

Data Path : Z:\HPCHEM1\BNA F\DATA\BF112217\
 Data File : BF100838.D
 Acq On : 22 Nov 2017 22:08
 Operator : SJ/JU
 Sample : I6043-19
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 N-T0A-BAJA-S001-0001-03

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

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

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110817.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	5.093	506	511	522	rBV	216608	322923	12.95%	1.497%
2	5.487	572	578	581	rBV	1833642	2095515	84.03%	9.714%
3	6.492	743	749	752	rBV	1889325	2232192	89.51%	10.348%
4	6.634	767	773	776	rBV	2172002	2217117	88.91%	10.278%
5	6.857	808	811	815	rBV	556709	450926	18.08%	2.090%
6	7.016	833	838	841	rBV	1900013	1689390	67.75%	7.831%
7	7.428	902	908	911	rBV	1272863	1421416	57.00%	6.589%
8	8.139	1025	1029	1033	rBV	711020	611716	24.53%	2.836%
9	8.692	1119	1123	1127	rBV	190123	156772	6.29%	0.727%
10	9.216	1207	1212	1215	rBV	2460542	2493696	100.00%	11.560%
11	9.551	1265	1269	1271	rBV	65766	50803	2.04%	0.236%
12	9.604	1275	1278	1282	rVV	188341	148055	5.94%	0.686%
13	9.898	1324	1328	1331	rBV	774249	669113	26.83%	3.102%
14	10.092	1357	1361	1365	rBV2	26298	26078	1.05%	0.121%
15	10.310	1391	1398	1403	rBV4	20089	37319	1.50%	0.173%
16	10.610	1445	1449	1453	rBV	883670	808114	32.41%	3.746%
17	10.686	1458	1462	1466	rBV	1439343	1493946	59.91%	6.925%
18	10.939	1499	1505	1509	rBV	53349	55452	2.22%	0.257%
19	10.992	1509	1514	1517	rBV	161109	159975	6.42%	0.742%
20	11.039	1517	1522	1525	rBV	179556	158505	6.36%	0.735%
21	11.286	1561	1564	1567	rVB	46463	34577	1.39%	0.160%
22	11.380	1577	1580	1584	rVB	725769	636611	25.53%	2.951%
23	11.904	1666	1669	1678	rBV2	193975	180451	7.24%	0.837%
24	12.604	1784	1788	1791	rBV5	18277	31572	1.27%	0.146%
25	12.686	1798	1802	1806	rBV	120567	130145	5.22%	0.603%
26	12.968	1845	1850	1853	rBV	1915592	1800822	72.21%	8.348%
27	13.851	1997	2000	2003	rBV	74266	72414	2.90%	0.336%
28	13.886	2003	2006	2010	rVB	26182	25127	1.01%	0.116%
29	13.992	2021	2024	2026	rBV	39683	35871	1.44%	0.166%
30	14.021	2026	2029	2035	rVB	447564	389313	15.61%	1.805%
31	15.492	2274	2279	2287	rVB	373670	455619	18.27%	2.112%
32	15.862	2338	2342	2348	rBV6	16577	29711	1.19%	0.138%
33	15.933	2350	2354	2360	rBV2	20012	28634	1.15%	0.133%
34	16.356	2422	2426	2432	rVB8	15098	25383	1.02%	0.118%

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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

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35	16.439	2435	2440	2448	rBV4	23935	46810	1.88%	0.217%
36	17.039	2537	2542	2549	rVB2	22957	38640	1.55%	0.179%
37	17.127	2551	2557	2565	rVB9	18897	43061	1.73%	0.200%
38	17.533	2620	2626	2636	rVB5	36816	84041	3.37%	0.390%
39	17.739	2657	2661	2669	rVB6	12267	26460	1.06%	0.123%
40	18.109	2720	2724	2730	rVB6	30972	55349	2.22%	0.257%
41	18.392	2765	2772	2780	rVB8	17538	46650	1.87%	0.216%
42	18.568	2794	2802	2810	rBV6	19512	55539	2.23%	0.257%

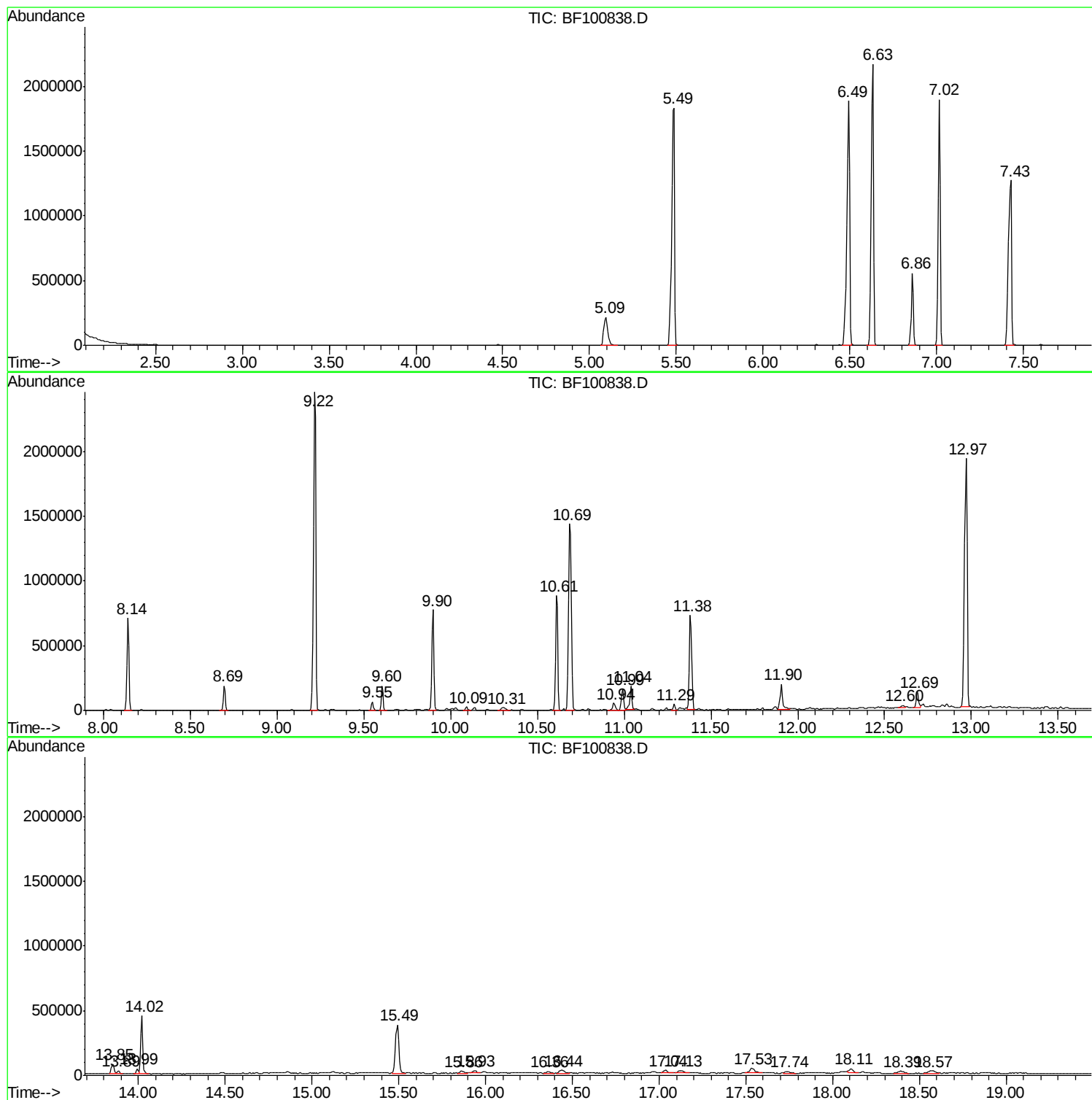
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110817.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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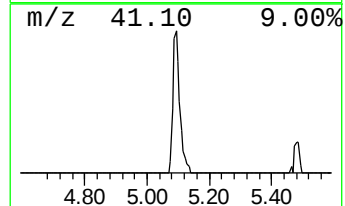
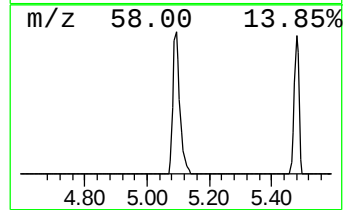
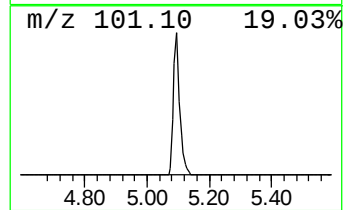
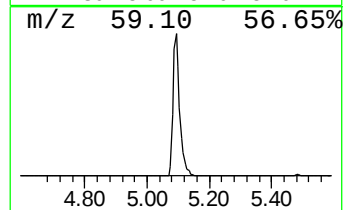
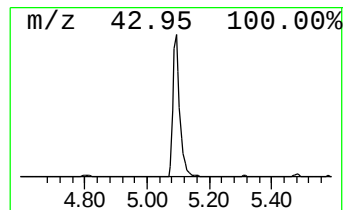
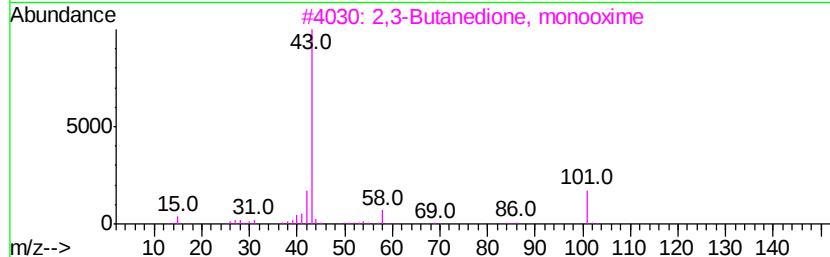
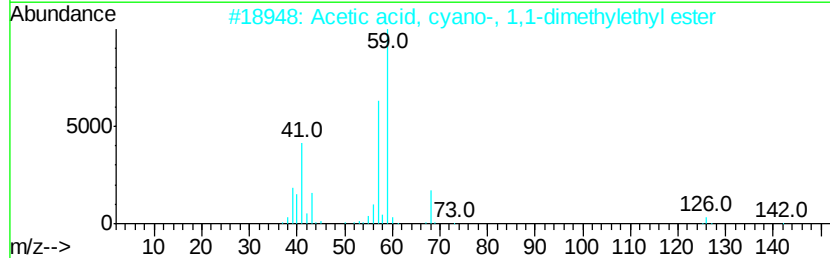
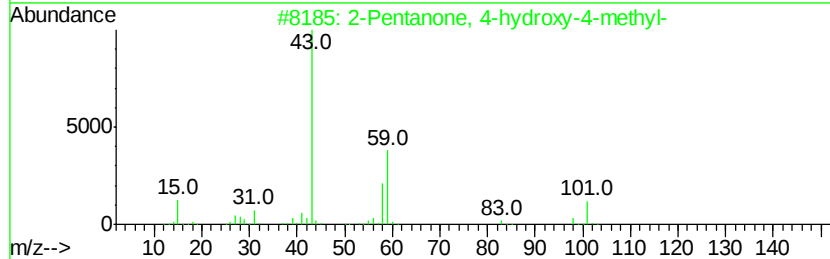
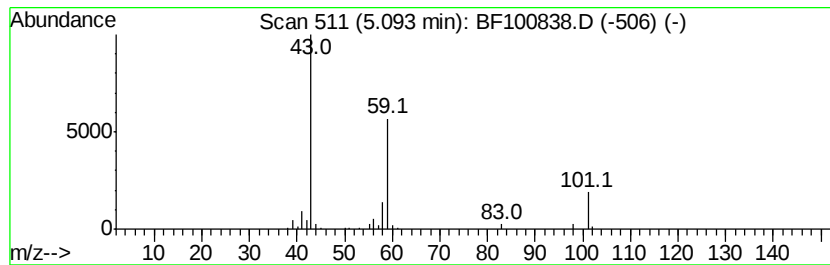
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110817.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
5.09	14.32 ng	322923	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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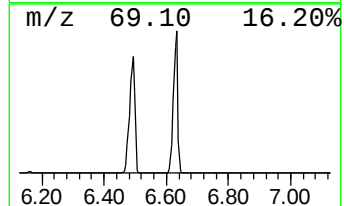
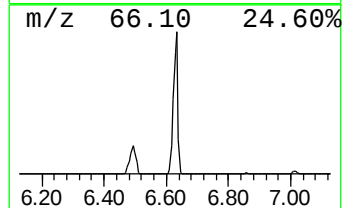
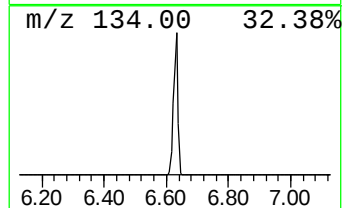
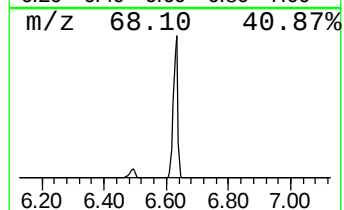
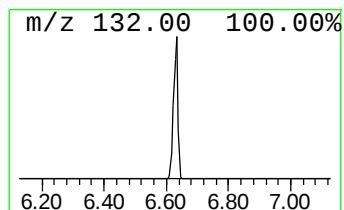
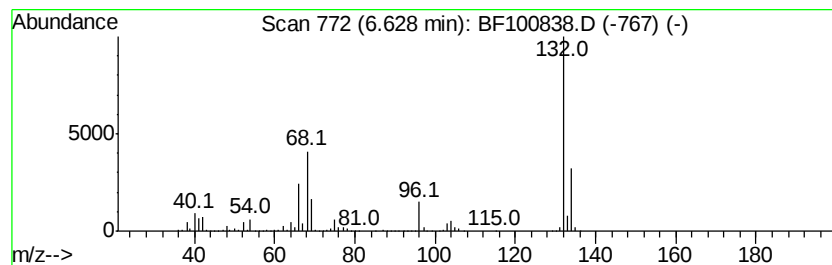
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110817.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 unknown6.63 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.63	98.34 ng	2217120	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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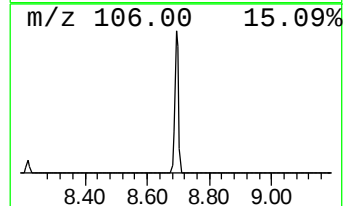
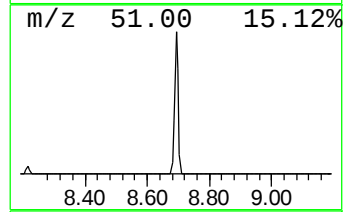
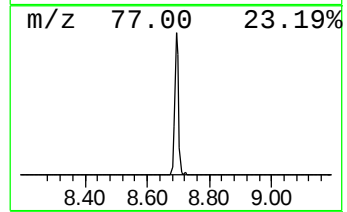
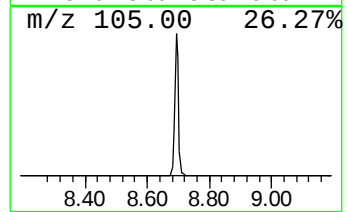
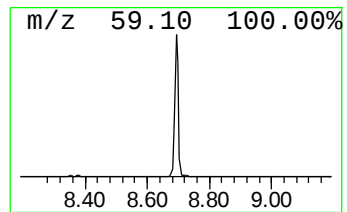
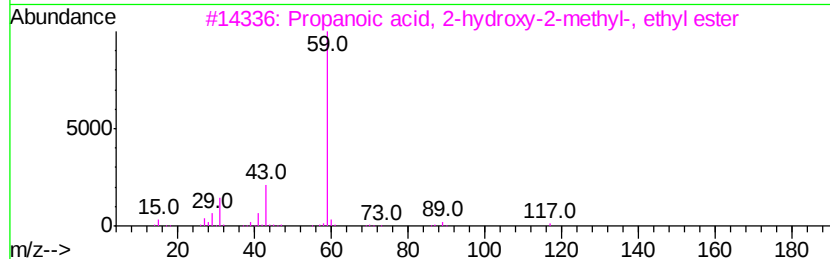
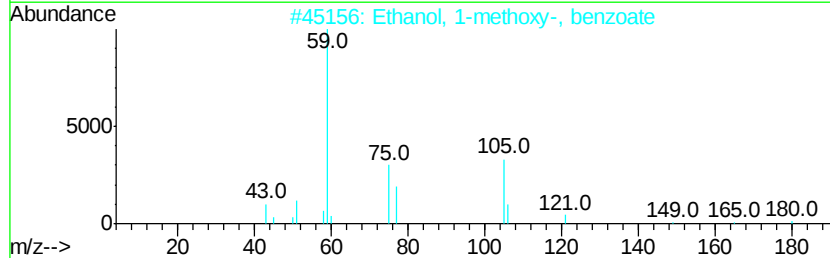
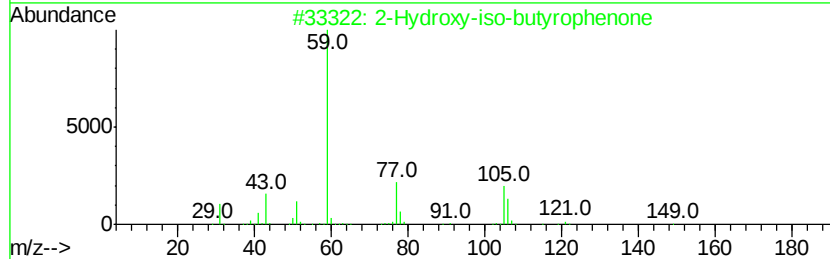
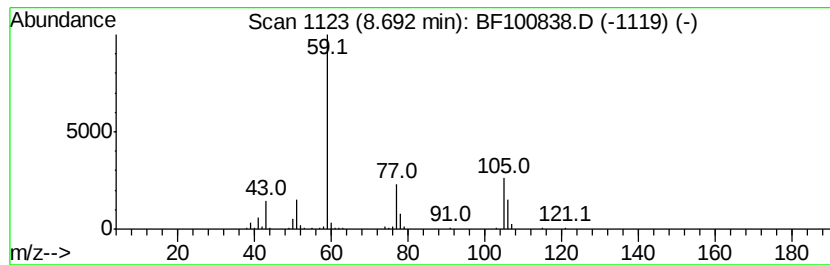
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 2-Hydroxy-iso-butyrophenone Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.	
8.69	5.13 ng	156772	Naphthalene-d8		8.14	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hydroxy-iso-butyrophenone		164	C10H12O2	007473-98-5	90
2	Ethanol, 1-methoxy-, benzoate		180	C10H12O3	051835-44-0	9
3	Propanoic acid, 2-hydroxy-2-meth...		132	C6H12O3	000080-55-7	9
4	Guanidine		59	CH5N3	000113-00-8	7
5	Formamide, N-methyl-		59	C2H5NO	000123-39-7	7



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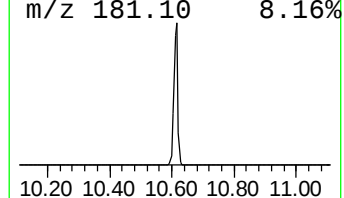
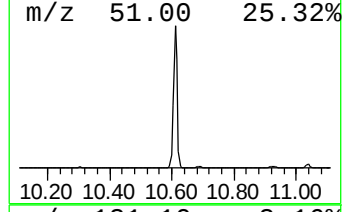
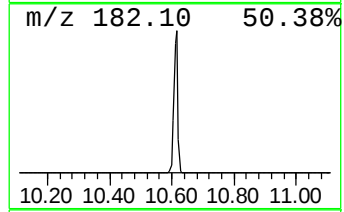
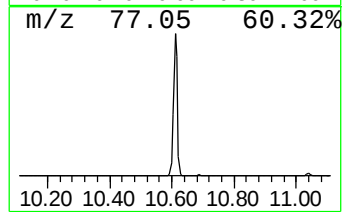
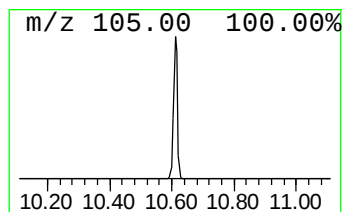
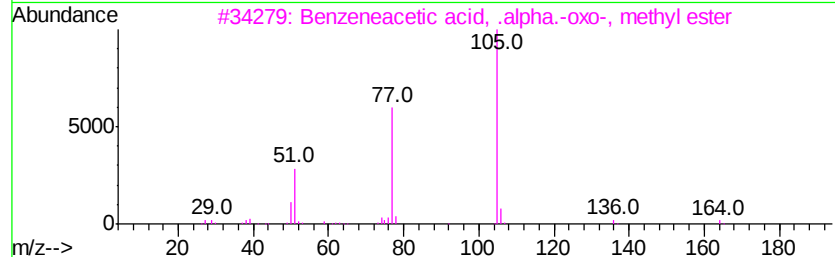
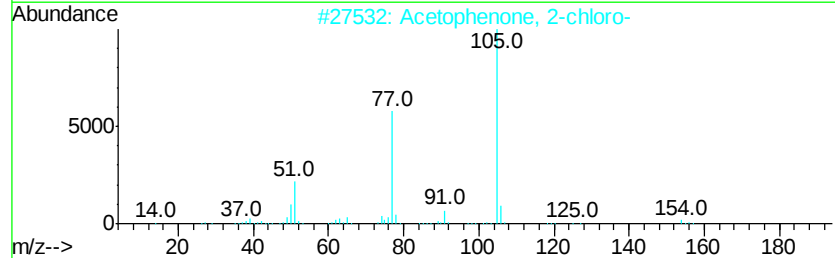
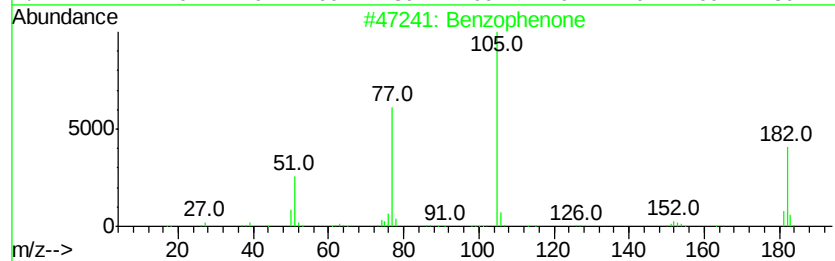
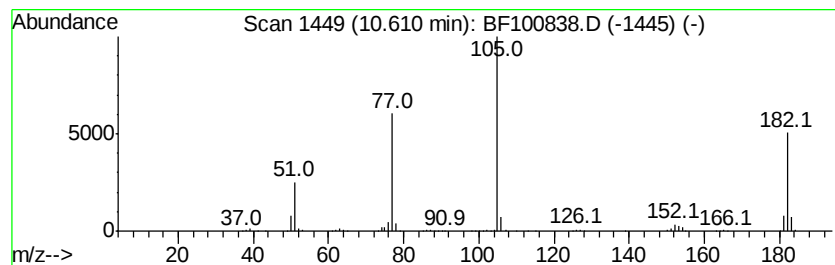
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzophenone Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.61	24.15 ng	808114	Acenaphthene-d10	9.90
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzophenone	182 C13H10O	000119-61-9 97
2		Acetophenone, 2-chloro-	154 C8H7ClO	000532-27-4 49
3		Benzeneacetic acid, .alpha.-oxo-...	164 C9H8O3	015206-55-0 47
4		Vinyl benzoate	148 C9H8O2	000769-78-8 47
5		N-Methoxy-N-methylbenzamide	165 C9H11NO2	006919-61-5 47



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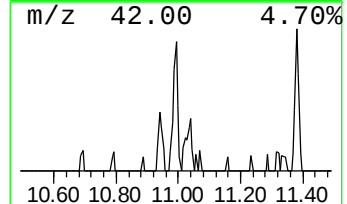
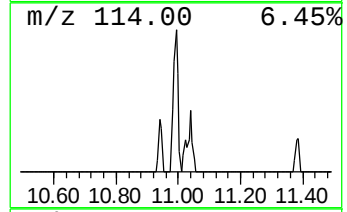
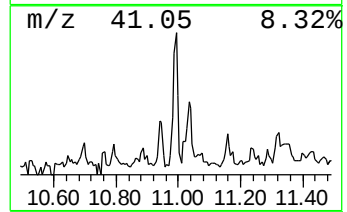
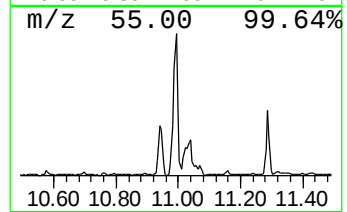
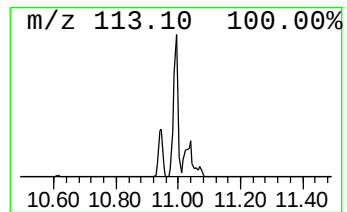
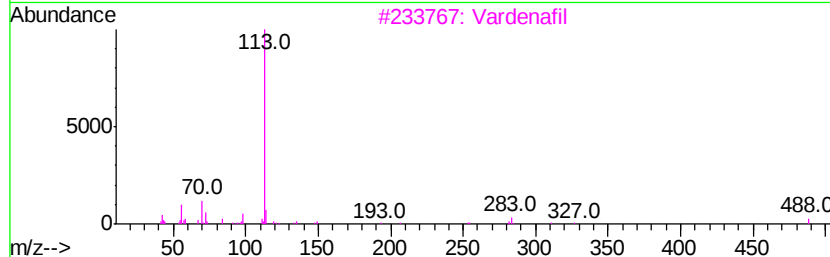
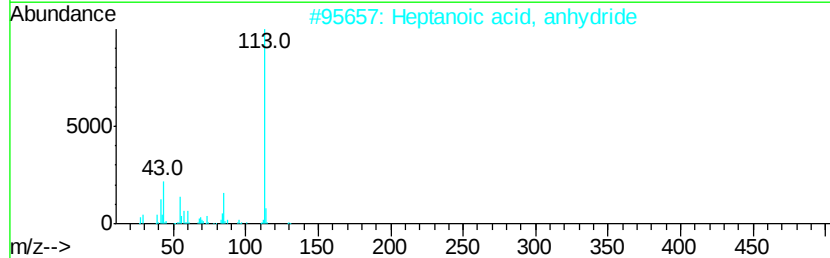
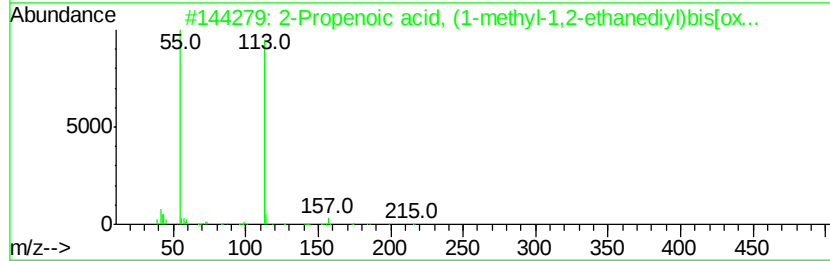
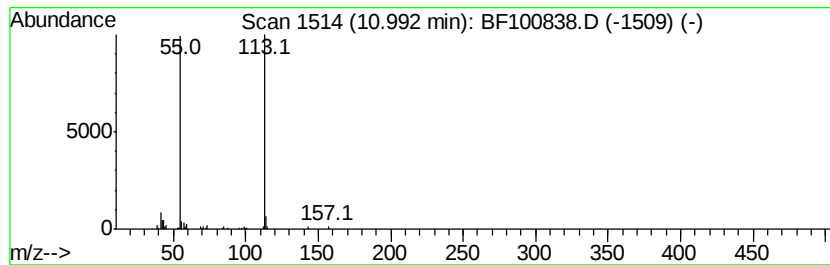
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 2-Propenoic acid, (1-methyl... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.99	5.03 ng	159975	Phenanthrene-d10	11.38

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Propenoic acid, (1-methyl-1,2-...	300	C15H24O6	042978-66-5	78
2		Heptanoic acid, anhydride	242	C14H26O3	000626-27-7	36
3		Vardenafil	488	C23H32N6O4S	224785-90-4	36
4		L-2-Aminobutyric acid, N-dimethy...	214	C11H22N2O2	1000375-52-5	36
5		2H-Pyrrrol-2-one, 1,5-dihydro-4-m...	113	C5H7NO2	069778-83-2	9



Data Path : Z:\HPCHEM1\BNA F\DATA\BF112217\
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Acq On : 22 Nov 2017 22:08
Operator : SJ/JU
Sample : I6043-19
Misc :
ALS Vial : 11 Sample Multiplier: 1

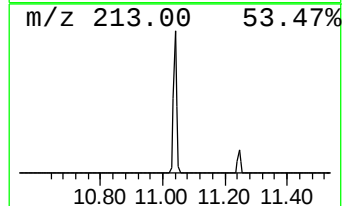
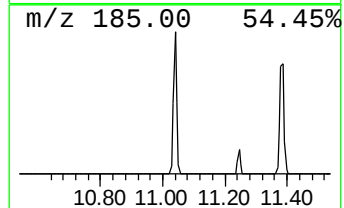
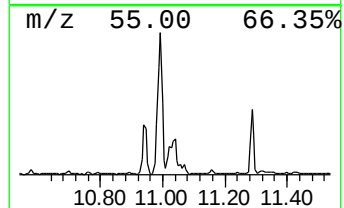
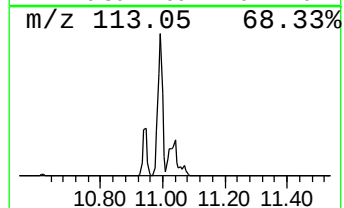
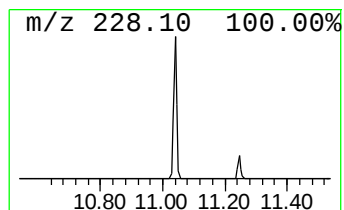
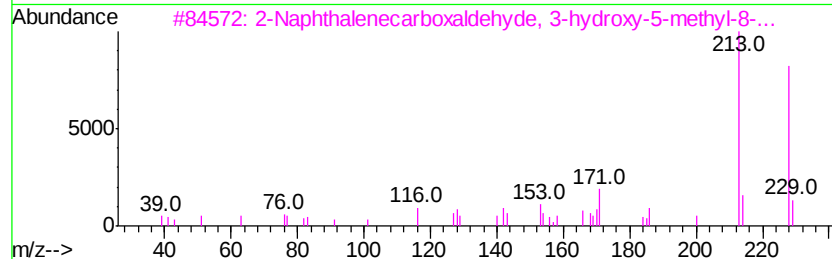
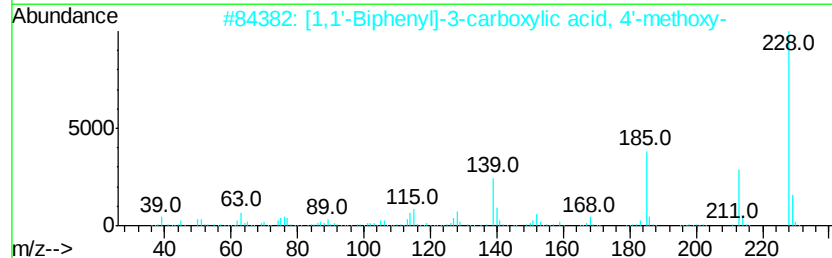
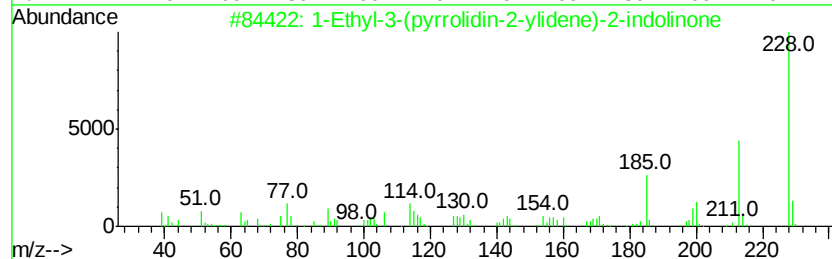
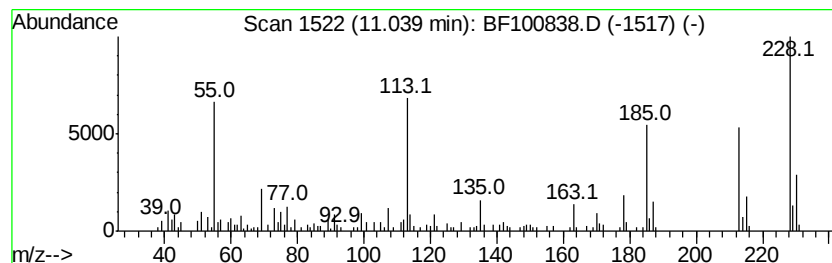
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 1-Ethyl-3-(pyrrolidin-2-ylidene) Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.04	4.98 ng	158505	Phenanthrene-d10		11.38	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Ethyl-3-(pyrrolidin-2-ylidene)...	228	C14H16N2O	1000260-86-3	62
2		[1,1'-Biphenyl]-3-carboxylic aci...	228	C14H12O3	1000338-15-5	47
3		2-Naphthalenecarboxaldehyde, 3-h...	228	C15H16O2	018478-73-4	43
4		1,2,3,4,5-Pentamethoxybenzene	228	C11H16O5	013909-75-6	43
5		1,4-Dimethylphenoxathin	228	C14H12OS	016314-44-6	27



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112217\
Data File : BF100838.D
Acq On : 22 Nov 2017 22:08
Operator : SJ/JU
Sample : I6043-19
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Instrument :
BNA_F
ClientSampleId :
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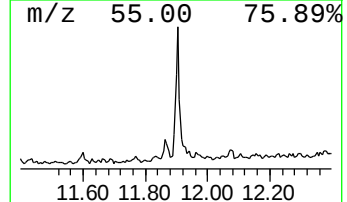
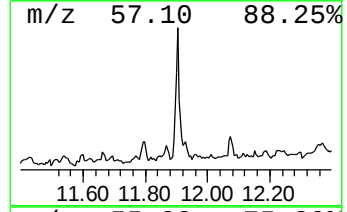
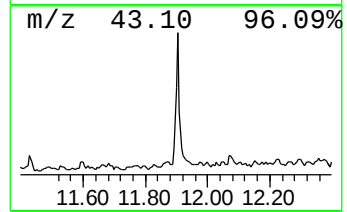
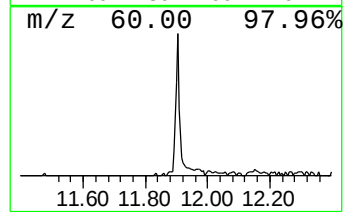
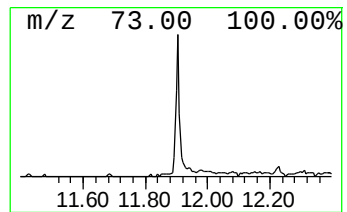
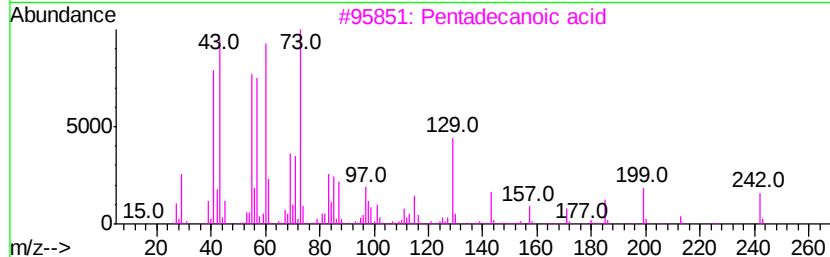
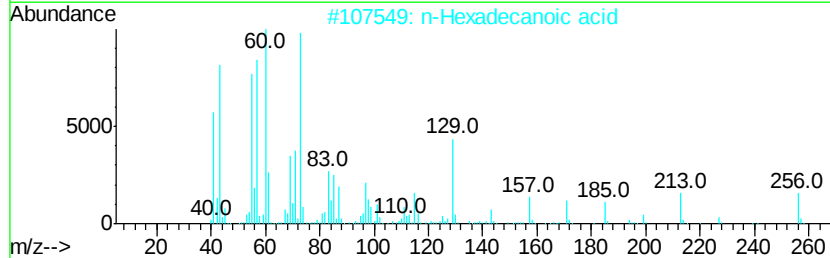
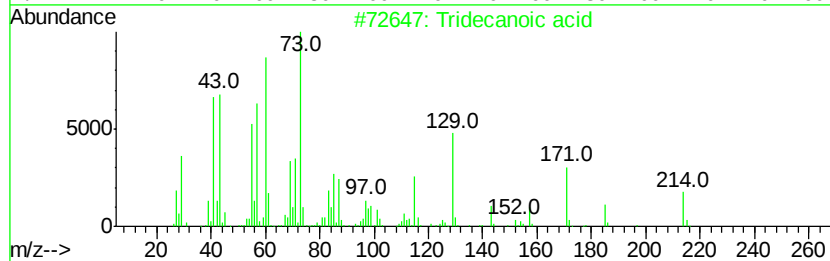
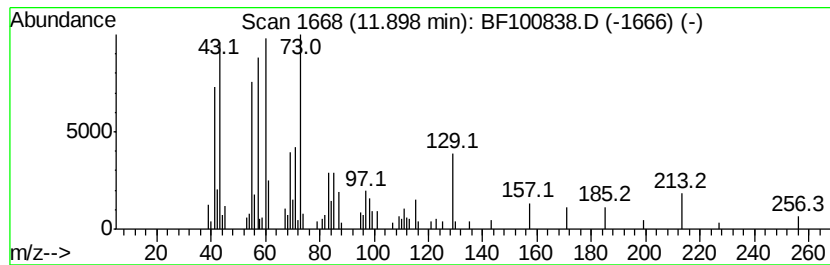
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Tridecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.90	5.67 ng	180451	Phenanthrene-d10	11.38

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecanoic acid	214	C13H26O2	000638-53-9	94
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	93
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	86
5	Undecanoic acid	186	C11H22O2	000112-37-8	83



Data Path : Z:\HPCHEM1\BNA F\DATA\BF112217\
Data File : BF100838.D
Acq On : 22 Nov 2017 22:08
Operator : SJ/JU
Sample : I6043-19
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
N-T0A-BAJA-S001-0001-03

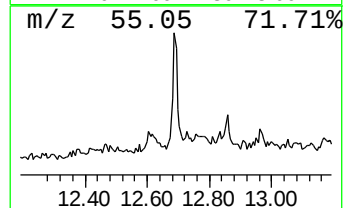
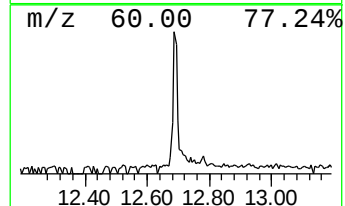
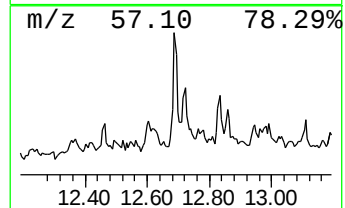
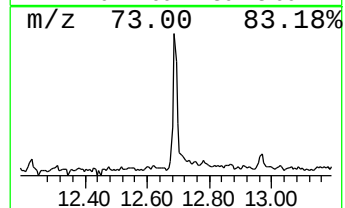
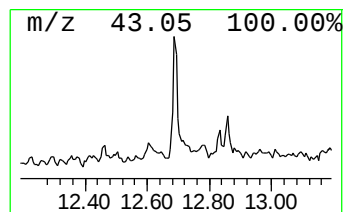
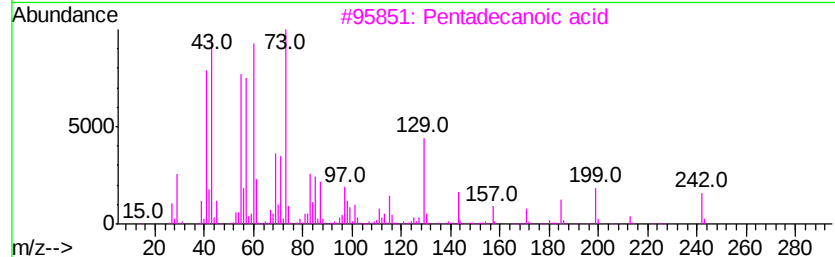
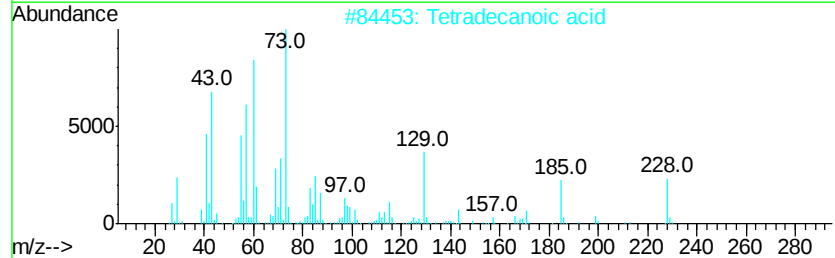
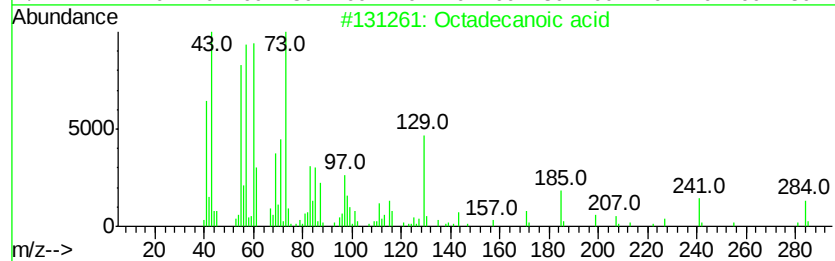
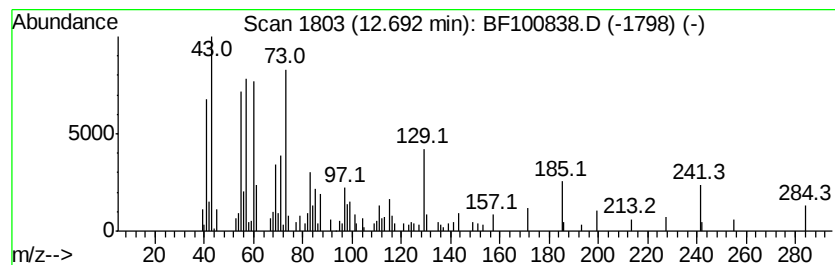
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Octadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.69	4.09 ng	130145	Phenanthrene-d10	11.38

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	98
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	87
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	68
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	68
5	n-Decanoic acid	172	C10H20O2	000334-48-5	49



Data Path : Z:\HPCHEM1\BNA F\DATA\BF112217\
Data File : BF100838.D
Acq On : 22 Nov 2017 22:08
Operator : SJ/JU
Sample : I6043-19
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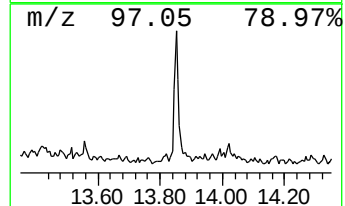
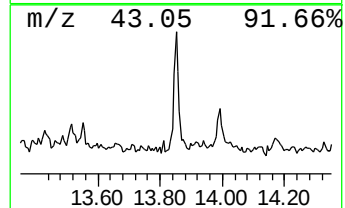
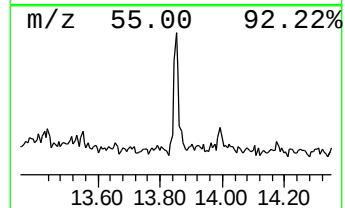
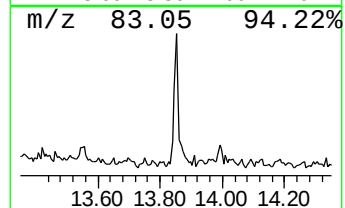
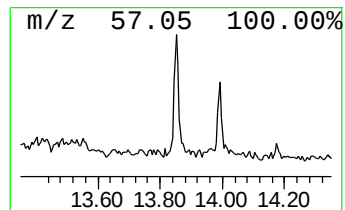
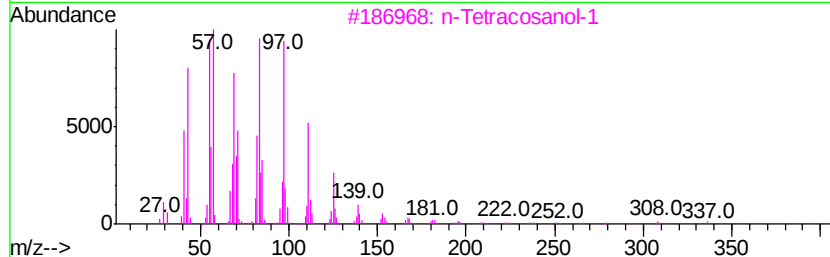
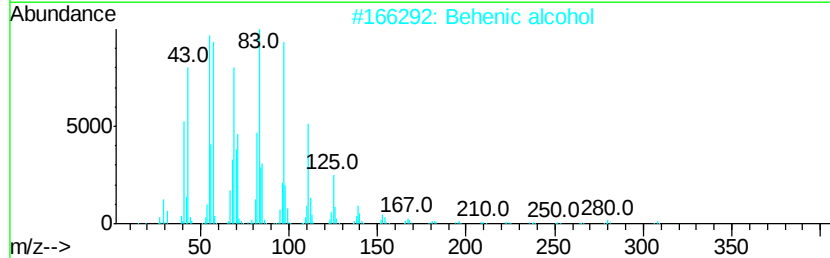
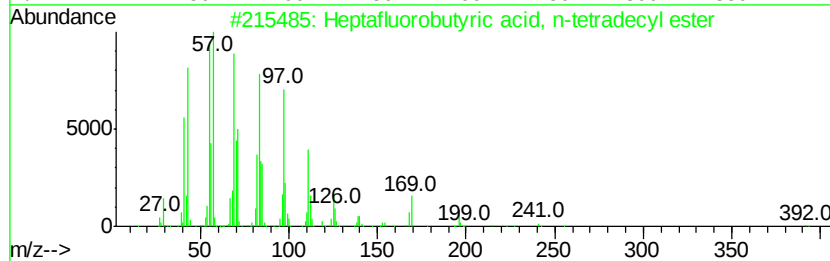
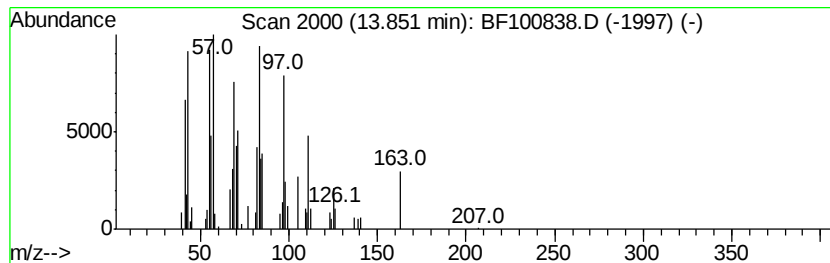
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Heptafluorobutyric acid, n-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.85	3.72 ng	72414	Chrysene-d12	14.02		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptafluorobutyric acid, n-tetra...	410	C18H29F7O2	007365-36-8	94
2		Behenic alcohol	326	C22H46O	000661-19-8	90
3		n-Tetracosanol-1	354	C24H50O	000506-51-4	90
4		3-Chloropropionic acid, heptadec...	346	C20H39ClO2	1000283-05-1	87
5		n-Nonadecanol-1	284	C19H40O	001454-84-8	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112217\
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Acq On : 22 Nov 2017 22:08
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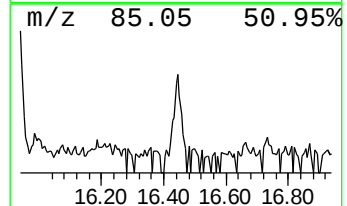
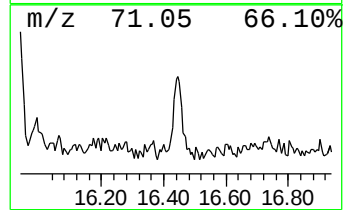
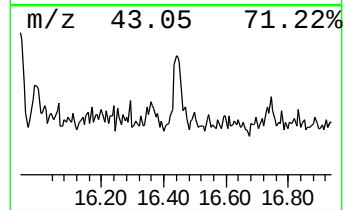
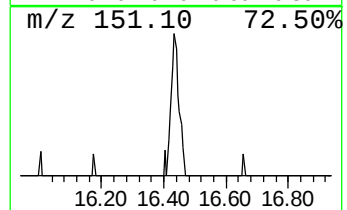
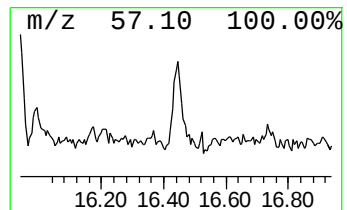
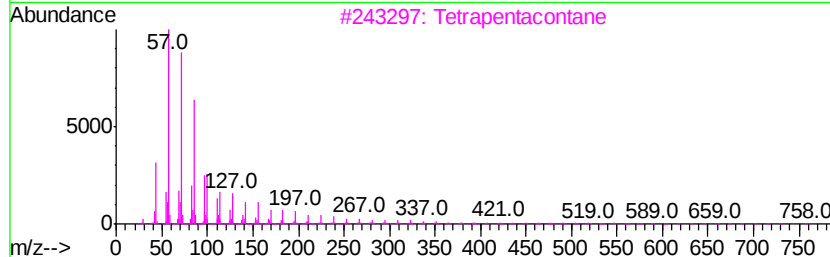
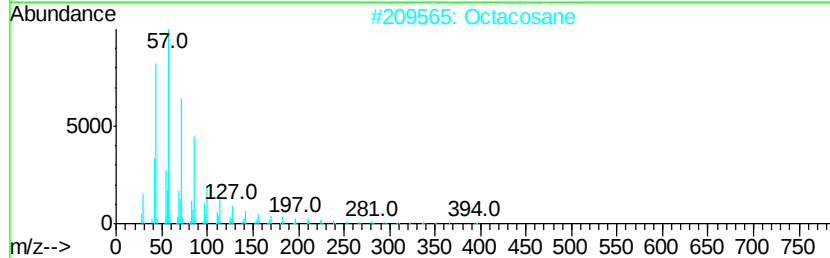
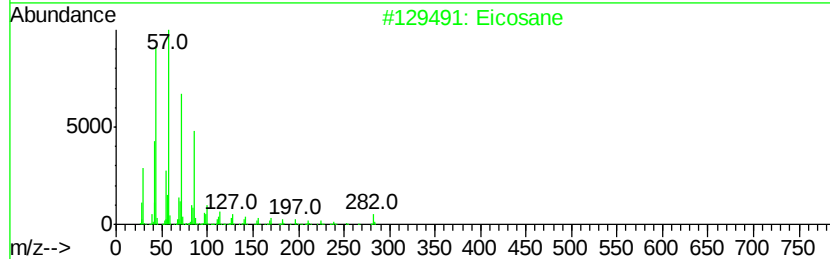
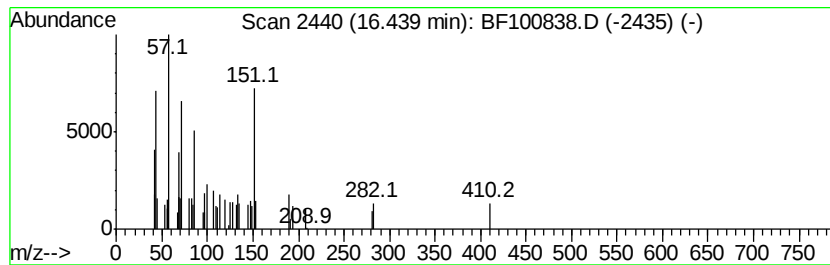
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown16.44 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD		R.T.		
16.44	2.05 ng	46810	Perylene-d12		15.49		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	46
2			Octacosane	394	C28H58	000630-02-4	38
3			Tetrapentacontane	759	C54H110	005856-66-6	38
4			Hexatriacontane	507	C36H74	000630-06-8	35
5			Heneicosane	296	C21H44	000629-94-7	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112217\
Data File : BF100838.D
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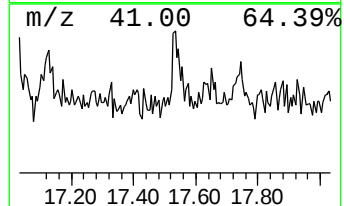
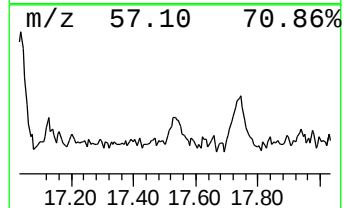
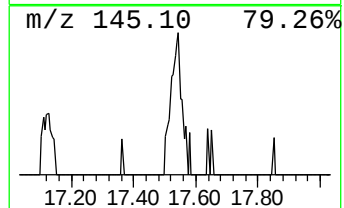
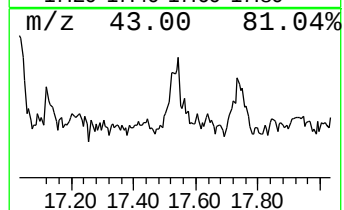
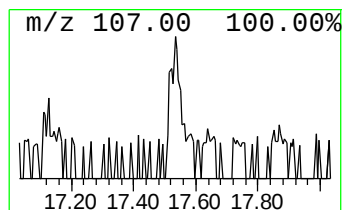
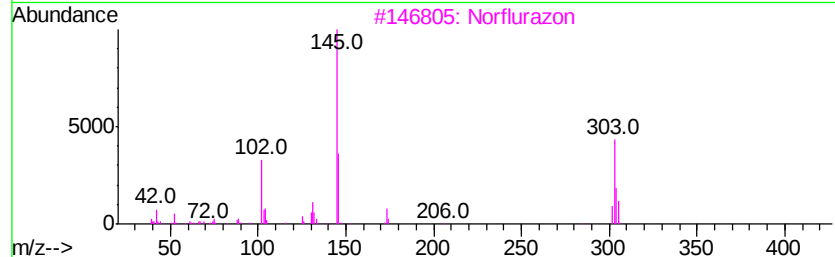
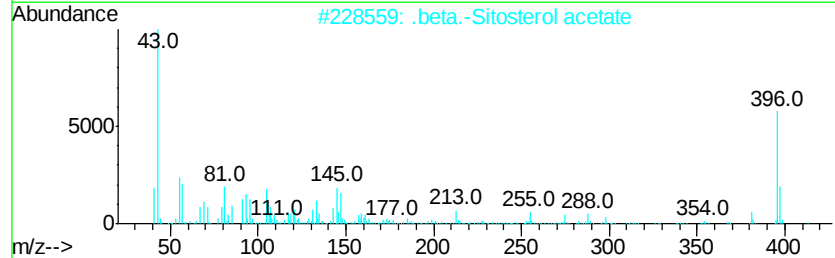
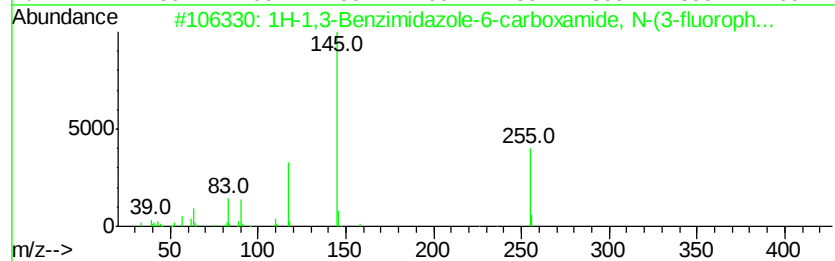
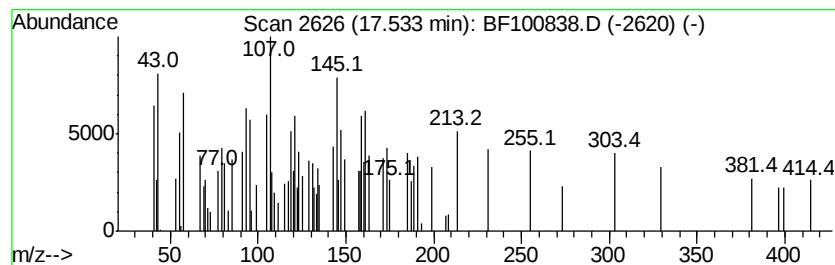
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown17.53 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.53	3.69 ng	84041	Perylene-d12	15.49

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-1,3-Benzimidazole-6-carboxami...	255	C14H10FN3O	1000350-37-6	14
2			.beta.-Sitosterol acetate	456	C31H52O2	000915-05-9	10
3			Norflurazon	303	C12H9ClF3N3O	027314-13-2	10
4			Imidazo[1,2-a]pyridine-2-methana...	255	C15H14FN3	1000319-52-3	9
5			Imidazo[1,2-a]pyridine-2-methana...	255	C15H14FN3	1000319-53-7	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112217\
Data File : BF100838.D
Acq On : 22 Nov 2017 22:08
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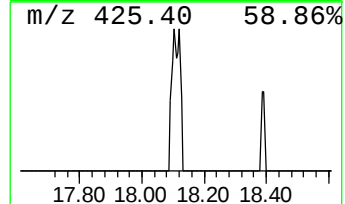
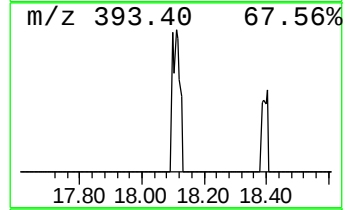
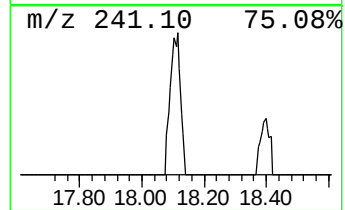
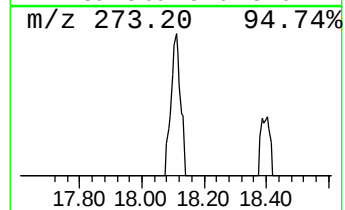
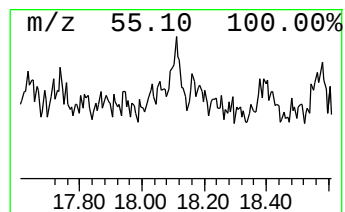
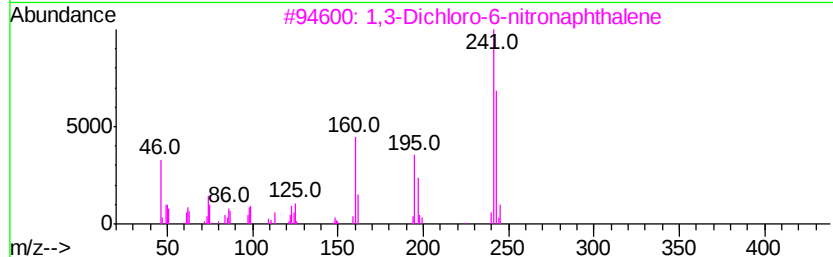
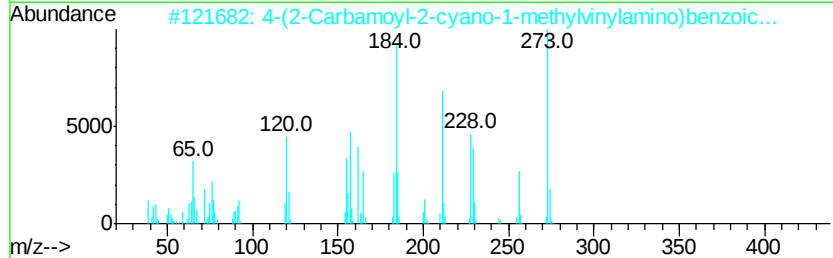
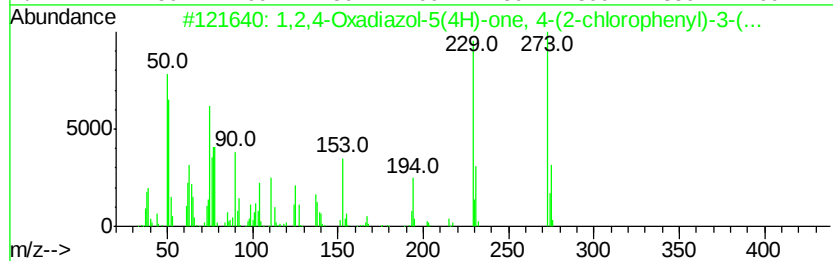
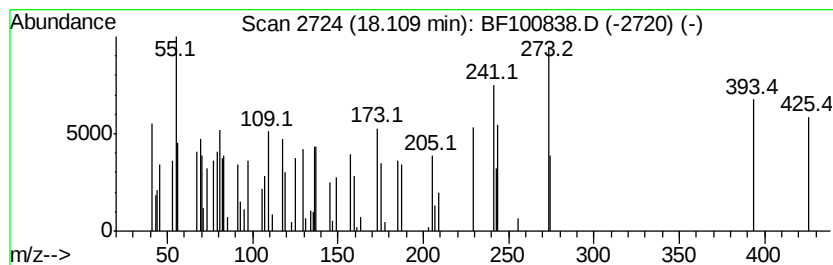
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown18.11 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.11	2.43 ng	55349	Perylene-d12	15.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,2,4-Oxadiazol-5(4H)-one, 4-(2-...	273	C13H8ClN3O2	288246-55-9	11
2		4-(2-Carbamoyl-2-cyano-1-methylv...	273	C14H15N3O3	1000311-98-1	9
3		1,3-Dichloro-6-nitronaphthalene	241	C10H5Cl2N02	023417-66-5	9
4		4-Ethyl-9-chloro-acridine	241	C15H12ClN	065753-76-6	9
5		CP 55,940	376	C24H40O3	083003-12-7	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112217\
Data File : BF100838.D
Acq On : 22 Nov 2017 22:08
Operator : SJ/JU
Sample : I6043-19
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
N-T0A-BAJA-S001-0001-03

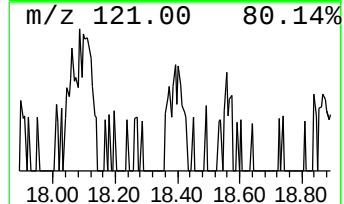
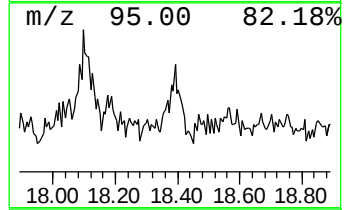
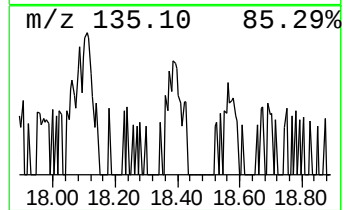
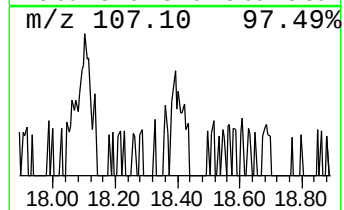
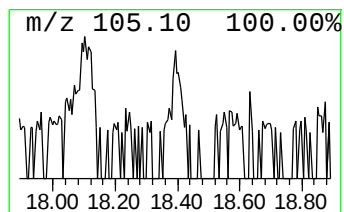
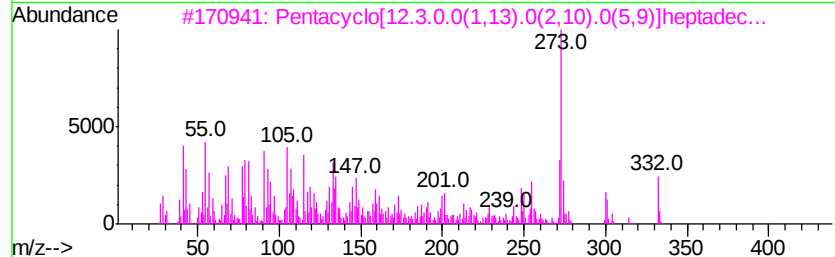
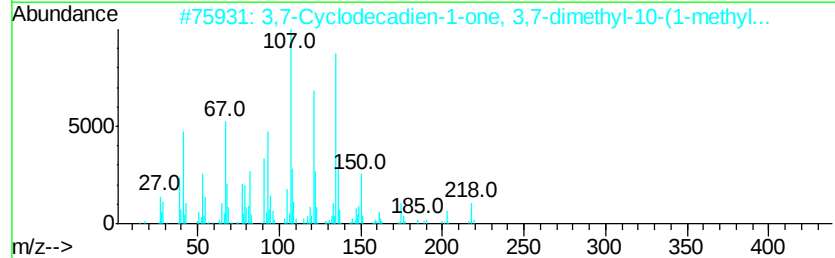
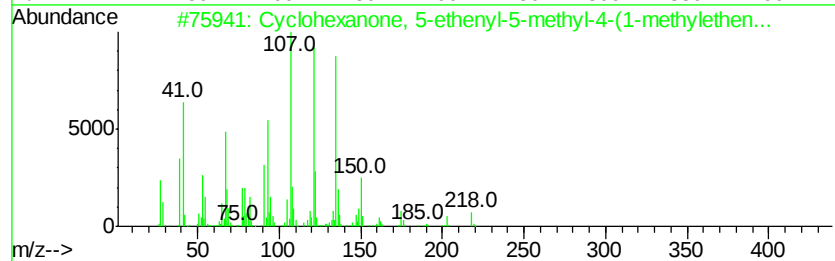
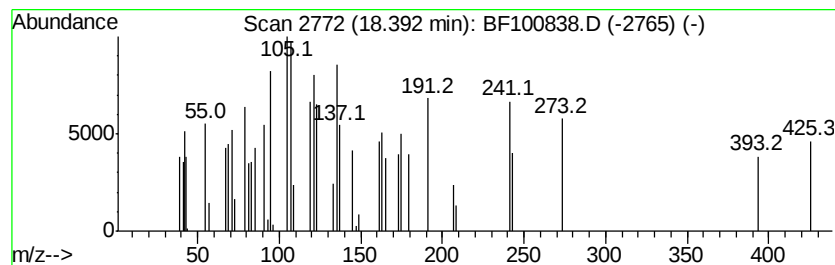
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown18.39 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.39	2.05 ng	46650	Perylene-d12	15.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexanone, 5-ethenyl-5-methyl...	218	C15H22O	032663-57-3	35
2		3,7-Cyclodecadien-1-one, 3,7-dim...	218	C15H22O	006902-91-6	25
3		Pentacyclo[12.3.0.0(1,13).0(2,10)...]	332	C20H28O4	1000192-08-1	23
4		1-Cyclohexene, 1,3,3-trimethyl-2...	206	C14H22O	1000197-08-4	17
5		.beta.-Elemenone	218	C15H22O	020303-60-0	17



Data Path : Z:\HPCHEM1\BNA F\DATA\BF112217\
Data File : BF100838.D
Acq On : 22 Nov 2017 22:08
Operator : SJ/JU
Sample : I6043-19
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
N-T0A-BAJA-S001-0001-03

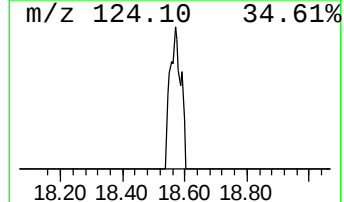
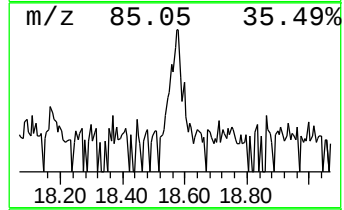
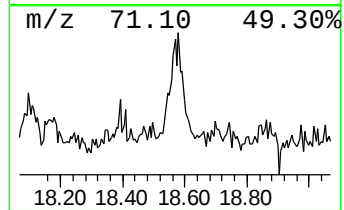
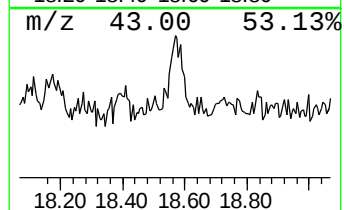
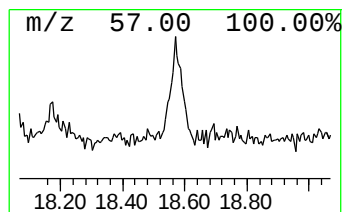
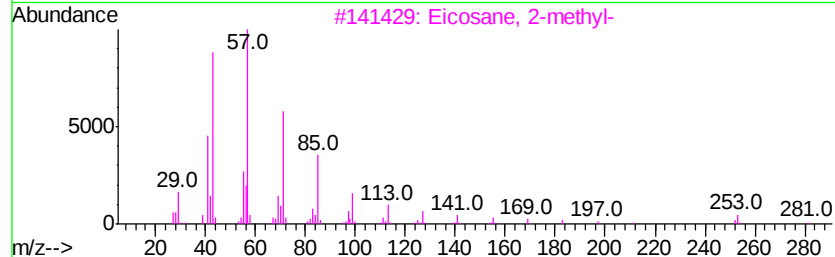
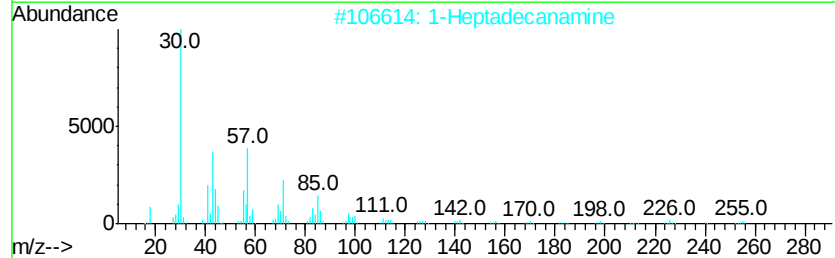
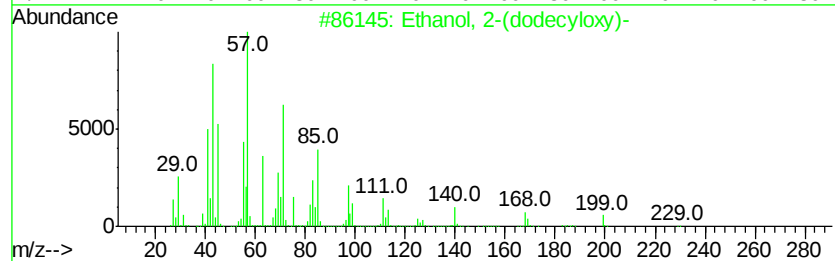
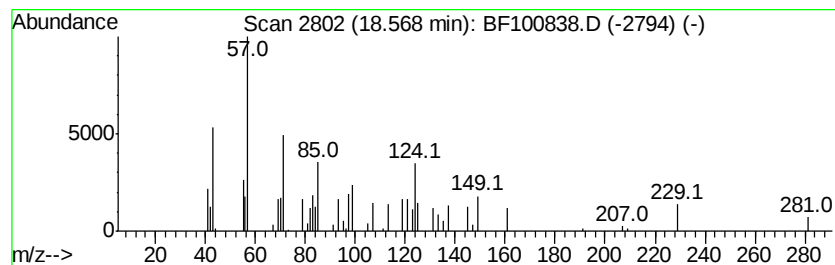
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 unknown18.57 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.57	2.44 ng	55539	Perylene-d12	15.49

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(dodecyloxy)-	230	C14H30O2	004536-30-5	43
2			1-Heptadecanamine	255	C17H37N	004200-95-7	38
3			Eicosane, 2-methyl-	296	C21H44	001560-84-5	38
4			Pentadecane, 2-methyl-	226	C16H34	001560-93-6	38
5			Nonadecane, 2-methyl-	282	C20H42	001560-86-7	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112217\
 Data File : BF100838.D
 Acq On : 22 Nov 2017 22:08
 Operator : SJ/JU
 Sample : I6043-19
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 N-T0A-BAJA-S001-0001-03

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110817.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...	5.09	14.3	ng	322923	1	6.86	450926	20.0
unknown6.63	6.63	98.3	ng	2217120	1	6.86	450926	20.0
2-Hydroxy-iso-but...	8.69	5.1	ng	156772	2	8.14	611716	20.0
Benzophenone	10.61	24.1	ng	808114	3	9.90	669113	20.0
2-Propenoic acid,...	10.99	5.0	ng	159975	4	11.38	636611	20.0
1-Ethyl-3-(pyrrol...	11.04	5.0	ng	158505	4	11.38	636611	20.0
Tridecanoic acid	11.90	5.7	ng	180451	4	11.38	636611	20.0
Octadecanoic acid	12.69	4.1	ng	130145	4	11.38	636611	20.0
Heptafluorobutyri...	13.85	3.7	ng	72414	5	14.02	389313	20.0
unknown16.44	16.44	2.0	ng	46810	6	15.49	455619	20.0
unknown17.53	17.53	3.7	ng	84041	6	15.49	455619	20.0
unknown18.11	18.11	2.4	ng	55349	6	15.49	455619	20.0
unknown18.39	18.39	2.0	ng	46650	6	15.49	455619	20.0
unknown18.57	18.57	2.4	ng	55539	6	15.49	455619	20.0