

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF010917\
 Data File : BF092161.D
 Acq On : 10 Jan 2017 1:57
 Operator : UM/SJ
 Sample : I1045-06 2X
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RAP-SP4

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-BF122116.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	4.744	248	250	258	rBV	280067	345179	25.29%	1.729%
2	5.247	291	294	302	rBV	958439	1128646	82.68%	5.654%
3	6.150	369	373	378	rVB5	9535	27173	1.99%	0.136%
4	6.287	382	385	386	rBV	120575	162023	11.87%	0.812%
5	6.310	386	387	394	rVB	786060	1163265	85.22%	5.828%
6	6.413	394	396	399	rBV	1023593	1133963	83.07%	5.681%
7	6.642	414	416	418	rBV	876371	908816	66.58%	4.553%
8	6.790	427	429	432	rVB	866288	853675	62.54%	4.277%
9	7.030	449	450	454	rBV3	16765	31238	2.29%	0.156%
10	7.202	463	465	468	rVV	497034	488591	35.79%	2.448%
11	7.259	468	470	471	rVV	24586	29097	2.13%	0.146%
12	7.293	471	473	475	rVB2	19056	28229	2.07%	0.141%
13	7.453	485	487	489	rVB2	30691	39034	2.86%	0.196%
14	7.568	493	497	498	rVV2	20702	30637	2.24%	0.153%
15	7.670	503	506	507	rBV2	20745	24963	1.83%	0.125%
16	7.762	511	514	516	rVB3	13130	28078	2.06%	0.141%
17	7.819	516	519	522	rBV4	9743	28516	2.09%	0.143%
18	7.888	522	525	526	rBV3	15348	26872	1.97%	0.135%
19	7.933	526	529	531	rBV	957340	1216958	89.15%	6.097%
20	8.139	542	547	549	rBV4	18195	47581	3.49%	0.238%
21	8.219	551	554	556	rBV3	19269	40433	2.96%	0.203%
22	8.368	564	567	571	rBV	55284	80345	5.89%	0.403%
23	8.436	571	573	575	rBV2	25826	31755	2.33%	0.159%
24	8.539	579	582	584	rVV	46497	69413	5.09%	0.348%
25	8.630	588	590	592	rVB2	23447	37929	2.78%	0.190%
26	8.699	592	596	598	rBV4	9245	25409	1.86%	0.127%
27	8.756	598	601	603	rVV2	24000	44035	3.23%	0.221%
28	8.893	610	613	616	rBV5	13587	31911	2.34%	0.160%
29	8.962	616	619	621	rBV3	77706	90390	6.62%	0.453%
30	9.008	621	623	625	rVB	1465856	1187052	86.96%	5.947%
31	9.099	627	631	633	rBV	90595	121727	8.92%	0.610%
32	9.156	633	636	639	rVB3	26853	51236	3.75%	0.257%
33	9.271	643	646	648	rBV2	38849	75006	5.49%	0.376%
34	9.339	651	652	654	rBV2	62073	85078	6.23%	0.426%

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35	9.408	656	658	660	rVB2	143255	176465	12.93%	0.884%
36	9.545	667	670	673	rBV4	26007	53156	3.89%	0.266%
37	9.591	673	674	676	rVB	38094	30463	2.23%	0.153%
38	9.625	676	677	679	rBV	105638	87356	6.40%	0.438%
39	9.682	679	682	683	rVV	1449293	1313357	96.21%	6.580%
40	9.716	683	685	687	rVV	75465	111843	8.19%	0.560%
41	9.785	690	691	694	rVB2	32689	44384	3.25%	0.222%
42	9.888	694	700	701	rBV3	88840	154565	11.32%	0.774%
43	9.922	701	703	704	rVV2	41799	67263	4.93%	0.337%
44	10.002	707	710	713	rBV3	35775	64321	4.71%	0.322%
45	10.082	715	717	719	rBV3	26796	49730	3.64%	0.249%
46	10.128	719	721	723	rVB	84324	86769	6.36%	0.435%
47	10.174	723	725	728	rBV4	15861	27888	2.04%	0.140%
48	10.219	728	729	733	rVB3	57675	99822	7.31%	0.500%
49	10.288	733	735	736	rBV	30199	28323	2.07%	0.142%
50	10.322	736	738	742	rBV2	124430	213831	15.66%	1.071%
51	10.425	746	747	748	rBV	28913	29224	2.14%	0.146%
52	10.471	748	751	753	rBV	562008	532426	39.00%	2.667%
53	10.562	757	759	760	rBV2	16825	32758	2.40%	0.164%
54	10.596	760	762	766	rVB	174266	256545	18.79%	1.285%
55	10.665	766	768	769	rBV	28931	34396	2.52%	0.172%
56	10.802	778	780	781	rVB2	38177	37139	2.72%	0.186%
57	10.974	793	795	797	rBV	21022	27606	2.02%	0.138%
58	11.042	800	801	802	rVV	143224	132909	9.74%	0.666%
59	11.076	802	804	806	rVB	100695	114939	8.42%	0.576%
60	11.156	809	811	815	rVV	997706	1365039	100.00%	6.838%
61	11.236	815	818	819	rVB	103543	109501	8.02%	0.549%
62	11.271	819	821	823	rBV3	22572	35458	2.60%	0.178%
63	11.396	829	832	833	rBV2	93007	120428	8.82%	0.603%
64	11.465	837	838	846	rVB3	103014	274893	20.14%	1.377%
65	11.625	846	852	853	rBV3	109613	280301	20.53%	1.404%
66	11.659	853	855	857	rBV2	53571	92367	6.77%	0.463%
67	11.774	863	865	870	rVB2	122581	212273	15.55%	1.063%
68	11.877	872	874	875	rBV	104917	82455	6.04%	0.413%
69	12.208	901	903	907	rBV4	101718	276631	20.27%	1.386%
70	12.265	907	908	910	rVV	144058	112229	8.22%	0.562%
71	12.299	910	911	913	rVV2	57143	76400	5.60%	0.383%

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72	12.368	915	917	919	rVV	506578	467101	34.22%	2.340%
73	12.597	934	937	939	rBV	358891	427925	31.35%	2.144%
74	12.745	948	950	954	rVB	883783	815863	59.77%	4.087%
75	12.997	970	972	973	rVB	154753	187900	13.77%	0.941%
76	13.031	973	975	978	rBV3	114266	253909	18.60%	1.272%
77	13.797	1040	1042	1046	rBV2	691037	941569	68.98%	4.717%
78	13.980	1056	1058	1060	rBV	260884	307232	22.51%	1.539%

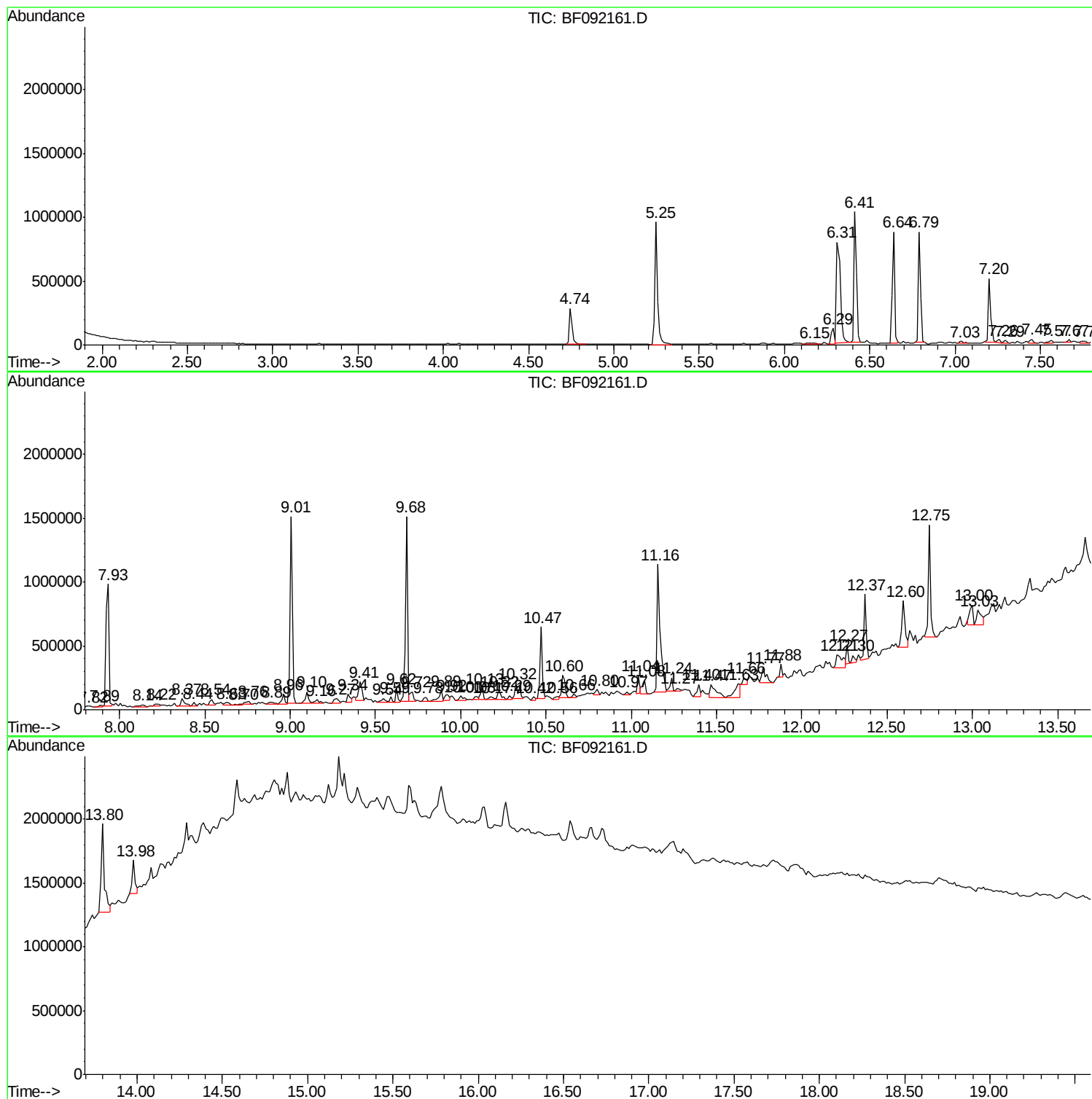
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Instrument :
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ClientSampleId :
RAP-SP4

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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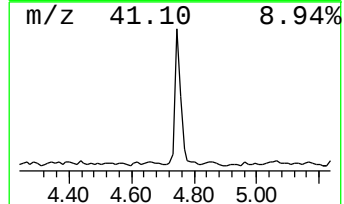
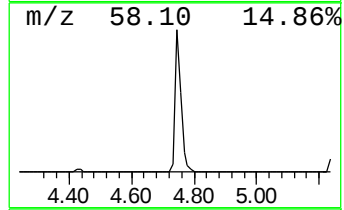
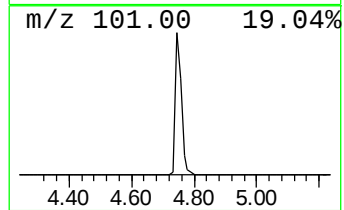
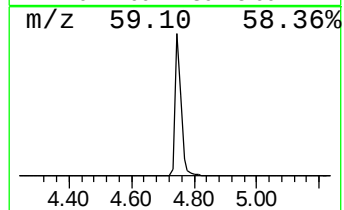
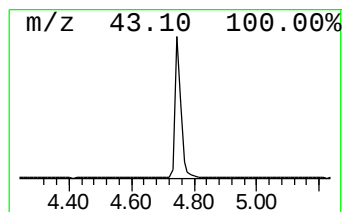
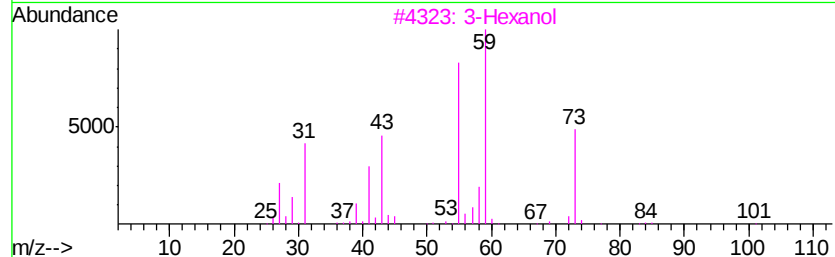
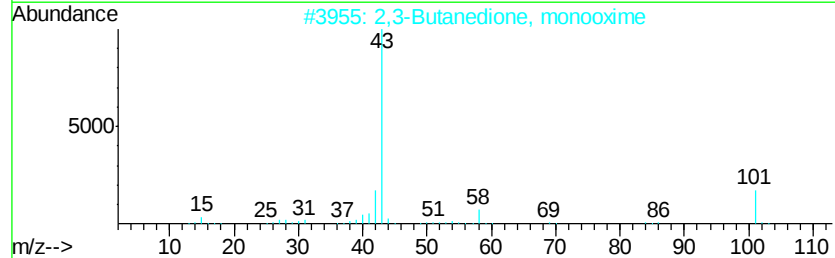
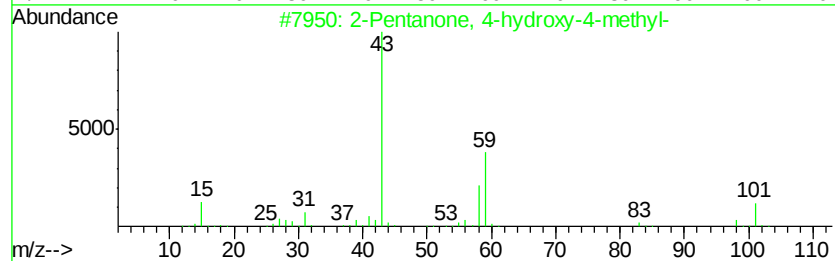
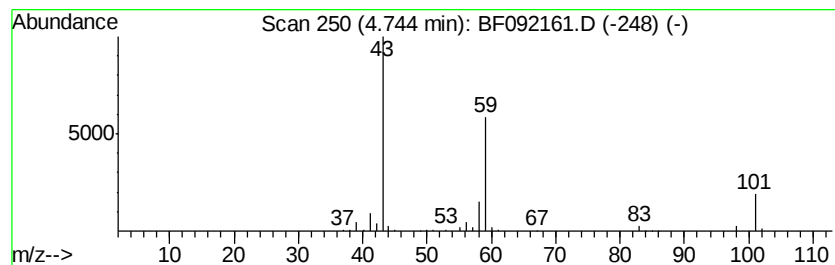
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.74	7.60 ng	345179	1,4-Dichlorobenzene-d4	6.64

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	35
3			3-Hexanol	102	C6H14O	000623-37-0	28
4			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17
5			(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9



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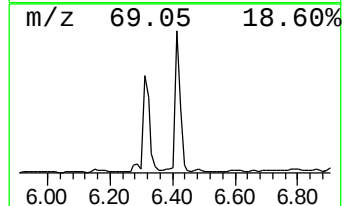
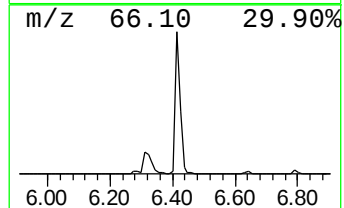
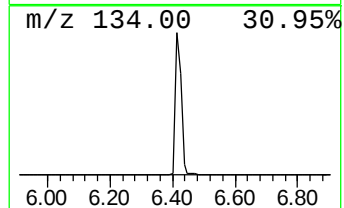
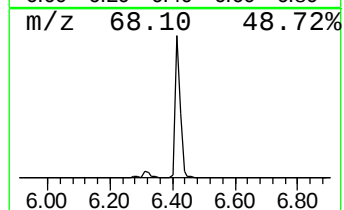
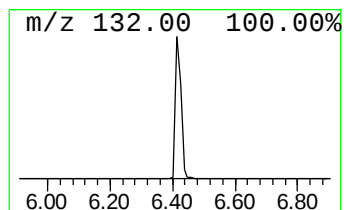
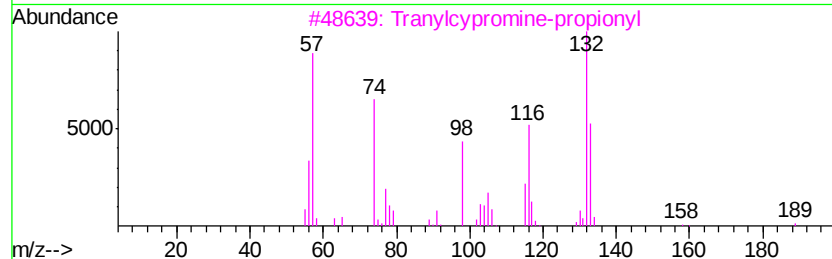
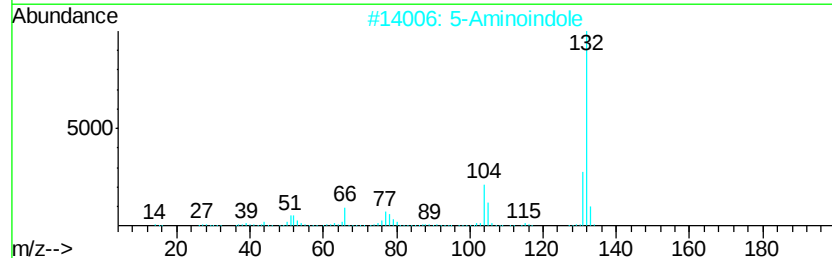
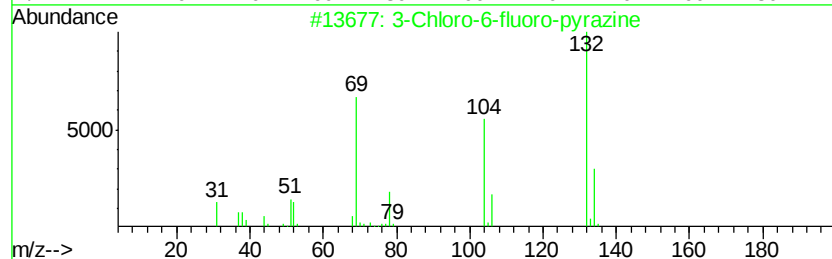
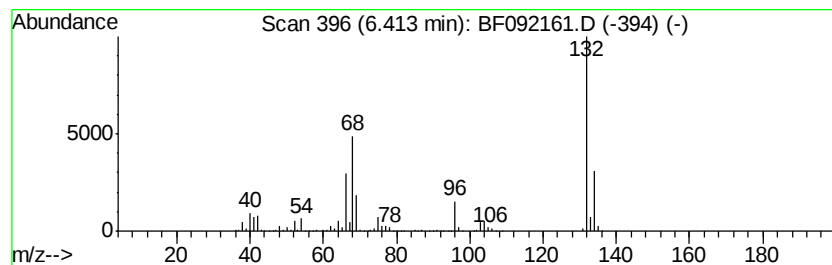
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.41 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.41	24.95 ng	1133960	1,4-Dichlorobenzene-d4	6.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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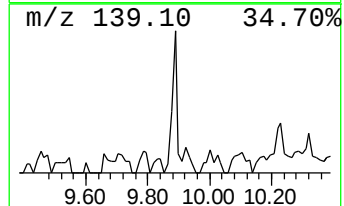
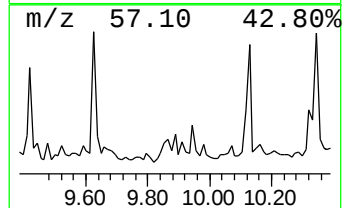
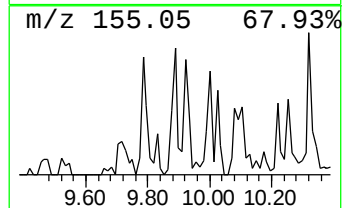
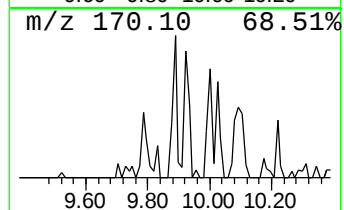
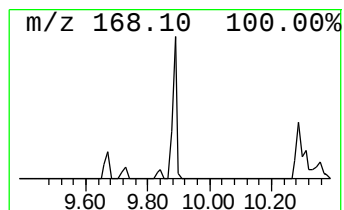
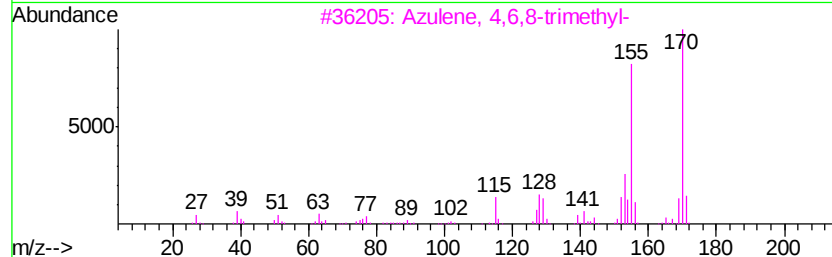
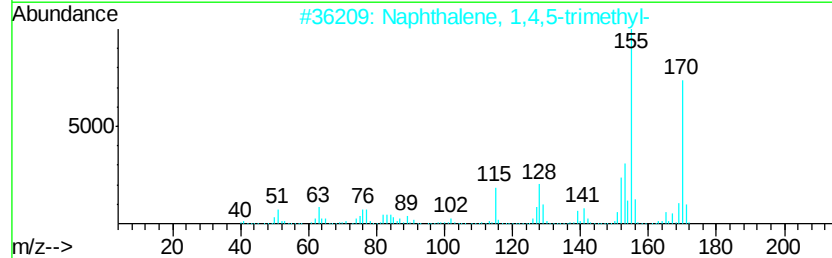
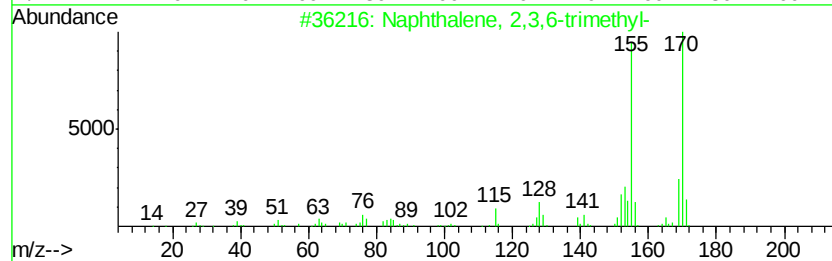
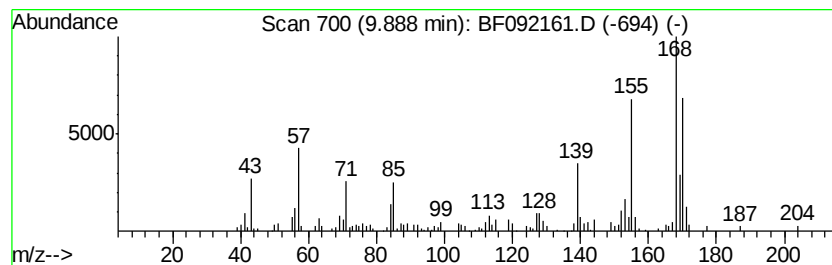
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Naphthalene, 2,3,6-trimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.89	2.35 ng	154565	Acenaphthene-d10	9.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	90
2		Naphthalene, 1,4,5-trimethyl-	170	C13H14	002131-41-1	86
3		Azulene, 4,6,8-trimethyl-	170	C13H14	000941-81-1	60
4		Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	50
5		1-(2,2-Dimethylcyclopropyl)-2-ph...	170	C13H14	1000294-91-1	46



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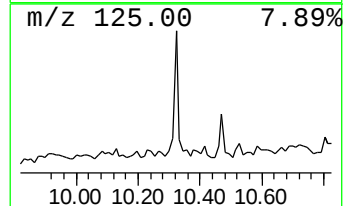
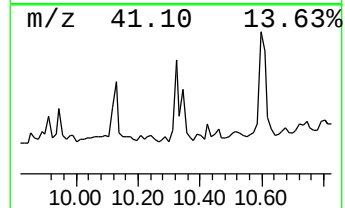
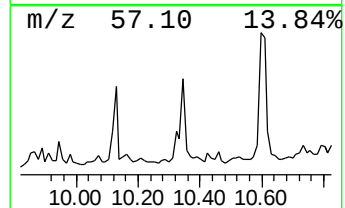
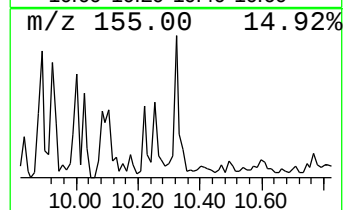
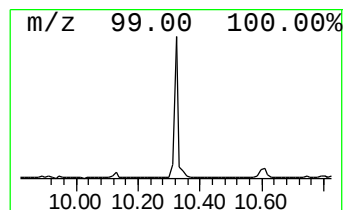
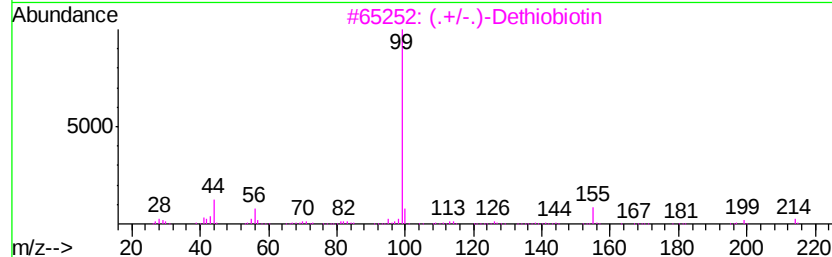
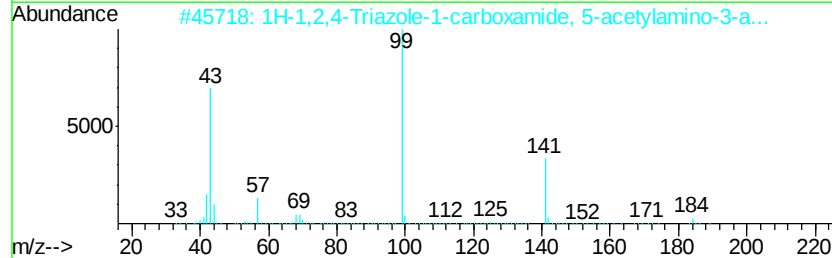
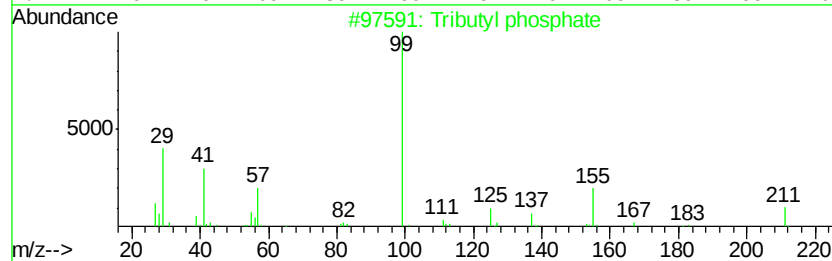
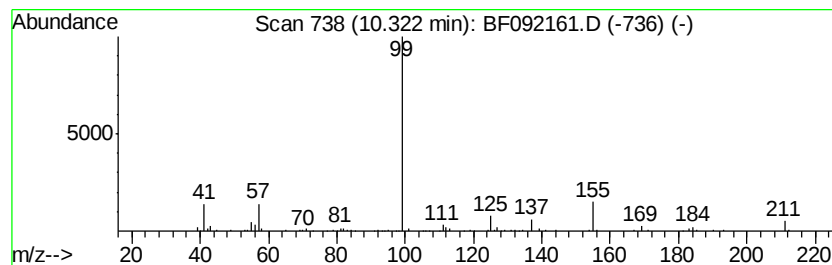
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 unknown10.32 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.32	3.26 ng	213831	Acenaphthene-d10	9.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tributyl phosphate	266	C12H27O4P	000126-73-8	47
2		1H-1,2,4-Triazole-1-carboxamide,...	184	C5H8N6O2	1000277-52-4	42
3		(.+/-)-Dethiobiotin	214	C10H18N2O3	000636-20-4	9
4		Spiro[1,3-dioxolane-2,2'(1'H)-na...	196	C12H20O2	006857-86-9	9
5		n-Heptyl methylphosphonofluoridate	196	C8H18FO2P	162085-82-7	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF010917\
Data File : BF092161.D
Acq On : 10 Jan 2017 1:57
Operator : UM/SJ
Sample : I1045-06 2X
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
RAP-SP4

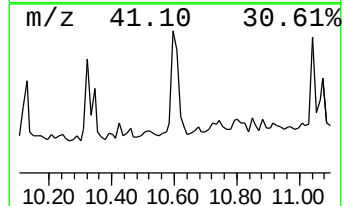
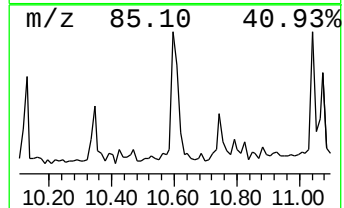
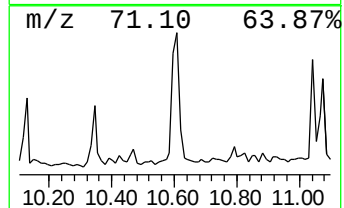
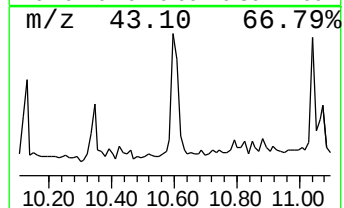
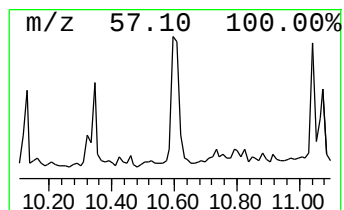
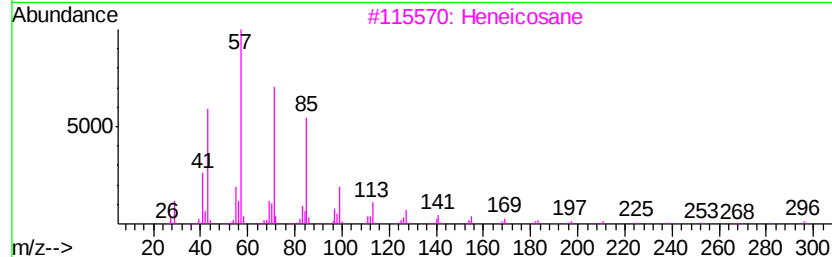
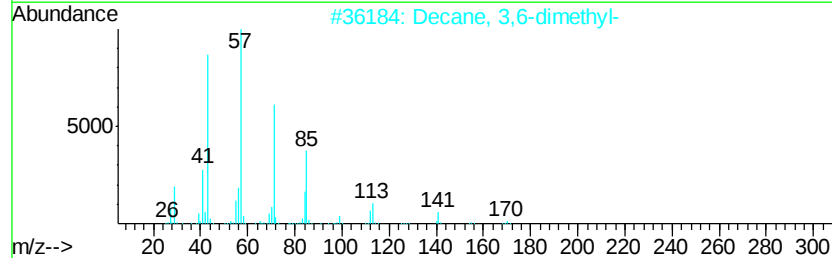
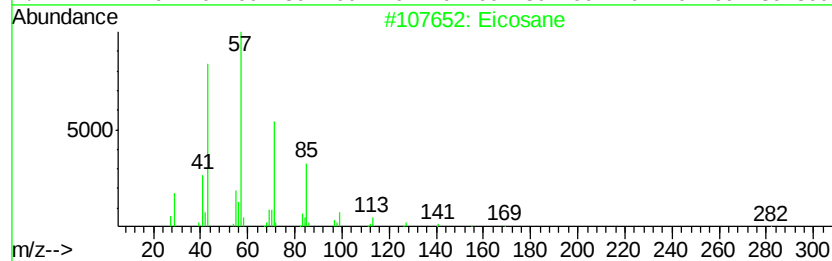
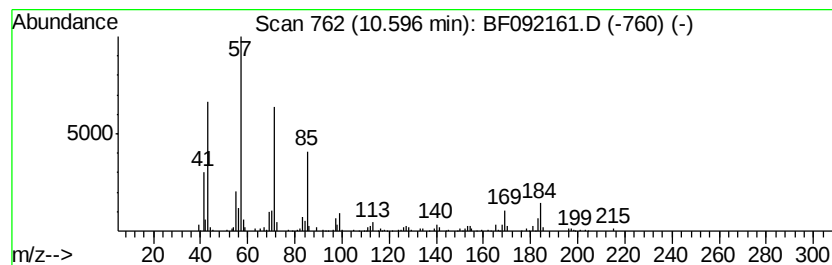
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Eicosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.60	3.76 ng	256545	Phenanthrene-d10	11.16

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	68
2	Decane, 3,6-dimethyl-		170	C12H26	017312-53-7	64
3	Heneicosane		296	C21H44	000629-94-7	64
4	2-Bromo dodecane		248	C12H25Br	013187-99-0	64
5	1-Iodo-2-methylundecane		296	C12H25I	073105-67-6	59



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF010917\
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Sample : I1045-06 2X
Misc :
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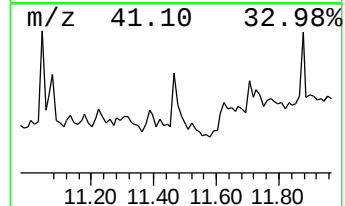
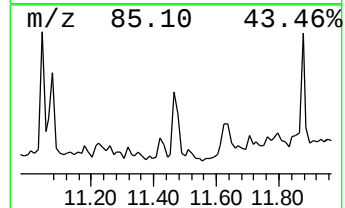
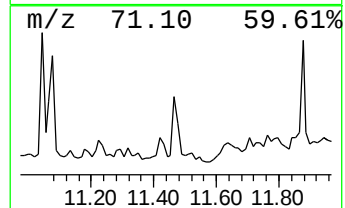
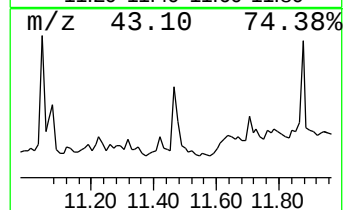
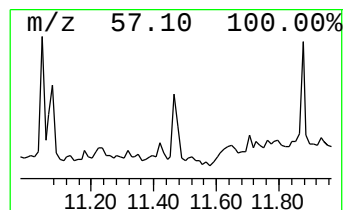
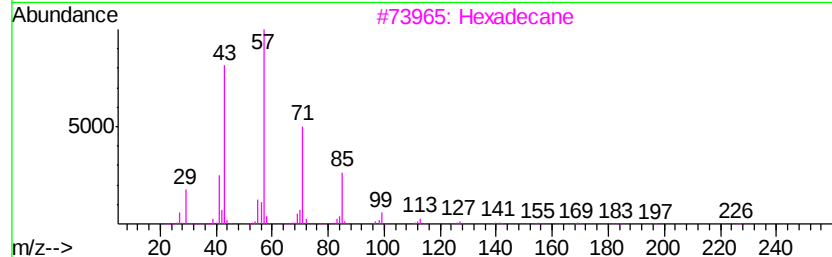
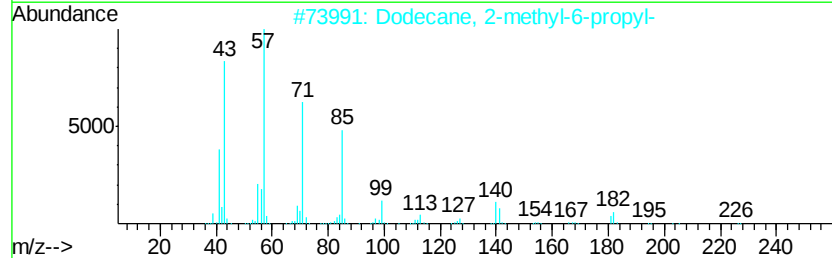
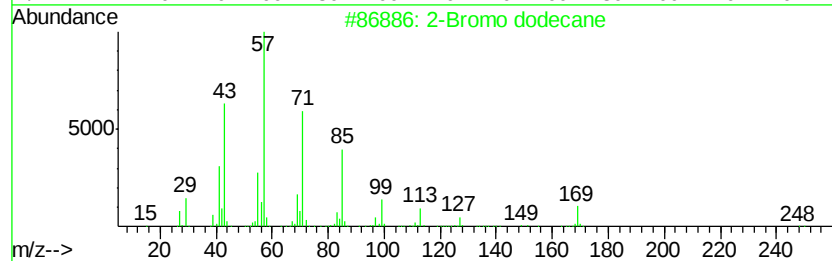
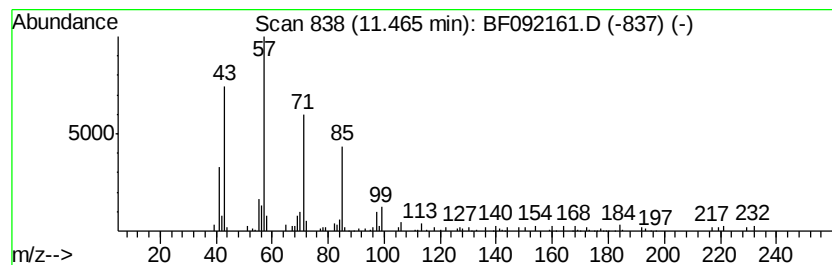
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 2-Bromo dodecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.47	4.03 ng	274893	Phenanthrene-d10		11.16	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Bromo dodecane		248	C12H25Br	013187-99-0	72
2	Dodecane, 2-methyl-6-propyl-		226	C16H34	055045-08-4	72
3	Hexadecane		226	C16H34	000544-76-3	52
4	Heptane, 4-ethyl-		128	C9H20	002216-32-2	46
5	Tridecane		184	C13H28	000629-50-5	45



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF010917\
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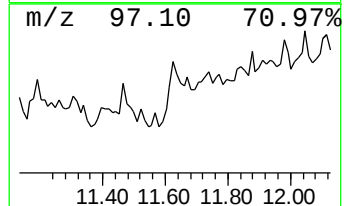
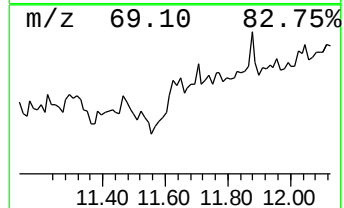
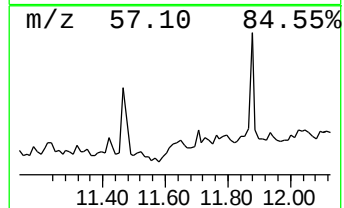
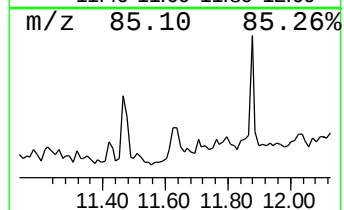
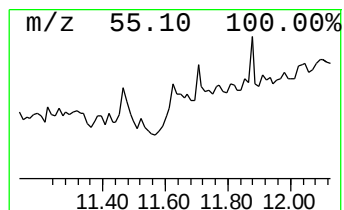
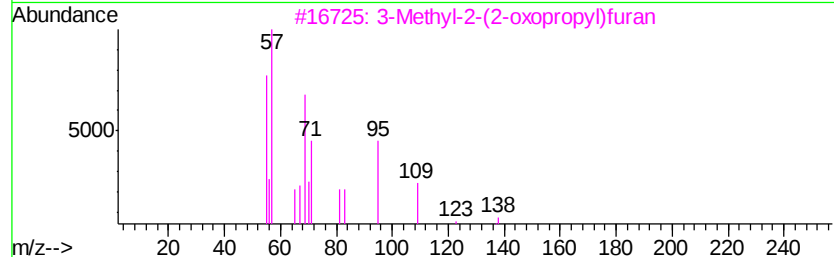
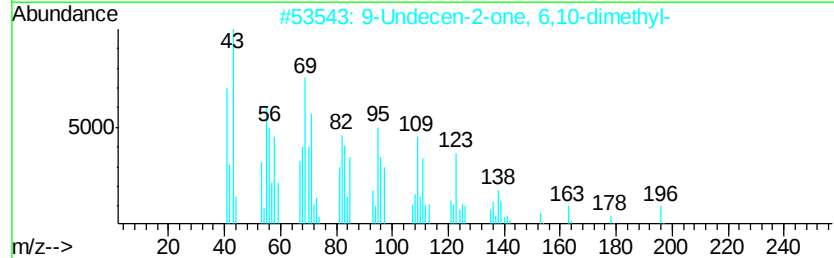
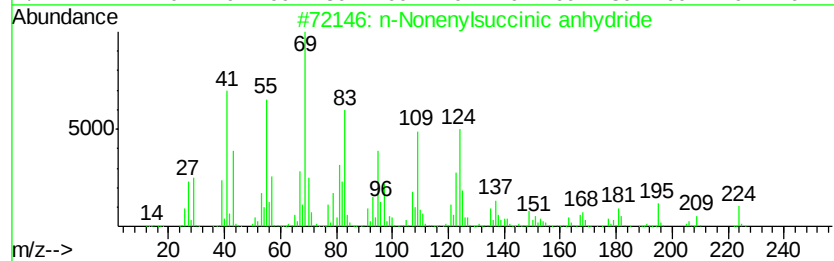
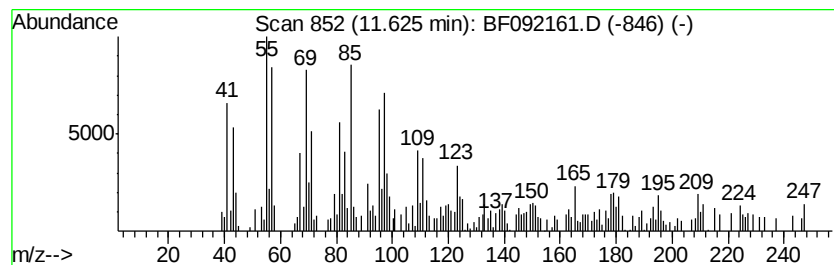
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 n-Nonenylsuccinic anhydride Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.63	4.11 ng	280301	Phenanthrene-d10		11.16
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Nonenylsuccinic anhydride	224	C13H20O3	028928-97-4	53
2	9-Undecen-2-one, 6,10-dimethyl-	196	C13H24O	004433-36-7	49
3	3-Methyl-2-(2-oxopropyl)furan	138	C8H10O2	087773-62-4	46
4	1-Tetradecene	196	C14H28	001120-36-1	46
5	Cyclopentadecane	210	C15H30	000295-48-7	42



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF010917\
Data File : BF092161.D
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Misc :
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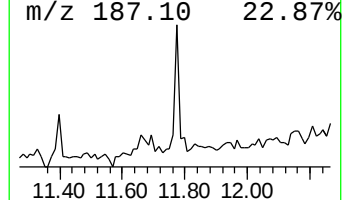
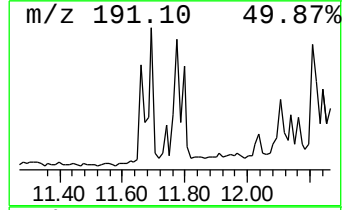
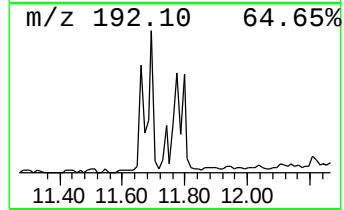
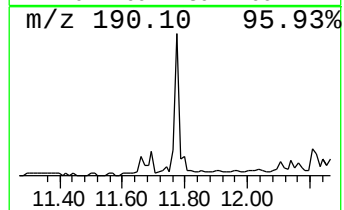
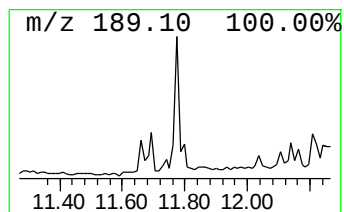
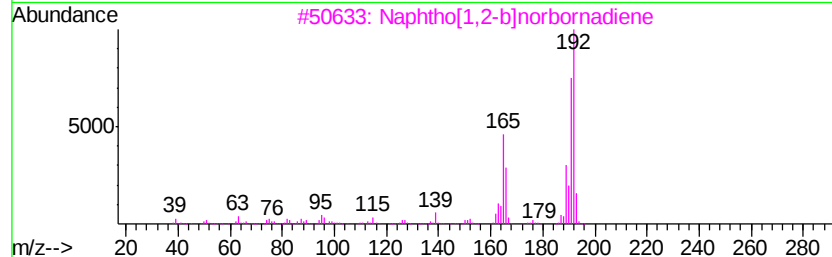
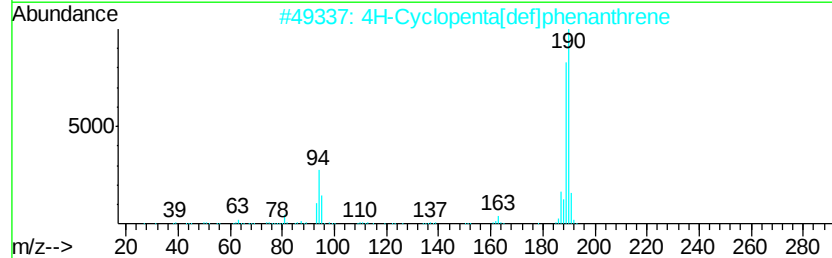
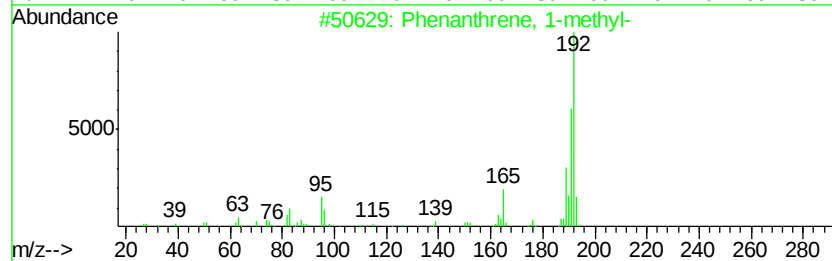
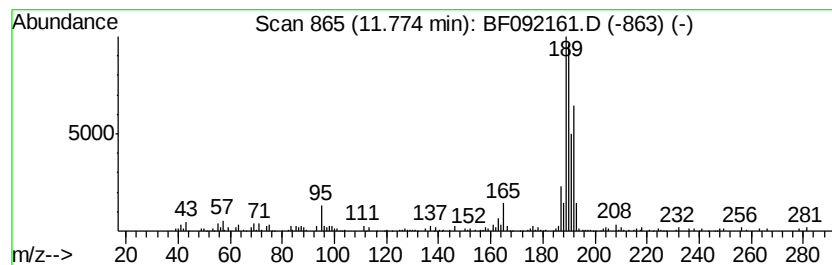
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Phenanthrene, 1-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.77	3.11 ng	212273	Phenanthrene-d10	11.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	74
2		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	64
3		Naphtho[1,2-b]norbornadiene	192	C15H12	1000210-14-8	38
4		Anthracene, 1-methyl-	192	C15H12	000610-48-0	38
5		1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	30



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF010917\
Data File : BF092161.D
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Misc :
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Instrument :
BNA_F
ClientSampled :
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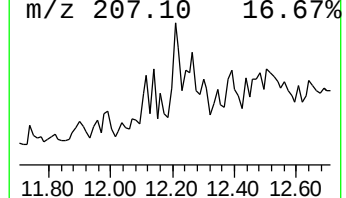
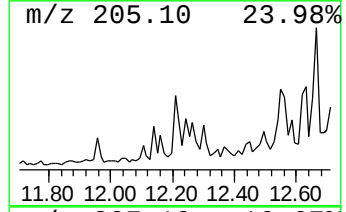
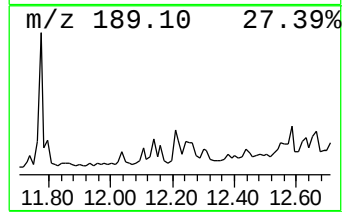
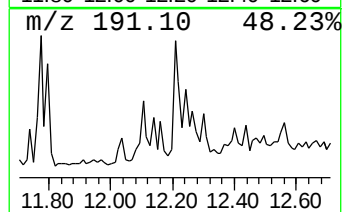
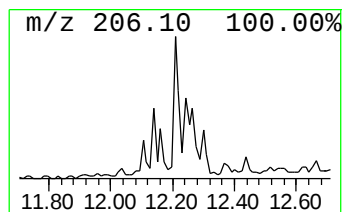
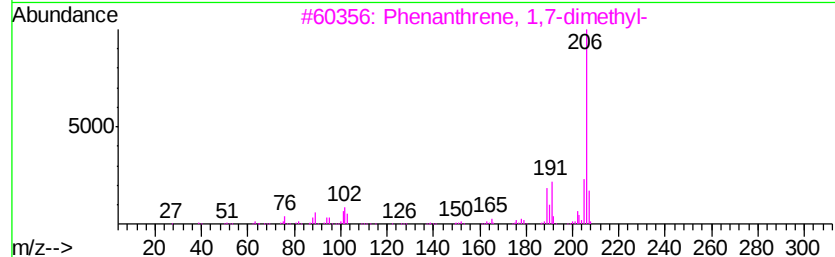
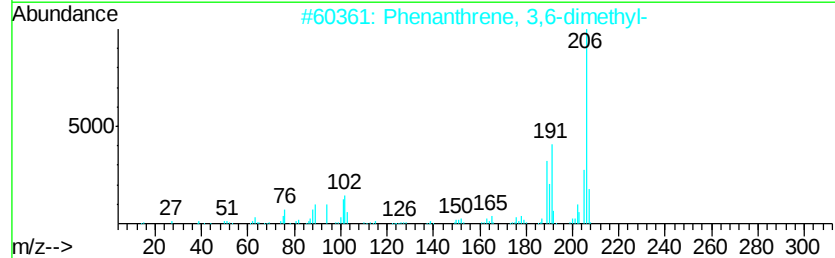
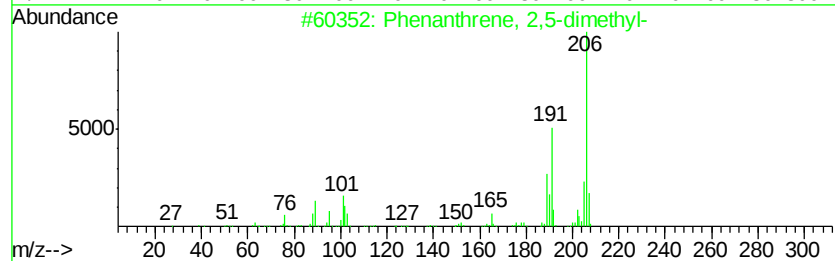
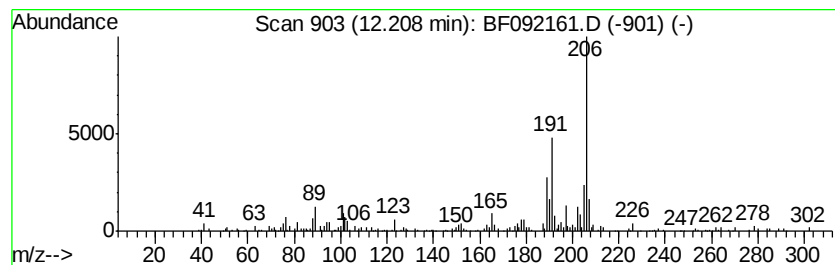
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Phenanthrene, 2,5-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.21	4.05 ng	276631	Phenanthrene-d10	11.16

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	98
2			Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	93
3			Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	91
4			9,10-Dimethylantracene	206	C16H14	000781-43-1	90
5			3-Methyl-1-phenyl-1H-indene	206	C16H14	022360-63-0	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF010917\
Data File : BF092161.D
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Misc :
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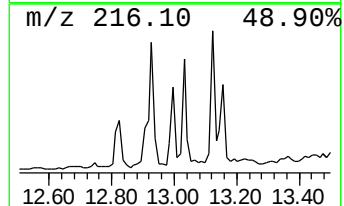
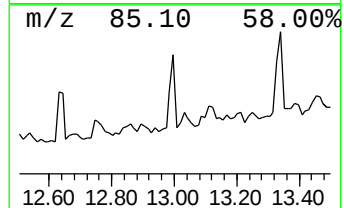
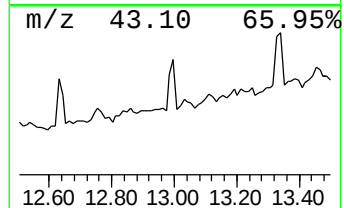
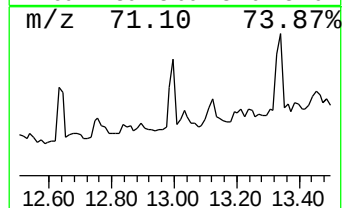
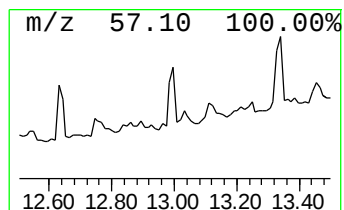
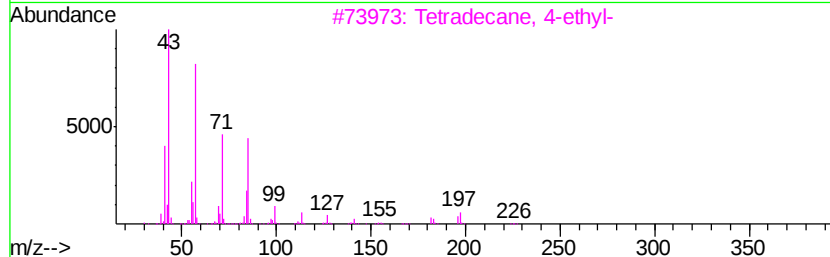
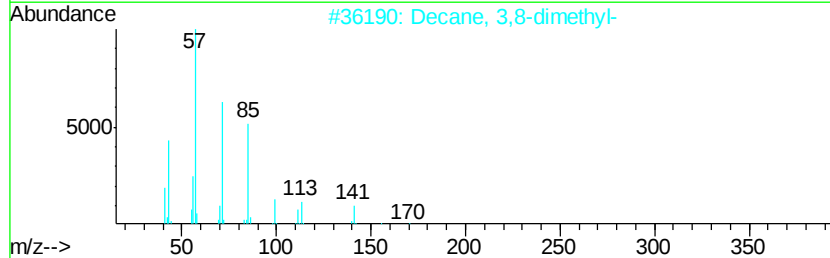
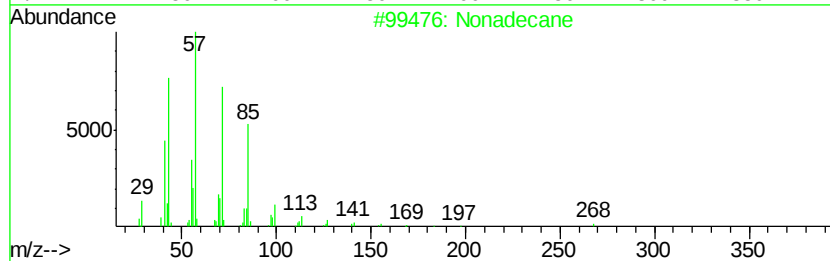
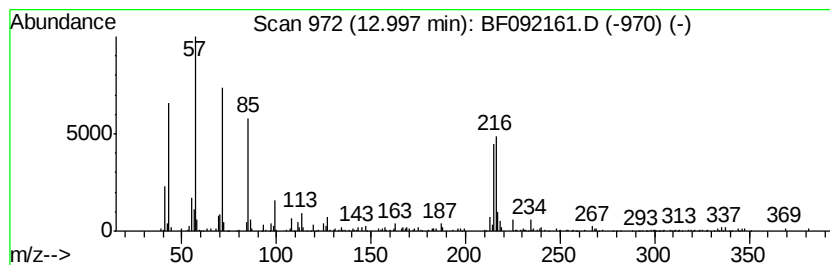
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown13.00 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.00	3.99 ng	187900	Chrysene-d12	13.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	49
2		Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	49
3		Tetradecane, 4-ethyl-	226	C16H34	055045-14-2	47
4		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	47
5		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	47



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF010917\
Data File : BF092161.D
Acq On : 10 Jan 2017 1:57
Operator : UM/SJ
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Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
RAP-SP4

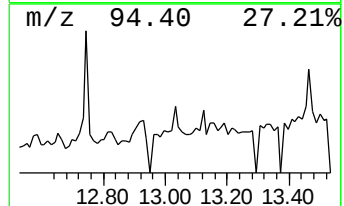
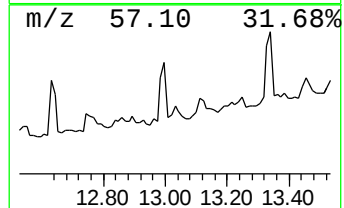
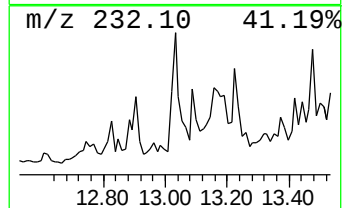
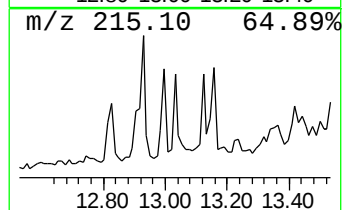
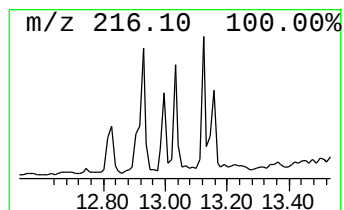
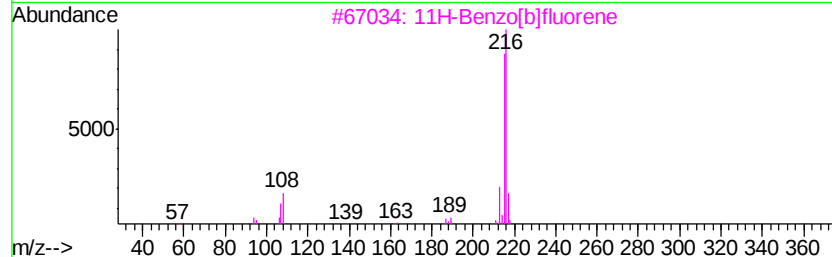
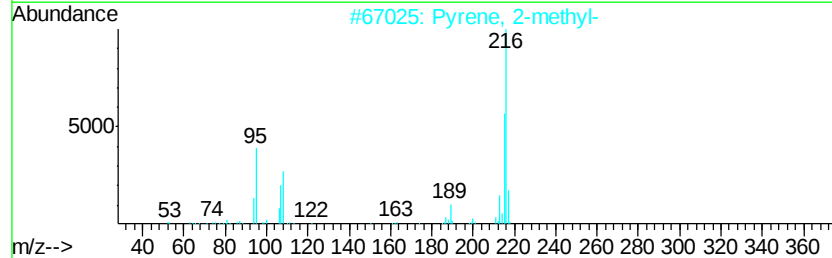
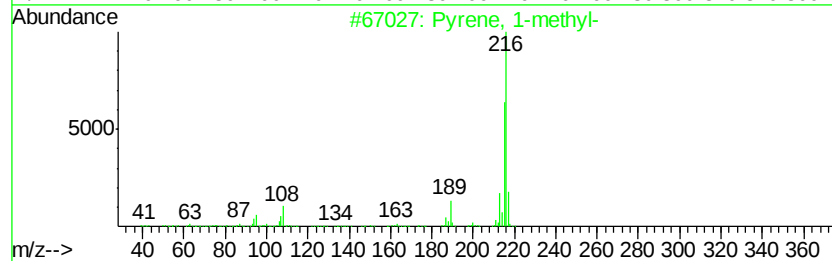
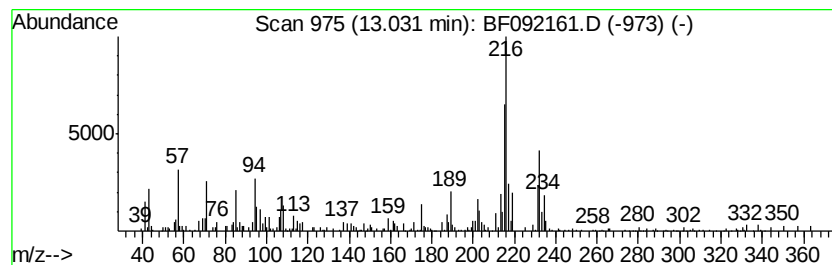
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Pyrene, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.03	5.39 ng	253909	Chrysene-d12	13.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pyrene, 1-methyl-	216	C17H12	002381-21-7	87
2		Pyrene, 2-methyl-	216	C17H12	003442-78-2	60
3		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	50
4		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	50
5		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	49



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF010917\
Data File : BF092161.D
Acq On : 10 Jan 2017 1:57
Operator : UM/SJ
Sample : I1045-06 2X
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
RAP-SP4

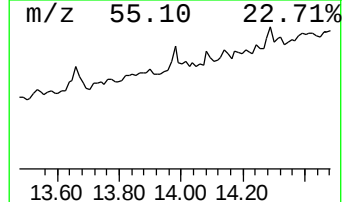
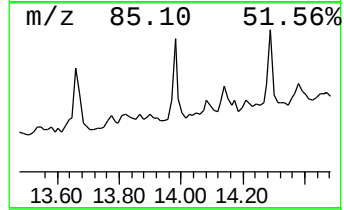
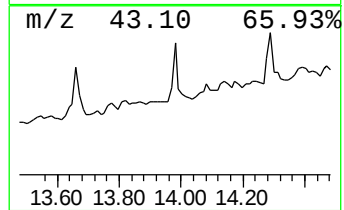
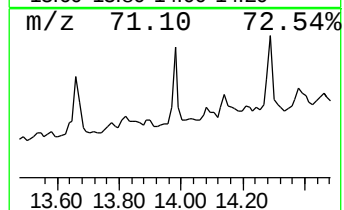
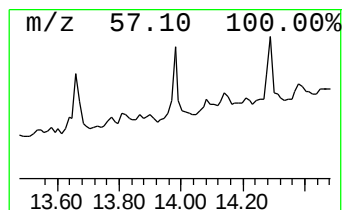
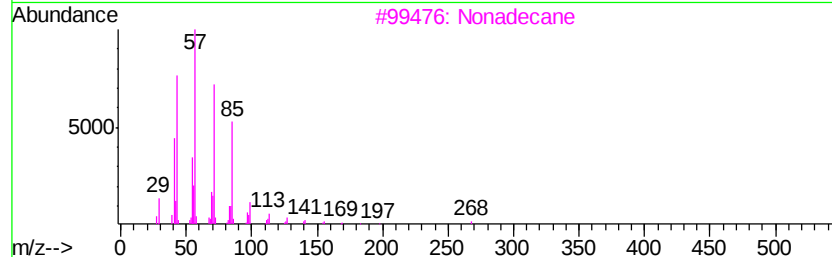
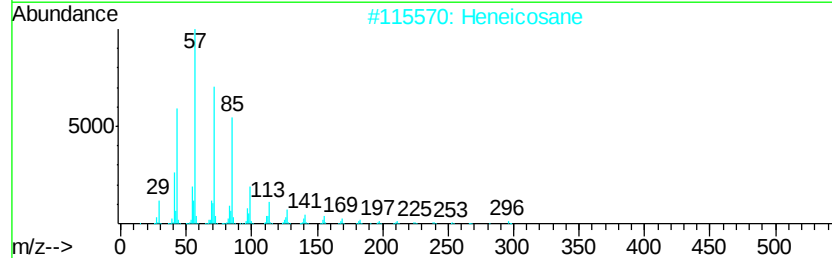
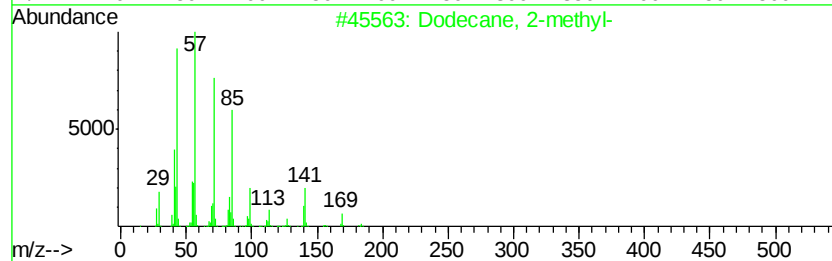
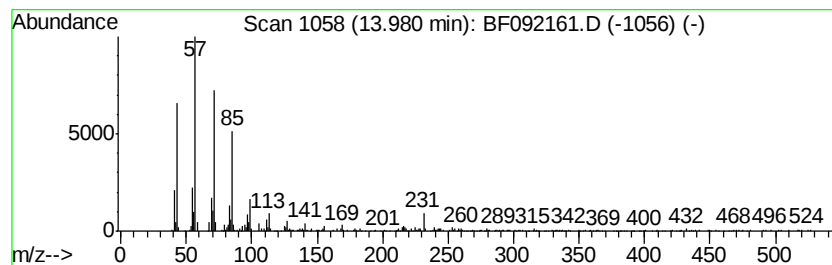
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Dodecane, 2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.98	6.53 ng	307232	Chrysene-d12	13.80

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane, 2-methyl-	184	C13H28	001560-97-0	81
2			Heneicosane	296	C21H44	000629-94-7	81
3			Nonadecane	268	C19H40	000629-92-5	81
4			Hentriacontane	437	C31H64	000630-04-6	78
5			Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	74



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF010917\
Data File : BF092161.D
Acq On : 10 Jan 2017 1:57
Operator : UM/SJ
Sample : I1045-06 2X
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
RAP-SP4

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.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...	4.74	7.6 ng		345179	1	6.64	908816	20.0
unknown6.41	6.41	24.9 ng		1133960	1	6.64	908816	20.0
Naphthalene, 2,3,...	9.89	2.4 ng		154565	3	9.68	1313360	20.0
unknown10.32	10.32	3.3 ng		213831	3	9.68	1313360	20.0
Eicosane	10.60	3.8 ng		256545	4	11.16	1365040	20.0
2-Bromo dodecane	11.47	4.0 ng		274893	4	11.16	1365040	20.0
n-Nonenylsuccinic...	11.63	4.1 ng		280301	4	11.16	1365040	20.0
Phenanthrene, 1-m...	11.77	3.1 ng		212273	4	11.16	1365040	20.0
Phenanthrene, 2,5...	12.21	4.0 ng		276631	4	11.16	1365040	20.0
unknown13.00	13.00	4.0 ng		187900	5	13.80	941569	20.0
Pyrene, 1-methyl-	13.03	5.4 ng		253909	5	13.80	941569	20.0
Dodecane, 2-methyl-	13.98	6.5 ng		307232	5	13.80	941569	20.0