

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110118\
 Data File : BF110378.D
 Acq On : 1 Nov 2018 20:07
 Operator : JU/SJ
 Sample : J5192-05 2X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AU-05-103118-A

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-BF102318.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	2.257	24	29	48	rVB	152859	276575	27.72%	2.160%
2	5.181	523	526	534	rBV	24546	29150	2.92%	0.228%
3	5.575	589	593	611	rBV	760300	704415	70.61%	5.502%
4	6.575	759	763	781	rBV	739524	731356	73.31%	5.712%
5	6.722	784	788	802	rBV	880293	751254	75.31%	5.868%
6	6.957	824	828	831	rBV	503199	443362	44.44%	3.463%
7	7.110	850	854	868	rBV	692266	590891	59.23%	4.615%
8	7.516	919	923	934	rBV	433604	436600	43.77%	3.410%
9	8.239	1042	1046	1066	rBV	542509	541534	54.28%	4.230%
10	9.310	1224	1228	1238	rBV	812349	723519	72.53%	5.651%
11	9.704	1290	1295	1301	rBV	70857	76378	7.66%	0.597%
12	9.998	1341	1345	1352	rVB	503033	485907	48.71%	3.795%
13	10.786	1475	1479	1488	rBV	337430	323186	32.40%	2.524%
14	10.892	1492	1497	1502	rVB2	11289	17483	1.75%	0.137%
15	11.486	1594	1598	1601	rBV	490597	422883	42.39%	3.303%
16	11.510	1601	1602	1607	rVB	71303	45070	4.52%	0.352%
17	11.857	1658	1661	1664	rBV	409622	315955	31.67%	2.468%
18	11.992	1680	1684	1687	rBV	41234	39002	3.91%	0.305%
19	12.021	1687	1689	1693	rVB2	15499	14882	1.49%	0.116%
20	12.104	1700	1703	1706	rBV2	18241	21440	2.15%	0.167%
21	12.545	1774	1778	1780	rBV	1060804	945247	94.75%	7.383%
22	12.568	1780	1782	1789	rVV	1167590	997583	100.00%	7.792%
23	12.651	1793	1796	1799	rVB	297953	234312	23.49%	1.830%
24	12.704	1799	1805	1809	rBV	156874	138509	13.88%	1.082%
25	12.933	1838	1844	1850	rVB	134581	153763	15.41%	1.201%
26	13.068	1859	1867	1870	rBV	690086	620700	62.22%	4.848%
27	13.151	1877	1881	1885	rBV5	10966	19002	1.90%	0.148%
28	13.245	1892	1897	1898	rBV4	15645	21333	2.14%	0.167%
29	13.263	1898	1900	1904	rVV2	22991	34781	3.49%	0.272%
30	13.298	1904	1906	1909	rVV2	21322	20981	2.10%	0.164%
31	13.374	1913	1919	1925	rBV4	40538	57330	5.75%	0.448%
32	13.486	1935	1938	1942	rBV4	29996	34467	3.46%	0.269%
33	13.527	1942	1945	1951	rVB2	64929	78507	7.87%	0.613%
34	13.615	1957	1960	1964	rBV3	14713	18583	1.86%	0.145%

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

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

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	13.774	1984	1987	1988	rBV	17461	16991	1.70%	0.133%
36	13.804	1990	1992	1996	rVB4	14911	19597	1.96%	0.153%
37	13.880	2002	2005	2008	rBV3	15404	19163	1.92%	0.150%
38	13.951	2013	2017	2020	rVV3	58932	68211	6.84%	0.533%
39	13.980	2020	2022	2029	rVB	25234	31862	3.19%	0.249%
40	14.039	2029	2032	2035	rBV2	36733	39872	4.00%	0.311%
41	14.133	2043	2048	2051	rBV	646287	658942	66.05%	5.147%
42	14.157	2051	2052	2060	rVB	102897	95805	9.60%	0.748%
43	14.239	2064	2066	2068	rBV	13153	13037	1.31%	0.102%
44	14.262	2068	2070	2074	rBV	32993	30199	3.03%	0.236%
45	14.515	2112	2113	2117	rVB	18064	15866	1.59%	0.124%
46	14.568	2119	2122	2131	rVB2	54673	68240	6.84%	0.533%
47	14.886	2172	2176	2179	rVB	74578	75980	7.62%	0.593%
48	15.033	2199	2201	2206	rVB5	16568	19072	1.91%	0.149%
49	15.204	2226	2230	2233	rBV	61764	101331	10.16%	0.791%
50	15.239	2233	2236	2240	rVB2	101023	122399	12.27%	0.956%
51	15.521	2281	2284	2289	rVB	42909	55045	5.52%	0.430%
52	15.586	2292	2295	2299	rVV	50039	60656	6.08%	0.474%
53	15.656	2299	2307	2311	rBV2	389052	607833	60.93%	4.747%
54	16.092	2376	2381	2386	rBV	74723	118439	11.87%	0.925%
55	16.374	2426	2429	2433	rVB4	20455	24333	2.44%	0.190%
56	16.621	2467	2471	2483	rVB2	50298	100262	10.05%	0.783%
57	17.992	2699	2704	2717	rVB9	27832	74256	7.44%	0.580%

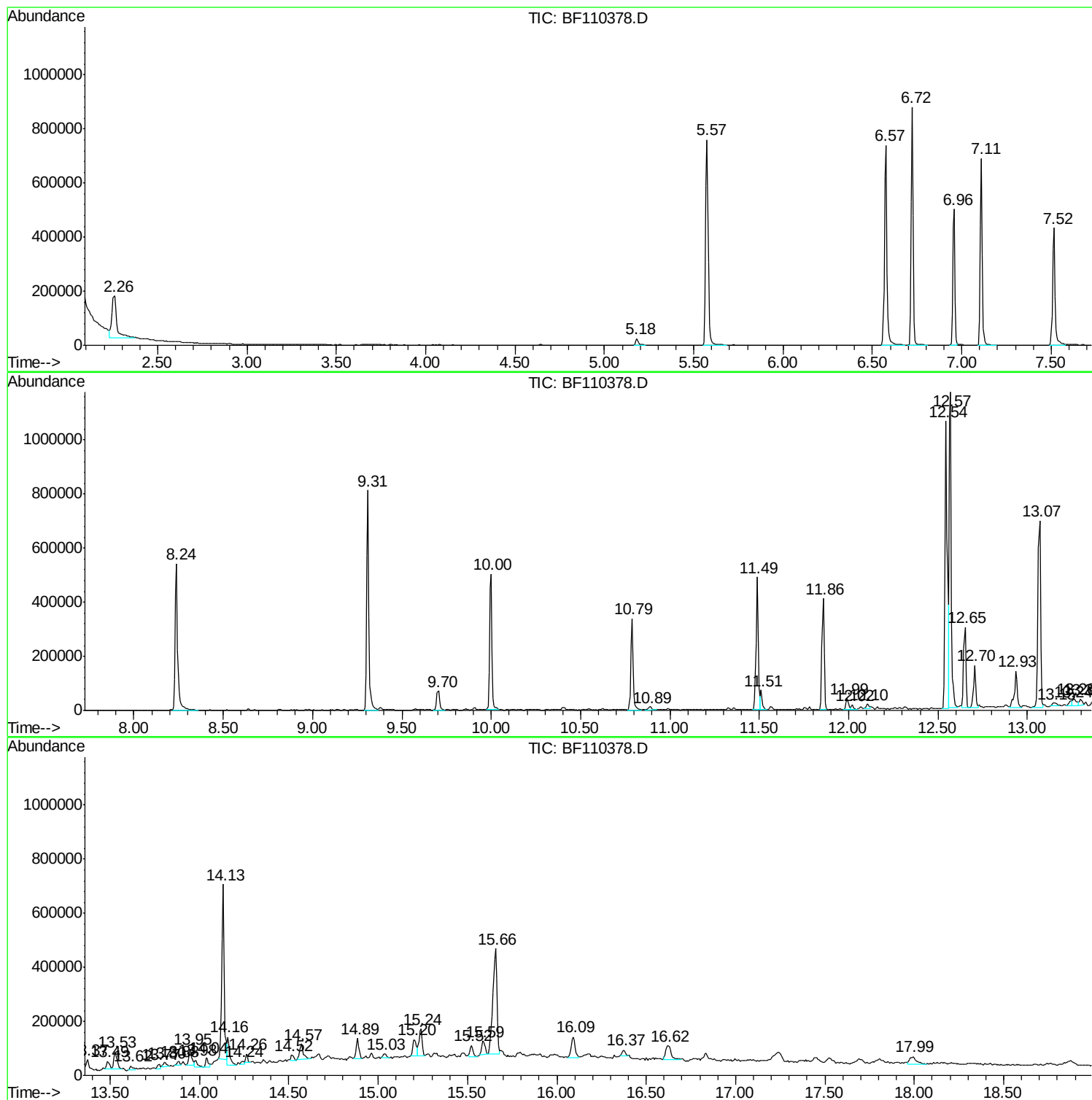
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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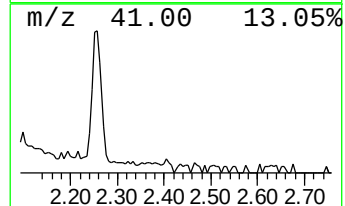
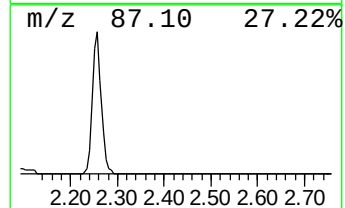
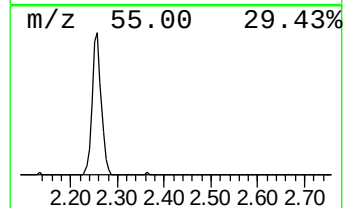
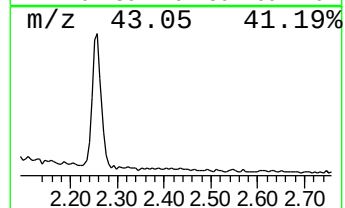
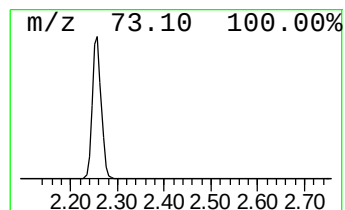
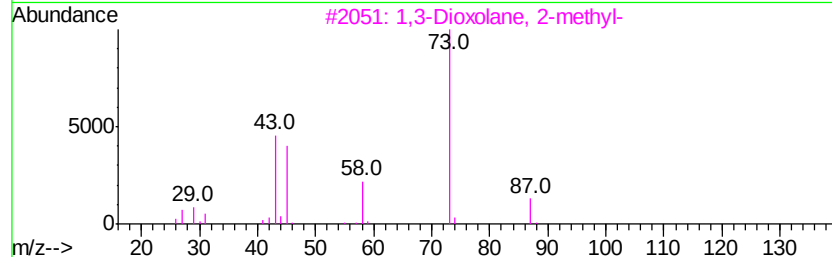
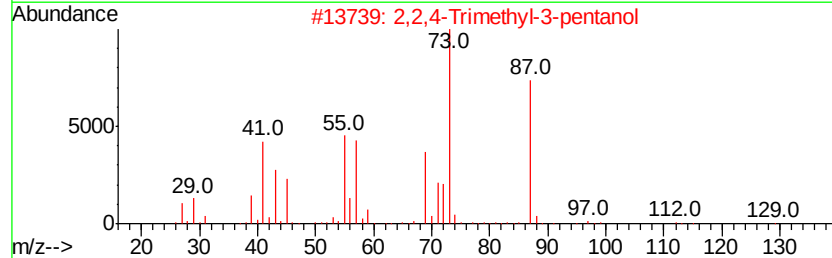
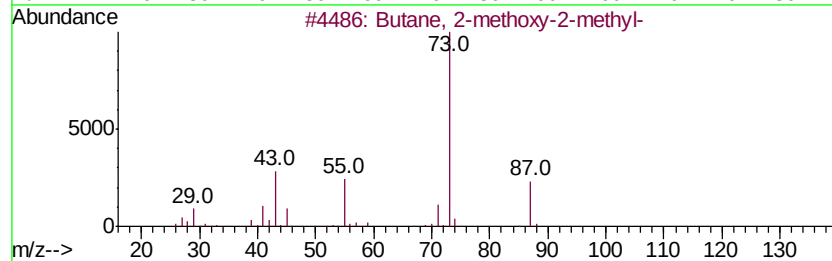
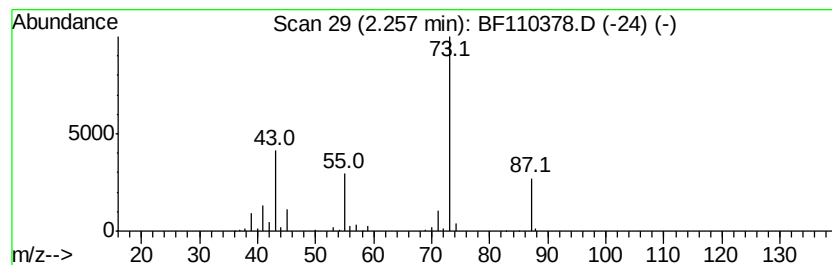
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.26	12.48 ng	276575	1,4-Dichlorobenzene-d4	6.96

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	37
4		Silane, tetramethyl-	88	C4H12Si	000075-76-3	25
5		Boronic acid, ethyl-, dimethyl e...	102	C4H11B02	007318-82-3	12



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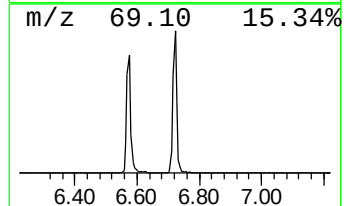
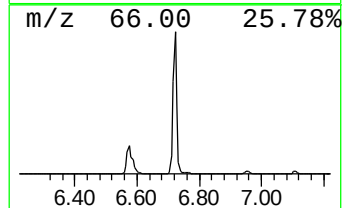
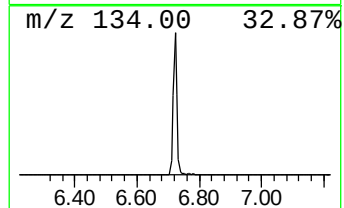
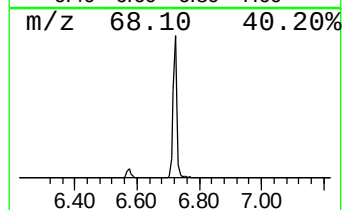
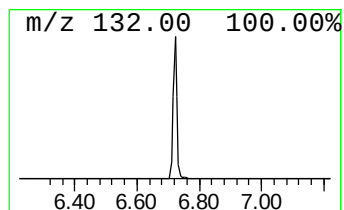
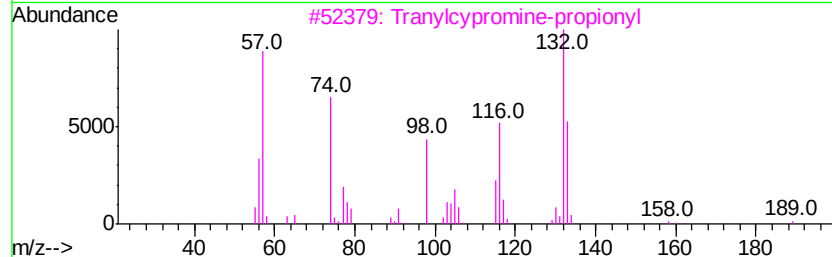
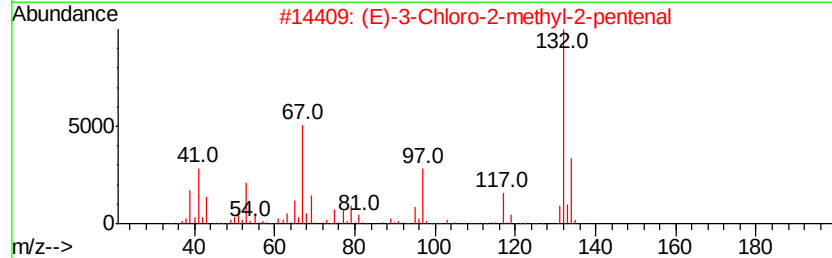
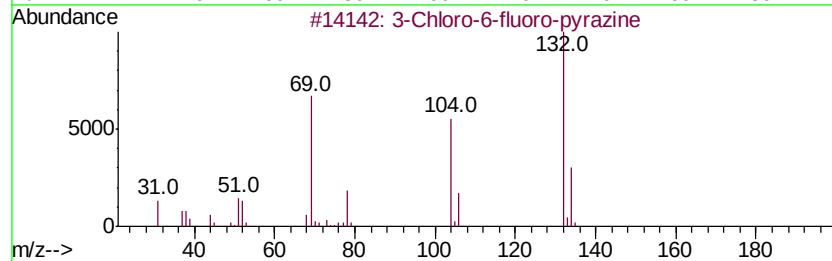
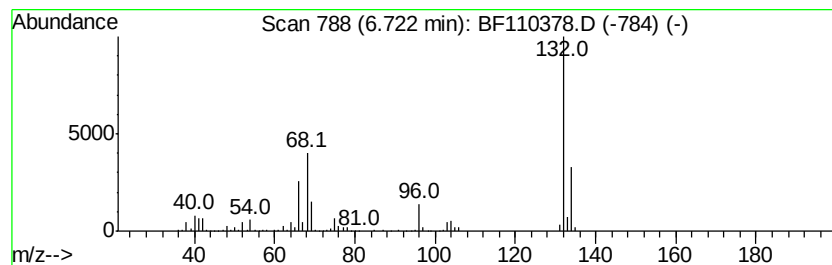
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.72 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.72	33.89 ng	751254	1,4-Dichlorobenzene-d4	6.96

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	17
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
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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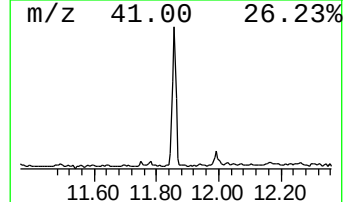
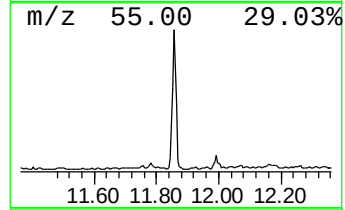
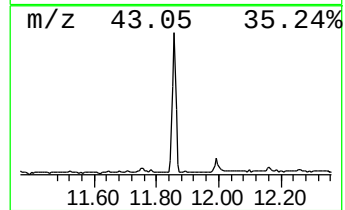
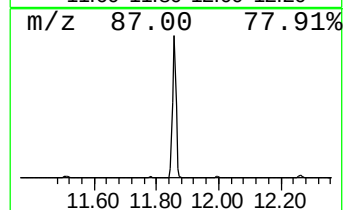
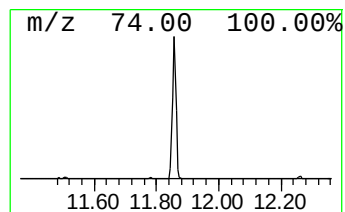
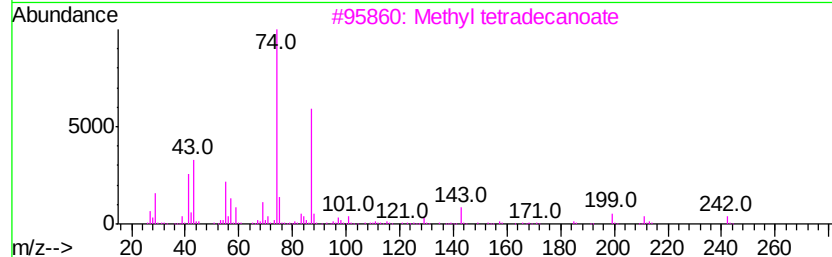
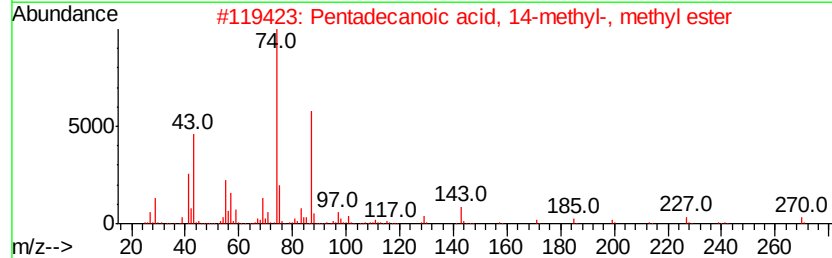
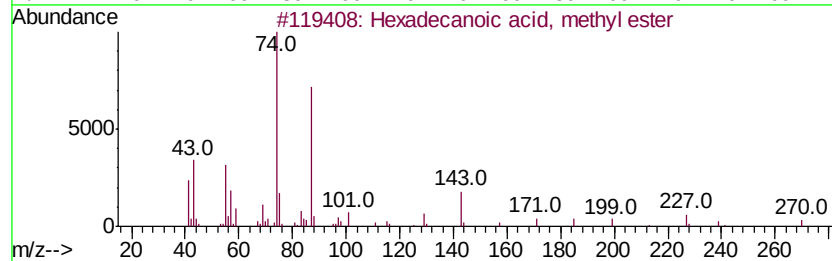
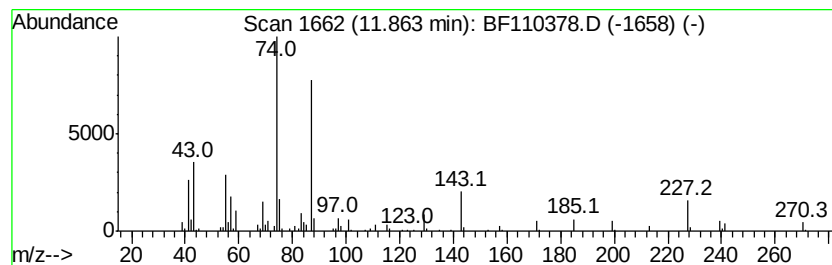
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Hexadecanoic acid, methyl e... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.		
11.86	14.94 ng	315955	Phenanthrene-d10	11.49		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	99
2		Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	93
3		Methyl tetradecanoate	242	C15H30O2	000124-10-7	87
4		Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	86
5		Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	80



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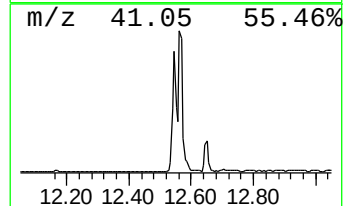
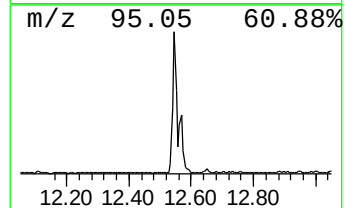
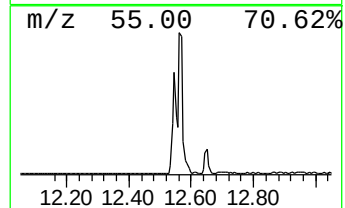
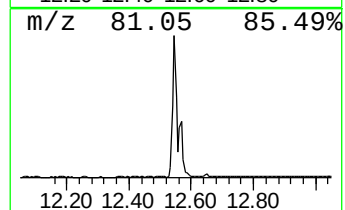
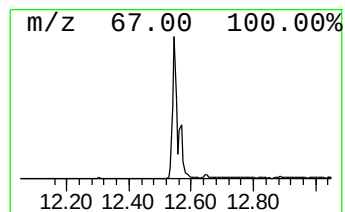
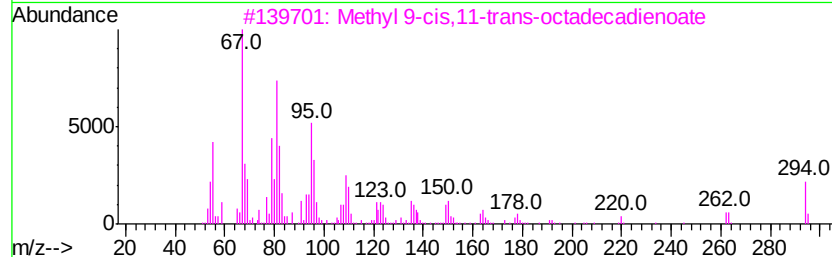
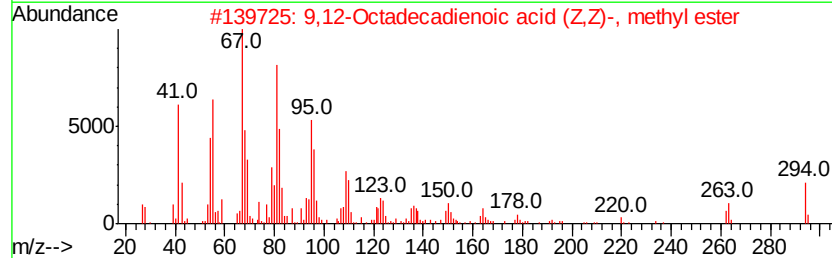
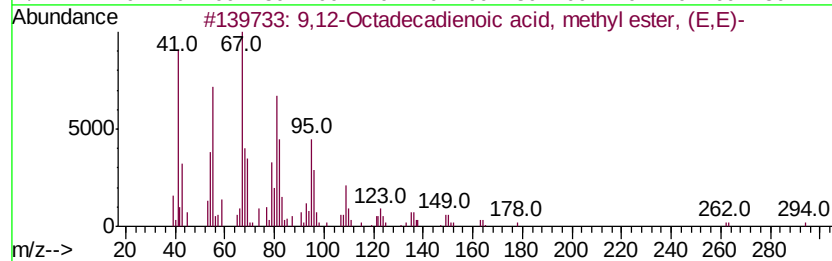
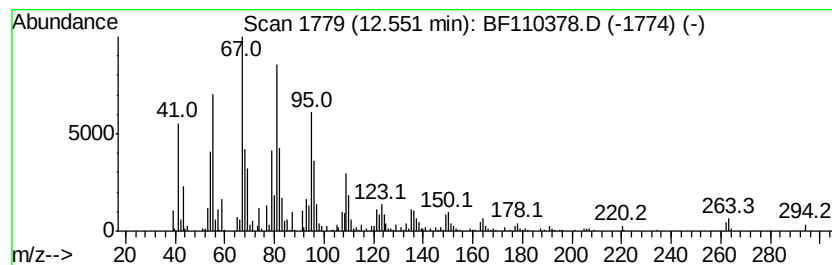
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 9,12-Octadecadienoic acid, ... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.55	44.70 ng	945247	Phenanthrene-d10	11.49

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	99
2		9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	99
3		Methyl 9-cis,11-trans-octadecadi...	294	C19H34O2	1000336-44-0	99
4		10,13-Octadecadienoic acid, meth...	294	C19H34O2	056554-62-2	99
5		9,12-Octadecadienoic acid, methy...	294	C19H34O2	002462-85-3	99



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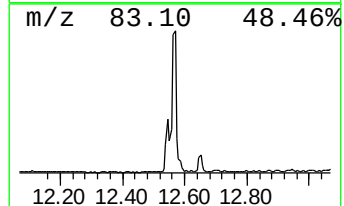
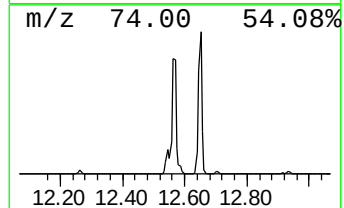
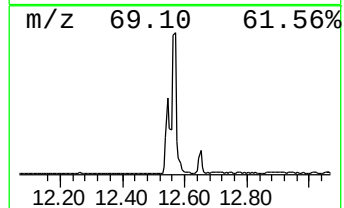
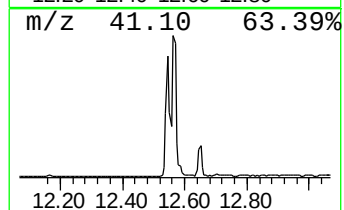
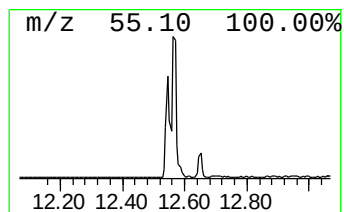
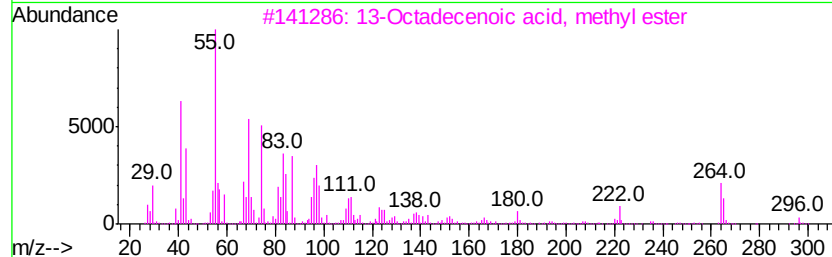
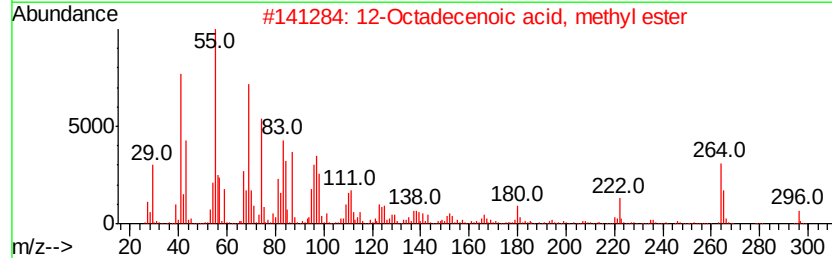
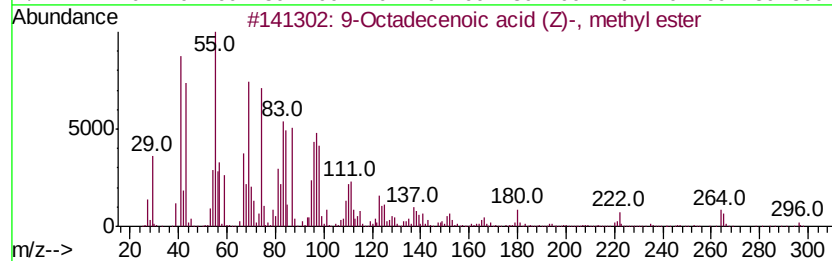
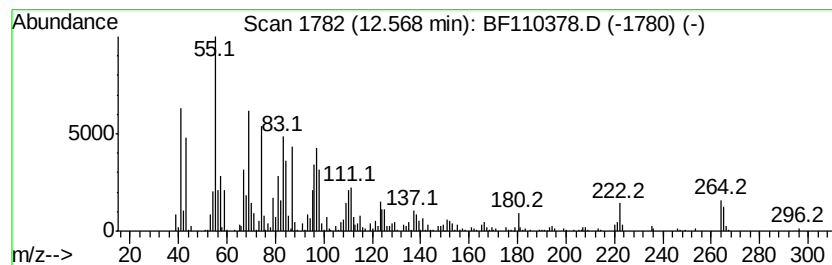
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 9-Octadecenoic acid (Z)-, m... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.57	47.18 ng	997583	Phenanthrene-d10	11.49	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	99
2	12-Octadecenoic acid, methyl ester	296	C19H36O2	056554-46-2	99
3	13-Octadecenoic acid, methyl ester	296	C19H36O2	056554-47-3	99
4	14-Octadecenoic acid, methyl ester	296	C19H36O2	056554-48-4	99
5	16-Octadecenoic acid, methyl ester	296	C19H36O2	056554-49-5	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110118\
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Operator : JU/SJ
Sample : J5192-05 2X
Misc :
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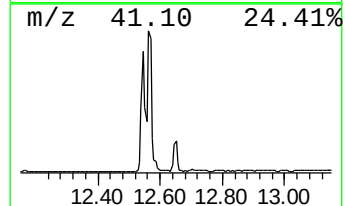
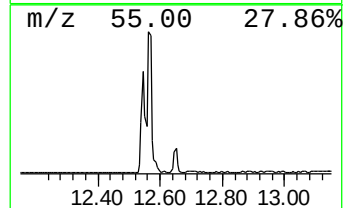
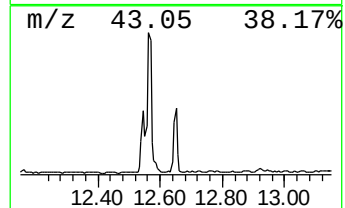
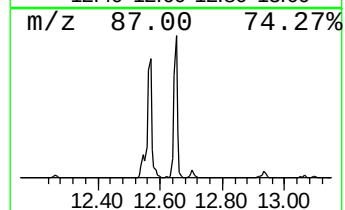
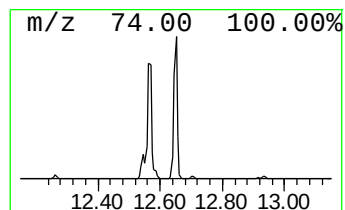
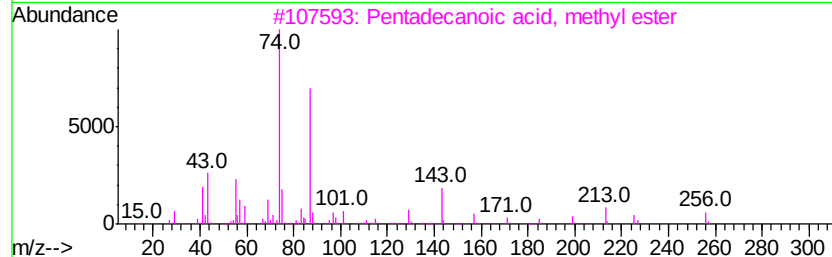
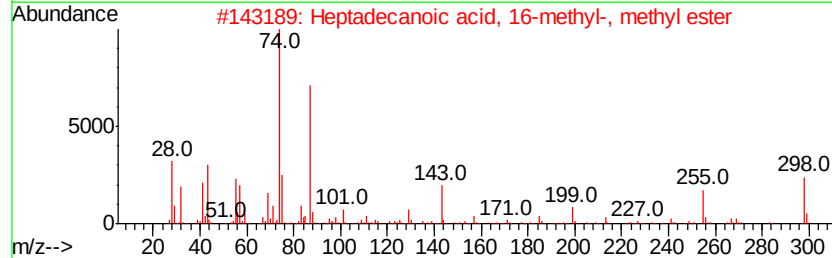
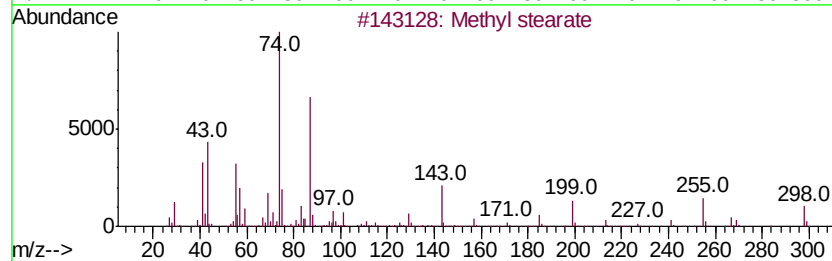
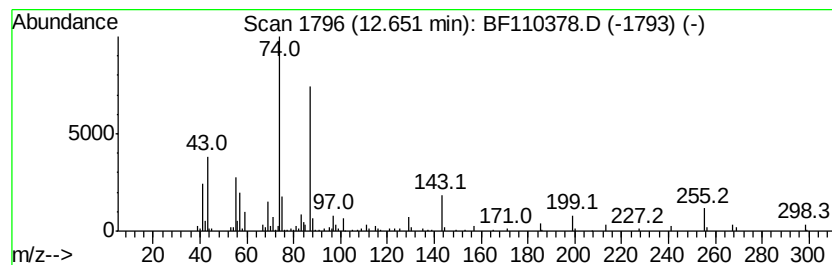
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Methyl stearate Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.65	11.08 ng	234312	Phenanthrene-d10	11.49
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Methyl stearate	298 C19H38O2	000112-61-8 98
2		Heptadecanoic acid, 16-methyl-, ...	298 C19H38O2	005129-61-3 96
3		Pentadecanoic acid, methyl ester	256 C16H32O2	007132-64-1 90
4		Decanoic acid, methyl ester	186 C11H22O2	000110-42-9 87
5		Hexadecanoic acid, methyl ester	270 C17H34O2	000112-39-0 87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110118\
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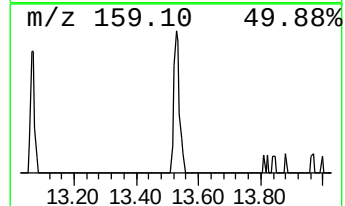
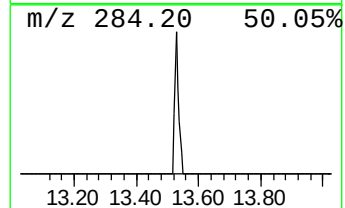
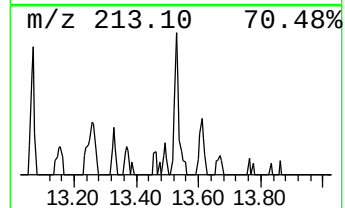
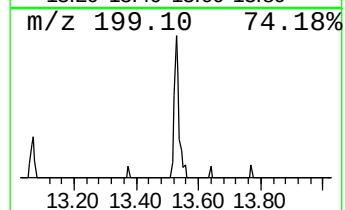
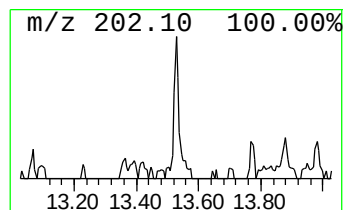
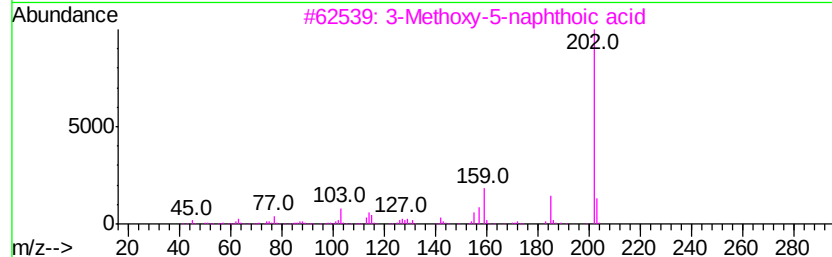
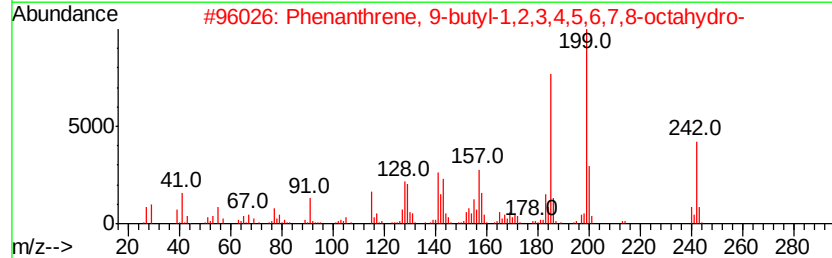
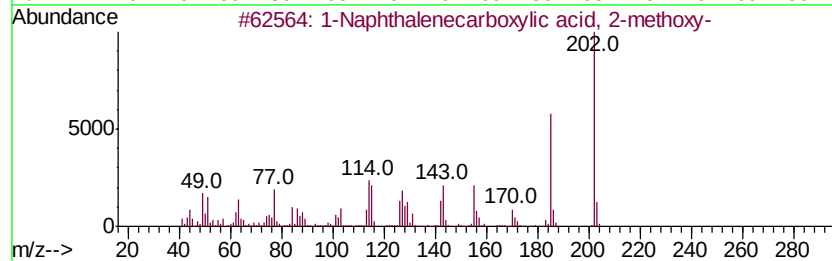
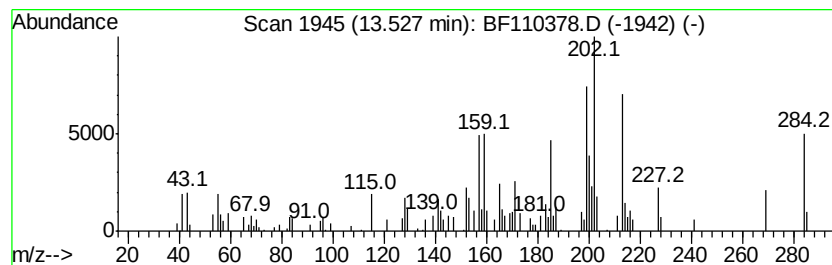
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown13.53 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.53	2.38 ng	78507	Chrysene-d12	14.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Naphthalenecarboxylic acid, 2-...	202	C12H10O3	000947-62-6	25
2		Phenanthrene, 9-butyl-1,2,3,4,5,...	242	C18H26	016703-29-0	20
3		3-Methoxy-5-naphthoic acid	202	C12H10O3	007498-58-0	20
4		9-Chloromethyl-S-octahydrophenan...	234	C15H19Cl	1000255-17-2	18
5		n-Heptanoic acid, methyl(tetramet...	228	C12H24O2Si	1000217-03-6	18



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110118\
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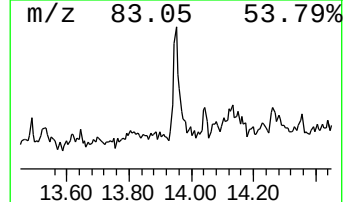
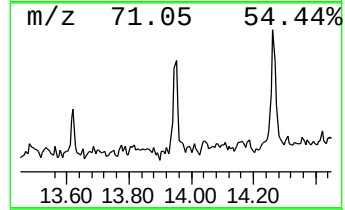
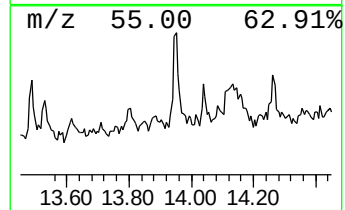
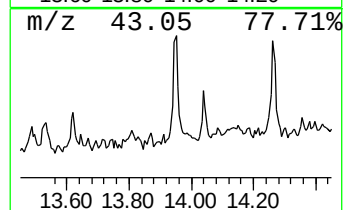
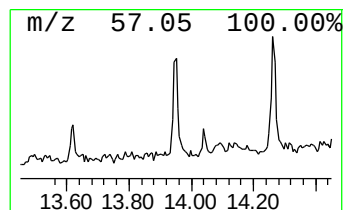
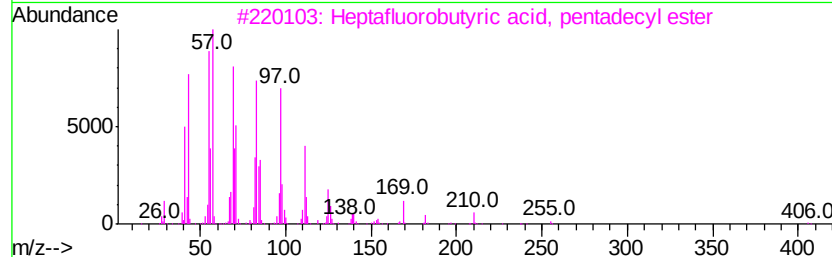
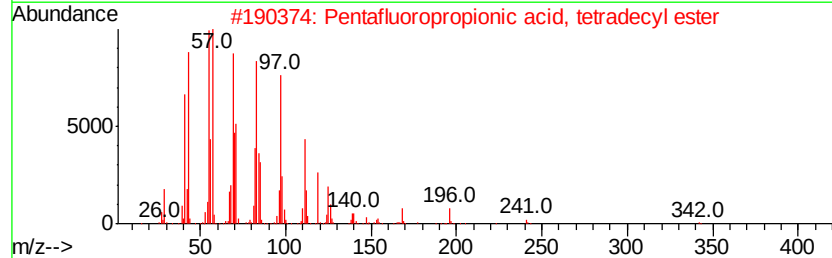
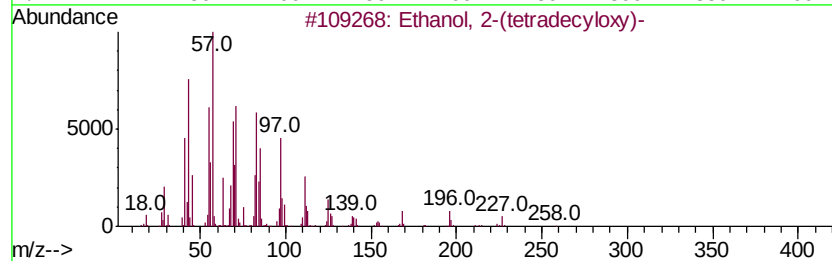
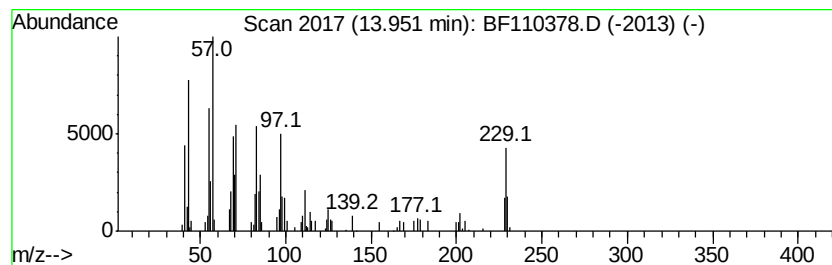
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Ethanol, 2-(tetradecyloxy)- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.95	2.07 ng	68211	Chrysene-d12	14.13
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Ethanol, 2-(tetradecyloxy)-	258 C16H34O2	002136-70-1 53
2		Pentafluoropropionic acid, tetra...	360 C17H29F5O2	006222-06-6 53
3		Heptafluorobutyric acid, pentade...	424 C19H31F7O2	959261-23-5 53
4		Heptafluorobutyric acid, n-tetra...	410 C18H29F7O2	007365-36-8 53
5		Pentafluoropropionic acid, penta...	374 C18H31F5O2	959092-08-1 53



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110118\
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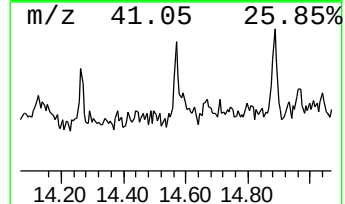
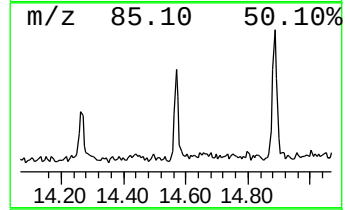
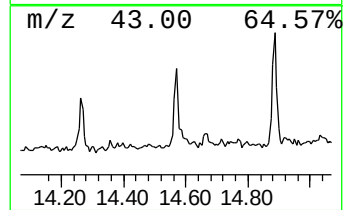
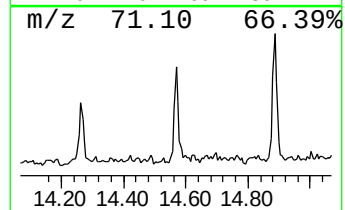
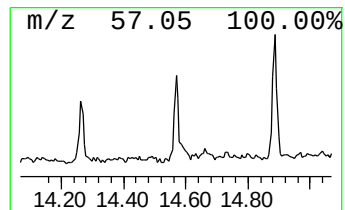
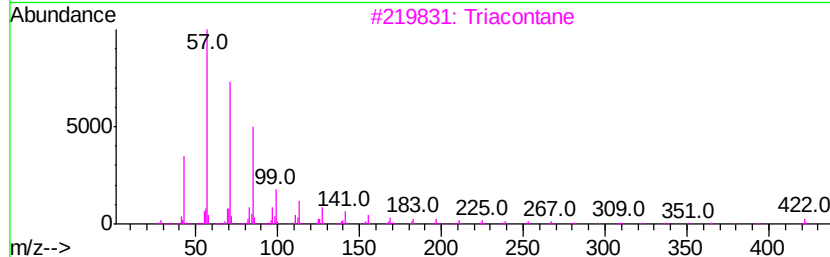
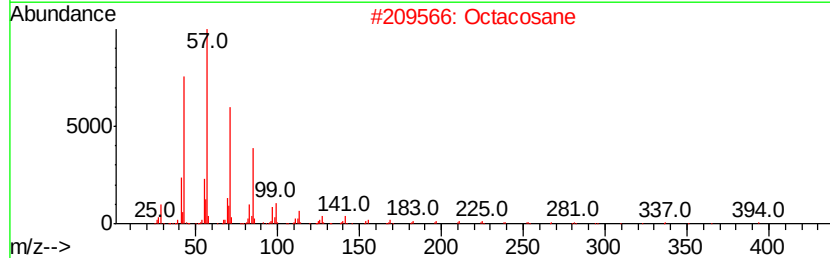
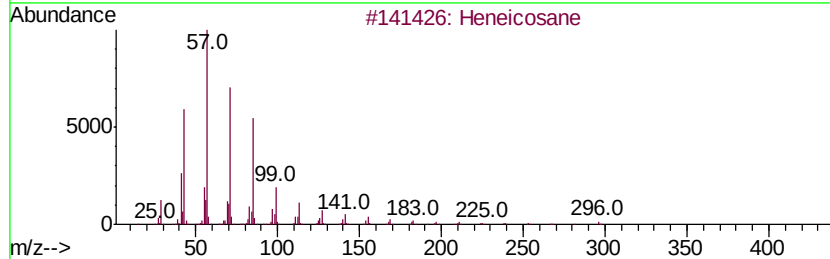
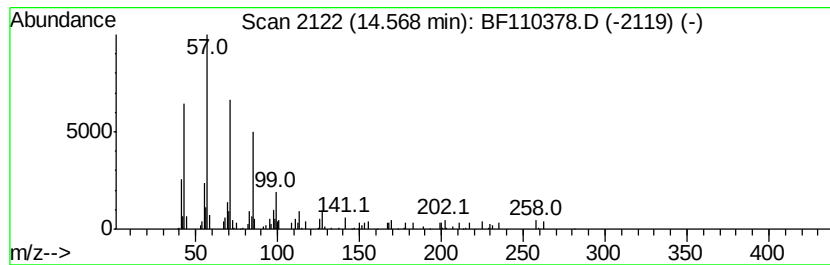
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Heneicosane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.57	2.07 ng	68240	Chrysene-d12	14.13
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Heneicosane	296 C21H44	000629-94-7	90
2	Octacosane	394 C28H58	000630-02-4	87
3	Triacontane	422 C30H62	000638-68-6	87
4	Hexacosane	366 C26H54	000630-01-3	87
5	Tetracosane	338 C24H50	000646-31-1	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110118\
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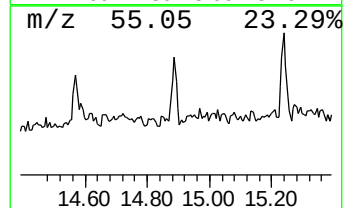
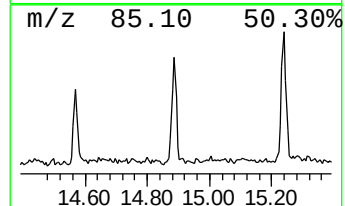
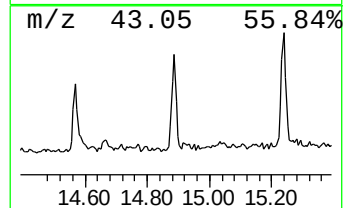
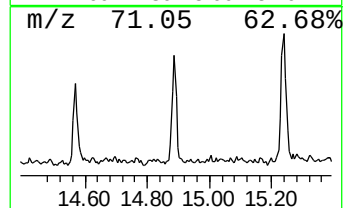
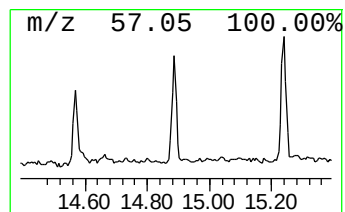
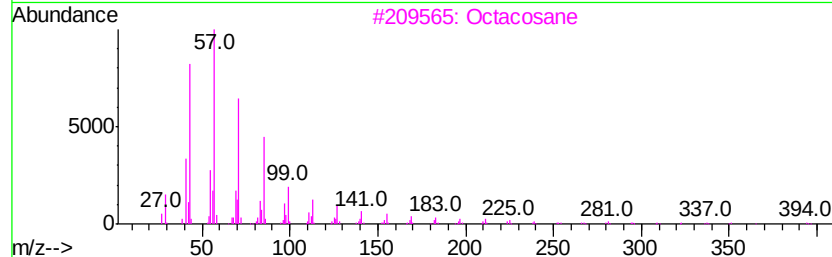
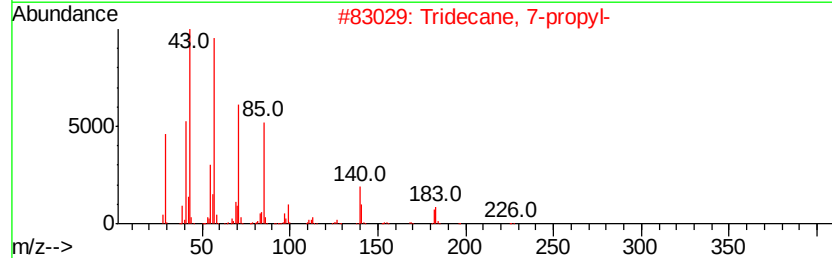
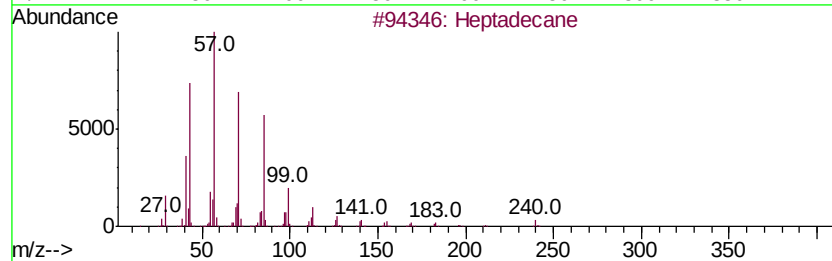
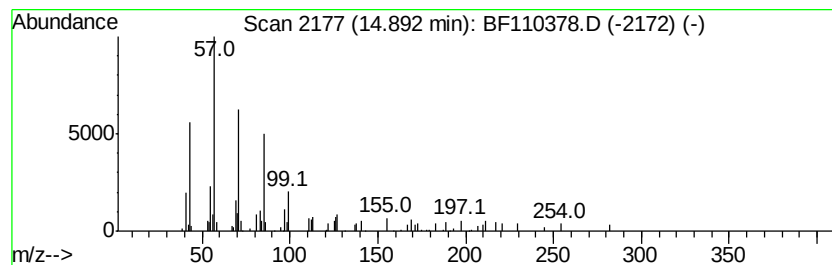
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Heptadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.89	2.31 ng	75980	Chrysene-d12	14.13
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Heptadecane	240 C17H36	000629-78-7	96
2	Tridecane, 7-propyl-	226 C16H34	055045-09-5	87
3	Octacosane	394 C28H58	000630-02-4	83
4	Tridecane, 1-iodo-	310 C13H27I	035599-77-0	83
5	Hexacosane	366 C26H54	000630-01-3	83



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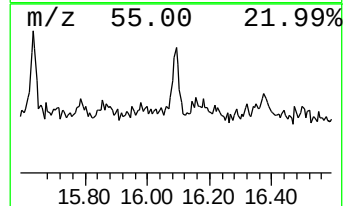
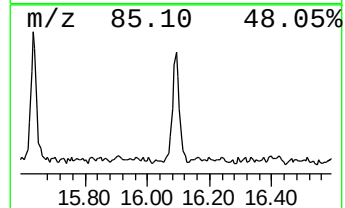
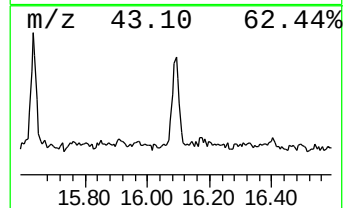
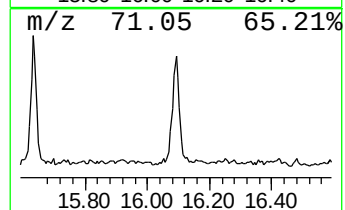
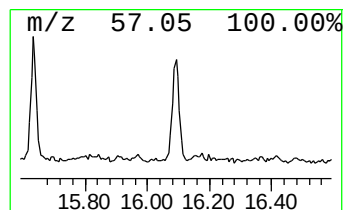
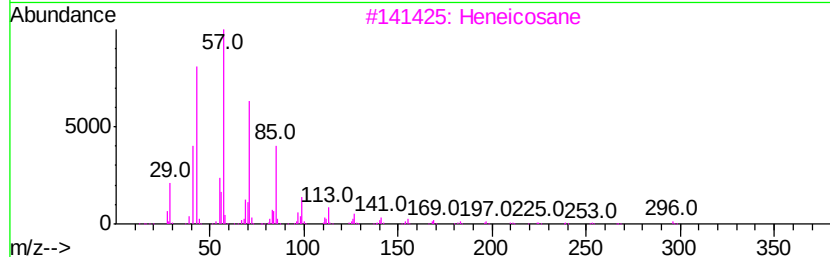
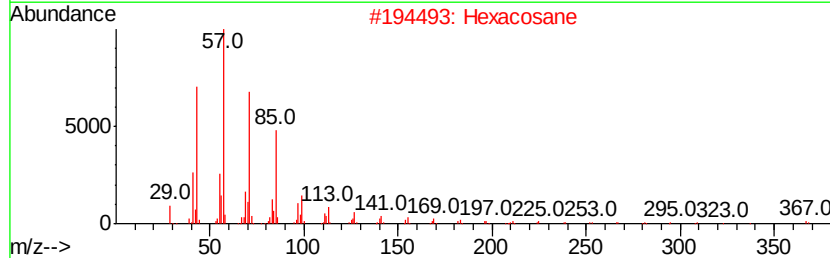
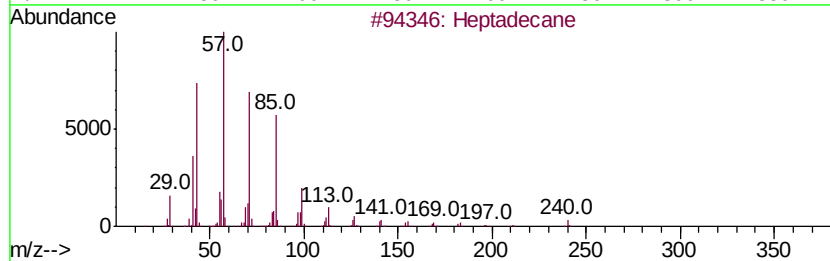
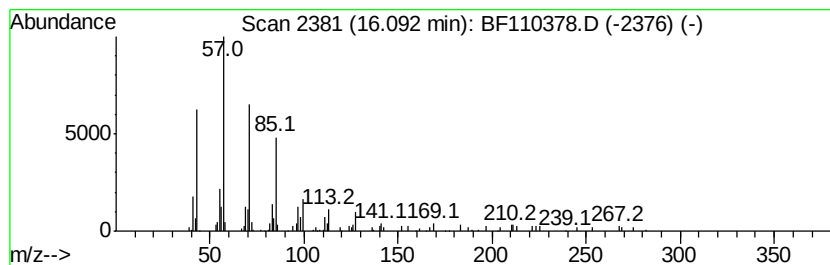
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Pentacosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.09	3.90 ng	118439	Perylene-d12	15.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane	240	C17H36	000629-78-7	91
2		Hexacosane	366	C26H54	000630-01-3	91
3		Heneicosane	296	C21H44	000629-94-7	91
4		Pentacosane	352	C25H52	000629-99-2	91
5		Hentriacontane	437	C31H64	000630-04-6	91



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Misc :
ALS Vial : 19 Sample Multiplier: 1

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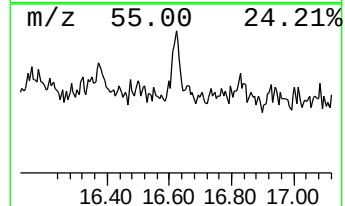
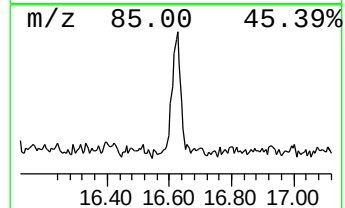
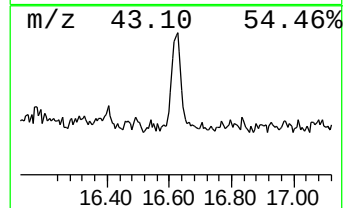
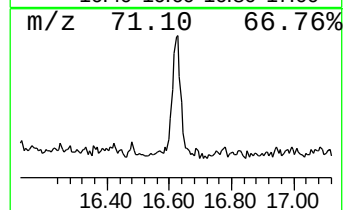
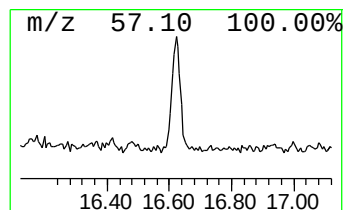
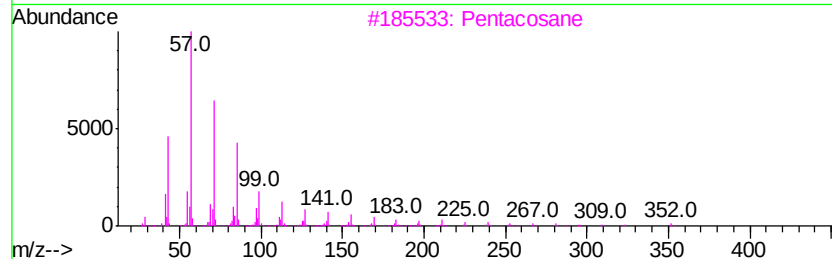
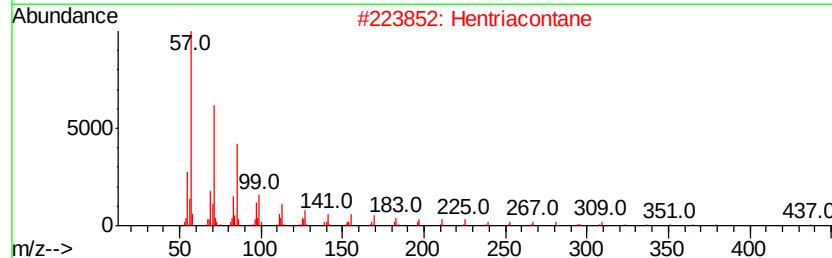
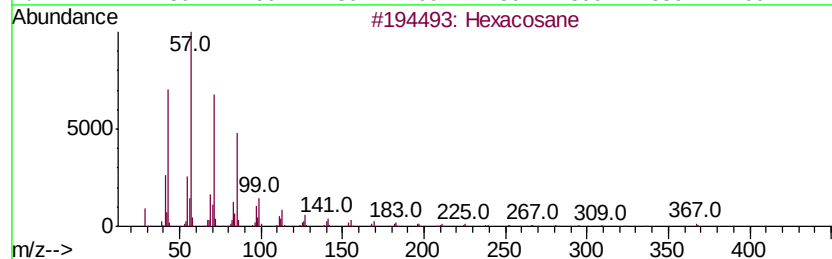
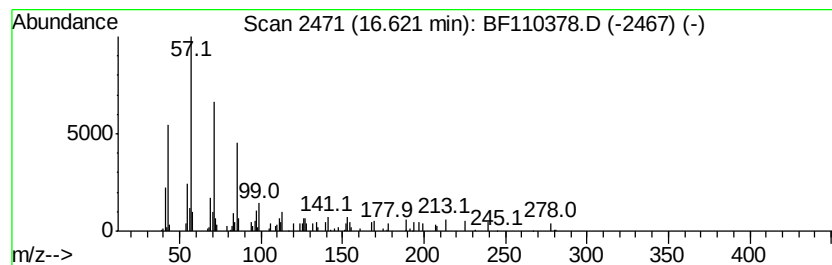
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Hexacosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.62	3.30 ng	100262	Perylene-d12	15.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	83
2		Hentriacontane	437	C31H64	000630-04-6	83
3		Pentacosane	352	C25H52	000629-99-2	83
4		Octacosane	394	C28H58	000630-02-4	80
5		Tetratetracontane	619	C44H90	007098-22-8	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110118\
Data File : BF110378.D
Acq On : 1 Nov 2018 20:07
Operator : JU/SJ
Sample : J5192-05 2X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
AU-05-103118-A

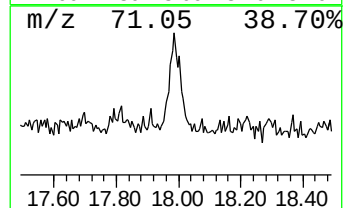
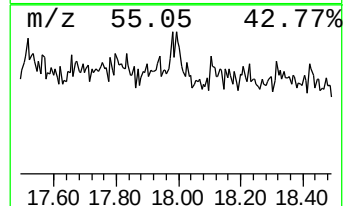
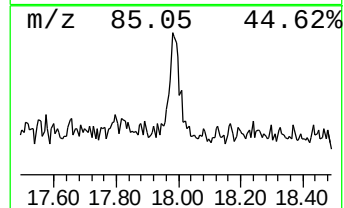
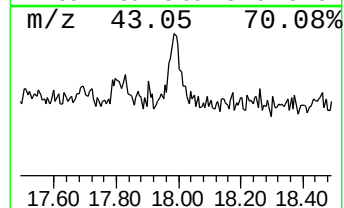
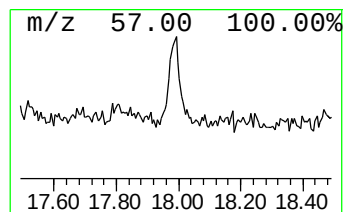
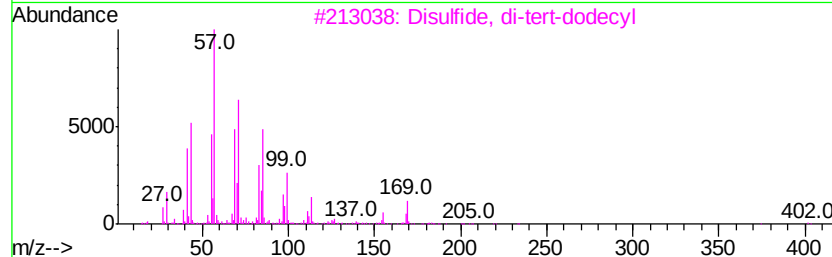
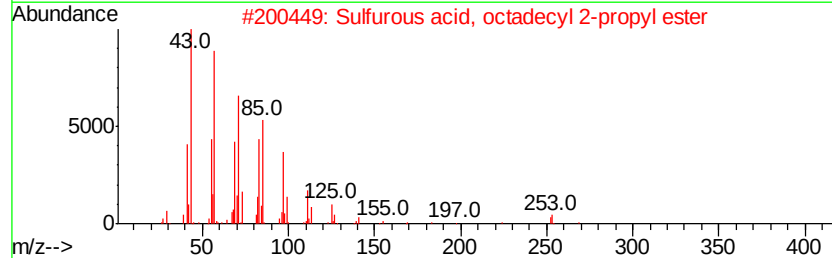
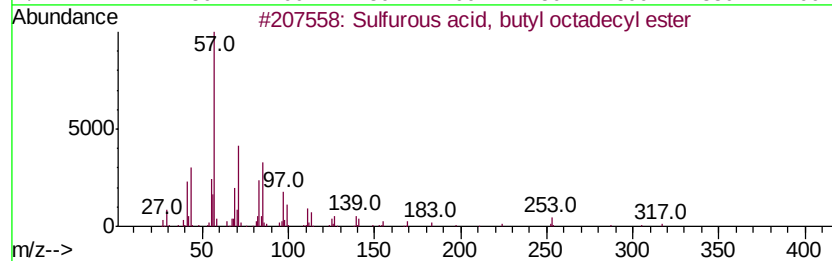
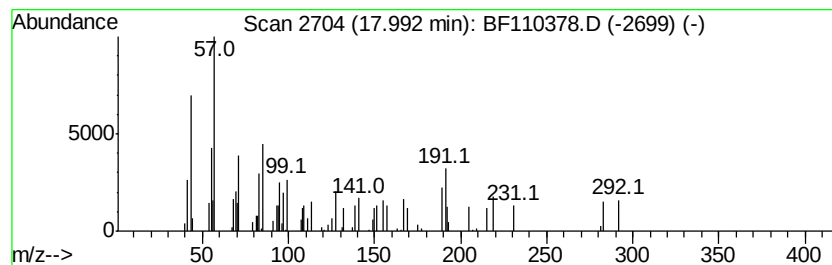
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown17.99 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.99	2.44 ng	74256	Perylene-d12	15.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Sulfurous acid, butyl octadecyl ...	390	C22H46O3S	1000309-18-5	32
2		Sulfurous acid, octadecyl 2-prop...	376	C21H44O3S	1000309-12-7	27
3		Disulfide, di-tert-dodecyl	402	C24H50S2	027458-90-8	22
4		iso-Valeraldehyde propylenealco...	144	C8H16O2	1000285-57-8	22
5		Oxalic acid, 6-ethyloct-3-yl hep...	328	C19H36O4	1000309-34-5	16



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF110118\
Data File : BF110378.D
Acq On : 1 Nov 2018 20:07
Operator : JU/SJ
Sample : J5192-05 2X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
AU-05-103118-A

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF102318.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...	2.26	12.5	ng	276575	1	6.96	443362	20.0
unknown6.72	6.72	33.9	ng	751254	1	6.96	443362	20.0
Hexadecanoic acid...	11.86	14.9	ng	315955	4	11.49	422883	20.0
9,12-Octadecadien...	12.55	44.7	ng	945247	4	11.49	422883	20.0
9-Octadecenoic ac...	12.57	47.2	ng	997583	4	11.49	422883	20.0
Methyl stearate	12.65	11.1	ng	234312	4	11.49	422883	20.0
unknown13.53	13.53	2.4	ng	78507	5	14.13	658942	20.0
Ethanol, 2-(tetra...	13.95	2.1	ng	68211	5	14.13	658942	20.0
Heneicosane	14.57	2.1	ng	68240	5	14.13	658942	20.0
Heptadecane	14.89	2.3	ng	75980	5	14.13	658942	20.0
Pentacosane	16.09	3.9	ng	118439	6	15.66	607833	20.0
Hexacosane	16.62	3.3	ng	100262	6	15.66	607833	20.0
unknown17.99	17.99	2.4	ng	74256	6	15.66	607833	20.0