

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102521\
 Data File : BP007732.D
 Acq On : 27 Oct 2021 22:19
 Operator : CG/JU
 Sample : M4215-18
 Misc :
 ALS Vial : 89 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 JDGF5

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-BP102521.M
 Title : SVOA CALIBRATION

Signal : TIC: BP007732.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.364	9	13	22	rVB	54319	68086	1.04%	0.102%
2	3.652	58	62	69	rBV2	8672	19071	0.29%	0.029%
3	3.828	89	92	94	rBV2	9196	12229	0.19%	0.018%
4	4.028	122	126	130	rVB	13515	19420	0.30%	0.029%
5	4.117	137	141	149	rVB	16399	27937	0.43%	0.042%
6	5.040	293	298	301	rVB2	7417	11101	0.17%	0.017%
7	6.022	462	465	472	rVB6	6576	9237	0.14%	0.014%
8	6.317	510	515	521	rVB6	6098	11044	0.17%	0.017%
9	7.093	641	647	660	rVB	1104442	1732534	26.50%	2.604%
10	7.258	669	675	685	rVB	865422	1384494	21.18%	2.081%
11	7.464	702	710	718	rBV	1123050	1769252	27.06%	2.659%
12	7.934	781	790	798	rBV	882159	1422588	21.76%	2.138%
13	8.005	798	802	807	rVB8	6326	11732	0.18%	0.018%
14	8.640	899	910	921	rBV	1363504	2163612	33.10%	3.252%
15	9.087	978	986	996	rBV	1025382	1656590	25.34%	2.490%
16	9.258	1010	1015	1020	rVB8	8132	13785	0.21%	0.021%
17	9.552	1060	1065	1073	rVB6	8526	16229	0.25%	0.024%
18	9.810	1100	1109	1120	rBV	782706	1309016	20.02%	1.967%
19	10.352	1193	1201	1217	rBV	1317768	2207784	33.77%	3.318%
20	10.516	1222	1229	1242	rBV	131877	237759	3.64%	0.357%
21	10.734	1256	1266	1276	rVB	1200195	2013683	30.80%	3.026%
22	10.869	1280	1289	1300	rBV	1111434	1968134	30.11%	2.958%
23	11.393	1371	1378	1381	rBV8	2986	6908	0.11%	0.010%
24	12.046	1483	1489	1494	rBV9	4396	8015	0.12%	0.012%
25	12.322	1530	1536	1545	rVB	21248	37369	0.57%	0.056%
26	13.181	1677	1682	1688	rVB8	7805	12315	0.19%	0.019%
27	13.304	1697	1703	1706	rBV8	4089	7758	0.12%	0.012%
28	13.475	1728	1732	1738	rVV5	6769	12958	0.20%	0.019%
29	13.540	1738	1743	1750	rVB7	6817	13995	0.21%	0.021%
30	13.981	1811	1818	1832	rBV	2276729	3192996	48.84%	4.799%
31	14.263	1856	1866	1879	rBV	2552035	3843465	58.79%	5.776%
32	14.569	1910	1918	1931	rBV	1852174	2741475	41.93%	4.120%
33	14.757	1941	1950	1961	rBV	986494	1471245	22.50%	2.211%
34	14.987	1984	1989	1993	rVV4	11905	18784	0.29%	0.028%
35	15.387	2052	2057	2061	rBV4	6776	11392	0.17%	0.017%
36	15.563	2079	2087	2099	rBV	3307274	4906933	75.06%	7.374%

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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-BP102521.M
 Title : SVOA CALIBRATION

37	15.675	2099	2106	2119	rVV	848139	1171756	17.92%	1.761%
38	15.804	2123	2128	2135	rVB	41074	66011	1.01%	0.099%
39	16.016	2158	2164	2169	rBV10	8289	14897	0.23%	0.022%
40	16.063	2169	2172	2175	rBV4	6239	7250	0.11%	0.011%
41	16.134	2180	2184	2188	rBV7	4859	8032	0.12%	0.012%
42	16.251	2199	2204	2209	rBV8	4401	8224	0.13%	0.012%
43	16.351	2218	2221	2228	rVB9	7571	14487	0.22%	0.022%
44	16.469	2236	2241	2243	rBV6	5671	6924	0.11%	0.010%
45	16.998	2327	2331	2334	rBV6	5860	7103	0.11%	0.011%
46	17.104	2337	2349	2353	rBV2	12568	28425	0.43%	0.043%
47	17.257	2370	2375	2379	rBV7	4581	7772	0.12%	0.012%
48	17.322	2379	2386	2395	rBV	2321305	3322859	50.83%	4.994%
49	17.422	2396	2403	2415	rVV2	3660891	5242992	80.20%	7.879%
50	17.522	2415	2420	2426	rVB8	15527	28482	0.44%	0.043%
51	17.692	2447	2449	2456	rVB6	7167	13570	0.21%	0.020%
52	18.004	2496	2502	2508	rBV	52193	67860	1.04%	0.102%
53	18.216	2534	2538	2545	rBV	94681	130934	2.00%	0.197%
54	18.287	2547	2550	2554	rVB	14873	20247	0.31%	0.030%
55	19.234	2707	2711	2715	rBV2	56571	74932	1.15%	0.113%
56	19.304	2719	2723	2727	rVV	39598	51917	0.79%	0.078%
57	19.451	2744	2748	2753	rBV4	32702	48657	0.74%	0.073%
58	19.598	2770	2773	2777	rBV5	15760	17215	0.26%	0.026%
59	19.657	2777	2783	2790	rBV	4926489	6424437	98.27%	9.655%
60	21.133	3029	3034	3043	rBV2	250407	450400	6.89%	0.677%
61	21.416	3076	3082	3095	rBV	2946315	4122737	63.06%	6.196%
62	23.698	3463	3470	3479	rVB2	3266866	6537460	100.00%	9.825%
63	23.857	3490	3497	3511	rVB2	1986620	4273901	65.38%	6.423%

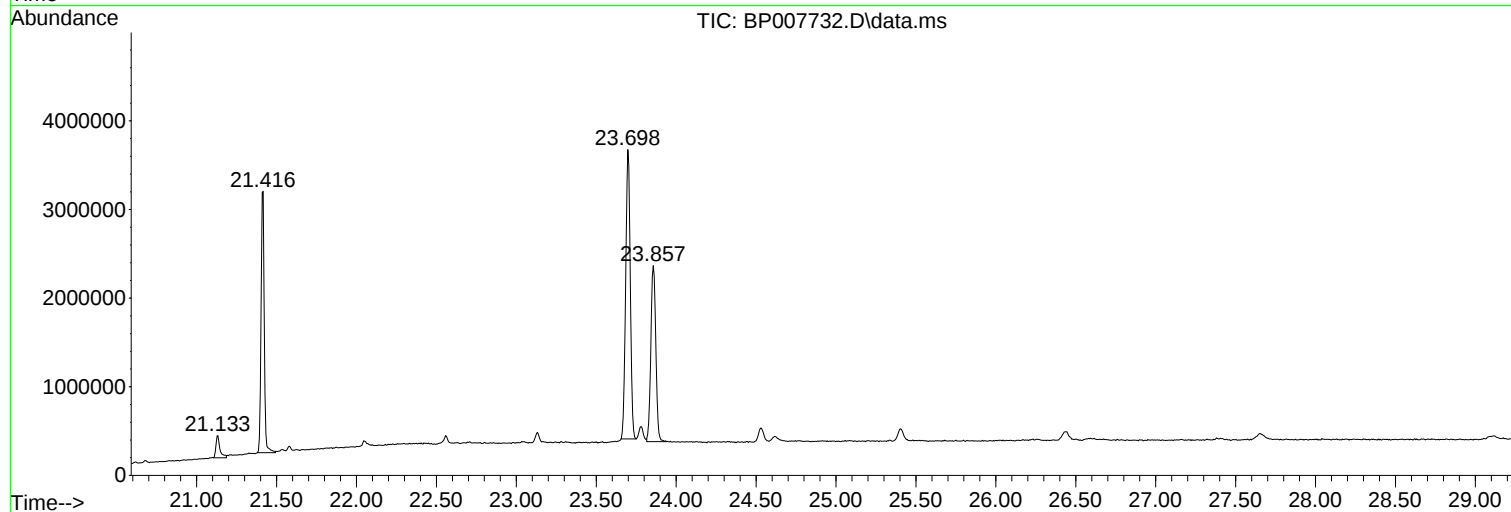
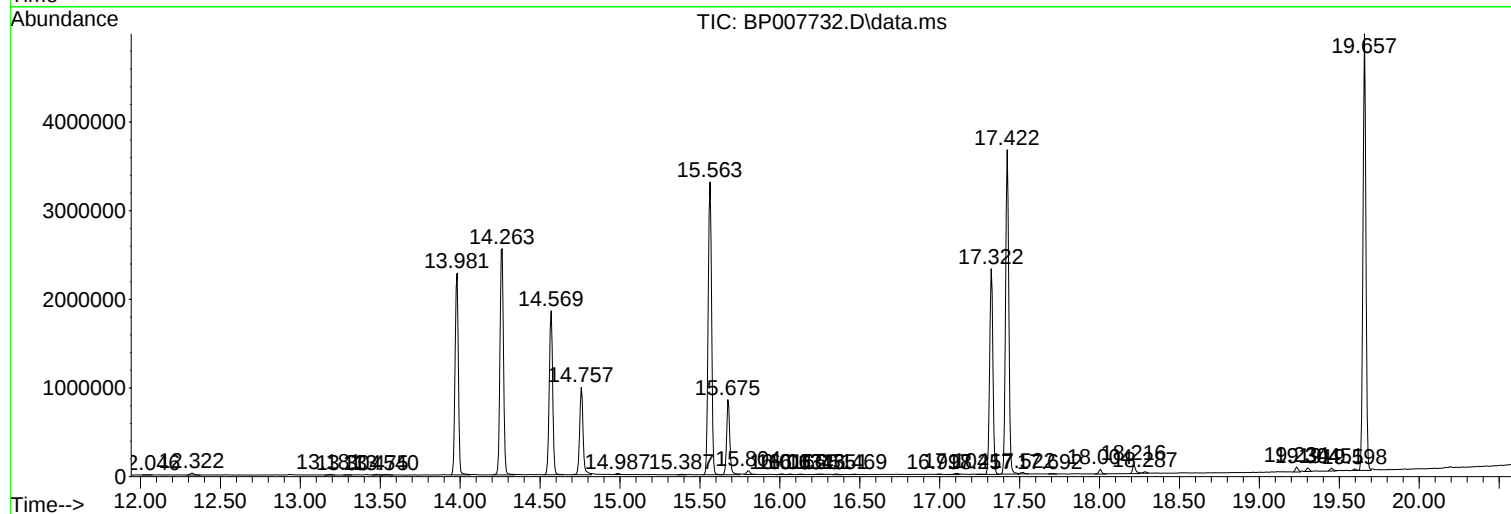
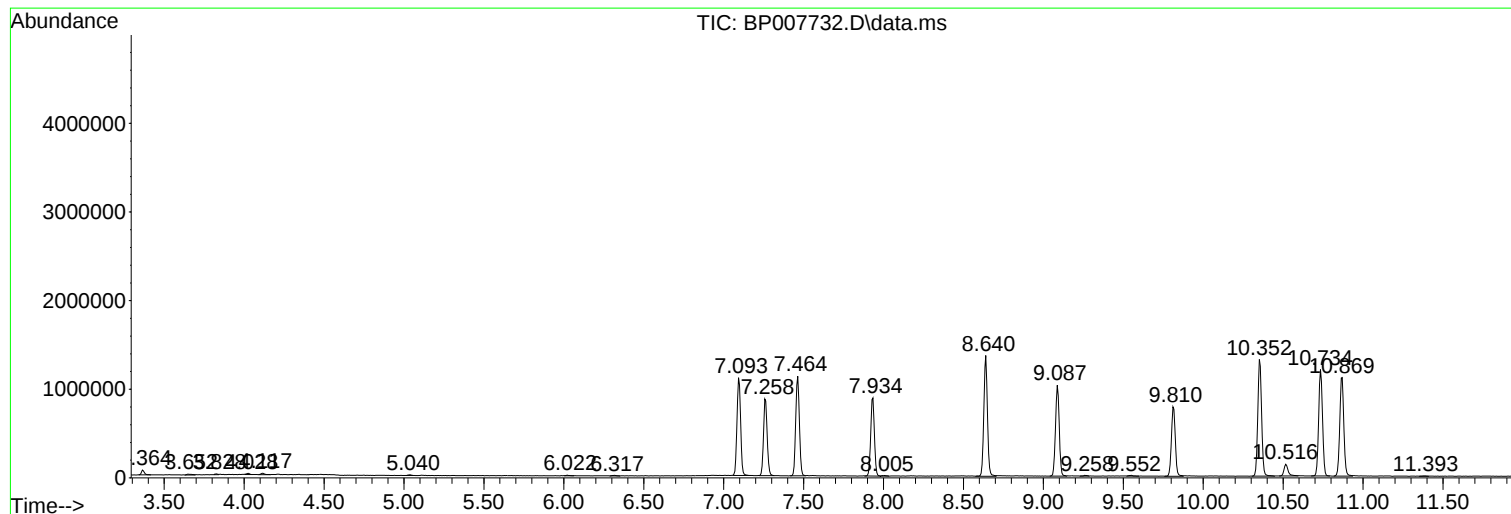
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
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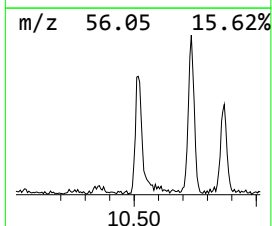
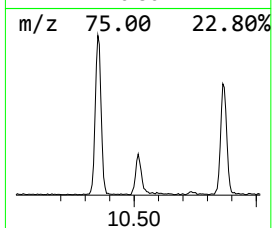
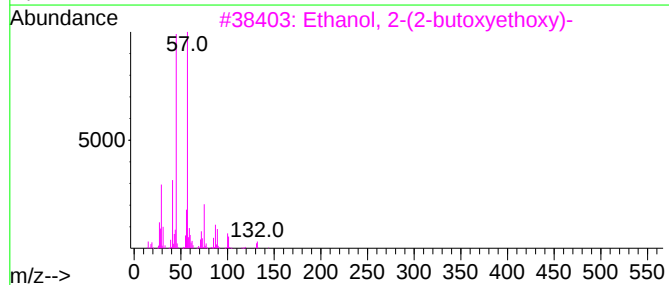
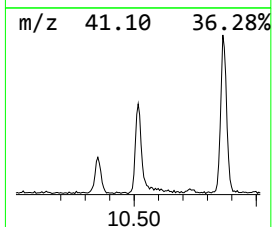
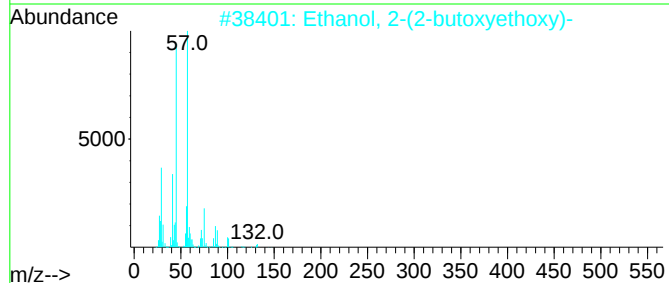
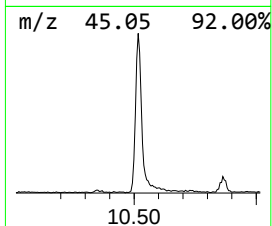
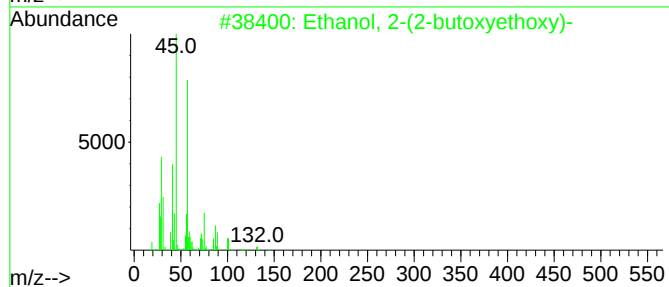
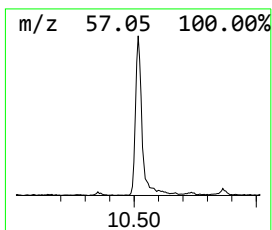
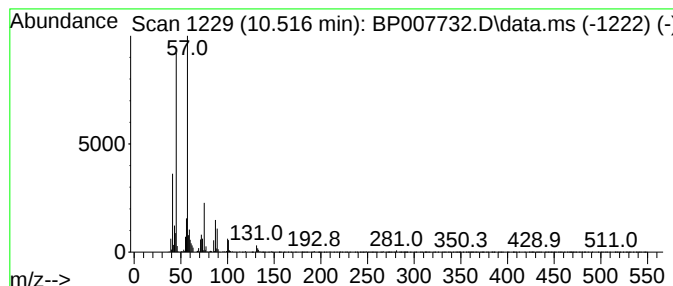
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.516	2.36 ng/ul	237759	Naphthalene-d8	10.734

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
3			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	83
4			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	83
5			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	83



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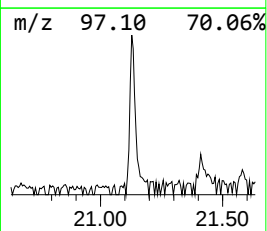
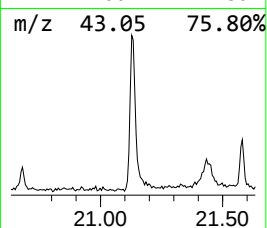
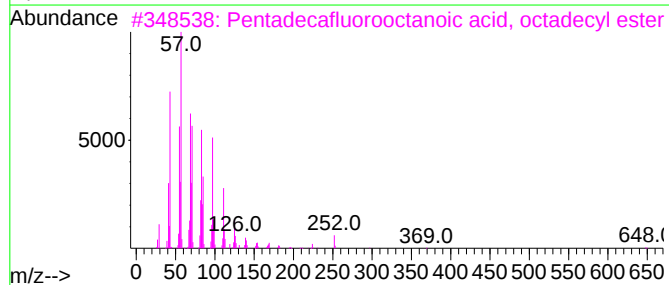
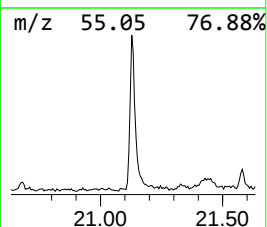
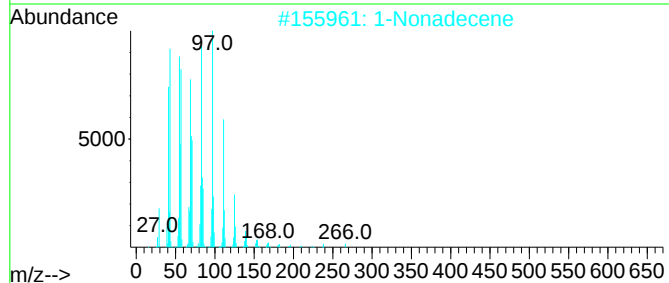
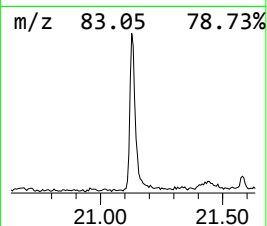
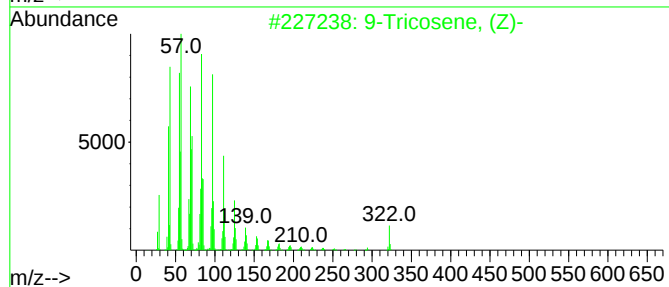
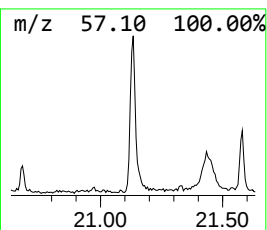
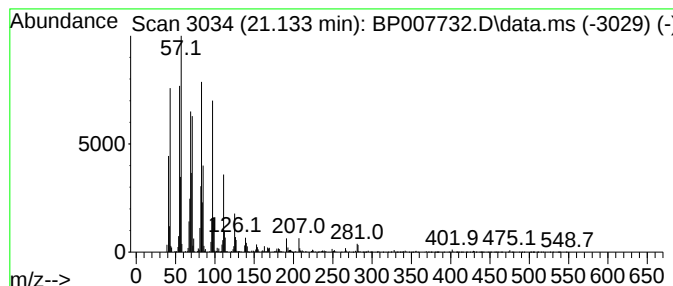
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

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.	
21.133	2.18 ng/ul	450400	Chrysene-d12	21.416	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Tricosene, (Z)-	322	C23H46	027519-02-4	95
2	1-Nonadecene	266	C19H38	018435-45-5	94
3	Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	94
4	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
5	Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	94



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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
Ethanol, 2-(2-b...	10.516	2.4	ng/u1	237759	2	10.734	2013680	20.0
(DEL) Alkane: S...	21.133	2.2	ng/u1	450400	5	21.416	4122740	20.0