

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF102817\
 Data File : BF099916.D
 Acq On : 28 Oct 2017 14:44
 Operator : SJ/JU
 Sample : I6071-03
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ENV-116-5(10-15)

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102317.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.010	494	497	514	rBV	183605	293968	12.20%	1.593%
2	5.410	562	565	592	rBV	1594947	1839594	76.31%	9.971%
3	6.433	735	739	757	rBV	1734821	1865496	77.39%	10.111%
4	6.563	757	761	768	rVB	2016512	1934457	80.25%	10.485%
5	6.786	796	799	808	rBV	647908	585193	24.28%	3.172%
6	6.939	821	825	844	rBB	1765029	1615450	67.02%	8.756%
7	7.357	892	896	909	rBV	1269102	1187506	49.26%	6.437%
8	8.069	1013	1017	1033	rBV	947697	816360	33.87%	4.425%
9	9.145	1195	1200	1203	rBV	2591933	2410540	100.00%	13.066%
10	9.251	1215	1218	1223	rVB	97276	78617	3.26%	0.426%
11	9.539	1264	1267	1273	rVB	323834	258248	10.71%	1.400%
12	9.822	1311	1315	1325	rVB2	988863	857331	35.57%	4.647%
13	10.239	1384	1386	1389	rVB	43084	29463	1.22%	0.160%
14	10.457	1417	1423	1425	rBV3	20308	25080	1.04%	0.136%
15	10.539	1434	1437	1441	rVB	88447	77733	3.22%	0.421%
16	10.616	1447	1450	1460	rBV	1007274	963031	39.95%	5.220%
17	10.710	1463	1466	1476	rVB	49742	67029	2.78%	0.363%
18	11.316	1565	1569	1578	rBV	799580	722444	29.97%	3.916%
19	11.827	1653	1656	1659	rBV	36305	34648	1.44%	0.188%
20	11.863	1659	1662	1668	rVB	66073	52161	2.16%	0.283%
21	12.898	1833	1838	1841	rBV	1701187	1541759	63.96%	8.357%
22	13.774	1983	1987	1993	rBV2	70669	77751	3.23%	0.421%
23	13.962	2015	2019	2026	rBV	529244	498943	20.70%	2.704%
24	15.427	2263	2268	2278	rVB	408145	537069	22.28%	2.911%
25	15.809	2328	2333	2338	rVB3	18697	29280	1.21%	0.159%
26	16.292	2411	2415	2419	rBV3	15937	25420	1.05%	0.138%
27	17.527	2621	2625	2631	rVB7	11621	24688	1.02%	0.134%

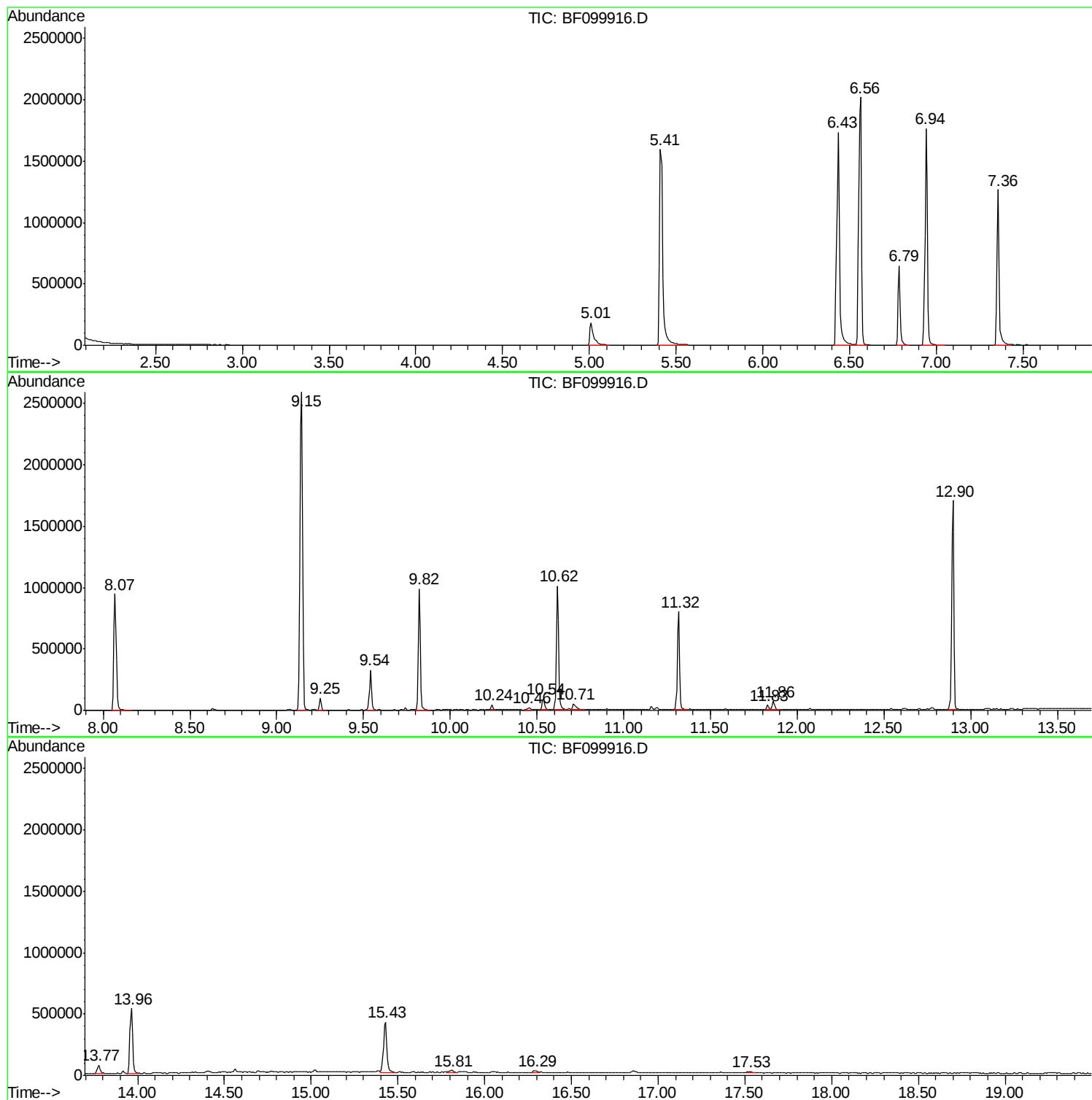
Sum of corrected areas: 18449259

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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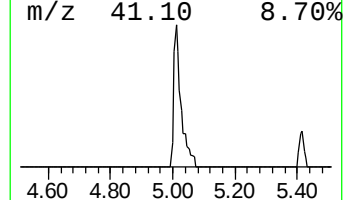
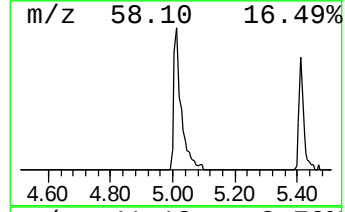
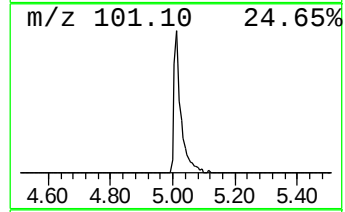
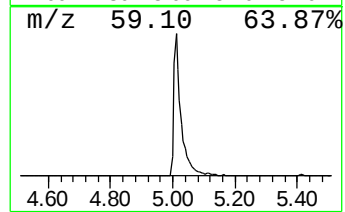
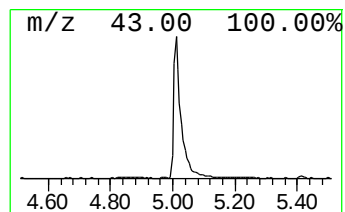
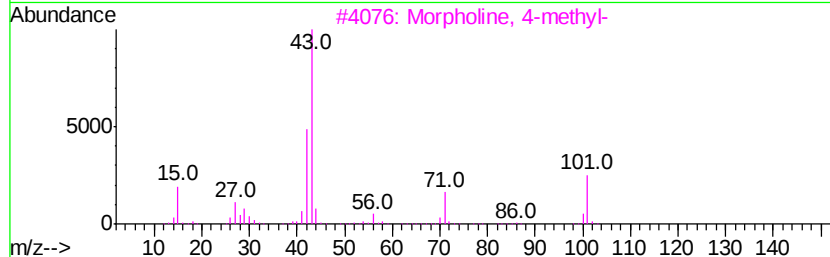
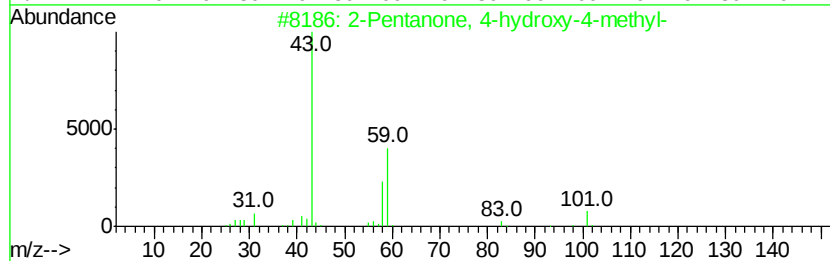
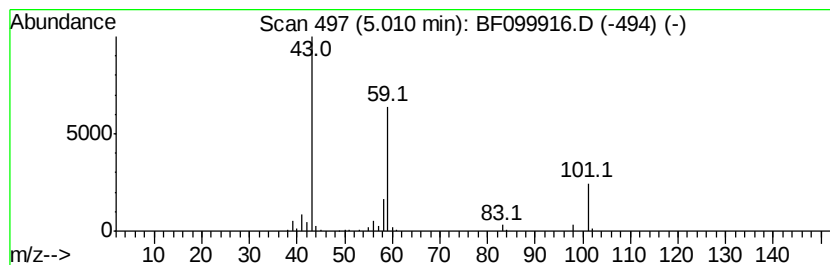
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.01	10.05 ng	293968	1,4-Dichlorobenzene-d4	6.79

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
3			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4			Guanidine	59	CH5N3	000113-00-8	9
5			(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9



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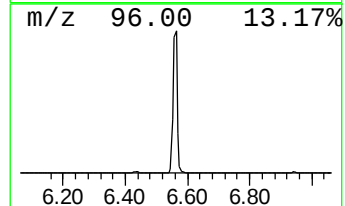
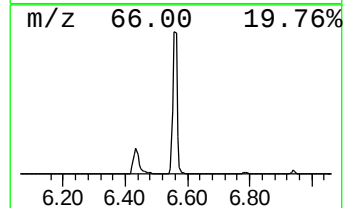
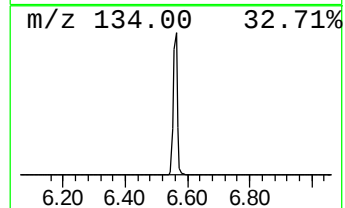
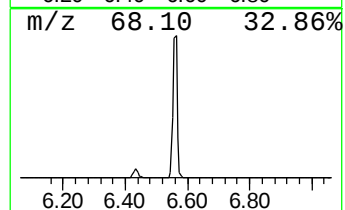
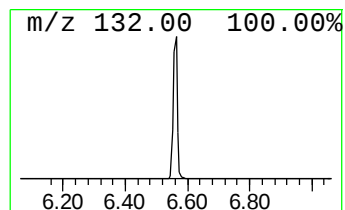
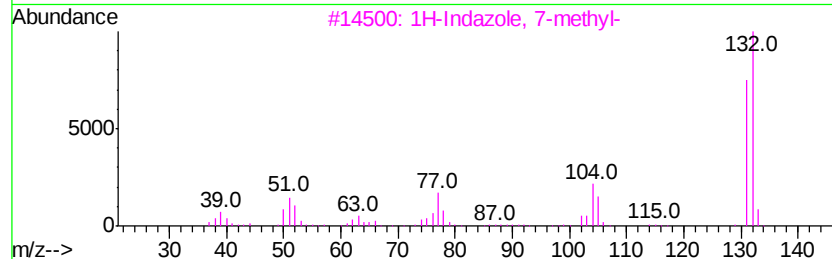
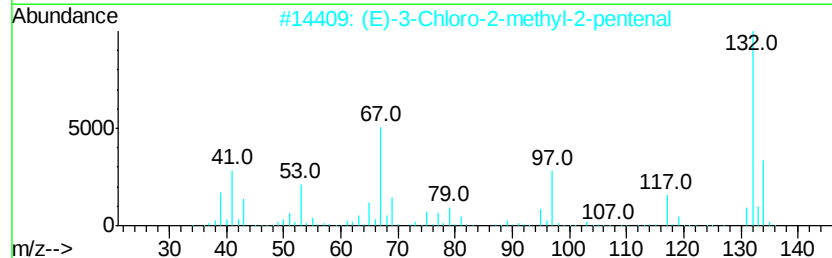
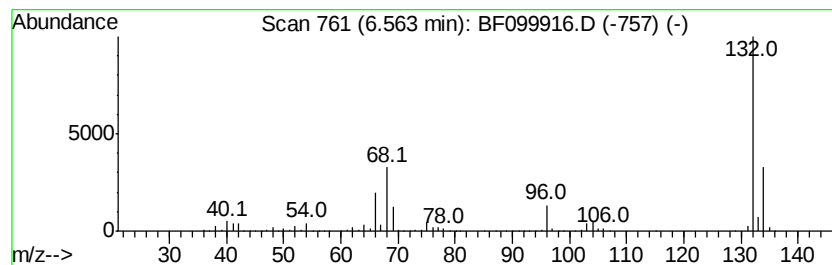
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Peak Number 2 unknown6.56 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.56	66.11 ng	1934460	1,4-Dichlorobenzene-d4	6.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		1H-Indazole, 7-methyl-	132	C8H8N2	1000316-00-7	10
4		Xenon	132	Xe	007440-63-3	9
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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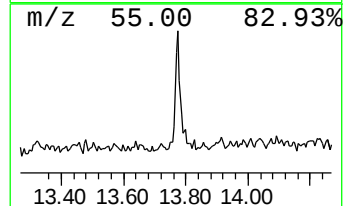
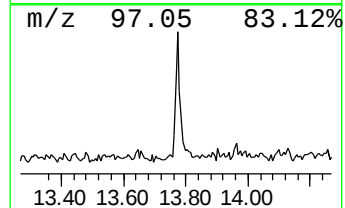
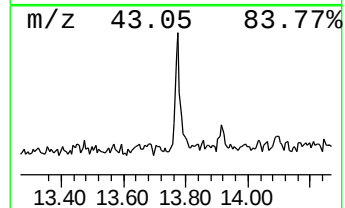
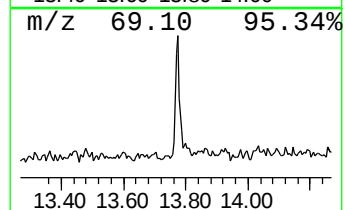
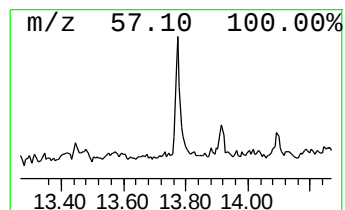
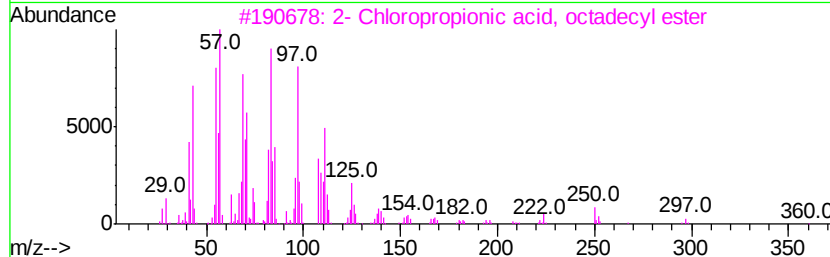
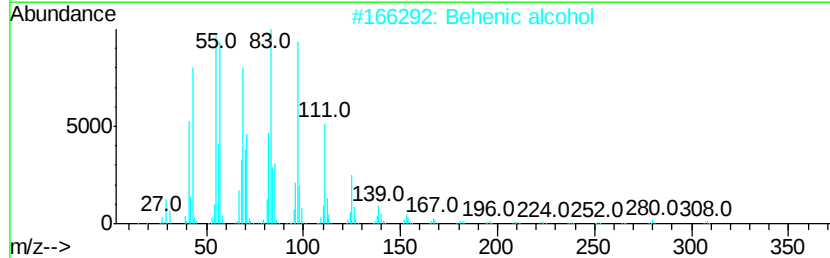
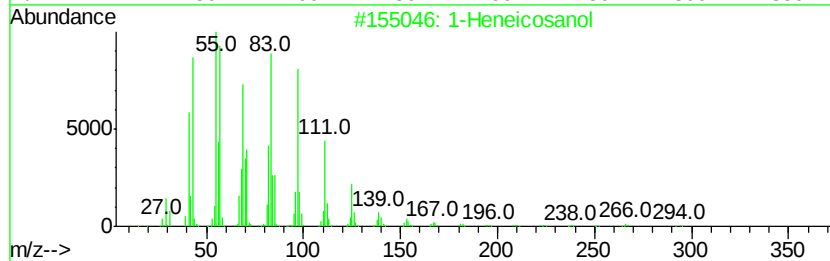
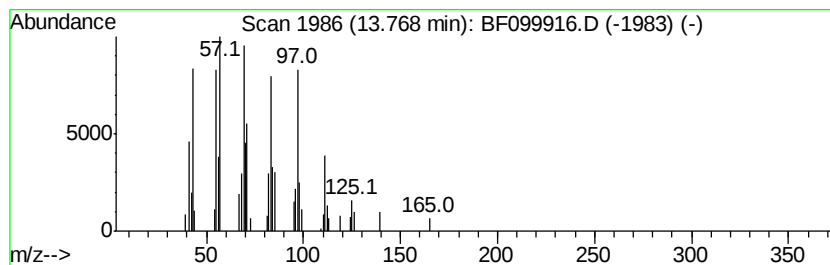
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Peak Number 4 1-Heneicosanol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.77	3.12 ng	77751	Chrysene-d12	13.96

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosanol		312	C21H44O	015594-90-8	95
2	Behenic alcohol		326	C22H46O	000661-19-8	94
3	2- Chloropropionic acid, octadec...		360	C21H41ClO2	088104-31-8	94
4	n-Tetracosanol-1		354	C24H50O	000506-51-4	93
5	Heptadecyl trifluoroacetate		352	C19H35F3O2	1000351-87-0	93



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.01	10.1	ng	293968	1	6.79	585193	20.0
unknown6.56	6.56	66.1	ng	1934460	1	6.79	585193	20.0
1-Heneicosanol	13.77	3.1	ng	77751	5	13.96	498943	20.0