

Data Path : Z:\HPCHEM1\BNA F\DATA\BF080517\
 Data File : BF097395.D
 Acq On : 5 Aug 2017 14:17
 Operator : SJ/JU
 Sample : I4584-01 5X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OK-01-080217

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-BF072517.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	4.551	466	470	475	rBV2	11095	18219	2.15%	0.259%
2	5.045	551	554	569	rBV	43343	96796	11.40%	1.378%
3	6.116	733	736	747	rBV	46921	102640	12.08%	1.461%
4	6.216	750	753	760	rVB	79971	114401	13.47%	1.628%
5	6.422	785	788	799	rBV	664229	590909	69.57%	8.411%
6	6.581	812	815	825	rBV	75611	79941	9.41%	1.138%
7	7.004	885	887	897	rBV	50074	81632	9.61%	1.162%
8	7.704	1003	1006	1017	rBV	862744	838856	98.76%	11.941%
9	8.786	1186	1190	1197	rBV	126476	118037	13.90%	1.680%
10	8.886	1202	1207	1211	rBV	12857	13942	1.64%	0.198%
11	9.192	1256	1259	1263	rBV2	12656	15264	1.80%	0.217%
12	9.410	1292	1296	1299	rBV	20565	18718	2.20%	0.266%
13	9.451	1299	1303	1312	rVB	914229	849359	100.00%	12.090%
14	9.880	1372	1376	1378	rBV	54127	45391	5.34%	0.646%
15	9.904	1378	1380	1383	rVB	24756	18188	2.14%	0.259%
16	10.121	1414	1417	1420	rBV	6909	9235	1.09%	0.131%
17	10.239	1434	1437	1442	rBV	56009	53089	6.25%	0.756%
18	10.374	1456	1460	1467	rVB	29226	35931	4.23%	0.511%
19	10.816	1532	1535	1538	rBV2	31807	29381	3.46%	0.418%
20	10.921	1549	1553	1562	rVB	825275	759514	89.42%	10.811%
21	11.239	1603	1607	1613	rBV	38690	36826	4.34%	0.524%
22	11.339	1621	1624	1627	rVB2	18722	14747	1.74%	0.210%
23	11.404	1632	1635	1638	rBV4	5165	10251	1.21%	0.146%
24	11.539	1655	1658	1660	rBV2	7382	8856	1.04%	0.126%
25	11.645	1673	1676	1681	rBV	48887	46184	5.44%	0.657%
26	11.810	1697	1704	1707	rBV7	6957	15140	1.78%	0.216%
27	11.892	1715	1718	1721	rBV4	6961	8854	1.04%	0.126%
28	11.921	1721	1723	1728	rVB5	7272	12420	1.46%	0.177%
29	12.033	1736	1742	1745	rBV2	74603	116000	13.66%	1.651%
30	12.133	1754	1759	1761	rBV4	14202	14825	1.75%	0.211%
31	12.357	1793	1797	1802	rBV6	7448	12135	1.43%	0.173%
32	12.398	1802	1804	1807	rVV	74677	65485	7.71%	0.932%
33	12.498	1818	1821	1825	rBV	128225	113907	13.41%	1.621%
34	12.751	1859	1864	1868	rBV	87390	91212	10.74%	1.298%

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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

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072517.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	12.874	1883	1885	1888	rVB4	13875	12201	1.44%	0.174%
36	13.004	1904	1907	1911	rVB5	10504	10410	1.23%	0.148%
37	13.092	1919	1922	1925	rBV	99287	89338	10.52%	1.272%
38	13.210	1938	1942	1944	rBV2	12677	14893	1.75%	0.212%
39	13.286	1953	1955	1956	rBV2	11173	9012	1.06%	0.128%
40	13.421	1974	1978	1983	rVB	105851	111788	13.16%	1.591%
41	13.539	1994	1998	2007	rBV	565546	600814	70.74%	8.552%
42	13.651	2015	2017	2022	rVB6	12703	14459	1.70%	0.206%
43	13.733	2025	2031	2040	rVV	110083	144303	16.99%	2.054%
44	13.886	2055	2057	2060	rBV4	10682	9519	1.12%	0.135%
45	13.915	2060	2062	2067	rVB6	15630	19833	2.34%	0.282%
46	14.039	2079	2083	2088	rVB	110962	132064	15.55%	1.880%
47	14.133	2096	2099	2100	rBV3	23427	25141	2.96%	0.358%
48	14.327	2129	2132	2135	rBV	111417	98024	11.54%	1.395%
49	14.609	2177	2180	2184	rBV	97857	98486	11.60%	1.402%
50	14.880	2222	2226	2230	rBV	442305	524804	61.79%	7.470%
51	14.915	2230	2232	2237	rVB	120419	132487	15.60%	1.886%
52	15.262	2287	2291	2295	rBV2	99558	152919	18.00%	2.177%
53	15.421	2315	2318	2325	rVB8	38337	67764	7.98%	0.965%
54	15.545	2337	2339	2341	rBV2	17208	13775	1.62%	0.196%
55	15.656	2354	2358	2364	rVB	78623	113512	13.36%	1.616%
56	16.121	2431	2437	2440	rBV4	54479	104063	12.25%	1.481%
57	16.615	2519	2521	2523	rBV3	14081	11352	1.34%	0.162%
58	18.039	2760	2763	2770	rBV7	22865	42930	5.05%	0.611%
59	18.686	2871	2873	2875	rBV3	13889	14975	1.76%	0.213%

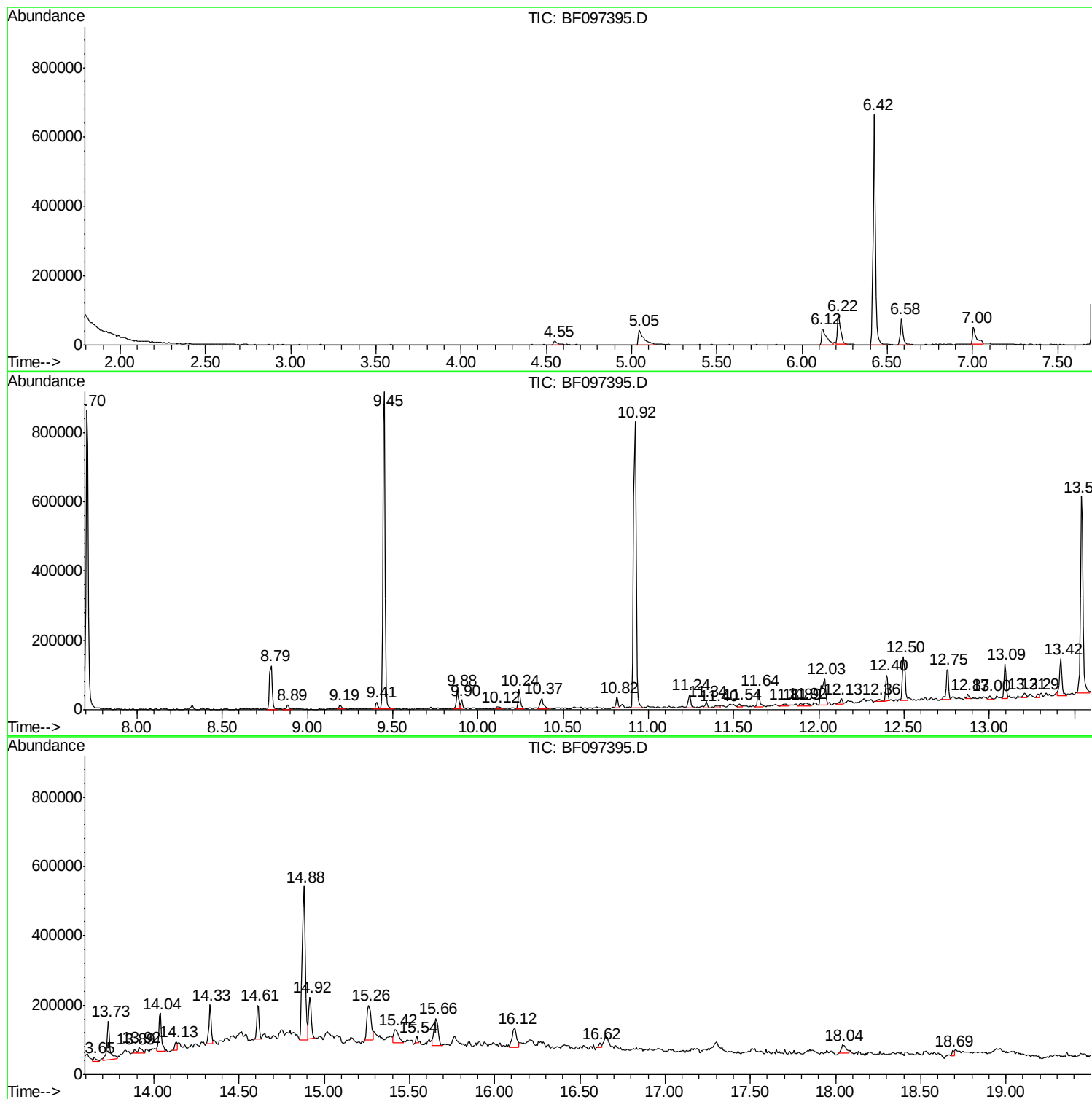
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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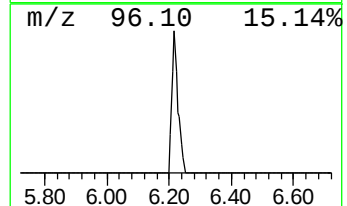
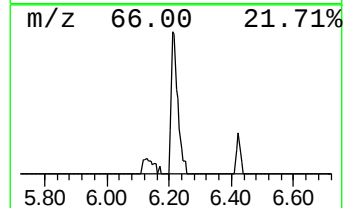
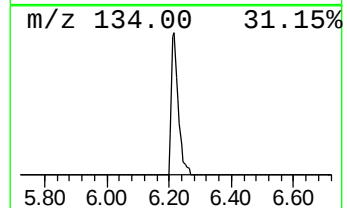
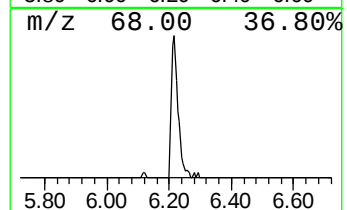
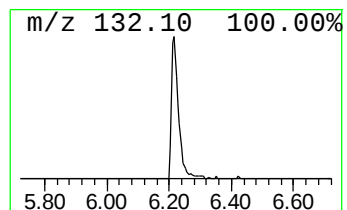
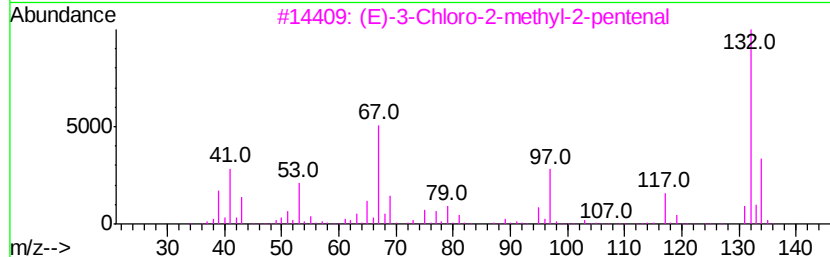
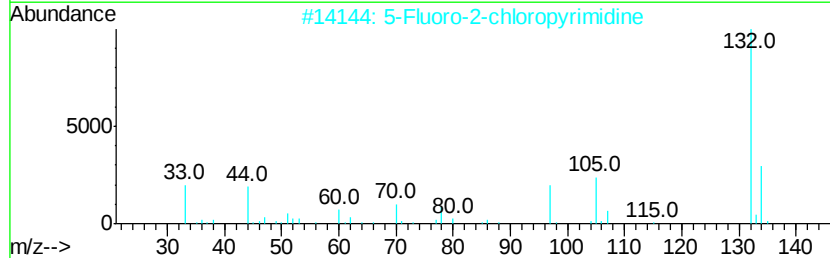
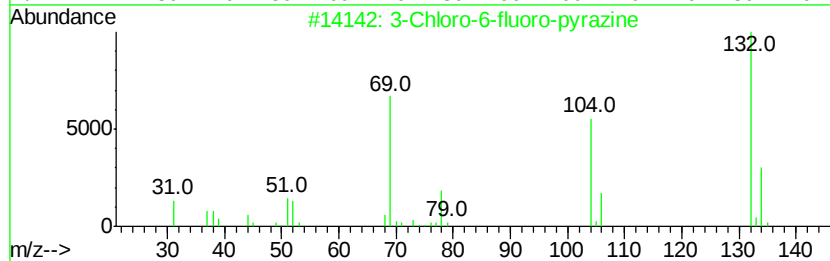
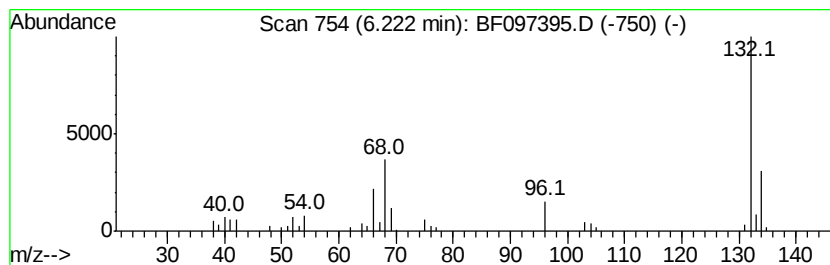
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown6.22 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.22	3.87 ng	114401	1,4-Dichlorobenzene-d4	6.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	30
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
4		5-Aminoindole	132	C8H8N2	005192-03-0	10
5		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10



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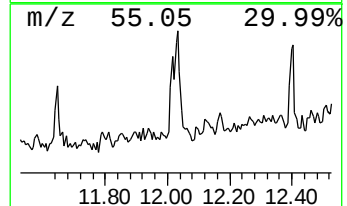
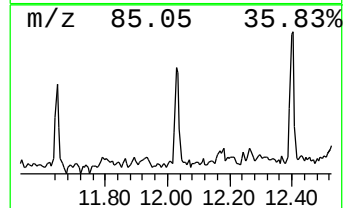
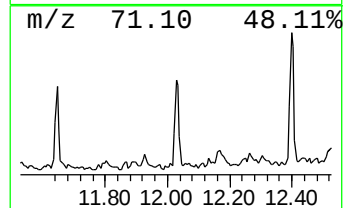
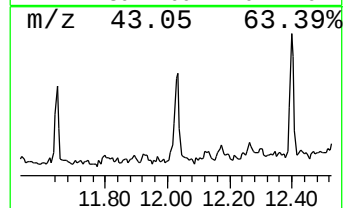
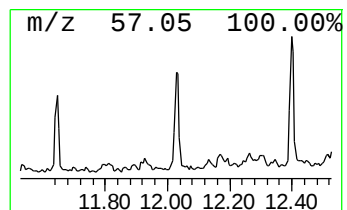
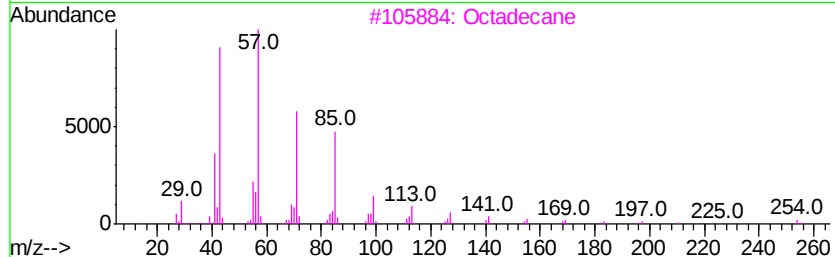
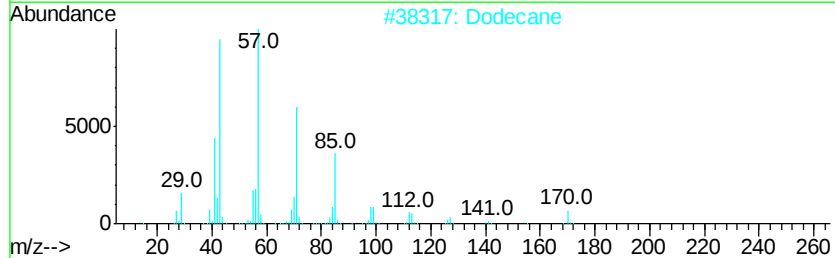
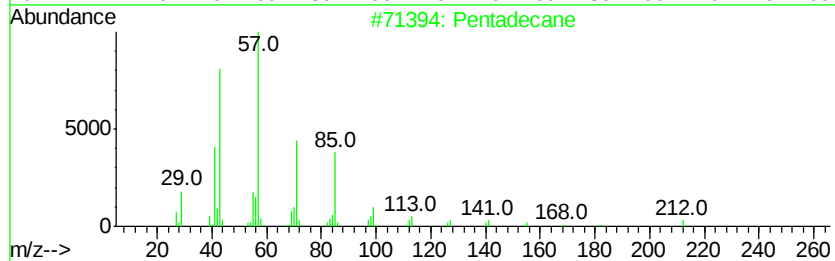
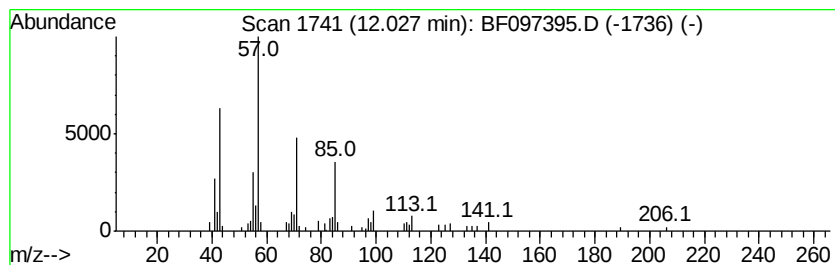
Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Pentadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.03	3.05 ng	116000	Phenanthrene-d10	10.92	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane	212	C15H32	000629-62-9	80
2	Dodecane	170	C12H26	000112-40-3	80
3	Octadecane	254	C18H38	000593-45-3	80
4	Sulfurous acid, butyl tridecyl e...	320	C17H36O3S	1000309-18-0	80
5	Sulfurous acid, butyl tetradecyl...	334	C18H38O3S	1000309-18-1	80



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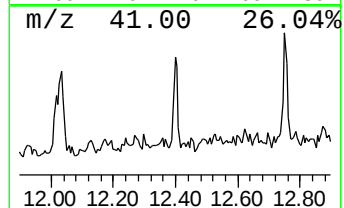
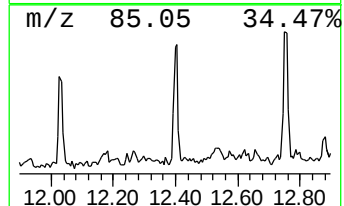
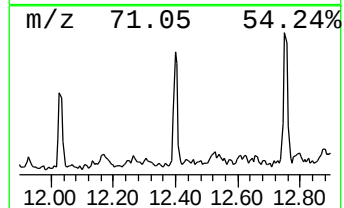
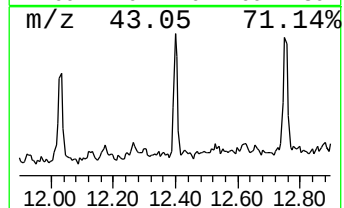
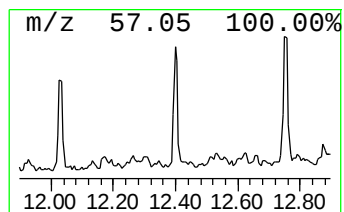
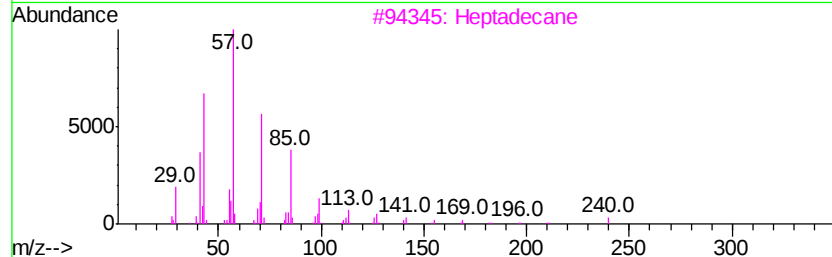
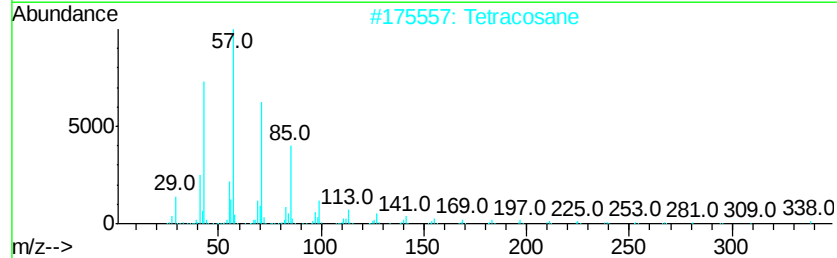
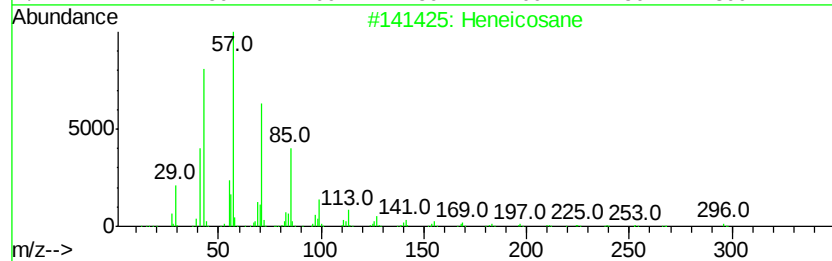
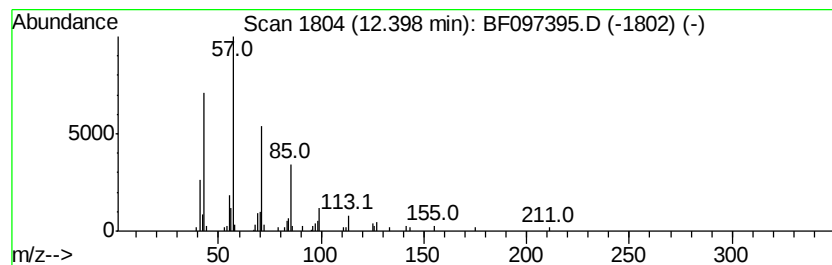
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Heneicosane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.40	2.18 ng	65485	Chrysene-d12	13.54

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	91
2	Tetracosane		338	C24H50	000646-31-1	91
3	Heptadecane		240	C17H36	000629-78-7	90
4	Silane, trichlorooctadecyl-		386	C18H37Cl3Si	000112-04-9	90
5	Hexadecane		226	C16H34	000544-76-3	86



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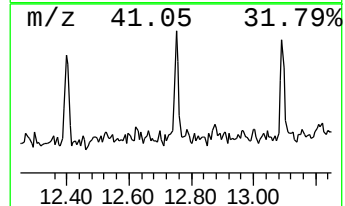
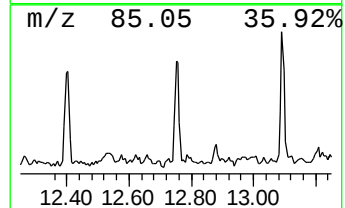
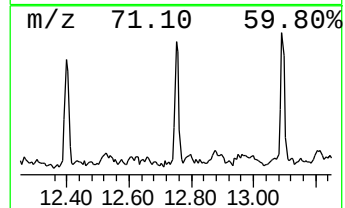
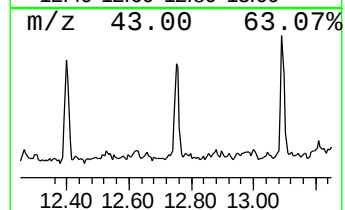
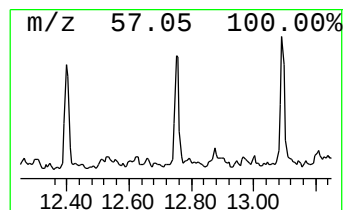
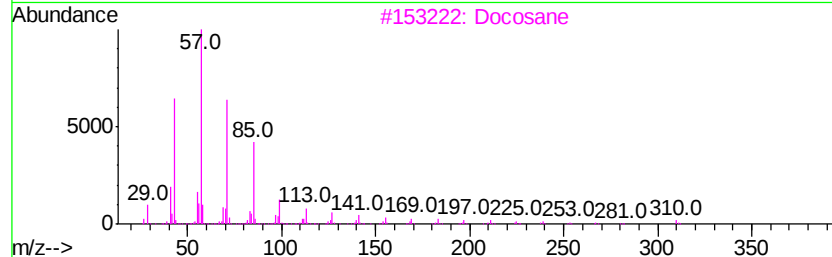
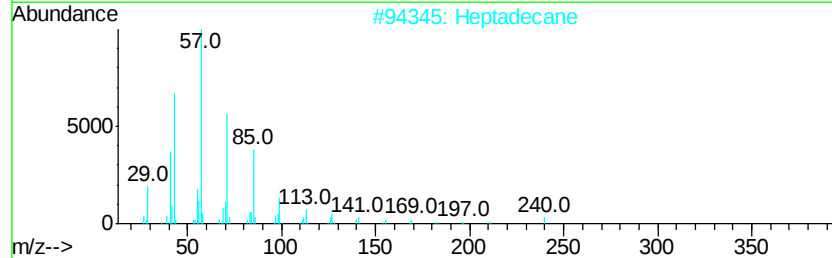
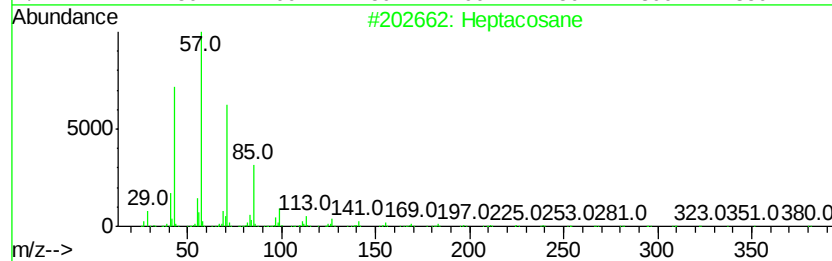
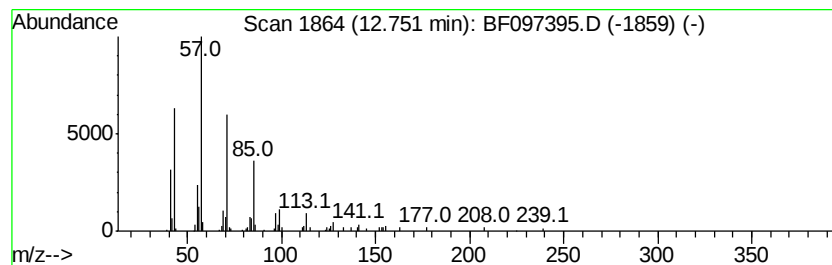
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Heptacosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.75	3.04 ng	91212	Chrysene-d12	13.54

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosane		380	C27H56	000593-49-7	91
2	Heptadecane		240	C17H36	000629-78-7	90
3	Docosane		310	C22H46	000629-97-0	90
4	Heneicosane		296	C21H44	000629-94-7	87
5	Eicosane		282	C20H42	000112-95-8	87



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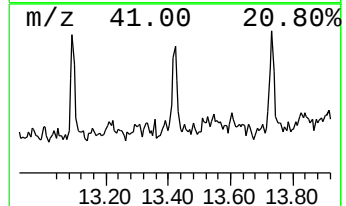
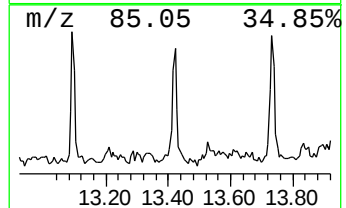
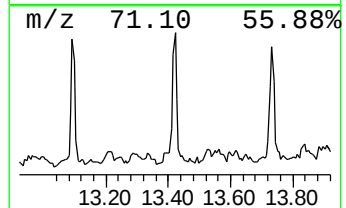
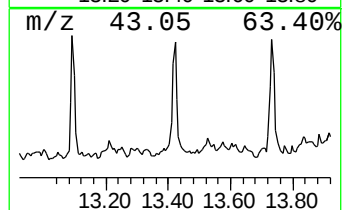
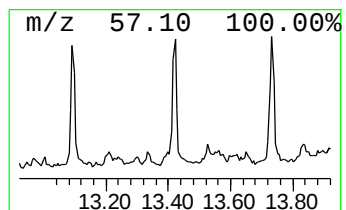
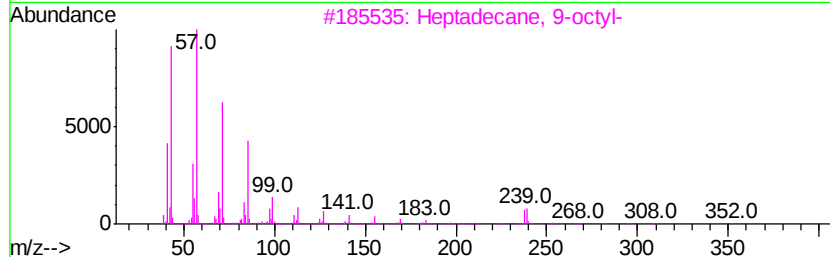
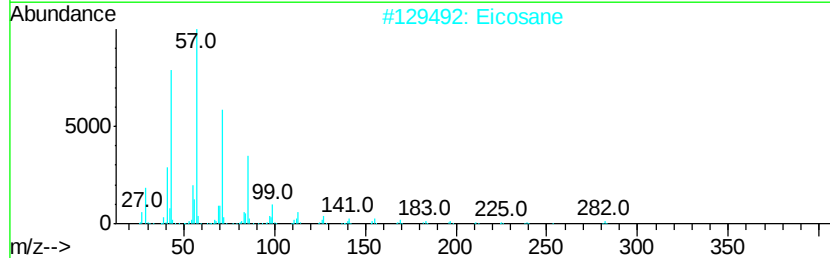
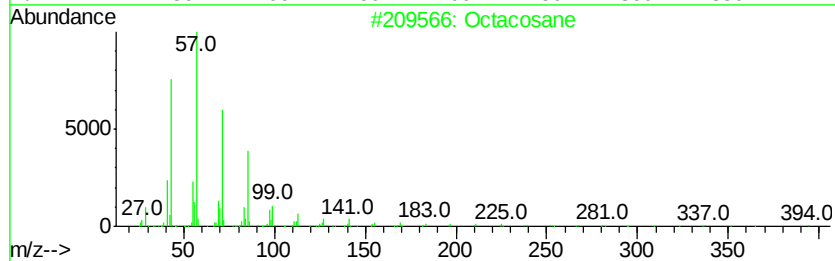
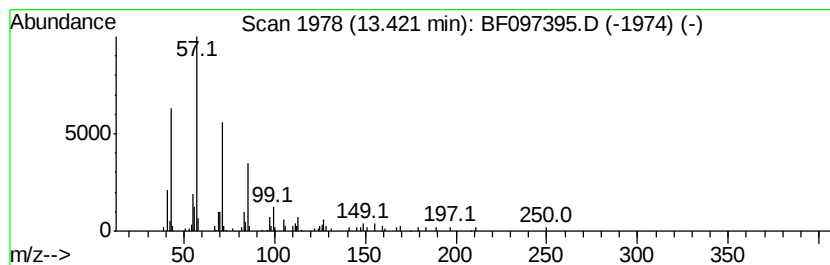
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Octacosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.42	3.72 ng	111788	Chrysene-d12	13.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	90
2		Eicosane	282	C20H42	000112-95-8	87
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87
4		Tetracosane	338	C24H50	000646-31-1	87
5		Heptacosane	380	C27H56	000593-49-7	87



Data Path : Z:\HPCHEM1\BNA F\DATA\BF080517\
Data File : BF097395.D
Acq On : 5 Aug 2017 14:17
Operator : SJ/JU
Sample : I4584-01 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
OK-01-080217

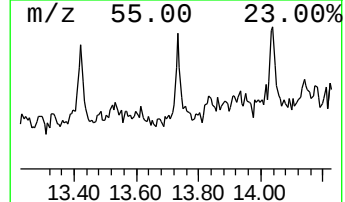
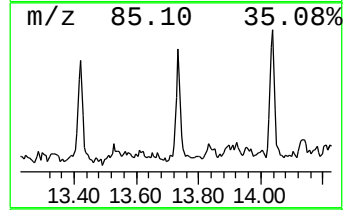
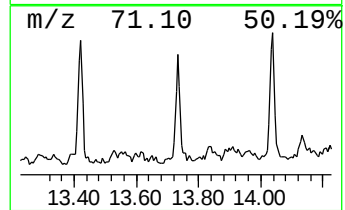
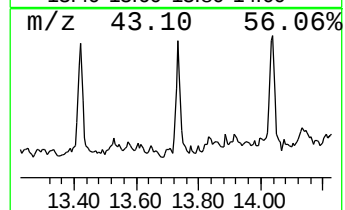
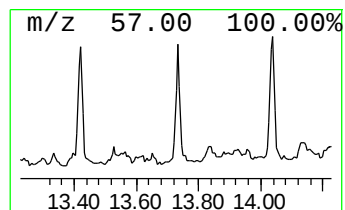
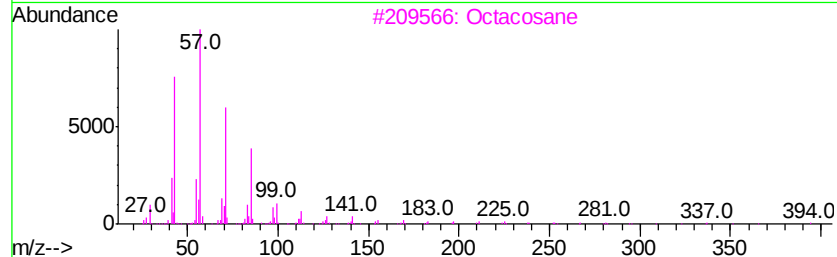
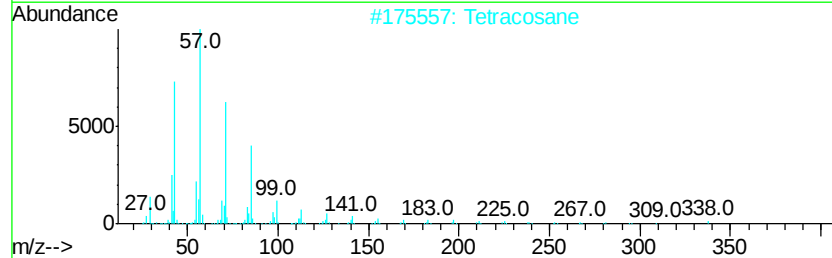
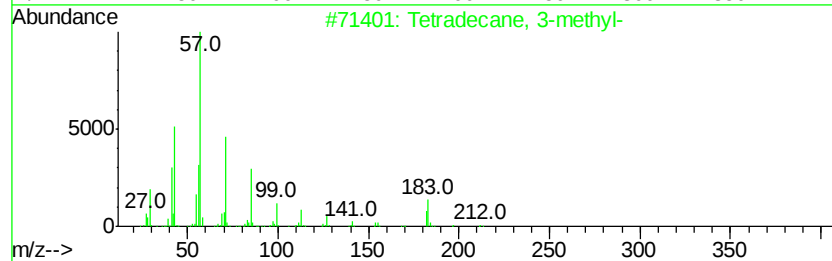
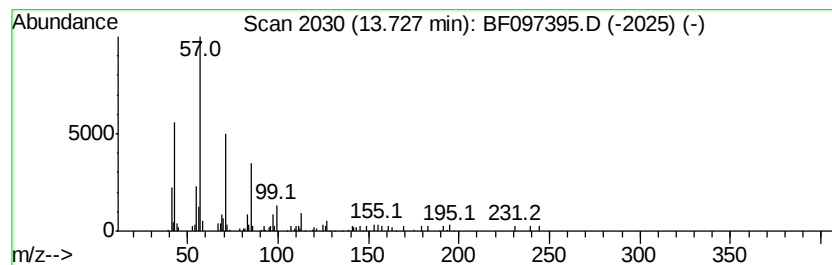
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Tetradecane, 3-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.73	4.80 ng	144303	Chrysene-d12	13.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetradecane, 3-methyl-	212	C15H32	018435-22-8	91
2		Tetracosane	338	C24H50	000646-31-1	90
3		Octacosane	394	C28H58	000630-02-4	90
4		Heptacosane	380	C27H56	000593-49-7	87
5		Heneicosane	296	C21H44	000629-94-7	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF080517\
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Acq On : 5 Aug 2017 14:17
Operator : SJ/JU
Sample : I4584-01 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
OK-01-080217

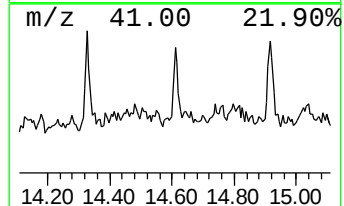
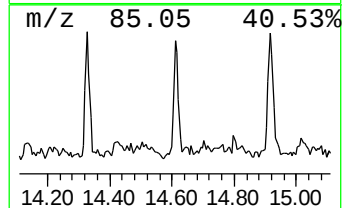
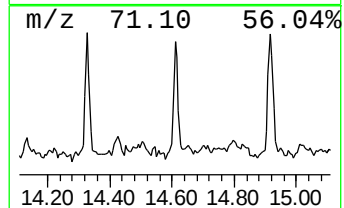
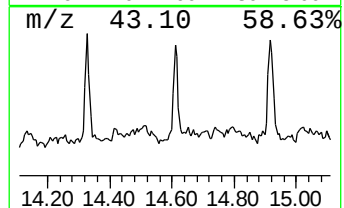
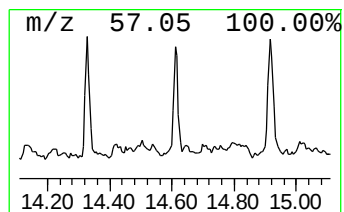
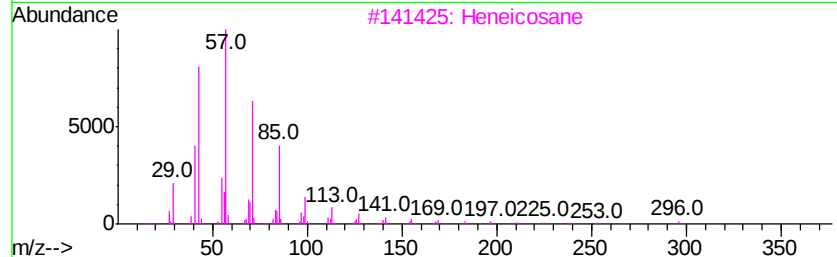
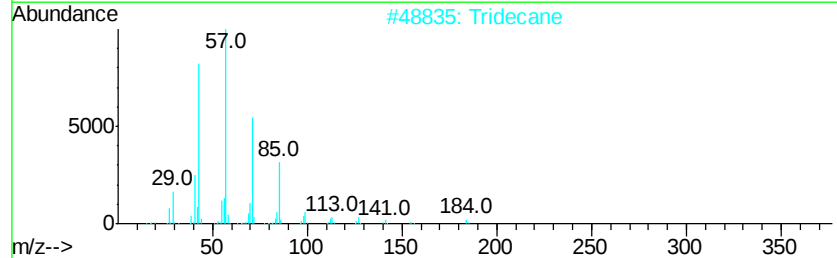
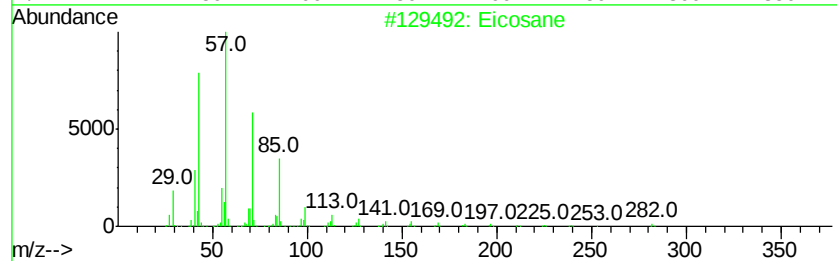
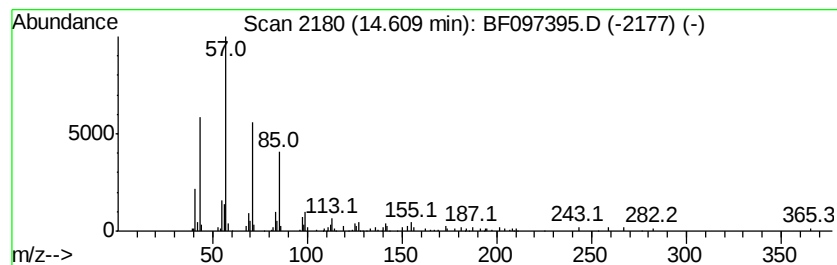
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Eicosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.61	3.75 ng	98486	Perylene-d12	14.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	93
2		Tridecane	184	C13H28	000629-50-5	86
3		Heneicosane	296	C21H44	000629-94-7	83
4		Heptacosane	380	C27H56	000593-49-7	83
5		Hexadecane	226	C16H34	000544-76-3	80



Data Path : Z:\HPCHEM1\BNA F\DATA\BF080517\
Data File : BF097395.D
Acq On : 5 Aug 2017 14:17
Operator : SJ/JU
Sample : I4584-01 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

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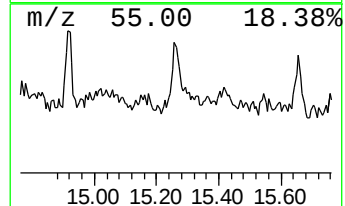
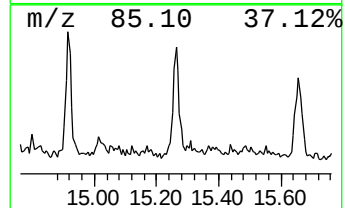
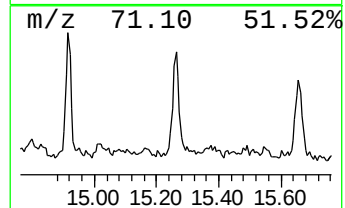
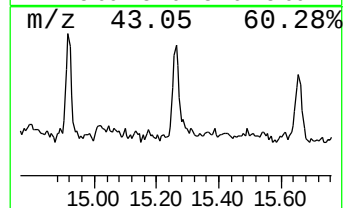
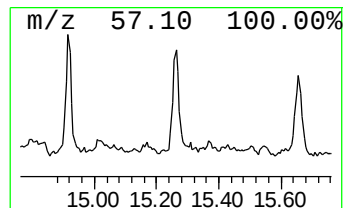
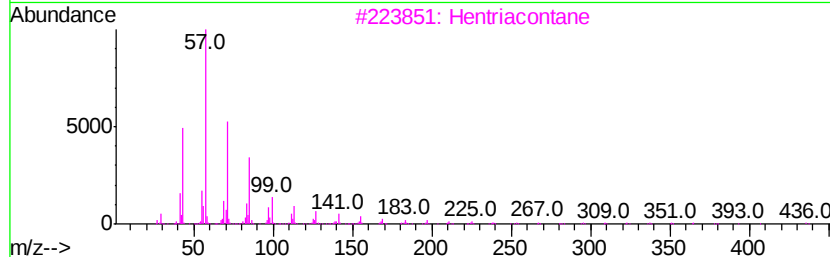
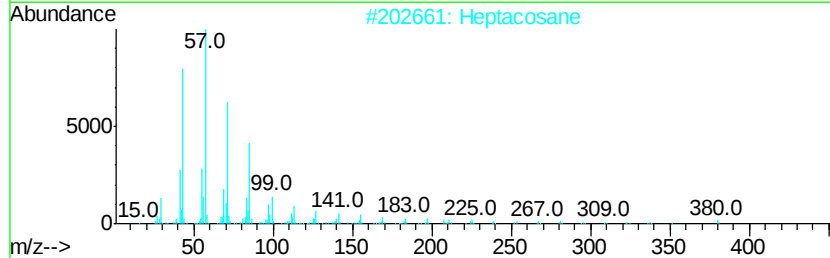
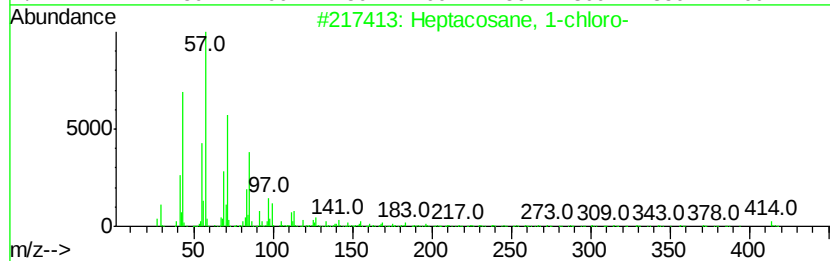
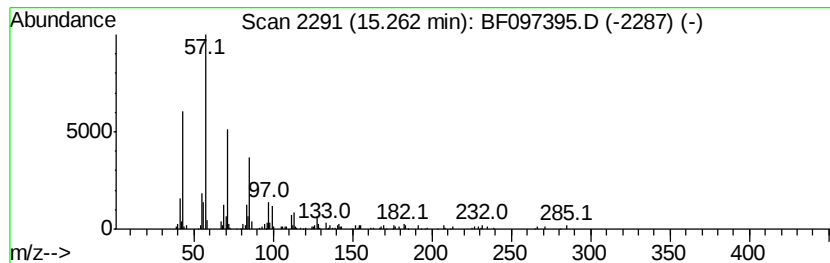
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Heptacosane, 1-chloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.26	5.83 ng	152919	Perylene-d12	14.88

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptacosane, 1-chloro-	414	C27H55Cl	062016-79-9	87
2		Heptacosane	380	C27H56	000593-49-7	86
3		Hentriacontane	437	C31H64	000630-04-6	86
4		Eicosane	282	C20H42	000112-95-8	83
5		Heneicosane	296	C21H44	000629-94-7	83



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF080517\
Data File : BF097395.D
Acq On : 5 Aug 2017 14:17
Operator : SJ/JU
Sample : I4584-01 5X
Misc :
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Instrument :
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ClientSampleId :
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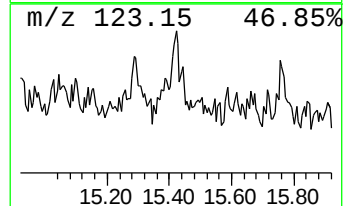
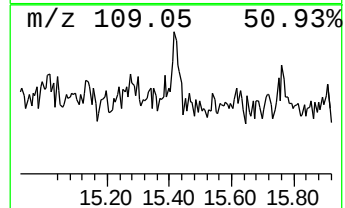
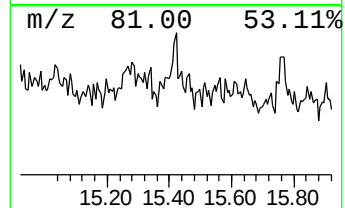
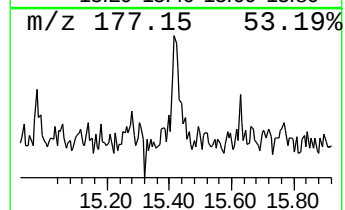
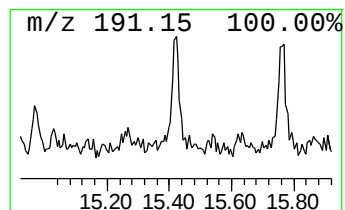
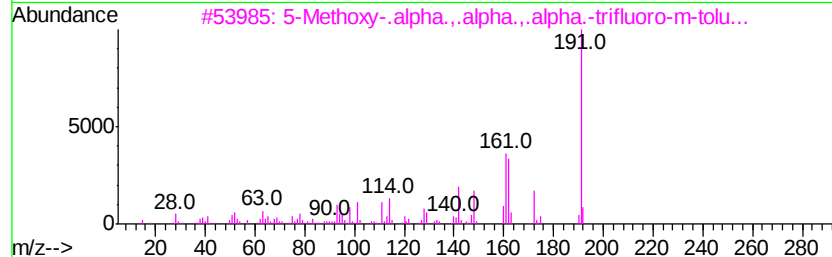
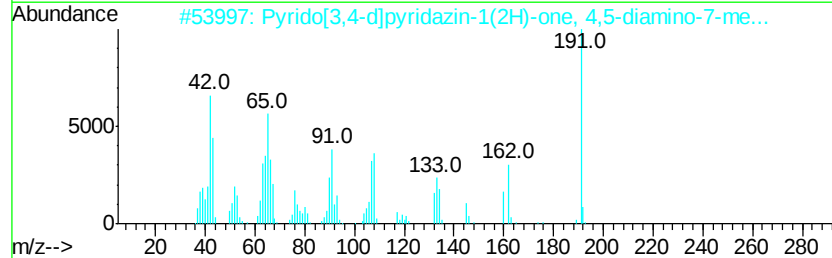
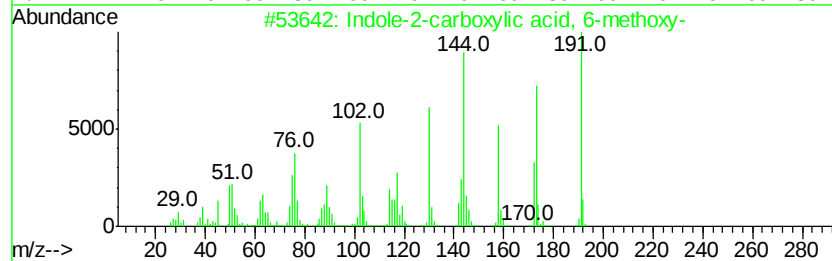
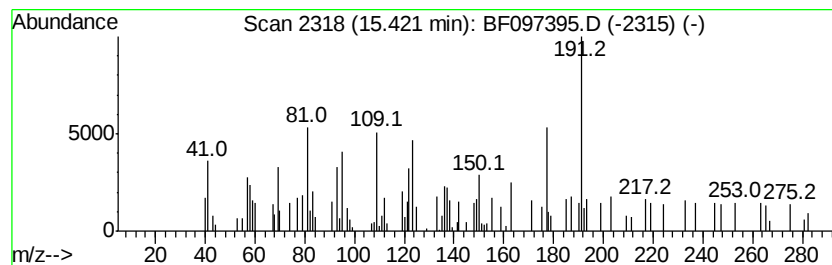
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown15.42 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.42	2.58 ng	67764	Perylene-d12	14.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indole-2-carboxylic acid, 6-meth...	191	C10H9NO3	016732-73-3	25
2		Pyrido[3,4-d]pyridazin-1(2H)-one...	191	C8H9N5O	1000263-69-7	15
3		5-Methoxy-.alpha.,.alpha.,.alpha...	191	C8H8F3NO	000349-55-3	11
4		4-Fluoro-3-trifluoromethylbenzoi...	390	C21H30F4O2	1000338-67-7	10
5		2-Fluoro-5-trifluoromethylbenzoi...	390	C21H30F4O2	1000338-62-3	10



Data Path : Z:\HPCHEM1\BNA F\DATA\BF080517\
Data File : BF097395.D
Acq On : 5 Aug 2017 14:17
Operator : SJ/JU
Sample : I4584-01 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
OK-01-080217

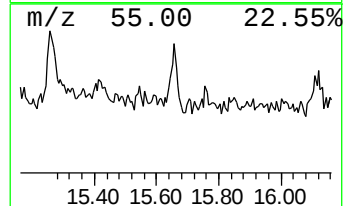
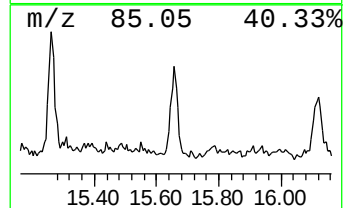
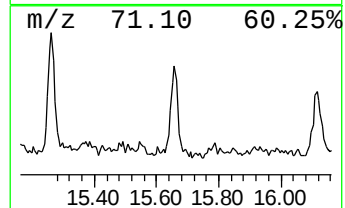
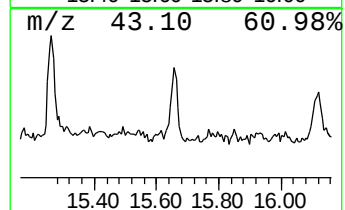
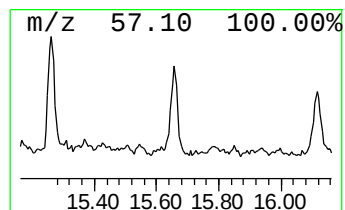
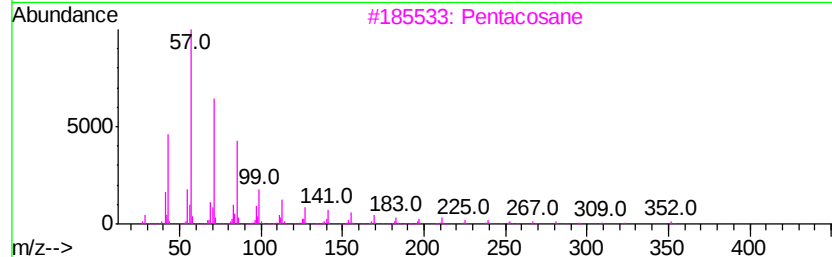
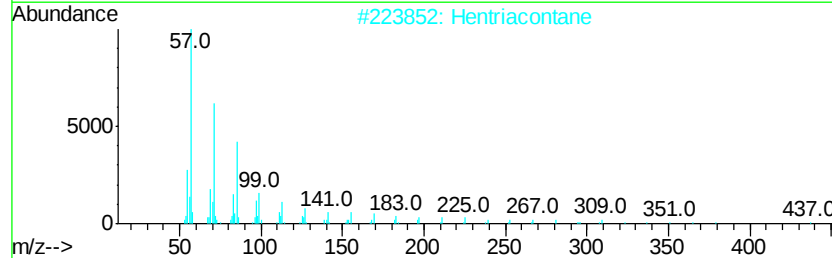
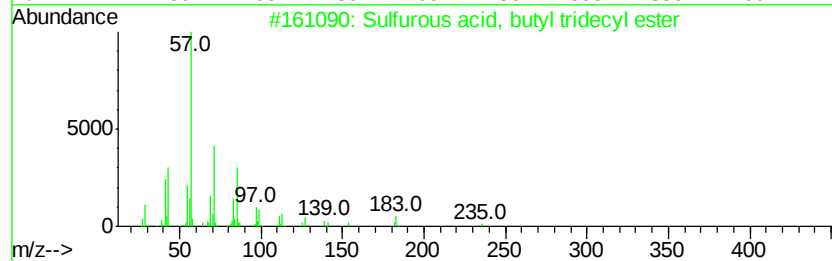
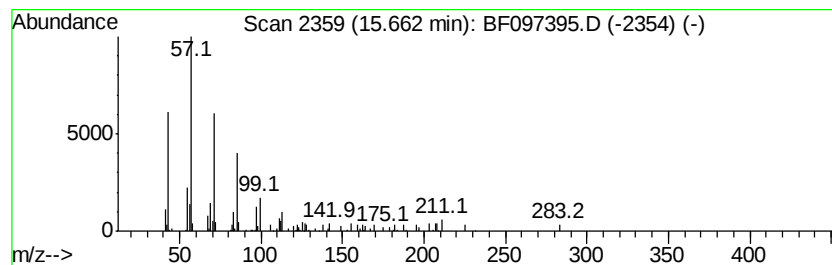
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Sulfurous acid, butyl tridecyl e... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.66	4.33 ng	113512	Perylene-d12	14.88

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfurous acid, butyl tridecyl e...	320	C17H36O3S	1000309-18-0	86
2			Hentriacontane	437	C31H64	000630-04-6	80
3			Pentacosane	352	C25H52	000629-99-2	80
4			Hexacosane	366	C26H54	000630-01-3	80
5			Tetracosane	338	C24H50	000646-31-1	72



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF080517\
Data File : BF097395.D
Acq On : 5 Aug 2017 14:17
Operator : SJ/JU
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Misc :
ALS Vial : 6 Sample Multiplier: 1

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ClientSampleId :
OK-01-080217

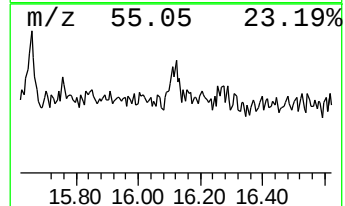
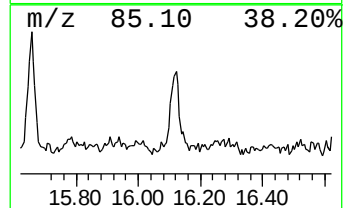
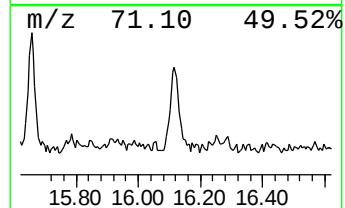
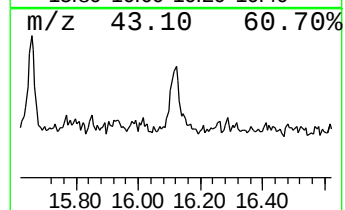
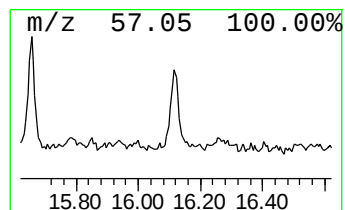
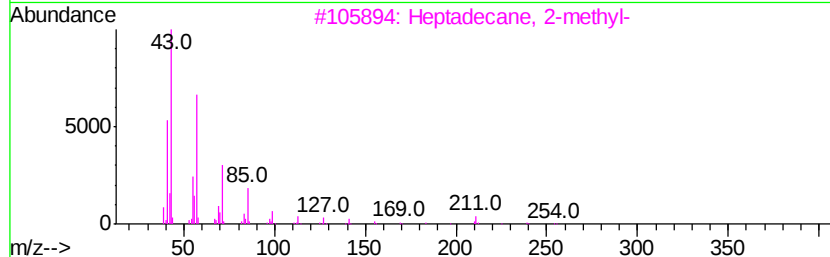
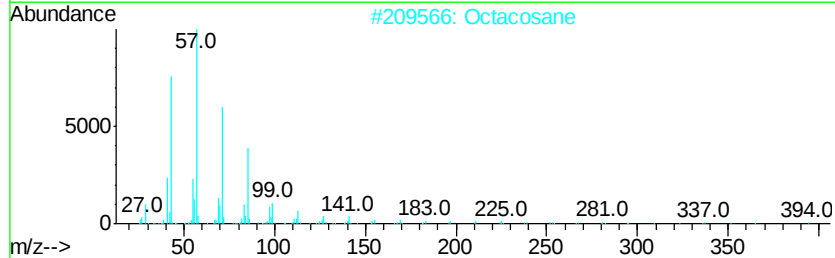
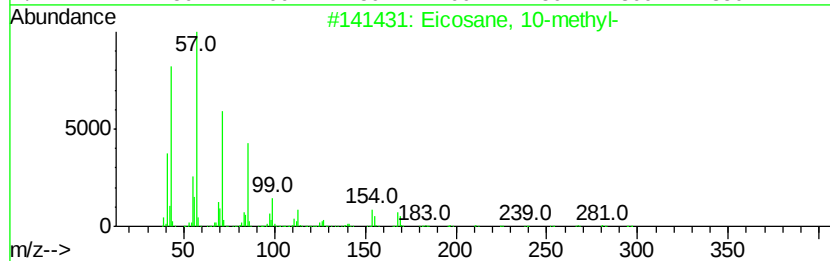
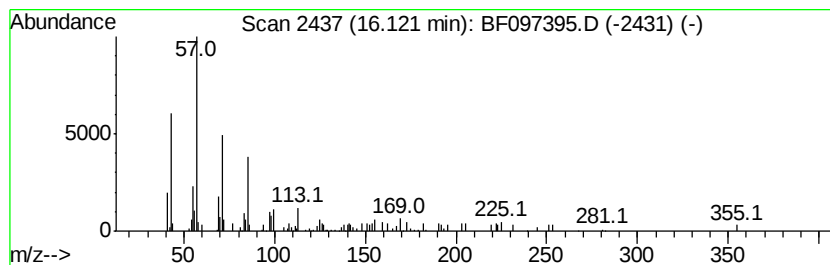
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Eicosane, 10-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.12	3.97 ng	104063	Perylene-d12	14.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane, 10-methyl-	296	C21H44	054833-23-7	86
2		Octacosane	394	C28H58	000630-02-4	80
3		Heptadecane, 2-methyl-	254	C18H38	001560-89-0	72
4		Tetratriacontane	479	C34H70	014167-59-0	72
5		Eicosane, 9-octyl-	394	C28H58	013475-77-9	72



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF080517\
 Data File : BF097395.D
 Acq On : 5 Aug 2017 14:17
 Operator : SJ/JU
 Sample : I4584-01 5X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OK-01-080217

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072517.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
unknown6.22	6.22	3.9 ng		114401	1	6.42	590909	20.0
Pentadecane	12.03	3.0 ng		116000	4	10.92	759514	20.0
Heneicosane	12.40	2.2 ng		65485	5	13.54	600814	20.0
Heptacosane	12.75	3.0 ng		91212	5	13.54	600814	20.0
Octacosane	13.42	3.7 ng		111788	5	13.54	600814	20.0
Tetradecane, 3-me...	13.73	4.8 ng		144303	5	13.54	600814	20.0
Eicosane	14.61	3.8 ng		98486	6	14.88	524804	20.0
Heptacosane, 1-ch...	15.26	5.8 ng		152919	6	14.88	524804	20.0
unknown15.42	15.42	2.6 ng		67764	6	14.88	524804	20.0
Sulfurous acid, b...	15.66	4.3 ng		113512	6	14.88	524804	20.0
Eicosane, 10-methyl-	16.12	4.0 ng		104063	6	14.88	524804	20.0