

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042318\
 Data File : BF104714.D
 Acq On : 23 Apr 2018 17:21
 Operator : JU/SJ
 Sample : J2500-01
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RBR-400028-2171

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:\HPCHEM1\BNA_F\METHODS\8270-BF040618.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	5.016	494	498	504	rBV	62190	81935	2.83%	0.227%
2	5.434	564	569	572	rBV	2232856	2132862	73.79%	5.916%
3	6.451	738	742	757	rBV	2069169	2235649	77.35%	6.201%
4	6.581	760	764	767	rBV	2455827	2178853	75.38%	6.044%
5	6.810	799	803	806	rBV	757170	667114	23.08%	1.850%
6	6.963	825	829	833	rBV	2088138	1733421	59.97%	4.808%
7	7.375	894	899	902	rBV	1379811	1176430	40.70%	3.263%
8	8.092	1017	1021	1024	rBV	1016508	817794	28.29%	2.268%
9	9.169	1200	1204	1207	rBV	2349676	2080197	71.97%	5.770%
10	9.563	1267	1271	1278	rVB	143181	142116	4.92%	0.394%
11	9.769	1303	1306	1310	rVB	85020	65689	2.27%	0.182%
12	9.851	1316	1320	1323	rBV	931211	813961	28.16%	2.258%
13	9.881	1323	1325	1334	rVB	93773	79970	2.77%	0.222%
14	10.057	1352	1355	1358	rVB2	44591	47065	1.63%	0.131%
15	10.092	1358	1361	1365	rBV2	39267	34309	1.19%	0.095%
16	10.245	1384	1387	1389	rBV	33731	37153	1.29%	0.103%
17	10.269	1389	1391	1394	rVB	114738	89201	3.09%	0.247%
18	10.398	1409	1413	1417	rVB2	60753	62427	2.16%	0.173%
19	10.551	1435	1439	1447	rVB3	36271	52679	1.82%	0.146%
20	10.639	1451	1454	1461	rBV	1123202	1114922	38.57%	3.093%
21	10.745	1469	1472	1481	rVB3	93471	120532	4.17%	0.334%
22	10.945	1500	1506	1509	rBV4	37652	51698	1.79%	0.143%
23	11.075	1523	1528	1530	rBV3	35986	46690	1.62%	0.130%
24	11.092	1530	1531	1537	rVV2	36271	40793	1.41%	0.113%
25	11.145	1537	1540	1544	rVV2	46726	42787	1.48%	0.119%
26	11.192	1544	1548	1551	rVV2	81219	87799	3.04%	0.244%
27	11.233	1551	1555	1561	rVB2	55641	70101	2.43%	0.194%
28	11.339	1569	1573	1575	rBV	897980	746816	25.84%	2.072%
29	11.363	1575	1577	1581	rVV	1048272	836648	28.95%	2.321%
30	11.416	1583	1586	1589	rVV	332881	269676	9.33%	0.748%
31	11.439	1589	1590	1597	rVB3	22136	35009	1.21%	0.097%
32	11.575	1606	1613	1618	rBV3	50687	88334	3.06%	0.245%
33	11.622	1618	1621	1627	rVV3	41184	69312	2.40%	0.192%
34	11.680	1627	1631	1635	rVB4	29966	43848	1.52%	0.122%

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Integrator: RTE

Smoothing : ON

Sampling : 1

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Stop Thrs : 0

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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:\HPCHEM1\BNA_F\METHODS\8270-BF040618.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	11.722	1635	1638	1642	rBV	1020966	844413	29.21%	2.342%
36	11.839	1654	1658	1661	rBV	141870	132074	4.57%	0.366%
37	11.869	1661	1663	1667	rVV	206948	172625	5.97%	0.479%
38	11.916	1669	1671	1674	rVV	97085	80961	2.80%	0.225%
39	11.957	1674	1678	1680	rVV	207878	232935	8.06%	0.646%
40	11.975	1680	1681	1687	rVB	101961	74653	2.58%	0.207%
41	12.027	1687	1690	1692	rBV2	47896	45999	1.59%	0.128%
42	12.139	1705	1709	1711	rBV	116825	107986	3.74%	0.300%
43	12.169	1711	1714	1718	rVB	54120	51148	1.77%	0.142%
44	12.286	1731	1734	1737	rVV	39756	32693	1.13%	0.091%
45	12.345	1741	1744	1746	rVB	35204	31740	1.10%	0.088%
46	12.416	1752	1756	1758	rBV	2950265	2890352	100.00%	8.017%
47	12.433	1758	1759	1765	rVB2	2031828	1552425	53.71%	4.306%
48	12.480	1765	1767	1770	rVB	130962	120994	4.19%	0.336%
49	12.516	1770	1773	1776	rVB	456145	355499	12.30%	0.986%
50	12.557	1776	1780	1784	rBV	1612474	1531701	52.99%	4.249%
51	12.622	1789	1791	1794	rVB	48203	43667	1.51%	0.121%
52	12.710	1798	1806	1808	rBV3	72096	110305	3.82%	0.306%
53	12.792	1812	1820	1823	rBV2	1377552	1606355	55.58%	4.456%
54	12.833	1825	1827	1832	rVB2	75239	86379	2.99%	0.240%
55	12.927	1836	1843	1846	rBV	1893124	1755845	60.75%	4.870%
56	13.004	1851	1856	1860	rBV2	114069	182737	6.32%	0.507%
57	13.116	1867	1875	1878	rBV2	162325	298686	10.33%	0.828%
58	13.180	1884	1886	1888	rVB	70510	48212	1.67%	0.134%
59	13.216	1888	1892	1895	rBV2	126829	157041	5.43%	0.436%
60	13.245	1895	1897	1900	rVB2	57335	47638	1.65%	0.132%
61	13.310	1905	1908	1911	rBV	72419	60450	2.09%	0.168%
62	13.339	1911	1913	1917	rBV2	70126	78358	2.71%	0.217%
63	13.492	1936	1939	1941	rBV4	35614	47049	1.63%	0.131%
64	13.621	1959	1961	1965	rVB	107559	104101	3.60%	0.289%
65	13.733	1976	1980	1982	rBV2	100850	108273	3.75%	0.300%
66	13.763	1982	1985	1987	rVB	96586	68308	2.36%	0.189%
67	13.786	1987	1989	1991	rBV	84007	77837	2.69%	0.216%
68	13.816	1991	1994	1996	rBV2	98911	94011	3.25%	0.261%
69	13.980	2015	2022	2026	rBV2	835641	1313360	45.44%	3.643%
70	14.016	2026	2028	2035	rVB	552271	468396	16.21%	1.299%
71	14.092	2035	2041	2044	rBV2	106403	133452	4.62%	0.370%

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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 >
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Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040618.M
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72	14.380	2087	2090	2093	rVB	95780	80120	2.77%	0.222%
73	14.416	2093	2096	2098	rBV	42998	43955	1.52%	0.122%
74	14.451	2099	2102	2103	rBV3	39884	41613	1.44%	0.115%
75	14.510	2109	2112	2113	rBV	37647	36326	1.26%	0.101%
76	15.039	2197	2202	2205	rBV	389872	611028	21.14%	1.695%
77	15.145	2217	2220	2224	rVB	75433	87374	3.02%	0.242%
78	15.339	2249	2253	2259	rVB	223890	291435	10.08%	0.808%
79	15.398	2259	2263	2269	rBV	335937	441504	15.28%	1.225%
80	15.463	2270	2274	2277	rVV	319310	418626	14.48%	1.161%
81	15.492	2277	2279	2286	rVB	108572	149410	5.17%	0.414%
82	16.927	2518	2523	2532	rVB2	140641	272380	9.42%	0.756%
83	17.368	2594	2598	2611	rVB	100964	235511	8.15%	0.653%

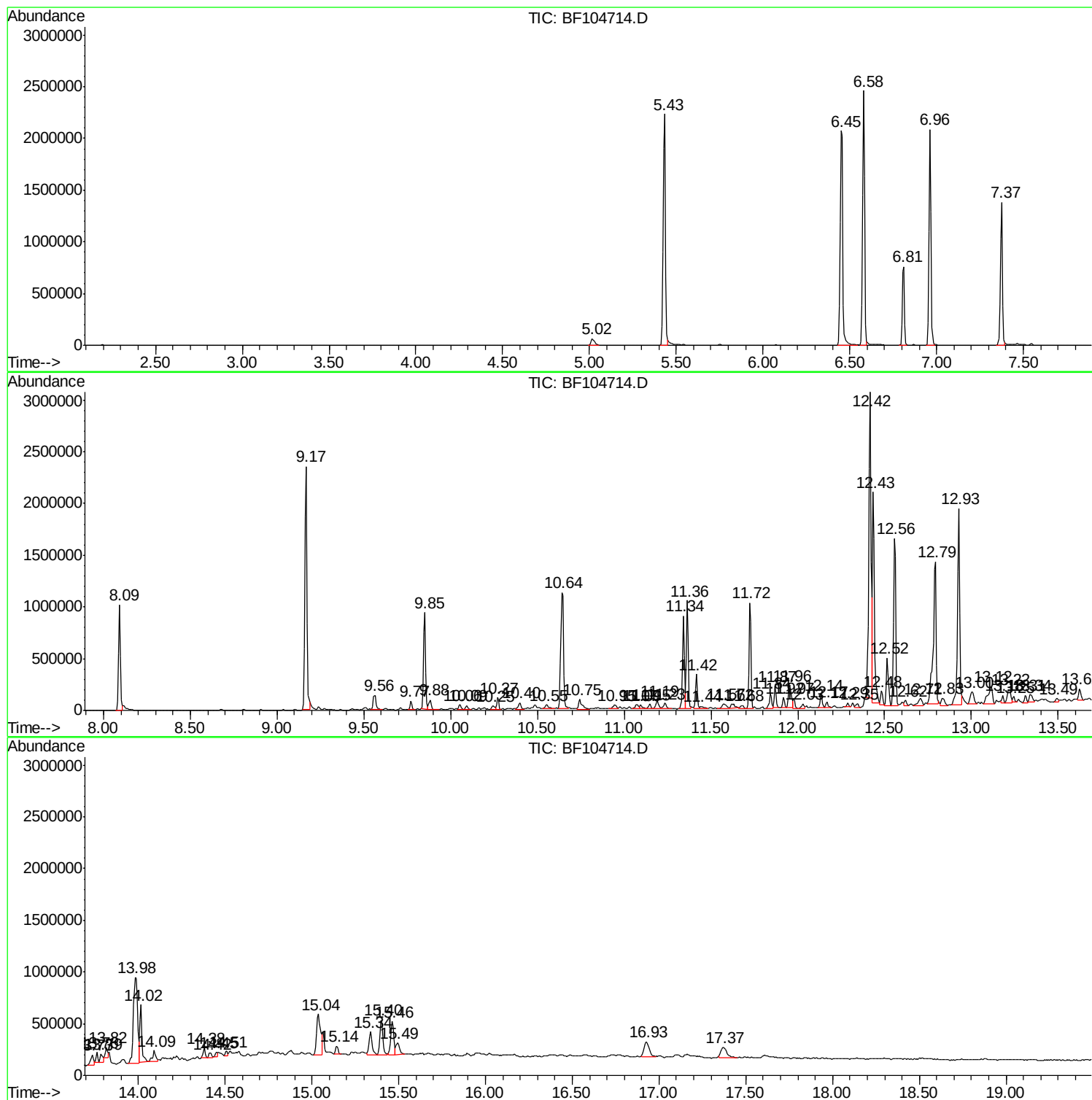
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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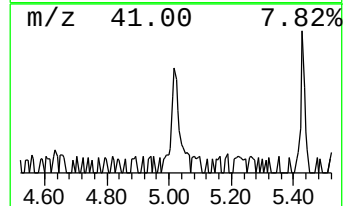
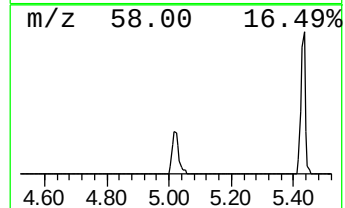
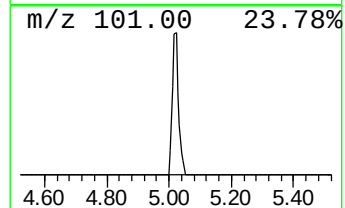
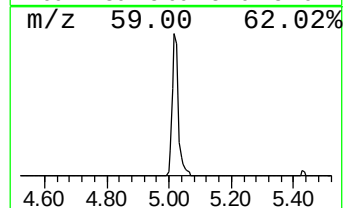
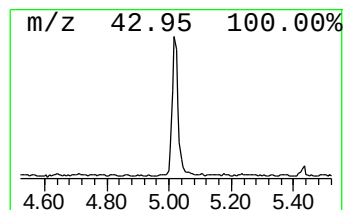
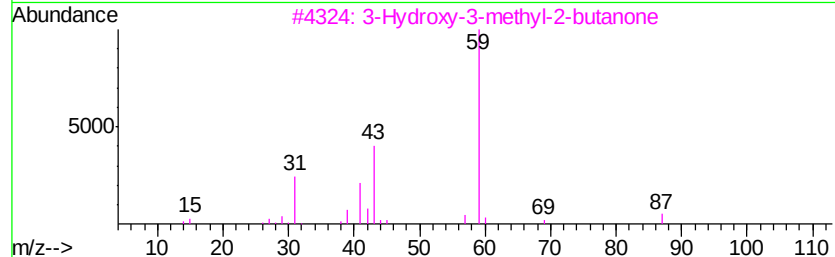
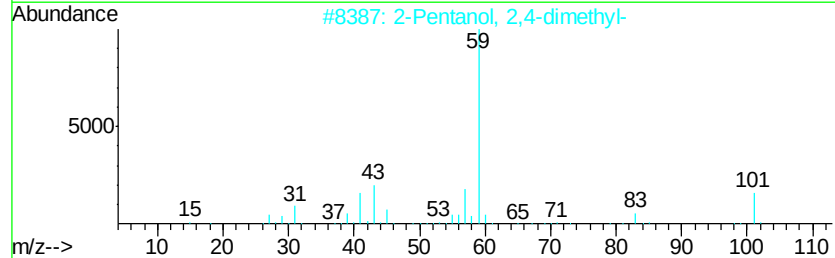
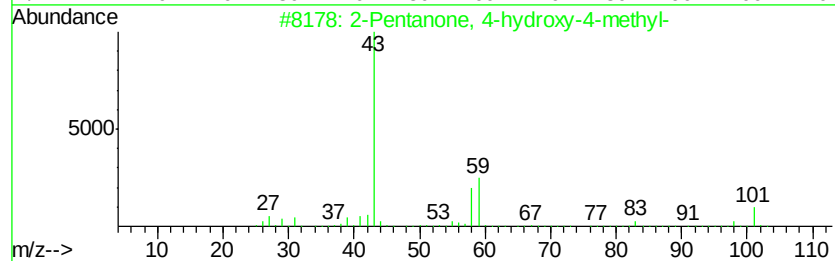
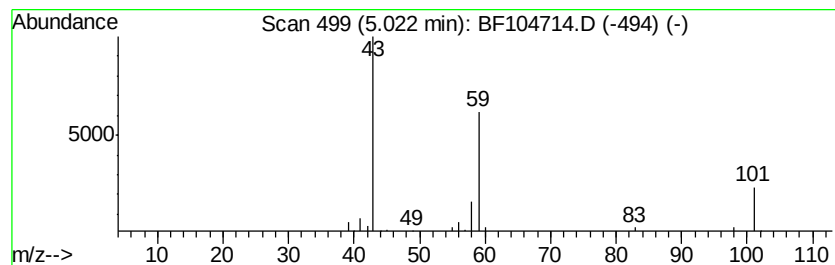
Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.	
5.02	2.46 ng	81935	1,4-Dichlorobenzene-d4	6.81	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2	2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	25
3	3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	17
4	Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	16
5	1-Butanol, 3-methoxy-	104	C5H12O2	002517-43-3	16



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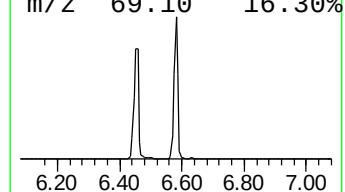
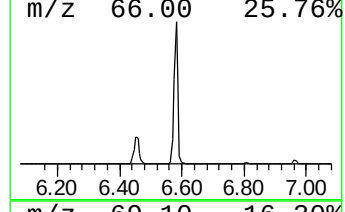
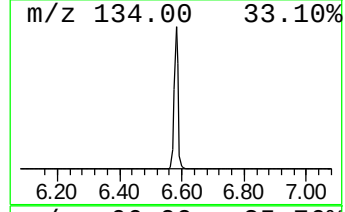
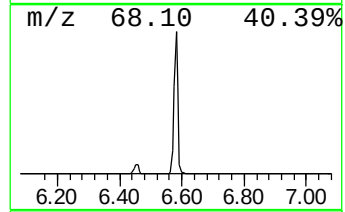
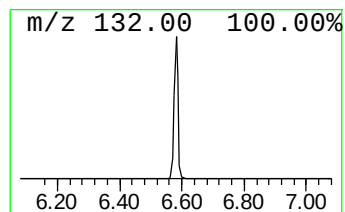
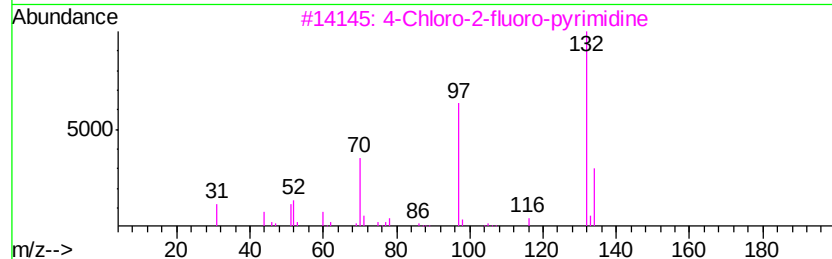
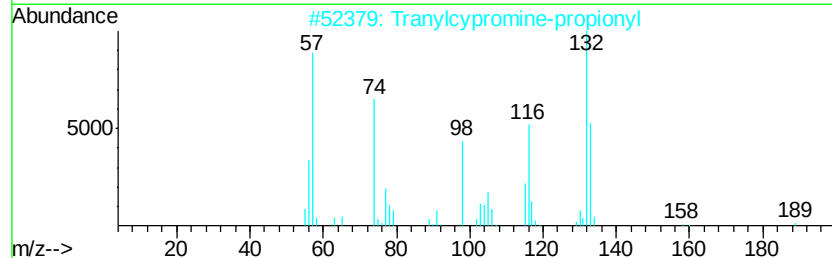
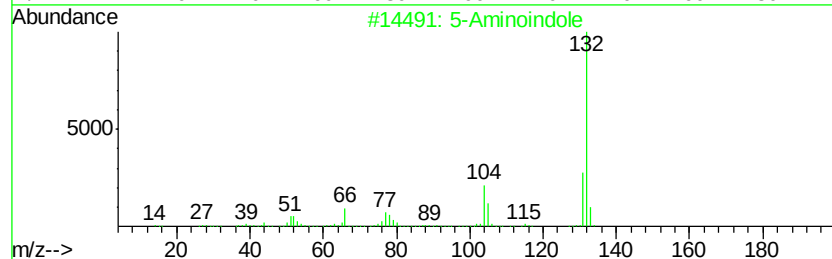
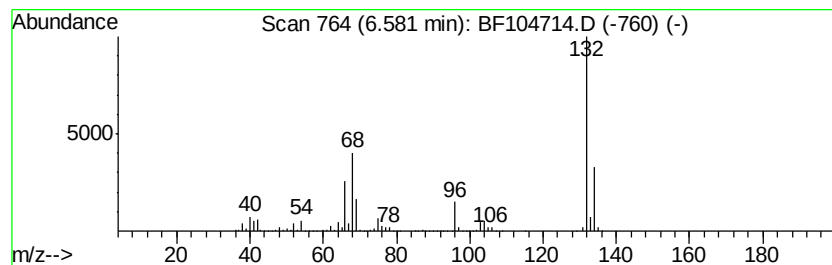
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040618.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.58 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.58	65.32 ng	2178850	1,4-Dichlorobenzene-d4	6.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Aminoindole	132	C8H8N2	005192-03-0	12
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
4		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9
5		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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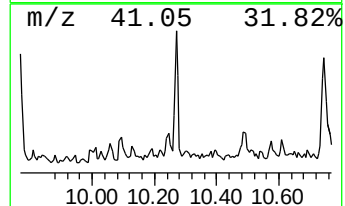
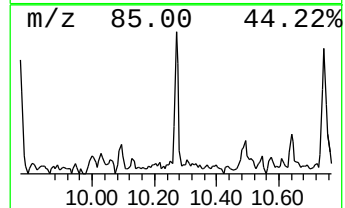
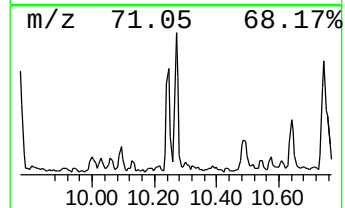
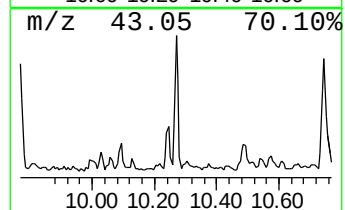
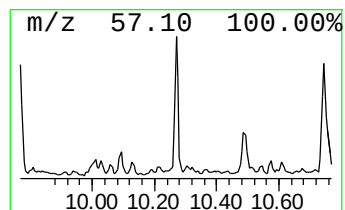
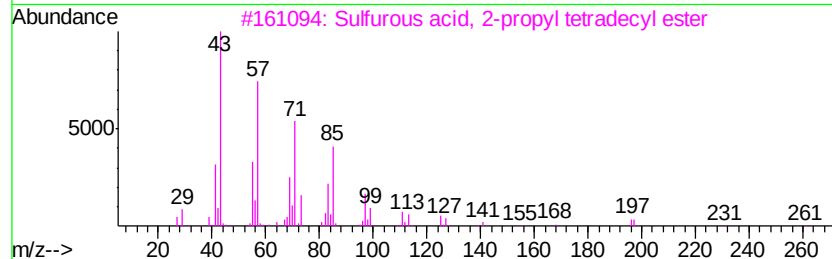
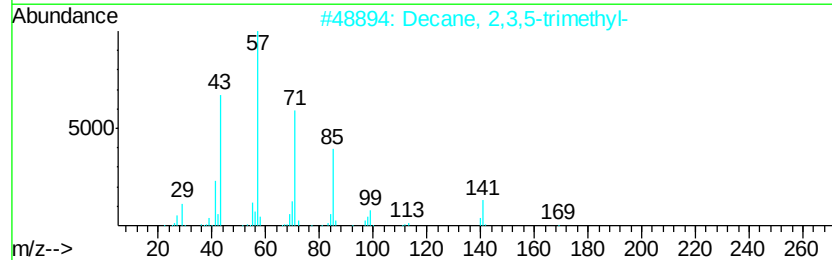
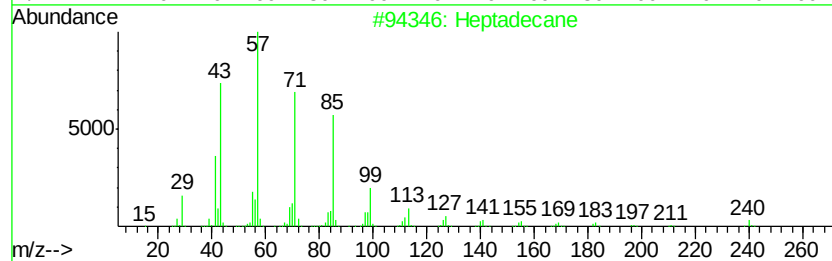
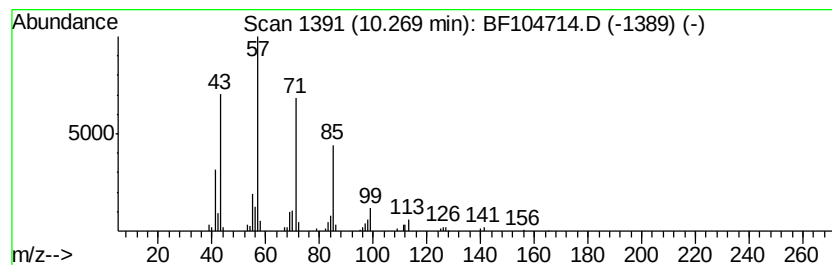
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Heptadecane Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.27	2.19 ng	89201	Acenaphthene-d10	9.85

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane		240	C17H36	000629-78-7	91
2	Decane, 2,3,5-trimethyl-		184	C13H28	062238-11-3	90
3	Sulfurous acid, 2-propyl tetradecyl-		320	C17H36O3S	1000309-12-5	86
4	Eicosane		282	C20H42	000112-95-8	86
5	Pentadecane		212	C15H32	000629-62-9	86



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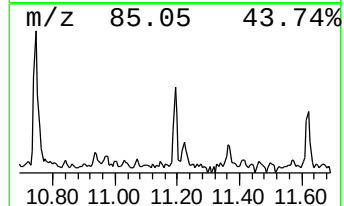
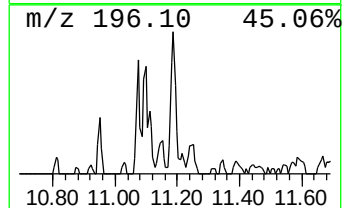
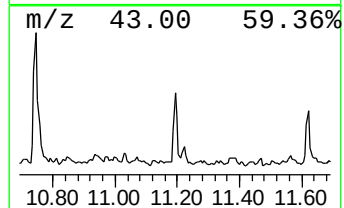
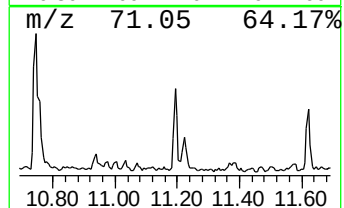
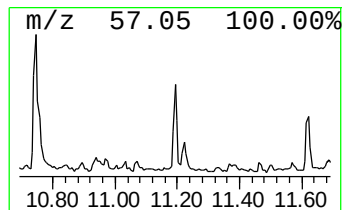
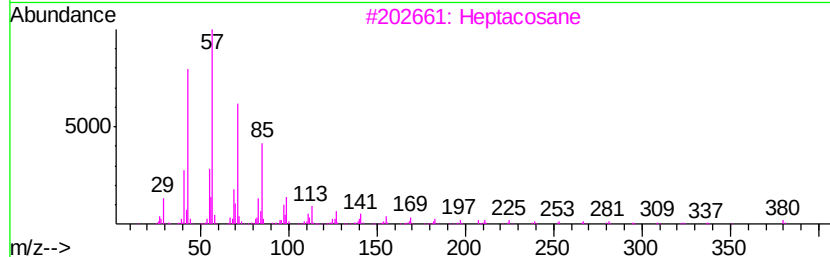
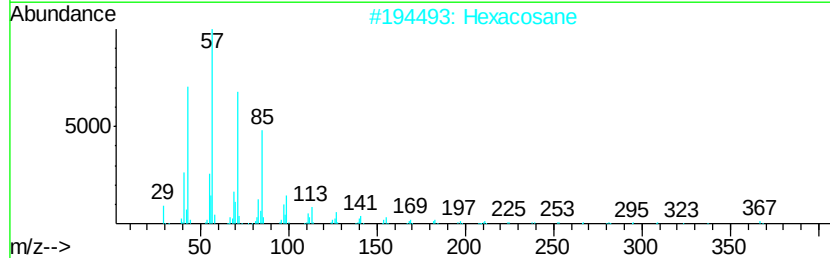
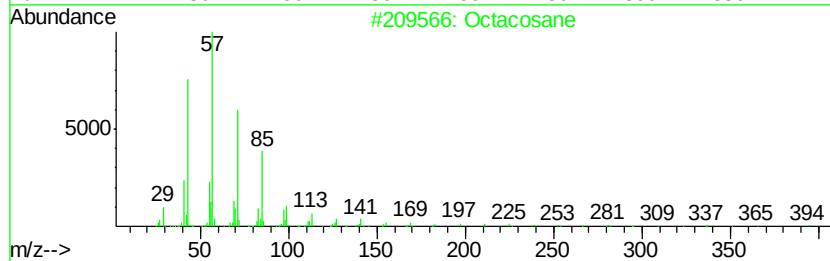
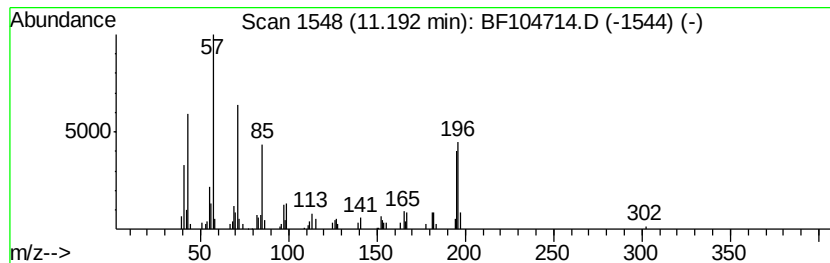
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 unknown11.19 Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.19	2.35 ng	87799	Phenanthrene-d10	11.34

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosane	394	C28H58	000630-02-4	49
2			Hexacosane	366	C26H54	000630-01-3	49
3			Heptacosane	380	C27H56	000593-49-7	49
4			Tetratetracontane	619	C44H90	007098-22-8	46
5			10-Methylnonadecane	282	C20H42	056862-62-5	46



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042318\
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Acq On : 23 Apr 2018 17:21
Operator : JU/SJ
Sample : J2500-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

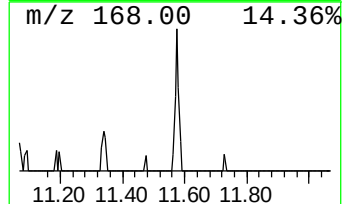
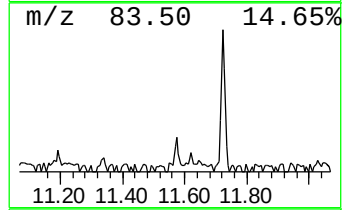
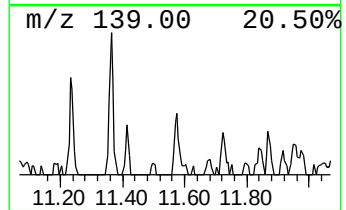
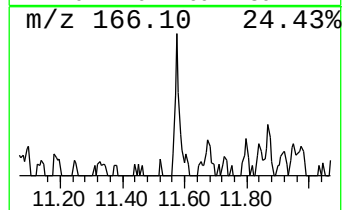
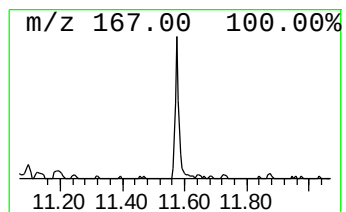
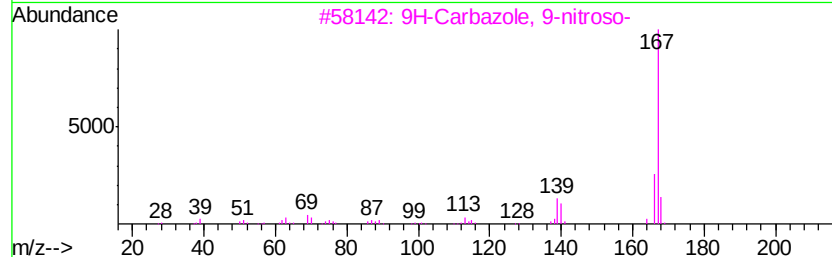
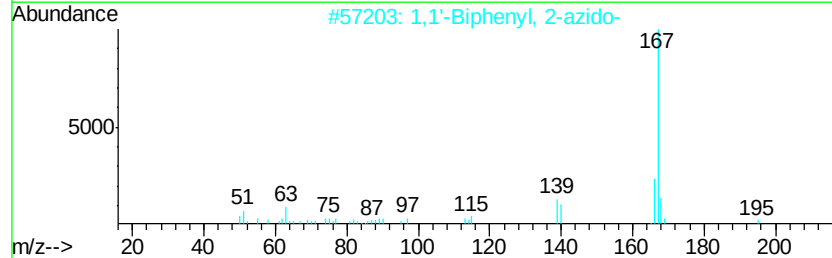
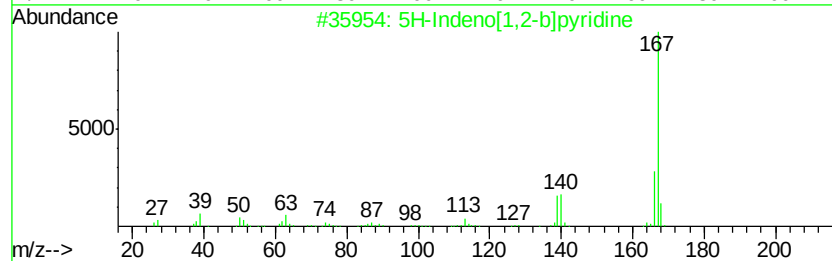
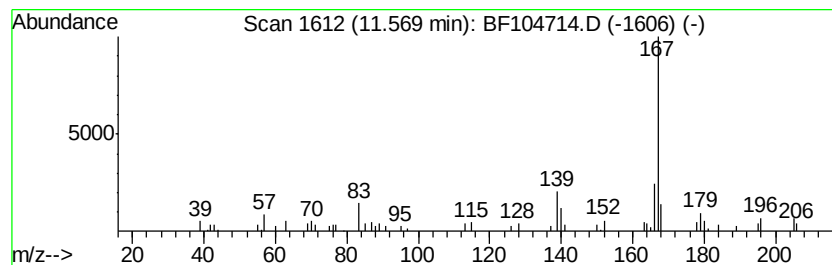
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 5H-Indeno[1,2-b]pyridine Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.57	2.37 ng	88334	Phenanthrene-d10		11.34
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5H-Indeno[1,2-b]pyridine	167	C12H9N	000244-99-5	87
2	1,1'-Biphenyl, 2-azido-	195	C12H9N3	007599-23-7	87
3	9H-Carbazole, 9-nitroso-	196	C12H8N2O	002788-23-0	83
4	9H-Carbazole-9-methanol	197	C13H11NO	002409-36-1	83
5	D-Galactitol, 2-(acetylmethylami...	363	C16H29NO8	074410-42-7	80



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042318\
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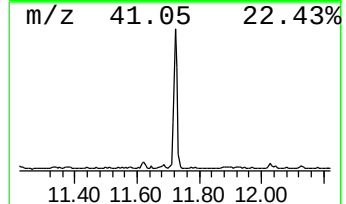
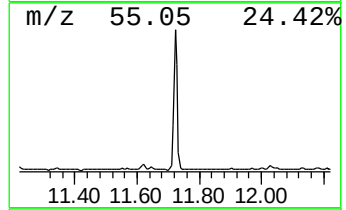
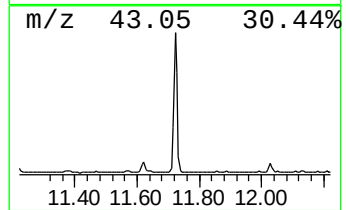
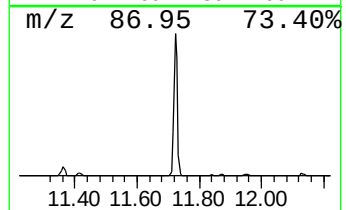
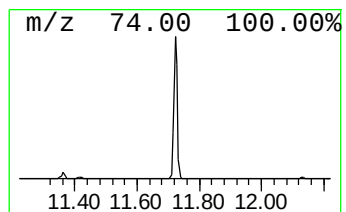
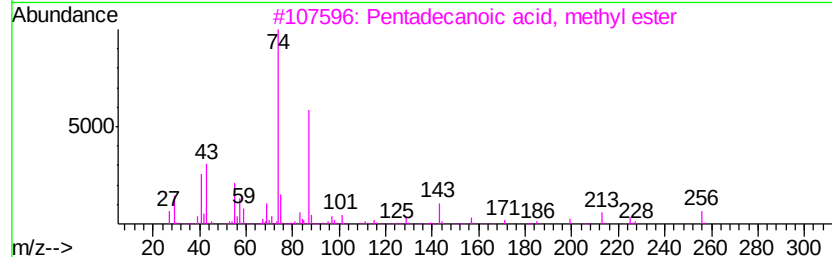
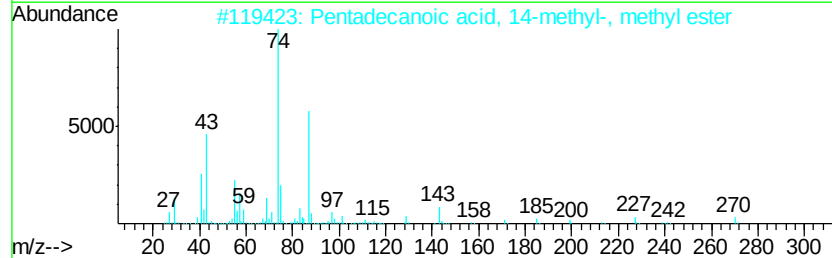
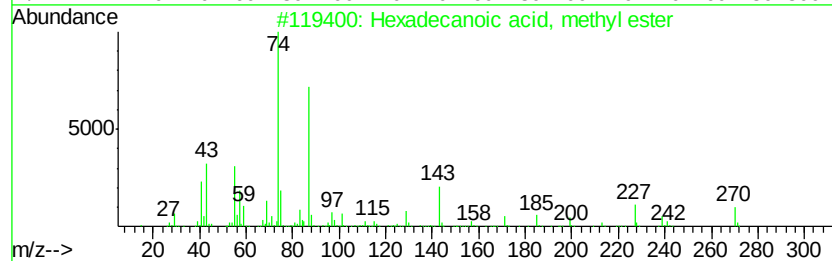
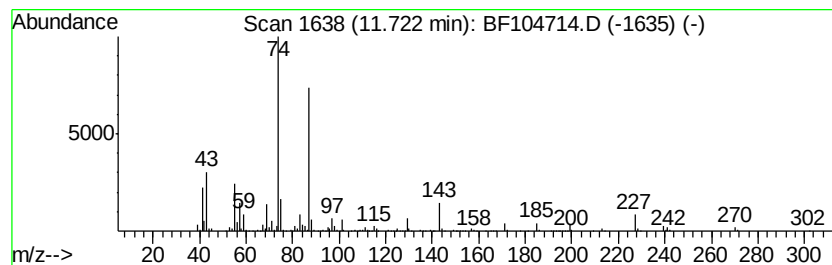
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Hexadecanoic acid, methyl e... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.72	22.61 ng	844413	Phenanthrene-d10	11.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	97
2		Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	91
3		Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	90
4		Hexadecanoic acid, 2-methyl-	270	C17H34O2	027147-71-3	90
5		Dodecanoic acid, methyl ester	214	C13H26O2	000111-82-0	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042318\
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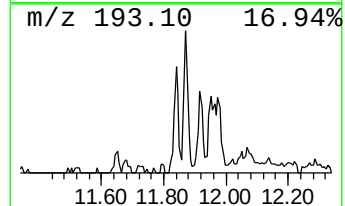
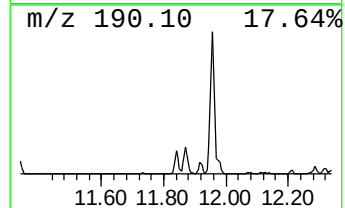
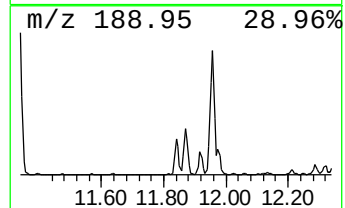
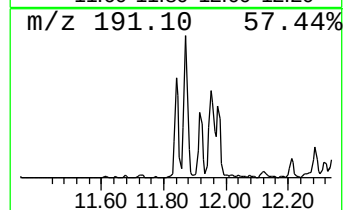
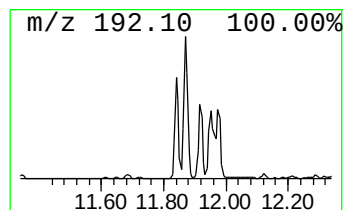
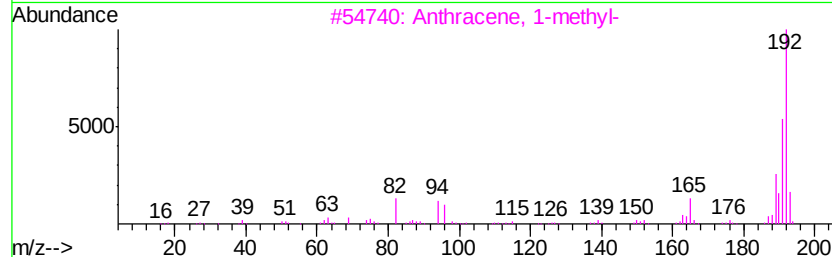
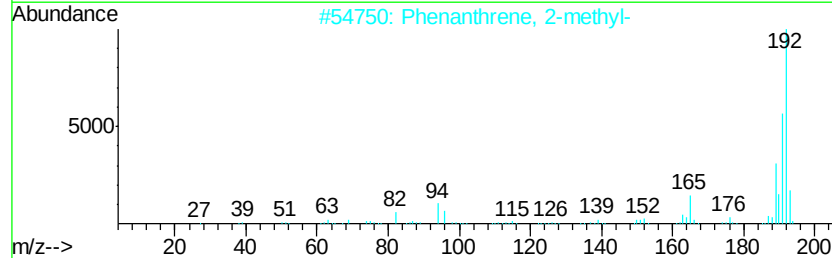
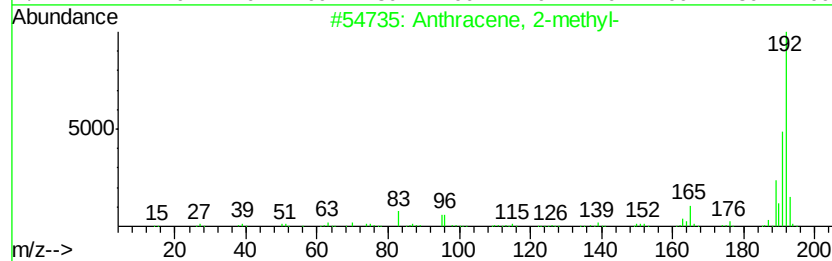
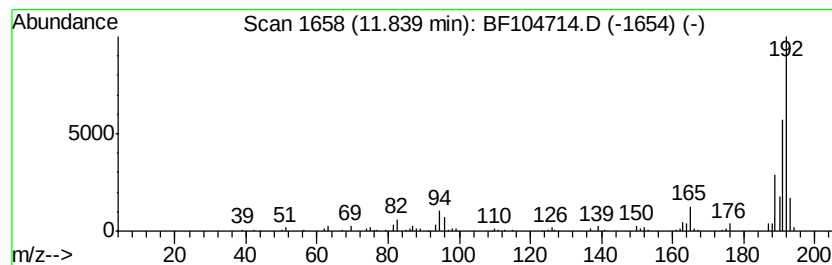
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Anthracene, 2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.84	3.54 ng	132074	Phenanthrene-d10		11.34
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
3	Anthracene, 1-methyl-	192	C15H12	000610-48-0	93
4	Anthracene, 9-methyl-	192	C15H12	000779-02-2	91
5	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	91



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042318\
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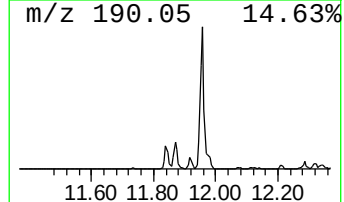
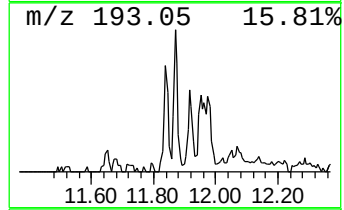
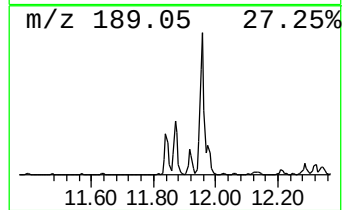
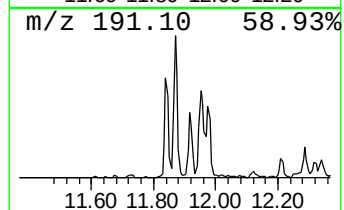
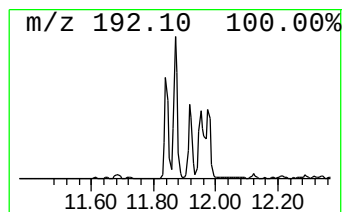
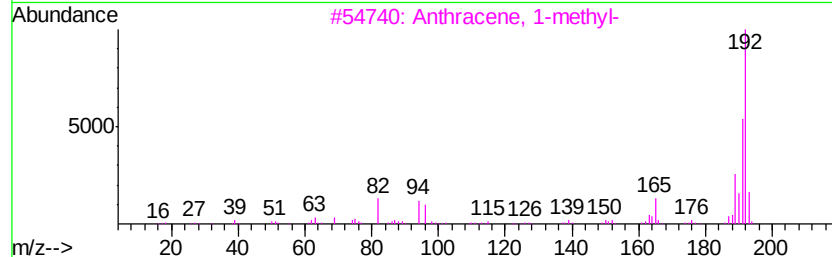
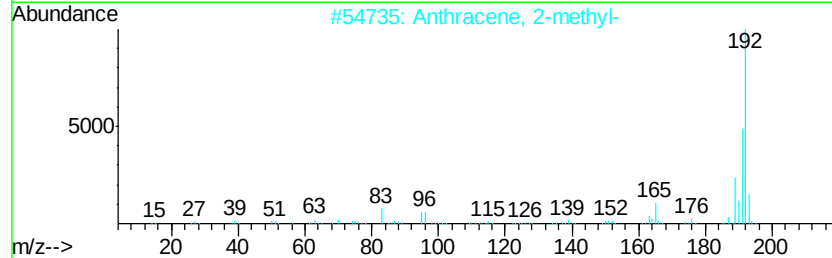
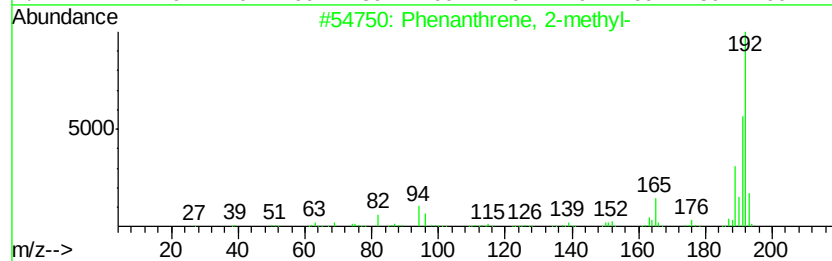
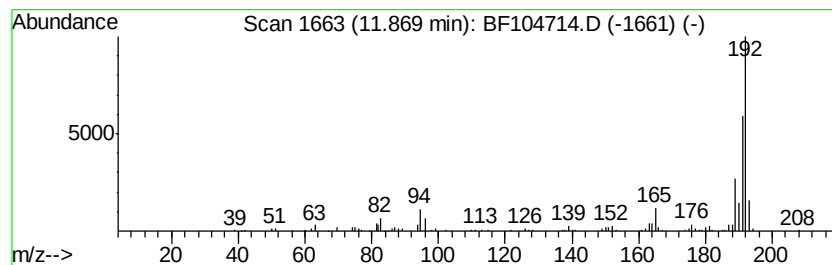
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Phenanthrene, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.87	4.62 ng	172625	Phenanthrene-d10	11.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	98
2		Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
3		Anthracene, 1-methyl-	192	C15H12	000610-48-0	95
4		Anthracene, 9-methyl-	192	C15H12	000779-02-2	95
5		1H-Cyclopropa[l]phenanthrene, 1a, ...	192	C15H12	000949-41-7	94



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042318\
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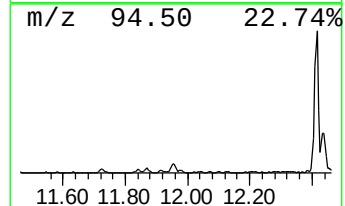
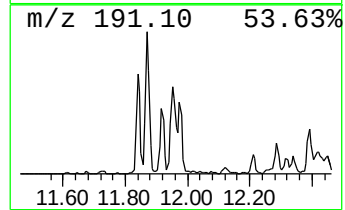
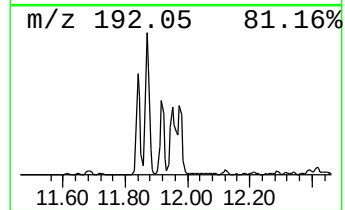
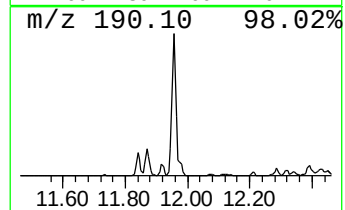
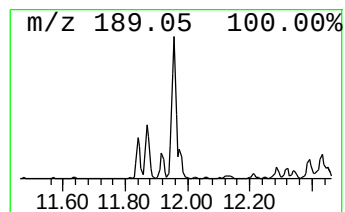
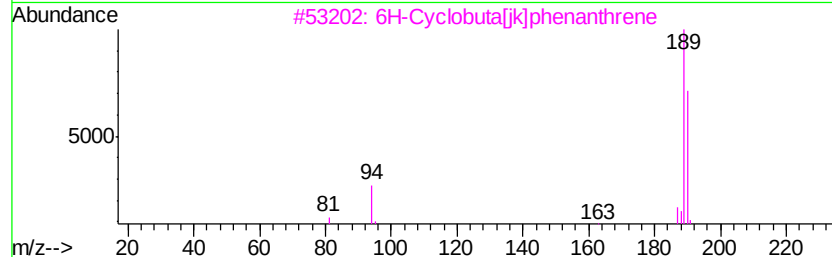
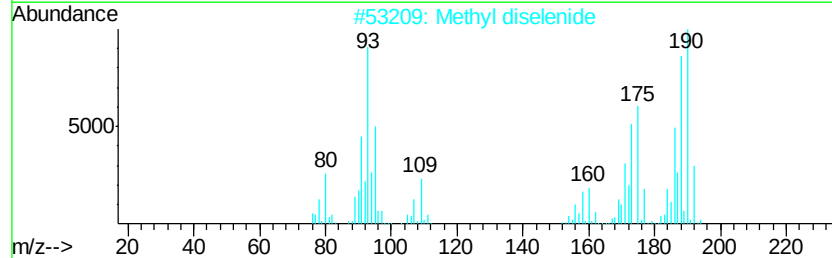
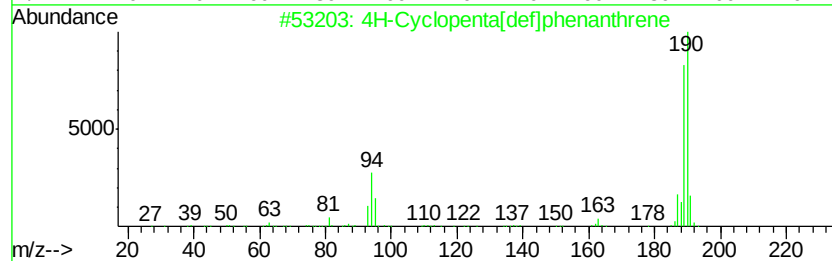
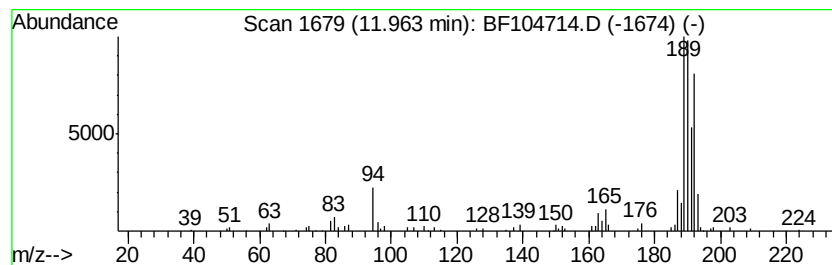
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 4H-Cyclopenta[def]phenanthrene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.96	6.24 ng	232935	Phenanthrene-d10		11.34
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	94
2	Methyl diselenide	190	C2H6Se2	007101-31-7	43
3	6H-Cyclobuta[f]phenanthrene	190	C15H10	083469-43-6	42
4	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	40
5	2,6-Dichlorobenzaldoxime	189	C7H5Cl2NO	025185-95-9	30



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042318\
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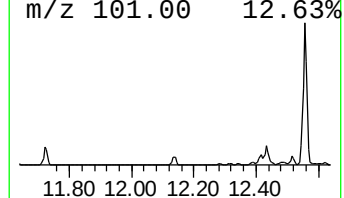
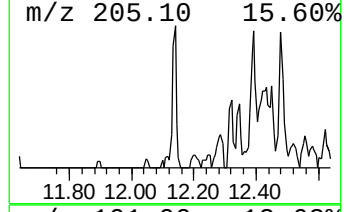
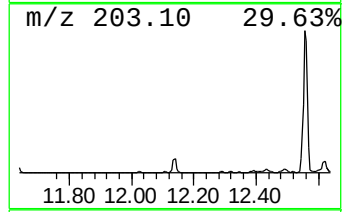
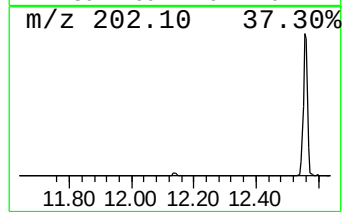
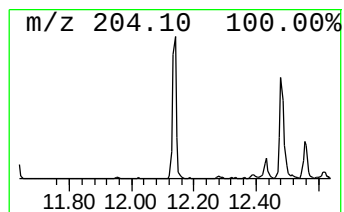
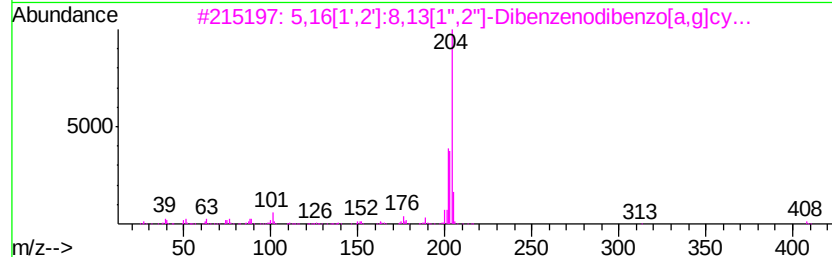
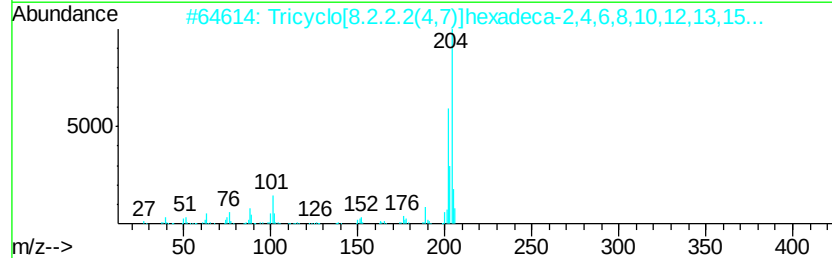
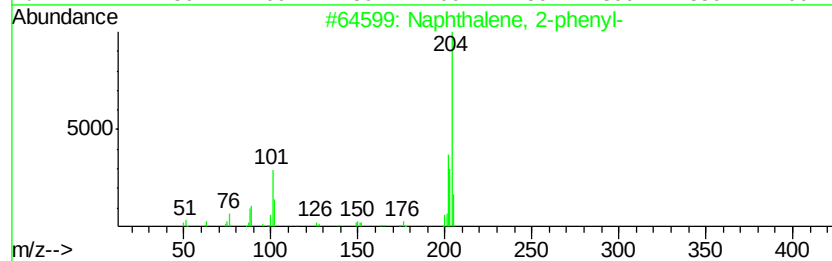
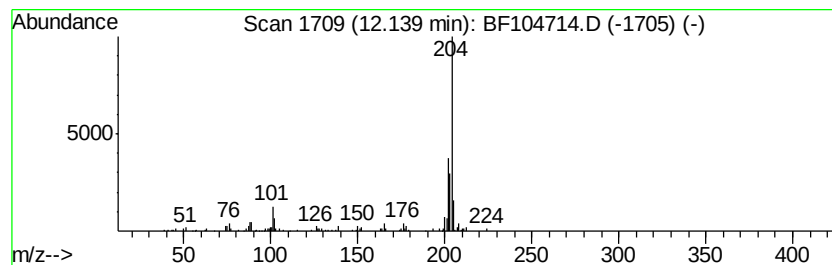
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Naphthalene, 2-phenyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.14	2.89 ng	107986	Phenanthrene-d10	11.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	90
2		Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	81
3		5,16[1',2']8,13[1',2']-Dibenz...	408	C32H24	005672-97-9	80
4		Benzene, 1,1'-(1-buten-3-yne-1,4...	204	C16H12	013141-45-2	60
5		9,10-Bis(bromomethyl)anthracene	362	C16H12Br2	034373-96-1	58



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042318\
Data File : BF104714.D
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ALS Vial : 17 Sample Multiplier: 1

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ClientSampleId :
RBR-400028-2171

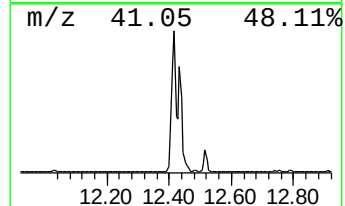
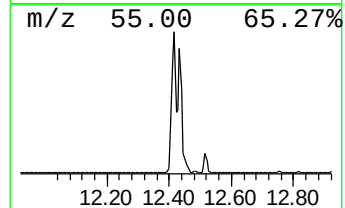
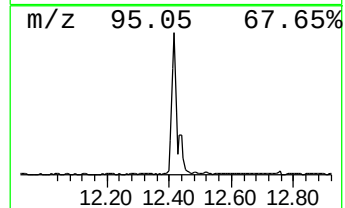
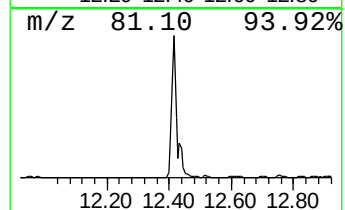
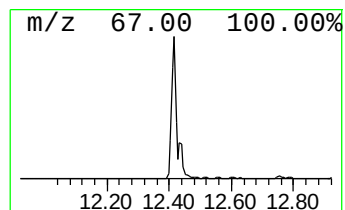
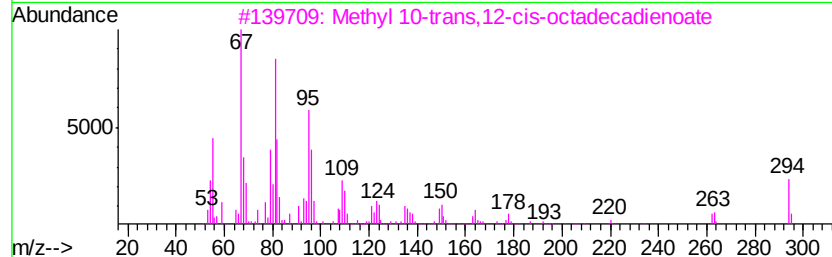
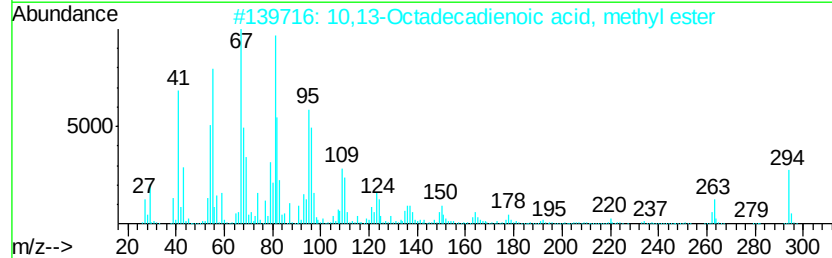
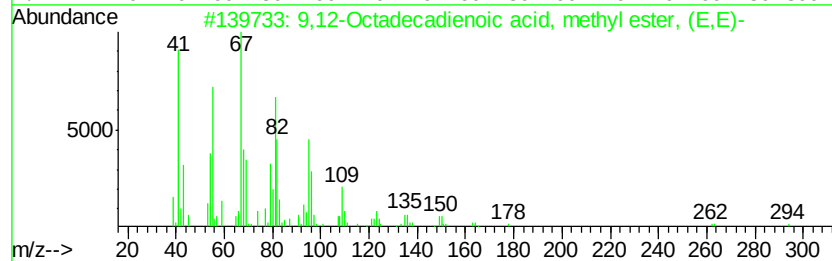
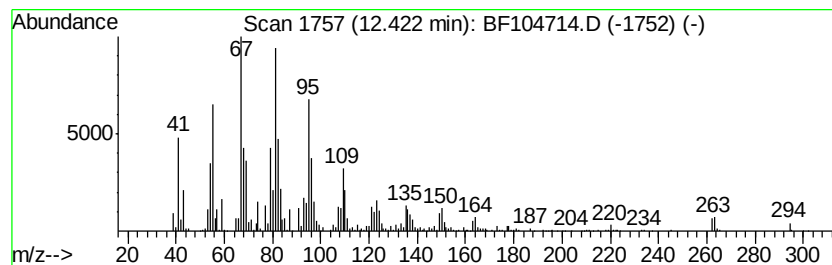
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 9,12-Octadecadienoic acid, ... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.42	77.40 ng	2890350	Phenanthrene-d10	11.34

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	99
2			10,13-Octadecadienoic acid, meth...	294	C19H34O2	056554-62-2	99
3			Methyl 10-trans,12-cis-octadecad...	294	C19H34O2	1000336-44-2	99
4			9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	99
5			9,15-Octadecadienoic acid, methy...	294	C19H34O2	017309-05-6	99



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042318\
Data File : BF104714.D
Acq On : 23 Apr 2018 17:21
Operator : JU/SJ
Sample : J2500-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
RBR-400028-2171

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 9-Octadecenoic acid (Z)-, m... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.43	41.57 ng	1552430	Phenanthrene-d10	11.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	99
2		10-Octadecenoic acid, methyl ester	296	C19H36O2	013481-95-3	99
3		9-Octadecenoic acid, methyl este...	296	C19H36O2	001937-62-8	99
4		cis-13-Octadecenoic acid, methyl...	296	C19H36O2	1000333-58-3	99
5		trans-13-Octadecenoic acid, meth...	296	C19H36O2	1000333-61-3	99

