

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092216\
 Data File : BG024073.D
 Acq On : 22 Sep 2016 17:47
 Operator : UM/SJ
 Sample : H4820-03
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 YORK-TP1-03

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-BG083116.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	2.916	8	13	28	rVB	3836196	5762755	100.00%	17.823%
2	5.113	381	387	397	rVB	214633	348161	6.04%	1.077%
3	5.595	463	469	485	rBV	1134136	1944382	33.74%	6.013%
4	7.222	739	746	762	rVB	1162395	2213617	38.41%	6.846%
5	7.580	799	807	818	rBV	1199144	2145881	37.24%	6.637%
6	8.044	880	886	895	rVB	295202	499071	8.66%	1.543%
7	8.373	935	942	949	rVB	870921	1529358	26.54%	4.730%
8	9.225	1078	1087	1098	rBV	750312	1459386	25.32%	4.513%
9	10.876	1361	1368	1375	rBV	377684	704939	12.23%	2.180%
10	12.544	1645	1652	1658	rBV2	41510	70082	1.22%	0.217%
11	13.331	1777	1786	1794	rBV	1891718	2967060	51.49%	9.176%
12	14.036	1900	1906	1911	rBV5	28938	59161	1.03%	0.183%
13	14.165	1923	1928	1938	rVB	392432	578418	10.04%	1.789%
14	14.723	2008	2023	2030	rBV	584288	989989	17.18%	3.062%
15	16.227	2272	2279	2289	rBV	1460375	2236115	38.80%	6.916%
16	16.433	2306	2314	2319	rBV6	52118	123296	2.14%	0.381%
17	17.208	2442	2446	2449	rBV4	50590	59479	1.03%	0.184%
18	17.490	2487	2494	2498	rBV2	627015	1013231	17.58%	3.134%
19	17.531	2498	2501	2507	rVB	158067	247326	4.29%	0.765%
20	18.189	2610	2613	2619	rVB2	75824	106484	1.85%	0.329%
21	18.366	2638	2643	2648	rVV2	247046	419814	7.28%	1.298%
22	18.418	2649	2652	2657	rVB2	74053	130661	2.27%	0.404%
23	18.565	2672	2677	2680	rBV3	47290	86894	1.51%	0.269%
24	19.282	2794	2799	2802	rBV4	71629	115044	2.00%	0.356%
25	19.546	2840	2844	2849	rVB	302279	422617	7.33%	1.307%
26	19.634	2855	2859	2870	rBV2	300284	517813	8.99%	1.601%
27	19.916	2902	2907	2913	rVB	309681	479703	8.32%	1.484%
28	20.099	2933	2938	2945	rBV	2544859	3450243	59.87%	10.671%
29	20.380	2983	2986	2994	rVB6	110626	176372	3.06%	0.545%
30	21.409	3157	3161	3166	rVB7	91009	158476	2.75%	0.490%
31	21.802	3221	3228	3233	rBV2	665815	1318328	22.88%	4.077%

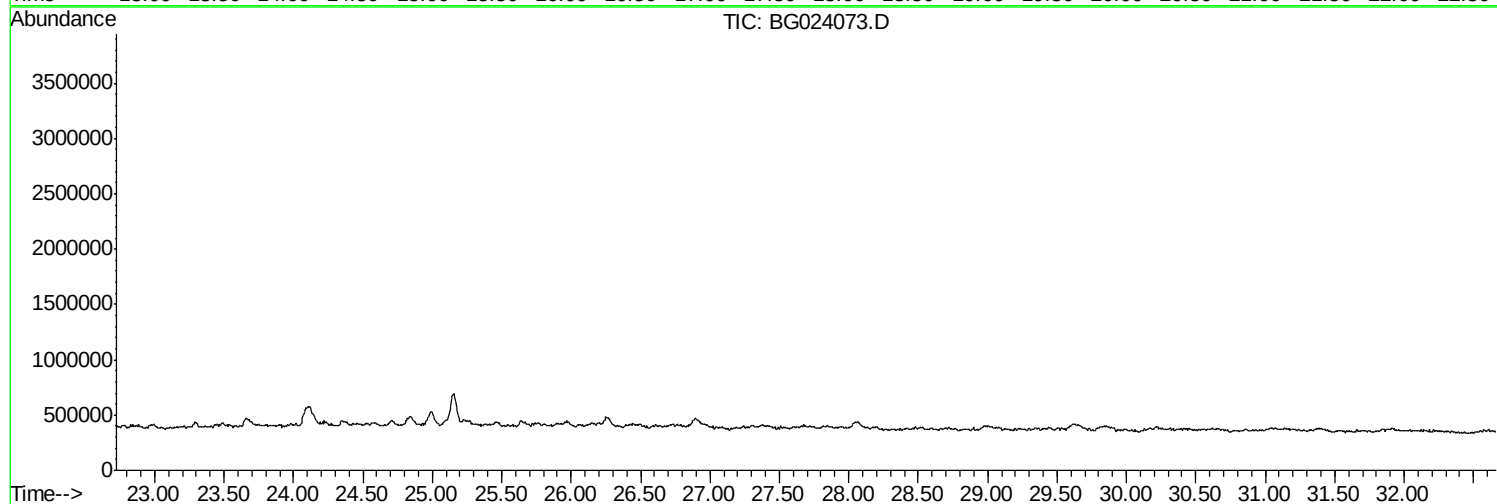
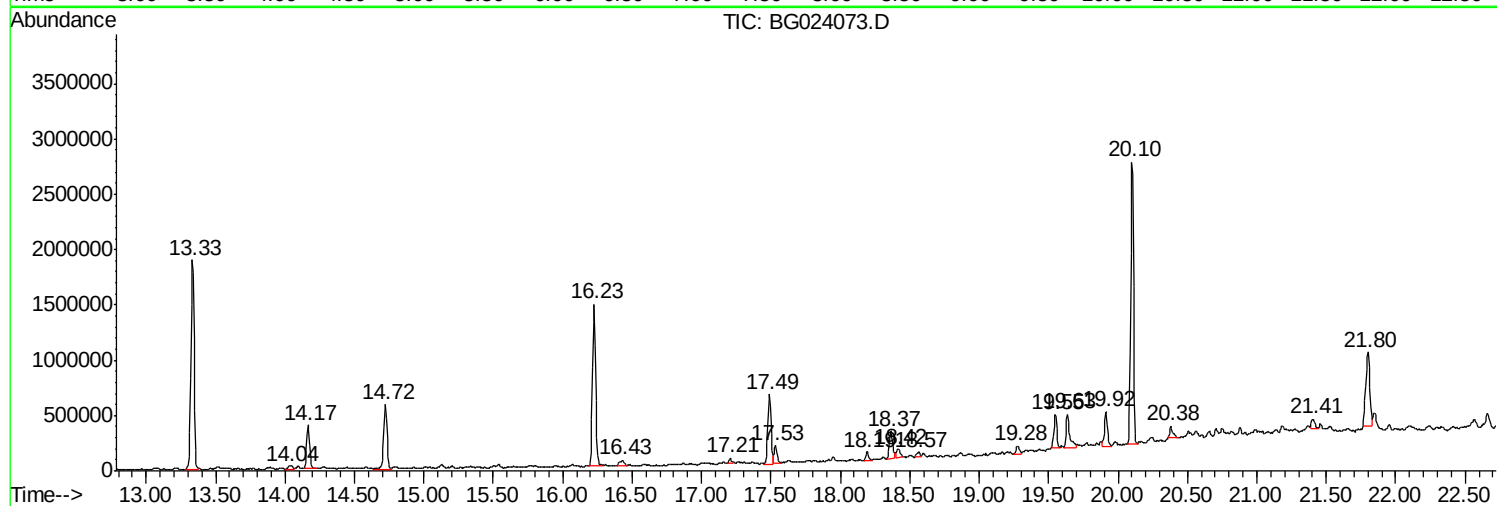
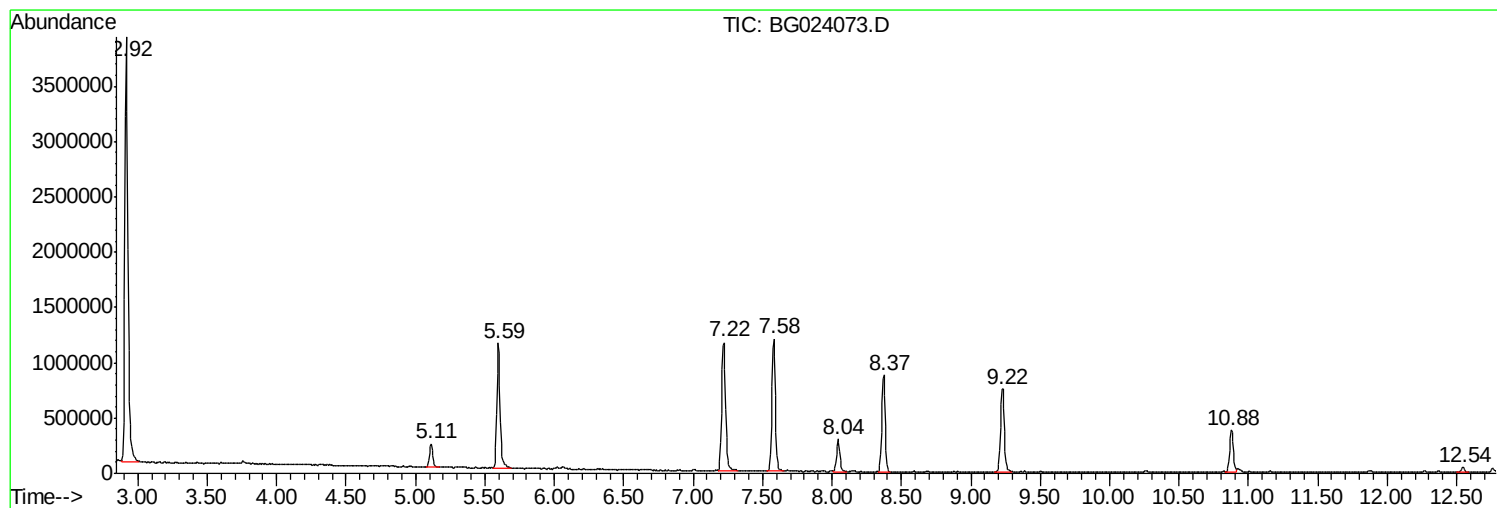
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG083116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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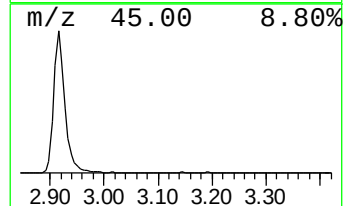
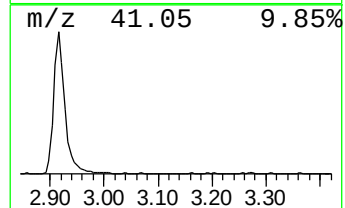
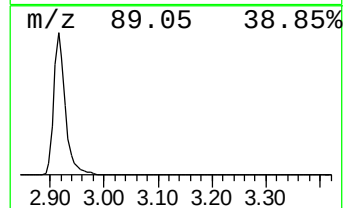
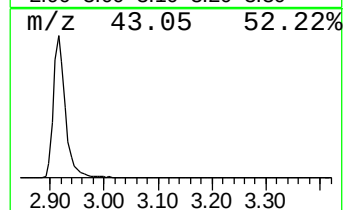
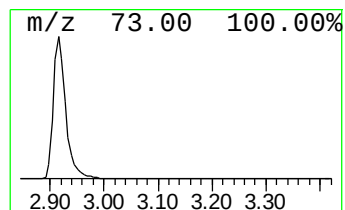
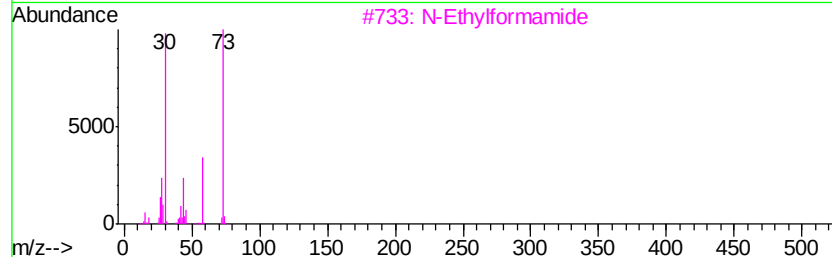
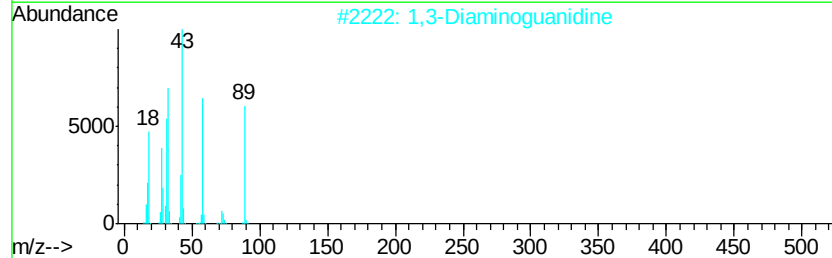
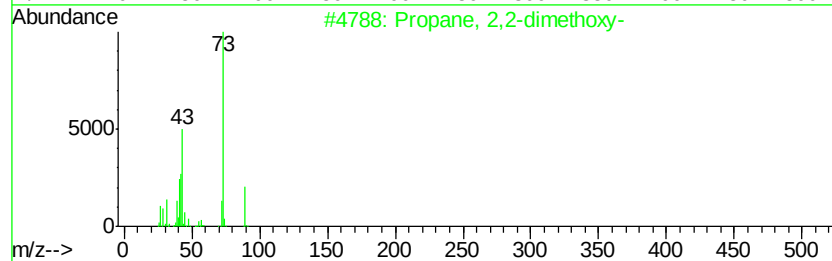
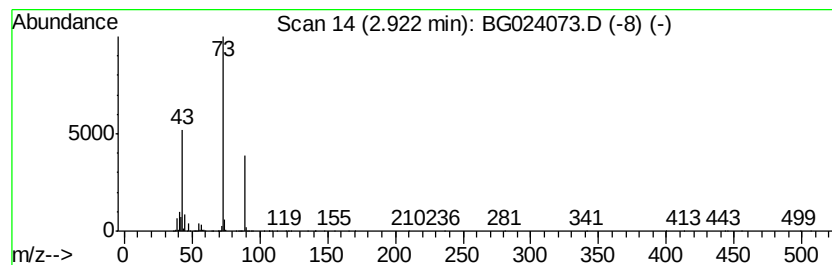
Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG083116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Propane, 2,2-dimethoxy- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
2.92	230.94 ng	5762760	1,4-Dichlorobenzene-d4	8.04	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	72
2	1,3-Diaminoguanidine	89	CH7N5	004364-78-7	25
3	N-Ethylformamide	73	C3H7NO	000627-45-2	9
4	Acetamide, N-methyl-	73	C3H7NO	000079-16-3	9
5	1-Propanol, 2-methyl-	74	C4H10O	000078-83-1	9



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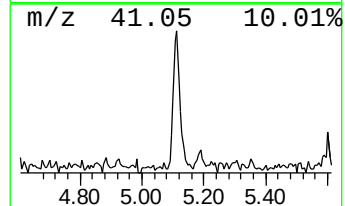
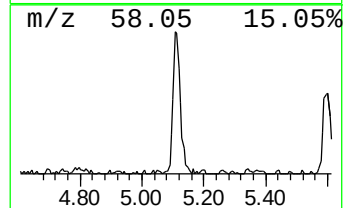
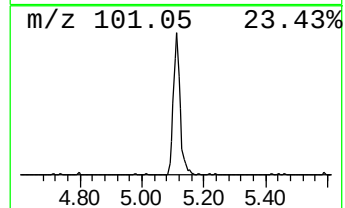
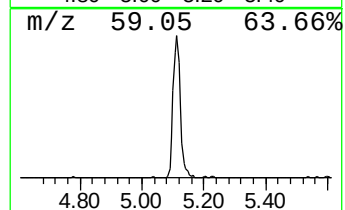
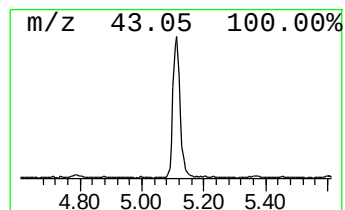
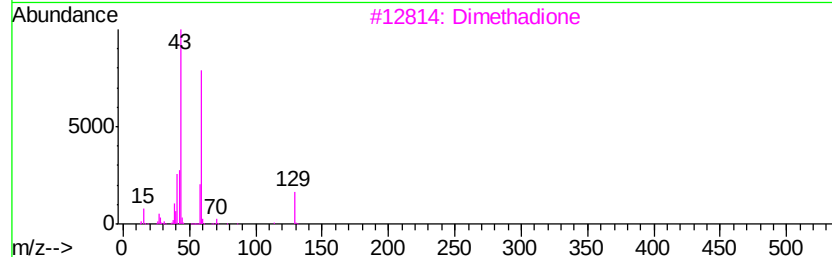
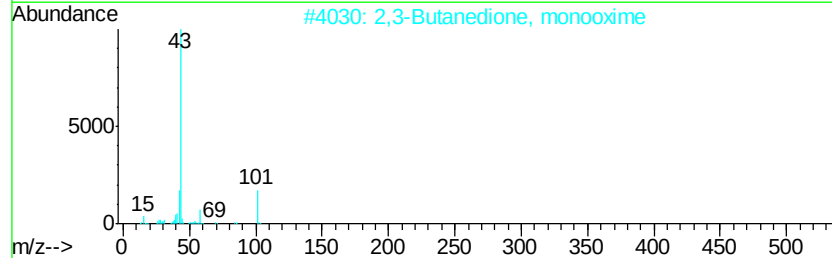
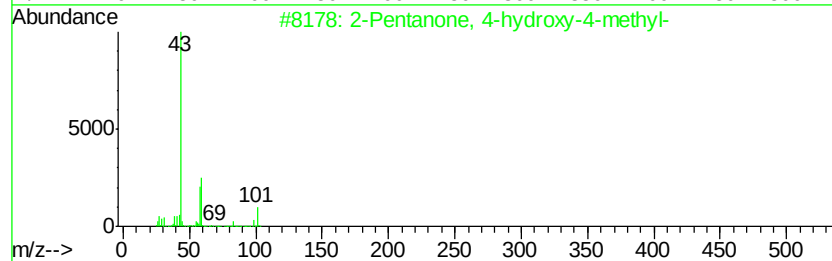
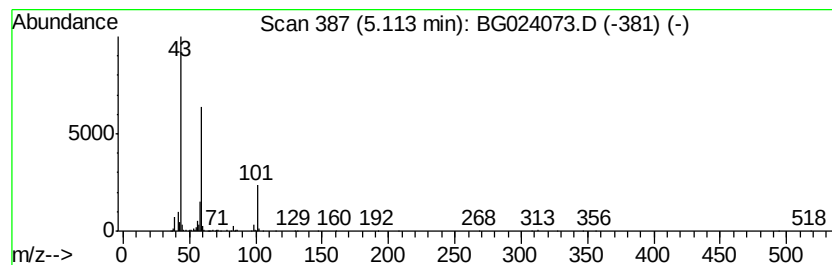
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.11	13.95 ng	348161	1,4-Dichlorobenzene-d4	8.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	35
3		Dimethadione	129	C5H7NO3	000695-53-4	33
4		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	25
5		4-Methoxy-1-pentene	100	C6H12O	098386-09-5	17



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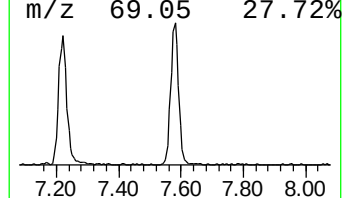
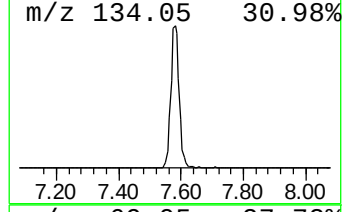
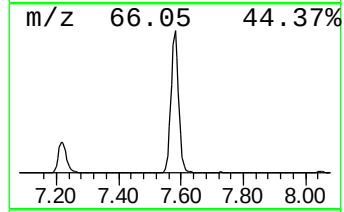
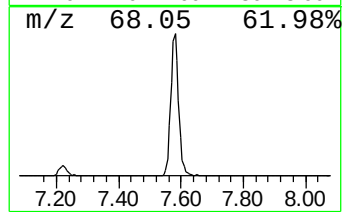
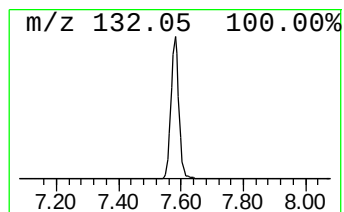
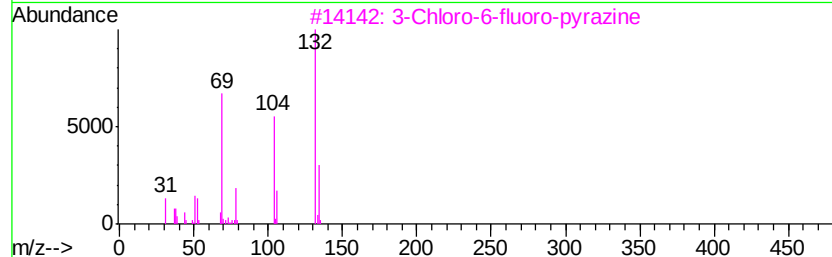
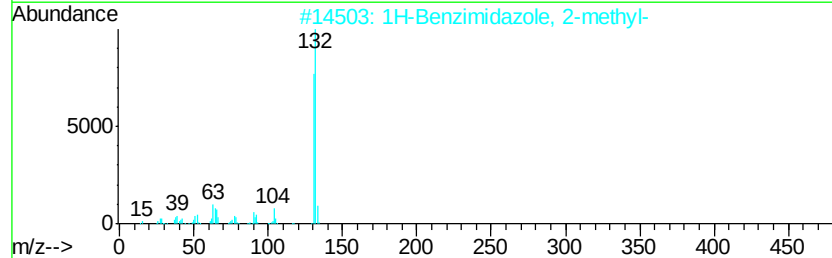
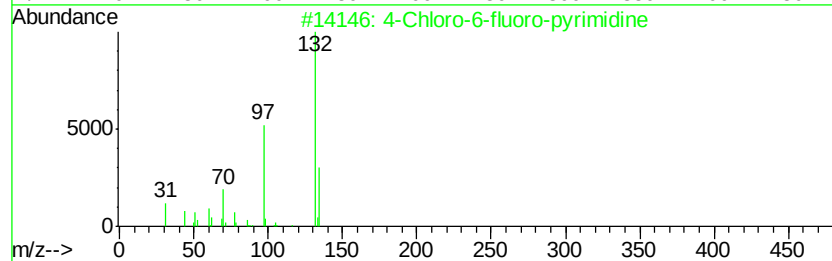
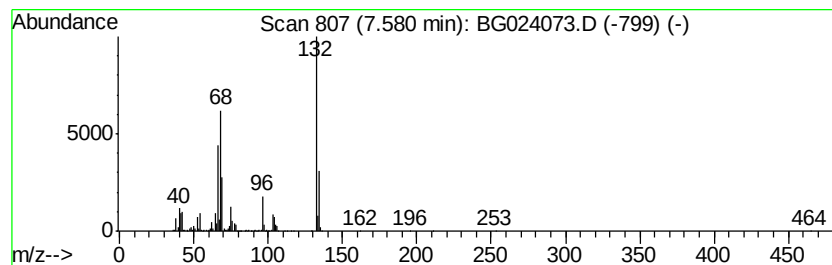
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 unknown7.58 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	86.00 ng	2145880	1,4-Dichlorobenzene-d4	8.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	14
2		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	14
3		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	14
4		3H-1,2-Dithiol-3-one, 5-methyl-	132	C4H4OS2	003620-08-4	10
5		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10



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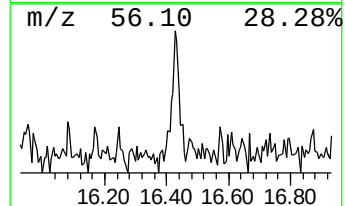
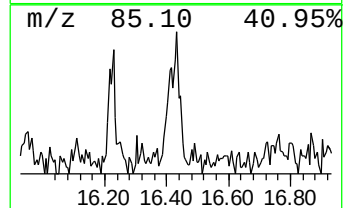
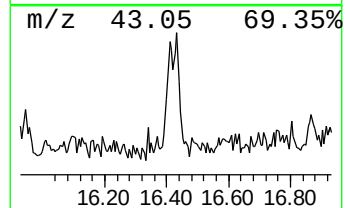
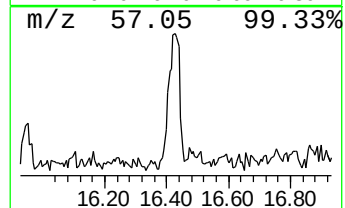
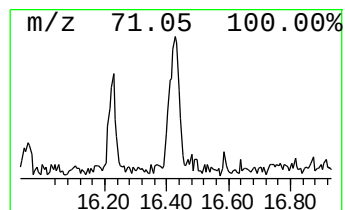
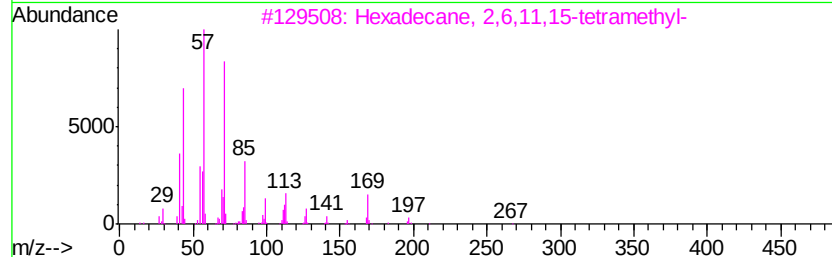
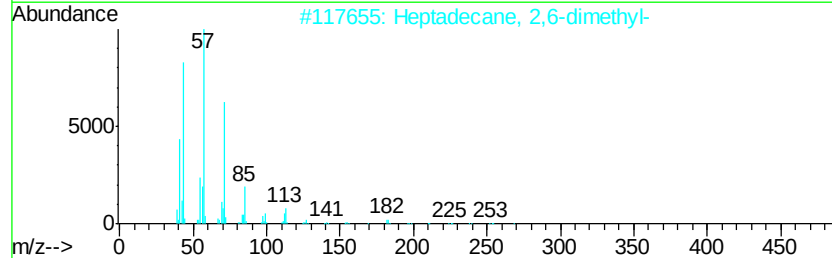
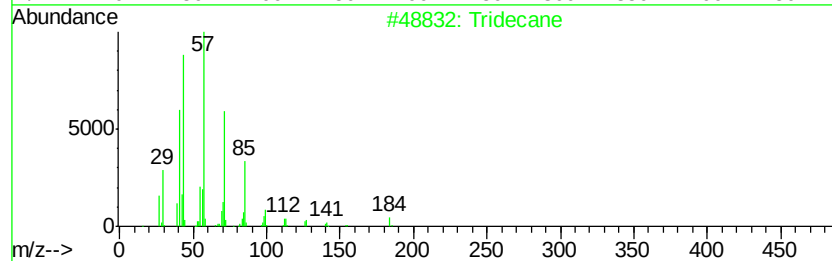
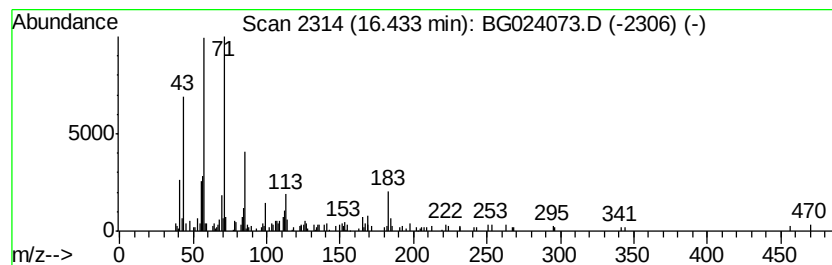
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 Tridecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.43	2.43 ng	123296	Phenanthrene-d10	17.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane	184	C13H28	000629-50-5	62
2		Heptadecane, 2,6-dimethyl-	268	C19H40	054105-67-8	62
3		Hexadecane, 2,6,11,15-tetramethyl-	282	C20H42	000504-44-9	58
4		Hexadecane, 2,6,10-trimethyl-	268	C19H40	055000-52-7	58
5		Heptadecane, 7-methyl-	254	C18H38	020959-33-5	53



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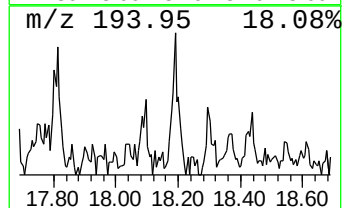
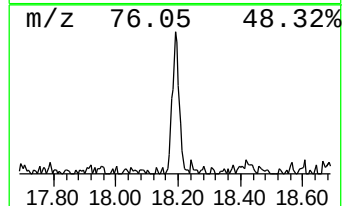
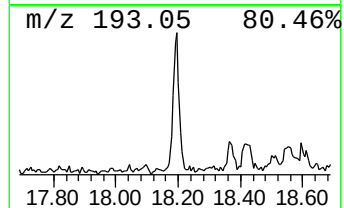
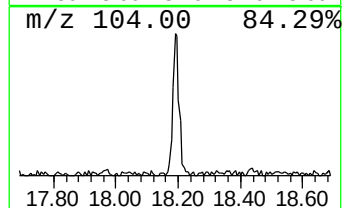
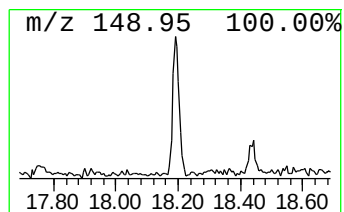
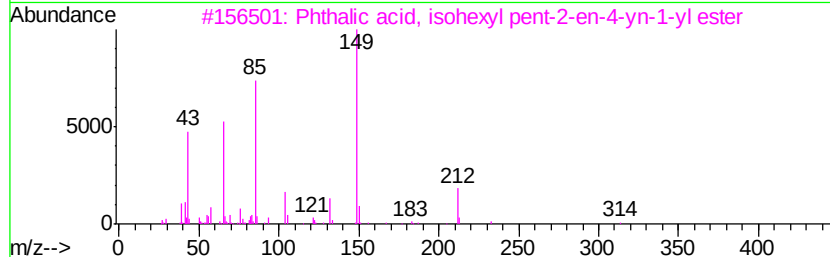
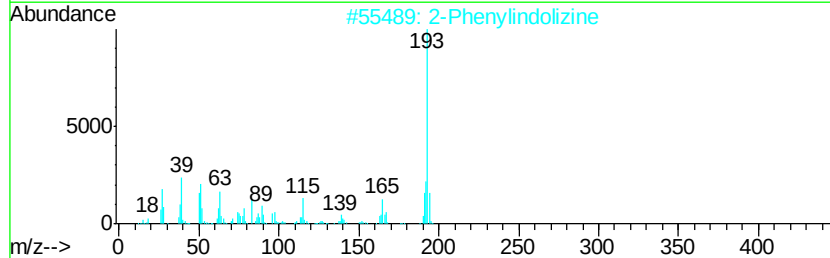
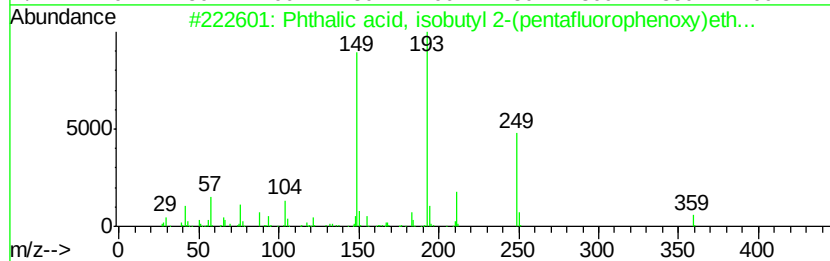
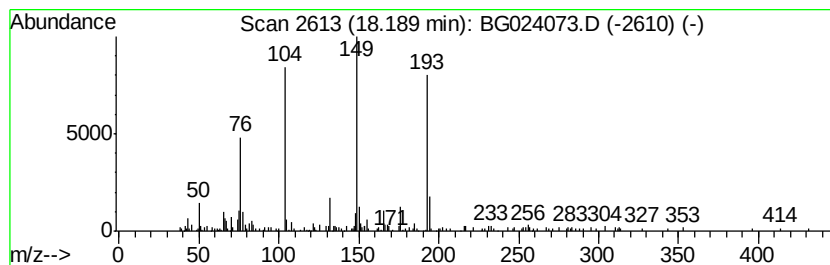
Instrument :
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ClientSampleId :
YORK-TP1-03

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

Peak Number 6 Phthalic acid, isobutyl 2-(... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.		
18.19	2.10 ng	106484	Phenanthrene-d10	17.49		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phthalic acid, isobutyl 2-(penta...	432	C20H17F5O5	1000378-00-4	50
2		2-Phenylindolizine	193	C14H11N	025379-20-8	18
3		Phthalic acid, isohexyl pent-2-e...	314	C19H22O4	1000315-45-7	18
4		4-Methylthieno[2,3-b]pyridine	149	C8H7NS	013362-81-7	14
5		Phthalic acid, hexyl 2-tert-buty...	396	C25H32O4	1000357-09-7	14



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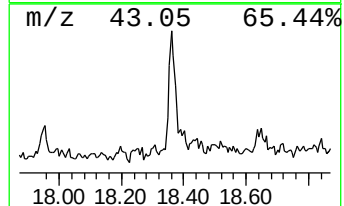
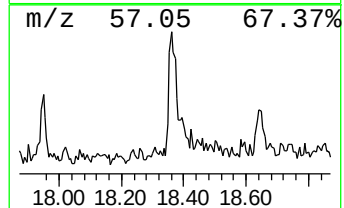
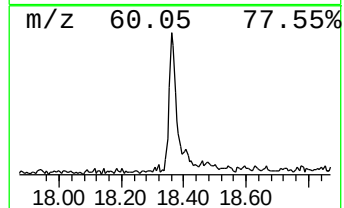
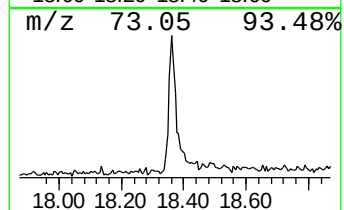
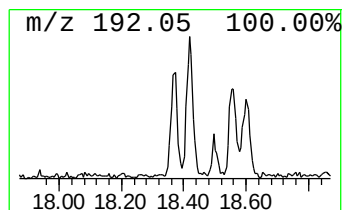
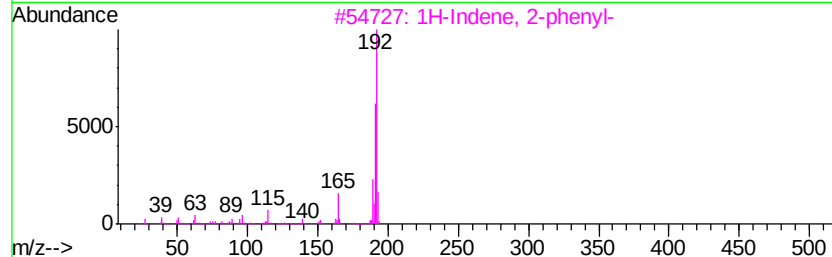
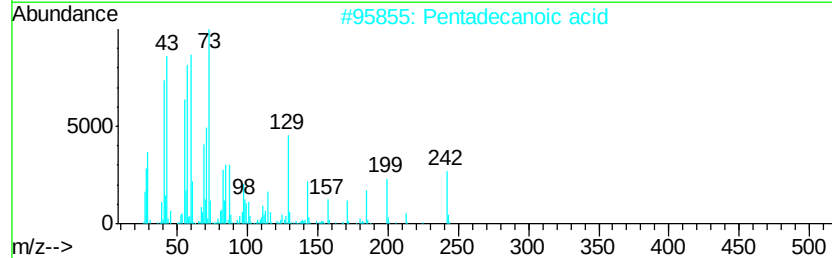
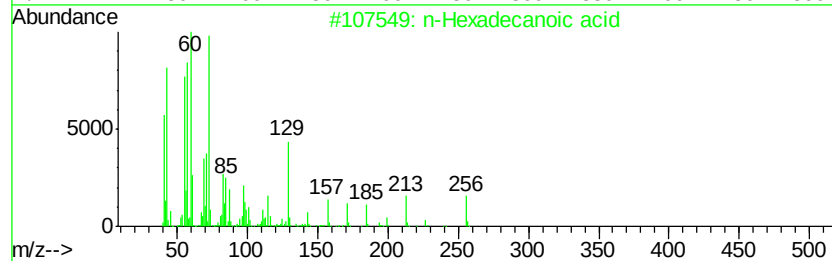
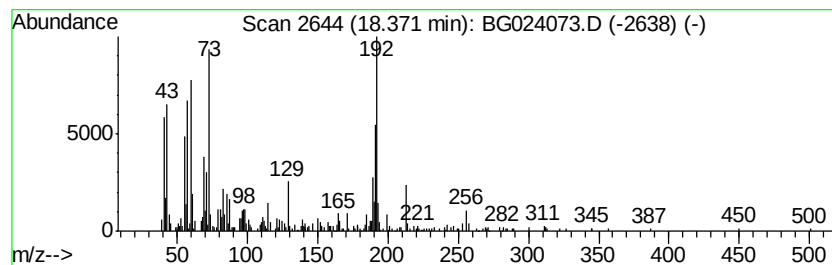
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ClientSampleId :
YORK-TP1-03

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG083116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.
18.37	8.29 ng	419814	Phenanthrene-d10		17.49
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	83
3	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	55
4	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	53
5	Tetradecanoic acid	228	C14H28O2	000544-63-8	50



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092216\
Data File : BG024073.D
Acq On : 22 Sep 2016 17:47
Operator : UM/SJ
Sample : H4820-03
Misc :
ALS Vial : 4 Sample Multiplier: 1

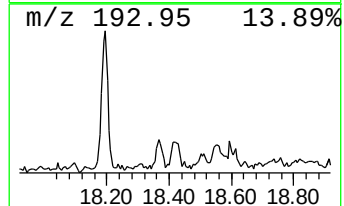
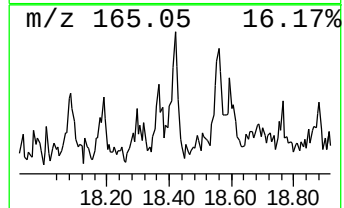
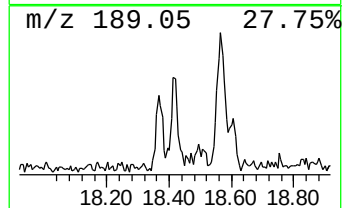
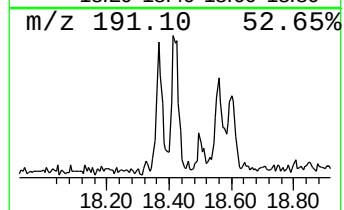
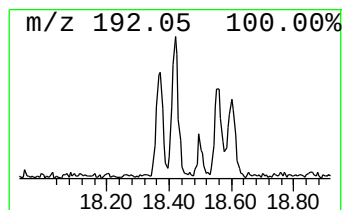
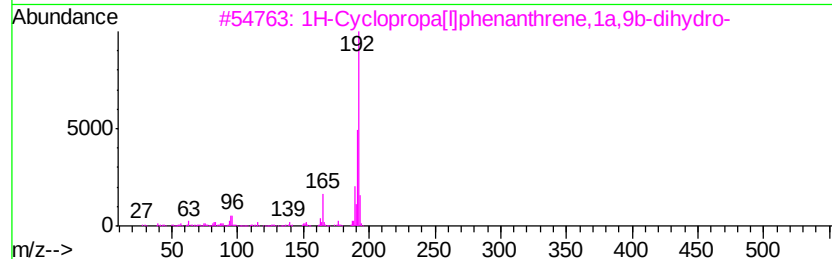
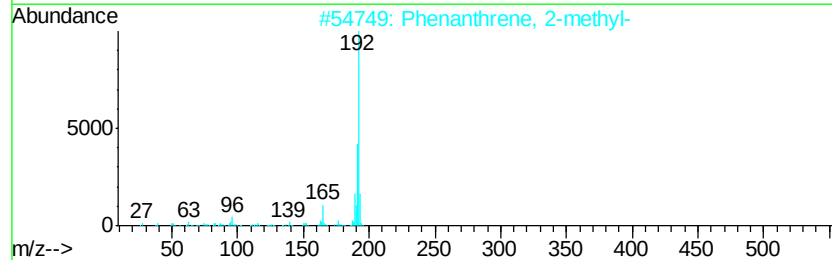
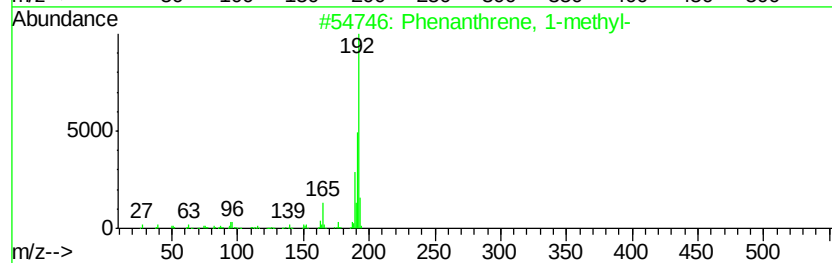
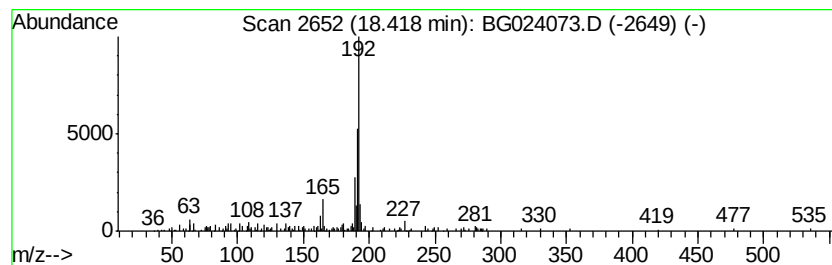
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG083116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Phenanthrene, 1-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.42	2.58 ng	130661	Phenanthrene-d10	17.49	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	81
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	81
3	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	81
4	Anthracene, 2-methyl-	192	C15H12	000613-12-7	81
5	Phenanthrene, 3-methyl-	192	C15H12	000832-71-3	81



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092216\
Data File : BG024073.D
Acq On : 22 Sep 2016 17:47
Operator : UM/SJ
Sample : H4820-03
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YORK-TP1-03

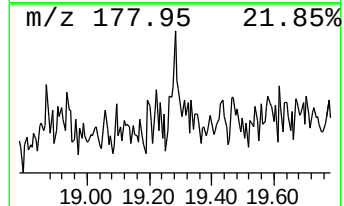
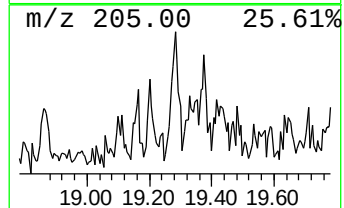
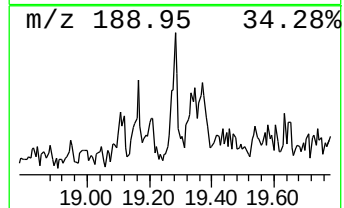
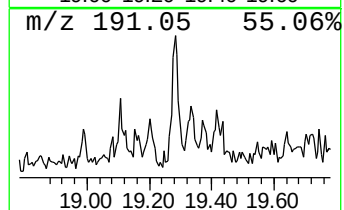
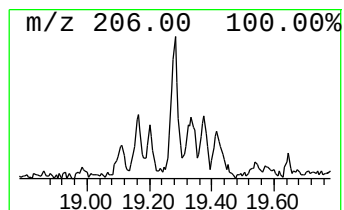
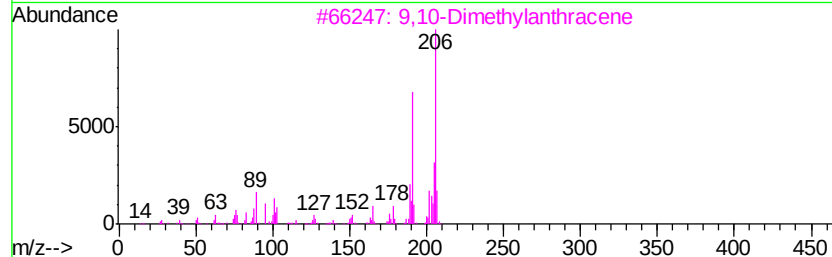
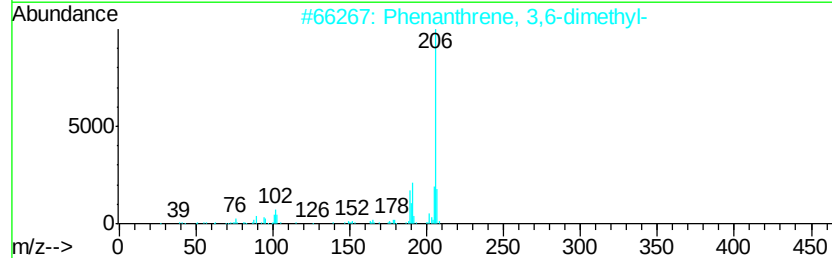
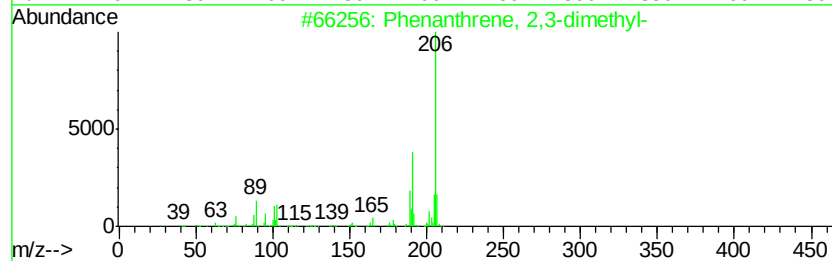
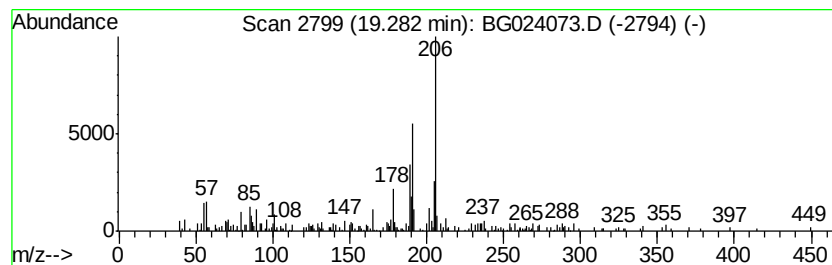
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Phenanthrene, 2,3-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.28	2.27 ng	115044	Phenanthrene-d10	17.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	81
2		Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	76
3		9,10-Dimethylantracene	206	C16H14	000781-43-1	74
4		Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	74
5		[14]Annulene, 1,6:8,13-bis(metha...	206	C16H14	085385-68-8	74



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092216\
Data File : BG024073.D
Acq On : 22 Sep 2016 17:47
Operator : UM/SJ
Sample : H4820-03
Misc :
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Instrument :
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ClientSampleId :
YORK-TP1-03

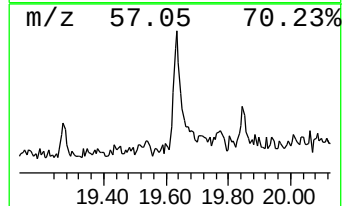
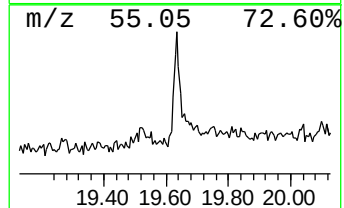
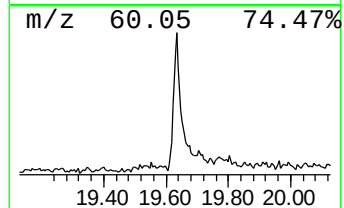
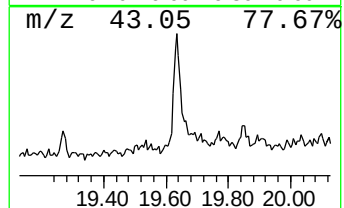
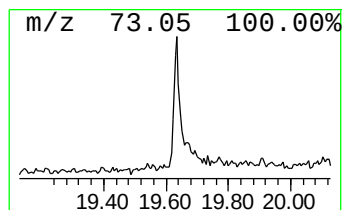
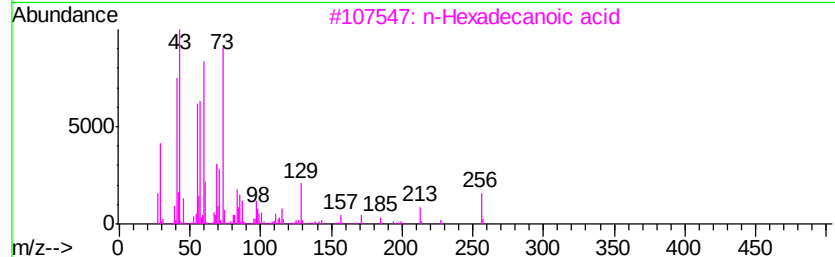
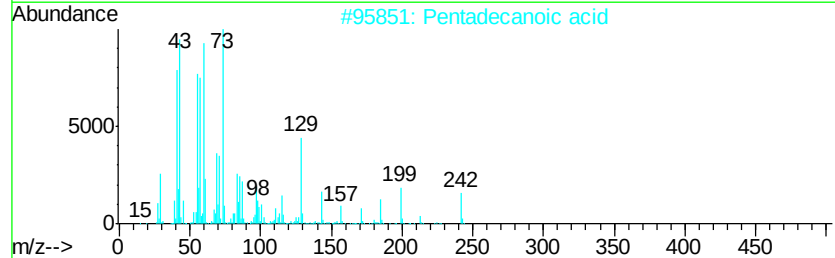
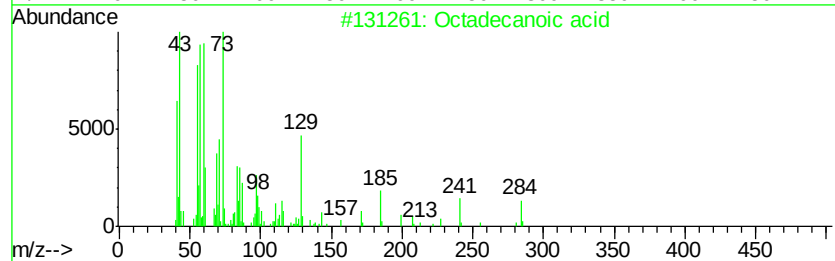
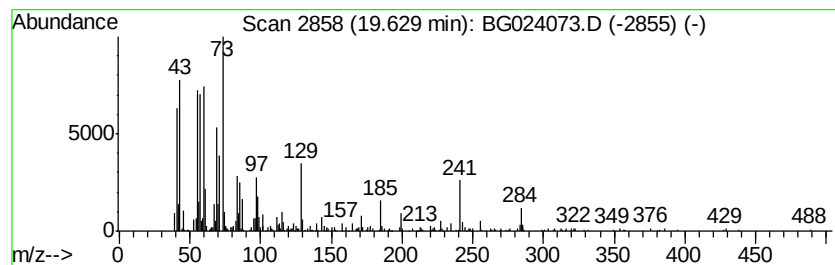
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Octadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.63	10.22 ng	517813	Phenanthrene-d10	17.49

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	94
3		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	70
5		Tridecanoic acid	214	C13H26O2	000638-53-9	47



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092216\
Data File : BG024073.D
Acq On : 22 Sep 2016 17:47
Operator : UM/SJ
Sample : H4820-03
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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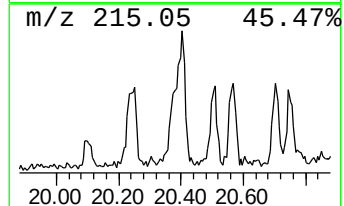
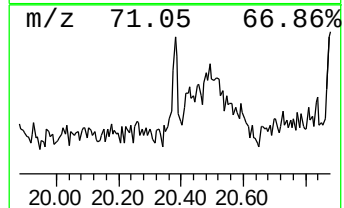
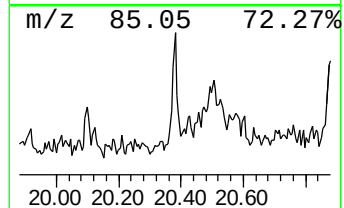
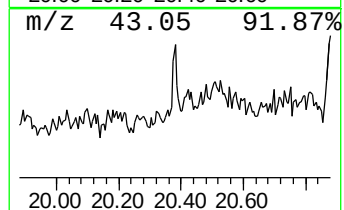
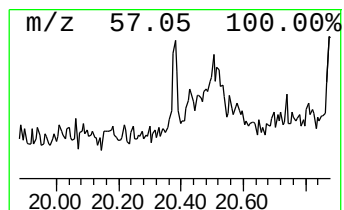
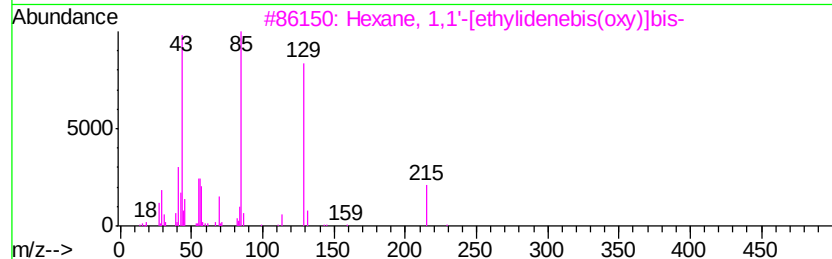
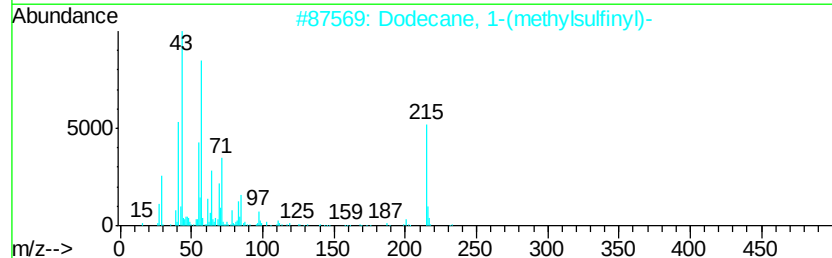
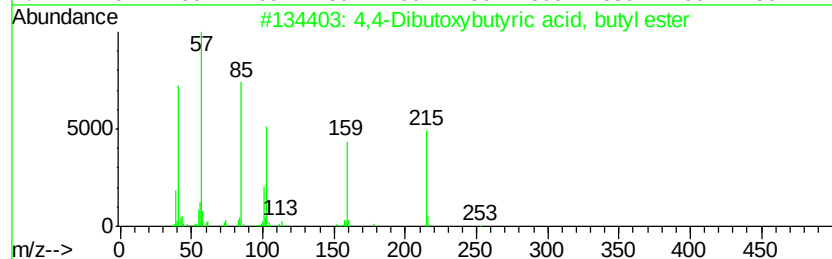
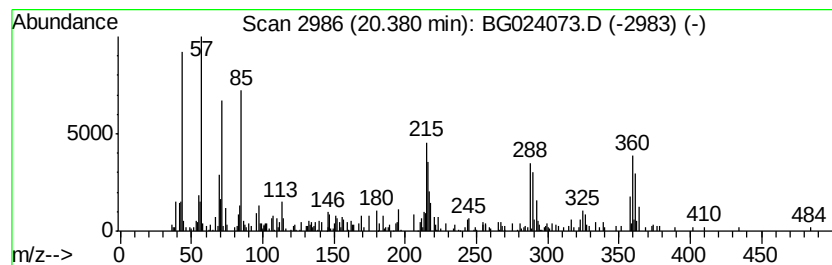
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown20.38 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.38	2.68 ng	176372	Chrysene-d12	21.80

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4,4-Dibutoxybutyric acid, butyl ...	288	C16H32O4	101443-57-6	12
2			Dodecane, 1-(methylsulfinyl)-	232	C13H28OS	003079-30-9	9
3			Hexane, 1,1'-[ethylidenebis(oxy)...	230	C14H30O2	005405-58-3	9
4			.alpha.-D-Xylofuranose, 1,2:3,5-...	230	C11H18O5	020881-04-3	9
5			5-(6,8-Dichloro-1-decylbenz[cd]i...	503	C27H31Cl2NO4	303099-58-3	8



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092216\
Data File : BG024073.D
Acq On : 22 Sep 2016 17:47
Operator : UM/SJ
Sample : H4820-03
Misc :
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ClientSampleId :
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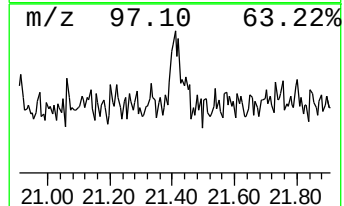
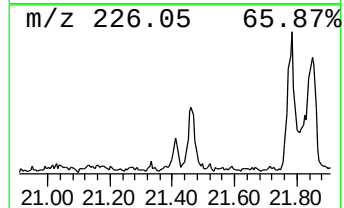
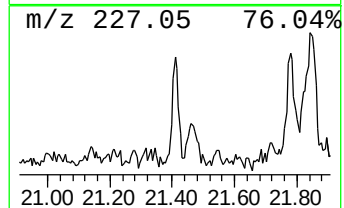
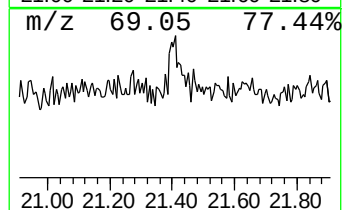
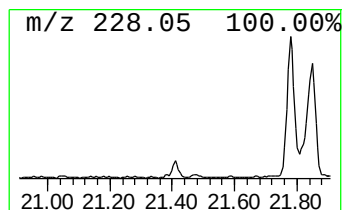
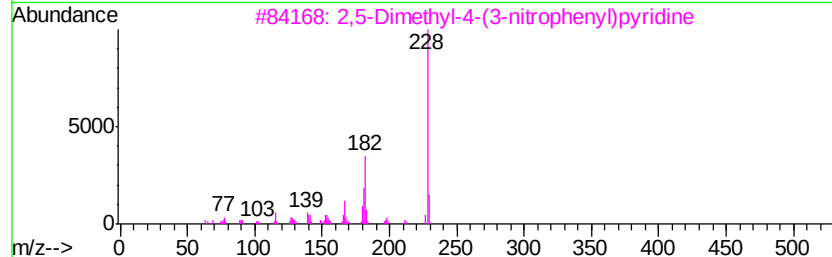
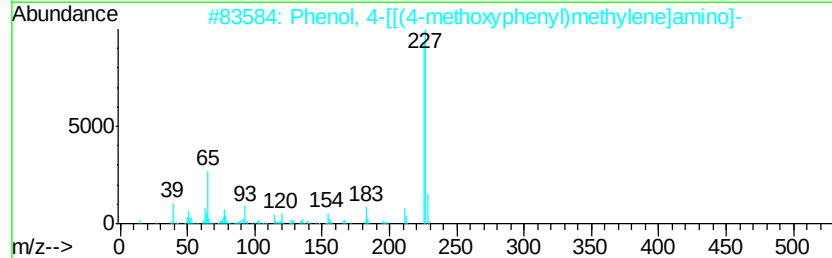
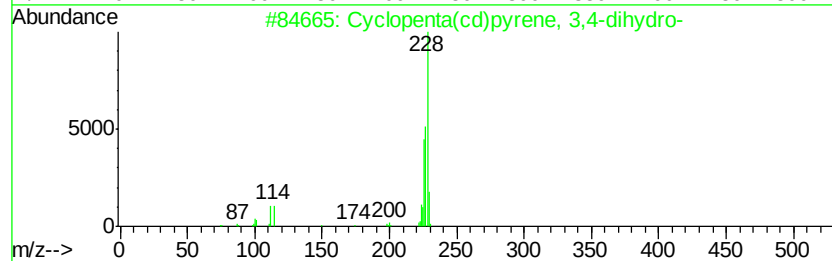
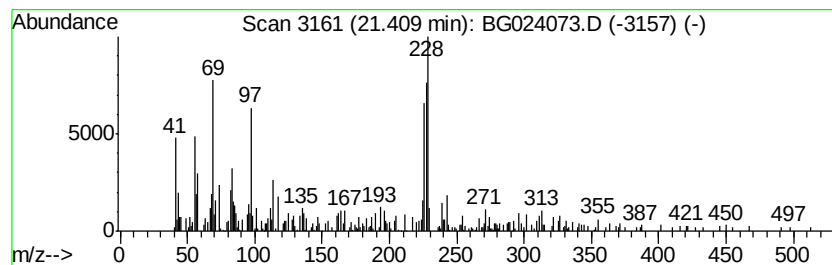
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown21.41 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.41	2.40 ng	158476	Chrysene-d12	21.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopenta(cd)pyrene, 3,4-dihydro-	228	C18H12	025732-74-5	27
2		Phenol, 4-[[4-methoxyphenyl)met...	227	C14H13NO2	003230-39-5	18
3		2,5-Dimethyl-4-(3-nitrophenyl)py...	228	C13H12N2O2	068018-39-3	15
4		2-Methoxy-5-phenyliminomethyl-ph...	227	C14H13NO2	1000294-90-8	14
5		Naphtho[2,3-d]pyrazole-4,9(4H,9H...	227	C12H9N3O2	1000263-96-1	14



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092216\
 Data File : BG024073.D
 Acq On : 22 Sep 2016 17:47
 Operator : UM/SJ
 Sample : H4820-03
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG083116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propane, 2,2-dime...	2.92	230.9	ng	5762760	1	8.04	499071	20.0
2-Pentanone, 4-hy...	5.11	13.9	ng	348161	1	8.04	499071	20.0
unknown7.58	7.58	86.0	ng	2145880	1	8.04	499071	20.0
Tridecane	16.43	2.4	ng	123296	4	17.49	1013230	20.0
Phthalic acid, is...	18.19	2.1	ng	106484	4	17.49	1013230	20.0
n-Hexadecanoic acid	18.37	8.3	ng	419814	4	17.49	1013230	20.0
Phenanthrene, 1-m...	18.42	2.6	ng	130661	4	17.49	1013230	20.0
Phenanthrene, 2,3...	19.28	2.3	ng	115044	4	17.49	1013230	20.0
Octadecanoic acid	19.63	10.2	ng	517813	4	17.49	1013230	20.0
unknown20.38	20.38	2.7	ng	176372	5	21.80	1318330	20.0
unknown21.41	21.41	2.4	ng	158476	5	21.80	1318330	20.0