

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
 Data File : BP013188.D
 Acq On : 21 Dec 2022 03:52
 Operator : CG/JU
 Sample : N5984-01
 Misc :
 ALS Vial : 70 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 COLC3

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-BP120722.M
 Title : SVOA CALIBRATION

Signal : TIC: BP013188.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.334	22	25	34	rVB	160465	204245	1.79%	0.142%
2	3.764	95	98	114	rBV	2127160	3090537	27.11%	2.143%
3	4.093	150	154	159	rVB	38253	49159	0.43%	0.034%
4	5.040	312	315	325	rVB	32965	58459	0.51%	0.041%
5	6.928	631	636	640	rBV5	14450	26979	0.24%	0.019%
6	7.111	657	667	680	rBV	2543467	3988201	34.98%	2.766%
7	7.275	689	695	704	rBV	2004695	3165116	27.76%	2.195%
8	7.481	723	730	741	rBV	2547544	3955273	34.69%	2.743%
9	7.952	803	810	817	rBV	2112025	3242128	28.44%	2.249%
10	8.011	817	820	826	rVB3	13828	22108	0.19%	0.015%
11	8.664	921	931	943	rBV	3076136	4964096	43.54%	3.443%
12	9.122	1000	1009	1021	rBV	2383995	3932953	34.49%	2.728%
13	9.228	1023	1027	1031	rVV6	13367	25412	0.22%	0.018%
14	9.275	1031	1035	1042	rVB10	12596	22973	0.20%	0.016%
15	9.522	1071	1077	1085	rVB	54067	85035	0.75%	0.059%
16	9.846	1124	1132	1145	rBV	2051052	3342447	29.32%	2.318%
17	10.063	1164	1169	1176	rBV2	36386	63304	0.56%	0.044%
18	10.381	1214	1223	1239	rBV	3110106	5259935	46.13%	3.648%
19	10.540	1247	1250	1257	rBV9	5800	12015	0.11%	0.008%
20	10.763	1281	1288	1298	rVB	2860659	4824707	42.32%	3.346%
21	10.905	1304	1312	1326	rBV	2914476	4935208	43.28%	3.423%
22	11.546	1418	1421	1428	rVB7	13371	24695	0.22%	0.017%
23	12.016	1496	1501	1506	rVB8	10415	17844	0.16%	0.012%
24	12.081	1506	1512	1517	rBV7	8052	15380	0.13%	0.011%
25	12.175	1523	1528	1531	rBV5	14880	22782	0.20%	0.016%
26	12.346	1551	1557	1563	rBV2	57108	96960	0.85%	0.067%
27	12.863	1639	1645	1651	rBV7	27306	51605	0.45%	0.036%
28	13.116	1686	1688	1696	rVB9	8143	15361	0.13%	0.011%
29	13.193	1696	1701	1706	rVB9	6137	12185	0.11%	0.008%
30	13.410	1732	1738	1744	rVB	42950	67149	0.59%	0.047%
31	13.487	1746	1751	1755	rBV7	9273	17213	0.15%	0.012%
32	13.563	1757	1764	1768	rVB6	12697	22450	0.20%	0.016%
33	13.916	1818	1824	1828	rBV9	8098	14521	0.13%	0.010%
34	13.999	1828	1838	1847	rBV	5258942	7350483	64.47%	5.098%
35	14.110	1853	1857	1864	rVB8	18156	32251	0.28%	0.022%
36	14.240	1875	1879	1880	rBV4	9748	14651	0.13%	0.010%

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Peak Location: TOP

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
 Title : SVOA CALIBRATION

37	14.287	1880	1887	1900	rVV	6250006	9114335	79.94%	6.321%
38	14.481	1916	1920	1925	rVB3	19790	32143	0.28%	0.022%
39	14.593	1928	1939	1949	rBV	4520856	6600838	57.89%	4.578%
40	14.787	1962	1972	1997	rBV	1811294	3025074	26.53%	2.098%
41	14.998	2003	2008	2022	rBV2	246903	376323	3.30%	0.261%
42	15.587	2101	2108	2120	rBV	7788814	11150919	97.80%	7.734%
43	15.704	2122	2128	2143	rVV	1847952	2563794	22.49%	1.778%
44	15.828	2145	2149	2154	rVB3	37522	62941	0.55%	0.044%
45	15.951	2164	2170	2176	rBV	1180871	1565739	13.73%	1.086%
46	16.369	2238	2241	2248	rVB8	12106	23169	0.20%	0.016%
47	16.522	2264	2267	2273	rBV8	7117	12432	0.11%	0.009%
48	16.734	2299	2303	2309	rBV8	30793	50334	0.44%	0.035%
49	17.140	2368	2372	2375	rVB4	11424	15538	0.14%	0.011%
50	17.351	2401	2408	2418	rBV	5267406	7290694	63.94%	5.057%
51	17.451	2418	2425	2437	rVV	7731734	11049770	96.91%	7.664%
52	17.540	2437	2440	2446	rVB6	67477	100680	0.88%	0.070%
53	17.728	2468	2472	2476	rBV6	29576	38955	0.34%	0.027%
54	18.222	2551	2556	2565	rBV	381130	574285	5.04%	0.398%
55	18.628	2622	2625	2631	rBV5	37765	49640	0.44%	0.034%
56	19.257	2728	2732	2736	rBV2	94006	128235	1.12%	0.089%
57	19.328	2739	2744	2748	rBV	96170	157509	1.38%	0.109%
58	19.451	2761	2765	2769	rBV4	70739	99368	0.87%	0.069%
59	19.681	2798	2804	2816	rBV	8772008	11401671	100.00%	7.908%
60	20.969	3020	3023	3028	rBV	217931	228626	2.01%	0.159%
61	21.128	3046	3050	3057	rBV	1057749	1494111	13.10%	1.036%
62	21.433	3097	3102	3112	rBV	4540734	6156089	53.99%	4.270%
63	23.757	3490	3497	3509	rVB	5342989	11004524	96.52%	7.632%
64	23.916	3518	3524	3535	rVB2	3342707	6768756	59.37%	4.695%

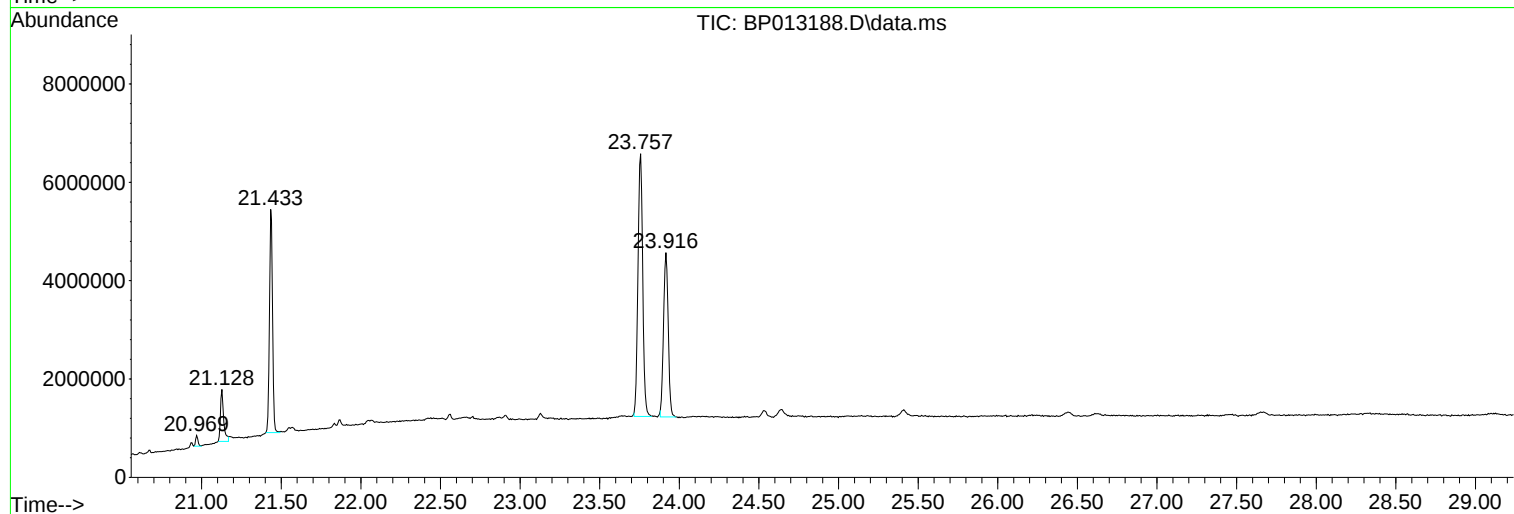
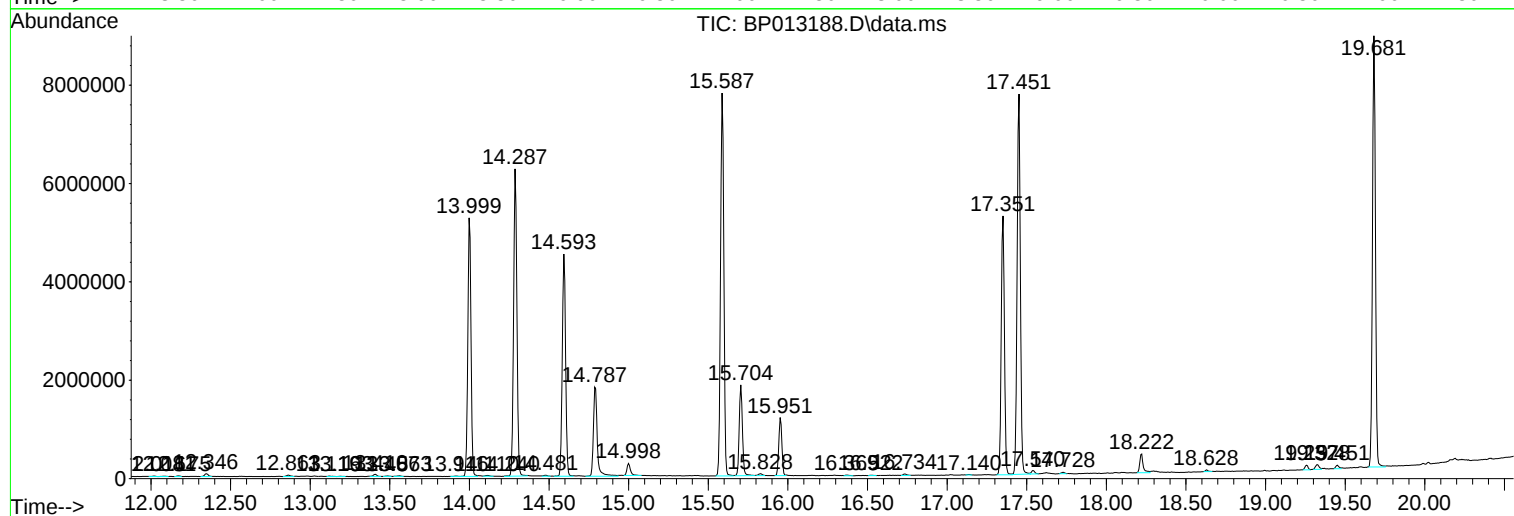
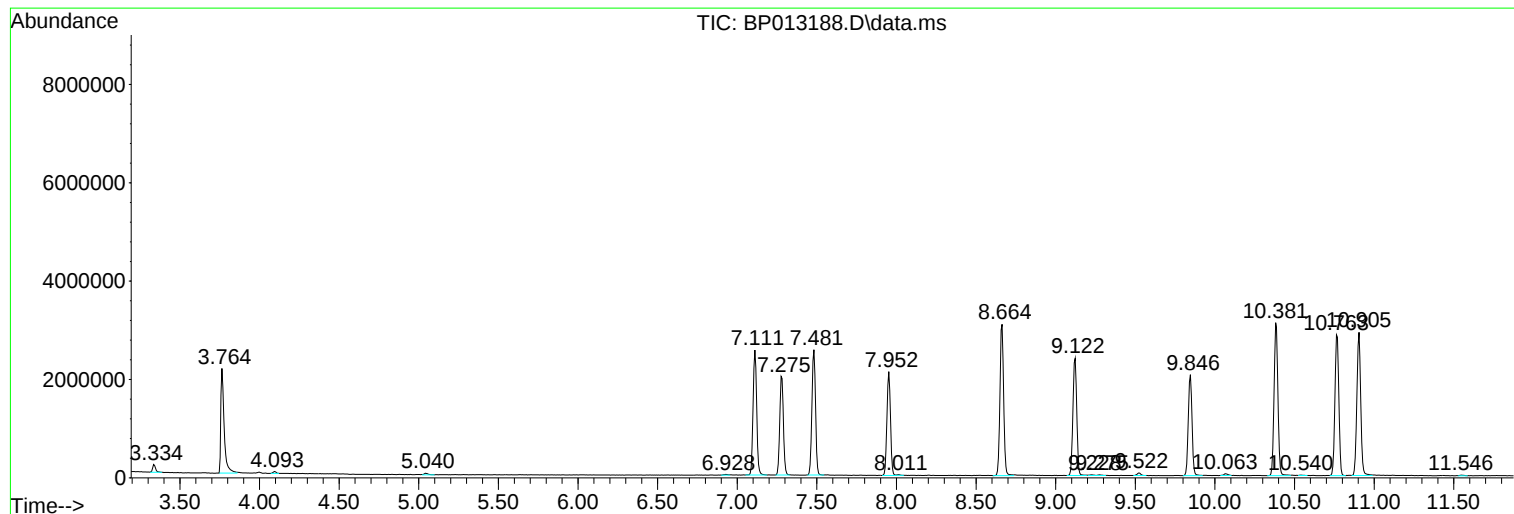
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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



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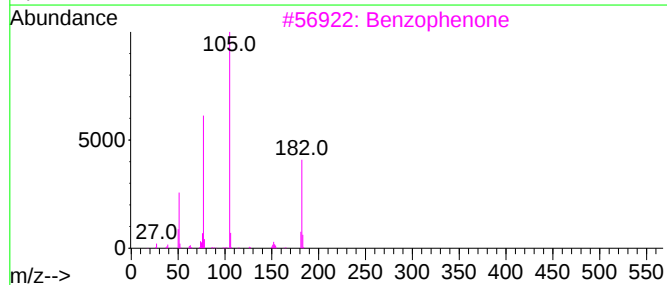
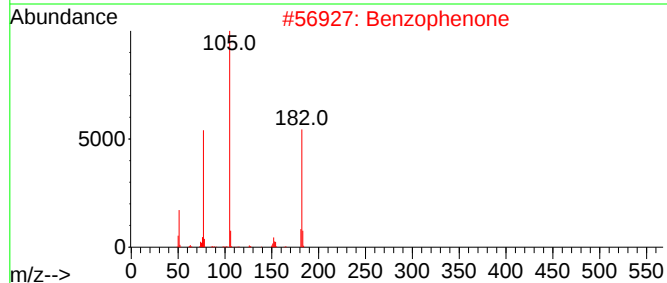
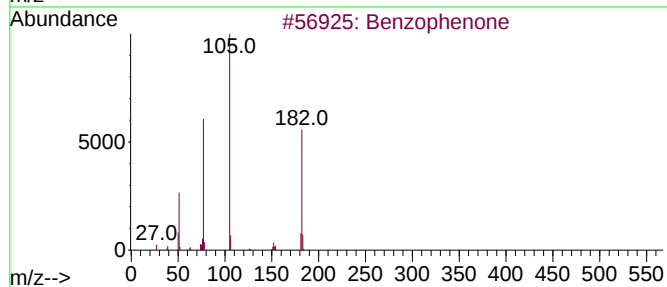
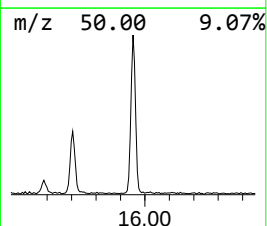
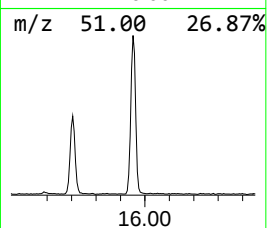
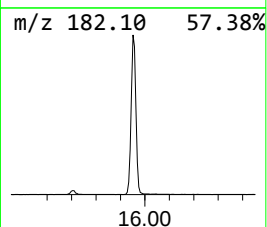
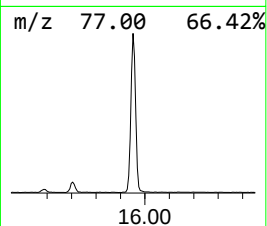
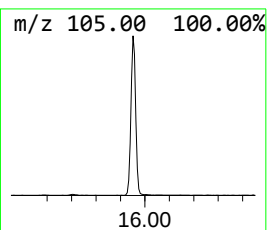
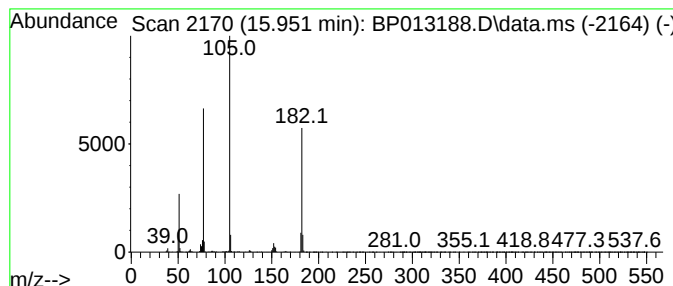
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzophenone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.951	4.74 ng/ul	1565740	Acenaphthene-d10	14.593

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzophenone	182	C13H10O	000119-61-9	95
3			Benzophenone	182	C13H10O	000119-61-9	94
4			Benzophenone	182	C13H10O	000119-61-9	90
5			Benzophenone	182	C13H10O	000119-61-9	90



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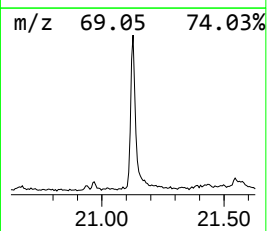
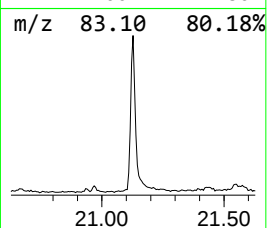
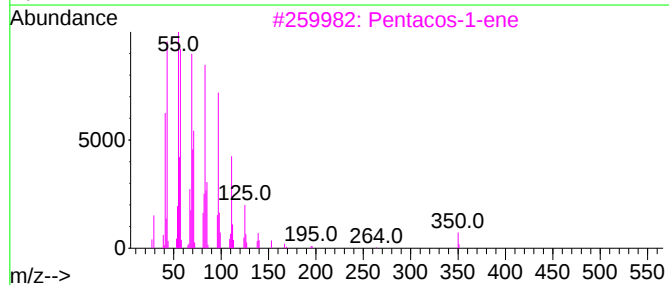
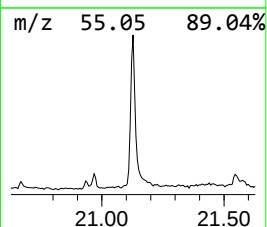
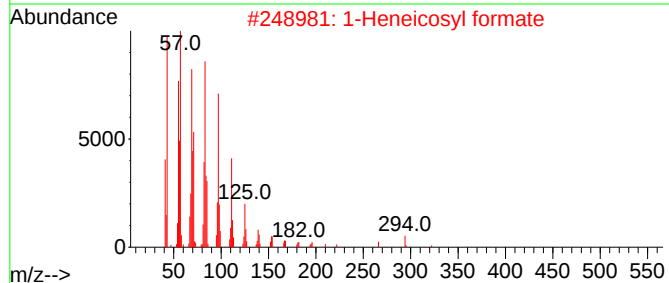
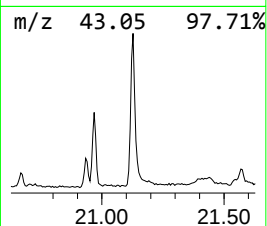
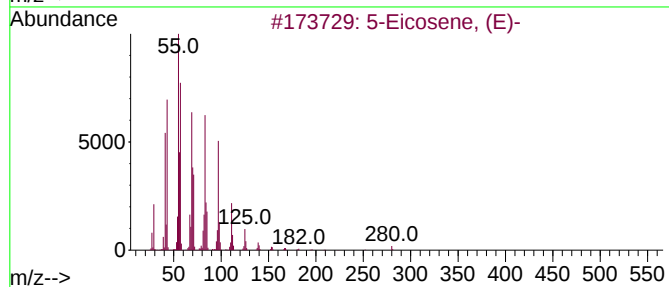
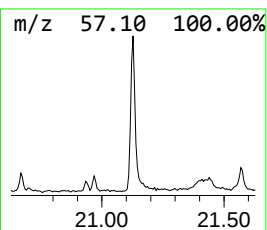
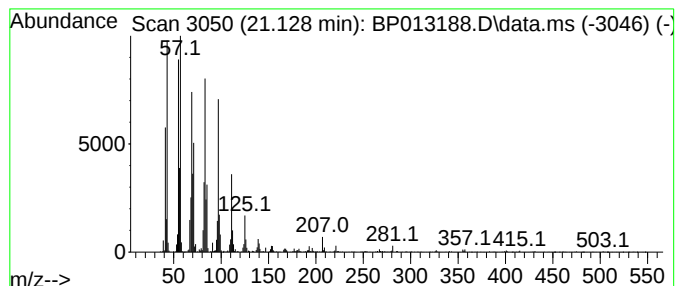
Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.128	4.85 ng/ul	1494110	Chrysene-d12	21.433	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Eicosene, (E)-	280	C20H40	074685-30-6	97
2	1-Heneicosyl formate	340	C22H44O2	077899-03-7	97
3	Pentacos-1-ene	350	C25H50	016980-85-1	94
4	Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	94
5	n-Tetracosanol-1	354	C24H50O	000506-51-4	94



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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
Benzophenone	15.951	4.7	ng/u1	1565740	3	14.593	6600840	20.0
(DEL) Alkane: S...	21.128	4.8	ng/u1	1494110	5	21.433	6156090	20.0