

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121518\
 Data File : BF111474.D
 Acq On : 15 Dec 2018 17:48
 Operator : JU/SJ
 Sample : J6361-01
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TN4803-121218

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-BF121418.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.998	491	495	507	rBV	280057	325512	8.35%	0.855%
2	5.422	562	567	570	rBV	3732777	3369972	86.47%	8.856%
3	6.439	735	740	757	rBV	3513854	3897256	100.00%	10.242%
4	6.569	757	762	765	rBV	4058969	3571986	91.65%	9.387%
5	6.792	797	800	804	rBV	1223141	959192	24.61%	2.521%
6	6.951	823	827	830	rBV	3444994	2814766	72.22%	7.397%
7	7.351	890	895	898	rBV	2459665	2268097	58.20%	5.961%
8	8.075	1014	1018	1025	rBV	1444066	1232028	31.61%	3.238%
9	9.151	1197	1201	1204	rBV	4238439	3494684	89.67%	9.184%
10	9.545	1264	1268	1277	rBV	203600	200052	5.13%	0.526%
11	9.828	1312	1316	1320	rBV	1354050	1209282	31.03%	3.178%
12	10.616	1446	1450	1461	rBV	1816252	1679745	43.10%	4.414%
13	10.751	1468	1473	1479	rVB4	41765	64575	1.66%	0.170%
14	10.804	1479	1482	1484	rBV	61680	53896	1.38%	0.142%
15	10.833	1484	1487	1491	rVB3	56321	57684	1.48%	0.152%
16	10.869	1491	1493	1495	rBV	55210	42063	1.08%	0.111%
17	10.980	1509	1512	1516	rVV2	65463	70699	1.81%	0.186%
18	11.016	1516	1518	1521	rVV	62553	62607	1.61%	0.165%
19	11.057	1523	1525	1531	rVB	39888	45282	1.16%	0.119%
20	11.310	1565	1568	1571	rBV	1212557	1132704	29.06%	2.977%
21	11.333	1571	1572	1576	rVV	366423	282762	7.26%	0.743%
22	11.386	1578	1581	1585	rVB	45059	48805	1.25%	0.128%
23	11.545	1603	1608	1614	rBV2	40151	59111	1.52%	0.155%
24	11.816	1651	1654	1656	rBV	48952	46799	1.20%	0.123%
25	11.845	1656	1659	1666	rVV2	195145	218584	5.61%	0.574%
26	11.927	1670	1673	1675	rBV	75618	73321	1.88%	0.193%
27	12.133	1706	1708	1714	rVB	51127	51173	1.31%	0.134%
28	12.369	1743	1748	1750	rBV3	56780	72486	1.86%	0.190%
29	12.521	1771	1774	1780	rBV	639864	576839	14.80%	1.516%
30	12.627	1788	1792	1798	rBV4	90553	155946	4.00%	0.410%
31	12.751	1807	1813	1816	rBV	582656	589241	15.12%	1.549%
32	12.898	1834	1838	1841	rBV	3291949	2828805	72.58%	7.434%
33	13.063	1862	1866	1867	rBV4	31673	41349	1.06%	0.109%
34	13.080	1867	1869	1872	rVB	80882	63700	1.63%	0.167%

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 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	13.145	1874	1880	1883	rBV2	59648	77618	1.99%	0.204%
36	13.186	1883	1887	1889	rBV2	44793	49007	1.26%	0.129%
37	13.310	1905	1908	1910	rBV	37388	47100	1.21%	0.124%
38	13.586	1952	1955	1959	rVB2	46877	64003	1.64%	0.168%
39	13.698	1971	1974	1977	rBV3	38446	48964	1.26%	0.129%
40	13.751	1981	1983	1988	rBV3	47291	52489	1.35%	0.138%
41	13.804	1988	1992	1999	rBV3	175435	255969	6.57%	0.673%
42	13.951	2011	2017	2019	rBV2	1322812	1475800	37.87%	3.878%
43	13.974	2019	2021	2028	rVB	335357	297732	7.64%	0.782%
44	14.057	2032	2035	2038	rBV	65626	58329	1.50%	0.153%
45	14.121	2043	2046	2051	rVB2	122830	121155	3.11%	0.318%
46	14.374	2086	2089	2092	rBV2	45182	58561	1.50%	0.154%
47	14.427	2095	2098	2101	rVB	257076	212261	5.45%	0.558%
48	14.727	2146	2149	2154	rVB	338716	315104	8.09%	0.828%
49	14.974	2187	2191	2201	rVB	257430	512430	13.15%	1.347%
50	15.057	2201	2205	2208	rBV	459572	491508	12.61%	1.292%
51	15.268	2238	2241	2246	rVB	131428	162504	4.17%	0.427%
52	15.327	2247	2251	2256	rVV	155249	179551	4.61%	0.472%
53	15.386	2257	2261	2265	rVV	648603	863002	22.14%	2.268%
54	15.427	2265	2268	2277	rVB2	354339	452594	11.61%	1.189%
55	15.851	2336	2340	2345	rVB	309550	433444	11.12%	1.139%
56	16.339	2419	2423	2428	rBV2	105800	160975	4.13%	0.423%

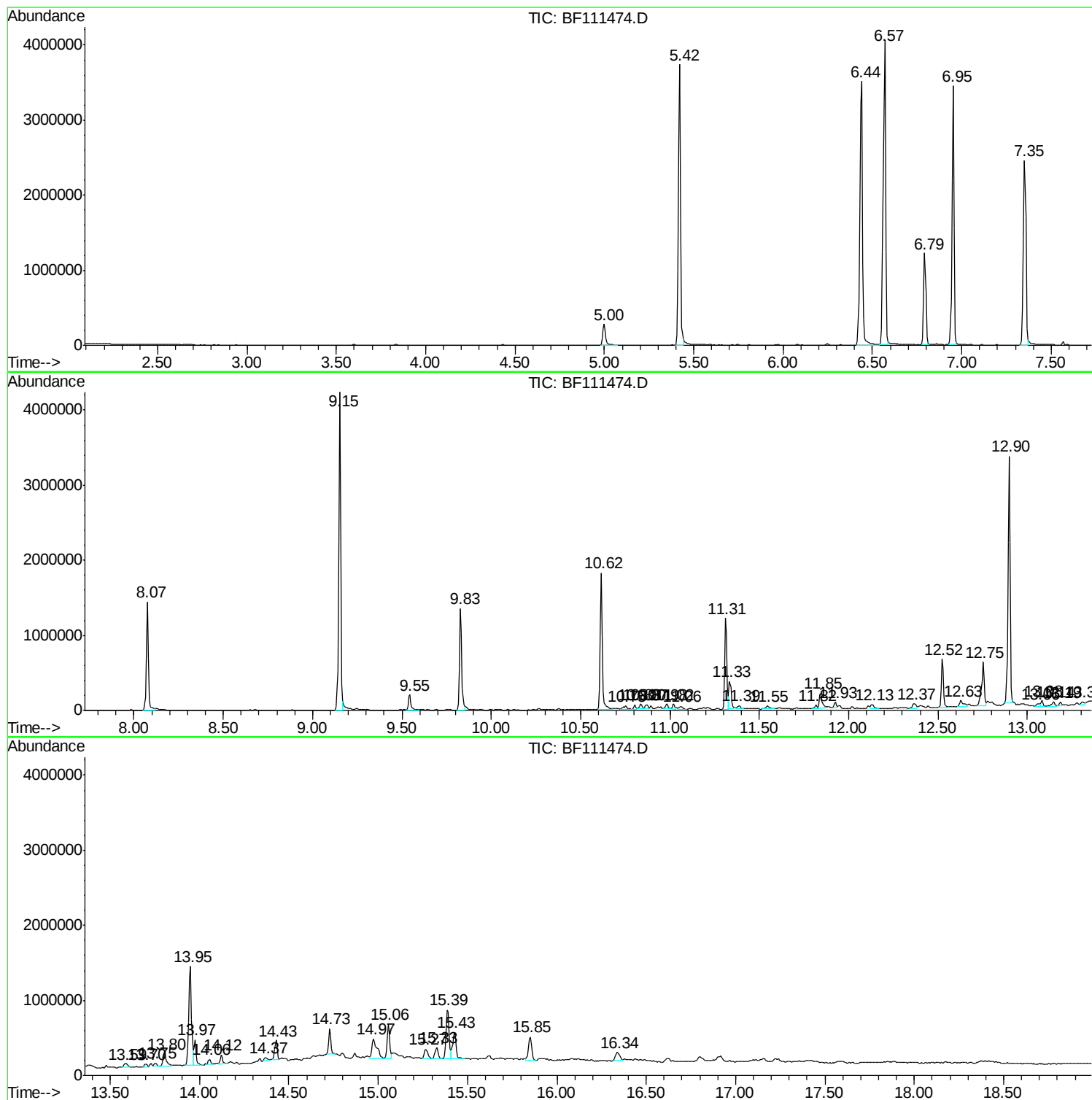
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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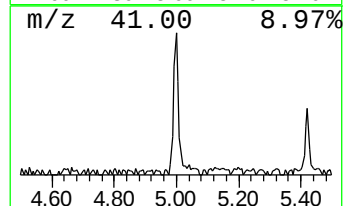
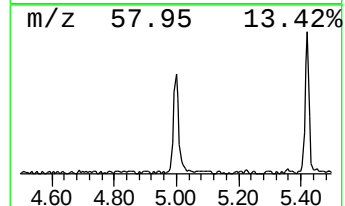
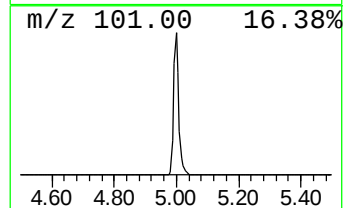
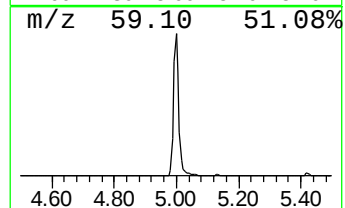
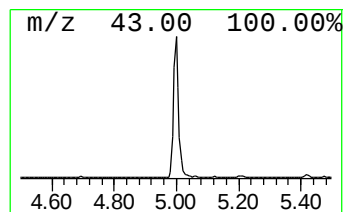
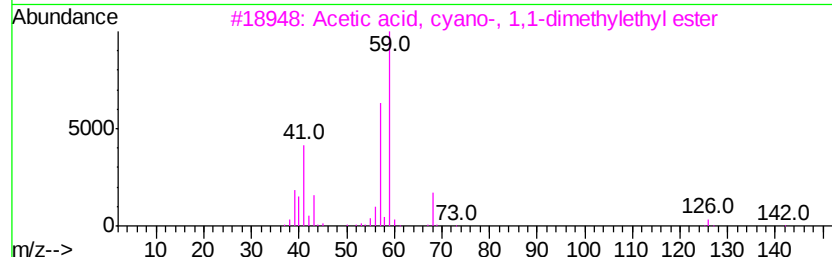
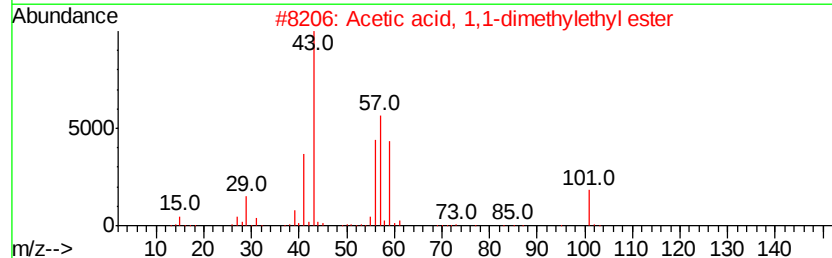
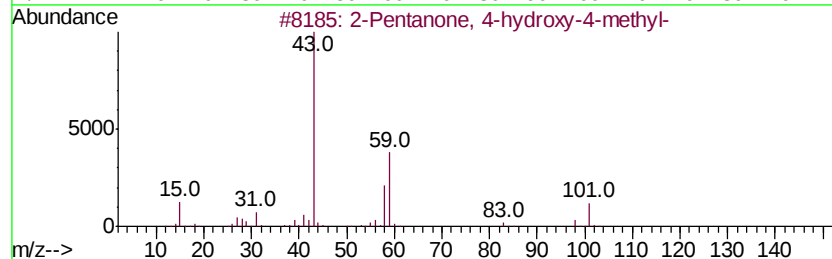
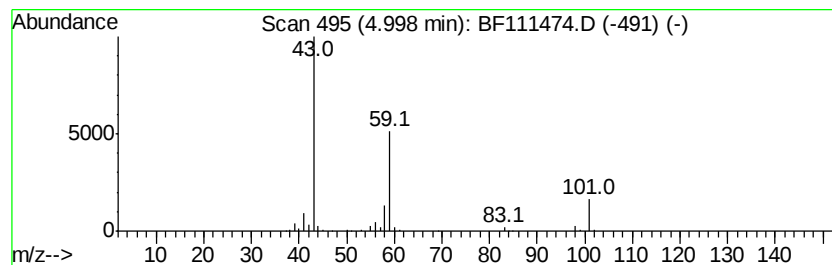
Instrument :
BNA_F
ClientSampleId :
TN4803-121218

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.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 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
5.00	6.79 ng	325512	1,4-Dichlorobenzene-d4	6.79	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
3	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5	3-Hexanol	102	C6H14O	000623-37-0	9



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Instrument :
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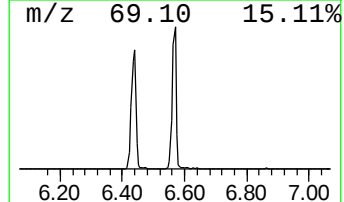
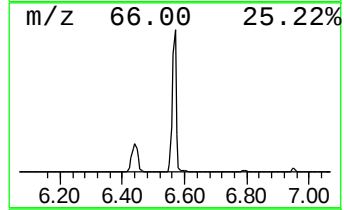
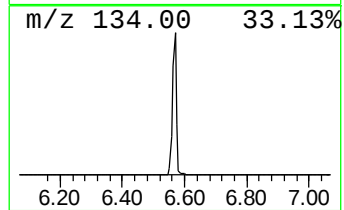
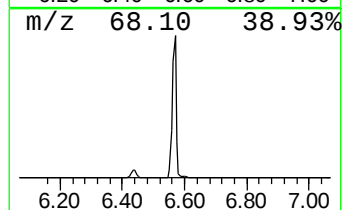
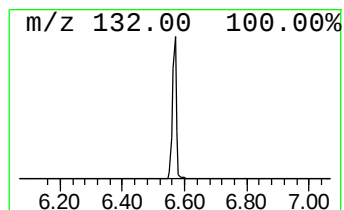
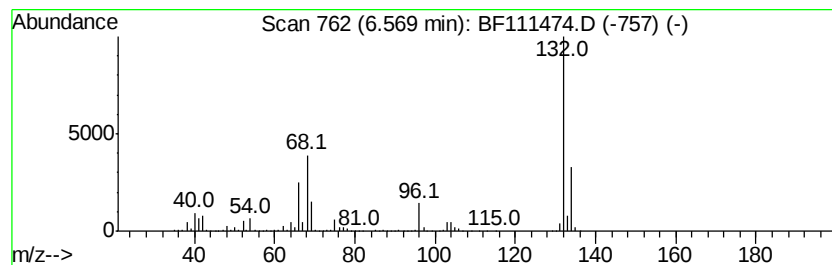
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 unknown6.57 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.57	74.48 ng	3571990	1,4-Dichlorobenzene-d4	6.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	12
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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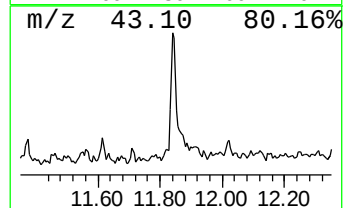
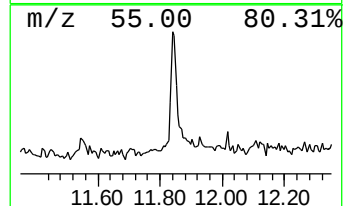
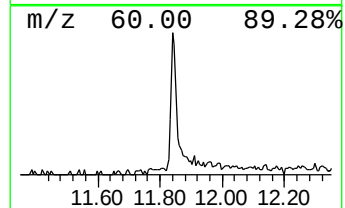
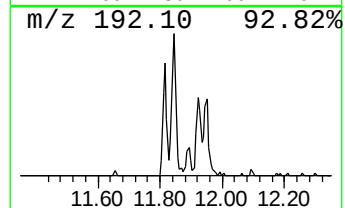
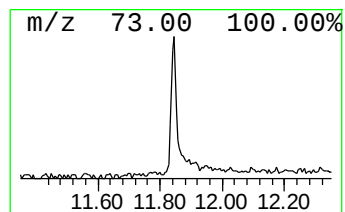
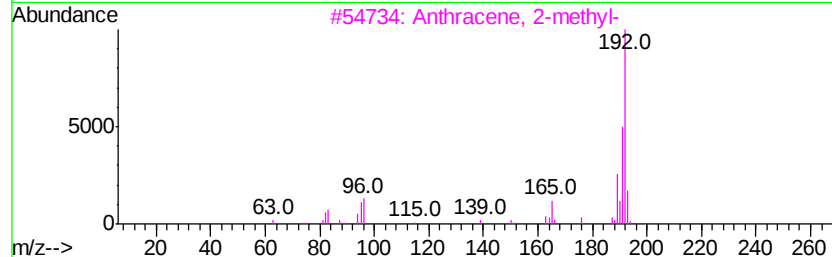
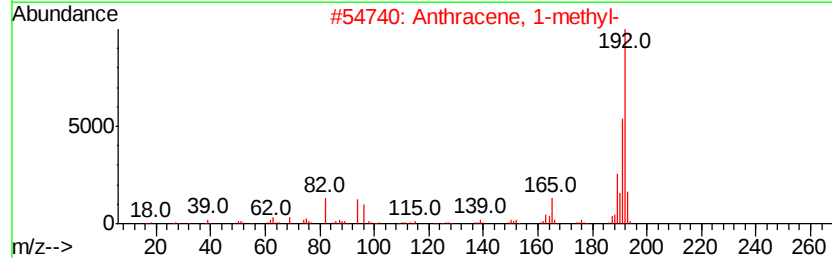
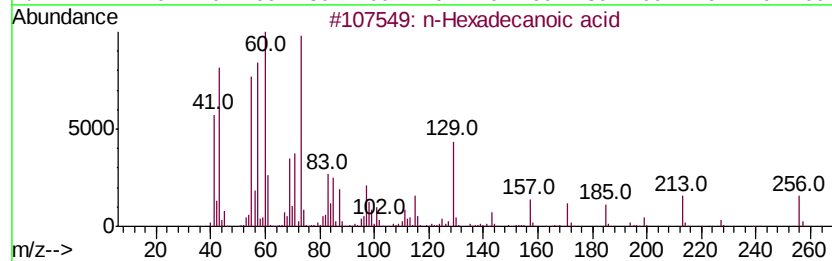
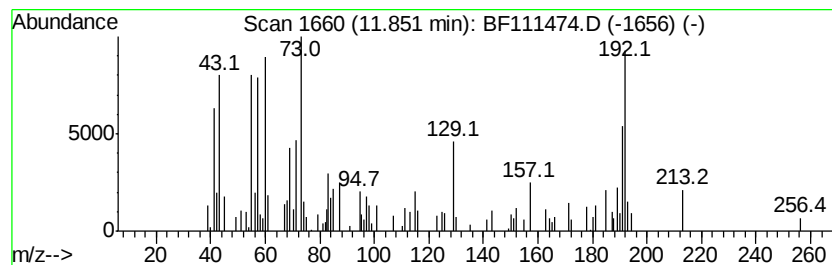
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.85	3.86 ng	218584	Phenanthrene-d10	11.31

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
2		Anthracene, 1-methyl-	192	C15H12	000610-48-0	80
3		Anthracene, 2-methyl-	192	C15H12	000613-12-7	50
4		Tridecanoic acid	214	C13H26O2	000638-53-9	47
5		Phenanthrene, 3-methyl-	192	C15H12	000832-71-3	42



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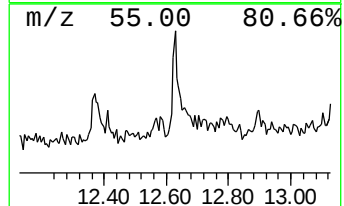
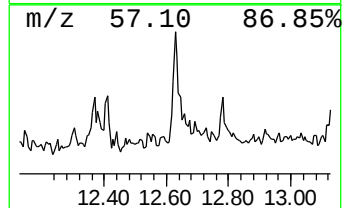
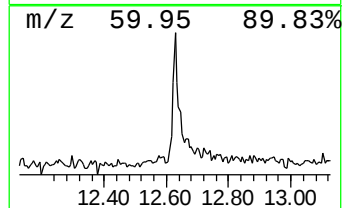
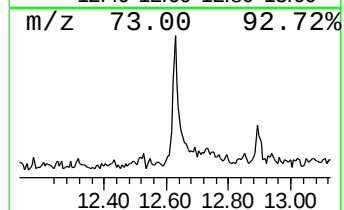
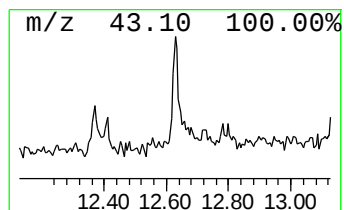
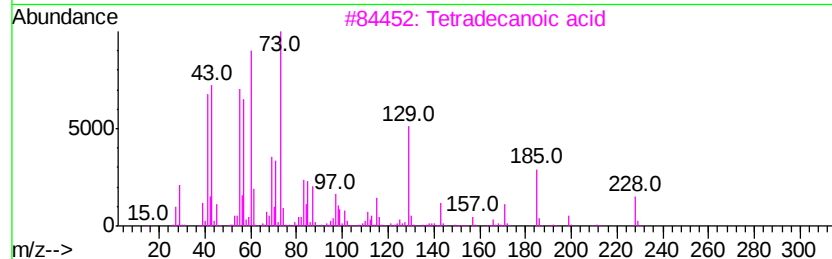
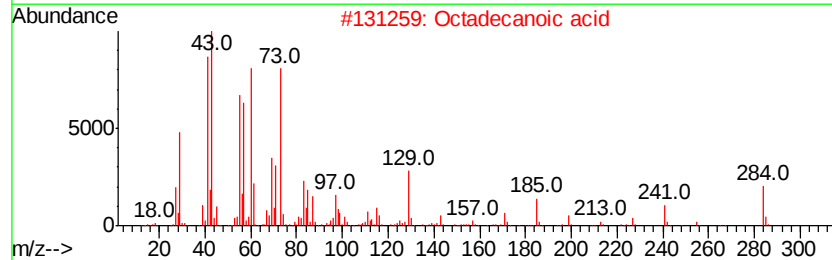
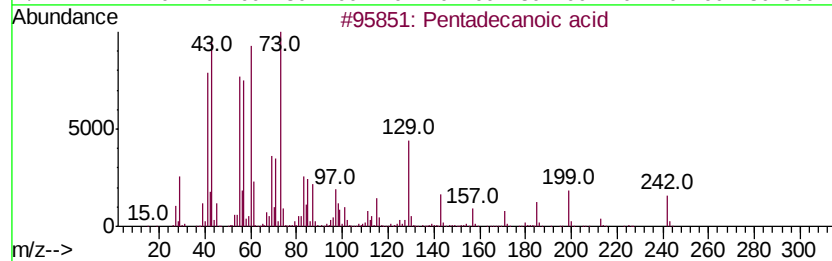
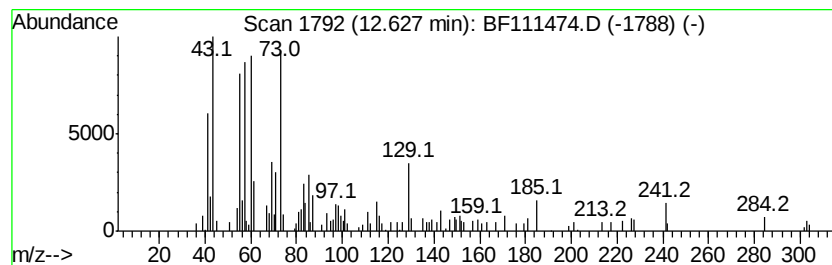
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Pentadecanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.63	2.75 ng	155946	Phenanthrene-d10	11.31

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentadecanoic acid	242	C15H30O2	001002-84-2	94
2		Octadecanoic acid	284	C18H36O2	000057-11-4	93
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	83
4		Tridecanoic acid	214	C13H26O2	000638-53-9	80
5		Dodecanoic acid	200	C12H24O2	000143-07-7	68



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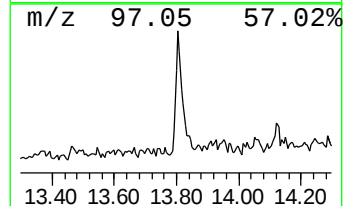
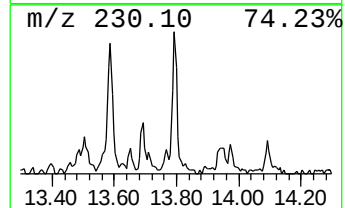
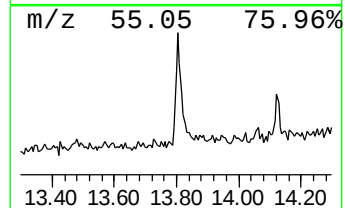
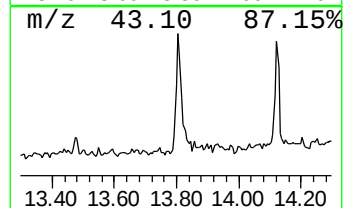
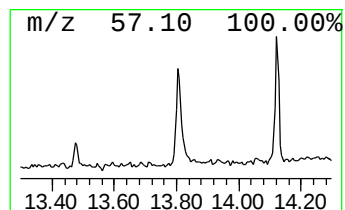
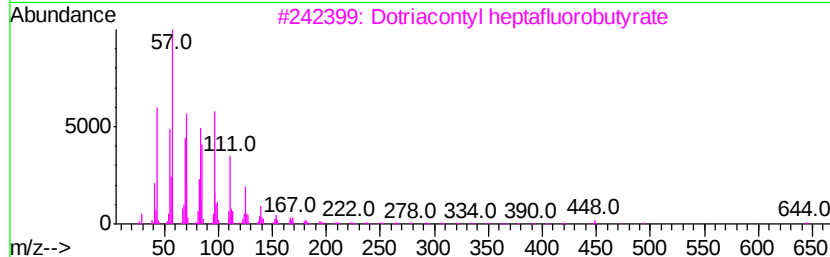
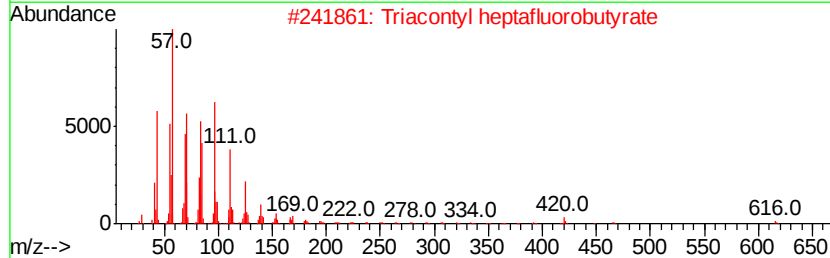
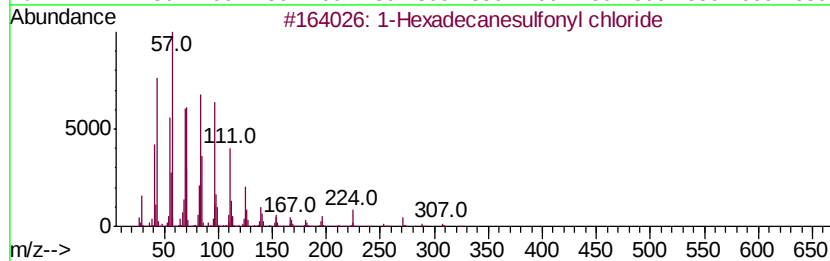
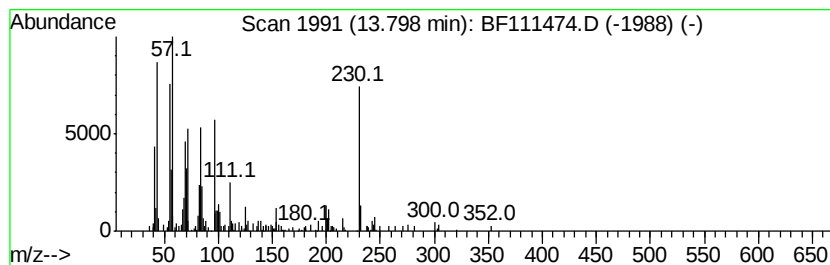
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 1-Hexadecanesulfonyl chloride Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.		
13.80	3.47 ng	255969	Chrysene-d12		13.95		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Hexadecanesulfonyl chloride	324	C16H33ClO2S	038775-38-1	91
2			Triacontyl heptafluorobutyrate	634	C34H61F7O2	1000351-83-2	91
3			Dotriacontyl heptafluorobutyrate	662	C36H65F7O2	1000351-84-2	91
4			Dotriacontyl pentafluoropropionate	612	C35H65F5O2	1000351-81-4	91
5			Tricosyl trifluoroacetate	436	C25H47F3O2	1000351-75-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121518\
Data File : BF111474.D
Acq On : 15 Dec 2018 17:48
Operator : JU/SJ
Sample : J6361-01
Misc :
ALS Vial : 18 Sample Multiplier: 1

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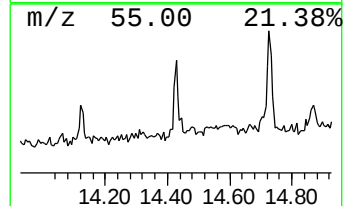
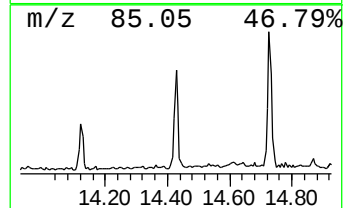
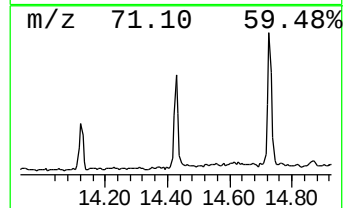
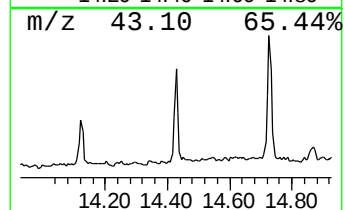
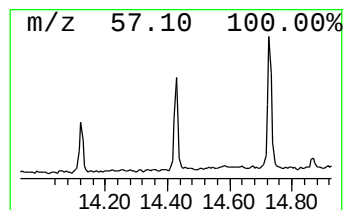
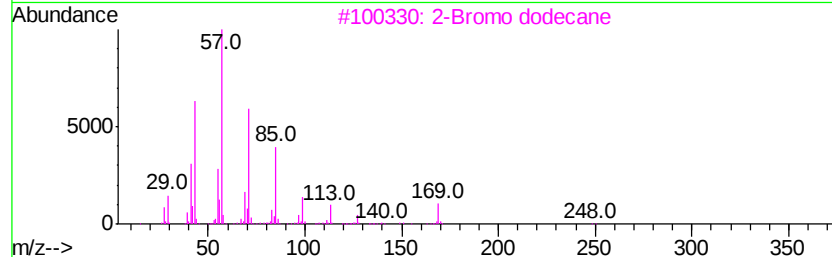
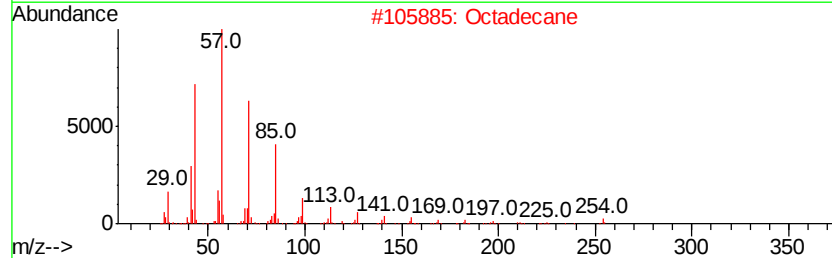
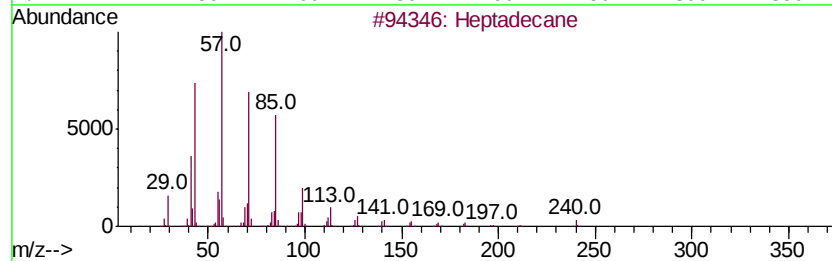
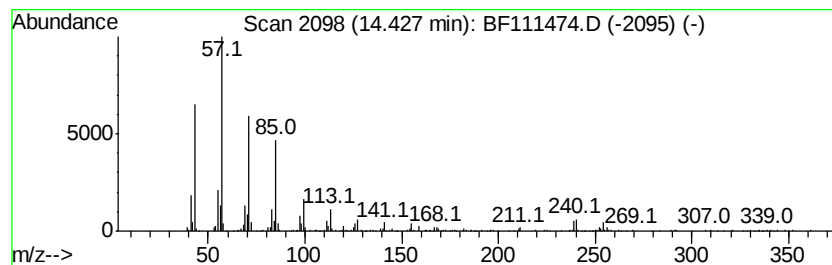
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Heptadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.43	2.88 ng	212261	Chrysene-d12	13.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane	240	C17H36	000629-78-7	96
2		Octadecane	254	C18H38	000593-45-3	93
3		2-Bromo dodecane	248	C12H25Br	013187-99-0	91
4		Hexacosane	366	C26H54	000630-01-3	87
5		Heptacosane	380	C27H56	000593-49-7	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121518\
Data File : BF111474.D
Acq On : 15 Dec 2018 17:48
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Misc :
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ClientSampleId :
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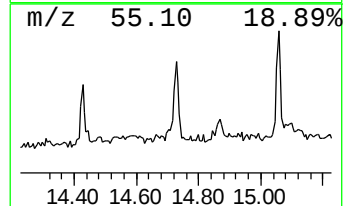
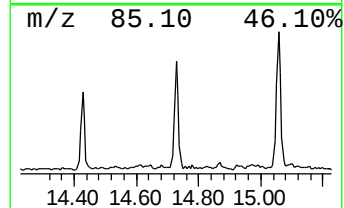
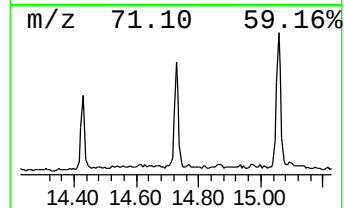
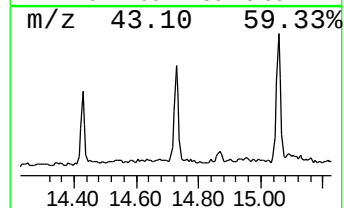
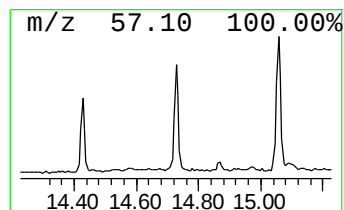
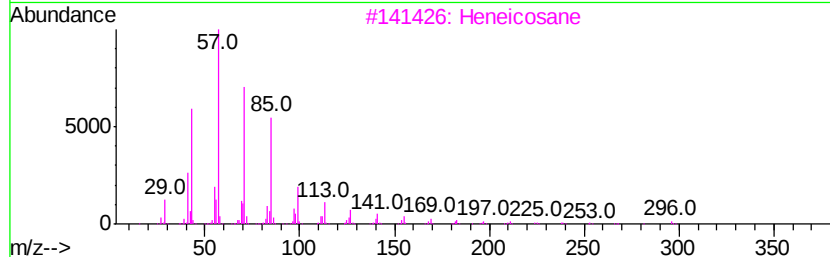
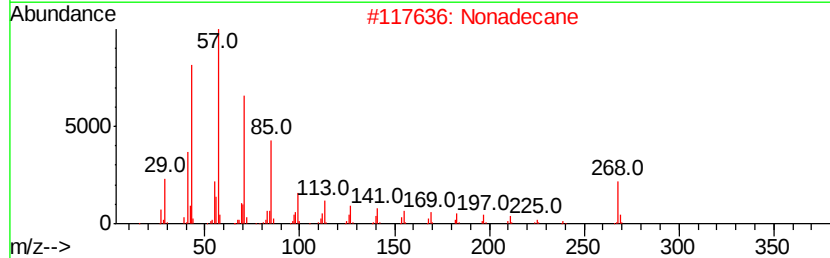
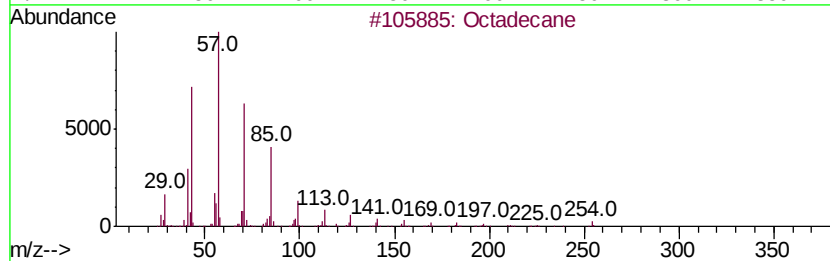
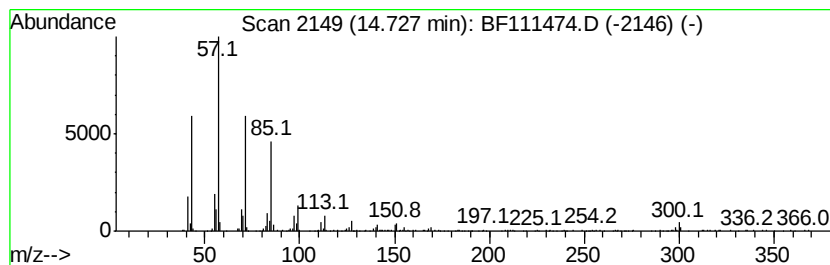
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Octadecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.73	7.30 ng	315104	Perylene-d12	15.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	91
2		Nonadecane	268	C19H40	000629-92-5	87
3		Heneicosane	296	C21H44	000629-94-7	83
4		Decane, 1-iodo-	268	C10H21I	002050-77-3	83
5		Tetracosane	338	C24H50	000646-31-1	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121518\
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ClientSampleId :
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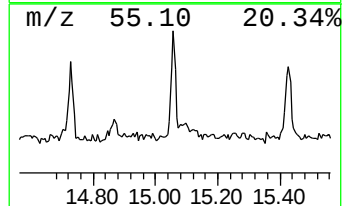
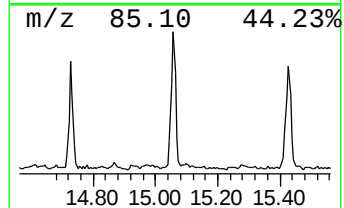
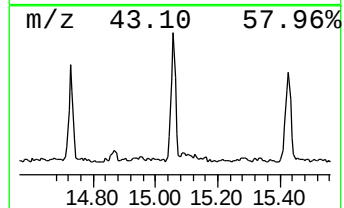
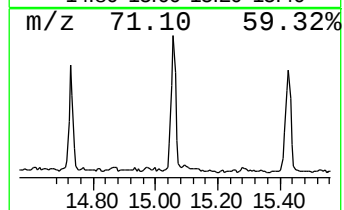
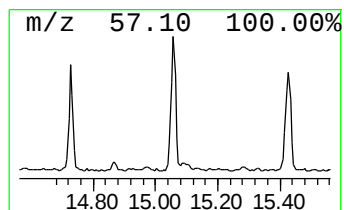
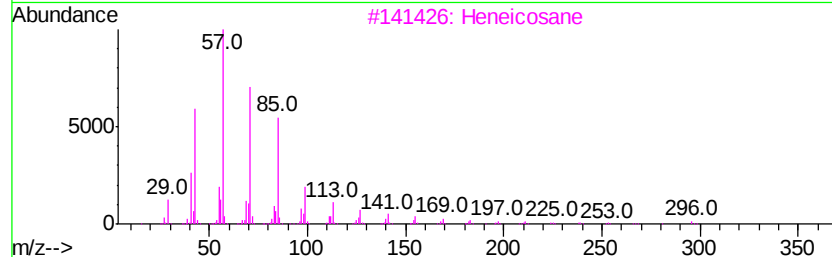
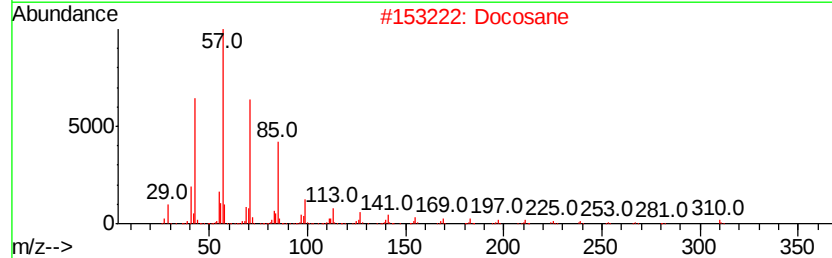
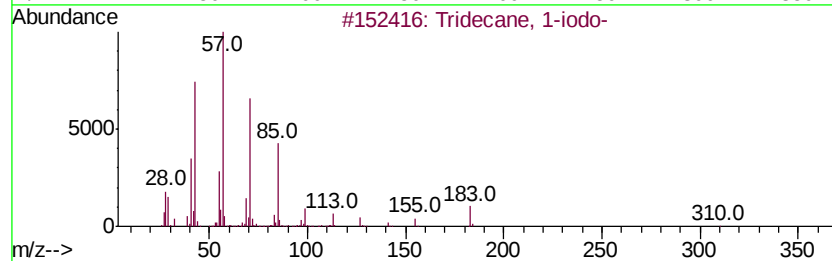
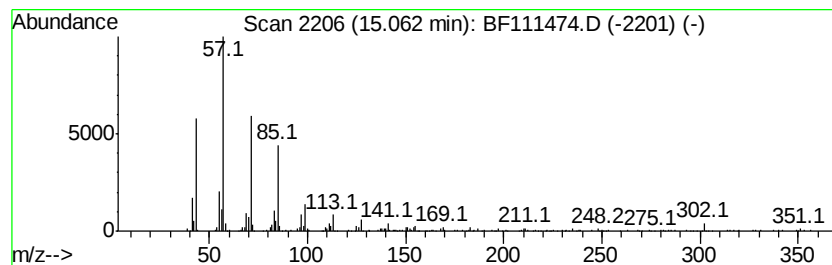
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Tridecane, 1-iodo- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.06	11.39 ng	491508	Perylene-d12	15.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane, 1-iodo-	310	C13H27I	035599-77-0	93
2		Docosane	310	C22H46	000629-97-0	90
3		Heneicosane	296	C21H44	000629-94-7	90
4		Pentacosane	352	C25H52	000629-99-2	87
5		Nonadecane	268	C19H40	000629-92-5	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121518\
Data File : BF111474.D
Acq On : 15 Dec 2018 17:48
Operator : JU/SJ
Sample : J6361-01
Misc :
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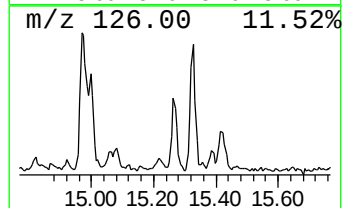
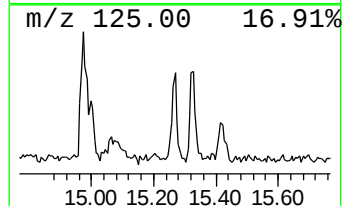
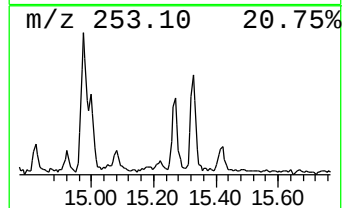
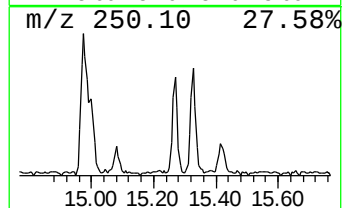
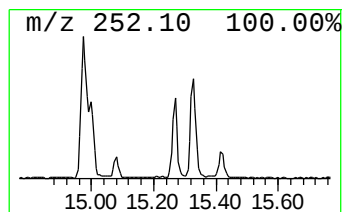
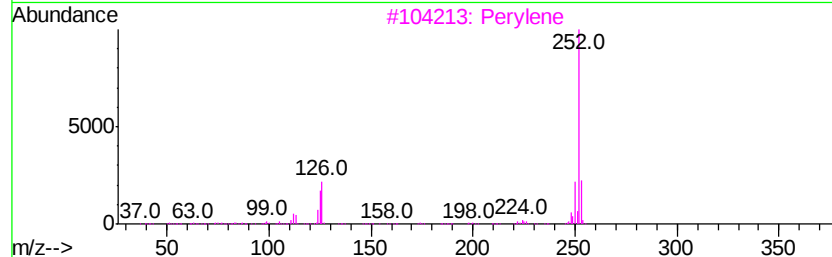
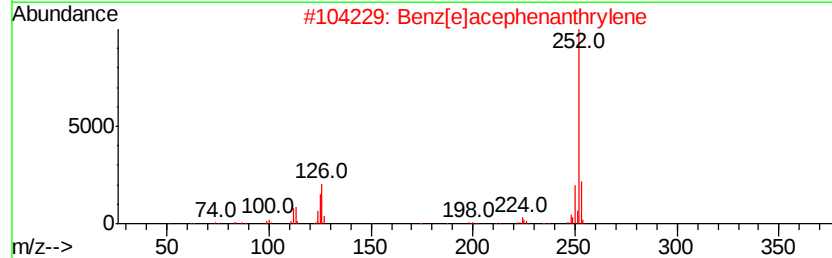
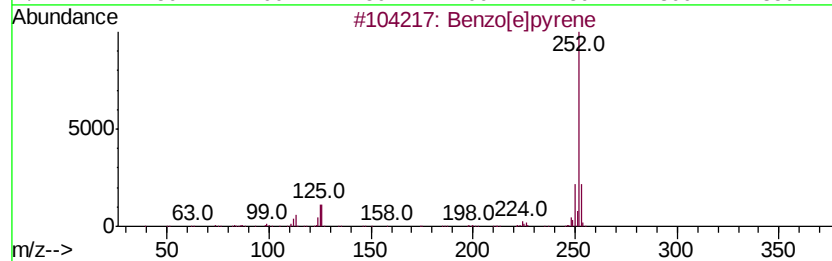
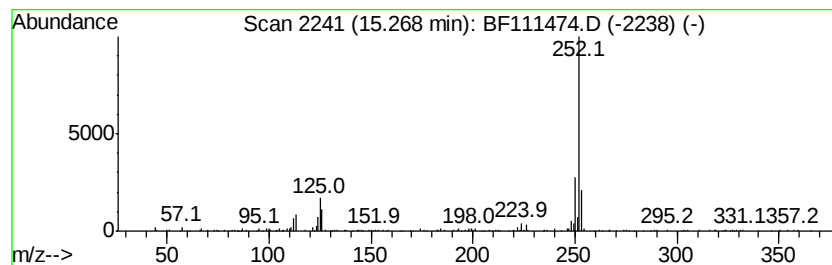
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Benzo[e]pyrene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.27	3.77 ng	162504	Perylene-d12	15.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[e]pyrene	252	C20H12	000192-97-2	96
2		Benz[e]acephenanthrylene	252	C20H12	000205-99-2	93
3		Perylene	252	C20H12	000198-55-0	93
4		Benzo[k]fluoranthene	252	C20H12	000207-08-9	89
5		Benzo[a]pyrene	252	C20H12	000050-32-8	81



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121518\
Data File : BF111474.D
Acq On : 15 Dec 2018 17:48
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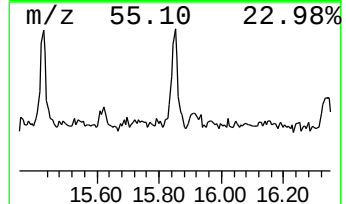
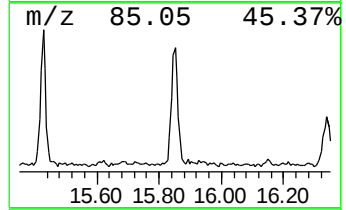
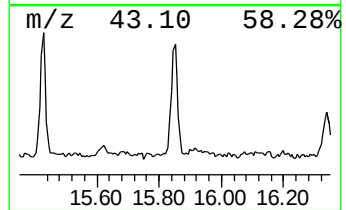
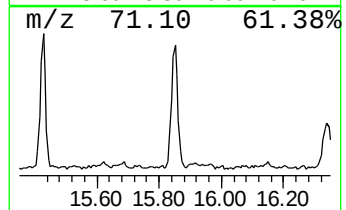
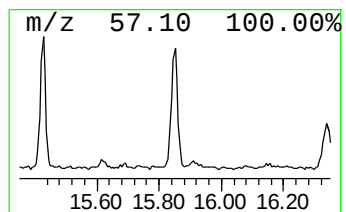
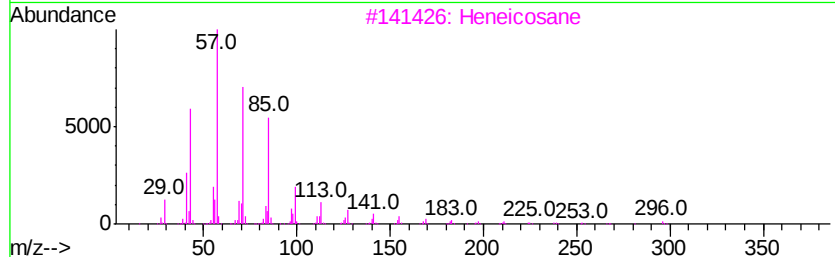
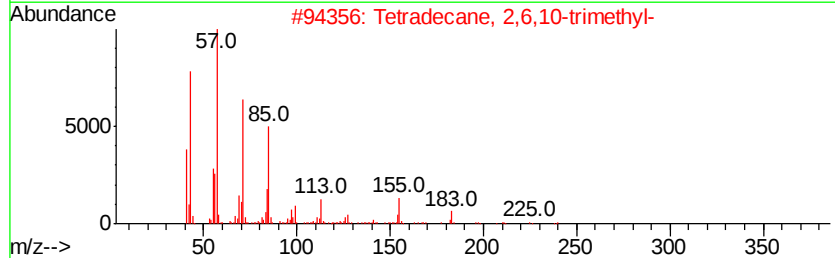
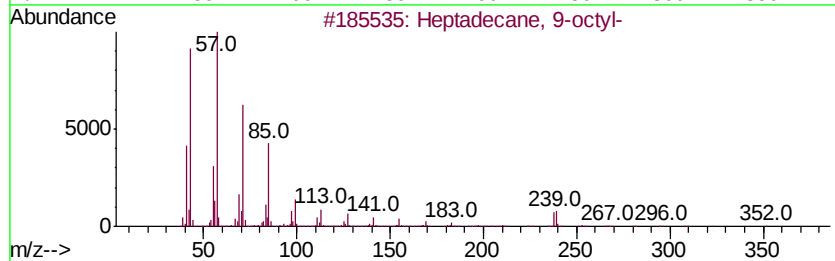
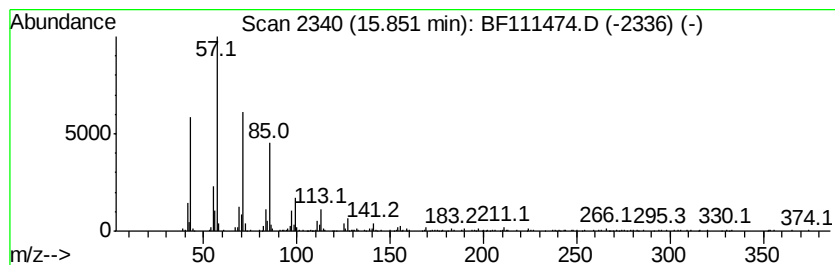
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Heptadecane, 9-octyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.85	10.05 ng	433444	Perylene-d12	15.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	95
2		Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	93
3		Heneicosane	296	C21H44	000629-94-7	91
4		Eicosane	282	C20H42	000112-95-8	91
5		Hexacosane	366	C26H54	000630-01-3	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF121518\
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Sample : J6361-01
Misc :
ALS Vial : 18 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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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unknown6.57	6.57	74.5	ng	3571990	1	6.79	959192	20.0
n-Hexadecanoic acid	11.85	3.9	ng	218584	4	11.31	1132700	20.0
Pentadecanoic acid	12.63	2.8	ng	155946	4	11.31	1132700	20.0
1-Hexadecanesulfo...	13.80	3.5	ng	255969	5	13.95	1475800	20.0
Heptadecane	14.43	2.9	ng	212261	5	13.95	1475800	20.0
Octadecane	14.73	7.3	ng	315104	6	15.39	863002	20.0
Tridecane, 1-iodo-	15.06	11.4	ng	491508	6	15.39	863002	20.0
Benzo[e]pyrene	15.27	3.8	ng	162504	6	15.39	863002	20.0
Heptadecane, 9-oc...	15.85	10.1	ng	433444	6	15.39	863002	20.0