

Data Path : Z:\HPCHEM1\BNA F\DATA\BF060217\
 Data File : BF095647.D
 Acq On : 2 Jun 2017 22:13
 Operator : SJ/MA
 Sample : I3440-01 5X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-02-RO-4146-060117

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	5.481	642	645	672	rBV2	72120	268596	9.36%	1.692%
2	7.122	921	924	936	rBV2	71023	216951	7.56%	1.367%
3	7.216	936	940	955	rVB	163579	391051	13.63%	2.464%
4	7.539	991	995	1006	rBV	326031	425809	14.85%	2.683%
5	7.775	1031	1035	1050	rBV	180821	279503	9.75%	1.761%
6	8.469	1149	1153	1166	rBV	54951	137557	4.80%	0.867%
7	9.569	1336	1340	1352	rBV	394011	535185	18.66%	3.372%
8	9.651	1352	1354	1365	rVB4	29107	48273	1.68%	0.304%
9	10.645	1517	1523	1529	rVB2	32081	45793	1.60%	0.289%
10	10.892	1559	1565	1570	rBV6	15752	29442	1.03%	0.185%
11	11.345	1638	1642	1654	rVB	274736	355261	12.39%	2.238%
12	11.563	1674	1679	1687	rVB3	39292	60659	2.11%	0.382%
13	12.045	1757	1761	1765	rBV	35615	56345	1.96%	0.355%
14	12.174	1779	1783	1789	rBV2	29229	41223	1.44%	0.260%
15	12.392	1816	1820	1825	rBV2	371500	497297	17.34%	3.133%
16	12.751	1875	1881	1887	rBV6	18407	39111	1.36%	0.246%
17	12.945	1911	1914	1920	rVB5	21535	29774	1.04%	0.188%
18	13.239	1958	1964	1970	rBV	49868	63273	2.21%	0.399%
19	13.298	1970	1974	1980	rVV	42813	60223	2.10%	0.379%
20	13.704	2040	2043	2052	rBV2	79878	129170	4.50%	0.814%
21	14.010	2092	2095	2098	rBV	40453	52204	1.82%	0.329%
22	14.745	2215	2220	2224	rBV	37433	47444	1.65%	0.299%
23	14.798	2224	2229	2232	rBV	345627	424291	14.79%	2.673%
24	14.833	2232	2235	2243	rVB	309116	402607	14.04%	2.537%
25	14.927	2246	2251	2262	rVB	82909	132478	4.62%	0.835%
26	15.415	2329	2334	2337	rBV2	28906	37632	1.31%	0.237%
27	15.568	2355	2360	2364	rBV	736790	853517	29.76%	5.378%
28	15.639	2368	2372	2380	rVV	56767	80453	2.81%	0.507%
29	15.710	2380	2384	2386	rVV	22977	29564	1.03%	0.186%
30	15.745	2386	2390	2394	rVV2	73796	108419	3.78%	0.683%
31	15.786	2394	2397	2404	rVB2	57765	89059	3.11%	0.561%
32	16.015	2432	2436	2438	rBV3	26552	31940	1.11%	0.201%
33	16.039	2438	2440	2445	rVB	26519	31012	1.08%	0.195%
34	16.398	2493	2501	2504	rBV	42723	65458	2.28%	0.412%

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35	16.439	2505	2508	2512	rVB4	19504	31037	1.08%	0.196%
36	16.545	2518	2526	2528	rBV	2294551	2868135	100.00%	18.070%
37	16.574	2528	2531	2534	rVV	1869263	2098109	73.15%	13.219%
38	16.598	2534	2535	2541	rVB	646975	527418	18.39%	3.323%
39	16.698	2548	2552	2556	rBV	362629	369299	12.88%	2.327%
40	16.739	2556	2559	2562	rBV3	62658	79291	2.76%	0.500%
41	16.768	2562	2564	2567	rVB	76639	73310	2.56%	0.462%
42	16.904	2581	2587	2594	rBV	583198	685480	23.90%	4.319%
43	17.015	2603	2606	2611	rVB5	36814	46920	1.64%	0.296%
44	17.145	2624	2628	2635	rVB	257233	293072	10.22%	1.846%
45	17.233	2636	2643	2647	rBV4	34103	66993	2.34%	0.422%
46	17.339	2657	2661	2664	rBV4	53468	77811	2.71%	0.490%
47	17.374	2664	2667	2672	rVB	66131	86116	3.00%	0.543%
48	17.456	2678	2681	2685	rBV	57737	79957	2.79%	0.504%
49	17.504	2685	2689	2693	rVV3	43212	60659	2.11%	0.382%
50	17.621	2706	2709	2713	rVB	29595	29557	1.03%	0.186%
51	17.662	2713	2716	2720	rVB	58851	63323	2.21%	0.399%
52	18.062	2782	2784	2791	rVB3	32117	37777	1.32%	0.238%
53	18.203	2805	2808	2811	rVB2	61445	62109	2.17%	0.391%
54	18.239	2811	2814	2817	rBV	34238	30979	1.08%	0.195%
55	18.492	2852	2857	2861	rBV2	515673	799559	27.88%	5.038%
56	18.527	2861	2863	2870	rVB2	221621	307577	10.72%	1.938%
57	18.627	2877	2880	2883	rVB2	50273	52701	1.84%	0.332%
58	19.003	2942	2944	2948	rVB2	41340	39727	1.39%	0.250%
59	19.768	3070	3074	3077	rBV	233194	309222	10.78%	1.948%
60	20.062	3121	3124	3128	rVB	114016	119698	4.17%	0.754%
61	20.115	3131	3133	3137	rBV	178087	191018	6.66%	1.203%
62	20.174	3139	3143	3146	rBV	252189	288513	10.06%	1.818%

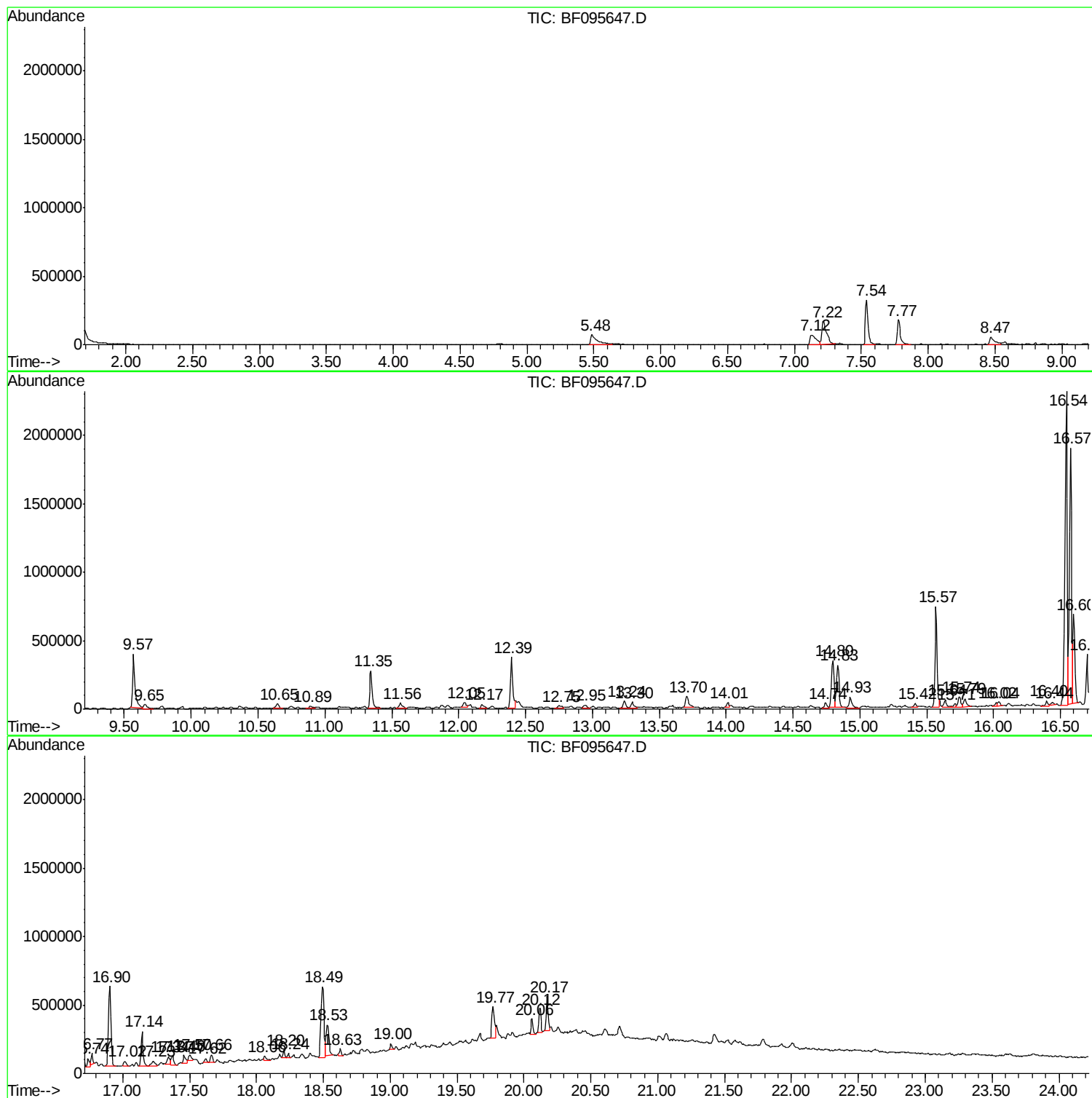
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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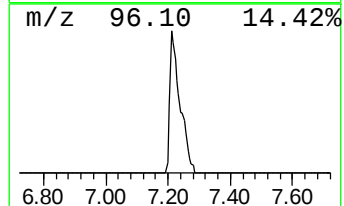
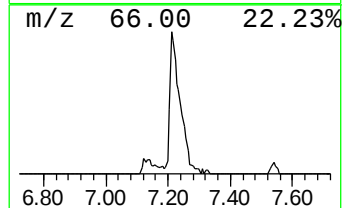
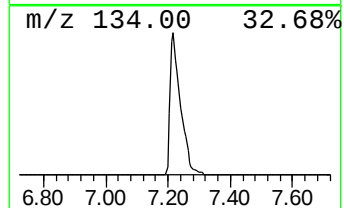
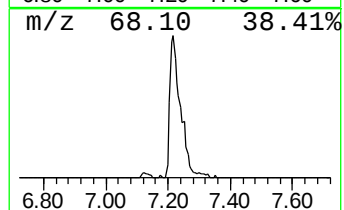
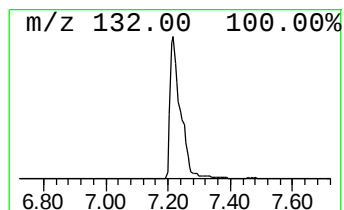
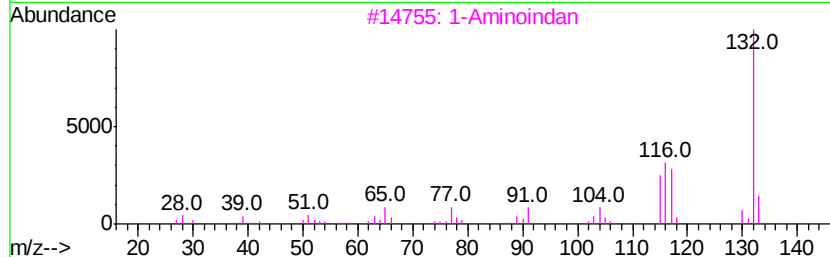
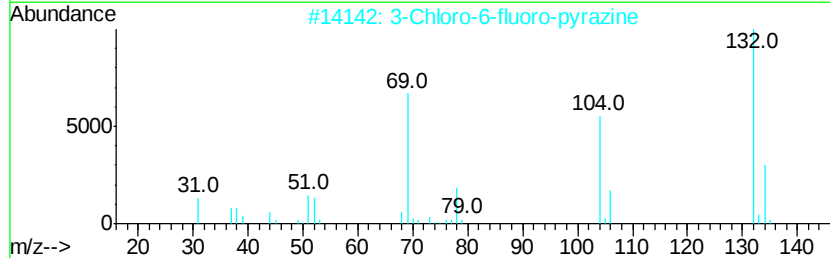
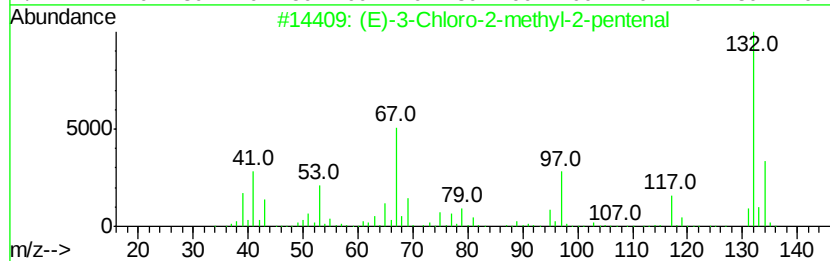
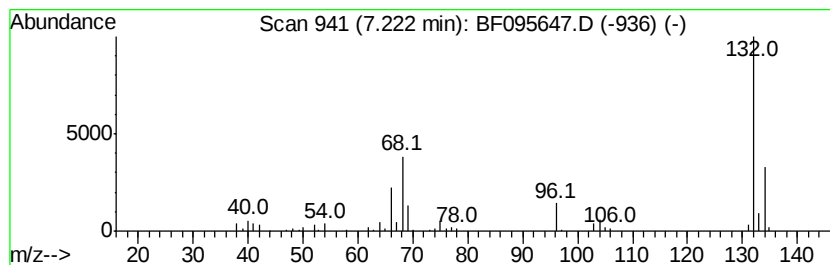
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown7.22 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.22	18.37 ng	391051	1,4-Dichlorobenzene-d4	7.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
3		1-Aminoindan	133	C9H11N	034698-41-4	9
4		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9
5		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9



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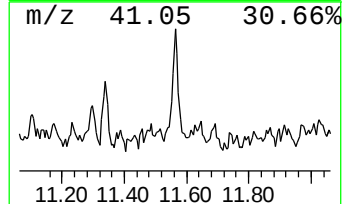
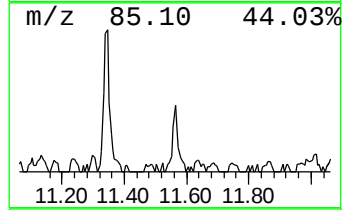
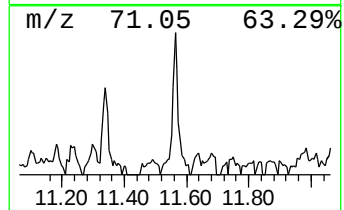
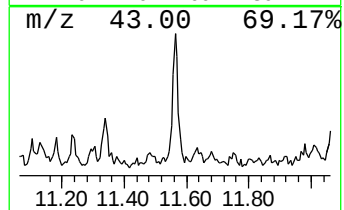
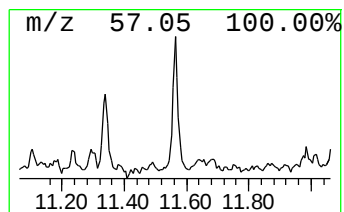
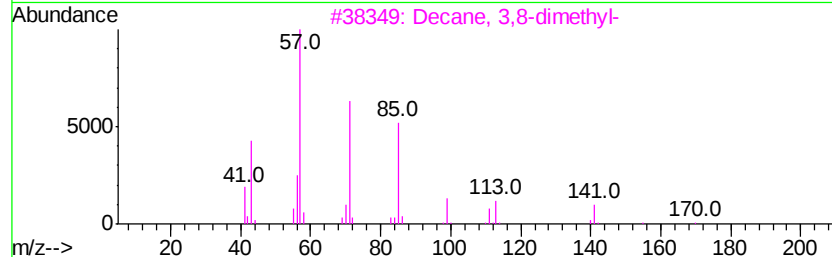
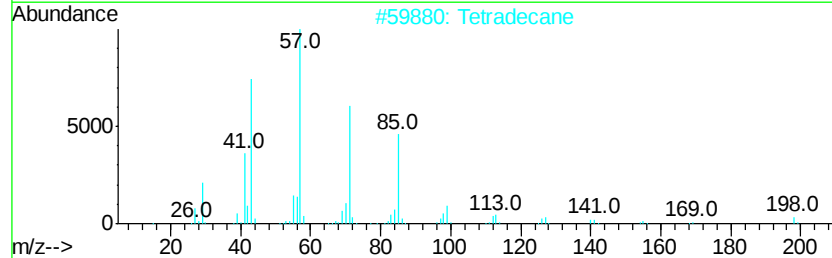
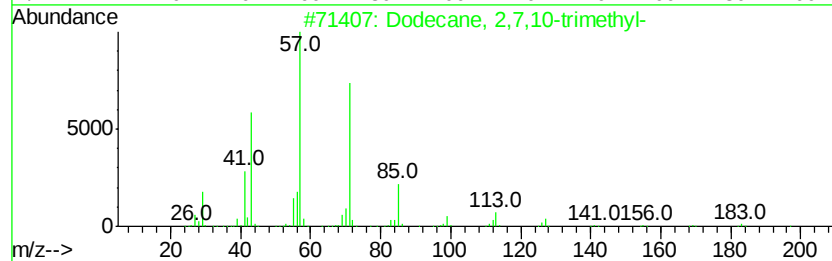
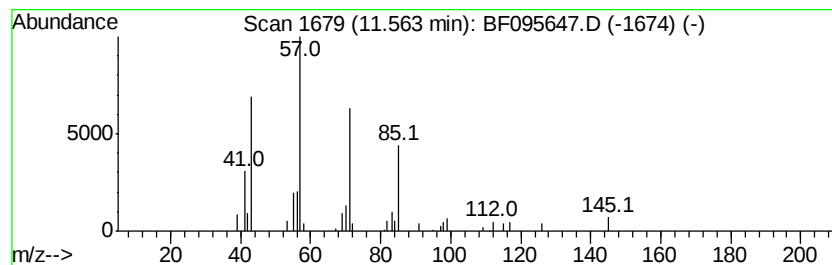
Instrument :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Dodecane, 2,7,10-trimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.56	2.44 ng	60659	Acenaphthene-d10		12.39	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane,	2,7,10-trimethyl-	212	C15H32	074645-98-0	64
2	Tetradecane		198	C14H30	000629-59-4	64
3	Decane,	3,8-dimethyl-	170	C12H26	017312-55-9	64
4	Decane,	2,3,5-trimethyl-	184	C13H28	062238-11-3	59
5	Tridecane		184	C13H28	000629-50-5	53



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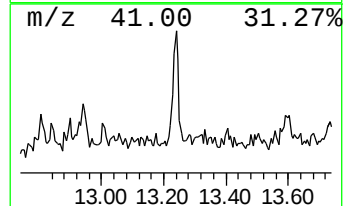
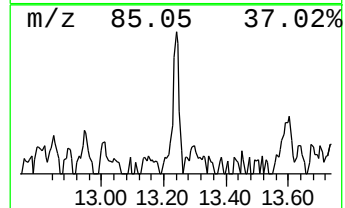
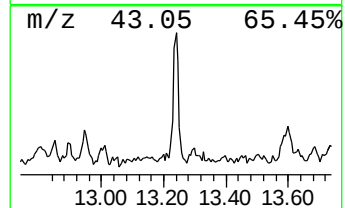
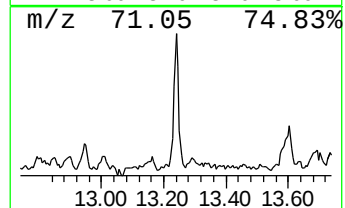
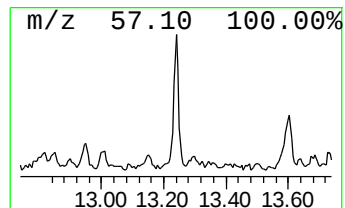
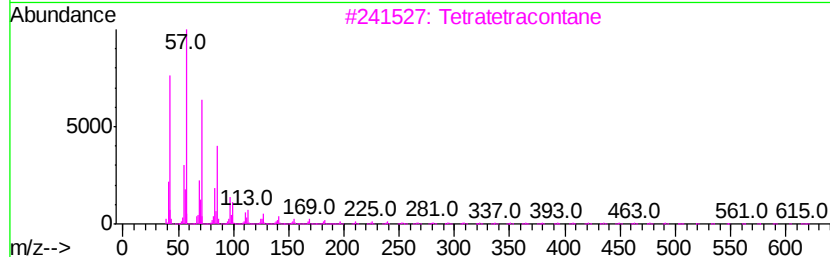
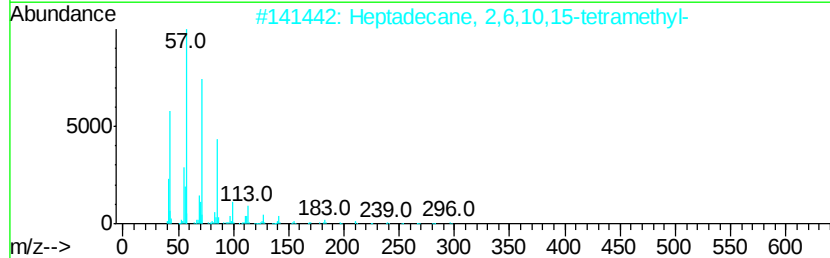
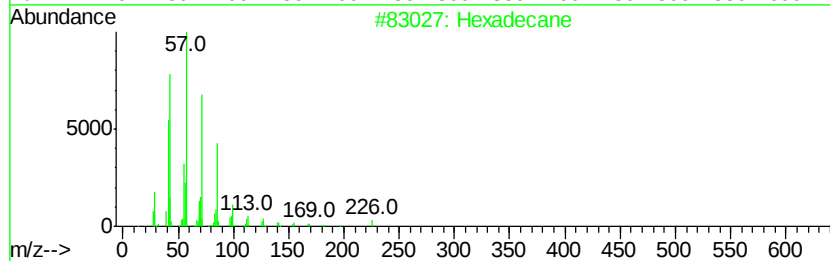
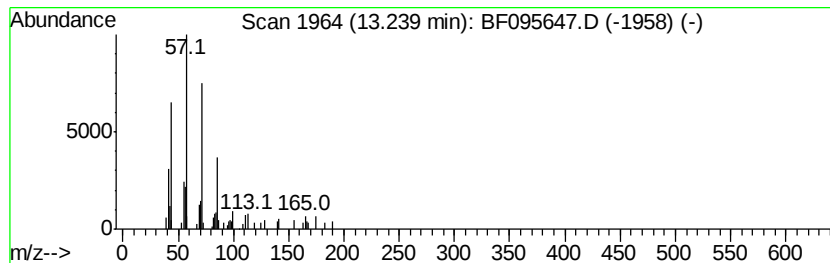
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Hexadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.24	2.54 ng	63273	Acenaphthene-d10	12.39
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Hexadecane	226 C16H34	000544-76-3	83
2	Heptadecane, 2,6,10,15-tetramethyl-	296 C21H44	054833-48-6	80
3	Tetratetracontane	619 C44H90	007098-22-8	80
4	Hexadecane, 2,6,11,15-tetramethyl-	282 C20H42	000504-44-9	80
5	Octane, 3-ethyl-2,7-dimethyl-	170 C12H26	062183-55-5	72



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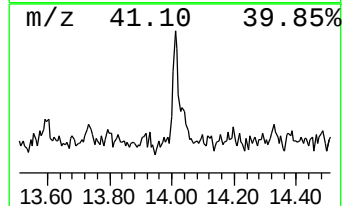
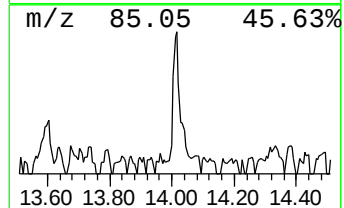
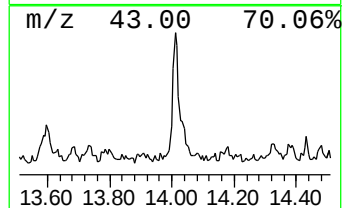
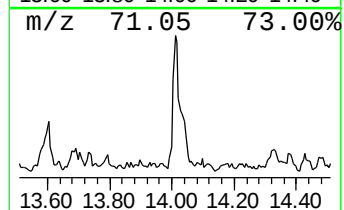
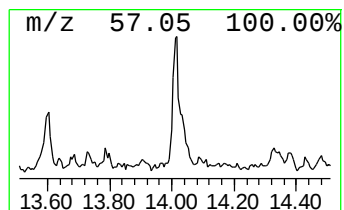
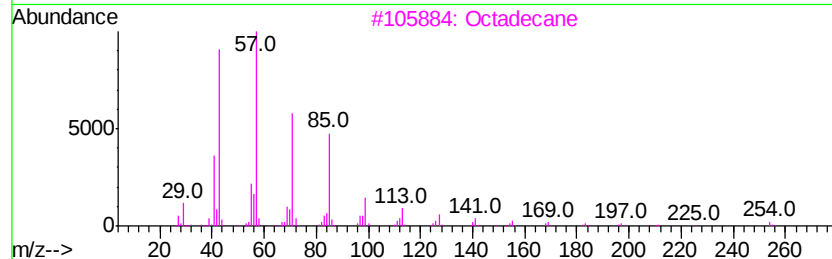
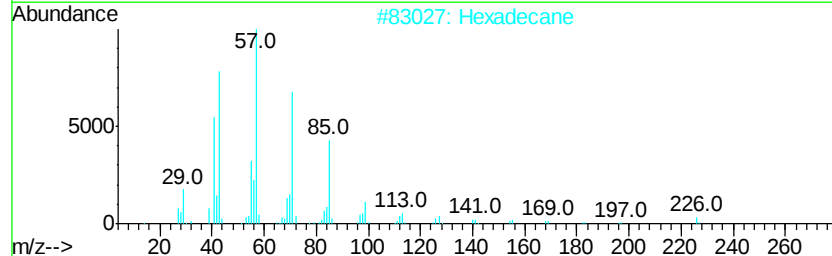
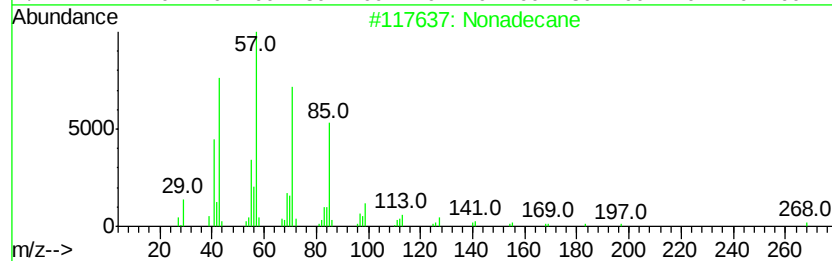
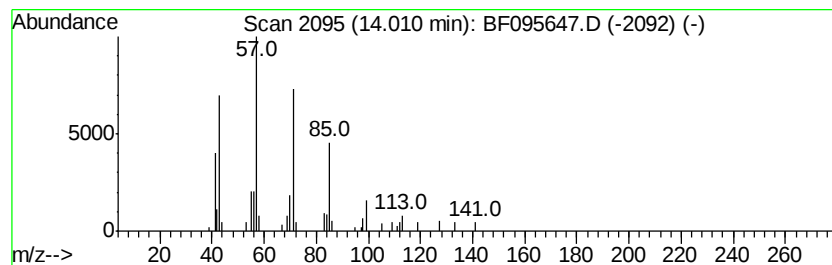
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Nonadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.01	2.46 ng	52204	Phenanthrene-d10	14.80

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane		268	C19H40	000629-92-5	87
2	Hexadecane		226	C16H34	000544-76-3	86
3	Octadecane		254	C18H38	000593-45-3	80
4	Tetradecane		198	C14H30	000629-59-4	80
5	Heptadecane		240	C17H36	000629-78-7	80



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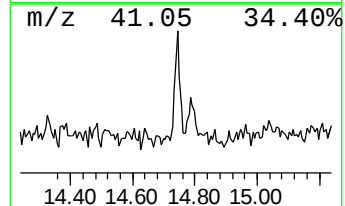
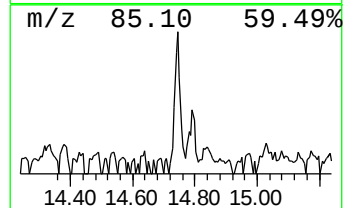
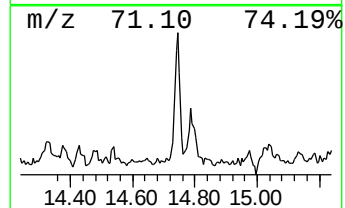
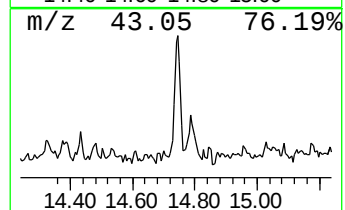
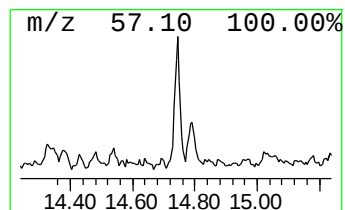
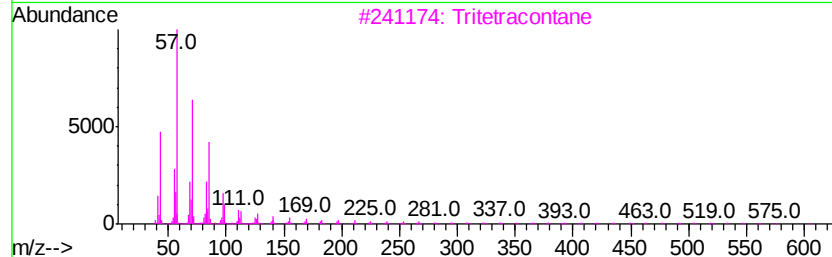
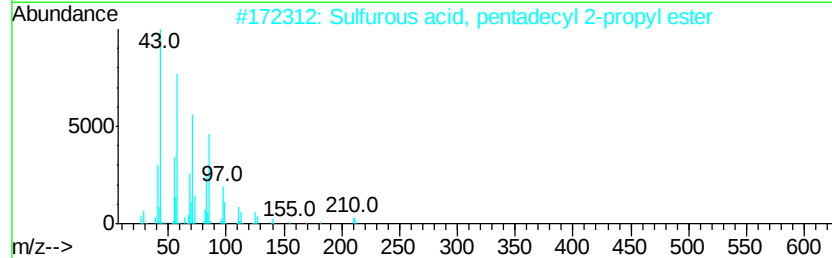
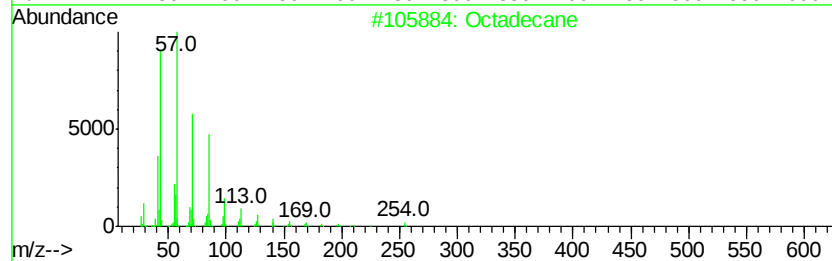
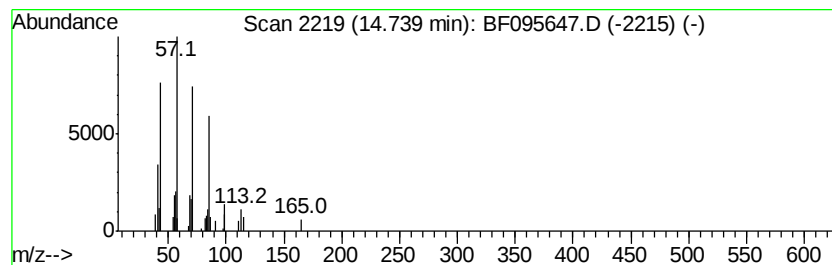
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Octadecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.74	2.24 ng	47444	Phenanthrene-d10	14.80

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane		254	C18H38	000593-45-3	91
2	Sulfurous acid, pentadecyl 2-pro...		334	C18H38O3S	1000309-12-6	86
3	Tritetracontane		605	C43H88	007098-21-7	86
4	Heptadecane		240	C17H36	000629-78-7	80
5	Hexadecane		226	C16H34	000544-76-3	78



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060217\
Data File : BF095647.D
Acq On : 2 Jun 2017 22:13
Operator : SJ/MA
Sample : I3440-01 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
EO-02-RO-4146-060117

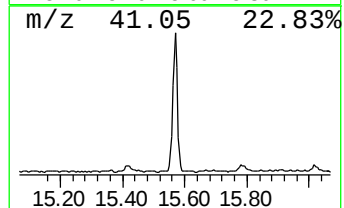
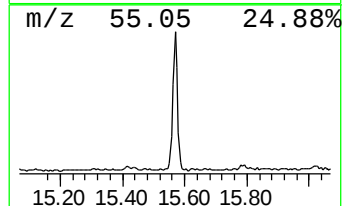
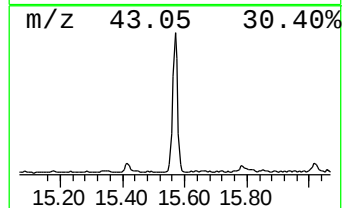
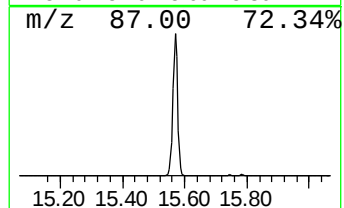
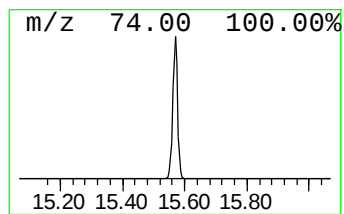
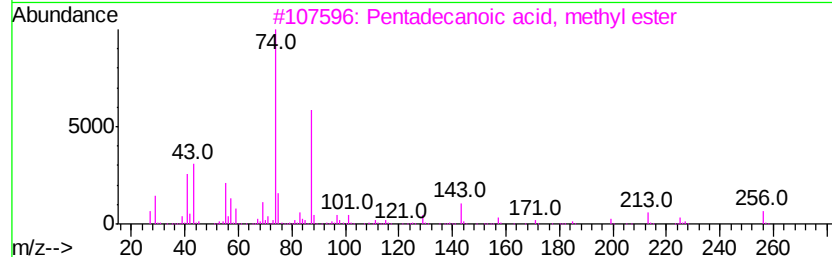
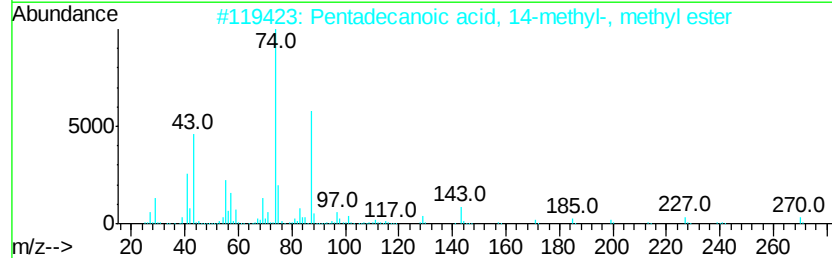
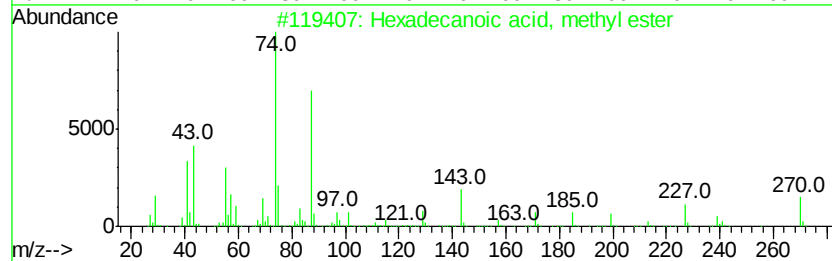
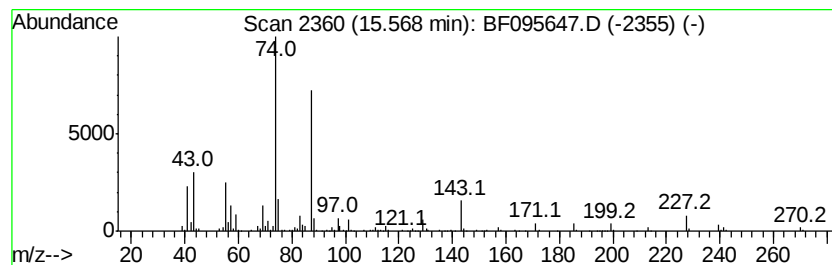
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Hexadecanoic acid, methyl e... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.57	40.23 ng	853517	Phenanthrene-d10	14.80

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	96
2		Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	96
3		Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	90
4		Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	87
5		Methyl tetradecanoate	242	C15H30O2	000124-10-7	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060217\
Data File : BF095647.D
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Operator : SJ/MA
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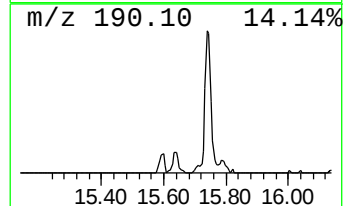
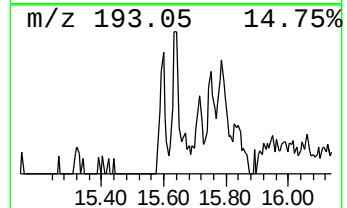
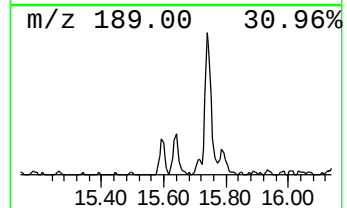
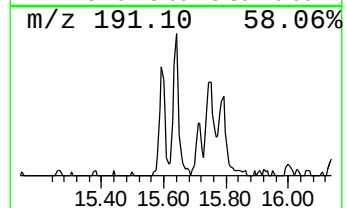
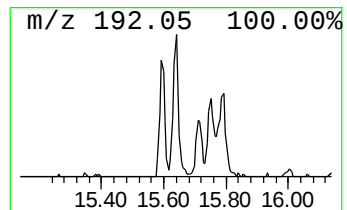
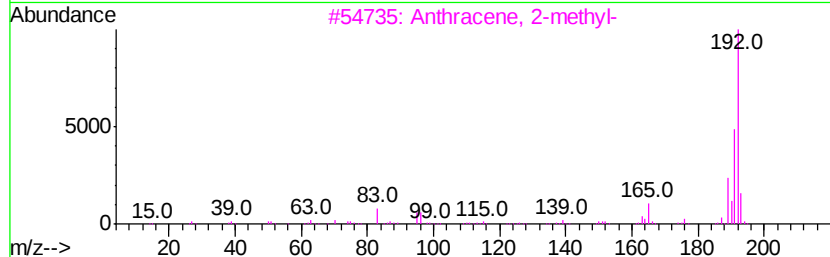
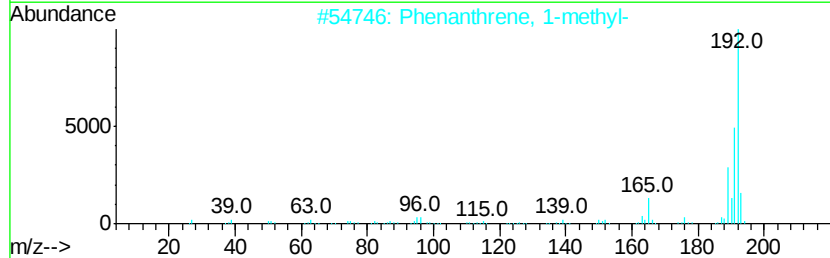
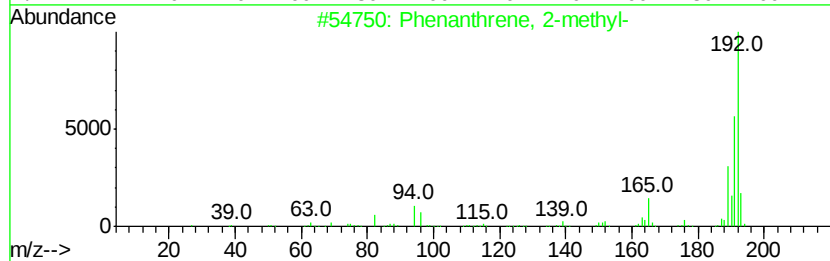
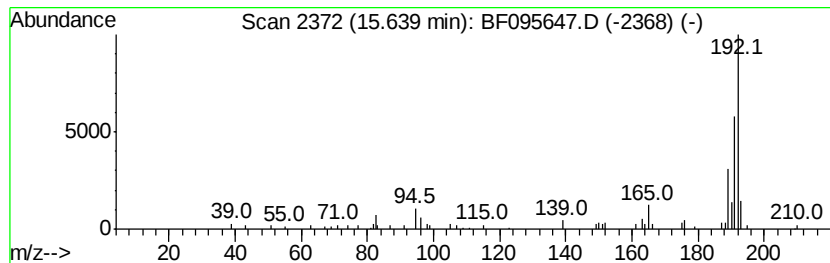
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Phenanthrene, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.64	3.79 ng	80453	Phenanthrene-d10	14.80

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
2		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
3		Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
4		Anthracene, 1-methyl-	192	C15H12	000610-48-0	92
5		Phenanthrene, 3-methyl-	192	C15H12	000832-71-3	92



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060217\
Data File : BF095647.D
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Operator : SJ/MA
Sample : I3440-01 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

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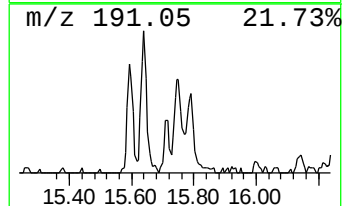
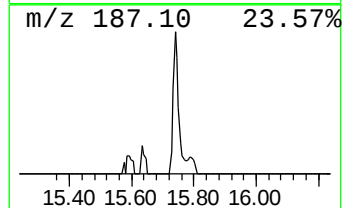
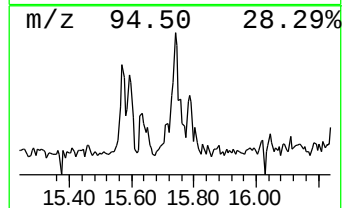
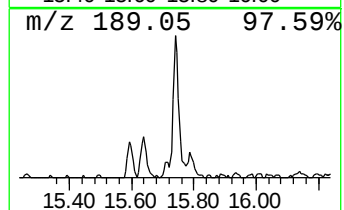
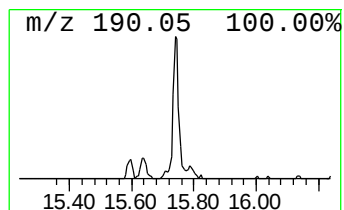
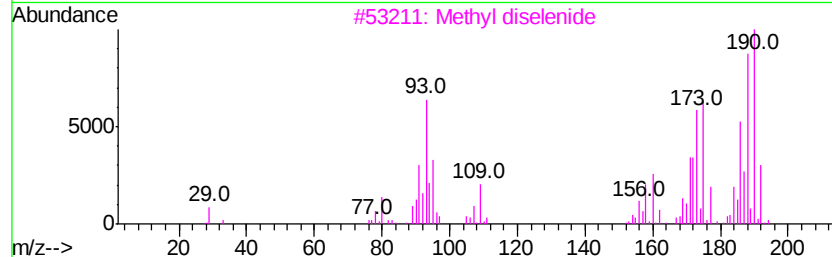
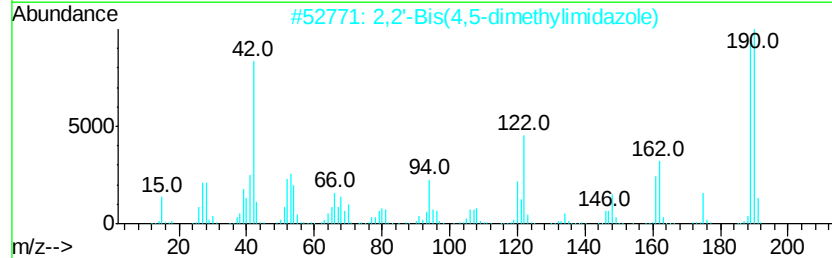
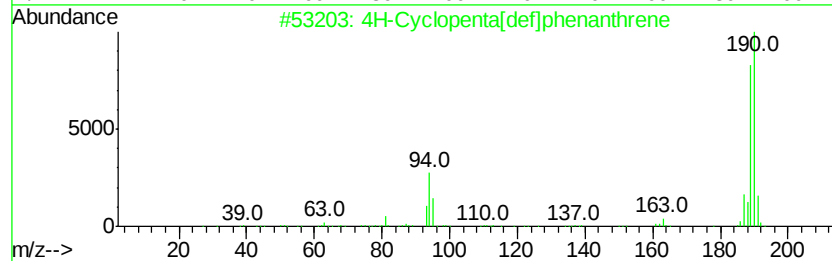
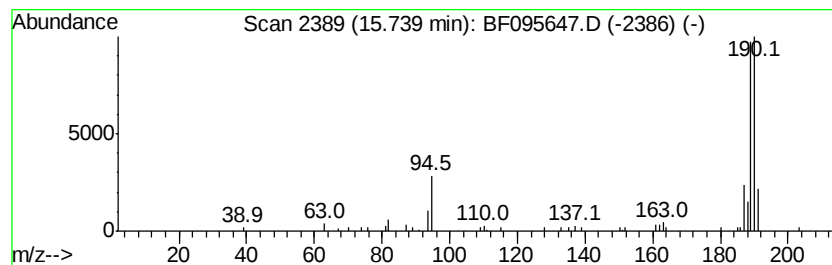
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 4H-Cyclopenta[def]phenanthrene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.74	5.11 ng	108419	Phenanthrene-d10	14.80

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	87
2		2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	45
3		Methyl diselenide	190	C2H6Se2	007101-31-7	43
4		6H-Cyclobuta[flk]phenanthrene	190	C15H10	083469-43-6	23
5		1,1'-Biphenyl, 4,4'-difluoro-	190	C12H8F2	000398-23-2	17



Data Path : Z:\HPCHEM1\BNA F\DATA\BF060217\
Data File : BF095647.D
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Sample : I3440-01 5X
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ClientSampleId :
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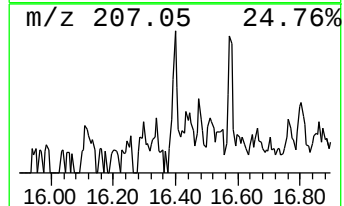
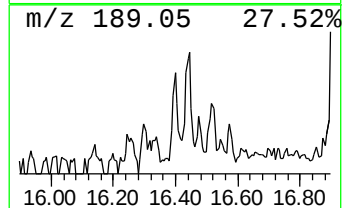
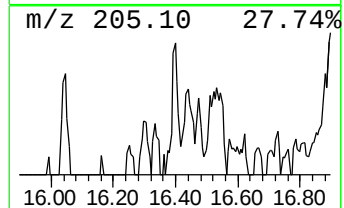
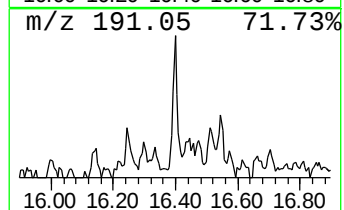
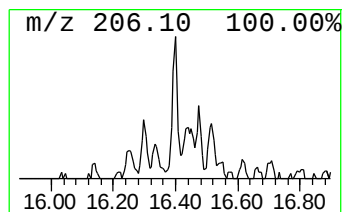
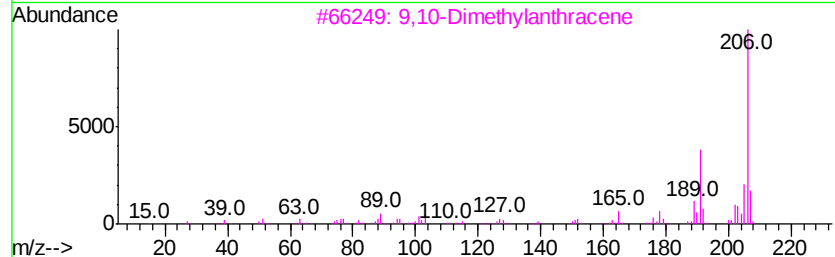
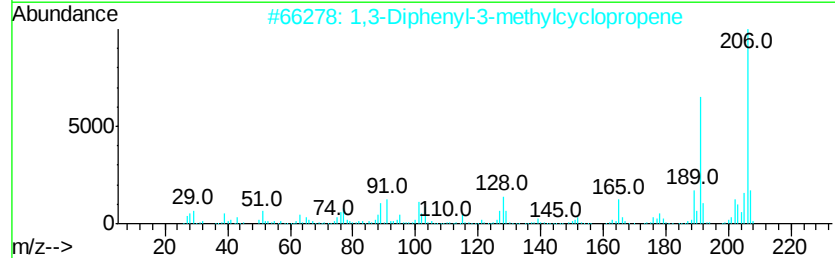
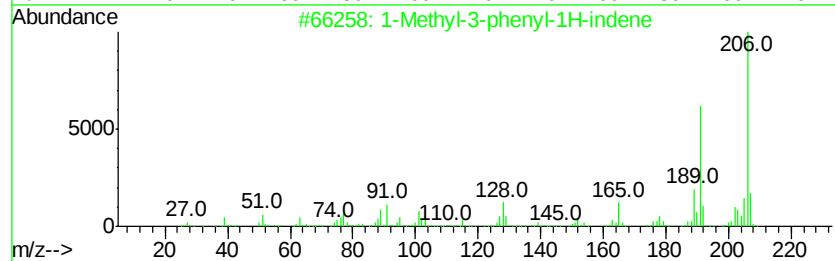
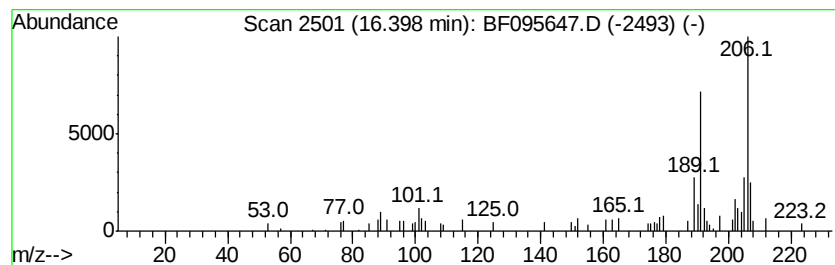
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 1-Methyl-3-phenyl-1H-indene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.40	3.09 ng	65458	Phenanthrene-d10	14.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Methyl-3-phenyl-1H-indene	206	C16H14	022360-62-9	93
2		1,3-Diphenyl-3-methylcyclopropene	206	C16H14	086544-79-8	91
3		9,10-Dimethylanthracene	206	C16H14	000781-43-1	90
4		Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	89
5		3-Methyl-1-phenyl-1H-indene	206	C16H14	022360-63-0	87



Data Path : Z:\HPCHEM1\BNA F\DATA\BF060217\
Data File : BF095647.D
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Operator : SJ/MA
Sample : I3440-01 5X
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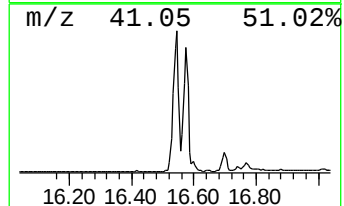
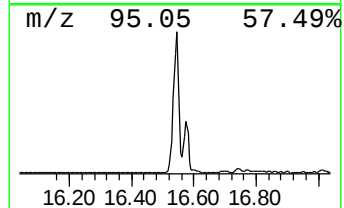
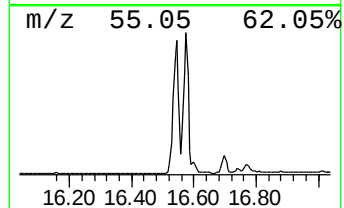
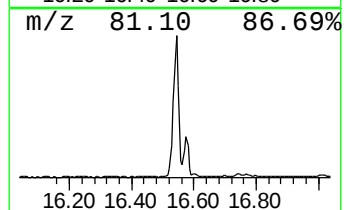
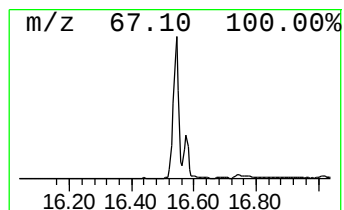
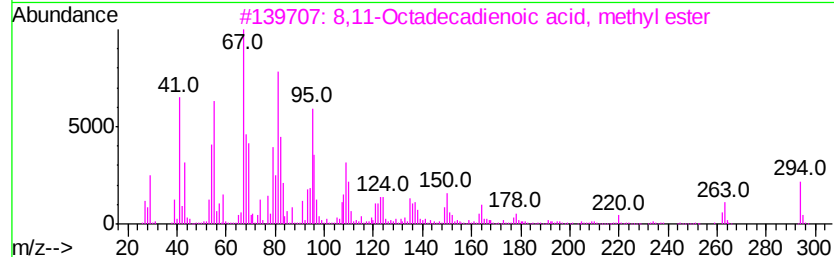
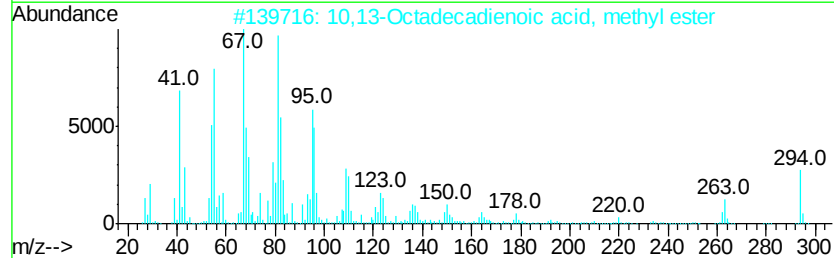
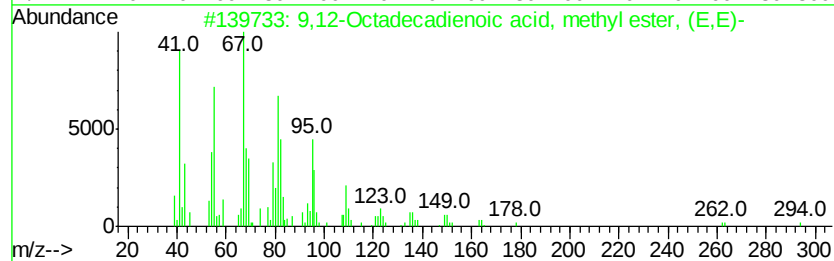
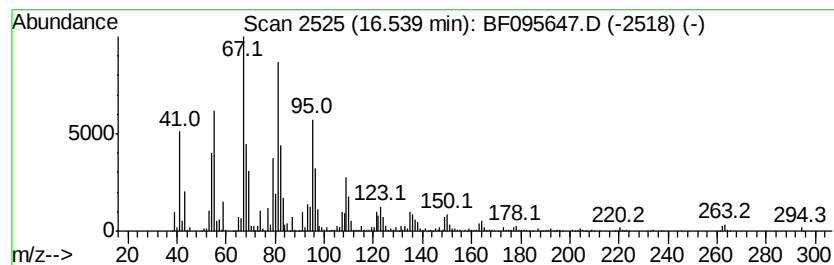
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 9,12-Octadecadienoic acid, ... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.54	135.20 ng	2868140	Phenanthrene-d10	14.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	99
2		10,13-Octadecadienoic acid, meth...	294	C19H34O2	056554-62-2	99
3		8,11-Octadecadienoic acid, methv...	294	C19H34O2	056599-58-7	99
4		9,15-Octadecadienoic acid, methy...	294	C19H34O2	017309-05-6	99
5		9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	99



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060217\
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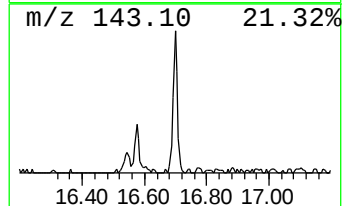
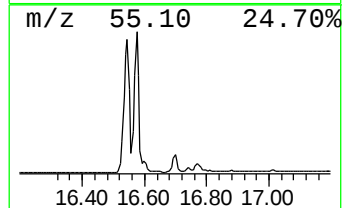
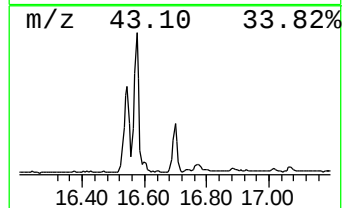
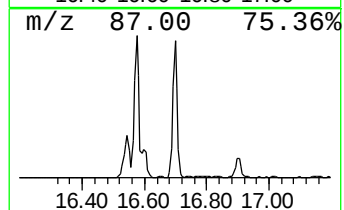
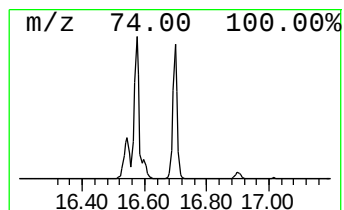
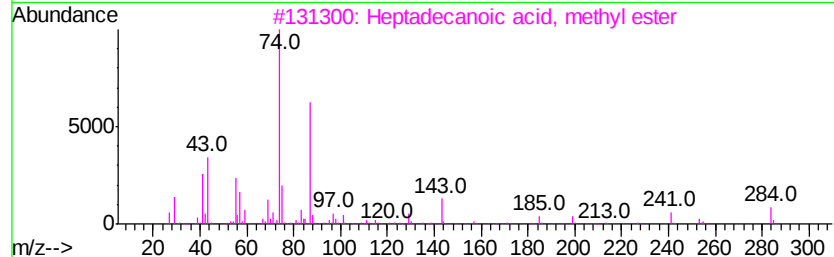
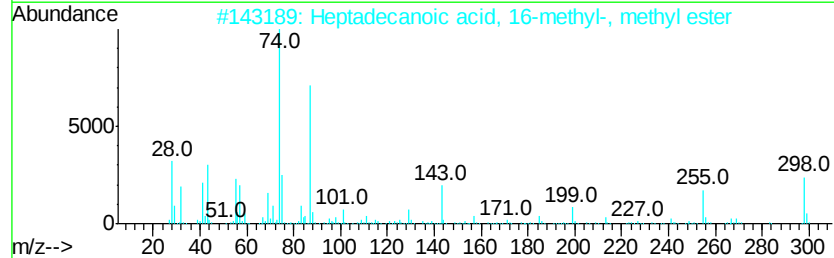
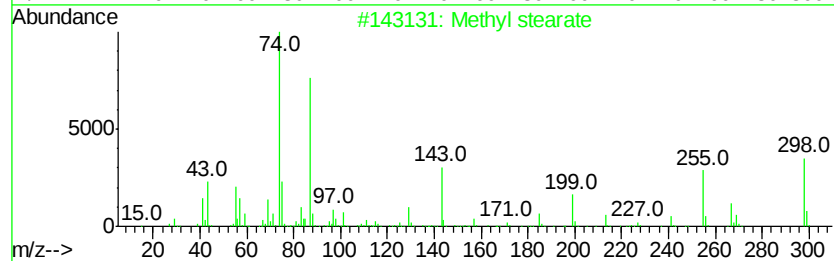
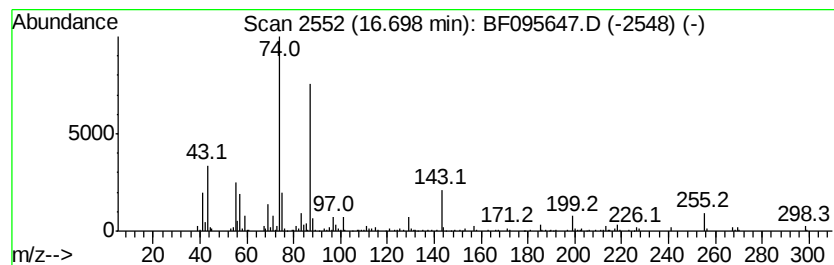
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Methyl stearate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.70	9.24 ng	369299	Chrysene-d12	18.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Methyl stearate	298	C19H38O2	000112-61-8	98
2		Heptadecanoic acid, 16-methyl-, ...	298	C19H38O2	005129-61-3	95
3		Heptadecanoic acid, methyl ester	284	C18H36O2	001731-92-6	86
4		Hexadecanoic acid, 15-methyl-, m...	284	C18H36O2	006929-04-0	83
5		Methyl tetradecanoate	242	C15H30O2	000124-10-7	83



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060217\
Data File : BF095647.D
Acq On : 2 Jun 2017 22:13
Operator : SJ/MA
Sample : I3440-01 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
EO-02-RO-4146-060117

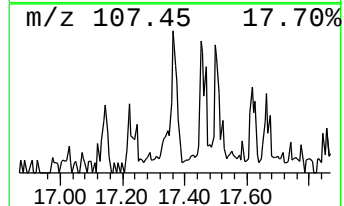
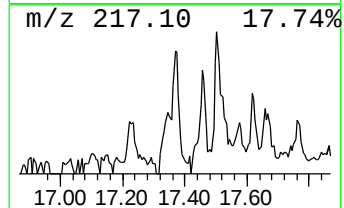
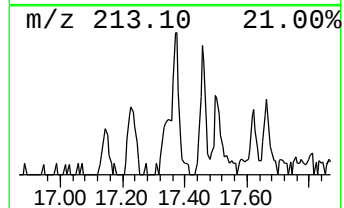
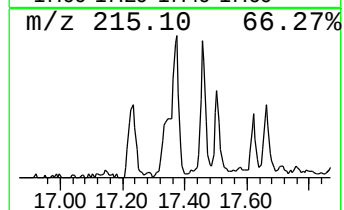
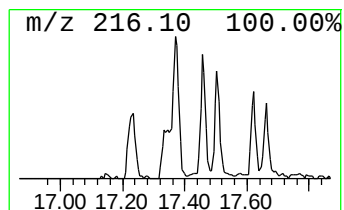
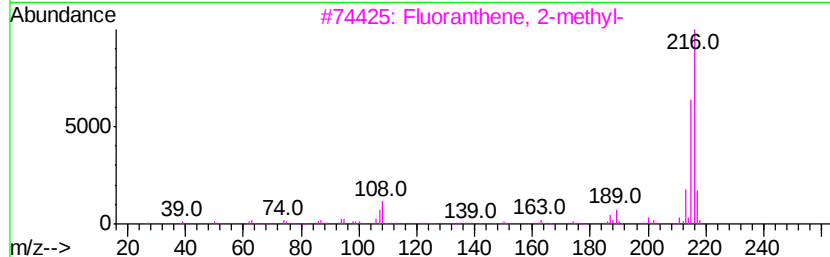
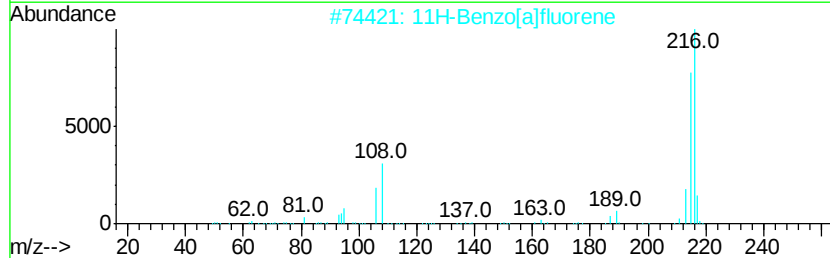
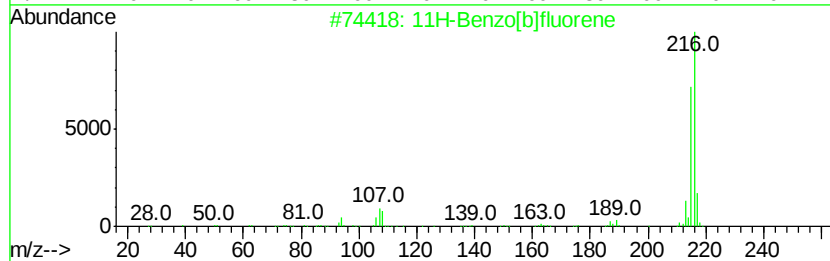
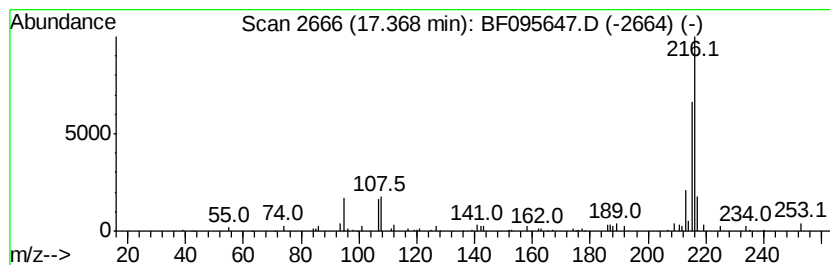
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 11H-Benzo[b]fluorene Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.37	2.15 ng	86116	Chrysene-d12	18.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	87
2		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	72
3		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	72
4		Pvrene, 1-methyl-	216	C17H12	002381-21-7	64
5		6H-Thiazolo[5,4-e]indole, 2,7,8-...	216	C12H12N2S	1000338-32-0	64



Data Path : Z:\HPCHEM1\BNA F\DATA\BF060217\
Data File : BF095647.D
Acq On : 2 Jun 2017 22:13
Operator : SJ/MA
Sample : I3440-01 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
EO-02-RO-4146-060117

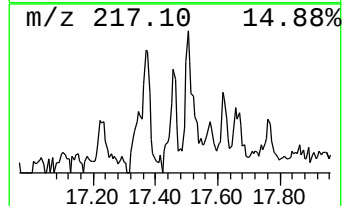
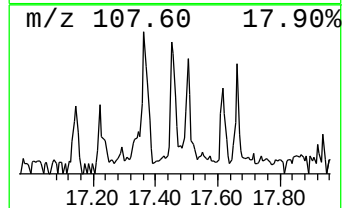
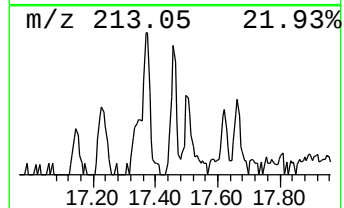
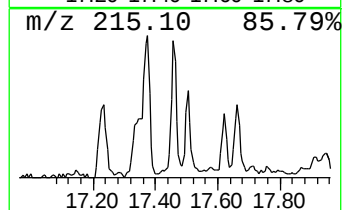
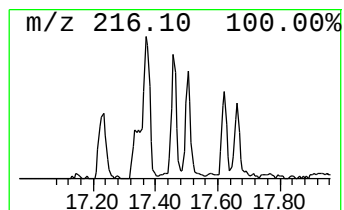
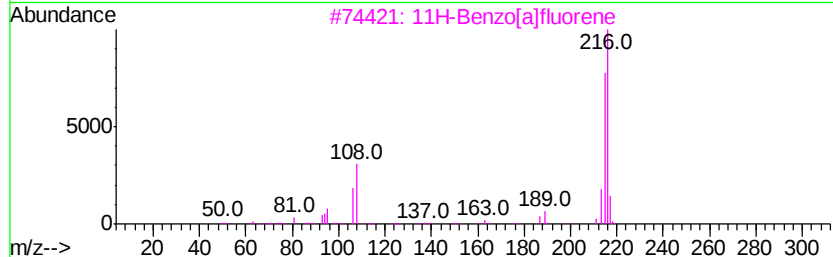
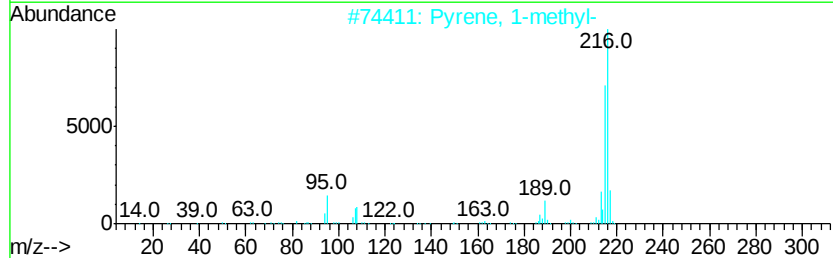
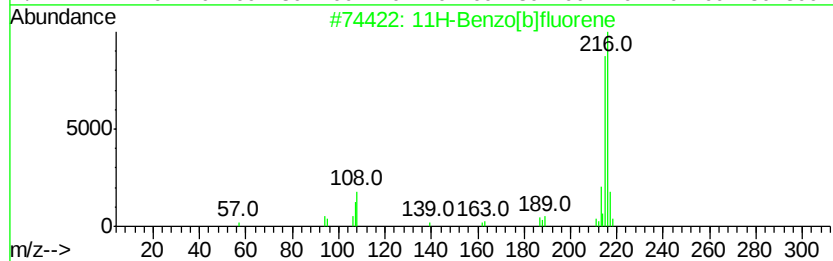
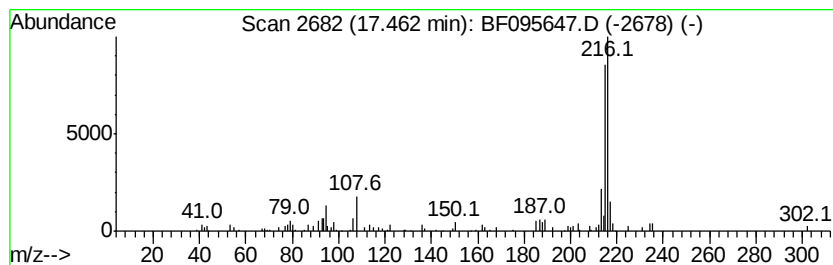
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 11H-Benzo[a]fluorene Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.46	2.00 ng	79957	Chrysene-d12	18.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	89
2		Pyrene, 1-methyl-	216	C17H12	002381-21-7	81
3		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	76
4		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	68
5		7H-Benzo[c]fluorene	216	C17H12	000205-12-9	58



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060217\
 Data File : BF095647.D
 Acq On : 2 Jun 2017 22:13
 Operator : SJ/MA
 Sample : I3440-01 5X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-02-RO-4146-060117

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
unknown7.22	7.22	18.4	ng	391051	1	7.54	425809	20.0
Dodecane, 2,7,10-...	11.56	2.4	ng	60659	3	12.39	497297	20.0
Hexadecane	13.24	2.5	ng	63273	3	12.39	497297	20.0
Nonadecane	14.01	2.5	ng	52204	4	14.80	424291	20.0
Octadecane	14.74	2.2	ng	47444	4	14.80	424291	20.0
Hexadecanoic acid...	15.57	40.2	ng	853517	4	14.80	424291	20.0
Phenanthrene, 2-m...	15.64	3.8	ng	80453	4	14.80	424291	20.0
4H-Cyclopenta[def...	15.74	5.1	ng	108419	4	14.80	424291	20.0
1-Methyl-3-phenyl...	16.40	3.1	ng	65458	4	14.80	424291	20.0
9,12-Octadecadien...	16.54	135.2	ng	2868140	4	14.80	424291	20.0
Methyl stearate	16.70	9.2	ng	369299	5	18.50	799559	20.0
11H-Benzo[b]fluorene	17.37	2.1	ng	86116	5	18.50	799559	20.0
11H-Benzo[a]fluorene	17.46	2.0	ng	79957	5	18.50	799559	20.0