

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063017\
 Data File : BF096368.D
 Acq On : 1 Jul 2017 6:07
 Operator : SJ/MA
 Sample : I3947-01
 Misc :
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP1-TP2-COMP

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-BF062717.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	4.951	483	487	494	rBV	360596	443739	14.47%	1.576%
2	5.375	554	559	562	rBV	3095232	2940835	95.90%	10.444%
3	6.393	727	732	742	rVB2	2801230	3066685	100.00%	10.891%
4	6.528	751	755	759	rBV	2864041	2760637	90.02%	9.804%
5	6.763	791	795	798	rBV	1682681	1486775	48.48%	5.280%
6	6.916	817	821	825	rBV	2007142	1762214	57.46%	6.259%
7	7.316	884	889	892	rBV	1734688	1563259	50.98%	5.552%
8	7.416	901	906	910	rBV	36547	40183	1.31%	0.143%
9	7.751	957	963	965	rBV	61978	72000	2.35%	0.256%
10	8.045	1008	1013	1016	rBV	2032297	1836305	59.88%	6.522%
11	9.116	1191	1195	1199	rBV	2103980	1963321	64.02%	6.973%
12	9.210	1206	1211	1214	rBV3	29611	34590	1.13%	0.123%
13	9.510	1259	1262	1265	rBV	462167	349475	11.40%	1.241%
14	9.739	1298	1301	1306	rVB	60645	48799	1.59%	0.173%
15	9.798	1306	1311	1314	rBV	1809307	1661200	54.17%	5.900%
16	10.239	1382	1386	1388	rVB	78024	55021	1.79%	0.195%
17	10.457	1419	1423	1429	rBV2	30184	44728	1.46%	0.159%
18	10.580	1439	1444	1447	rBV	1653644	1486163	48.46%	5.278%
19	10.680	1457	1461	1463	rBV	129095	103152	3.36%	0.366%
20	10.710	1463	1466	1473	rVB	82558	96602	3.15%	0.343%
21	11.157	1538	1542	1545	rBV2	51942	47742	1.56%	0.170%
22	11.275	1558	1562	1565	rBV2	1695540	1539874	50.21%	5.469%
23	11.298	1565	1566	1569	rVV	122780	65313	2.13%	0.232%
24	11.345	1569	1574	1577	rVB2	27073	32376	1.06%	0.115%
25	11.810	1649	1653	1657	rBV	146363	150842	4.92%	0.536%
26	12.322	1737	1740	1744	rBV5	35349	42786	1.40%	0.152%
27	12.480	1764	1767	1770	rBV	178628	162222	5.29%	0.576%
28	12.710	1800	1806	1809	rBV	175533	196997	6.42%	0.700%
29	12.857	1827	1831	1835	rBV	1980361	1874650	61.13%	6.658%
30	12.933	1840	1844	1847	rBV6	18362	36578	1.19%	0.130%
31	13.757	1981	1984	1990	rBV2	142898	139467	4.55%	0.495%
32	13.904	2005	2009	2012	rBV	1231924	1227646	40.03%	4.360%
33	15.327	2247	2251	2256	rVB	658308	824883	26.90%	2.930%

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Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % 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:\HPCHEM1\BNA_F\METHODS\8270-BF062717.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

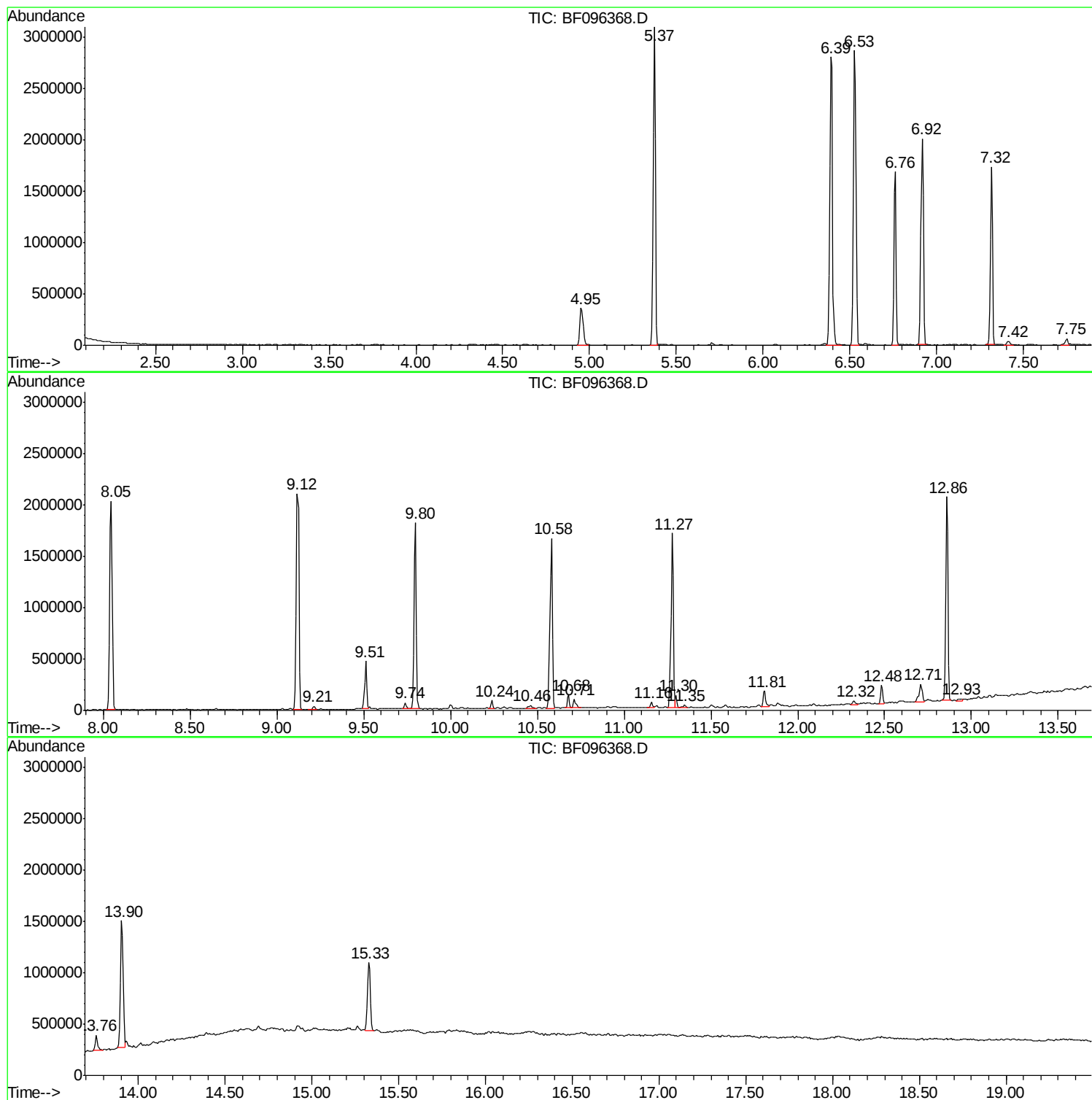
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF062717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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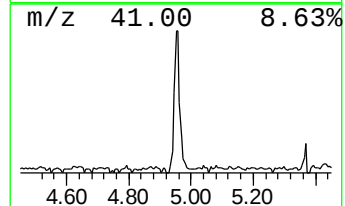
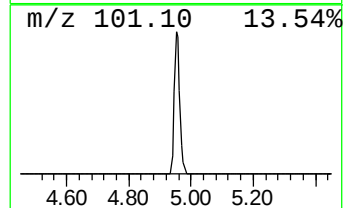
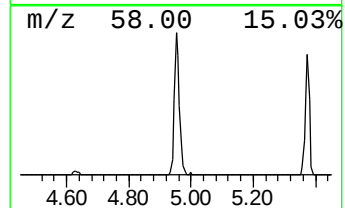
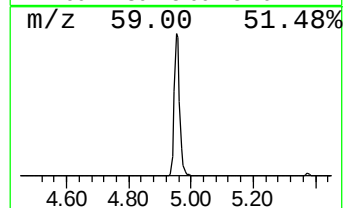
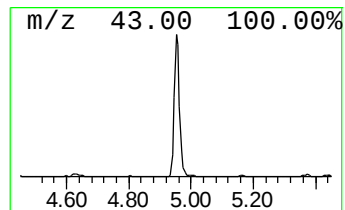
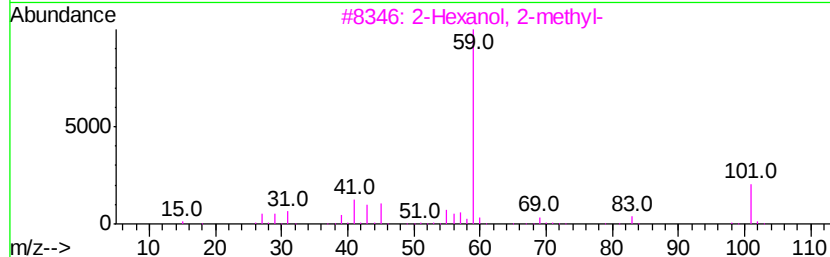
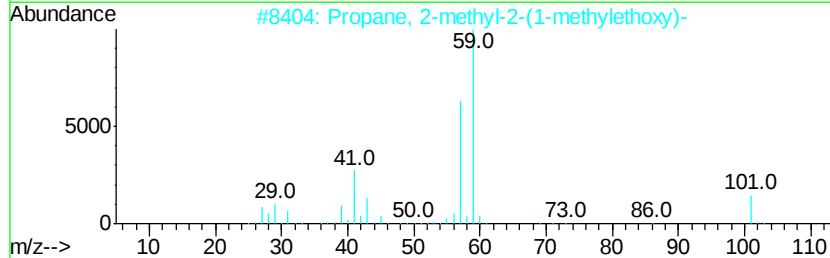
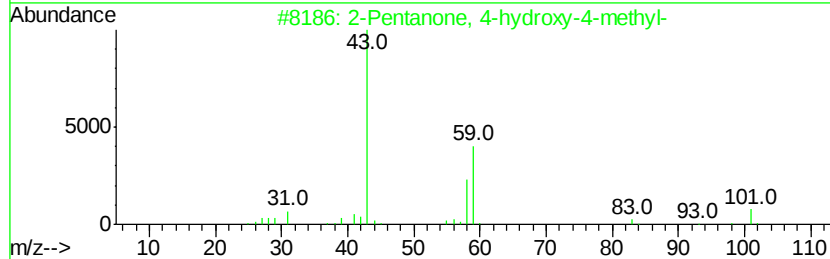
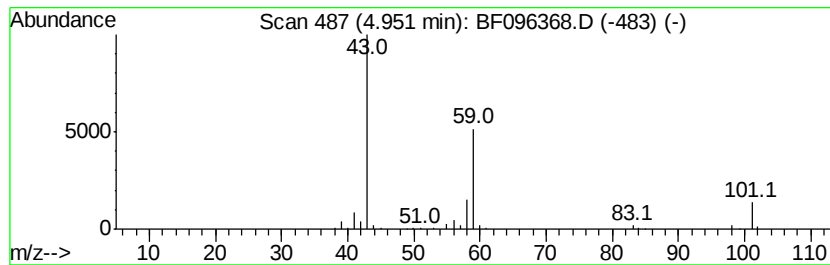
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.95	5.97 ng	443739	1,4-Dichlorobenzene-d4	6.76

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	42
3			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	9



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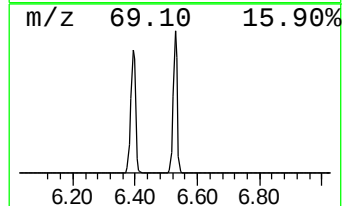
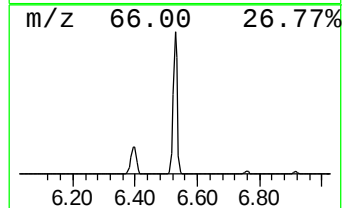
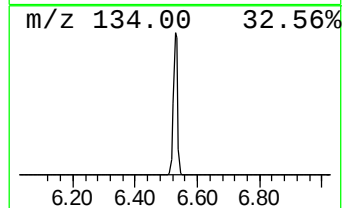
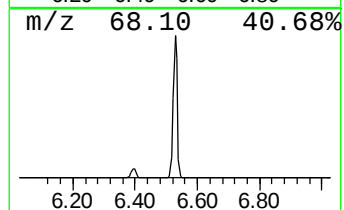
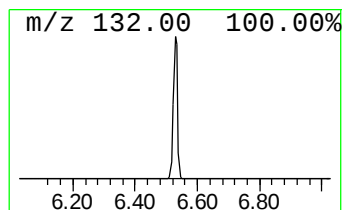
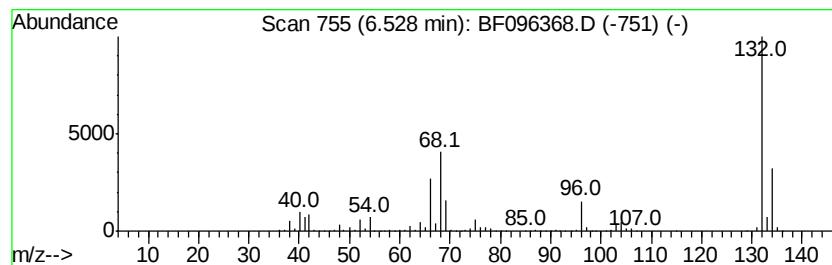
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Peak Number 2 unknown6.53 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.53	37.14 ng	2760640	1,4-Dichlorobenzene-d4	6.76

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		1-Aminoindan	133	C9H11N	034698-41-4	9



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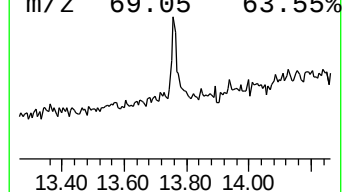
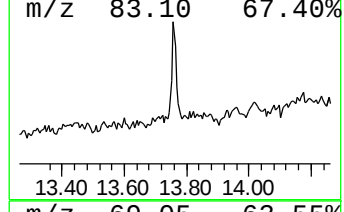
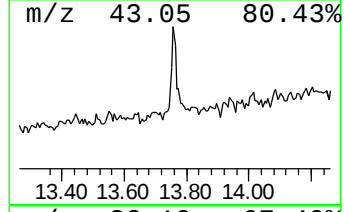
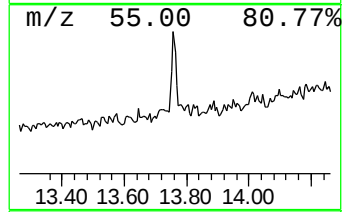
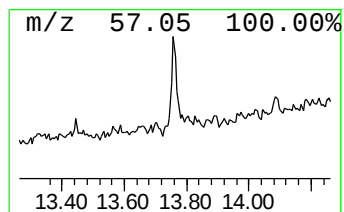
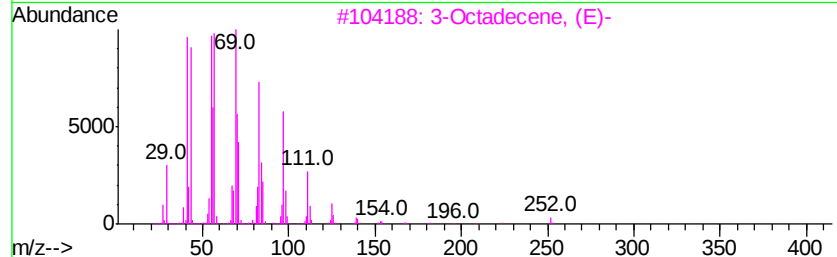
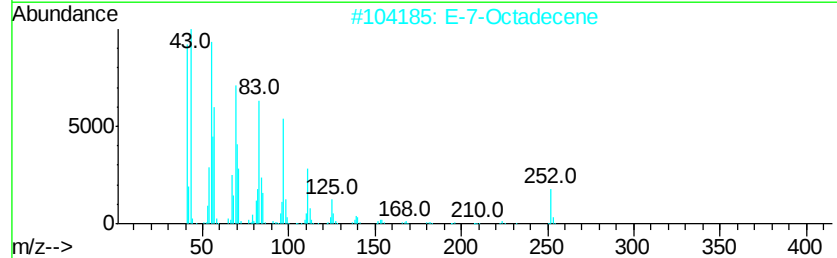
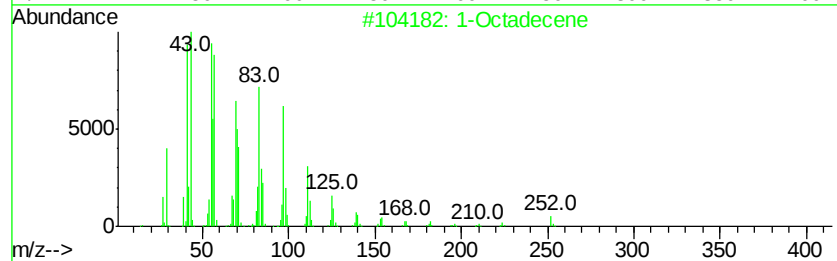
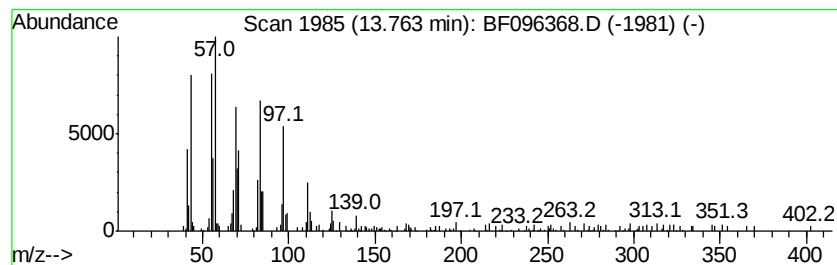
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TIC Integration Parameters: LSCINT.P

Peak Number 5 1-Octadecene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.76	2.27 ng	139467	Chrysene-d12	13.90	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Octadecene	252	C18H36	000112-88-9	96
2	E-7-Octadecene	252	C18H36	1000130-92-0	93
3	3-Octadecene, (E)-	252	C18H36	007206-19-1	93
4	1-Docosene	308	C22H44	001599-67-3	93
5	8-Heptadecene	238	C17H34	002579-04-6	91



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.95	6.0	ng	443739	1	6.76	1486780	20.0
unknown6.53	6.53	37.1	ng	2760640	1	6.76	1486780	20.0
1-Octadecene	13.76	2.3	ng	139467	5	13.90	1227650	20.0