

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG082516\
 Data File : BG023906.D
 Acq On : 25 Aug 2016 22:58
 Operator : UM/SJ
 Sample : H4588-03
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MW-04

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-BG080116.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	5.152	389	394	403	rBV	184333	296123	4.36%	0.718%
2	5.634	467	476	486	rBV	911436	1533653	22.57%	3.720%
3	7.249	743	751	764	rBV	600486	1145309	16.86%	2.778%
4	7.619	806	814	827	rBV	1623230	2867577	42.20%	6.955%
5	8.089	886	894	900	rBV	454112	799078	11.76%	1.938%
6	8.148	900	904	916	rVB9	25955	79548	1.17%	0.193%
7	8.418	942	950	964	rBV	1412900	2550239	37.53%	6.186%
8	9.164	1072	1077	1087	rVB	87399	177633	2.61%	0.431%
9	9.270	1088	1095	1105	rBV	1272352	2345987	34.52%	5.690%
10	10.926	1369	1377	1384	rBV	691631	1245184	18.32%	3.020%
11	11.502	1466	1475	1481	rBV4	70623	169355	2.49%	0.411%
12	12.325	1609	1615	1620	rVB9	38963	96750	1.42%	0.235%
13	12.560	1648	1655	1665	rBV9	40315	145270	2.14%	0.352%
14	13.376	1786	1794	1800	rBV	3235705	5135546	75.58%	12.456%
15	13.429	1800	1803	1808	rVB7	70212	102908	1.51%	0.250%
16	13.764	1856	1860	1865	rVB7	44277	88230	1.30%	0.214%
17	13.987	1894	1898	1902	rVB7	43845	70250	1.03%	0.170%
18	14.093	1913	1916	1926	rVB6	95034	168766	2.48%	0.409%
19	14.204	1931	1935	1941	rBV	338670	472837	6.96%	1.147%
20	14.580	1996	1999	2004	rVB	78230	122178	1.80%	0.296%
21	14.680	2013	2016	2019	rBV4	64888	76116	1.12%	0.185%
22	14.762	2025	2030	2036	rBV2	972230	1652289	24.32%	4.008%
23	15.467	2146	2150	2153	rBV5	120145	173378	2.55%	0.421%
24	15.585	2167	2170	2175	rBV4	90538	122997	1.81%	0.298%
25	16.155	2261	2267	2274	rVB	778982	1392670	20.50%	3.378%
26	16.266	2280	2286	2294	rBV	2552342	3918385	57.67%	9.504%
27	17.330	2464	2467	2475	rVB	348090	482887	7.11%	1.171%
28	17.529	2497	2501	2507	rVB2	1315560	1940329	28.55%	4.706%
29	18.410	2648	2651	2657	rVB5	253656	302155	4.45%	0.733%
30	19.673	2863	2866	2872	rVB2	191909	276851	4.07%	0.672%
31	20.137	2940	2945	2954	rVB	4869043	6795059	100.00%	16.482%
32	21.847	3230	3236	3243	rVB2	1340228	2295166	33.78%	5.567%
33	25.231	3803	3812	3825	rVB3	725295	2187587	32.19%	5.306%

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Integration Parameters: rteint.p
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_G\METHODS\8270-BG080116.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

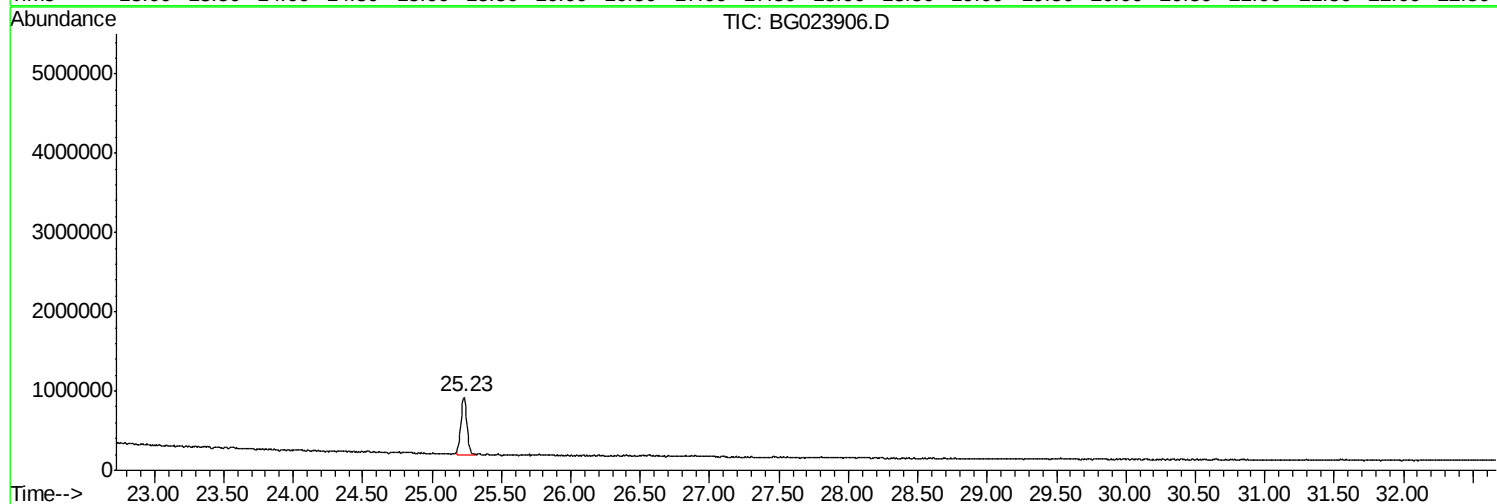
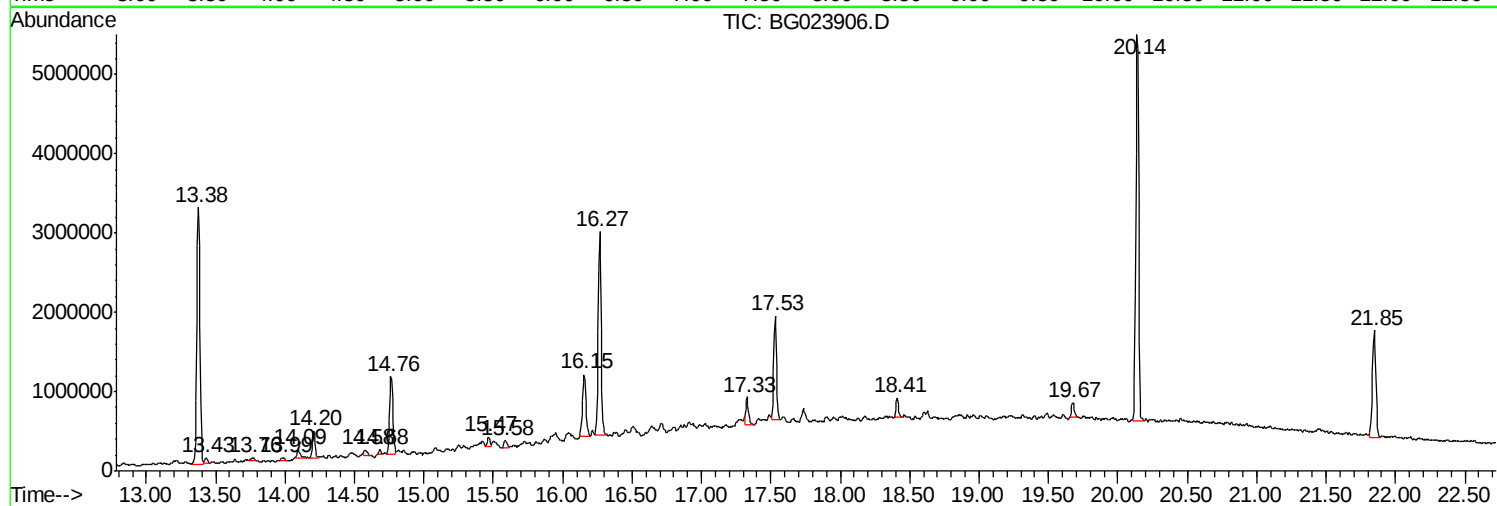
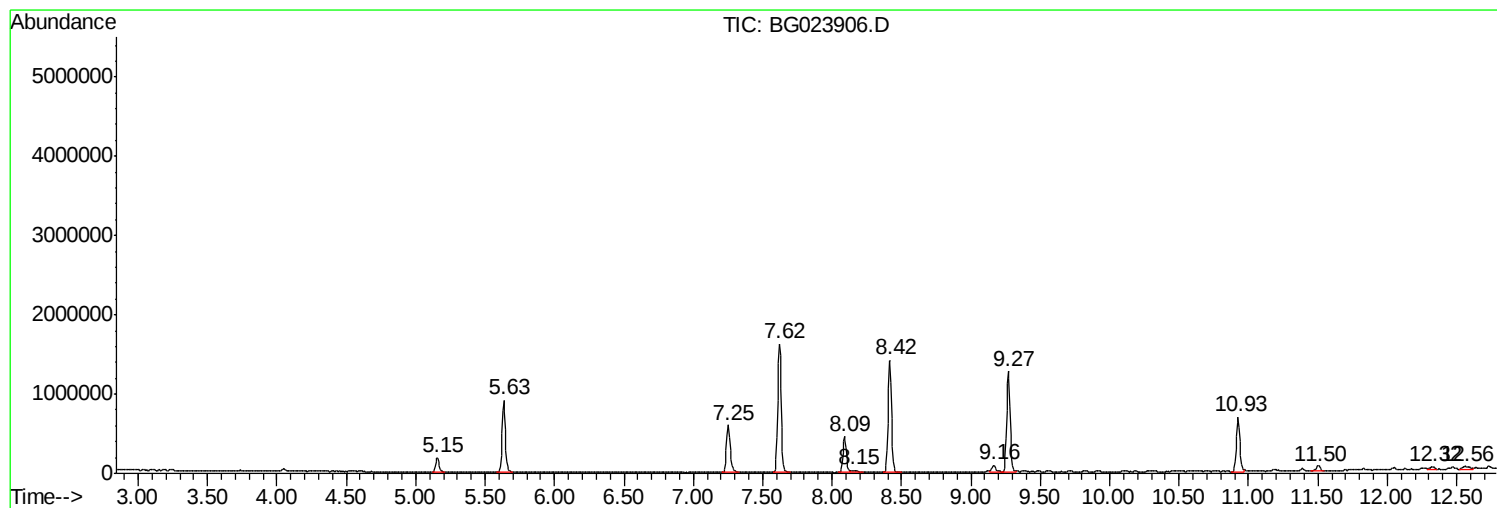
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG080116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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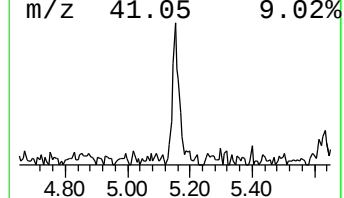
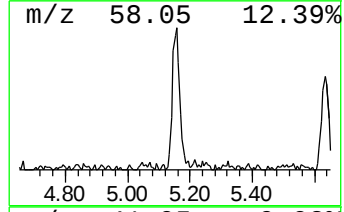
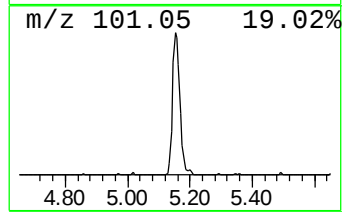
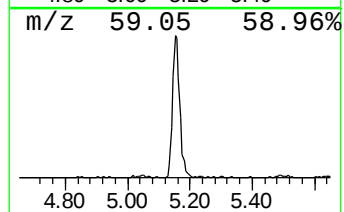
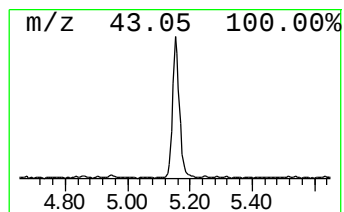
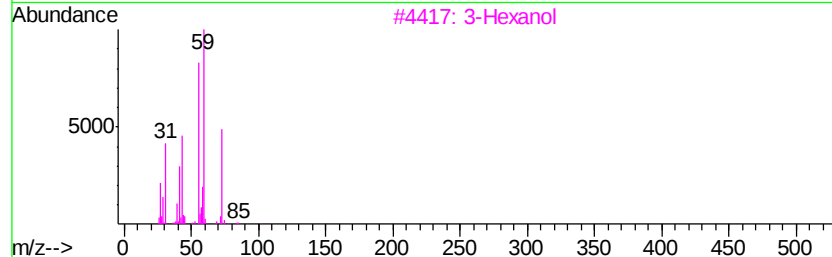
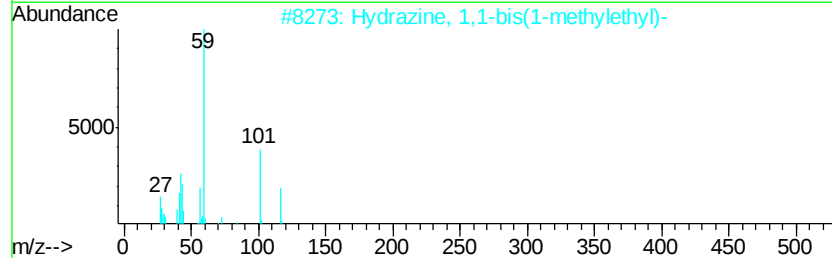
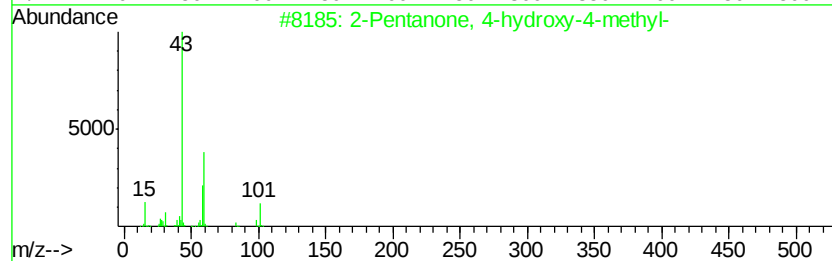
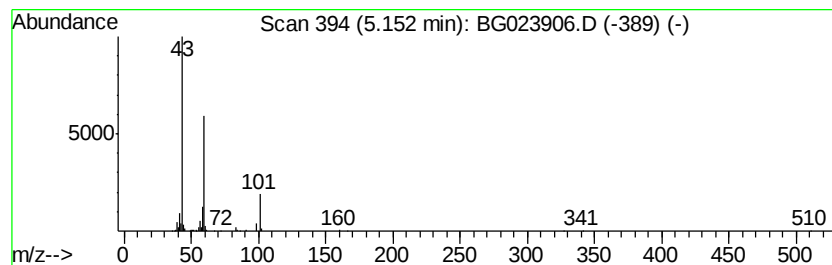
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.15	7.41 ng	296123	1,4-Dichlorobenzene-d4	8.09

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	38
3			3-Hexanol	102	C6H14O	000623-37-0	36
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	35
5			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	32



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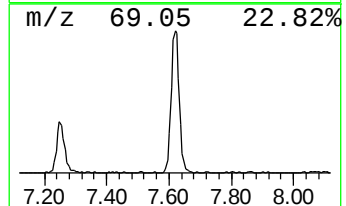
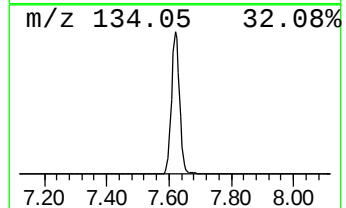
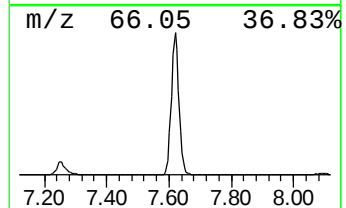
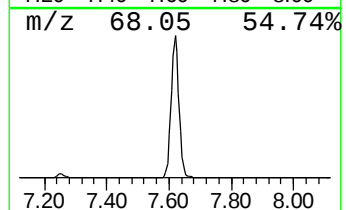
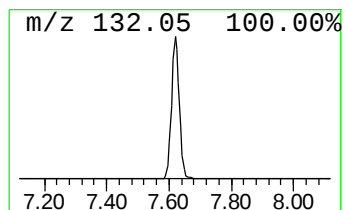
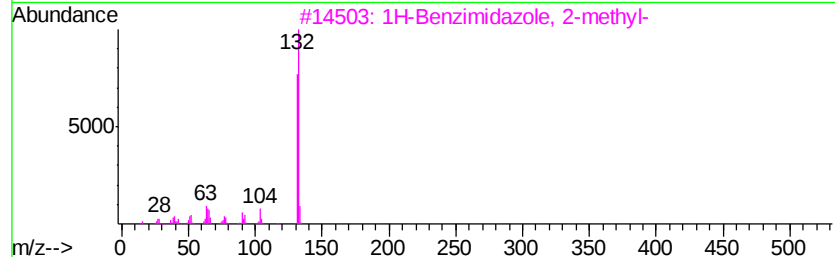
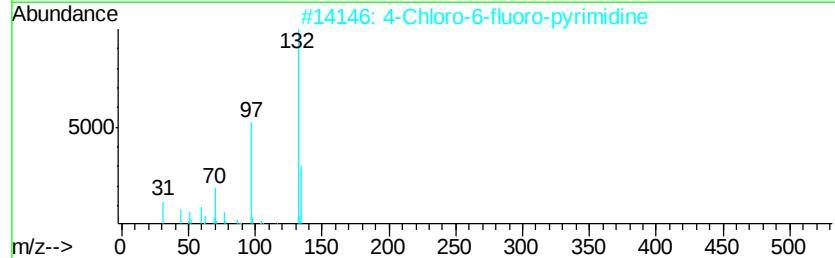
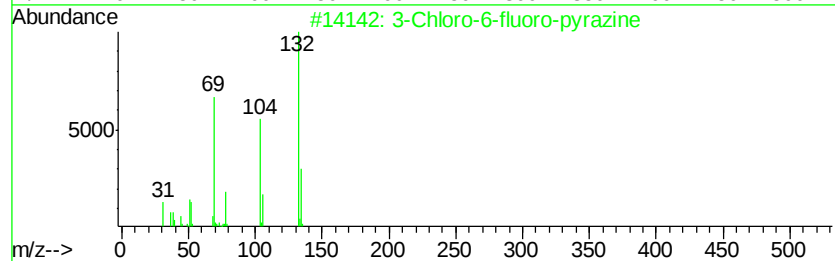
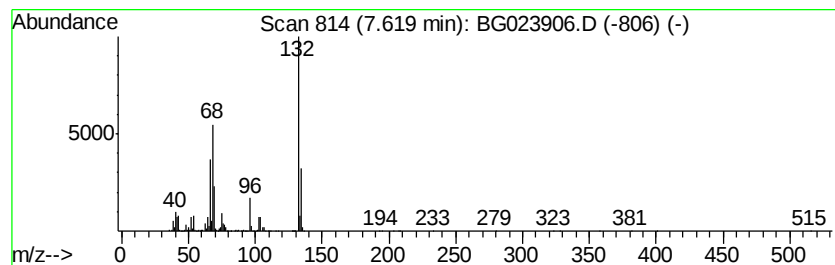
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 unknown7.62 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.62	71.77 ng	2867580	1,4-Dichlorobenzene-d4	8.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	22
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	16



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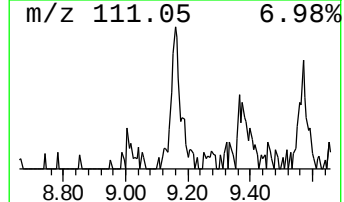
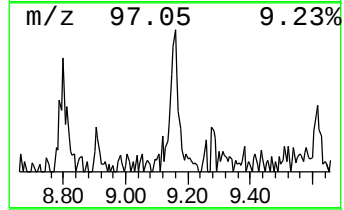
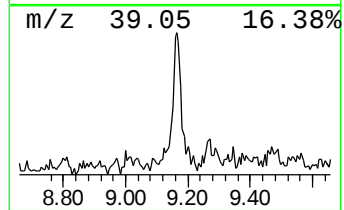
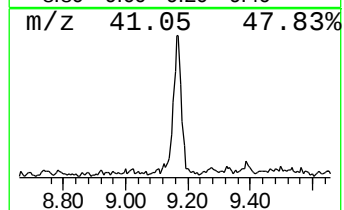
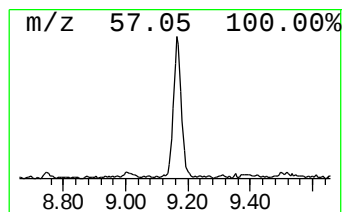
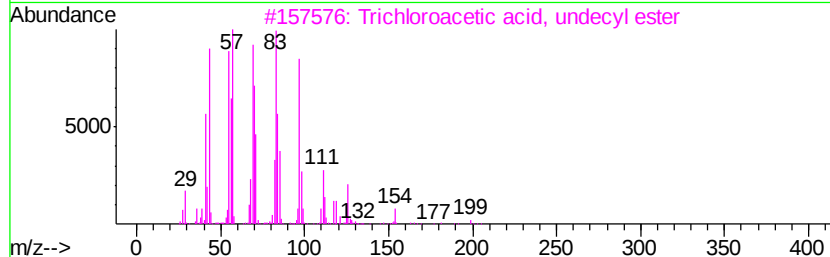
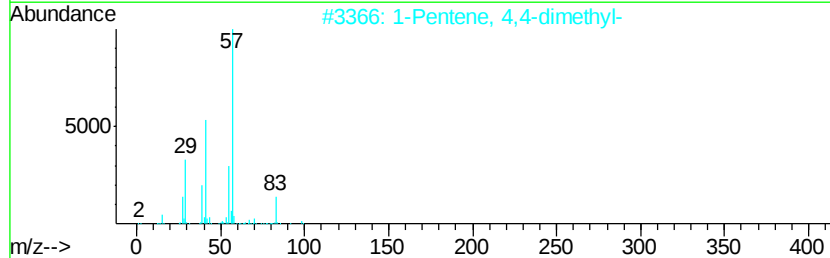
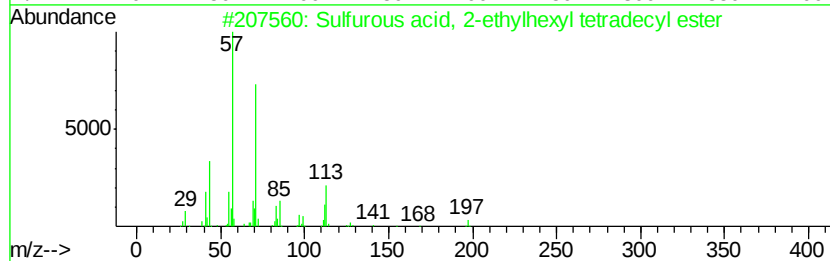
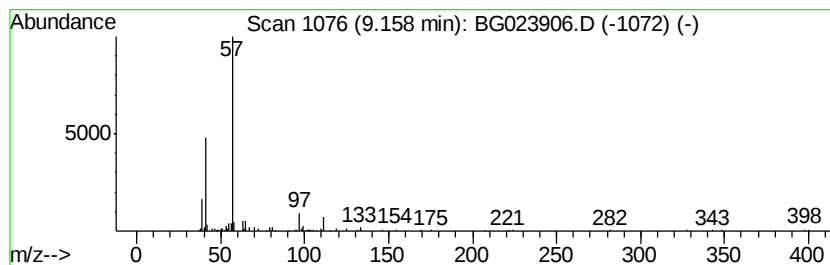
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown9.16 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.16	4.45 ng	177633	1,4-Dichlorobenzene-d4	8.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Sulfurous acid, 2-ethylhexyl tet...	390	C22H46O3S	1000309-19-7	28
2		1-Pentene, 4,4-dimethyl-	98	C7H14	000762-62-9	9
3		Trichloroacetic acid, undecyl ester	316	C13H23Cl3O2	074339-49-4	9
4		Propane, 2-bromo-2-methyl-	136	C4H9Br	000507-19-7	9
5		tert-Butyl iodide	184	C4H9I	000558-17-8	9



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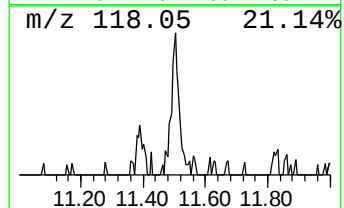
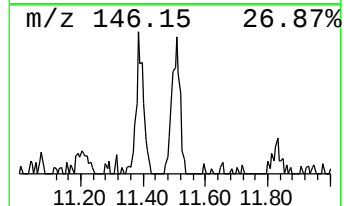
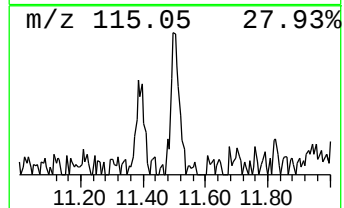
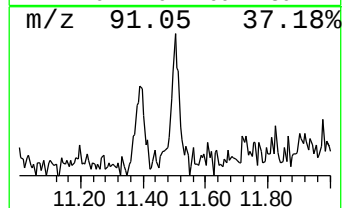
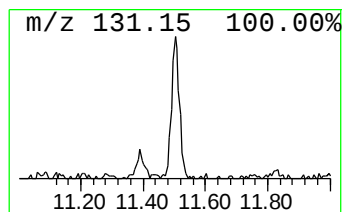
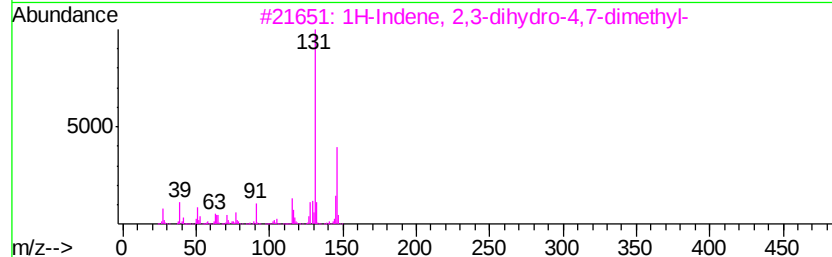
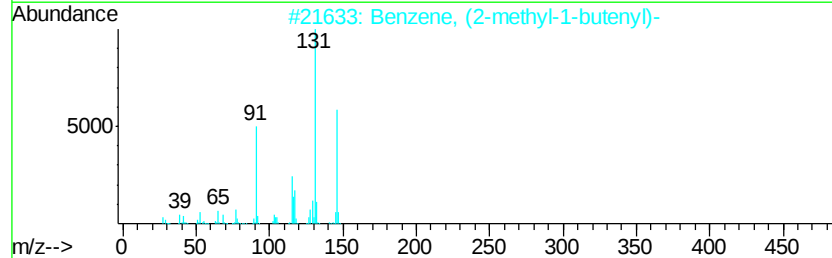
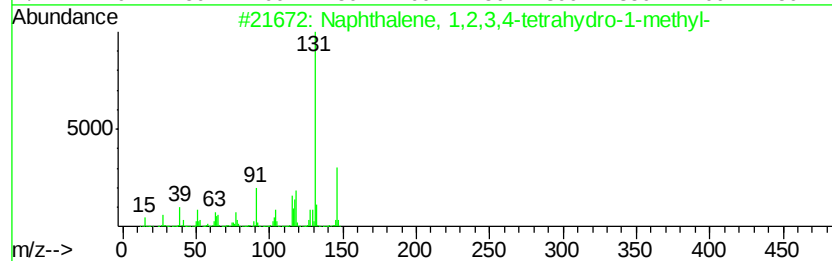
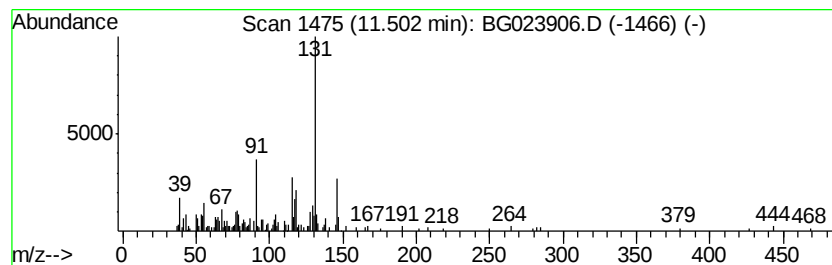
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.50	2.72 ng	169355	Naphthalene-d8	10.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	93
2		Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	90
3		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	76
4		1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	76
5		Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	74



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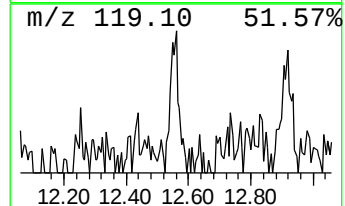
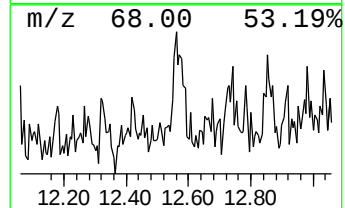
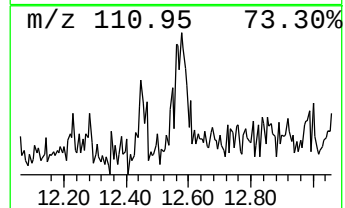
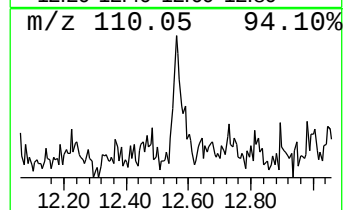
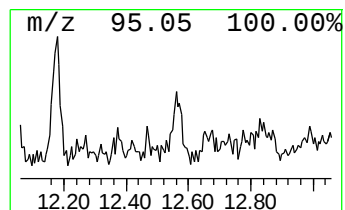
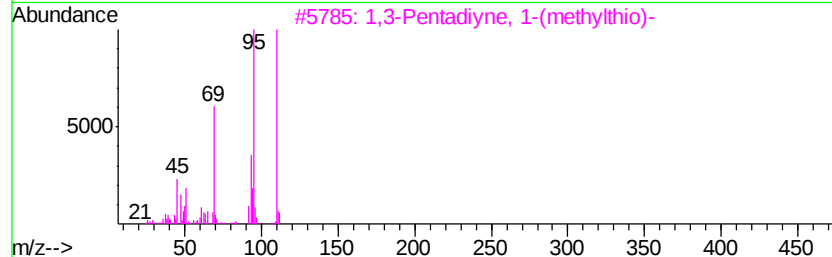
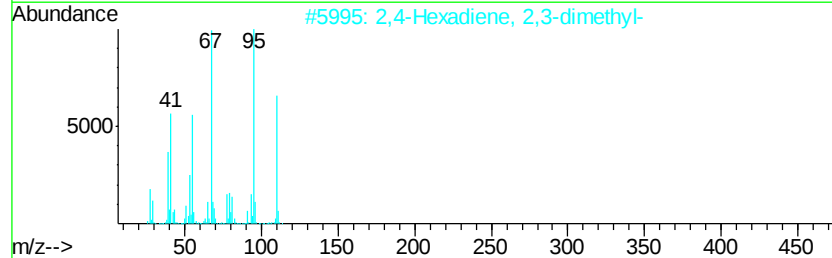
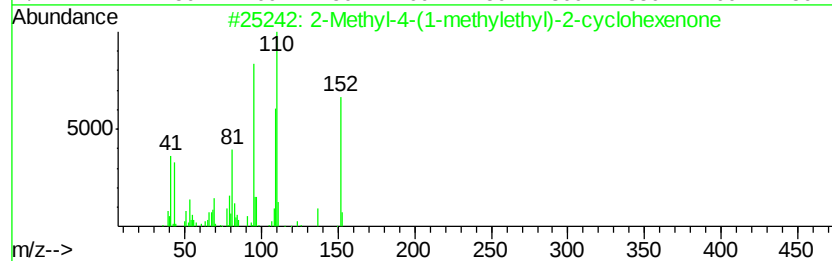
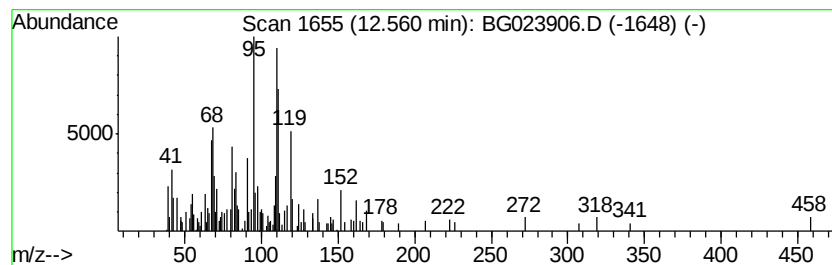
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Peak Number 6 unknown12.56 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.56	2.33 ng	145270	Naphthalene-d8	10.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Methyl-4-(1-methylethyl)-2-cyc...	152	C10H16O	041469-46-9	49
2		2,4-Hexadiene, 2,3-dimethyl-	110	C8H14	005678-98-8	35
3		1,3-Pentadiyne, 1-(methylthio)-	110	C6H6S	042875-80-9	30
4		Bicyclo[2.2.1]heptan-2-one, 5,5,...	152	C10H16O	003649-86-3	25
5		2(4H)-Benzofuranone, 5,6,7,7a-te...	180	C11H16O2	017092-92-1	18



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG082516\
Data File : BG023906.D
Acq On : 25 Aug 2016 22:58
Operator : UM/SJ
Sample : H4588-03
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleID :
MW-04

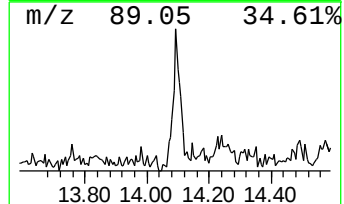
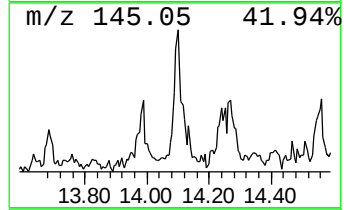
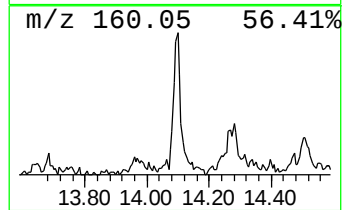
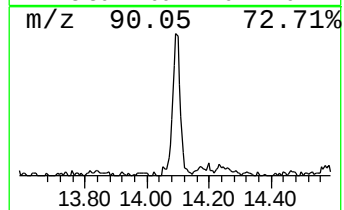
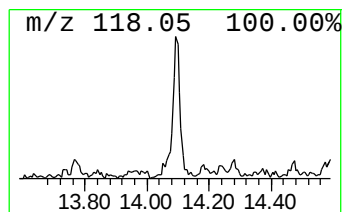
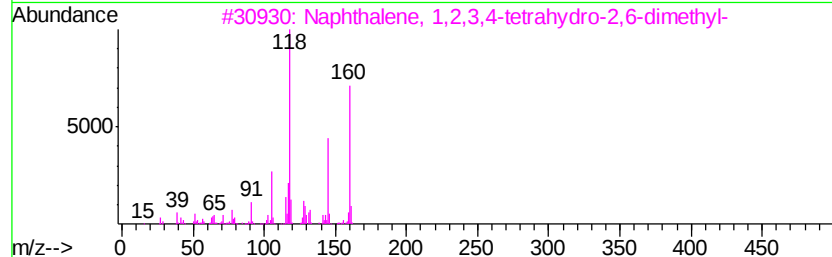
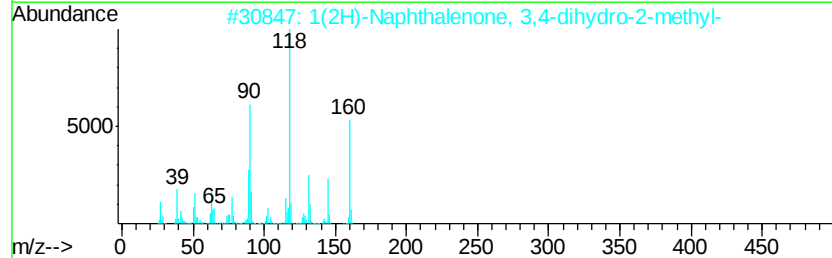
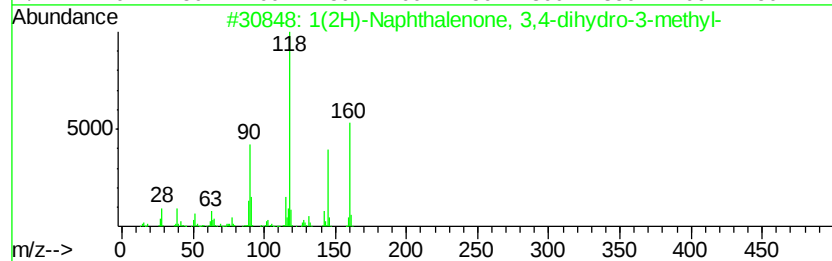
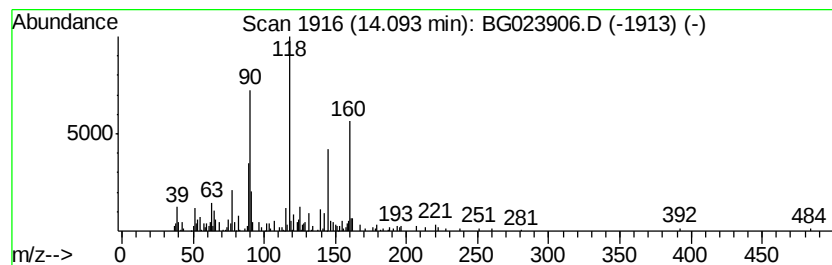
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 1(2H)-Naphthalenone, 3,4-di... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.09	2.04 ng	168766	Acenaphthene-d10	14.76

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1(2H)-Naphthalenone, 3,4-dihydro...	160	C11H12O	014944-23-1	80
2		1(2H)-Naphthalenone, 3,4-dihydro...	160	C11H12O	001590-08-5	64
3		Naphthalene, 1,2,3,4-tetrahydro...	160	C12H16	007524-63-2	52
4		1(2H)-Naphthalenone, 3,4-dihydro-	146	C10H10O	000529-34-0	43
5		Benzofuran	118	C8H6O	000271-89-6	38



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG082516\
Data File : BG023906.D
Acq On : 25 Aug 2016 22:58
Operator : UM/SJ
Sample : H4588-03
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
MW-04

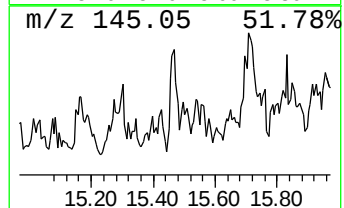
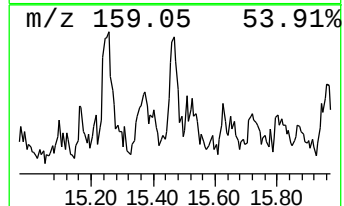
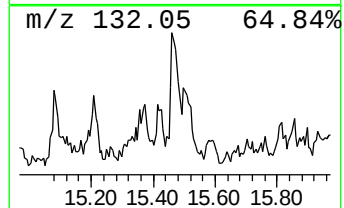
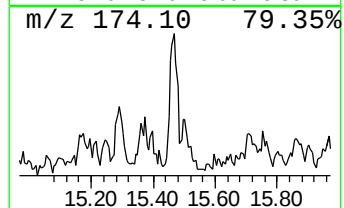
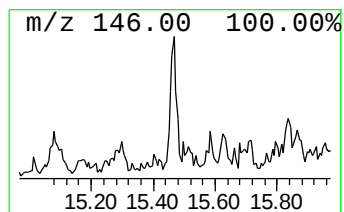
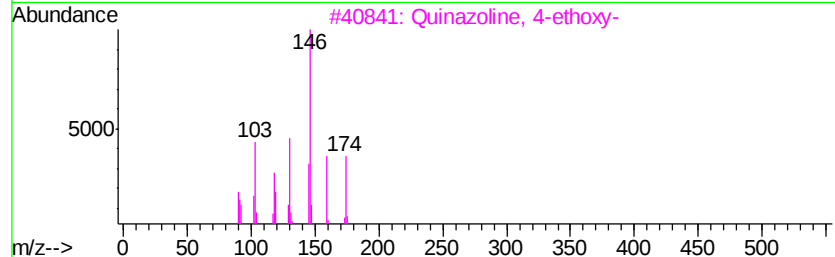
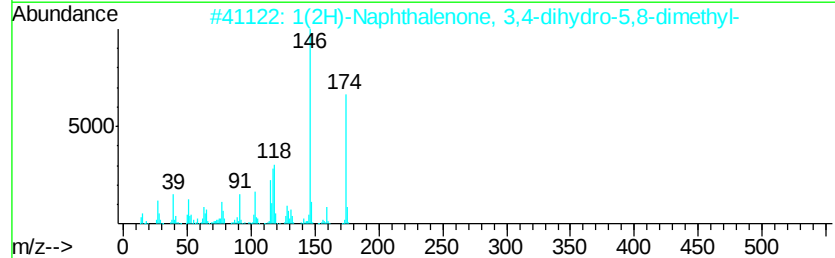
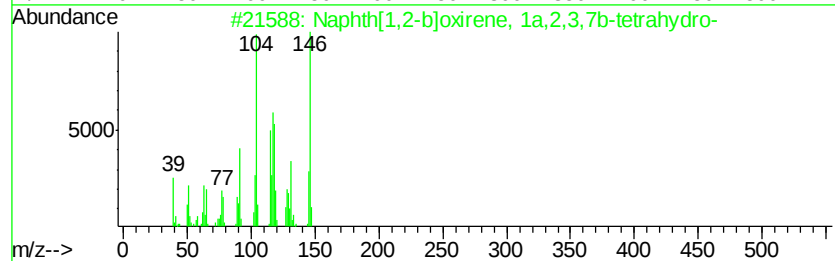
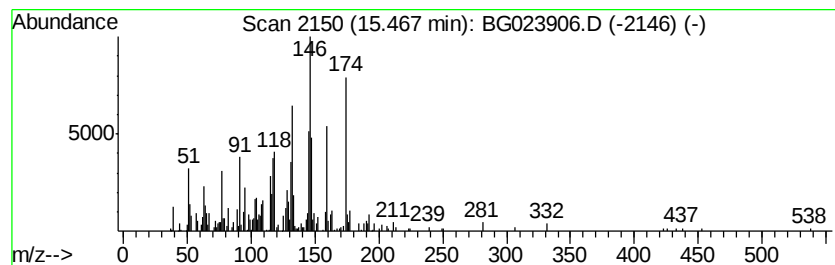
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Naphth[1,2-b]oxirene, 1a,2,... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.47	2.10 ng	173378	Acenaphthene-d10	14.76

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphth[1,2-b]oxirene, 1a,2,3,7b-...	146	C10H10O	002461-34-9	60
2		1(2H)-Naphthalenone, 3,4-dihydro...	174	C12H14O	005037-63-8	53
3		Quinazoline, 4-ethoxy-	174	C10H10N2O	016347-96-9	49
4		Benzene, 1,3-diisocyanato-2-methyl-	174	C9H6N2O2	000091-08-7	46
5		1(2H)-Naphthalenone, 3,4-dihydro...	174	C12H14O	032281-65-5	46



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG082516\
Data File : BG023906.D
Acq On : 25 Aug 2016 22:58
Operator : UM/SJ
Sample : H4588-03
Misc :
ALS Vial : 15 Sample Multiplier: 1

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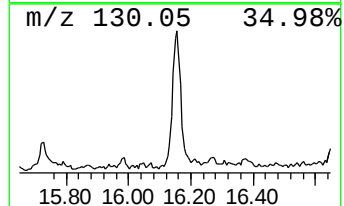
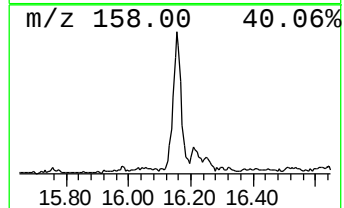
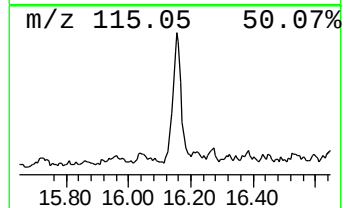
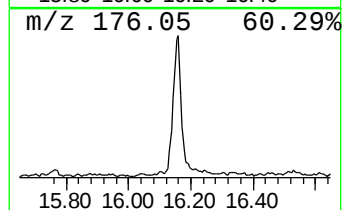
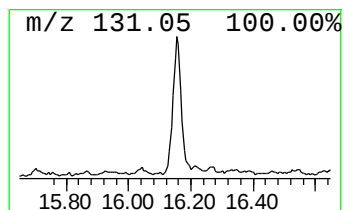
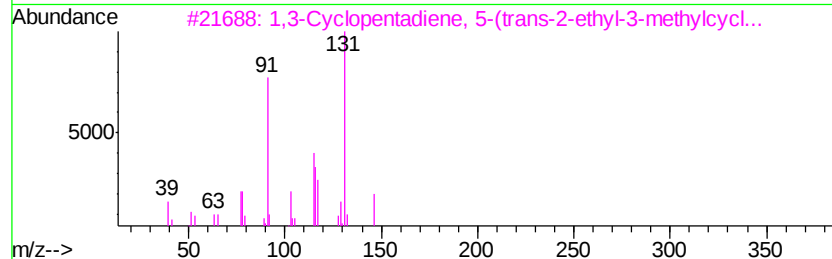
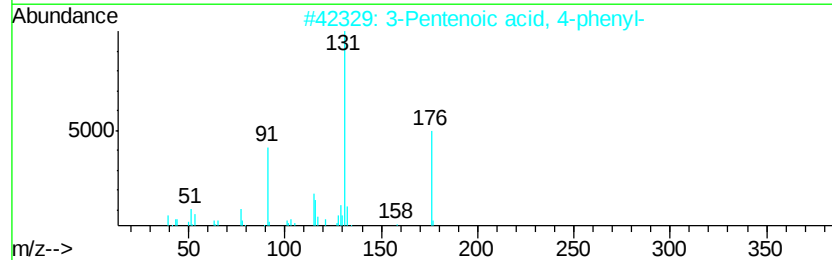
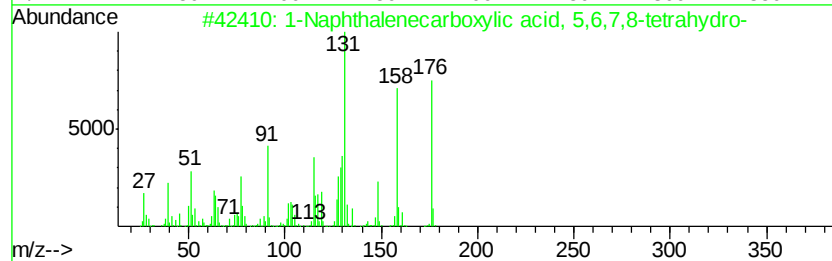
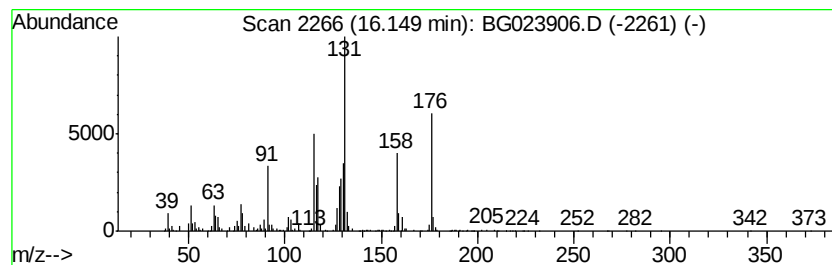
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 1-Naphthalenecarboxylic aci... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.15	14.36 ng	1392670	Phenanthrene-d10	17.53

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Naphthalenecarboxylic acid, 5,...	176	C11H12O2	004242-18-6	93
2		3-Pentenoic acid, 4-phenyl-	176	C11H12O2	053774-19-9	46
3		1,3-Cyclopentadiene, 5-(trans-2-...	146	C11H14	079209-36-2	43
4		Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	38
5		Cinnamic acid, .alpha.,.beta.-di...	176	C11H12O2	1000151-56-4	38



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG082516\
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Acq On : 25 Aug 2016 22:58
Operator : UM/SJ
Sample : H4588-03
Misc :
ALS Vial : 15 Sample Multiplier: 1

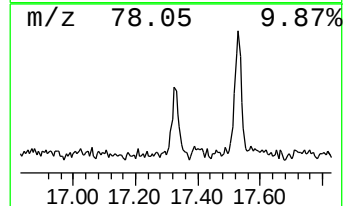
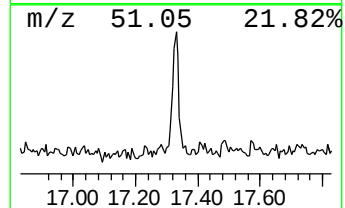
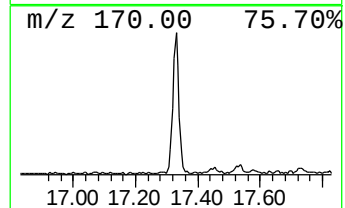
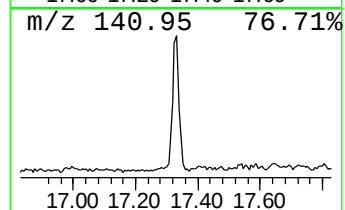
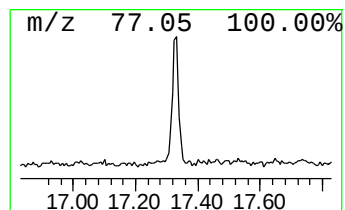
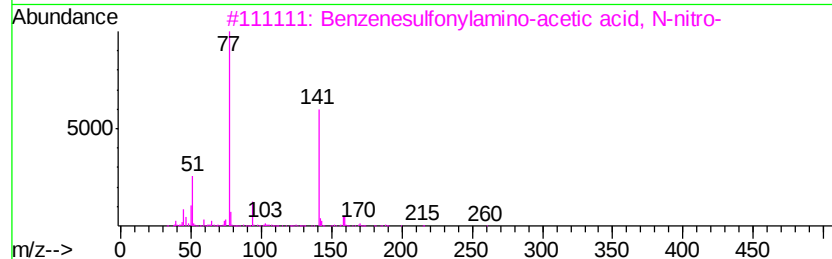
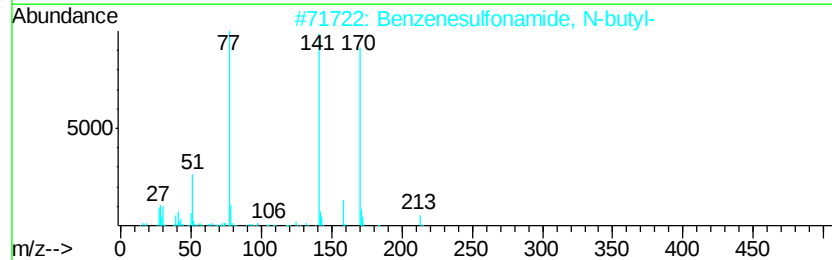
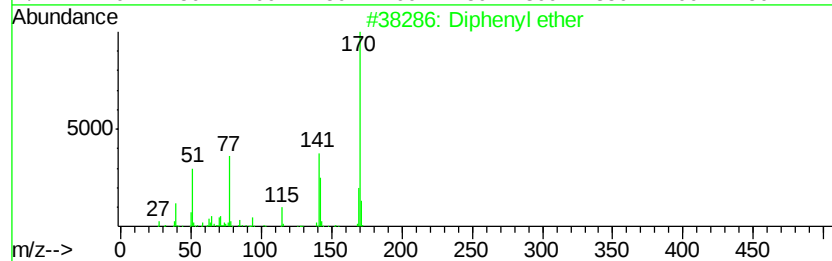
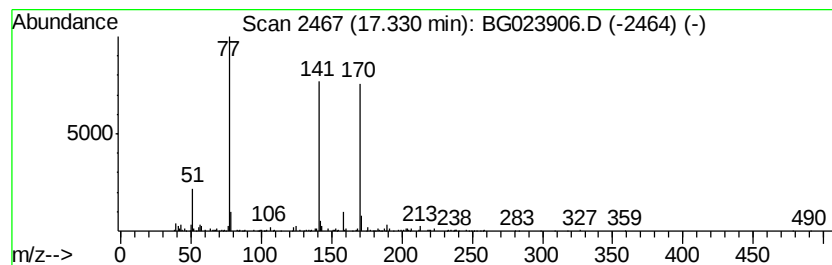
Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG080116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Diphenyl ether Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.33	4.98 ng	482887	Phenanthrene-d10	17.53	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diphenyl ether	170	C12H10O	000101-84-8	64
2	Benzenesulfonamide, N-butyl-	213	C10H15NO2S	003622-84-2	58
3	Benzenesulfonylamino-acetic acid...	260	C8H8N2O6S	1000301-09-0	53
4	N,N-Dichlorobenzenesulfonamide	225	C6H5Cl2NO2S	000473-29-0	50
5	3-Hydroxybiphenyl	170	C12H10O	000580-51-8	43



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG082516\
Data File : BG023906.D
Acq On : 25 Aug 2016 22:58
Operator : UM/SJ
Sample : H4588-03
Misc :
ALS Vial : 15 Sample Multiplier: 1

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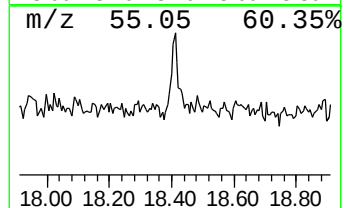
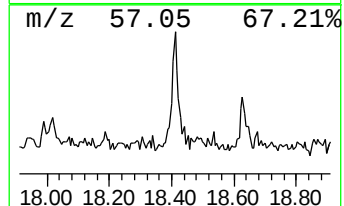
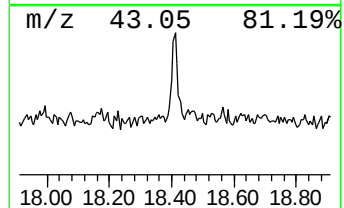
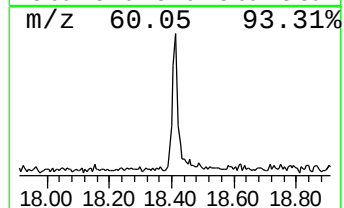
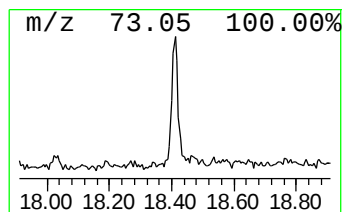
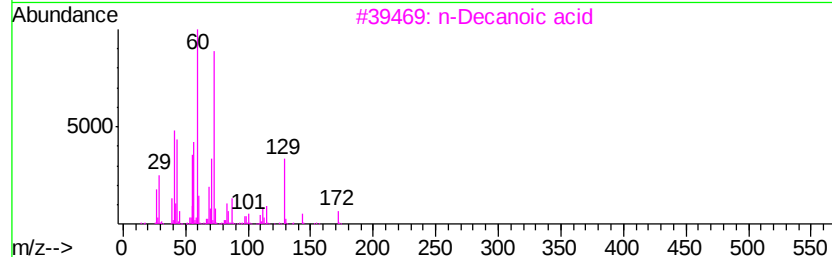
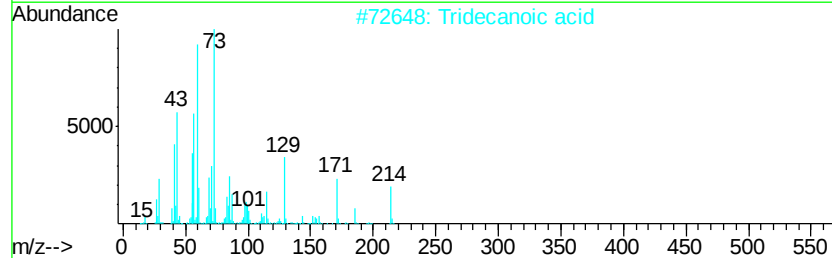
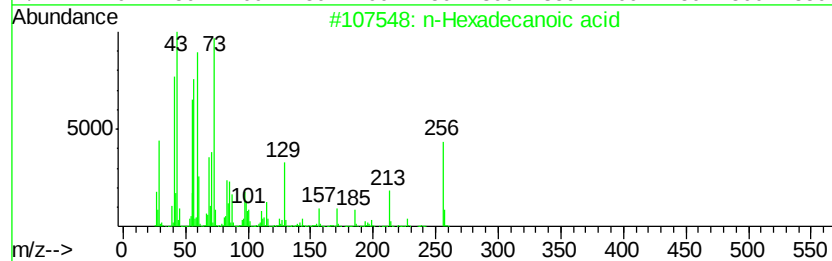
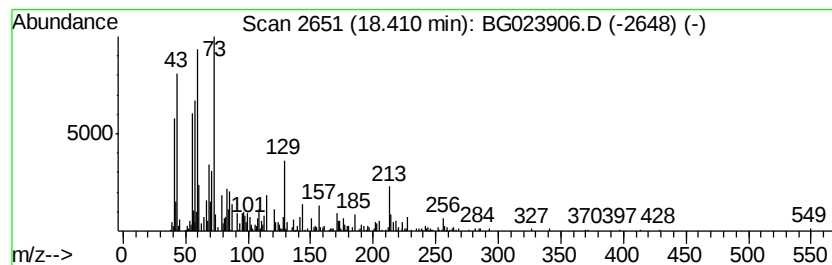
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.41	3.11 ng	302155	Phenanthrene-d10	17.53

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	83
2			Tridecanoic acid	214	C13H26O2	000638-53-9	74
3			n-Decanoic acid	172	C10H20O2	000334-48-5	64
4			Pentadecanoic acid	242	C15H30O2	001002-84-2	64
5			Dodecanoic acid	200	C12H24O2	000143-07-7	59



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG082516\
Data File : BG023906.D
Acq On : 25 Aug 2016 22:58
Operator : UM/SJ
Sample : H4588-03
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
MW-04

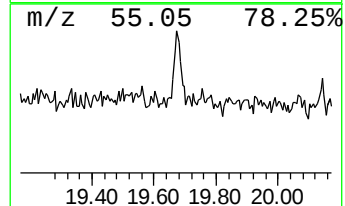
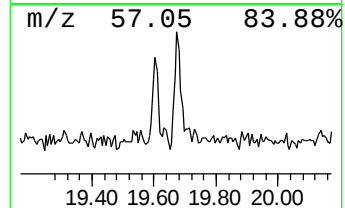
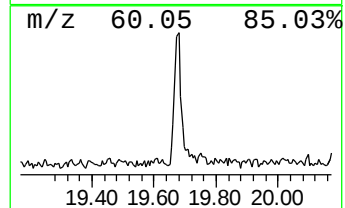
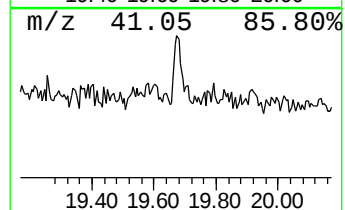
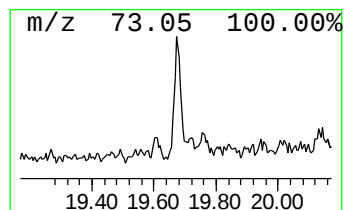
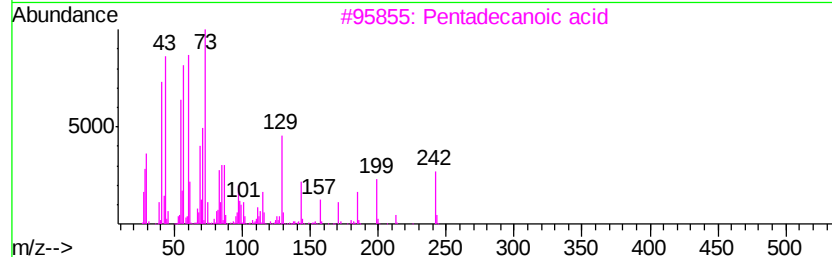
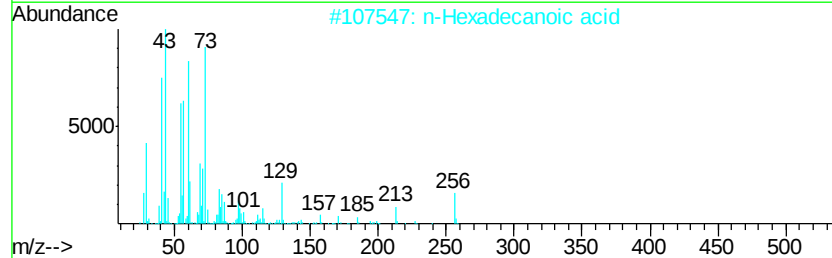
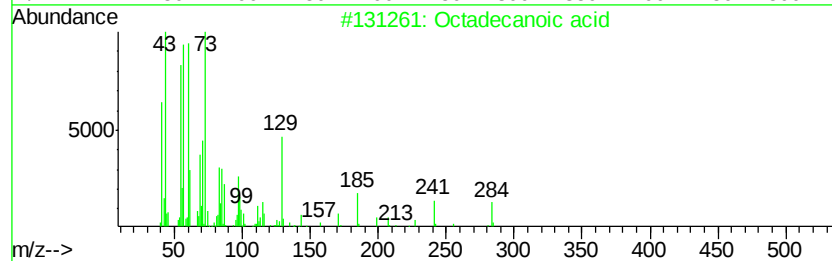
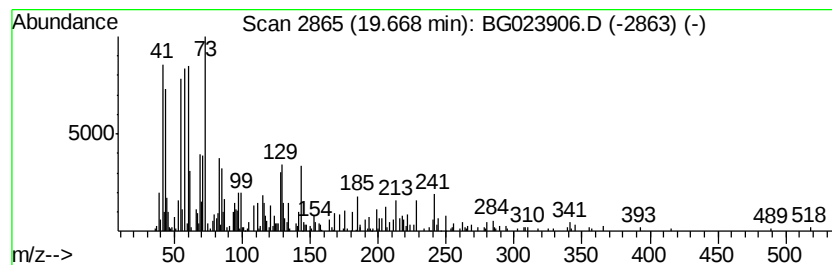
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Octadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.67	2.85 ng	276851	Phenanthrene-d10	17.53

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	93
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	90
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	72
5		Hexadecane, 1-chloro-	260	C16H33Cl	004860-03-1	20



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG082516\
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Acq On : 25 Aug 2016 22:58
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Sample : H4588-03
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG080116.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
2-Pentanone, 4-hy...	5.15	7.4	ng	296123	1	8.09	799078	20.0
unknown7.62	7.62	71.8	ng	2867580	1	8.09	799078	20.0
unknown9.16	9.16	4.5	ng	177633	1	8.09	799078	20.0
Naphthalene, 1,2,...	11.50	2.7	ng	169355	2	10.93	1245180	20.0
unknown12.56	12.56	2.3	ng	145270	2	10.93	1245180	20.0
1(2H)-Naphthaleno...	14.09	2.0	ng	168766	3	14.76	1652290	20.0
Naphth[1,2-b]oxir...	15.47	2.1	ng	173378	3	14.76	1652290	20.0
1-Naphthalenecarb...	16.15	14.4	ng	1392670	4	17.53	1940330	20.0
Diphenyl ether	17.33	5.0	ng	482887	4	17.53	1940330	20.0
n-Hexadecanoic acid	18.41	3.1	ng	302155	4	17.53	1940330	20.0
Octadecanoic acid	19.67	2.9	ng	276851	4	17.53	1940330	20.0