

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062619\
 Data File : BF115216.D
 Acq On : 26 Jun 2019 13:57
 Operator : HP/JU
 Sample : K3164-10
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BAT-B05

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-BF062119.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.454	423	428	433	rBV	550984	659882	8.07%	1.052%
2	5.207	551	556	559	rBV	3581166	3426089	41.87%	5.463%
3	6.242	728	732	735	rBV	2670143	2304764	28.17%	3.675%
4	6.278	735	738	742	rVB	464262	379246	4.64%	0.605%
5	6.378	750	755	758	rBV	5768291	5845341	71.44%	9.321%
6	6.607	791	794	798	rBV	1704288	1477814	18.06%	2.357%
7	6.766	817	821	825	rBV	5646664	5464872	66.79%	8.715%
8	7.172	885	890	893	rBV	4129448	4428239	54.12%	7.062%
9	7.413	925	931	934	rBV	257786	227082	2.78%	0.362%
10	7.889	1008	1012	1015	rBV	2285082	1902370	23.25%	3.034%
11	8.454	1105	1108	1111	rBV	162305	137120	1.68%	0.219%
12	8.972	1190	1196	1199	rBV	7537044	7839728	95.82%	12.502%
13	9.360	1259	1262	1267	rVB2	184035	212923	2.60%	0.340%
14	9.601	1299	1303	1305	rBV2	82436	99967	1.22%	0.159%
15	9.636	1305	1309	1312	rVV	2246216	2062544	25.21%	3.289%
16	9.666	1312	1314	1319	rVB	111914	121983	1.49%	0.195%
17	10.172	1394	1400	1405	rBV5	161999	327852	4.01%	0.523%
18	10.419	1437	1442	1445	rBV	4706912	4988929	60.98%	7.956%
19	11.107	1555	1559	1562	rBV	2690249	2294148	28.04%	3.658%
20	11.666	1650	1654	1658	rBV	1075481	967640	11.83%	1.543%
21	12.348	1768	1770	1771	rBV	140222	109709	1.34%	0.175%
22	12.372	1771	1774	1781	rVB	499571	481413	5.88%	0.768%
23	12.448	1784	1787	1795	rVV	1028347	962405	11.76%	1.535%
24	12.530	1797	1801	1804	rVB	1817637	1466953	17.93%	2.339%
25	12.695	1824	1829	1832	rBV	7042250	8181856	100.00%	13.047%
26	12.742	1834	1837	1840	rVB	450791	405147	4.95%	0.646%
27	13.542	1970	1973	1979	rVB	285332	268703	3.28%	0.428%
28	13.648	1988	1991	1998	rVB	244086	303933	3.71%	0.485%
29	13.724	2000	2004	2013	rVB	2380082	2275962	27.82%	3.629%
30	14.230	2088	2090	2093	rVB	101870	97034	1.19%	0.155%
31	14.524	2137	2140	2143	rBV	164470	148439	1.81%	0.237%
32	14.824	2188	2191	2200	rVB	230292	247783	3.03%	0.395%
33	15.089	2231	2236	2244	rVV	1681362	1835346	22.43%	2.927%
34	15.160	2244	2248	2258	rVB	243398	312614	3.82%	0.499%

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ClientSampleId :
BAT-B05

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	15.542	2309	2313	2320	rVB	173029	237160	2.90%	0.378%
36	15.983	2384	2388	2397	rVB2	110156	206105	2.52%	0.329%

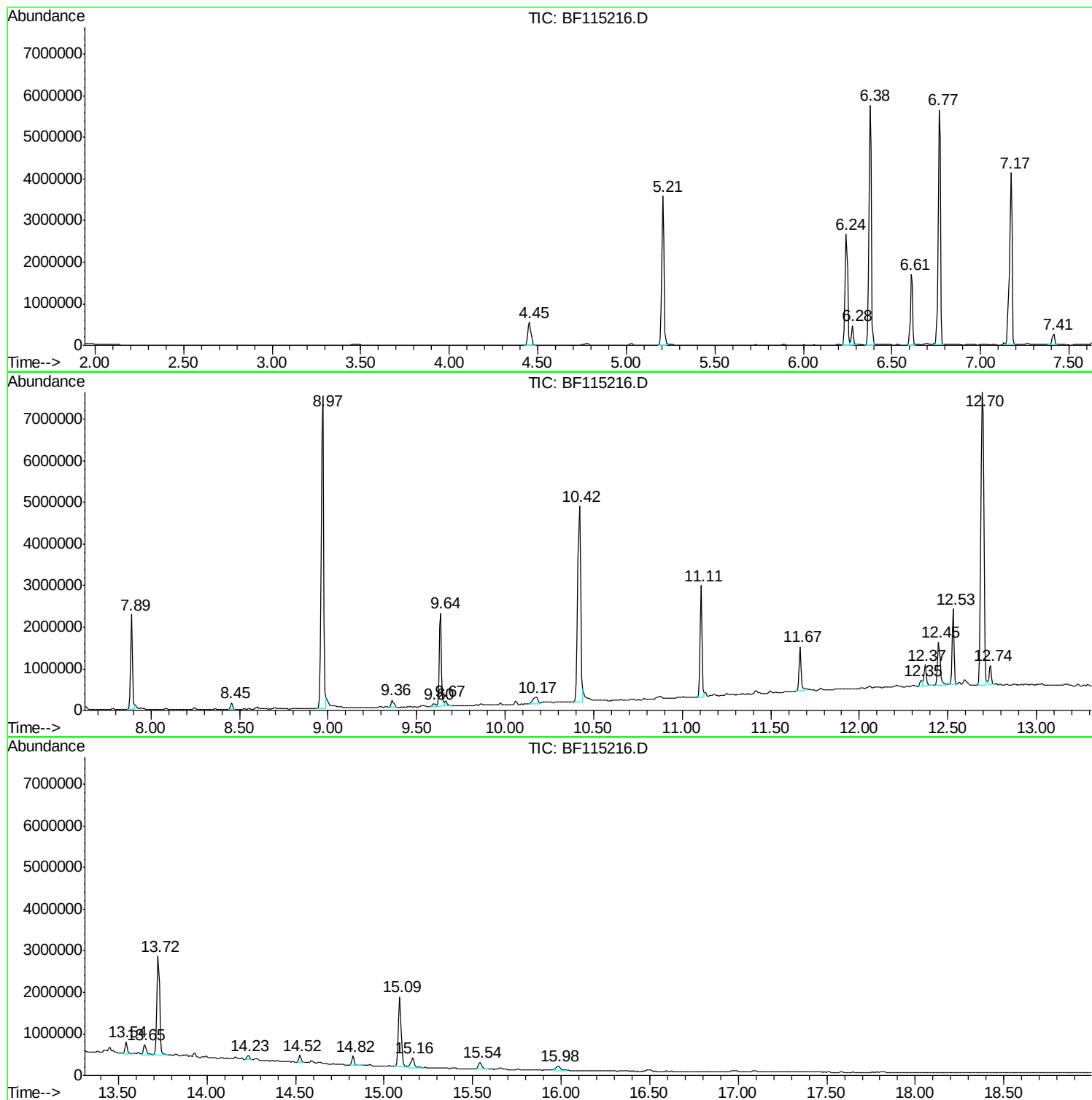
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF062119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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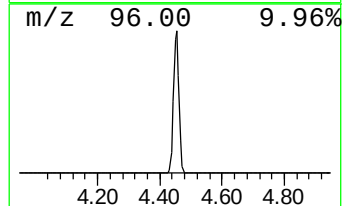
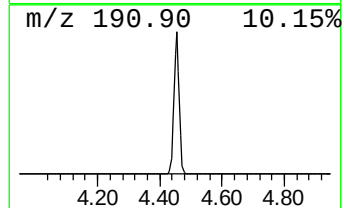
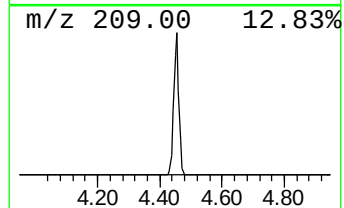
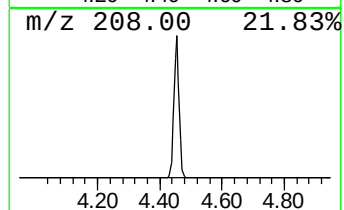
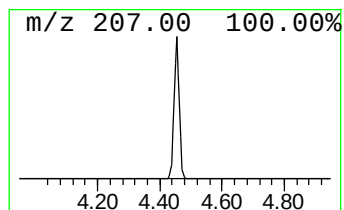
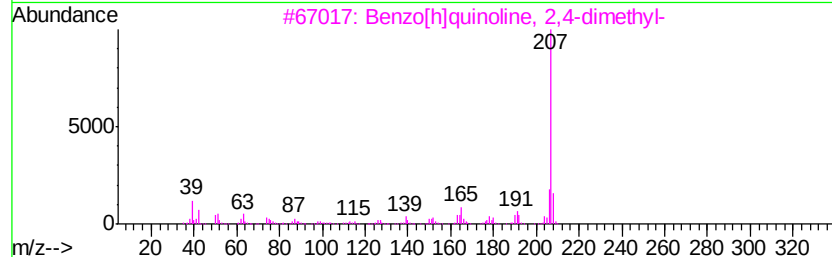
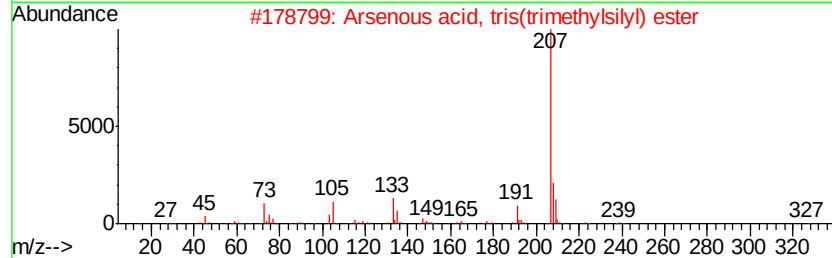
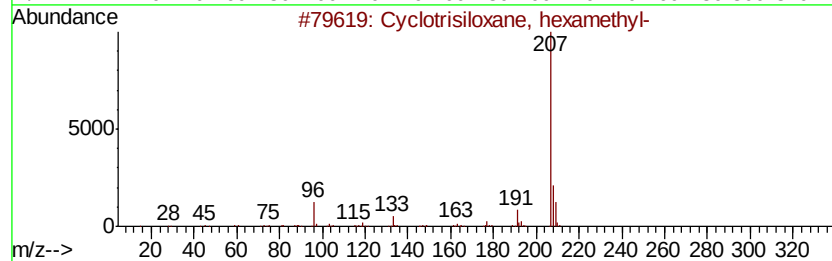
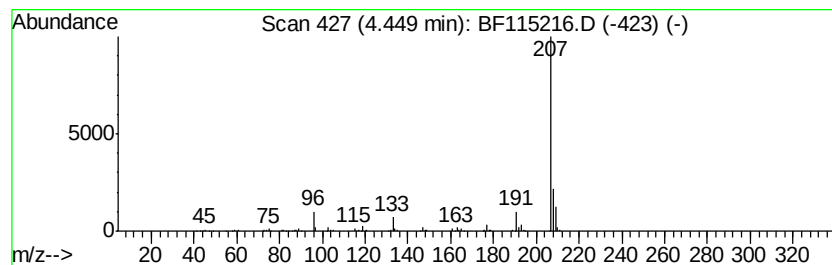
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF062119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Cyclotrisiloxane, hexamethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.45	8.93 ng	659882	1,4-Dichlorobenzene-d4	6.61

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclotrisiloxane, hexamethyl-	222	C6H18O3Si3	000541-05-9	91
2		Arsenous acid, tris(trimethylsil...	342	C9H27AsO3Si3	055429-29-3	64
3		Benzo[h]quinoline, 2,4-dimethyl-	207	C15H13N	000605-67-4	39
4		1H-Indole, 1-methyl-2-phenyl-	207	C15H13N	003558-24-5	39
5		1H-Indole, 5-methyl-2-phenyl-	207	C15H13N	013228-36-9	9



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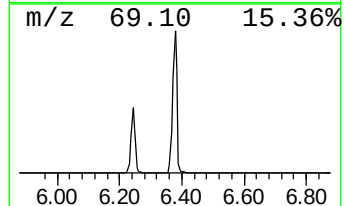
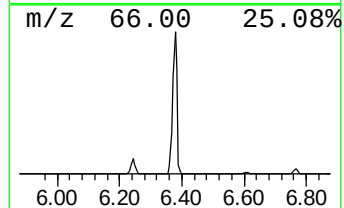
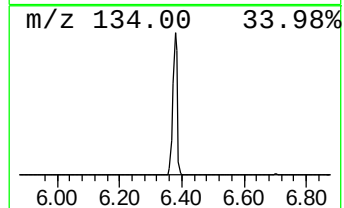
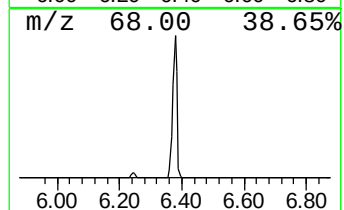
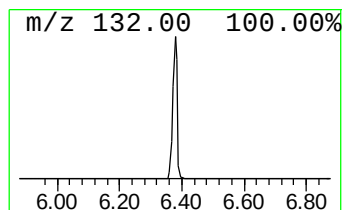
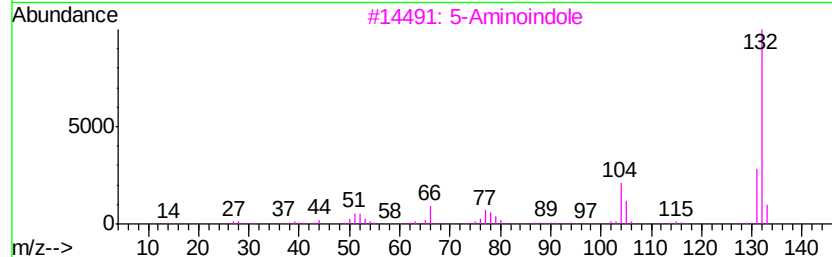
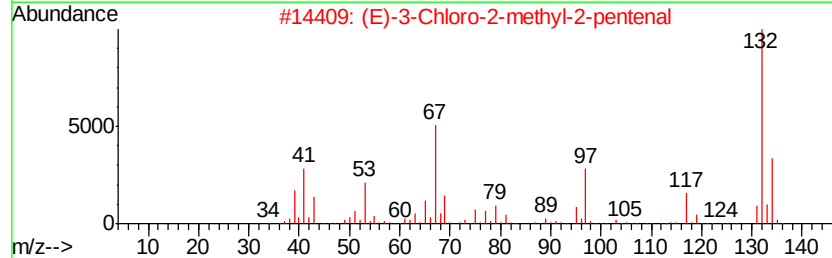
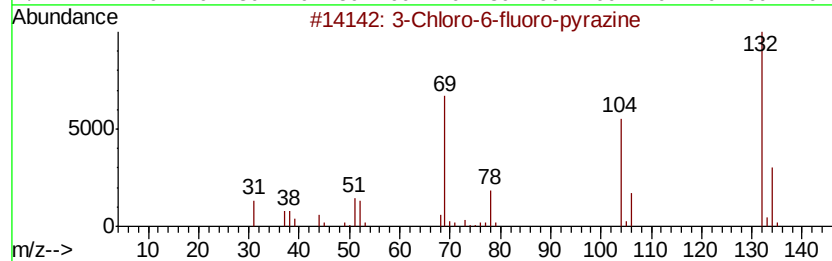
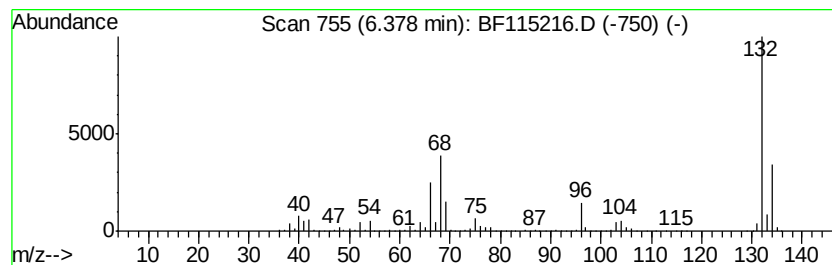
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 unknown6.38 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.38	79.11 ng	5845340	1,4-Dichlorobenzene-d4	6.61

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
5		Xenon	132	Xe	007440-63-3	9



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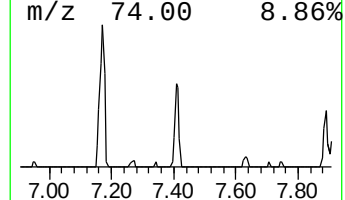
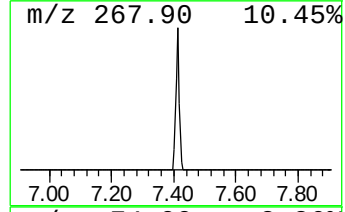
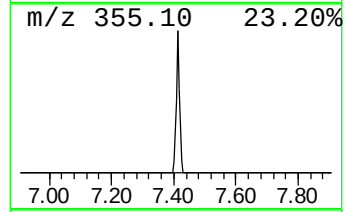
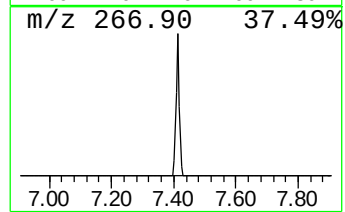
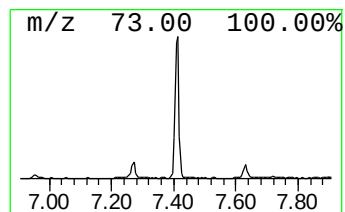
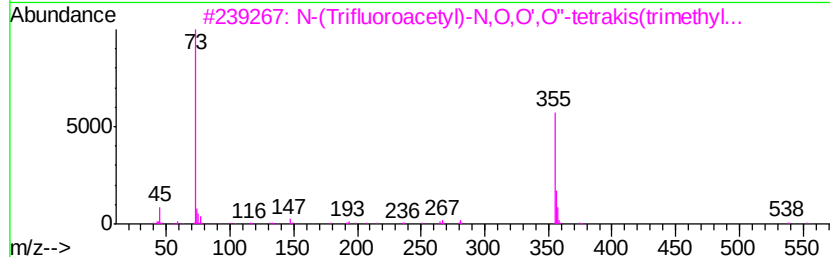
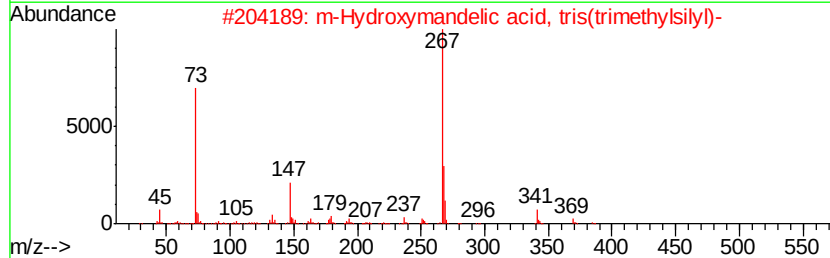
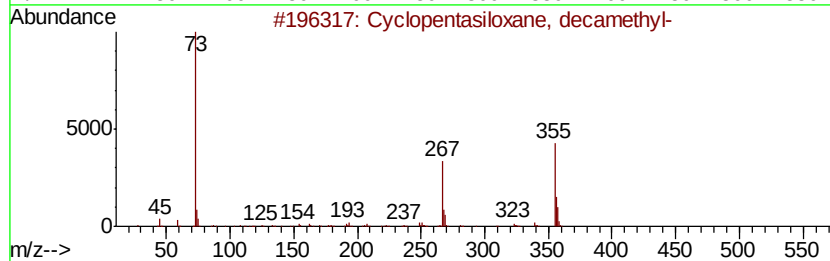
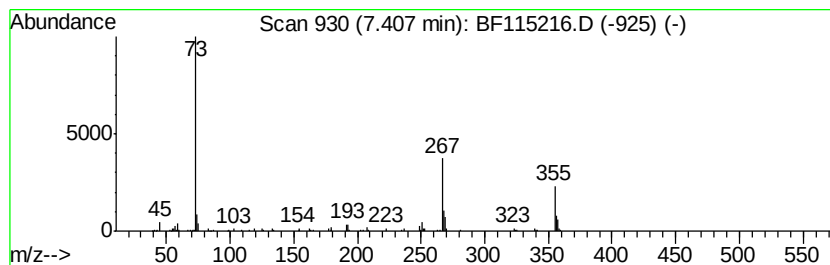
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Cyclopentasiloxane, decamet... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.41	2.39 ng	227082	Naphthalene-d8	7.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	81
2		m-Hydroxymandelic acid, tris(tri...	384	C17H32O4Si3	068595-69-7	43
3		N-(Trifluoroacetyl)-N,O,O',O''-t...	553	C22H42F3N04Si4	1000072-26-7	43
4		4-(2-Acetylamino-1-(trimethylsil...	339	C16H29N03Si2	1000373-43-5	38
5		Silanamine, 1,1,1-trimethyl-N-[2...	369	C17H35N02Si3	068595-60-8	38



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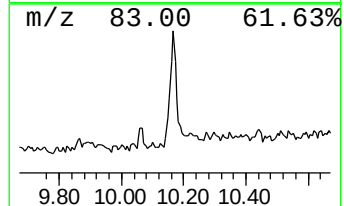
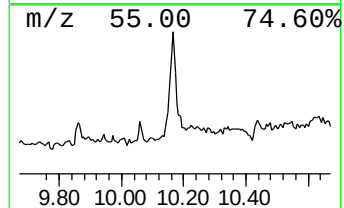
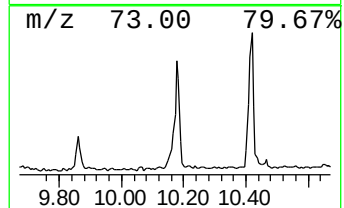
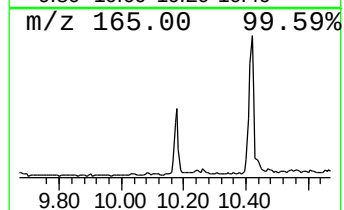
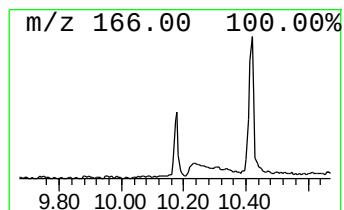
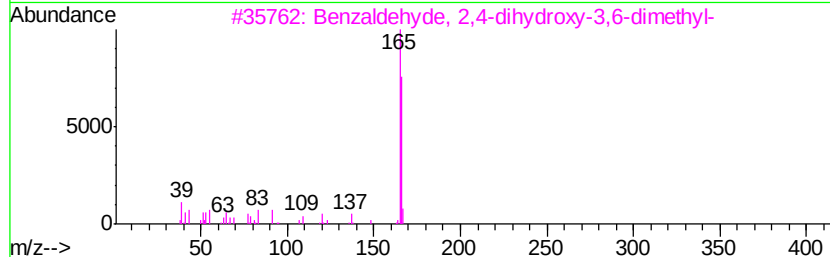
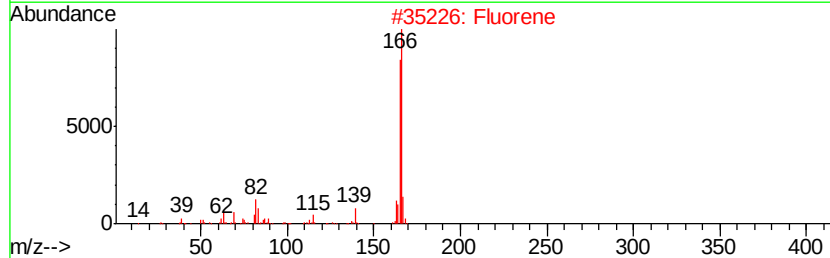
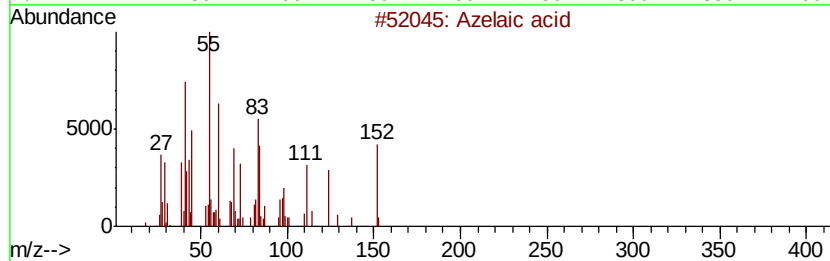
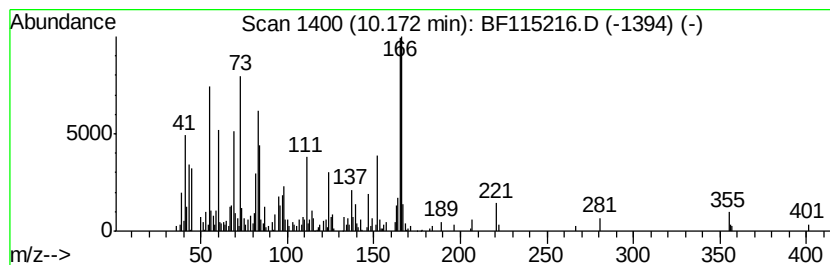
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 unknown10.17 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.17	3.18 ng	327852	Acenaphthene-d10	9.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Azelaic acid	188	C9H16O4	000123-99-9	38
2		Fluorene	166	C13H10	000086-73-7	38
3		Benzaldehyde, 2,4-dihydroxy-3,6-dimethyl-	166	C9H10O3	034883-14-2	30
4		1,3-Benzodioxole-5-carboxylic acid	166	C8H6O4	000094-53-1	27
5		Naphthalene, 1-(2-nitro-2-propenyl)-	213	C13H11NO2	080255-13-6	22



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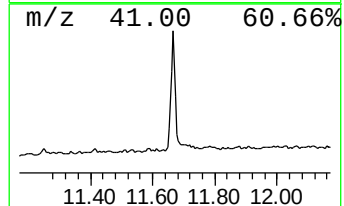
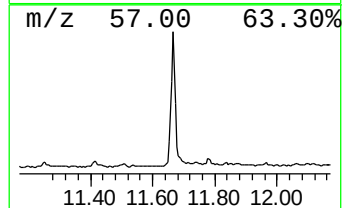
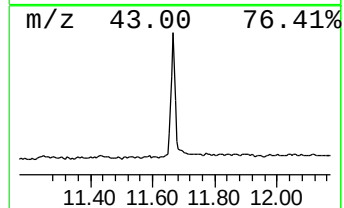
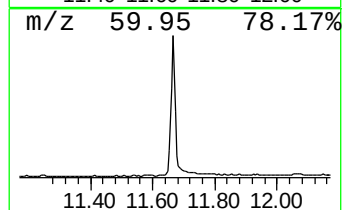
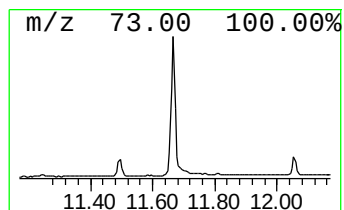
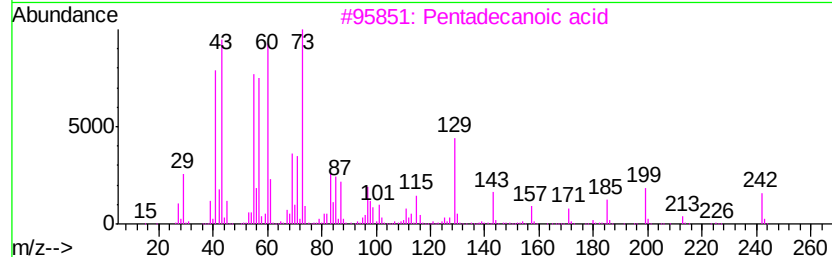
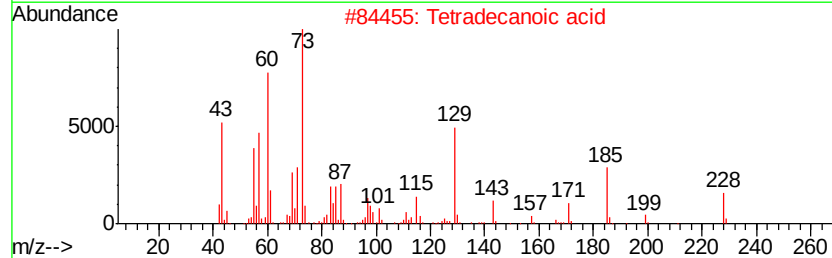
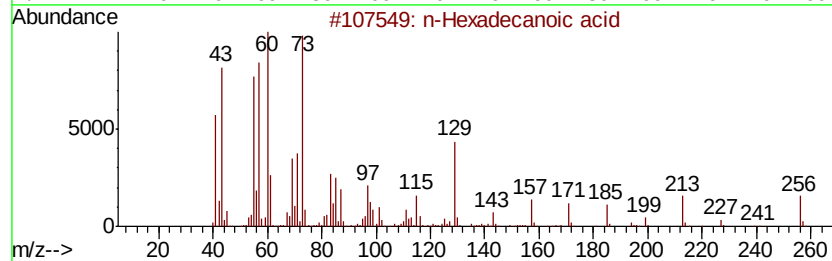
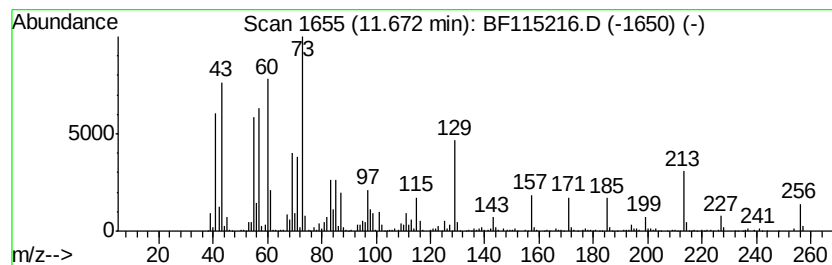
Instrument :
BNA_F
ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.67	8.44 ng	967640	Phenanthrene-d10	11.11	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	91
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	90
4	Tridecanoic acid	214	C13H26O2	000638-53-9	64
5	n-Decanoic acid	172	C10H20O2	000334-48-5	62



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062619\
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Acq On : 26 Jun 2019 13:57
Operator : HP/JU
Sample : K3164-10
Misc :
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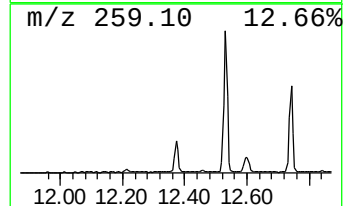
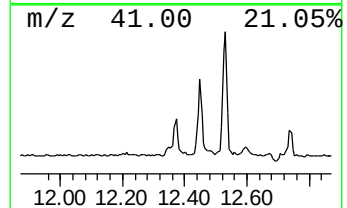
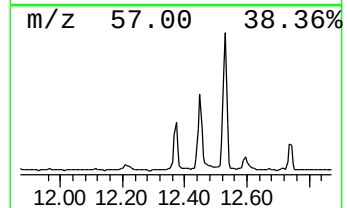
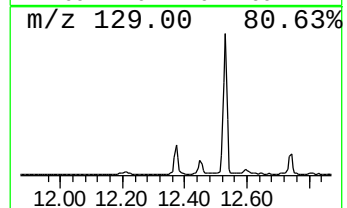
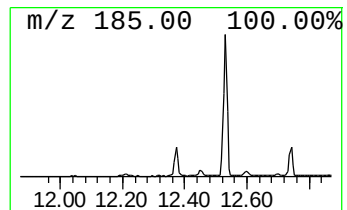
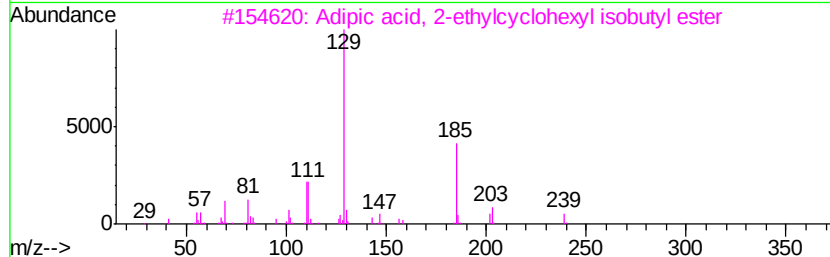
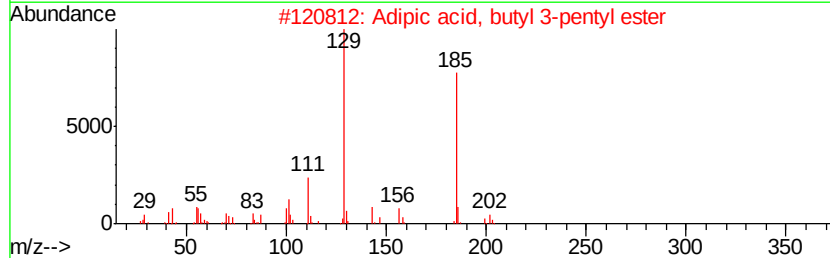
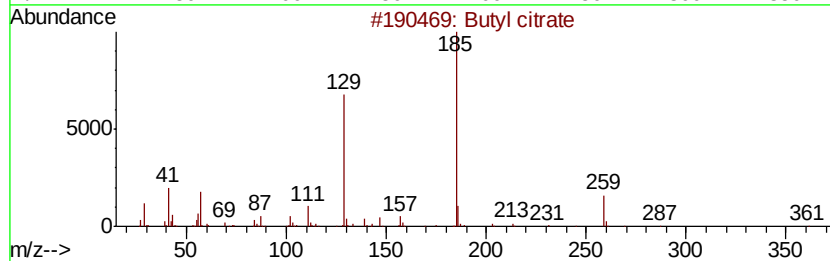
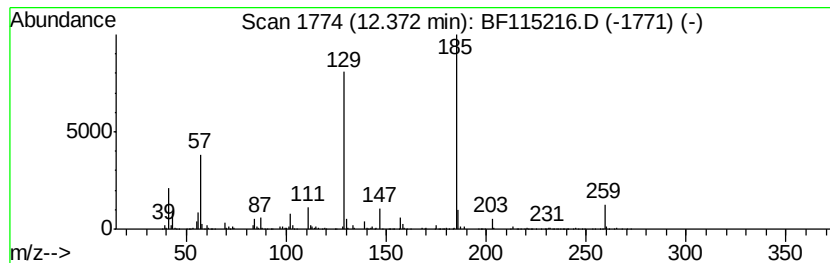
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Adipic acid, butyl 3-pentyl... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.37	4.20 ng	481413	Phenanthrene-d10	11.11
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Butyl citrate	360 C18H32O7	000077-94-1 87
2		Adipic acid, butyl 3-pentyl ester	272 C15H28O4	1000324-60-5 72
3		Adipic acid, 2-ethylcyclohexyl i...	312 C18H32O4	1000324-22-5 56
4		Adipic acid, isobutyl 3-pentyl e...	272 C15H28O4	1000324-60-4 50
5		Adipic acid, 3-hexyl isobutyl ester	286 C16H30O4	1000353-62-2 40



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062619\
Data File : BF115216.D
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Sample : K3164-10
Misc :
ALS Vial : 10 Sample Multiplier: 1

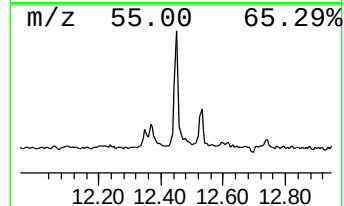
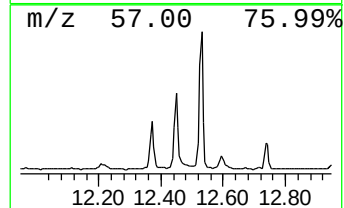
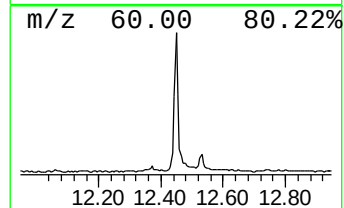
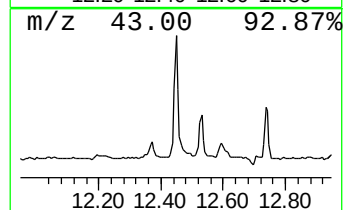
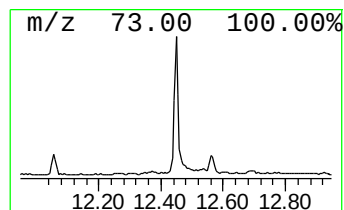
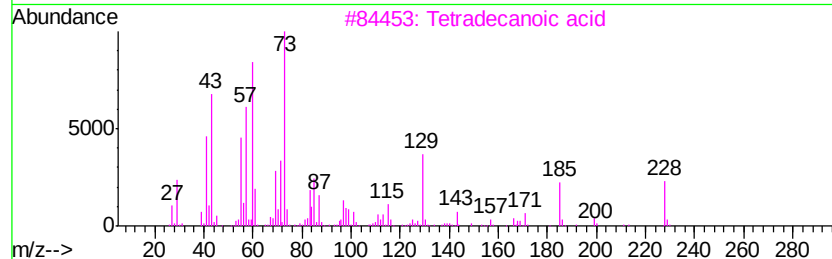
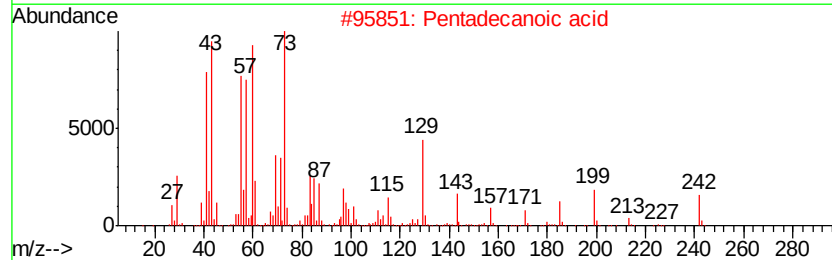
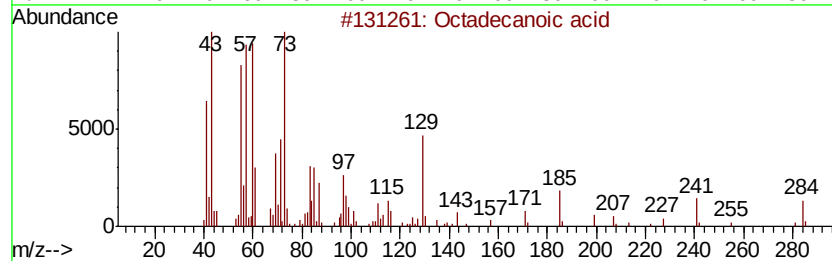
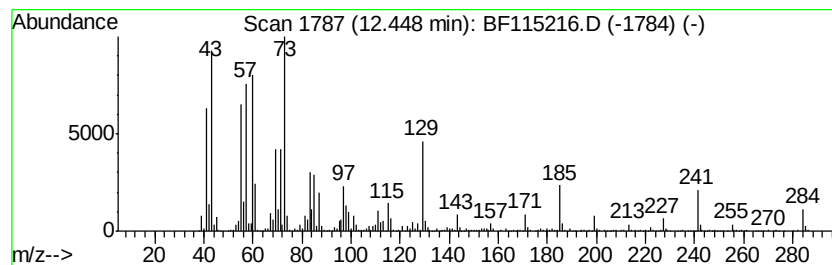
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Octadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.45	8.46 ng	962405	Chrysene-d12	13.72	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	89
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	83
4	Tridecanoic acid	214	C13H26O2	000638-53-9	74
5	n-Decanoic acid	172	C10H20O2	000334-48-5	70



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062619\
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Sample : K3164-10
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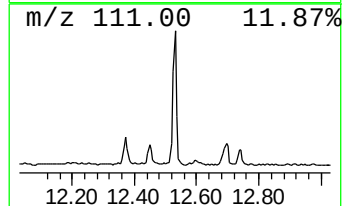
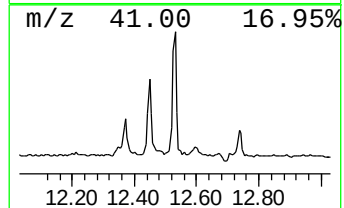
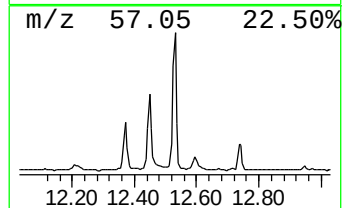
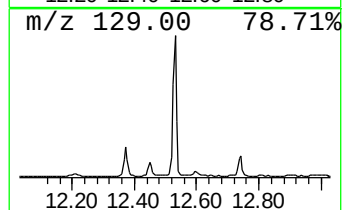
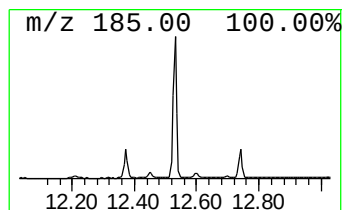
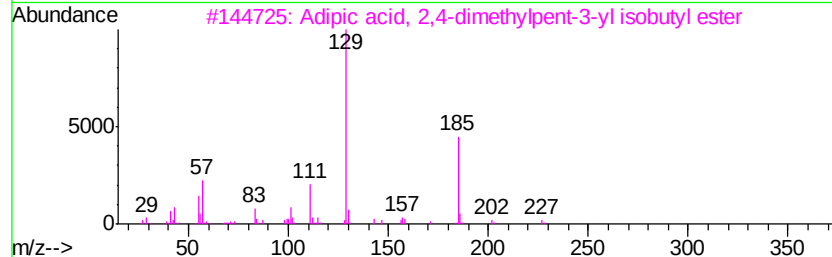
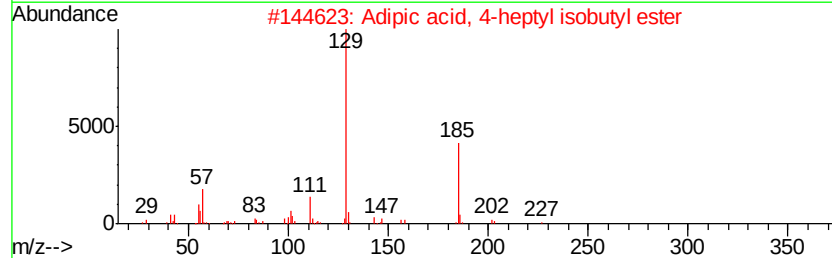
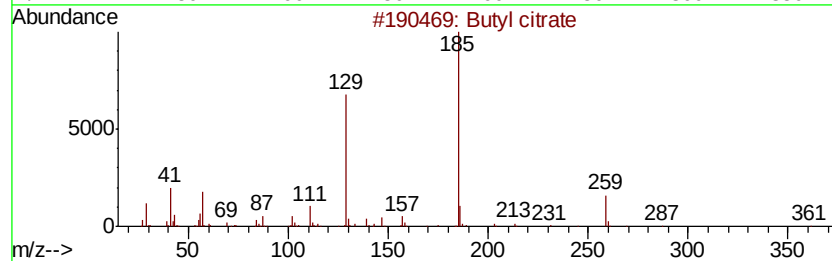
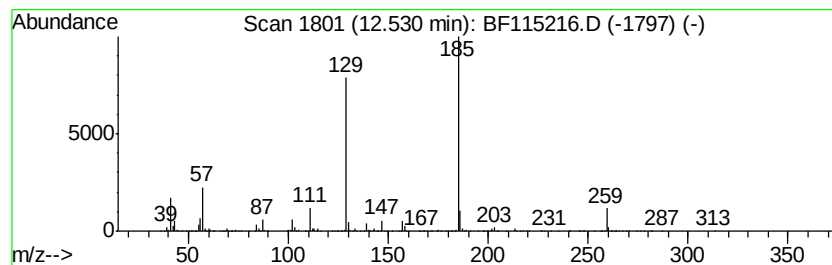
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF062119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Butyl citrate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.		
12.53	12.89 ng	1466950	Chrysene-d12	13.72		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Butyl citrate	360	C18H32O7	000077-94-1	91
2		Adipic acid, 4-heptyl isobutyl e...	300	C17H32O4	1000349-73-6	78
3		Adipic acid, 2,4-dimethylpent-3-...	300	C17H32O4	1000349-77-6	64
4		Adipic acid, 2-hexyl isobutyl ester	286	C16H30O4	1000353-70-9	64
5		Dibutyl adipate	258	C14H26O4	000105-99-7	56



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062619\
Data File : BF115216.D
Acq On : 26 Jun 2019 13:57
Operator : HP/JU
Sample : K3164-10
Misc :
ALS Vial : 10 Sample Multiplier: 1

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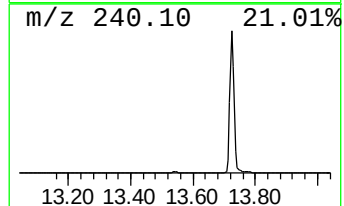
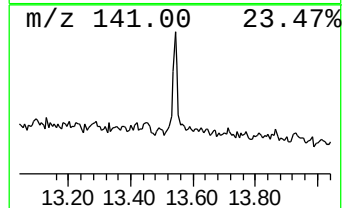
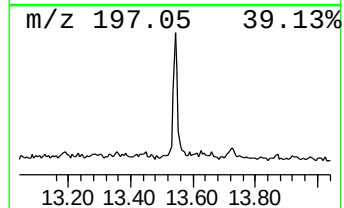
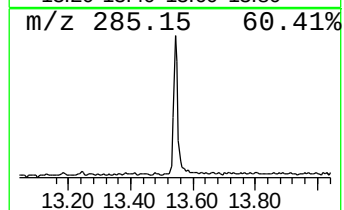
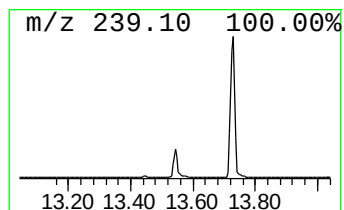
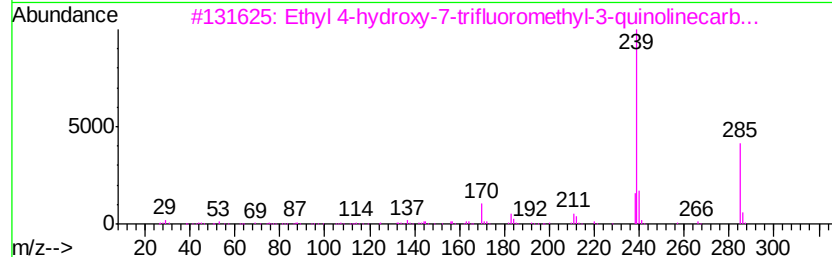
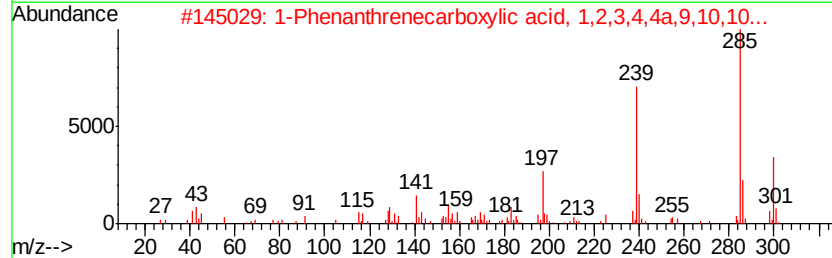
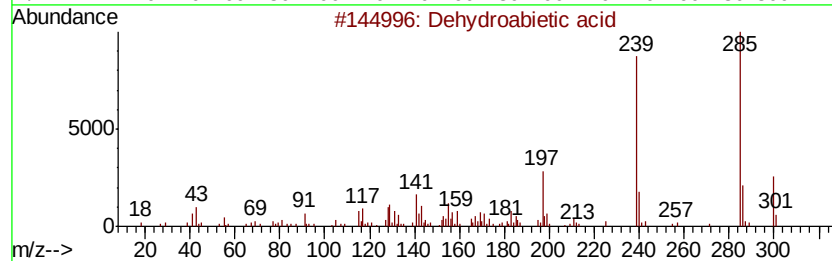
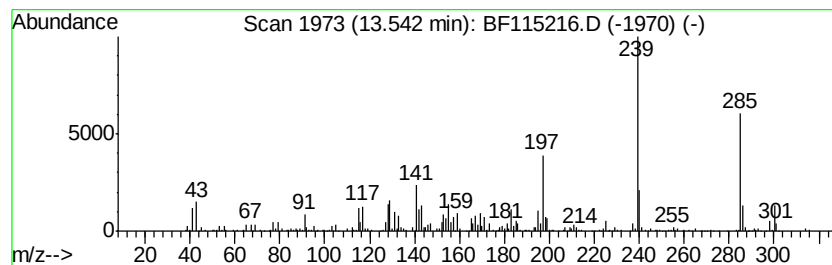
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Dehydroabietic acid Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.54	2.36 ng	268703	Chrysene-d12	13.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dehydroabietic acid	300	C20H28O2	001740-19-8	97
2		1-Phenanthrenecarboxylic acid, 1...	300	C20H28O2	005155-70-4	94
3		Ethyl 4-hydroxy-7-trifluoromethyl...	285	C13H10F3NO3	000391-02-6	49
4		2-Ethylidene-hydrazono-3-methyl-...	239	C10H10ClN3S	1000148-08-4	43
5		Imidazolo[2,1-b]thiazole, 5-(3-i...	239	C13H9N3S	327097-37-0	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062619\
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Sample : K3164-10
Misc :
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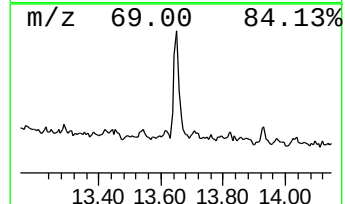
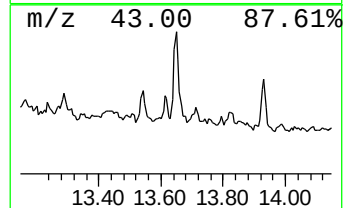
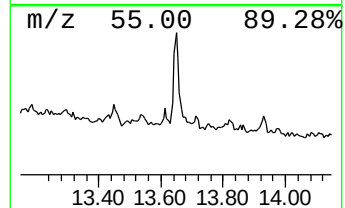
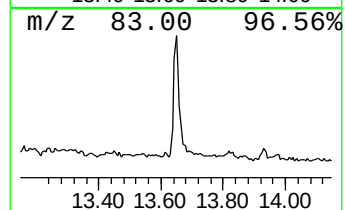
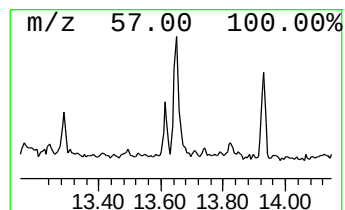
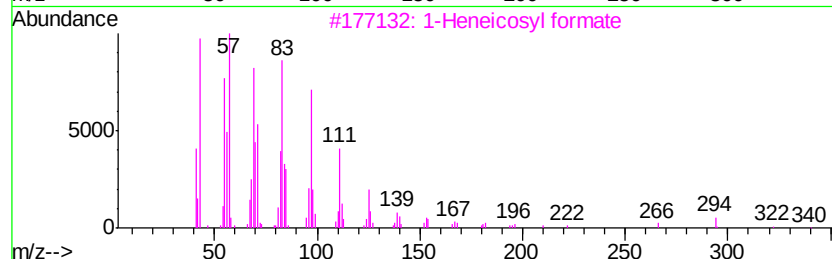
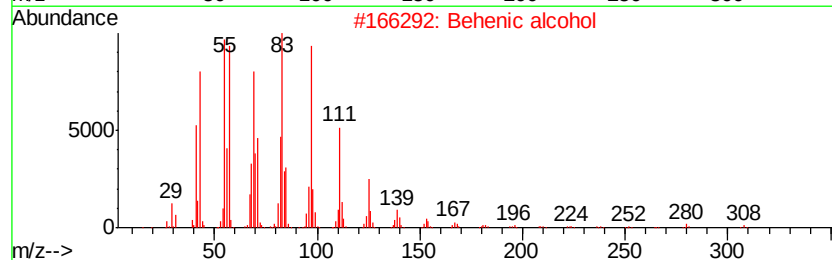
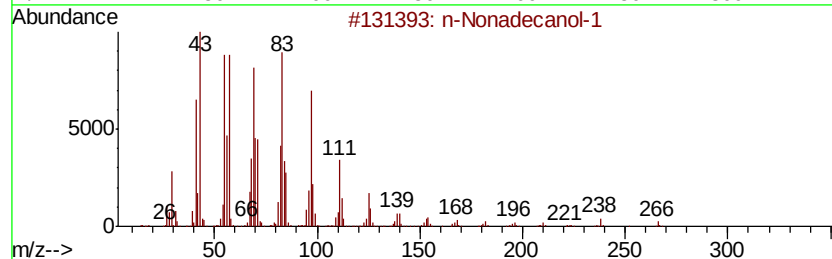
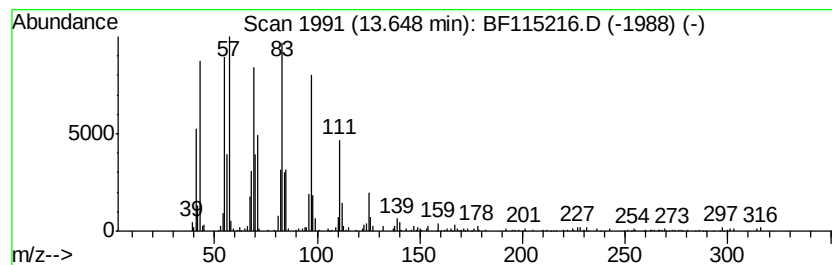
Instrument :
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 n-Nonadecanol-1 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.
13.65	2.67 ng	303933	Chrysene-d12		13.72
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Nonadecanol-1	284	C19H40O	001454-84-8	94
2	Behenic alcohol	326	C22H46O	000661-19-8	94
3	1-Heneicosyl formate	340	C22H44O2	077899-03-7	94
4	n-Tetracosanol-1	354	C24H50O	000506-51-4	94
5	5-Eicosene, (E)-	280	C20H40	074685-30-6	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062619\
Data File : BF115216.D
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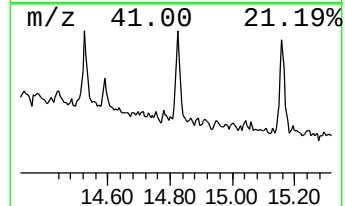
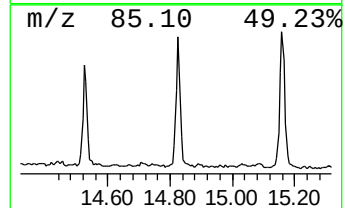
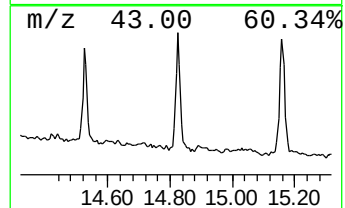
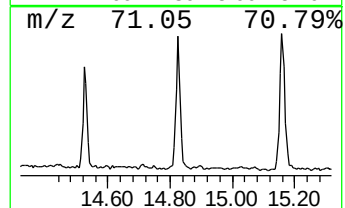
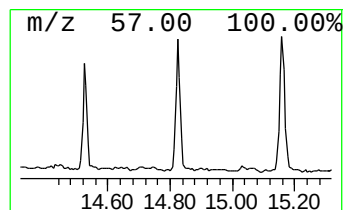
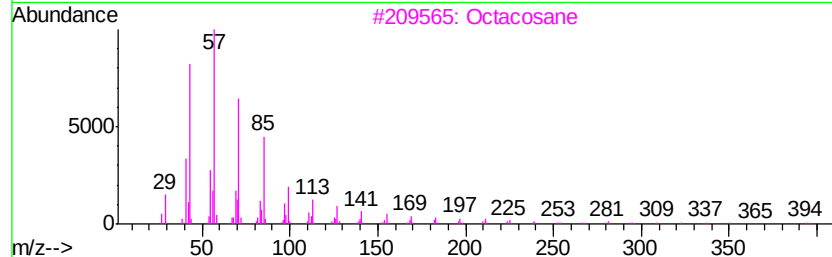
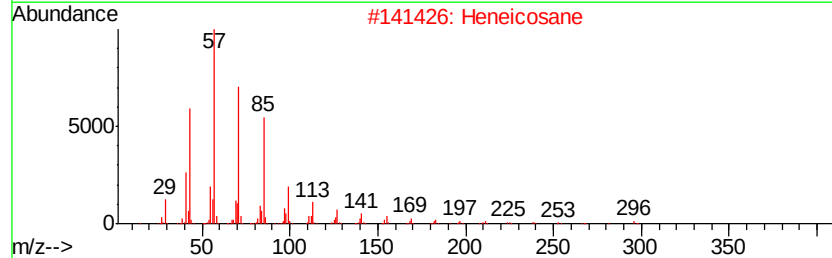
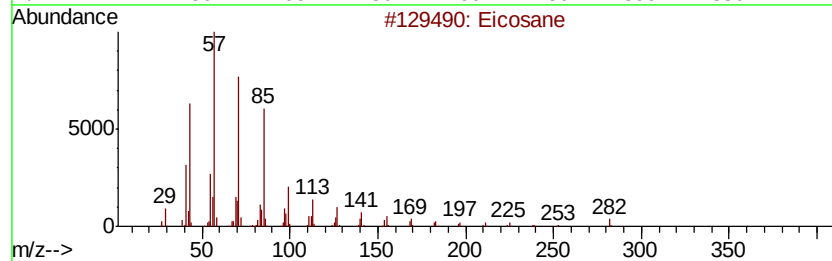
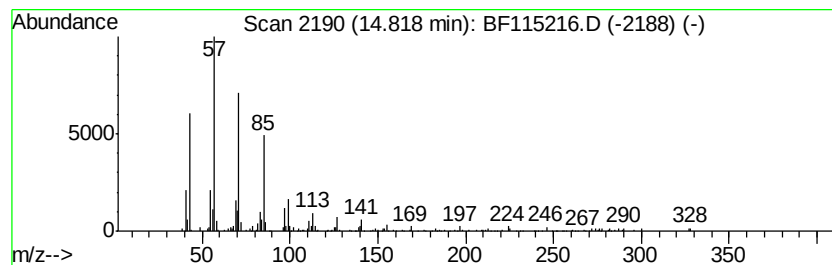
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Eicosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.82	2.70 ng	247783	Perylene-d12	15.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	97
2		Heneicosane	296	C21H44	000629-94-7	91
3		Octacosane	394	C28H58	000630-02-4	91
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
5		2-methyloctacosane	408	C29H60	1000376-72-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062619\
Data File : BF115216.D
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ALS Vial : 10 Sample Multiplier: 1

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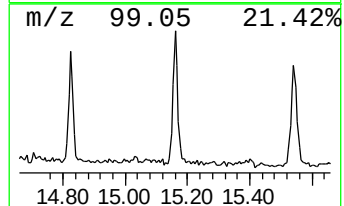
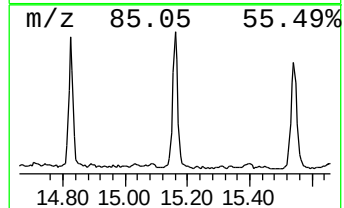
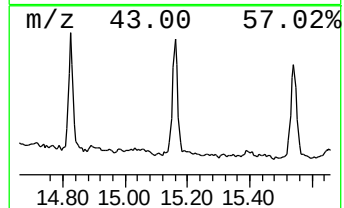
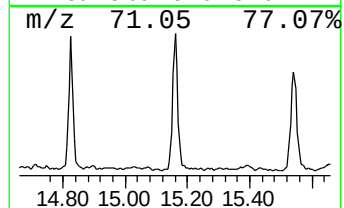
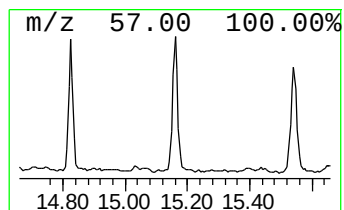
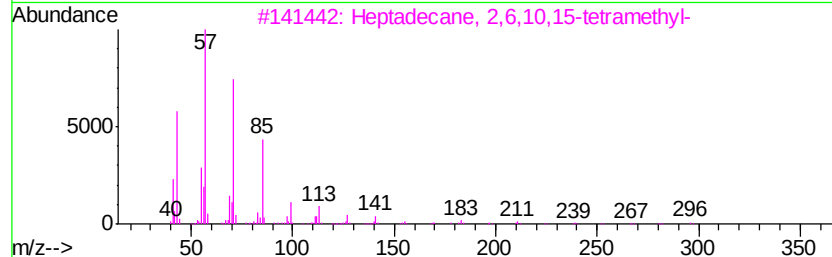
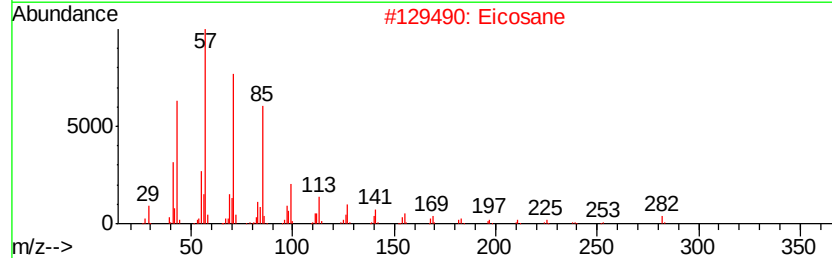
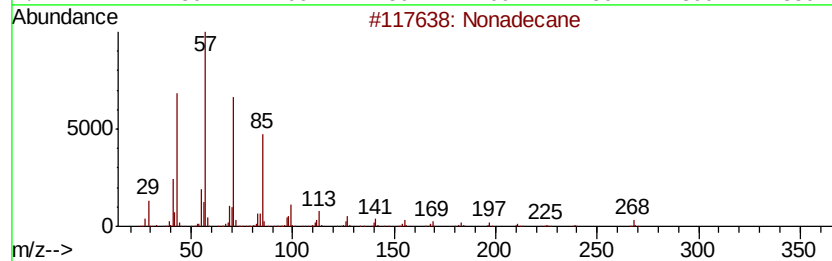
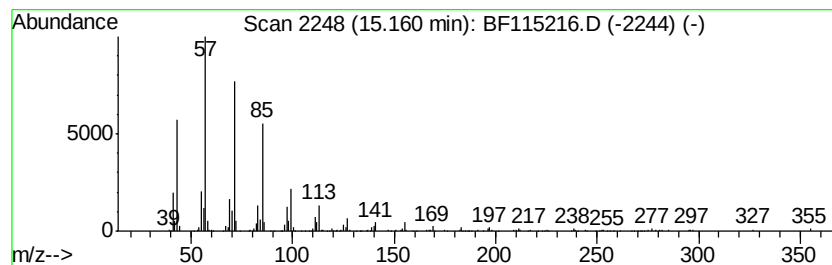
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 Nonadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.16	3.41 ng	312614	Perylene-d12	15.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	95
2		Eicosane	282	C20H42	000112-95-8	91
3		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	91
4		Heneicosane	296	C21H44	000629-94-7	91
5		Hexadecane	226	C16H34	000544-76-3	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062619\
Data File : BF115216.D
Acq On : 26 Jun 2019 13:57
Operator : HP/JU
Sample : K3164-10
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
BAT-B05

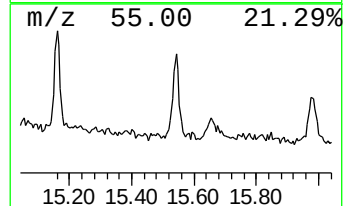
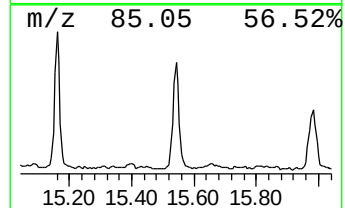
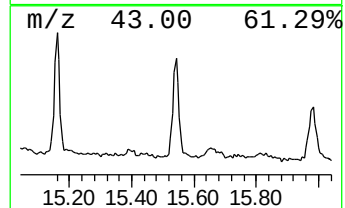
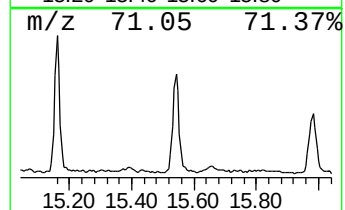
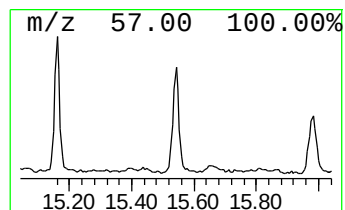
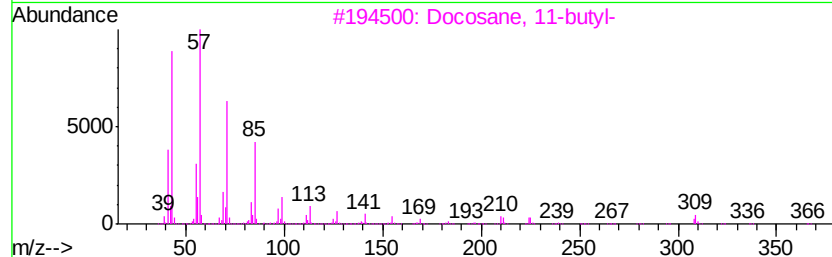
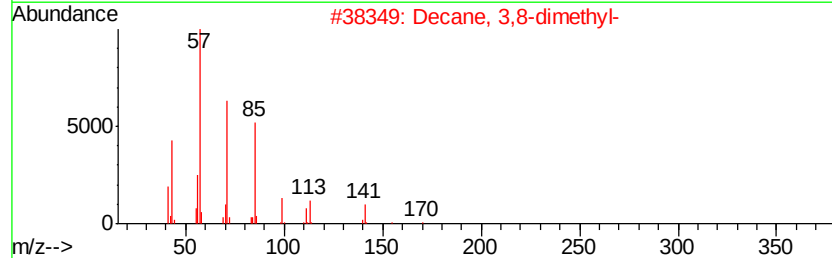
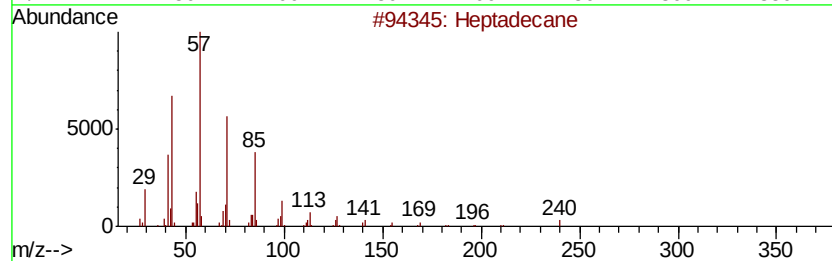
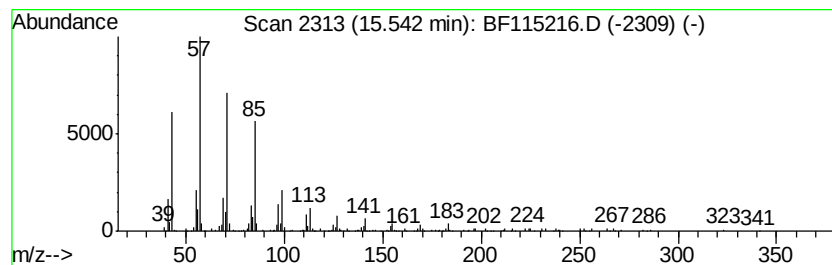
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Heptadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.54	2.58 ng	237160	Perylene-d12	15.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane	240	C17H36	000629-78-7	87
2		Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	81
3		Docosane, 11-butyl-	366	C26H54	013475-76-8	80
4		Hexacosane	366	C26H54	000630-01-3	80
5		Octadecane	254	C18H38	000593-45-3	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062619\
Data File : BF115216.D
Acq On : 26 Jun 2019 13:57
Operator : HP/JU
Sample : K3164-10
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
BAT-B05

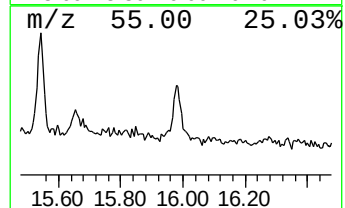
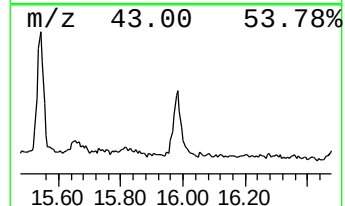
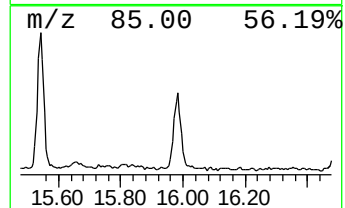
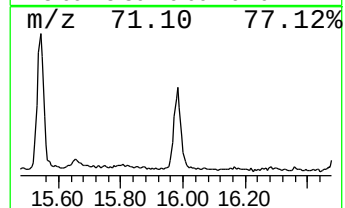
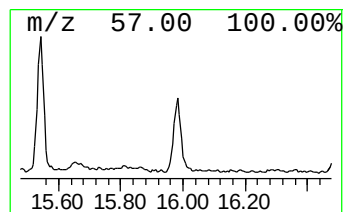
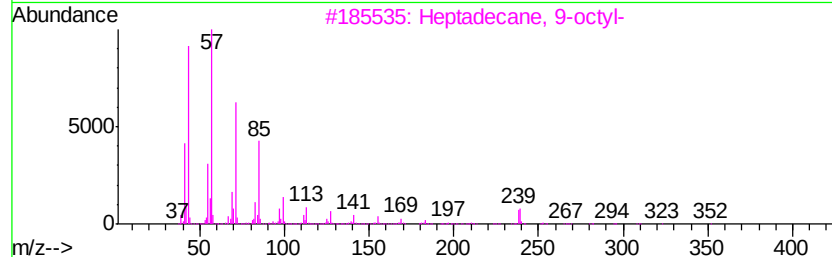
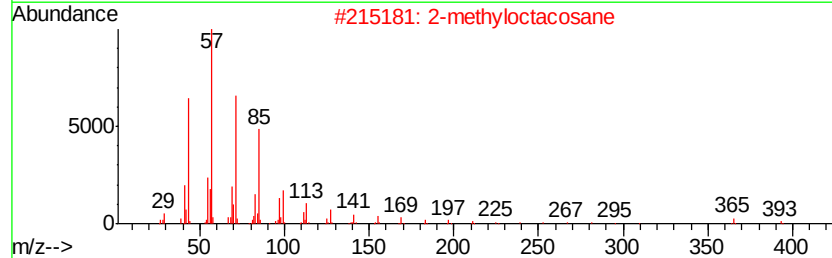
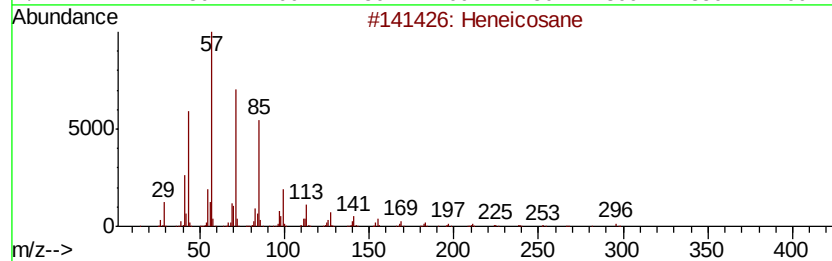
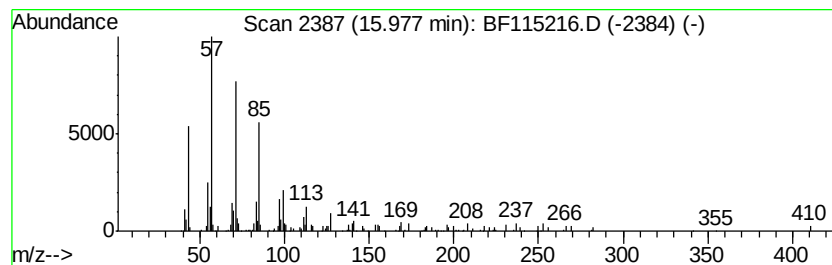
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Heneicosane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.98	2.25 ng	206105	Perylene-d12	15.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		2-methyloctacosane	408	C29H60	1000376-72-8	91
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	90
4		Docosane, 11-butyl-	366	C26H54	013475-76-8	90
5		Pentacosane	352	C25H52	000629-99-2	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF062619\
 Data File : BF115216.D
 Acq On : 26 Jun 2019 13:57
 Operator : HP/JU
 Sample : K3164-10
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BAT-B05

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF062119.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
Cyclotrisiloxane,...	4.45	8.9	ng	659882	1	6.61	1477810	20.0
unknown6.38	6.38	79.1	ng	5845340	1	6.61	1477810	20.0
Cyclopentasiloxan...	7.41	2.4	ng	227082	2	7.89	1902370	20.0
unknown10.17	10.17	3.2	ng	327852	3	9.64	2062540	20.0
n-Hexadecanoic acid	11.67	8.4	ng	967640	4	11.11	2294150	20.0
Adipic acid, buty...	12.37	4.2	ng	481413	4	11.11	2294150	20.0
Octadecanoic acid	12.45	8.5	ng	962405	5	13.72	2275960	20.0
Butyl citrate	12.53	12.9	ng	1466950	5	13.72	2275960	20.0
Dehydroabietic acid	13.54	2.4	ng	268703	5	13.72	2275960	20.0
n-Nonadecanol-1	13.65	2.7	ng	303933	5	13.72	2275960	20.0
Eicosane	14.82	2.7	ng	247783	6	15.09	1835350	20.0
Nonadecane	15.16	3.4	ng	312614	6	15.09	1835350	20.0
Heptadecane	15.54	2.6	ng	237160	6	15.09	1835350	20.0
Heneicosane	15.98	2.3	ng	206105	6	15.09	1835350	20.0