

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120617\
 Data File : BF101169.D
 Acq On : 6 Dec 2017 15:57
 Operator : SJ/JU
 Sample : I6303-01 10X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HD-02-120417

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-BF113017.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	5.457	570	573	586	rBB	43992	49314	13.90%	1.500%
2	6.475	743	746	758	rVB	41761	48858	13.77%	1.487%
3	6.616	767	770	777	rBB	66846	60804	17.14%	1.850%
4	6.845	805	809	817	rBB	320124	266258	75.05%	8.101%
5	6.998	832	835	841	rVB	59292	49468	13.94%	1.505%
6	7.410	902	905	911	rBV	33477	34703	9.78%	1.056%
7	8.128	1023	1027	1033	rBV	420856	336969	94.98%	10.252%
8	8.710	1123	1126	1129	rBV	5794	5120	1.44%	0.156%
9	9.198	1206	1209	1214	rBV	84144	69660	19.63%	2.119%
10	9.275	1218	1222	1225	rVB	9851	8609	2.43%	0.262%
11	9.469	1250	1255	1258	rBV3	2962	4326	1.22%	0.132%
12	9.592	1273	1276	1284	rVB3	11013	15047	4.24%	0.458%
13	9.804	1310	1312	1317	rVB2	12875	10082	2.84%	0.307%
14	9.880	1322	1325	1335	rBV	350604	343787	96.90%	10.460%
15	10.080	1358	1359	1364	rVB4	3370	4144	1.17%	0.126%
16	10.128	1364	1367	1371	rBV4	6464	7605	2.14%	0.231%
17	10.304	1388	1397	1399	rBV2	17650	17769	5.01%	0.541%
18	10.522	1428	1434	1436	rBV2	10749	12883	3.63%	0.392%
19	10.639	1453	1454	1457	rVB2	5314	4100	1.16%	0.125%
20	10.675	1457	1460	1464	rBV	25456	24966	7.04%	0.760%
21	10.775	1474	1477	1483	rBV3	23016	31218	8.80%	0.950%
22	10.869	1490	1493	1496	rVB3	3492	4017	1.13%	0.122%
23	10.998	1513	1515	1519	rVB4	4924	5275	1.49%	0.160%
24	11.104	1528	1533	1534	rBV3	3172	4299	1.21%	0.131%
25	11.222	1549	1553	1556	rBV	20245	18197	5.13%	0.554%
26	11.251	1556	1558	1564	rVB3	18421	19954	5.62%	0.607%
27	11.374	1575	1579	1582	rBV	402316	354793	100.00%	10.795%
28	11.492	1597	1599	1603	rVB4	3637	4815	1.36%	0.146%
29	11.533	1603	1606	1608	rBV3	4761	5506	1.55%	0.168%
30	11.604	1615	1618	1620	rBV3	5166	4464	1.26%	0.136%
31	11.651	1623	1626	1628	rVV	25749	18706	5.27%	0.569%
32	11.704	1632	1635	1638	rVB3	5951	6428	1.81%	0.196%
33	11.792	1646	1650	1652	rBV3	5001	4943	1.39%	0.150%
34	11.886	1662	1666	1667	rBV3	10176	9720	2.74%	0.296%

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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-BF113017.M
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35	11.904	1667	1669	1672	rVB3	5978	5924	1.67%	0.180%
36	12.057	1693	1695	1698	rBV	23327	18574	5.24%	0.565%
37	12.104	1701	1703	1708	rVB3	6173	7569	2.13%	0.230%
38	12.204	1716	1720	1723	rVB5	6249	6804	1.92%	0.207%
39	12.280	1728	1733	1735	rBV5	5656	8590	2.42%	0.261%
40	12.298	1735	1736	1740	rVB4	7205	6047	1.70%	0.184%
41	12.345	1741	1744	1747	rBV5	4486	5203	1.47%	0.158%
42	12.421	1755	1757	1759	rBV	8750	9789	2.76%	0.298%
43	12.445	1759	1761	1766	rVB3	34665	34841	9.82%	1.060%
44	12.516	1770	1773	1775	rBV4	4579	4629	1.30%	0.141%
45	12.545	1775	1778	1783	rVB5	14664	16904	4.76%	0.514%
46	12.586	1783	1785	1787	rBV	13184	13037	3.67%	0.397%
47	12.627	1791	1792	1794	rVB2	8051	5316	1.50%	0.162%
48	12.686	1794	1802	1805	rBV7	11616	22943	6.47%	0.698%
49	12.739	1808	1811	1813	rBV4	10012	10646	3.00%	0.324%
50	12.763	1813	1815	1819	rVB5	12045	12242	3.45%	0.372%
51	12.816	1819	1824	1831	rBV4	45408	62764	17.69%	1.910%
52	12.874	1831	1834	1837	rBV2	11748	12924	3.64%	0.393%
53	12.927	1841	1843	1845	rBV2	10077	8998	2.54%	0.274%
54	12.951	1845	1847	1850	rBV	86853	67468	19.02%	2.053%
55	13.051	1862	1864	1867	rBV4	6963	7319	2.06%	0.223%
56	13.133	1875	1878	1882	rBV5	10965	13347	3.76%	0.406%
57	13.174	1882	1885	1887	rBV	32585	28176	7.94%	0.857%
58	13.251	1894	1898	1902	rBV6	10935	21341	6.02%	0.649%
59	13.374	1916	1919	1920	rBV3	7816	8020	2.26%	0.244%
60	13.516	1940	1943	1946	rBV2	37070	34390	9.69%	1.046%
61	13.568	1950	1952	1953	rVB2	10223	5749	1.62%	0.175%
62	13.751	1981	1983	1984	rBV2	8498	6925	1.95%	0.211%
63	13.845	1996	1999	2002	rVB	38873	33203	9.36%	1.010%
64	14.015	2024	2028	2035	rBV	298313	298419	84.11%	9.080%
65	14.163	2049	2053	2055	rBV2	46938	52914	14.91%	1.610%
66	14.468	2102	2105	2108	rVB	45855	35758	10.08%	1.088%
67	14.774	2154	2157	2159	rBV	38917	37235	10.49%	1.133%
68	15.110	2211	2214	2220	rVB3	47705	58132	16.38%	1.769%
69	15.492	2275	2279	2284	rVB	229690	303440	85.53%	9.232%
70	16.609	2466	2469	2475	rVB3	45365	67631	19.06%	2.058%
71	17.192	2562	2568	2572	rBV8	28694	61185	17.25%	1.862%

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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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72	19.462	2952	2954	2955	rBV2	6259	5518	1.56%	0.168%
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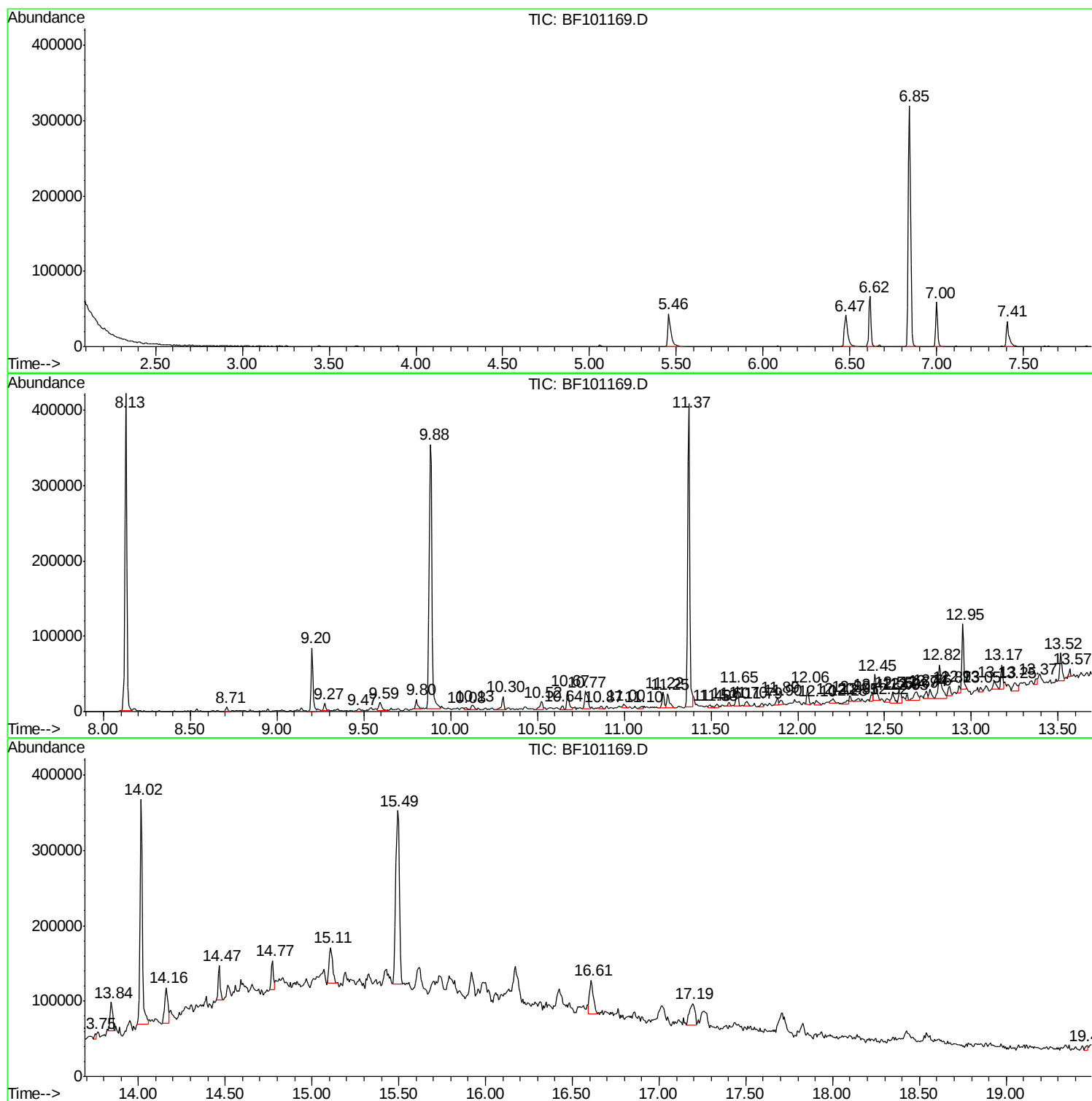
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Instrument :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF113017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Acq On : 6 Dec 2017 15:57
Operator : SJ/JU
Sample : I6303-01 10X
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Instrument :
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ClientSampleId :
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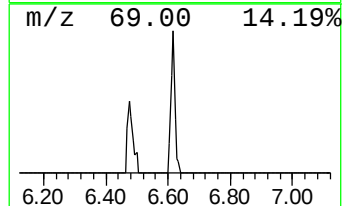
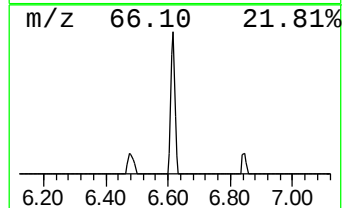
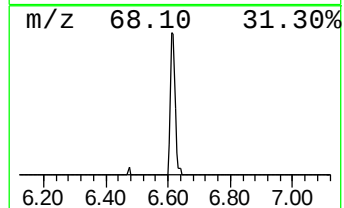
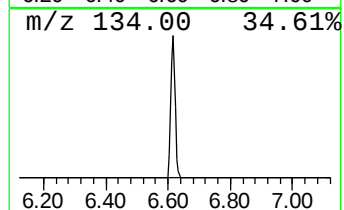
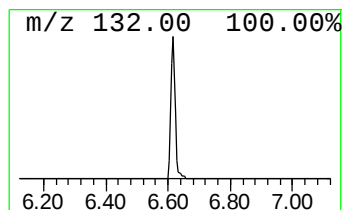
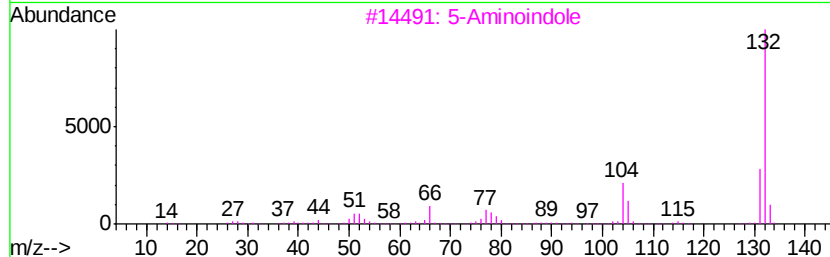
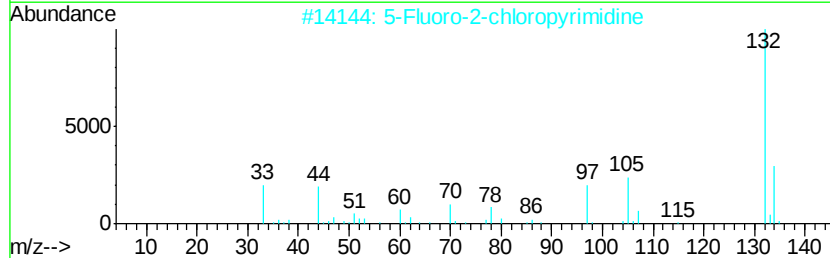
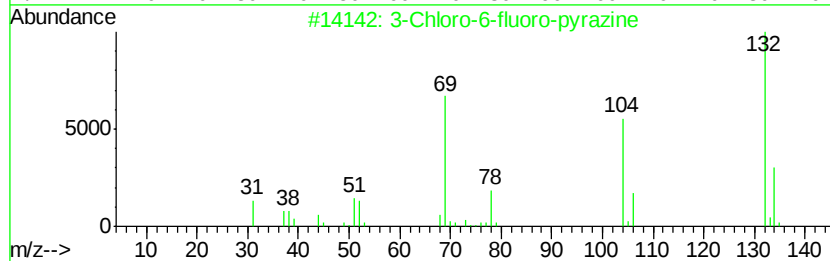
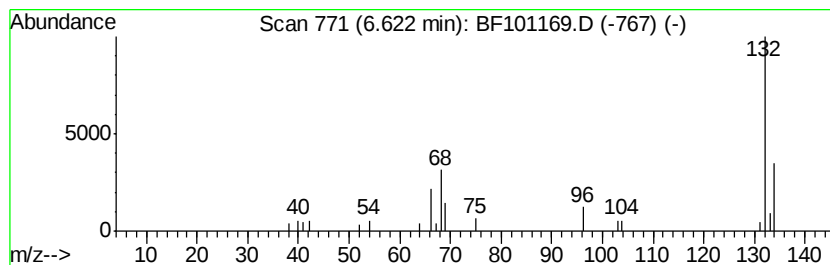
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF113017.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown6.62 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.62	4.57 ng	60804	1,4-Dichlorobenzene-d4	6.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
3		5-Aminoindole	132	C8H8N2	005192-03-0	9
4		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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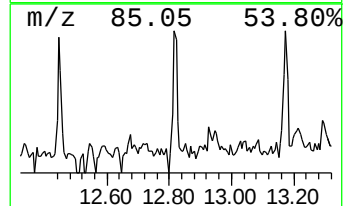
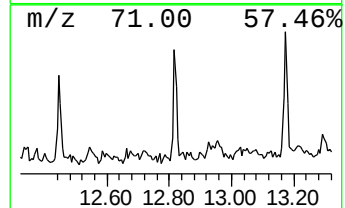
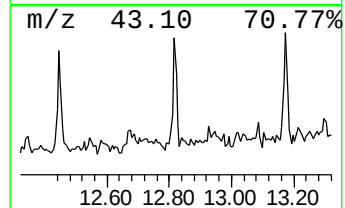
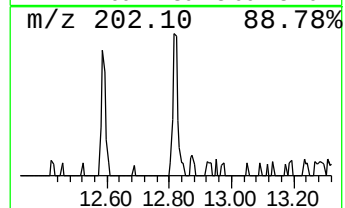
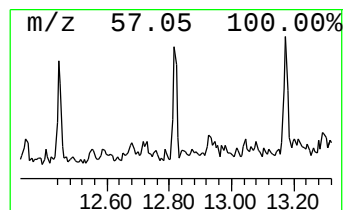
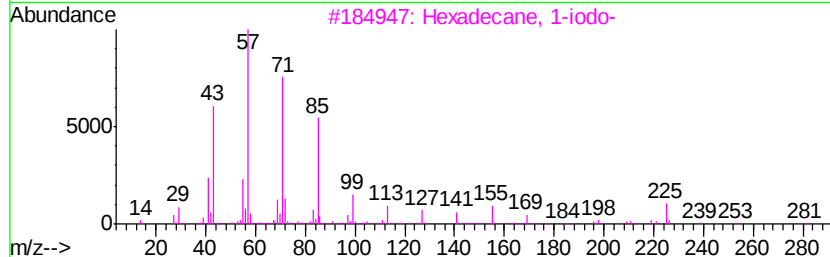
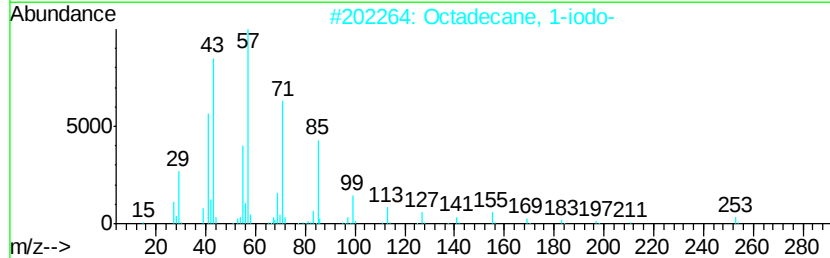
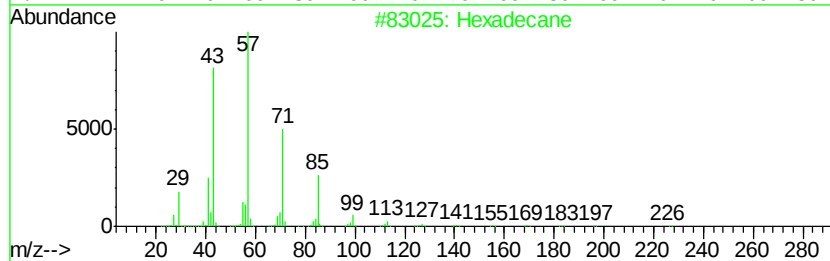
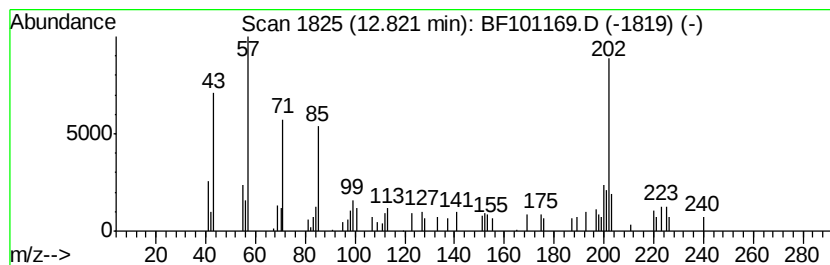
Instrument :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF113017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Hexadecane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.82	4.21 ng	62764	Chrysene-d12	14.02
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Hexadecane	226 C16H34	000544-76-3	90
2	Octadecane, 1-iodo-	380 C18H37I	000629-93-6	43
3	Hexadecane, 1-iodo-	352 C16H33I	000544-77-4	43
4	Eicosane, 10-methyl-	296 C21H44	054833-23-7	43
5	Silane, trichlorooctadecyl-	386 C18H37Cl3Si	000112-04-9	43



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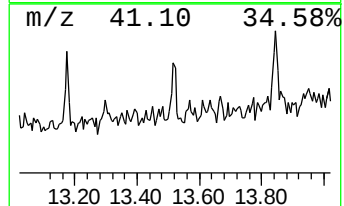
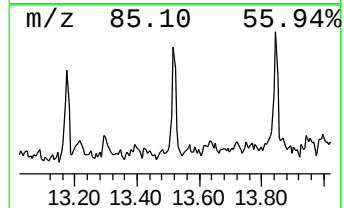
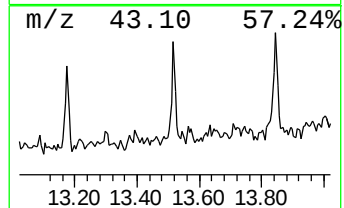
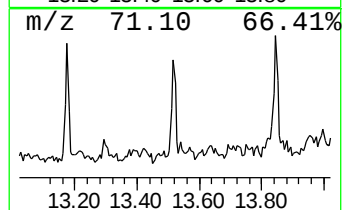
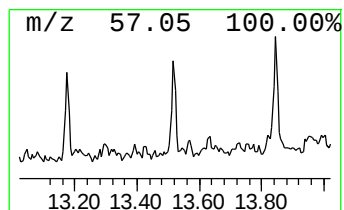
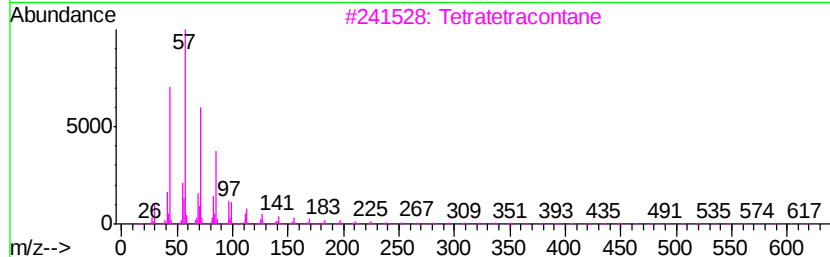
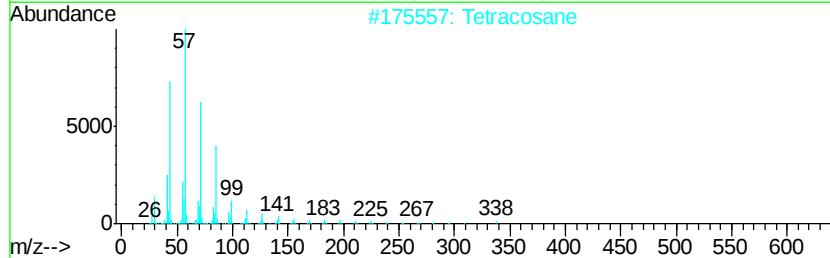
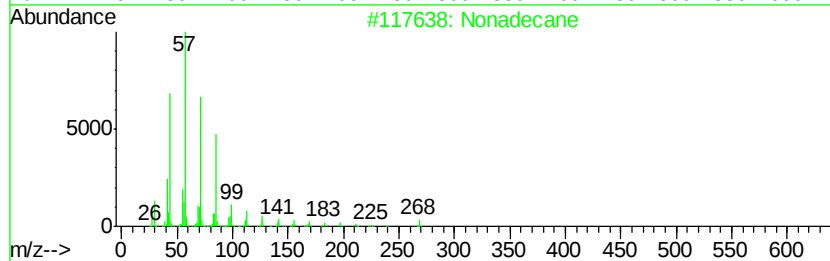
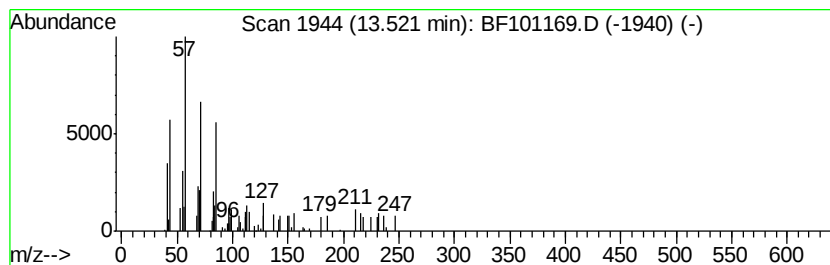
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF113017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Nonadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.52	2.30 ng	34390	Chrysene-d12	14.02

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane		268	C19H40	000629-92-5	96
2	Tetracosane		338	C24H50	000646-31-1	86
3	Tetratetracontane		619	C44H90	007098-22-8	86
4	Octadecane		254	C18H38	000593-45-3	83
5	Hexatriacontane		507	C36H74	000630-06-8	83



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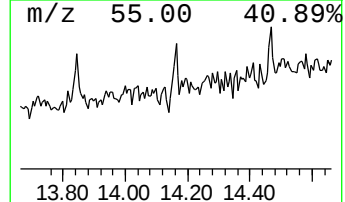
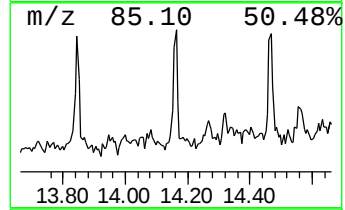
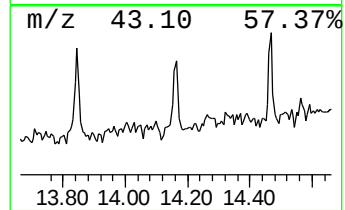
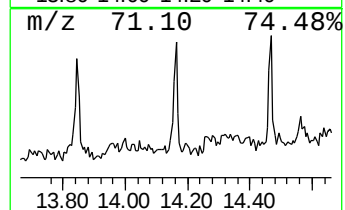
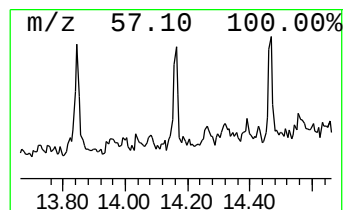
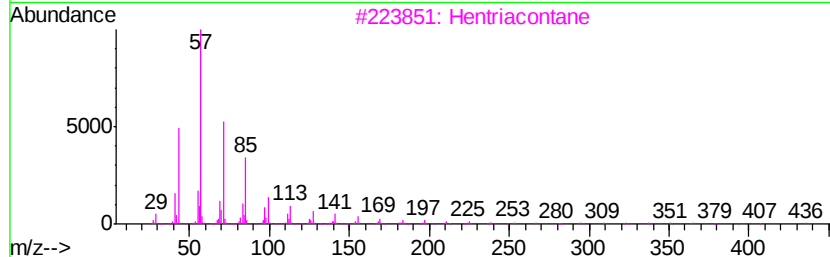
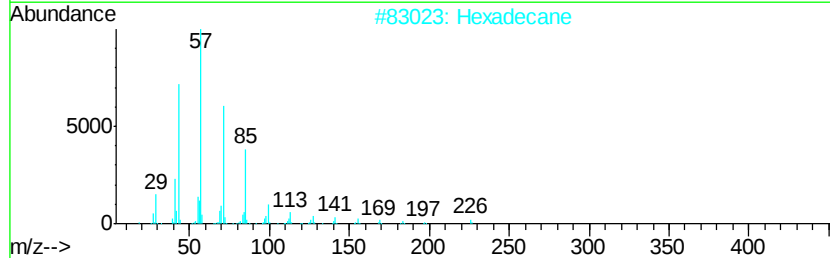
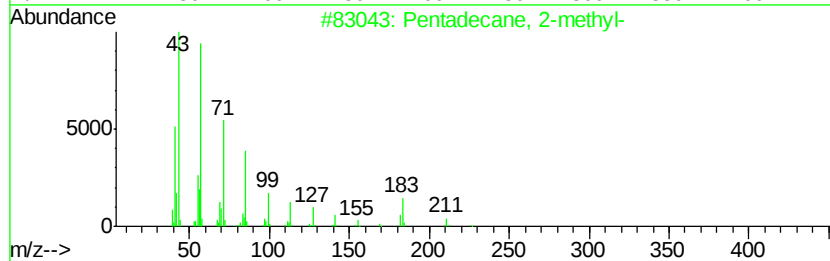
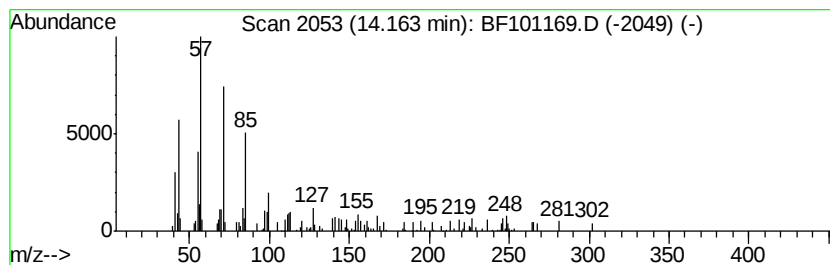
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF113017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Pentadecane, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.16	3.55 ng	52914	Chrysene-d12	14.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane, 2-methyl-	226	C16H34	001560-93-6	81
2		Hexadecane	226	C16H34	000544-76-3	76
3		Hentriacontane	437	C31H64	000630-04-6	74
4		Heneicosane	296	C21H44	000629-94-7	74
5		Pentacosane	352	C25H52	000629-99-2	74



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120617\
Data File : BF101169.D
Acq On : 6 Dec 2017 15:57
Operator : SJ/JU
Sample : I6303-01 10X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
HD-02-120417

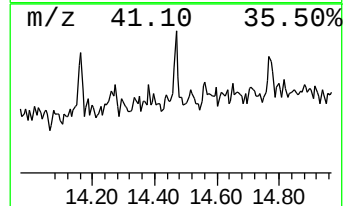
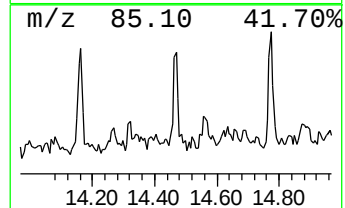
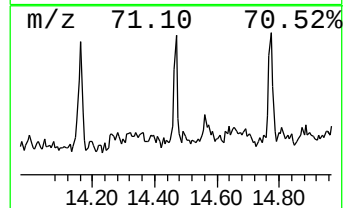
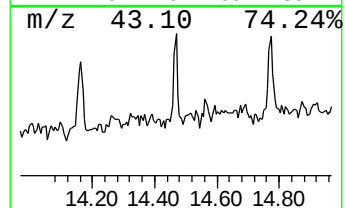
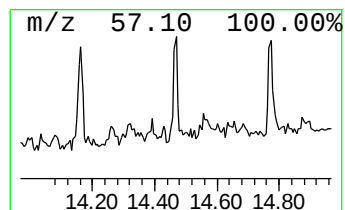
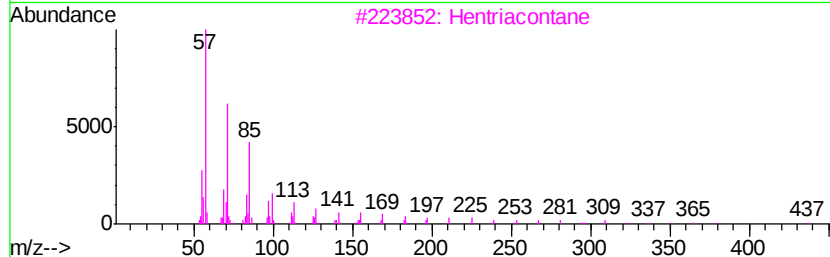
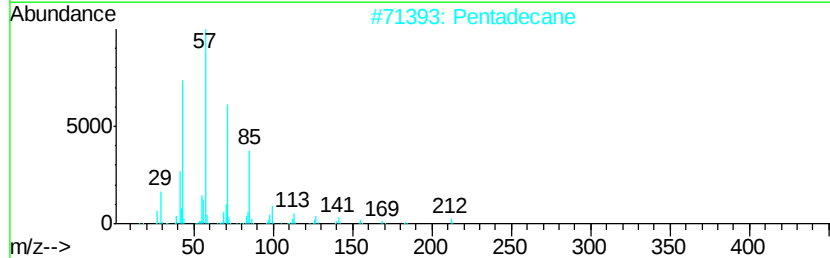
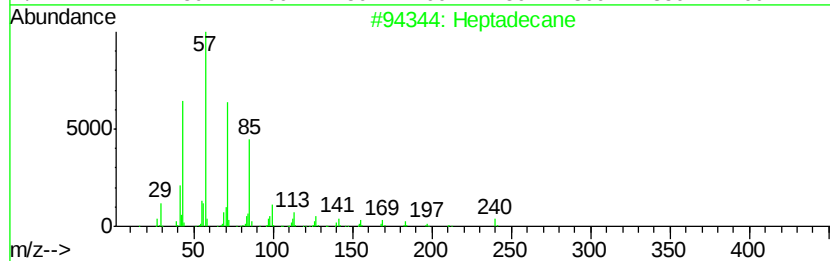
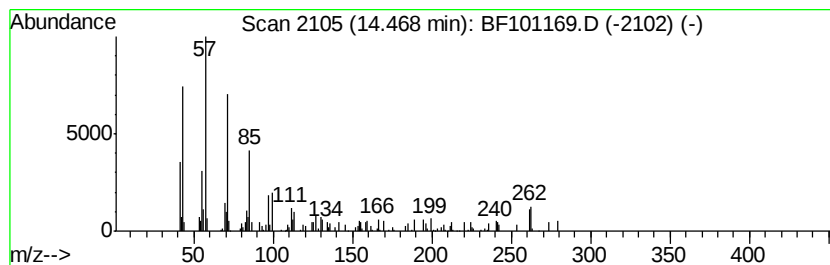
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Heptadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.47	2.40 ng	35758	Chrysene-d12	14.02

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane		240	C17H36	000629-78-7	91
2	Pentadecane		212	C15H32	000629-62-9	70
3	Hentriacontane		437	C31H64	000630-04-6	60
4	Octadecane, 2-methyl-		268	C19H40	001560-88-9	58
5	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	53



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120617\
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ClientSampleId :
HD-02-120417

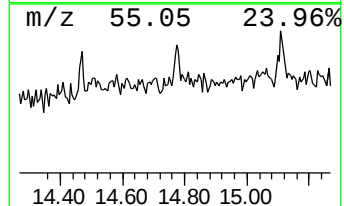
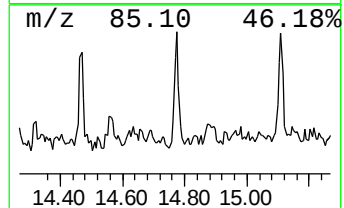
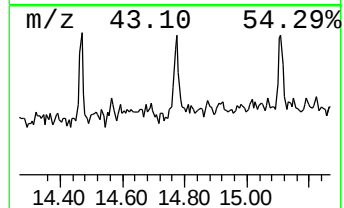
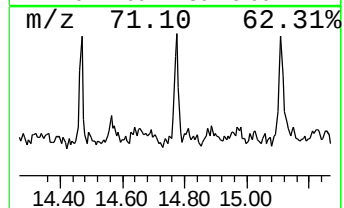
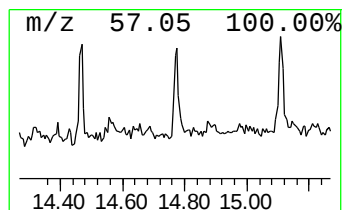
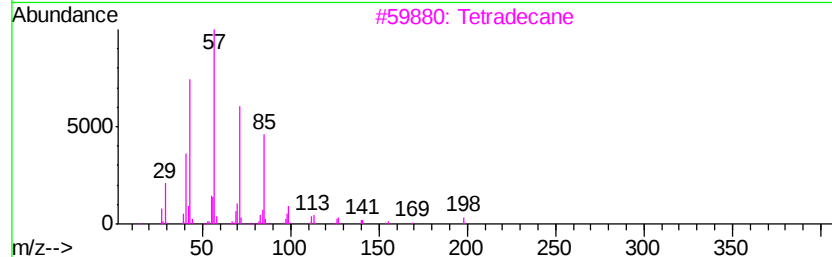
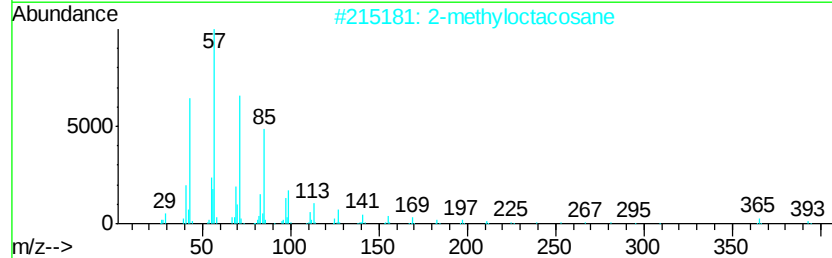
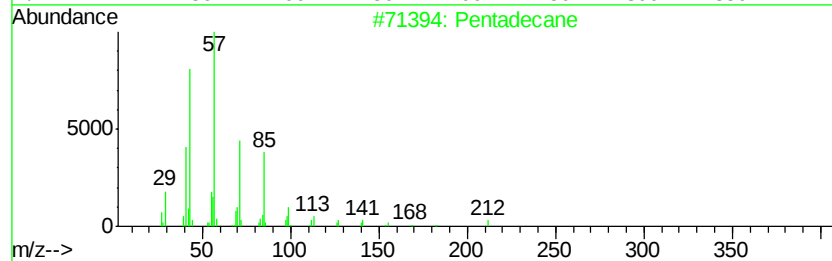
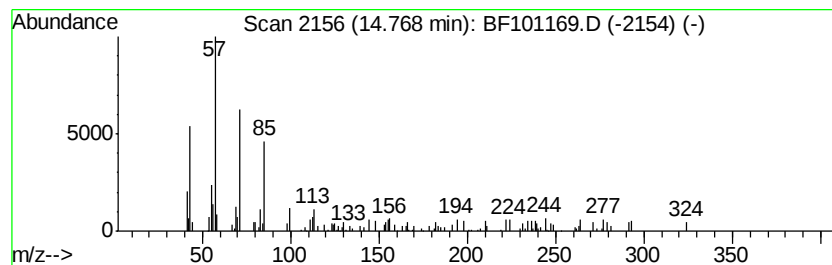
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Pentadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.77	2.45 ng	37235	Perylene-d12	15.50

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane		212	C15H32	000629-62-9	68
2	2-methyloctacosane		408	C29H60	1000376-72-8	58
3	Tetradecane		198	C14H30	000629-59-4	58
4	Heptadecane, 2,6,10,15-tetramethyl-		296	C21H44	054833-48-6	52
5	Tetracosane		338	C24H50	000646-31-1	52



Data Path : Z:\HPCHEM1\BNA F\DATA\BF120617\
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Sample : I6303-01 10X
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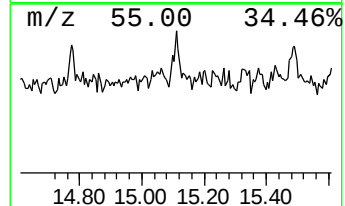
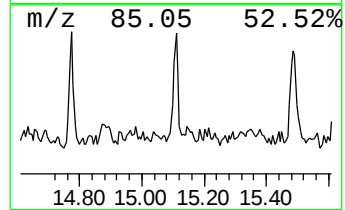
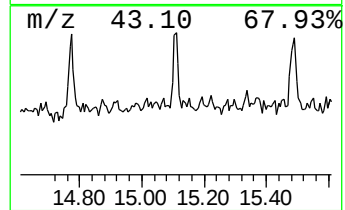
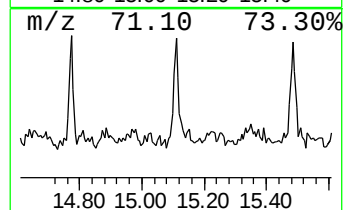
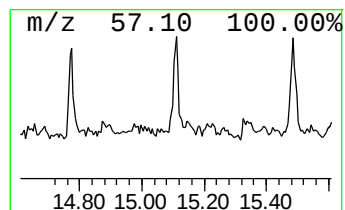
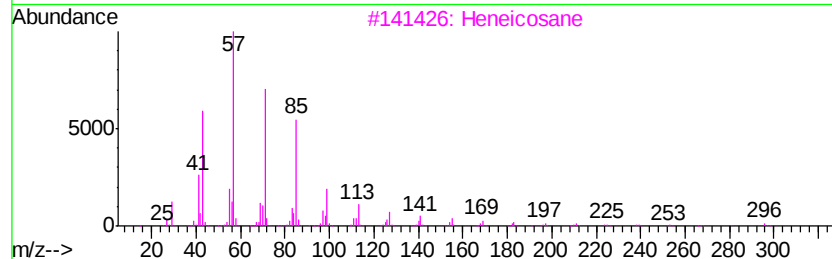
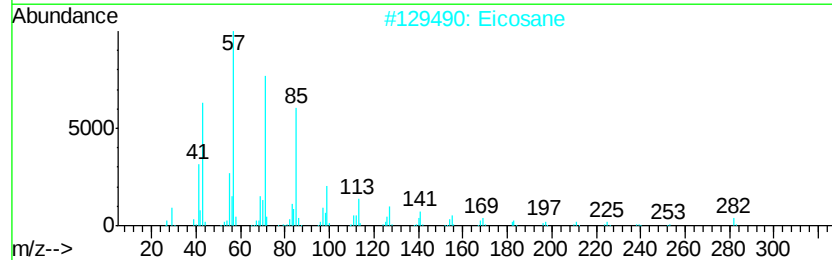
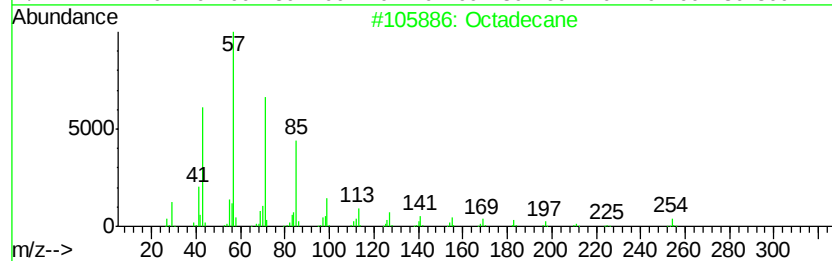
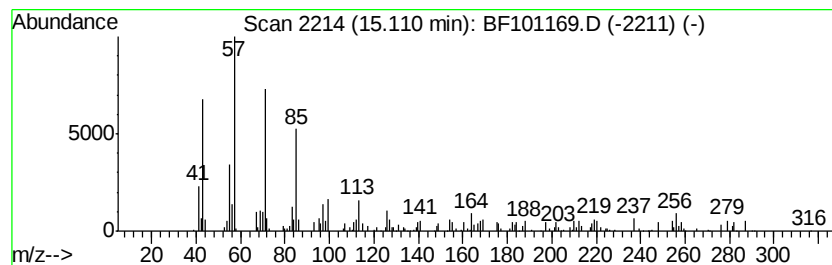
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF113017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Octadecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.
15.11	3.83 ng	58132	Perylene-d12		15.50
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane	254	C18H38	000593-45-3	98
2	Eicosane	282	C20H42	000112-95-8	93
3	Heneicosane	296	C21H44	000629-94-7	70
4	Hexacosane	366	C26H54	000630-01-3	70
5	5-Ethyl-5-methylnonadecane	310	C22H46	1000360-42-7	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120617\
Data File : BF101169.D
Acq On : 6 Dec 2017 15:57
Operator : SJ/JU
Sample : I6303-01 10X
Misc :
ALS Vial : 11 Sample Multiplier: 1

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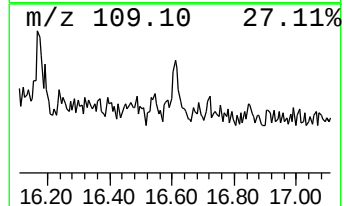
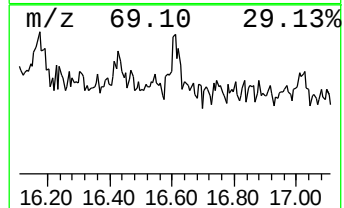
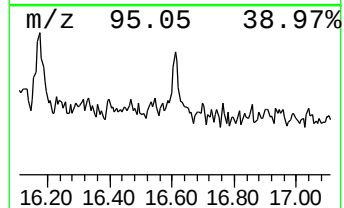
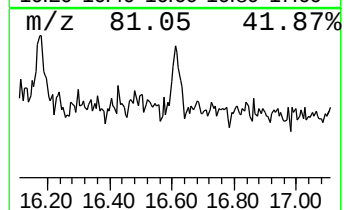
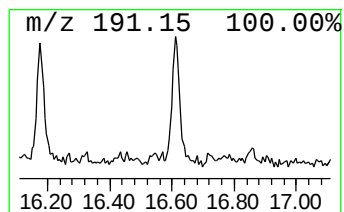
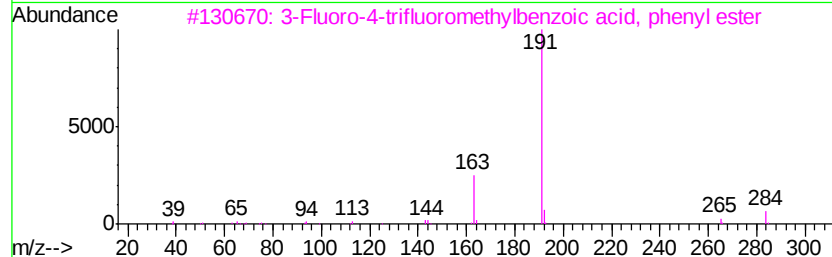
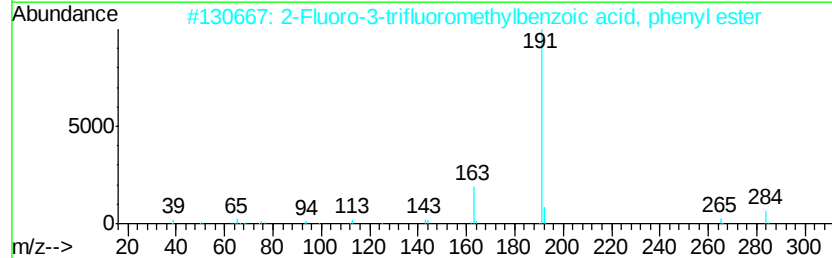
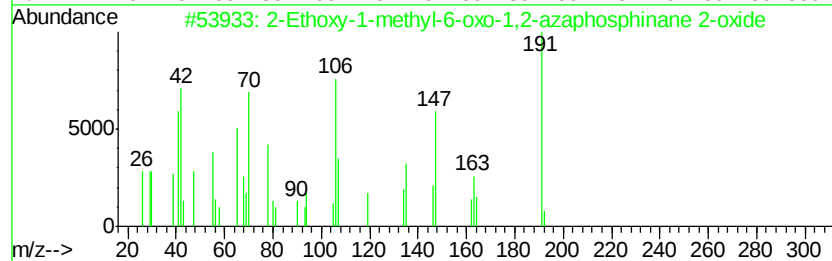
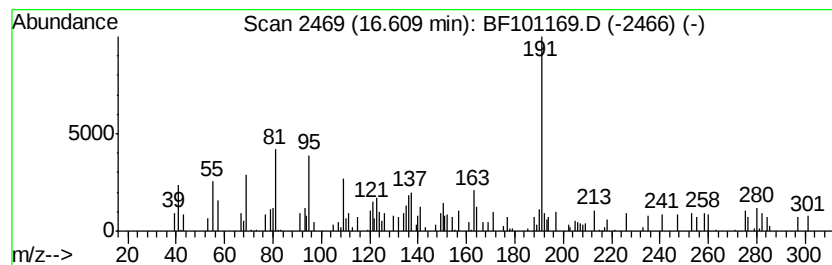
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown16.61 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.61	4.46 ng	67631	Perylene-d12	15.50

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Ethoxy-1-methyl-6-oxo-1,2-azap...	191	C7H14N03P	093989-00-5	38		
2	2-Fluoro-3-trifluoromethylbenzoi...	284	C14H8F4O2	1000357-64-3	38		
3	3-Fluoro-4-trifluoromethylbenzoi...	284	C14H8F4O2	1000357-93-7	38		
4	5-Fluoro-2-trifluoromethylbenzoi...	284	C14H8F4O2	1000357-62-7	38		
5	2-Fluoro-6-trifluoromethylbenzoi...	284	C14H8F4O2	1000357-69-4	38		



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120617\
Data File : BF101169.D
Acq On : 6 Dec 2017 15:57
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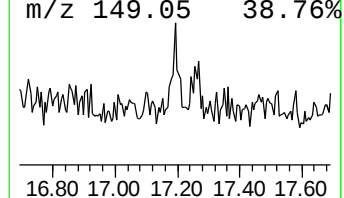
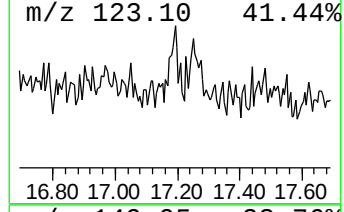
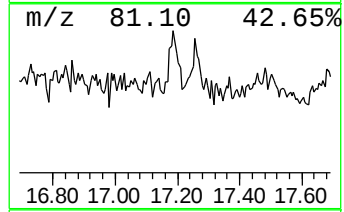
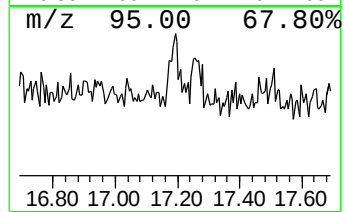
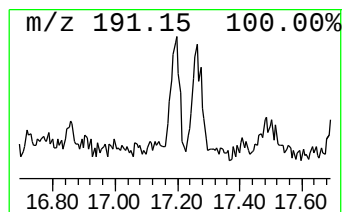
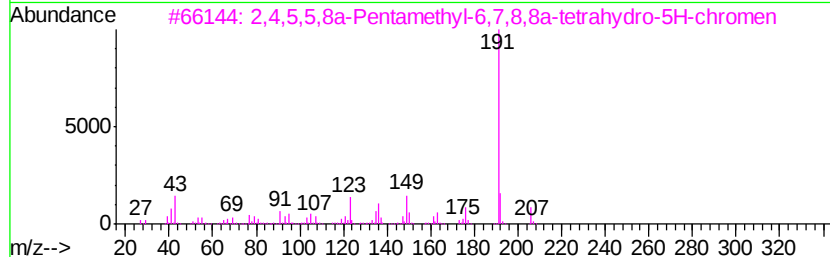
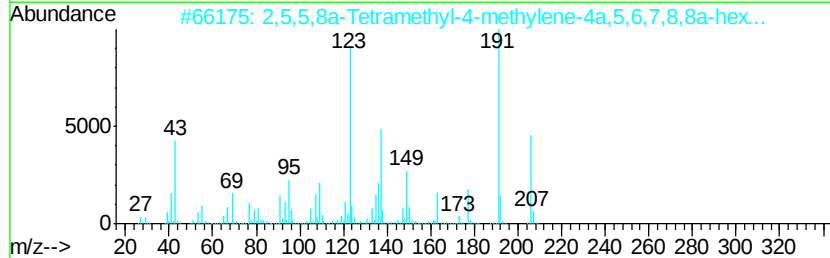
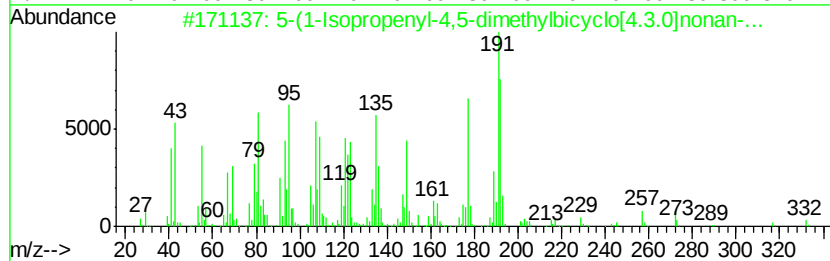
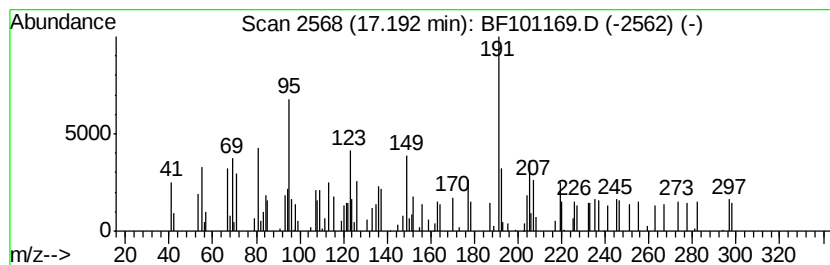
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown17.19 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.19	4.03 ng	61185	Perylene-d12	15.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-(1-Isopropenyl-4,5-dimethylbic...	332	C22H36O2	1000195-69-8	45
2		2,5,5,8a-Tetramethyl-4-methylene...	206	C14H22O	093175-76-9	43
3		2,4,5,5,8a-Pentamethyl-6,7,8,8a-...	206	C14H22O	1000195-40-9	38
4		Isophthalic acid, 3-methylphenyl...	298	C18H18O4	1000344-51-5	38
5		2-Phenylamino-5,6(4H)dihydro-1,3...	192	C10H12N2S	1000147-35-4	38



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

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					#	RT	Resp	Conc
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Nonadecane	13.52	2.3	ng	34390	5	14.02	298419	20.0
Pentadecane, 2-me...	14.16	3.5	ng	52914	5	14.02	298419	20.0
Heptadecane	14.47	2.4	ng	35758	5	14.02	298419	20.0
Pentadecane	14.77	2.5	ng	37235	6	15.50	303440	20.0
Octadecane	15.11	3.8	ng	58132	6	15.50	303440	20.0
unknown16.61	16.61	4.5	ng	67631	6	15.50	303440	20.0
unknown17.19	17.19	4.0	ng	61185	6	15.50	303440	20.0