

Data Path : Z:\HPCHEM1\BNA M\DATA\BM051117\
 Data File : BM010008.D
 Acq On : 12 May 2017 05:18
 Operator : SJ/MA
 Sample : I3079-05
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DUP-0399-170509

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_M\METHODS\8270-BM051117.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	4.893	303	307	318	rBV	215637	314113	6.15%	0.819%
2	5.357	378	386	394	rBV	426079	665439	13.03%	1.734%
3	5.469	397	405	419	rVV	2146655	3209144	62.86%	8.365%
4	5.957	484	488	498	rVB	34995	54513	1.07%	0.142%
5	6.946	648	656	668	rBV	360210	633416	12.41%	1.651%
6	7.293	709	715	724	rBV	1109226	1815655	35.56%	4.732%
7	7.757	788	794	803	rBV	257157	423488	8.29%	1.104%
8	8.075	841	848	860	rBV	1000756	1645364	32.23%	4.289%
9	8.934	986	994	1007	rBV	1016913	1713832	33.57%	4.467%
10	10.557	1262	1270	1272	rBV	365908	654098	12.81%	1.705%
11	11.439	1409	1420	1432	rBV	892309	1617626	31.68%	4.216%
12	13.028	1683	1690	1699	rVB	2303495	3325564	65.14%	8.668%
13	13.539	1770	1777	1790	rBV2	240341	407149	7.97%	1.061%
14	13.880	1829	1835	1842	rBV	66993	103621	2.03%	0.270%
15	14.416	1919	1926	1935	rVB2	468415	733777	14.37%	1.913%
16	15.816	2157	2164	2169	rBV2	61205	97784	1.92%	0.255%
17	15.874	2169	2174	2176	rVV	117518	170951	3.35%	0.446%
18	15.916	2176	2181	2195	rVB	1951740	2764216	54.14%	7.205%
19	16.098	2208	2212	2221	rBV2	32067	55017	1.08%	0.143%
20	17.168	2389	2394	2404	rBV	595133	885316	17.34%	2.308%
21	18.057	2538	2545	2564	rBV	1460140	2193545	42.97%	5.717%
22	19.192	2731	2738	2740	rBV	922967	1235916	24.21%	3.221%
23	19.221	2740	2743	2746	rVV	1107316	1577286	30.89%	4.111%
24	19.262	2746	2750	2759	rVV	1776541	3101850	60.76%	8.085%
25	19.333	2759	2762	2775	rVB	635983	979046	19.18%	2.552%
26	19.474	2782	2786	2796	rBV5	104609	250005	4.90%	0.652%
27	19.698	2819	2824	2834	rBV2	224672	378743	7.42%	0.987%
28	19.792	2835	2840	2847	rVB	3942843	5105422	100.00%	13.307%
29	20.327	2928	2931	2938	rBV3	113532	165640	3.24%	0.432%
30	20.427	2946	2948	2956	rVB7	36921	52070	1.02%	0.136%
31	21.045	3050	3053	3058	rVB4	87794	106293	2.08%	0.277%
32	21.356	3101	3106	3111	rBV	809049	1063622	20.83%	2.772%
33	23.650	3490	3496	3504	rVB3	443701	866220	16.97%	2.258%

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ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DUP-0399-170509

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_M\METHODS\8270-BM051117.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

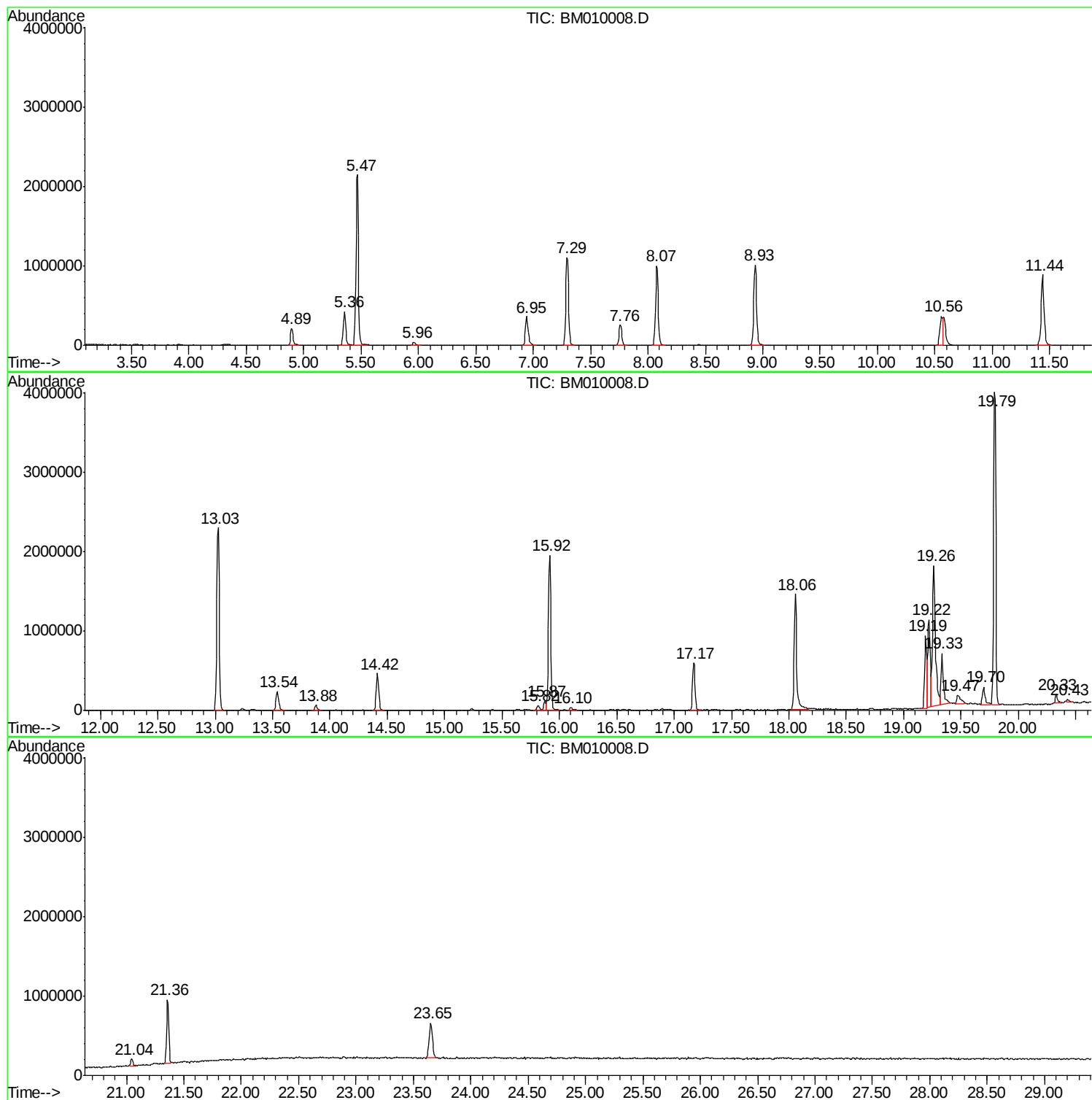
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Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM051117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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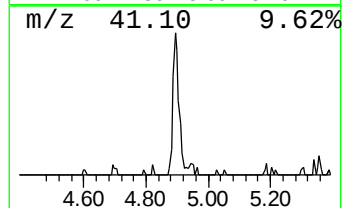
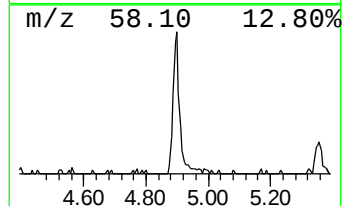
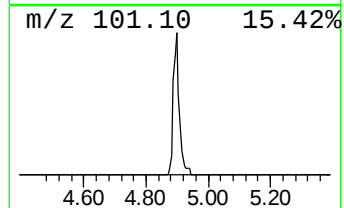
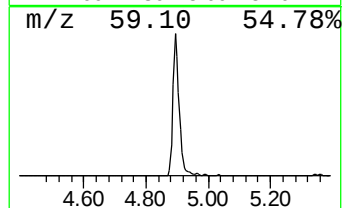
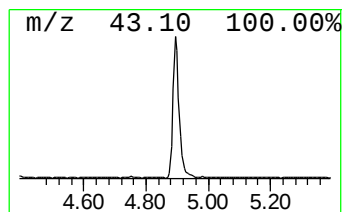
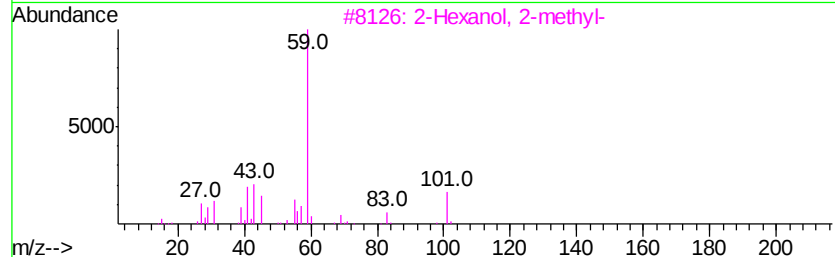
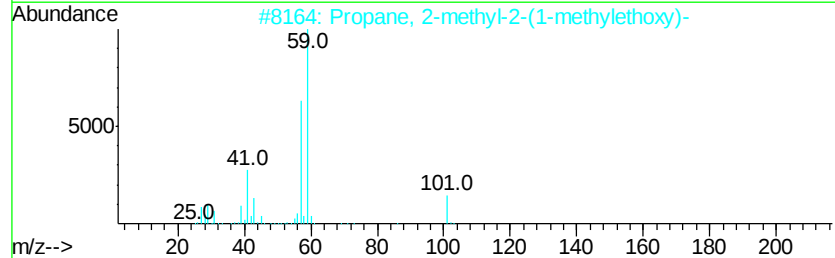
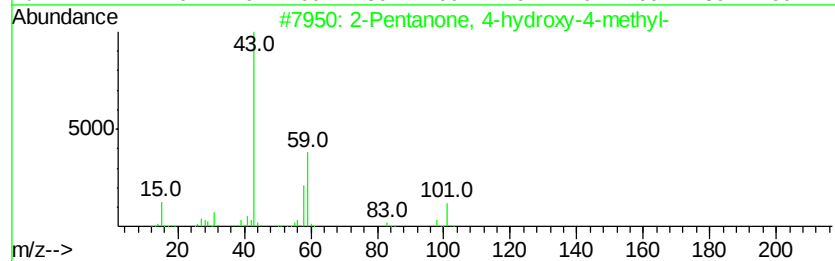
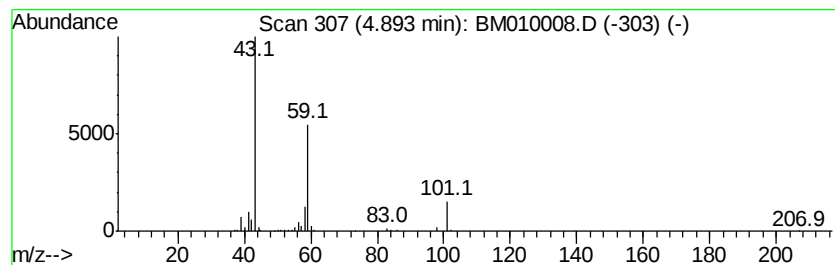
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.89	14.83 ng	314113	1,4-Dichlorobenzene-d4	7.76

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
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			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
5			Diacetamide	101	C4H7NO2	000625-77-4	9



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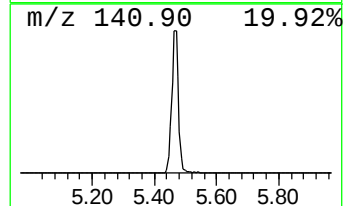
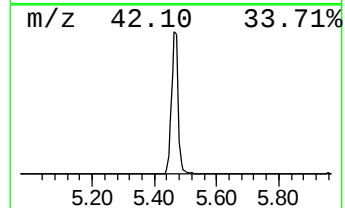
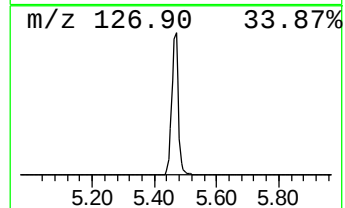
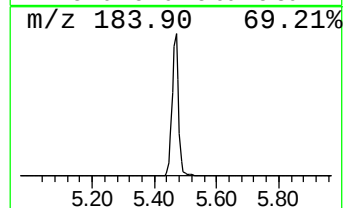
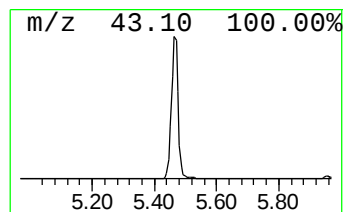
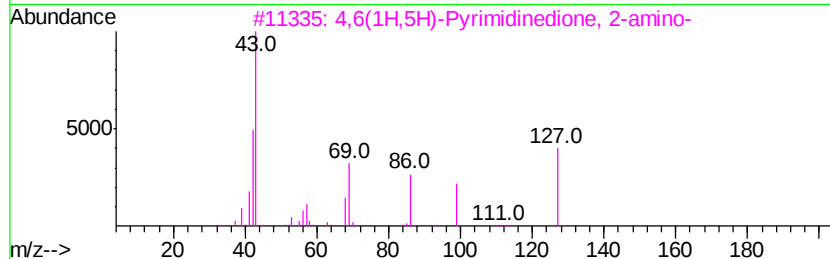
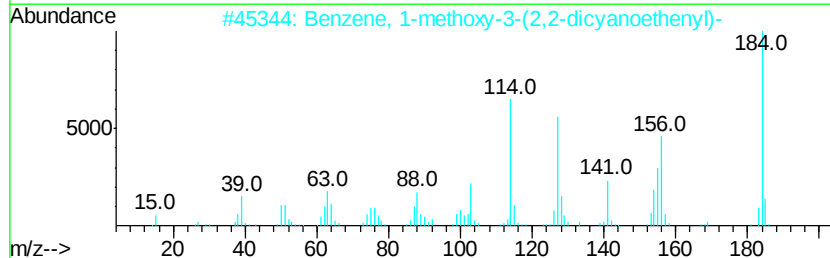
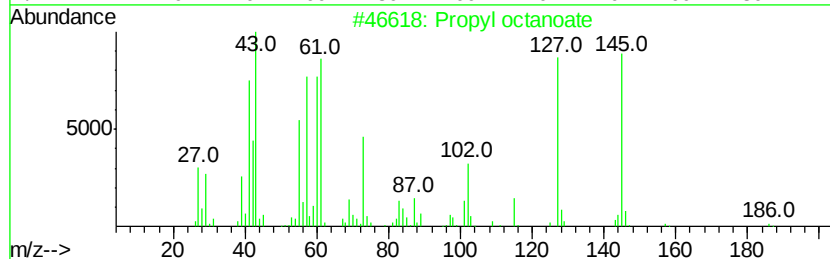
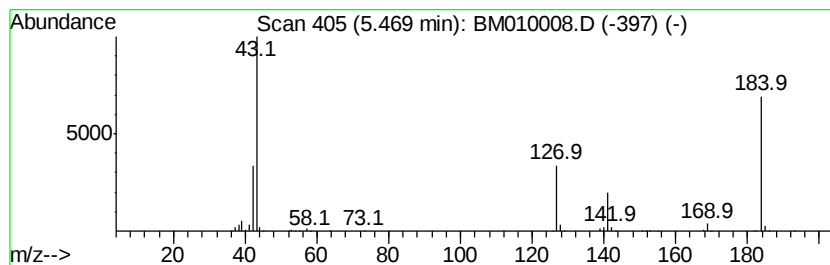
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown5.47 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.47	151.56 ng	3209140	1,4-Dichlorobenzene-d4	7.76

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propyl octanoate	186	C11H22O2	000624-13-5	35
2		Benzene, 1-methoxy-3-(2,2-dicyan...	184	C11H8N2O	002972-72-7	28
3		4,6(1H,5H)-Pyrimidinedione, 2-am...	127	C4H5N3O2	004425-67-6	17
4		6-Cyano-5-methoxyquinoline	184	C11H8N2O	1000241-27-7	9
5		1,1'-Biphenyl, 3-methoxy-	184	C13H12O	002113-56-6	9



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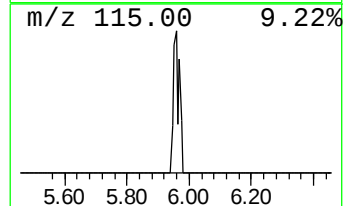
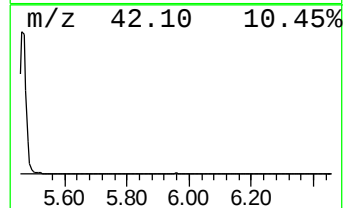
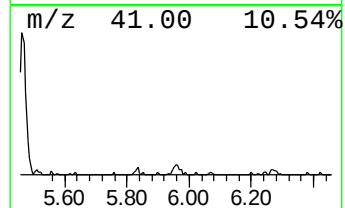
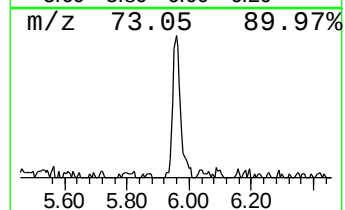
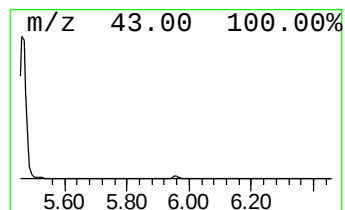
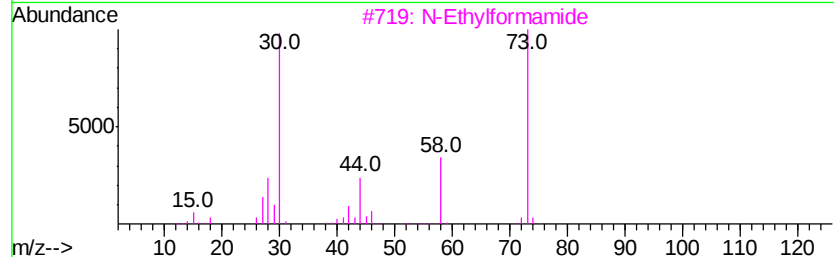
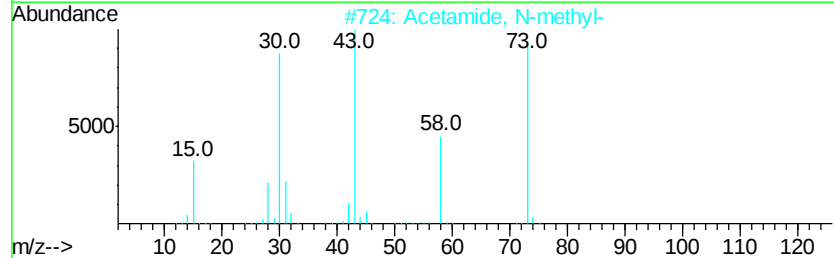
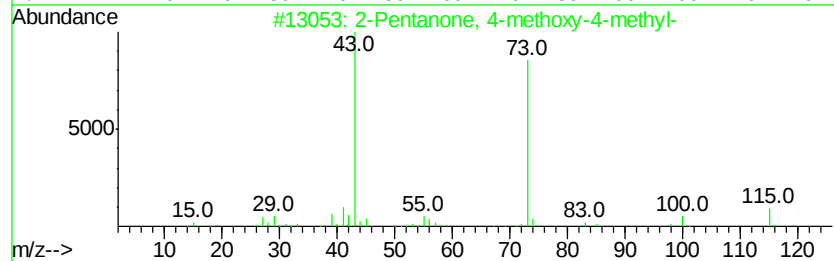
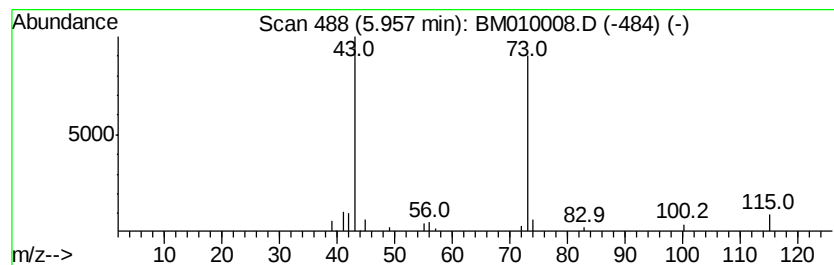
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 2-Pentanone, 4-methoxy-4-me... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.96	2.57 ng	54513	1,4-Dichlorobenzene-d4	7.76

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-methoxy-4-methyl-	130	C7H14O2	000107-70-0	83
2		Acetamide, N-methyl-	73	C3H7NO	000079-16-3	38
3		N-Ethylformamide	73	C3H7NO	000627-45-2	25
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	9
5		Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	9



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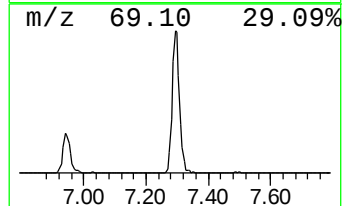
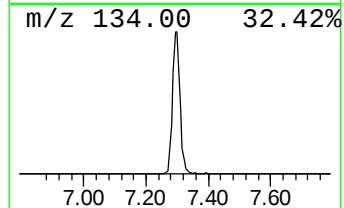
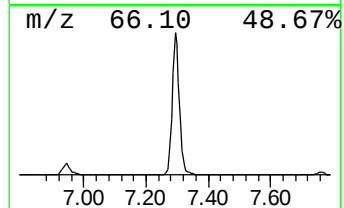
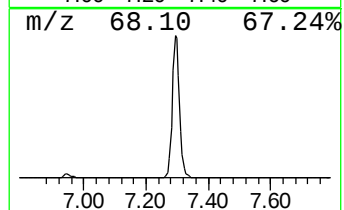
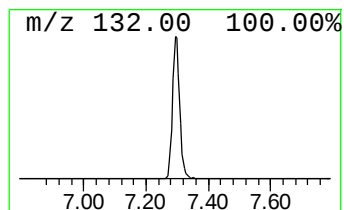
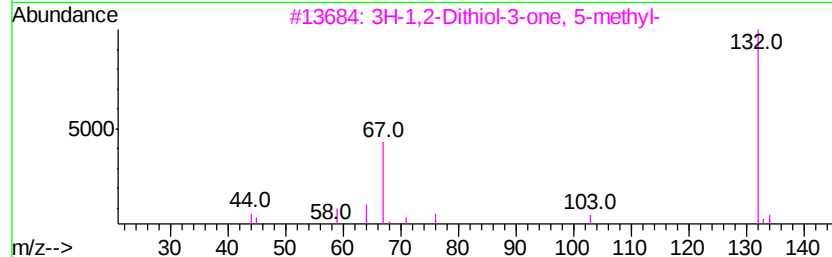
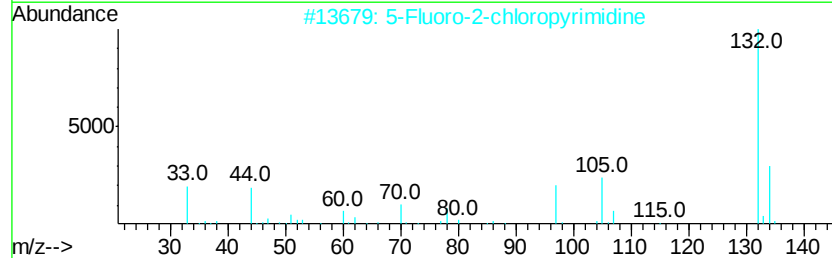
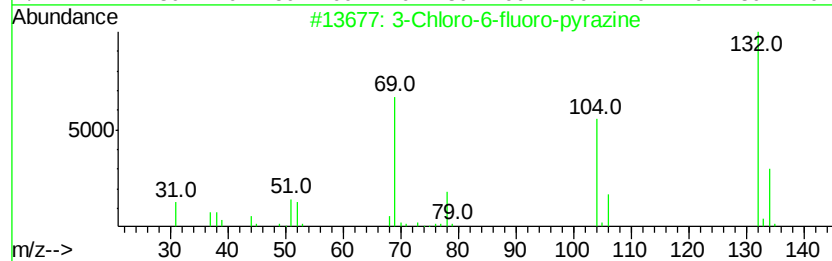
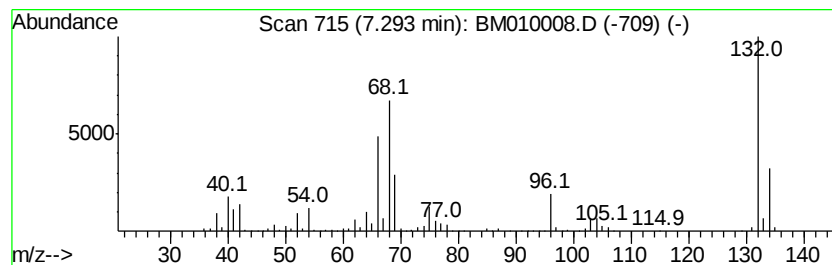
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown7.29 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.29	85.75 ng	1815660	1,4-Dichlorobenzene-d4	7.76

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	14
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	14
3		3H-1,2-Dithiol-3-one, 5-methyl-	132	C4H4OS2	003620-08-4	14
4		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	10
5		Xenon	132	Xe	007440-63-3	9



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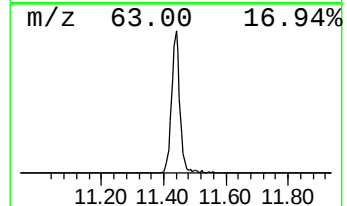
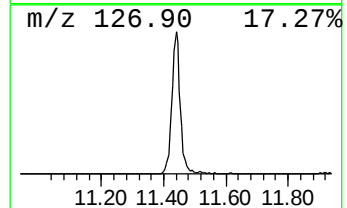
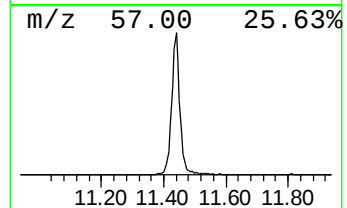
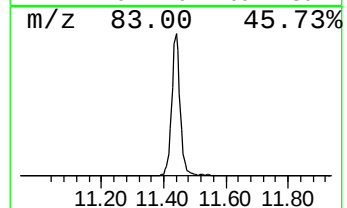
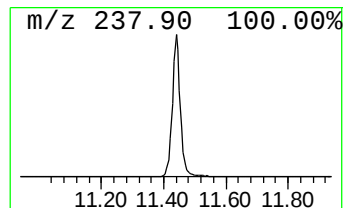
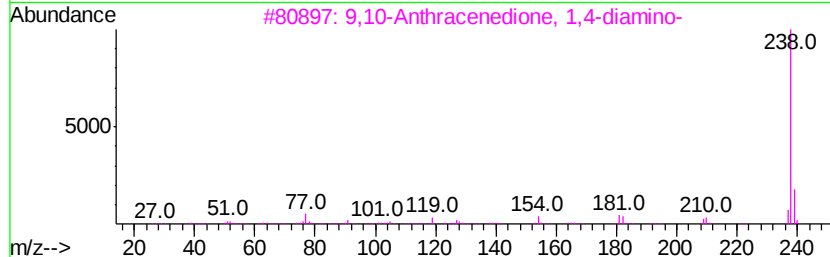
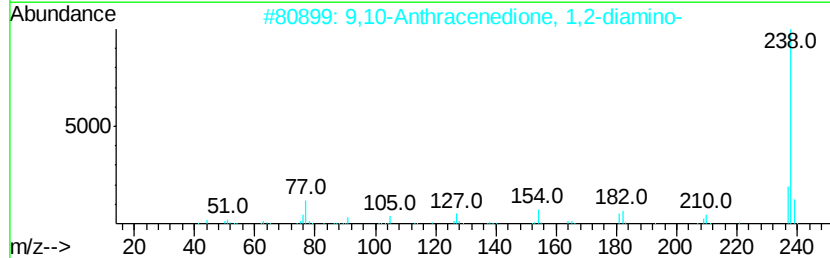
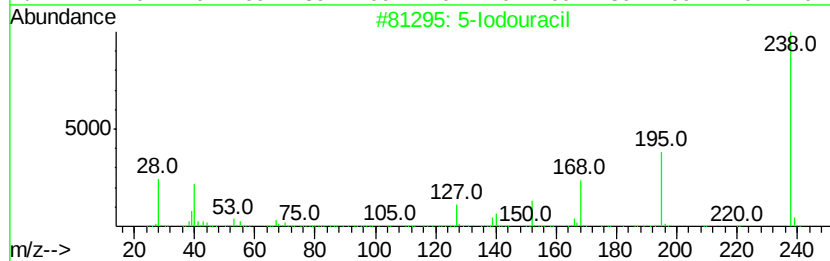
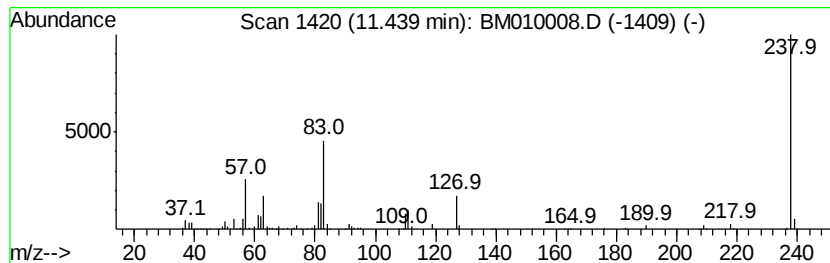
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 6 unknown11.44 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.44	49.46 ng	1617630	Naphthalene-d8	10.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Iodouracil	238	C4H3IN2O2	000696-07-1	32
2		9,10-Anthracenedione, 1,2-diamino-	238	C14H10N2O2	001758-68-5	9
3		9,10-Anthracenedione, 1,4-diamino-	238	C14H10N2O2	000128-95-0	9
4		6-Hydroxyflavone	238	C15H10O3	006665-83-4	5
5		Imidazo[4,5-e][1,4]diazepine-5,8...	238	C10H14N4O3	077246-18-5	5



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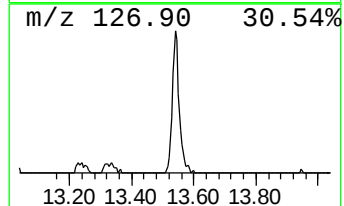
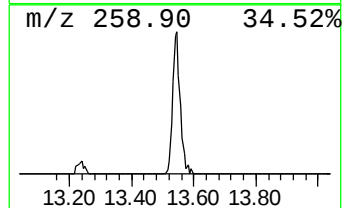
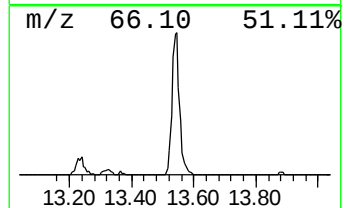
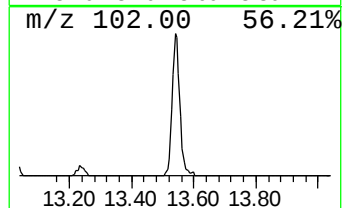
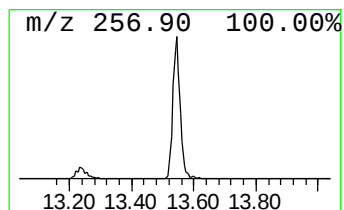
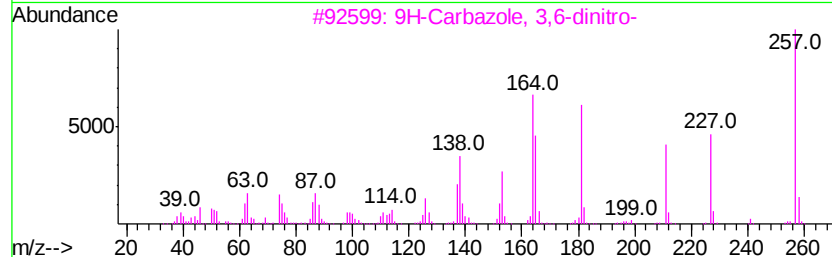
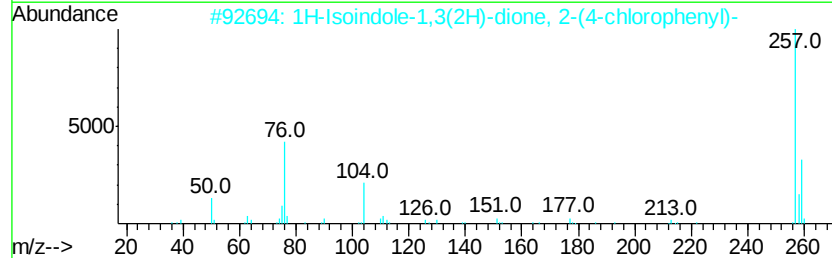
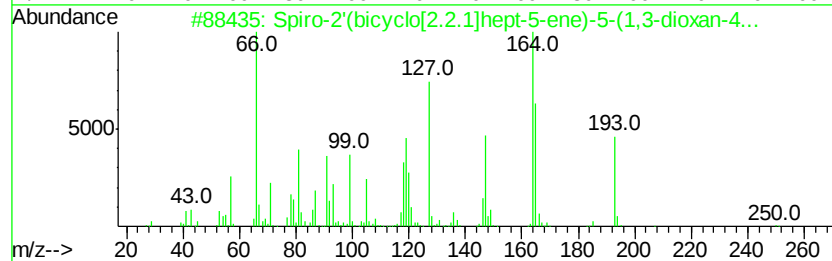
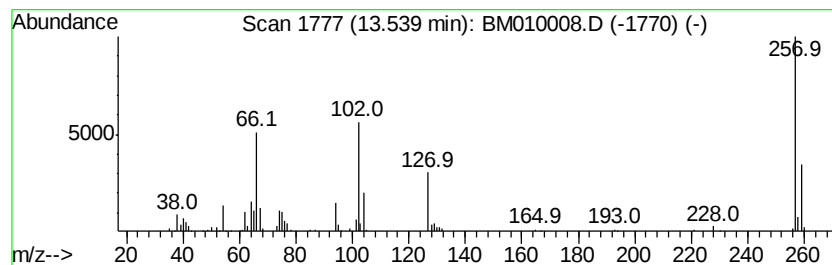
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM051117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown13.54 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.54	11.10 ng	407149	Acenaphthene-d10	14.42
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Spiro-2'(bicyclo[2.2.1]hept-5-en...	250 C15H22O3	1000188-54-9 25
2		1H-Isoindole-1,3(2H)-dione, 2-(4...	257 C14H8ClN02	007386-21-2 14
3		9H-Carbazole, 3,6-dinitro-	257 C12H7N3O4	003244-54-0 14
4		1-(2-Cyanovinyl)aziridine-2-carb...	119 C6H5N3	1000191-13-8 12
5		Acetic acid, 2-phenoxy-, (1,3,3a...	327 C18H17NO5	1000265-61-8 12



Data Path : Z:\HPCHEM1\BNA M\DATA\BM051117\
Data File : BM010008.D
Acq On : 12 May 2017 05:18
Operator : SJ/MA
Sample : I3079-05
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
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ClientSampleId :
DUP-0399-170509

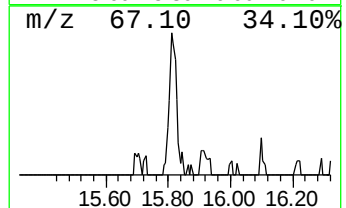
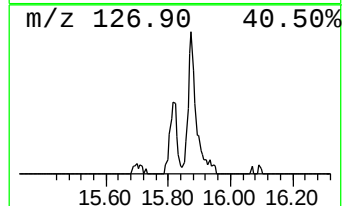
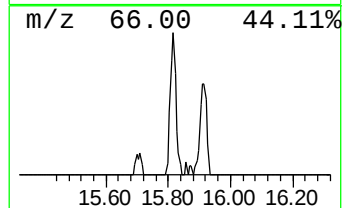
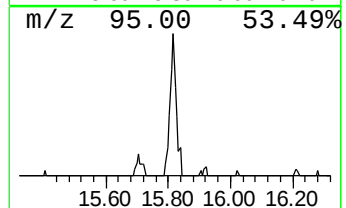
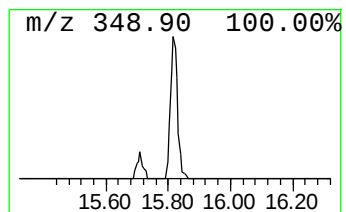
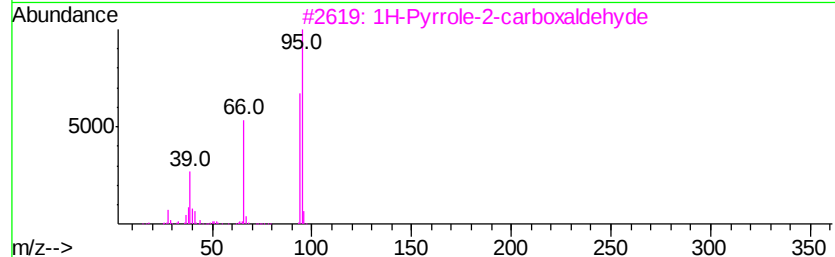
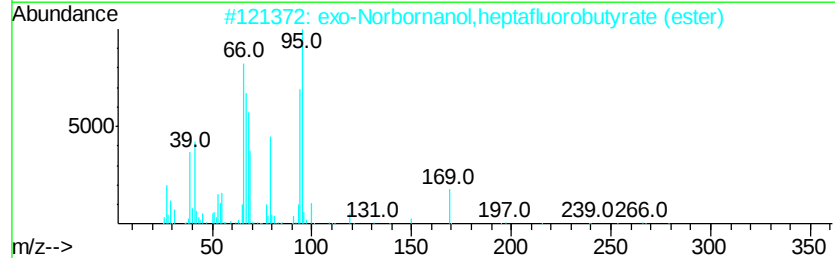
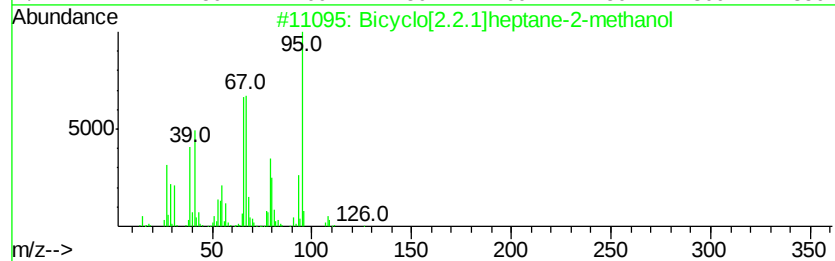
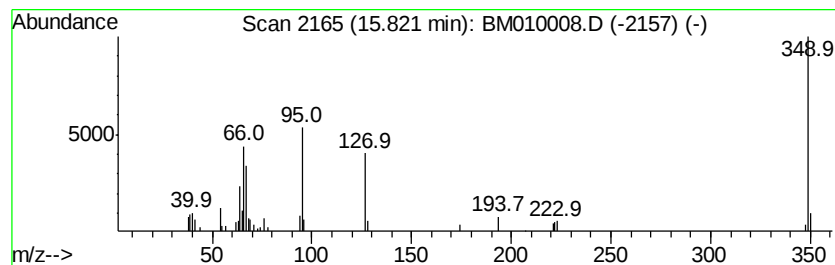
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown15.82 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.82	2.21 ng	97784	Phenanthrene-d10	17.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bicyclo[2.2.1]heptane-2-methanol	126	C8H14O	005240-72-2	23
2		exo-Norbornanol,heptafluorobutyr...	308	C11H11F7O2	1000245-88-2	23
3		1H-Pyrrole-2-carboxaldehyde	95	C5H5NO	001003-29-8	23
4		N-(2-Ethylhexyl)-5-norbornene-2,...	275	C17H25NO2	000113-48-4	10
5		4H-Thiopyran-3,5-dicarbonitrile,...	344	C16H16N4O3S	312265-39-7	10



Data Path : Z:\HPCHEM1\BNA M\DATA\BM051117\
Data File : BM010008.D
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Misc :
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ClientSampleId :
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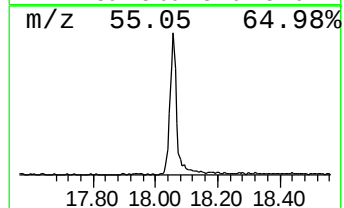
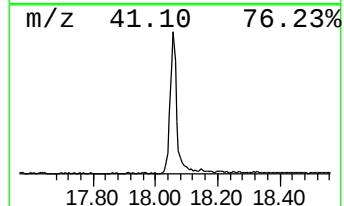
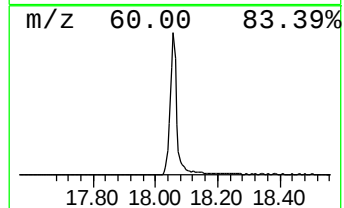
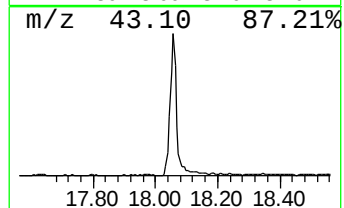
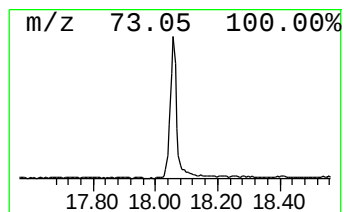
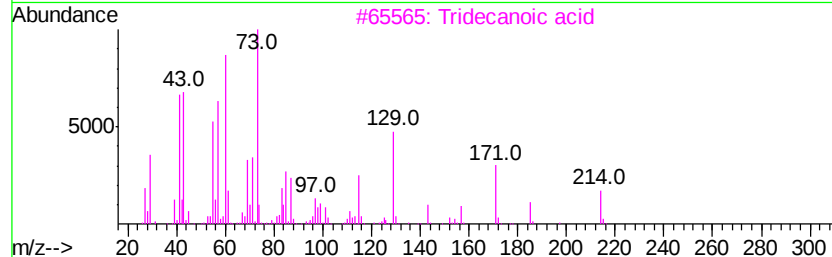
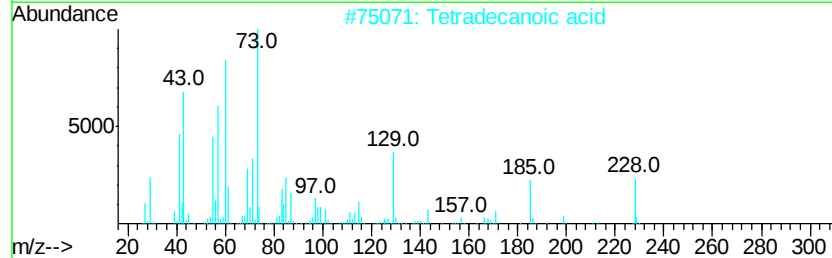
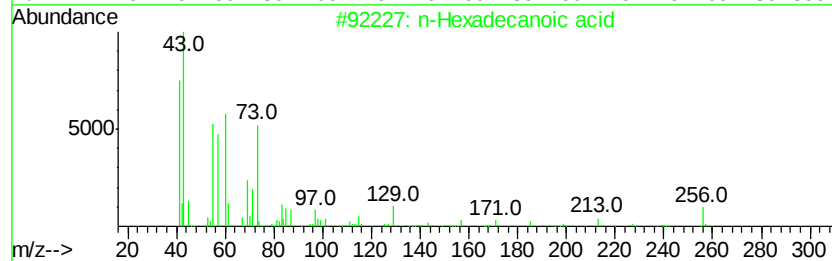
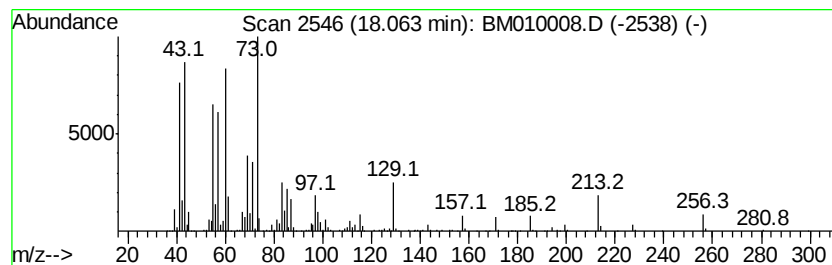
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.06	49.55 ng	2193550	Phenanthrene-d10	17.17

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	94
3		Tridecanoic acid	214	C13H26O2	000638-53-9	90
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	87
5		Undecanoic acid	186	C11H22O2	000112-37-8	80



Data Path : Z:\HPCHEM1\BNA M\DATA\BM051117\
Data File : BM010008.D
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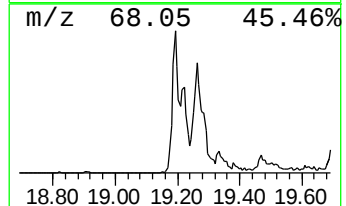
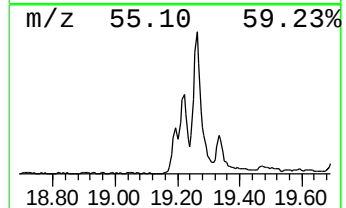
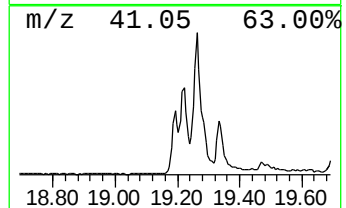
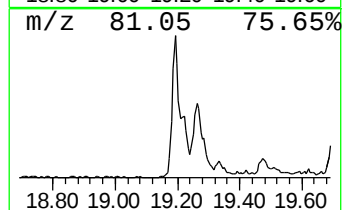
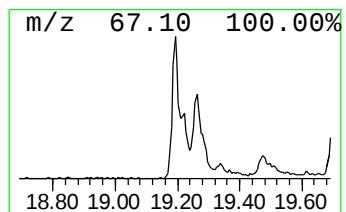
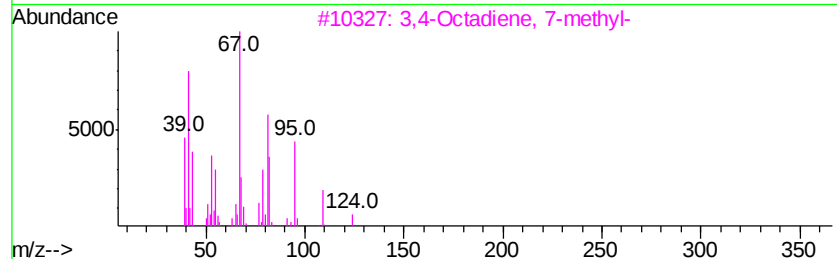
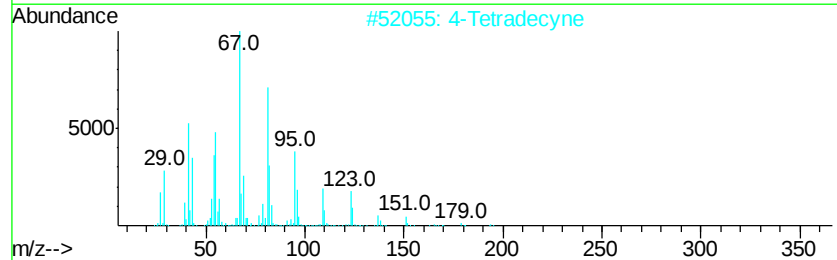
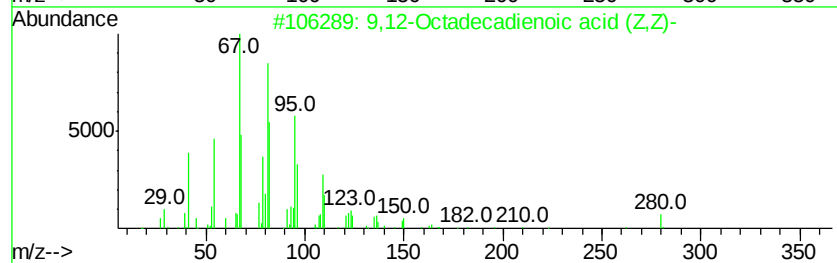
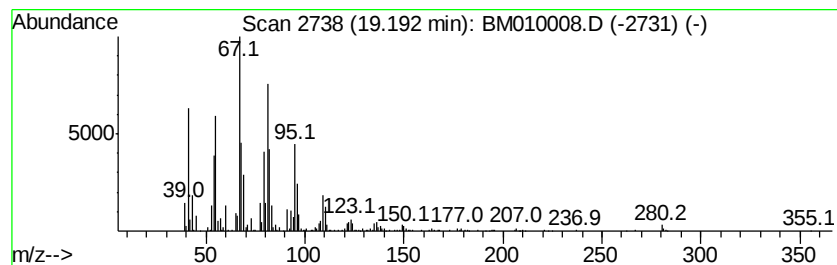
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 9,12-Octadecadienoic acid (... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.19	27.92 ng	1235920	Phenanthrene-d10	17.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9,12-Octadecadienoic acid (Z,Z)-	280	C18H32O2	000060-33-3	98
2		4-Tetradecyne	194	C14H26	060212-33-1	93
3		3,4-Octadiene, 7-methyl-	124	C9H16	037050-05-8	76
4		9-Octadecynoic acid, methyl ester	294	C19H34O2	001120-32-7	72
5		Bicyclo[2.2.2]octane, 2-methyl-	124	C9H16	000766-53-0	72



Data Path : Z:\HPCHEM1\BNA M\DATA\BM051117\
Data File : BM010008.D
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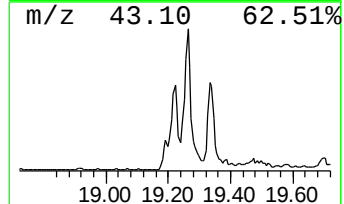
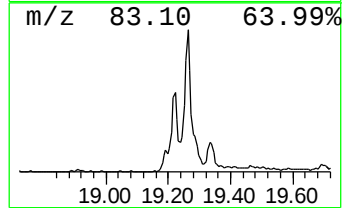
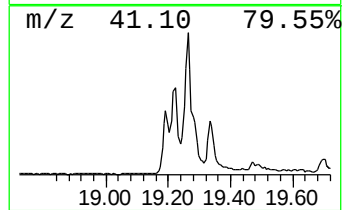
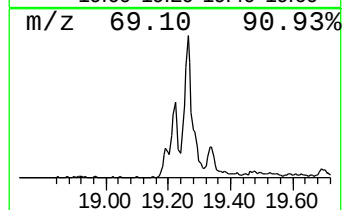
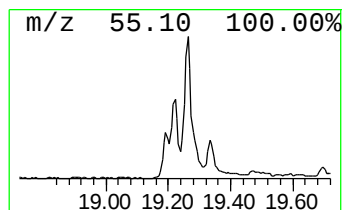
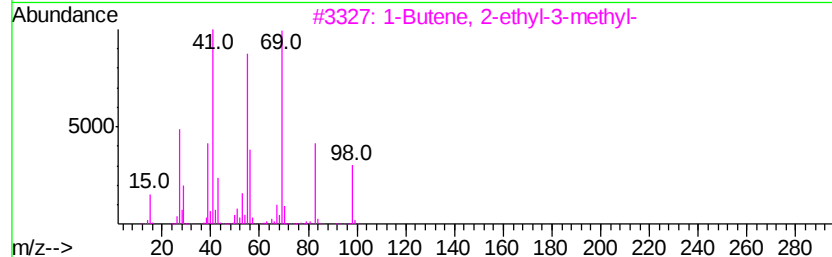
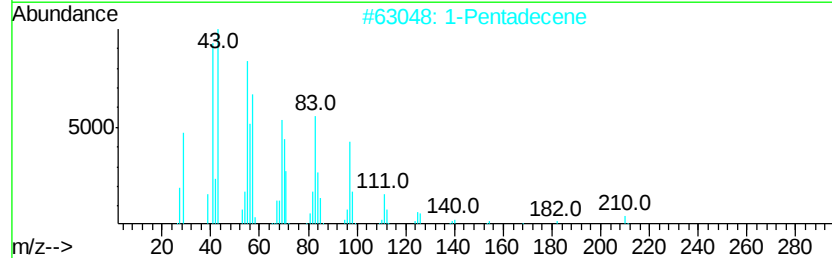
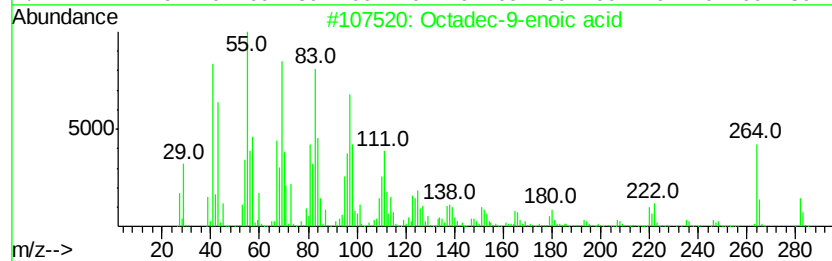
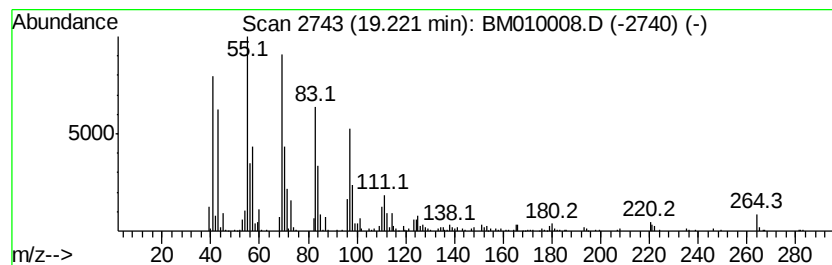
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Octadec-9-enoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.22	35.63 ng	1577290	Phenanthrene-d10	17.17

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadec-9-enoic acid	282	C18H34O2	1000190-13-7	86
2	1-Pentadecene	210	C15H30	013360-61-7	47
3	1-Butene, 2-ethyl-3-methyl-	98	C7H14	007357-93-9	38
4	1-Hexene, 3,3-dimethyl-	112	C8H16	003404-77-1	38
5	Cyclohexadecane	224	C16H32	000295-65-8	38



Data Path : Z:\HPCHEM1\BNA M\DATA\BM051117\
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Acq On : 12 May 2017 05:18
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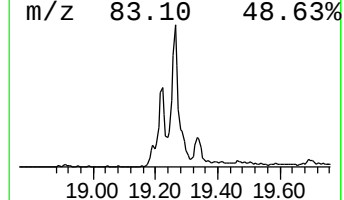
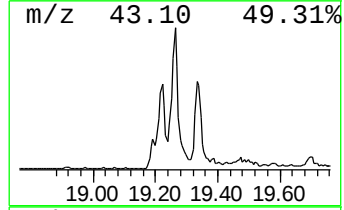
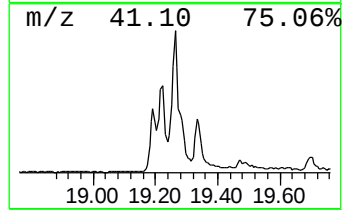
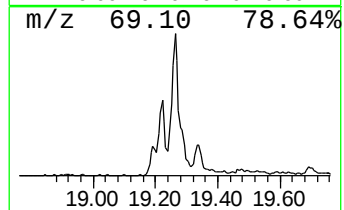
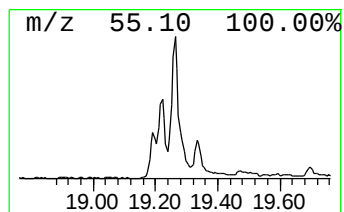
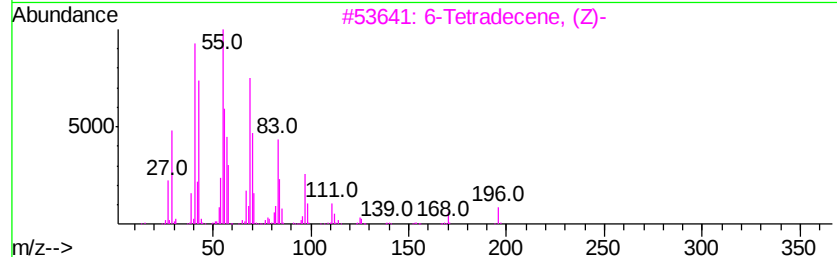
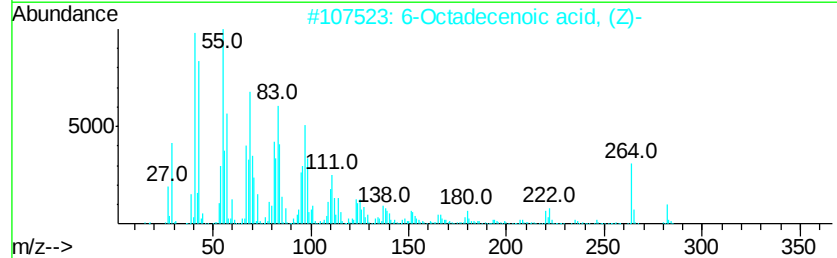
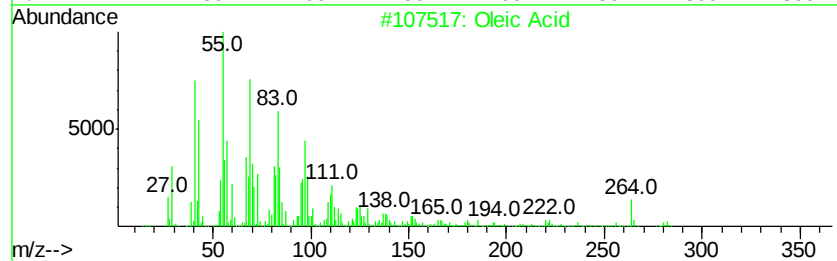
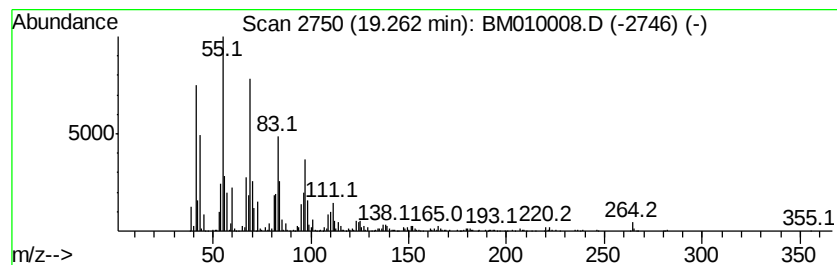
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Oleic Acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.26	70.07 ng	3101850	Phenanthrene-d10	17.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Oleic Acid	282	C18H34O2	000112-80-1	94
2		6-Octadecenoic acid, (Z)-	282	C18H34O2	000593-39-5	83
3		6-Tetradecene, (Z)-	196	C14H28	041446-61-1	64
4		2,6-Octadienal, 3,7-dimethyl-, (E)-	152	C10H16O	000141-27-5	41
5		2,6-Octadienal, 3,7-dimethyl-	152	C10H16O	005392-40-5	38



Data Path : Z:\HPCHEM1\BNA M\DATA\BM051117\
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ALS Vial : 23 Sample Multiplier: 1

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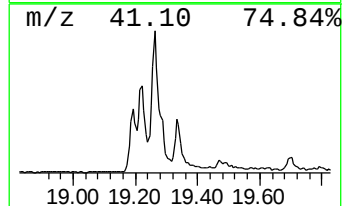
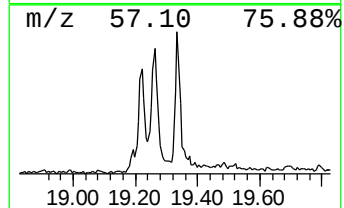
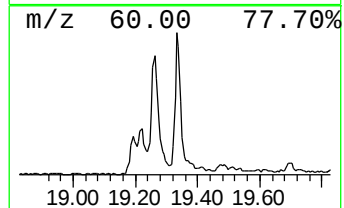
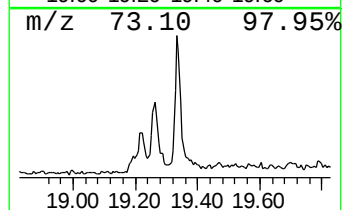
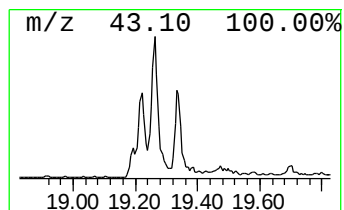
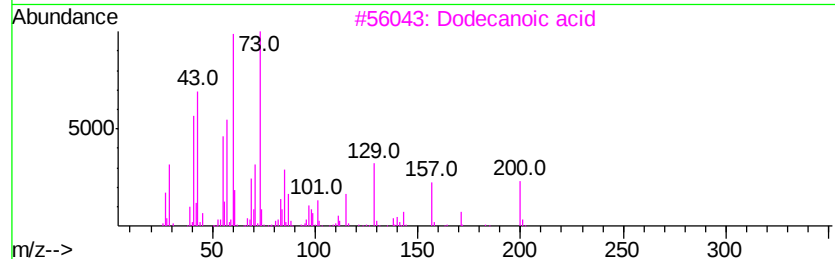
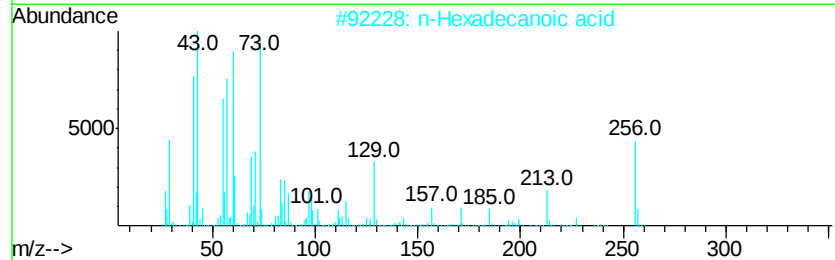
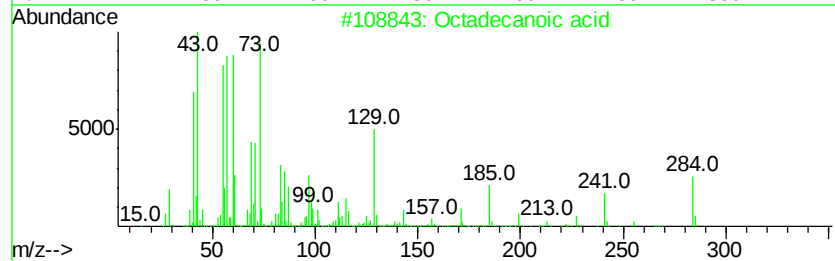
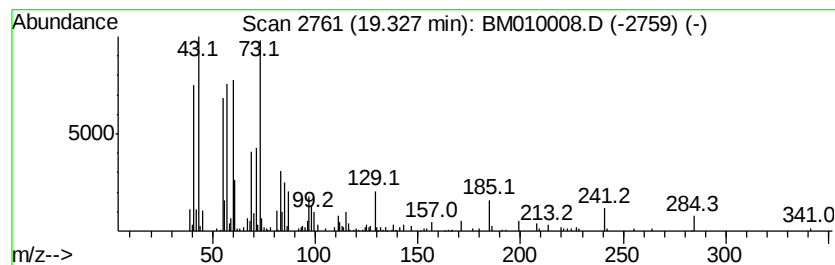
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Octadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.33	18.41 ng	979046	Chrysene-d12	21.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	72
3		Dodecanoic acid	200	C12H24O2	000143-07-7	64
4		Tridecanoic acid	214	C13H26O2	000638-53-9	59
5		Tetradecanoic acid	228	C14H28O2	000544-63-8	58



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Sample : I3079-05
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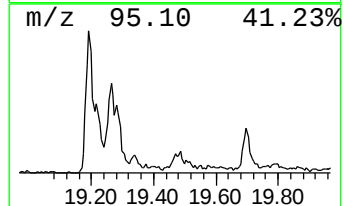
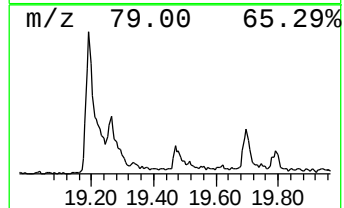
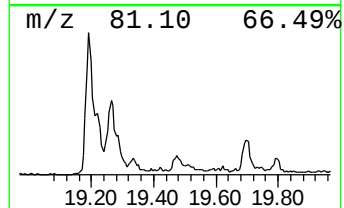
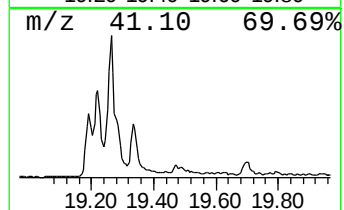
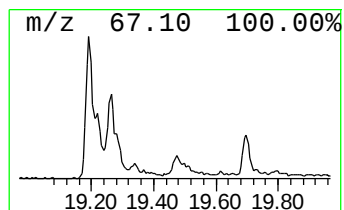
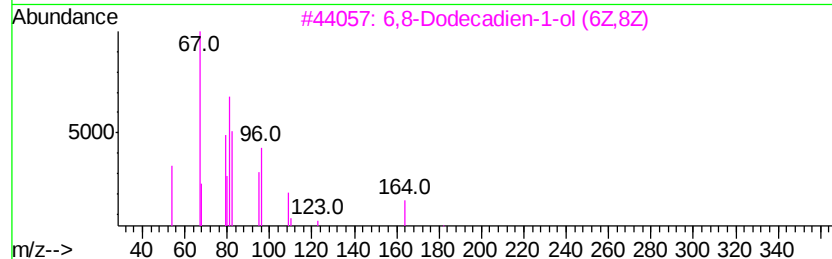
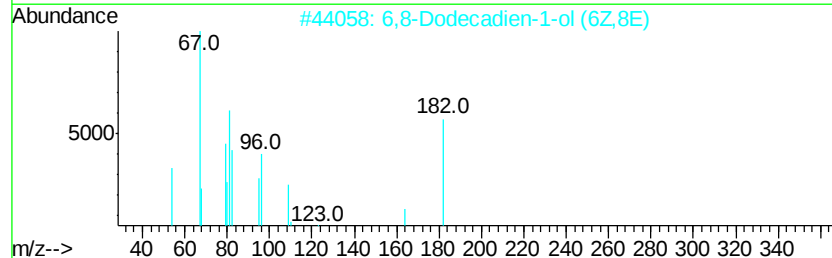
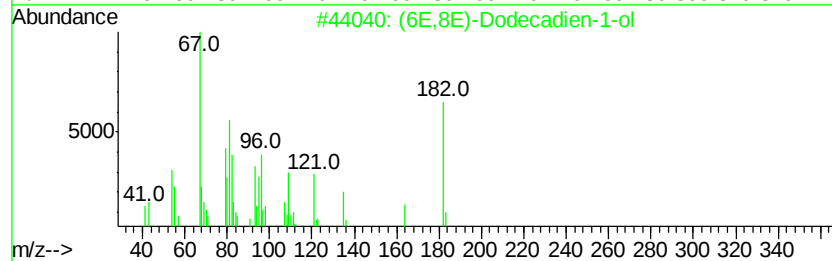
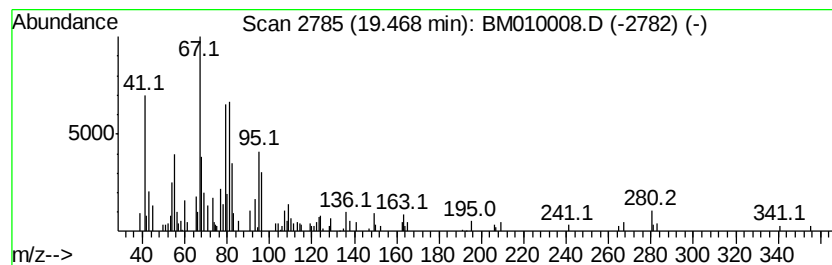
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 (6E,8E)-Dodecadien-1-ol Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.		
19.47	4.70 ng	250005	Chrysene-d12	21.36		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		(6E,8E)-Dodecadien-1-ol	182	C12H22O	072648-89-6	72
2		6,8-Dodecadien-1-ol (6Z,8E)	182	C12H22O	1000222-02-7	53
3		6,8-Dodecadien-1-ol (6Z,8Z)	182	C12H22O	1000222-02-6	49
4		4,5-Nonadiene	124	C9H16	000821-74-9	47
5		3,4-Octadiene, 7-methyl-	124	C9H16	037050-05-8	38



Data Path : Z:\HPCHEM1\BNA M\DATA\BM051117\
Data File : BM010008.D
Acq On : 12 May 2017 05:18
Operator : SJ/MA
Sample : I3079-05
Misc :
ALS Vial : 23 Sample Multiplier: 1

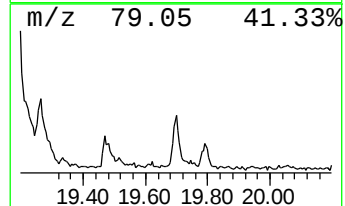
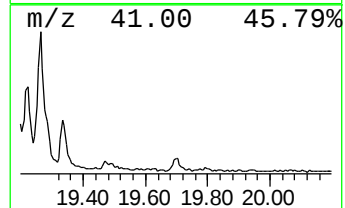
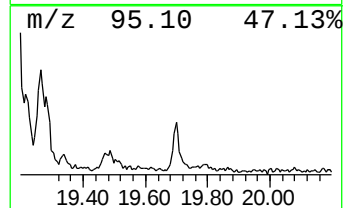
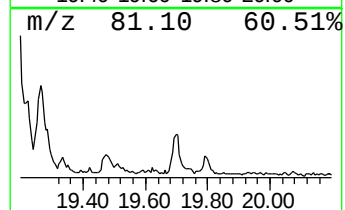
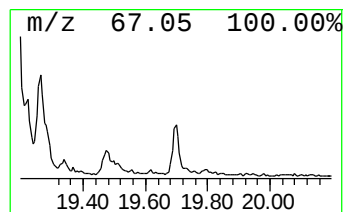
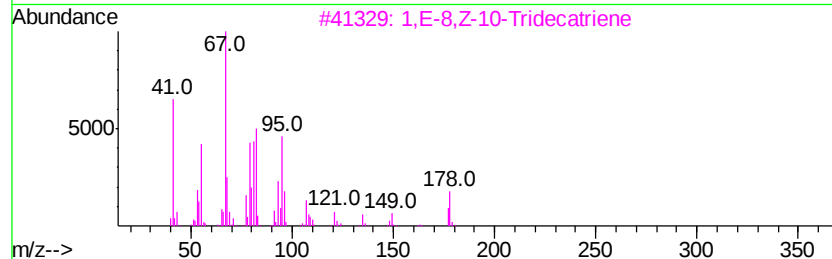
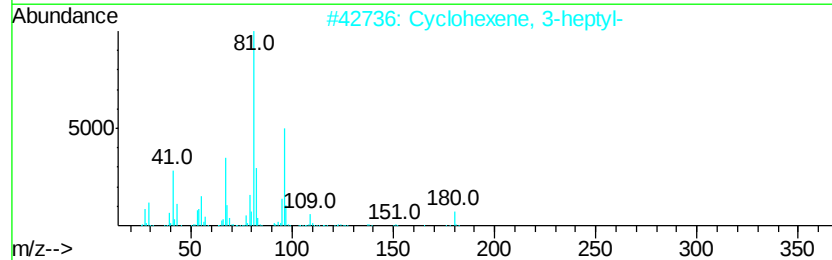
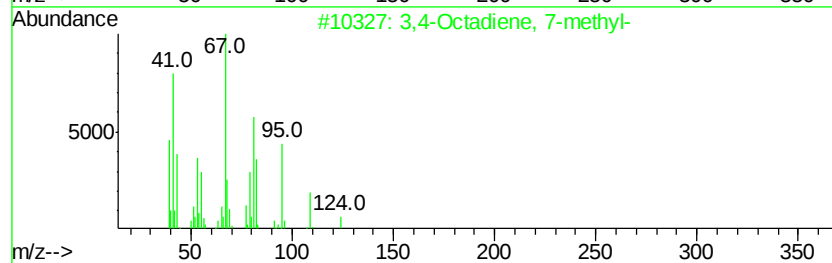
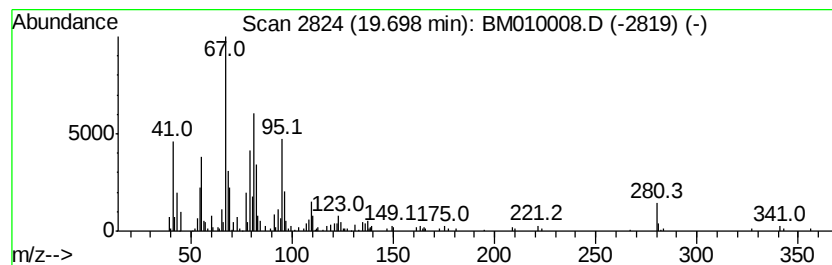
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 3,4-Octadiene, 7-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.70	7.12 ng	378743	Chrysene-d12	21.36	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,4-Octadiene, 7-methyl-	124	C9H16	037050-05-8	72
2	Cyclohexene, 3-heptyl-	180	C13H24	015232-93-6	59
3	1,E-8,Z-10-Tridecatriene	178	C13H22	080625-30-5	53
4	(Z,Z)-3,6-Nonadienal	138	C9H14O	021944-83-2	49
5	7-Hexadecyne	222	C16H30	074685-28-2	43



Data Path : Z:\HPCHEM1\BNA M\DATA\BM051117\
Data File : BM010008.D
Acq On : 12 May 2017 05:18
Operator : SJ/MA
Sample : I3079-05
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DUP-0399-170509

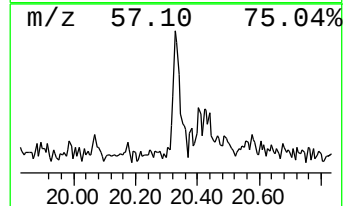
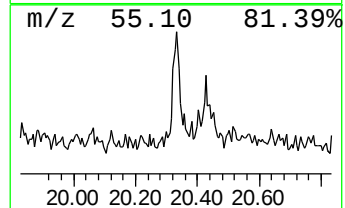
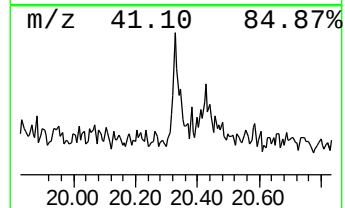
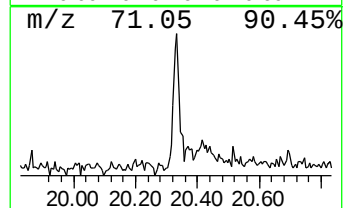
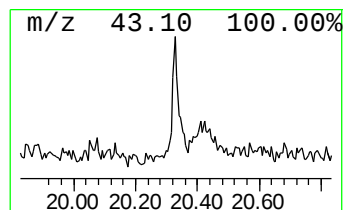
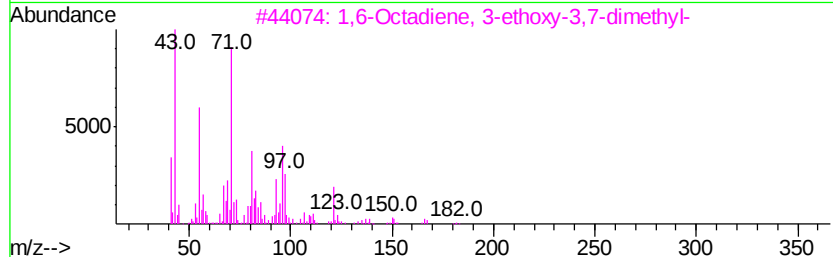
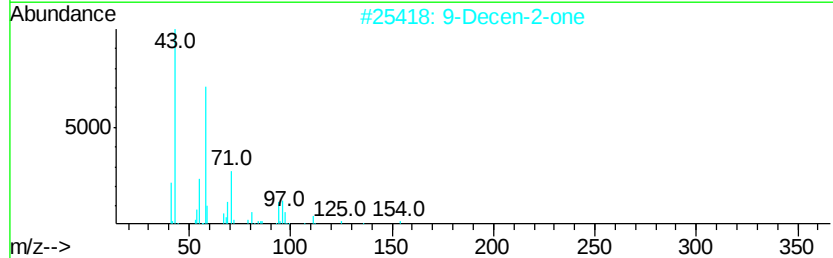
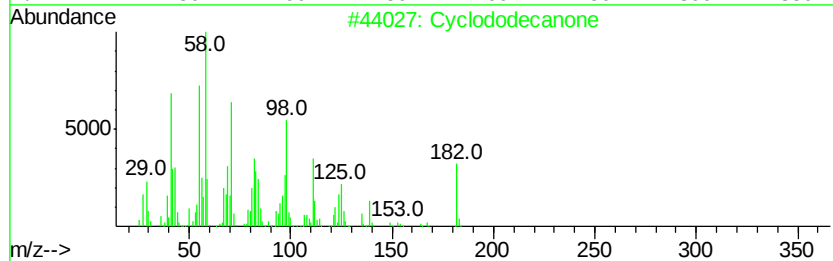
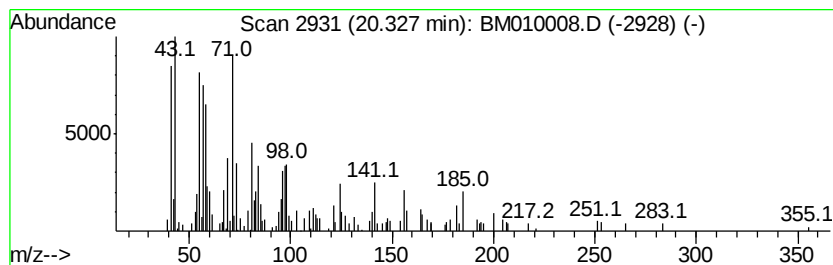
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 unknown20.33 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.33	3.11 ng	165640	Chrysene-d12	21.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclododecanone	182	C12H22O	000830-13-7	47
2		9-Decen-2-one	154	C10H18O	035194-30-0	38
3		1,6-Octadiene, 3-ethoxy-3,7-dime...	182	C12H22O	072845-33-1	35
4		4-Bromobenzoic acid, 2-tetrahydr...	284	C12H13BrO3	1000282-24-9	35
5		9-Heptadecanone	254	C17H34O	000540-08-9	35



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 Acq On : 12 May 2017 05:18
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 Sample : I3079-05
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-BM051117.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
2-Pentanone, 4-hy...	4.89	14.8	ng	314113	1	7.76	423488	20.0
unknown5.47	5.47	151.6	ng	3209140	1	7.76	423488	20.0
2-Pentanone, 4-me...	5.96	2.6	ng	54513	1	7.76	423488	20.0
unknown7.29	7.29	85.8	ng	1815660	1	7.76	423488	20.0
unknown11.44	11.44	49.5	ng	1617630	2	10.56	654098	20.0
unknown13.54	13.54	11.1	ng	407149	3	14.42	733777	20.0
unknown15.82	15.82	2.2	ng	97784	4	17.17	885316	20.0
n-Hexadecanoic acid	18.06	49.5	ng	2193550	4	17.17	885316	20.0
9,12-Octadecadien...	19.19	27.9	ng	1235920	4	17.17	885316	20.0
Octadec-9-enoic acid	19.22	35.6	ng	1577290	4	17.17	885316	20.0
Oleic Acid	19.26	70.1	ng	3101850	4	17.17	885316	20.0
Octadecanoic acid	19.33	18.4	ng	979046	5	21.36	1063620	20.0
(6E,8E)-Dodecadie...	19.47	4.7	ng	250005	5	21.36	1063620	20.0
3,4-Octadiene, 7-...	19.70	7.1	ng	378743	5	21.36	1063620	20.0
unknown20.33	20.33	3.1	ng	165640	5	21.36	1063620	20.0