

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100124\
 Data File : BP022134.D
 Acq On : 01 Oct 2024 13:48
 Operator : RC/JU
 Sample : P4010-06
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AW0

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-BP092424.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP022134.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.217	36	39	45	rVB5	5658	8255	0.30%	0.053%
2	3.658	111	114	119	rVV5	3524	6019	0.22%	0.038%
3	3.734	123	127	131	rVB	9291	11823	0.43%	0.076%
4	3.952	160	164	172	rVB2	7493	10612	0.39%	0.068%
5	4.858	313	318	323	rBV	12133	19699	0.72%	0.126%
6	5.722	461	465	470	rBV6	2458	4470	0.16%	0.029%
7	6.410	576	582	587	rBV5	5012	9379	0.34%	0.060%
8	6.881	652	662	673	rBV2	85813	167725	6.12%	1.072%
9	7.034	682	688	696	rBV	73733	110045	4.02%	0.703%
10	7.240	716	723	733	rVB	91443	147563	5.39%	0.943%
11	7.328	733	738	744	rBV7	2884	5278	0.19%	0.034%
12	7.693	793	800	808	rBV	423069	720005	26.29%	4.602%
13	8.393	913	919	930	rBV	43188	83726	3.06%	0.535%
14	8.505	932	938	947	rBV2	68737	117486	4.29%	0.751%
15	8.840	988	995	1004	rBV	76247	136339	4.98%	0.871%
16	9.393	1082	1089	1098	rBV8	6573	13570	0.50%	0.087%
17	9.546	1107	1115	1123	rBV2	62633	116192	4.24%	0.743%
18	9.710	1137	1143	1148	rBV	13352	25933	0.95%	0.166%
19	9.904	1166	1176	1183	rVB3	8411	16406	0.60%	0.105%
20	10.087	1200	1207	1219	rBV	100903	173358	6.33%	1.108%
21	10.451	1261	1269	1282	rBV	568733	948397	34.63%	6.061%
22	10.593	1282	1293	1305	rBV	56207	112175	4.10%	0.717%
23	10.999	1354	1362	1369	rBV10	7947	17618	0.64%	0.113%
24	11.240	1399	1403	1404	rBV2	4127	4678	0.17%	0.030%
25	11.963	1522	1526	1529	rBV6	2078	2991	0.11%	0.019%
26	12.046	1534	1540	1543	rVV7	2307	4089	0.15%	0.026%
27	13.634	1805	1810	1815	rVB7	5716	9525	0.35%	0.061%
28	13.722	1817	1825	1833	rBV	219081	302291	11.04%	1.932%
29	13.887	1849	1853	1858	rVB2	5035	6760	0.25%	0.043%
30	13.998	1865	1872	1882	rBV	195255	315666	11.53%	2.017%
31	14.087	1882	1887	1892	rVB9	2823	5499	0.20%	0.035%
32	14.304	1916	1924	1933	rBV	903441	1359378	49.64%	8.688%
33	14.375	1933	1936	1941	rVB7	2636	4878	0.18%	0.031%
34	14.522	1956	1961	1967	rBV2	39256	77582	2.83%	0.496%
35	14.634	1977	1980	1983	rVB4	3385	4471	0.16%	0.029%
36	14.681	1985	1988	1993	rVV5	2574	3615	0.13%	0.023%

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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-BP092424.MA.M
 Title : SVOA CALIBRATION

37	14.745	1995	1999	2005	rVB8	5957	10127	0.37%	0.065%
38	15.122	2057	2063	2064	rBV6	2654	3566	0.13%	0.023%
39	15.316	2089	2096	2103	rBV	290740	438804	16.02%	2.804%
40	15.445	2112	2118	2125	rBV	58568	84978	3.10%	0.543%
41	15.692	2154	2160	2168	rBV2	27642	49433	1.81%	0.316%
42	16.504	2295	2298	2302	rVV6	3404	4662	0.17%	0.030%
43	16.634	2312	2320	2325	rBV2	10138	17897	0.65%	0.114%
44	16.757	2337	2341	2346	rBV8	9453	15606	0.57%	0.100%
45	16.922	2367	2369	2376	rVB8	3332	3717	0.14%	0.024%
46	17.116	2393	2402	2407	rBV	1135897	1780002	65.00%	11.376%
47	17.157	2407	2409	2413	rVV	72879	84169	3.07%	0.538%
48	17.210	2413	2418	2429	rVB	332351	464468	16.96%	2.968%
49	17.539	2464	2474	2477	rBV2	9730	20524	0.75%	0.131%
50	17.622	2486	2488	2490	rBV3	3902	3090	0.11%	0.020%
51	17.757	2506	2511	2516	rVB9	12540	19060	0.70%	0.122%
52	17.816	2517	2521	2524	rBV6	8212	10583	0.39%	0.068%
53	17.904	2534	2536	2545	rVB10	5774	12475	0.46%	0.080%
54	18.051	2551	2561	2568	rBV3	81675	167326	6.11%	1.069%
55	18.222	2584	2590	2593	rBV5	12309	20484	0.75%	0.131%
56	18.628	2656	2659	2667	rVB2	23634	29887	1.09%	0.191%
57	18.875	2698	2701	2706	rBV2	15487	23387	0.85%	0.149%
58	19.239	2759	2763	2773	rVB	162779	309617	11.31%	1.979%
59	19.375	2783	2786	2791	rVB5	27754	37090	1.35%	0.237%
60	19.522	2806	2811	2816	rBV	230391	316439	11.56%	2.022%
61	19.586	2817	2822	2826	rBV	384917	601613	21.97%	3.845%
62	19.757	2848	2851	2856	rBV4	50181	64684	2.36%	0.413%
63	21.539	3146	3154	3159	rBV	1438594	2526783	92.27%	16.149%
64	23.968	3563	3567	3574	rVB	173281	334806	12.23%	2.140%
65	24.586	3667	3672	3680	rVB	175315	369554	13.49%	2.362%
66	24.810	3701	3710	3723	rVB	951406	2738482	100.00%	17.502%

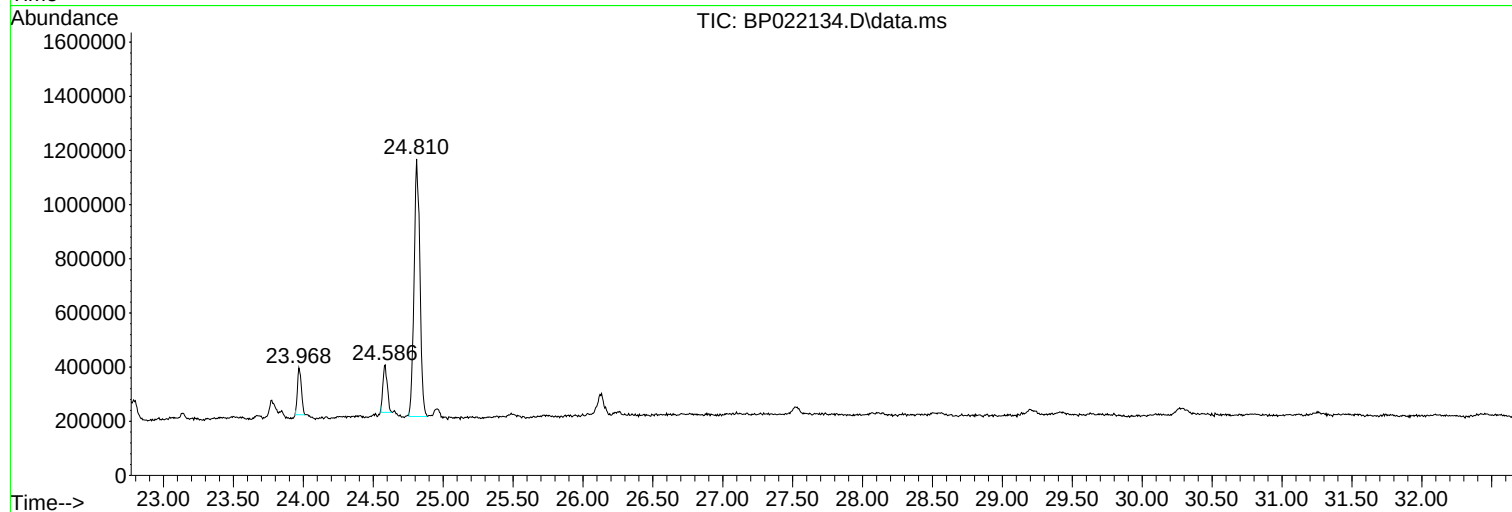
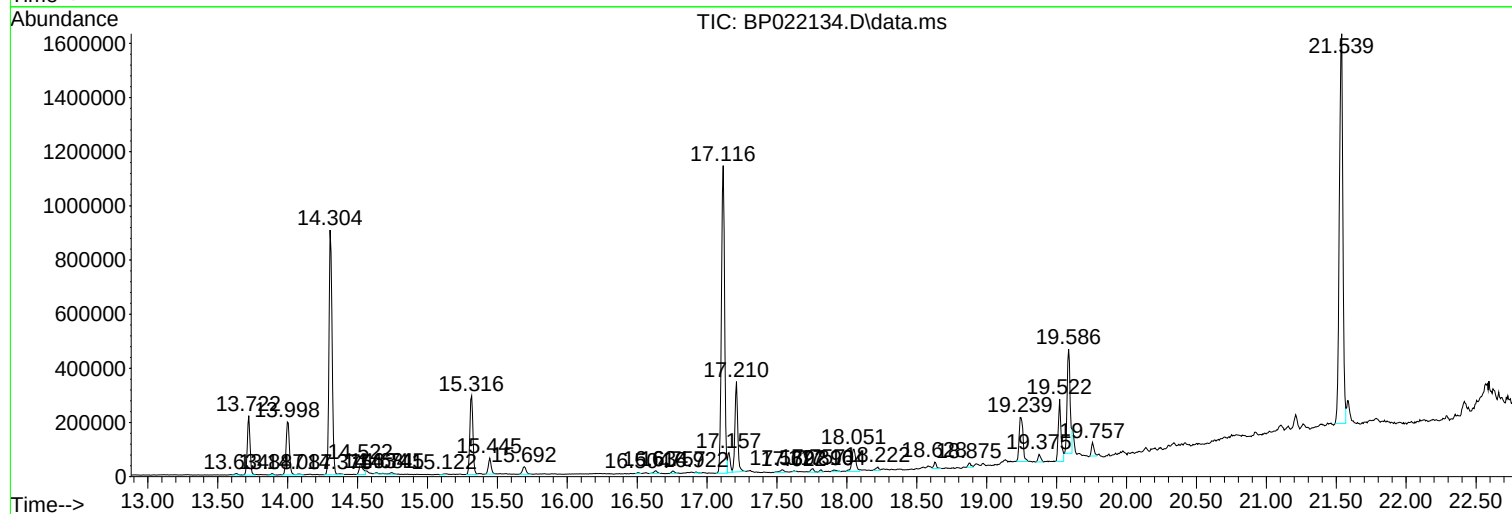
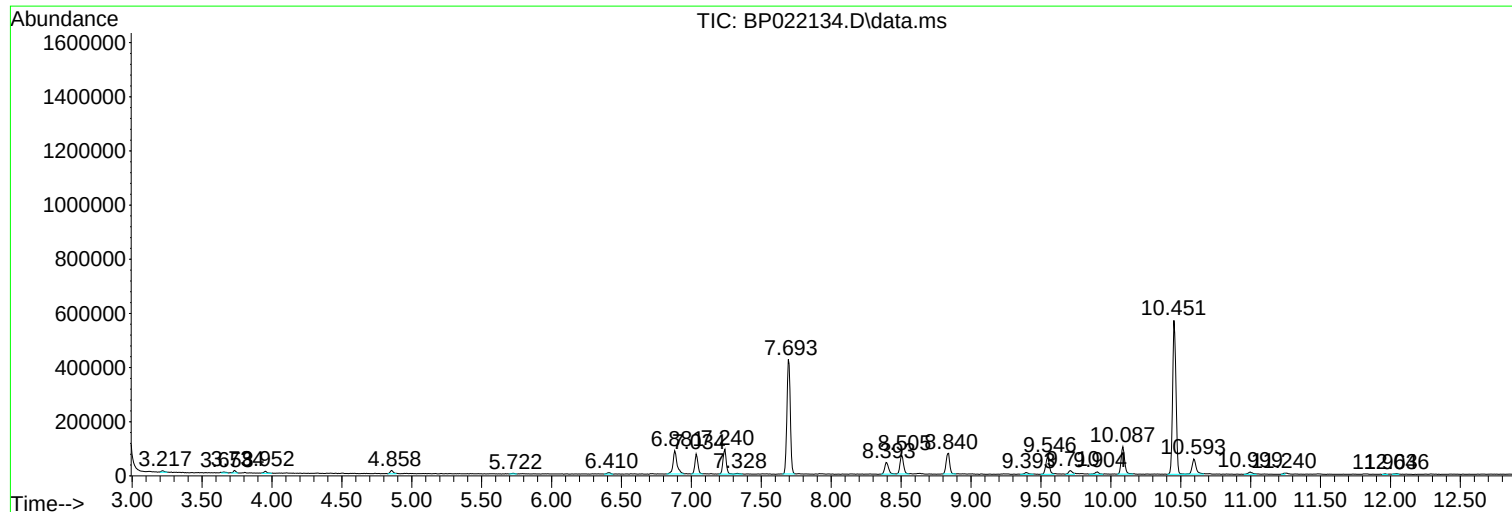
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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



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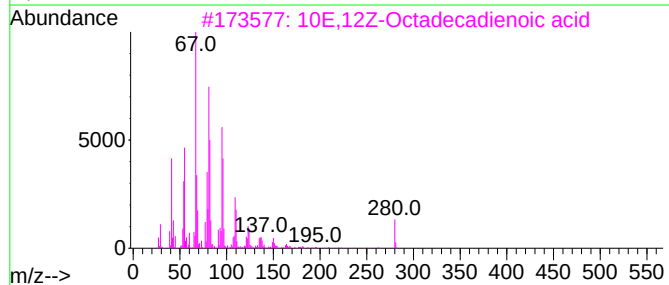
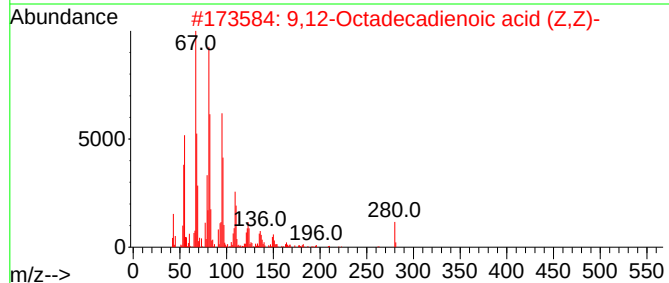
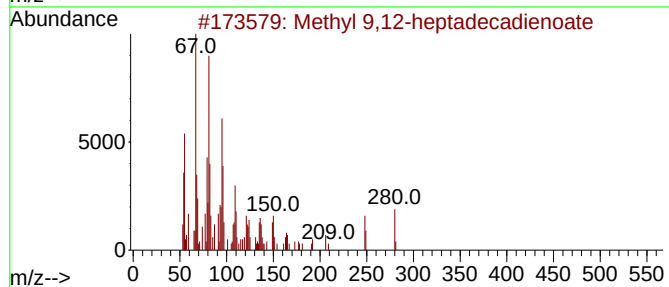
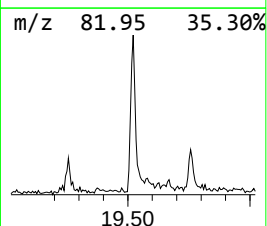
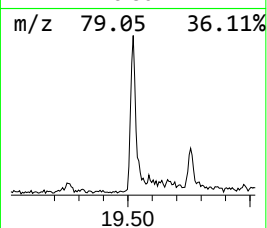
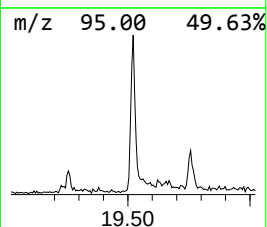
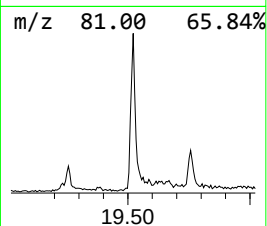
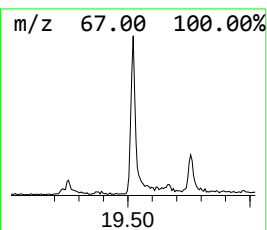
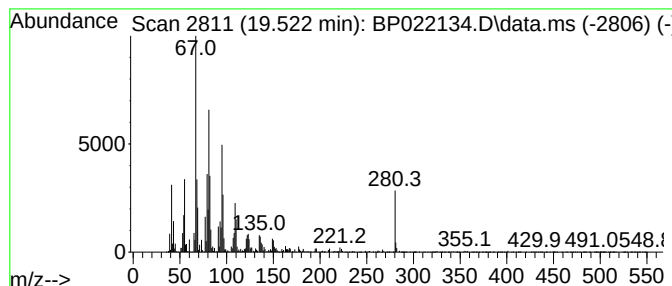
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Methyl 9,12-heptadecadienoate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.522	2.50 ng/ul	316439	Chrysene-d12	21.539

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl 9,12-heptadecadienoate	280	C18H32O2	1000336-36-2	86
2			9,12-Octadecadienoic acid (Z,Z)-	280	C18H32O2	000060-33-3	86
3			10E,12Z-Octadecadienoic acid	280	C18H32O2	002420-56-6	76
4			3,4-Octadiene, 7-methyl-	124	C9H16	037050-05-8	70
5			(9E,11E)-Octadecadienoic acid	280	C18H32O2	000544-71-8	68



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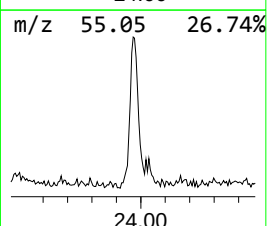
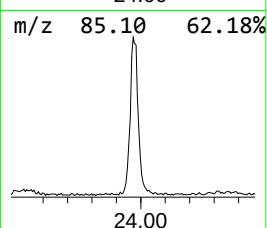
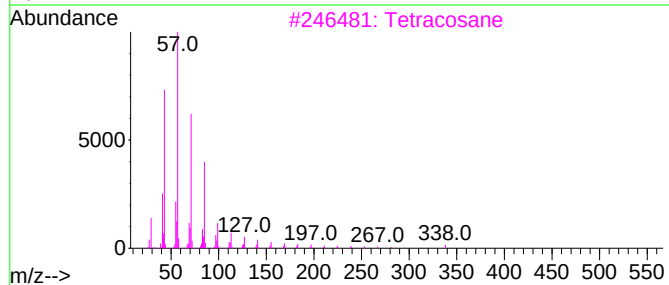
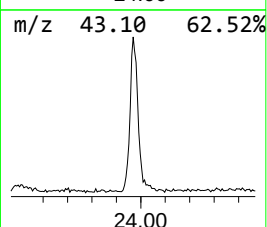
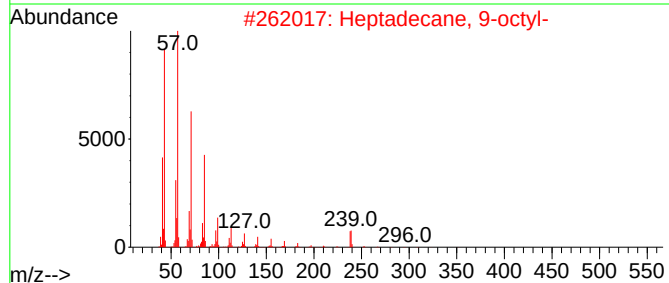
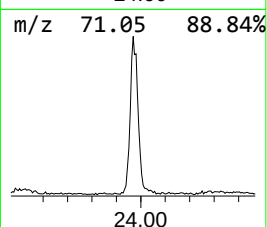
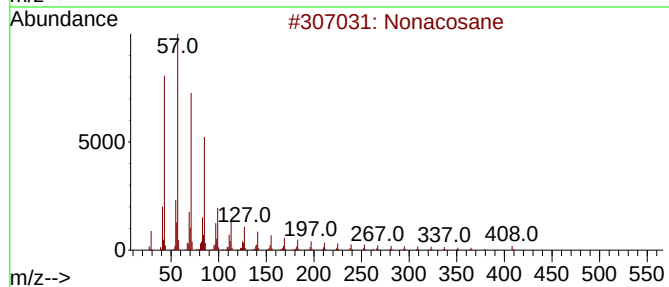
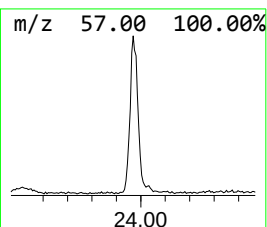
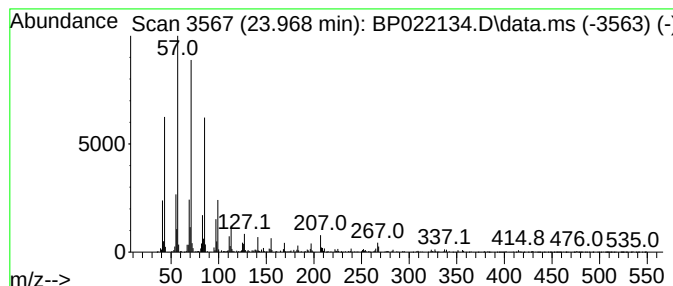
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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.
23.968	2.45 ng/ul	334806	Perylene-d12	24.810

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonacosane	408	C29H60	000630-03-5	94
2			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
3			Tetracosane	338	C24H50	000646-31-1	91
4			Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
5			Heneicosane, 11-decyl-	437	C31H64	055320-06-4	91



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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Methyl 9,12-hep...	19.522	2.5	ng/u1	316439	5	21.539	2526780	20.0
(DEL) Alkane: S...	23.968	2.5	ng/u1	334806	6	24.810	2738480	20.0