

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG071915\
 Data File : BG017971.D
 Acq On : 19 Jul 2015 23:30
 Operator : TP/UM
 Sample : G2966-08
 Misc :
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TEPMW-PCB2

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_G\METHODS\8270-BG071715.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	2.788	12	17	25	rBV	58909	94864	1.91%	0.376%
2	2.864	25	30	39	rVB	1009029	1340446	27.00%	5.306%
3	4.733	342	348	354	rBV2	55899	82297	1.66%	0.326%
4	5.191	420	426	434	rBV	483105	682461	13.75%	2.701%
5	6.766	687	694	703	rBV	311850	488037	9.83%	1.932%
6	7.118	747	754	764	rBV	900698	1418988	28.59%	5.617%
7	7.582	827	833	840	rBV	337235	521633	10.51%	2.065%
8	7.900	880	887	893	rBV	814133	1282895	25.84%	5.078%
9	8.734	1022	1029	1040	rVB	657657	1120111	22.57%	4.434%
10	10.362	1298	1306	1313	rBV	484381	815833	16.44%	3.229%
11	11.930	1567	1573	1578	rBV2	36965	66540	1.34%	0.263%
12	12.259	1620	1629	1635	rVB3	50498	114250	2.30%	0.452%
13	12.582	1677	1684	1690	rBV6	32852	74278	1.50%	0.294%
14	12.859	1721	1731	1738	rBV	1896699	2893601	58.29%	11.454%
15	13.288	1800	1804	1811	rVB3	29916	59772	1.20%	0.237%
16	13.564	1845	1851	1856	rBV10	21641	51644	1.04%	0.204%
17	13.705	1870	1875	1880	rBV	114473	167106	3.37%	0.661%
18	13.957	1914	1918	1923	rVB4	44065	67900	1.37%	0.269%
19	14.239	1961	1966	1974	rBV2	743085	1177401	23.72%	4.661%
20	15.614	2196	2200	2208	rVB4	76376	142333	2.87%	0.563%
21	15.738	2215	2221	2233	rBV	1555276	2202360	44.37%	8.718%
22	16.037	2268	2272	2277	rVB	82305	113831	2.29%	0.451%
23	16.995	2429	2435	2443	rVB2	1013965	1399399	28.19%	5.539%
24	17.912	2587	2591	2600	rBV2	171987	280671	5.65%	1.111%
25	19.204	2807	2811	2820	rBV2	109155	162789	3.28%	0.644%
26	19.645	2880	2886	2895	rBV	3898164	4963930	100.00%	19.649%
27	20.526	3033	3036	3041	rVB	59833	69807	1.41%	0.276%
28	20.920	3100	3103	3111	rBV	127074	171627	3.46%	0.679%
29	21.196	3145	3150	3157	rVB	1261063	1617083	32.58%	6.401%
30	23.411	3520	3527	3536	rBV2	845891	1618582	32.61%	6.407%

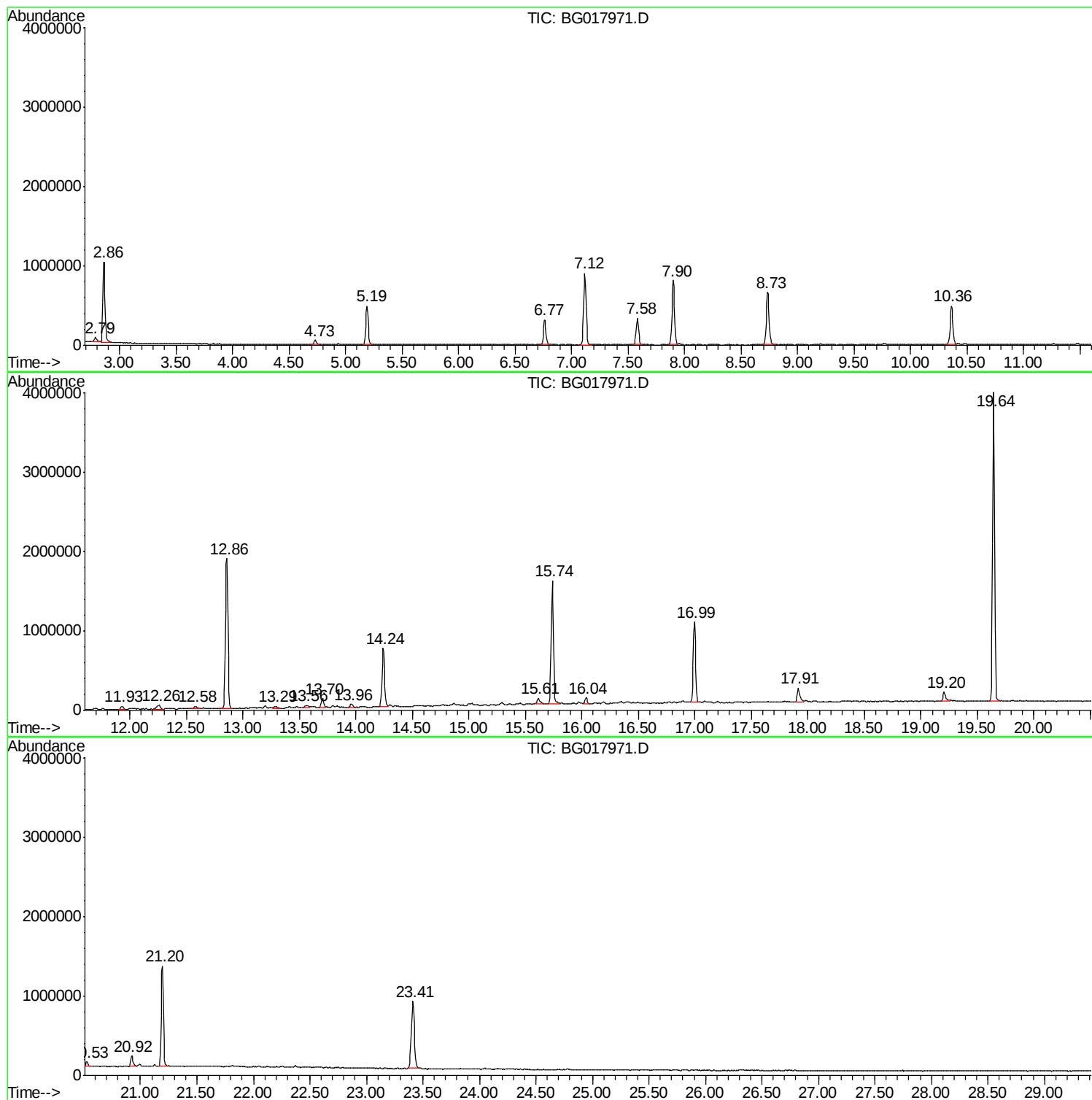
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Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG071715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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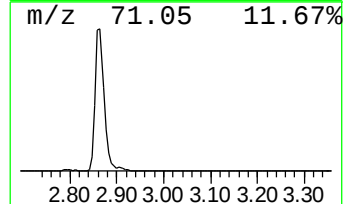
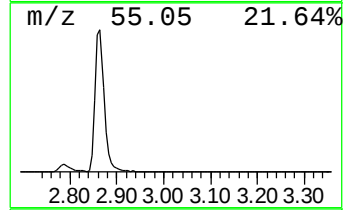
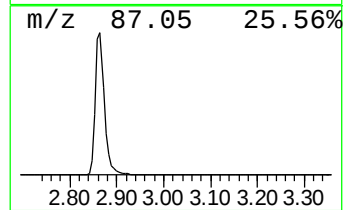
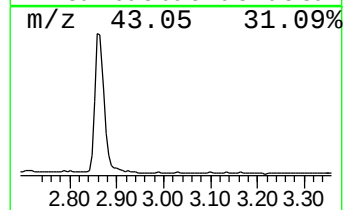
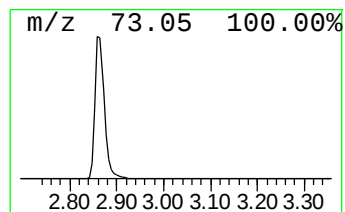
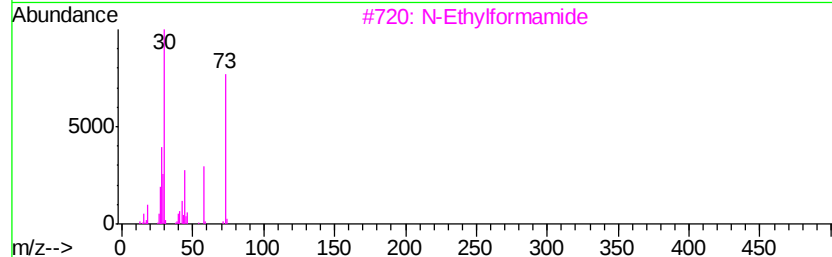
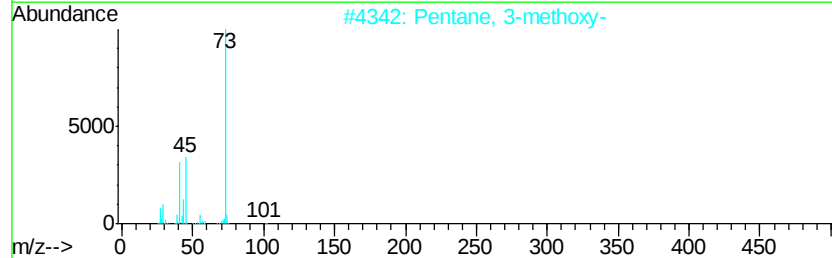
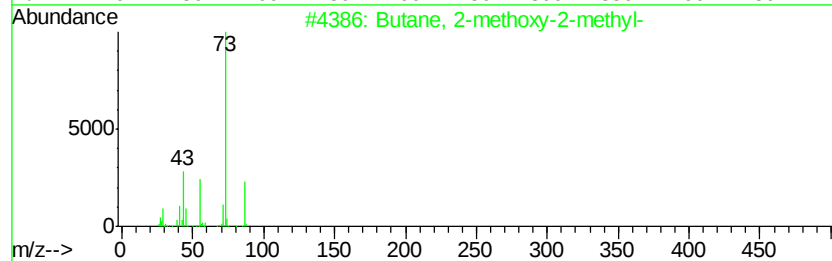
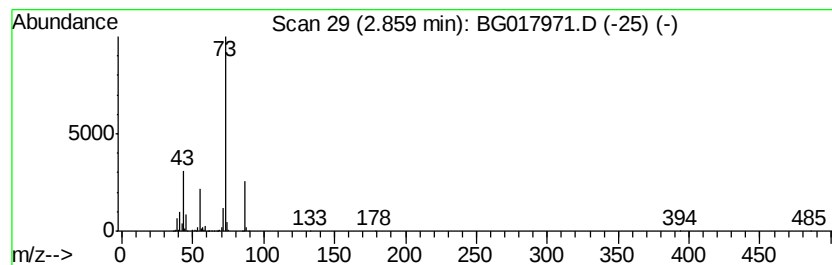
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.86	51.39 ng	1340450	1,4-Dichlorobenzene-d4	7.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	86
2		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	25
3		N-Ethylformamide	73	C3H7NO	000627-45-2	9
4		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9
5		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9



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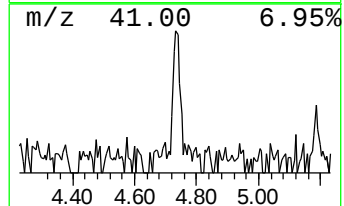
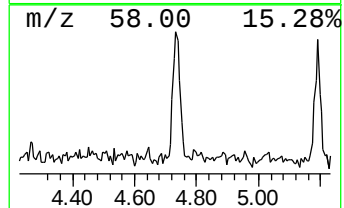
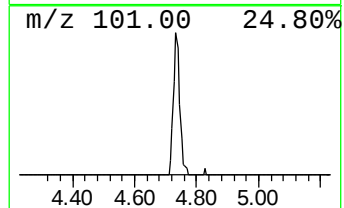
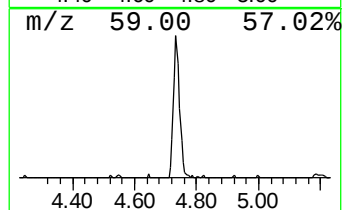
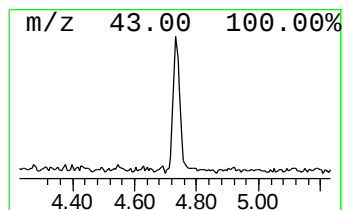
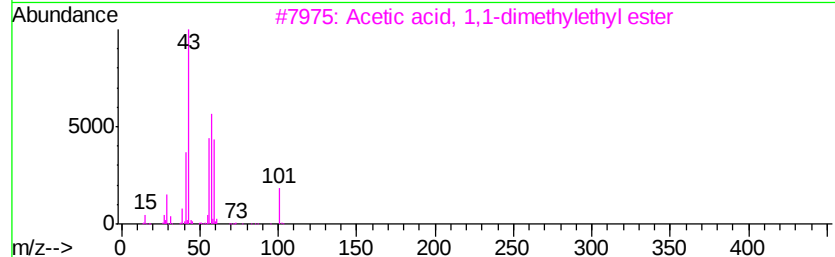
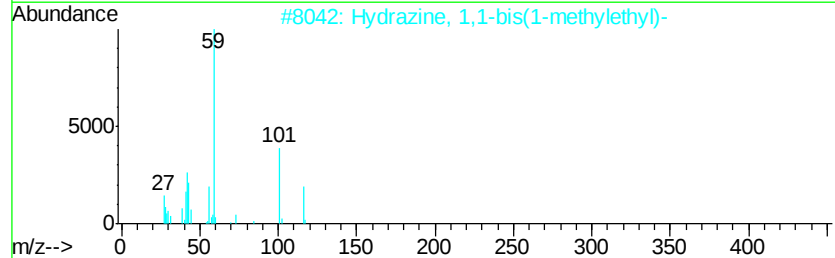
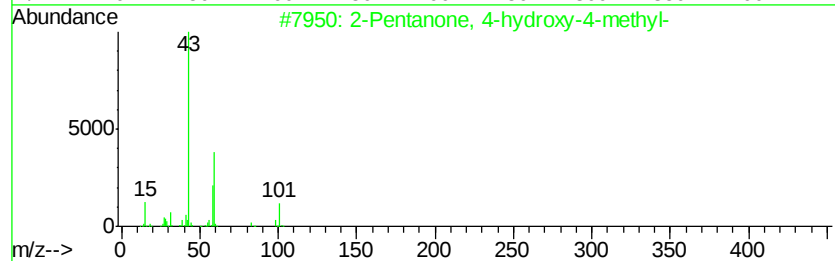
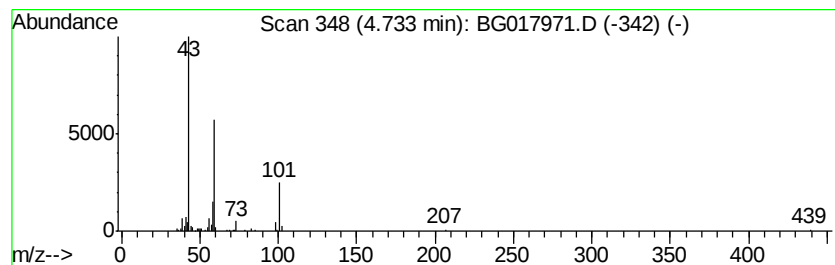
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Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.73	3.16 ng	82297	1,4-Dichlorobenzene-d4	7.58

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2			Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	38
3			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	35
5			4-Ethyl-3-octanol	158	C10H22O	019781-26-1	32



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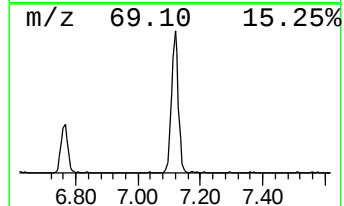
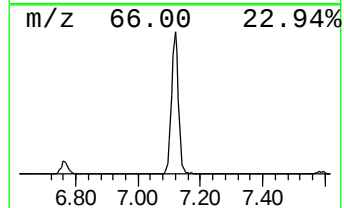
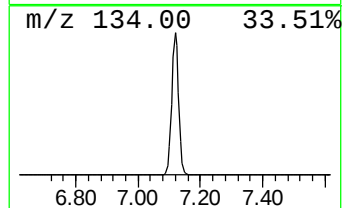
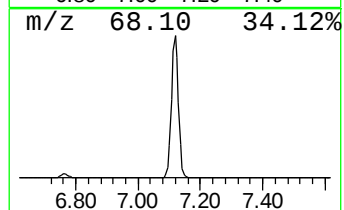
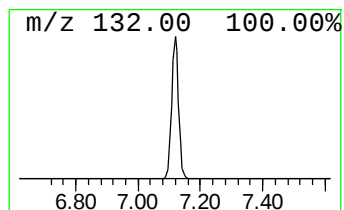
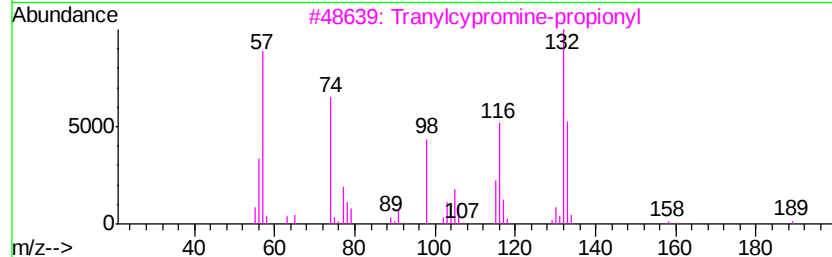
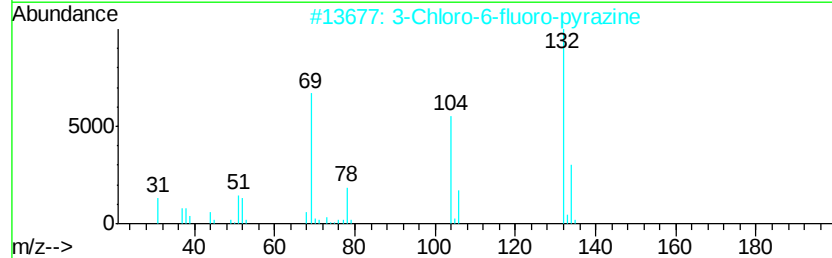
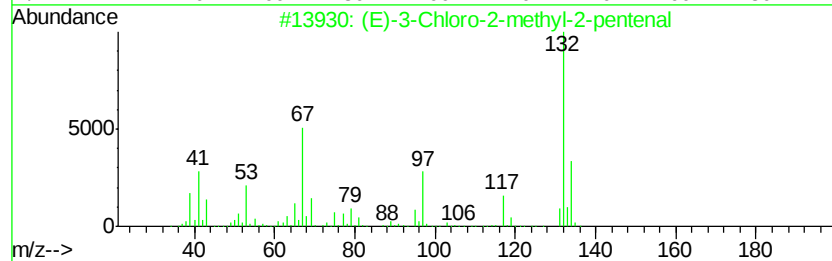
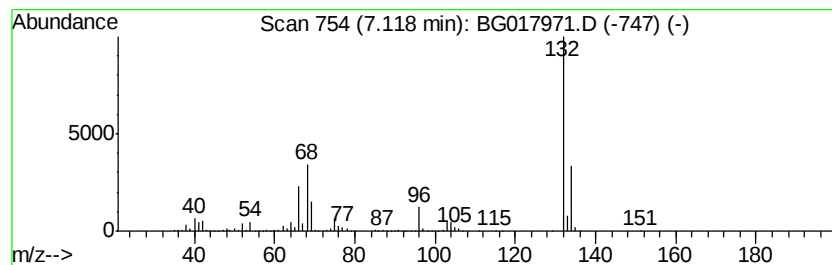
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Peak Number 4 unknown7.12 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.12	54.41 ng	1418990	1,4-Dichlorobenzene-d4	7.58

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	35	
2	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25	
3	Tranlylcypromine-propionyl	189	C12H15NO	1000123-86-3	12	
4	2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9	
5	4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9	



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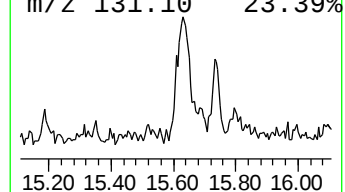
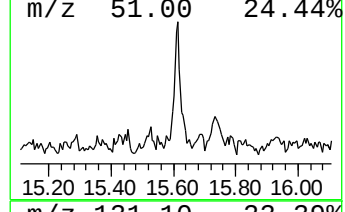
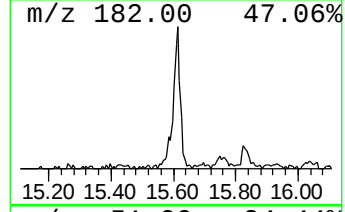
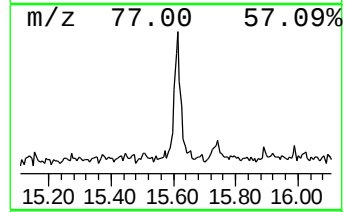
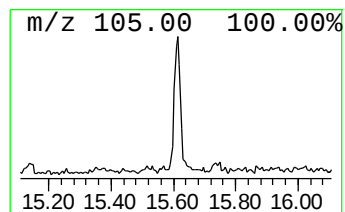
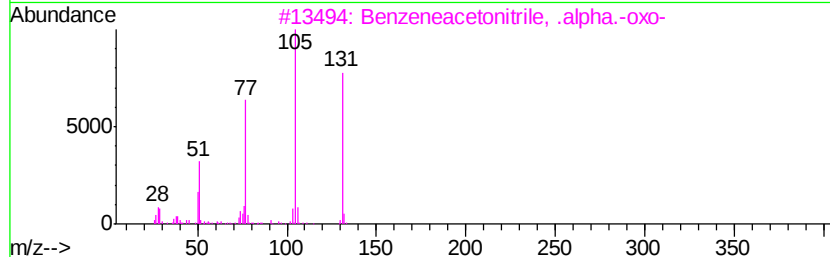
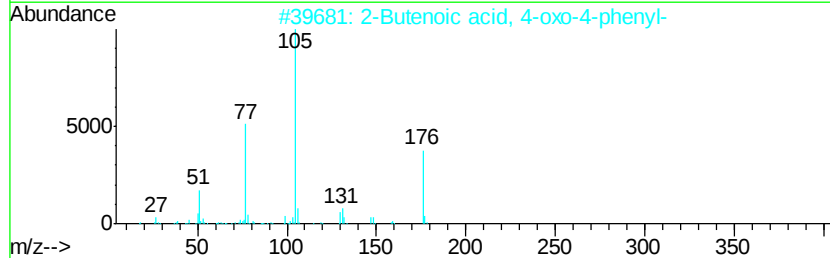
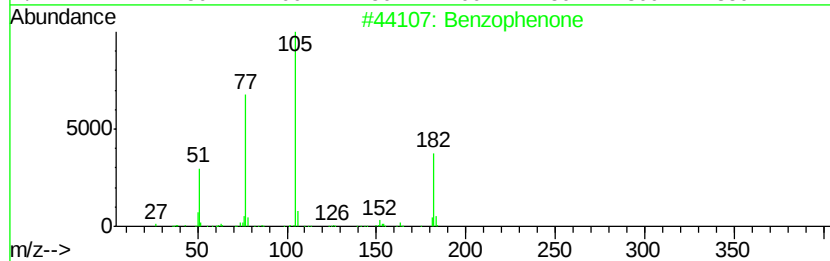
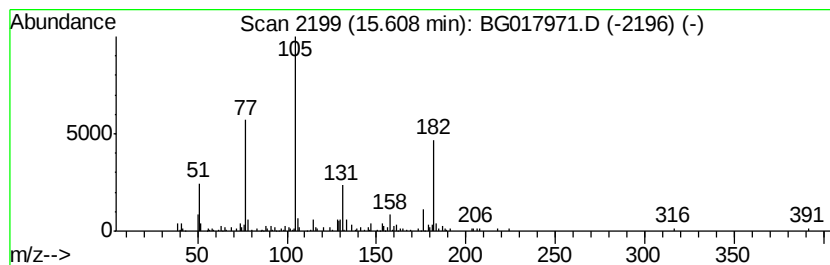
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 7 Benzophenone Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.61	2.42 ng	142333	Acenaphthene-d10	14.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzophenone	182	C13H10O	000119-61-9	58
2		2-Butenoic acid, 4-oxo-4-phenyl-	176	C10H8O3	000583-06-2	43
3		Benzeneacetonitrile, .alpha.-oxo-	131	C8H5NO	000613-90-1	43
4		Butanenitrile, 2,3-bis(benzoylox...	335	C18H13N3O4	1000269-66-6	38
5		7-Benzoyloxybicyclo[2.2.1]hepta...	212	C14H12O2	004796-68-3	38



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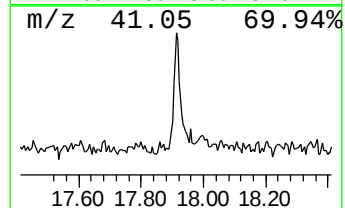
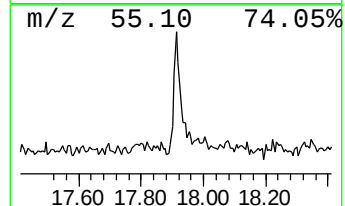
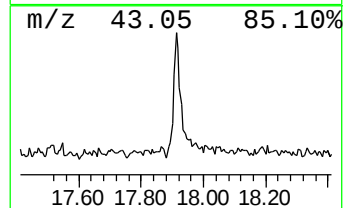
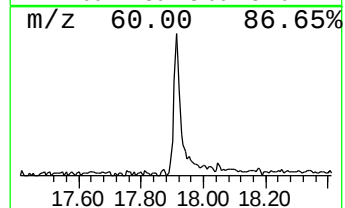
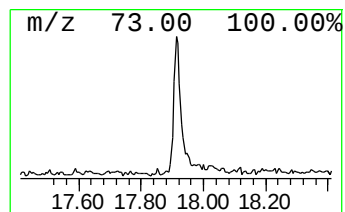
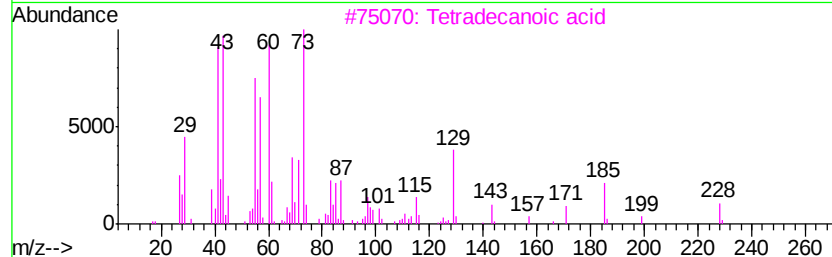
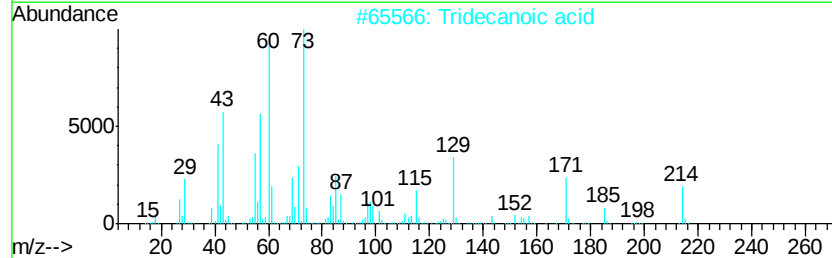
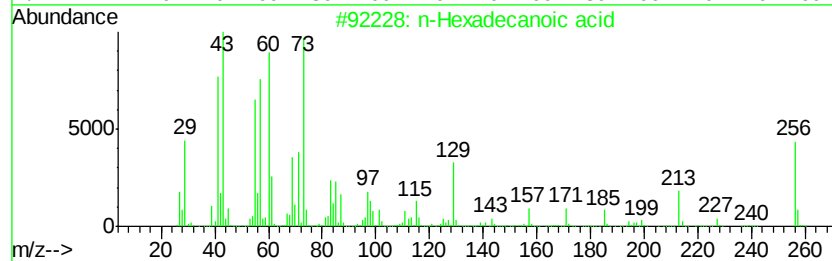
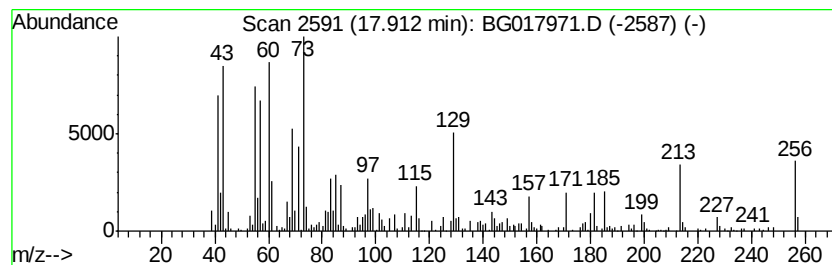
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 8 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.91	4.01 ng	280671	Phenanthrene-d10	16.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
2		Tridecanoic acid	214	C13H26O2	000638-53-9	64
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	62
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	50
5		Dodecanoic acid	200	C12H24O2	000143-07-7	49



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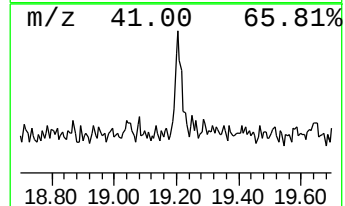
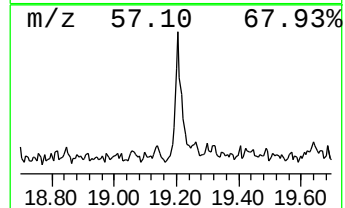
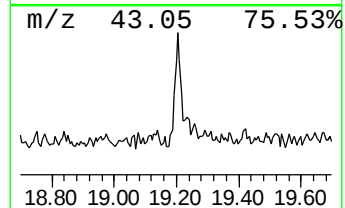
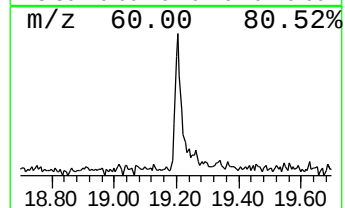
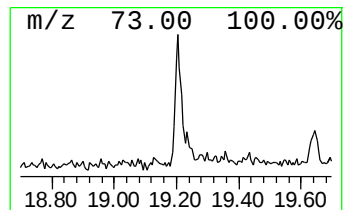
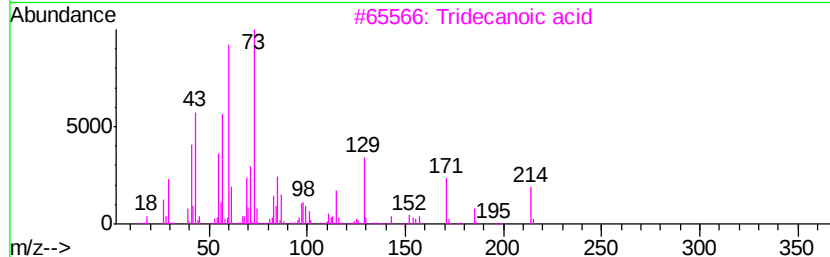
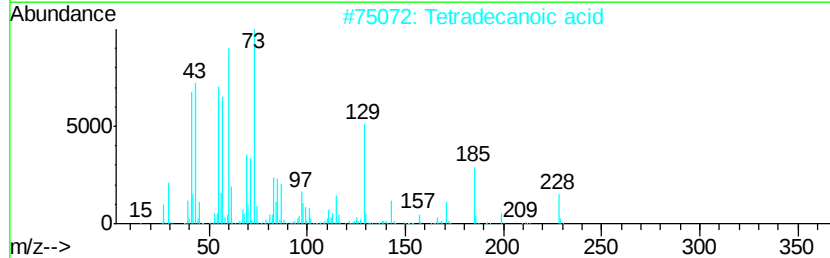
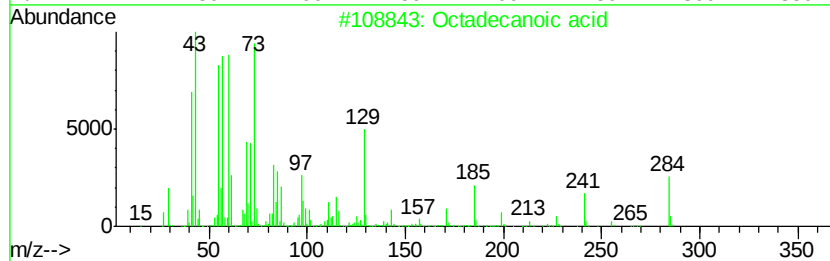
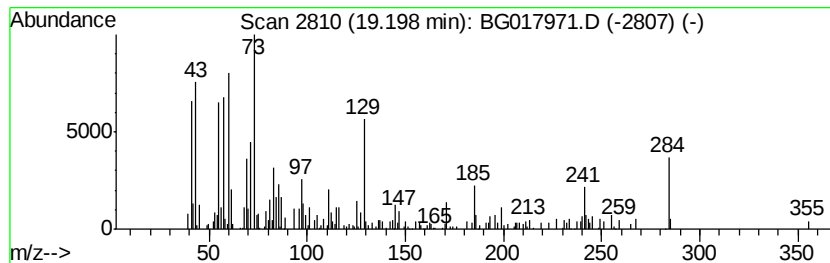
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Octadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.20	2.01 ng	162789	Chrysene-d12	21.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	98
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	62
3		Tridecanoic acid	214	C13H26O2	000638-53-9	58
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	55
5		Undecanoic acid	186	C11H22O2	000112-37-8	49



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG071915\
Data File : BG017971.D
Acq On : 19 Jul 2015 23:30
Operator : TP/UM
Sample : G2966-08
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
TEPMW-PCB2

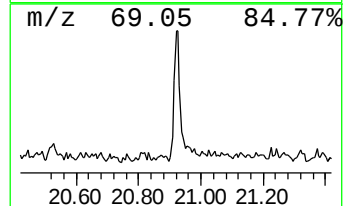
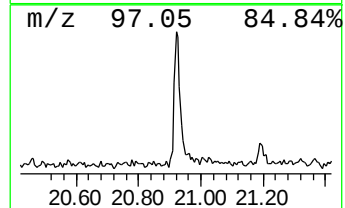
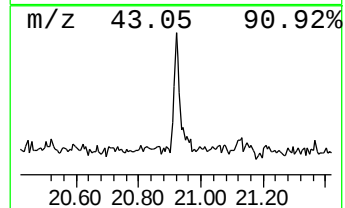
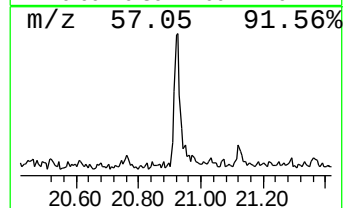
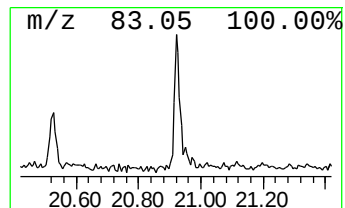
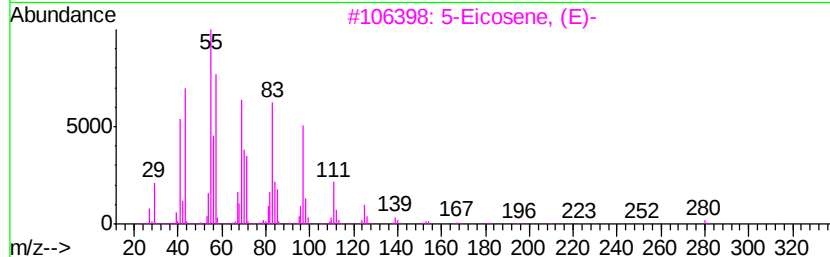
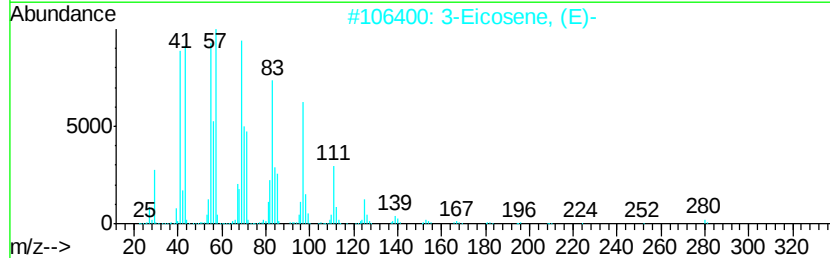
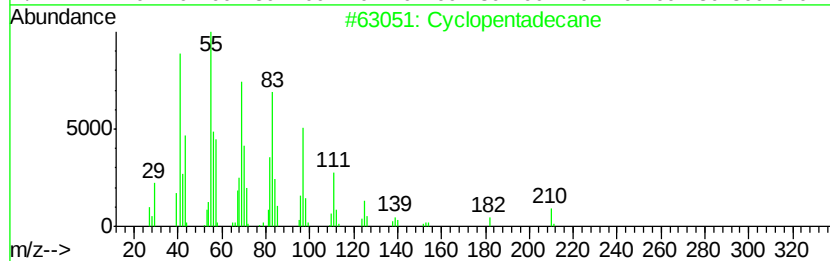
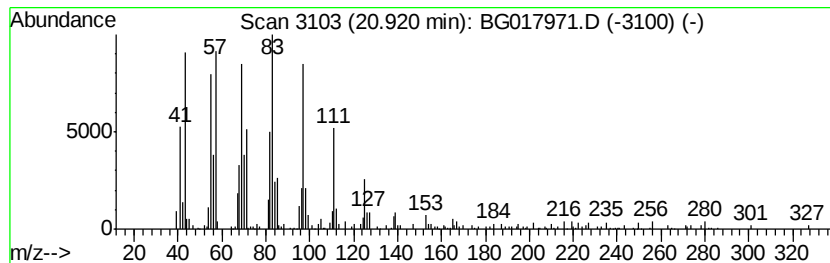
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Cyclopentadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.92	2.12 ng	171627	Chrysene-d12	21.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentadecane	210	C15H30	000295-48-7	93
2		3-Eicosene, (E)-	280	C20H40	074685-33-9	92
3		5-Eicosene, (E)-	280	C20H40	074685-30-6	91
4		Bromoacetic acid, octadecyl ester	390	C20H39BrO2	018992-03-5	87
5		1-Eicosene	280	C20H40	003452-07-1	86



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG071915\
Data File : BG017971.D
Acq On : 19 Jul 2015 23:30
Operator : TP/UM
Sample : G2966-08
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
TEPMW-PCB2

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG071715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	2.86	51.4	ng	1340450	1	7.58	521633	20.0
2-Pentanone, 4-hy...	4.73	3.2	ng	82297	1	7.58	521633	20.0
unknown7.12	7.12	54.4	ng	1418990	1	7.58	521633	20.0
Benzophenone	15.61	2.4	ng	142333	3	14.24	1177400	20.0
n-Hexadecanoic acid	17.91	4.0	ng	280671	4	16.99	1399400	20.0
Octadecanoic acid	19.20	2.0	ng	162789	5	21.20	1617080	20.0
Cyclopentadecane	20.92	2.1	ng	171627	5	21.20	1617080	20.0