

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN082021\
 Data File : BN015997.D
 Acq On : 20 Aug 2021 22:03
 Operator : CG/JU
 Sample : M3448-09
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBKE8

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_N\Methods\SFAM-EPA-BN081621MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BN015997.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.740	109	111	116	rBV2	142465	198731	3.25%	0.347%
2	7.034	666	671	683	rVB	202459	353902	5.79%	0.617%
3	7.205	694	700	710	rBV	674706	1146534	18.75%	1.999%
4	7.405	728	734	744	rVB	722491	1200141	19.63%	2.093%
5	7.875	807	814	822	rVB	858509	1404618	22.97%	2.449%
6	8.575	927	933	942	rVB	553349	929698	15.21%	1.621%
7	8.705	953	955	957	rBV3	6022	6362	0.10%	0.011%
8	9.034	1004	1011	1019	rVB	910884	1546832	25.30%	2.697%
9	9.757	1126	1134	1146	rVB	667625	1180618	19.31%	2.059%
10	10.293	1218	1225	1235	rBV	987243	1722658	28.18%	3.004%
11	10.452	1246	1252	1262	rBV	109452	185507	3.03%	0.323%
12	10.675	1282	1290	1301	rVB	1125348	2005939	32.81%	3.498%
13	10.810	1305	1313	1328	rVV	674615	1186656	19.41%	2.069%
14	11.457	1421	1423	1426	rBV3	4616	6663	0.11%	0.012%
15	11.493	1426	1429	1434	rVB5	6575	10994	0.18%	0.019%
16	11.957	1505	1508	1514	rVB6	11545	20111	0.33%	0.035%
17	12.263	1554	1560	1566	rVB5	18219	31328	0.51%	0.055%
18	12.757	1641	1644	1649	rVB5	9719	15248	0.25%	0.027%
19	12.875	1660	1664	1671	rBV6	9575	19930	0.33%	0.035%
20	13.128	1701	1707	1710	rBV6	5591	10368	0.17%	0.018%
21	13.922	1835	1842	1853	rBV	2117138	3086674	50.49%	5.382%
22	14.157	1877	1882	1883	rBV4	10180	12041	0.20%	0.021%
23	14.204	1883	1890	1900	rVB	2245996	3376998	55.23%	5.888%
24	14.416	1923	1926	1930	rBV5	4931	6947	0.11%	0.012%
25	14.510	1935	1942	1952	rVV	1695418	2603126	42.58%	4.539%
26	14.704	1966	1975	1983	rBV	124488	204218	3.34%	0.356%
27	14.840	1995	1998	2001	rBV3	5393	7102	0.12%	0.012%
28	15.322	2076	2080	2084	rVV6	14253	23622	0.39%	0.041%
29	15.504	2104	2111	2124	rBV	2928298	4498693	73.58%	7.844%
30	15.616	2124	2130	2137	rBV	470850	689753	11.28%	1.203%
31	15.740	2148	2151	2159	rVB3	34719	54638	0.89%	0.095%
32	15.940	2180	2185	2187	rBV3	8143	11915	0.19%	0.021%
33	16.075	2204	2208	2214	rBV6	4994	10964	0.18%	0.019%
34	16.292	2241	2245	2247	rVB4	8355	11340	0.19%	0.020%
35	16.387	2257	2261	2262	rBV4	7634	9731	0.16%	0.017%
36	16.563	2289	2291	2295	rBV5	6364	7700	0.13%	0.013%

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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_N\Methods\SFAM-EPA-BN081621MA.M
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37	16.745	2318	2322	2328	rBV4	8594	16678	0.27%	0.029%
38	16.934	2351	2354	2357	rVB5	6366	6493	0.11%	0.011%
39	17.045	2369	2373	2381	rVB7	10771	19733	0.32%	0.034%
40	17.181	2393	2396	2398	rBV3	6966	8272	0.14%	0.014%
41	17.257	2403	2409	2416	rBV	2089039	3062464	50.09%	5.340%
42	17.357	2419	2426	2435	rVV	3406434	5081879	83.12%	8.861%
43	17.628	2470	2472	2475	rBV4	5170	7521	0.12%	0.013%
44	17.757	2492	2494	2496	rBV3	4697	6174	0.10%	0.011%
45	17.810	2502	2503	2508	rVB5	6786	8139	0.13%	0.014%
46	17.934	2519	2524	2528	rVB7	16219	26491	0.43%	0.046%
47	17.986	2528	2533	2534	rBV4	4966	7834	0.13%	0.014%
48	18.051	2538	2544	2548	rVB6	11451	18350	0.30%	0.032%
49	18.122	2551	2556	2558	rBV6	5723	8890	0.15%	0.016%
50	18.157	2558	2562	2570	rBV9	10526	23688	0.39%	0.041%
51	18.228	2570	2574	2578	rBV3	11817	19062	0.31%	0.033%
52	18.404	2600	2604	2605	rBV3	5386	6711	0.11%	0.012%
53	18.616	2637	2640	2641	rBV3	7762	8026	0.13%	0.014%
54	18.692	2650	2653	2654	rBV3	7383	7842	0.13%	0.014%
55	19.069	2714	2717	2720	rBV4	6074	7169	0.12%	0.013%
56	19.157	2729	2732	2734	rBV4	6184	7752	0.13%	0.014%
57	19.216	2738	2742	2748	rVV4	29604	45348	0.74%	0.079%
58	19.286	2750	2754	2758	rVV	43841	63120	1.03%	0.110%
59	19.539	2793	2797	2801	rBV4	20199	26784	0.44%	0.047%
60	19.651	2809	2816	2822	rBV	4362463	6046297	98.89%	10.543%
61	19.792	2838	2840	2844	rVB5	10227	10033	0.16%	0.017%
62	19.828	2844	2846	2848	rBV3	8744	8631	0.14%	0.015%
63	19.963	2867	2869	2870	rBV2	10410	7018	0.11%	0.012%
64	20.651	2981	2986	2992	rBV	690034	852302	13.94%	1.486%
65	21.157	3068	3072	3078	rBV7	88587	143909	2.35%	0.251%
66	21.439	3115	3120	3131	rBV	2504432	3460183	56.60%	6.033%
67	23.110	3402	3404	3409	rVB2	108867	148680	2.43%	0.259%
68	23.686	3494	3502	3509	rBV	3006590	6113934	100.00%	10.661%
69	23.839	3521	3528	3537	rVB	1701446	3514860	57.49%	6.129%
70	24.463	3628	3634	3643	rBV4	172077	413067	6.76%	0.720%
71	25.310	3773	3778	3786	rVB5	137500	338816	5.54%	0.591%

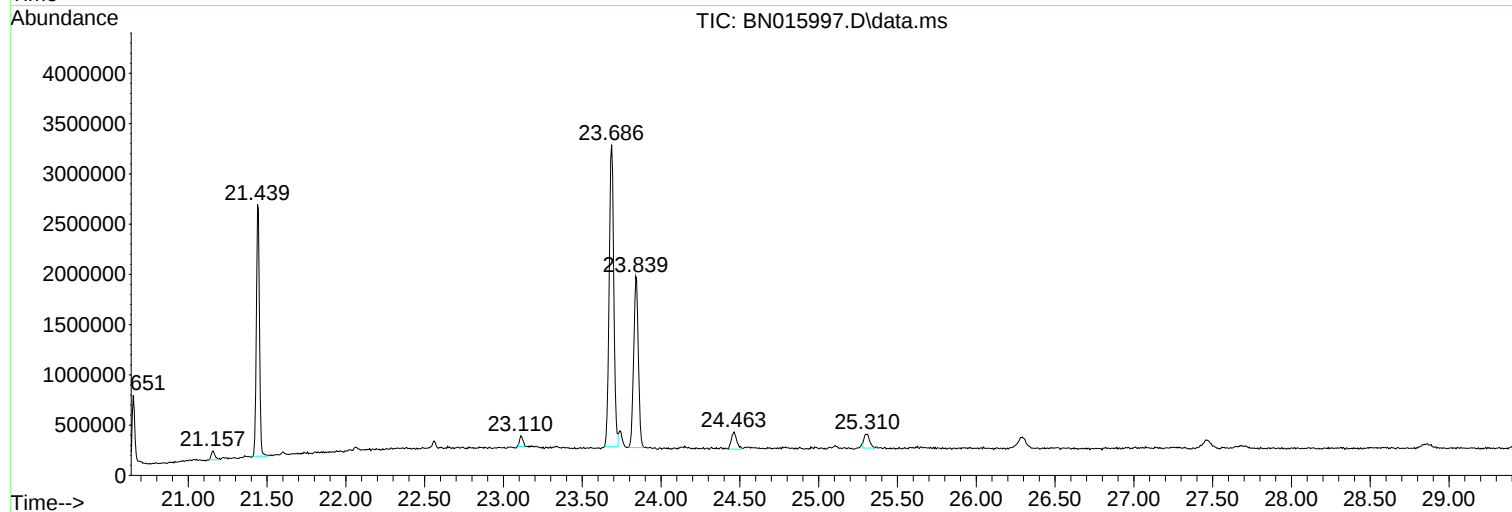
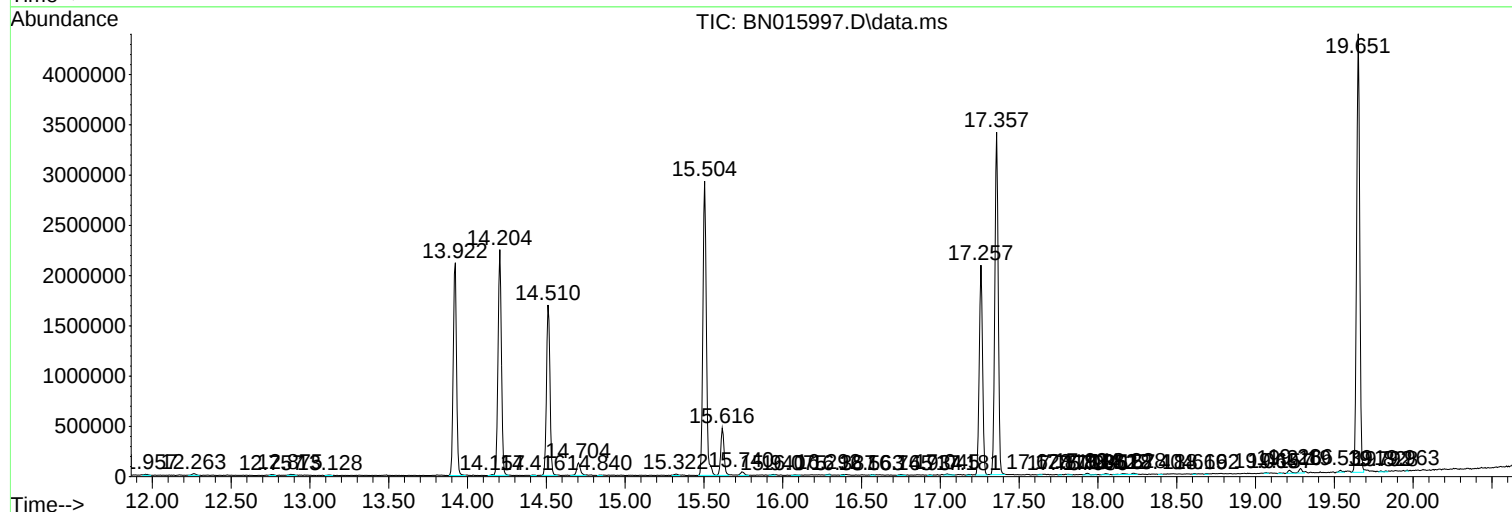
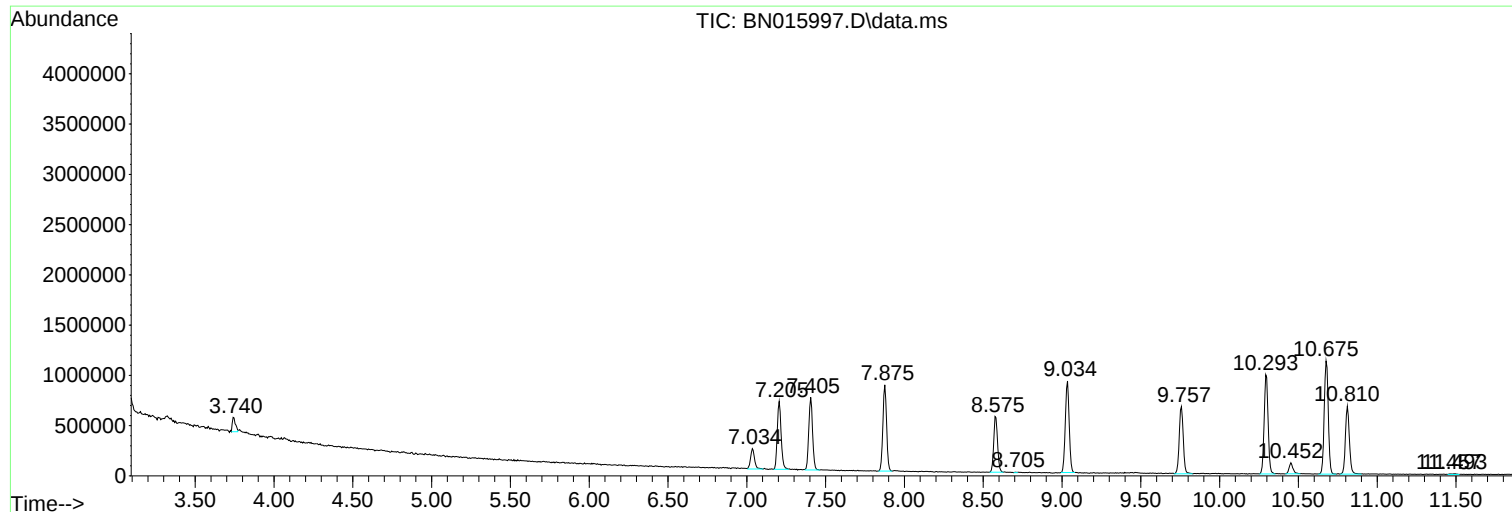
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081621MA.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
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ClientSampleId :
DBKE8

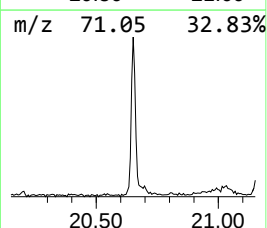
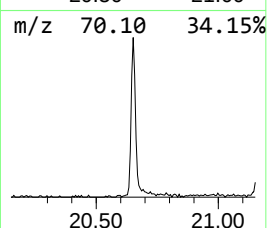
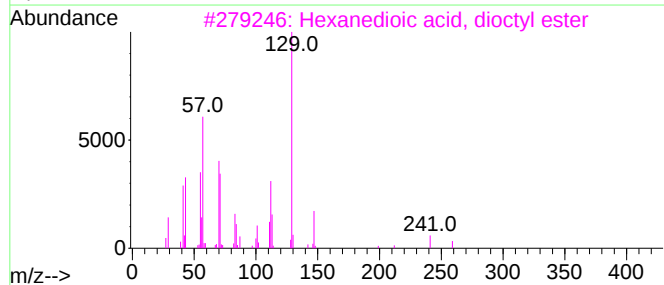
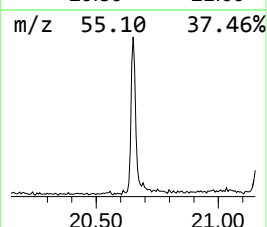
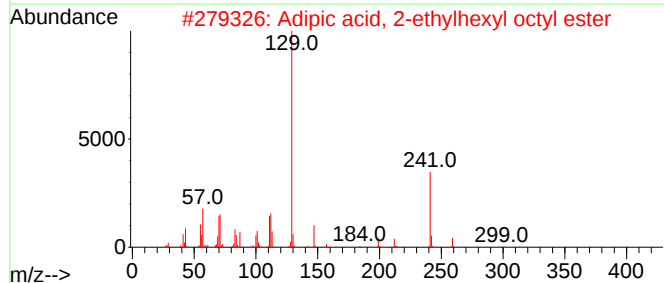
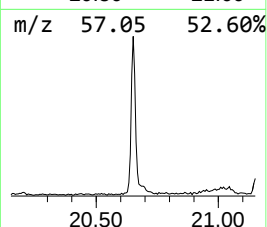
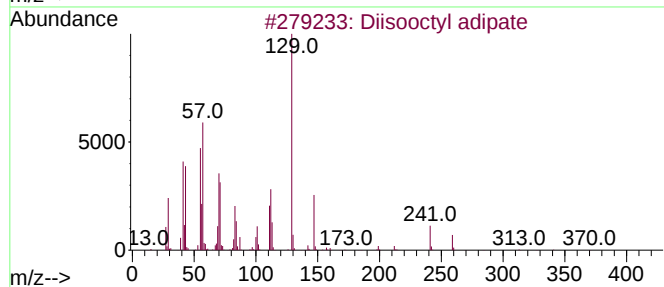
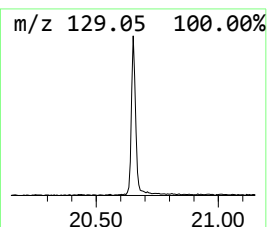
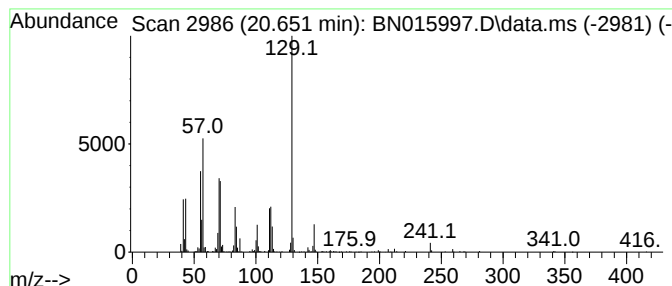
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081621MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Diisooctyl adipate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.651	4.93 ng/ul	852302	Chrysene-d12	21.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Diisooctyl adipate	370	C22H42O4	001330-86-5	91
2			Adipic acid, 2-ethylhexyl octyl ...	370	C22H42O4	1000324-48-6	90
3			Hexanedioic acid, dioctyl ester	370	C22H42O4	000123-79-5	87
4			Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	87
5			Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	83



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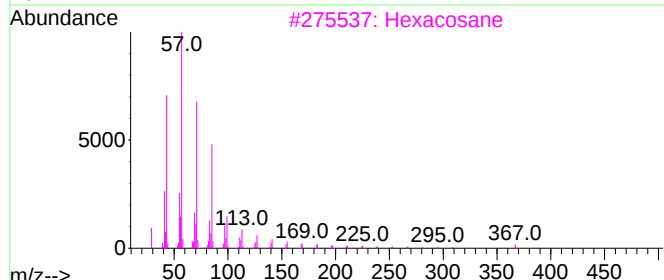
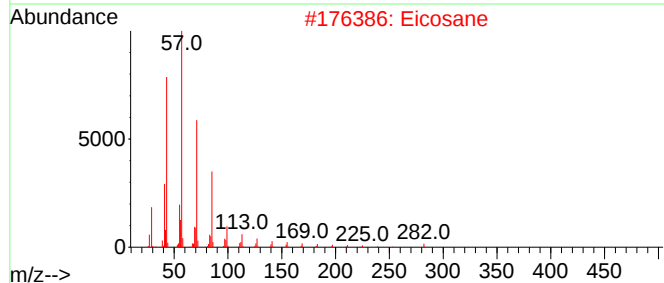
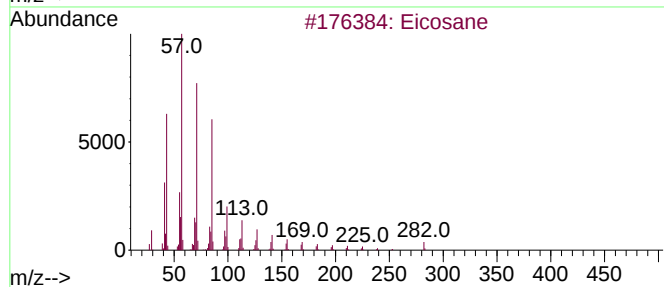
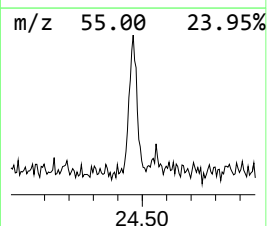
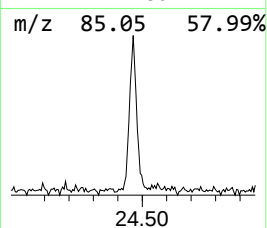
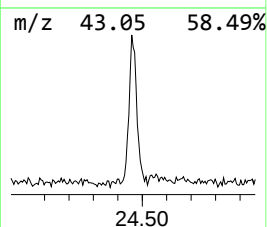
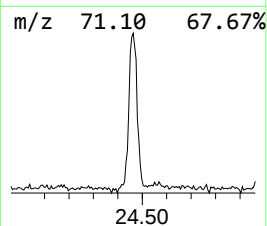
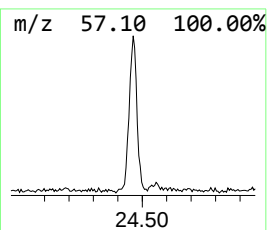
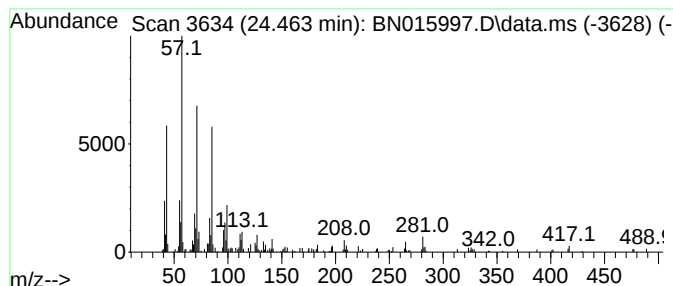
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081621MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.463	2.35 ng/ul	413067	Perylene-d12	23.839

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Eicosane	282	C20H42	000112-95-8	76
2		Eicosane	282	C20H42	000112-95-8	76
3		Hexacosane	366	C26H54	000630-01-3	76
4		Eicosane, 1-iodo-	408	C20H41I	1000406-31-8	74
5		Hexacosane, 1-iodo-	492	C26H53I	1000406-32-1	74



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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
Diisooctyl adipate	20.651	4.9	ng/u1	852302	5	21.445	3460180	20.0
(DEL) Alkane: S...	24.463	2.4	ng/u1	413067	6	23.839	3514860	20.0