

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111324\
 Data File : BP023129.D
 Acq On : 15 Nov 2024 07:58
 Operator : RC/JU
 Sample : P4784-04
 Misc :
 ALS Vial : 53 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E27N2

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: BP023129.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.099	16	19	22	rBV4	3187	4079	0.12%	0.015%
2	3.134	22	25	32	rVB	16580	21606	0.62%	0.080%
3	3.558	93	97	108	rBV	58270	110170	3.15%	0.406%
4	3.858	145	148	152	rBV5	3918	5149	0.15%	0.019%
5	4.281	216	220	226	rVB3	3133	5484	0.16%	0.020%
6	6.722	625	635	641	rBV10	2818	6011	0.17%	0.022%
7	6.793	641	647	664	rBV	80787	181623	5.20%	0.669%
8	6.922	664	669	685	rVB	210045	329262	9.42%	1.212%
9	7.128	698	704	720	rBV	260082	448056	12.82%	1.650%
10	7.575	774	780	790	rBV	262881	439300	12.57%	1.617%
11	8.287	895	901	920	rBV	245894	479958	13.74%	1.767%
12	8.722	969	975	992	rBV	297024	491420	14.07%	1.809%
13	9.116	1036	1042	1049	rBV7	3174	5500	0.16%	0.020%
14	9.434	1089	1096	1118	rBV	242436	457470	13.09%	1.684%
15	9.975	1181	1188	1215	rBV	324728	737587	21.11%	2.715%
16	10.169	1215	1221	1232	rVB8	5212	16094	0.46%	0.059%
17	10.334	1241	1249	1262	rBV	371466	613414	17.56%	2.258%
18	10.481	1267	1274	1301	rBV	209668	497301	14.23%	1.831%
19	11.175	1385	1392	1394	rBV7	3755	6745	0.19%	0.025%
20	11.940	1515	1522	1529	rBV4	4576	8057	0.23%	0.030%
21	12.504	1613	1618	1622	rBV7	3247	6421	0.18%	0.024%
22	12.799	1662	1668	1674	rBV9	2296	4443	0.13%	0.016%
23	13.081	1706	1716	1723	rBV2	15860	31695	0.91%	0.117%
24	13.163	1726	1730	1736	rBV8	4814	9930	0.28%	0.037%
25	13.604	1799	1805	1818	rBV	828113	1161629	33.25%	4.277%
26	13.893	1842	1854	1867	rBV	690779	1296890	37.12%	4.775%
27	14.145	1891	1897	1900	rBV	144134	215451	6.17%	0.793%
28	14.198	1900	1906	1912	rVV	539453	883415	25.29%	3.252%
29	14.251	1912	1915	1916	rVV2	25880	32088	0.92%	0.118%
30	14.281	1916	1920	1925	rVV	67842	134956	3.86%	0.497%
31	14.487	1949	1955	1976	rVB4	38456	143660	4.11%	0.529%
32	14.869	2013	2020	2026	rBV2	31639	83838	2.40%	0.309%
33	15.040	2047	2049	2055	rVB	7288	9470	0.27%	0.035%
34	15.198	2069	2076	2089	rBV	1053901	1720436	49.24%	6.334%
35	15.363	2096	2104	2117	rBV	376640	647006	18.52%	2.382%
36	15.445	2117	2118	2127	rVB7	9643	15398	0.44%	0.057%

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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
 Title : SVOA CALIBRATION

37	15.587	2138	2142	2148	rVB5	8656	12642	0.36%	0.047%
38	15.787	2170	2176	2183	rBV3	26147	51931	1.49%	0.191%
39	15.845	2184	2186	2190	rVB5	5408	5783	0.17%	0.021%
40	16.004	2208	2213	2219	rBV	79788	105180	3.01%	0.387%
41	16.081	2219	2226	2240	rBV	265751	605851	17.34%	2.230%
42	16.475	2290	2293	2299	rVB8	5077	8548	0.24%	0.031%
43	16.769	2336	2343	2346	rVV9	3854	7847	0.22%	0.029%
44	16.804	2346	2349	2354	rVB5	2910	5059	0.14%	0.019%
45	16.934	2364	2371	2376	rBV	100617	134324	3.84%	0.495%
46	16.998	2376	2382	2393	rVV	852059	1327881	38.01%	4.889%
47	17.116	2395	2402	2416	rVV	1520080	2283588	65.36%	8.407%
48	17.245	2418	2424	2430	rVV	240211	321357	9.20%	1.183%
49	17.304	2430	2434	2444	rVB2	16762	30278	0.87%	0.111%
50	17.439	2452	2457	2459	rBV6	4826	8888	0.25%	0.033%
51	17.469	2459	2462	2467	rVB2	15370	20354	0.58%	0.075%
52	17.586	2477	2482	2485	rBV	12425	17391	0.50%	0.064%
53	17.628	2485	2489	2495	rVV	21899	43274	1.24%	0.159%
54	17.698	2497	2501	2508	rVB6	19375	40184	1.15%	0.148%
55	17.945	2538	2543	2550	rBV5	25488	48416	1.39%	0.178%
56	18.010	2552	2554	2560	rVB2	9699	12526	0.36%	0.046%
57	18.386	2614	2618	2624	rVB	149736	174166	4.99%	0.641%
58	18.439	2625	2627	2631	rVB4	6419	5420	0.16%	0.020%
59	18.481	2631	2634	2636	rBV4	5141	7354	0.21%	0.027%
60	18.692	2665	2670	2674	rBV	77907	103147	2.95%	0.380%
61	18.928	2707	2710	2714	rVB5	8522	11799	0.34%	0.043%
62	19.033	2724	2728	2732	rBV2	28059	43913	1.26%	0.162%
63	19.116	2738	2742	2747	rVB	14549	23314	0.67%	0.086%
64	19.286	2767	2771	2774	rBV6	10595	16103	0.46%	0.059%
65	19.492	2799	2806	2815	rBV	1962476	3017107	86.36%	11.108%
66	21.445	3132	3138	3150	rBV	1093664	1833933	52.49%	6.752%
67	24.439	3637	3647	3660	rVB	1401566	3493687	100.00%	12.862%
68	24.663	3676	3685	3706	rVB	680168	2079942	59.53%	7.657%

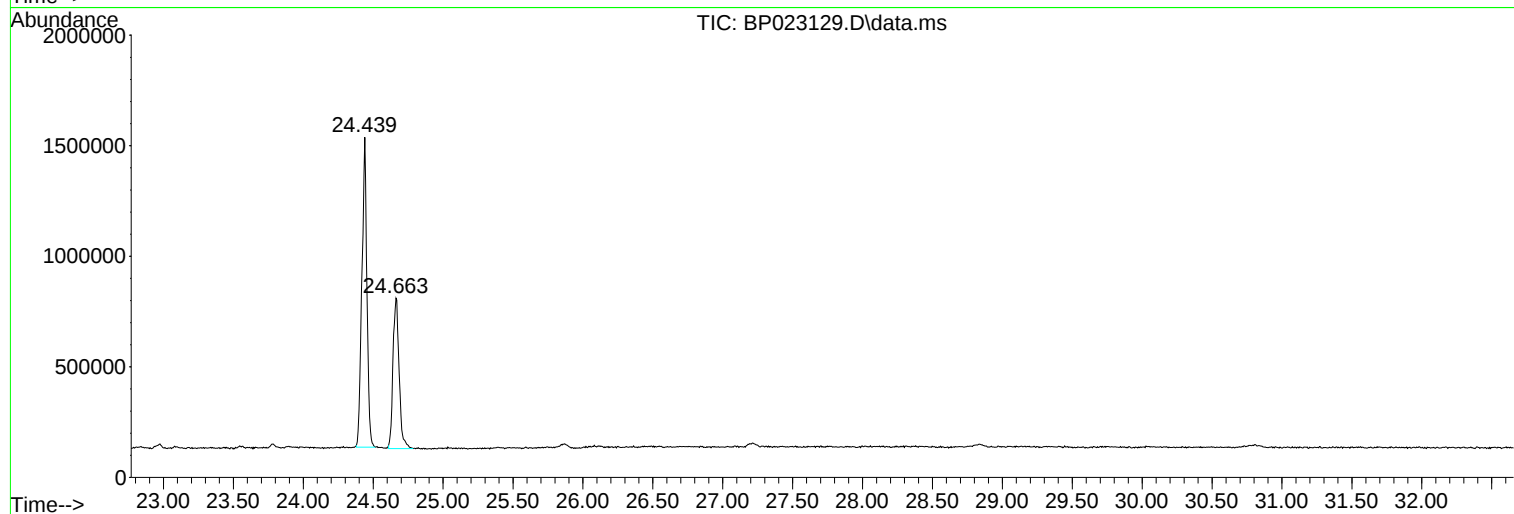
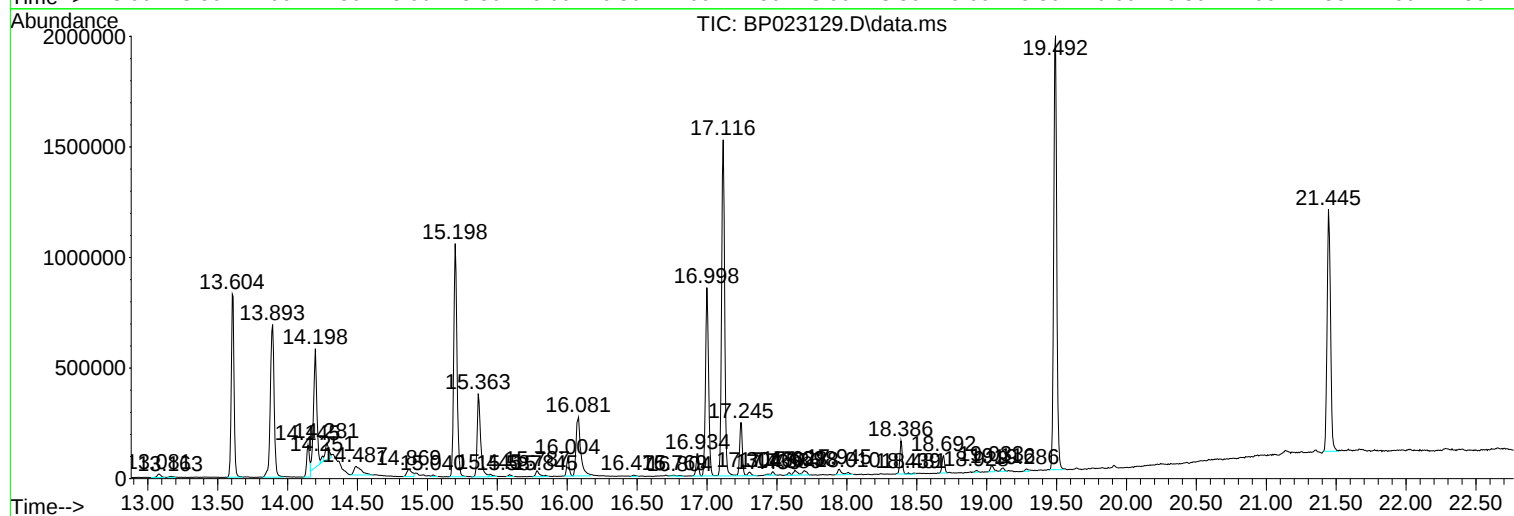
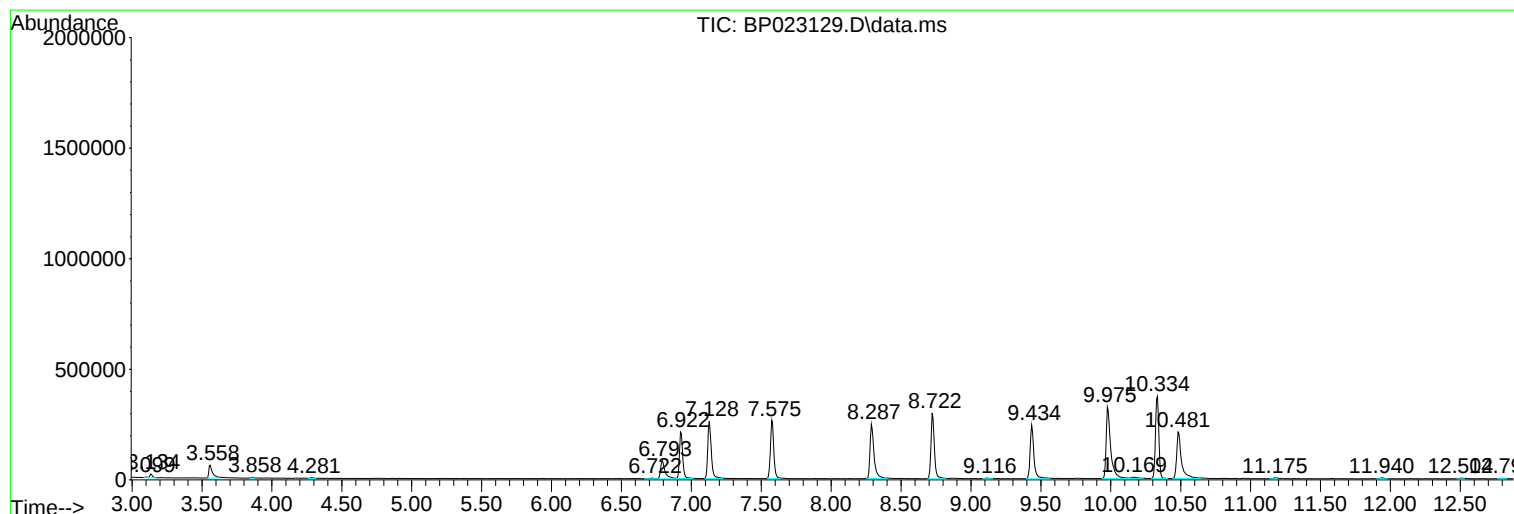
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
Quant Title : SVOA CALIBRATION

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



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Sample : P4784-04
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ALS Vial : 53 Sample Multiplier: 1

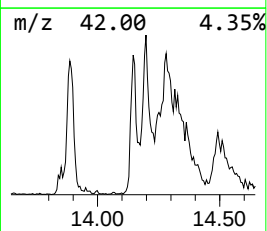
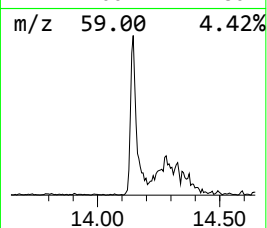
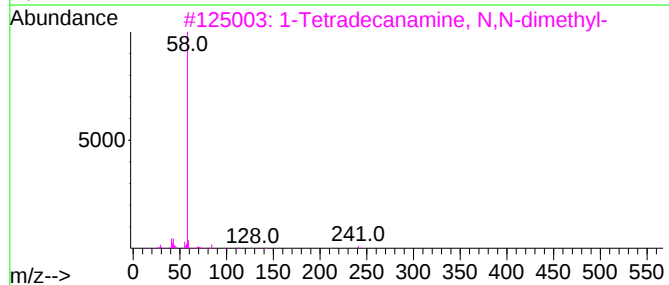
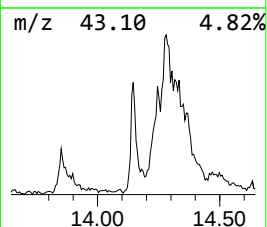
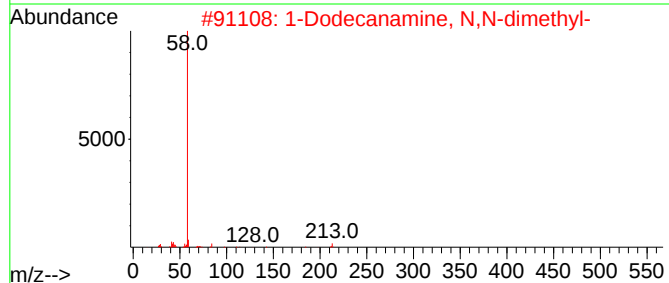
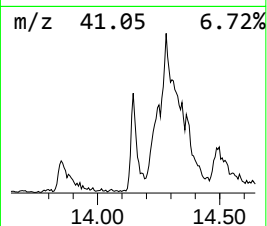
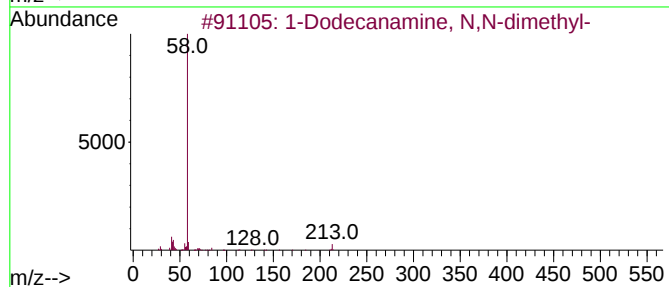
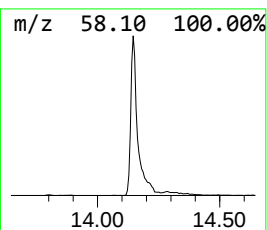
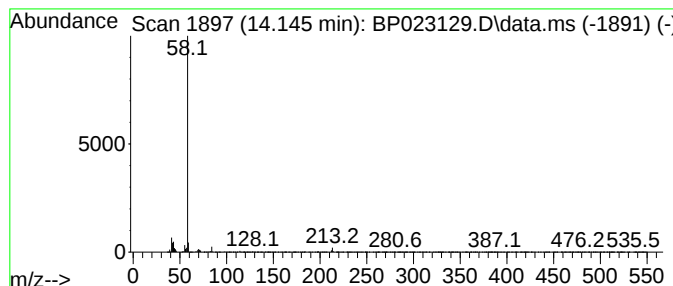
Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 1-Dodecanamine, N,N-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.146	4.88 ng/ul	215451	Acenaphthene-d10	14.198	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Dodecanamine, N,N-dimethyl-	213	C14H31N	000112-18-5	94
2	1-Dodecanamine, N,N-dimethyl-	213	C14H31N	000112-18-5	91
3	1-Tetradecanamine, N,N-dimethyl-	241	C16H35N	000112-75-4	78
4	3(N,N-Dimethylmyristylammonio)pr...	363	C19H41NO3S	014933-09-6	78
5	Dimethyl palmitamine	269	C18H39N	000112-69-6	72



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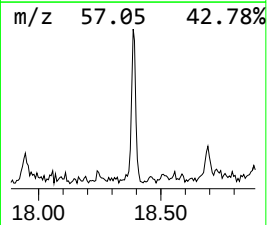
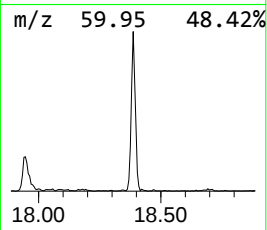
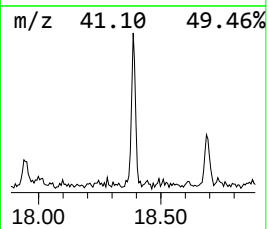
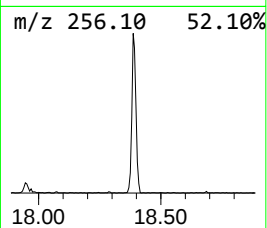
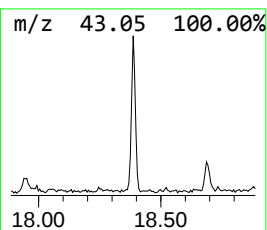
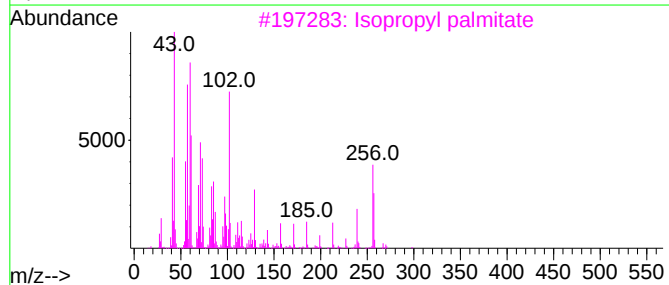
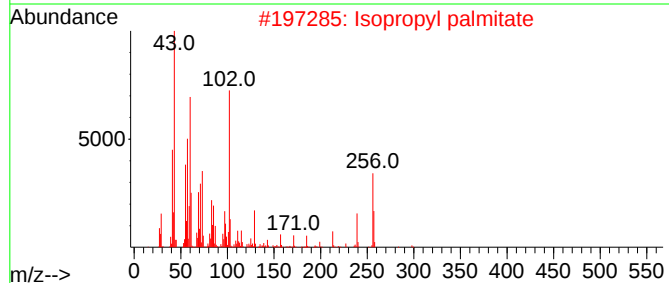
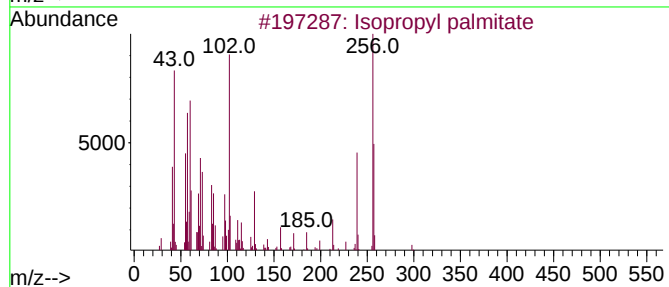
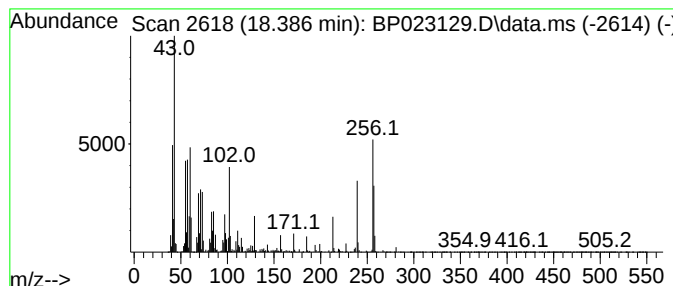
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 Isopropyl palmitate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.386	2.62 ng/ul	174166	Phenanthrene-d10	16.998

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Isopropyl palmitate	298	C19H38O2	000142-91-6	91
2			Isopropyl palmitate	298	C19H38O2	000142-91-6	86
3			Isopropyl palmitate	298	C19H38O2	000142-91-6	52
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	45
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	27



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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
1-Dodecanamine,...	14.146	4.9	ng/ul	215451	3	14.198	883415	20.0
Isopropyl palmi...	18.386	2.6	ng/ul	174166	4	16.998	1327880	20.0