

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030519\
 Data File : BF112876.D
 Acq On : 6 Mar 2019 2:28
 Operator : JU/SJ
 Sample : K1705-12
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EXT-027-SW014-022719

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-BF022719.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	5.351	550	555	576	rBV	3014145	3026716	94.14%	8.813%
2	6.375	724	729	732	rBV	3113590	3147693	97.90%	9.165%
3	6.510	747	752	755	rBV	3648246	3215098	100.00%	9.361%
4	6.739	787	791	795	rBV	1229569	979753	30.47%	2.853%
5	6.898	814	818	821	rBV	2934054	2494067	77.57%	7.262%
6	7.298	881	886	889	rBV	2044420	1911074	59.44%	5.564%
7	8.022	1005	1009	1021	rBV	1182696	1209083	37.61%	3.520%
8	8.733	1124	1130	1136	rBV	37891	40832	1.27%	0.119%
9	9.098	1188	1192	1206	rBV	3436814	3192939	99.31%	9.297%
10	9.357	1230	1236	1241	rBV3	24809	60443	1.88%	0.176%
11	9.486	1256	1258	1264	rVB	126101	129637	4.03%	0.377%
12	9.633	1279	1283	1289	rVB	64174	63054	1.96%	0.184%
13	9.769	1302	1306	1316	rBV	1253544	1191316	37.05%	3.469%
14	9.980	1335	1342	1345	rBV5	25566	37876	1.18%	0.110%
15	10.557	1435	1440	1455	rBV	1930757	2024408	62.97%	5.894%
16	11.251	1554	1558	1560	rBV	1158525	1087268	33.82%	3.166%
17	11.269	1560	1561	1566	rVV	243609	202340	6.29%	0.589%
18	11.321	1568	1570	1575	rVB	51642	48112	1.50%	0.140%
19	11.710	1632	1636	1639	rBV4	31524	45572	1.42%	0.133%
20	11.751	1640	1643	1645	rBV	55260	51416	1.60%	0.150%
21	11.786	1645	1649	1653	rBV2	504490	518671	16.13%	1.510%
22	11.857	1658	1661	1664	rVV	92114	126629	3.94%	0.369%
23	11.880	1664	1665	1670	rVB	57933	54563	1.70%	0.159%
24	12.068	1695	1697	1703	rVB4	37588	47561	1.48%	0.138%
25	12.298	1732	1736	1739	rBV	71896	79166	2.46%	0.231%
26	12.380	1748	1750	1755	rVB4	34377	41113	1.28%	0.120%
27	12.457	1759	1763	1767	rBV	295345	289637	9.01%	0.843%
28	12.504	1767	1771	1776	rVB8	35034	71116	2.21%	0.207%
29	12.568	1779	1782	1788	rVV2	160692	170937	5.32%	0.498%
30	12.680	1796	1801	1804	rBV	316193	344795	10.72%	1.004%
31	12.827	1823	1826	1830	rBV	2855269	2804724	87.24%	8.167%
32	12.904	1834	1839	1842	rBV3	54578	78433	2.44%	0.228%
33	13.010	1855	1857	1861	rVB	65056	60922	1.89%	0.177%
34	13.051	1861	1864	1866	rBV4	31448	41600	1.29%	0.121%

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 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 >
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35	13.074	1866	1868	1871	rVB2	63937	48590	1.51%	0.141%
36	13.115	1871	1875	1877	rBV2	67484	72574	2.26%	0.211%
37	13.204	1888	1890	1893	rBV	39903	37980	1.18%	0.111%
38	13.239	1893	1896	1899	rVB	46725	45544	1.42%	0.133%
39	13.315	1906	1909	1914	rVB5	76400	88687	2.76%	0.258%
40	13.415	1923	1926	1929	rBV	55500	49763	1.55%	0.145%
41	13.515	1940	1943	1948	rVB	49200	69923	2.17%	0.204%
42	13.621	1958	1961	1965	rBV2	39064	59761	1.86%	0.174%
43	13.733	1976	1980	1988	rBV3	324669	452831	14.08%	1.319%
44	13.874	1999	2004	2007	rBV	1478534	1544565	48.04%	4.497%
45	14.057	2032	2035	2039	rVB	218257	187702	5.84%	0.547%
46	14.262	2068	2070	2073	rVB	53484	39755	1.24%	0.116%
47	14.362	2084	2087	2093	rBV	427890	503887	15.67%	1.467%
48	14.657	2134	2137	2143	rVB	322055	320504	9.97%	0.933%
49	14.886	2173	2176	2183	rVB	95220	181211	5.64%	0.528%
50	14.974	2188	2191	2194	rBV	311896	376103	11.70%	1.095%
51	15.227	2231	2234	2238	rVB	67261	75790	2.36%	0.221%
52	15.286	2240	2244	2248	rVV	678007	773323	24.05%	2.252%
53	15.333	2249	2252	2258	rVB	236573	285199	8.87%	0.830%
54	15.739	2318	2321	2326	rVB	177131	241851	7.52%	0.704%

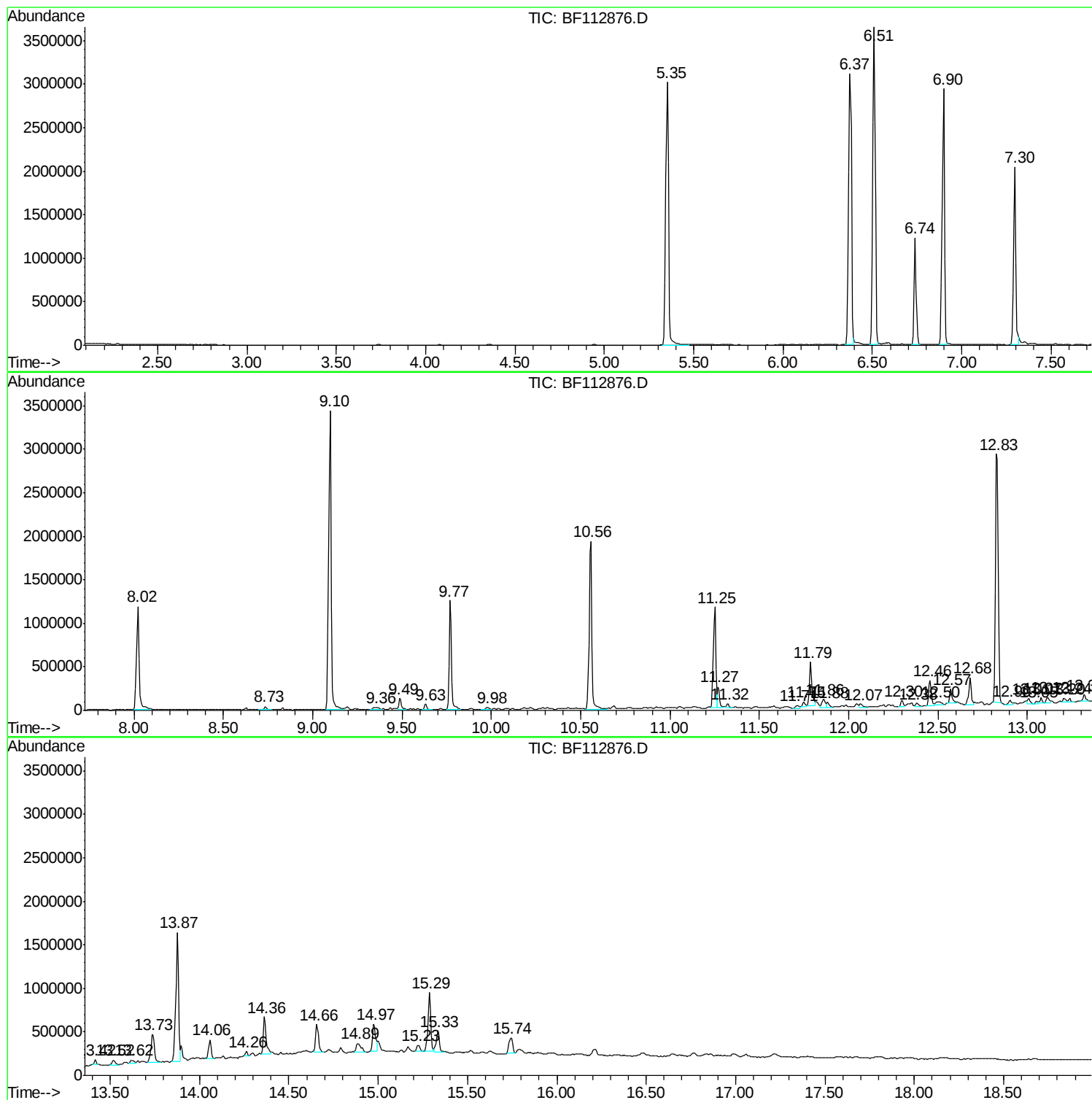
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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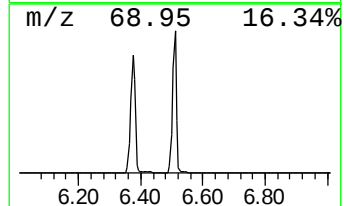
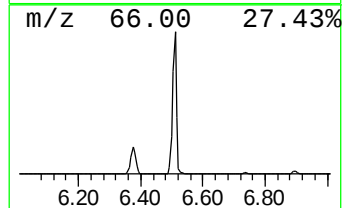
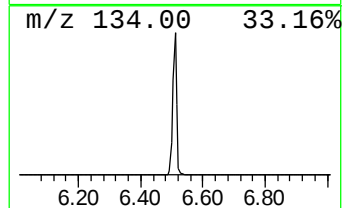
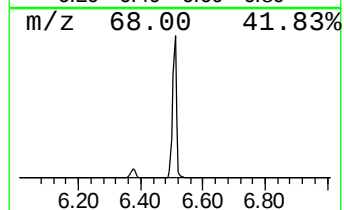
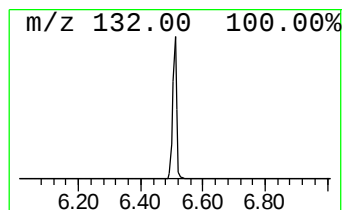
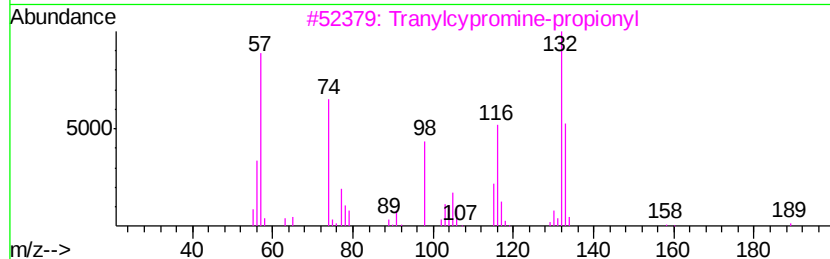
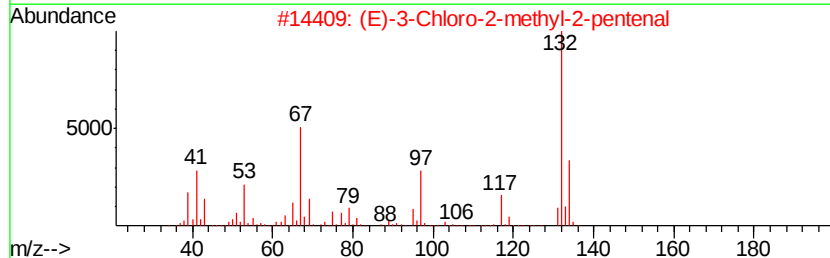
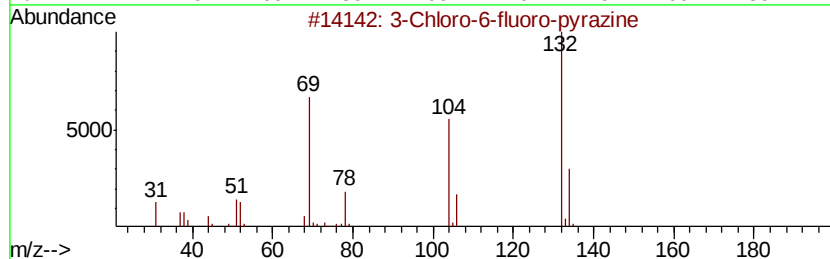
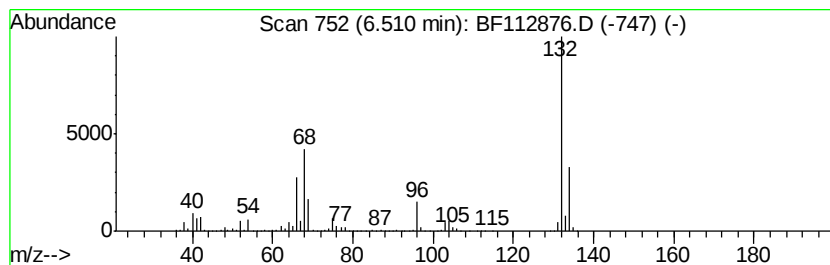
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown6.51 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.51	65.63 ng	3215100	1,4-Dichlorobenzene-d4	6.74

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		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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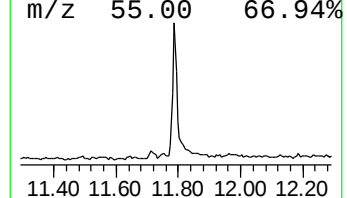
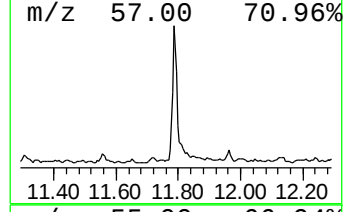
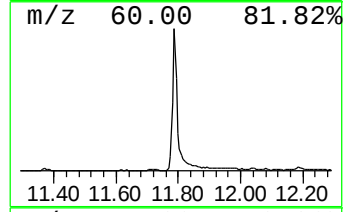
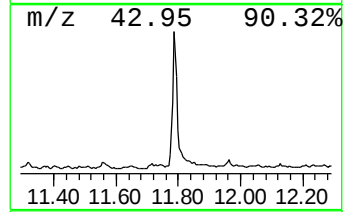
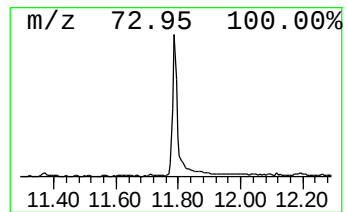
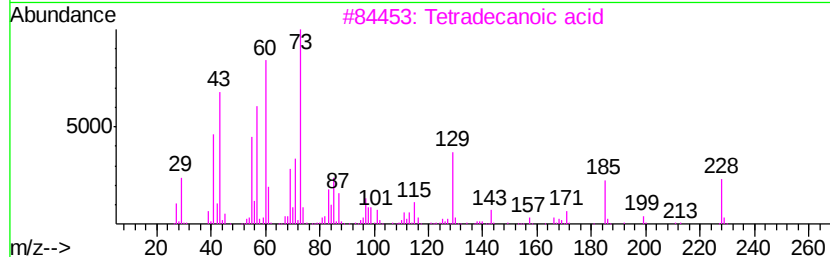
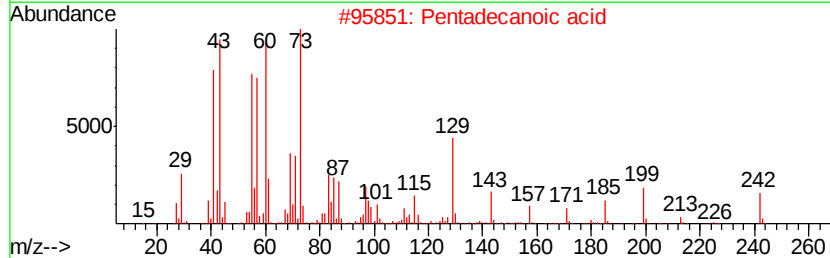
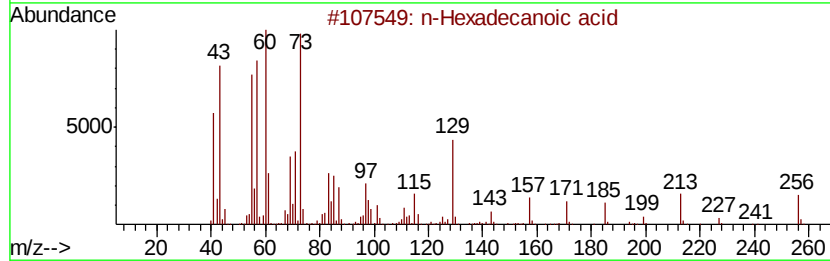
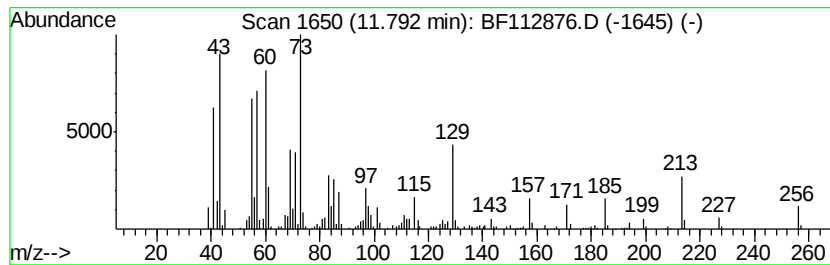
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.79	9.54 ng	518671	Phenanthrene-d10	11.25

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	93
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	90
4		Tridecanoic acid	214	C13H26O2	000638-53-9	86
5		n-Decanoic acid	172	C10H20O2	000334-48-5	76



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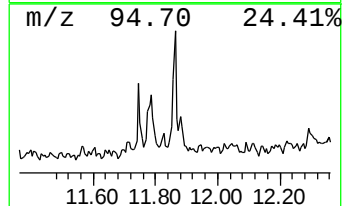
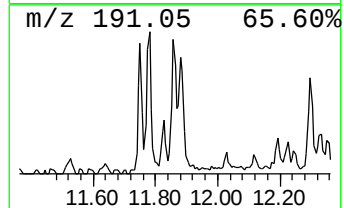
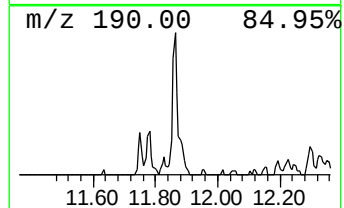
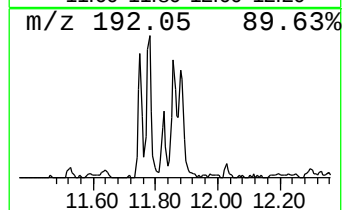
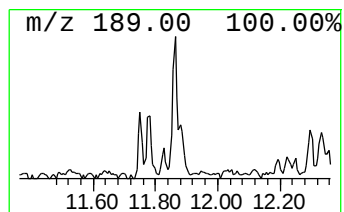
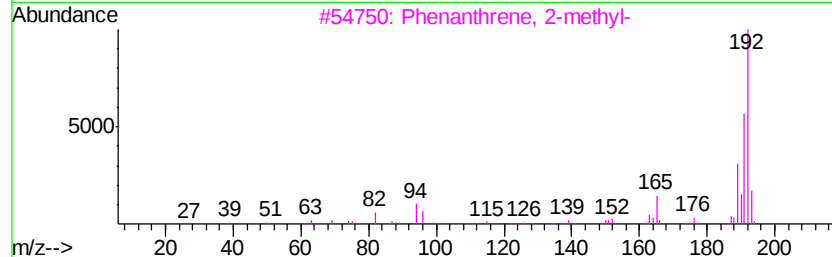
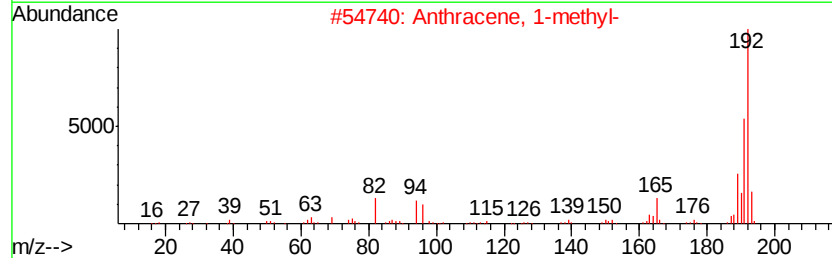
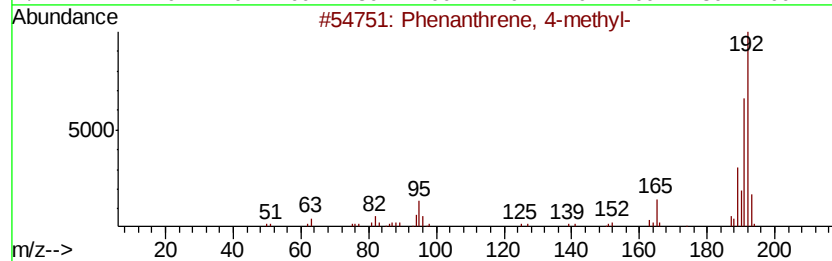
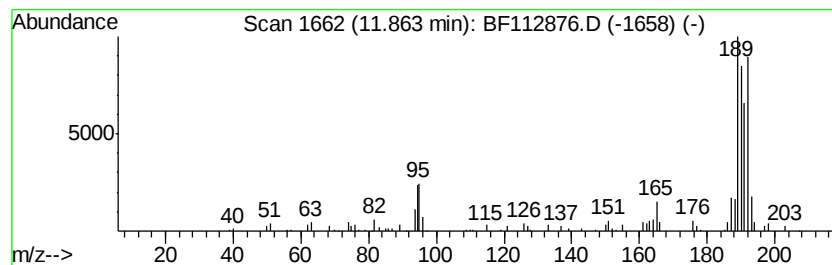
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Phenanthrene, 4-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.86	2.33 ng	126629	Phenanthrene-d10	11.25	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	92
2	Anthracene, 1-methyl-	192	C15H12	000610-48-0	90
3	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	60
4	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	60
5	Anthracene, 2-methyl-	192	C15H12	000613-12-7	60



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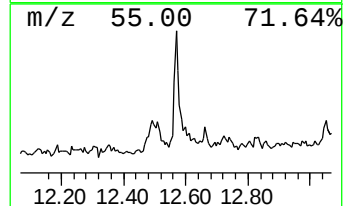
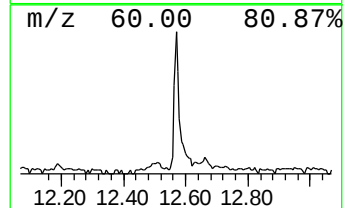
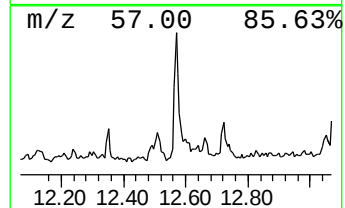
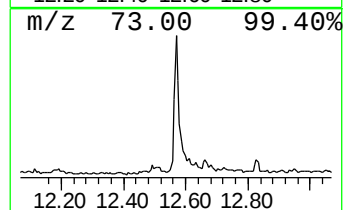
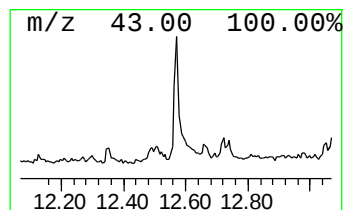
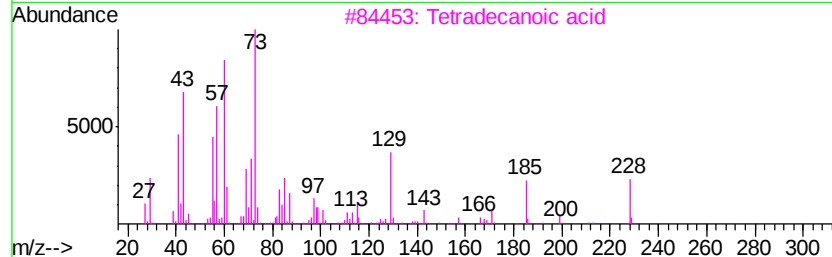
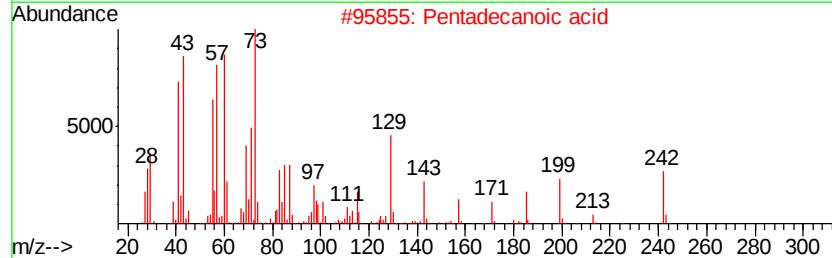
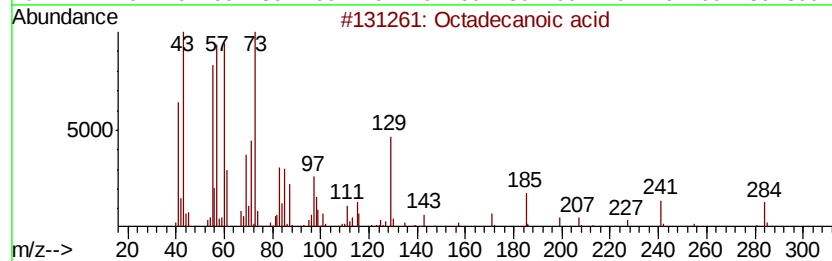
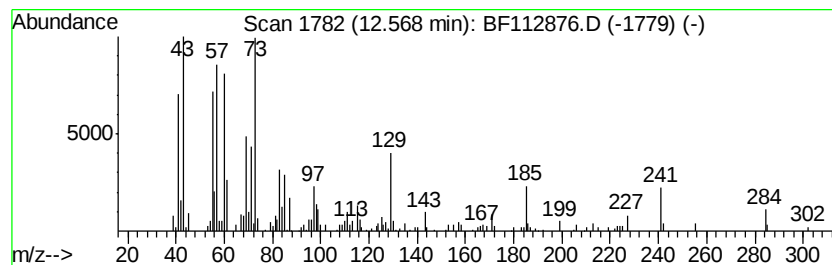
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Octadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.57	2.21 ng	170937	Chrysene-d12	13.87

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	87
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	87
4		n-Decanoic acid	172	C10H20O2	000334-48-5	64
5		Undecanoic acid	186	C11H22O2	000112-37-8	55



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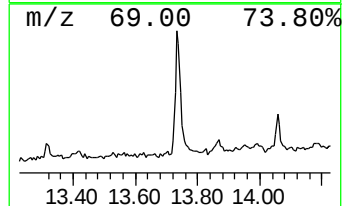
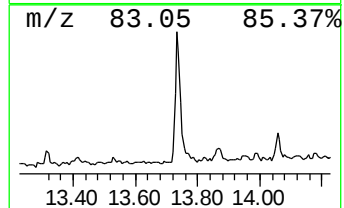
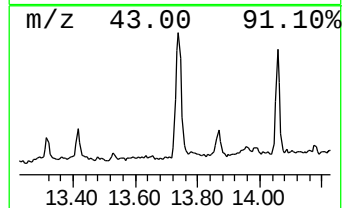
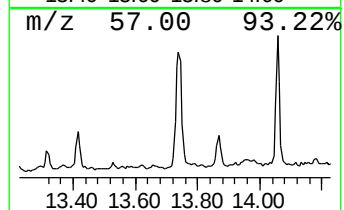
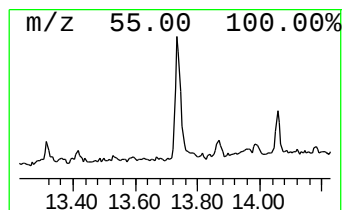
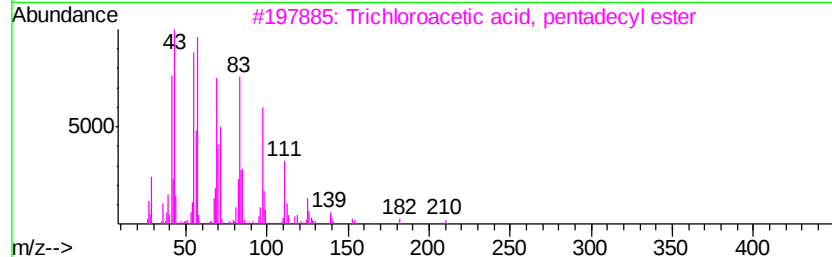
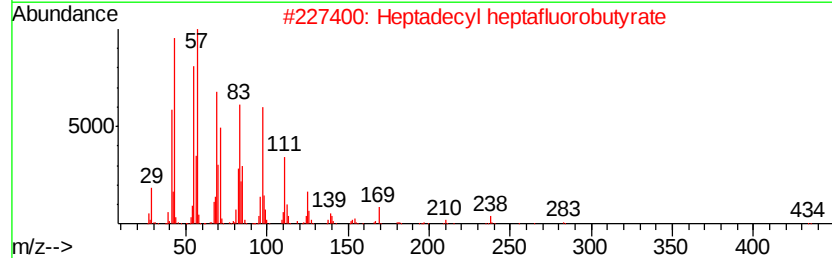
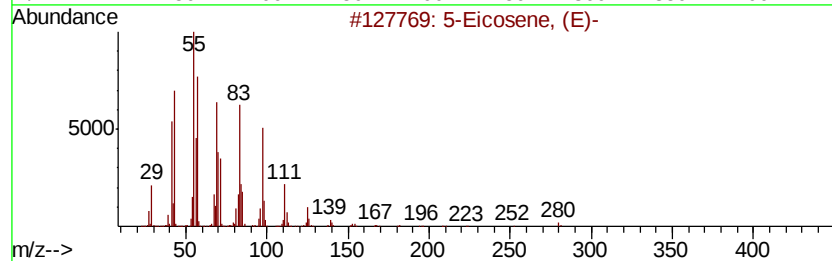
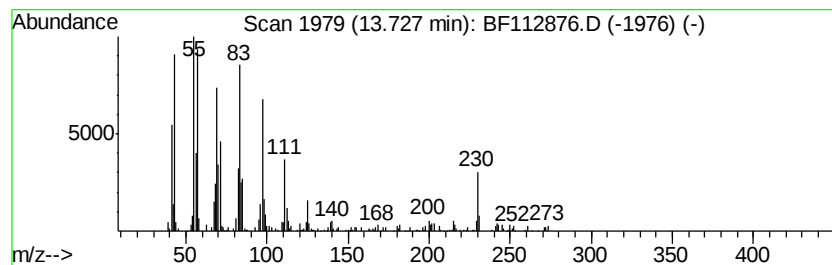
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 6 5-Eicosene, (E)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.73	5.86 ng	452831	Chrysene-d12	13.87
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	5-Eicosene, (E)-	280 C20H40	074685-30-6	97
2	Heptadecyl heptafluorobutyrate	452 C21H35F7O2	959085-66-6	94
3	Trichloroacetic acid, pentadecyl...	372 C17H31Cl3O2	074339-53-0	94
4	n-Tetracosanol-1	354 C24H50O	000506-51-4	94
5	E-7-Octadecene	252 C18H36	1000130-92-0	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030519\
Data File : BF112876.D
Acq On : 6 Mar 2019 2:28
Operator : JU/SJ
Sample : K1705-12
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
EXT-027-SW014-022719

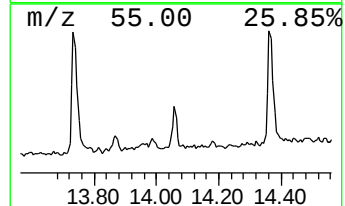
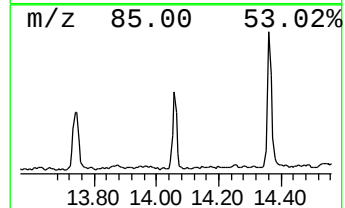
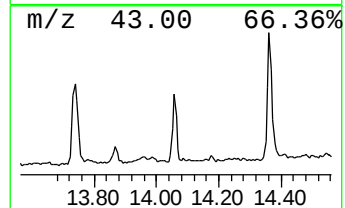
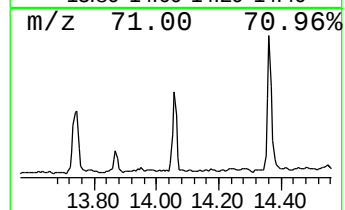
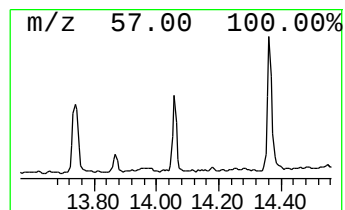
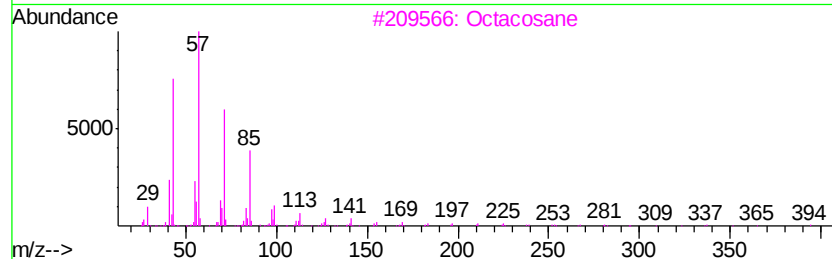
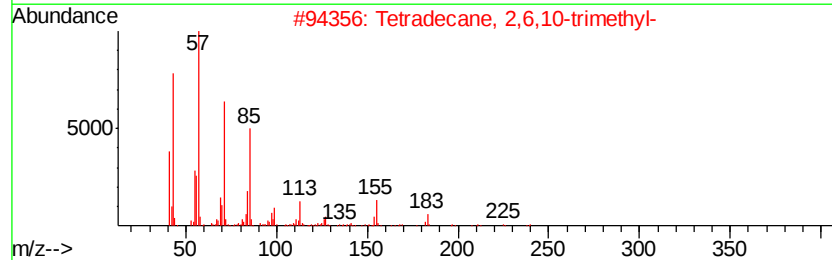
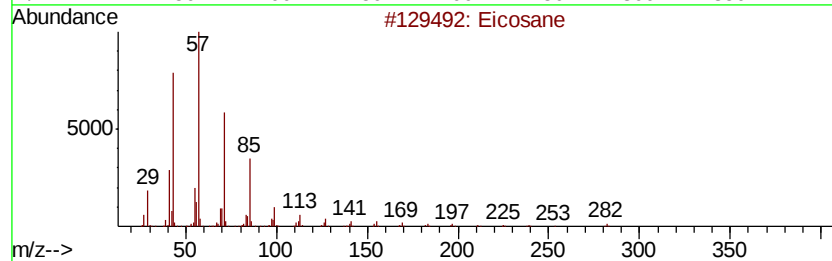
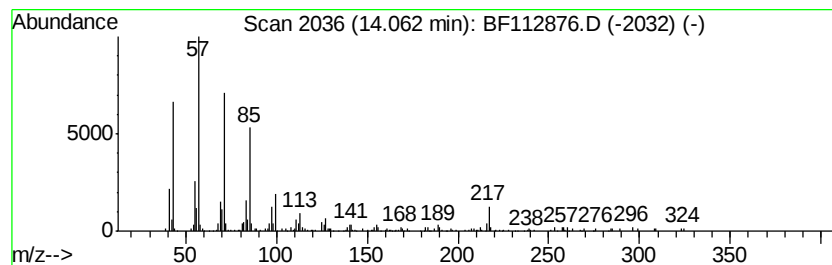
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Eicosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.06	2.43 ng	187702	Chrysene-d12	13.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	93
2		Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	91
3		Octacosane	394	C28H58	000630-02-4	90
4		Hexacosane	366	C26H54	000630-01-3	90
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030519\
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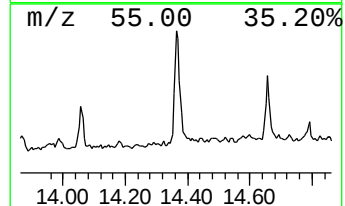
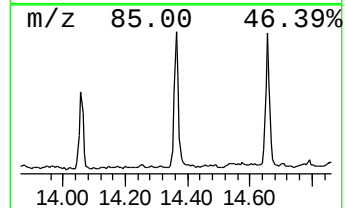
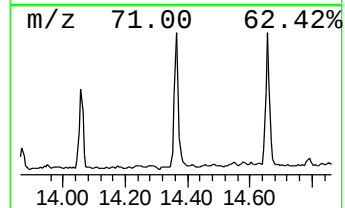
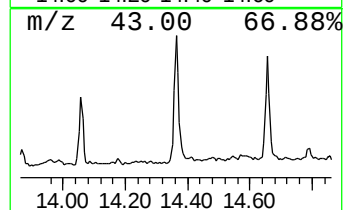
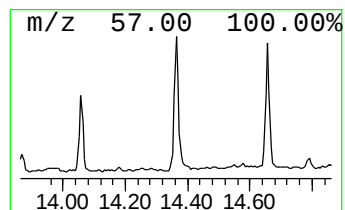
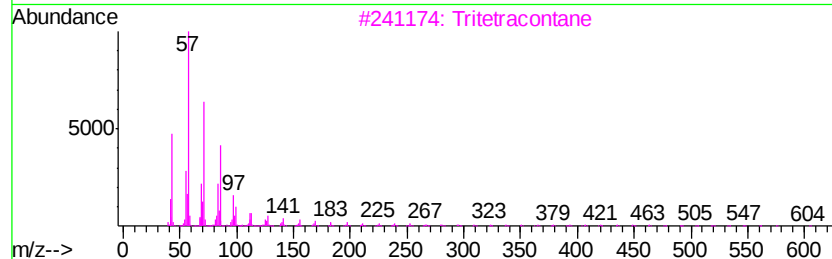
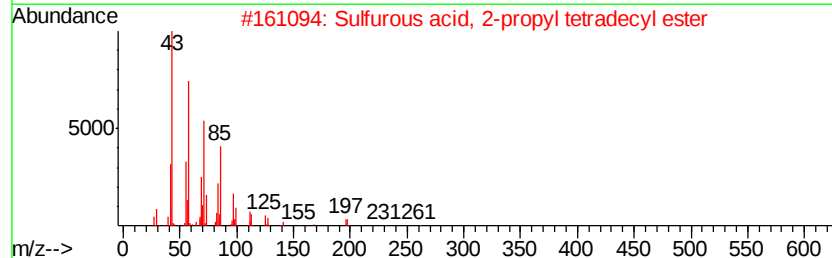
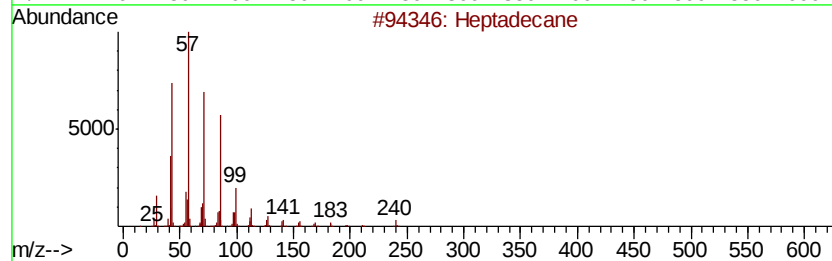
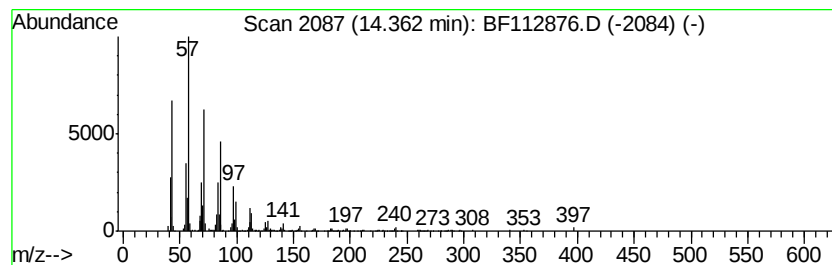
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 8 Heptadecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.36	6.52 ng	503887	Chrysene-d12	13.87	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	96
2	Sulfurous acid, 2-propyl tetradec...	320	C17H36O3S	1000309-12-5	91
3	Tritetracontane	605	C43H88	007098-21-7	87
4	Sulfurous acid, dodecyl 2-propyl...	292	C15H32O3S	1000309-12-3	87
5	Heptacosane, 1-chloro-	414	C27H55Cl	062016-79-9	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030519\
Data File : BF112876.D
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Operator : JU/SJ
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Instrument :
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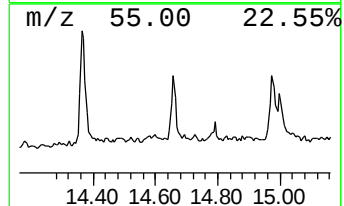
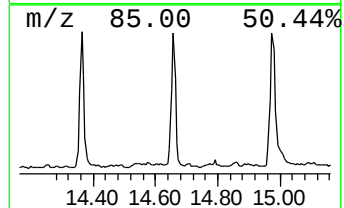
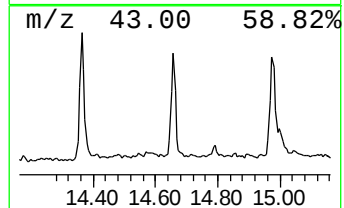
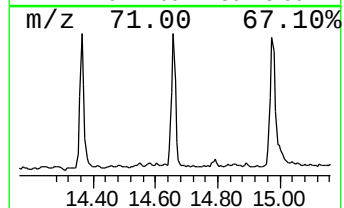
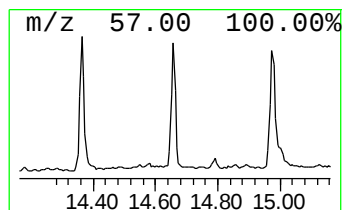
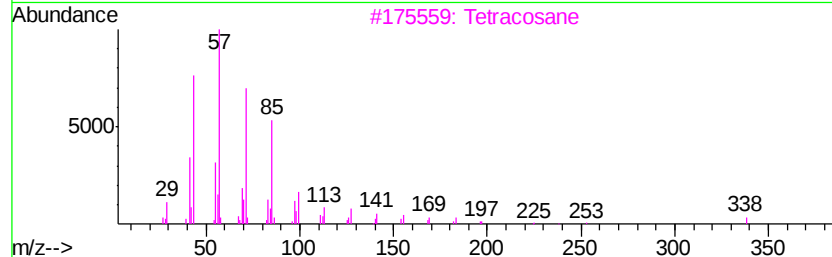
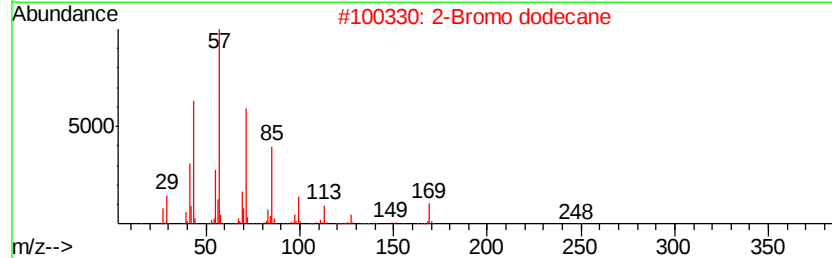
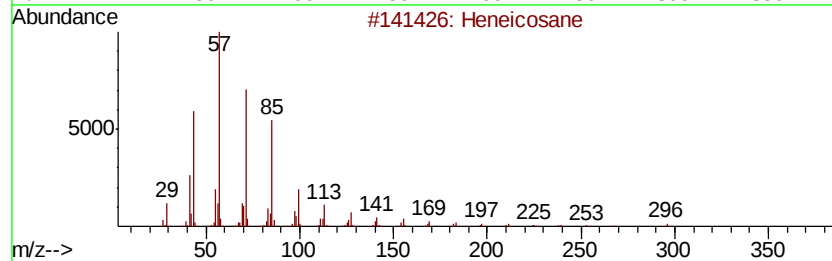
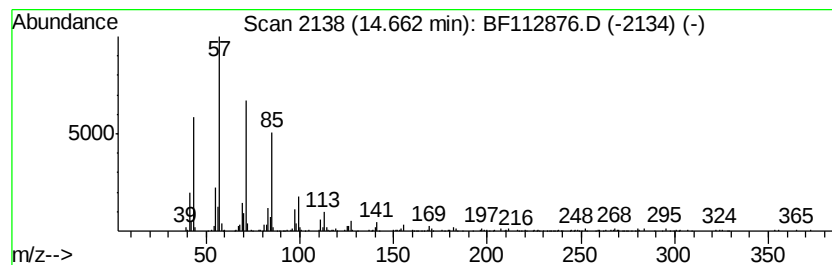
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Heneicosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.66	8.29 ng	320504	Perylene-d12	15.29

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	91
2	2-Bromo dodecane		248	C12H25Br	013187-99-0	90
3	Tetracosane		338	C24H50	000646-31-1	90
4	Eicosane		282	C20H42	000112-95-8	87
5	Heptadecane		240	C17H36	000629-78-7	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030519\
 Data File : BF112876.D
 Acq On : 6 Mar 2019 2:28
 Operator : JU/SJ
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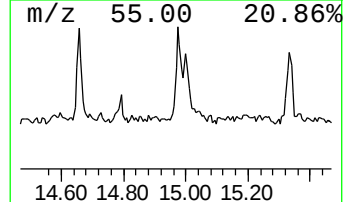
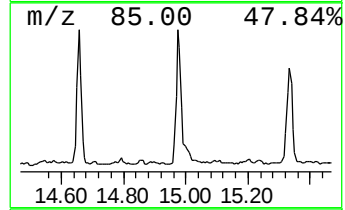
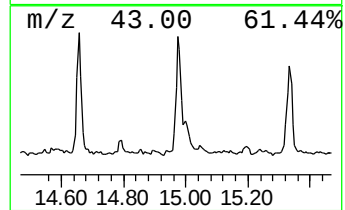
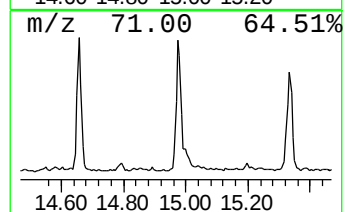
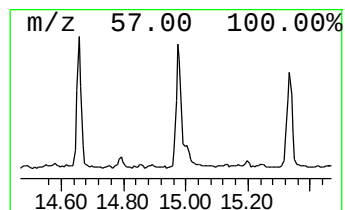
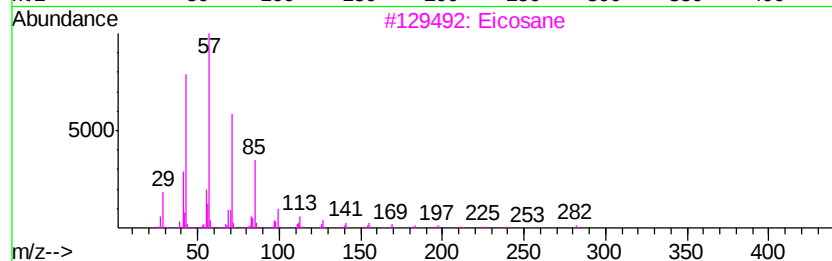
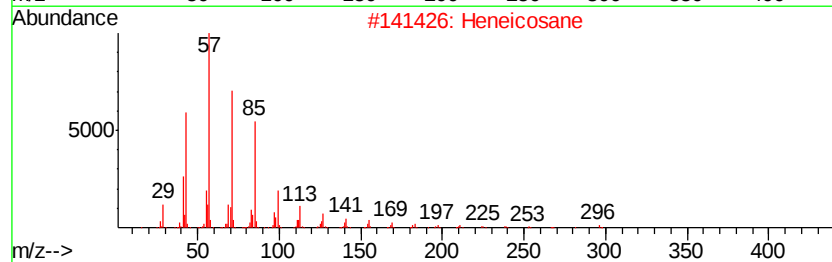
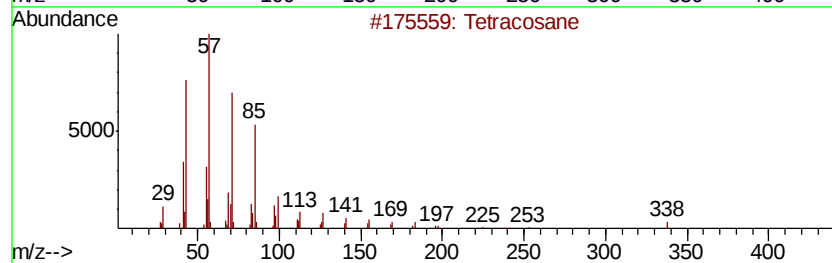
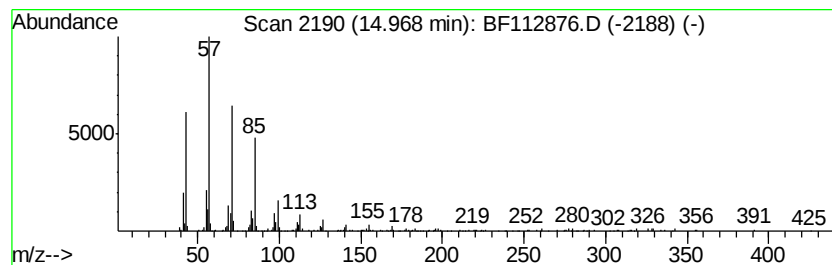
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 10 Tetracosane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.97	9.73 ng	376103	Perylene-d12	15.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane	338	C24H50	000646-31-1	91
2		Heneicosane	296	C21H44	000629-94-7	90
3		Eicosane	282	C20H42	000112-95-8	90
4		Tridecane, 6-propyl-	226	C16H34	055045-10-8	90
5		Octadecane	254	C18H38	000593-45-3	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030519\
Data File : BF112876.D
Acq On : 6 Mar 2019 2:28
Operator : JU/SJ
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ClientSampleId :
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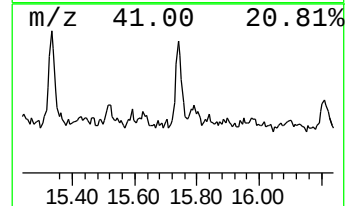
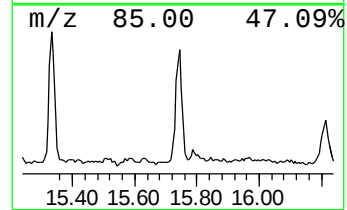
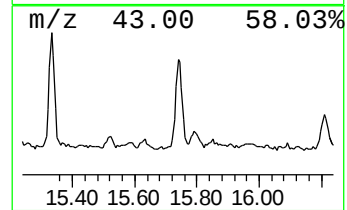
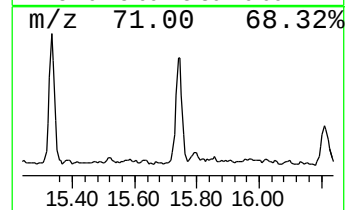
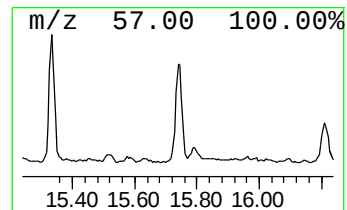
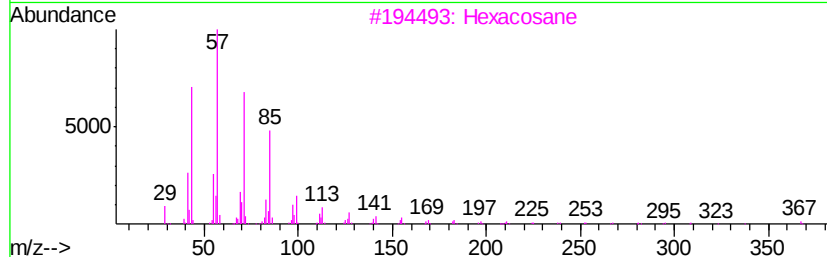
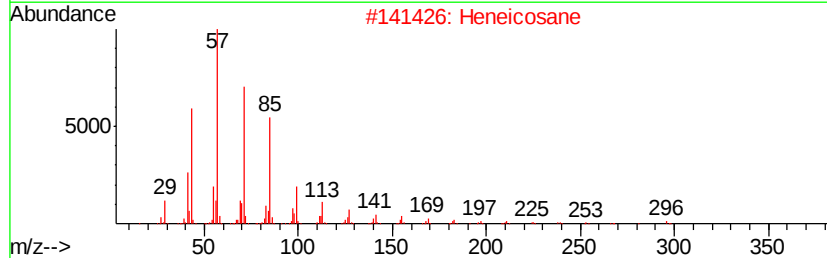
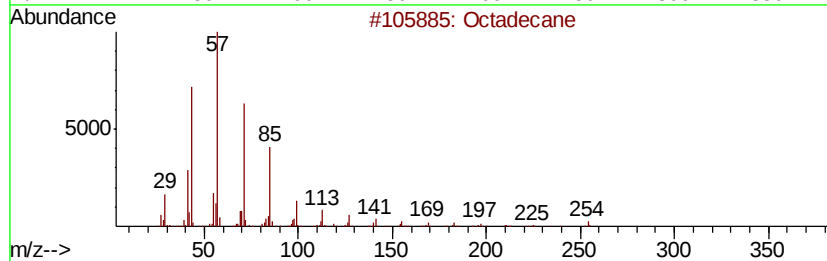
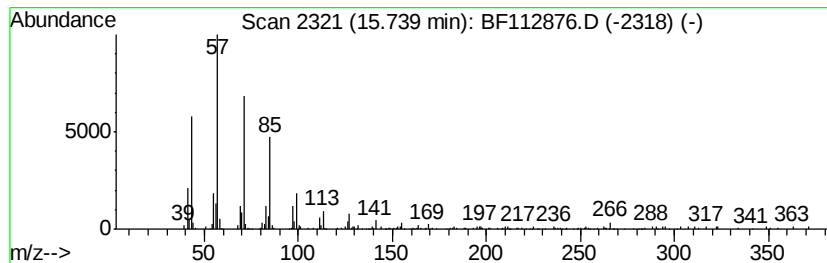
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Octadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.74	6.25 ng	241851	Perylene-d12	15.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	91
2		Heneicosane	296	C21H44	000629-94-7	90
3		Hexacosane	366	C26H54	000630-01-3	90
4		Tetracosane	338	C24H50	000646-31-1	87
5		triacontane	422	C30H62	000638-68-6	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF030519\
Data File : BF112876.D
Acq On : 6 Mar 2019 2:28
Operator : JU/SJ
Sample : K1705-12
Misc :
ALS Vial : 21 Sample Multiplier: 1

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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n-Hexadecanoic acid	11.79	9.5	ng	518671	4	11.25	1087270	20.0
Phenanthrene, 4-m...	11.86	2.3	ng	126629	4	11.25	1087270	20.0
Octadecanoic acid	12.57	2.2	ng	170937	5	13.87	1544570	20.0
5-Eicosene, (E)-	13.73	5.9	ng	452831	5	13.87	1544570	20.0
Eicosane	14.06	2.4	ng	187702	5	13.87	1544570	20.0
Heptadecane	14.36	6.5	ng	503887	5	13.87	1544570	20.0
Heneicosane	14.66	8.3	ng	320504	6	15.29	773323	20.0
Tetracosane	14.97	9.7	ng	376103	6	15.29	773323	20.0
Octadecane	15.74	6.3	ng	241851	6	15.29	773323	20.0