

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092615\
 Data File : BF081825.D
 Acq On : 26 Sep 2015 13:53
 Operator : UM/IZ
 Sample : G3754-04
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-6B

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-BF092415.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	3.954	156	159	162	rVB	156084	215544	5.15%	0.382%
2	5.143	260	263	265	rVB	1096684	1678091	40.13%	2.973%
3	5.406	284	286	288	rBV	216155	195797	4.68%	0.347%
4	5.531	293	297	300	rBV2	1772676	2242038	53.62%	3.972%
5	5.771	315	318	321	rBV	476830	555899	13.29%	0.985%
6	6.457	376	378	380	rBV	376095	312382	7.47%	0.553%
7	6.491	380	381	383	rVV	126053	88765	2.12%	0.157%
8	6.560	383	387	389	rVV	1753899	2149752	51.41%	3.808%
9	6.617	390	392	393	rVV	116451	93738	2.24%	0.166%
10	6.697	397	399	401	rVV	1754577	2003941	47.92%	3.550%
11	6.766	401	405	407	rVV	566517	665699	15.92%	1.179%
12	6.811	407	409	411	rVB	193675	171258	4.10%	0.303%
13	6.937	418	420	421	rBV	933217	764850	18.29%	1.355%
14	6.960	421	422	424	rVV	213673	301894	7.22%	0.535%
15	6.994	424	425	429	rVV2	300209	404899	9.68%	0.717%
16	7.097	431	434	435	rVV	2533023	2454076	58.69%	4.347%
17	7.200	440	443	445	rVB	1094626	1015926	24.29%	1.800%
18	7.257	445	448	449	rBV	113293	93512	2.24%	0.166%
19	7.474	465	467	468	rBV	178043	238675	5.71%	0.423%
20	7.509	468	470	471	rVB	1644066	1933284	46.23%	3.425%
21	7.543	471	473	475	rVB2	91261	152392	3.64%	0.270%
22	7.577	475	476	479	rVB3	93826	100783	2.41%	0.179%
23	7.863	499	501	503	rBV2	97045	98798	2.36%	0.175%
24	7.966	507	510	512	rBV2	418425	593275	14.19%	1.051%
25	8.000	512	513	517	rVB	299011	535223	12.80%	0.948%
26	8.092	520	521	523	rVB	155820	133904	3.20%	0.237%
27	8.252	531	535	537	rBV2	3337815	4181734	100.00%	7.408%
28	8.457	549	553	556	rVB3	130138	236222	5.65%	0.418%
29	8.572	559	563	566	rBV2	129221	226245	5.41%	0.401%
30	8.629	566	568	569	rBV2	50225	93855	2.24%	0.166%
31	8.812	582	584	586	rBV	446022	629478	15.05%	1.115%
32	8.857	586	588	589	rBV	265422	295423	7.06%	0.523%
33	8.937	594	595	597	rVB	1214363	910466	21.77%	1.613%
34	9.040	602	604	606	rVV	1188507	1120182	26.79%	1.984%

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35	9.075	606	607	611	rVB3	86774	146996	3.52%	0.260%
36	9.223	617	620	621	rBV2	131539	229200	5.48%	0.406%
37	9.246	621	622	624	rVV	145516	145633	3.48%	0.258%
38	9.303	624	627	629	rVV	3727257	3423452	81.87%	6.065%
39	9.372	631	633	634	rVV	126633	126802	3.03%	0.225%
40	9.406	634	636	638	rVV	254545	387449	9.27%	0.686%
41	9.463	638	641	643	rVV4	357547	588857	14.08%	1.043%
42	9.497	643	644	646	rVV	223528	273426	6.54%	0.484%
43	9.543	646	648	649	rVV	248557	374865	8.96%	0.664%
44	9.635	655	656	657	rVV	211077	210479	5.03%	0.373%
45	9.669	657	659	660	rVV2	215401	303737	7.26%	0.538%
46	9.703	660	662	664	rVV2	452266	451223	10.79%	0.799%
47	9.760	664	667	670	rVB4	68810	157250	3.76%	0.279%
48	9.817	670	672	673	rBV	253759	315161	7.54%	0.558%
49	9.840	673	674	677	rVV	1525432	1431515	34.23%	2.536%
50	9.943	681	683	684	rVV	115248	175591	4.20%	0.311%
51	9.977	684	686	688	rVV2	928580	1097995	26.26%	1.945%
52	10.012	688	689	693	rVV	470900	437956	10.47%	0.776%
53	10.069	693	694	697	rVV2	162842	179561	4.29%	0.318%
54	10.138	698	700	703	rVV4	188572	336443	8.05%	0.596%
55	10.183	703	704	711	rVB2	215753	384375	9.19%	0.681%
56	10.286	711	713	716	rBV3	117858	208441	4.98%	0.369%
57	10.378	719	721	724	rVB3	100578	126479	3.02%	0.224%
58	10.435	724	726	727	rVB	158765	123698	2.96%	0.219%
59	10.469	727	729	730	rBV	104475	121945	2.92%	0.216%
60	10.503	730	732	733	rVV	111827	148958	3.56%	0.264%
61	10.526	733	734	737	rVB	589050	511986	12.24%	0.907%
62	10.595	737	740	742	rBV3	69712	126723	3.03%	0.224%
63	10.675	746	747	749	rVV2	81666	88810	2.12%	0.157%
64	10.766	752	755	757	rVV2	1539382	1651519	39.49%	2.926%
65	10.892	763	766	769	rVB2	204847	234412	5.61%	0.415%
66	10.960	769	772	776	rVB5	55086	120279	2.88%	0.213%
67	11.063	776	781	783	rBV2	88595	181507	4.34%	0.322%
68	11.269	796	799	802	rVB2	90917	179379	4.29%	0.318%
69	11.463	815	816	817	rBV	736396	966168	23.10%	1.712%
70	11.498	817	819	820	rVB	1579953	1634024	39.08%	2.895%
71	11.543	820	823	825	rBV	449459	380486	9.10%	0.674%

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72	11.772	840	843	845	rVV3	74653	146972	3.51%	0.260%
73	11.818	845	847	849	rVV3	52098	104434	2.50%	0.185%
74	11.989	858	862	865	rBV2	308887	590341	14.12%	1.046%
75	12.081	868	870	874	rVV2	339988	602830	14.42%	1.068%
76	12.263	882	886	891	rVB2	153610	213586	5.11%	0.378%
77	12.412	896	899	903	rBV3	157979	246808	5.90%	0.437%
78	12.515	906	908	910	rVV	110374	157027	3.76%	0.278%
79	12.572	912	913	915	rVV	197967	174138	4.16%	0.308%
80	12.686	921	923	925	rVV	455389	636593	15.22%	1.128%
81	12.721	925	926	929	rVV2	111529	143259	3.43%	0.254%
82	12.778	929	931	933	rVV	228675	231614	5.54%	0.410%
83	12.846	935	937	940	rBV3	48456	90443	2.16%	0.160%
84	12.915	940	943	945	rVV	759695	971244	23.23%	1.721%
85	12.972	945	948	950	rVB	109371	124634	2.98%	0.221%
86	13.064	953	956	958	rVV	2250137	3229612	77.23%	5.721%
87	13.121	960	961	964	rBV2	71910	134174	3.21%	0.238%
88	13.189	965	967	969	rVV	327083	299333	7.16%	0.530%
89	13.235	969	971	974	rVV	95674	123137	2.94%	0.218%
90	13.304	974	977	979	rVV	129099	174710	4.18%	0.309%
91	13.338	979	980	984	rVB2	64597	99732	2.38%	0.177%
92	13.441	984	989	990	rBV2	85922	146681	3.51%	0.260%
93	13.909	1027	1030	1032	rBV2	68875	113670	2.72%	0.201%
94	13.944	1032	1033	1035	rVB	87865	86721	2.07%	0.154%
95	14.104	1044	1047	1052	rVV2	802506	1187478	28.40%	2.104%
96	15.144	1135	1138	1143	rBV	75219	168018	4.02%	0.298%
97	15.509	1168	1170	1173	rVB	125497	149402	3.57%	0.265%
98	15.578	1173	1176	1184	rBV	549558	708581	16.94%	1.255%
99	17.475	1338	1342	1347	rVB	41966	91647	2.19%	0.162%
100	21.476	1684	1692	1698	rBV	97652	431960	10.33%	0.765%

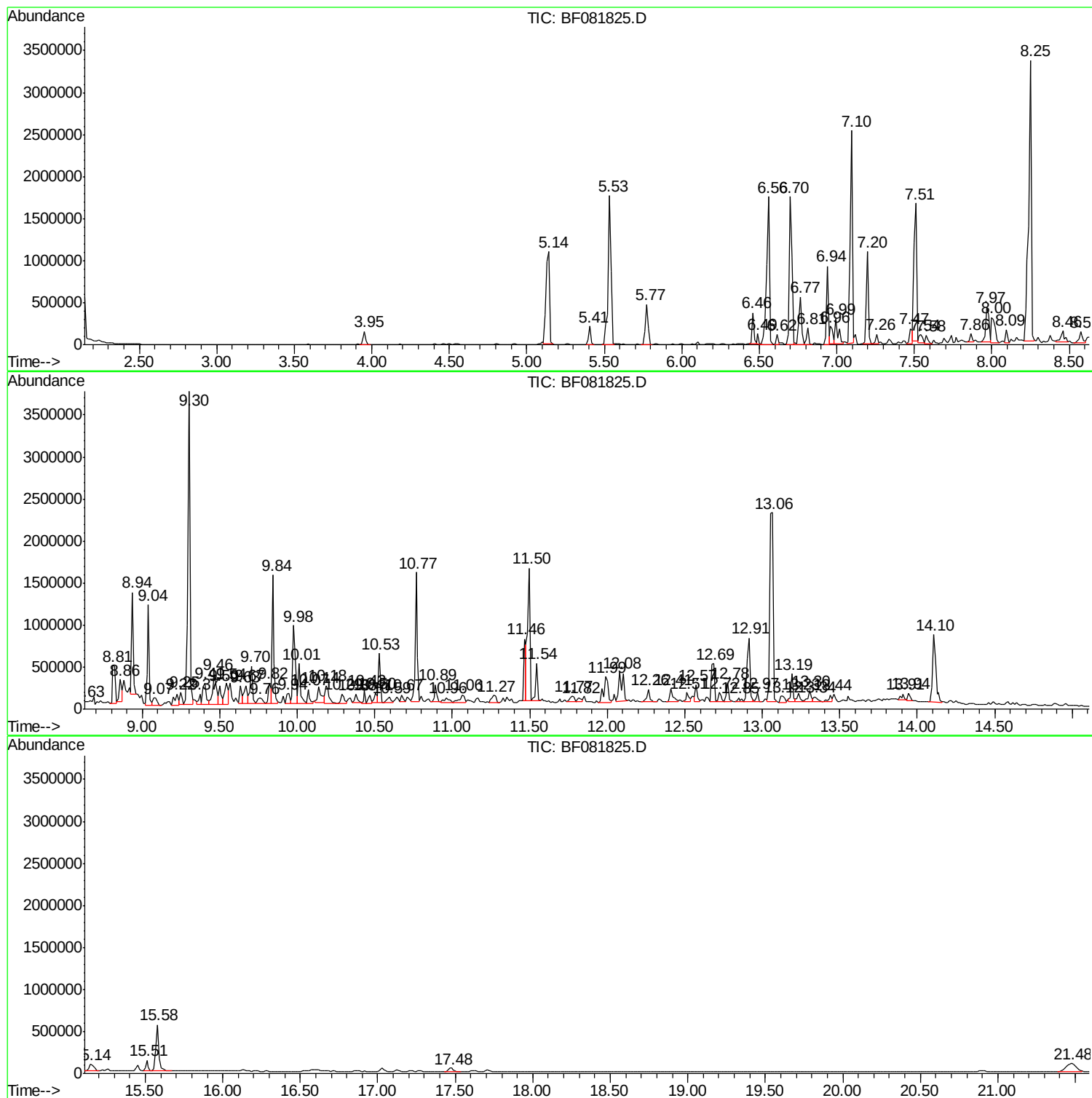
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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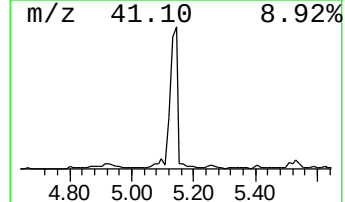
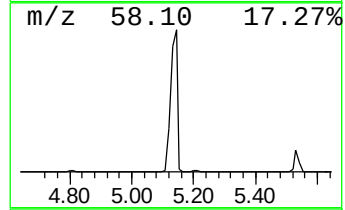
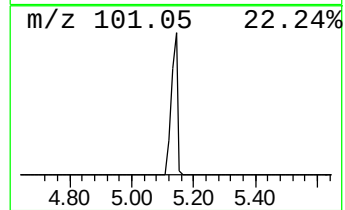
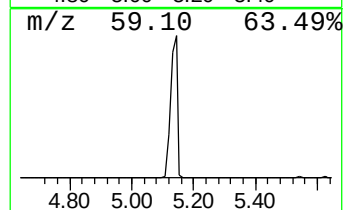
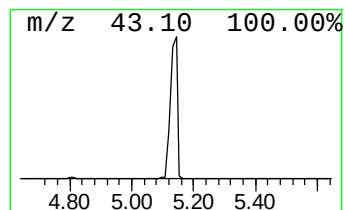
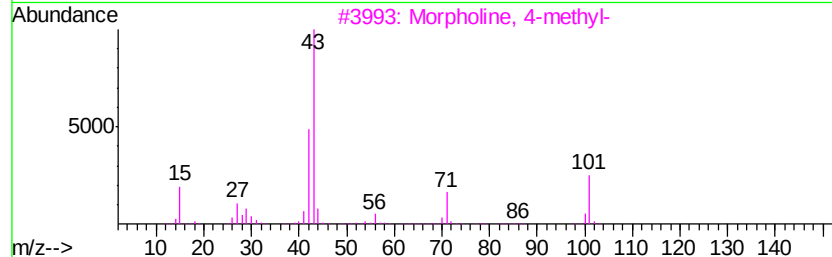
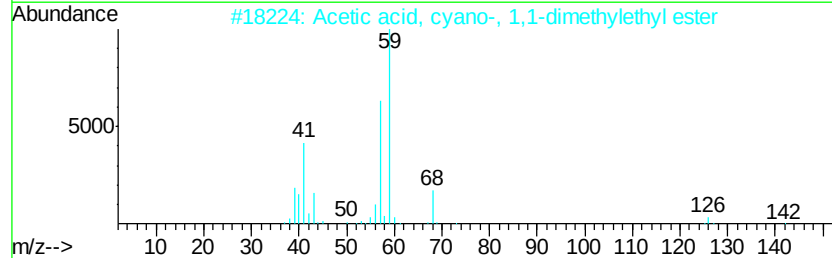
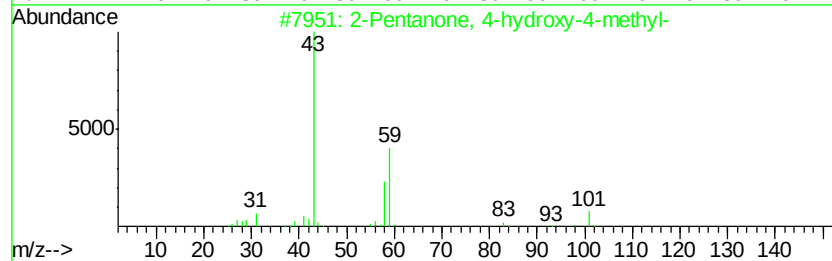
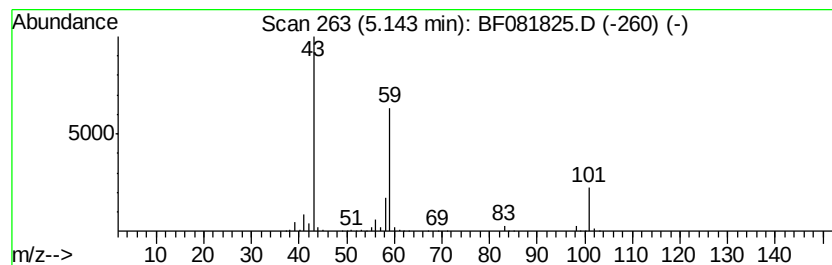
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.14	43.88 ng	1678090	1,4-Dichlorobenzene-d4	6.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
3		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4		Acetone	58	C3H6O	000067-64-1	9
5		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	9



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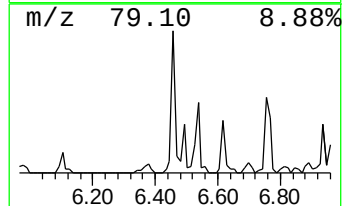
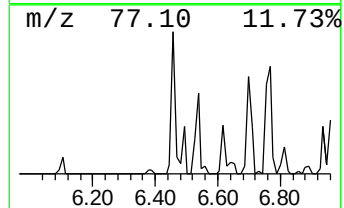
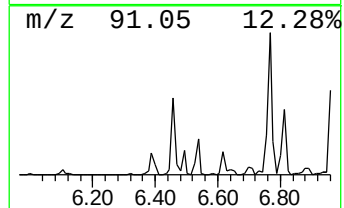
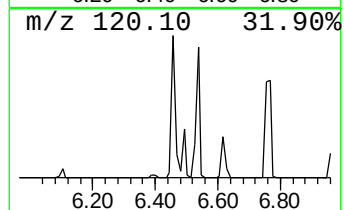
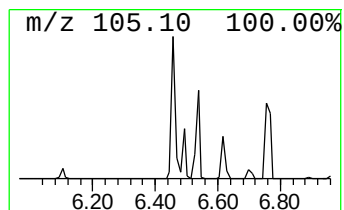
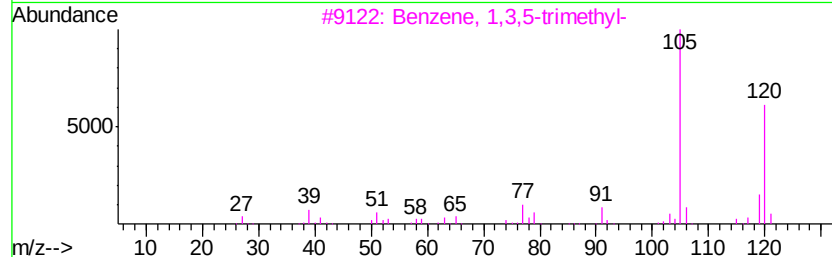
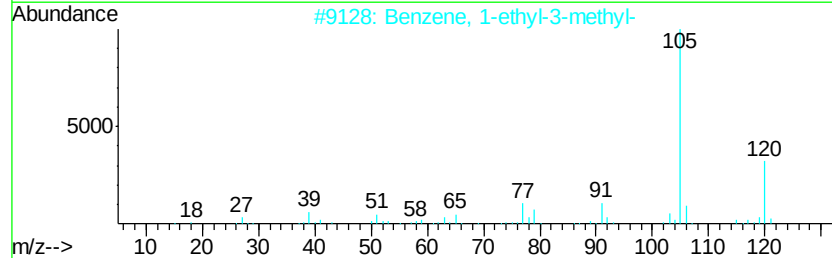
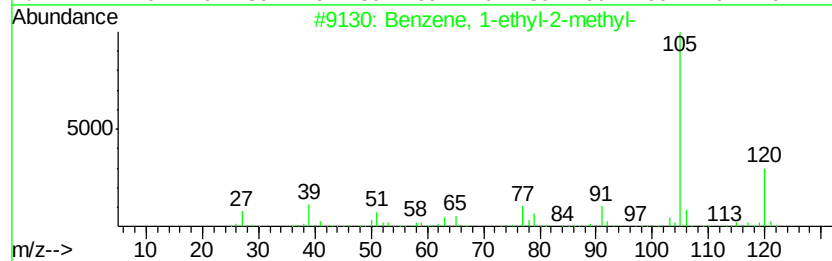
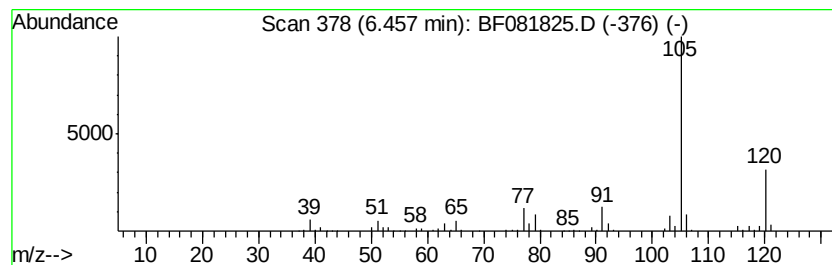
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, 1-ethyl-2-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.46	8.17 ng	312382	1,4-Dichlorobenzene-d4	6.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
3		Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	91
4		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



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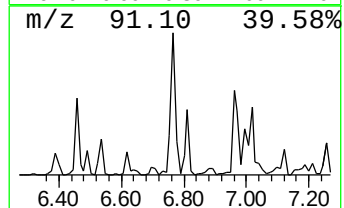
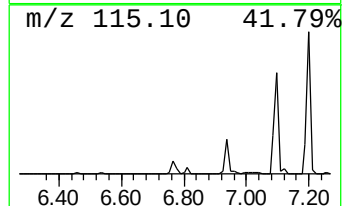
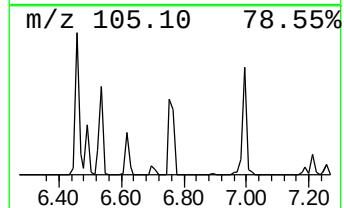
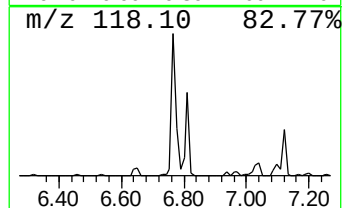
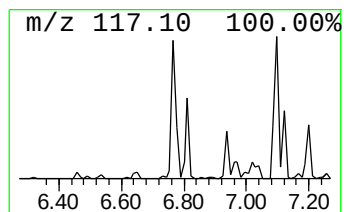
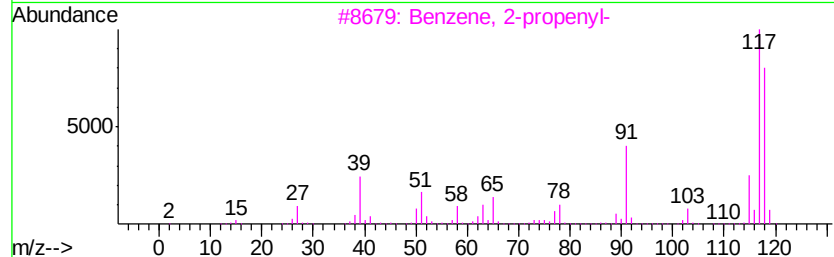
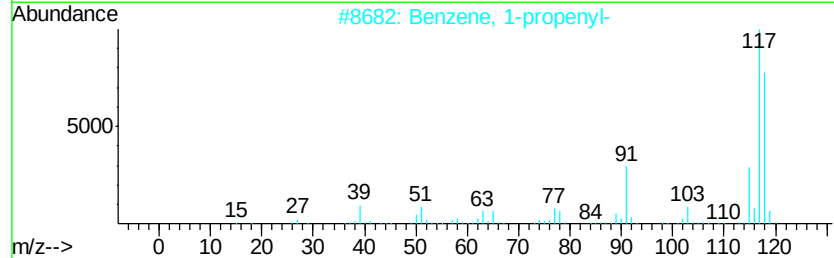
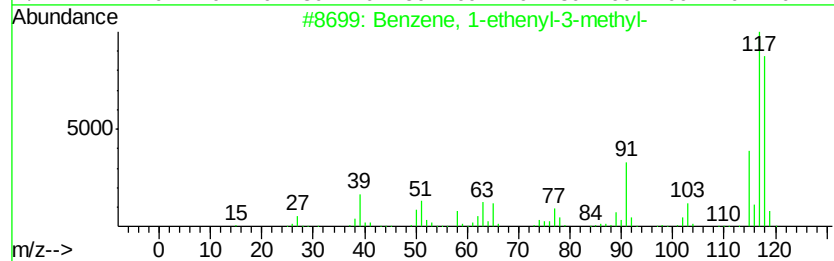
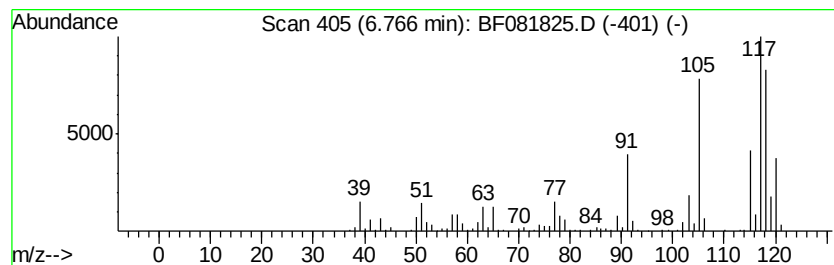
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Peak Number 7 Benzene, 1-ethenyl-3-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.77	17.41 ng	665699	1,4-Dichlorobenzene-d4	6.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	92
2		Benzene, 1-propenyl-	118	C9H10	000637-50-3	91
3		Benzene, 2-propenyl-	118	C9H10	000300-57-2	90
4		Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	64
5		Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	64



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Operator : UM/IZ
Sample : G3754-04
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-6B

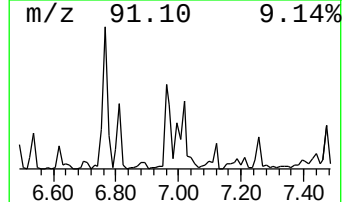
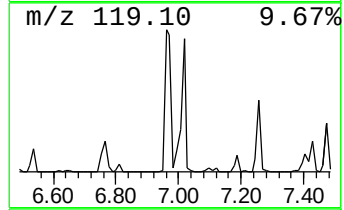
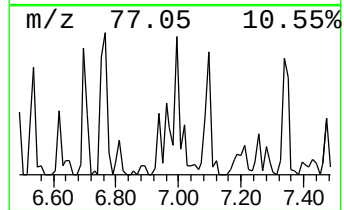
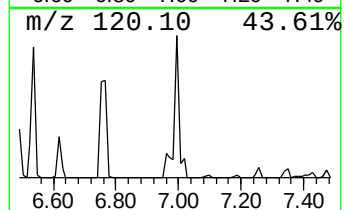
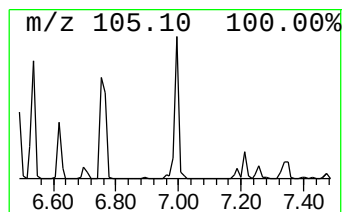
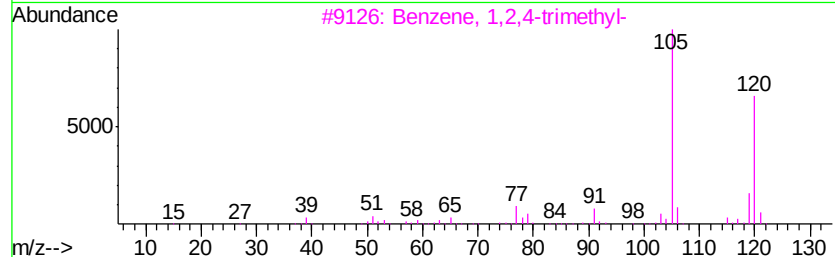
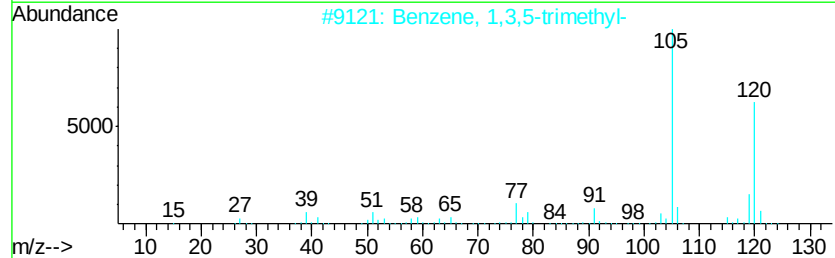
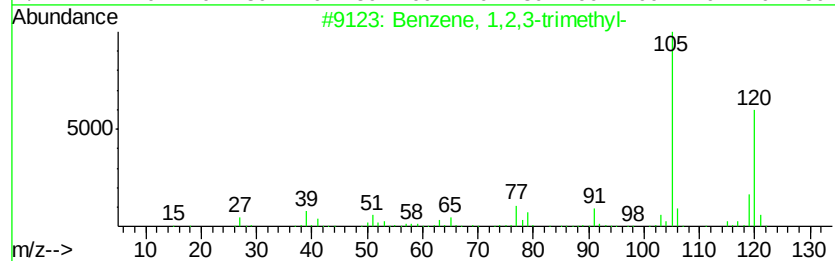
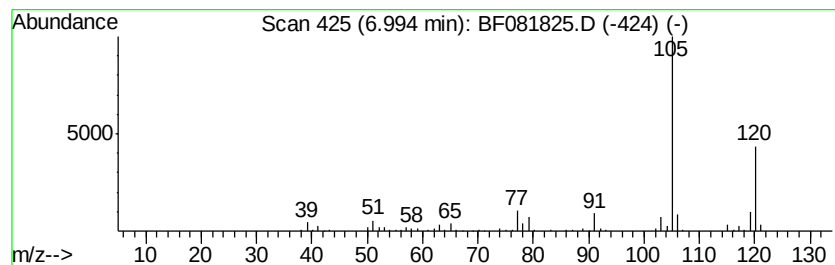
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Benzene, 1,2,3-trimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.99	10.59 ng	404899	1,4-Dichlorobenzene-d4	6.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2		Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	94
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90
5		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092615\
Data File : BF081825.D
Acq On : 26 Sep 2015 13:53
Operator : UM/IZ
Sample : G3754-04
Misc :
ALS Vial : 39 Sample Multiplier: 1

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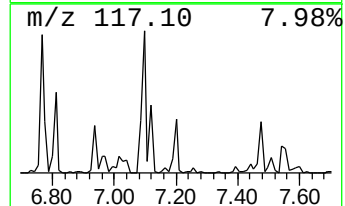
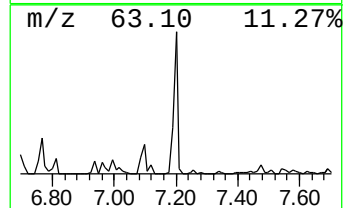
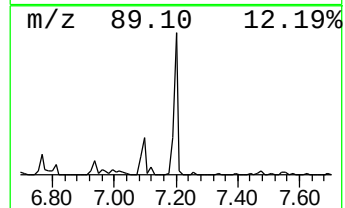
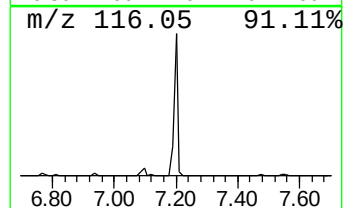
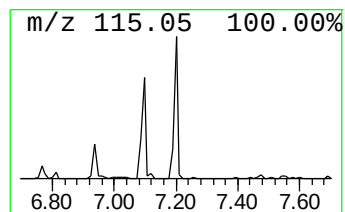
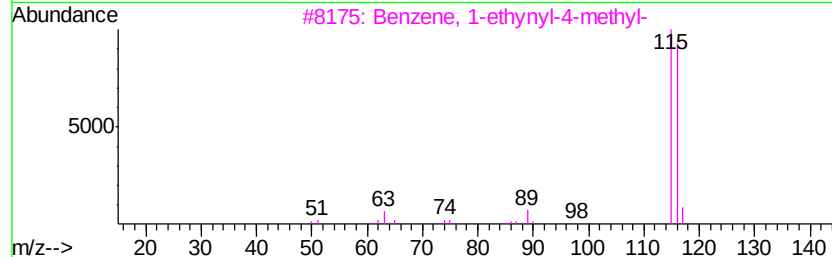
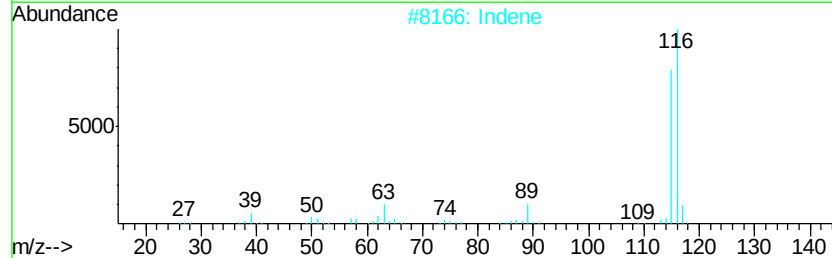
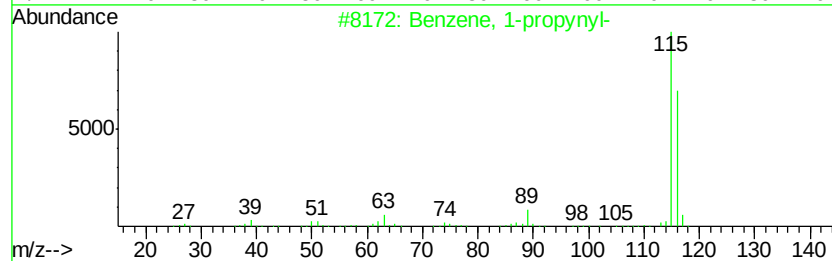
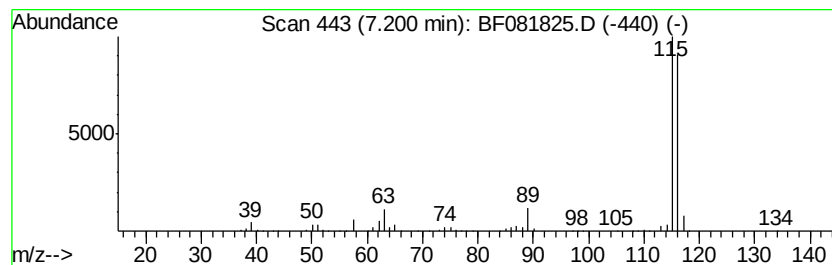
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Benzene, 1-propynyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.20	26.57 ng	1015930	1,4-Dichlorobenzene-d4	6.94

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-propynyl-	116	C9H8	000673-32-5	94
2	Indene	116	C9H8	000095-13-6	93
3	Benzene, 1-ethynyl-4-methyl-	116	C9H8	000766-97-2	91
4	Benzene, 1-ethynyl-2-methyl-	116	C9H8	000766-47-2	91
5	Benzene, 1,2-propadienyl-	116	C9H8	002327-99-3	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092615\
Data File : BF081825.D
Acq On : 26 Sep 2015 13:53
Operator : UM/IZ
Sample : G3754-04
Misc :
ALS Vial : 39 Sample Multiplier: 1

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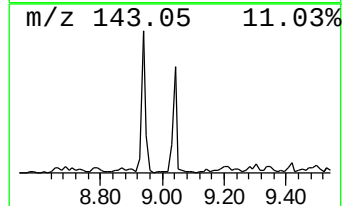
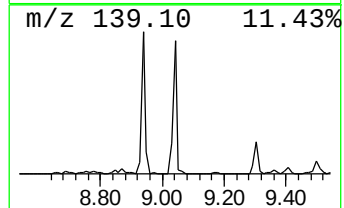
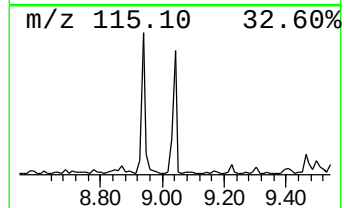
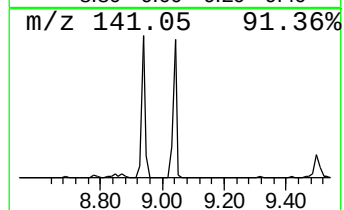
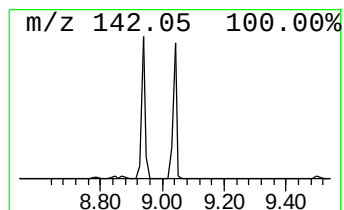
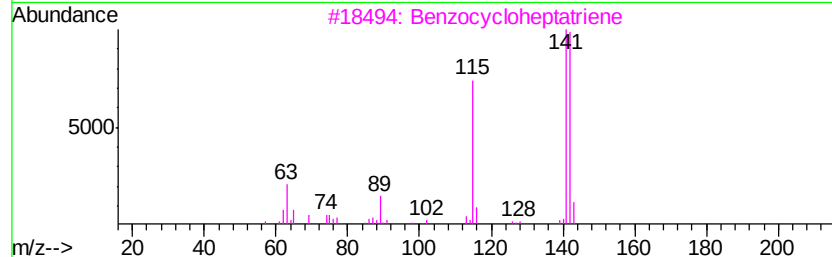
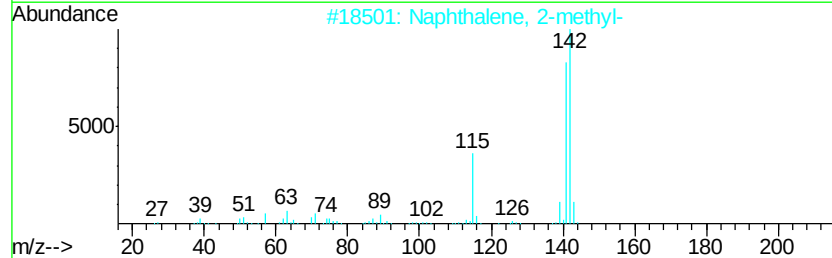
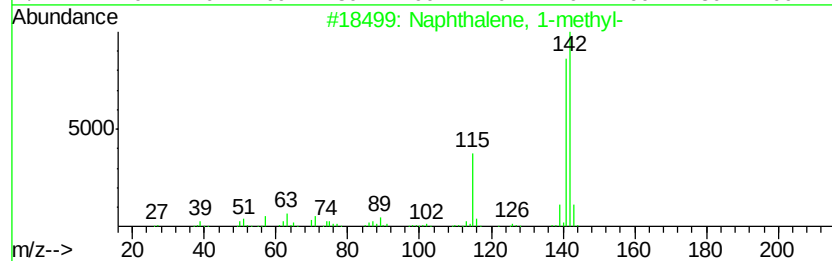
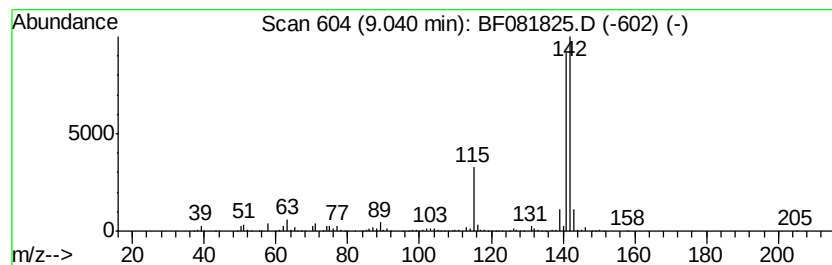
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Naphthalene, 1-methyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.04	5.36 ng	1120180	Naphthalene-d8	8.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3		Benzocycloheptatriene	142	C11H10	000264-09-5	91
4		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
5		1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092615\
Data File : BF081825.D
Acq On : 26 Sep 2015 13:53
Operator : UM/IZ
Sample : G3754-04
Misc :
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Instrument :
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ClientSampleId :
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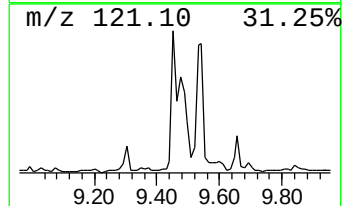
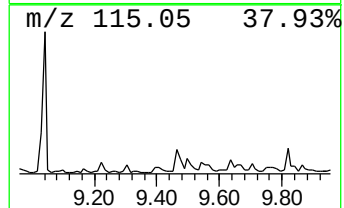
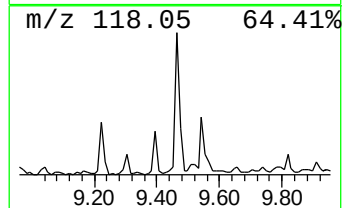
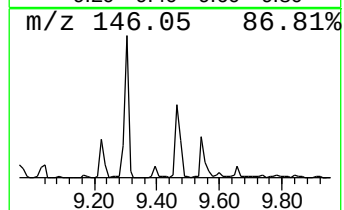
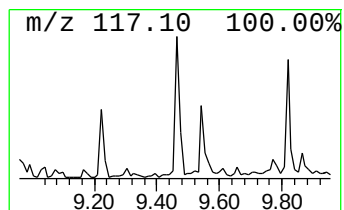
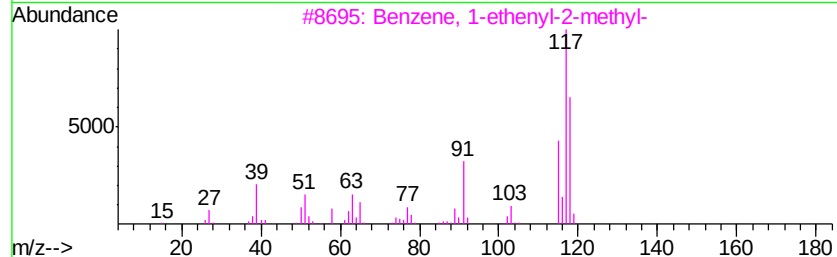
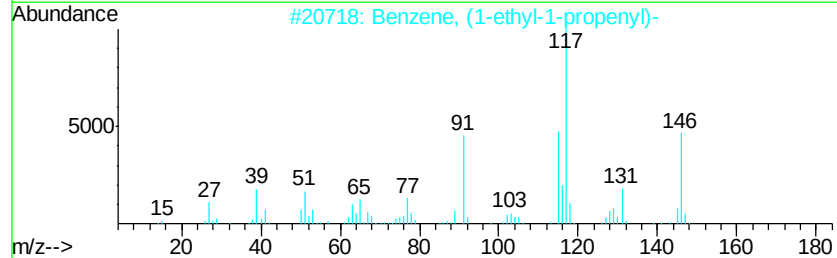
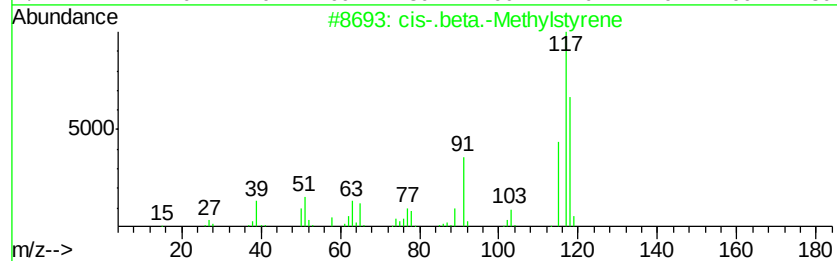
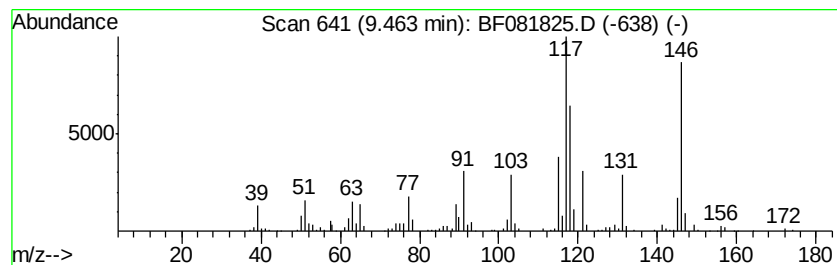
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 cis-.beta.-Methylstyrene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.46	10.73 ng	588857	Acenaphthene-d10	9.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	cis-.beta.-Methylstyrene	118	C9H10	000766-90-5	86
2		Benzene, (1-ethyl-1-propenyl)-	146	C11H14	004701-36-4	76
3		Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	64
4		Benzene, 2-propenyl-	118	C9H10	000300-57-2	62
5		Benzene, 1-pentenyl-	146	C11H14	000826-18-6	58



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092615\
Data File : BF081825.D
Acq On : 26 Sep 2015 13:53
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Sample : G3754-04
Misc :
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Instrument :
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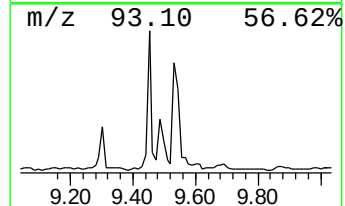
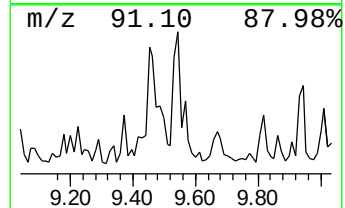
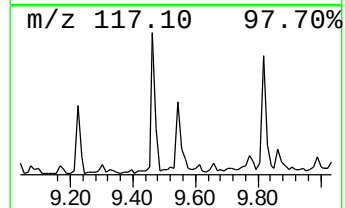
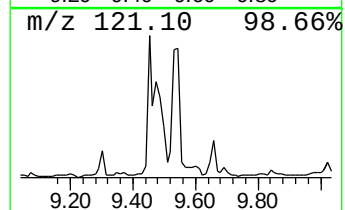
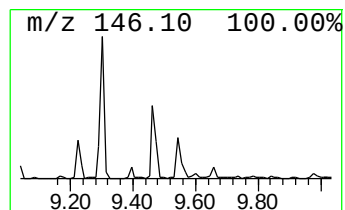
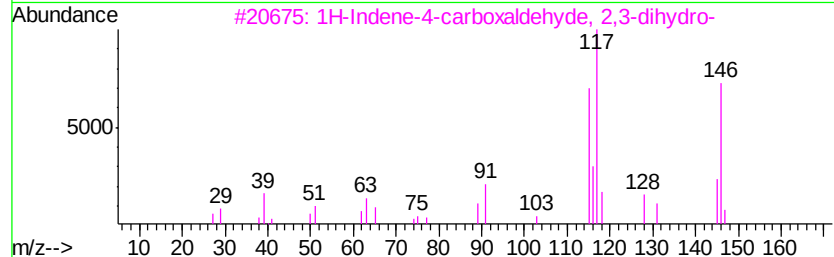
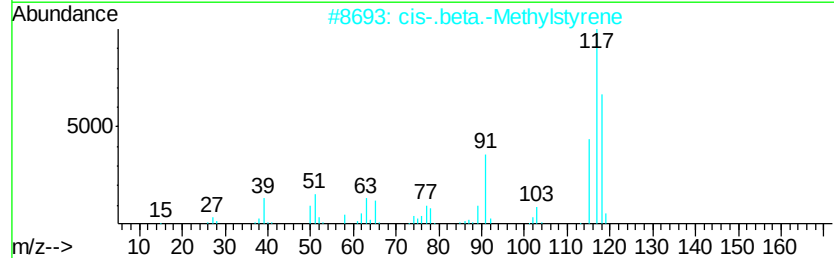
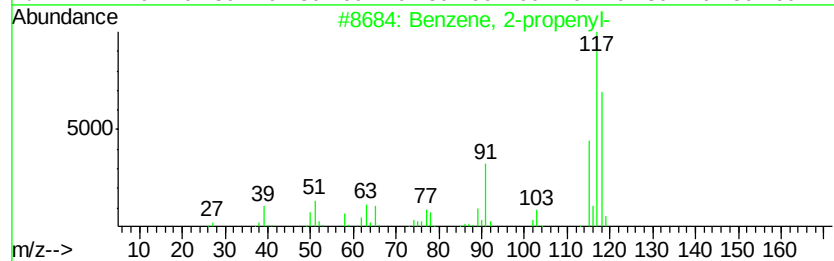
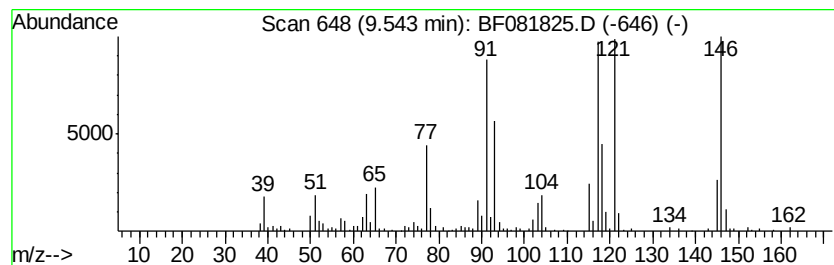
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 unknown9.54 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.54	6.83 ng	374865	Acenaphthene-d10	9.98

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-propenyl-	118	C9H10	000300-57-2	43
2			cis-.beta.-Methylstyrene	118	C9H10	000766-90-5	43
3			1H-Indene-4-carboxaldehyde, 2,3-...	146	C10H10O	051932-70-8	38
4			7-Methylindan-1-one	146	C10H10O	039627-61-7	38
5			1H-1,3,2-Benzodiazaborole, 2-eth...	146	C8H11BN2	074663-81-3	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092615\
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Acq On : 26 Sep 2015 13:53
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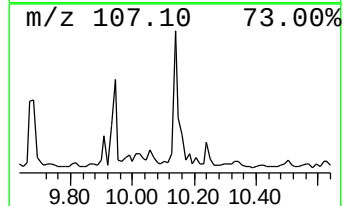
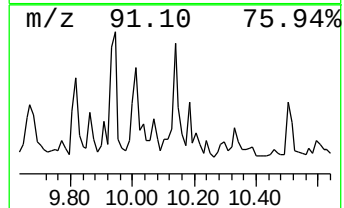
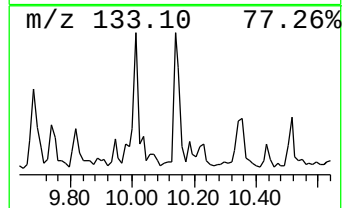
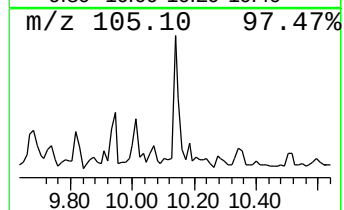
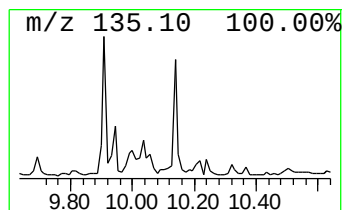
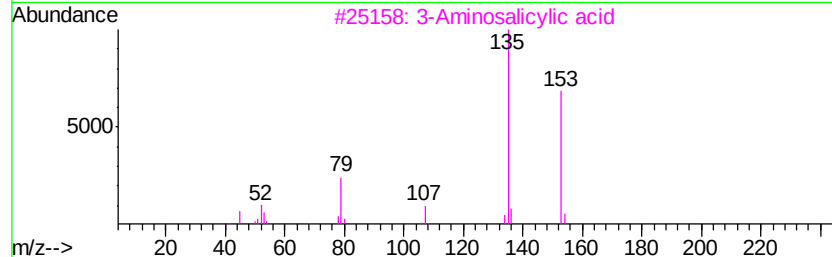
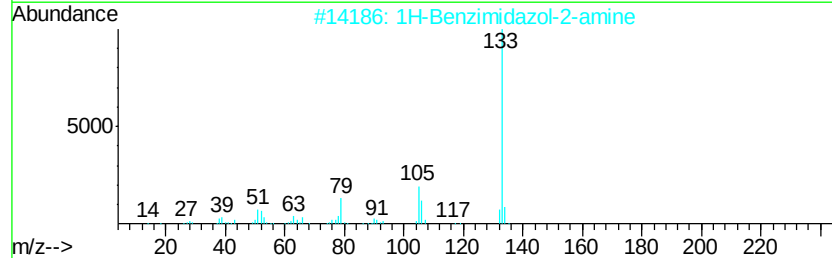
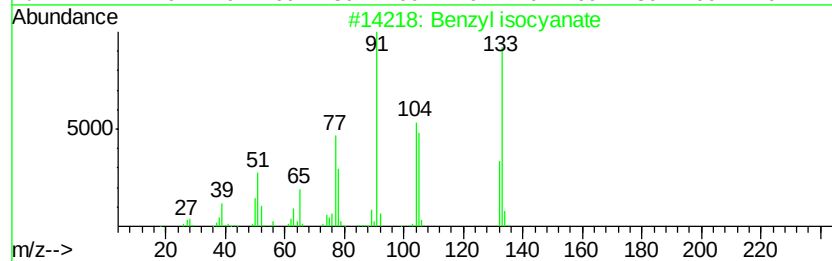
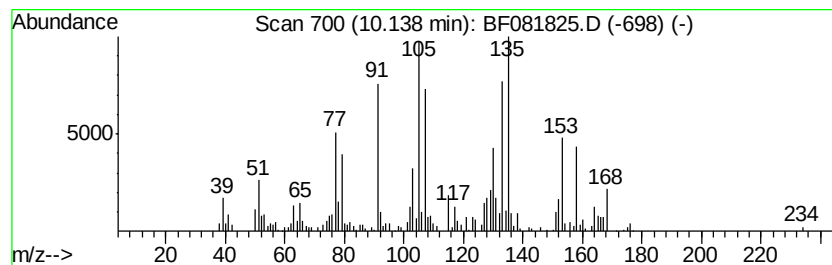
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ClientSampled :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 unknown10.14 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.14	6.13 ng	336443	Acenaphthene-d10	9.98
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzyl isocyanate	133 C8H7NO	003173-56-6 25
2		1H-Benzimidazol-2-amine	133 C7H7N3	000934-32-7 20
3		3-Aminosalicylic acid	153 C7H7NO3	000570-23-0 18
4		Aminosalicylic Acid	153 C7H7NO3	000065-49-6 18
5		1H-Indol-4-ol	133 C8H7NO	002380-94-1 11



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092615\
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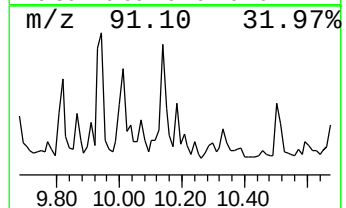
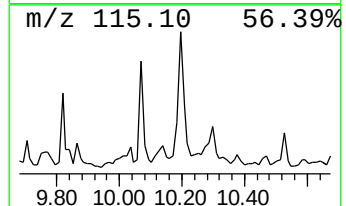
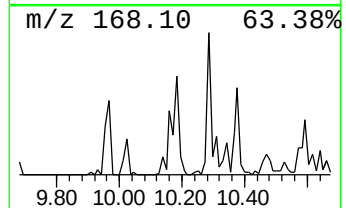
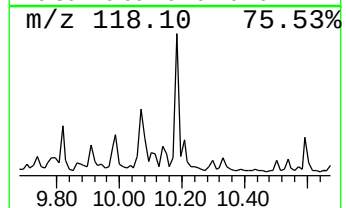
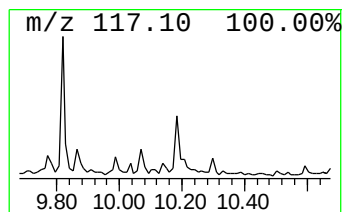
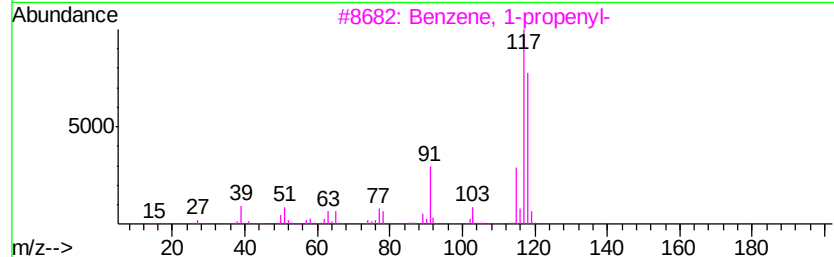
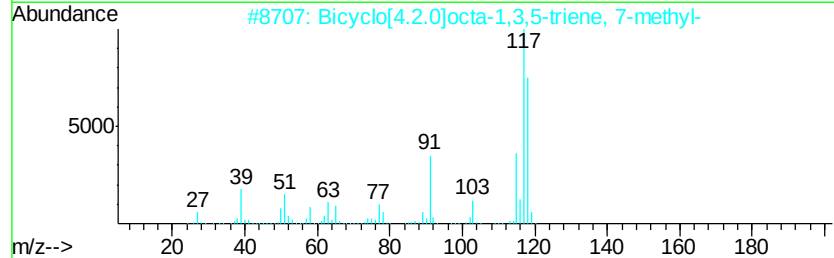
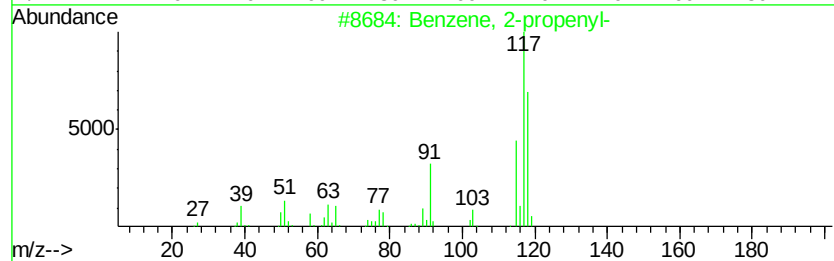
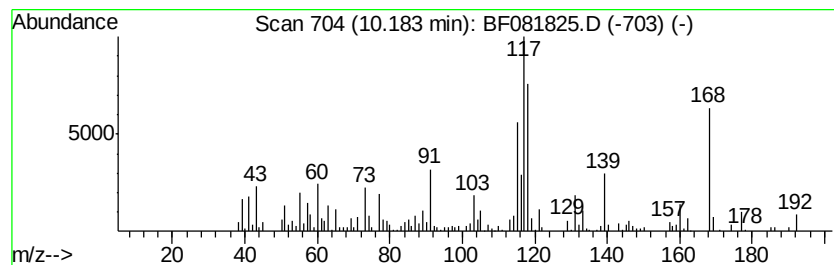
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 Bicyclo[4.2.0]octa-1,3,5-tr... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.18	7.00 ng	384375	Acenaphthene-d10	9.98
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 2-propenyl-	118 C9H10	000300-57-2 86
2		Bicyclo[4.2.0]octa-1,3,5-triene, ...	118 C9H10	055337-80-9 64
3		Benzene, 1-propenyl-	118 C9H10	000637-50-3 50
4		Benzene, 1-ethenyl-4-methyl-	118 C9H10	000622-97-9 50
5		Benzene, 1-ethenyl-2-methyl-	118 C9H10	000611-15-4 50



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092615\
Data File : BF081825.D
Acq On : 26 Sep 2015 13:53
Operator : UM/IZ
Sample : G3754-04
Misc :
ALS Vial : 39 Sample Multiplier: 1

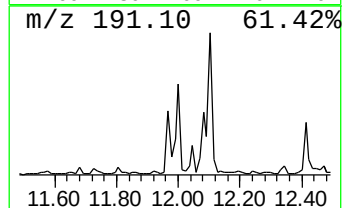
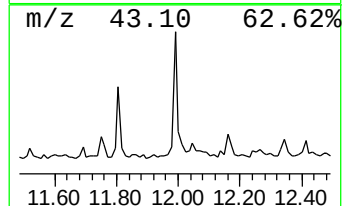
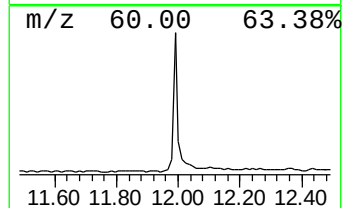
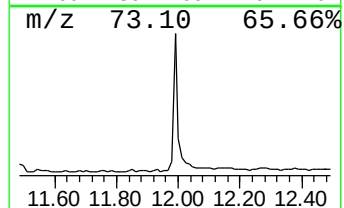
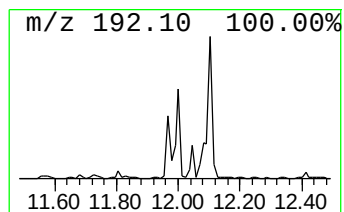
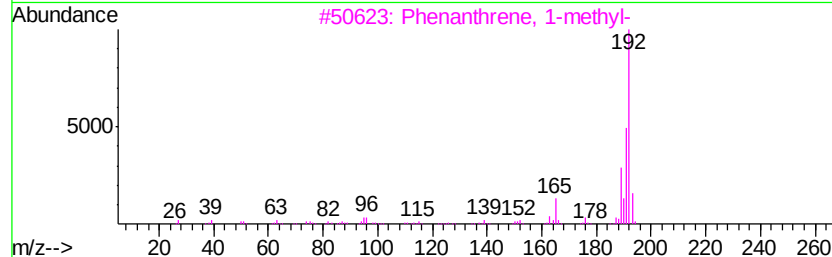
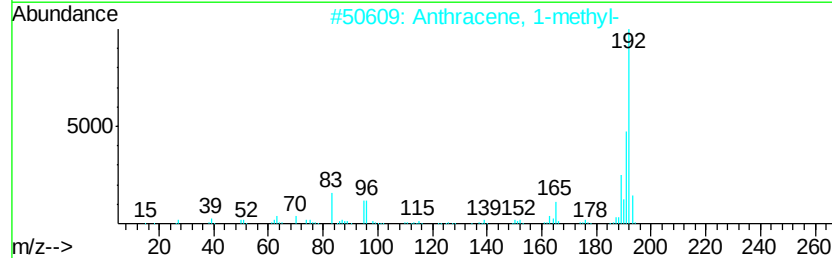
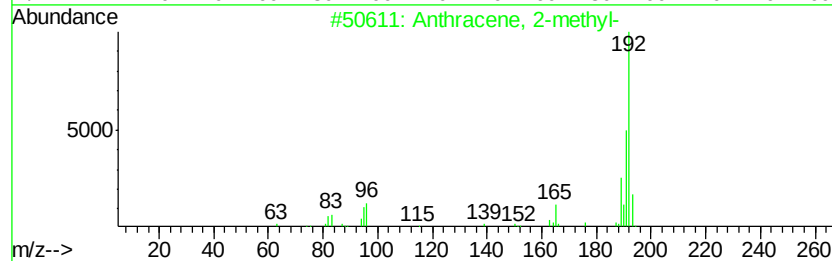
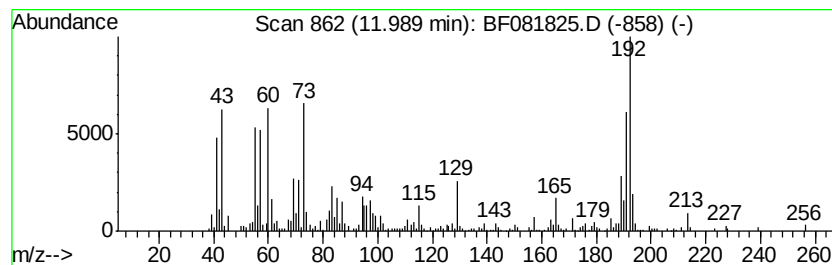
Instrument :
BNA_F
ClientSampleId :
MW-6B

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 Anthracene, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.99	12.22 ng	590341	Phenanthrene-d10	11.47	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
2	Anthracene, 1-methyl-	192	C15H12	000610-48-0	96
3	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	92
4	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	92
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092615\
Data File : BF081825.D
Acq On : 26 Sep 2015 13:53
Operator : UM/IZ
Sample : G3754-04
Misc :
ALS Vial : 39 Sample Multiplier: 1

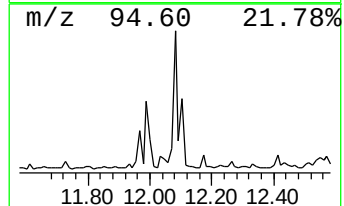
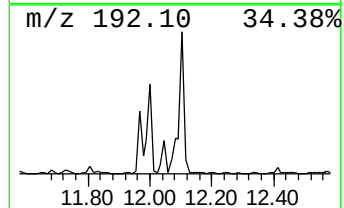
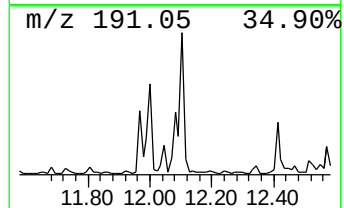
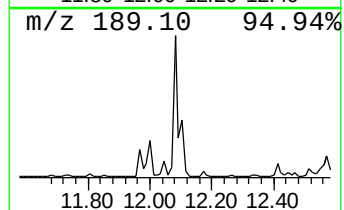
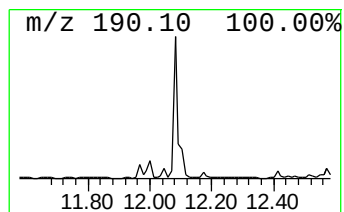
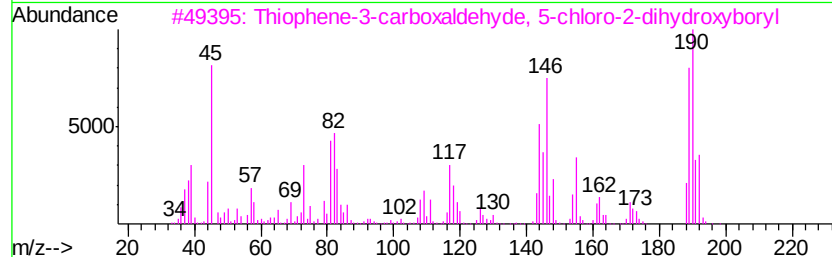
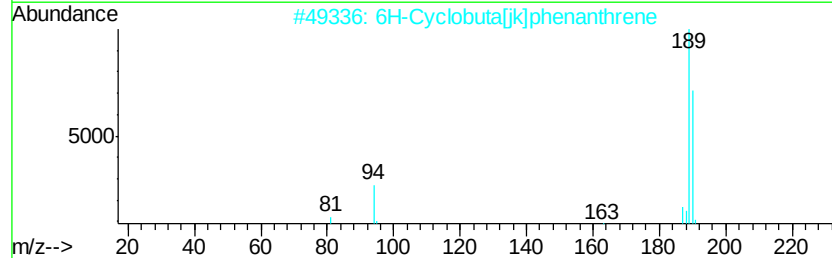
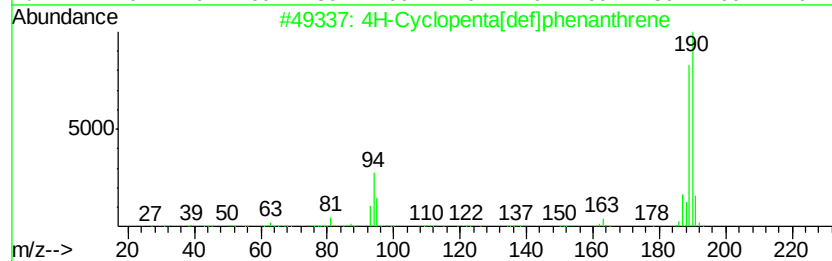
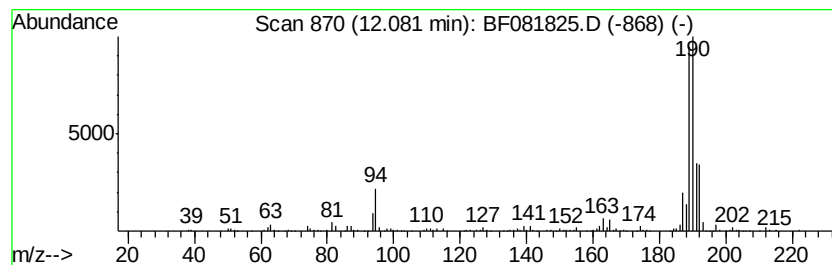
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 17 4H-Cyclopenta[def]phenanthrene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.		
12.08	12.48 ng	602830	Phenanthrene-d10	11.47		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	93
2		6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	56
3		Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	53
4		2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	50
5		2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	45



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092615\
Data File : BF081825.D
Acq On : 26 Sep 2015 13:53
Operator : UM/IZ
Sample : G3754-04
Misc :
ALS Vial : 39 Sample Multiplier: 1

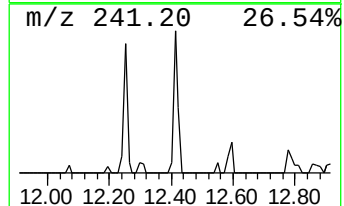
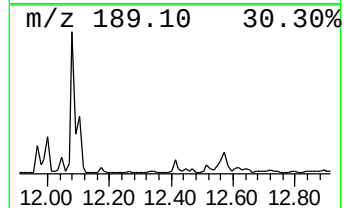
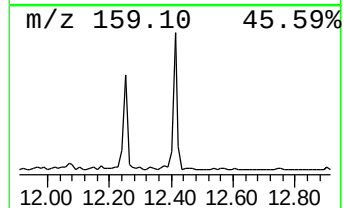
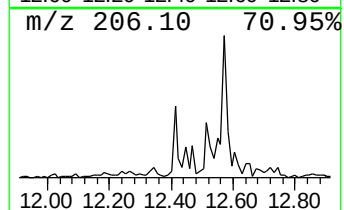
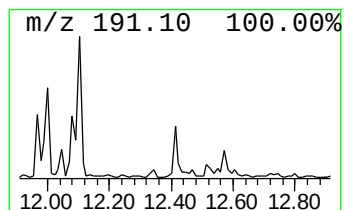
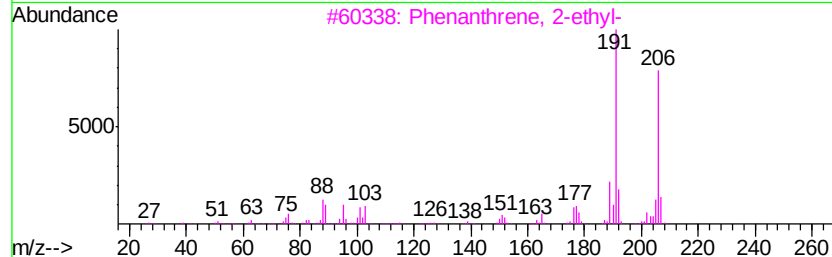
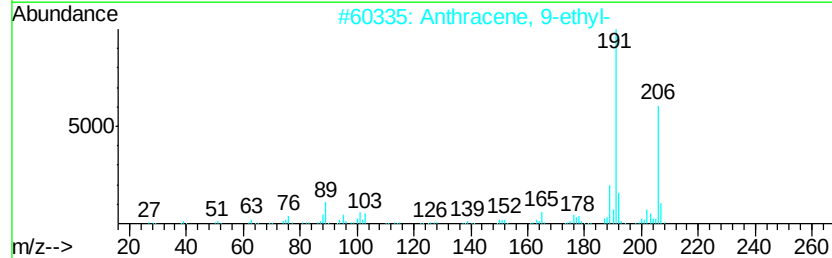
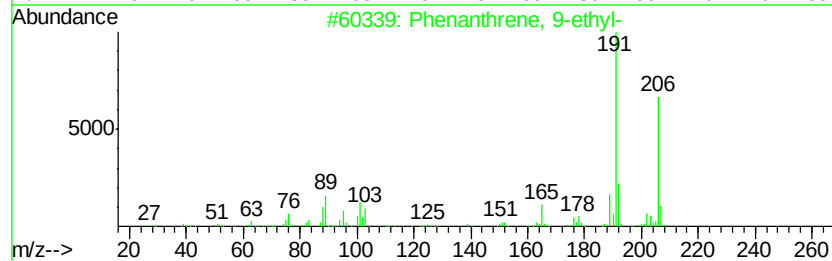
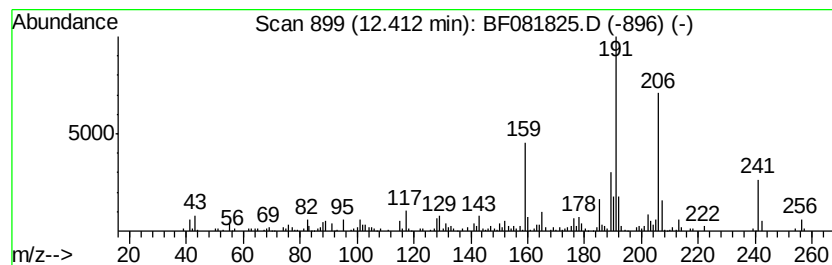
Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 18 Phenanthrene, 9-ethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD		R.T.
12.41	5.11 ng	246808	Phenanthrene-d10		11.47
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 9-ethyl-	206	C16H14	003674-75-7	93
2	Anthracene, 9-ethyl-	206	C16H14	000605-83-4	70
3	Phenanthrene, 2-ethyl-	206	C16H14	003674-74-6	64
4	Phenanthrene, 4,5-dimethyl-	206	C16H14	003674-69-9	58
5	Anthracene, 2-ethyl-	206	C16H14	052251-71-5	58



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092615\
Data File : BF081825.D
Acq On : 26 Sep 2015 13:53
Operator : UM/IZ
Sample : G3754-04
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-6B

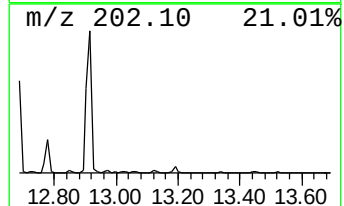
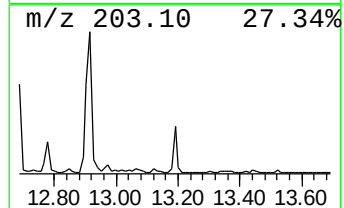
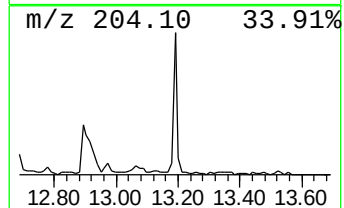
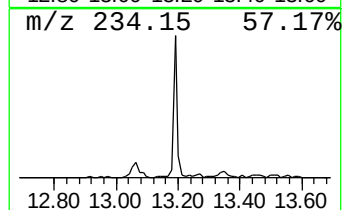
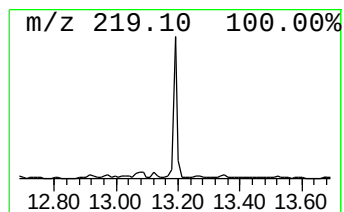
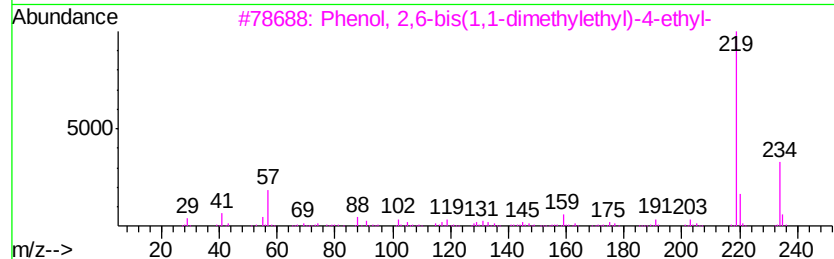
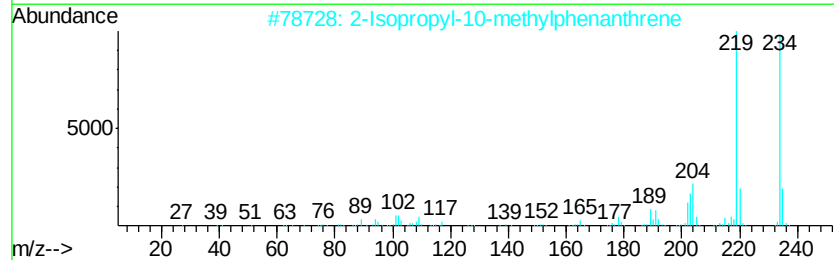
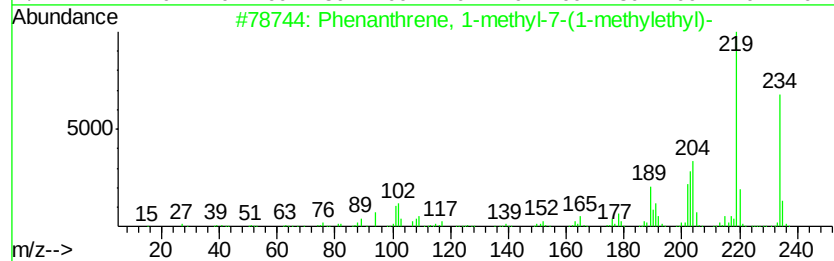
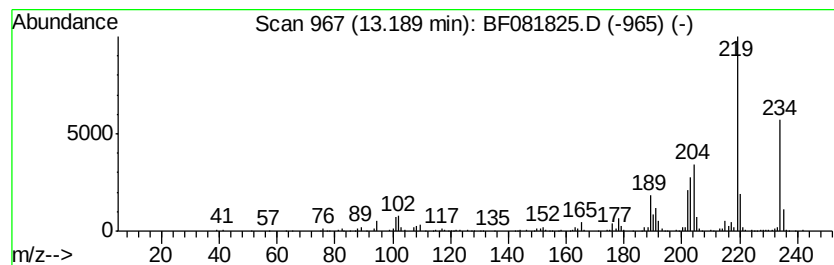
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 19 Phenanthrene, 1-methyl-7-(1... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.19	5.04 ng	299333	Chrysene-d12	14.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 1-methyl-7-(1-meth...	234	C18H18	000483-65-8	95
2		2-Isopropyl-10-methylphenanthrene	234	C18H18	066552-97-4	95
3		Phenol, 2,6-bis(1,1-dimethylethy...	234	C16H26O	004130-42-1	52
4		Benzene, 1,3-bis(1,1-dimethyleth...	234	C16H26O	001518-53-2	52
5		Ethanone, 1-(4,6-dihydroxy-2,3,5...	234	C13H14O4	021987-07-5	47



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092615\
Data File : BF081825.D
Acq On : 26 Sep 2015 13:53
Operator : UM/IZ
Sample : G3754-04
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-6B

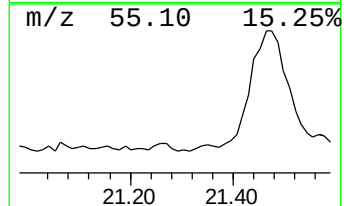
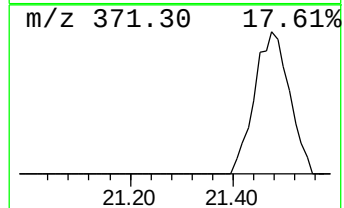
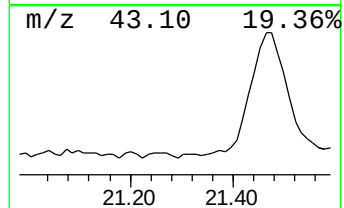
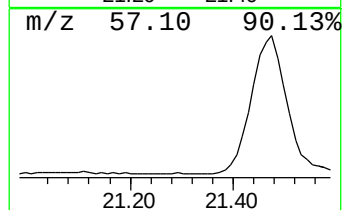
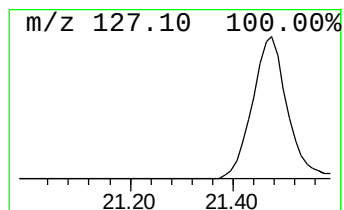
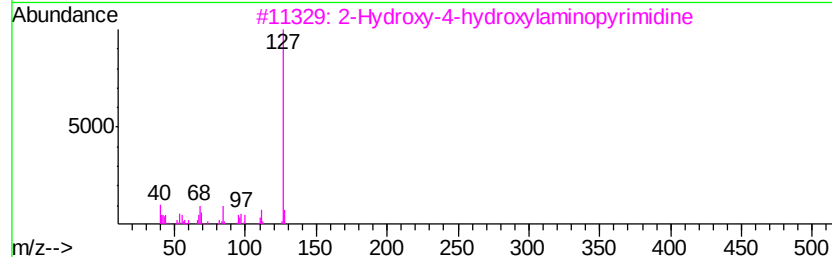
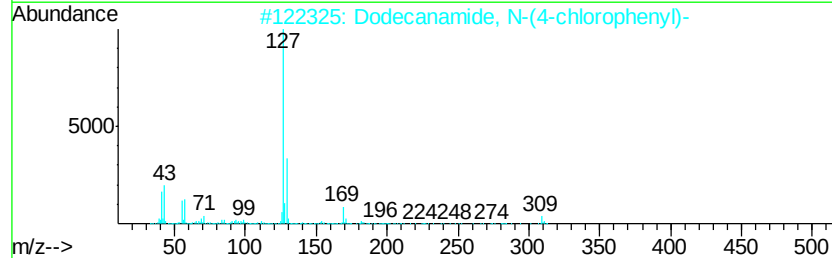
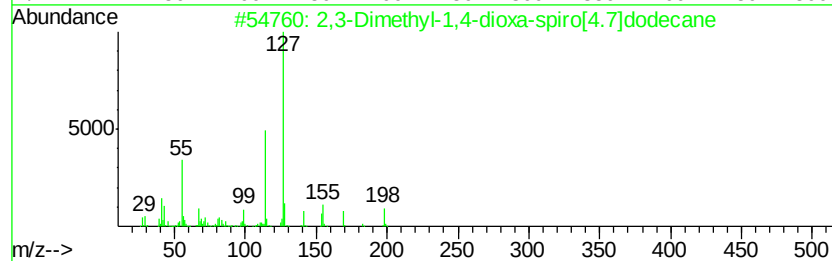
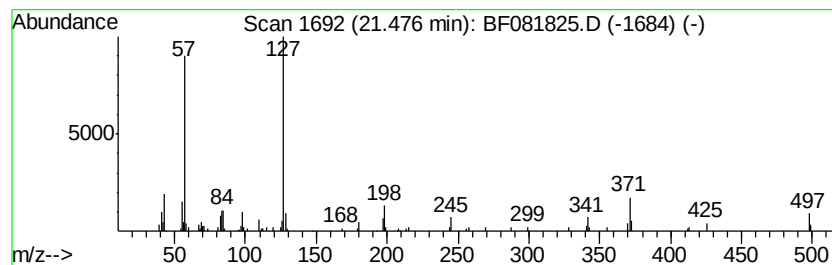
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 20 2,3-Dimethyl-1,4-dioxo-spiro... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.48	12.19 ng	431960	Perylene-d12	15.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,3-Dimethyl-1,4-dioxo-spiro[4.7]dodecane	198	C12H22O2	062406-91-1	50
2		Dodecanamide, N-(4-chlorophenyl)-	309	C18H28ClNO	099002-89-8	47
3		2-Hydroxy-4-hydroxylaminopyrimidine	127	C4H5N3O2	020555-88-8	47
4		1H-Imidazole, 4-methyl-5-nitro-	127	C4H5N3O2	014003-66-8	47
5		Propylphosphonic acid, fluoroanh...	224	C10H22F02P	333416-19-6	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092615\
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 Acq On : 26 Sep 2015 13:53
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Instrument :
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 ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.14	43.9	ng	1678090	1	6.94	764850	20.0
Benzene, 1-ethyl-...	6.46	8.2	ng	312382	1	6.94	764850	20.0
Benzene, 1-etheny...	6.77	17.4	ng	665699	1	6.94	764850	20.0
Benzene, 1,2,3-tr...	6.99	10.6	ng	404899	1	6.94	764850	20.0
Benzene, 1-propynyl-	7.20	26.6	ng	1015930	1	6.94	764850	20.0
Naphthalene, 1-me...	9.04	5.4	ng	1120180	2	8.23	4181730	20.0
cis-.beta.-Methyl...	9.46	10.7	ng	588857	3	9.98	1098000	20.0
unknown9.54	9.54	6.8	ng	374865	3	9.98	1098000	20.0
unknown10.14	10.14	6.1	ng	336443	3	9.98	1098000	20.0
Bicyclo[4.2.0]oct...	10.18	7.0	ng	384375	3	9.98	1098000	20.0
Anthracene, 2-met...	11.99	12.2	ng	590341	4	11.47	966168	20.0
4H-Cyclopenta[def...	12.08	12.5	ng	602830	4	11.47	966168	20.0
Phenanthrene, 9-e...	12.41	5.1	ng	246808	4	11.47	966168	20.0
Phenanthrene, 1-m...	13.19	5.0	ng	299333	5	14.10	1187480	20.0
2,3-Dimethyl-1,4-...	21.48	12.2	ng	431960	6	15.58	708581	20.0