

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092717\
 Data File : BF098877.D
 Acq On : 27 Sep 2017 22:00
 Operator : SJ/JU
 Sample : I5423-04
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SA-06-092517-B

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-BF092317.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	2.199	15	19	34	rVB	86047	156728	5.29%	0.410%
2	5.140	515	519	530	rVB	470858	545185	18.39%	1.426%
3	5.534	581	586	589	rBV	2858022	2640531	89.06%	6.908%
4	6.539	751	757	766	rVB	2697288	2733109	92.19%	7.150%
5	6.681	777	781	785	rBV	2780922	2737263	92.33%	7.161%
6	6.916	817	821	824	rVB	1133054	929849	31.36%	2.433%
7	7.069	843	847	851	rBV	2271025	2027683	68.39%	5.305%
8	7.475	911	916	919	rBV	1886003	1681575	56.72%	4.399%
9	8.198	1035	1039	1042	rBV	1447211	1266314	42.71%	3.313%
10	9.269	1216	1221	1224	rBV	3399544	2964797	100.00%	7.756%
11	9.345	1230	1234	1236	rBV2	35768	34527	1.16%	0.090%
12	9.604	1276	1278	1281	rBV	33320	35572	1.20%	0.093%
13	9.657	1284	1287	1295	rVB	449699	414682	13.99%	1.085%
14	9.810	1310	1313	1317	rBV	175034	157820	5.32%	0.413%
15	9.875	1319	1324	1326	rVB2	59813	69142	2.33%	0.181%
16	9.951	1331	1337	1341	rVV2	1508570	1343094	45.30%	3.514%
17	9.986	1341	1343	1346	rVB	62003	49640	1.67%	0.130%
18	10.151	1366	1371	1375	rVB3	58204	74385	2.51%	0.195%
19	10.192	1375	1378	1382	rBV2	39768	34933	1.18%	0.091%
20	10.263	1387	1390	1393	rBV3	33713	41647	1.40%	0.109%
21	10.375	1402	1409	1411	rVB	113510	127216	4.29%	0.333%
22	10.498	1427	1430	1435	rVB	98978	93304	3.15%	0.244%
23	10.586	1442	1445	1452	rVB2	43977	71967	2.43%	0.188%
24	10.739	1467	1471	1474	rBV	1708465	1487634	50.18%	3.892%
25	10.845	1486	1489	1490	rBV	137872	108528	3.66%	0.284%
26	11.045	1518	1523	1525	rBV3	38190	49976	1.69%	0.131%
27	11.292	1562	1565	1567	rBV	91730	74153	2.50%	0.194%
28	11.327	1567	1571	1576	rVB3	55438	93229	3.14%	0.244%
29	11.439	1586	1590	1592	rBV	1423426	1202640	40.56%	3.146%
30	11.463	1592	1594	1597	rVV	605488	450137	15.18%	1.178%
31	11.516	1600	1603	1605	rVB	179113	161154	5.44%	0.422%
32	11.545	1605	1608	1611	rBV2	33240	33218	1.12%	0.087%
33	11.663	1625	1628	1632	rBV	56944	53606	1.81%	0.140%
34	11.721	1635	1638	1642	rVB2	45945	53031	1.79%	0.139%

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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-BF092317.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	11.821	1652	1655	1658	rBV	56716	47387	1.60%	0.124%
36	11.939	1672	1675	1678	rBV2	105759	151677	5.12%	0.397%
37	11.969	1678	1680	1684	rVB	131842	109093	3.68%	0.285%
38	12.016	1685	1688	1691	rBV	48862	44461	1.50%	0.116%
39	12.051	1691	1694	1697	rVV2	209704	235709	7.95%	0.617%
40	12.074	1697	1698	1702	rVB	84339	43801	1.48%	0.115%
41	12.127	1704	1707	1712	rBV2	36572	36476	1.23%	0.095%
42	12.233	1722	1725	1727	rBV	67281	52344	1.77%	0.137%
43	12.257	1727	1729	1732	rVB2	46743	40022	1.35%	0.105%
44	12.486	1765	1768	1770	rBV	105871	108531	3.66%	0.284%
45	12.527	1773	1775	1780	rVB2	143717	153712	5.18%	0.402%
46	12.574	1780	1783	1787	rBV	96363	109323	3.69%	0.286%
47	12.651	1792	1796	1800	rBV	996940	859120	28.98%	2.248%
48	12.745	1809	1812	1814	rBV	63527	53642	1.81%	0.140%
49	12.880	1829	1835	1841	rBV	1080610	1173125	39.57%	3.069%
50	12.933	1841	1844	1847	rVB2	51658	65491	2.21%	0.171%
51	13.021	1855	1859	1862	rBV	2481506	2191785	73.93%	5.734%
52	13.098	1868	1872	1875	rBV3	94686	136536	4.61%	0.357%
53	13.210	1884	1891	1894	rVB	160861	261735	8.83%	0.685%
54	13.274	1899	1902	1905	rVV	116584	92051	3.10%	0.241%
55	13.310	1905	1908	1911	rVV2	141635	143437	4.84%	0.375%
56	13.404	1921	1924	1927	rBV	112681	92953	3.14%	0.243%
57	13.433	1927	1929	1932	rVB	104768	90557	3.05%	0.237%
58	13.710	1974	1976	1981	rVB	78135	90411	3.05%	0.237%
59	13.827	1994	1996	1999	rVB	79032	65148	2.20%	0.170%
60	13.857	1999	2001	2002	rBV	73924	48541	1.64%	0.127%
61	13.880	2002	2005	2007	rVV	117181	112234	3.79%	0.294%
62	13.904	2007	2009	2017	rVB4	248005	296619	10.00%	0.776%
63	14.045	2031	2033	2035	rBV	156605	168570	5.69%	0.441%
64	14.080	2035	2039	2041	rVV2	1199617	1568068	52.89%	4.102%
65	14.104	2041	2043	2048	rVB	575124	518921	17.50%	1.358%
66	14.186	2054	2057	2060	rVB	128158	95790	3.23%	0.251%
67	14.292	2073	2075	2079	rBV2	56455	76175	2.57%	0.199%
68	14.462	2102	2104	2107	rVB	149901	128871	4.35%	0.337%
69	14.498	2108	2110	2113	rVB2	81112	62942	2.12%	0.165%
70	14.586	2123	2125	2127	rBV	95651	78594	2.65%	0.206%
71	15.133	2214	2218	2221	rBV	696547	1032194	34.81%	2.700%

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72	15.239	2234	2236	2239	rVB	141319	145510	4.91%	0.381%
73	15.445	2267	2271	2277	rVB	457990	580722	19.59%	1.519%
74	15.509	2278	2282	2288	rVV	622224	826011	27.86%	2.161%
75	15.574	2289	2293	2296	rVV	621736	833761	28.12%	2.181%
76	15.604	2296	2298	2305	rVB	234062	299289	10.09%	0.783%
77	17.080	2544	2549	2556	rVB2	255706	514705	17.36%	1.347%
78	17.539	2622	2627	2635	rVB	196285	390179	13.16%	1.021%

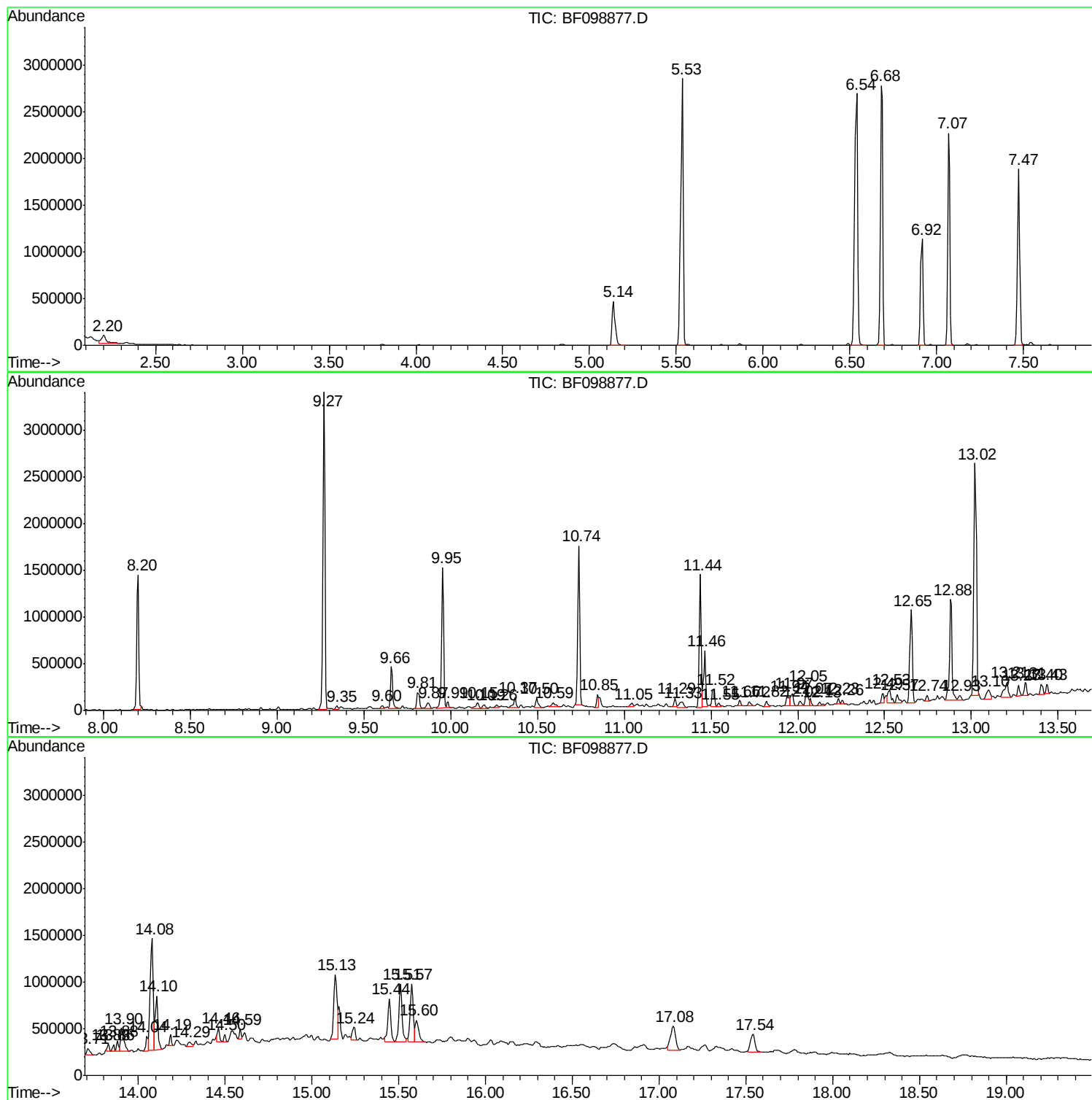
Sum of corrected areas: 38224492

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Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.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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Instrument :
BNA_F
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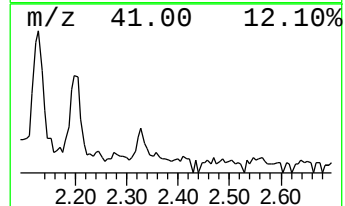
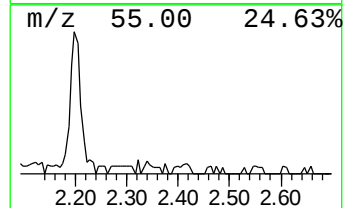
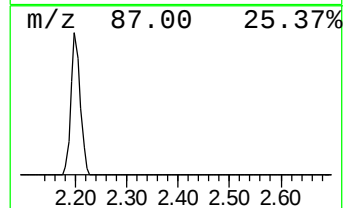
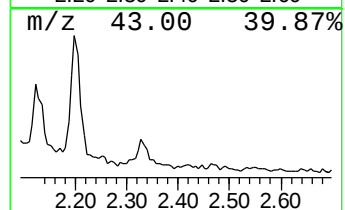
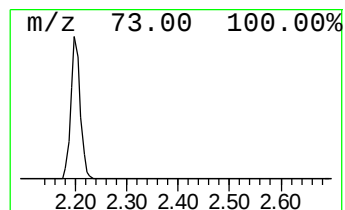
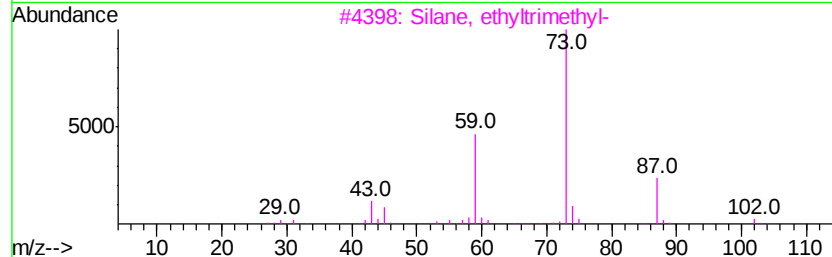
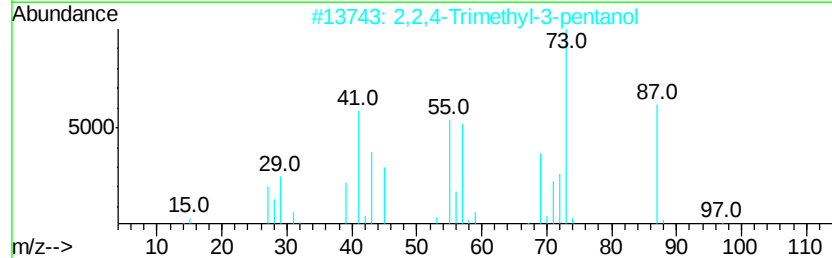
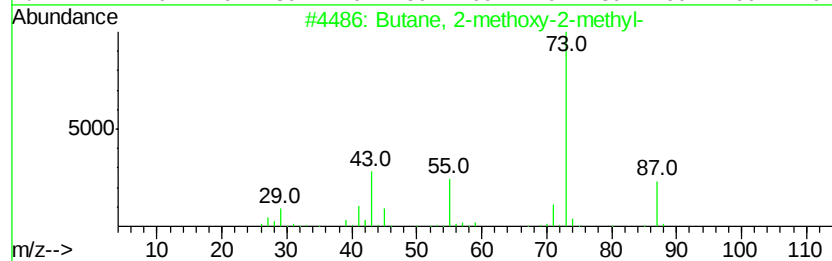
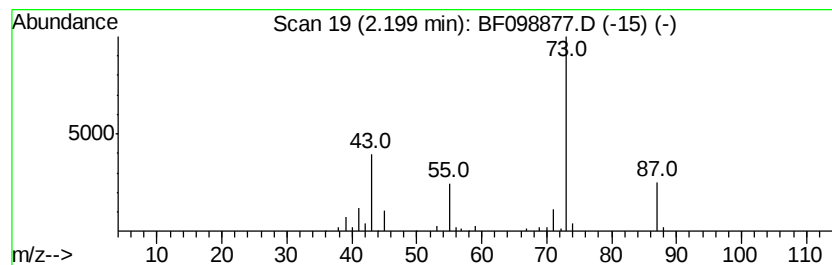
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.20	3.37 ng	156728	1,4-Dichlorobenzene-d4	6.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3		Silane, ethyltrimethyl-	102	C5H14Si	003439-38-1	33
4		3-Hexanol, 2-methyl-	116	C7H16O	000617-29-8	25
5		Boronic acid, ethyl-, dimethyl e...	102	C4H11BO2	007318-82-3	17



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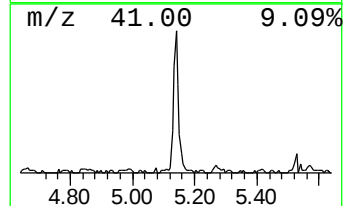
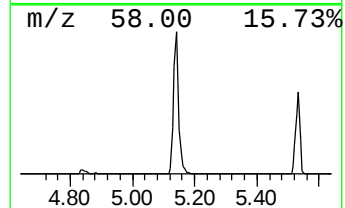
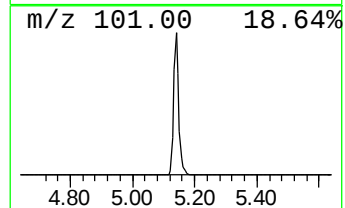
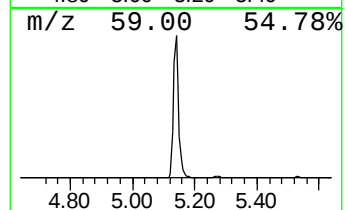
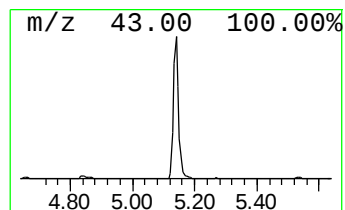
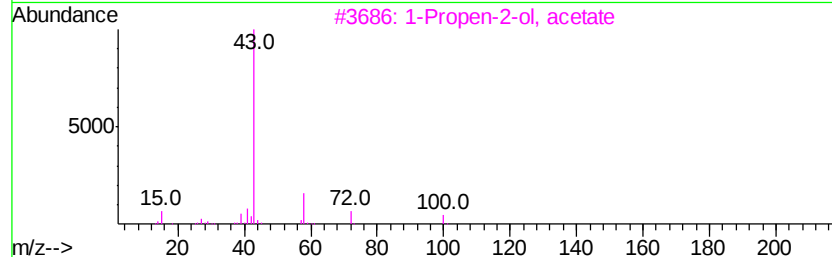
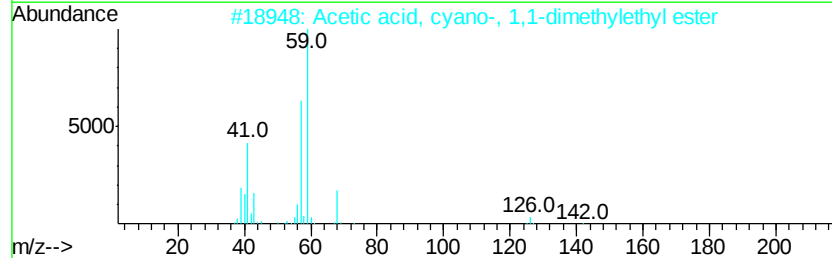
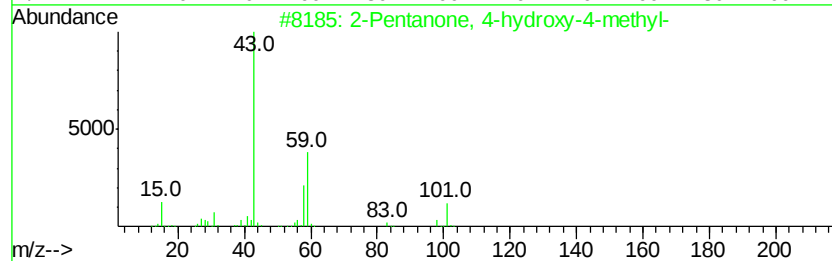
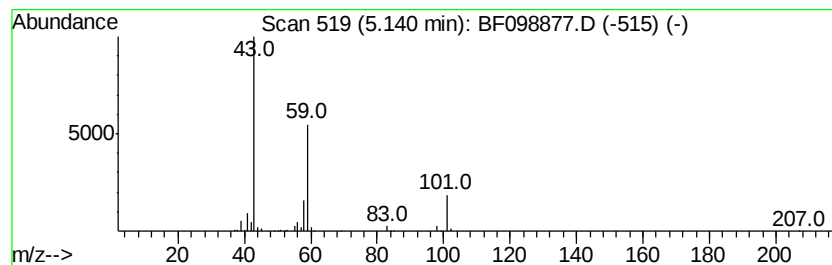
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.14	11.73 ng	545185	1,4-Dichlorobenzene-d4	6.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
3		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	9



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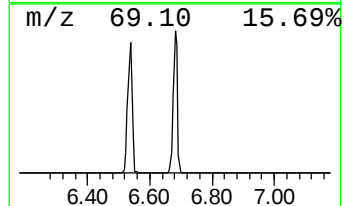
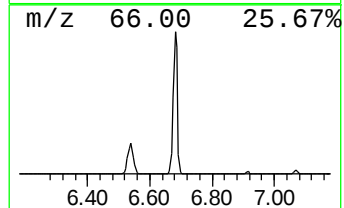
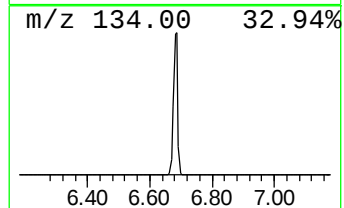
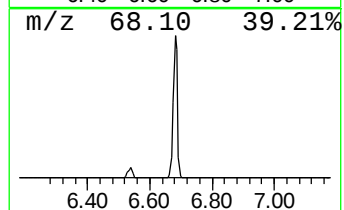
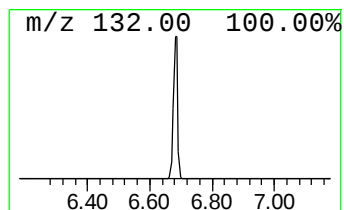
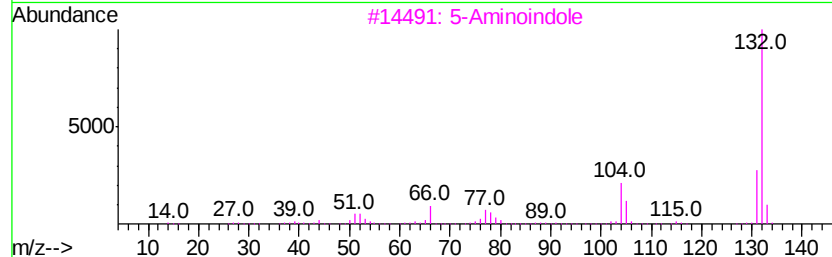
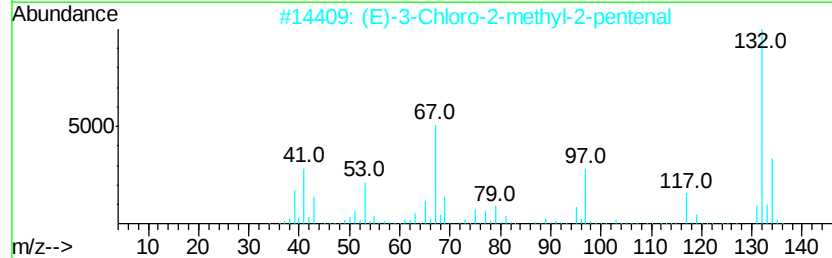
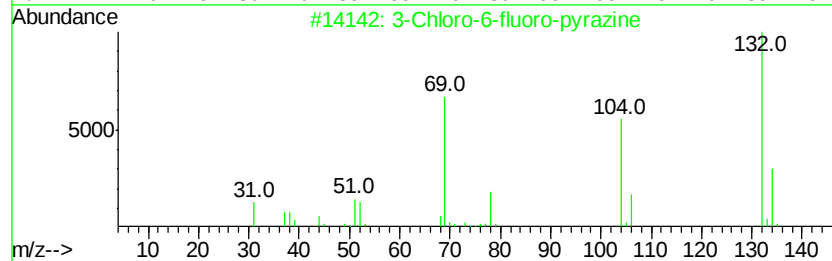
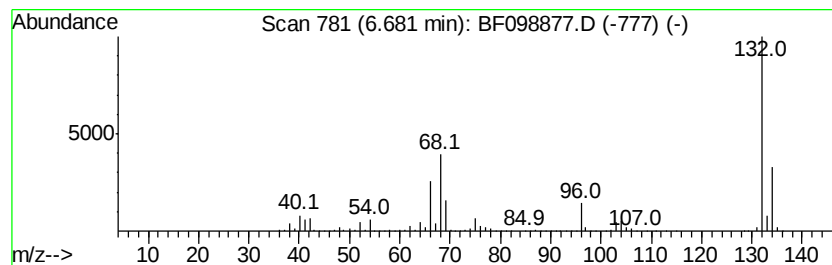
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 unknown6.68 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.68	58.88 ng	2737260	1,4-Dichlorobenzene-d4	6.92

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		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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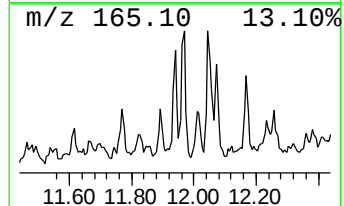
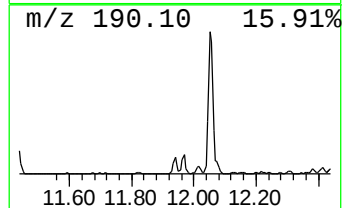
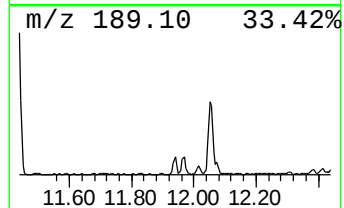
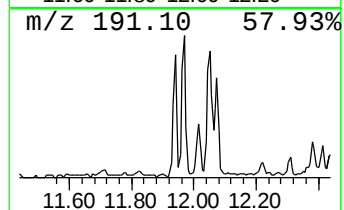
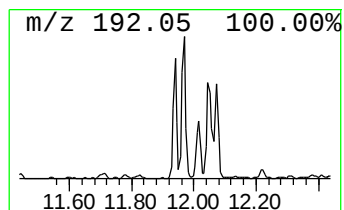
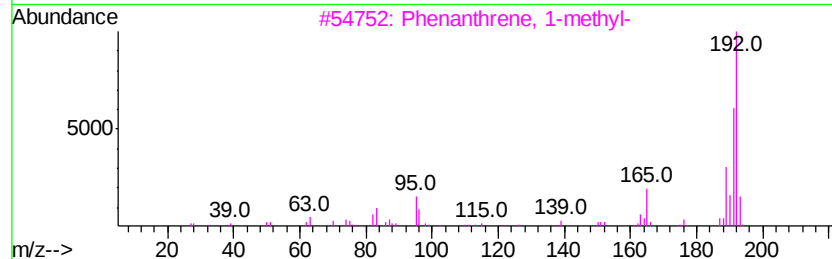
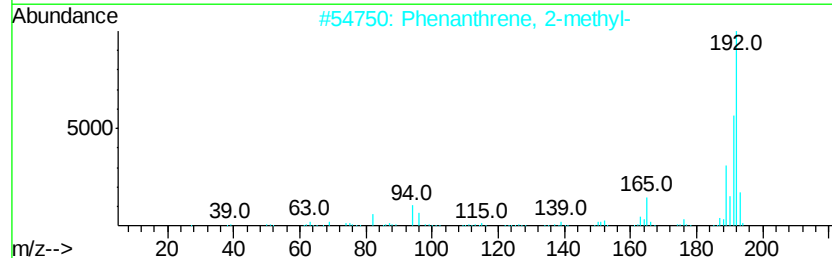
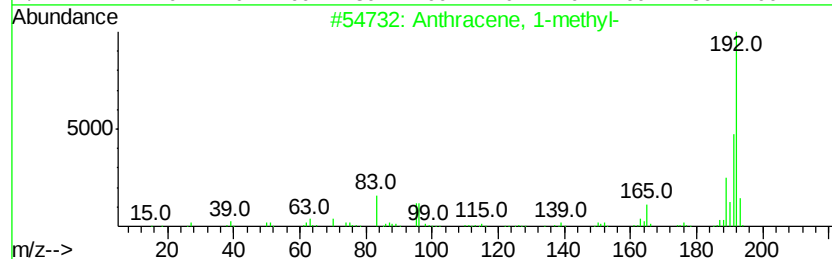
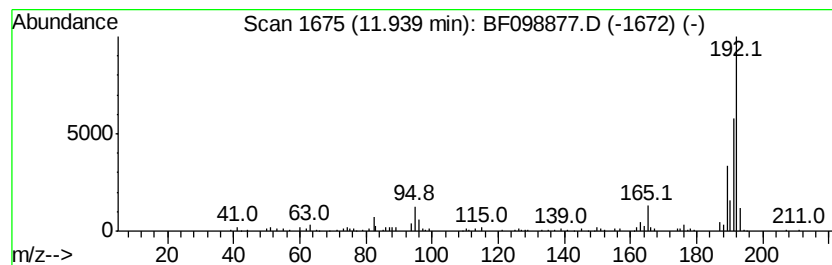
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Anthracene, 1-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.94	2.52 ng	151677	Phenanthrene-d10	11.44

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Anthracene, 1-methyl-	192	C15H12	000610-48-0	93
2		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	91
3		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	91
4		Anthracene, 9-methyl-	192	C15H12	000779-02-2	91
5		1H-Cyclopropa[1]phenanthrene, 1a, ...	192	C15H12	000949-41-7	89



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092717\
Data File : BF098877.D
Acq On : 27 Sep 2017 22:00
Operator : SJ/JU
Sample : I5423-04
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SA-06-092517-B

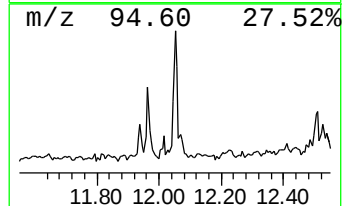
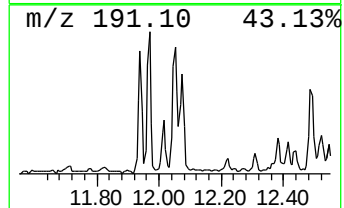
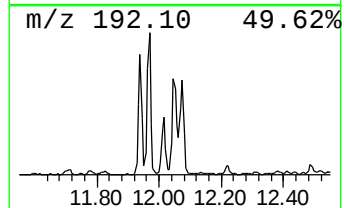
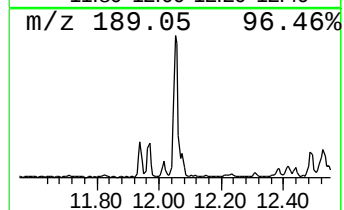
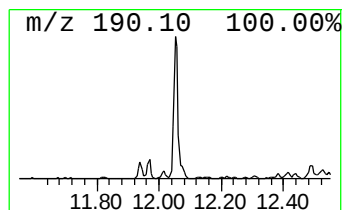
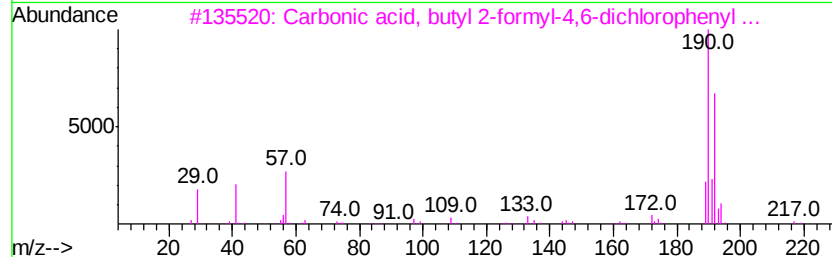
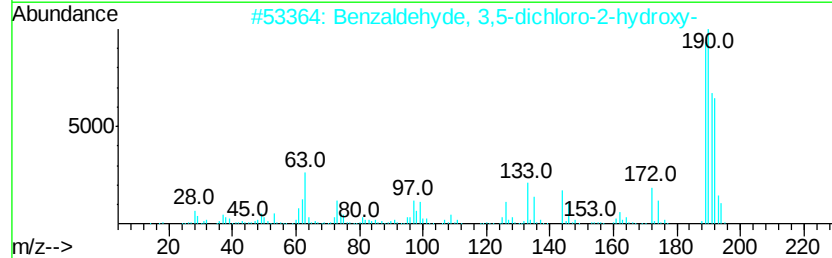
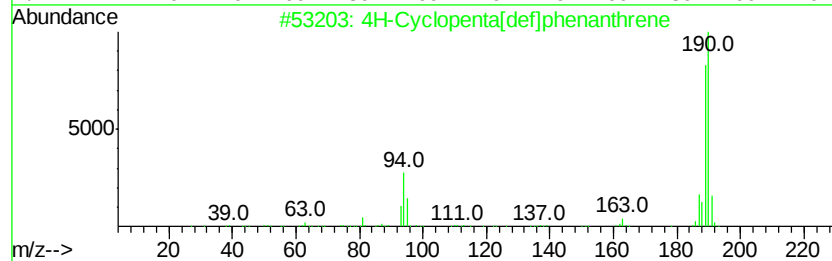
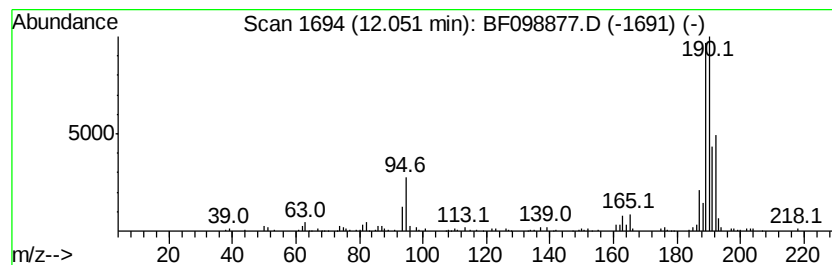
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 4H-Cyclopenta[def]phenanthrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.05	3.92 ng	235709	Phenanthrene-d10	11.44

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	76
2		Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	28
3		Carbonic acid, butyl 2-formyl-4,...	290	C12H12Cl2O4	1000331-42-2	25
4		1,1'-Biphenyl, 4,4'-difluoro-	190	C12H8F2	000398-23-2	23
5		2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	23



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092717\
Data File : BF098877.D
Acq On : 27 Sep 2017 22:00
Operator : SJ/JU
Sample : I5423-04
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ALS Vial : 13 Sample Multiplier: 1

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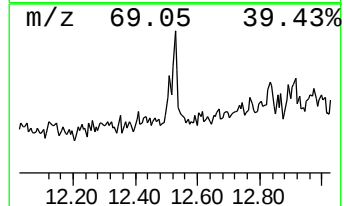
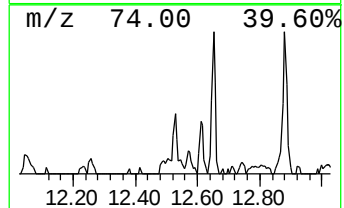
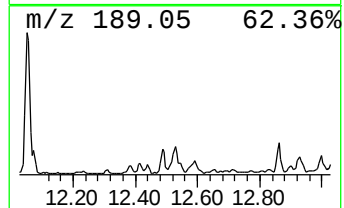
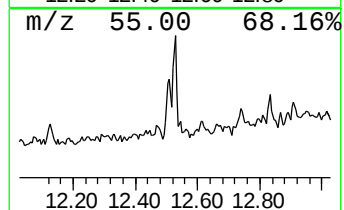
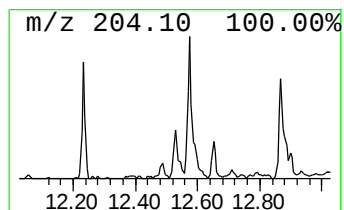
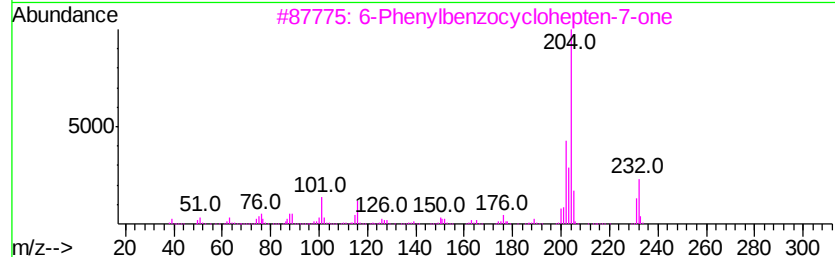
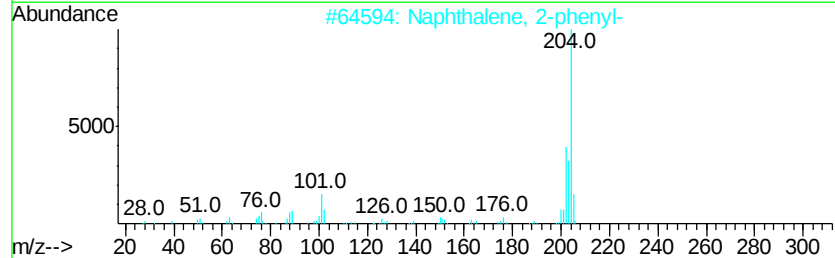
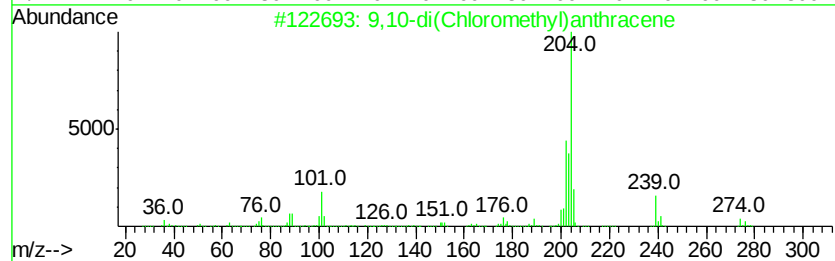
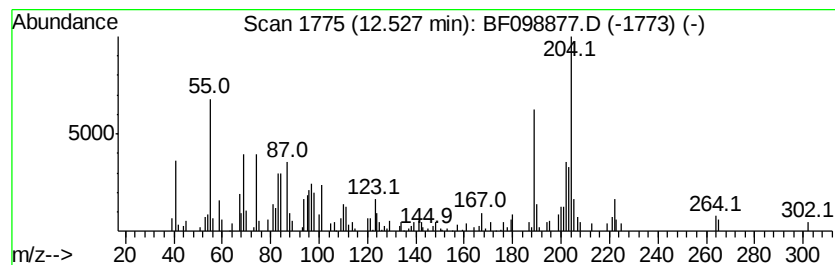
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown12.53 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.53	2.56 ng	153712	Phenanthrene-d10	11.44

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-di(Chloromethyl)anthracene	274	C16H12Cl2	010387-13-0	43		
2	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	38		
3	6-Phenylbenzocyclohepten-7-one	232	C17H12O	093327-56-1	38		
4	Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	38		
5	Cyclododecane, 1,2-dithia-	204	C10H20S2	000294-67-7	38		



Data Path : Z:\HPCHEM1\BNA F\DATA\BF092717\
Data File : BF098877.D
Acq On : 27 Sep 2017 22:00
Operator : SJ/JU
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Misc :
ALS Vial : 13 Sample Multiplier: 1

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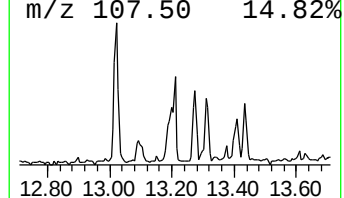
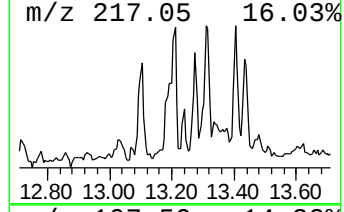
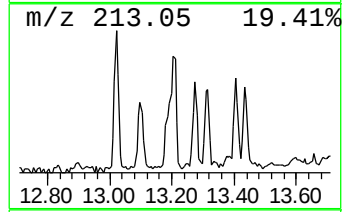
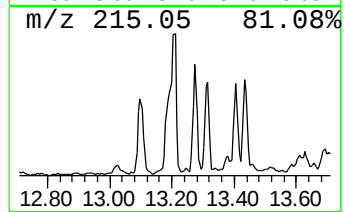
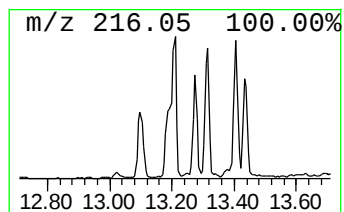
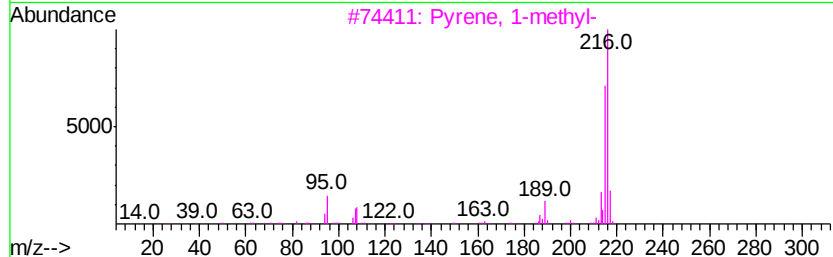
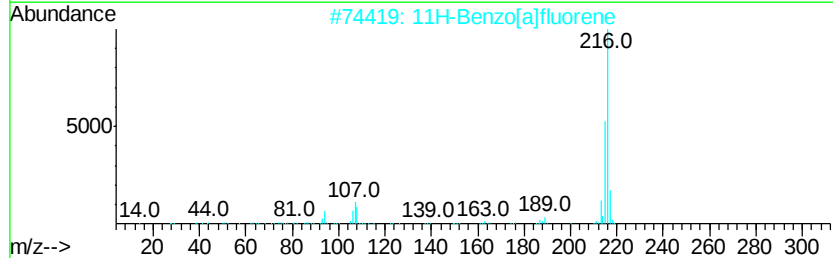
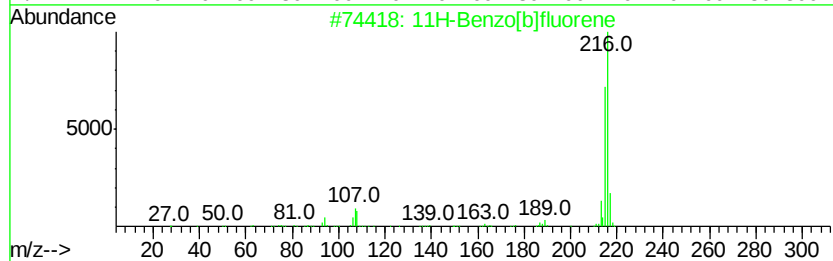
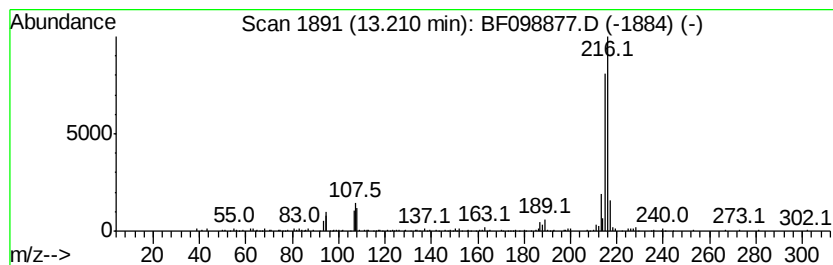
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 11H-Benzo[b]fluorene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.21	3.34 ng	261735	Chrysene-d12	14.08

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	93
2		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	93
3		Pvrene, 1-methyl-	216	C17H12	002381-21-7	91
4		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	90
5		7H-Benzo[c]fluorene	216	C17H12	000205-12-9	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092717\
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Sample : I5423-04
Misc :
ALS Vial : 13 Sample Multiplier: 1

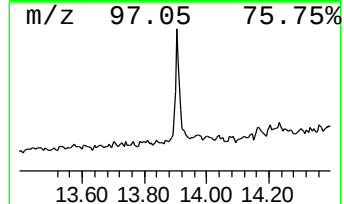
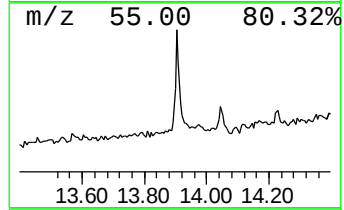
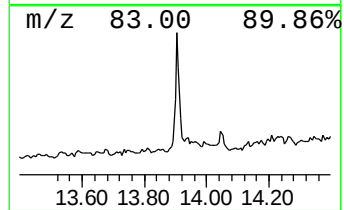
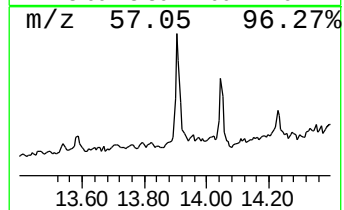
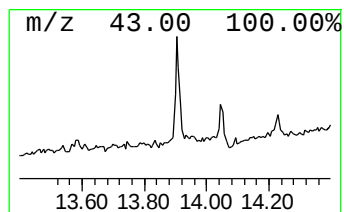
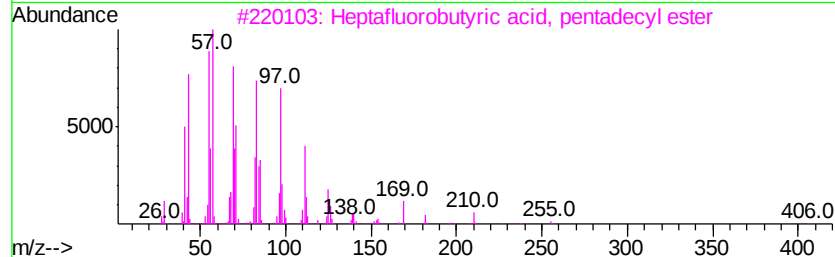
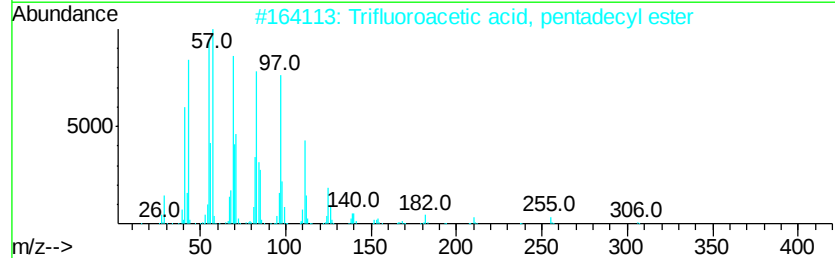
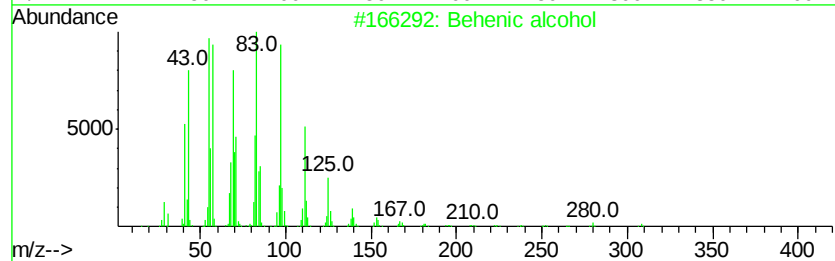
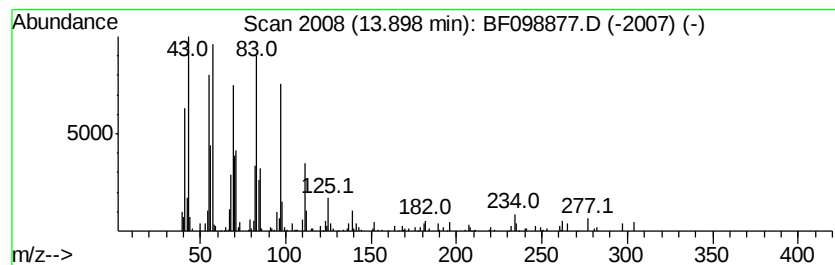
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Behenic alcohol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.	
13.90	3.78 ng	296619	Chrysene-d12		14.08	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Behenic alcohol	326	C22H46O	000661-19-8	90
2		Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	90
3		Heptafluorobutvric acid, pentade...	424	C19H31F7O2	959261-23-5	90
4		1-Nonadecene	266	C19H38	018435-45-5	90
5		Heptafluorobutyric acid, hexadec...	438	C20H33F7O2	006385-15-5	87



Data Path : Z:\HPCHEM1\BNA F\DATA\BF092717\
Data File : BF098877.D
Acq On : 27 Sep 2017 22:00
Operator : SJ/JU
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Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampled :
SA-06-092517-B

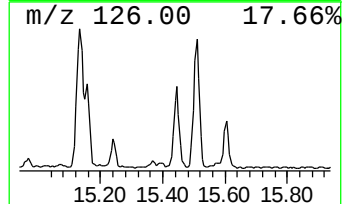
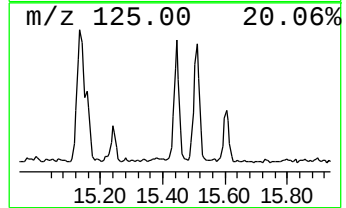
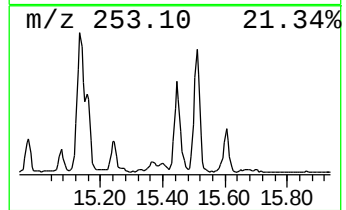
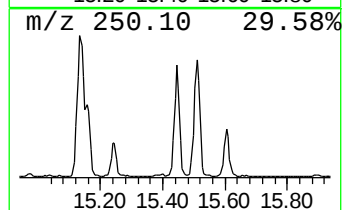
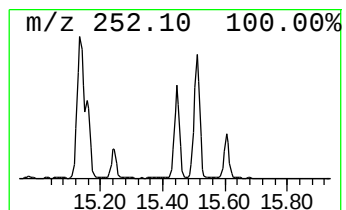
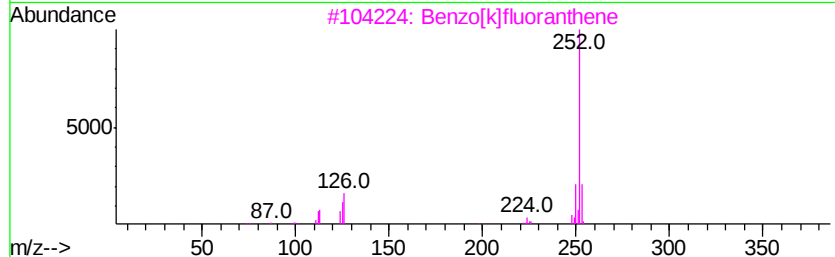
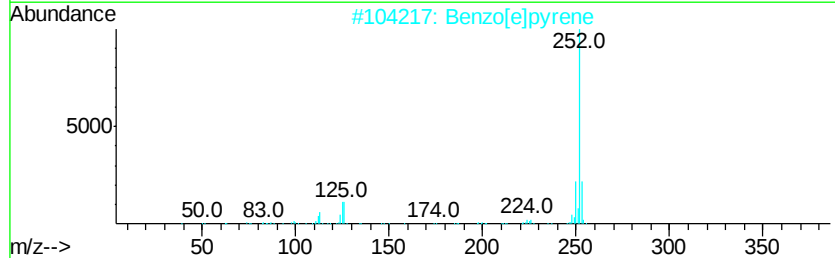
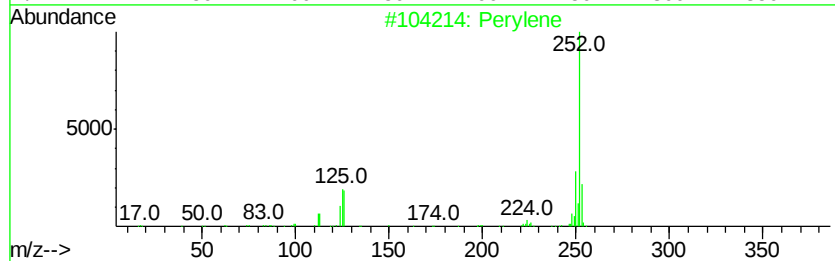
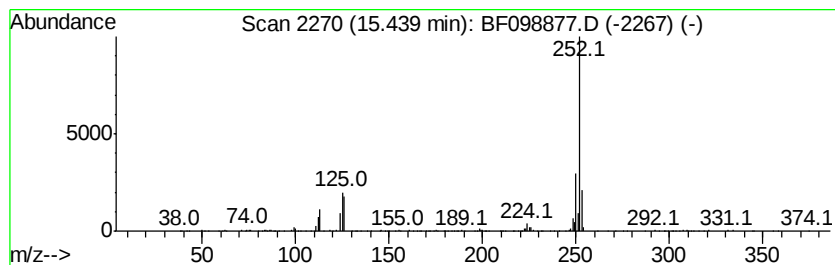
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Perylene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.44	13.93 ng	580722	Perylene-d12	15.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Perylene	252	C20H12	000198-55-0	99
2		Benzo[e]pyrene	252	C20H12	000192-97-2	98
3		Benzo[k]fluoranthene	252	C20H12	000207-08-9	97
4		Benzo[e]acephenanthrylene	252	C20H12	000205-99-2	96
5		Benzo[a]pyrene	252	C20H12	000050-32-8	96



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092717\
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Acq On : 27 Sep 2017 22:00
Operator : SJ/JU
Sample : I5423-04
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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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
Butane, 2-methoxy...	2.20	3.4	ng	156728	1	6.92	929849	20.0
2-Pentanone, 4-hy...	5.14	11.7	ng	545185	1	6.92	929849	20.0
unknown6.68	6.68	58.9	ng	2737260	1	6.92	929849	20.0
Anthracene, 1-met...	11.94	2.5	ng	151677	4	11.44	1202640	20.0
4H-Cyclopenta[def...	12.05	3.9	ng	235709	4	11.44	1202640	20.0
unknown12.53	12.53	2.6	ng	153712	4	11.44	1202640	20.0
11H-Benzo[b]fluorene	13.21	3.3	ng	261735	5	14.08	1568070	20.0
Behenic alcohol	13.90	3.8	ng	296619	5	14.08	1568070	20.0
Perylene	15.44	13.9	ng	580722	6	15.57	833761	20.0