

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF090319\
 Data File : BF116773.D
 Acq On : 3 Sep 2019 15:38
 Operator : HP/JU
 Sample : K4554-57
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-1-(0-2)

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-BF083019.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	5.463	569	574	577	rBV	2275370	2085860	78.93%	7.726%
2	6.475	741	746	749	rBV	2566654	2226979	84.26%	8.248%
3	6.616	766	770	773	rBV	2658923	2304175	87.19%	8.534%
4	6.845	806	809	813	rBV	903470	695021	26.30%	2.574%
5	7.004	832	836	839	rBV	2201203	1734524	65.63%	6.424%
6	7.404	899	904	907	rBV	1730208	1419295	53.70%	5.257%
7	8.128	1023	1027	1030	rBV	1201902	922240	34.90%	3.416%
8	9.204	1206	1210	1213	rBV	2887475	2551252	96.53%	9.449%
9	9.592	1272	1276	1279	rBV	269468	224473	8.49%	0.831%
10	9.881	1321	1325	1329	rBV	1343236	1090557	41.26%	4.039%
11	10.663	1454	1458	1462	rBV	1724860	1629901	61.67%	6.037%
12	11.363	1573	1577	1580	rBV	1332875	1099637	41.61%	4.073%
13	11.886	1662	1666	1671	rBV	472918	458248	17.34%	1.697%
14	12.439	1756	1760	1766	rBV	119930	132369	5.01%	0.490%
15	12.569	1779	1782	1784	rBV	33923	34925	1.32%	0.129%
16	12.669	1796	1799	1806	rBV	175719	179444	6.79%	0.665%
17	12.745	1809	1812	1814	rBV	28974	30627	1.16%	0.113%
18	12.822	1823	1825	1831	rVB	62373	66808	2.53%	0.247%
19	12.945	1841	1846	1849	rBV	2843487	2642838	100.00%	9.789%
20	13.045	1861	1863	1872	rVB10	28194	48492	1.83%	0.180%
21	13.304	1903	1907	1912	rBV	791299	738647	27.95%	2.736%
22	13.416	1924	1926	1931	rVB6	34137	42968	1.63%	0.159%
23	13.674	1967	1970	1973	rBV	131755	133036	5.03%	0.493%
24	13.727	1976	1979	1986	rVB	261393	254150	9.62%	0.941%
25	13.839	1995	1998	2003	rBV	298214	392478	14.85%	1.454%
26	13.992	2020	2024	2027	rVB	1088109	932174	35.27%	3.453%
27	14.157	2050	2052	2055	rVB	55073	46337	1.75%	0.172%
28	14.321	2077	2080	2083	rBV	365658	345191	13.06%	1.279%
29	14.463	2101	2104	2109	rBV2	116181	139992	5.30%	0.519%
30	14.768	2152	2156	2159	rBV	145035	153605	5.81%	0.569%
31	15.098	2209	2212	2223	rVB	234244	378987	14.34%	1.404%
32	15.445	2266	2271	2274	rBV	691701	839521	31.77%	3.109%
33	15.480	2274	2277	2281	rVB	212630	270888	10.25%	1.003%
34	15.910	2346	2350	2355	rVB	193308	290590	11.00%	1.076%

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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	16.410	2431	2435	2445	rvB4	155011	311452	11.78%	1.154%
36	16.504	2447	2451	2457	rvB2	89621	151688	5.74%	0.562%

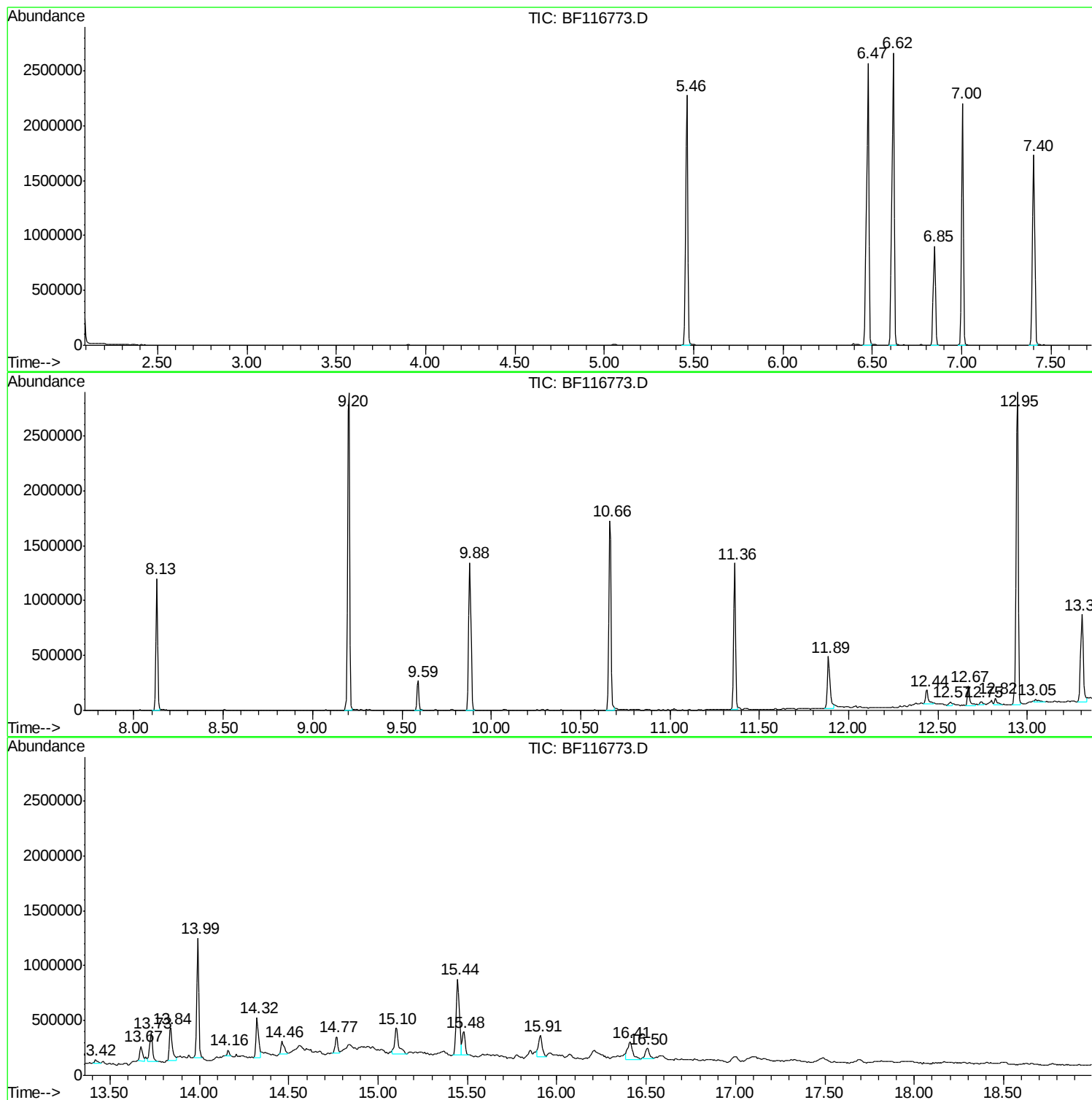
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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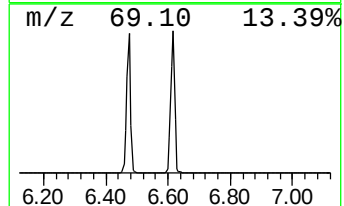
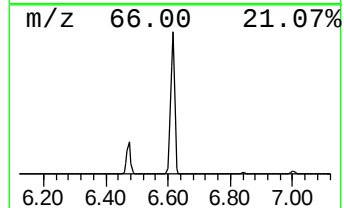
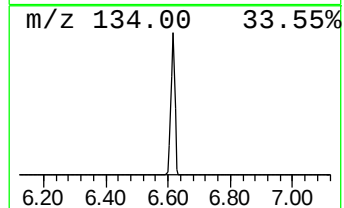
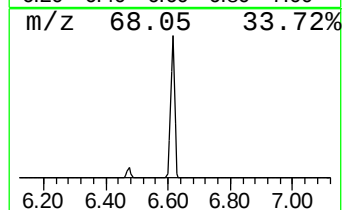
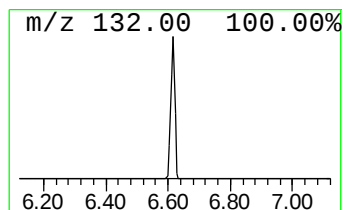
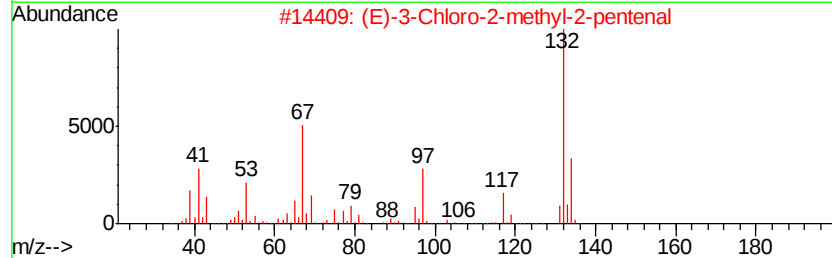
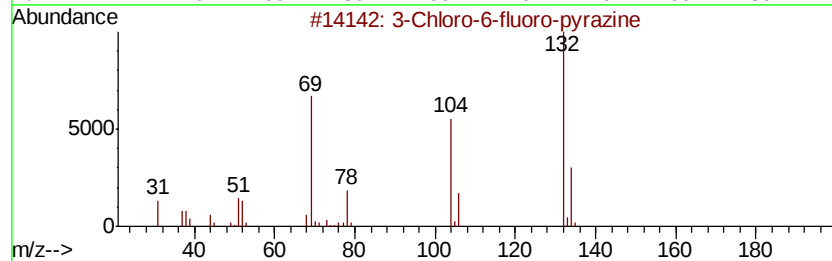
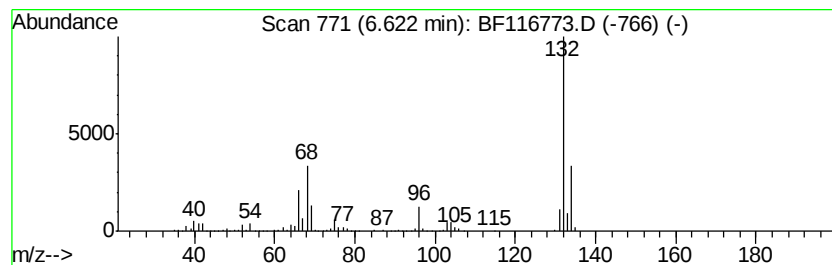
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown6.62 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.62	66.31 ng	2304180	1,4-Dichlorobenzene-d4	6.85

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		Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	12
4		2H-Cyclopentadlpyridazine, 2-me...	132	C8H8N2	022291-85-6	9
5		Xenon	132	Xe	007440-63-3	9



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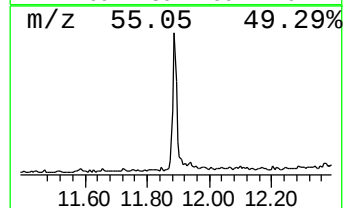
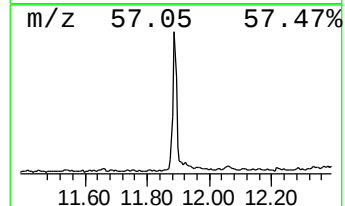
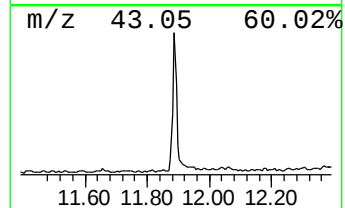
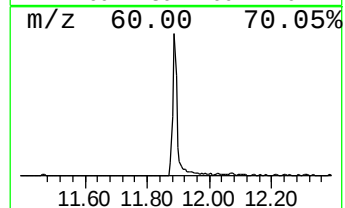
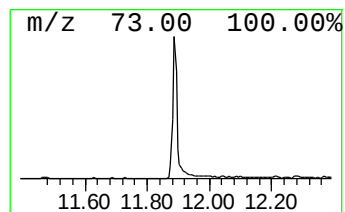
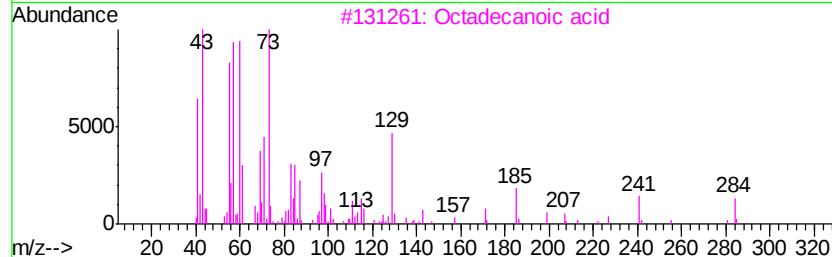
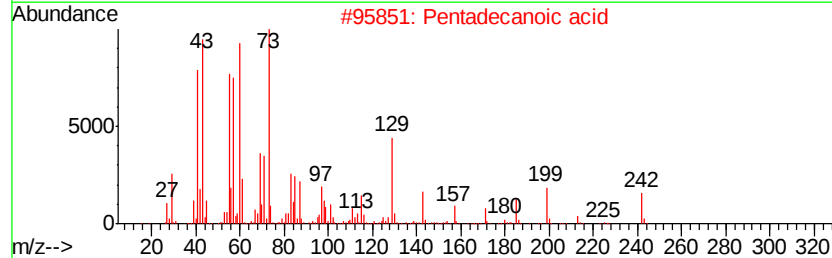
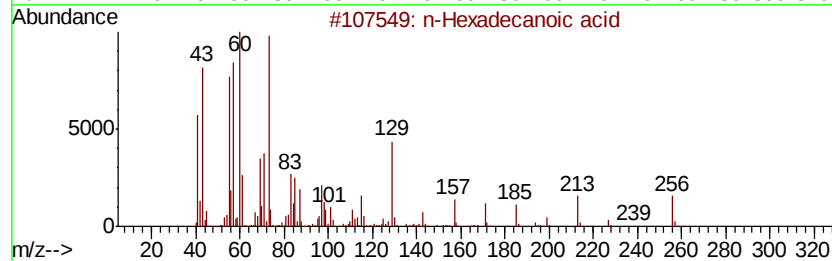
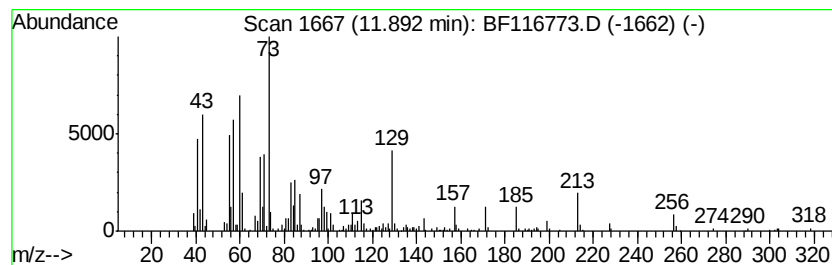
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.89	8.33 ng	458248	Phenanthrene-d10	11.36

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	87
3		Octadecanoic acid	284	C18H36O2	000057-11-4	83
4		Tridecanoic acid	214	C13H26O2	000638-53-9	72
5		Tetradecanoic acid	228	C14H28O2	000544-63-8	72



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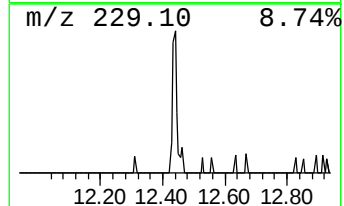
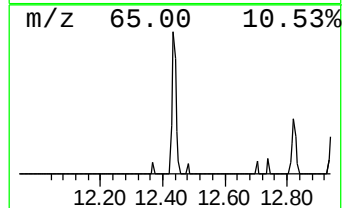
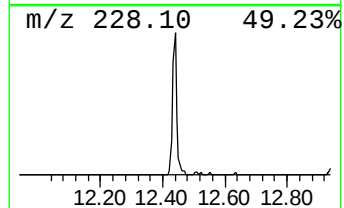
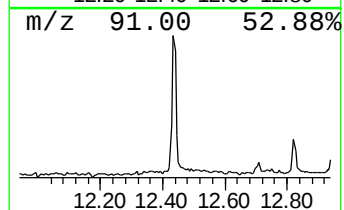
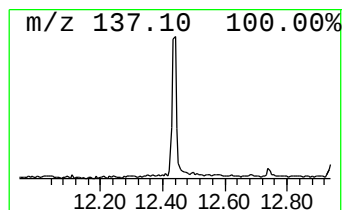
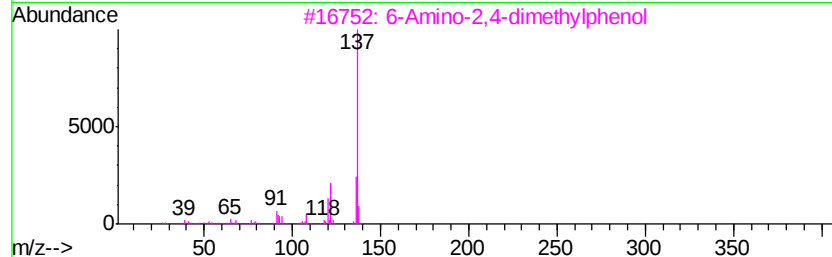
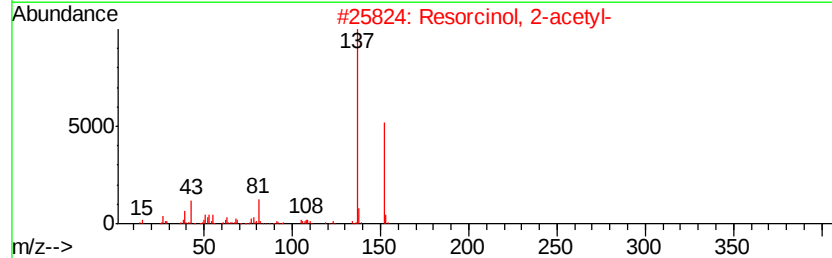
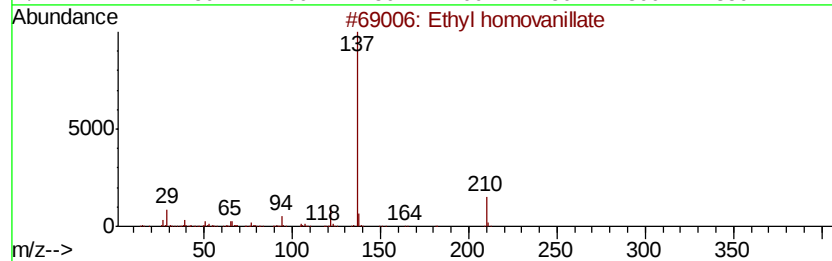
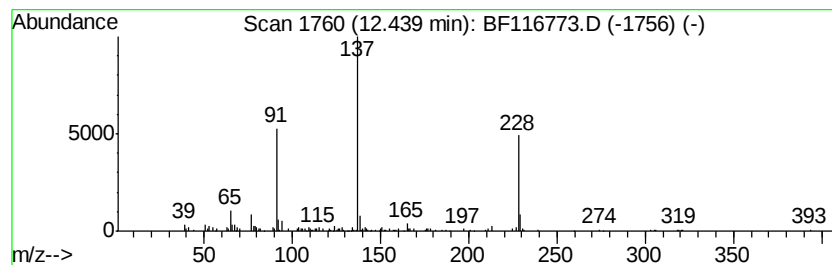
Instrument :
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown12.44 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.44	2.41 ng	132369	Phenanthrene-d10	11.36	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethyl homovanillate	210	C11H14O4	060563-13-5	43
2	Resorcinol, 2-acetyl-	152	C8H8O3	000699-83-2	38
3	6-Amino-2,4-dimethylphenol	137	C8H11NO	041458-65-5	38
4	(4-Methoxy-phenyl)-(2-nitrocyclo...	265	C14H19NO4	103077-68-5	37
5	Benzenemethanol, 4-methoxy-.alph...	251	C13H17NO4	103130-05-8	37



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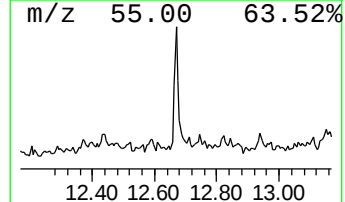
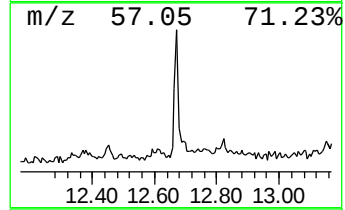
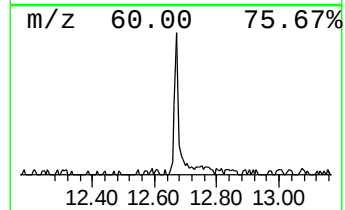
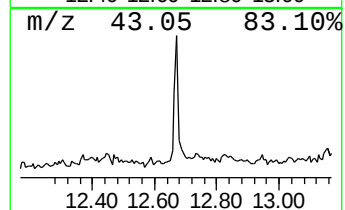
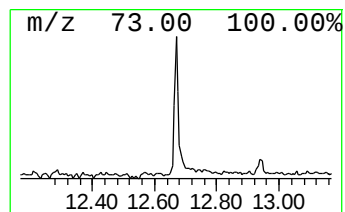
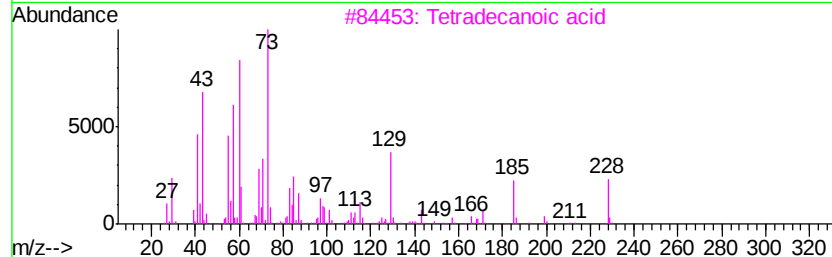
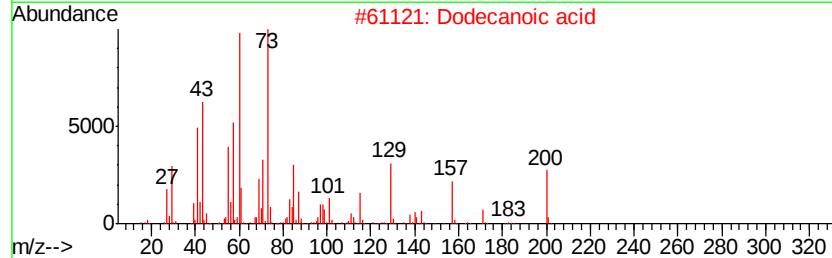
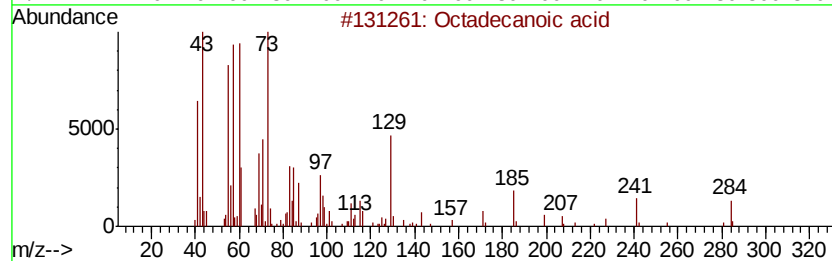
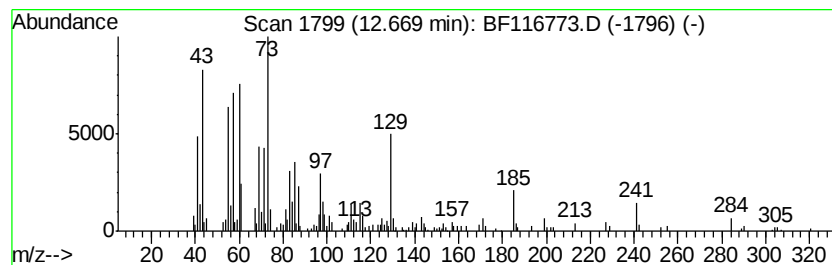
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Octadecanoic acid Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.67	3.26 ng	179444	Phenanthrene-d10	11.36

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		Dodecanoic acid	200	C12H24O2	000143-07-7	87
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	87
4		n-Decanoic acid	172	C10H20O2	000334-48-5	76
5		Undecanoic acid	186	C11H22O2	000112-37-8	46



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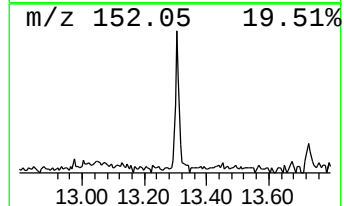
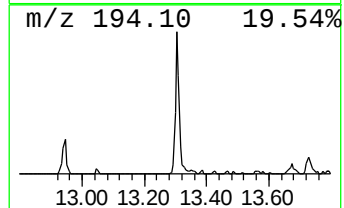
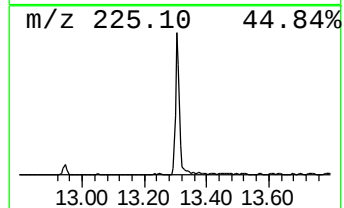
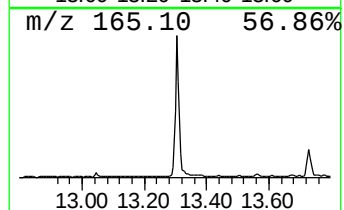
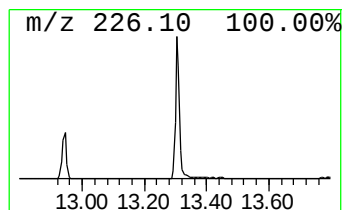
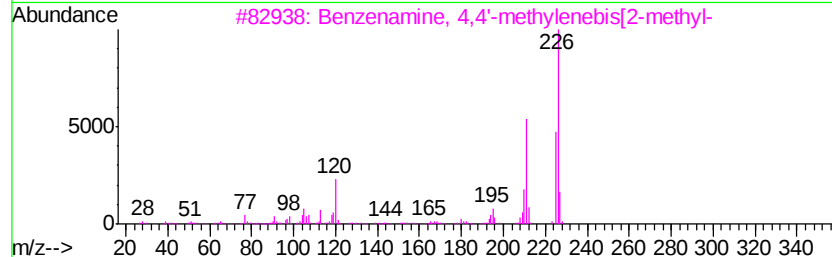
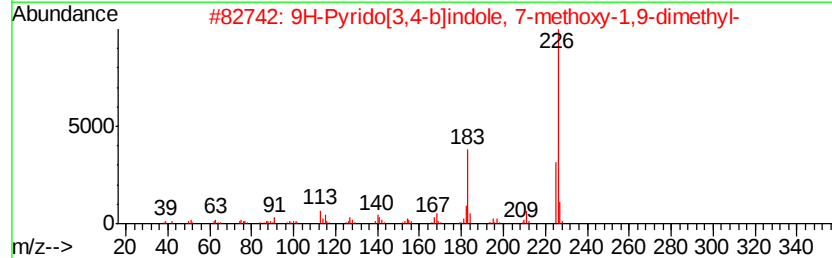
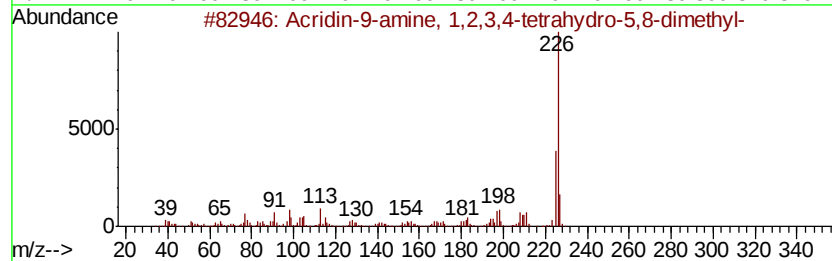
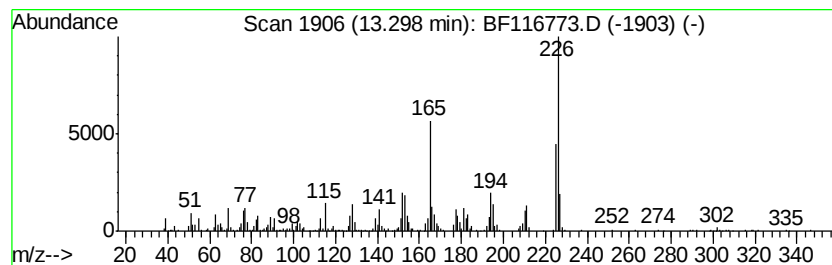
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Peak Number 6 Acridin-9-amine, 1,2,3,4-te... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.30	15.85 ng	738647	Chrysene-d12	13.99		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acridin-9-amine, 1,2,3,4-tetrahy...	226	C15H18N2	297758-19-1	70	
2	9H-Pyrido[3,4-b]indole, 7-methox...	226	C14H14N2O	143502-37-8	50	
3	Benzenamine, 4,4'-methylenebis[2...	226	C15H18N2	000838-88-0	49	
4	Acridin-9-amine, 1,2,3,4-tetrahy...	226	C15H18N2	1000300-57-6	49	
5	Benzenamine, N-[(4-nitrophenyl)m...	226	C13H10N2O2	000785-80-8	46	



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090319\
Data File : BF116773.D
Acq On : 3 Sep 2019 15:38
Operator : HP/JU
Sample : K4554-57
Misc :
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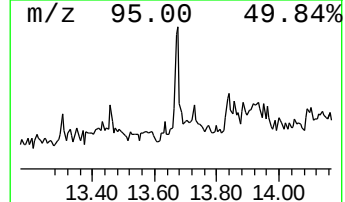
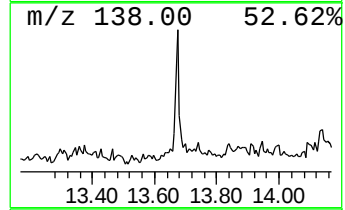
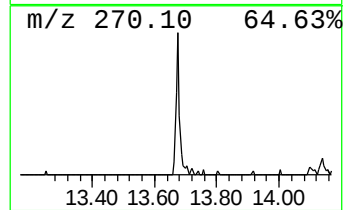
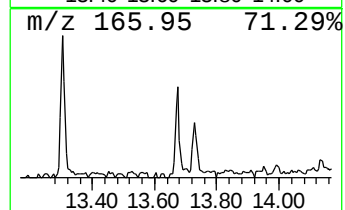
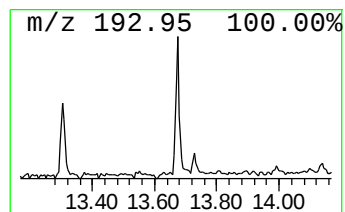
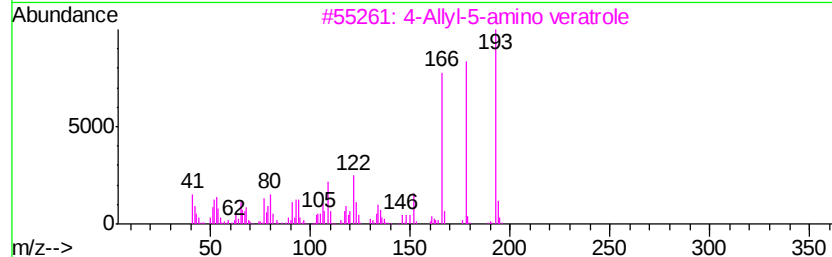
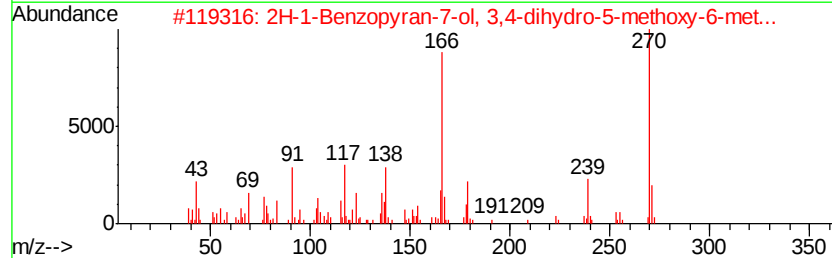
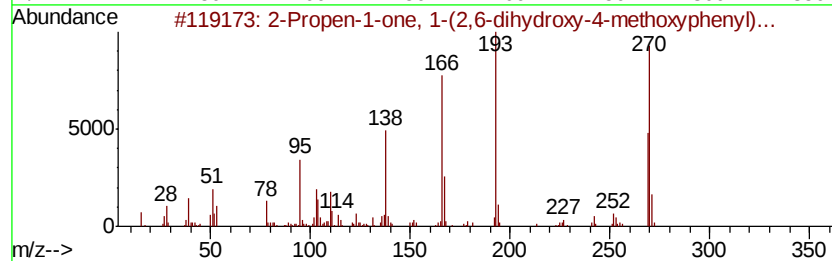
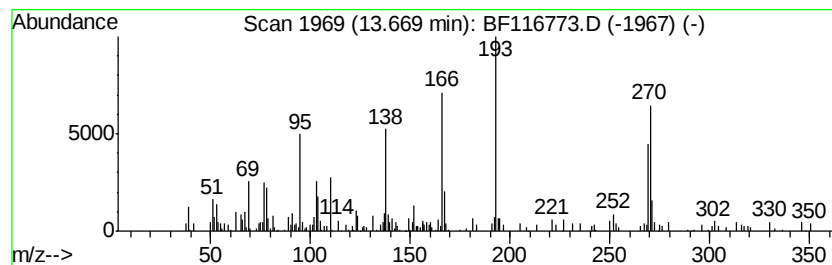
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ClientSampleId :
SB-1-(0-2)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF083019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 2-Propen-1-one, 1-(2,6-dihy... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.67	2.85 ng	133036	Chrysene-d12	13.99		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propen-1-one, 1-(2,6-dihydroxyv...	270	C16H14O4	018956-15-5	97	
2	2H-1-Benzopyran-7-ol, 3,4-dihydr...	270	C17H18O3	055670-26-3	30	
3	4-Allyl-5-amino veratrole	193	C11H15NO2	030058-51-6	30	
4	Benzene, 1,2-methylenedioxy-4-(3...	270	C14H10N2O4	191595-07-0	30	
5	3-(Benzylideneamino)carbazole	270	C19H14N2	014384-36-2	30	



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090319\
Data File : BF116773.D
Acq On : 3 Sep 2019 15:38
Operator : HP/JU
Sample : K4554-57
Misc :
ALS Vial : 11 Sample Multiplier: 1

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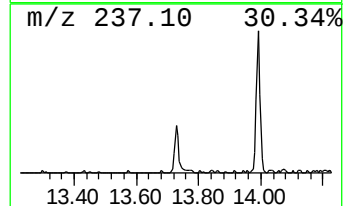
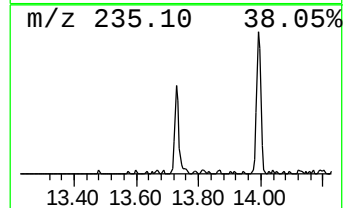
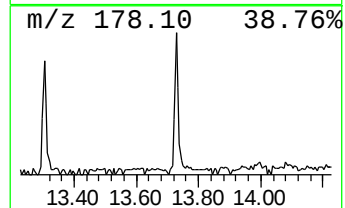
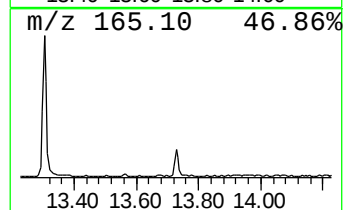
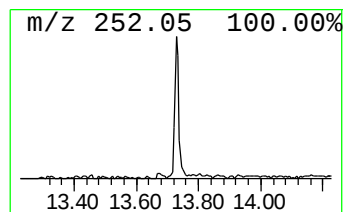
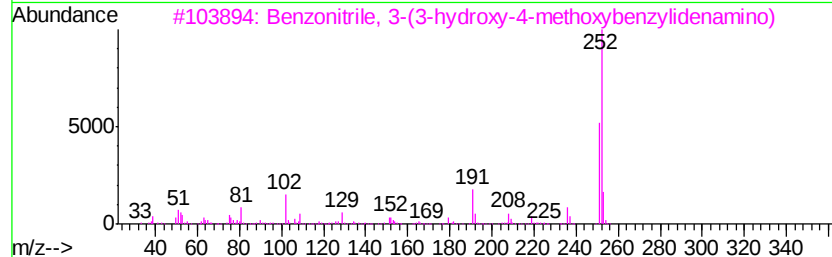
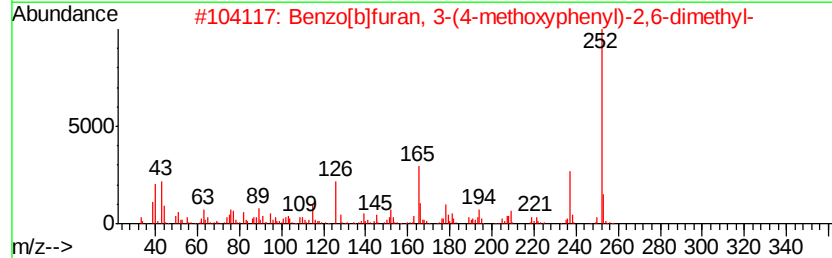
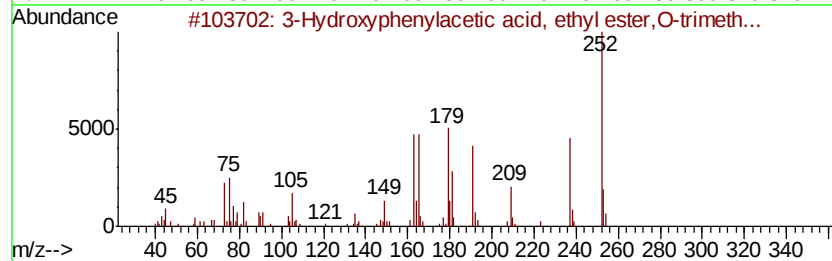
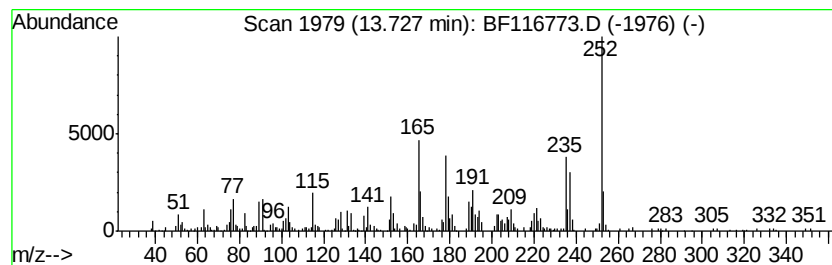
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 3-Hydroxyphenylacetic acid,... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.73	5.45 ng	254150	Chrysene-d12	13.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Hydroxyphenylacetic acid, ethyl...	252	C13H20O3Si	067903-48-4	64
2		Benzo[b]furan, 3-(4-methoxyphenyl)-2,6-dimethyl-	252	C17H16O2	074228-98-1	58
3		Benzonitrile, 3-(3-hydroxy-4-methoxybenzylideneamino)	252	C15H12N2O2	304683-23-6	46
4		4H-Benzopyran-4-one, 3-(4-methoxyphenyl)-2,6-dimethyl-	252	C16H12O3	001621-58-5	42
5		3H-Pyrrolo[3,2-f]quinoline, 1,2,3,4-tetrahydro-	252	C17H20N2	1000302-73-4	41



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090319\
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Operator : HP/JU
Sample : K4554-57
Misc :
ALS Vial : 11 Sample Multiplier: 1

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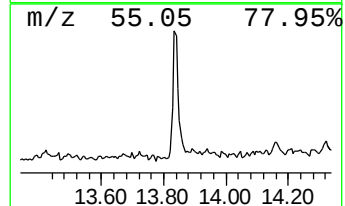
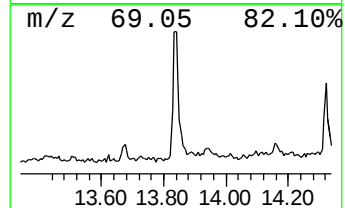
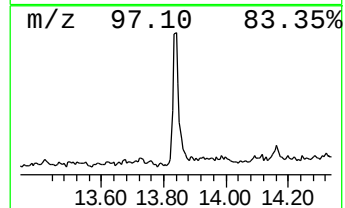
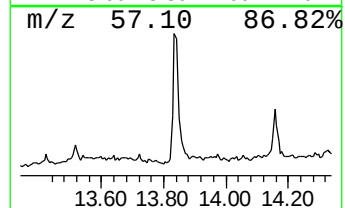
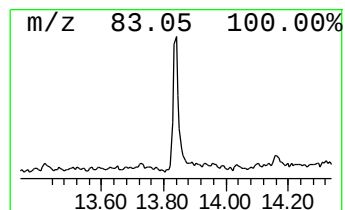
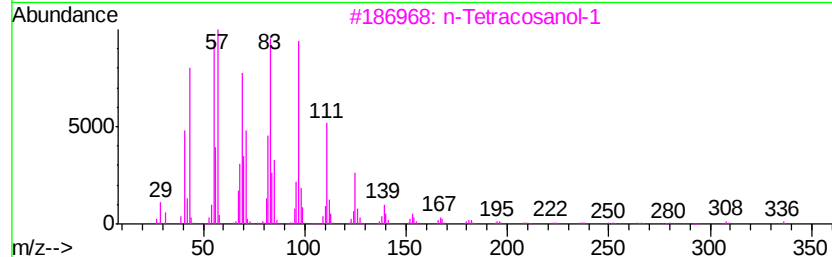
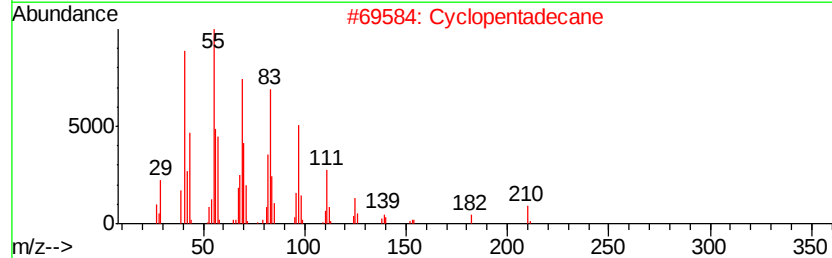
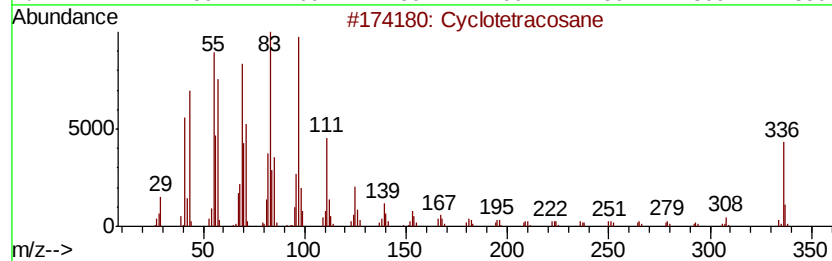
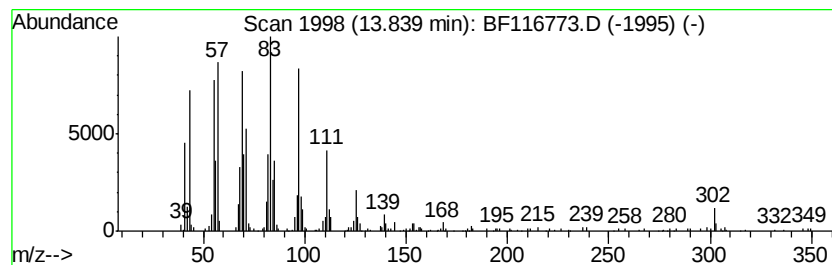
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Cyclotetracosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.84	8.42 ng	392478	Chrysene-d12	13.99

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclotetracosane	336	C24H48	000297-03-0	99
2		Cyclopentadecane	210	C15H30	000295-48-7	95
3		n-Tetracosanol-1	354	C24H50O	000506-51-4	94
4		1-Heptacosanol	396	C27H56O	002004-39-9	93
5		1-Heneicosanol	312	C21H44O	015594-90-8	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090319\
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ClientSampleId :
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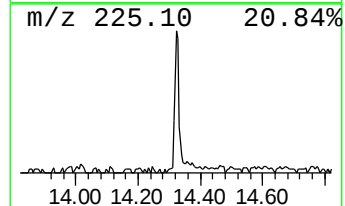
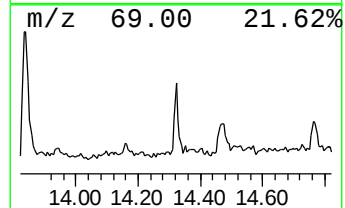
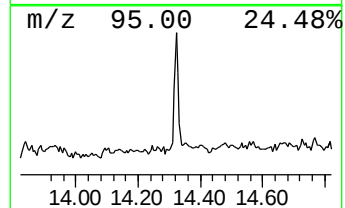
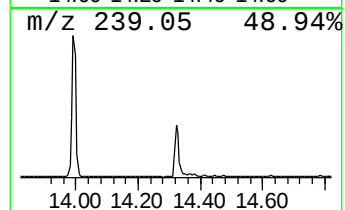
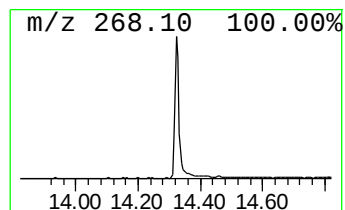
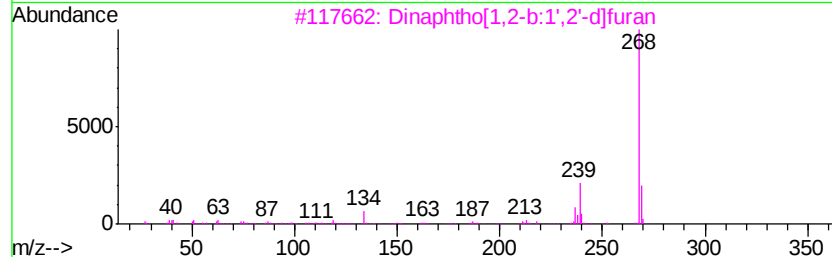
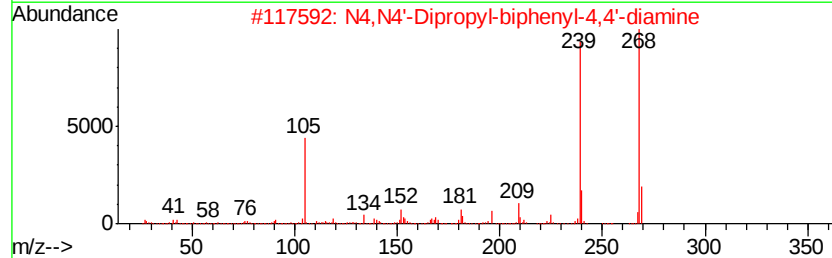
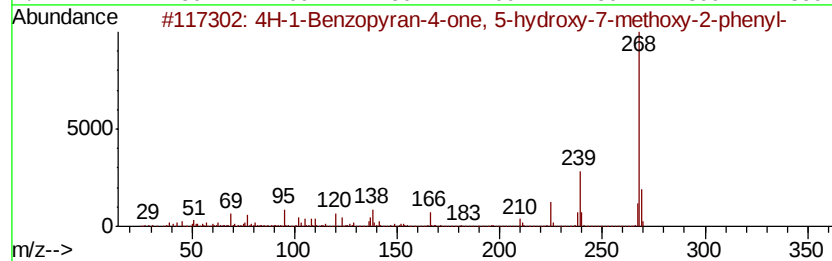
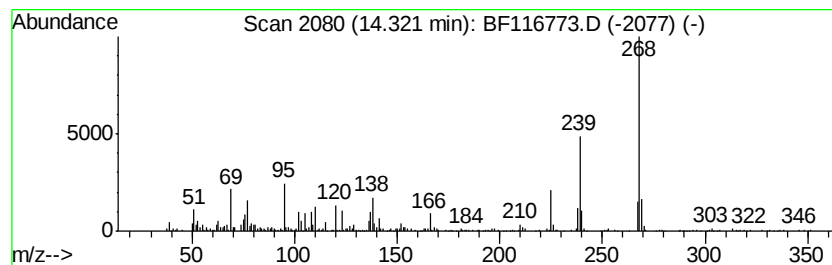
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 4H-1-Benzopyran-4-one, 5-hy... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.32	7.41 ng	345191	Chrysene-d12	13.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4H-1-Benzopyran-4-one, 5-hydroxy...	268	C16H12O4	000520-28-5	93
2		N4,N4'-Dipropyl-biphenyl-4,4'-di...	268	C18H24N2	017576-22-6	53
3		Dinaphtho[1,2-b:1',2'-d]furan	268	C20H12O	000207-93-2	53
4		Benzo(a)pyren-7-ol	268	C20H12O	037994-82-4	52
5		3-Hydroxy-cis-D-homoestra-1,3,5(...	268	C18H20O2	021547-08-0	49



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090319\
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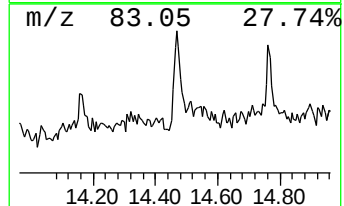
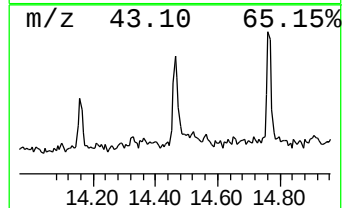
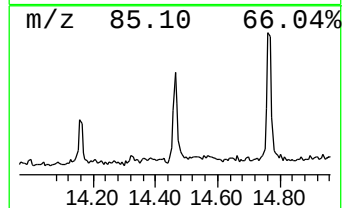
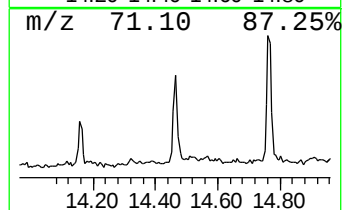
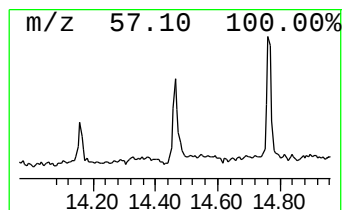
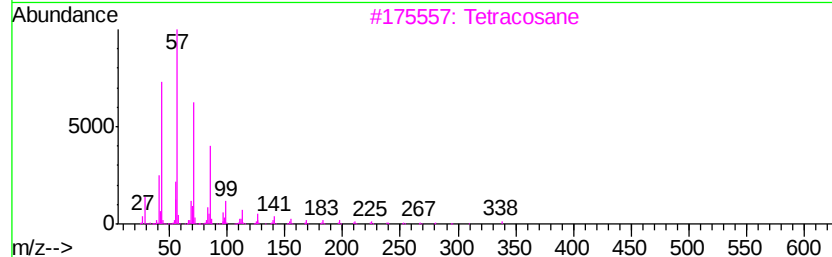
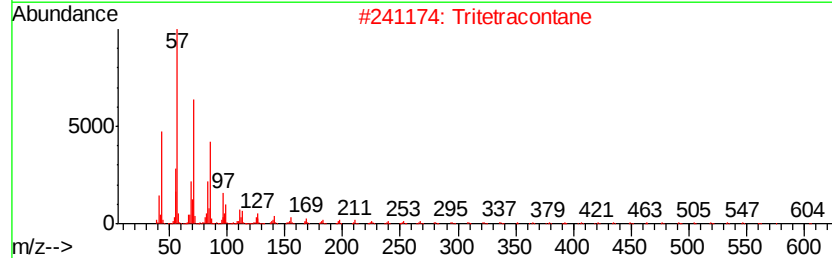
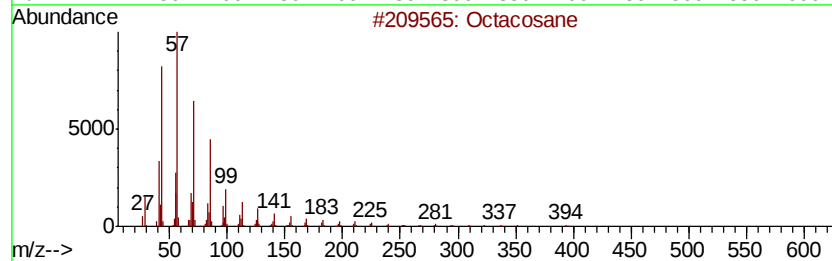
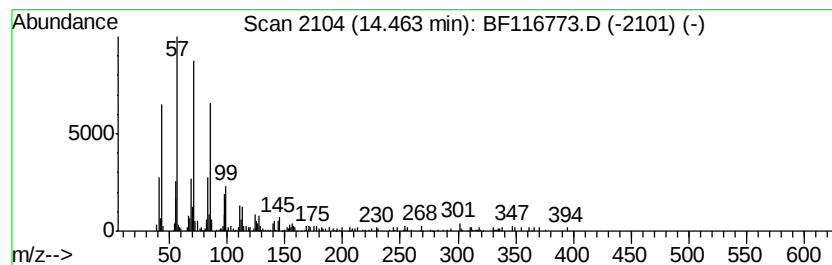
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Octacosane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.46	3.00 ng	139992	Chrysene-d12	13.99

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosane	394	C28H58	000630-02-4	91
2			Tritetracontane	605	C43H88	007098-21-7	86
3			Tetracosane	338	C24H50	000646-31-1	83
4			Heptadecane, 8-methyl-	254	C18H38	013287-23-5	76
5			Tricosane, 2-methyl-	338	C24H50	001928-30-9	74



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090319\
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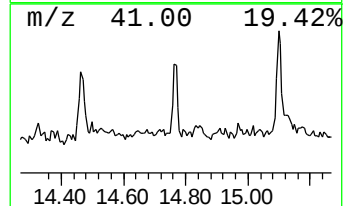
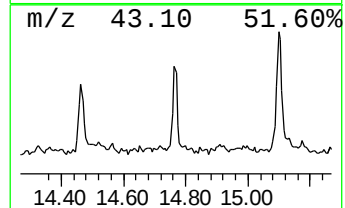
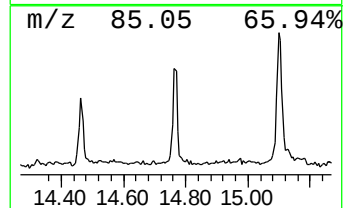
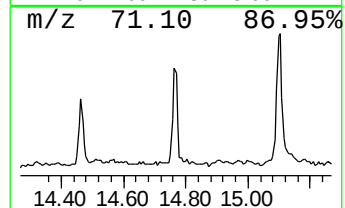
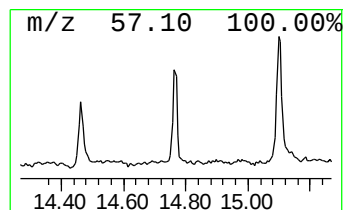
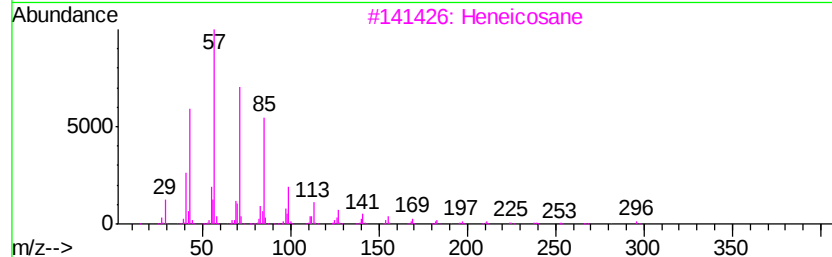
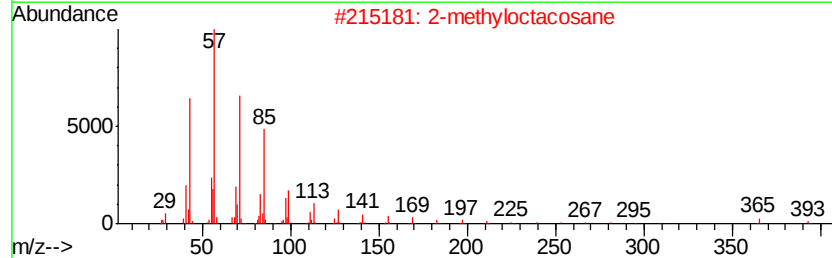
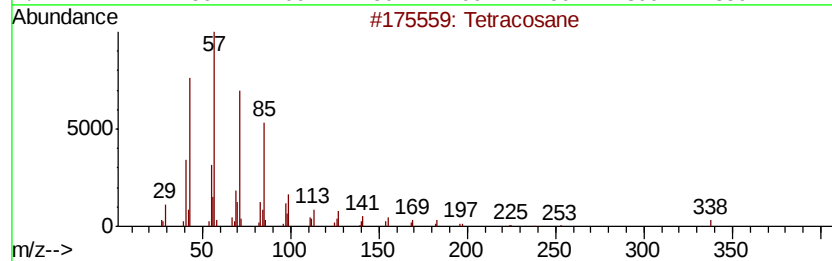
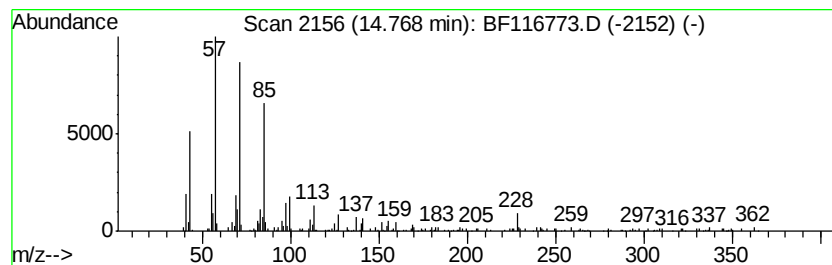
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 12 Tetracosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.77	3.66 ng	153605	Perylene-d12	15.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane	338	C24H50	000646-31-1	90
2		2-methyloctacosane	408	C29H60	1000376-72-8	87
3		Heneicosane	296	C21H44	000629-94-7	87
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	86
5		Eicosane	282	C20H42	000112-95-8	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090319\
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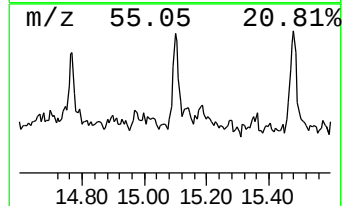
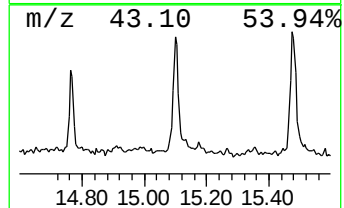
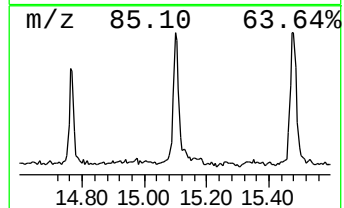
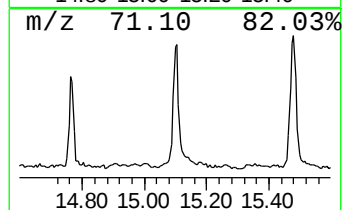
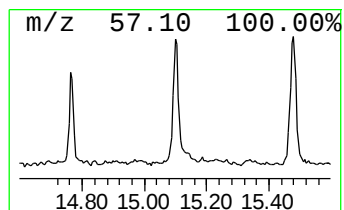
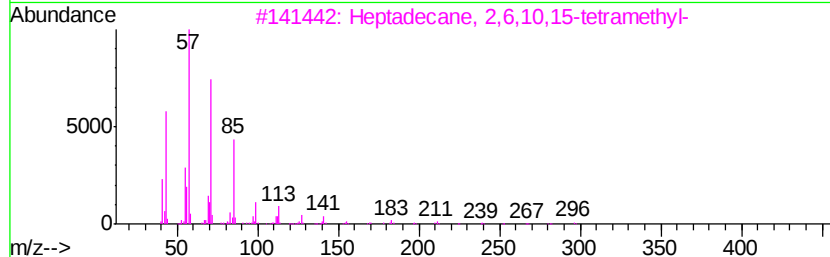
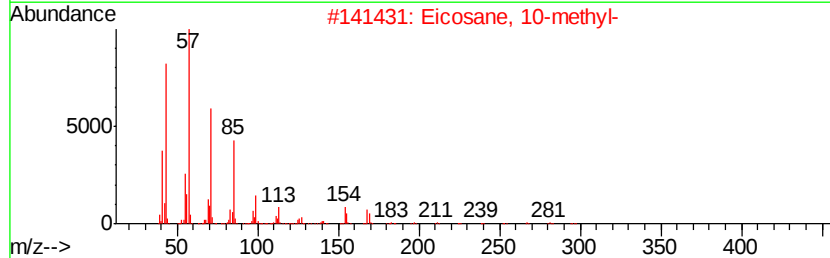
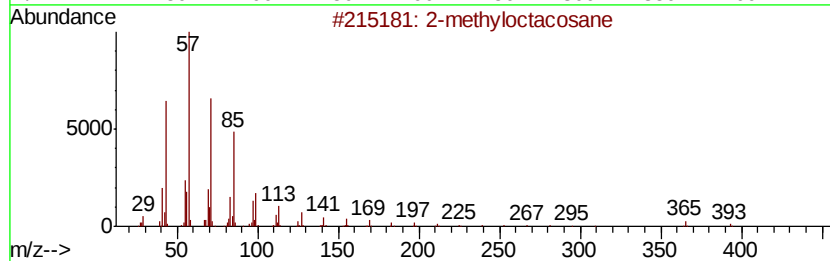
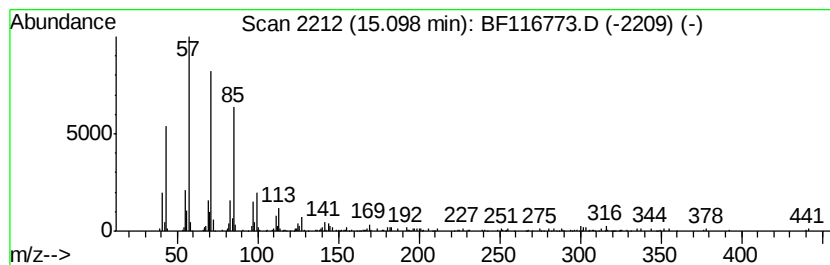
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 2-methyloctacosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.10	9.03 ng	378987	Perylene-d12	15.44

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-methyloctacosane	408	C29H60	1000376-72-8	91
2			Eicosane, 10-methyl-	296	C21H44	054833-23-7	90
3			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	89
4			Octadecane, 1-iodo-	380	C18H37I	000629-93-6	87
5			Hexacosane	366	C26H54	000630-01-3	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090319\
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 Acq On : 3 Sep 2019 15:38
 Operator : HP/JU
 Sample : K4554-57
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

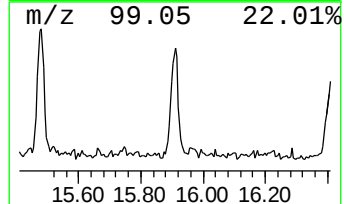
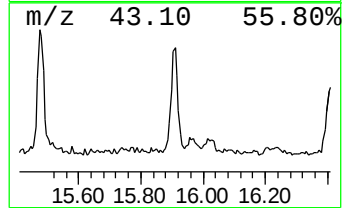
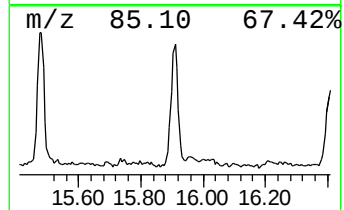
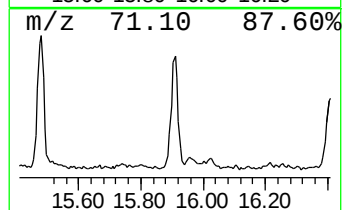
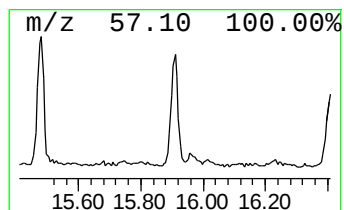
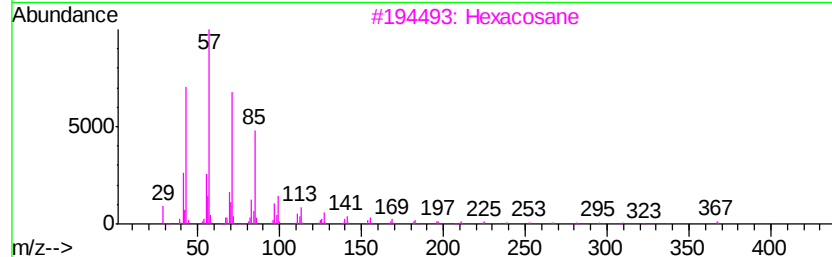
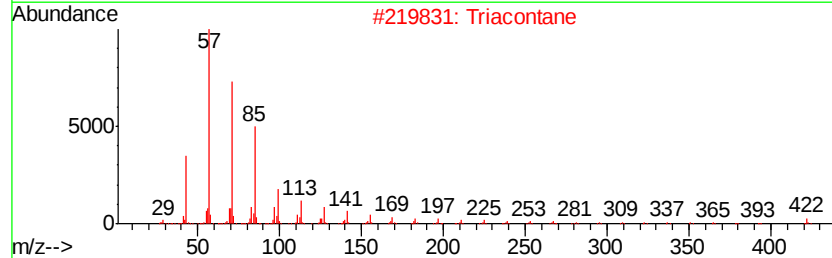
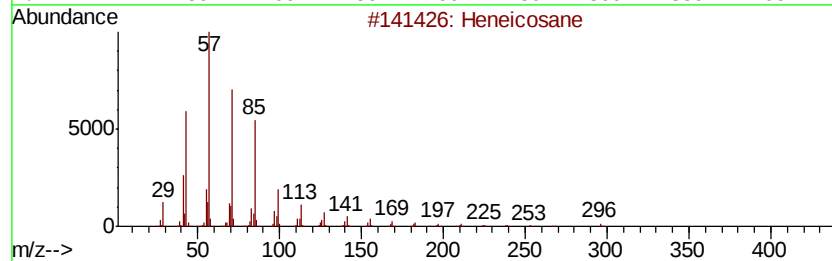
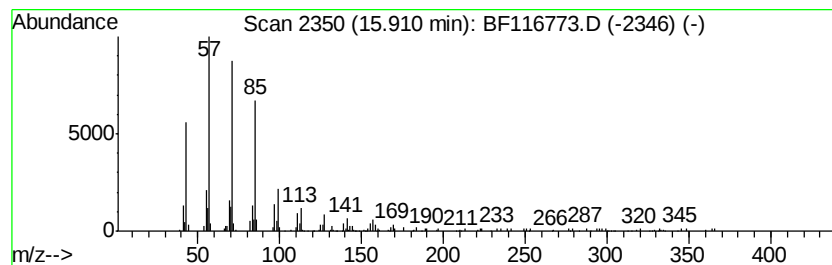
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 14 Heneicosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.
15.91	6.92 ng	290590	Perylene-d12		15.44
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	90
2	triacontane	422	C30H62	000638-68-6	90
3	Hexacosane	366	C26H54	000630-01-3	87
4	2-Bromotetradecane	276	C14H29Br	074036-95-6	86
5	Octacosane	394	C28H58	000630-02-4	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090319\
Data File : BF116773.D
Acq On : 3 Sep 2019 15:38
Operator : HP/JU
Sample : K4554-57
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-1-(0-2)

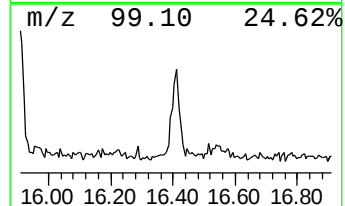
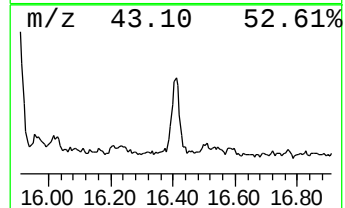
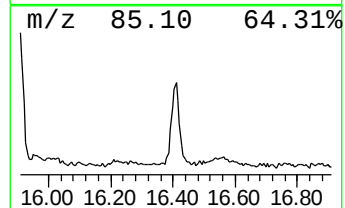
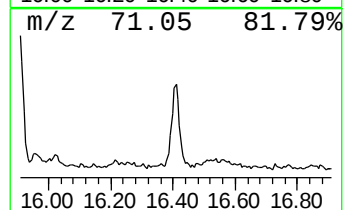
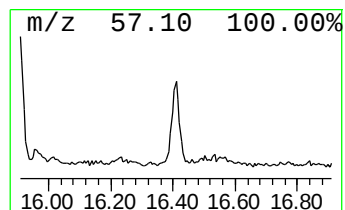
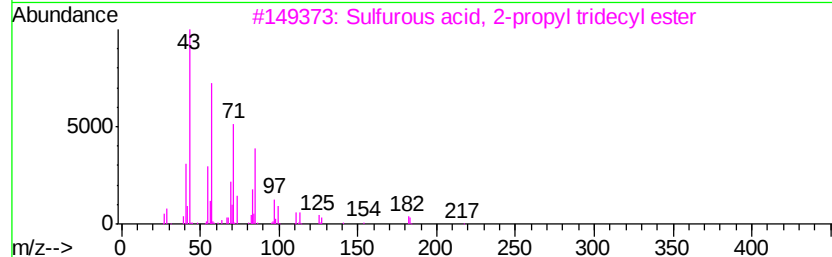
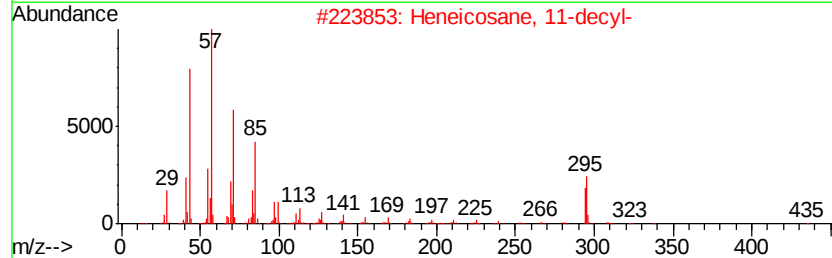
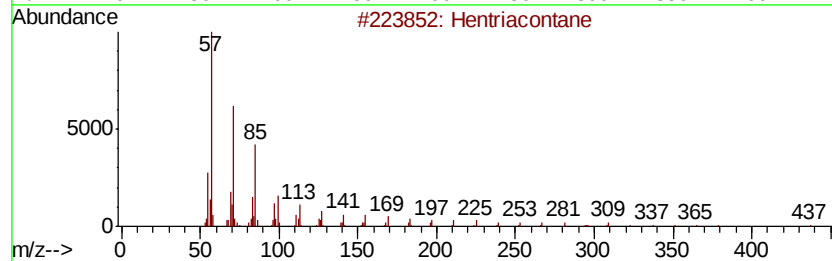
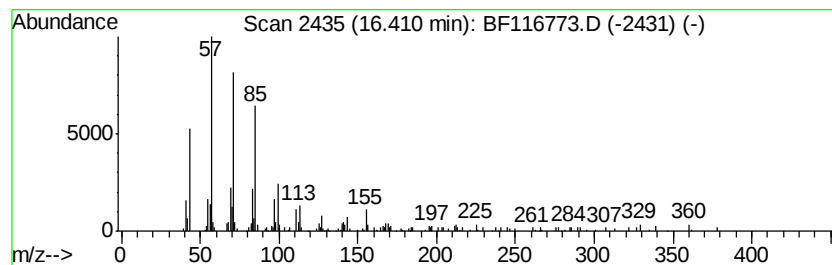
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Hentriacontane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.41	7.42 ng	311452	Perylene-d12	15.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hentriacontane	437	C31H64	000630-04-6	91
2		Heneicosane, 11-decyl-	437	C31H64	055320-06-4	86
3		Sulfurous acid, 2-propyl tridecyl...	306	C16H34O3S	1000309-12-4	86
4		Heneicosane	296	C21H44	000629-94-7	80
5		2-methyloctacosane	408	C29H60	1000376-72-8	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090319\
Data File : BF116773.D
Acq On : 3 Sep 2019 15:38
Operator : HP/JU
Sample : K4554-57
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-1-(0-2)

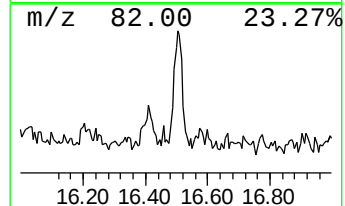
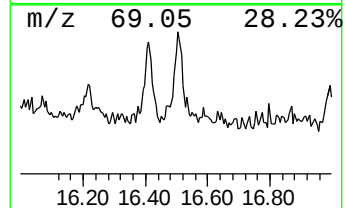
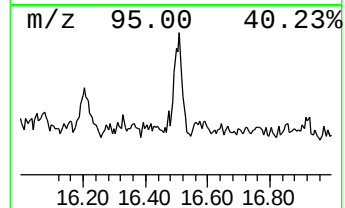
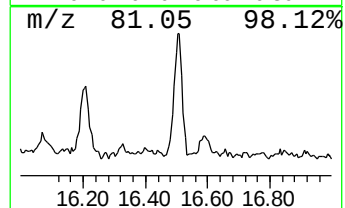
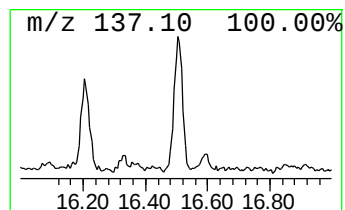
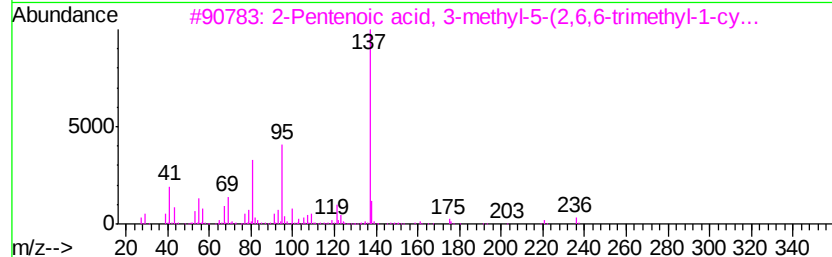
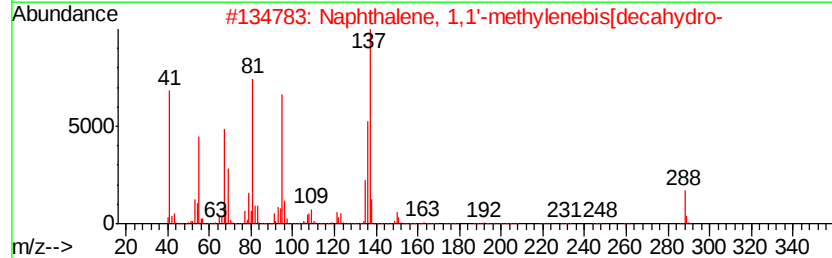
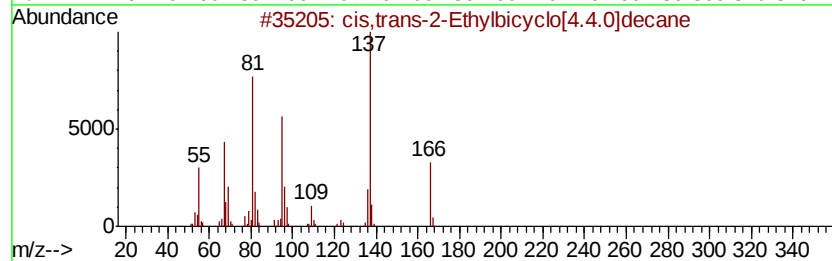
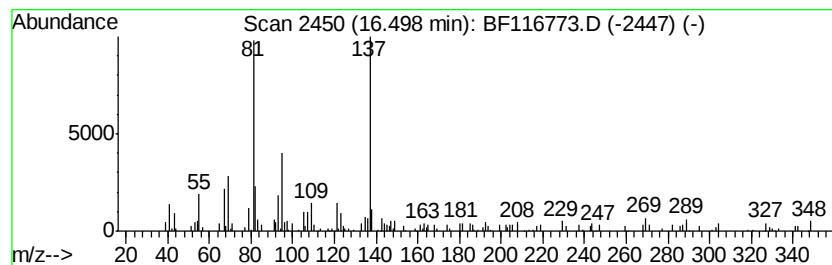
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 cis,trans-2-Ethylbicyclo[4.4.0]d... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.50	3.61 ng	151688	Perylene-d12	15.44

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			cis,trans-2-Ethylbicyclo[4.4.0]d...	166	C12H22	066660-38-6	59
2			Naphthalene, 1,1'-methylenebis[d...	288	C21H36	055125-02-5	50
3			2-Pentenoic acid, 3-methyl-5-(2,...	236	C15H24O2	1000196-81-2	43
4			Cyclohexene, 1-butyl-	138	C10H18	003282-53-9	38
5			(4-Methoxy-phenyl)-(2-nitrocyclo...	265	C14H19NO4	103077-68-5	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF090319\
 Data File : BF116773.D
 Acq On : 3 Sep 2019 15:38
 Operator : HP/JU
 Sample : K4554-57
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-1-(0-2)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF083019.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
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n-Hexadecanoic acid	11.89	8.3	ng	458248	4	11.36	1099640	20.0
unknown12.44	12.44	2.4	ng	132369	4	11.36	1099640	20.0
Octadecanoic acid	12.67	3.3	ng	179444	4	11.36	1099640	20.0
Acridin-9-amine, ...	13.30	15.9	ng	738647	5	13.99	932174	20.0
2-Propen-1-one, 1...	13.67	2.9	ng	133036	5	13.99	932174	20.0
3-Hydroxyphenylac...	13.73	5.5	ng	254150	5	13.99	932174	20.0
Cyclotetracosane	13.84	8.4	ng	392478	5	13.99	932174	20.0
4H-1-Benzopyran-4...	14.32	7.4	ng	345191	5	13.99	932174	20.0
Octacosane	14.46	3.0	ng	139992	5	13.99	932174	20.0
Tetracosane	14.77	3.7	ng	153605	6	15.44	839521	20.0
2-methyloctacosane	15.10	9.0	ng	378987	6	15.44	839521	20.0
Heneicosane	15.91	6.9	ng	290590	6	15.44	839521	20.0
Hentriacontane	16.41	7.4	ng	311452	6	15.44	839521	20.0
cis,trans-2-Ethyl...	16.50	3.6	ng	151688	6	15.44	839521	20.0