

Data Path : Z:\HPCHEM1\BNA F\DATA\BF052217\
 Data File : BF095325.D
 Acq On : 22 May 2017 21:06
 Operator : SJ/MA
 Sample : I3261-01
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NCC-BUCK-3-1

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.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	5.119	537	541	557	rBV	72257	186960	9.01%	0.986%
2	5.725	640	644	667	rBV	742187	1709571	82.37%	9.017%
3	7.301	908	912	928	rBV	841195	1659649	79.97%	8.754%
4	7.425	928	933	951	rVB	1233376	2075432	100.00%	10.947%
5	7.760	986	990	1002	rVB	369768	462878	22.30%	2.442%
6	7.995	1025	1030	1042	rBV	1168453	1468620	70.76%	7.746%
7	8.660	1138	1143	1174	rBV	576649	1069144	51.51%	5.639%
8	9.777	1329	1333	1362	rBV	350291	628170	30.27%	3.313%
9	11.548	1629	1634	1656	rBV	1304821	2010298	96.86%	10.604%
10	12.254	1749	1754	1764	rBV	113378	194159	9.36%	1.024%
11	12.607	1810	1814	1826	rBV2	413837	689314	33.21%	3.636%
12	13.442	1949	1956	1960	rBV	29342	33270	1.60%	0.175%
13	13.906	2030	2035	2047	rBV	477136	854162	41.16%	4.505%
14	14.212	2084	2087	2090	rBV2	25947	31509	1.52%	0.166%
15	14.418	2118	2122	2129	rBV3	17800	24934	1.20%	0.132%
16	15.012	2217	2223	2233	rBV	299767	482122	23.23%	2.543%
17	15.201	2251	2255	2258	rBV4	16026	20829	1.00%	0.110%
18	15.236	2258	2261	2268	rVB3	20098	25860	1.25%	0.136%
19	15.965	2381	2385	2395	rBV2	73440	132157	6.37%	0.697%
20	16.042	2395	2398	2405	rVB7	18387	27825	1.34%	0.147%
21	17.053	2566	2570	2574	rBV5	20778	33685	1.62%	0.178%
22	17.324	2612	2616	2628	rBV	976038	1218368	58.70%	6.426%
23	18.571	2824	2828	2833	rBV4	41063	62475	3.01%	0.330%
24	18.683	2843	2847	2861	rBV	378536	564243	27.19%	2.976%
25	19.359	2958	2962	2965	rBV	49034	63913	3.08%	0.337%
26	19.800	3034	3037	3040	rVB	40767	41543	2.00%	0.219%
27	19.971	3055	3066	3072	rBV	82083	314747	15.17%	1.660%
28	20.083	3081	3085	3089	rBV	261199	286322	13.80%	1.510%
29	20.306	3120	3123	3128	rVB5	35886	67351	3.25%	0.355%
30	20.359	3128	3132	3141	rBV	307273	473258	22.80%	2.496%
31	20.700	3186	3190	3199	rVB	72021	126915	6.12%	0.669%
32	20.812	3205	3209	3215	rVB	237236	304190	14.66%	1.604%
33	21.024	3242	3245	3253	rVB2	80719	131429	6.33%	0.693%
34	22.153	3429	3437	3447	rVB5	191575	554275	26.71%	2.924%

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35	22.571	3504	3508	3519	rVB6	135666	373462	17.99%	1.970%
36	22.782	3538	3544	3547	rBV6	75801	162849	7.85%	0.859%
37	23.841	3716	3724	3736	rVB6	135999	392844	18.93%	2.072%

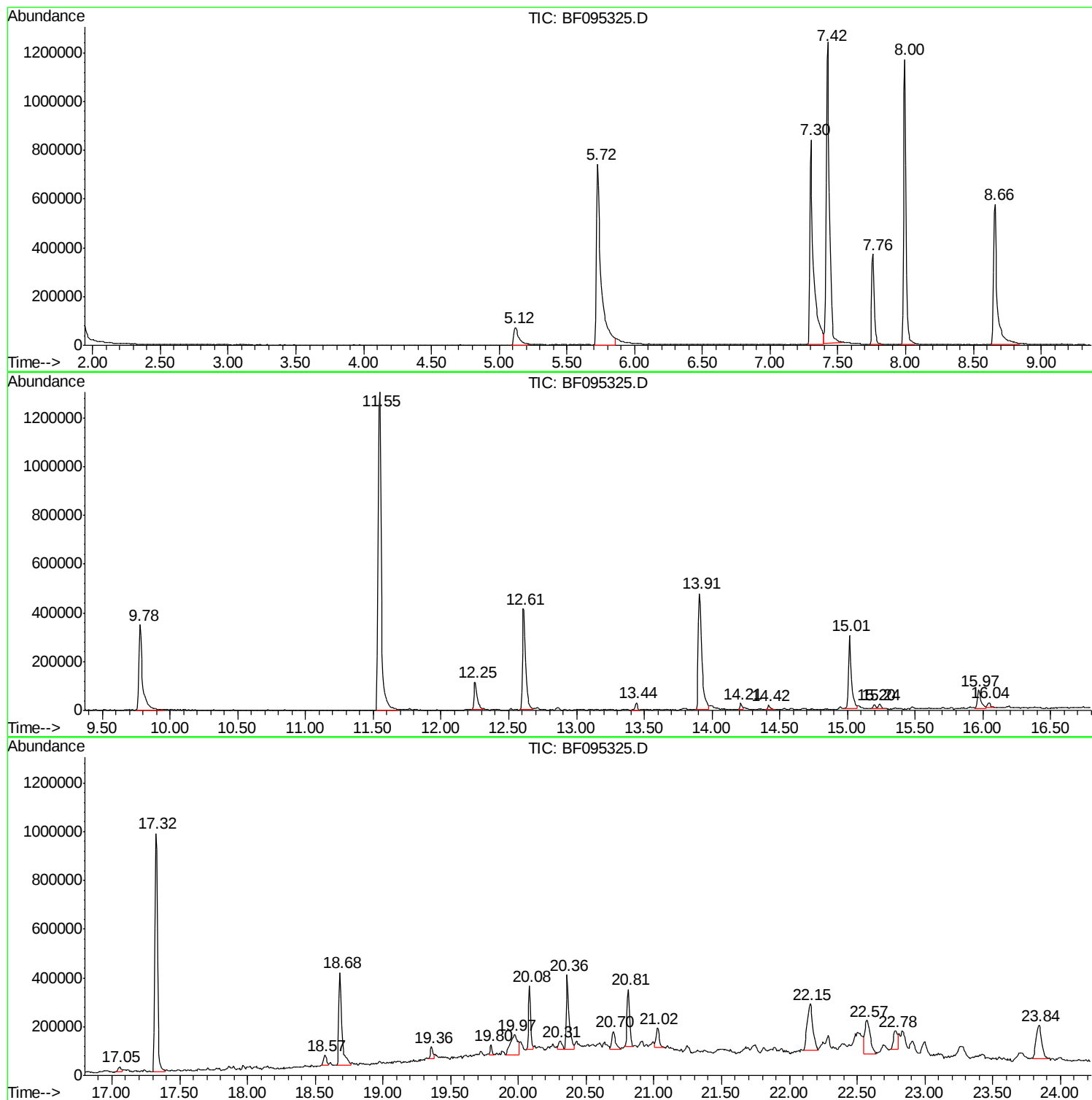
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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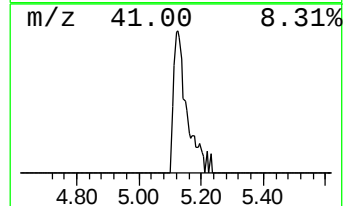
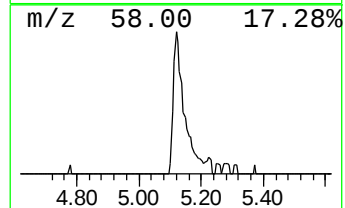
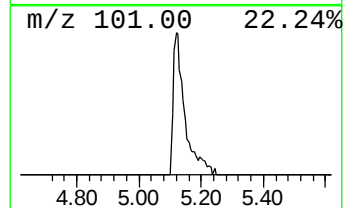
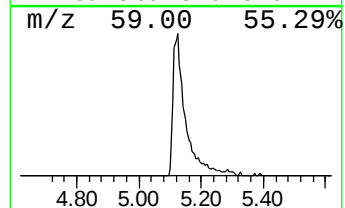
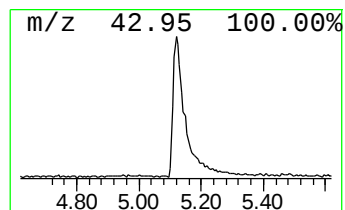
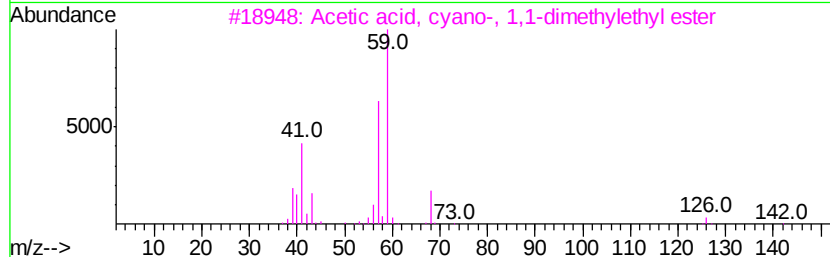
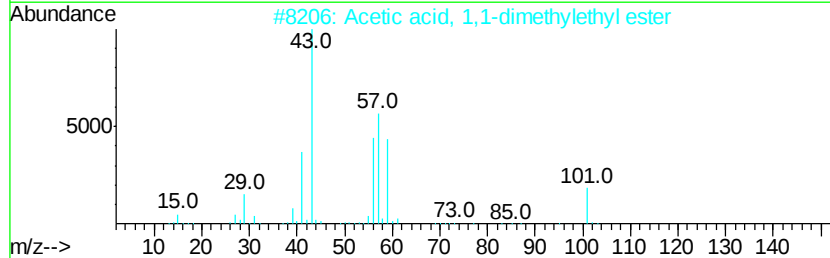
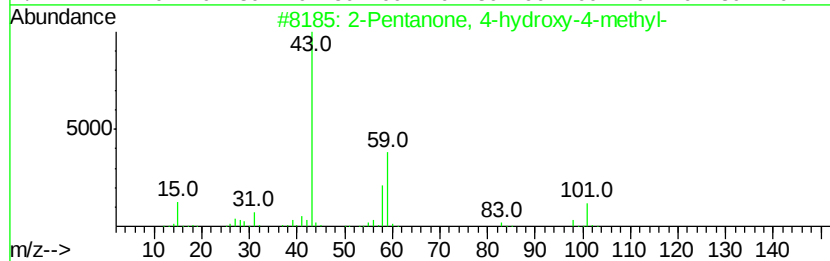
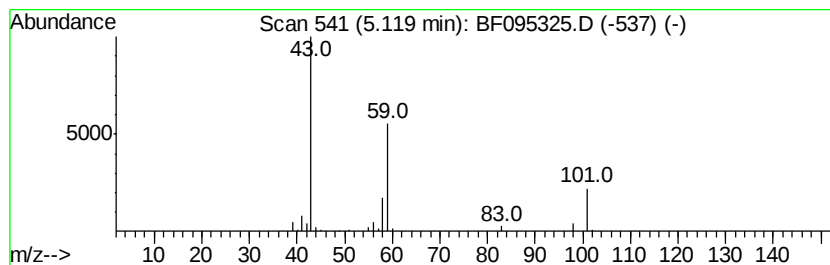
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.12	8.08 ng	186960	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	50
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
4			5-Hexen-2-one	98	C6H10O	000109-49-9	9
5			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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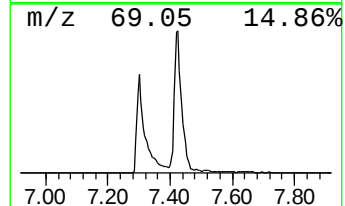
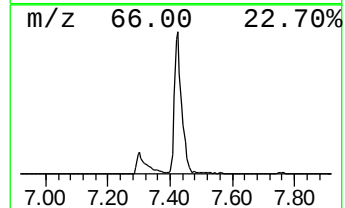
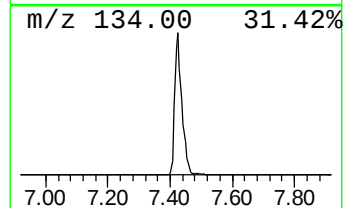
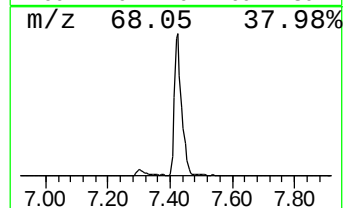
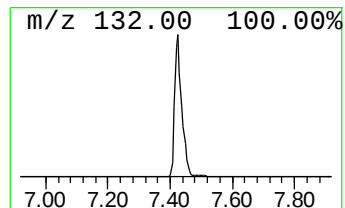
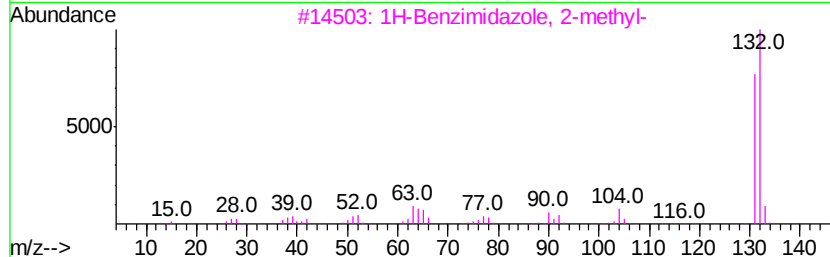
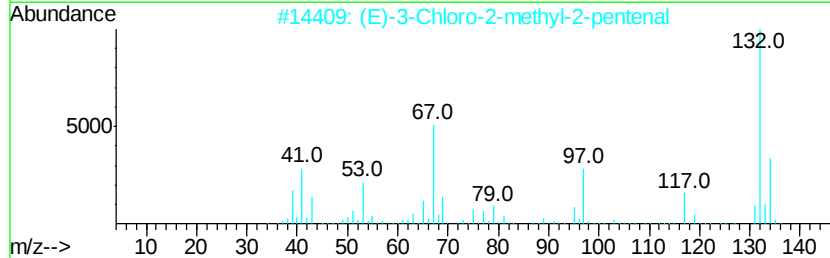
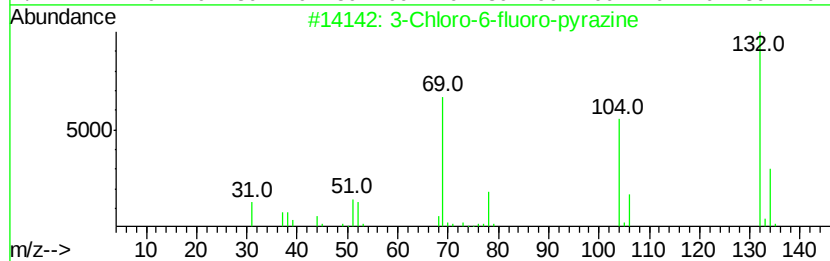
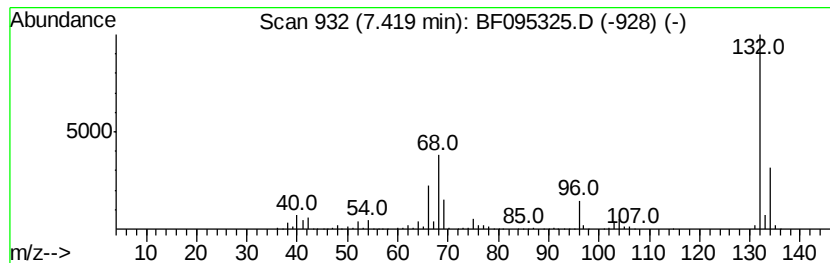
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Peak Number 2 unknown7.42 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.42	89.68 ng	2075430	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	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	35
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	10
4		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
5		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9



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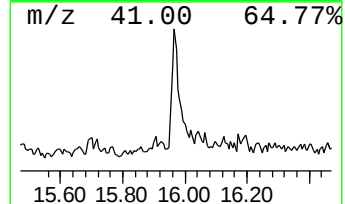
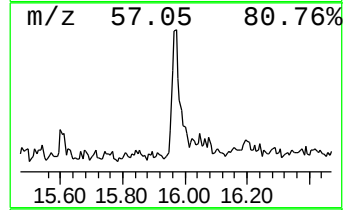
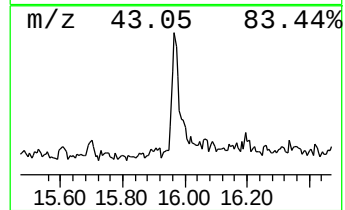
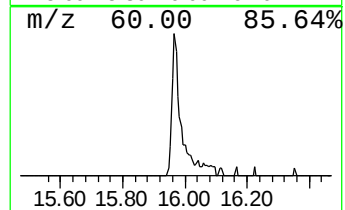
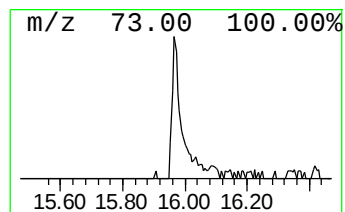
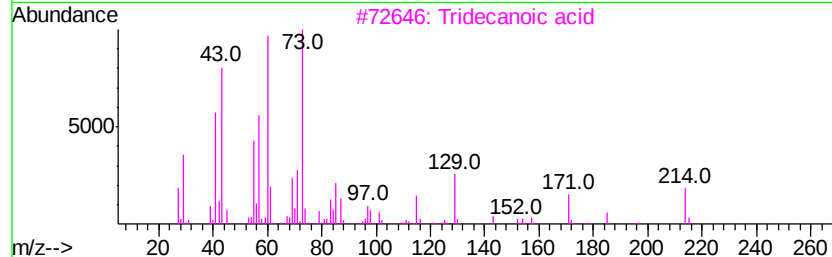
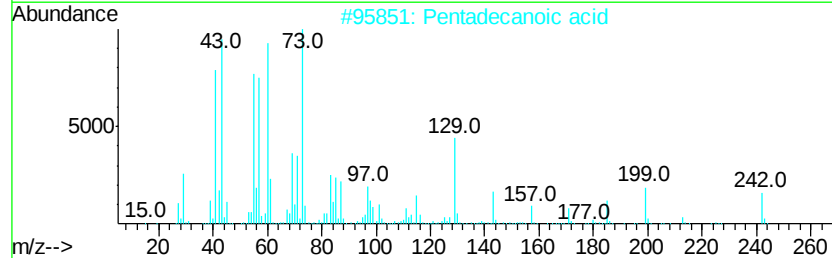
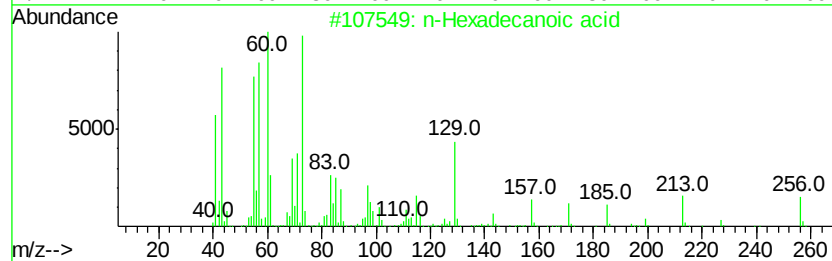
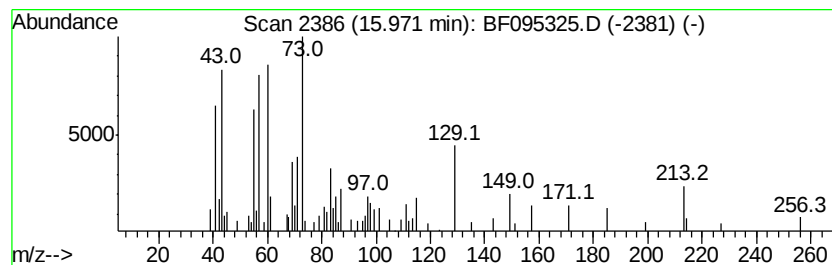
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 n-Hexadecanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.97	5.48 ng	132157	Phenanthrene-d10	15.01

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	87
3	Tridecanoic acid	214	C13H26O2	000638-53-9	80
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	80
5	Dodecanoic acid	200	C12H24O2	000143-07-7	53



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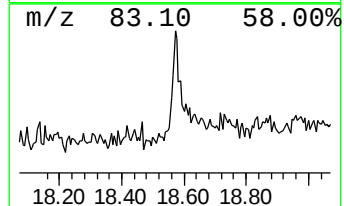
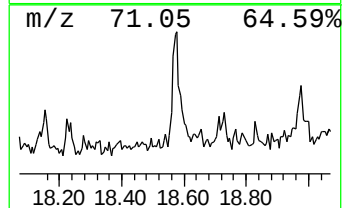
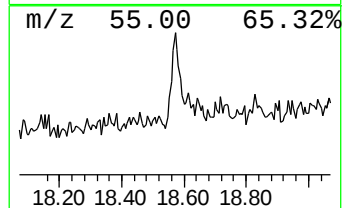
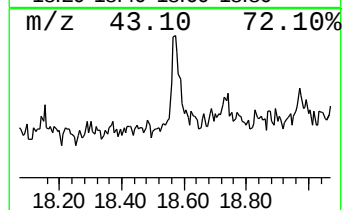
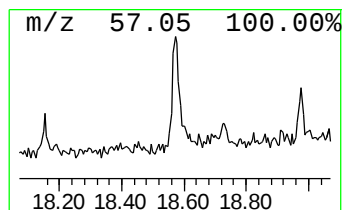
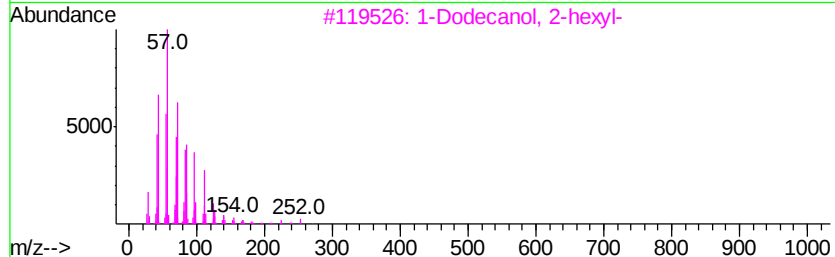
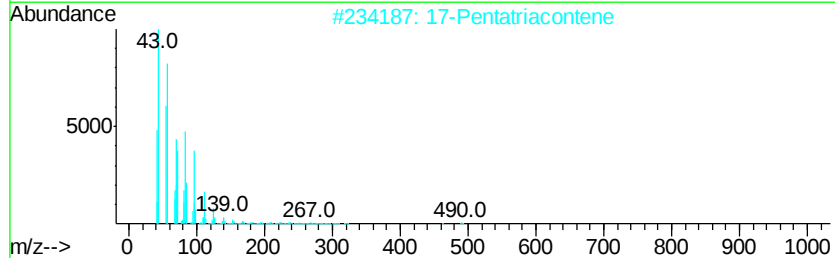
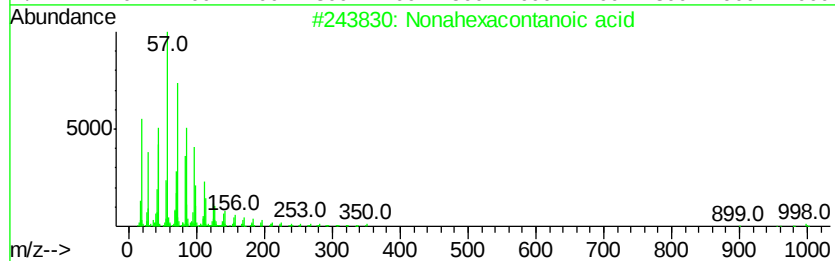
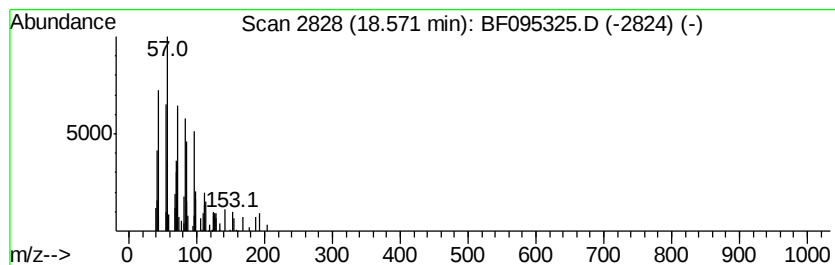
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 Nonaheptacontanoic acid Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.57	2.21 ng	62475	Chrysene-d12	18.68

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonaheptacontanoic acid	999	C69H138O2	040710-32-5	58
2		17-Pentatriacontene	491	C35H70	006971-40-0	52
3		1-Dodecanol, 2-hexyl-	270	C18H38O	110225-00-8	49
4		1-Decanol, 2-octyl-	270	C18H38O	045235-48-1	49
5		1-Decanol, 2-hexyl-	242	C16H34O	002425-77-6	49



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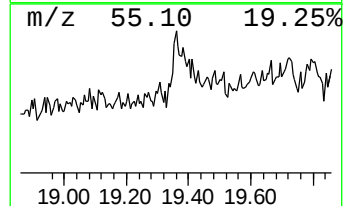
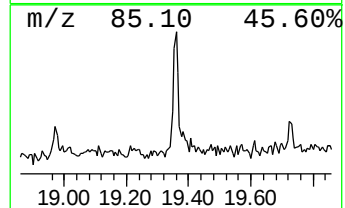
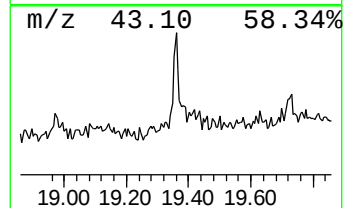
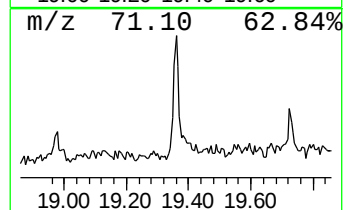
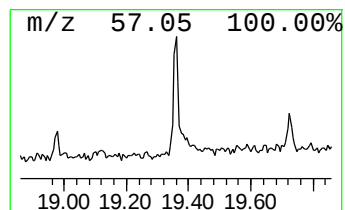
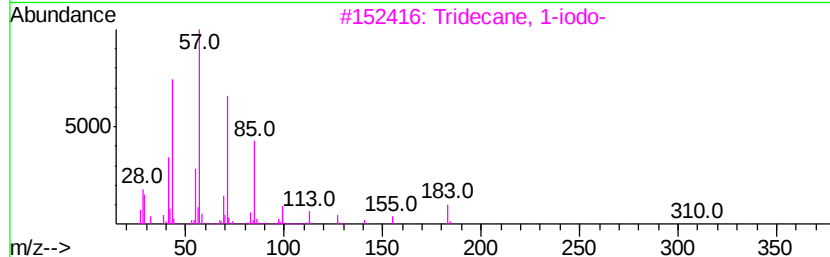
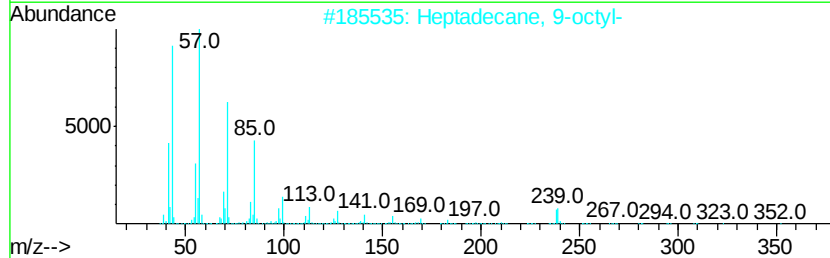
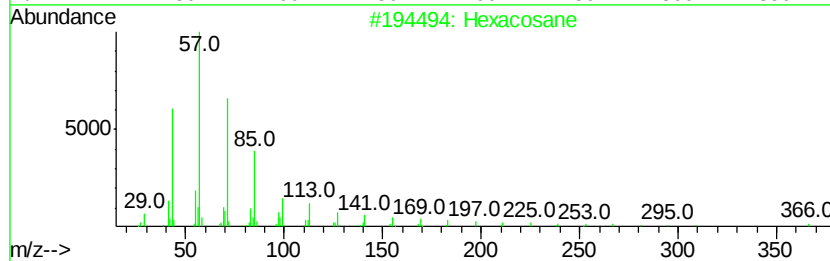
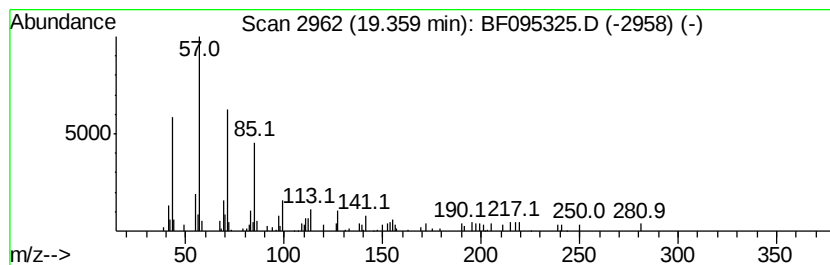
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Peak Number 6 Hexacosane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.36	2.27 ng	63913	Chrysene-d12	18.68

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexacosane		366	C26H54	000630-01-3	87
2	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	83
3	Tridecane, 1-iodo-		310	C13H27I	035599-77-0	80
4	Tetracosane		338	C24H50	000646-31-1	80
5	Octadecane, 2-methyl-		268	C19H40	001560-88-9	80



Data Path : Z:\HPCHEM1\BNA F\DATA\BF052217\
Data File : BF095325.D
Acq On : 22 May 2017 21:06
Operator : SJ/MA
Sample : I3261-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

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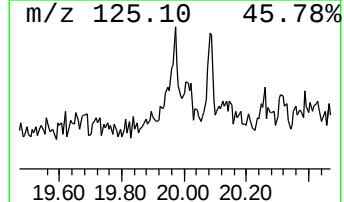
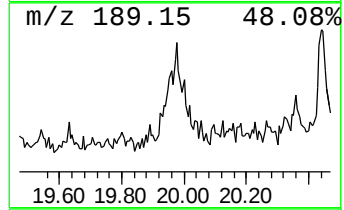
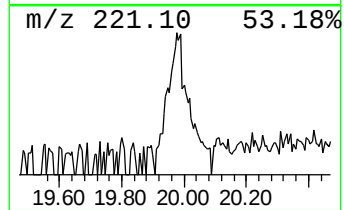
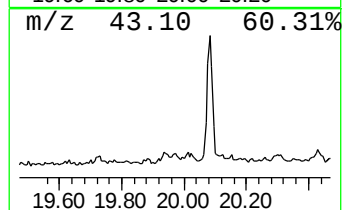
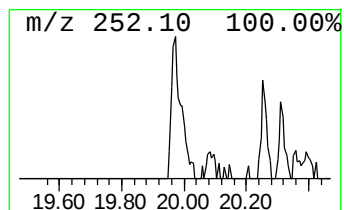
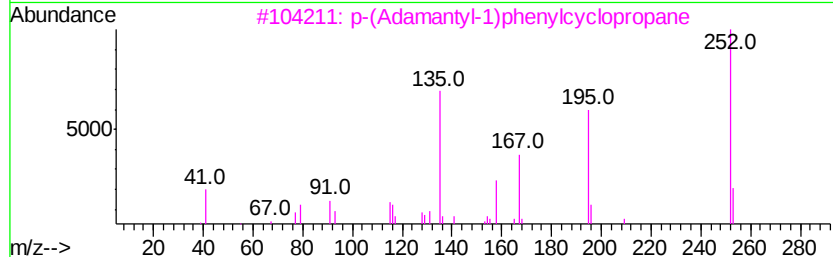
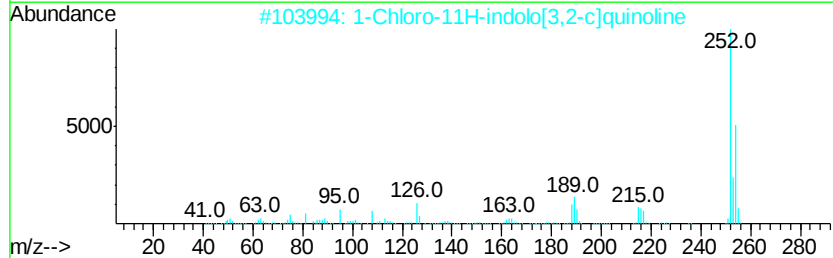
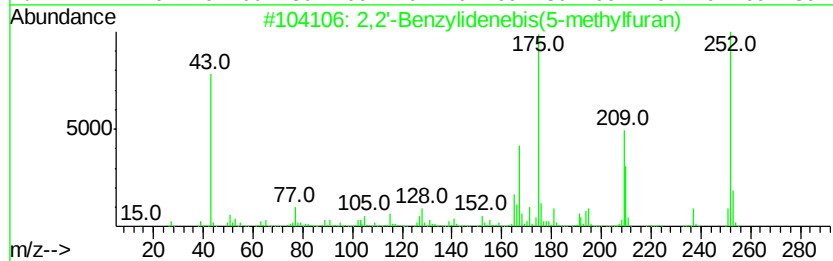
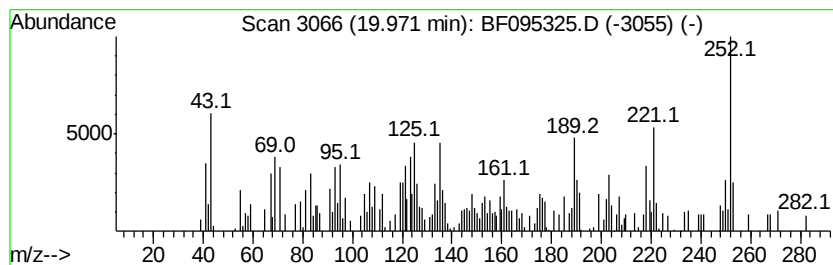
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown19.97 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.97	13.30 ng	314747	Perylene-d12	20.36

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2'	-Benzylidenebis(5-methylfuran)	252	C17H16O2	086694-47-5	35	
2	1-Chloro-11H-indolo[3,2-c]quinoline	252	C15H9ClN2	155249-51-7	30		
3	p-(Adamantyl-1)phenylcyclopropane	252	C19H24	1000140-46-1	30		
4	Benzo[e]pyrene	252	C20H12	000192-97-2	27		
5	Benz[e]acephenanthrylene	252	C20H12	000205-99-2	22		



Data Path : Z:\HPCHEM1\BNA F\DATA\BF052217\
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Misc :
ALS Vial : 19 Sample Multiplier: 1

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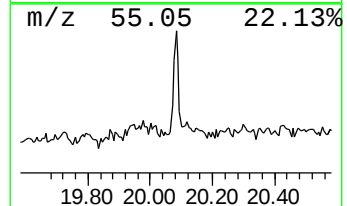
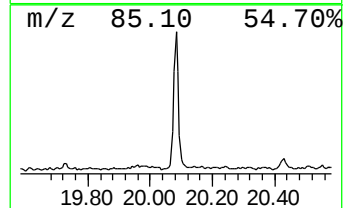
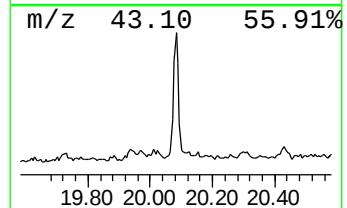
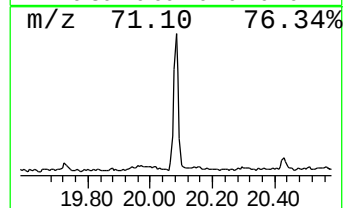
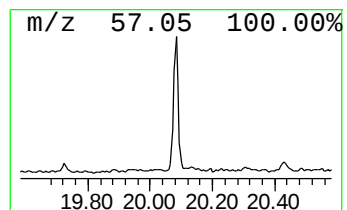
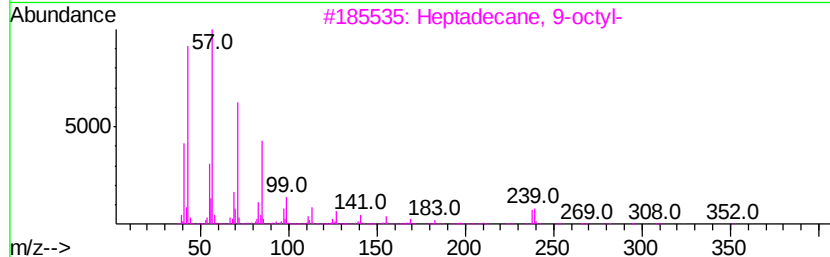
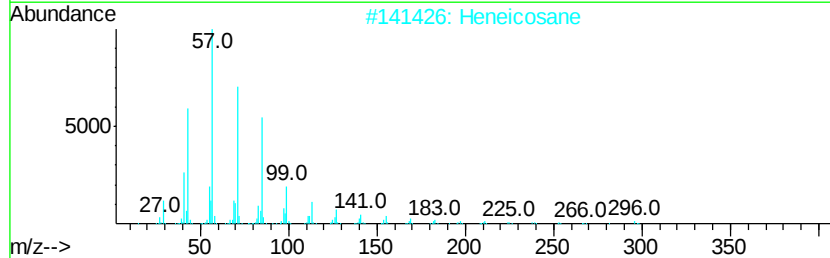
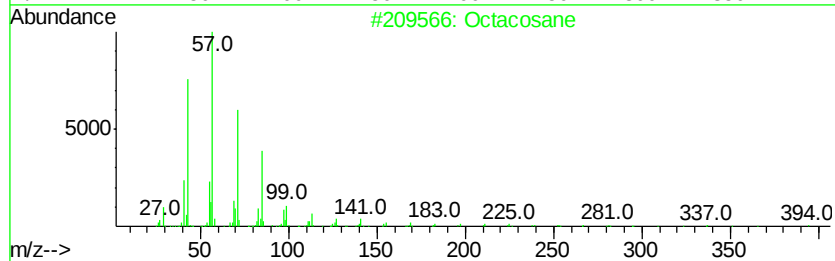
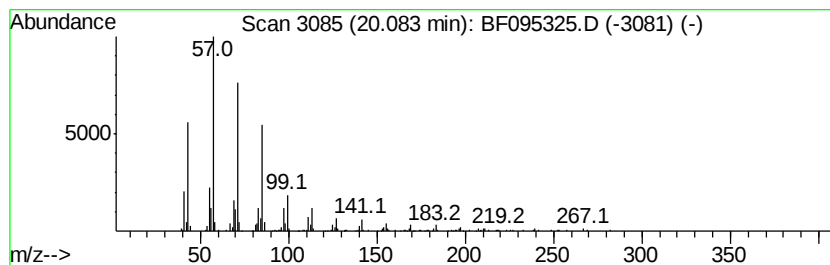
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Octacosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.08	12.10 ng	286322	Perylene-d12	20.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	91
2		Heneicosane	296	C21H44	000629-94-7	91
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4		2-methyloctacosane	408	C29H60	1000376-72-8	91
5		triacontane	422	C30H62	000638-68-6	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052217\
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Acq On : 22 May 2017 21:06
Operator : SJ/MA
Sample : I3261-01
Misc :
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Instrument :
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ClientSampleId :
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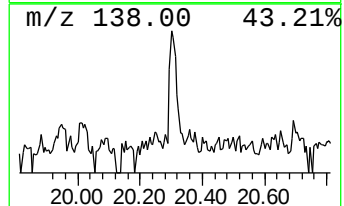
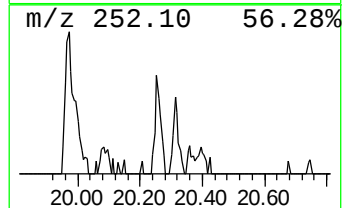
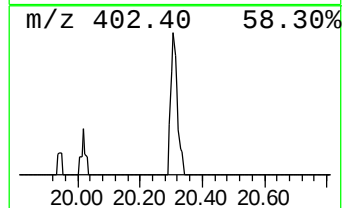
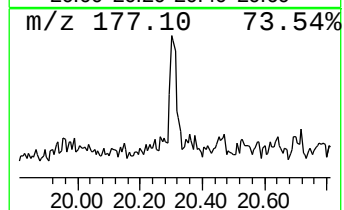
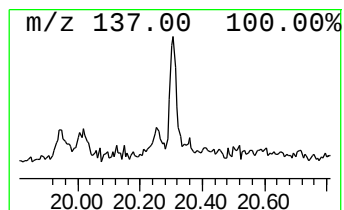
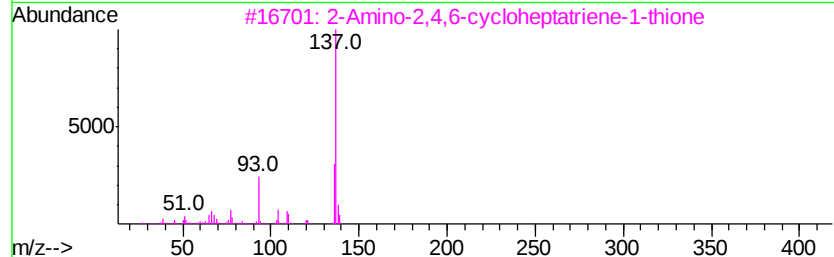
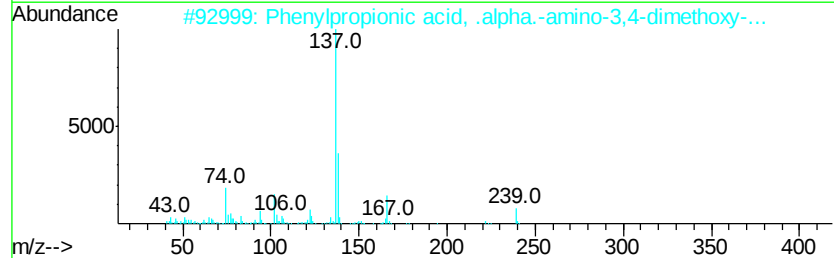
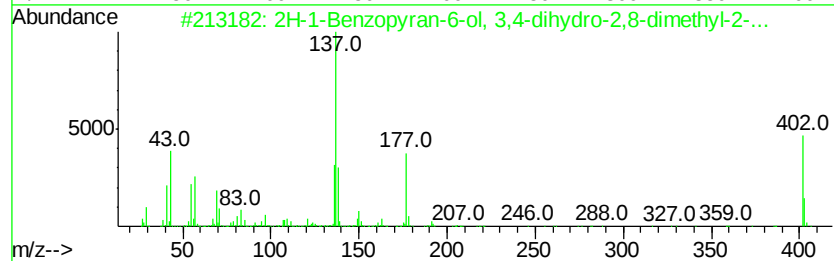
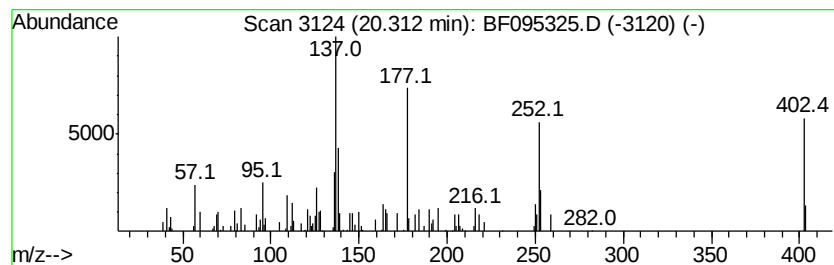
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 2H-1-Benzopyran-6-ol, 3,4-d... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.31	2.85 ng	67351	Perylene-d12	20.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2H-1-Benzopyran-6-ol, 3,4-dihydr...	402	C27H46O2	000119-13-1	94
2		Phenylpropionic acid, .alpha.-am...	239	C12H17NO4	067409-05-6	27
3		2-Amino-2,4,6-cycloheptatriene-1...	137	C7H7NS	003336-99-0	27
4		6-Amino-2,4-dimethylphenol	137	C8H11NO	041458-65-5	27
5		4-Dimethylaminopyridin-2-amine	137	C7H11N3	050426-31-8	27



Data Path : Z:\HPCHEM1\BNA F\DATA\BF052217\
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Sample : I3261-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

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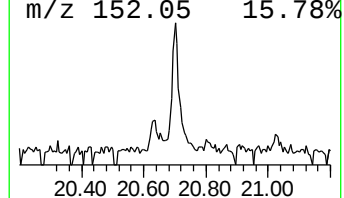
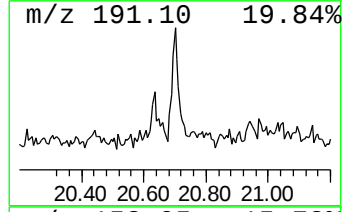
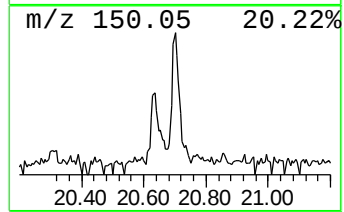
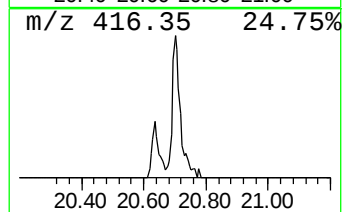
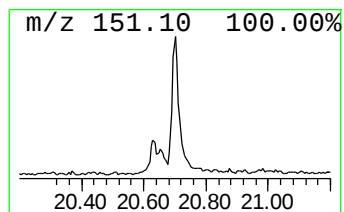
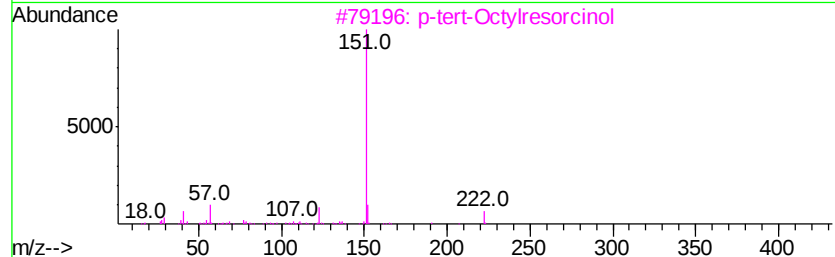
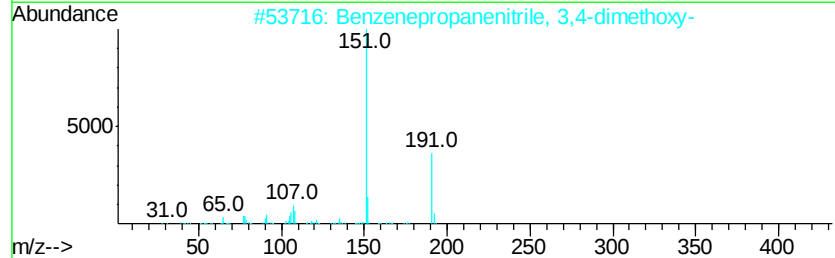
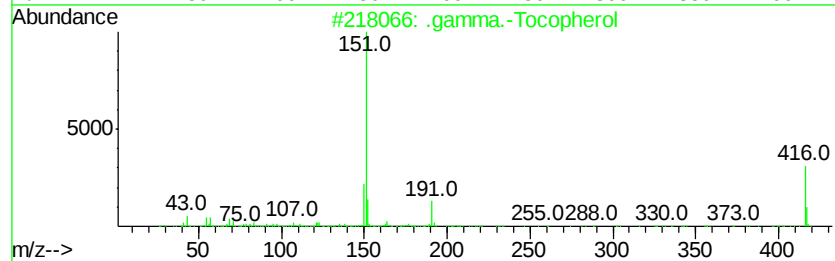
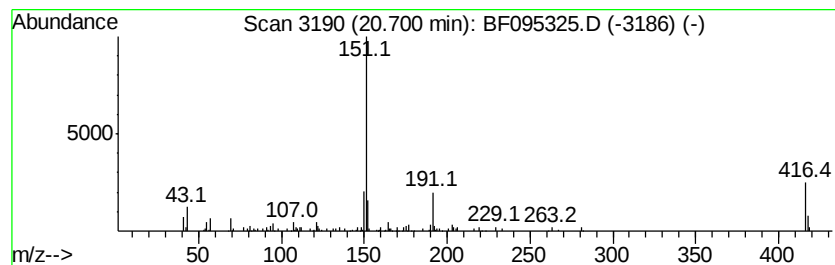
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 .gamma.-Tocopherol Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.70	5.36 ng	126915	Perylene-d12	20.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	.gamma.-Tocopherol	416	C28H48O2	007616-22-0	95
2		Benzenepropanenitrile, 3,4-dimet...	191	C11H13N02	049621-56-9	50
3		p-tert-Octylresorcinol	222	C14H22O2	028122-52-3	43
4		4H-1,3,2-Dioxaborin, 6-ethenyl-2...	208	C12H21B02	074630-05-0	40
5		Benzene, 1-[(2-bromophenoxy)meth...	322	C15H15Br03	1000115-65-9	37



Data Path : Z:\HPCHEM1\BNA F\DATA\BF052217\
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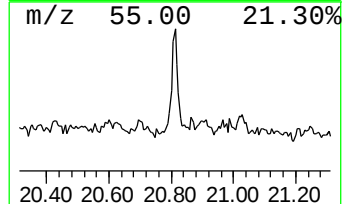
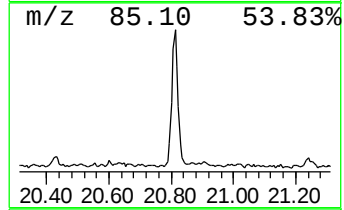
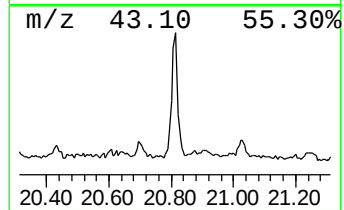
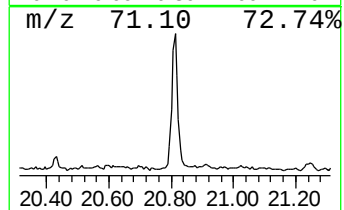
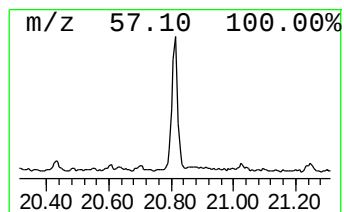
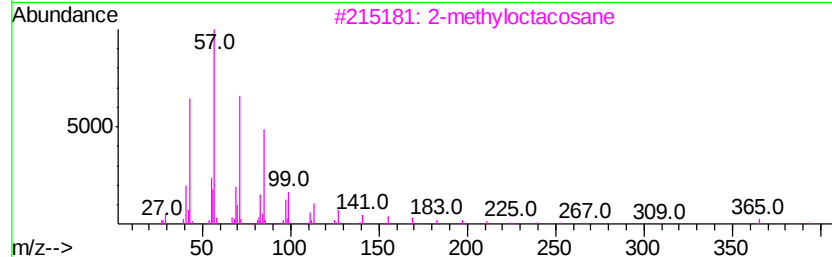
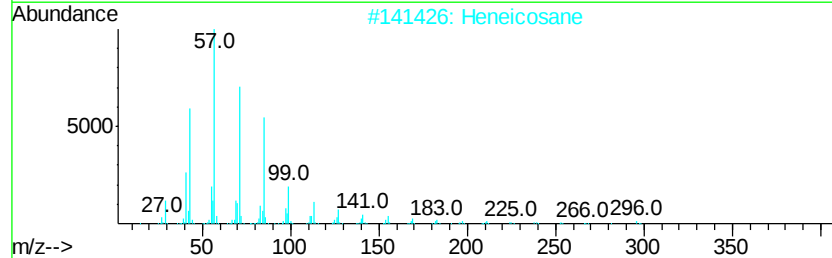
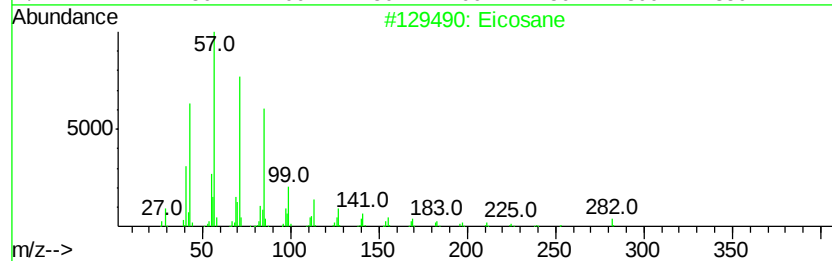
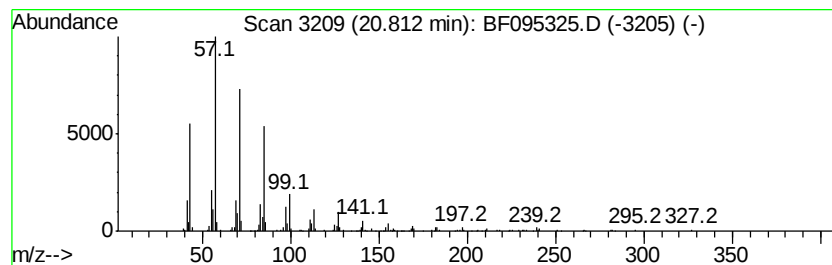
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Eicosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.81	12.86 ng	304190	Perylene-d12	20.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	95
2		Heneicosane	296	C21H44	000629-94-7	91
3		2-methyloctacosane	408	C29H60	1000376-72-8	91
4		Hexacosane	366	C26H54	000630-01-3	91
5		Octacosane	394	C28H58	000630-02-4	91



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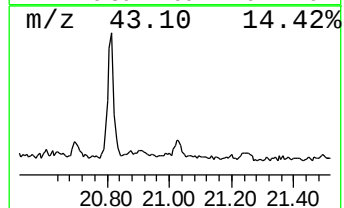
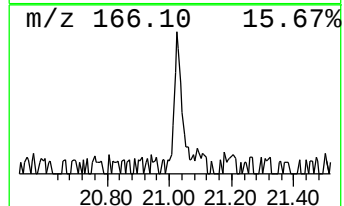
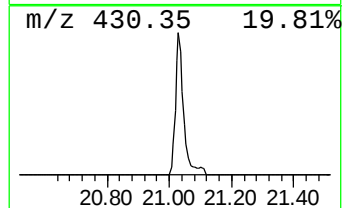
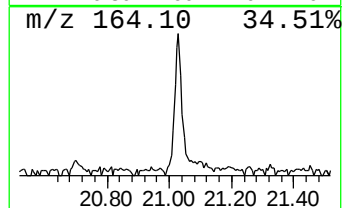
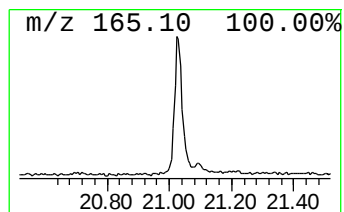
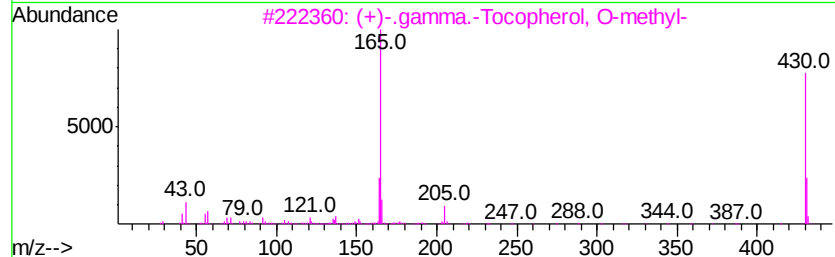
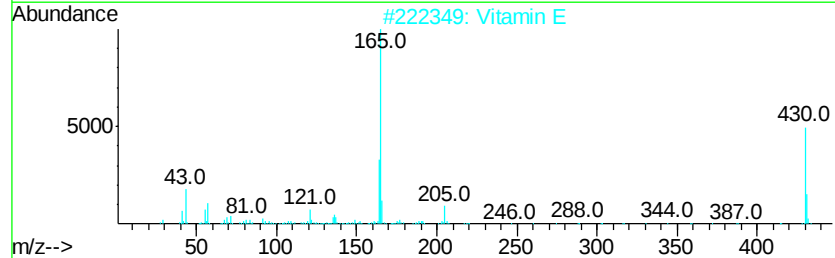
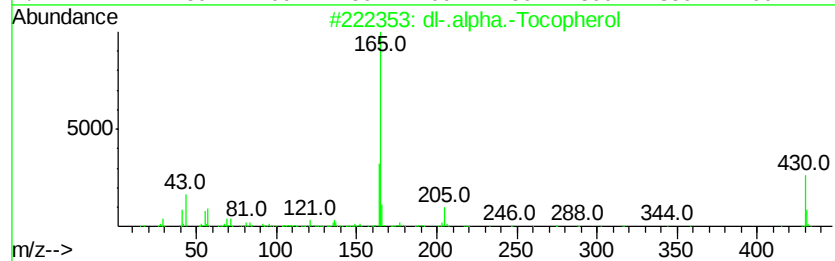
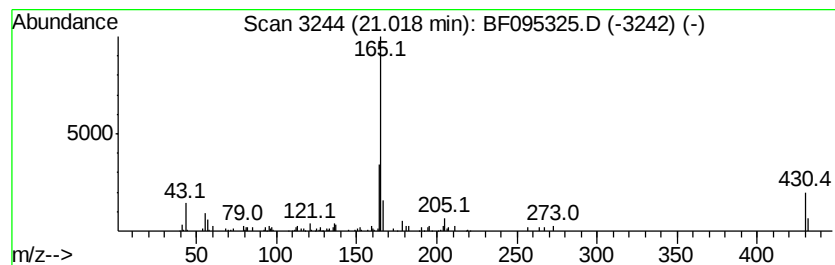
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 dl-.alpha.-Tocopherol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.02	5.55 ng	131429	Perylene-d12	20.36

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			dl-.alpha.-Tocopherol	430	C29H50O2	010191-41-0	93
2			Vitamin E	430	C29H50O2	000059-02-9	93
3			(+)-.gamma.-Tocopherol, 0-methyl-	430	C29H50O2	1000374-72-2	90
4			Benzenamine, 3-methoxy-2,4,6-tri...	165	C10H15NO	034874-88-9	50
5			9,10-Anthracenedione, 1,2-dimethyl-	236	C16H12O2	003285-98-1	43



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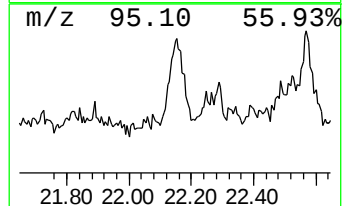
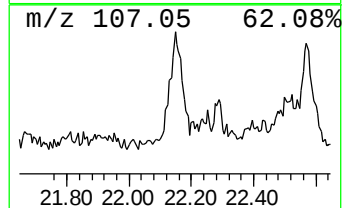
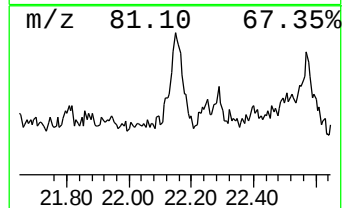
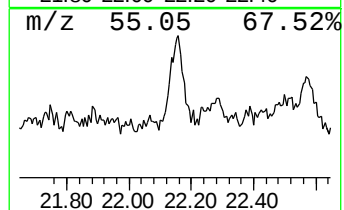
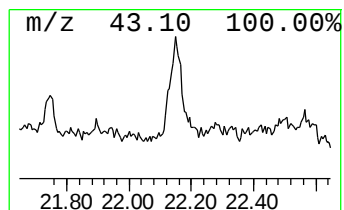
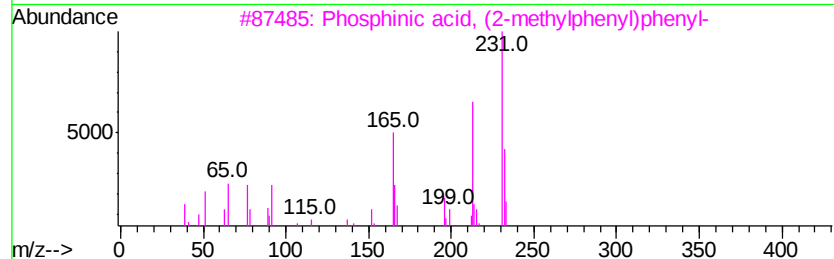
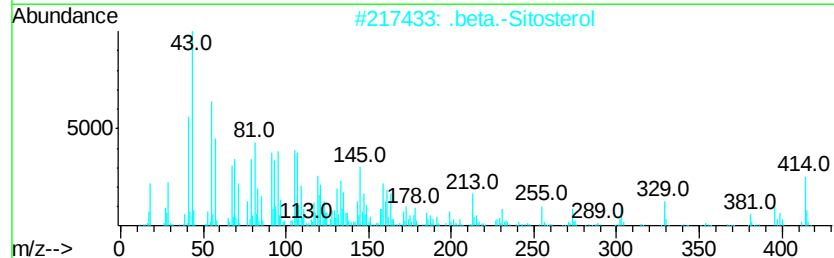
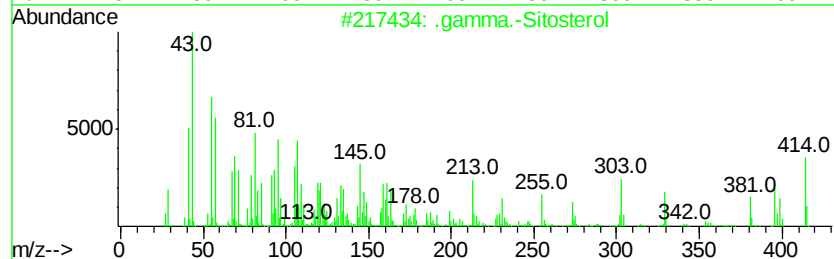
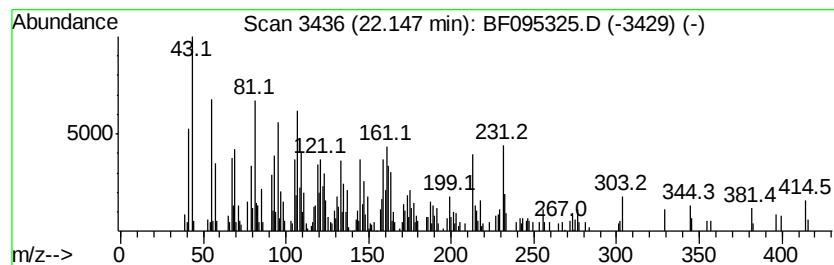
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown22.15 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.15	23.42 ng	554275	Perylene-d12	20.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	.gamma.-Sitosterol	414	C29H50O	000083-47-6	45
2		.beta.-Sitosterol	414	C29H50O	000083-46-5	42
3		Phosphinic acid, (2-methylphenyl)phenyl-	232	C13H13O2P	018593-18-5	18
4		Ethanone, 1-(1,2,3,4,7,7a-hexahydro-1H-benzofluoren-1-yl)-	246	C17H26O	032388-58-2	11
5		1,2,3,4,5,6-Hexahydro-[2,3']bipyridine	344	C19H24N2O2S	1000304-05-3	11



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052217\
Data File : BF095325.D
Acq On : 22 May 2017 21:06
Operator : SJ/MA
Sample : I3261-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampled :
NCC-BUCK-3-1

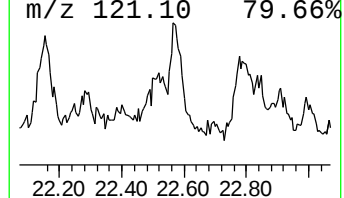
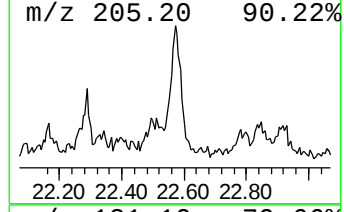
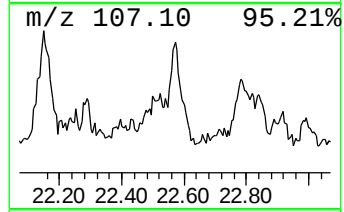
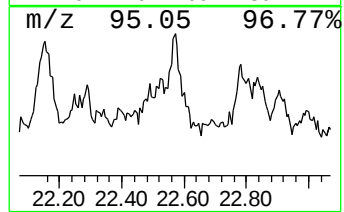
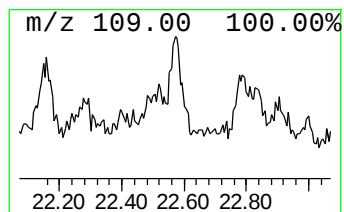
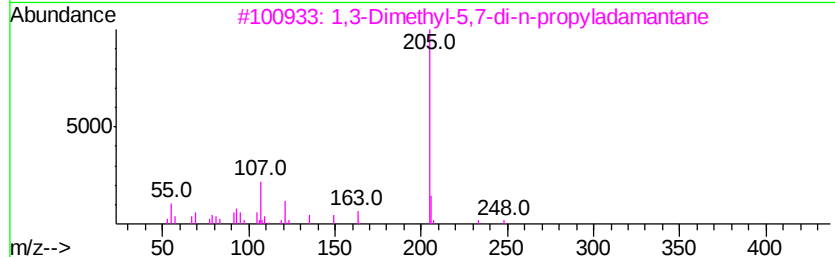
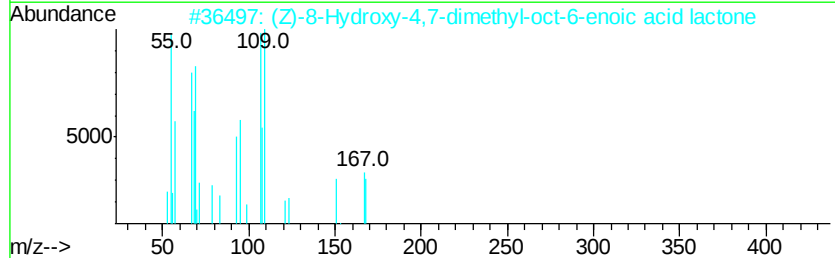
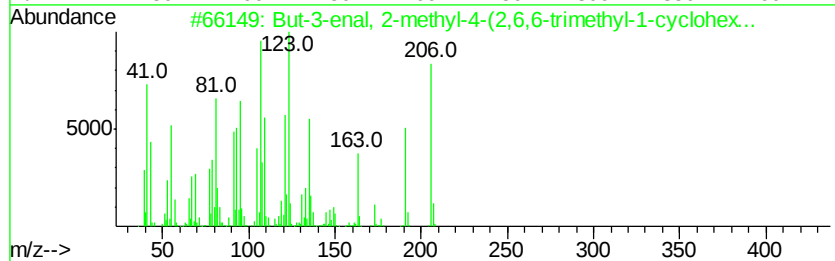
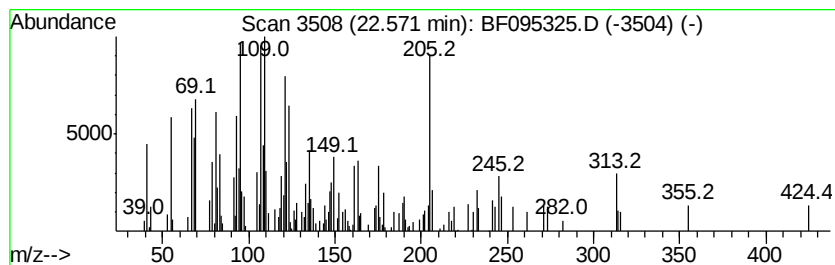
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown22.57 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.57	15.78 ng	373462	Perylene-d12	20.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	But-3-enal, 2-methyl-4-(2,6,6-tr...	206	C14H22O	104900-59-6	46
2		(Z)-8-Hydroxy-4,7-dimethyl-oct-6...	168	C10H16O2	106542-12-5	27
3		1,3-Dimethyl-5,7-di-n-propyladam...	248	C18H32	1000215-00-5	27
4		1,5-Cyclodecadiene, 1,5-dimethyl...	204	C15H24	015423-57-1	25
5		.gamma.-Elemene	204	C15H24	029873-99-2	25



Data Path : Z:\HPCHEM1\BNA F\DATA\BF052217\
Data File : BF095325.D
Acq On : 22 May 2017 21:06
Operator : SJ/MA
Sample : I3261-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampled :
NCC-BUCK-3-1

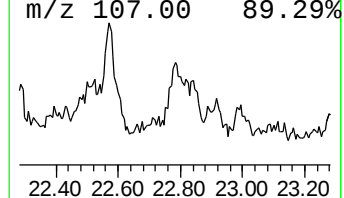
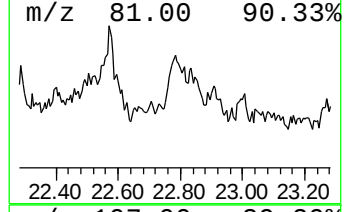
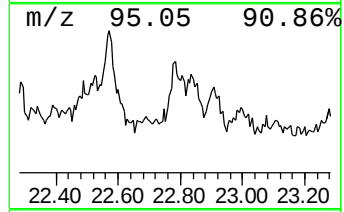
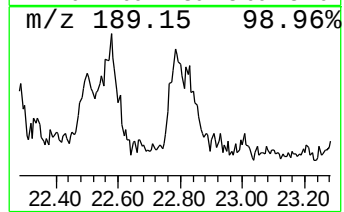
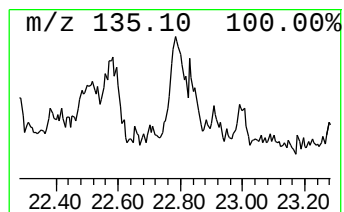
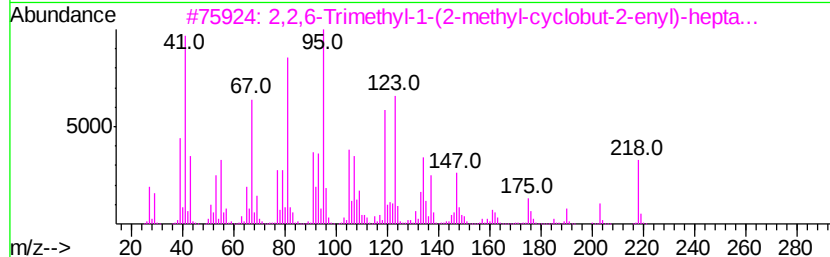
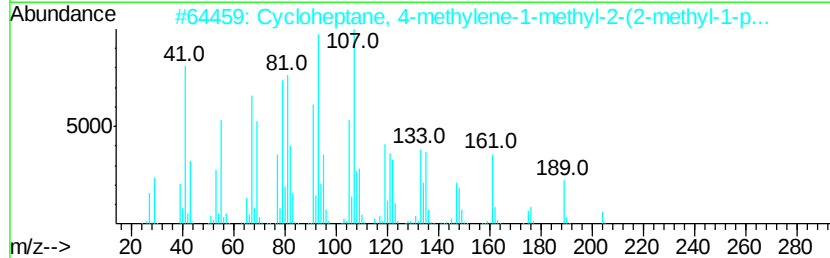
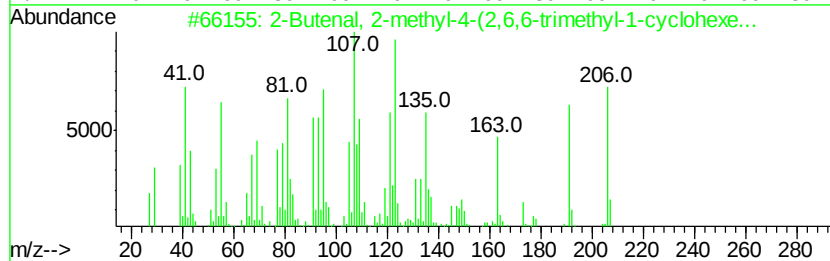
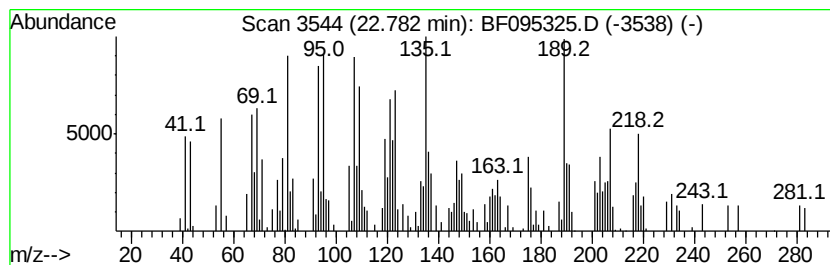
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 unknown22.78 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.78	6.88 ng	162849	Perylene-d12	20.36

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Butenal, 2-methyl-4-(2,6,6-tri...	206	C14H22O	003155-71-3	38
2			Cycloheptane, 4-methylene-1-meth...	204	C15H24	1000159-38-5	38
3			2,2,6-Trimethyl-1-(2-methyl-cycl...	218	C15H22O	1000188-72-8	35
4			(-)-Isolongifolol, methyl ether	236	C16H28O	1000333-80-8	25
5			.alpha.-Bulnesene	204	C15H24	1000374-19-9	25



Data Path : Z:\HPCHEM1\BNA F\DATA\BF052217\
Data File : BF095325.D
Acq On : 22 May 2017 21:06
Operator : SJ/MA
Sample : I3261-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

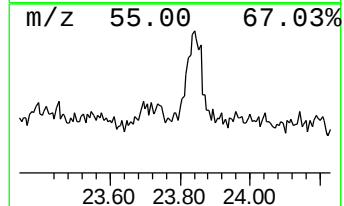
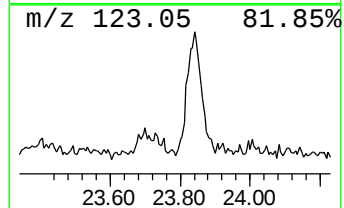
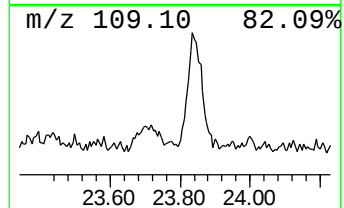
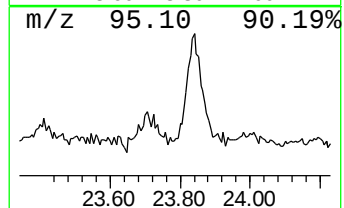
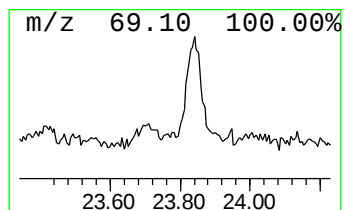
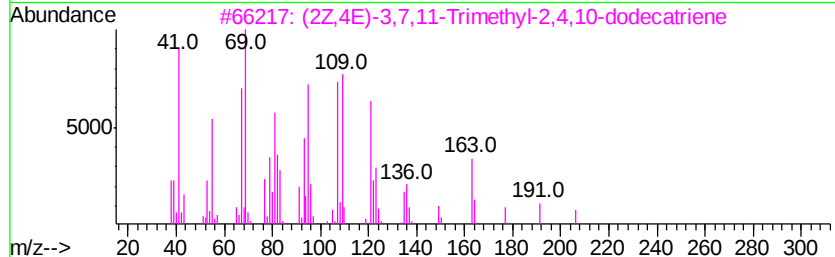
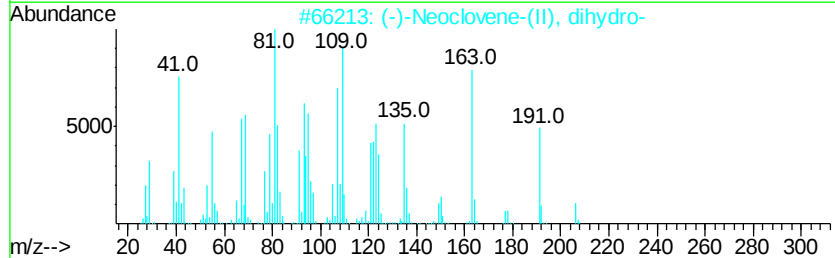
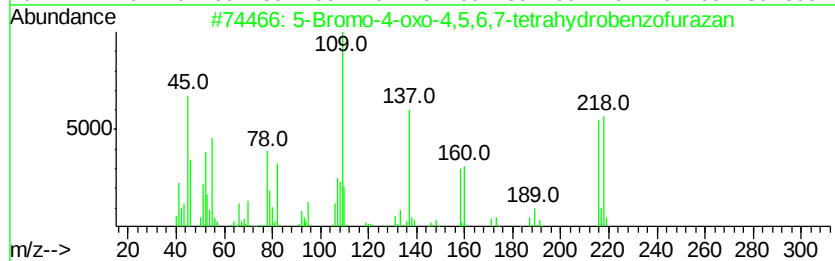
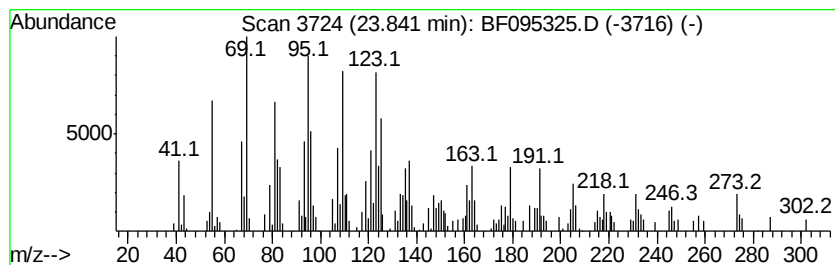
Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 5-Bromo-4-oxo-4,5,6,7-tetra... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
23.84	16.60 ng	392844	Perylene-d12	20.36	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Bromo-4-oxo-4,5,6,7-tetrahydro...	216	C6H5BrN2O2	300574-36-1	60
2	(-)-Neoclovene-(II), dihydro-	206	C15H26	1000152-83-6	47
3	(2Z,4E)-3,7,11-Trimethyl-2,4,10-...	206	C15H26	172549-29-0	43
4	1,4-Methanoazulene-9-methanol, d...	222	C15H26O	001139-17-9	42
5	d-Norandrostane (5.alpha.,14.alp...	246	C18H30	1000281-13-8	25



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052217\
 Data File : BF095325.D
 Acq On : 22 May 2017 21:06
 Operator : SJ/MA
 Sample : I3261-01
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NCC-BUCK-3-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.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...	5.12	8.1 ng		186960	1	7.76	462878	20.0
unknown7.42	7.42	89.7 ng		2075430	1	7.76	462878	20.0
n-Hexadecanoic acid	15.97	5.5 ng		132157	4	15.01	482122	20.0
Nonahexacontanoic...	18.57	2.2 ng		62475	5	18.68	564243	20.0
Hexacosane	19.36	2.3 ng		63913	5	18.68	564243	20.0
unknown19.97	19.97	13.3 ng		314747	6	20.36	473258	20.0
Octacosane	20.08	12.1 ng		286322	6	20.36	473258	20.0
2H-1-Benzopyran-6...	20.31	2.9 ng		67351	6	20.36	473258	20.0
.gamma.-Tocopherol	20.70	5.4 ng		126915	6	20.36	473258	20.0
Eicosane	20.81	12.9 ng		304190	6	20.36	473258	20.0
dl-.alpha.-Tocoph...	21.02	5.5 ng		131429	6	20.36	473258	20.0
unknown22.15	22.15	23.4 ng		554275	6	20.36	473258	20.0
unknown22.57	22.57	15.8 ng		373462	6	20.36	473258	20.0
unknown22.78	22.78	6.9 ng		162849	6	20.36	473258	20.0
5-Bromo-4-oxo-4,5...	23.84	16.6 ng		392844	6	20.36	473258	20.0