

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040319\
 Data File : BF113552.D
 Acq On : 3 Apr 2019 19:18
 Operator : JU/SJ
 Sample : K2213-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S9A(2-10)COMP

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-BF040319.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	5.298	542	546	570	rBV	2756953	2799902	83.01%	8.265%
2	6.334	718	722	725	rBV	2800469	2839856	84.20%	8.383%
3	6.457	739	743	753	rVB	3203281	2944796	87.31%	8.693%
4	6.687	778	782	792	rBV	823592	777333	23.05%	2.295%
5	6.840	804	808	822	rBV	2541900	2301893	68.25%	6.795%
6	7.251	873	878	881	rBV	1930848	1788998	53.04%	5.281%
7	7.969	996	1000	1015	rBV	1005627	993573	29.46%	2.933%
8	9.045	1178	1183	1186	rBV	3755939	3372794	100.00%	9.956%
9	9.439	1247	1250	1255	rBV	202822	168650	5.00%	0.498%
10	9.722	1294	1298	1302	rBV	1267648	1068426	31.68%	3.154%
11	10.510	1428	1432	1442	rBV	2189708	2034093	60.31%	6.004%
12	11.204	1546	1550	1553	rBV	1231510	1092879	32.40%	3.226%
13	11.228	1553	1554	1557	rVV	107644	57839	1.71%	0.171%
14	11.280	1557	1563	1567	rVB	116990	114944	3.41%	0.339%
15	11.439	1583	1590	1594	rBV	50693	61926	1.84%	0.183%
16	11.745	1638	1642	1651	rBV3	224472	370685	10.99%	1.094%
17	11.816	1651	1654	1663	rVB	54316	92408	2.74%	0.273%
18	12.298	1732	1736	1740	rBV3	51434	70264	2.08%	0.207%
19	12.345	1741	1744	1750	rVB	49452	52609	1.56%	0.155%
20	12.416	1752	1756	1759	rBV	545512	506994	15.03%	1.497%
21	12.527	1770	1775	1779	rBV	412969	520106	15.42%	1.535%
22	12.639	1789	1794	1797	rVB	496715	490001	14.53%	1.446%
23	12.669	1797	1799	1801	rVV	52732	40399	1.20%	0.119%
24	12.692	1801	1803	1808	rVB3	43128	42807	1.27%	0.126%
25	12.792	1813	1820	1822	rBV	3026065	3147566	93.32%	9.291%
26	12.810	1822	1823	1828	rVB2	122917	109367	3.24%	0.323%
27	12.974	1844	1851	1854	rVB2	73529	119080	3.53%	0.352%
28	13.039	1854	1862	1865	rBV3	76471	176985	5.25%	0.522%
29	13.369	1916	1918	1921	rVB	65931	45512	1.35%	0.134%
30	13.592	1953	1956	1958	rBV	53886	61166	1.81%	0.181%
31	13.698	1971	1974	1975	rBV2	144719	150480	4.46%	0.444%
32	13.839	1993	1998	2001	rBV2	1105822	1224396	36.30%	3.614%
33	13.863	2001	2002	2010	rVB	314913	251535	7.46%	0.743%
34	14.010	2024	2027	2030	rBV	183644	169839	5.04%	0.501%

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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	14.316	2076	2079	2082	rVB	240358	228198	6.77%	0.674%
36	14.610	2125	2129	2134	rBV	440727	456625	13.54%	1.348%
37	14.851	2167	2170	2177	rVB	217682	376163	11.15%	1.110%
38	14.921	2178	2182	2186	rVB	545315	577188	17.11%	1.704%
39	15.133	2216	2218	2225	rVB	138345	173237	5.14%	0.511%
40	15.192	2226	2228	2233	rBV	82381	102536	3.04%	0.303%
41	15.251	2234	2238	2240	rBV	666907	886728	26.29%	2.618%
42	15.668	2305	2309	2314	rVB	434146	644531	19.11%	1.903%
43	16.127	2383	2387	2393	rBV	230203	370893	11.00%	1.095%

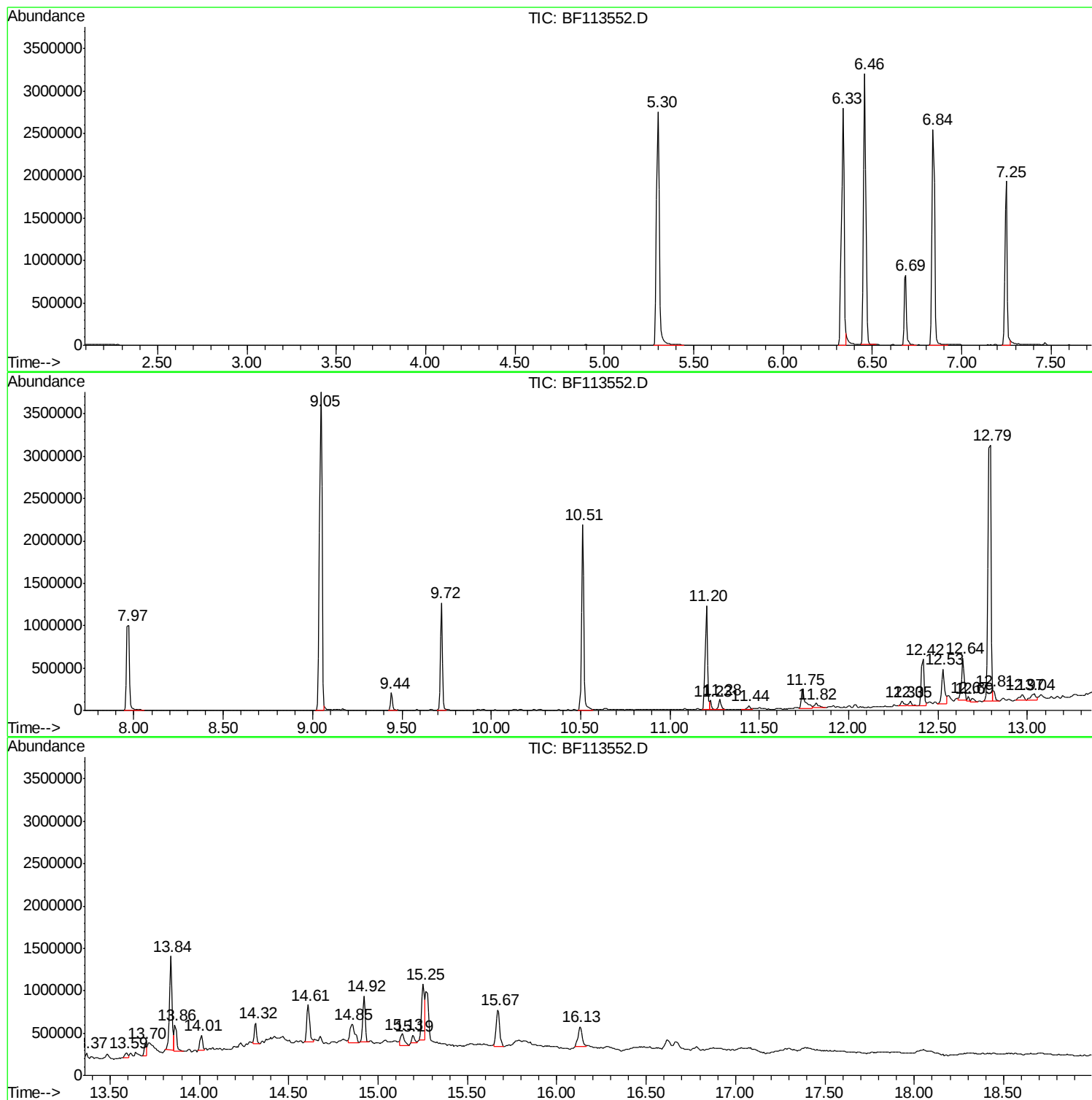
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF040319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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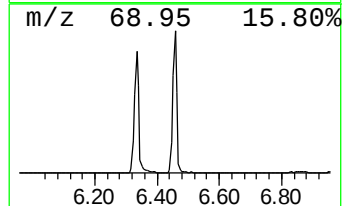
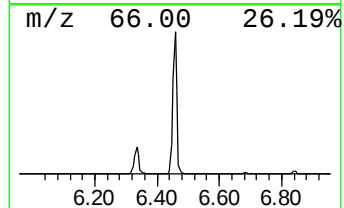
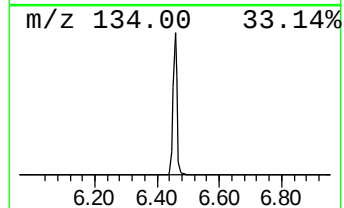
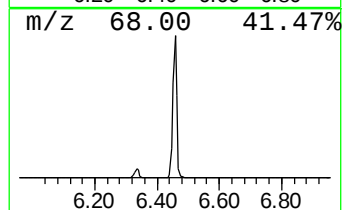
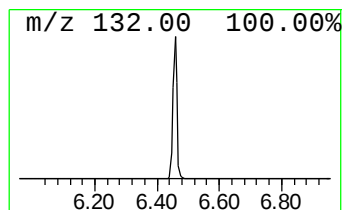
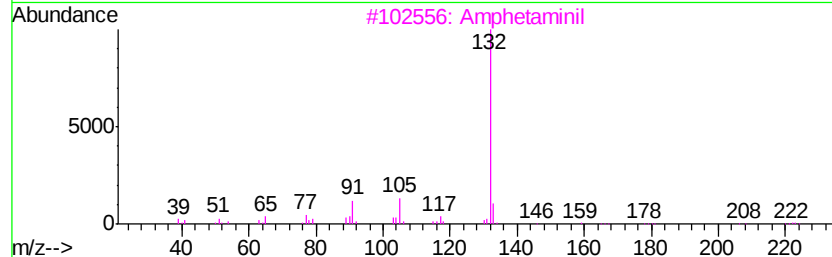
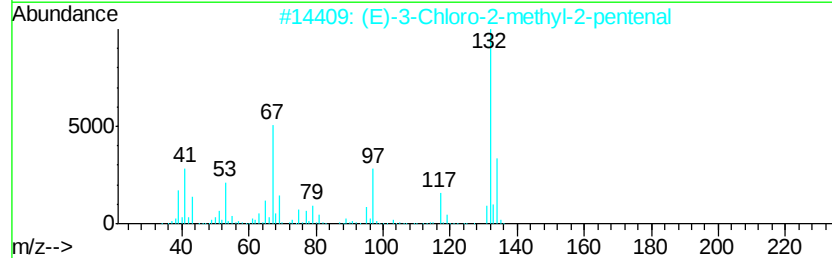
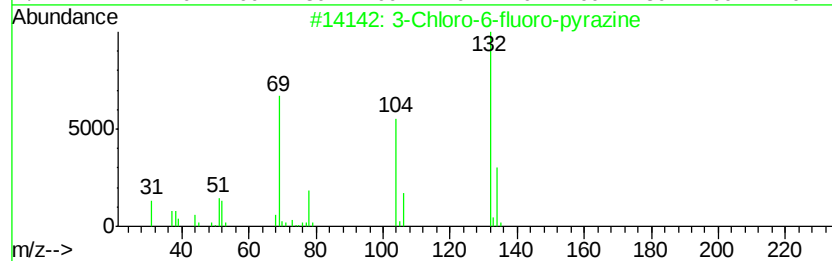
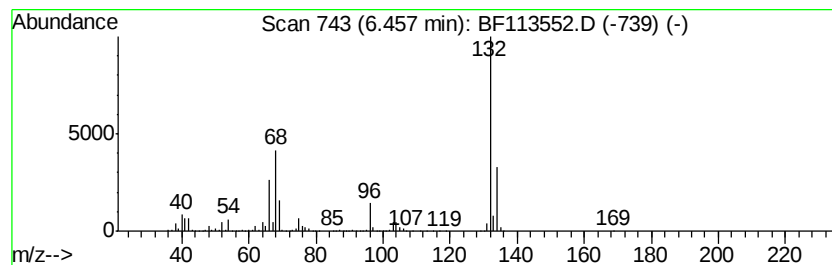
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown6.46 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.46	75.77 ng	2944800	1,4-Dichlorobenzene-d4	6.69

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	12
3		Amphetaminil	250	C17H18N2	017590-01-1	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		Xenon	132	Xe	007440-63-3	9



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ClientSampleId :
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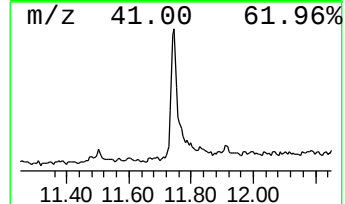
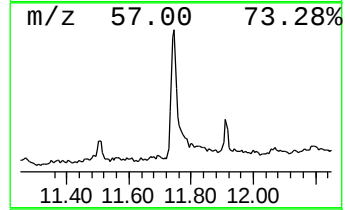
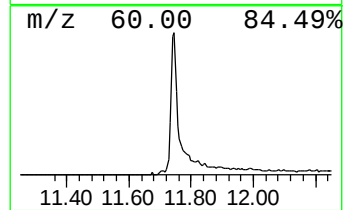
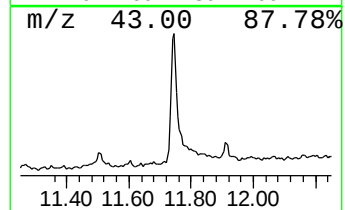
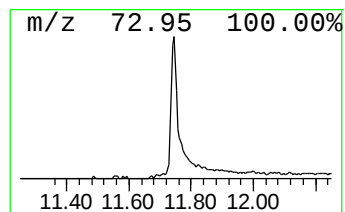
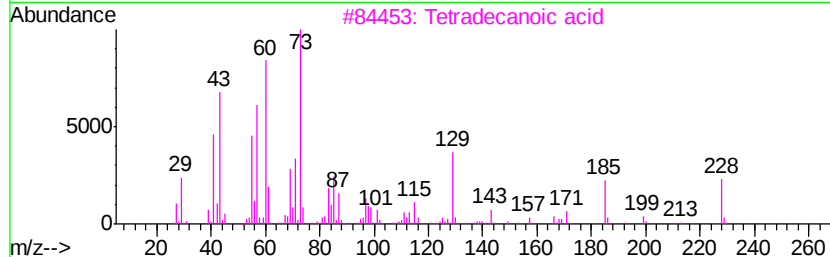
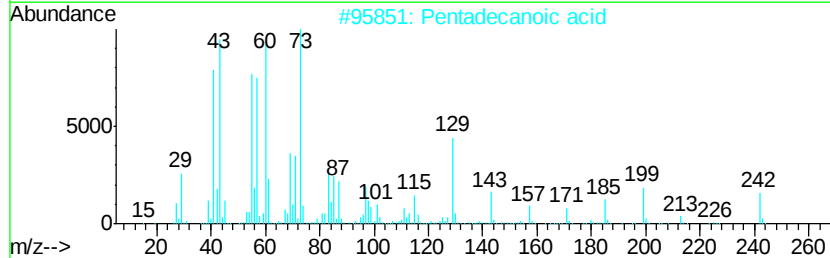
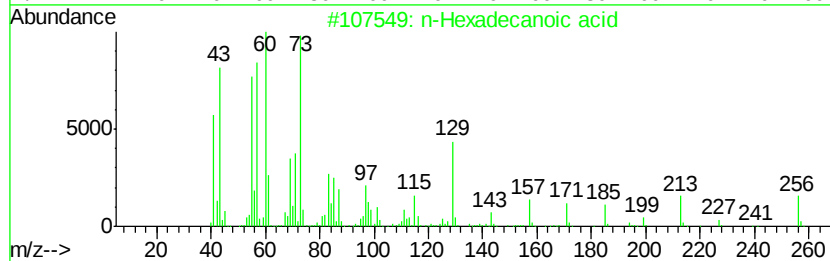
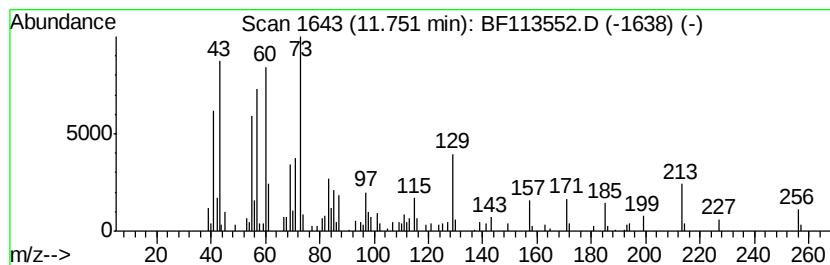
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.75	6.78 ng	370685	Phenanthrene-d10	11.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	90
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	90
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	80
4		Tridecanoic acid	214	C13H26O2	000638-53-9	76
5		n-Decanoic acid	172	C10H20O2	000334-48-5	76



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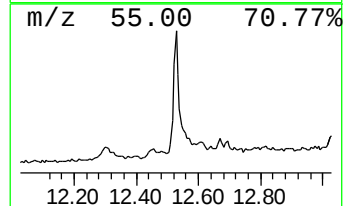
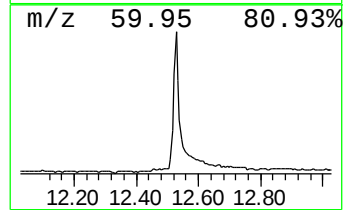
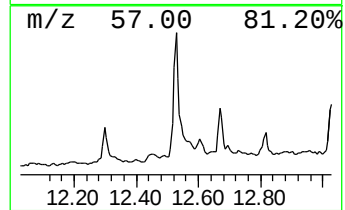
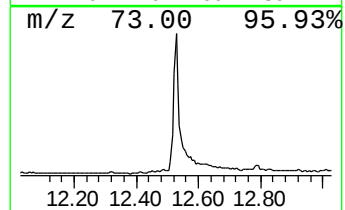
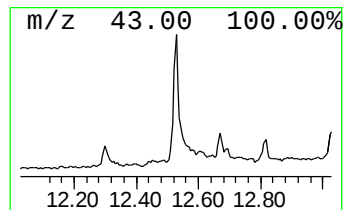
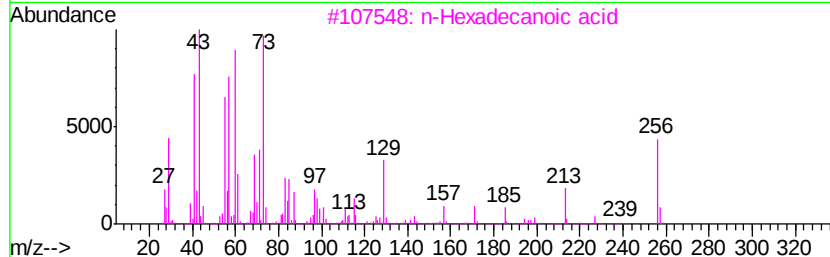
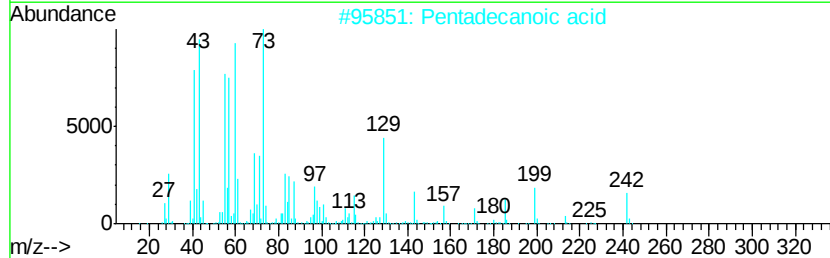
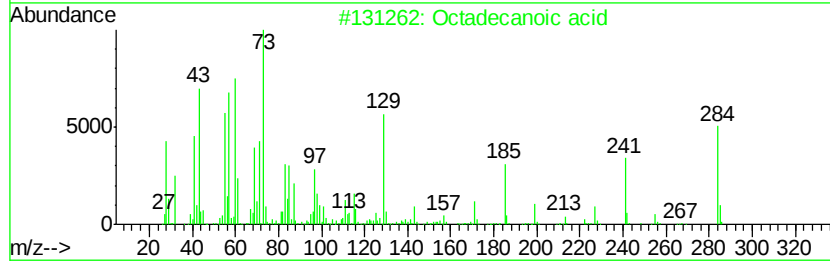
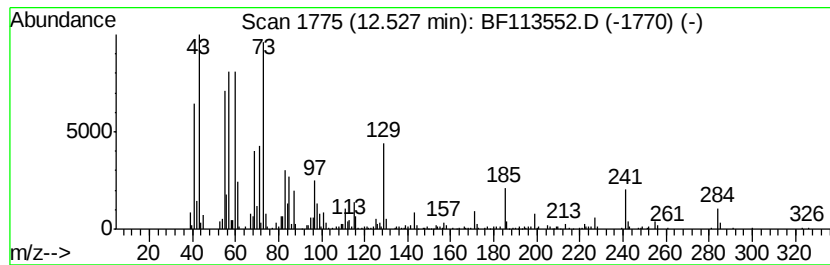
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Octadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.53	8.50 ng	520106	Chrysene-d12	13.84

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	93
3		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	87
5		n-Decanoic acid	172	C10H20O2	000334-48-5	62



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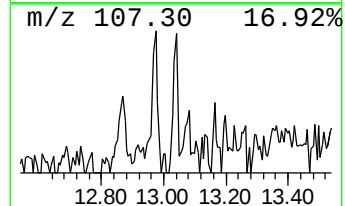
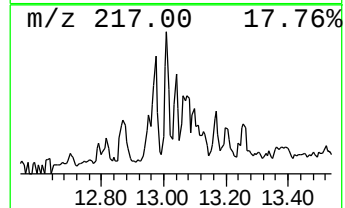
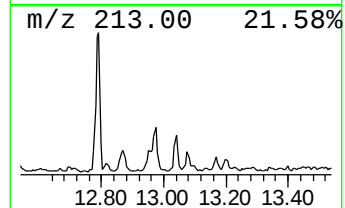
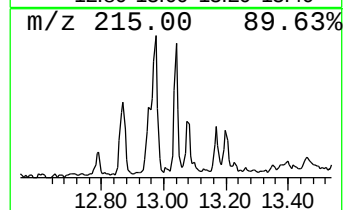
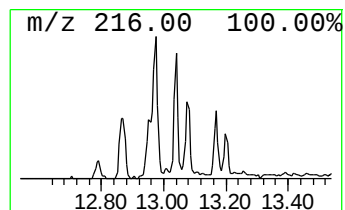
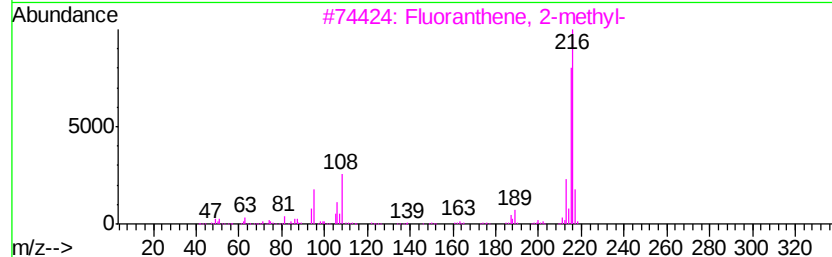
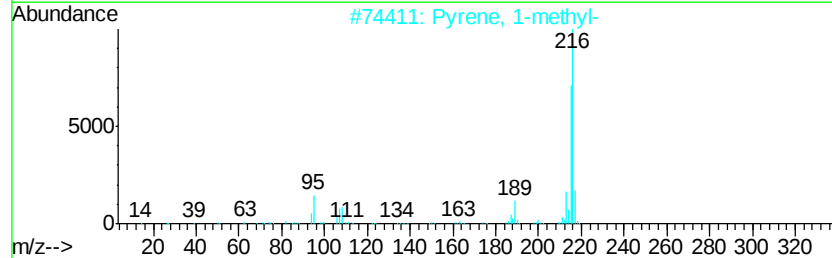
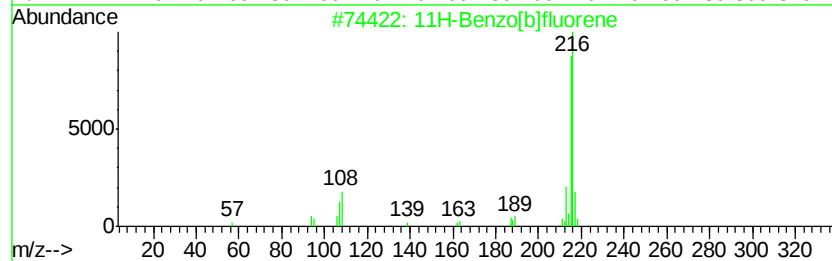
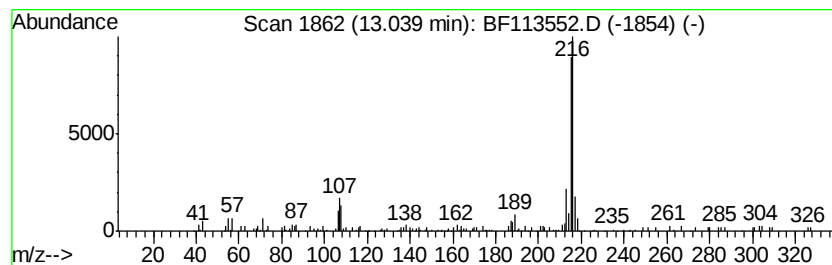
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 11H-Benzo[b]fluorene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.04	2.89 ng	176985	Chrysene-d12	13.84

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	93
2		Pyrene, 1-methyl-	216	C17H12	002381-21-7	91
3		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	91
4		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	81
5		Pyrene, 2-methyl-	216	C17H12	003442-78-2	70



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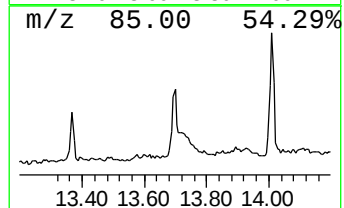
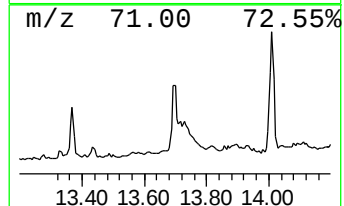
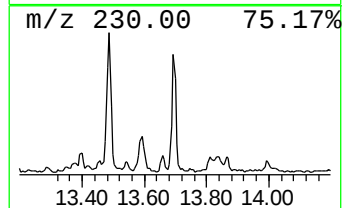
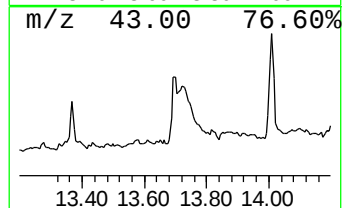
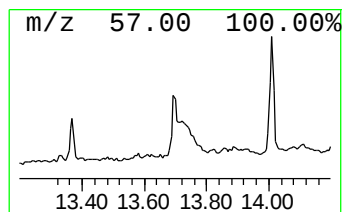
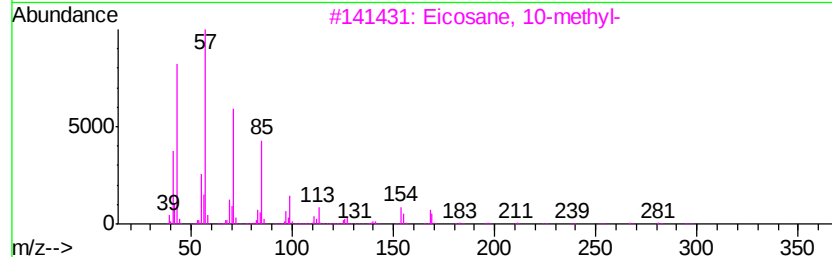
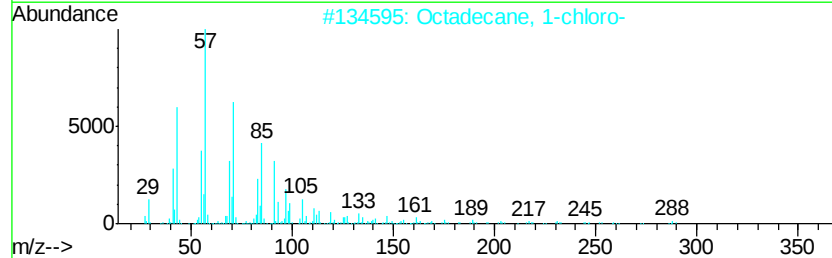
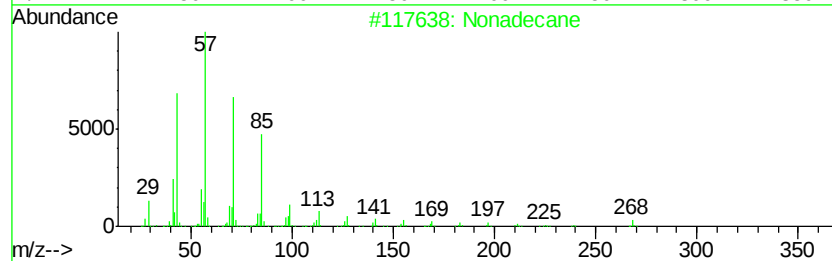
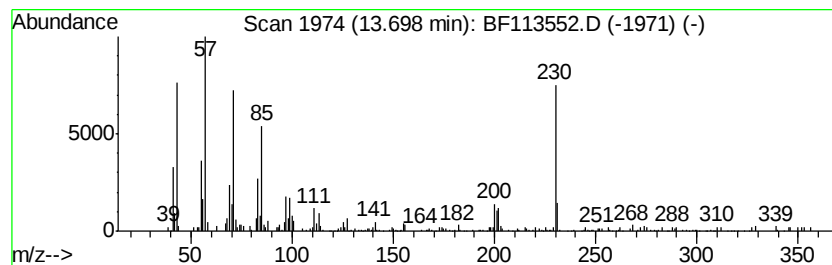
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TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 6 Nonadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.70	2.46 ng	150480	Chrysene-d12	13.84

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane		268	C19H40	000629-92-5	86
2	Octadecane, 1-chloro-		288	C18H37Cl	003386-33-2	80
3	Eicosane, 10-methyl-		296	C21H44	054833-23-7	60
4	Docosane		310	C22H46	000629-97-0	59
5	Hexadecane		226	C16H34	000544-76-3	52



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040319\
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Acq On : 3 Apr 2019 19:18
Operator : JU/SJ
Sample : K2213-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
S9A(2-10)COMP

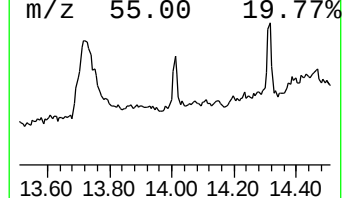
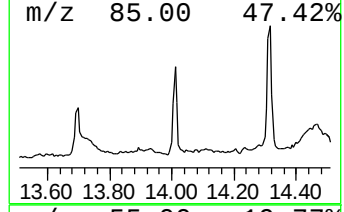
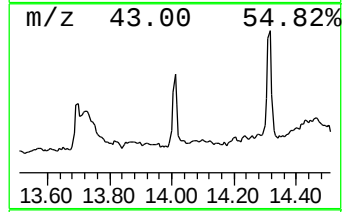
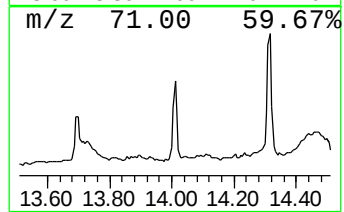
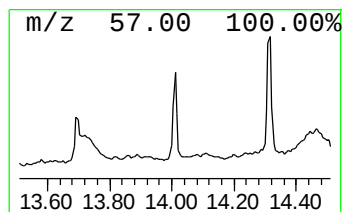
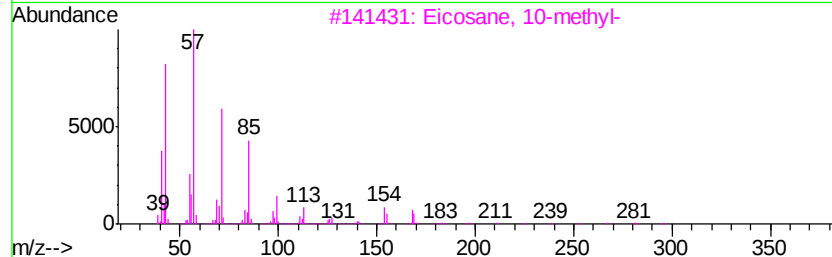
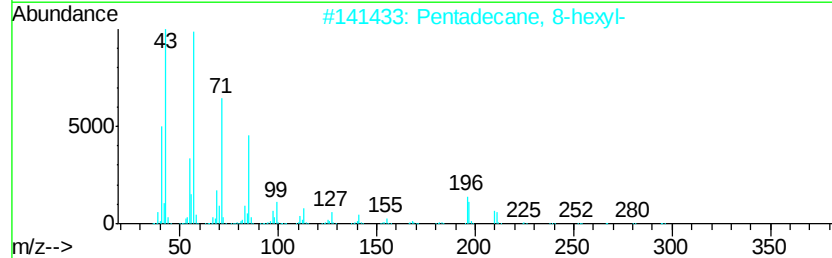
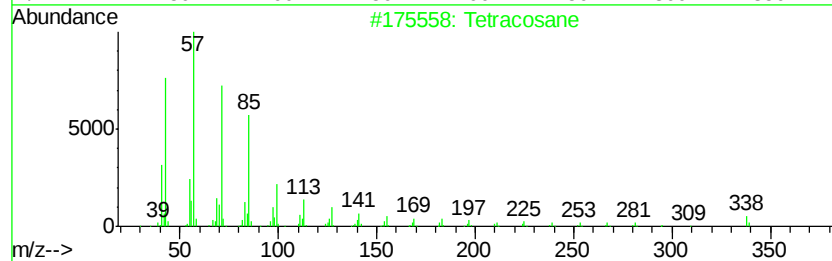
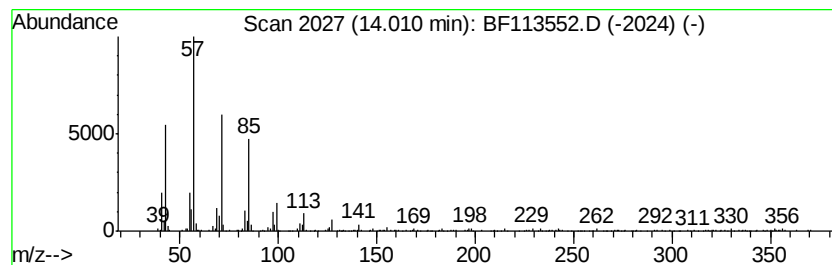
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Tetracosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.01	2.77 ng	169839	Chrysene-d12	13.84

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetracosane	338	C24H50	000646-31-1	95
2		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	93
3		Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
4		Octacosane	394	C28H58	000630-02-4	86
5		Tricosane, 2-methyl-	338	C24H50	001928-30-9	81



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040319\
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Operator : JU/SJ
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Instrument :
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ClientSampleId :
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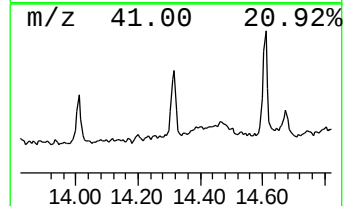
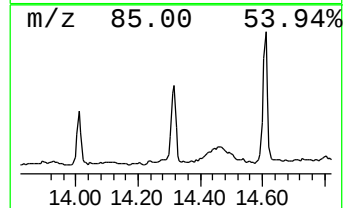
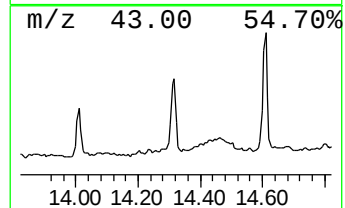
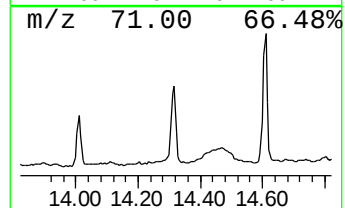
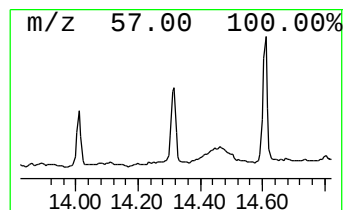
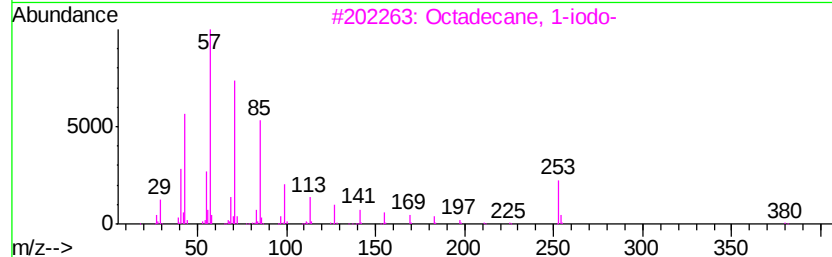
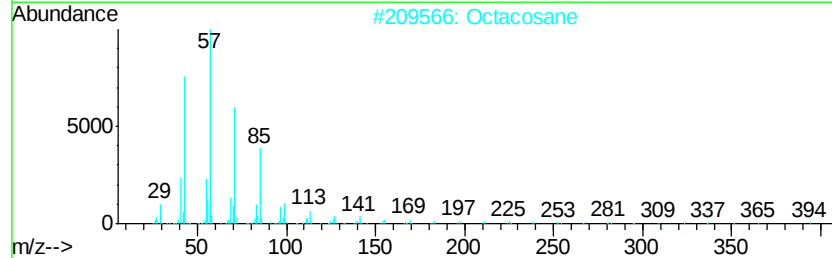
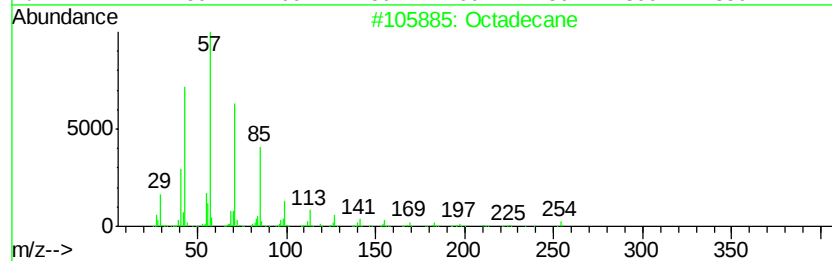
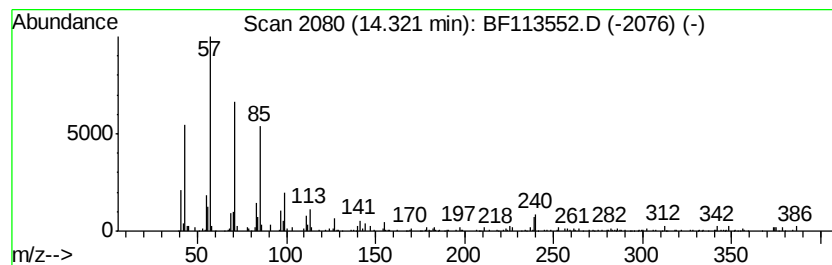
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Octadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.32	3.73 ng	228198	Chrysene-d12	13.84

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane		254	C18H38	000593-45-3	92
2	Octacosane		394	C28H58	000630-02-4	90
3	Octadecane, 1-iodo-		380	C18H37I	000629-93-6	87
4	Docosane		310	C22H46	000629-97-0	87
5	Eicosane		282	C20H42	000112-95-8	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040319\
Data File : BF113552.D
Acq On : 3 Apr 2019 19:18
Operator : JU/SJ
Sample : K2213-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
S9A(2-10)COMP

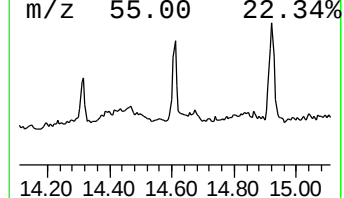
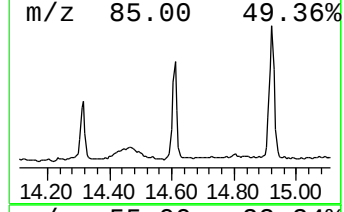
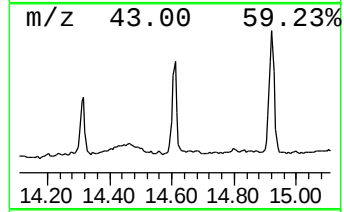
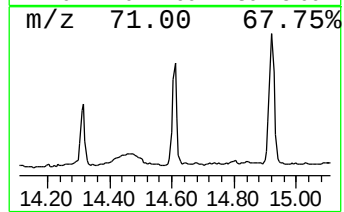
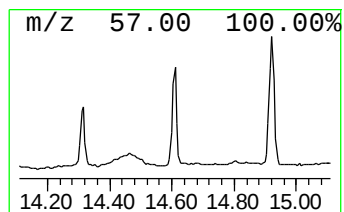
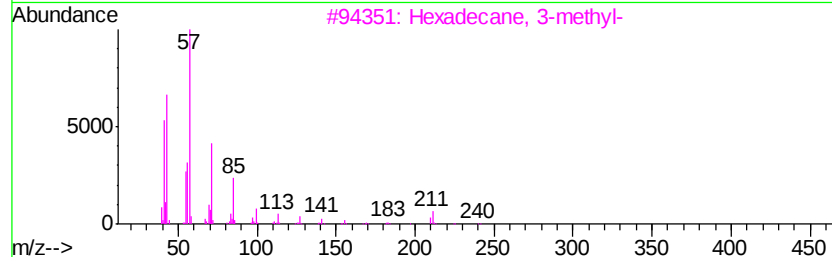
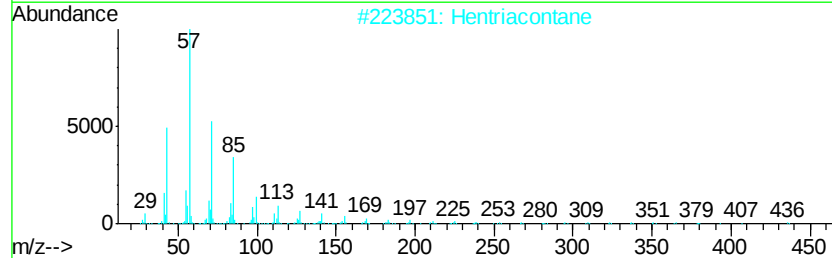
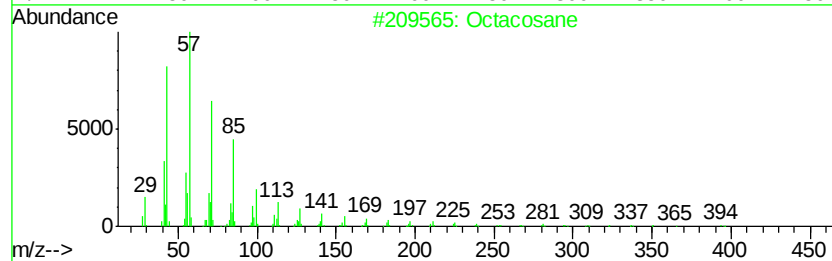
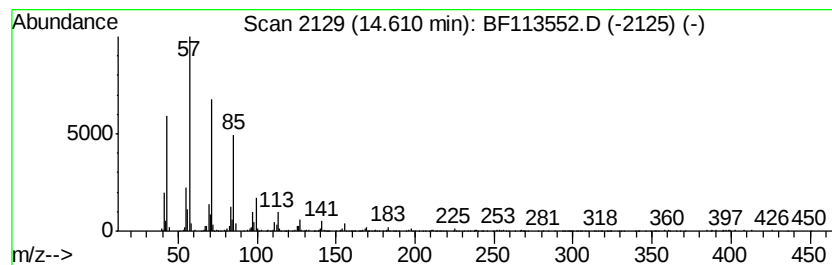
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Octacosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.61	10.30 ng	456625	Perylene-d12	15.25

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	96
2		Hentriacontane	437	C31H64	000630-04-6	95
3		Hexadecane, 3-methyl-	240	C17H36	006418-43-5	93
4		Eicosane, 9-octyl-	394	C28H58	013475-77-9	91
5		Octadecane	254	C18H38	000593-45-3	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040319\
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Acq On : 3 Apr 2019 19:18
Operator : JU/SJ
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ClientSampleId :
S9A(2-10)COMP

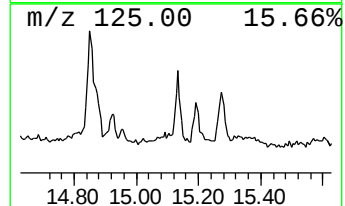
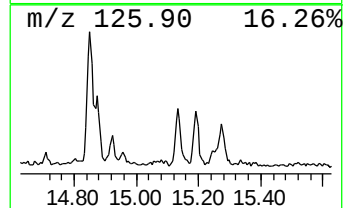
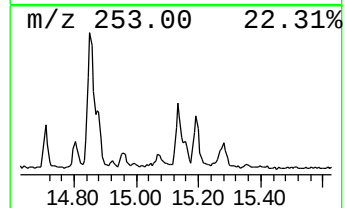
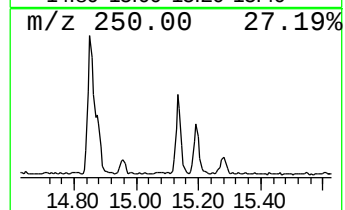
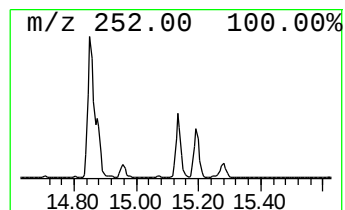
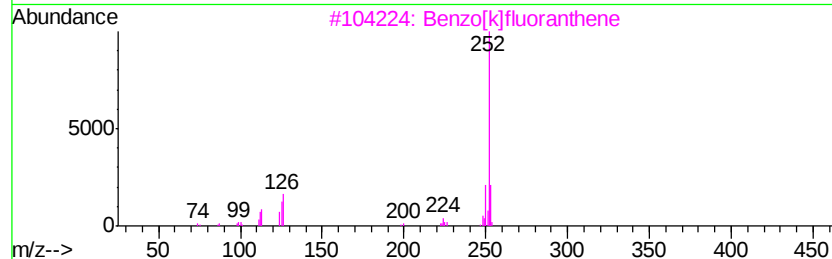
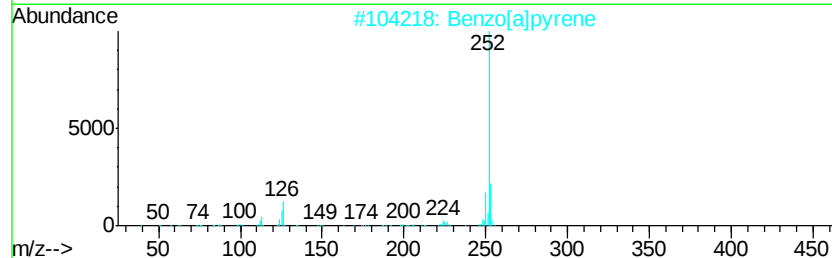
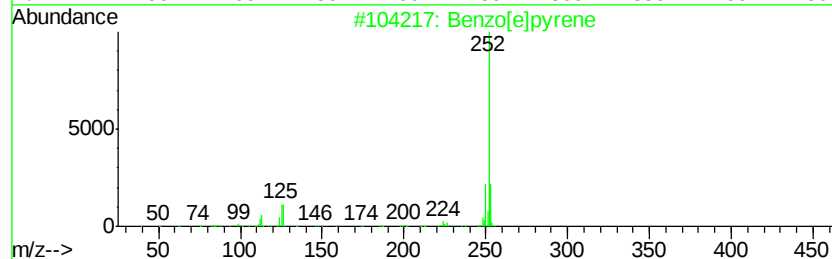
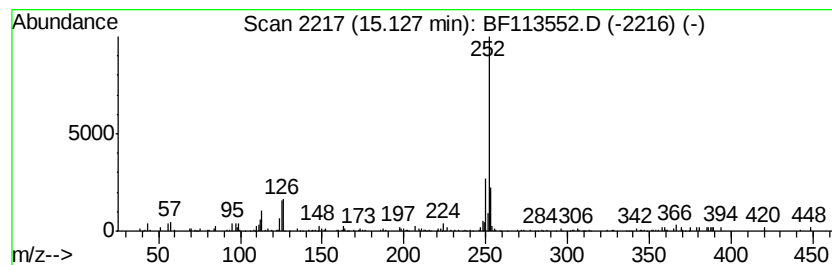
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
15.13	3.91 ng	173237	Perylene-d12	15.25

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	97
2			Benzo[a]pyrene	252	C20H12	000050-32-8	96
3			Benzo[k]fluoranthene	252	C20H12	000207-08-9	93
4			Benz[e]acephenanthrylene	252	C20H12	000205-99-2	92
5			Perylene	252	C20H12	000198-55-0	89



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040319\
Data File : BF113552.D
Acq On : 3 Apr 2019 19:18
Operator : JU/SJ
Sample : K2213-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
S9A(2-10)COMP

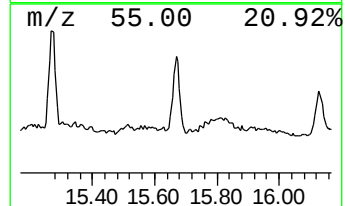
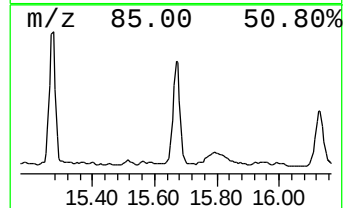
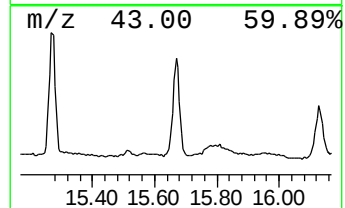
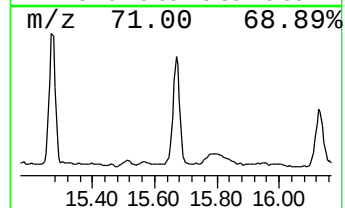
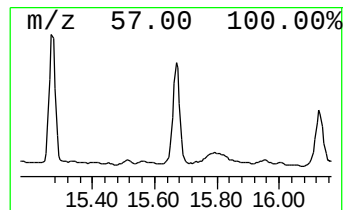
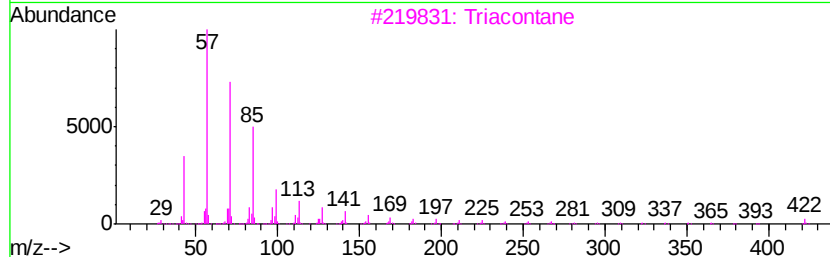
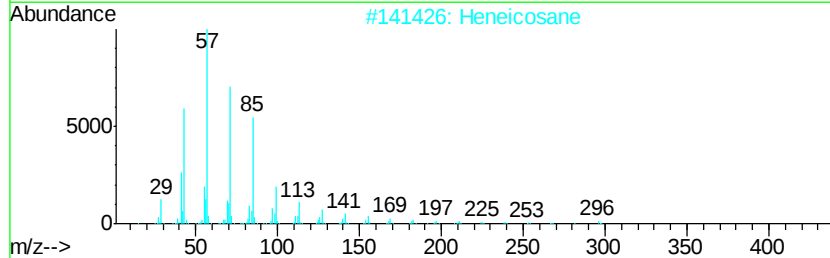
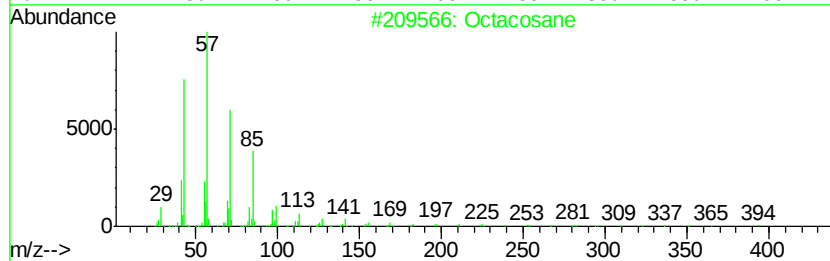
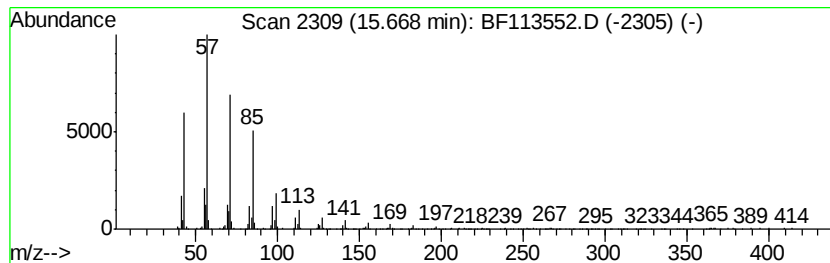
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Triacontane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.67	14.54 ng	644531	Perylene-d12	15.25

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	98
2		Heneicosane	296	C21H44	000629-94-7	91
3		Triacontane	422	C30H62	000638-68-6	91
4		Heptacosane	380	C27H56	000593-49-7	91
5		Hexacosane	366	C26H54	000630-01-3	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040319\
Data File : BF113552.D
Acq On : 3 Apr 2019 19:18
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ClientSampleId :
S9A(2-10)COMP

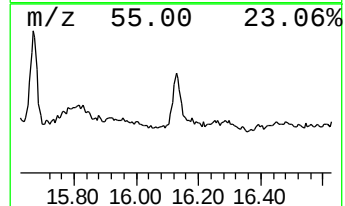
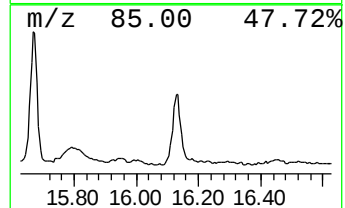
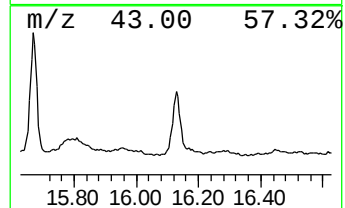
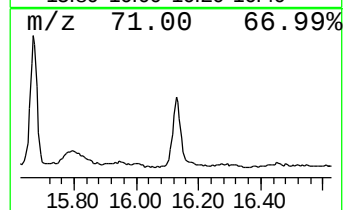
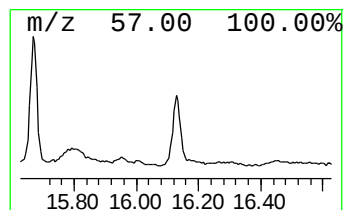
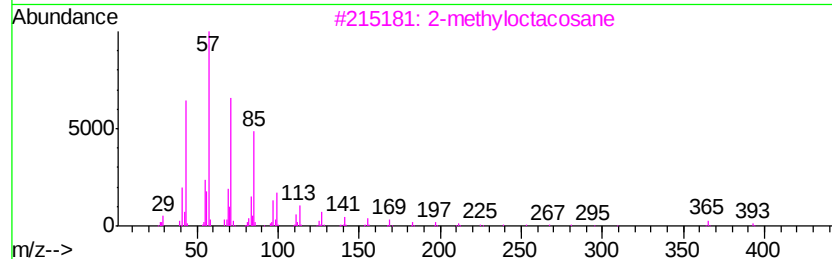
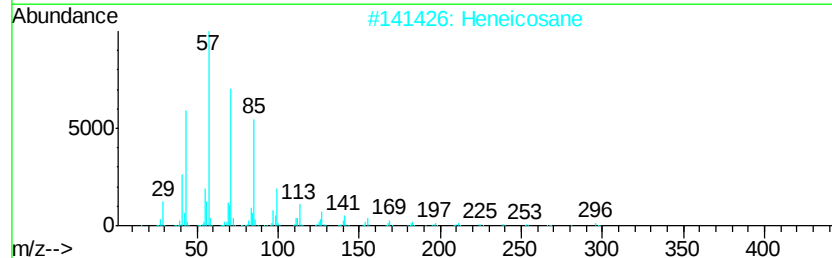
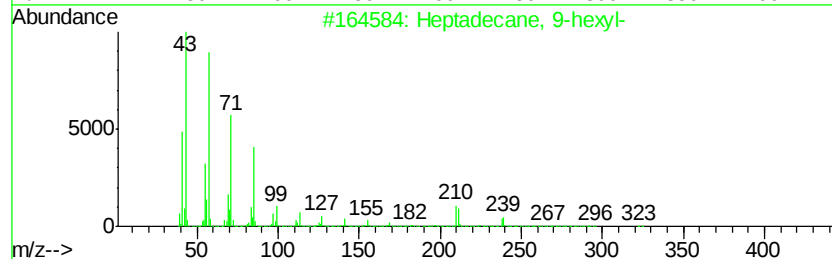
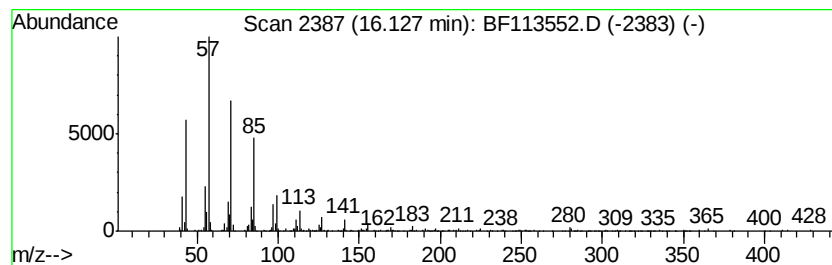
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Heptadecane, 9-hexyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.13	8.37 ng	370893	Perylene-d12	15.25

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	90
2		Heneicosane	296	C21H44	000629-94-7	90
3		2-methyloctacosane	408	C29H60	1000376-72-8	87
4		Octacosane	394	C28H58	000630-02-4	87
5		1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040319\
Data File : BF113552.D
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Sample : K2213-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
S9A(2-10)COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF040319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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n-Hexadecanoic acid	11.75	6.8	ng	370685	4	11.20	1092880	20.0
Octadecanoic acid	12.53	8.5	ng	520106	5	13.84	1224400	20.0
11H-Benzo[b]fluorene	13.04	2.9	ng	176985	5	13.84	1224400	20.0
Nonadecane	13.70	2.5	ng	150480	5	13.84	1224400	20.0
Tetracosane	14.01	2.8	ng	169839	5	13.84	1224400	20.0
Octadecane	14.32	3.7	ng	228198	5	13.84	1224400	20.0
Octacosane	14.61	10.3	ng	456625	6	15.25	886728	20.0
Benzo[e]pyrene	15.13	3.9	ng	173237	6	15.25	886728	20.0
Triacontane	15.67	14.5	ng	644531	6	15.25	886728	20.0
Heptadecane, 9-he...	16.13	8.4	ng	370893	6	15.25	886728	20.0