

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110819\
 Data File : BF117733.D
 Acq On : 8 Nov 2019 16:16
 Operator : JU
 Sample : K5722-13
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 1-C

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:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110619.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.228	17	24	30	rVB	1205109	1521069	19.69%	1.765%
2	5.145	515	520	528	rVB	1182978	1349807	17.47%	1.566%
3	5.540	582	587	590	rBV	5815788	5861704	75.88%	6.800%
4	6.545	752	758	761	rBV	6183195	6255086	80.97%	7.256%
5	6.692	779	783	787	rBV	6354377	6394824	82.78%	7.418%
6	6.928	819	823	826	rBV	2672234	2089476	27.05%	2.424%
7	7.087	846	850	853	rVB	5974011	5045537	65.31%	5.853%
8	7.487	913	918	921	rBV	4428977	4002993	51.82%	4.644%
9	8.216	1036	1042	1045	rBV	3332289	2832713	36.67%	3.286%
10	9.292	1220	1225	1228	rBV	8542861	7724944	100.00%	8.961%
11	9.680	1289	1291	1296	rVB	297985	255552	3.31%	0.296%
12	9.833	1314	1317	1321	rBV	131362	109064	1.41%	0.127%
13	9.975	1334	1341	1344	rBV	3944235	3418429	44.25%	3.966%
14	10.004	1344	1346	1350	rVB	147523	136963	1.77%	0.159%
15	10.175	1372	1375	1379	rVB2	91563	102266	1.32%	0.119%
16	10.522	1431	1434	1440	rBV2	143001	167335	2.17%	0.194%
17	10.763	1471	1475	1479	rBV	5167623	5204562	67.37%	6.038%
18	10.892	1493	1497	1501	rBV3	146886	218965	2.83%	0.254%
19	11.104	1530	1533	1535	rBV2	88414	115548	1.50%	0.134%
20	11.269	1559	1561	1564	rBV	143793	120682	1.56%	0.140%
21	11.363	1575	1577	1579	rBV	147917	103882	1.34%	0.121%
22	11.469	1591	1595	1597	rBV	3923333	3438165	44.51%	3.989%
23	11.492	1597	1599	1602	rVV	2254537	1625625	21.04%	1.886%
24	11.539	1605	1607	1610	rVV	251439	229253	2.97%	0.266%
25	11.698	1631	1634	1637	rBV2	303688	343633	4.45%	0.399%
26	11.969	1678	1680	1682	rBV	398409	376710	4.88%	0.437%
27	11.998	1682	1685	1688	rVB	1519141	1291340	16.72%	1.498%
28	12.080	1697	1699	1702	rBV2	430425	425787	5.51%	0.494%
29	12.686	1799	1802	1808	rVB	2621769	2250211	29.13%	2.610%
30	12.916	1839	1841	1846	rVB	1988836	1887636	24.44%	2.190%
31	13.063	1862	1866	1869	rBV	6836570	6596886	85.40%	7.653%
32	13.957	2015	2018	2030	rVB2	1365052	1667936	21.59%	1.935%
33	14.115	2040	2045	2048	rBV2	2703862	3513120	45.48%	4.075%
34	14.145	2048	2050	2058	rVB	1040923	949091	12.29%	1.101%

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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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35	14.586	2123	2125	2130	rVB3	349915	429830	5.56%	0.499%
36	14.904	2177	2179	2184	rVB	371235	357450	4.63%	0.415%
37	14.986	2190	2193	2197	rBV	635501	653140	8.45%	0.758%
38	15.174	2221	2225	2229	rBV2	668054	1112091	14.40%	1.290%
39	15.262	2237	2240	2243	rBV	325534	351489	4.55%	0.408%
40	15.492	2276	2279	2285	rVB	436584	516938	6.69%	0.600%
41	15.557	2286	2290	2294	rVV	379650	466277	6.04%	0.541%
42	15.621	2297	2301	2305	rVV	1711761	2274136	29.44%	2.638%
43	15.662	2305	2308	2315	rVB3	460663	641817	8.31%	0.745%
44	16.121	2382	2386	2392	rBV	328178	520913	6.74%	0.604%
45	16.662	2473	2478	2482	rBV2	187766	345407	4.47%	0.401%
46	17.145	2555	2560	2568	rVB2	203734	418612	5.42%	0.486%
47	17.292	2581	2585	2591	rBV3	74680	148280	1.92%	0.172%
48	17.604	2634	2638	2647	rVB2	164688	338668	4.38%	0.393%

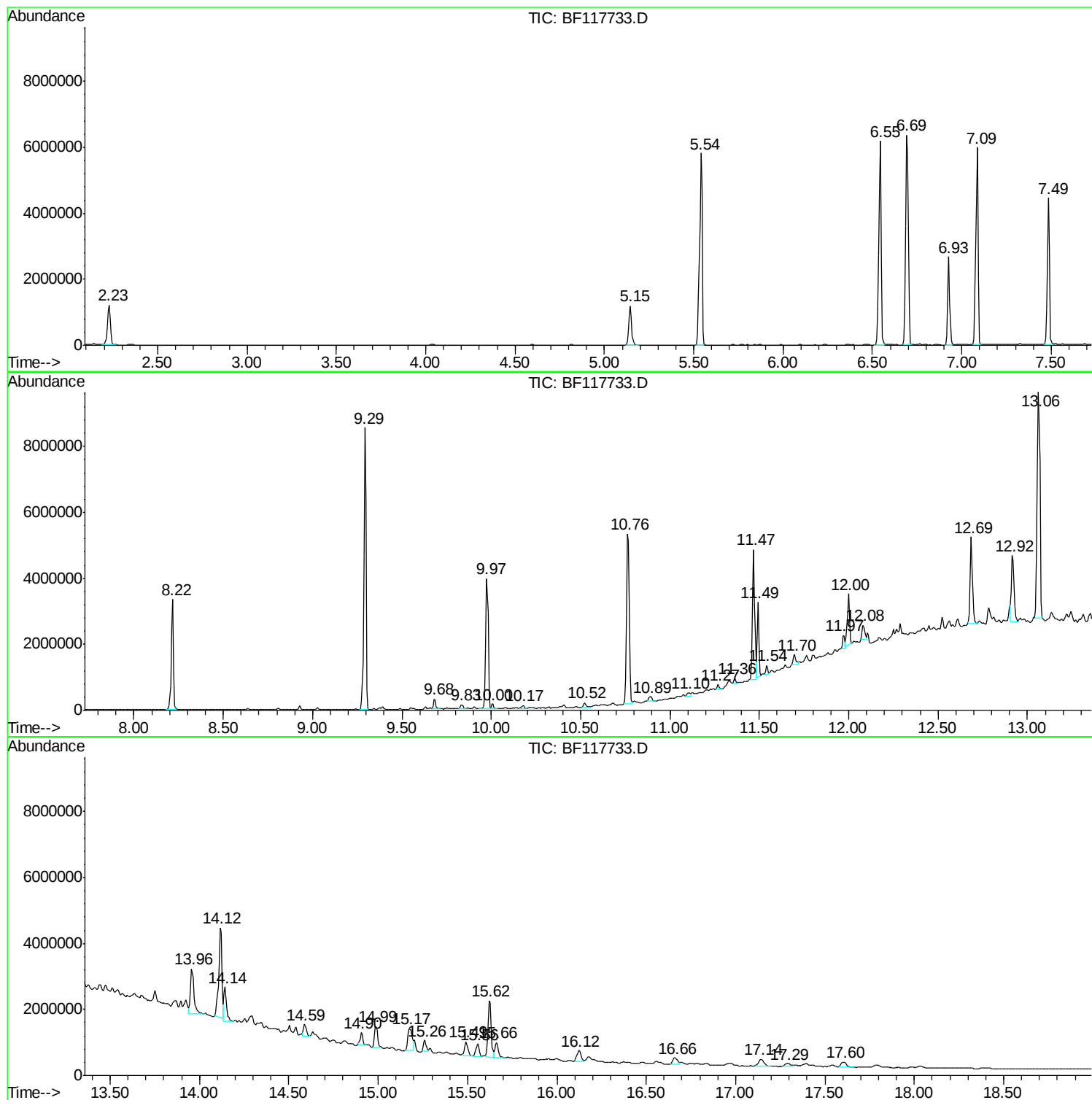
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ClientSampled :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Sample : K5722-13
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Instrument :
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ClientSampleId :
1-C

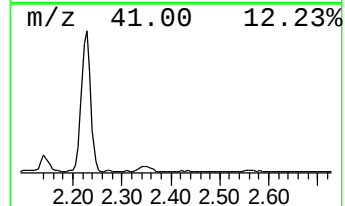
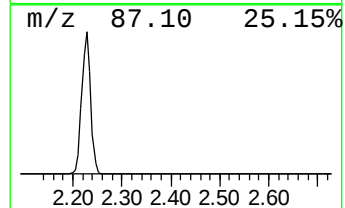
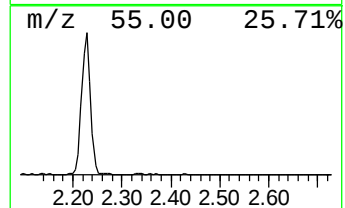
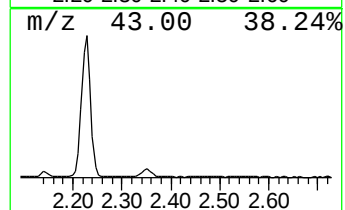
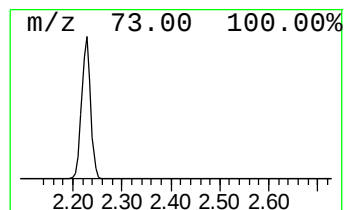
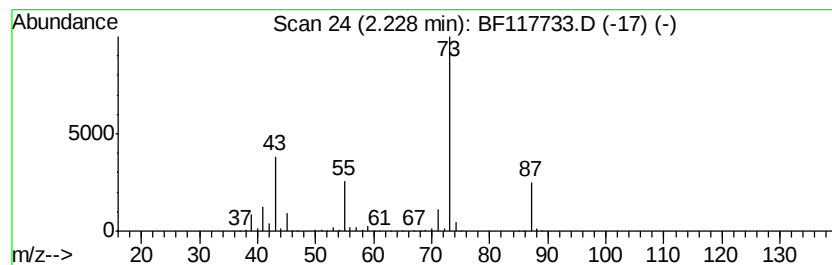
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.23	14.56 ng	1521070	1,4-Dichlorobenzene-d4	6.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12
5		Silane, tetramethyl-	88	C4H12Si	000075-76-3	10



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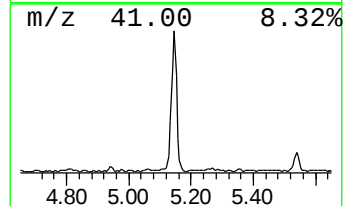
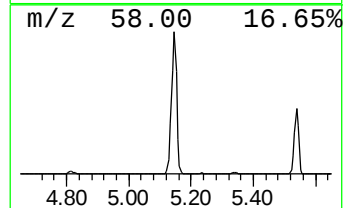
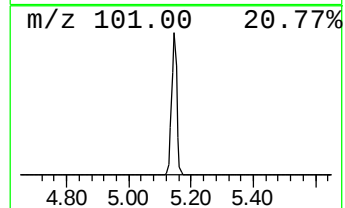
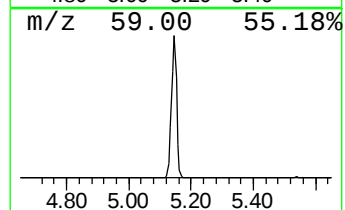
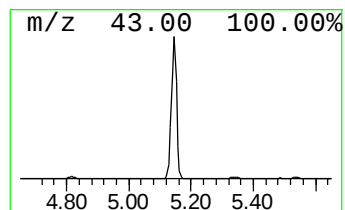
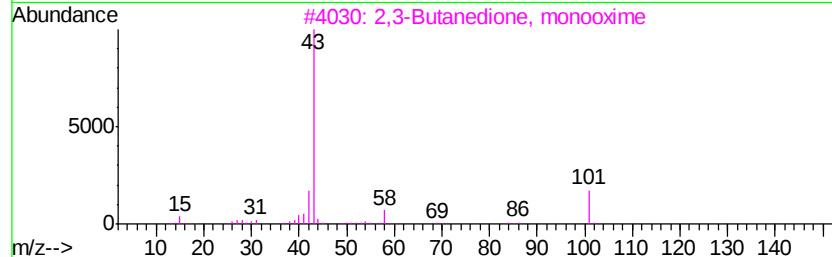
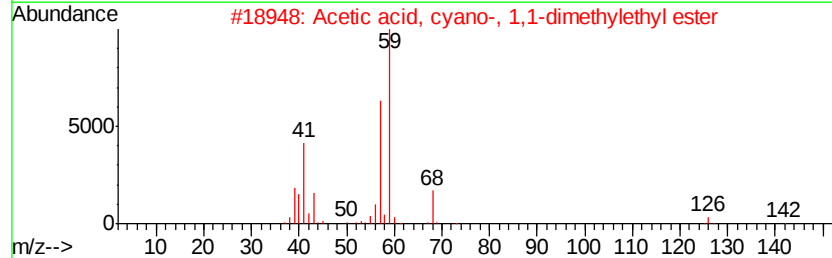
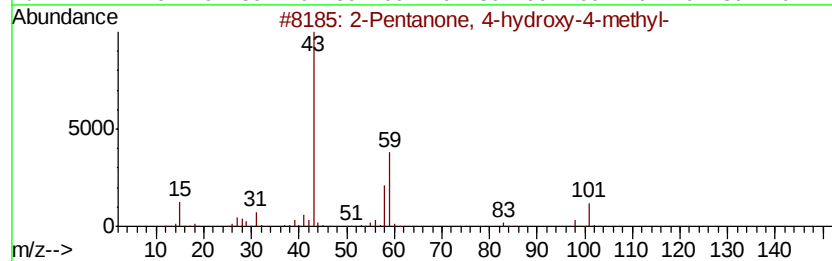
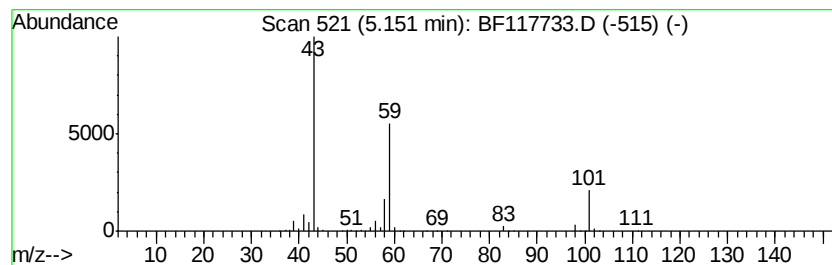
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.15	12.92 ng	1349810	1,4-Dichlorobenzene-d4	6.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
4		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
5		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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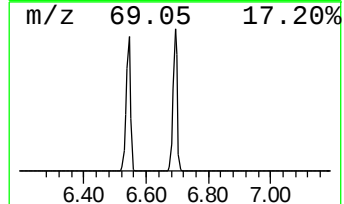
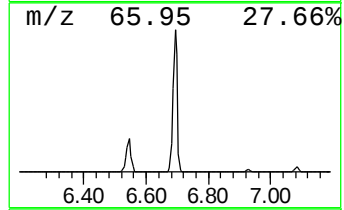
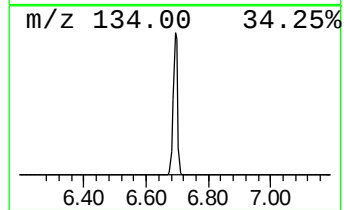
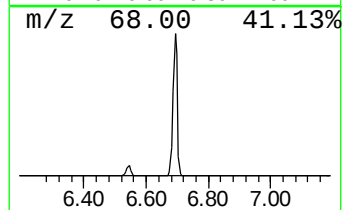
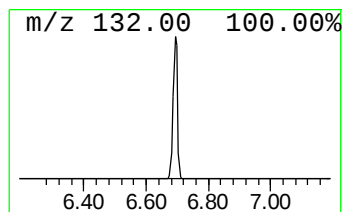
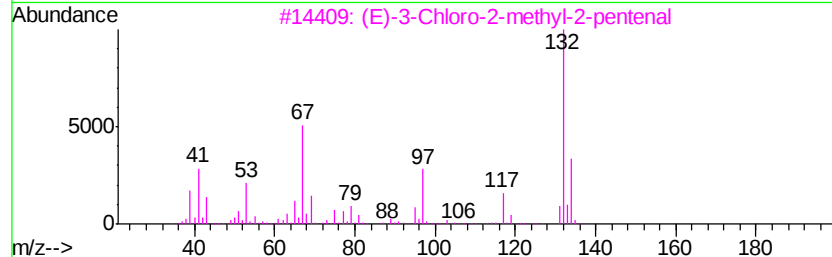
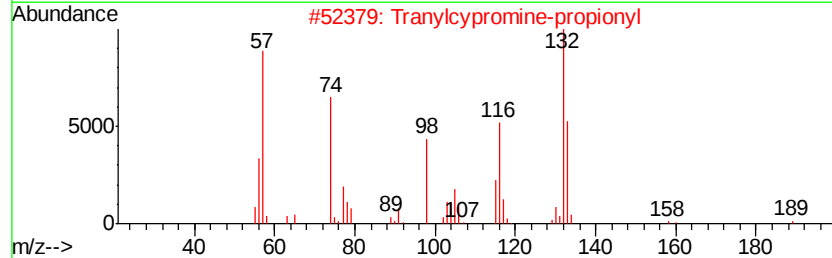
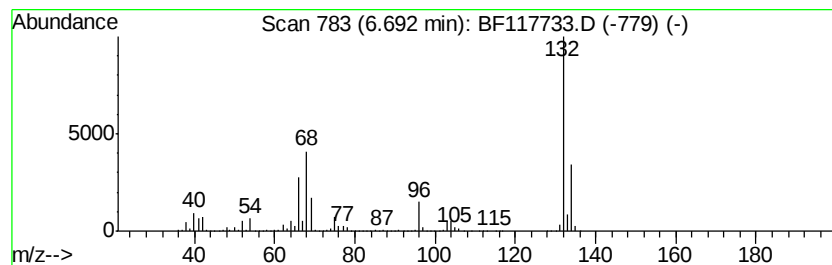
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 unknown6.69 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.69	61.21 ng	6394820	1,4-Dichlorobenzene-d4	6.93

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
3			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
4			1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
5			Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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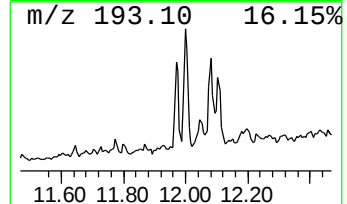
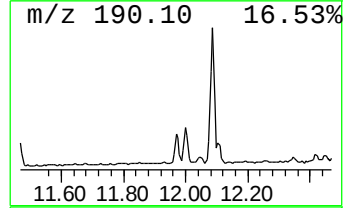
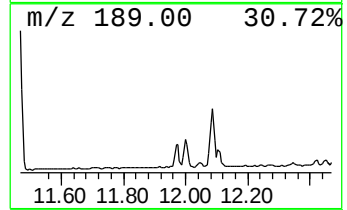
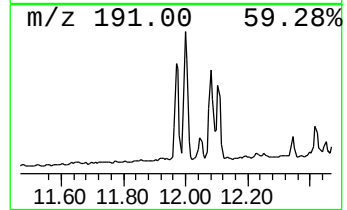
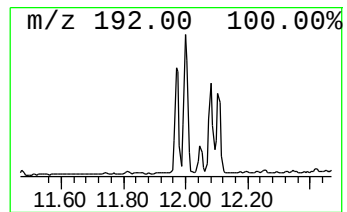
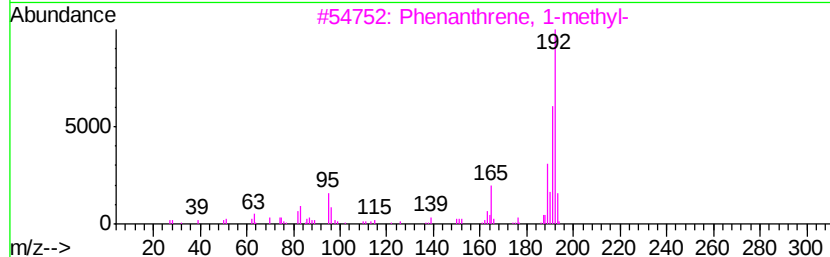
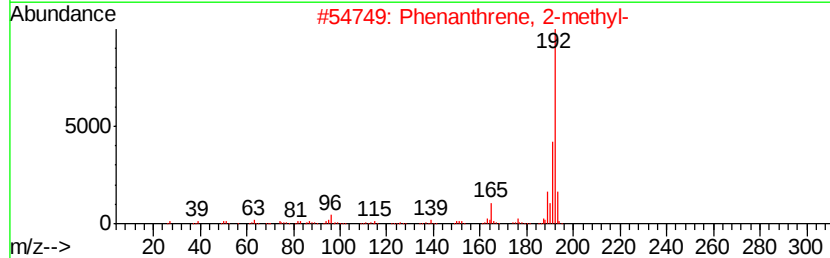
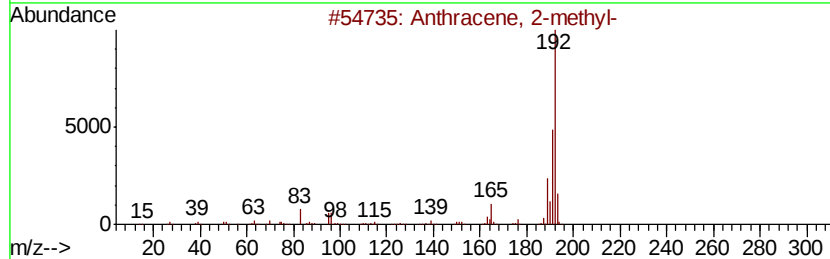
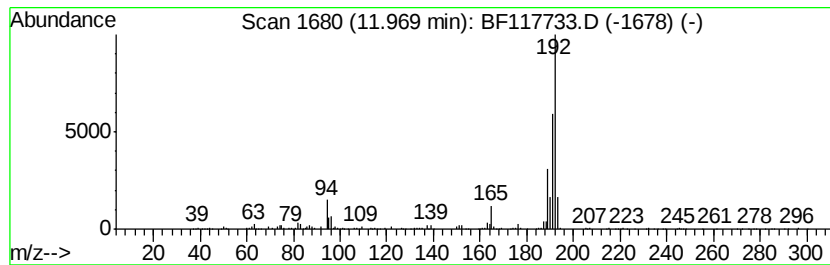
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Anthracene, 2-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.97	2.19 ng	376710	Phenanthrene-d10	11.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
2		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
3		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
4		1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	93
5		Anthracene, 9-methyl-	192	C15H12	000779-02-2	93



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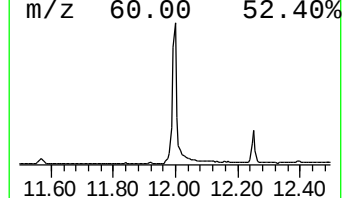
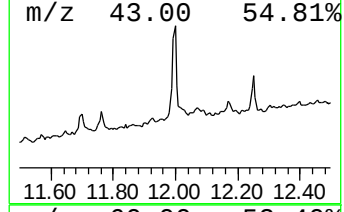
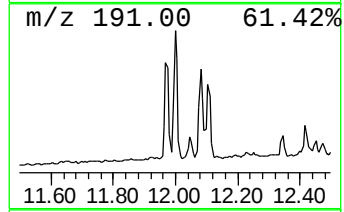
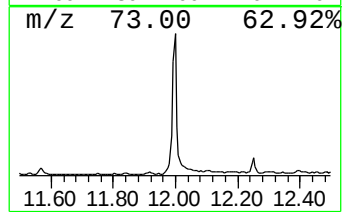
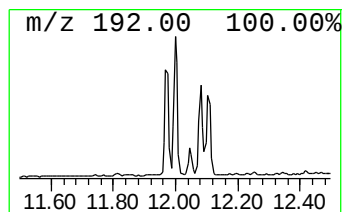
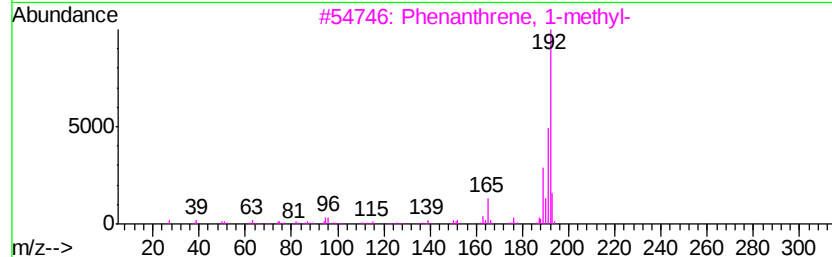
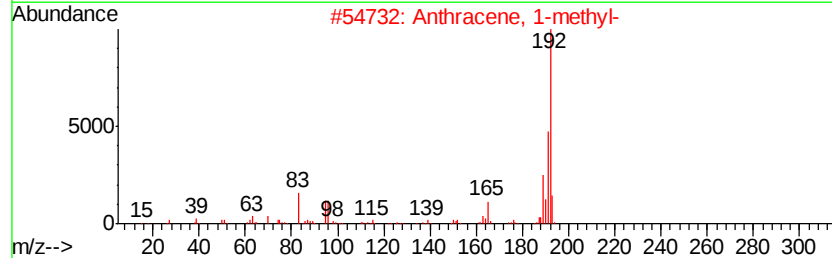
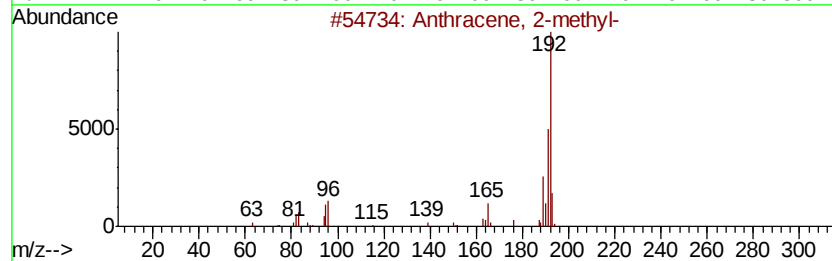
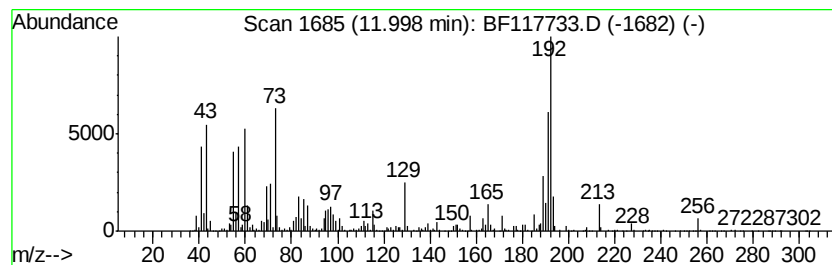
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Anthracene, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.00	7.51 ng	1291340	Phenanthrene-d10	11.47

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
2			Anthracene, 1-methyl-	192	C15H12	000610-48-0	95
3			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	92
4			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	92
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91



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Data File : BF117733.D
Acq On : 8 Nov 2019 16:16
Operator : JU
Sample : K5722-13
Misc :
ALS Vial : 6 Sample Multiplier: 1

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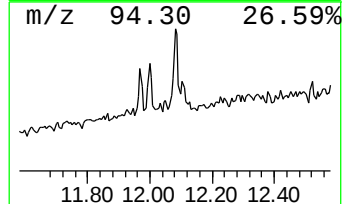
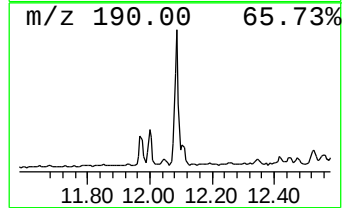
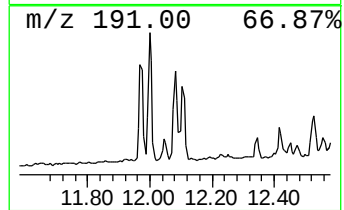
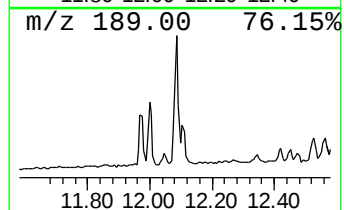
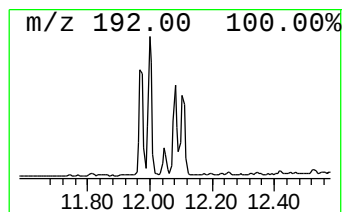
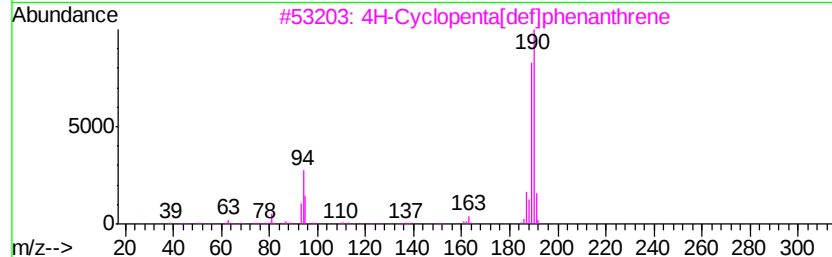
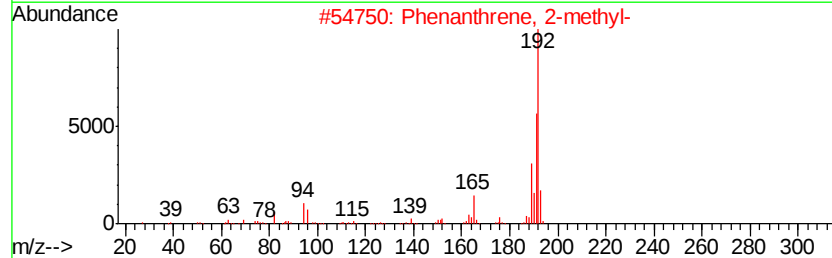
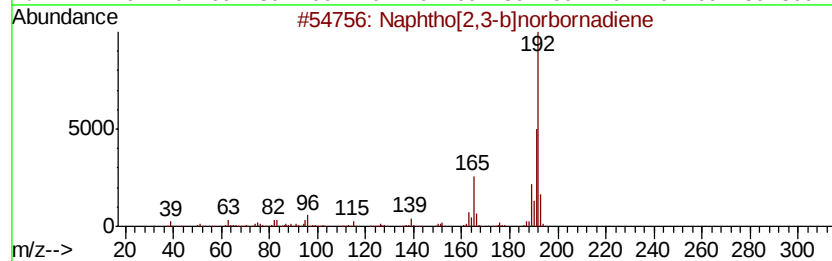
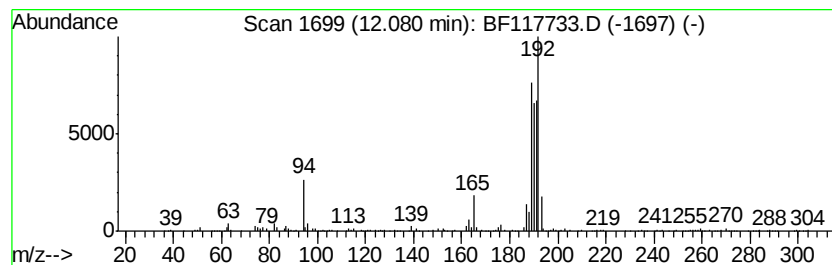
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Naphtho[2,3-b]norbornadiene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.08	2.48 ng	425787	Phenanthrene-d10	11.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphtho[2,3-b]norbornadiene	192	C15H12	107426-38-0	50
2		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	49
3		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	46
4		6H-Cyclobuta[fik]phenanthrene	190	C15H10	083469-43-6	46
5		Anthracene, 2-methyl-	192	C15H12	000613-12-7	45



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110819\
Data File : BF117733.D
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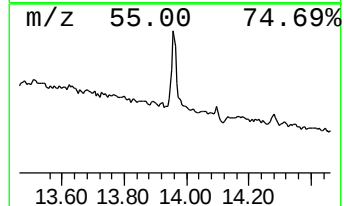
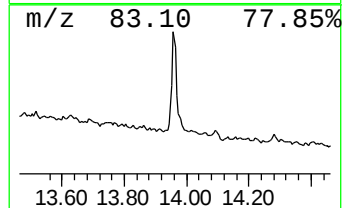
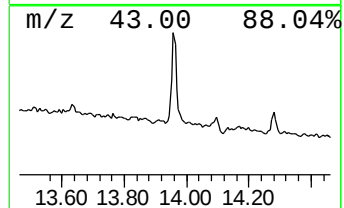
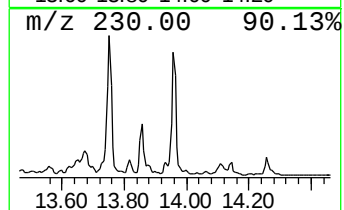
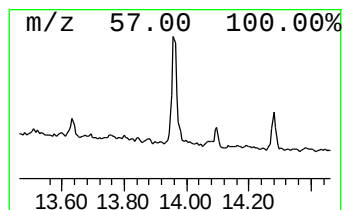
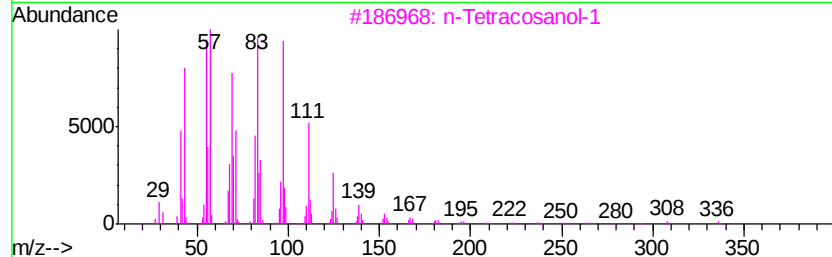
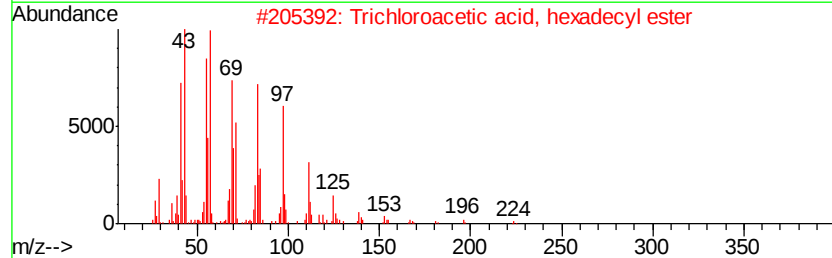
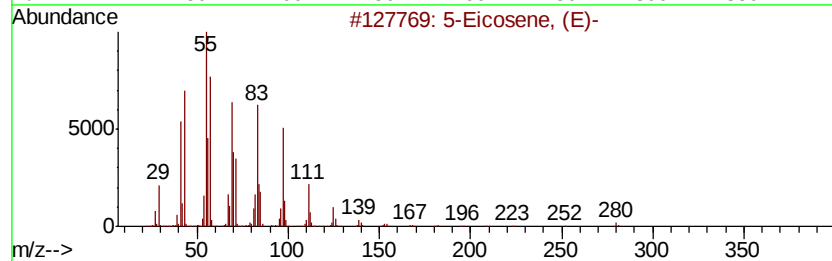
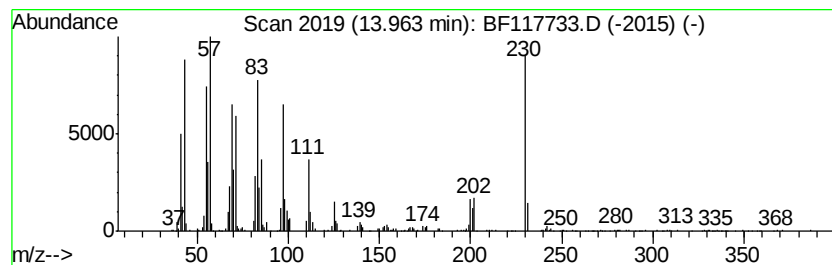
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 5-Eicosene, (E)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.96	9.50 ng	1667940	Chrysene-d12	14.12
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	5-Eicosene, (E)-		280 C20H40	074685-30-6 96
2	Trichloroacetic acid, hexadecyl ...		386 C18H33Cl3O2	074339-54-1 86
3	n-Tetracosanol-1		354 C24H50O	000506-51-4 86
4	1-Heneicosanol		312 C21H44O	015594-90-8 83
5	Nonadecyl trifluoroacetate		380 C21H39F3O2	1000351-76-3 70



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110819\
Data File : BF117733.D
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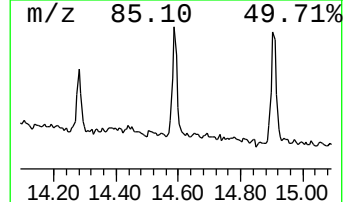
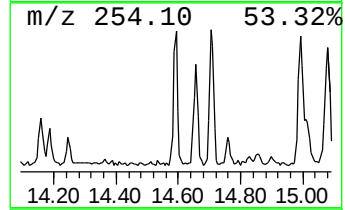
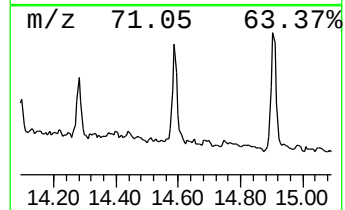
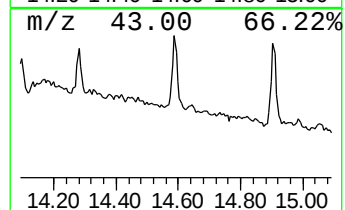
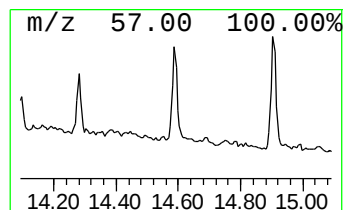
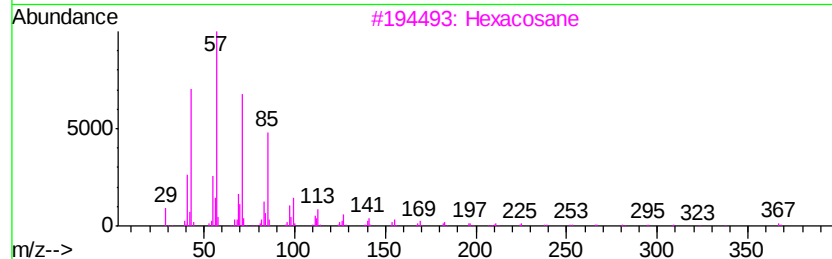
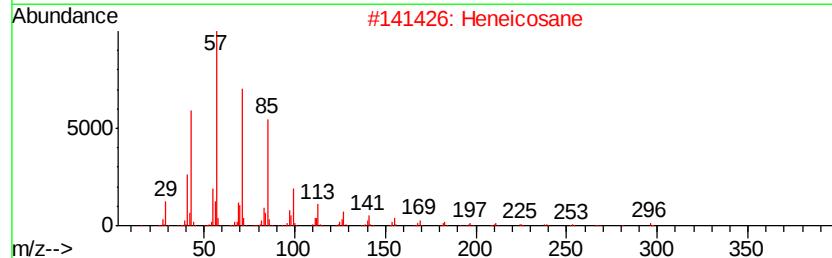
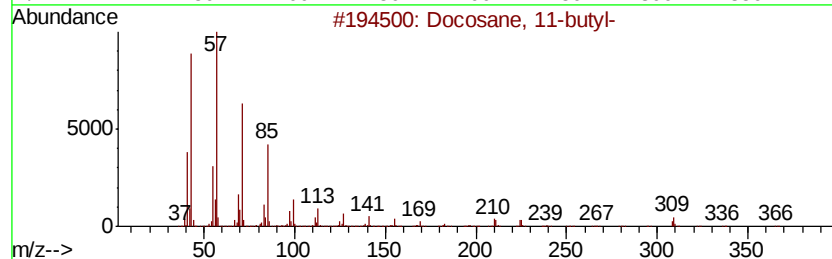
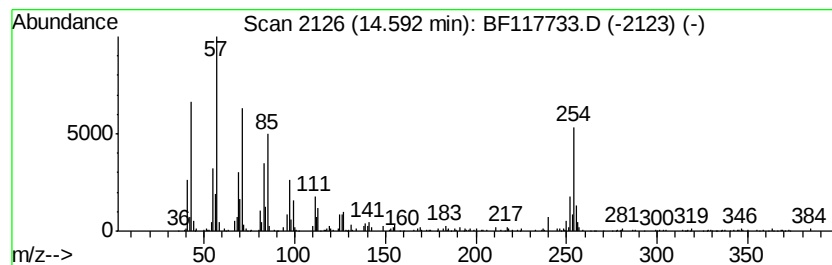
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Docosane, 11-butyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.59	2.45 ng	429830	Chrysene-d12	14.12
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Docosane, 11-butyl-	366 C26H54	013475-76-8 90
2		Heneicosane	296 C21H44	000629-94-7 81
3		Hexacosane	366 C26H54	000630-01-3 81
4		Hexatriacontane	507 C36H74	000630-06-8 74
5		Tetracosane	338 C24H50	000646-31-1 74



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110819\
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Misc :
ALS Vial : 6 Sample Multiplier: 1

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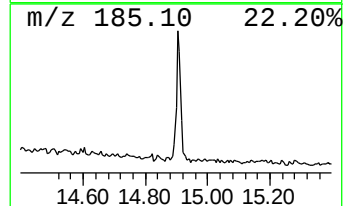
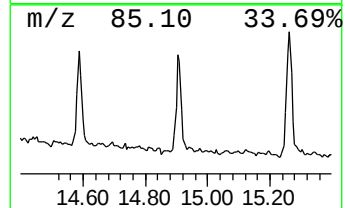
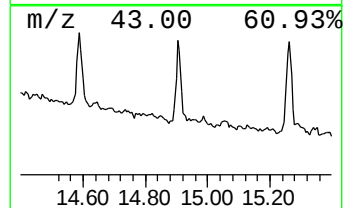
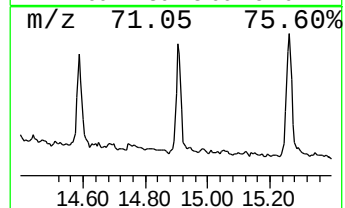
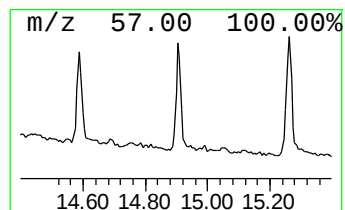
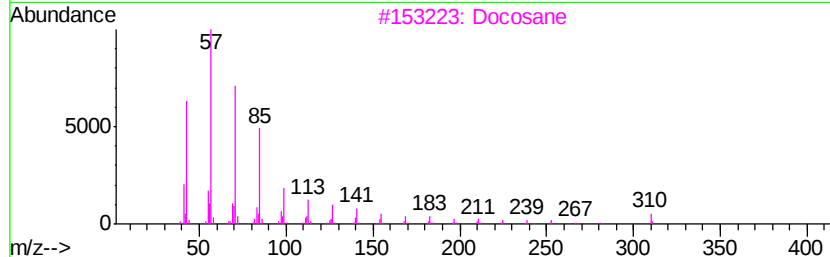
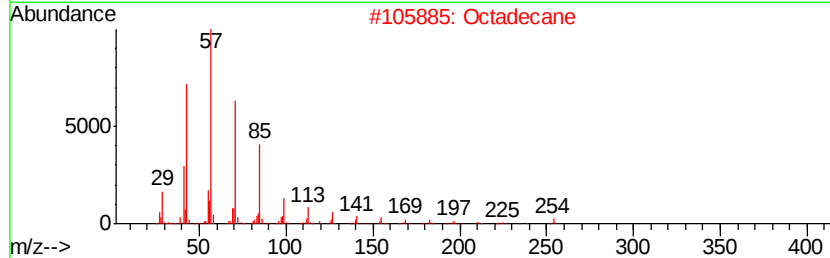
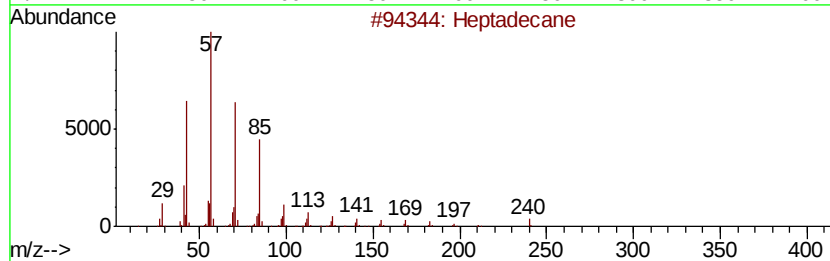
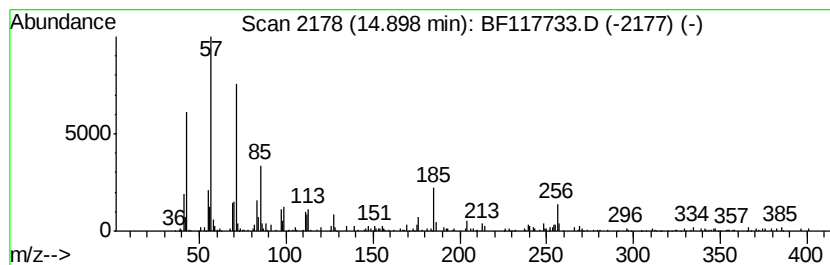
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Heptadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.90	3.14 ng	357450	Perylene-d12	15.63

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane		240	C17H36	000629-78-7	92
2	Octadecane		254	C18H38	000593-45-3	91
3	Docosane		310	C22H46	000629-97-0	89
4	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	86
5	Octacosane		394	C28H58	000630-02-4	76



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110819\
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ClientSampleId :
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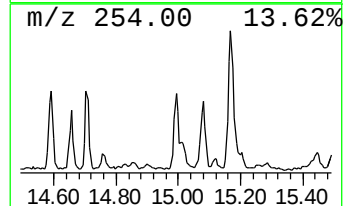
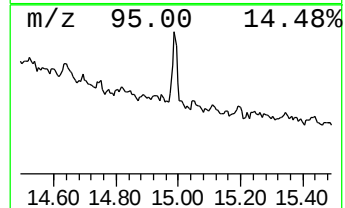
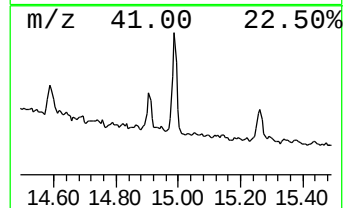
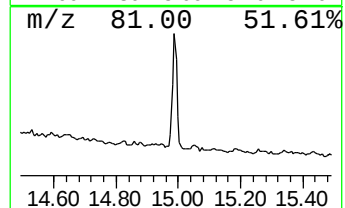
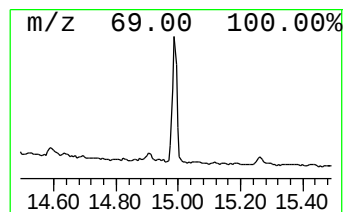
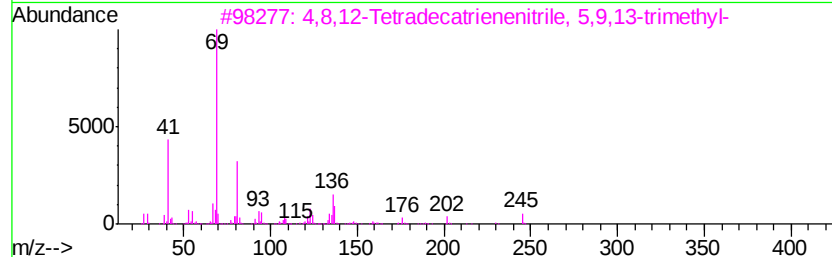
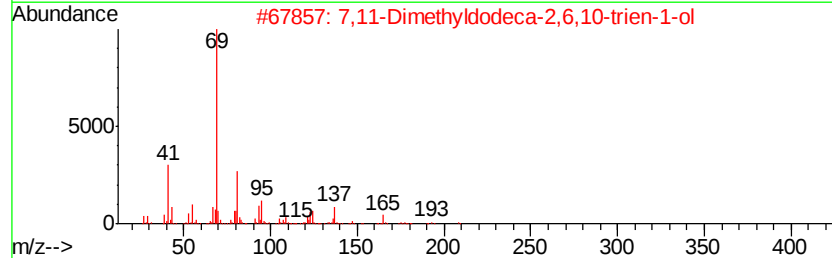
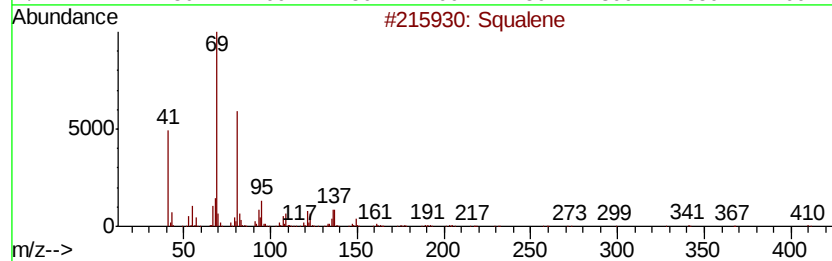
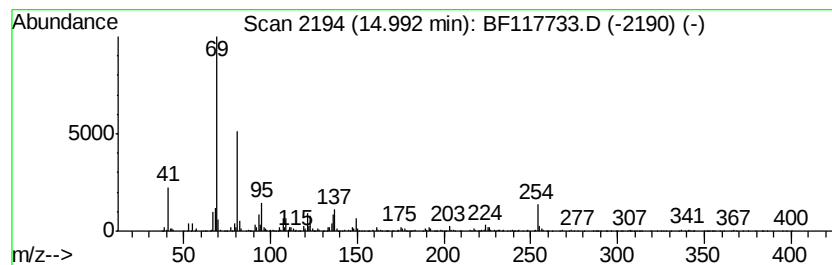
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Squalene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.99	5.74 ng	653140	Perylene-d12	15.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Squalene	410	C30H50	000111-02-4	95
2		7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24O	125529-10-4	72
3		4,8,12-Tetradecatrienenitrile, 5...	245	C17H27N	005981-31-7	64
4		2,6-Octadiene, 2,6-dimethyl-	138	C10H18	002792-39-4	53
5		Trideca-1,7,11-triene-1,1-dicarb...	270	C18H26N2	1000161-46-0	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110819\
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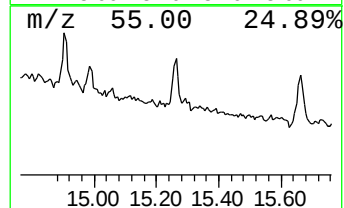
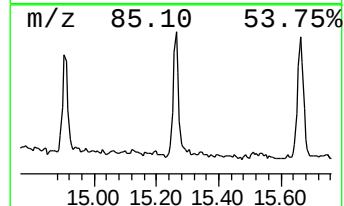
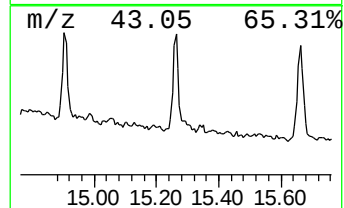
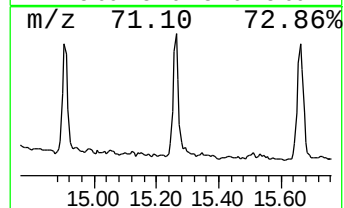
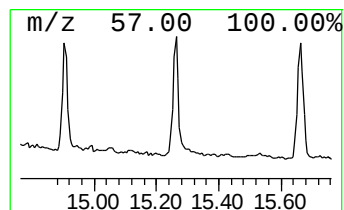
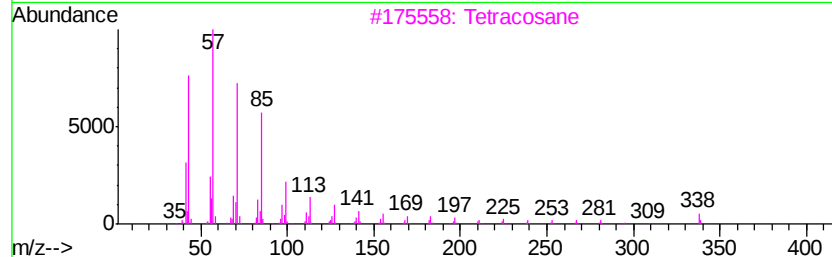
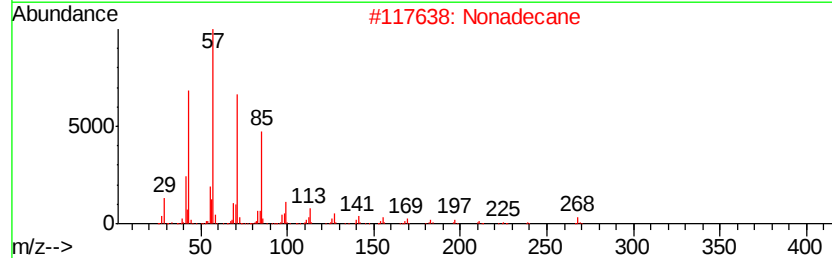
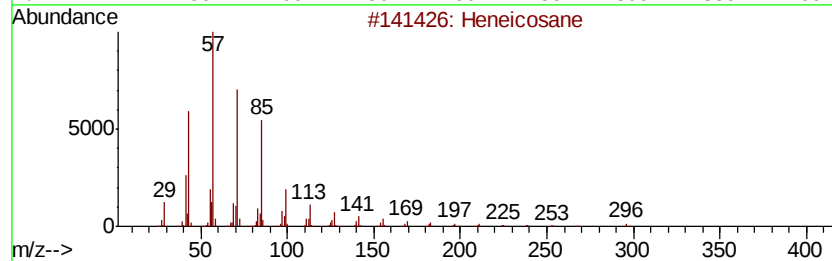
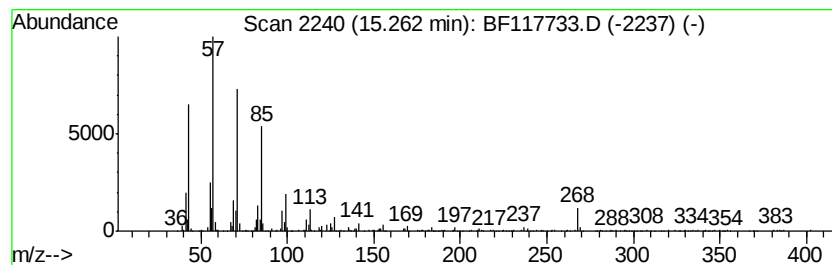
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Heneicosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.26	3.09 ng	351489	Perylene-d12	15.63
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Heneicosane	296	C21H44	000629-94-7 97
2	Nonadecane	268	C19H40	000629-92-5 91
3	Tetracosane	338	C24H50	000646-31-1 87
4	Eicosane, 7-hexyl-	366	C26H54	055333-99-8 86
5	Heptadecane, 9-octyl-	352	C25H52	007225-64-1 83



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110819\
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ClientSampleId :
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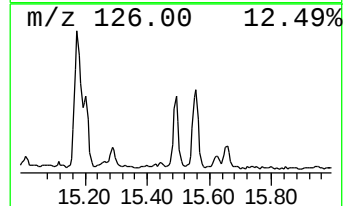
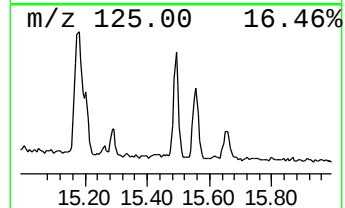
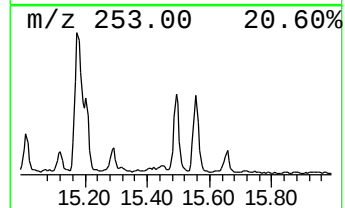
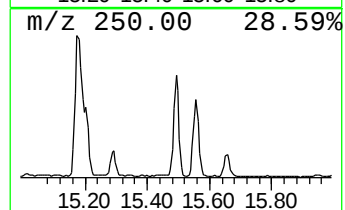
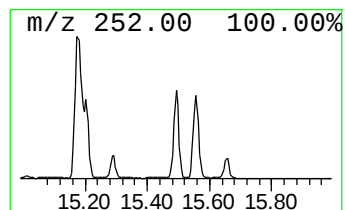
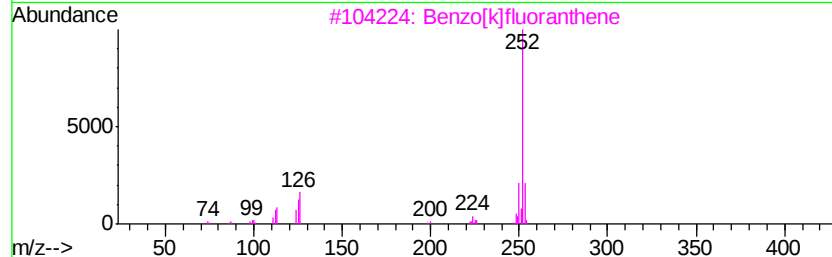
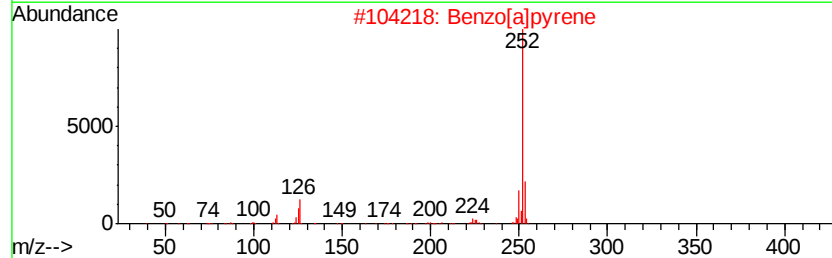
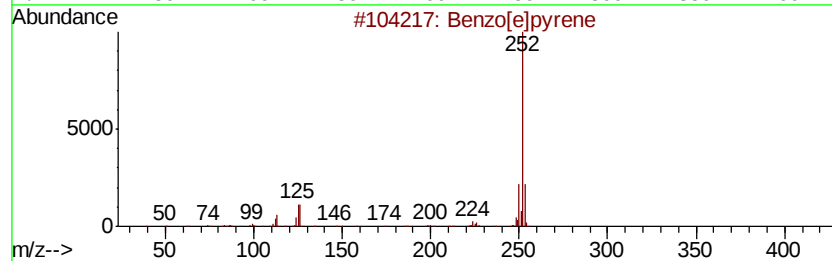
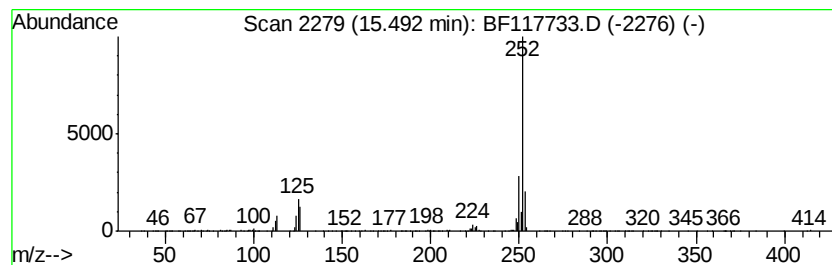
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Benzo[e]pyrene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.49	4.55 ng	516938	Perylene-d12	15.63

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	98
2			Benzo[a]pyrene	252	C20H12	000050-32-8	96
3			Benzo[k]fluoranthene	252	C20H12	000207-08-9	95
4			Benzo[e]acephenanthrylene	252	C20H12	000205-99-2	94
5			Perylene	252	C20H12	000198-55-0	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110819\
Data File : BF117733.D
Acq On : 8 Nov 2019 16:16
Operator : JU
Sample : K5722-13
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
1-C

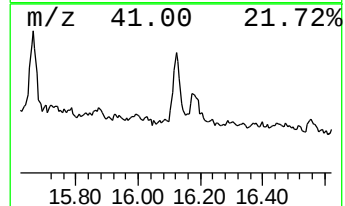
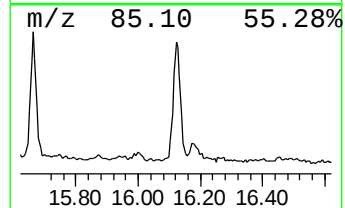
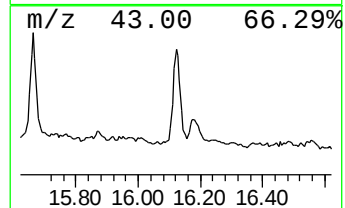
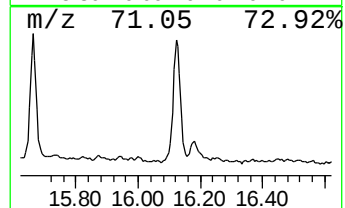
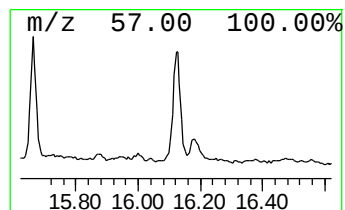
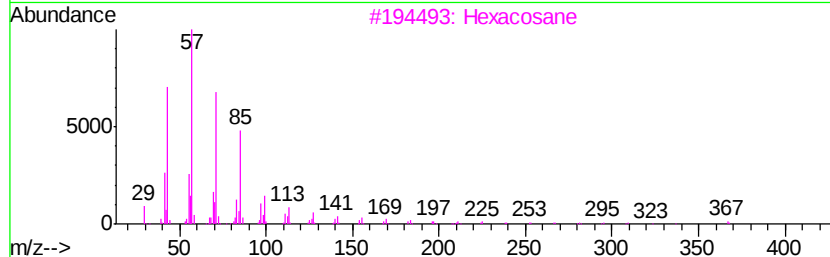
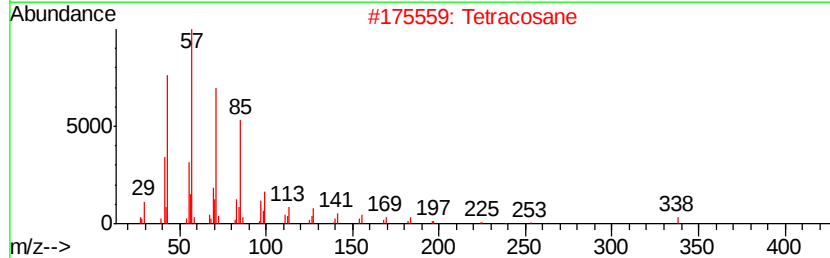
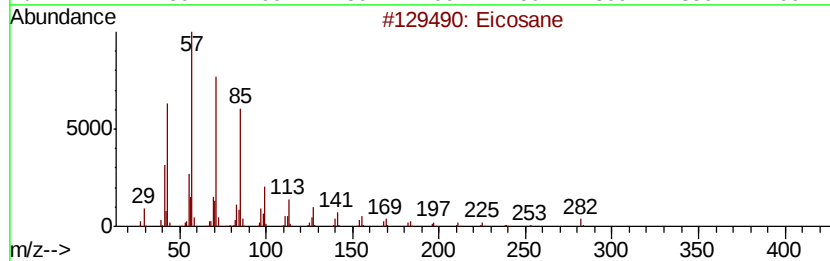
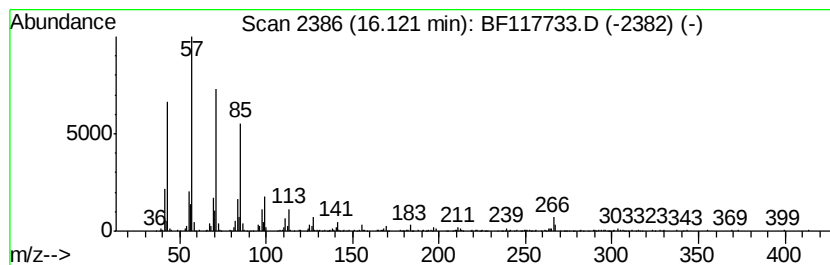
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Eicosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.12	4.58 ng	520913	Perylene-d12	15.63

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane			282	C20H42	000112-95-8	93
2	Tetracosane			338	C24H50	000646-31-1	90
3	Hexacosane			366	C26H54	000630-01-3	90
4	Heneicosane			296	C21H44	000629-94-7	90
5	2-methyloctacosane			408	C29H60	1000376-72-8	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110819\
 Data File : BF117733.D
 Acq On : 8 Nov 2019 16:16
 Operator : JU
 Sample : K5722-13
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 1-C

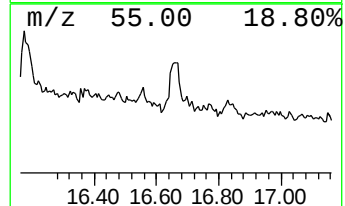
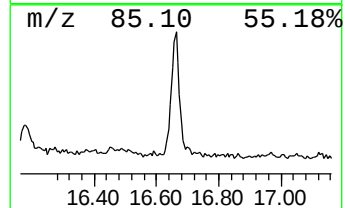
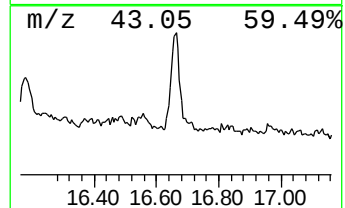
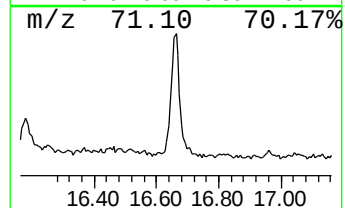
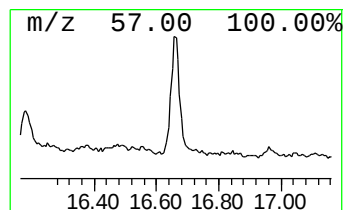
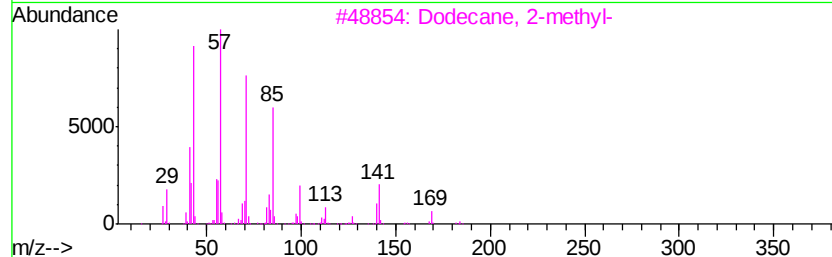
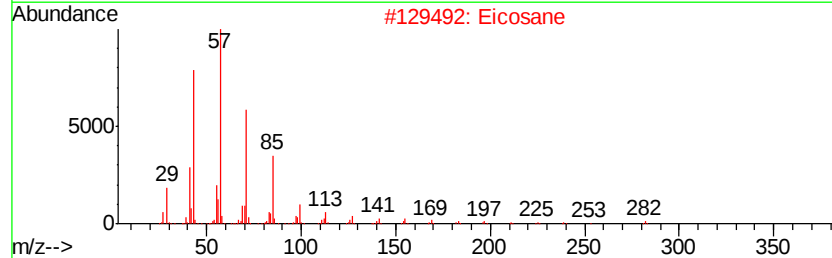
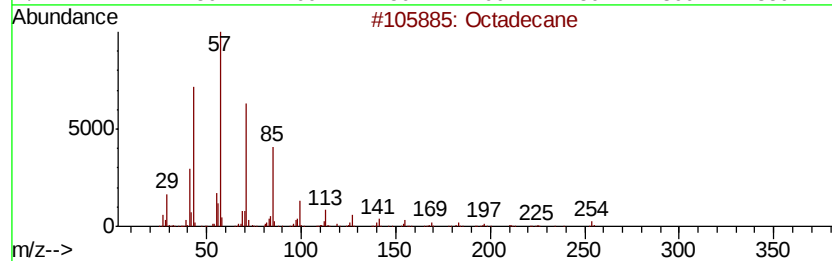
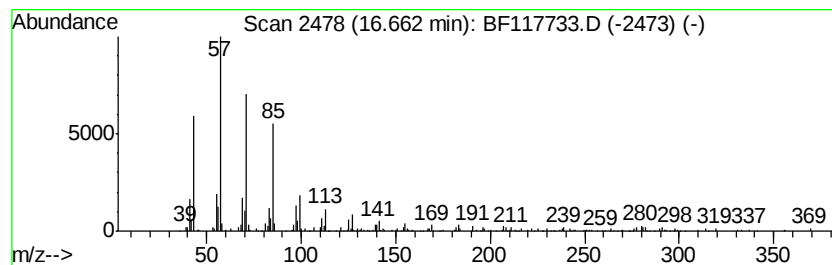
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 15 Octadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.66	3.04 ng	345407	Perylene-d12	15.63

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane		254	C18H38	000593-45-3	95
2	Eicosane		282	C20H42	000112-95-8	95
3	Dodecane, 2-methyl-		184	C13H28	001560-97-0	90
4	Triacontane		422	C30H62	000638-68-6	90
5	Heneicosane		296	C21H44	000629-94-7	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF110819\
 Data File : BF117733.D
 Acq On : 8 Nov 2019 16:16
 Operator : JU
 Sample : K5722-13
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 1-C

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110619.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
Butane, 2-methoxy...	2.23	14.6	ng	1521070	1	6.93	2089480	20.0
2-Pentanone, 4-hy...	5.15	12.9	ng	1349810	1	6.93	2089480	20.0
unknown6.69	6.69	61.2	ng	6394820	1	6.93	2089480	20.0
Anthracene, 2-met...	11.97	2.2	ng	376710	4	11.47	3438170	20.0
Anthracene, 1-met...	12.00	7.5	ng	1291340	4	11.47	3438170	20.0
Naphtho[2,3-b]nor...	12.08	2.5	ng	425787	4	11.47	3438170	20.0
5-Eicosene, (E)-	13.96	9.5	ng	1667940	5	14.12	3513120	20.0
Docosane, 11-butyl-	14.59	2.5	ng	429830	5	14.12	3513120	20.0
Heptadecane	14.90	3.1	ng	357450	6	15.63	2274140	20.0
Squalene	14.99	5.7	ng	653140	6	15.63	2274140	20.0
Heneicosane	15.26	3.1	ng	351489	6	15.63	2274140	20.0
Benzo[e]pyrene	15.49	4.5	ng	516938	6	15.63	2274140	20.0
Eicosane	16.12	4.6	ng	520913	6	15.63	2274140	20.0
Octadecane	16.66	3.0	ng	345407	6	15.63	2274140	20.0