

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF071218\
 Data File : BF107345.D
 Acq On : 12 Jul 2018 18:01
 Operator : JU/SJ
 Sample : J3828-03 5X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-01-071018-B

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-BF061918.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	5.149	475	478	492	rBB	36997	43050	1.49%	0.349%
2	5.566	546	549	572	rBV	283831	319396	11.02%	2.592%
3	6.572	717	720	733	rBV	288627	341003	11.76%	2.767%
4	6.707	739	743	748	rBV	351599	338839	11.69%	2.750%
5	6.925	777	780	788	rBV	396434	356843	12.31%	2.896%
6	7.084	803	807	816	rVB	309017	267565	9.23%	2.171%
7	7.495	873	877	886	rBV	191074	211428	7.29%	1.716%
8	8.213	995	999	1004	rBV	519412	447464	15.44%	3.631%
9	8.784	1091	1096	1102	rBV2	30221	32875	1.13%	0.267%
10	9.284	1178	1181	1189	rBV	404584	399644	13.79%	3.243%
11	9.348	1189	1192	1196	rBV	48525	44325	1.53%	0.360%
12	9.684	1243	1249	1252	rBV	83578	99081	3.42%	0.804%
13	9.883	1278	1283	1287	rBV	63949	61500	2.12%	0.499%
14	9.978	1295	1299	1303	rVB	520191	463758	16.00%	3.763%
15	10.383	1364	1368	1370	rBV	87780	69522	2.40%	0.564%
16	10.601	1400	1405	1408	rBV	39376	44902	1.55%	0.364%
17	10.778	1431	1435	1440	rBV	178000	180782	6.24%	1.467%
18	10.854	1445	1448	1453	rBV	88021	101482	3.50%	0.824%
19	11.307	1521	1525	1527	rBV	76309	62358	2.15%	0.506%
20	11.336	1527	1530	1536	rVB2	33667	36771	1.27%	0.298%
21	11.472	1550	1553	1557	rBV	457703	439362	15.16%	3.565%
22	11.730	1595	1597	1601	rBV	57339	49868	1.72%	0.405%
23	11.842	1612	1616	1619	rVB	1278940	1102126	38.02%	8.944%
24	12.142	1664	1667	1669	rBV	52198	42083	1.45%	0.342%
25	12.536	1729	1734	1736	rBV	2003688	2898762	100.00%	23.524%
26	12.560	1736	1738	1744	rVV2	1875690	1859463	64.15%	15.090%
27	12.636	1747	1751	1754	rVV	578709	519387	17.92%	4.215%
28	12.701	1759	1762	1767	rBV	19695	36106	1.25%	0.293%
29	12.907	1794	1797	1799	rBV2	32498	31499	1.09%	0.256%
30	12.930	1799	1801	1807	rVB2	41545	43892	1.51%	0.356%
31	13.060	1819	1823	1825	rBV	395837	346429	11.95%	2.811%
32	13.236	1848	1853	1855	rBV4	38721	56323	1.94%	0.457%
33	13.360	1871	1874	1877	rBV2	67991	64764	2.23%	0.526%
34	14.030	1985	1988	1991	rBV	58702	55763	1.92%	0.453%

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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	14.136	2002	2006	2009	rBV	467671	464070	16.01%	3.766%
36	14.560	2076	2078	2082	rBV5	36197	50816	1.75%	0.412%
37	15.677	2264	2268	2277	rVB	264729	339532	11.71%	2.755%

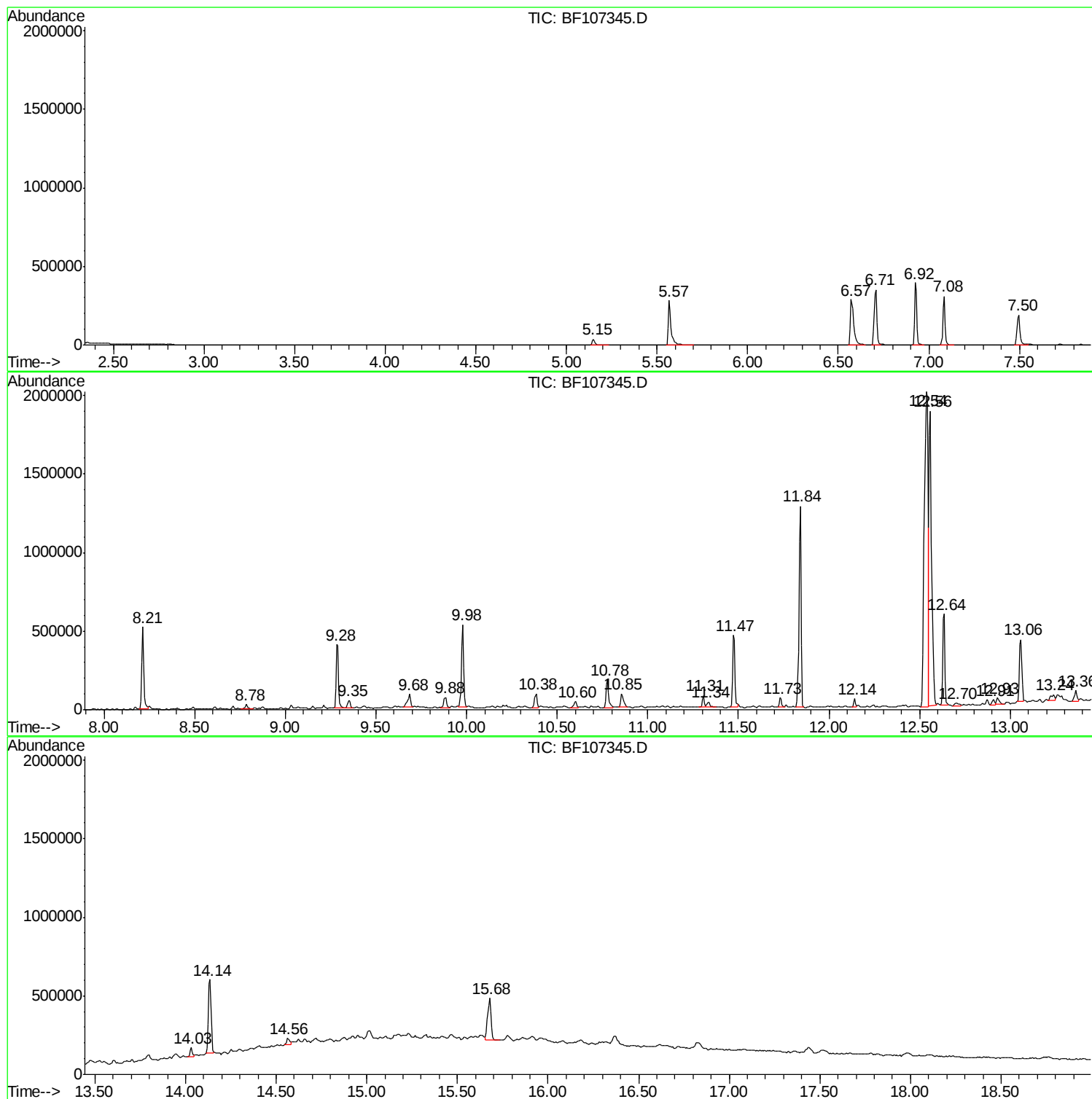
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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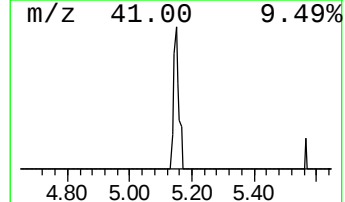
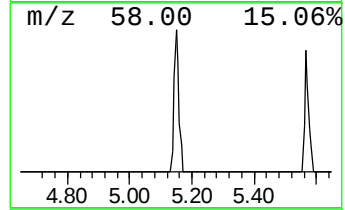
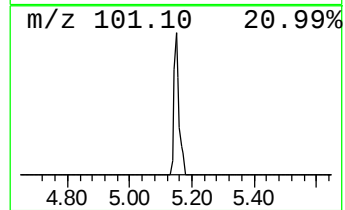
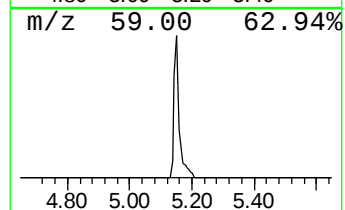
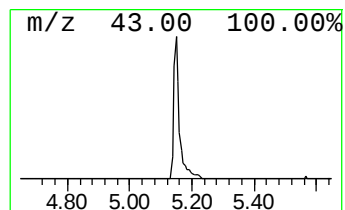
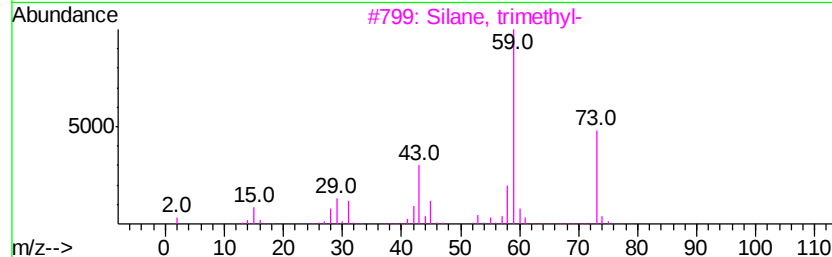
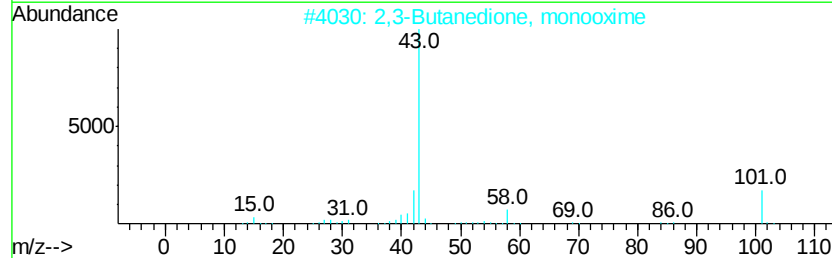
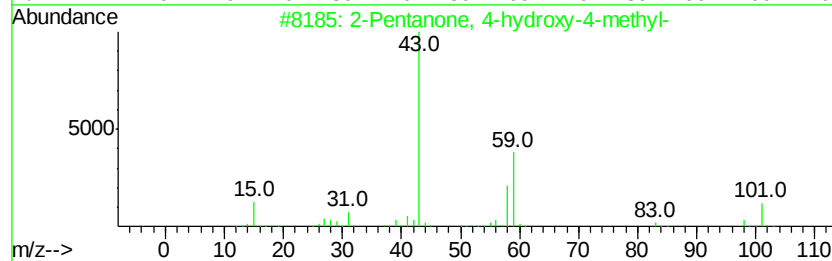
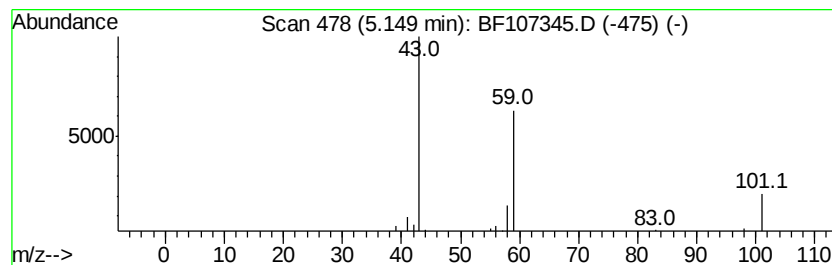
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.15	2.41 ng	43050	1,4-Dichlorobenzene-d4	6.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
3		Silane, trimethyl-	74	C3H10Si	000993-07-7	9
4		Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	9
5		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9



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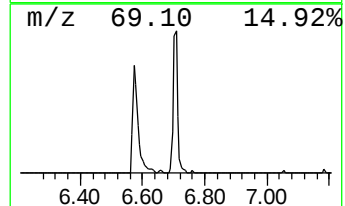
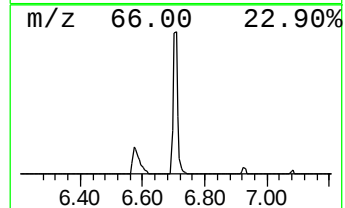
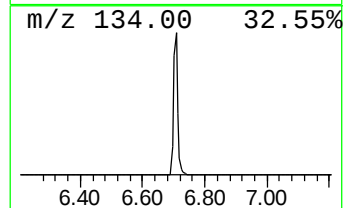
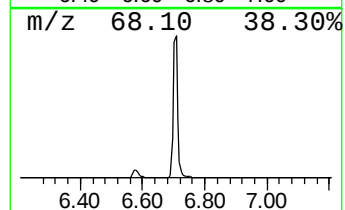
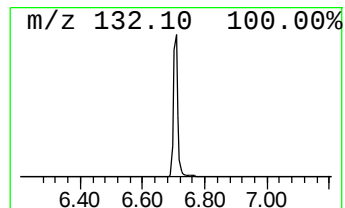
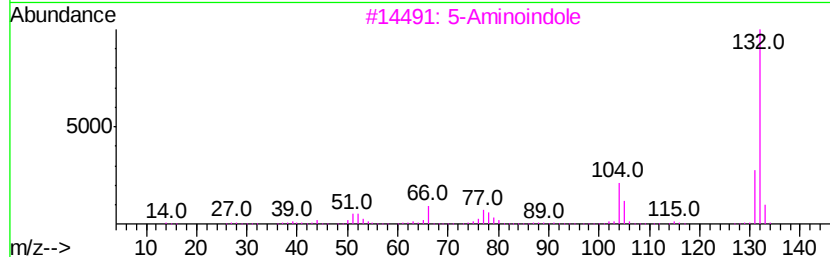
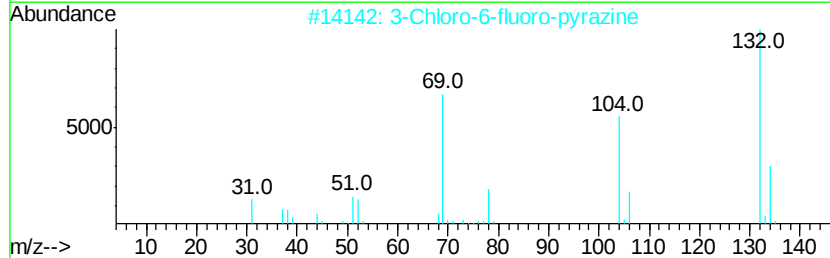
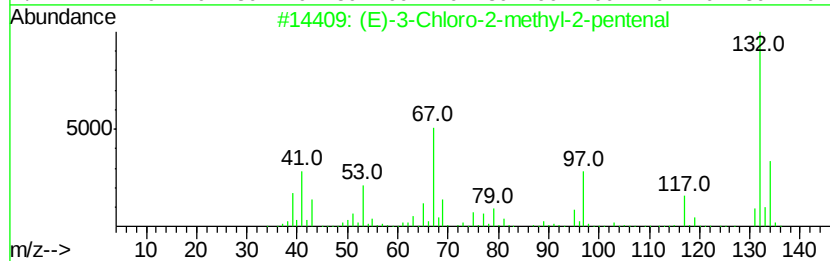
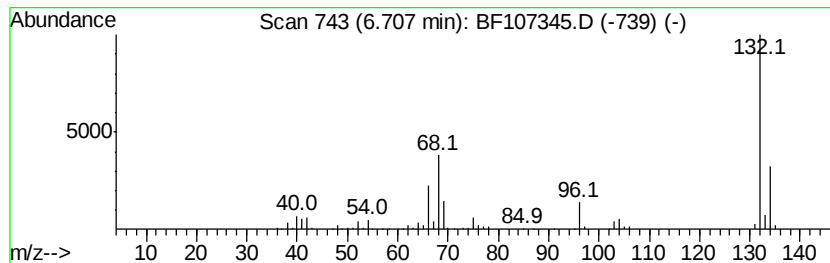
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 unknown6.71 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.71	18.99 ng	338839	1,4-Dichlorobenzene-d4	6.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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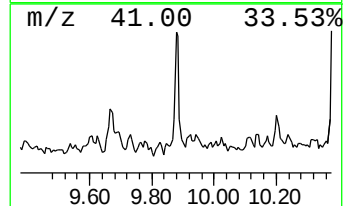
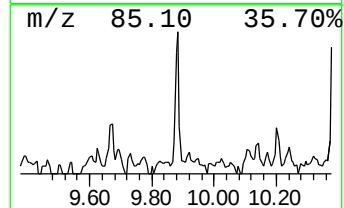
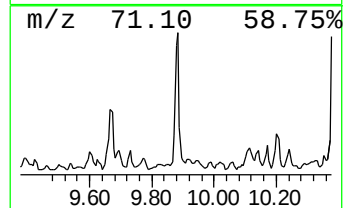
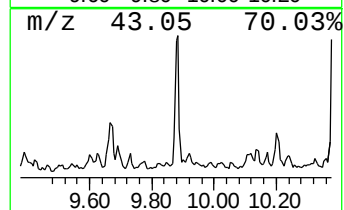
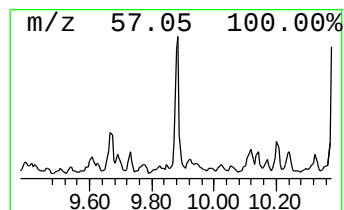
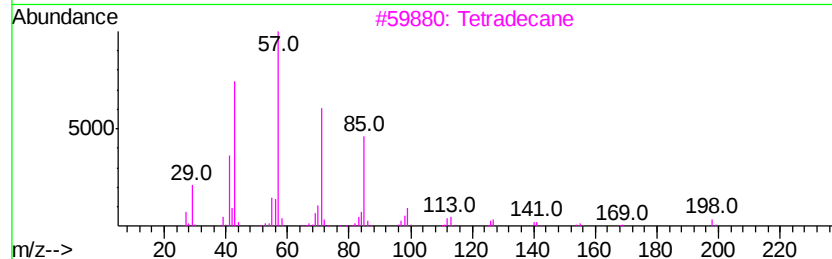
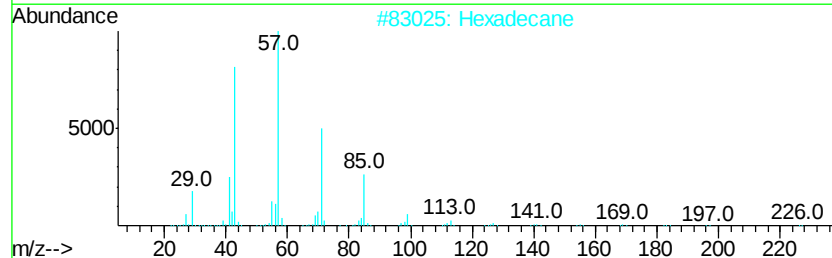
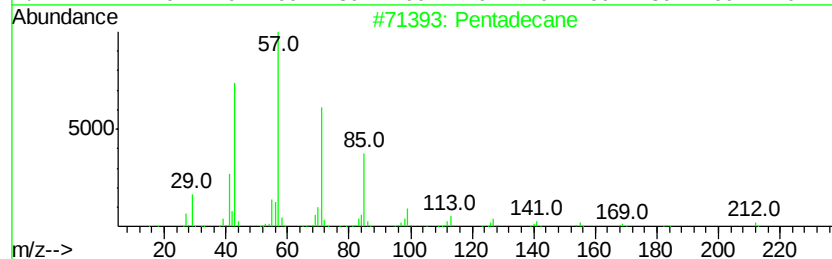
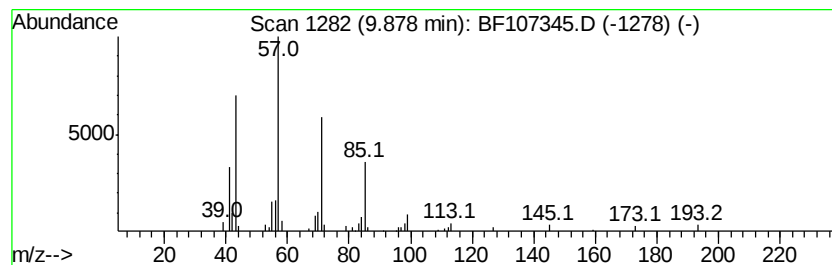
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Pentadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.88	2.65 ng	61500	Acenaphthene-d10	9.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane	212	C15H32	000629-62-9	86
2		Hexadecane	226	C16H34	000544-76-3	86
3		Tetradecane	198	C14H30	000629-59-4	80
4		Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	78
5		Decane, 2,3,6-trimethyl-	184	C13H28	062238-12-4	64



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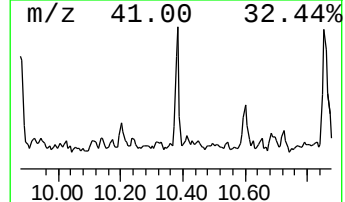
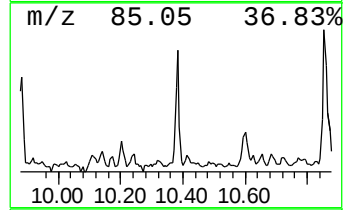
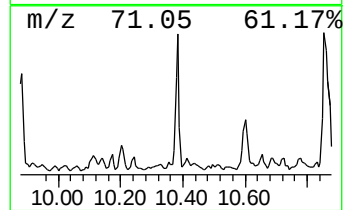
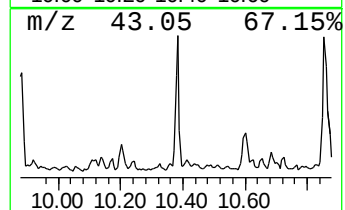
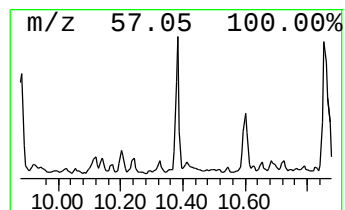
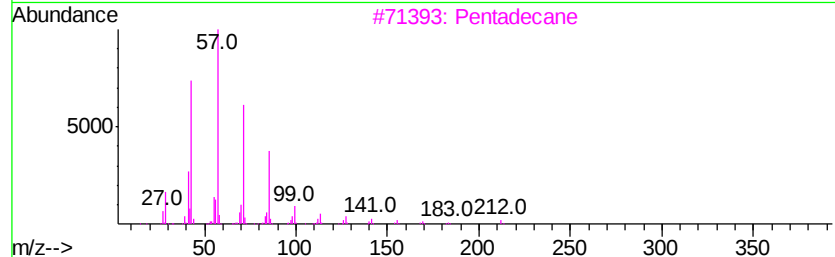
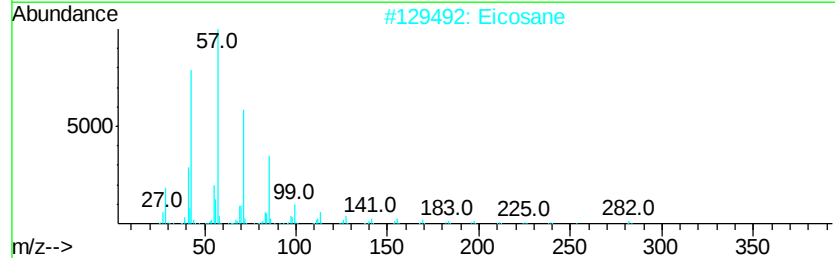
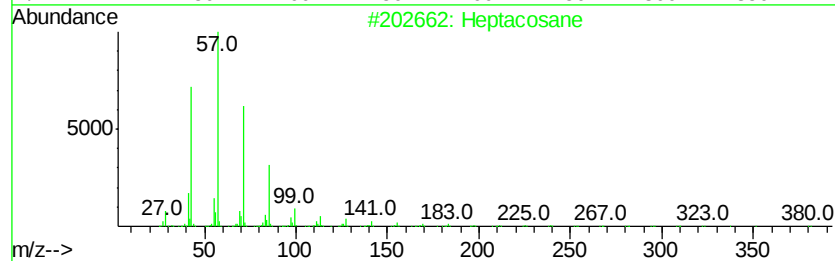
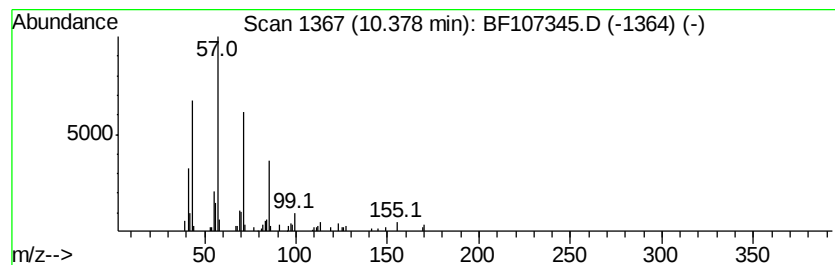
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 Heptacosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.38	3.00 ng	69522	Acenaphthene-d10	9.98

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosane		380	C27H56	000593-49-7	90
2	Eicosane		282	C20H42	000112-95-8	87
3	Pentadecane		212	C15H32	000629-62-9	87
4	Hexadecane		226	C16H34	000544-76-3	86
5	Tridecane		184	C13H28	000629-50-5	86



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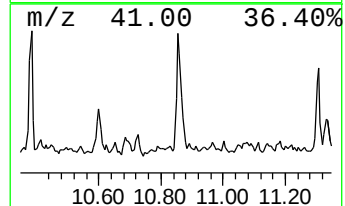
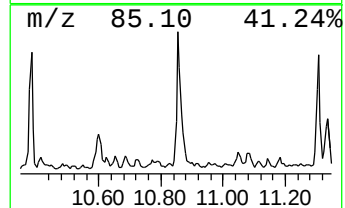
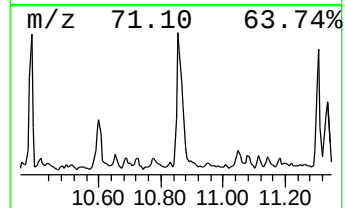
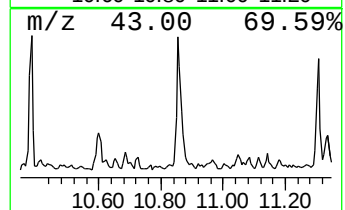
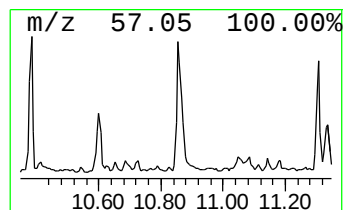
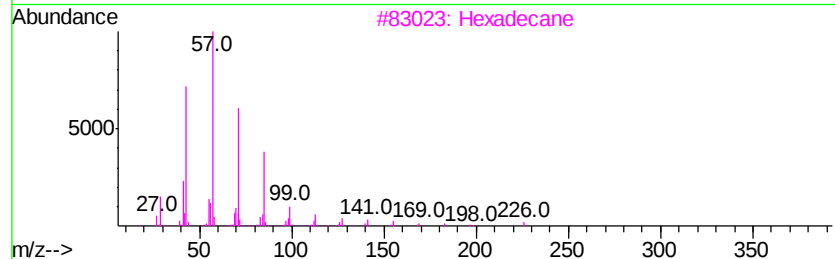
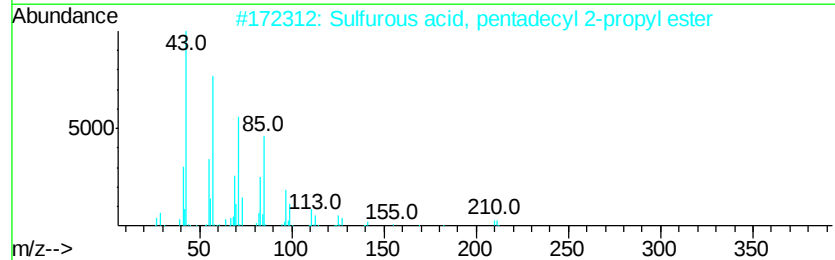
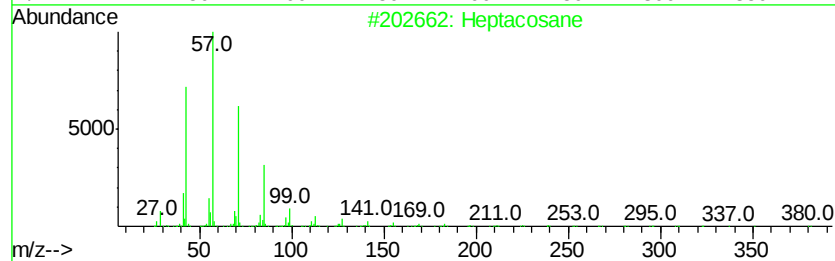
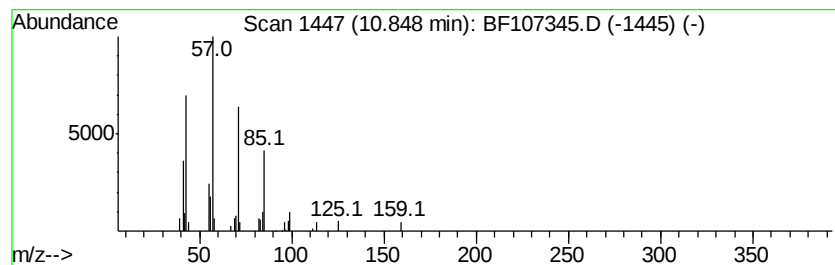
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Peak Number 6 Sulfurous acid, pentadecyl ... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.85	4.62 ng	101482	Phenanthrene-d10	11.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptacosane	380	C27H56	000593-49-7	90
2		Sulfurous acid, pentadecyl 2-pro...	334	C18H38O3S	1000309-12-6	86
3		Hexadecane	226	C16H34	000544-76-3	86
4		Sulfurous acid, 2-propyl tetra...	320	C17H36O3S	1000309-12-5	86
5		Dodecane, 1-iodo-	296	C12H25I	004292-19-7	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071218\
Data File : BF107345.D
Acq On : 12 Jul 2018 18:01
Operator : JU/SJ
Sample : J3828-03 5X
Misc :
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ClientSampleId :
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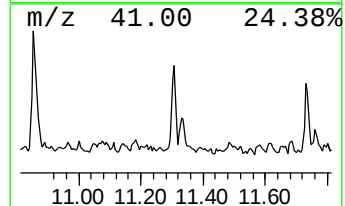
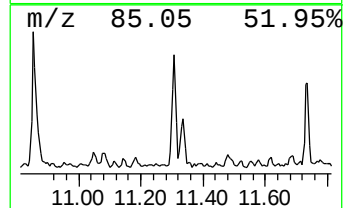
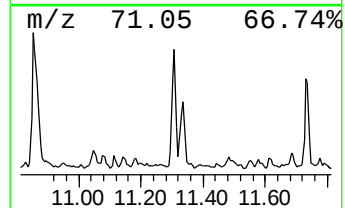
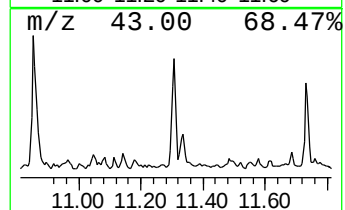
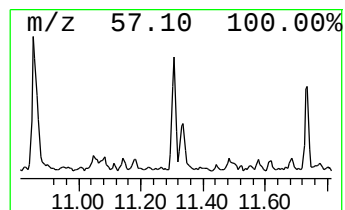
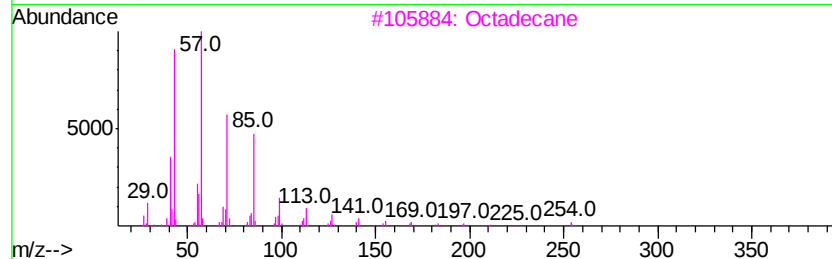
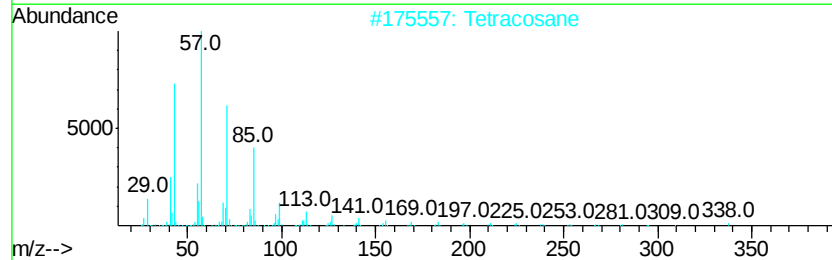
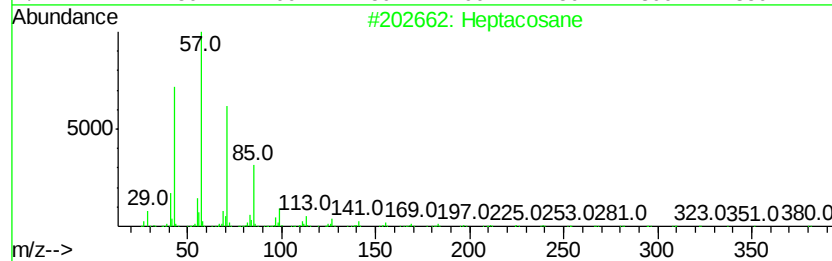
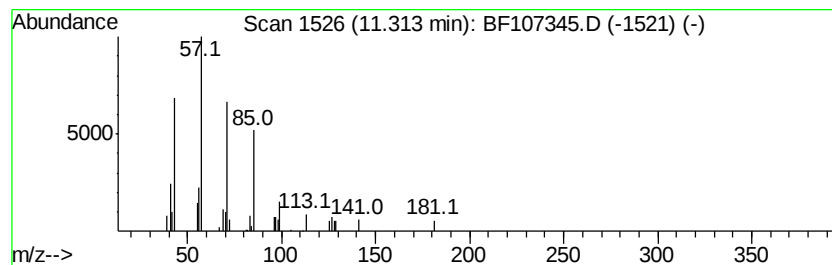
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Tetracosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.31	2.84 ng	62358	Phenanthrene-d10	11.47

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptacosane	380	C27H56	000593-49-7	91
2			Tetracosane	338	C24H50	000646-31-1	91
3			Octadecane	254	C18H38	000593-45-3	91
4			Pentadecane	212	C15H32	000629-62-9	91
5			Tritetracontane	605	C43H88	007098-21-7	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071218\
Data File : BF107345.D
Acq On : 12 Jul 2018 18:01
Operator : JU/SJ
Sample : J3828-03 5X
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ClientSampleId :
EO-01-071018-B

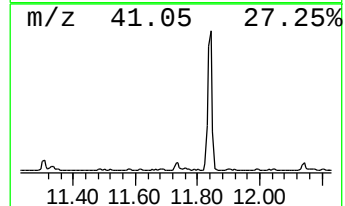
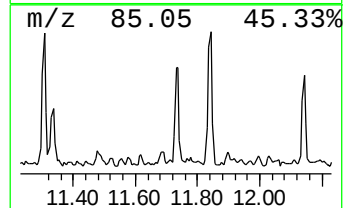
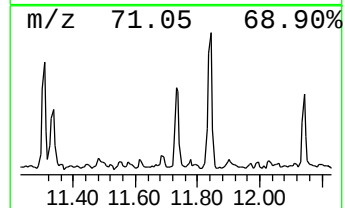
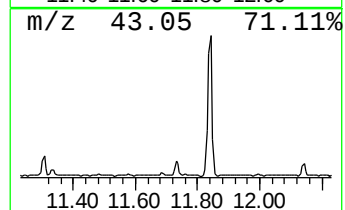
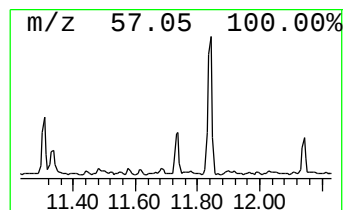
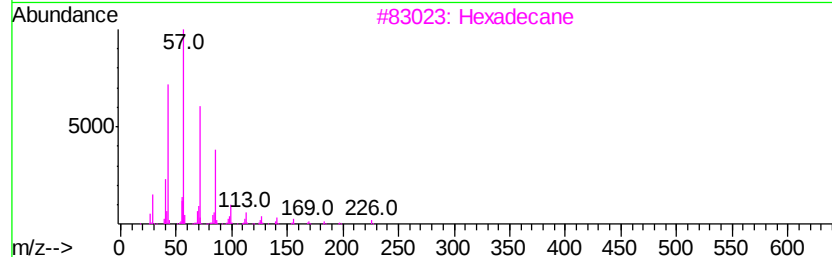
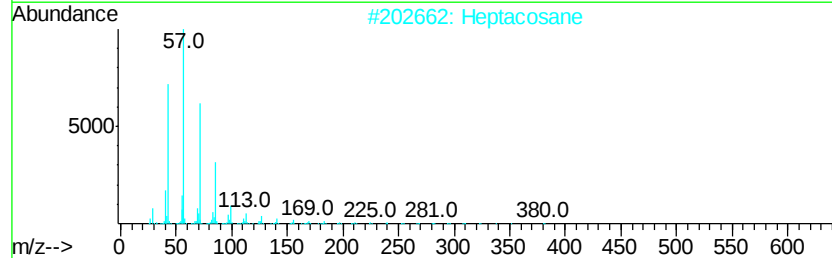
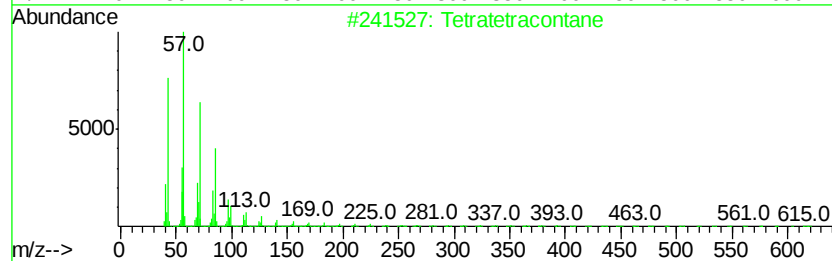
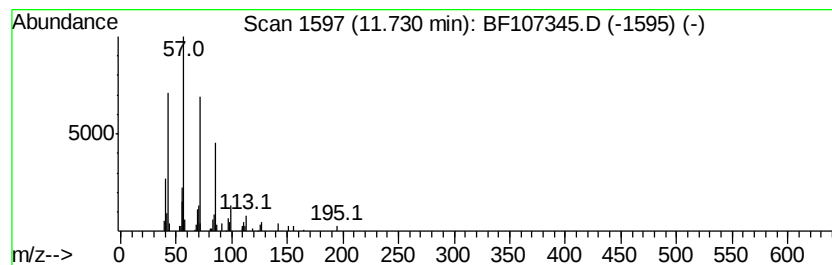
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Tetratetracontane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.73	2.27 ng	49868	Phenanthrene-d10	11.47

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetratetracontane	619	C44H90	007098-22-8	90
2		Heptacosane	380	C27H56	000593-49-7	90
3		Hexadecane	226	C16H34	000544-76-3	90
4		Eicosane, 10-methyl-	296	C21H44	054833-23-7	87
5		Heneicosane	296	C21H44	000629-94-7	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071218\
Data File : BF107345.D
Acq On : 12 Jul 2018 18:01
Operator : JU/SJ
Sample : J3828-03 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

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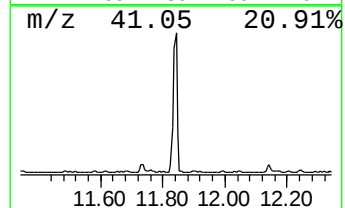
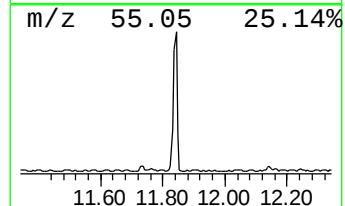
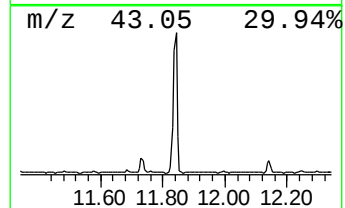
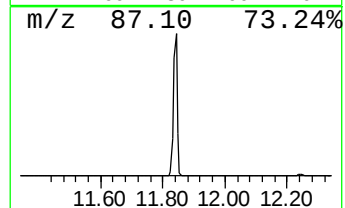
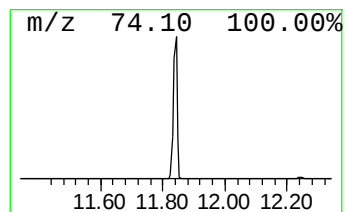
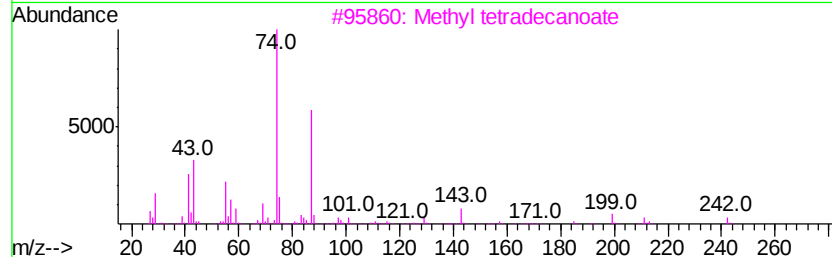
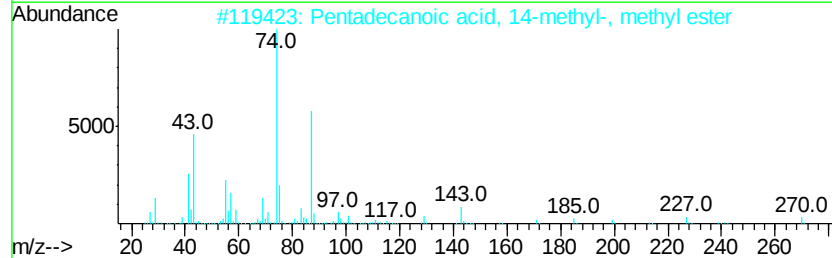
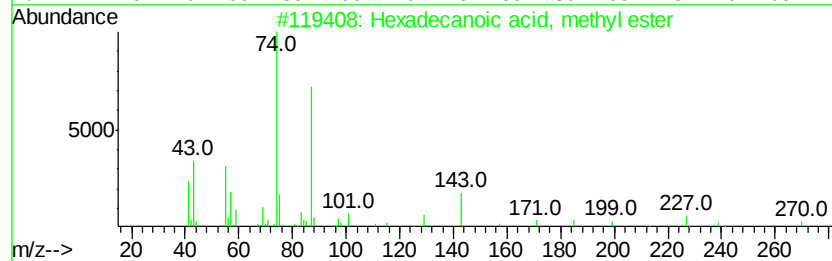
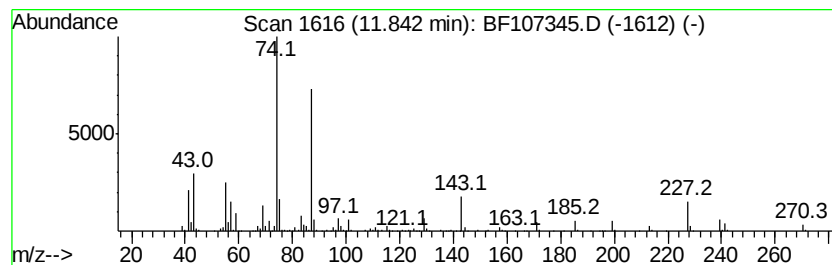
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Hexadecanoic acid, methyl e... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.84	50.17 ng	1102130	Phenanthrene-d10	11.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	98
2		Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	95
3		Methyl tetradecanoate	242	C15H30O2	000124-10-7	93
4		Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	90
5		Hexadecanoic acid, 15-methyl-, m...	284	C18H36O2	006929-04-0	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071218\
Data File : BF107345.D
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Misc :
ALS Vial : 17 Sample Multiplier: 1

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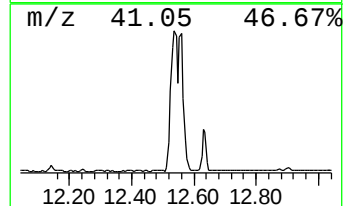
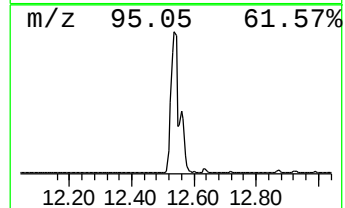
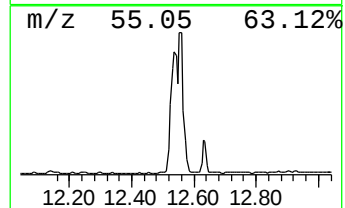
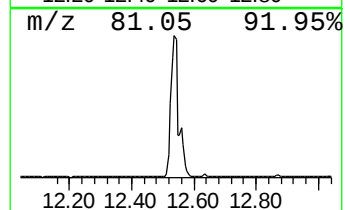
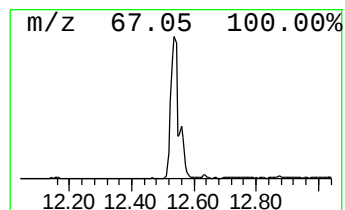
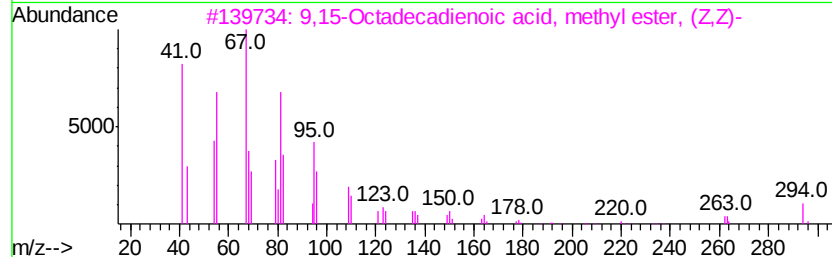
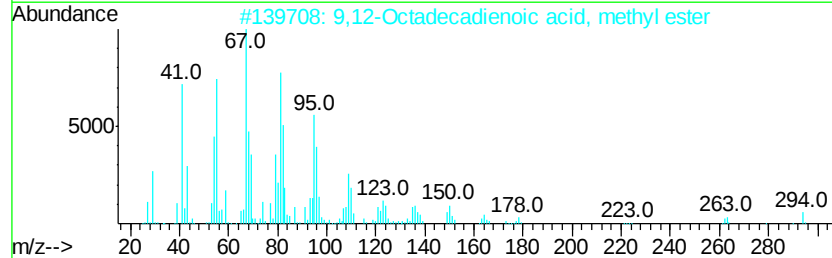
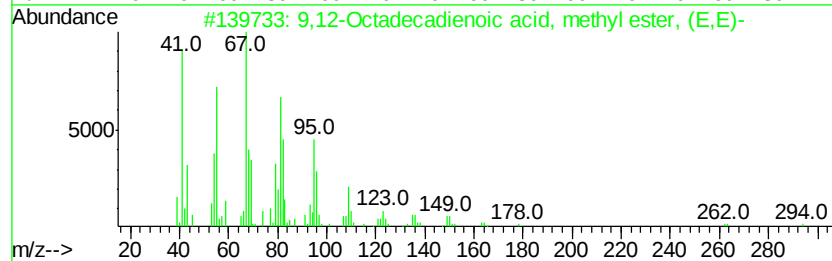
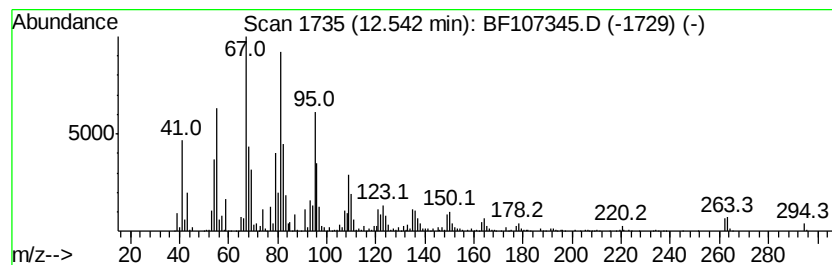
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 9,12-Octadecadienoic acid, ... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.54	131.95 ng	2898760	Phenanthrene-d10	11.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	99
2		9,12-Octadecadienoic acid, methy...	294	C19H34O2	002462-85-3	99
3		9,15-Octadecadienoic acid, methy...	294	C19H34O2	017309-05-6	99
4		10,13-Octadecadienoic acid, meth...	294	C19H34O2	056554-62-2	97
5		Methyl 9-cis,11-trans-octadecadi...	294	C19H34O2	1000336-44-0	97



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071218\
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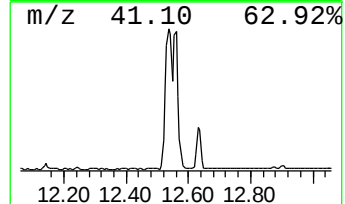
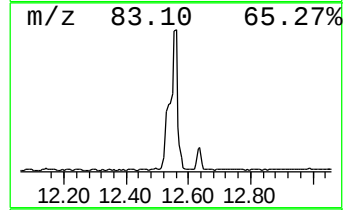
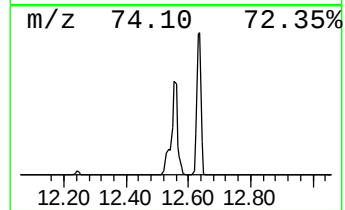
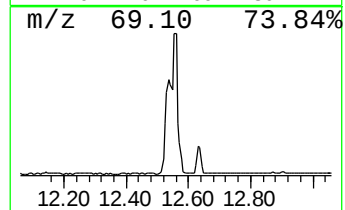
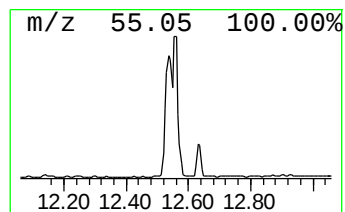
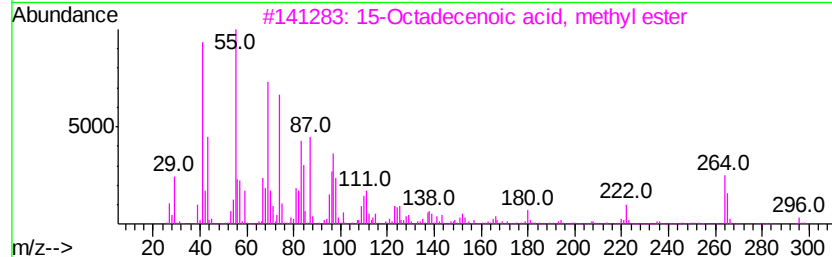
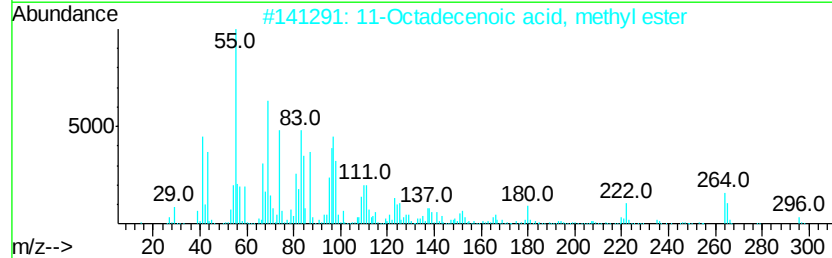
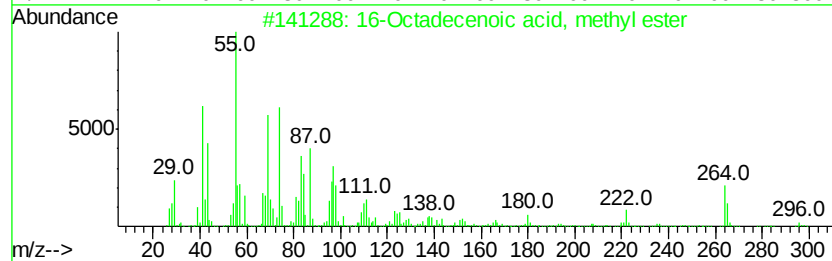
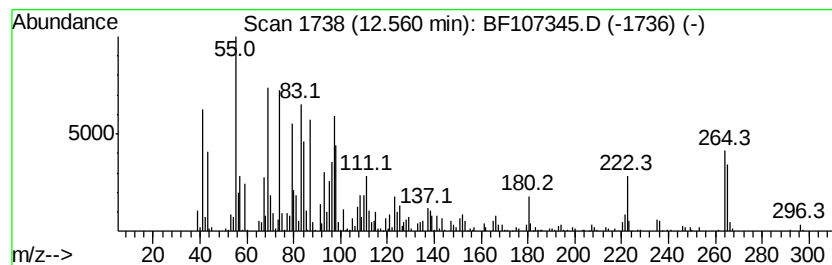
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 16-Octadecenoic acid, methyl ester Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.56	84.64 ng	1859460	Phenanthrene-d10	11.47

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			16-Octadecenoic acid, methyl ester	296	C19H36O2	056554-49-5	99
2			11-Octadecenoic acid, methyl ester	296	C19H36O2	052380-33-3	97
3			15-Octadecenoic acid, methyl ester	296	C19H36O2	004764-72-1	95
4			14-Octadecenoic acid, methyl ester	296	C19H36O2	056554-48-4	91
5			6-Octadecenoic acid, methyl ester	296	C19H36O2	052355-31-4	89



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071218\
Data File : BF107345.D
Acq On : 12 Jul 2018 18:01
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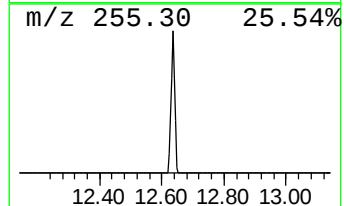
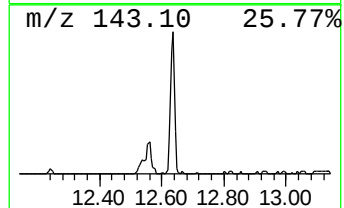
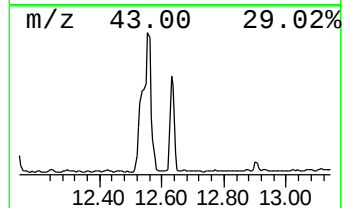
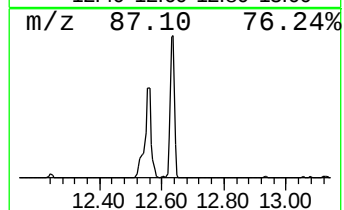
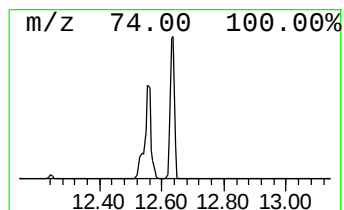
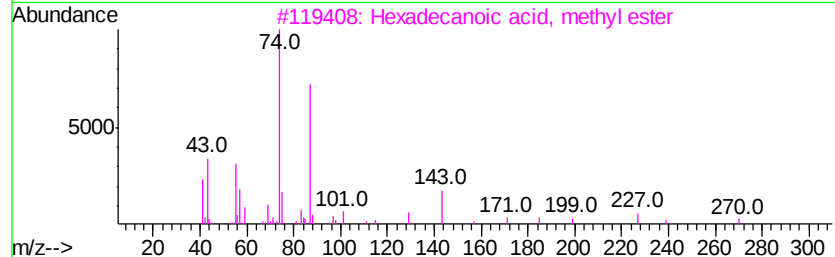
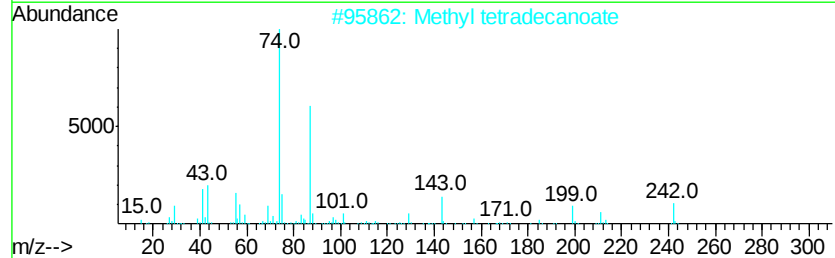
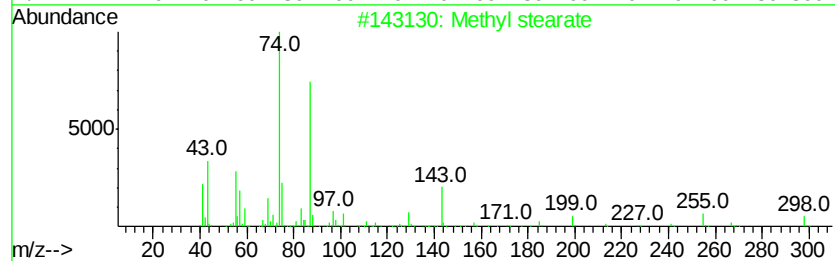
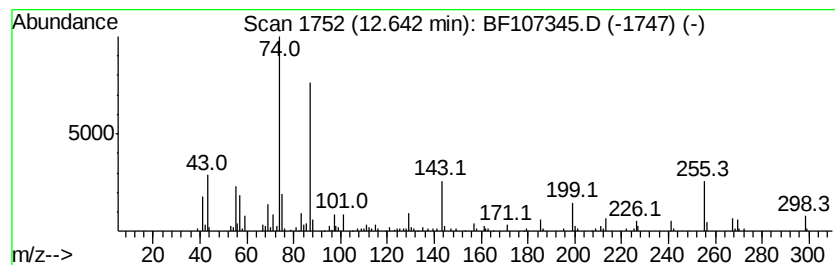
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Methyl stearate Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.64	23.64 ng	519387	Phenanthrene-d10	11.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Methyl stearate	298	C19H38O2	000112-61-8	97
2		Methyl tetradecanoate	242	C15H30O2	000124-10-7	93
3		Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	91
4		Heptadecanoic acid, 16-methyl-, ...	298	C19H38O2	005129-61-3	91
5		Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071218\
Data File : BF107345.D
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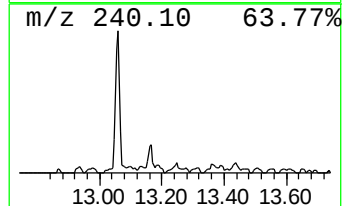
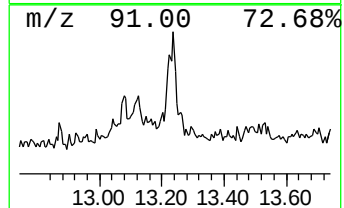
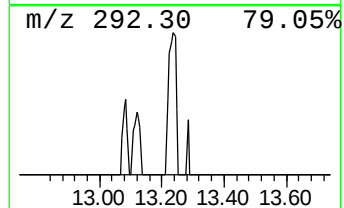
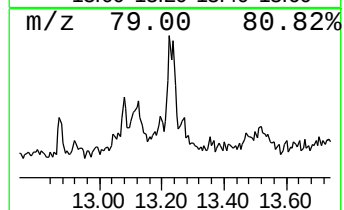
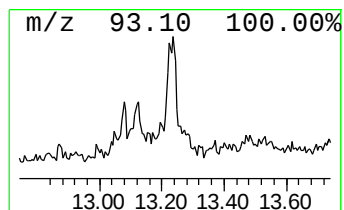
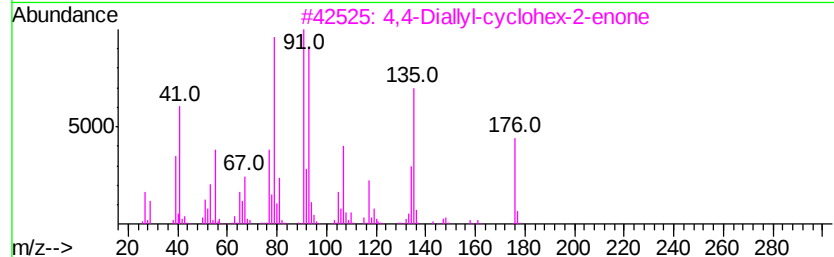
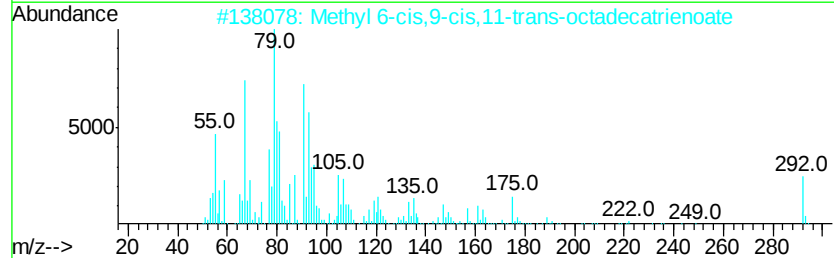
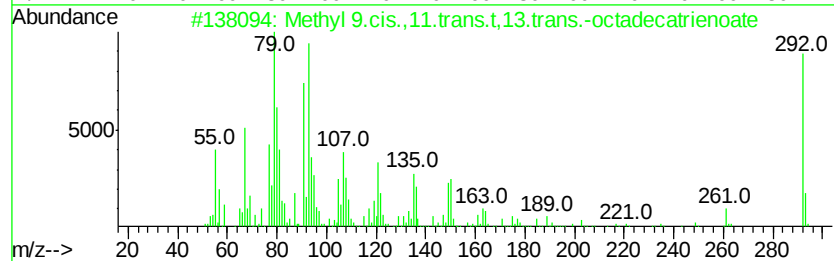
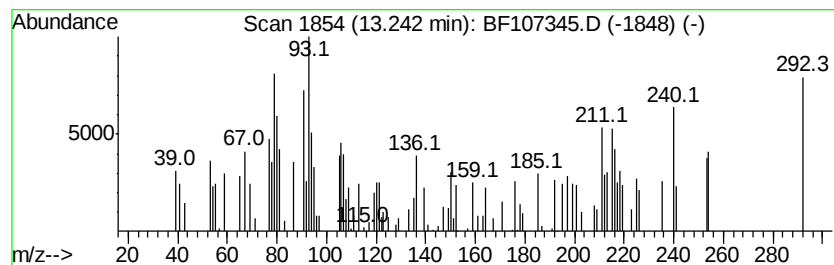
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Methyl 9.cis.,11.trans.t,13... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.24	2.43 ng	56323	Chrysene-d12	14.14

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl 9.cis.,11.trans.t,13.tran...	292	C19H32O2	1000336-42-6	90
2			Methyl 6-cis,9-cis,11-trans-octa...	292	C19H32O2	1000336-37-7	83
3			4,4-Diallyl-cyclohex-2-enone	176	C12H16O	1000186-50-1	55
4			1(2H)-Pentalenone, hexahydro-5-m...	150	C10H14O	120587-85-1	53
5			4-Hexadecen-6-yne, (E)-	220	C16H28	074744-51-7	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071218\
Data File : BF107345.D
Acq On : 12 Jul 2018 18:01
Operator : JU/SJ
Sample : J3828-03 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
EO-01-071018-B

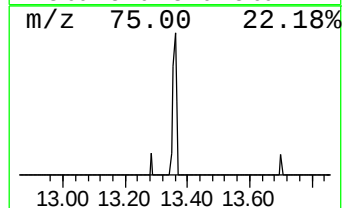
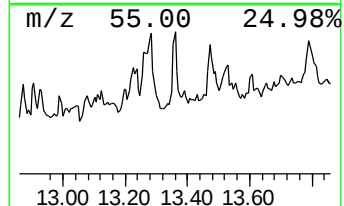
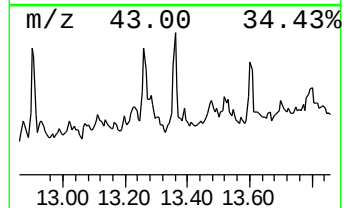
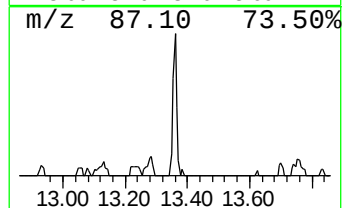
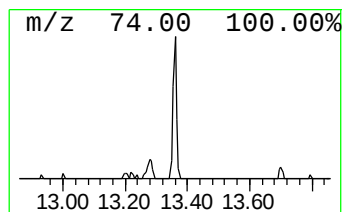
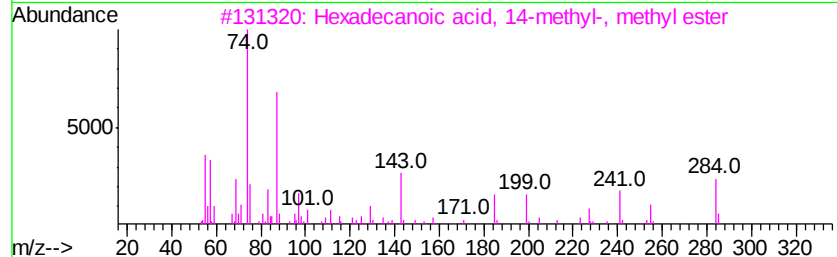
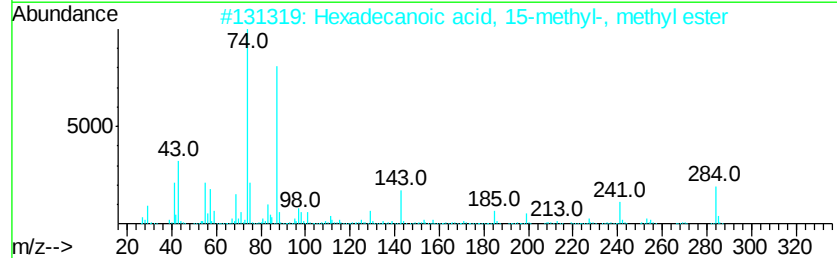
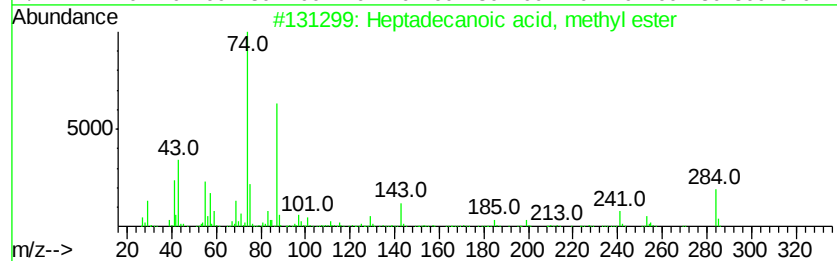
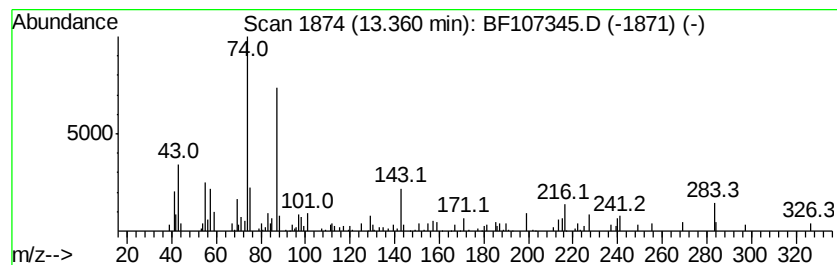
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Heptadecanoic acid, methyl ... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.36	2.79 ng	64764	Chrysene-d12	14.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecanoic acid, methyl ester	284	C18H36O2	001731-92-6	94
2		Hexadecanoic acid, 15-methyl-, methyl ester	284	C18H36O2	006929-04-0	93
3		Hexadecanoic acid, 14-methyl-, methyl ester	284	C18H36O2	002490-49-5	87
4		Eicosanoic acid, methyl ester	326	C21H42O2	001120-28-1	83
5		Decanoic acid, methyl ester	186	C11H22O2	000110-42-9	81



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071218\
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Sample : J3828-03 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
EO-01-071018-B

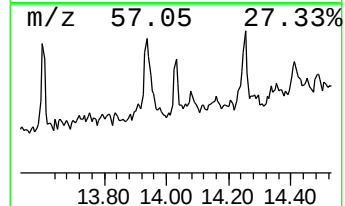
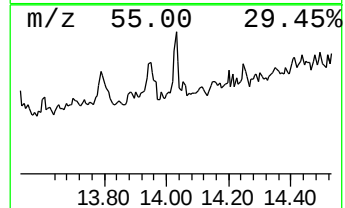
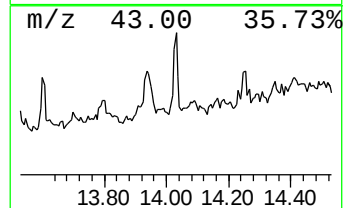
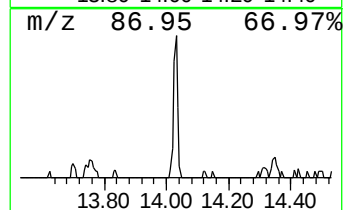
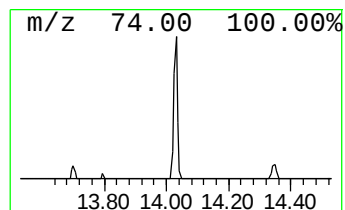
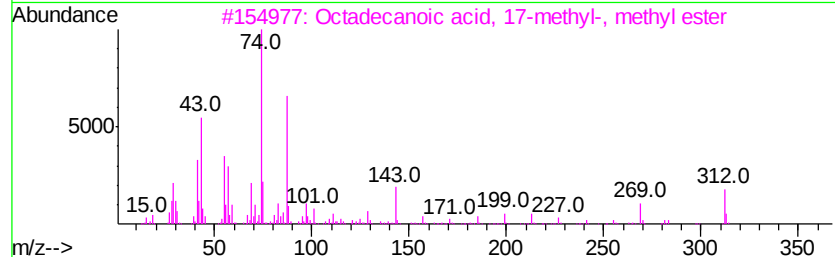
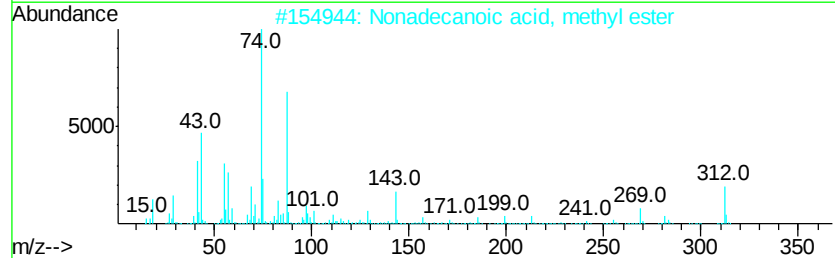
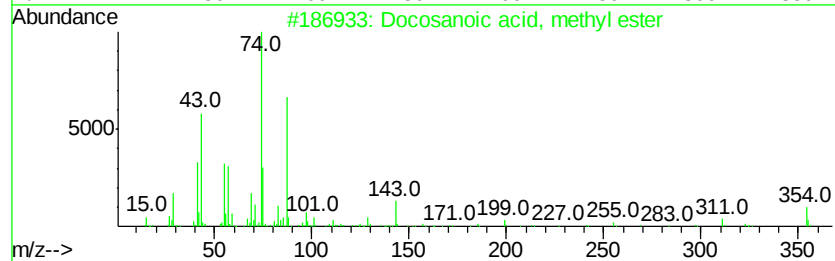
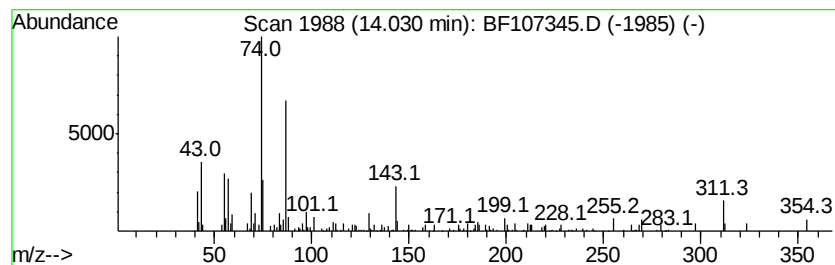
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Docosanoic acid, methyl ester Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.03	2.40 ng	55763	Chrysene-d12	14.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Docosanoic acid, methyl ester	354	C23H46O2	000929-77-1	97
2		Nonadecanoic acid, methyl ester	312	C20H40O2	001731-94-8	90
3		Octadecanoic acid, 17-methyl-, m...	312	C20H40O2	055124-97-5	90
4		Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	81
5		Methyl stearate	298	C19H38O2	000112-61-8	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071218\
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Sample : J3828-03 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

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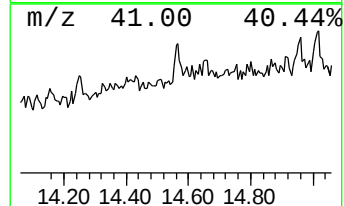
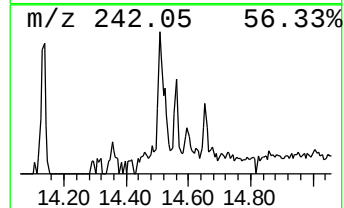
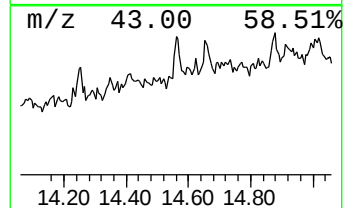
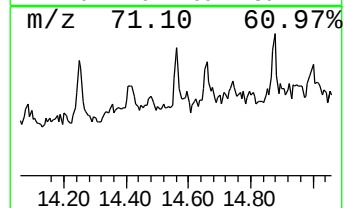
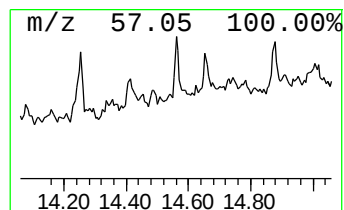
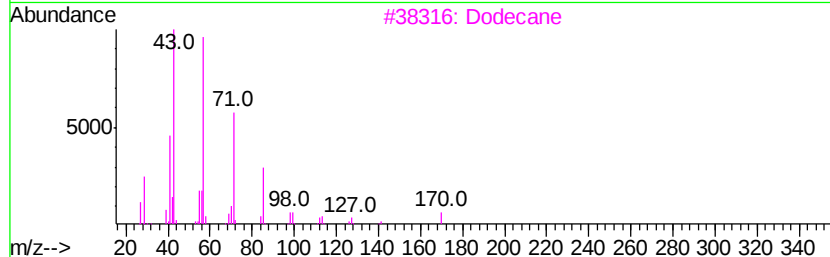
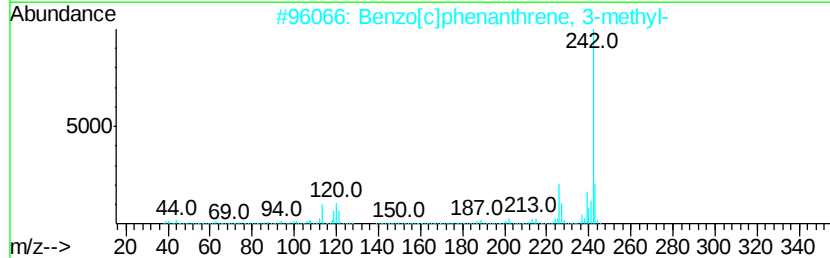
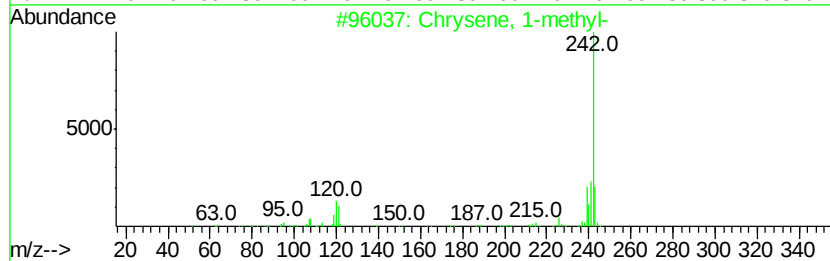
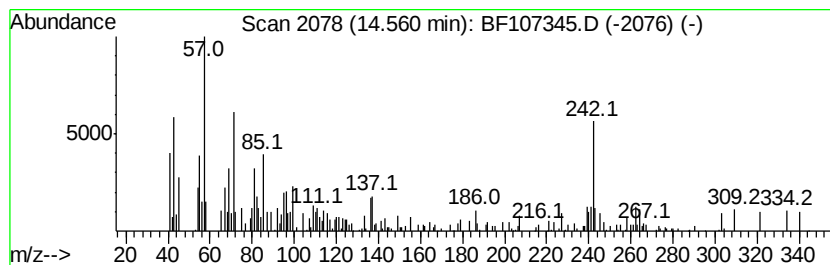
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 unknown14.56 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.56	2.19 ng	50816	Chrysene-d12	14.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Chrysene, 1-methyl-	242	C19H14	003351-28-8	47
2		Benzo[c]phenanthrene, 3-methyl-	242	C19H14	002381-19-3	25
3		Dodecane	170	C12H26	000112-40-3	25
4		Nonanoic acid, 2,2,3,3,4,4,4-hep...	340	C13H19F7O2	1000376-65-8	18
5		Hexadecane	226	C16H34	000544-76-3	15



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF071218\
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 Acq On : 12 Jul 2018 18:01
 Operator : JU/SJ
 Sample : J3828-03 5X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-01-071018-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.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
2-Pentanone, 4-hy...	5.15	2.4	ng	43050	1	6.92	356843	20.0
unknown6.71	6.71	19.0	ng	338839	1	6.92	356843	20.0
Pentadecane	9.88	2.6	ng	61500	3	9.98	463758	20.0
Heptacosane	10.38	3.0	ng	69522	3	9.98	463758	20.0
Sulfurous acid, p...	10.85	4.6	ng	101482	4	11.47	439362	20.0
Tetracosane	11.31	2.8	ng	62358	4	11.47	439362	20.0
Tetratetracontane	11.73	2.3	ng	49868	4	11.47	439362	20.0
Hexadecanoic acid...	11.84	50.2	ng	1102130	4	11.47	439362	20.0
9,12-Octadecadien...	12.54	131.9	ng	2898760	4	11.47	439362	20.0
16-Octadecenoic a...	12.56	84.6	ng	1859460	4	11.47	439362	20.0
Methyl stearate	12.64	23.6	ng	519387	4	11.47	439362	20.0
Methyl 9.cis.,11....	13.24	2.4	ng	56323	5	14.14	464070	20.0
Heptadecanoic aci...	13.36	2.8	ng	64764	5	14.14	464070	20.0
Docosanoic acid, ...	14.03	2.4	ng	55763	5	14.14	464070	20.0
unknown14.56	14.56	2.2	ng	50816	5	14.14	464070	20.0