

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011217\
 Data File : BF092230.D
 Acq On : 13 Jan 2017 2:52
 Operator : UM/SJ
 Sample : I1131-02
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SW-001

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	1.967	5	7	12	rVB2	20391	50893	1.04%	0.091%
2	3.990	180	184	189	rBV	32593	74425	1.53%	0.134%
3	4.721	245	248	265	rVB	1576026	2430883	49.88%	4.363%
4	5.190	287	289	301	rBV	3675504	3914817	80.33%	7.026%
5	6.264	380	383	385	rBV	4281086	4873544	100.00%	8.746%
6	6.367	390	392	395	rVB	5203326	4848902	99.49%	8.702%
7	6.596	409	412	414	rBV	936480	969592	19.90%	1.740%
8	6.745	423	425	428	rBV	2212014	2657680	54.53%	4.770%
9	7.167	459	462	464	rBV	2836333	2738695	56.20%	4.915%
10	7.876	522	524	527	rBV	938769	1133450	23.26%	2.034%
11	8.962	616	619	621	rBV	4177436	3785810	77.68%	6.794%
12	9.362	652	654	656	rBV	977328	819696	16.82%	1.471%
13	9.625	675	677	679	rBV	1682314	1749154	35.89%	3.139%
14	9.659	679	680	683	rVB	73565	60757	1.25%	0.109%
15	9.831	690	695	697	rBV4	37324	72178	1.48%	0.130%
16	10.048	711	714	715	rBV	87079	90412	1.86%	0.162%
17	10.082	715	717	718	rVB	102844	119034	2.44%	0.214%
18	10.173	723	725	729	rVB	75019	90140	1.85%	0.162%
19	10.299	733	736	738	rBV	44042	59286	1.22%	0.106%
20	10.413	744	746	749	rBV2	2221220	3002910	61.62%	5.389%
21	10.551	756	758	762	rBV	176358	222366	4.56%	0.399%
22	10.745	771	775	777	rBV4	25094	52916	1.09%	0.095%
23	10.996	795	797	799	rBV3	207425	228724	4.69%	0.410%
24	11.099	804	806	811	rVV2	1243236	2594419	53.23%	4.656%
25	11.179	811	813	815	rVB	295238	280998	5.77%	0.504%
26	11.351	825	828	832	rBV4	99260	264660	5.43%	0.475%
27	11.419	832	834	836	rVB	184847	161951	3.32%	0.291%
28	11.602	845	850	852	rBV	112130	200732	4.12%	0.360%
29	11.636	852	853	854	rVV	154466	147238	3.02%	0.264%
30	11.659	854	855	858	rVV2	392133	442084	9.07%	0.793%
31	11.716	858	860	865	rVB2	192722	304106	6.24%	0.546%
32	11.819	868	869	873	rVV	81512	120866	2.48%	0.217%
33	11.899	874	876	878	rVV	83717	99456	2.04%	0.178%
34	11.934	878	879	881	rVB	66398	50217	1.03%	0.090%

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 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	12.151	896	898	900	rBV	61994	86710	1.78%	0.156%
36	12.208	900	903	905	rVV2	55415	128119	2.63%	0.230%
37	12.242	905	906	910	rVB	90707	86753	1.78%	0.156%
38	12.311	910	912	914	rBV	1521908	1341554	27.53%	2.408%
39	12.379	914	918	919	rVV3	39831	70145	1.44%	0.126%
40	12.402	919	920	922	rVV	64097	85610	1.76%	0.154%
41	12.459	922	925	926	rVV2	89947	177459	3.64%	0.318%
42	12.539	929	932	934	rVV	1188090	1389353	28.51%	2.493%
43	12.585	934	936	939	rVB2	68777	114253	2.34%	0.205%
44	12.688	943	945	948	rVB	3354393	3469402	71.19%	6.226%
45	12.757	948	951	953	rBV2	55088	96917	1.99%	0.174%
46	12.871	957	961	963	rVB	106182	196461	4.03%	0.353%
47	12.939	963	967	968	rBV	90251	135155	2.77%	0.243%
48	12.974	968	970	974	rVB	101659	133179	2.73%	0.239%
49	13.088	979	980	985	rVB2	33539	51534	1.06%	0.092%
50	13.179	985	988	992	rVB2	65551	92978	1.91%	0.167%
51	13.282	992	997	1001	rBV4	43907	127525	2.62%	0.229%
52	13.374	1004	1005	1009	rVB	72812	93822	1.93%	0.168%
53	13.488	1012	1015	1016	rBV	73632	119366	2.45%	0.214%
54	13.602	1022	1025	1032	rVB3	162362	400260	8.21%	0.718%
55	13.728	1033	1036	1041	rVV	1292934	2211254	45.37%	3.969%
56	13.831	1043	1045	1047	rBV2	61428	80921	1.66%	0.145%
57	13.888	1047	1050	1051	rBV2	29549	51505	1.06%	0.092%
58	13.922	1051	1053	1056	rVB3	41941	79252	1.63%	0.142%
59	14.117	1068	1070	1072	rVB	76524	92461	1.90%	0.166%
60	14.220	1075	1079	1080	rBV3	53997	113856	2.34%	0.204%
61	14.242	1080	1081	1086	rVB4	70215	126110	2.59%	0.226%
62	14.437	1095	1098	1102	rBV5	36630	104981	2.15%	0.188%
63	14.517	1102	1105	1108	rBV4	60653	112953	2.32%	0.203%
64	14.585	1109	1111	1115	rBV2	59932	84196	1.73%	0.151%
65	14.642	1115	1116	1118	rBV2	99022	102129	2.10%	0.183%
66	14.722	1121	1123	1128	rBV	446244	934090	19.17%	1.676%
67	14.814	1128	1131	1133	rBV3	123893	207762	4.26%	0.373%
68	14.985	1144	1146	1149	rVB	290339	344382	7.07%	0.618%
69	15.042	1149	1151	1153	rBV	403990	485213	9.96%	0.871%
70	15.088	1153	1155	1157	rBV	843154	1136493	23.32%	2.040%
71	15.305	1171	1174	1178	rVB5	34732	63159	1.30%	0.113%

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72	15.500	1186	1191	1193	rBV2	98262	181657	3.73%	0.326%
73	15.568	1194	1197	1200	rVV3	41951	90708	1.86%	0.163%
74	15.923	1225	1228	1229	rBV	36949	65376	1.34%	0.117%
75	16.060	1235	1240	1244	rBV7	37377	157102	3.22%	0.282%
76	16.163	1246	1249	1251	rBV2	50651	96914	1.99%	0.174%
77	16.334	1261	1264	1267	rBV	226488	399927	8.21%	0.718%
78	16.403	1267	1270	1274	rVB	77715	158917	3.26%	0.285%
79	16.483	1274	1277	1279	rBV	45068	93808	1.92%	0.168%
80	16.700	1293	1296	1303	rVV	154730	328945	6.75%	0.590%
81	16.837	1304	1308	1312	rVV6	34602	117457	2.41%	0.211%
82	16.917	1312	1315	1318	rVV	45872	120530	2.47%	0.216%
83	16.974	1318	1320	1328	rVB3	45663	131546	2.70%	0.236%
84	17.660	1375	1380	1385	rBV3	46396	166855	3.42%	0.299%
85	18.471	1447	1451	1456	rVB3	21891	73543	1.51%	0.132%
86	18.597	1459	1462	1467	rVB2	33258	96531	1.98%	0.173%

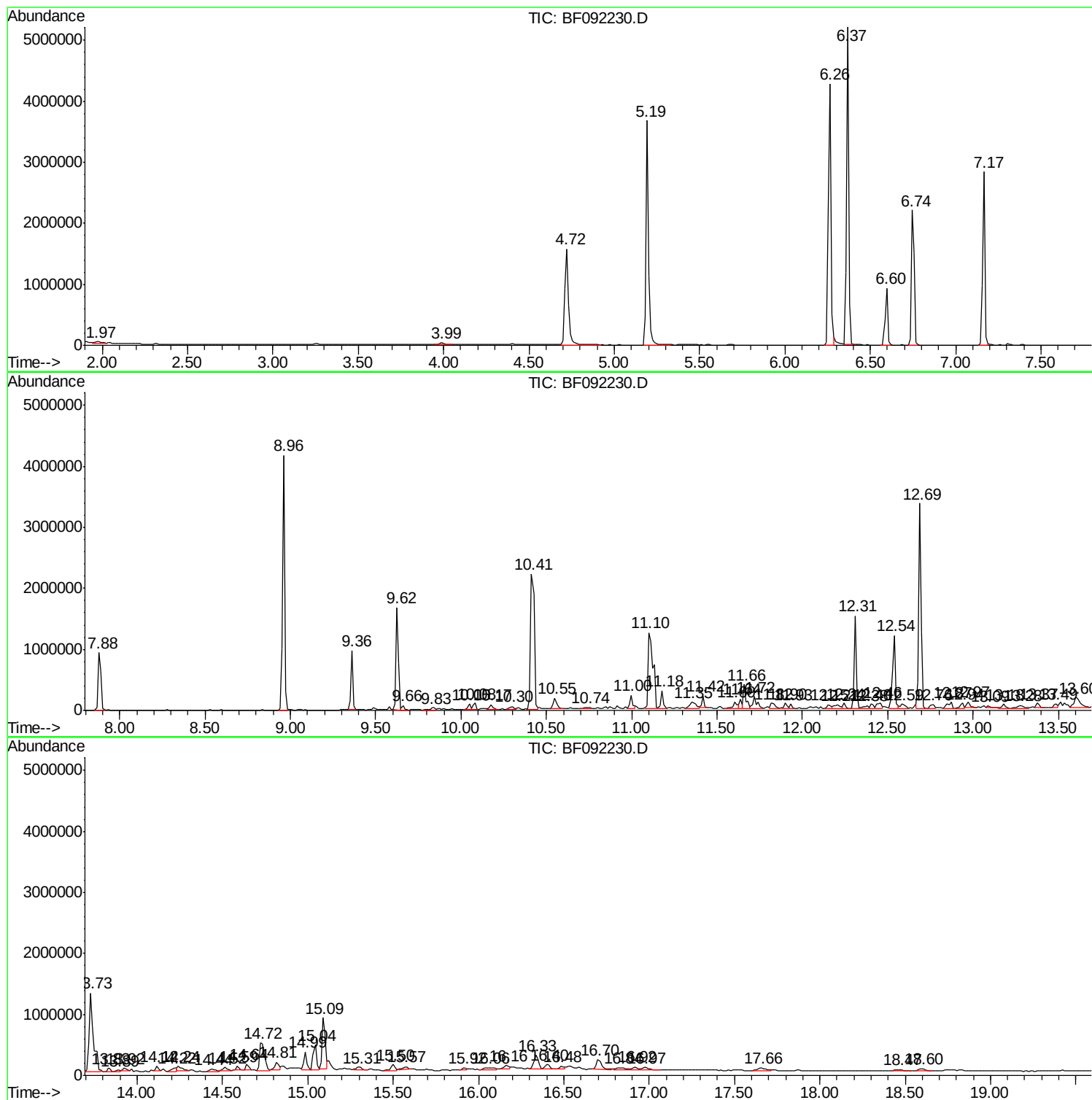
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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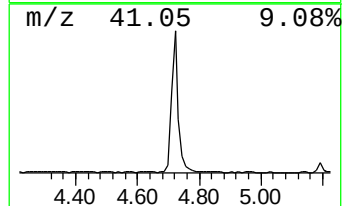
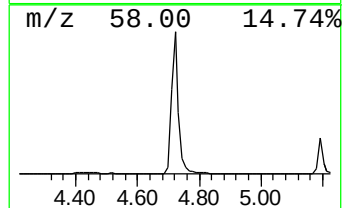
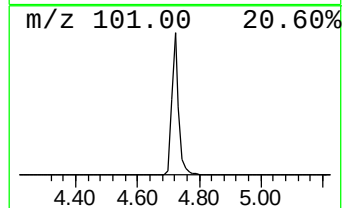
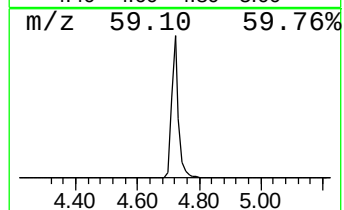
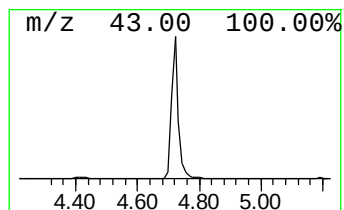
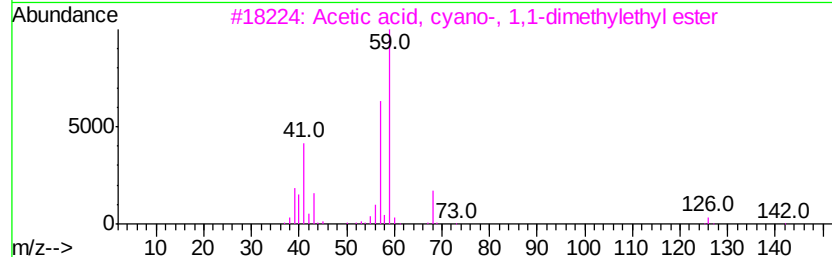
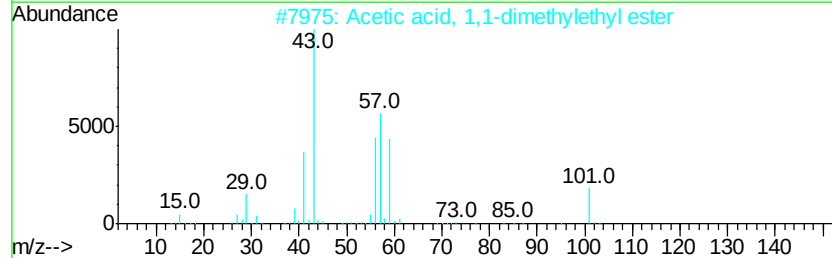
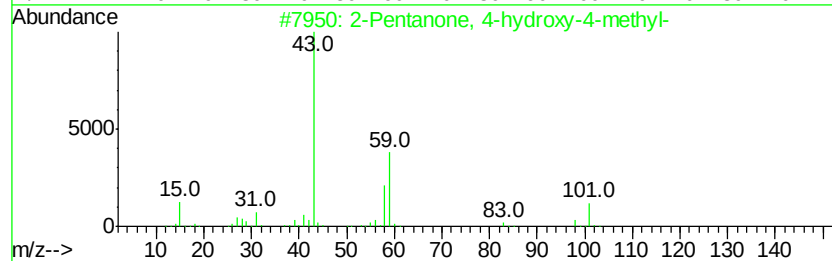
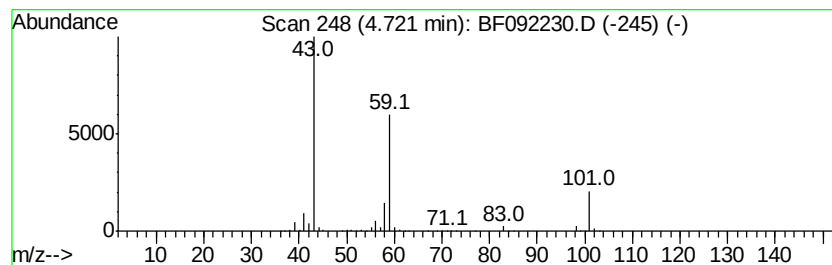
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
4.72	50.14 ng	2430880	1,4-Dichlorobenzene-d4	6.60

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72	
2	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39	
3	Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17	
4	Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9	
5	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9	



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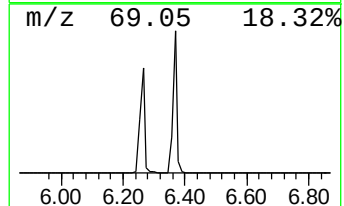
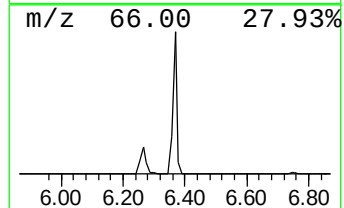
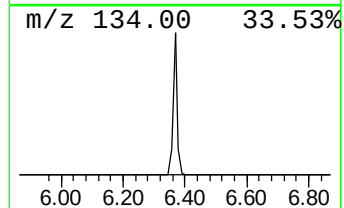
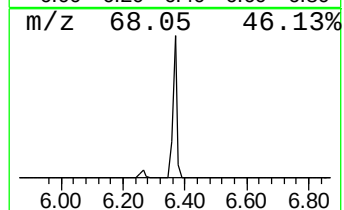
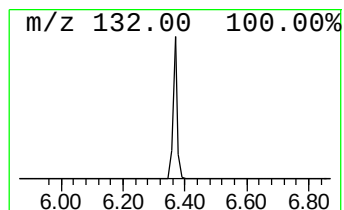
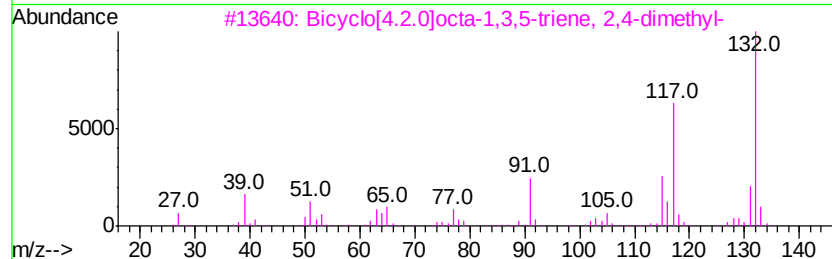
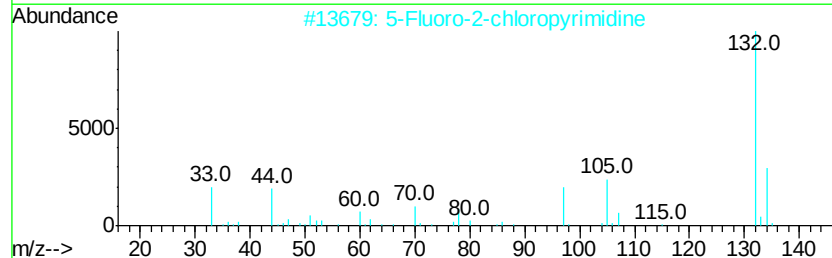
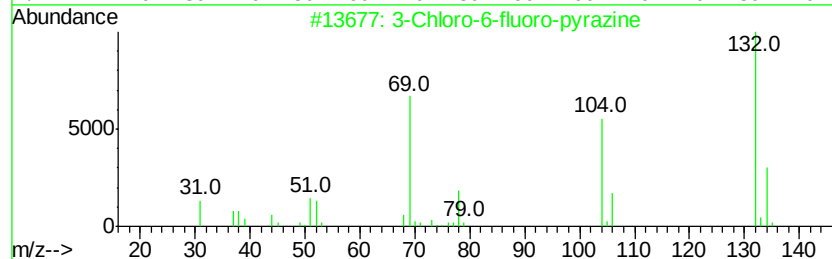
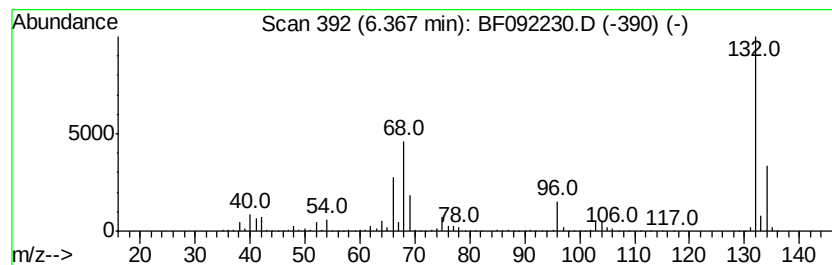
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 2 unknown6.37 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.37	100.02 ng	4848900	1,4-Dichlorobenzene-d4	6.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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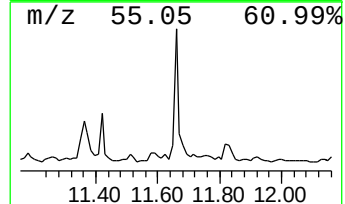
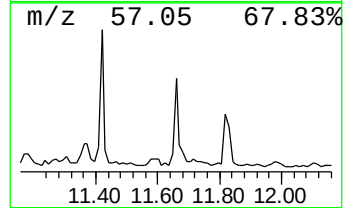
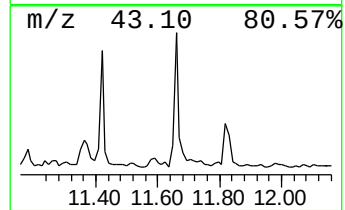
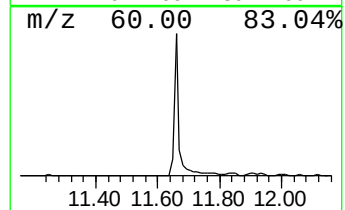
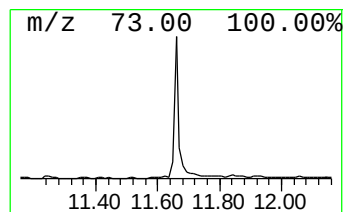
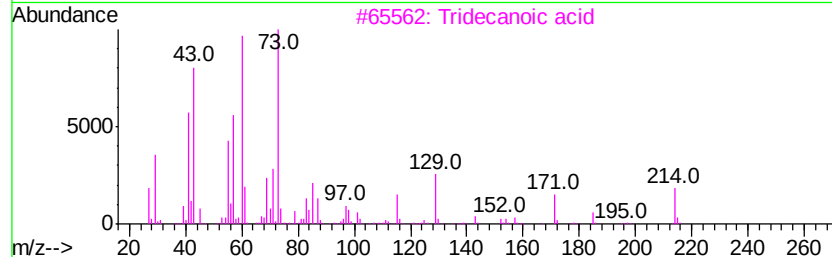
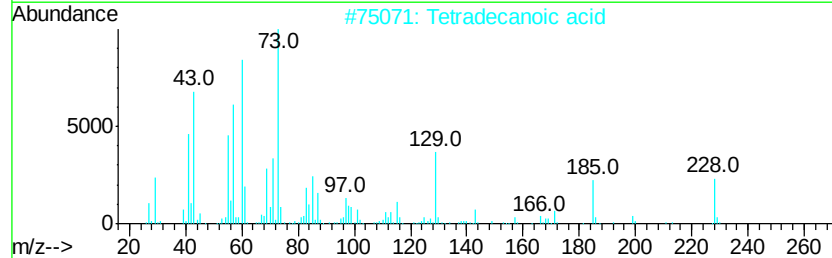
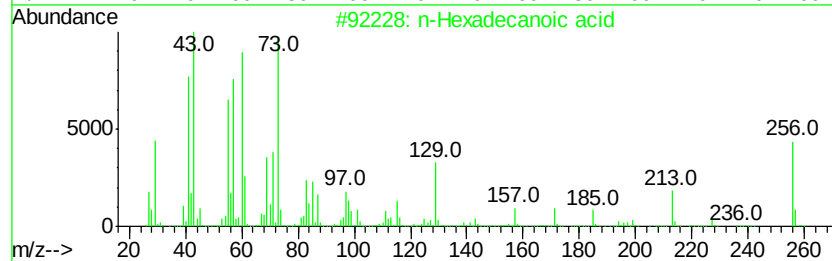
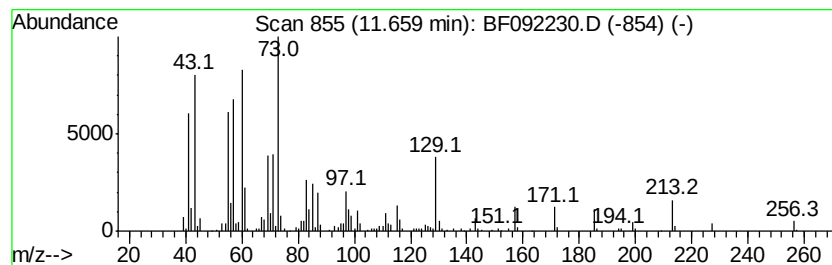
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.66	3.41 ng	442084	Phenanthrene-d10	11.11

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	86
3		Tridecanoic acid	214	C13H26O2	000638-53-9	81
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	80
5		n-Decanoic acid	172	C10H20O2	000334-48-5	68



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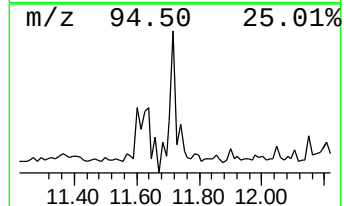
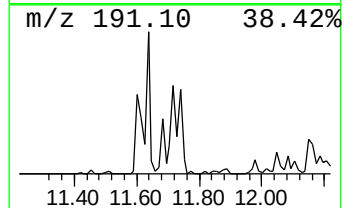
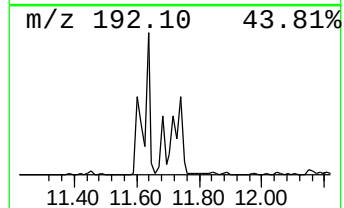
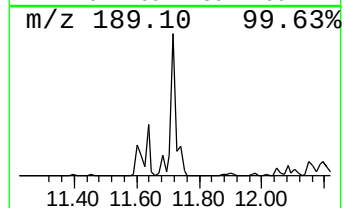
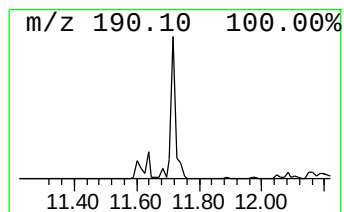
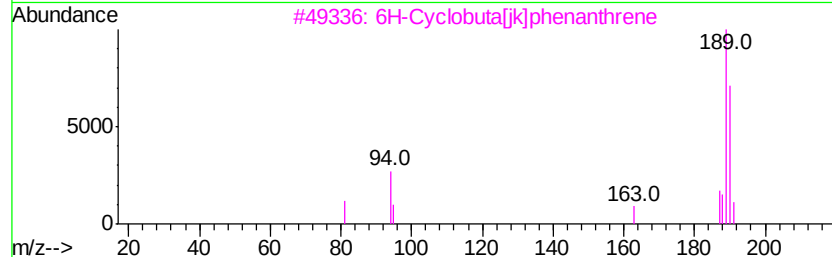
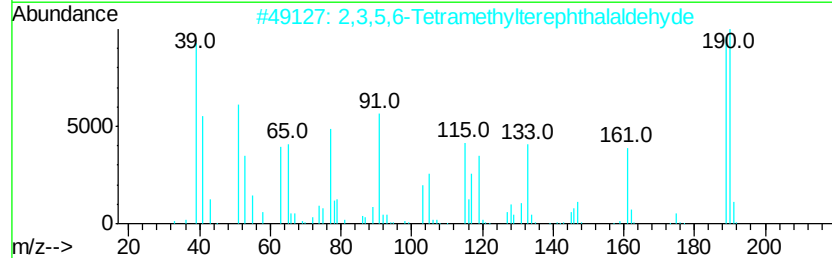
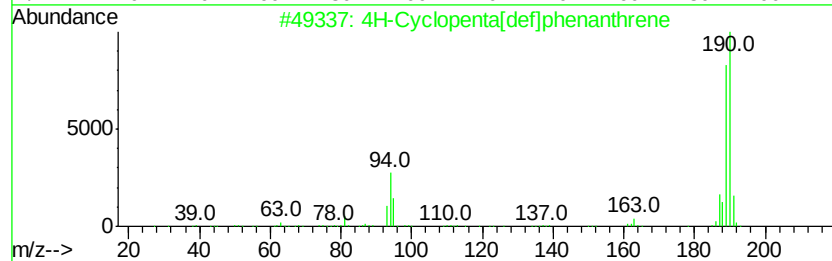
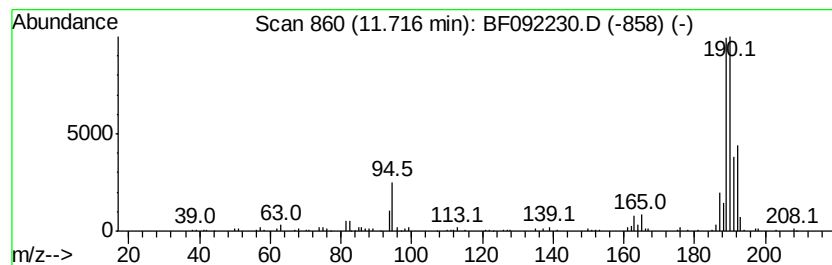
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 4H-Cyclopenta[def]phenanthrene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.		
11.72	2.34 ng	304106	Phenanthrene-d10	11.11		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	93
2		2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	43
3		6H-Cyclobuta[ik]phenanthrene	190	C15H10	083469-43-6	42
4		2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	40
5		1,4-Naphthalenedione, 2,5-dihydr...	190	C10H6O4	004923-55-1	25



Data Path : Z:\HPCHEM1\BNA F\DATA\BF011217\
Data File : BF092230.D
Acq On : 13 Jan 2017 2:52
Operator : UM/SJ
Sample : I1131-02
Misc :
ALS Vial : 25 Sample Multiplier: 1

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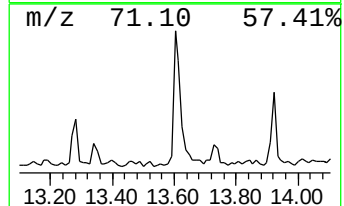
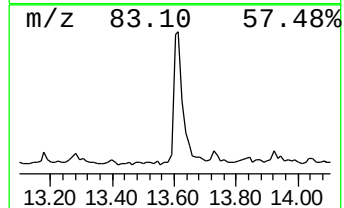
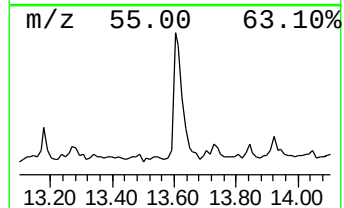
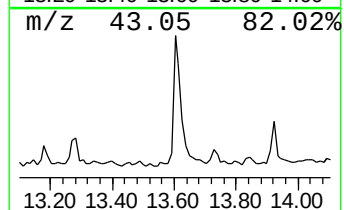
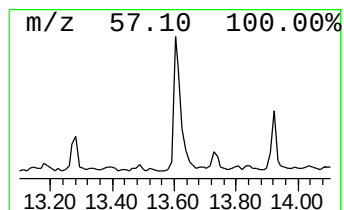
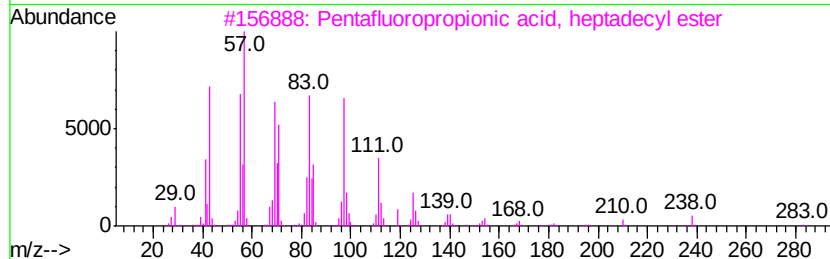
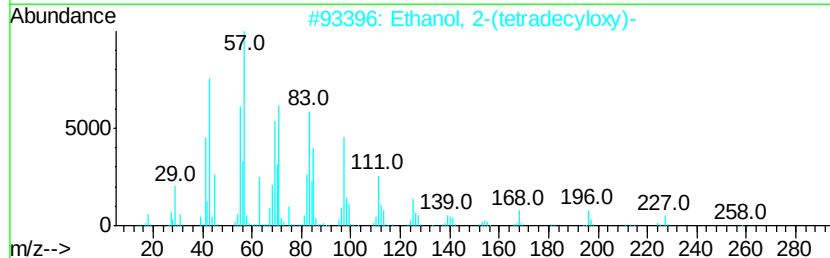
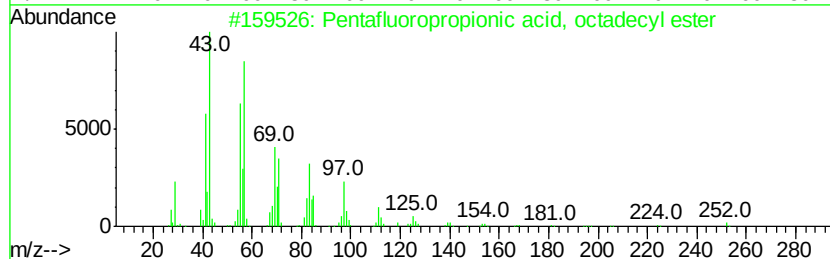
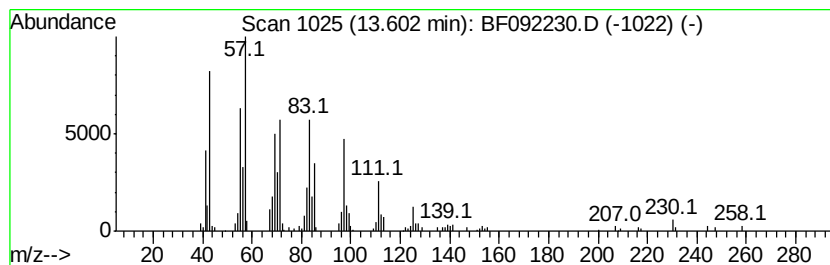
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Pentafluoropropionic acid, ... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.60	3.62 ng	400260	Chrysene-d12	13.73

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentafluoropropionic acid, octad...	416	C21H37F5O2	1000280-07-7	91
2		Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	91
3		Pentafluoropropionic acid, hepta...	402	C20H35F5O2	1000283-04-2	90
4		17-Pentatriacontene	491	C35H70	006971-40-0	83
5		Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	81



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011217\
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Acq On : 13 Jan 2017 2:52
Operator : UM/SJ
Sample : I1131-02
Misc :
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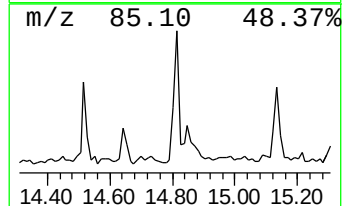
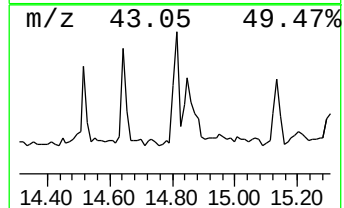
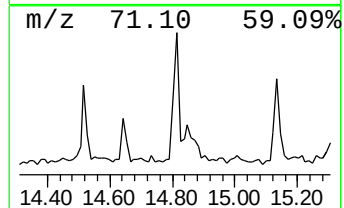
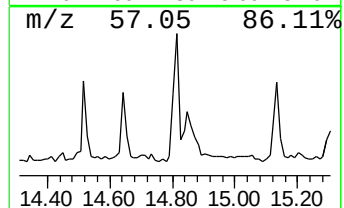
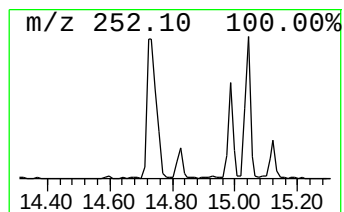
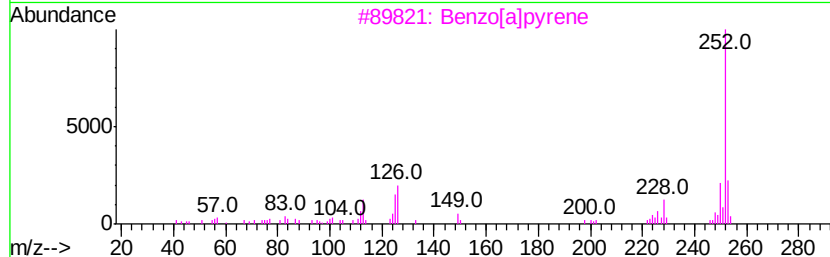
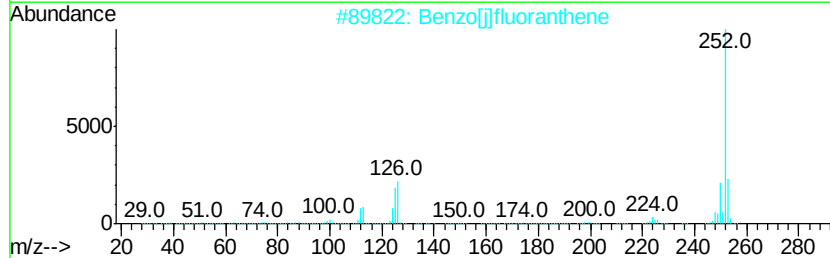
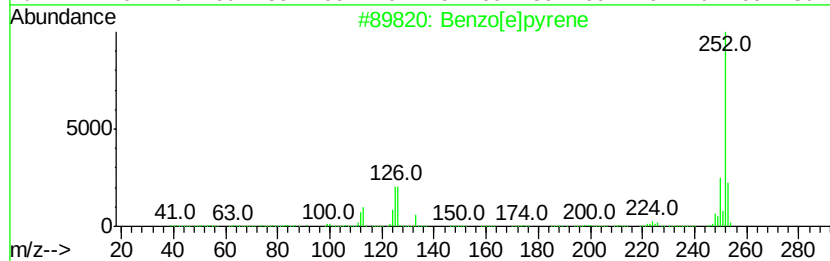
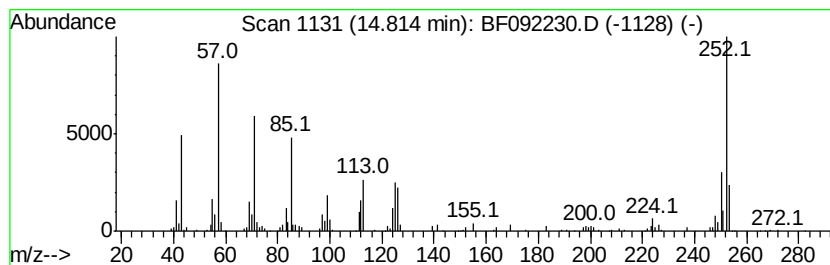
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Benzo[e]pyrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.81	3.66 ng	207762	Perylene-d12	15.09	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	90
2	Benzo[i]fluoranthene	252	C20H12	000205-82-3	87
3	Benzo[a]pyrene	252	C20H12	000050-32-8	78
4	Perylene	252	C20H12	000198-55-0	70
5	Benzo[k]fluoranthene	252	C20H12	000207-08-9	70



Data Path : Z:\HPCHEM1\BNA F\DATA\BF011217\
Data File : BF092230.D
Acq On : 13 Jan 2017 2:52
Operator : UM/SJ
Sample : I1131-02
Misc :
ALS Vial : 25 Sample Multiplier: 1

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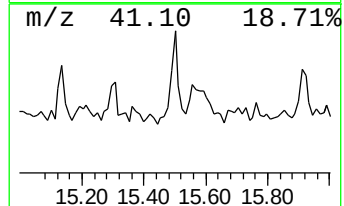
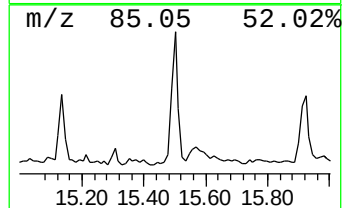
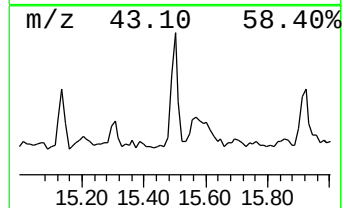
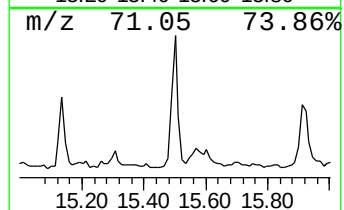
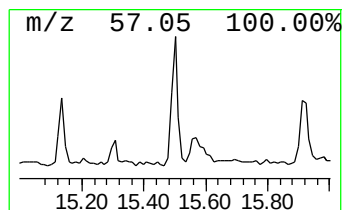
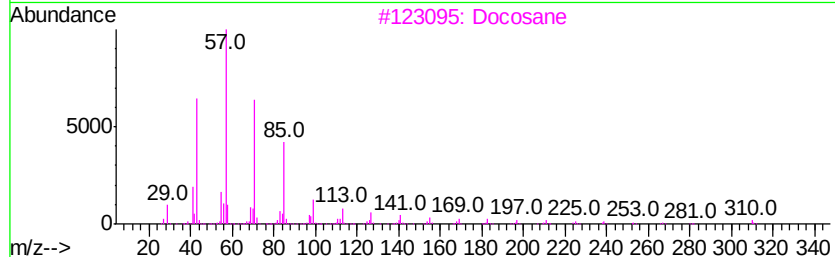
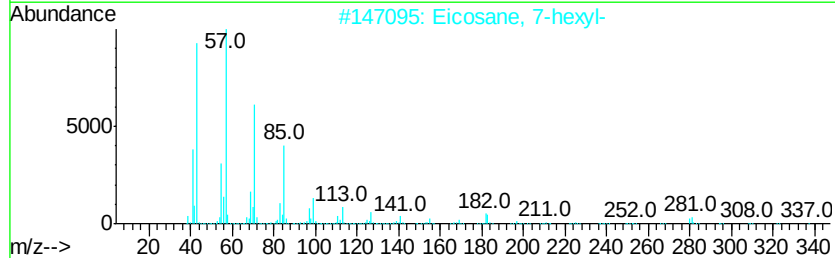
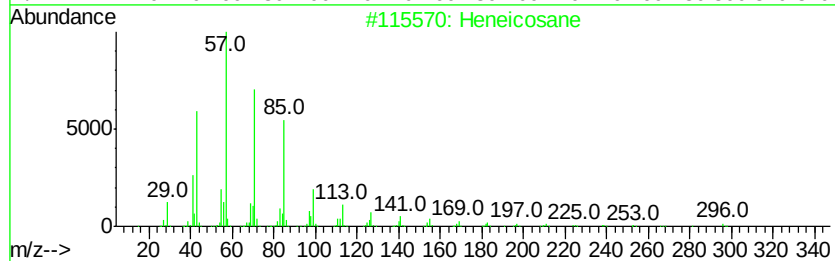
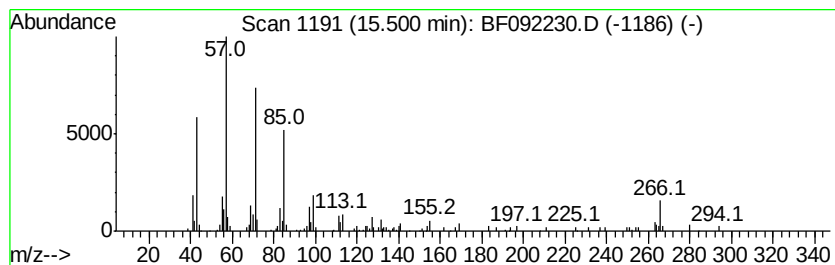
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Heneicosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.50	3.20 ng	181657	Perylene-d12	15.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	87
2		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	86
3		Docosane	310	C22H46	000629-97-0	83
4		Octacosane	394	C28H58	000630-02-4	83
5		Tridecane, 7-propyl-	226	C16H34	055045-09-5	83



Data Path : Z:\HPCHEM1\BNA F\DATA\BF011217\
Data File : BF092230.D
Acq On : 13 Jan 2017 2:52
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Sample : I1131-02
Misc :
ALS Vial : 25 Sample Multiplier: 1

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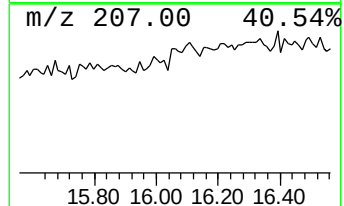
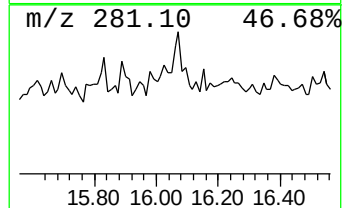
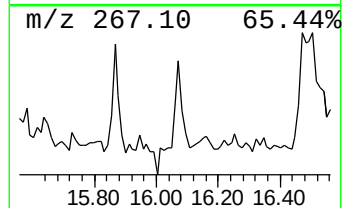
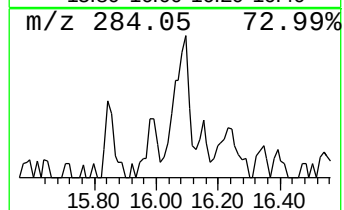
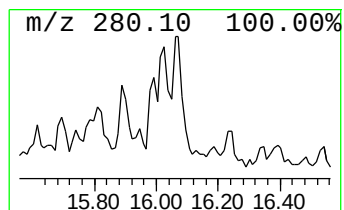
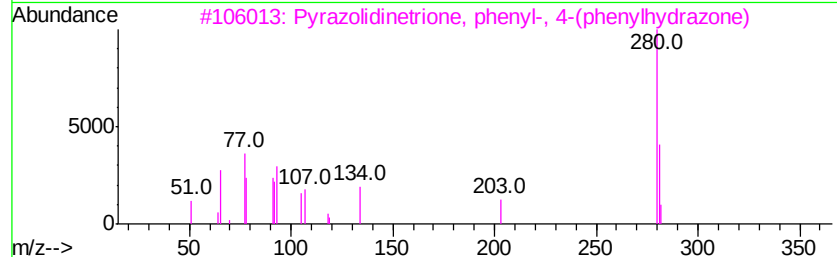
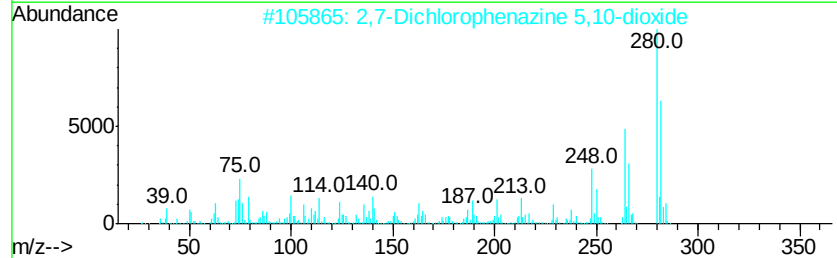
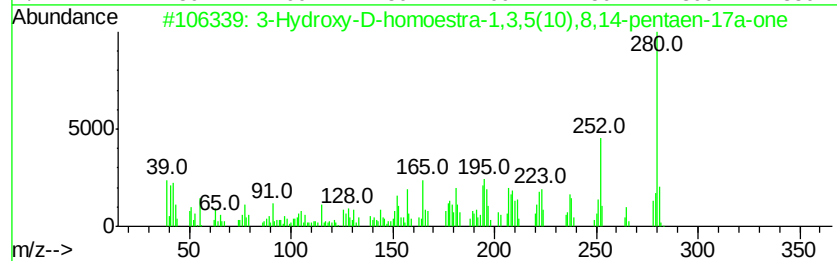
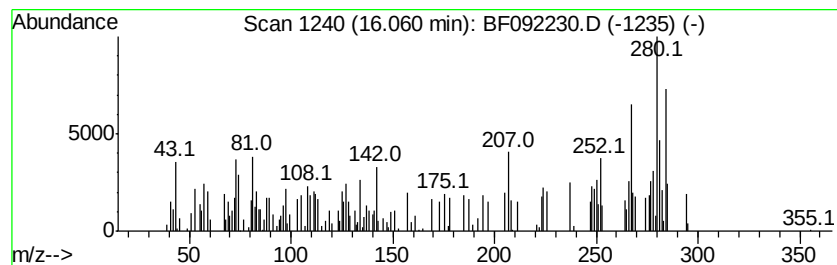
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown16.06 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.06	2.76 ng	157102	Perylene-d12	15.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Hydroxy-D-homoestra-1,3,5(10),...	280	C19H20O2	1000215-90-6	44
2		2,7-Dichlorophenazine 5,10-dioxide	280	C12H6Cl2N2O2	002089-15-8	27
3		Pyrazolidinetrione, phenyl-, 4-(...	280	C15H12N4O2	021272-26-4	27
4		3-Methoxy-D-homo-18-norestra-1,3...	280	C19H20O2	021070-83-7	22
5		Pentacene, 6,13-dihydro-	280	C22H16	013579-08-3	22



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011217\
Data File : BF092230.D
Acq On : 13 Jan 2017 2:52
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Sample : I1131-02
Misc :
ALS Vial : 25 Sample Multiplier: 1

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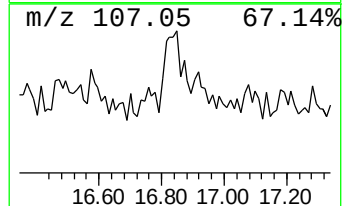
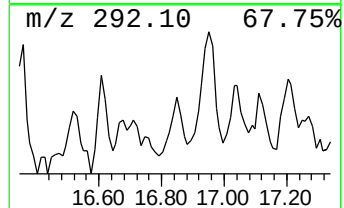
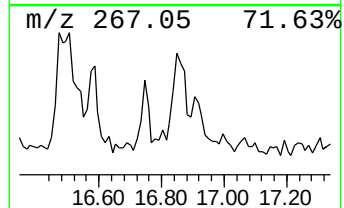
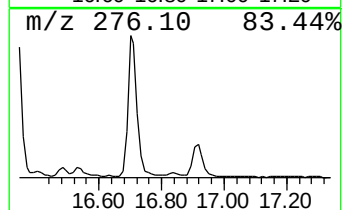
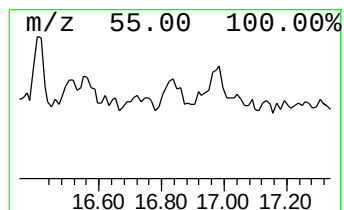
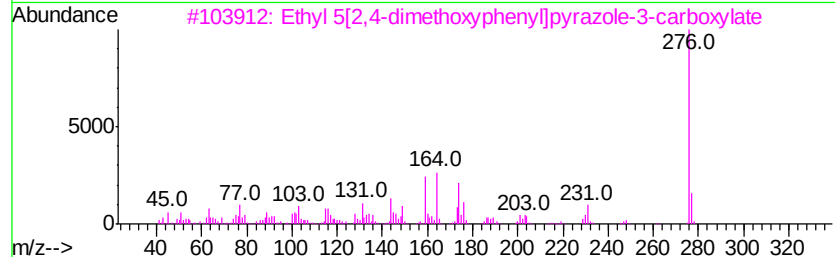
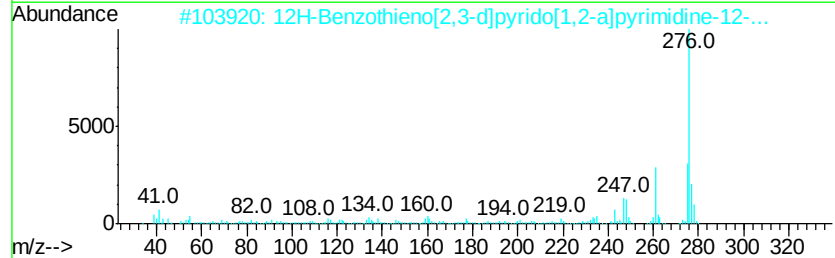
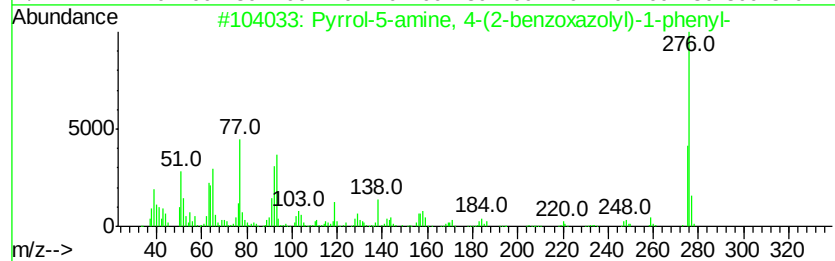
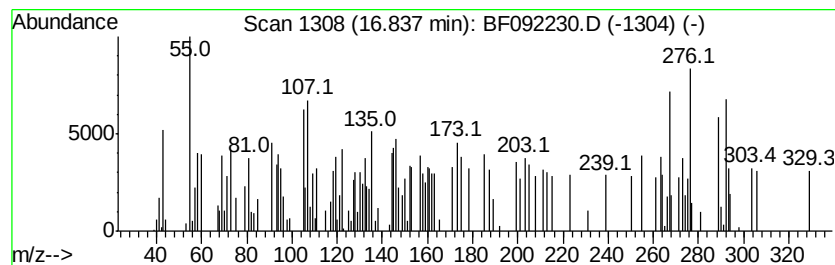
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown16.84 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.84	2.07 ng	117457	Perylene-d12	15.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pyrrol-5-amine, 4-(2-benzoxazolyl)-	276	C16H12N4O	039116-39-7	18
2		12H-Benzothieno[2,3-d]pyrido[1,2-...	276	C14H16N2S2	102254-63-7	14
3		Ethyl 5[2,4-dimethoxyphenyl]pyra...	276	C14H16N2O4	1000211-63-5	11
4		Thiambutene	291	C16H21NS2	000086-14-6	11
5		Diphenyl(4-tolyl)phosphine	276	C19H17P	001031-93-2	10



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011217\
Data File : BF092230.D
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Misc :
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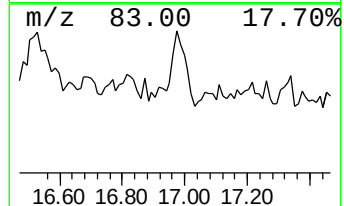
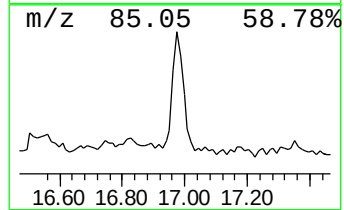
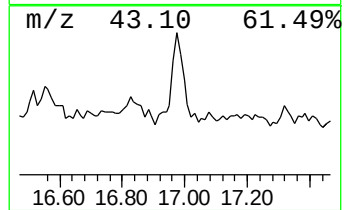
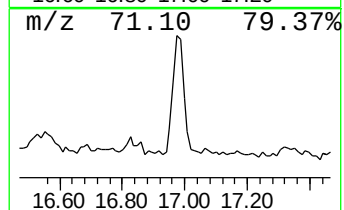
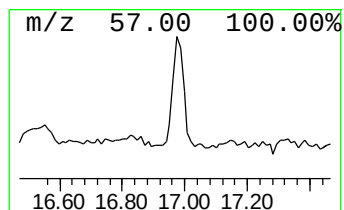
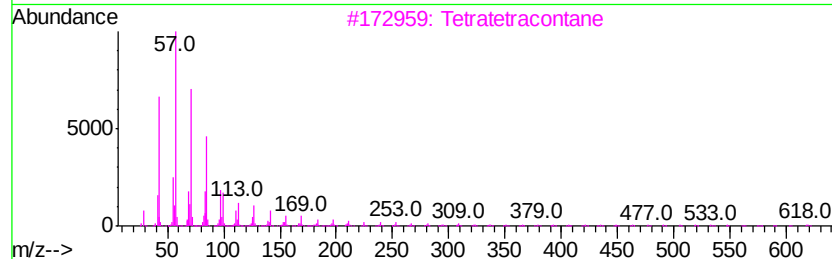
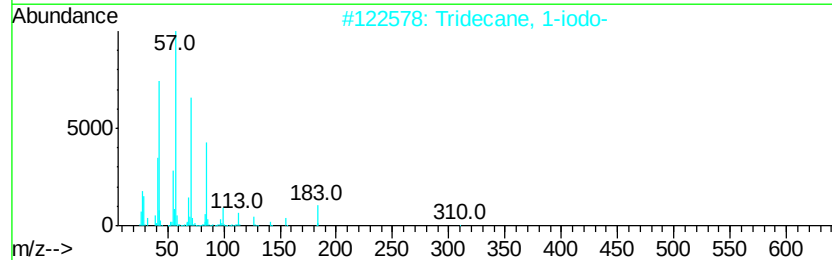
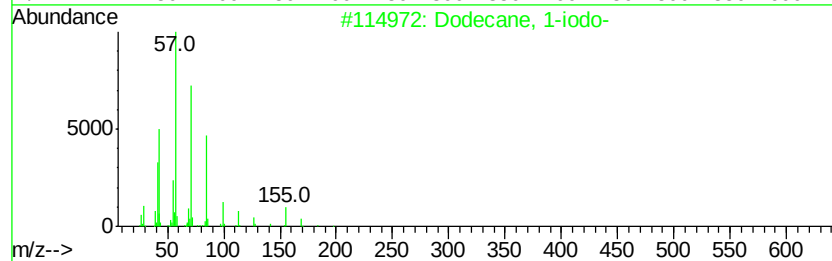
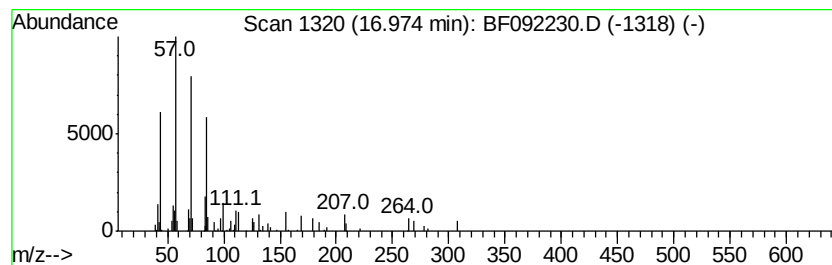
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Dodecane, 1-iodo- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.97	2.31 ng	131546	Perylene-d12	15.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dodecane, 1-iodo-	296	C12H25I	004292-19-7	83
2		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	72
3		Tetratetracontane	619	C44H90	007098-22-8	64
4		Undecane	156	C11H24	001120-21-4	64
5		Undecane, 2,7-dimethyl-	184	C13H28	017301-24-5	59



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011217\
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Sample : I1131-02
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SW-001

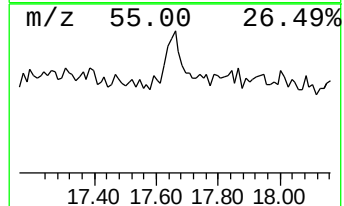
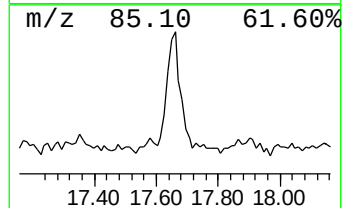
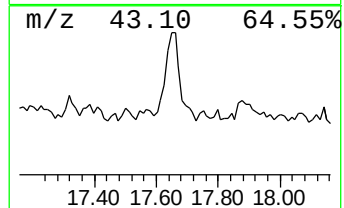
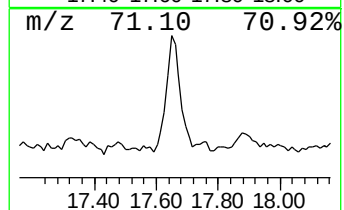
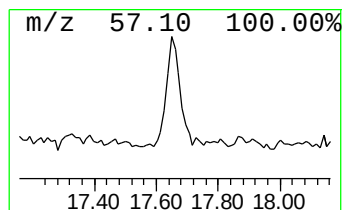
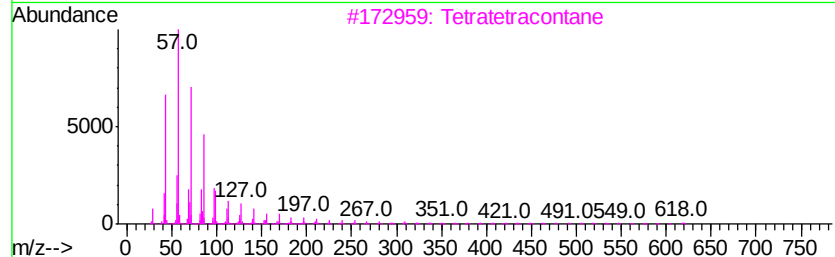
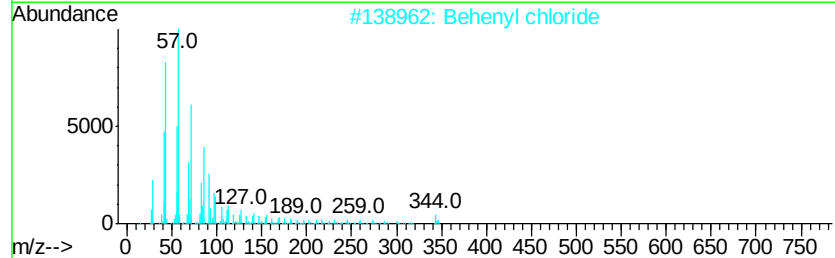
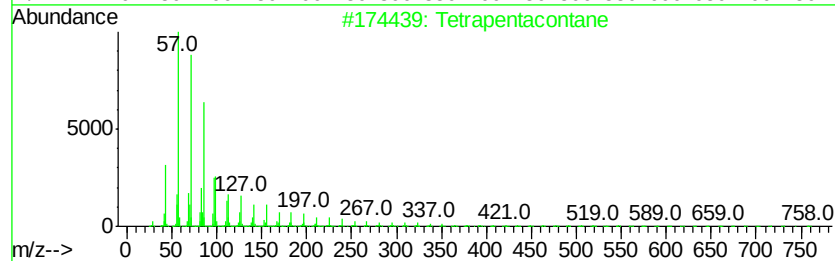
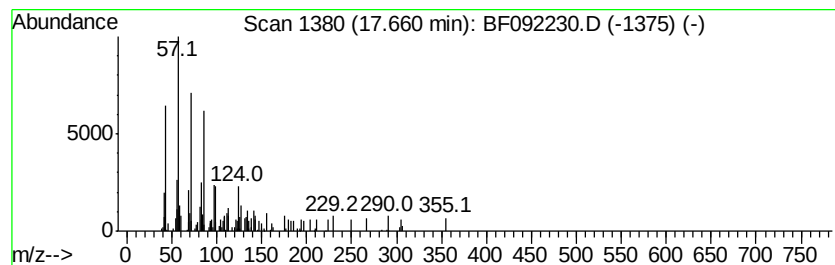
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Tetrapentacontane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.66	2.94 ng	166855	Perylene-d12	15.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetrapentacontane	759	C54H110	005856-66-6	58
2		Behenyl chloride	344	C22H45Cl	042217-03-8	58
3		Tetratetracontane	619	C44H90	007098-22-8	58
4		Docosane	310	C22H46	000629-97-0	53
5		Tetracosane, 11-decyl-	479	C34H70	055429-84-0	52



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011217\
Data File : BF092230.D
Acq On : 13 Jan 2017 2:52
Operator : UM/SJ
Sample : I1131-02
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SW-001

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...	4.72	50.1 ng		2430880	1	6.60	969592	20.0
unknown6.37	6.37	100.0 ng		4848900	1	6.60	969592	20.0
n-Hexadecanoic acid	11.66	3.4 ng		442084	4	11.11	2594420	20.0
4H-Cyclopenta[def...	11.72	2.3 ng		304106	4	11.11	2594420	20.0
Pentafluoropropio...	13.60	3.6 ng		400260	5	13.73	2211250	20.0
Benzo[e]pyrene	14.81	3.7 ng		207762	6	15.09	1136490	20.0
Heneicosane	15.50	3.2 ng		181657	6	15.09	1136490	20.0
unknown16.06	16.06	2.8 ng		157102	6	15.09	1136490	20.0
unknown16.84	16.84	2.1 ng		117457	6	15.09	1136490	20.0
Dodecane, 1-iodo-	16.97	2.3 ng		131546	6	15.09	1136490	20.0
Tetrapentacontane	17.66	2.9 ng		166855	6	15.09	1136490	20.0