

Data Path : U:\HPCHEM1\BNA F\DATA\BF082217\
 Data File : BF097988.D
 Acq On : 23 Aug 2017 7:57
 Operator : SJ/JU
 Sample : I4913-01
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TEC-BE2-1-3-COMP

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-BF081517.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	3.128	174	177	193	rBV	49866	109096	1.79%	0.257%
2	4.951	484	487	496	rVB	198703	297185	4.88%	0.700%
3	5.363	551	557	560	rBV	3208802	3492561	57.36%	8.225%
4	5.575	590	593	598	rVB	96734	79991	1.31%	0.188%
5	6.393	726	732	741	rBV	3182805	3745555	61.51%	8.820%
6	6.522	749	754	757	rBV	3356138	3461010	56.84%	8.150%
7	6.751	789	793	796	rBV	1004170	863153	14.17%	2.033%
8	6.910	815	820	823	rBV	2777905	2727519	44.79%	6.423%
9	7.316	883	889	892	rBV	2214191	2215204	36.38%	5.217%
10	8.034	1006	1011	1014	rBV	1237708	1046986	17.19%	2.466%
11	8.798	1137	1141	1147	rBV	125104	156373	2.57%	0.368%
12	9.110	1189	1194	1197	rBV	3686305	3573169	58.68%	8.414%
13	9.504	1257	1261	1264	rBV	557178	463064	7.60%	1.090%
14	9.786	1303	1309	1312	rBV	1077383	947256	15.56%	2.231%
15	10.569	1438	1442	1446	rBV	1513153	1559871	25.62%	3.673%
16	10.692	1460	1463	1471	rVB	91009	91560	1.50%	0.216%
17	11.263	1557	1560	1563	rBV	860769	826921	13.58%	1.947%
18	11.286	1563	1564	1568	rVV	309008	210738	3.46%	0.496%
19	11.339	1568	1573	1576	rVB	94215	90554	1.49%	0.213%
20	11.763	1642	1645	1648	rBV	63985	64055	1.05%	0.151%
21	11.792	1648	1650	1653	rVV	89019	68729	1.13%	0.162%
22	11.833	1653	1657	1661	rVB	575990	477920	7.85%	1.125%
23	11.875	1661	1664	1667	rBV	115354	119113	1.96%	0.280%
24	12.316	1735	1739	1742	rBV2	92631	93053	1.53%	0.219%
25	12.475	1762	1766	1770	rBV	1086361	904206	14.85%	2.129%
26	12.627	1784	1792	1794	rBV2	38402	61594	1.01%	0.145%
27	12.704	1797	1805	1808	rBV	868106	915859	15.04%	2.157%
28	12.857	1826	1831	1834	rVV	2440351	2535451	41.64%	5.971%
29	12.922	1837	1842	1845	rBV2	58986	97773	1.61%	0.230%
30	13.033	1858	1861	1864	rVB	131990	126867	2.08%	0.299%
31	13.098	1868	1872	1874	rVV	103822	86164	1.41%	0.203%
32	13.133	1874	1878	1882	rBV	104319	111879	1.84%	0.263%
33	13.333	1908	1912	1914	rBV2	103518	116183	1.91%	0.274%
34	13.533	1943	1946	1951	rVB	76984	93059	1.53%	0.219%

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Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	13.645	1961	1965	1968	rBV2	101929	132068	2.17%	0.311%
36	13.674	1968	1970	1972	rVV	94492	89797	1.47%	0.211%
37	13.710	1972	1976	1979	rVV2	157731	208411	3.42%	0.491%
38	13.751	1979	1983	1988	rVB2	185248	283998	4.66%	0.669%
39	13.898	2003	2008	2011	rBV	4958216	6089331	100.00%	14.340%
40	13.927	2011	2013	2018	rVB	509559	443624	7.29%	1.045%
41	14.004	2023	2026	2029	rVB	136173	107901	1.77%	0.254%
42	14.286	2071	2074	2077	rVB2	88426	77418	1.27%	0.182%
43	14.380	2084	2090	2093	rBV7	90934	153645	2.52%	0.362%
44	14.468	2103	2105	2109	rVB4	66582	81886	1.34%	0.193%
45	14.745	2149	2152	2157	rBV2	137839	158630	2.61%	0.374%
46	14.916	2177	2181	2190	rBV	385616	731812	12.02%	1.723%
47	14.998	2192	2195	2197	rBV2	62681	78317	1.29%	0.184%
48	15.204	2226	2230	2235	rVB	194400	229214	3.76%	0.540%
49	15.263	2236	2240	2246	rVV	286102	347530	5.71%	0.818%
50	15.321	2246	2250	2253	rVV	482819	552102	9.07%	1.300%
51	15.351	2253	2255	2262	rVB2	104600	172571	2.83%	0.406%
52	15.774	2323	2327	2332	rBV2	81454	129261	2.12%	0.304%
53	16.257	2405	2409	2412	rBV2	51690	75359	1.24%	0.177%
54	16.698	2480	2484	2494	rVB2	103018	198252	3.26%	0.467%
55	16.810	2500	2503	2509	rVB3	38176	64241	1.05%	0.151%
56	17.115	2550	2555	2561	rVB	67745	124655	2.05%	0.294%
57	17.468	2612	2615	2626	rVB5	45822	104926	1.72%	0.247%

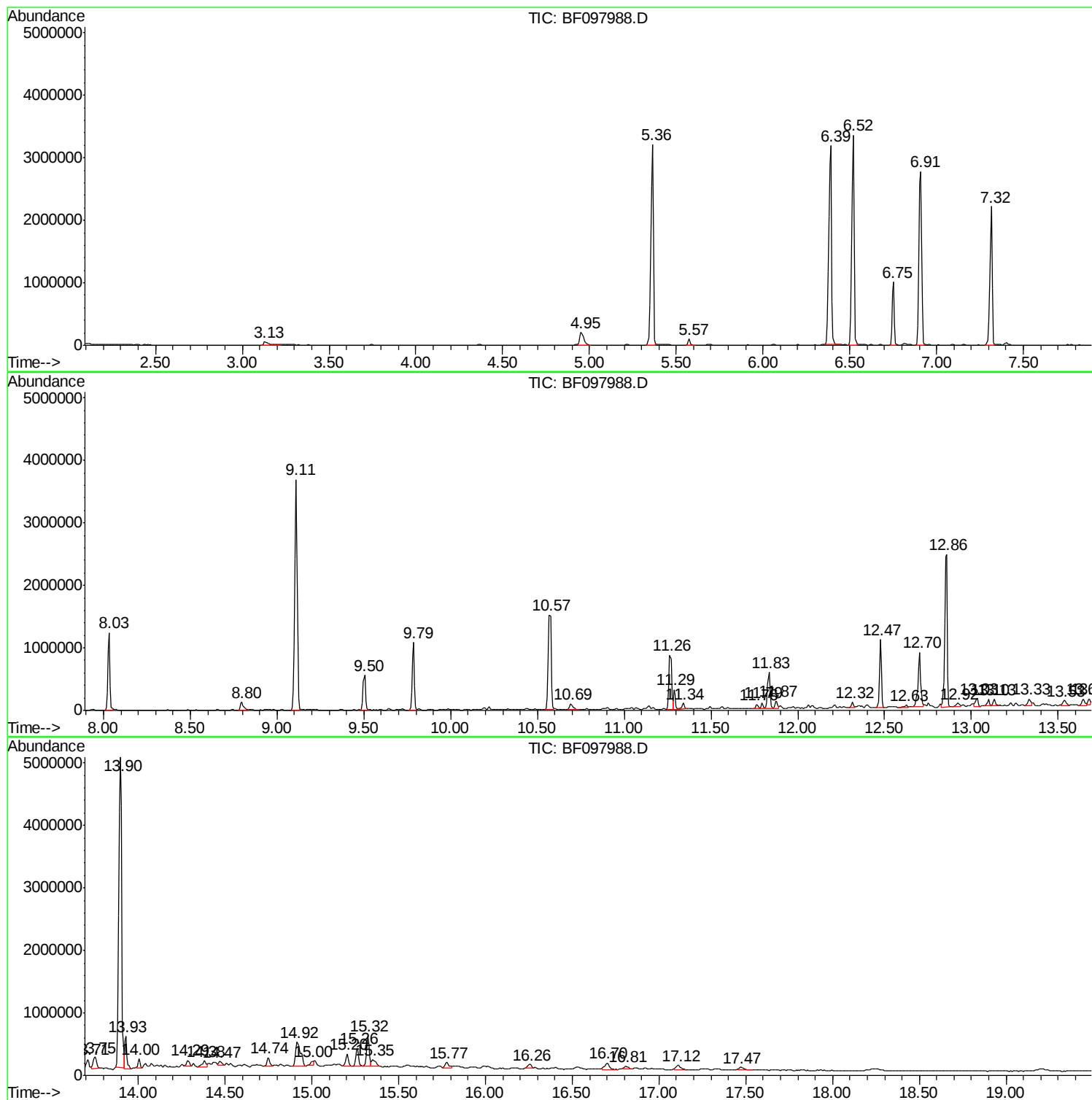
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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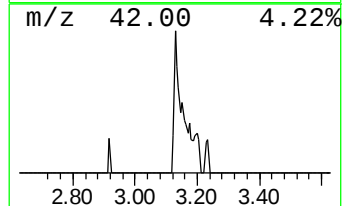
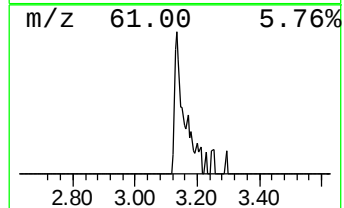
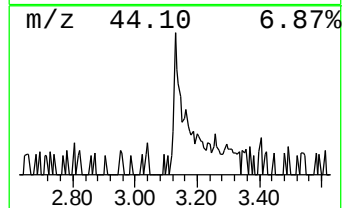
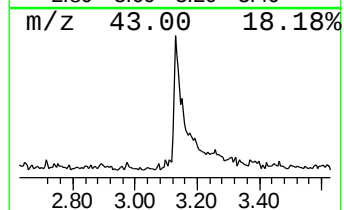
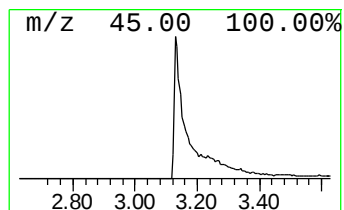
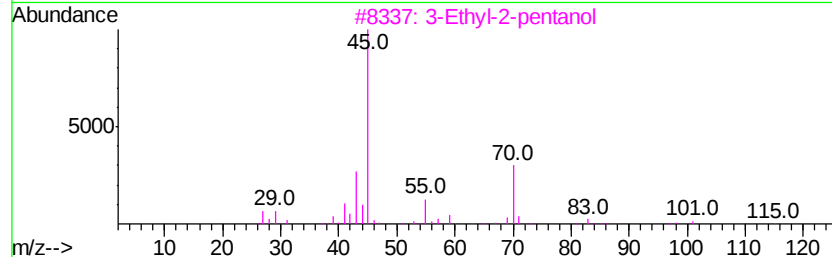
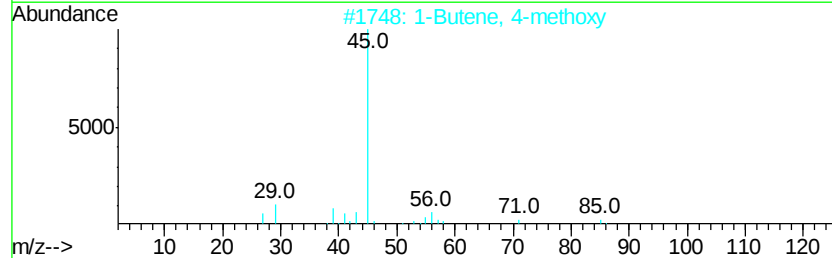
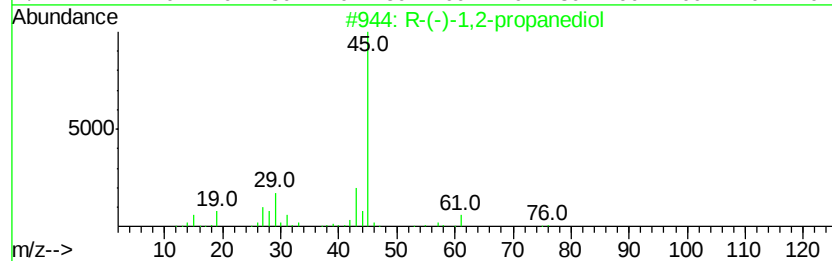
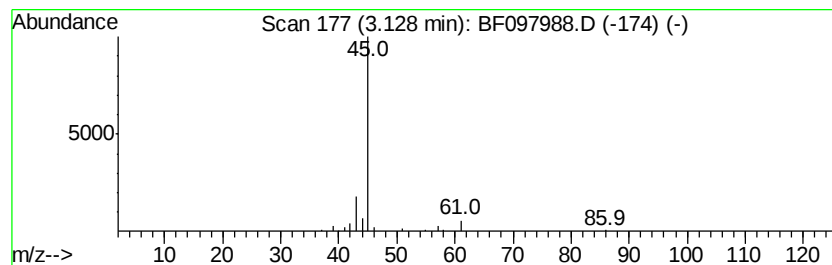
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown3.13 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.13	2.53 ng	109096	1,4-Dichlorobenzene-d4	6.75

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	43
2			1-Butene, 4-methoxy	86	C5H10O	004696-30-4	9
3			3-Ethyl-2-pentanol	116	C7H16O	000609-27-8	9
4			2-Hexanol, 3-methyl-	116	C7H16O	002313-65-7	9
5			Isopropyl Alcohol	60	C3H8O	000067-63-0	9



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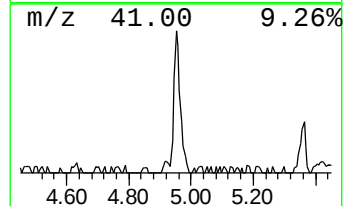
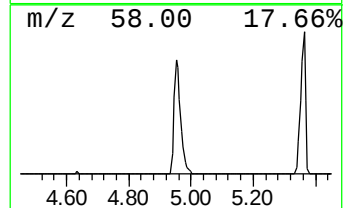
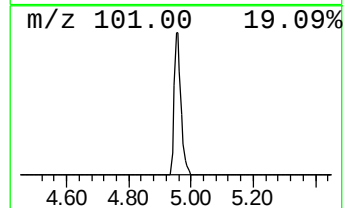
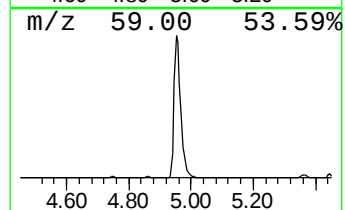
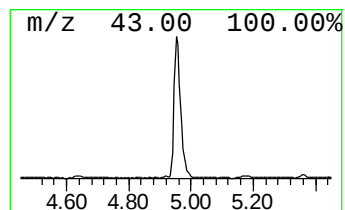
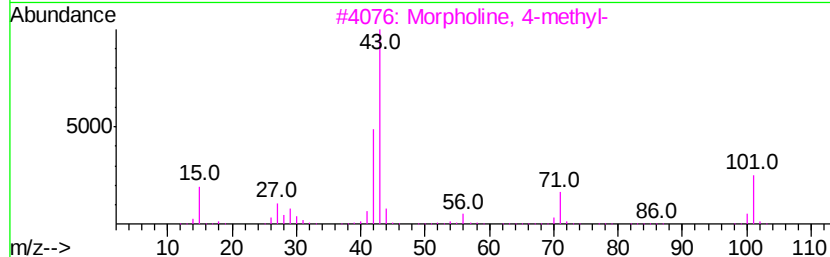
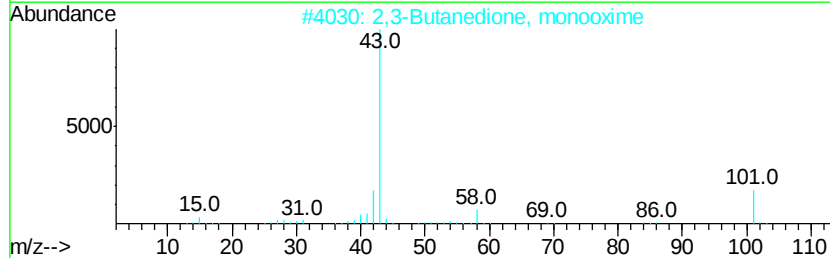
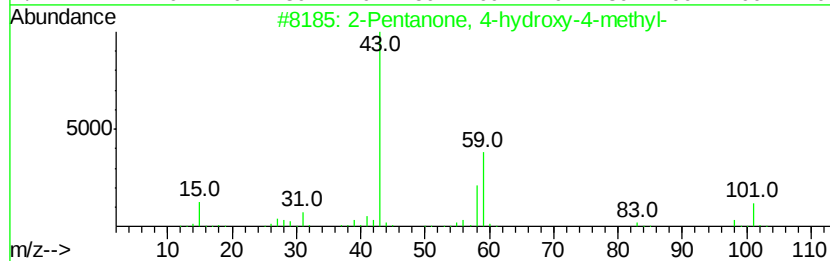
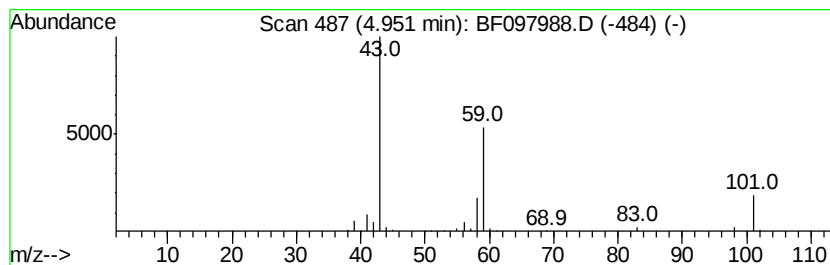
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.95	6.89 ng	297185	1,4-Dichlorobenzene-d4	6.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
3		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4		3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	9
5		Butanamide, 2-hydroxy-2,N-dimethyl-	117	C5H11NO2	039961-72-3	9



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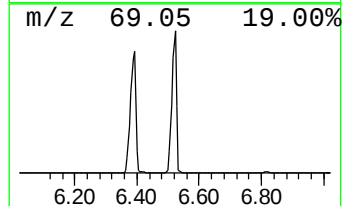
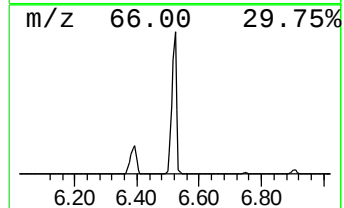
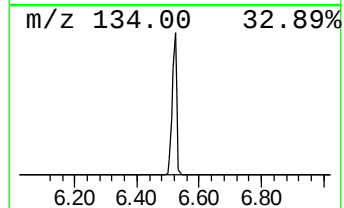
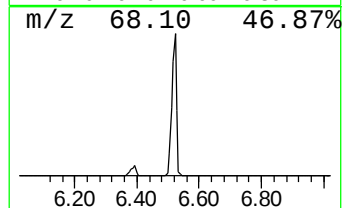
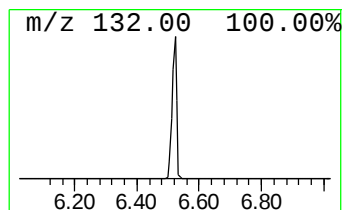
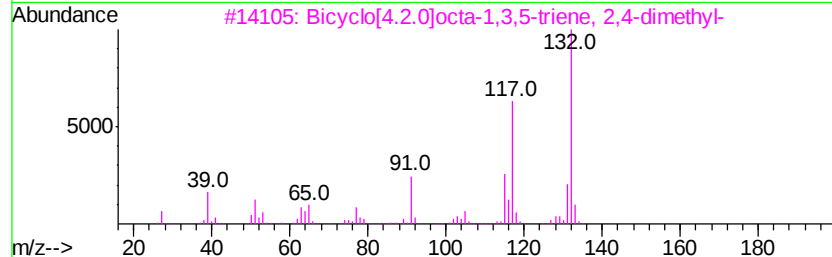
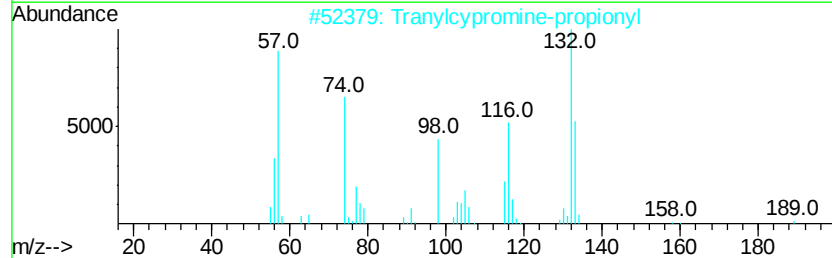
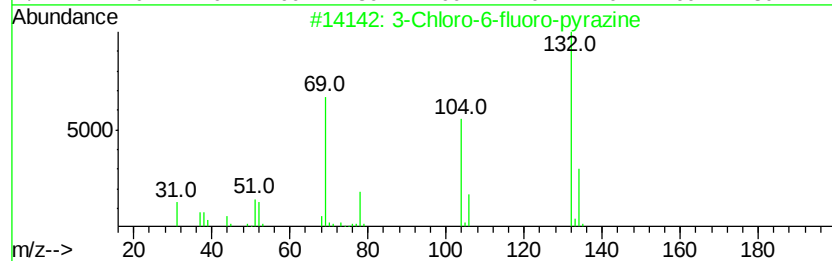
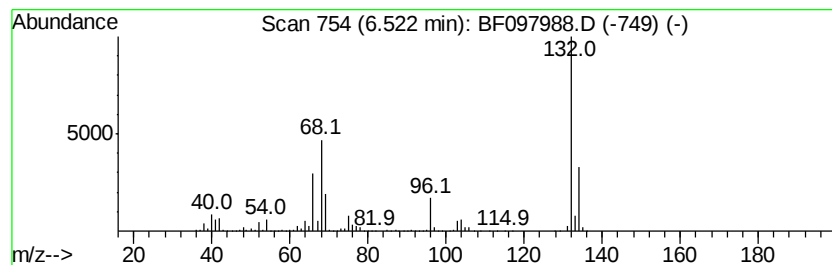
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.52 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.52	80.19 ng	3461010	1,4-Dichlorobenzene-d4	6.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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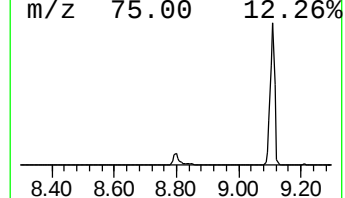
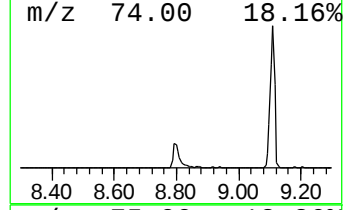
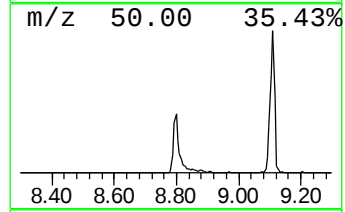
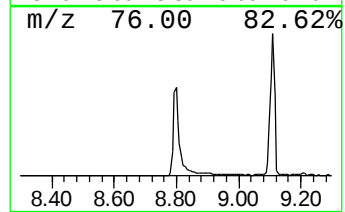
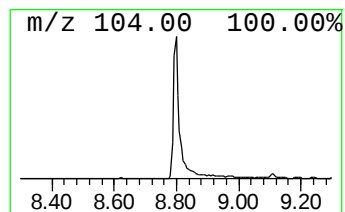
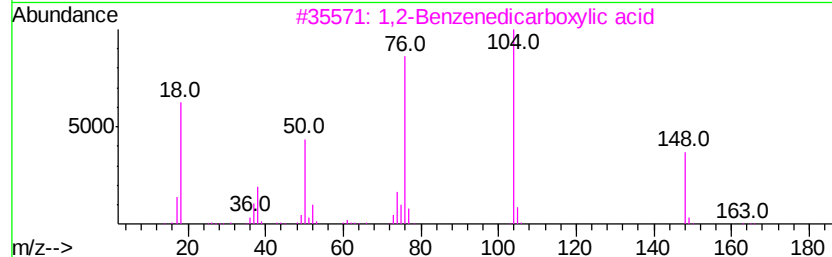
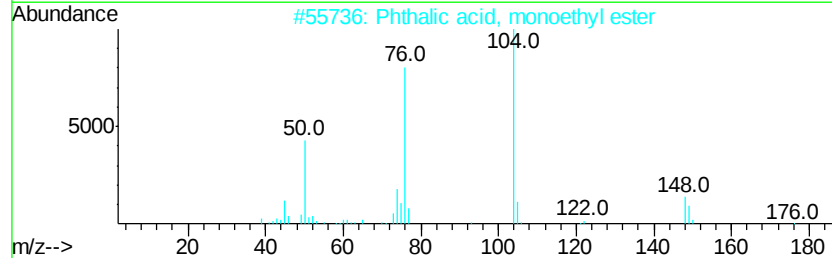
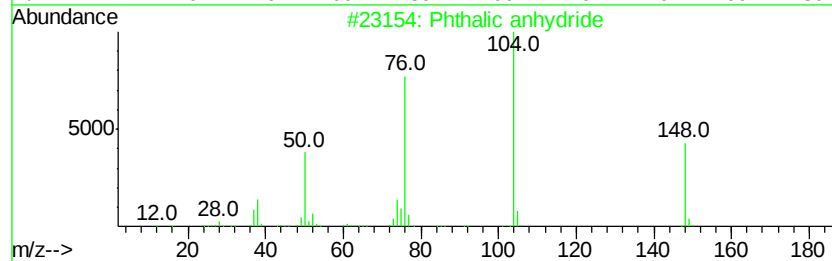
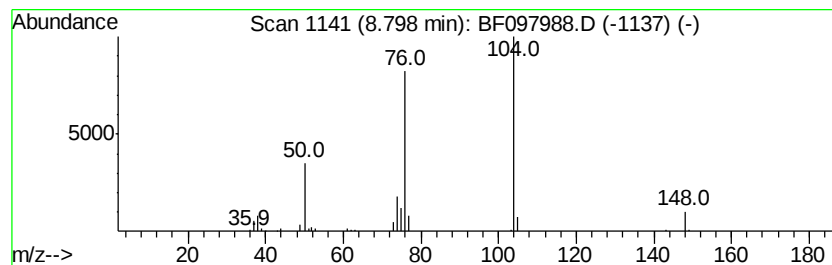
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Phthalic anhydride Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.80	2.99 ng	156373	Naphthalene-d8	8.03

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phthalic anhydride	148	C8H4O3	000085-44-9	95
2		Phthalic acid, monoethyl ester	194	C10H10O4	002306-33-4	90
3		1,2-Benzenedicarboxylic acid	166	C8H6O4	000088-99-3	83
4		Bicyclo[4.2.0]octa-1,3,5-triene...	132	C8H4O2	006383-11-5	83
5		Phthalimide	147	C8H5NO2	000085-41-6	78



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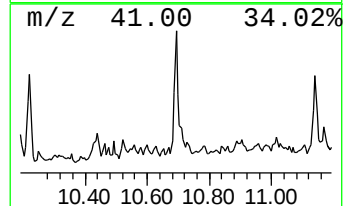
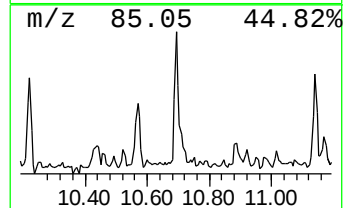
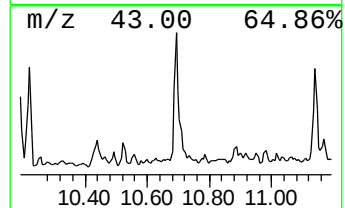
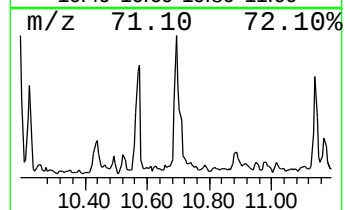
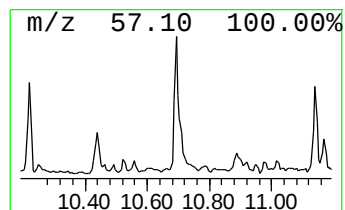
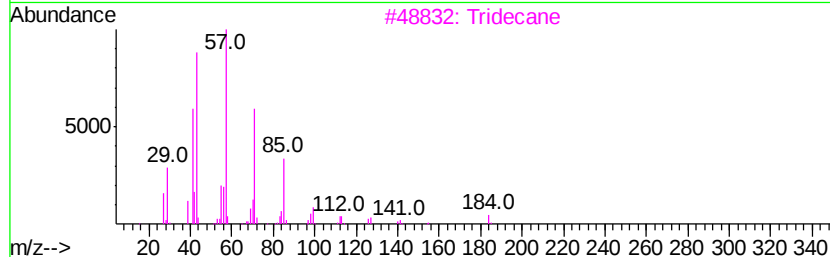
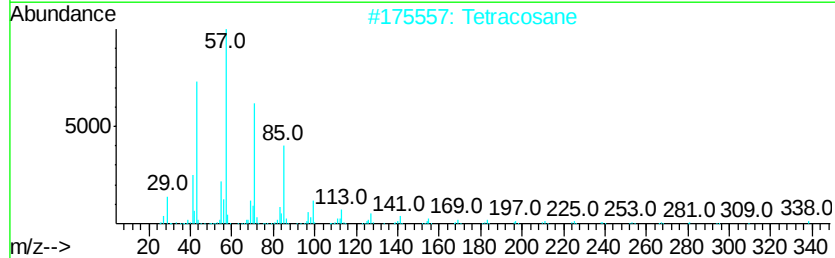
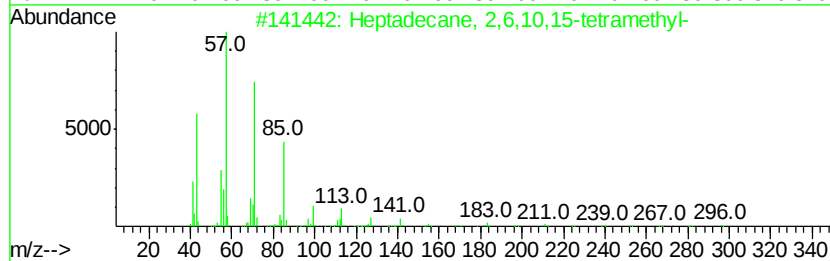
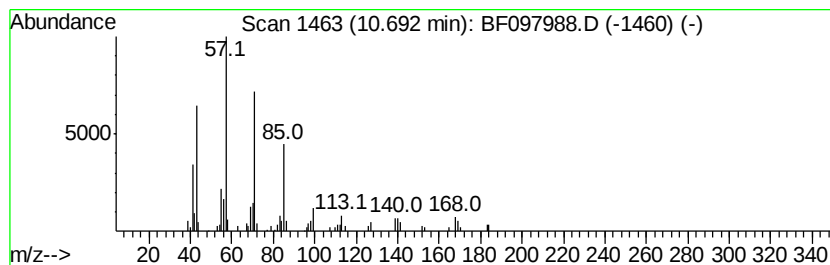
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Heptadecane, 2,6,10,15-tetr... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.69	2.21 ng	91560	Phenanthrene-d10		11.26
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	83
2	Tetracosane	338	C24H50	000646-31-1	83
3	Tridecane	184	C13H28	000629-50-5	83
4	Eicosane, 10-methyl-	296	C21H44	054833-23-7	80
5	Dodecane, 1-iodo-	296	C12H25I	004292-19-7	80



Data Path : U:\HPCHEM1\BNA F\DATA\BF082217\
Data File : BF097988.D
Acq On : 23 Aug 2017 7:57
Operator : SJ/JU
Sample : I4913-01
Misc :
ALS Vial : 30 Sample Multiplier: 1

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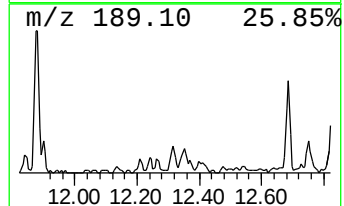
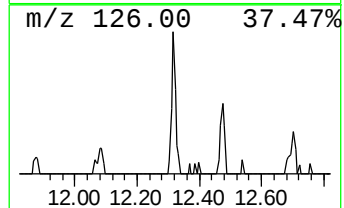
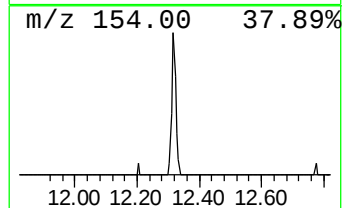
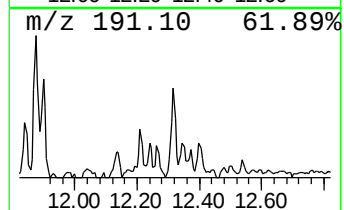
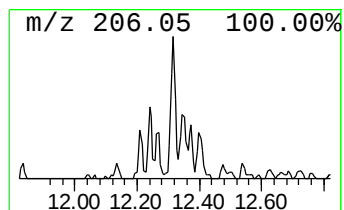
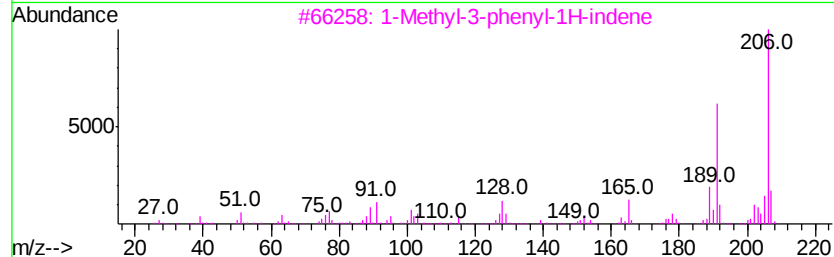
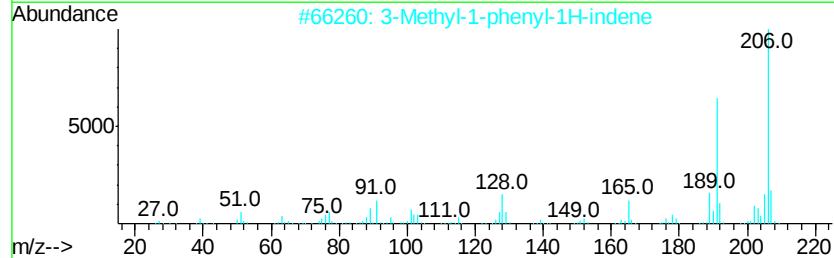
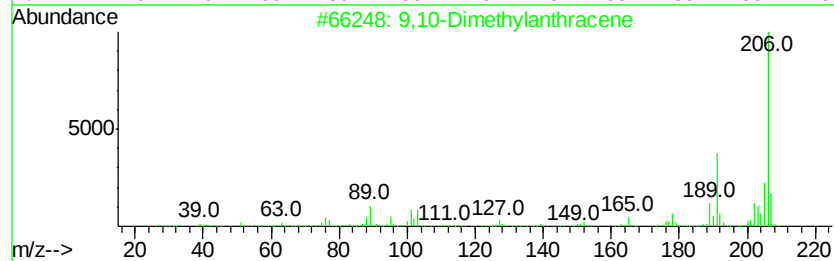
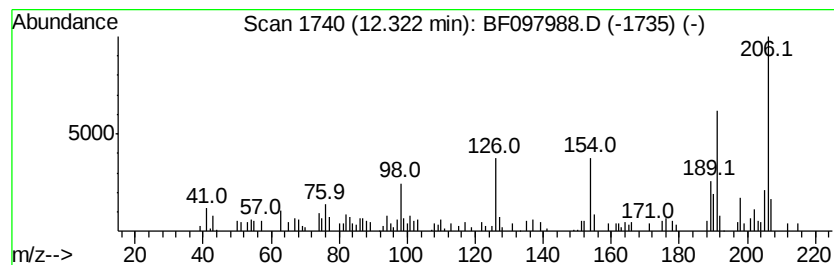
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 9,10-Dimethylantracene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.32	2.25 ng	93053	Phenanthrene-d10	11.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9,10-Dimethylantracene	206	C16H14	000781-43-1	91
2		3-Methyl-1-phenyl-1H-indene	206	C16H14	022360-63-0	64
3		1-Methyl-3-phenyl-1H-indene	206	C16H14	022360-62-9	64
4		Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	64
5		1H-Indene, 2-methyl-3-phenyl-	206	C16H14	035099-60-6	58



Data Path : U:\HPCHEM1\BNA F\DATA\BF082217\
Data File : BF097988.D
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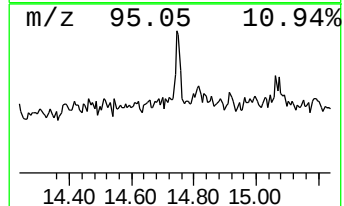
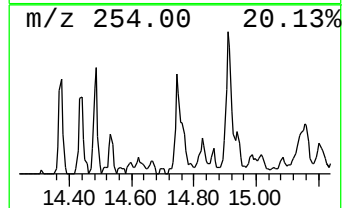
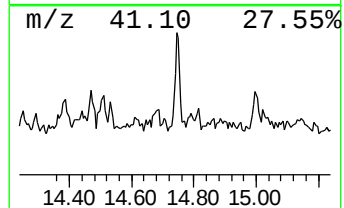
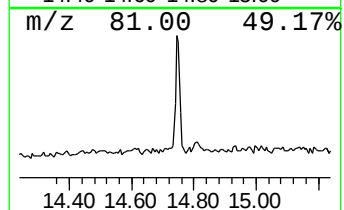
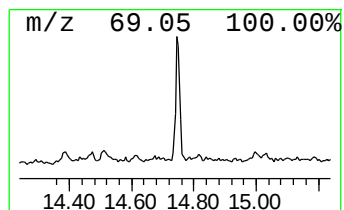
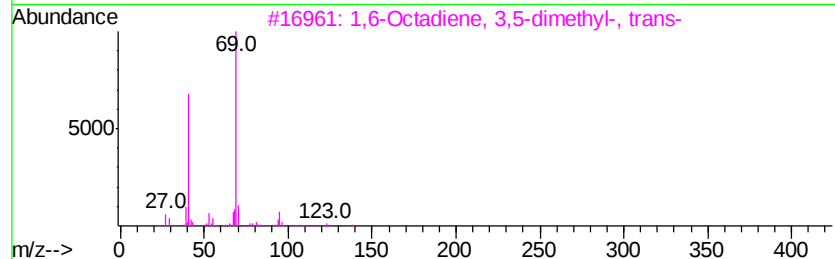
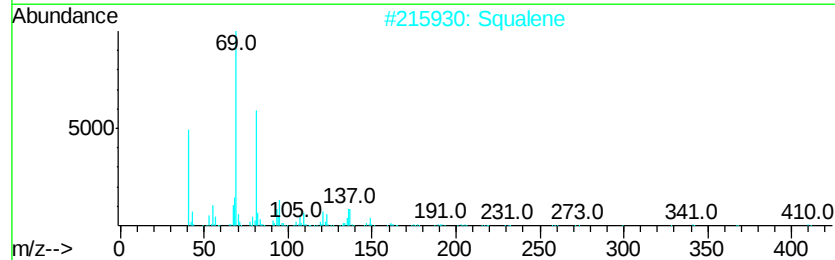
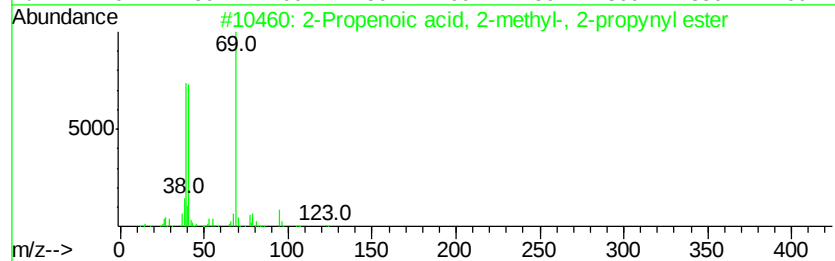
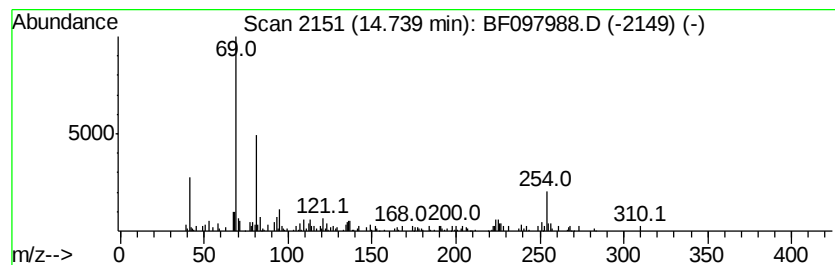
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 2-Propenoic acid, 2-methyl-... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.74	5.75 ng	158630	Perylene-d12	15.32

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Propenoic acid, 2-methyl-, 2-p...	124	C7H8O2	013861-22-8	58
2		Squalene	410	C30H50	000111-02-4	58
3		1,6-Octadiene, 3,5-dimethyl-, tr...	138	C10H18	074630-87-8	53
4		Trideca-1,3,7,11-tetraene-1,1-di...	268	C18H24N2	1000159-88-9	50
5		1,5-Heptadiene, 3,4-dimethyl-	124	C9H16	1000061-77-9	50



Data Path : U:\HPCHEM1\BNA_F\DATA\BF082217\
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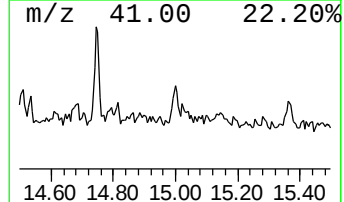
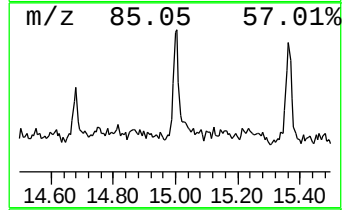
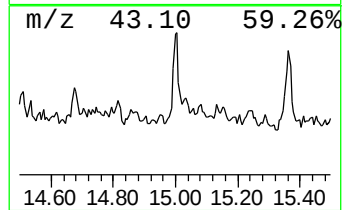
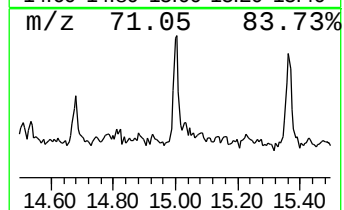
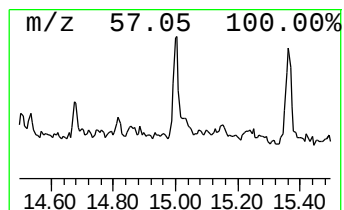
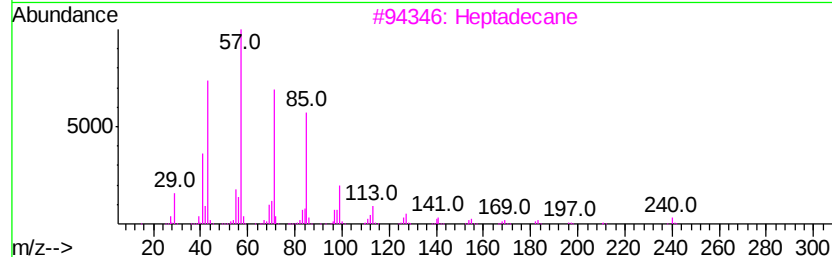
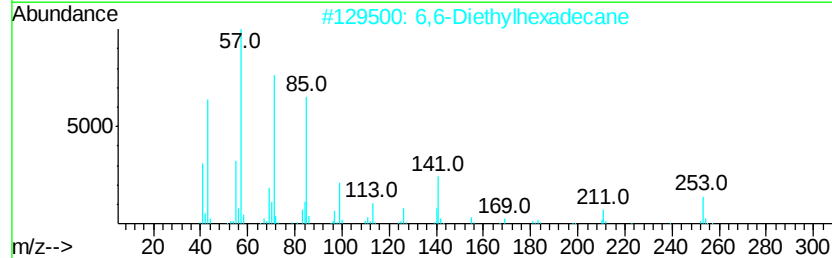
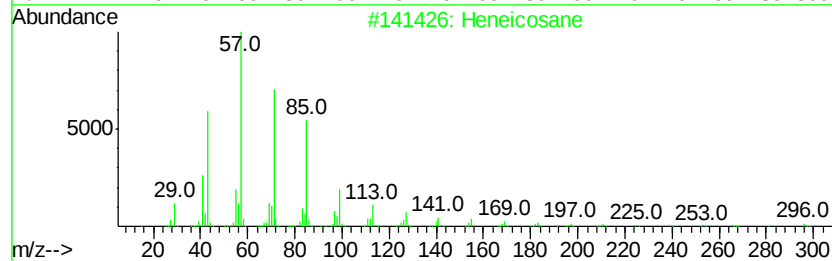
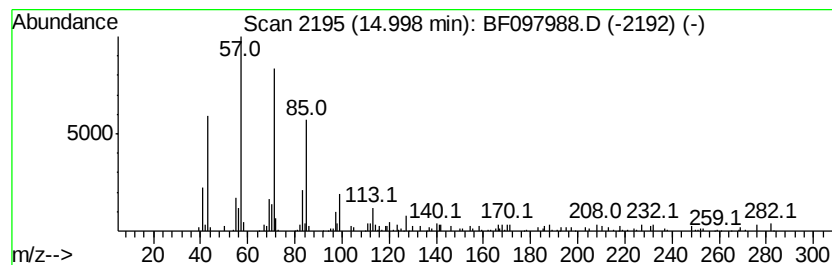
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Heneicosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.00	2.84 ng	78317	Perylene-d12	15.32
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Heneicosane	296 C21H44	000629-94-7	80
2	6,6-Diethylhexadecane	282 C20H42	1000360-41-4	80
3	Heptadecane	240 C17H36	000629-78-7	80
4	Dodecane	170 C12H26	000112-40-3	74
5	Heptadecane, 2,6,10,15-tetramethyl-	296 C21H44	054833-48-6	72



Data Path : U:\HPCHEM1\BNA F\DATA\BF082217\
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Operator : SJ/JU
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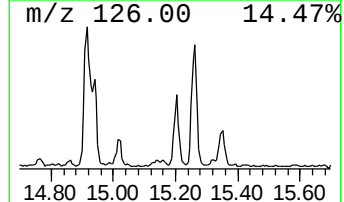
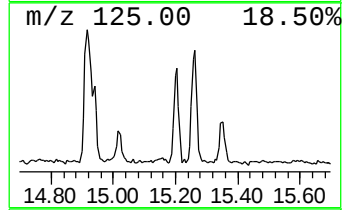
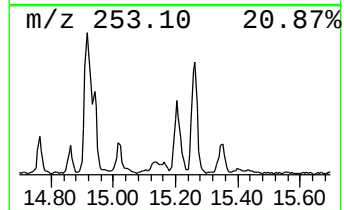
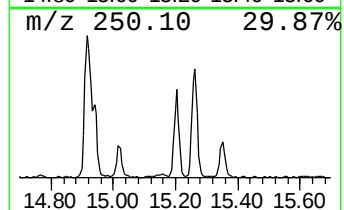
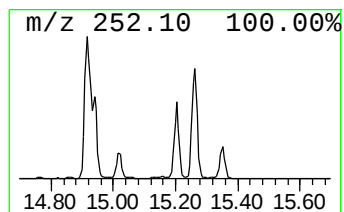
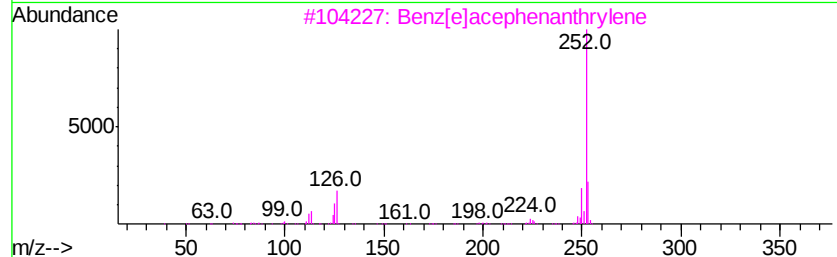
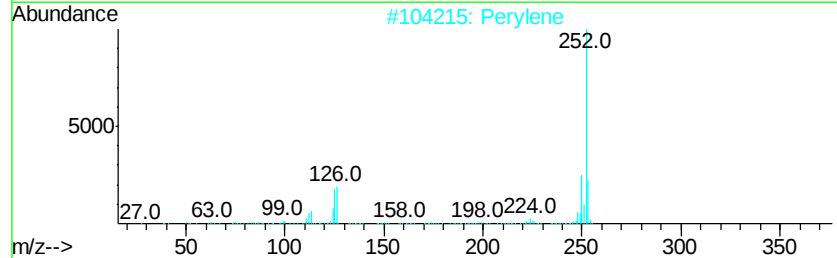
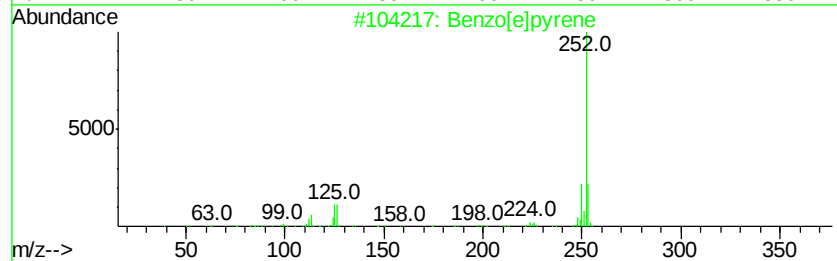
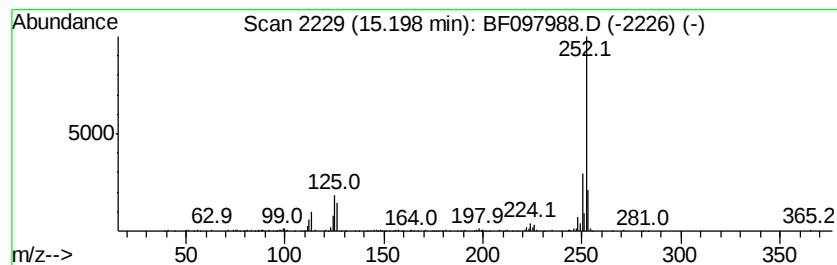
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Benzo[e]pyrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.20	8.30 ng	229214	Perylene-d12	15.32

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	98
2			Perylene	252	C20H12	000198-55-0	98
3			Benzo[e]acephenanthrylene	252	C20H12	000205-99-2	97
4			Benzo[k]fluoranthene	252	C20H12	000207-08-9	93
5			Benzo[a]pyrene	252	C20H12	000050-32-8	91



Data Path : U:\HPCHEM1\BNA_F\DATA\BF082217\
Data File : BF097988.D
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Operator : SJ/JU
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Instrument :
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ClientSampleId :
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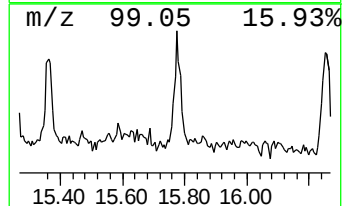
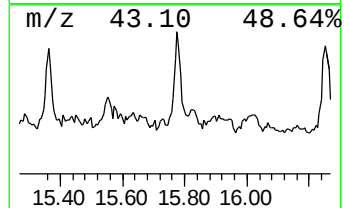
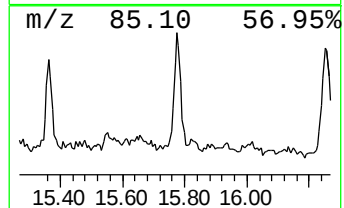
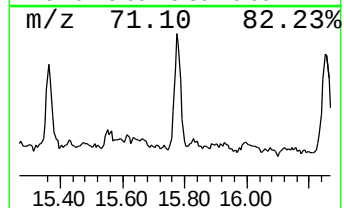
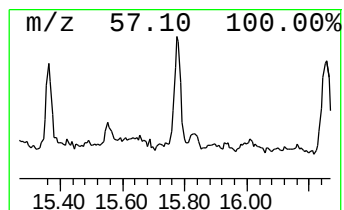
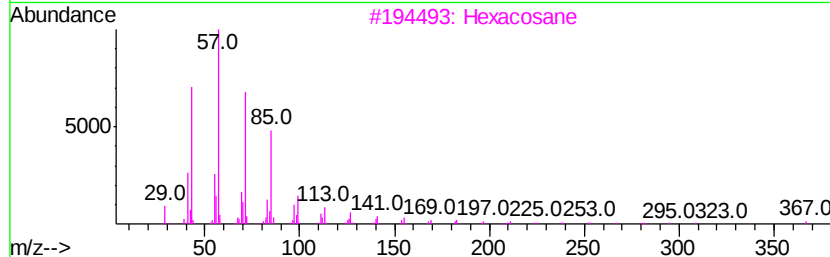
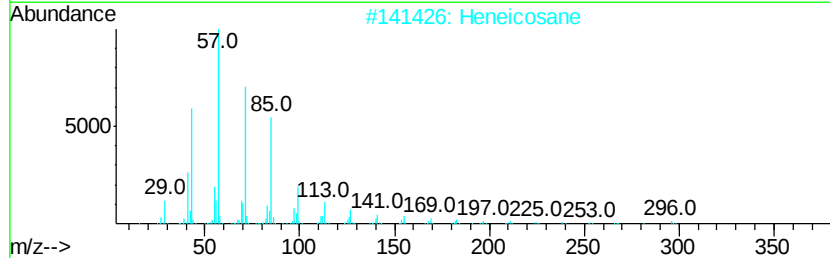
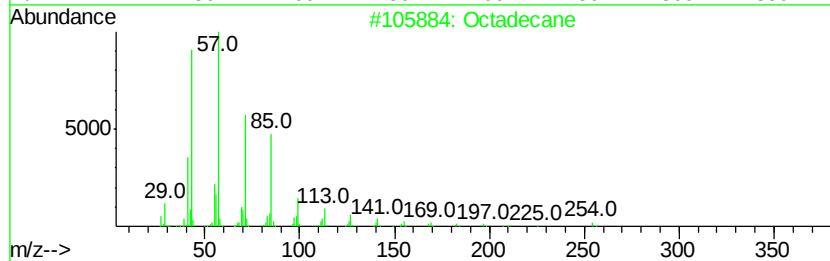
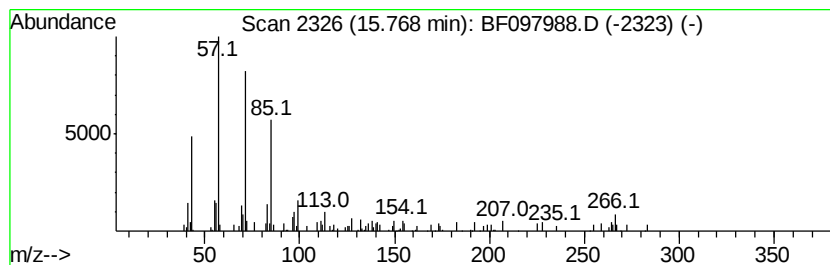
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Octadecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.77	4.68 ng	129261	Perylene-d12	15.32

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecane	254	C18H38	000593-45-3	93
2			Heneicosane	296	C21H44	000629-94-7	87
3			Hexacosane	366	C26H54	000630-01-3	86
4			Octacosane	394	C28H58	000630-02-4	83
5			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	80



Data Path : U:\HPCHEM1\BNA_F\DATA\BF082217\
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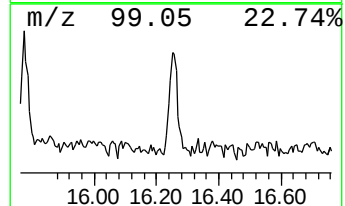
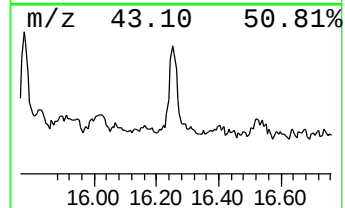
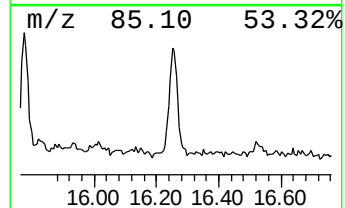
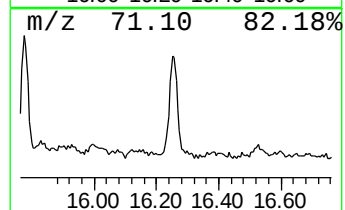
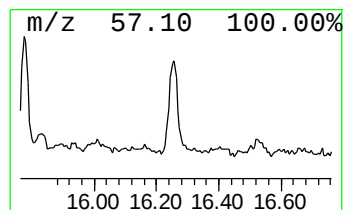
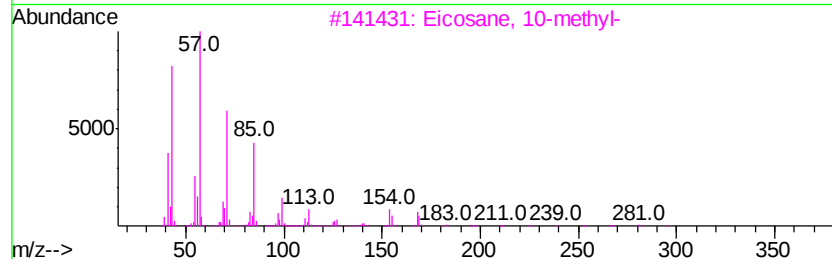
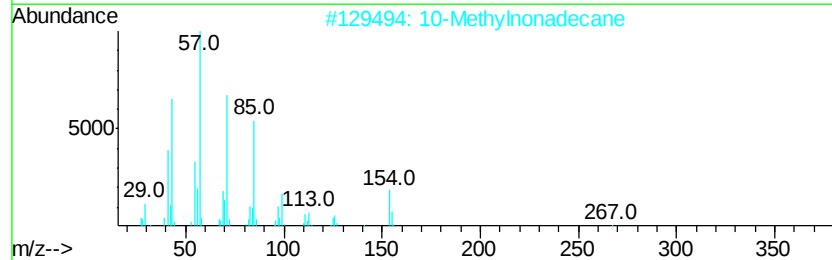
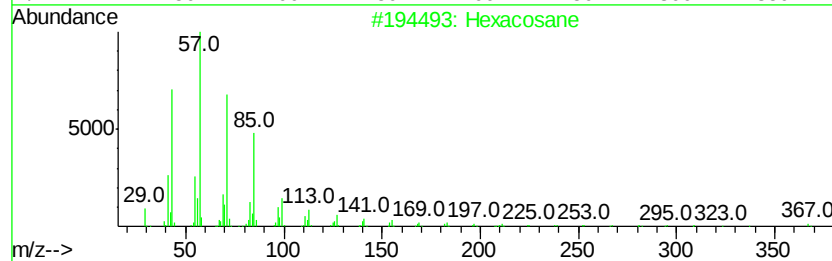
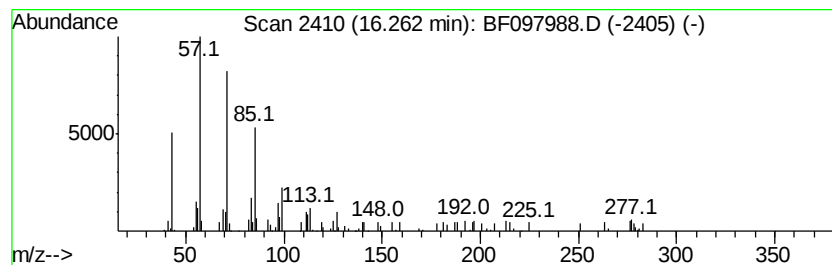
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Hexacosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.26	2.73 ng	75359	Perylene-d12	15.32
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Hexacosane	366 C26H54	000630-01-3	87
2	10-Methylnonadecane	282 C20H42	056862-62-5	86
3	Eicosane, 10-methyl-	296 C21H44	054833-23-7	86
4	Heptadecane	240 C17H36	000629-78-7	80
5	Eicosane, 9-octyl-	394 C28H58	013475-77-9	80



Data Path : U:\HPCHEM1\BNA_F\DATA\BF082217\
Data File : BF097988.D
Acq On : 23 Aug 2017 7:57
Operator : SJ/JU
Sample : I4913-01
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ALS Vial : 30 Sample Multiplier: 1

Instrument :
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ClientSampleId :
TEC-BE2-1-3-COMP

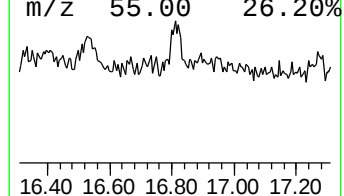
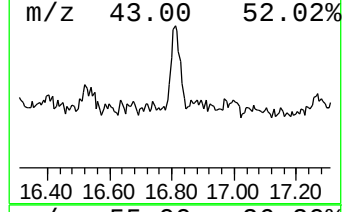
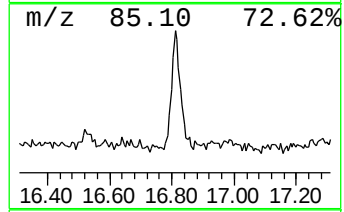
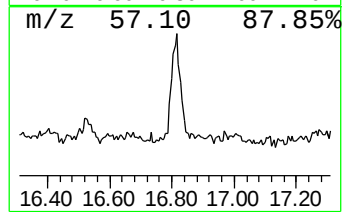
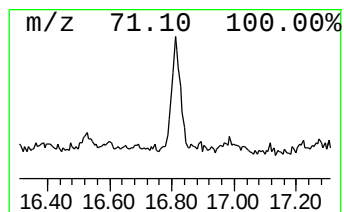
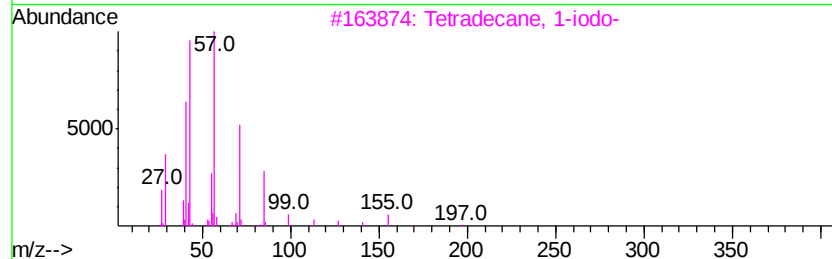
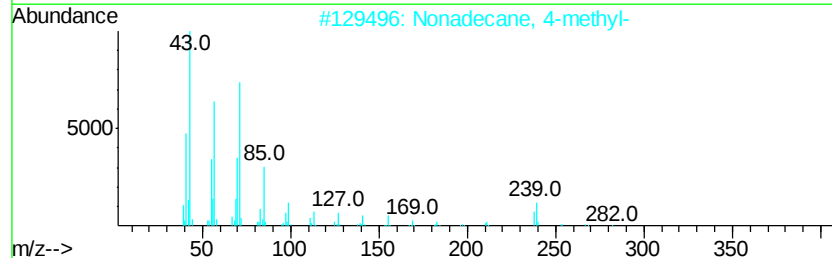
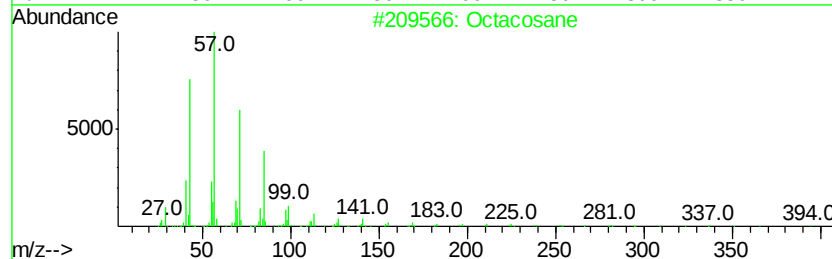
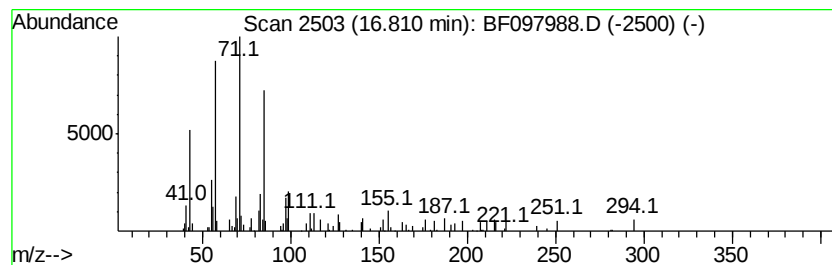
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Octacosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.81	2.33 ng	64241	Perylene-d12	15.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	72
2		Nonadecane, 4-methyl-	282	C20H42	025117-27-5	64
3		Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	53
4		Nonadecane, 9-methyl-	282	C20H42	013287-24-6	53
5		3-Ethyl-3-methylheptane	142	C10H22	017302-01-1	50



Data Path : U:\HPCHEM1\BNA_F\DATA\BF082217\
Data File : BF097988.D
Acq On : 23 Aug 2017 7:57
Operator : SJ/JU
Sample : I4913-01
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TEC-BE2-1-3-COMP

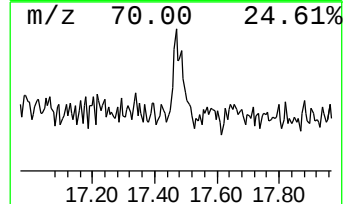
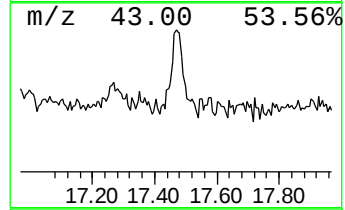
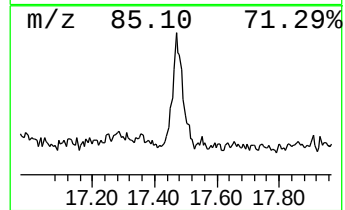
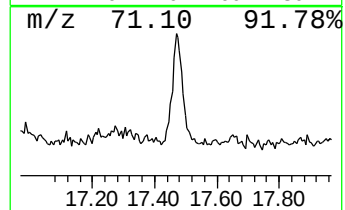
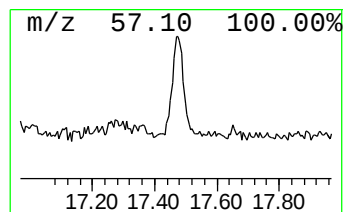
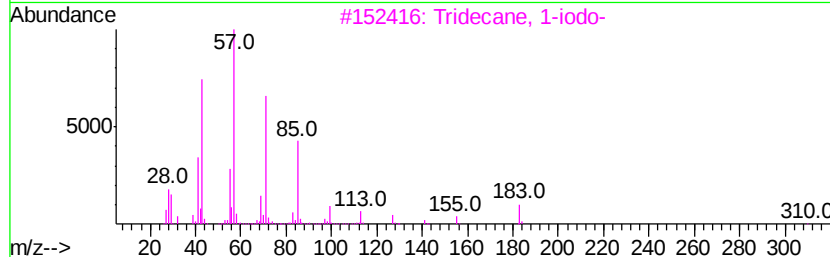
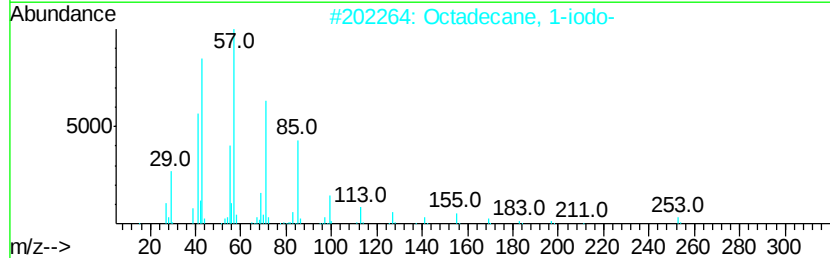
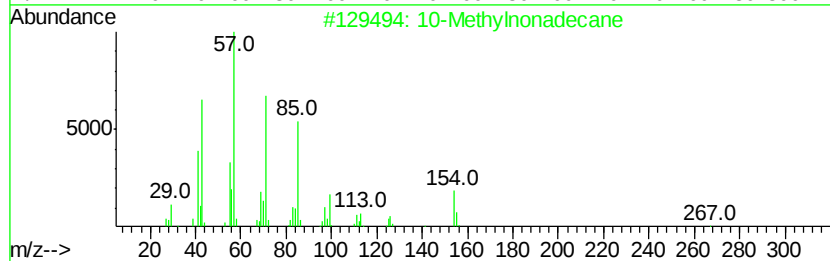
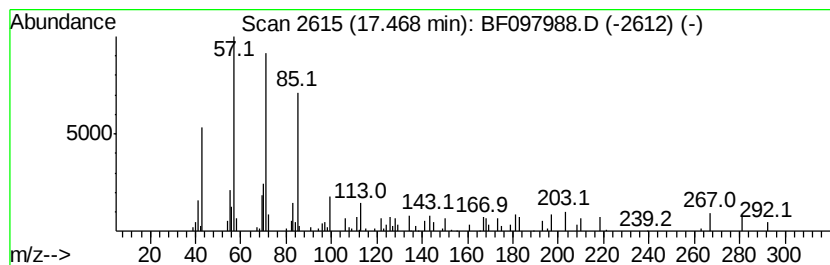
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 10-Methylnonadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.47	3.80 ng	104926	Perylene-d12	15.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	10-Methylnonadecane	282	C20H42	056862-62-5	59
2		Octadecane, 1-iodo-	380	C18H37I	000629-93-6	59
3		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	59
4		Tetradecane	198	C14H30	000629-59-4	59
5		Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	53



Data Path : U:\HPCHEM1\BNA_F\DATA\BF082217\
 Data File : BF097988.D
 Acq On : 23 Aug 2017 7:57
 Operator : SJ/JU
 Sample : I4913-01
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TEC-BE2-1-3-COMP

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown3.13	3.13	2.5	ng	109096	1	6.75	863153	20.0
2-Pentanone, 4-hy...	4.95	6.9	ng	297185	1	6.75	863153	20.0
unknown6.52	6.52	80.2	ng	3461010	1	6.75	863153	20.0
Phthalic anhydride	8.80	3.0	ng	156373	2	8.03	1046990	20.0
Heptadecane, 2,6,...	10.69	2.2	ng	91560	4	11.26	826921	20.0
9,10-Dimethylanthe...	12.32	2.3	ng	93053	4	11.26	826921	20.0
2-Propenoic acid,...	14.74	5.8	ng	158630	6	15.32	552102	20.0
Heneicosane	15.00	2.8	ng	78317	6	15.32	552102	20.0
Benzo[e]pyrene	15.20	8.3	ng	229214	6	15.32	552102	20.0
Octadecane	15.77	4.7	ng	129261	6	15.32	552102	20.0
Hexacosane	16.26	2.7	ng	75359	6	15.32	552102	20.0
Octacosane	16.81	2.3	ng	64241	6	15.32	552102	20.0
10-Methylnonadecane	17.47	3.8	ng	104926	6	15.32	552102	20.0