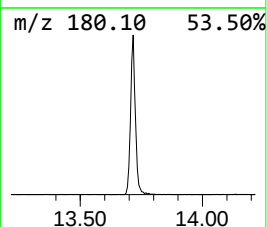
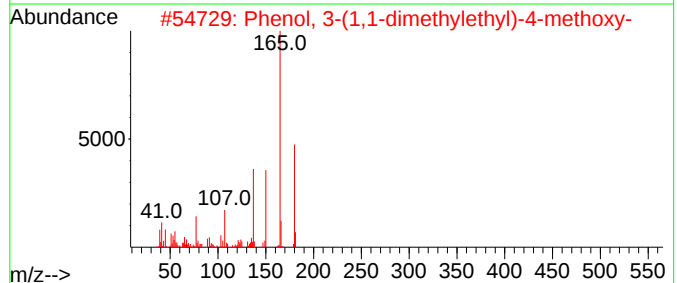
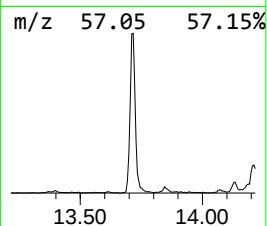
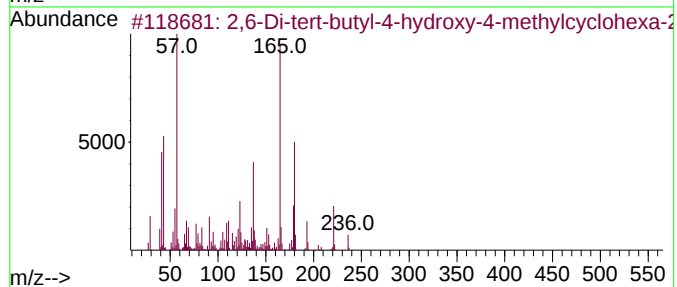
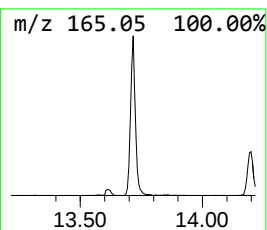
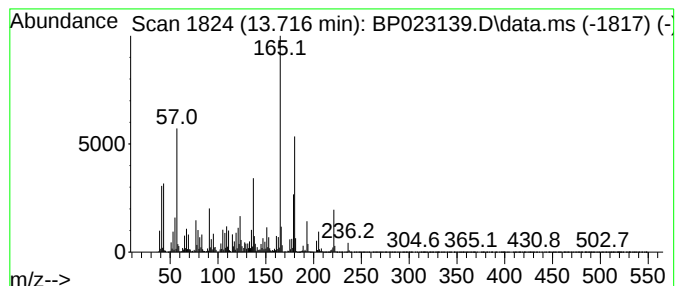
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 2,6-Di-tert-butyl-4-hydroxy... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.716	14.09 ng/ul	719950	Acenaphthene-d10	14.198

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,6-Di-tert-butyl-4-hydroxy-4-me...	236	C15H24O2	010396-80-2	95		
2	Phenol, 3-(1,1-dimethylethyl)-4-...	180	C11H16O2	000088-32-4	70		
3	Phenol, 3-(1,1-dimethylethyl)-4-...	180	C11H16O2	000088-32-4	70		
4	3-tert-Butyl-4-hydroxyanisole	180	C11H16O2	000121-00-6	64		
5	3H-Benz[e]indene, 2-methyl-	180	C14H12	150096-60-9	64		



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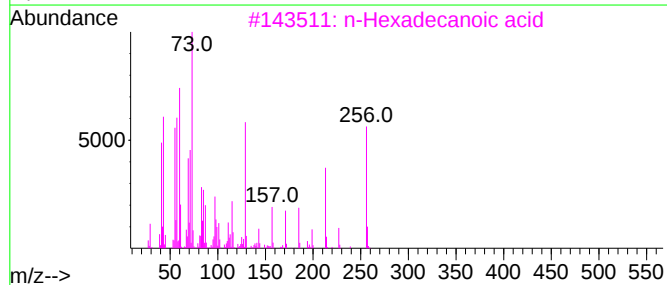
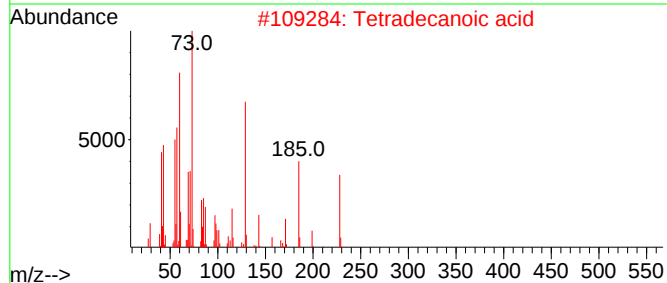
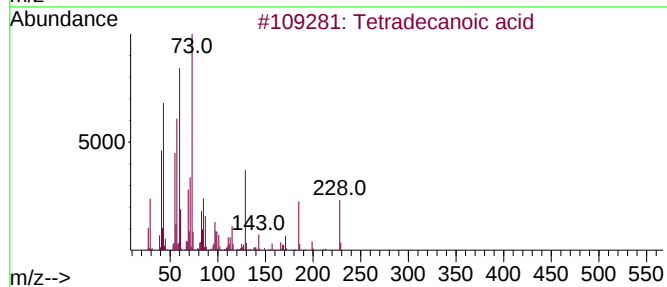
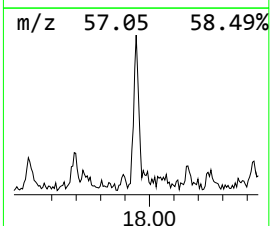
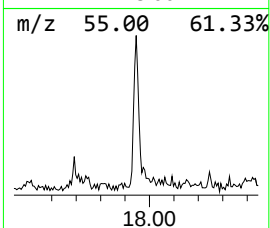
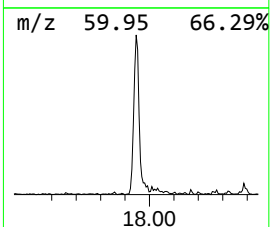
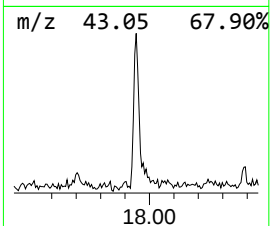
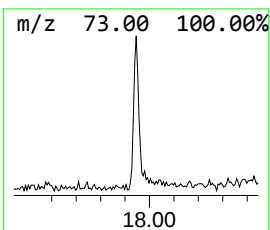
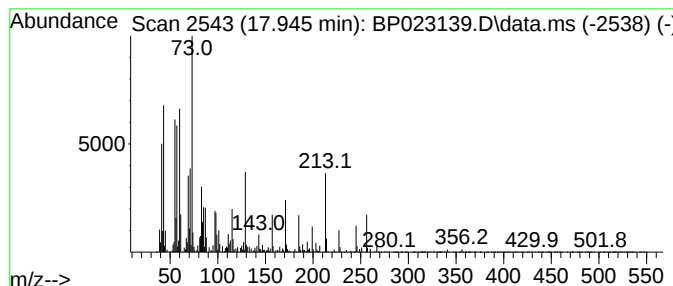
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Tetradecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.945	2.13 ng/ul	125814	Phenanthrene-d10	16.998

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecanoic acid	228	C14H28O2	000544-63-8	94
2			Tetradecanoic acid	228	C14H28O2	000544-63-8	93
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
4			Tridecanoic acid	214	C13H26O2	000638-53-9	93
5			Tridecanoic acid	214	C13H26O2	000638-53-9	93



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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
2,6-Di-tert-but...	13.716	14.1	ng/ul	719950	3	14.198	1021740	20.0
Tetradecanoic acid	17.945	2.1	ng/ul	125814	4	16.998	1181490	20.0