

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051622\
 Data File : BP010367.D
 Acq On : 16 May 2022 20:54
 Operator : CG/JU
 Sample : N2746-03
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBSK6

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-BP051322.M
 Title : SVOA CALIBRATION

Signal : TIC: BP010367.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.264	27	30	37	rVB2	18415	24963	0.68%	0.064%
2	3.346	40	44	50	rBV	25646	34623	0.94%	0.088%
3	3.652	93	96	99	rBV2	4191	4701	0.13%	0.012%
4	3.764	112	115	135	rBV	340509	690498	18.82%	1.759%
5	4.087	166	170	174	rVB	10830	14290	0.39%	0.036%
6	4.358	213	216	220	rVB6	3048	4399	0.12%	0.011%
7	4.511	238	242	246	rBV6	2624	3805	0.10%	0.010%
8	5.017	322	328	336	rBV2	10785	24692	0.67%	0.063%
9	5.893	474	477	482	rBV7	2314	3794	0.10%	0.010%
10	7.064	669	676	686	rBV	542031	919428	25.06%	2.342%
11	7.228	697	704	714	rBV	433468	715107	19.49%	1.821%
12	7.311	714	718	725	rVB	12104	21507	0.59%	0.055%
13	7.428	731	738	752	rBV	555795	926371	25.25%	2.359%
14	7.899	810	818	827	rBV	466770	772827	21.06%	1.968%
15	7.964	827	829	837	rVB5	8471	14978	0.41%	0.038%
16	8.605	931	938	958	rBV	724133	1206132	32.87%	3.072%
17	9.058	1005	1015	1026	rBV	583568	1010452	27.54%	2.573%
18	9.222	1039	1043	1051	rVB3	13532	23224	0.63%	0.059%
19	9.499	1083	1090	1096	rVB2	10281	17251	0.47%	0.044%
20	9.781	1129	1138	1154	rBV	487805	827024	22.54%	2.106%
21	10.216	1205	1212	1217	rBV	2237	5528	0.15%	0.014%
22	10.322	1222	1230	1247	rBV	677575	1221632	33.29%	3.111%
23	10.487	1254	1258	1264	rVB9	4778	7893	0.22%	0.020%
24	10.699	1284	1294	1309	rBV	682934	1173835	31.99%	2.990%
25	10.834	1309	1317	1340	rVV	646684	1269542	34.60%	3.233%
26	11.387	1408	1411	1414	rVB5	3113	3843	0.10%	0.010%
27	12.134	1534	1538	1543	rVB5	7395	10851	0.30%	0.028%
28	12.293	1559	1565	1573	rVB4	11195	22482	0.61%	0.057%
29	12.934	1672	1674	1681	rVB7	3372	4982	0.14%	0.013%
30	13.081	1694	1699	1704	rVV9	3807	6346	0.17%	0.016%
31	13.157	1707	1712	1716	rVB8	2556	4123	0.11%	0.011%
32	13.363	1742	1747	1755	rVB	15546	24741	0.67%	0.063%
33	13.499	1766	1770	1772	rVB5	3082	4693	0.13%	0.012%
34	13.940	1837	1845	1857	rBV	1345819	1933584	52.70%	4.925%
35	14.222	1883	1893	1905	rBV	1440426	2187482	59.62%	5.571%
36	14.434	1926	1929	1934	rVB7	3259	5200	0.14%	0.013%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
 Title : SVOA CALIBRATION

37	14.528	1938	1945	1957	rBV	1109094	1646819	44.88%	4.194%
38	14.610	1957	1959	1964	rVB5	3430	4242	0.12%	0.011%
39	14.728	1971	1979	2004	rBV	394211	856281	23.34%	2.181%
40	14.951	2013	2017	2023	rBV9	8541	15060	0.41%	0.038%
41	15.193	2054	2058	2062	rVB7	3866	6262	0.17%	0.016%
42	15.346	2080	2084	2088	rVB7	2726	5314	0.14%	0.014%
43	15.522	2107	2114	2127	rBV	1957479	2809781	76.58%	7.156%
44	15.645	2129	2135	2144	rBV	439387	666668	18.17%	1.698%
45	15.763	2151	2155	2162	rVB2	25266	34442	0.94%	0.088%
46	16.087	2204	2210	2215	rBV9	3415	8389	0.23%	0.021%
47	16.310	2243	2248	2252	rBV6	10202	14604	0.40%	0.037%
48	16.692	2310	2313	2318	rVB7	4188	4362	0.12%	0.011%
49	16.963	2355	2359	2362	rBV6	3506	4982	0.14%	0.013%
50	17.045	2370	2373	2374	rBV3	3946	4371	0.12%	0.011%
51	17.281	2406	2413	2423	rBV	1400798	2016683	54.96%	5.136%
52	17.381	2423	2430	2443	rVV	2141987	3055016	83.26%	7.781%
53	17.481	2444	2447	2451	rVB5	12756	18069	0.49%	0.046%
54	17.681	2478	2481	2487	rBV8	2558	4860	0.13%	0.012%
55	17.957	2524	2528	2533	rVB8	7900	10931	0.30%	0.028%
56	18.169	2559	2564	2573	rBV	121368	194860	5.31%	0.496%
57	19.192	2734	2738	2744	rBV2	31724	45549	1.24%	0.116%
58	19.263	2746	2750	2758	rBV	33516	56290	1.53%	0.143%
59	19.398	2770	2773	2781	rBV3	60119	101248	2.76%	0.258%
60	19.616	2804	2810	2815	rBV	2745322	3669313	100.00%	9.345%
61	20.139	2896	2899	2904	rVB3	32576	35345	0.96%	0.090%
62	20.504	2959	2961	2968	rBV3	33020	39668	1.08%	0.101%
63	20.622	2977	2981	2985	rBV5	57779	80698	2.20%	0.206%
64	21.075	3054	3058	3066	rBV	191786	295740	8.06%	0.753%
65	21.363	3102	3107	3121	rBV	1633622	2340924	63.80%	5.962%
66	21.516	3131	3133	3139	rVB2	62843	71652	1.95%	0.182%
67	21.980	3210	3212	3217	rVB2	58111	69846	1.90%	0.178%
68	22.480	3293	3297	3306	rVB2	202459	331820	9.04%	0.845%
69	23.639	3487	3494	3501	rBV	1721926	3366539	91.75%	8.574%
70	23.792	3513	3520	3532	rVB2	1070298	2230445	60.79%	5.681%

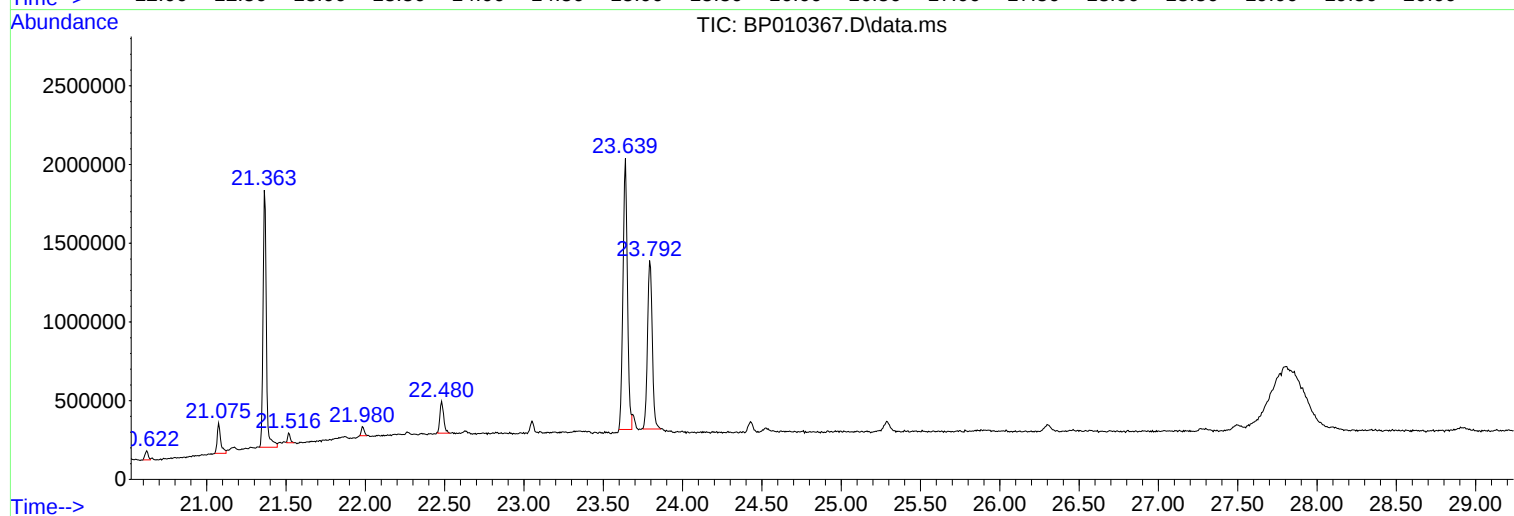
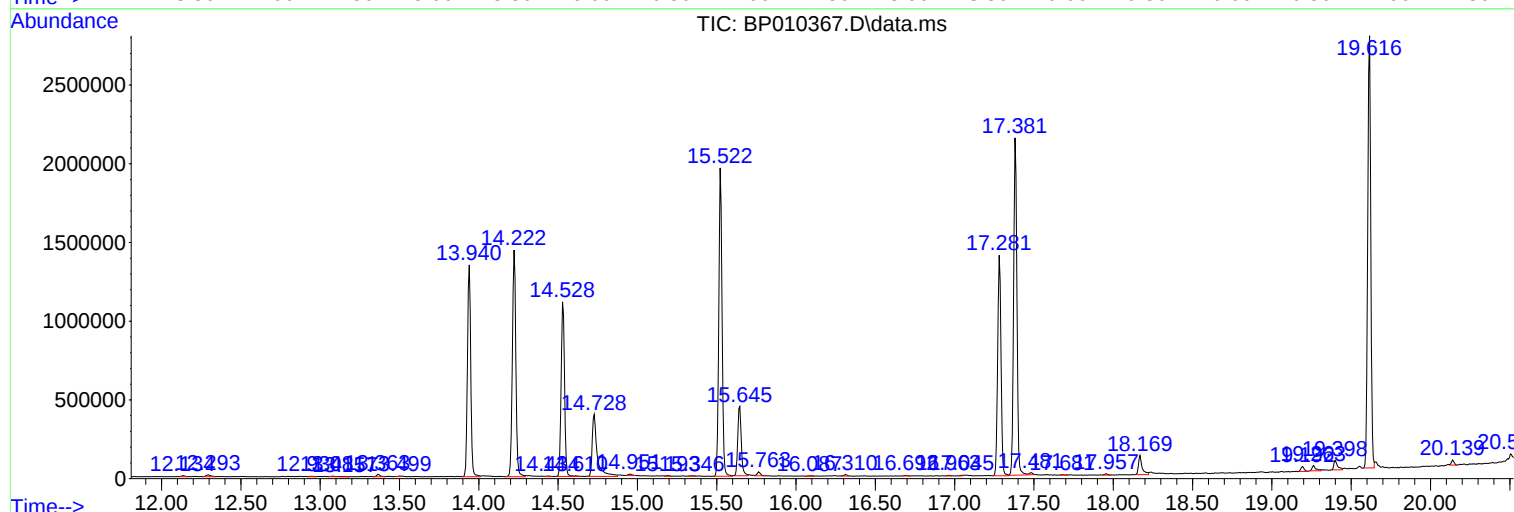
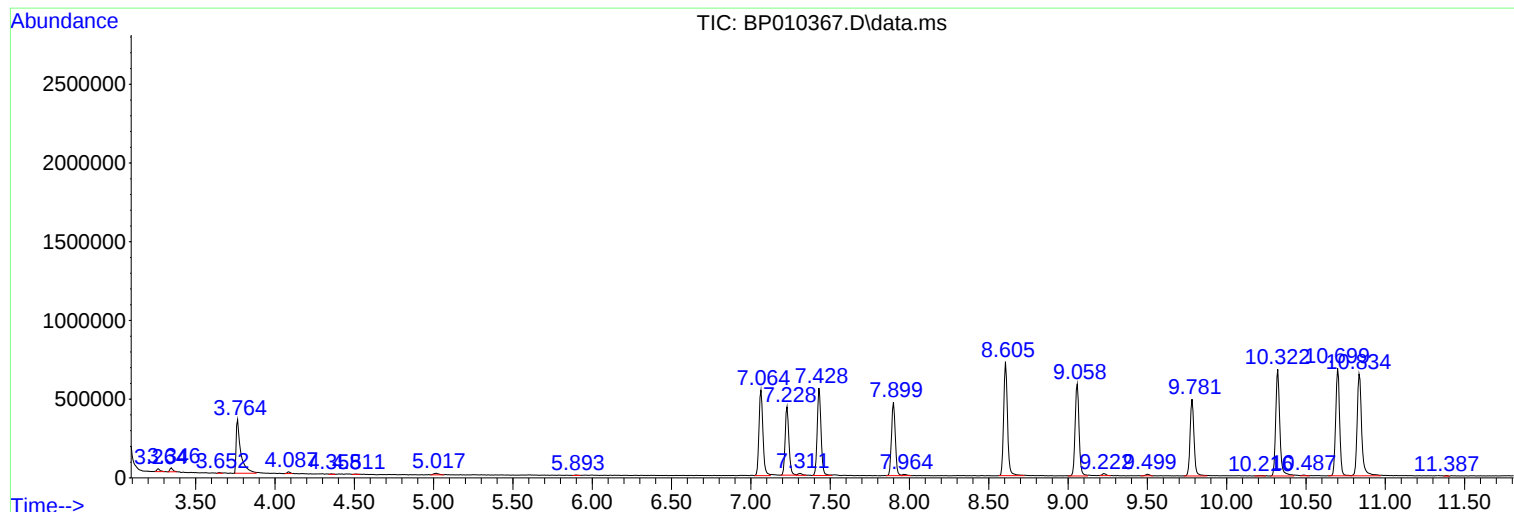
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION

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



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Operator : CG/JU
Sample : N2746-03
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBSK6

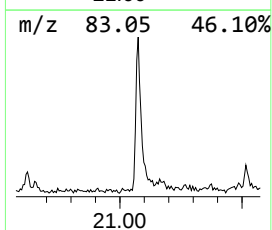
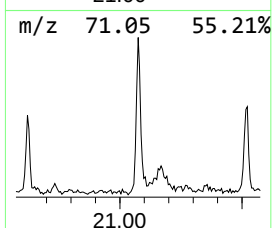
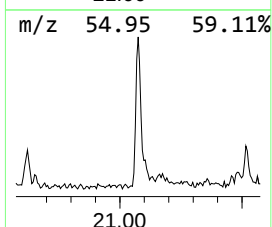
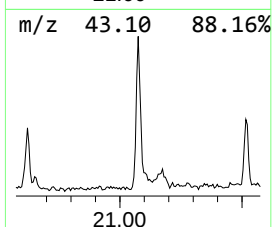
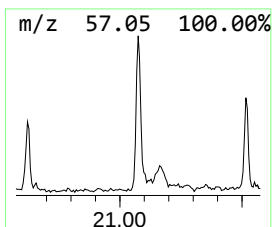
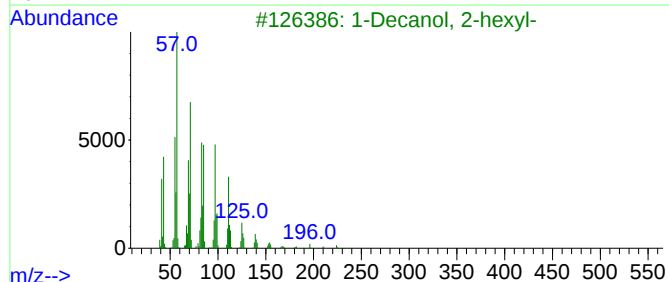
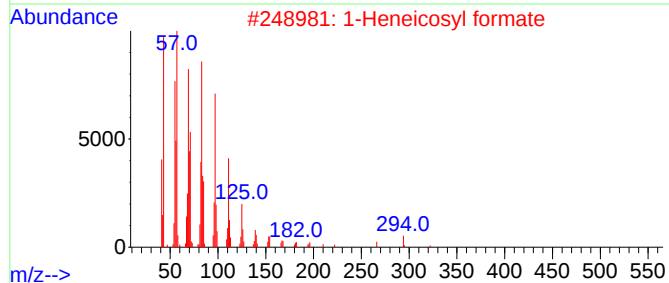
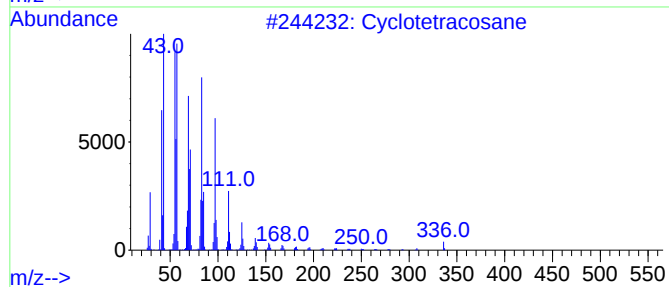
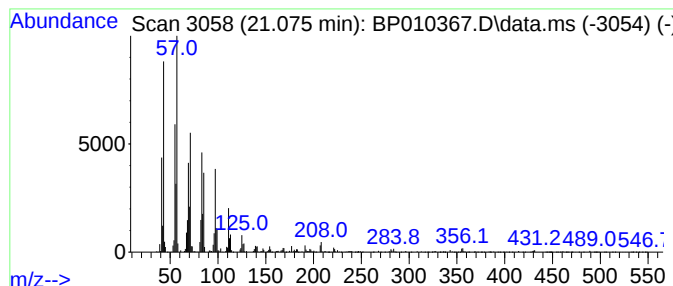
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Cyclic21.075 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.075	2.53 ng/ul	295740	Chrysene-d12	21.363

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclotetracosane	336	C24H48	000297-03-0	95
2			1-Heneicosyl formate	340	C22H44O2	077899-03-7	94
3			1-Decanol, 2-hexyl-	242	C16H34O	002425-77-6	91
4			1-Docosene	308	C22H44	001599-67-3	90
5			Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	90



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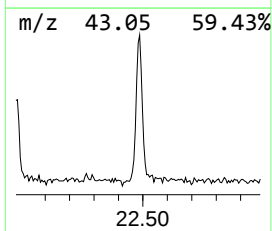
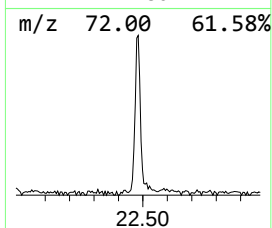
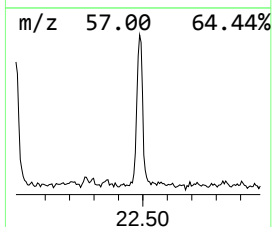
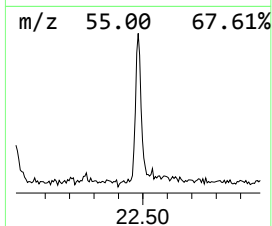
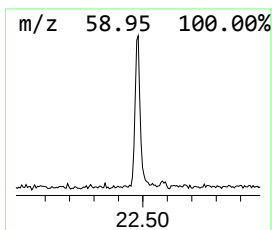
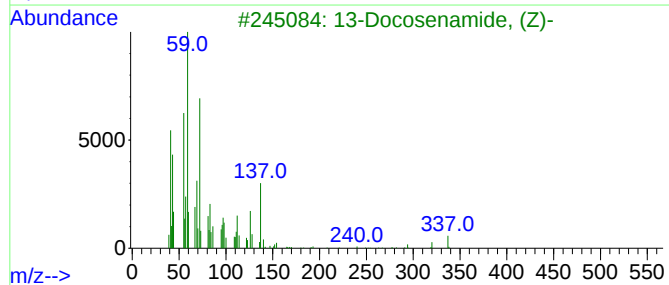
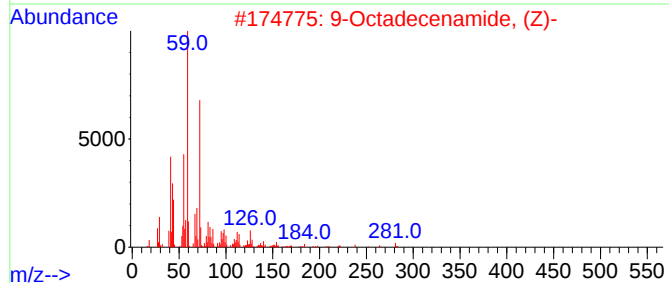
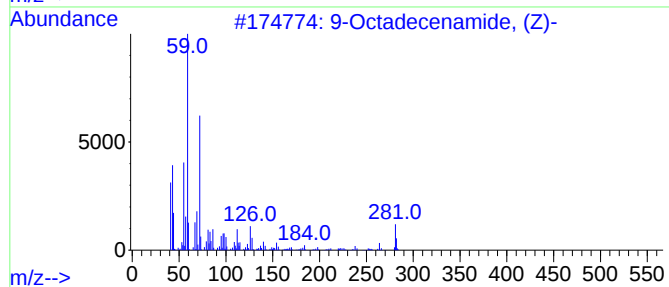
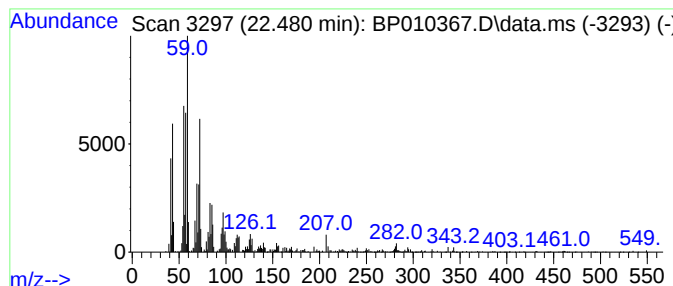
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 9-Octadecenamide, (Z)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.480	2.83 ng/ul	331820	Chrysene-d12	21.363

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	92
2			9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	90
3			13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	89
4			9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	86
5			9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	84



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: C...	21.075	2.5	ng/u1	295740	5	21.363	2340920	20.0
9-Octadecenamid...	22.480	2.8	ng/u1	331820	5	21.363	2340920	20.0