

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112118\
 Data File : BF110929.D
 Acq On : 21 Nov 2018 15:37
 Operator : JU/SJ
 Sample : J6000-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FS-12072-11929

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-BF110818.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	2.240	22	26	36	rVB	80451	114723	6.62%	0.753%
2	5.181	523	526	543	rBV	30315	54674	3.15%	0.359%
3	5.563	587	591	612	rBV	1323352	1394920	80.49%	9.158%
4	6.569	758	762	782	rBV	1414719	1500521	86.59%	9.851%
5	6.710	782	786	801	rVB	1741320	1513833	87.35%	9.939%
6	6.939	822	825	832	rBV	451837	396636	22.89%	2.604%
7	7.098	848	852	865	rBV	1213403	1165439	67.25%	7.651%
8	7.510	918	922	939	rBV	846216	934245	53.91%	6.134%
9	8.222	1040	1043	1060	rBV	432483	522645	30.16%	3.431%
10	9.298	1222	1226	1235	rBV	1723803	1732982	100.00%	11.378%
11	9.692	1287	1293	1298	rBV	343534	393986	22.73%	2.587%
12	9.833	1314	1317	1319	rBV2	31186	27968	1.61%	0.184%
13	9.880	1322	1325	1328	rVB	46244	38724	2.23%	0.254%
14	9.980	1339	1342	1350	rBV2	583678	573081	33.07%	3.762%
15	10.304	1392	1397	1399	rBV5	20536	27930	1.61%	0.183%
16	10.380	1407	1410	1413	rVB2	51577	45480	2.62%	0.299%
17	10.610	1445	1449	1451	rBV4	16719	22112	1.28%	0.145%
18	10.780	1474	1478	1490	rBV	819599	961598	55.49%	6.313%
19	10.874	1492	1494	1497	rVB2	28985	28213	1.63%	0.185%
20	11.480	1593	1597	1605	rBV	516788	526046	30.35%	3.454%
21	11.980	1679	1682	1687	rBV3	67532	91088	5.26%	0.598%
22	12.698	1801	1804	1811	rVB3	79726	103796	5.99%	0.681%
23	13.063	1862	1866	1869	rBV	1315757	1200536	69.28%	7.882%
24	13.604	1955	1958	1960	rBV2	17931	17472	1.01%	0.115%
25	13.933	2011	2014	2019	rBV2	137374	163486	9.43%	1.073%
26	14.133	2044	2048	2057	rVB	428777	437393	25.24%	2.872%
27	14.251	2065	2068	2073	rVB	30179	34773	2.01%	0.228%
28	14.551	2116	2119	2123	rBV	55336	62425	3.60%	0.410%
29	14.868	2169	2173	2177	rBV	85823	87542	5.05%	0.575%
30	15.215	2228	2232	2244	rVB	134195	186313	10.75%	1.223%
31	15.609	2294	2299	2304	rBV	130892	183353	10.58%	1.204%
32	15.668	2304	2309	2319	rVB	235064	347411	20.05%	2.281%
33	16.062	2371	2376	2381	rVV	111099	170819	9.86%	1.121%
34	16.133	2385	2388	2396	rVB9	13623	30304	1.75%	0.199%

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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	16.592	2460	2466	2471	rBV	53161	100543	5.80%	0.660%
36	17.209	2566	2571	2578	rVB4	20568	38619	2.23%	0.254%

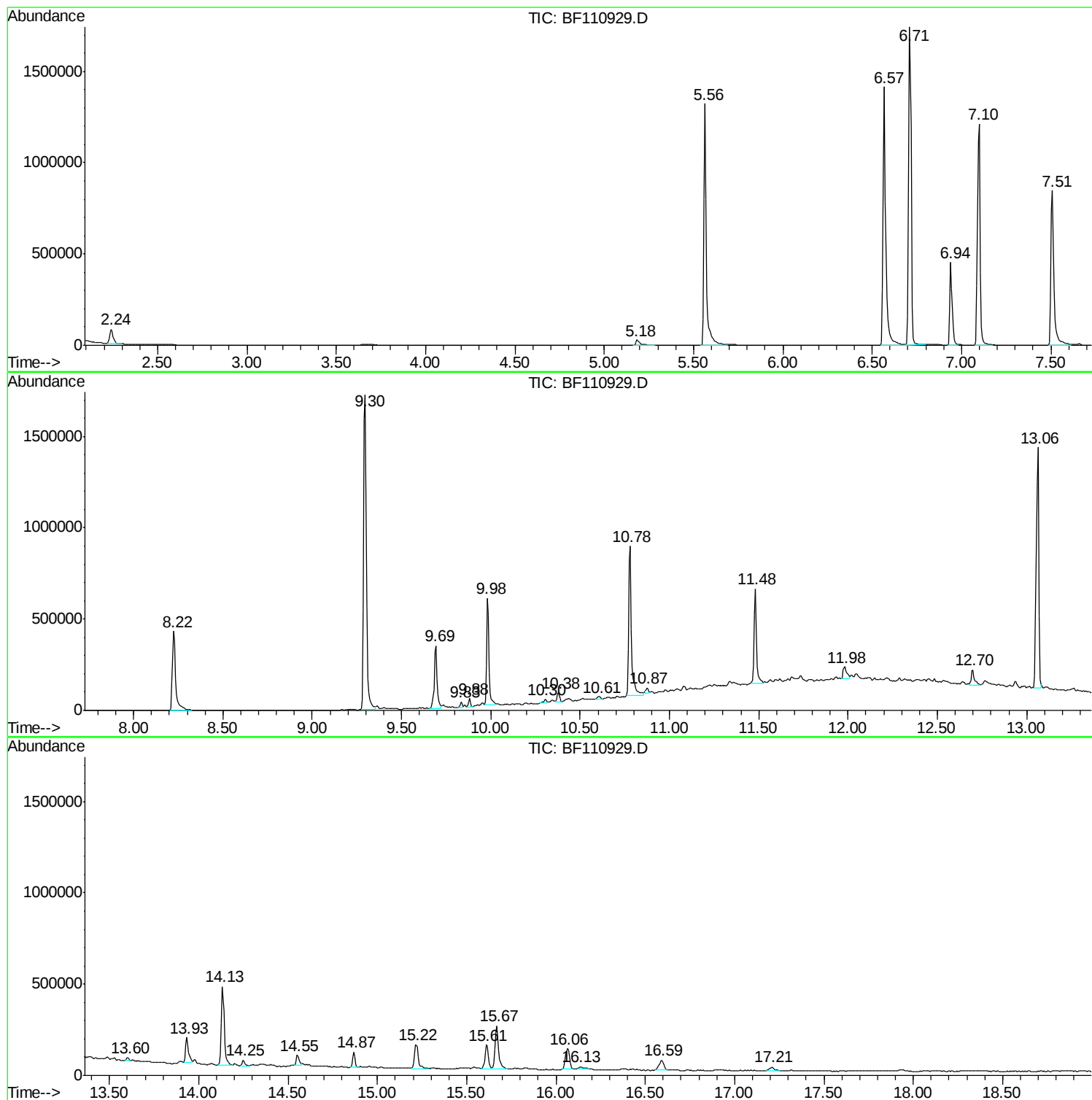
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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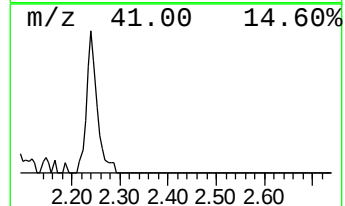
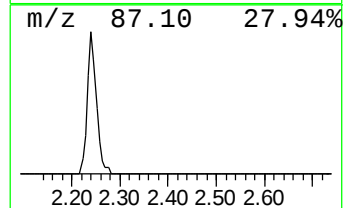
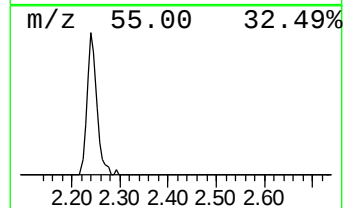
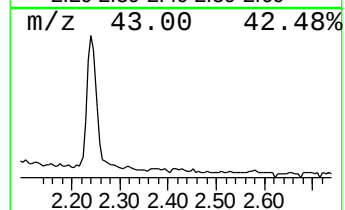
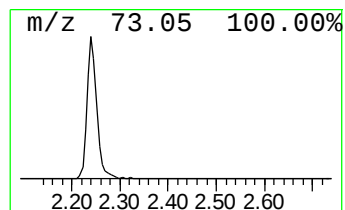
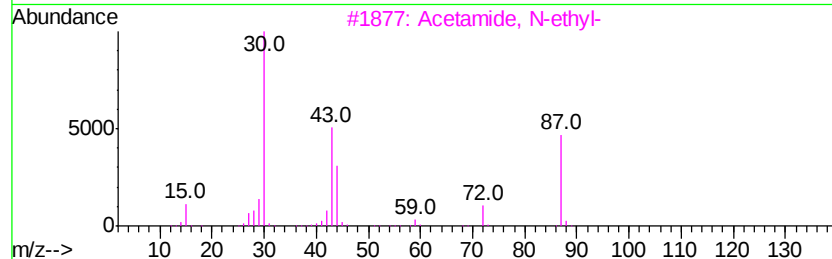
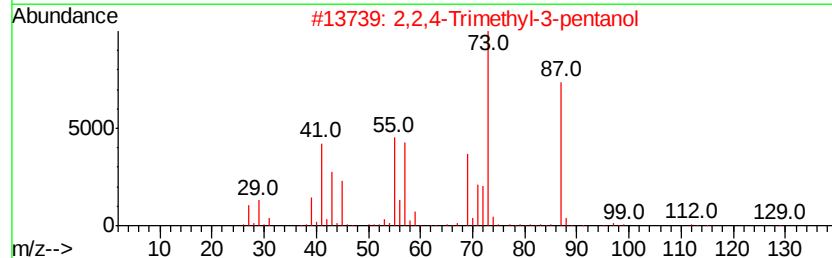
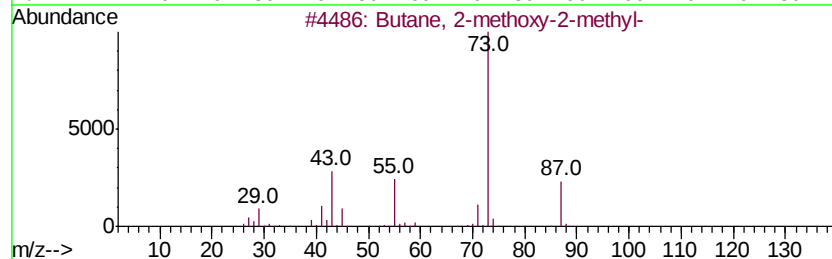
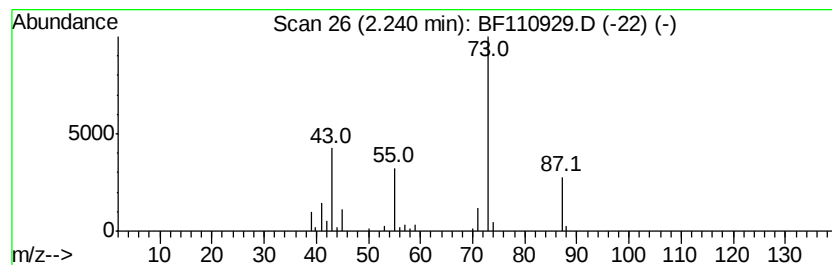
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.24	5.78 ng	114723	1,4-Dichlorobenzene-d4	6.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	38
3		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	25
4		2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	25
5		N-Ethylformamide	73	C3H7NO	000627-45-2	9



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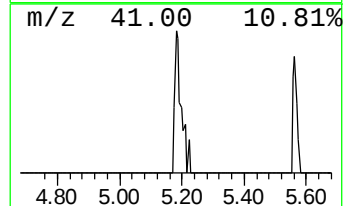
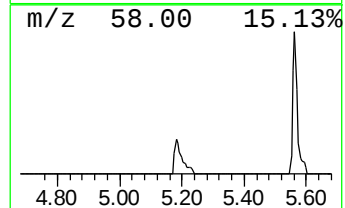
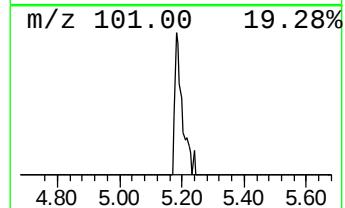
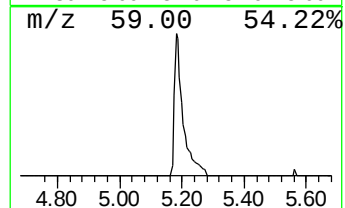
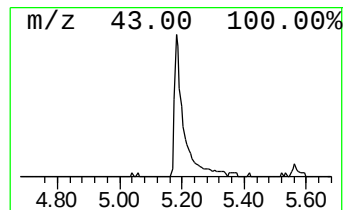
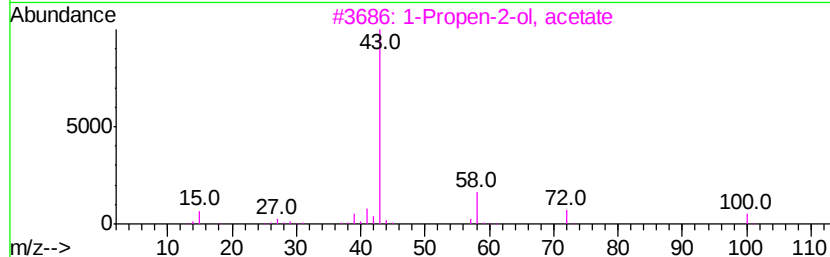
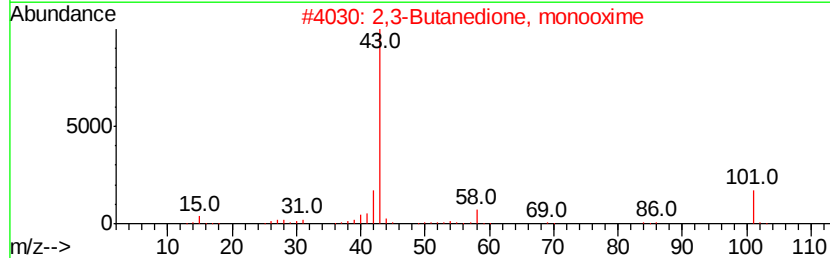
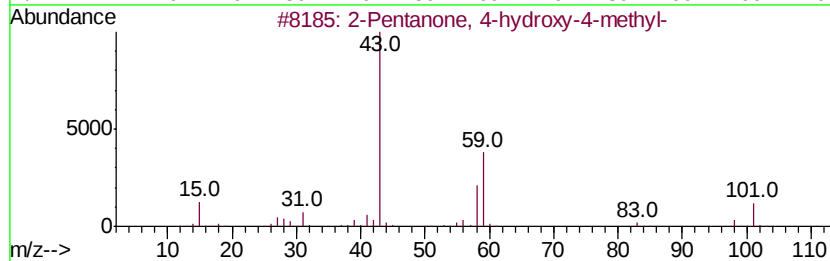
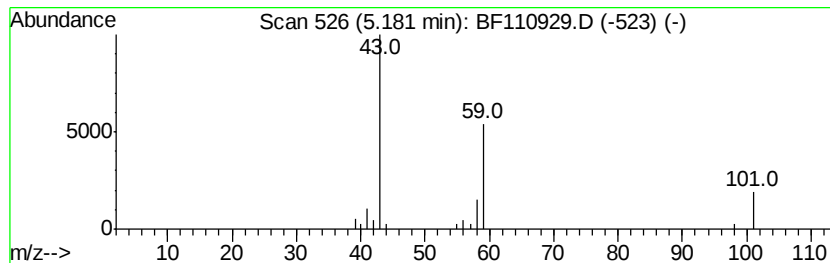
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.18	2.76 ng	54674	1,4-Dichlorobenzene-d4	6.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
3		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9
4		Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	9
5		Butane	58	C4H10	000106-97-8	7



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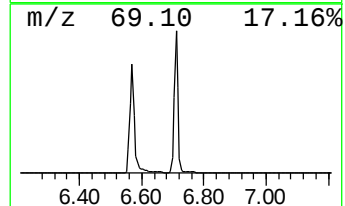
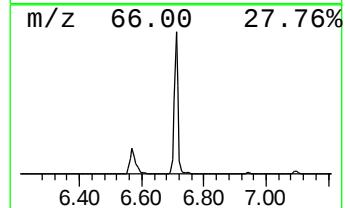
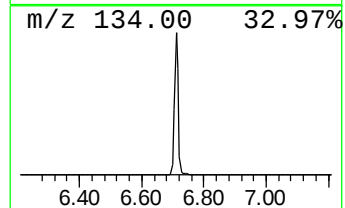
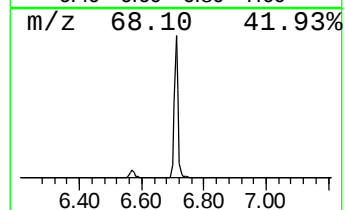
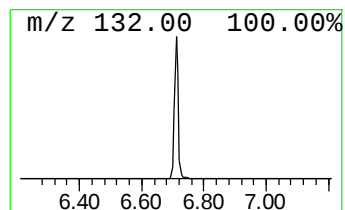
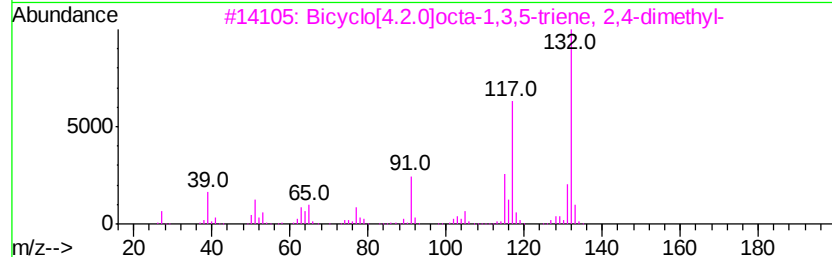
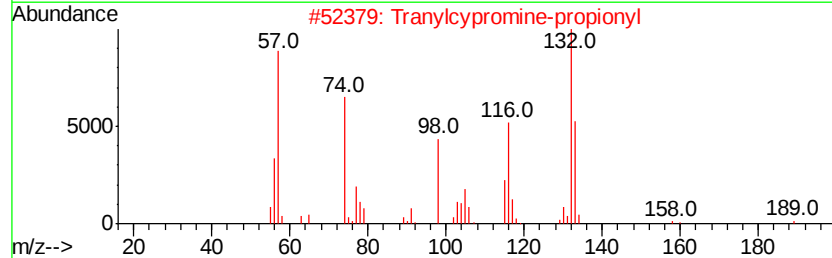
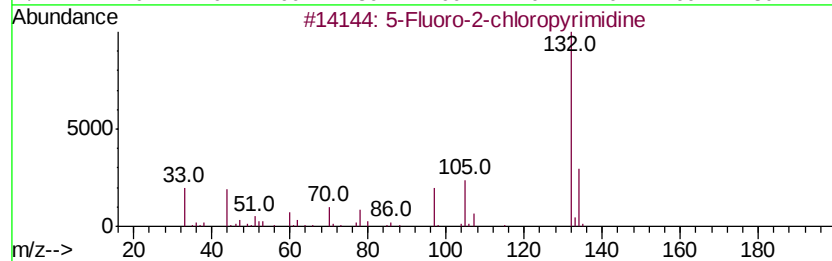
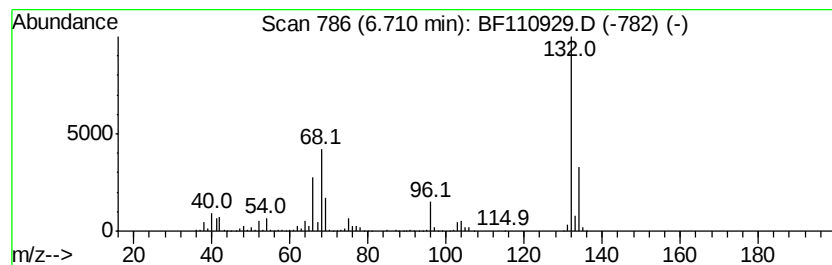
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.71 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.71	76.33 ng	1513830	1,4-Dichlorobenzene-d4	6.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
5		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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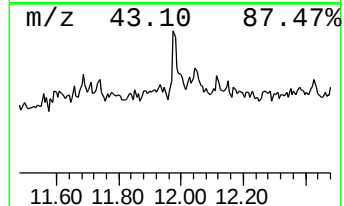
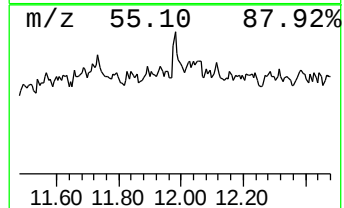
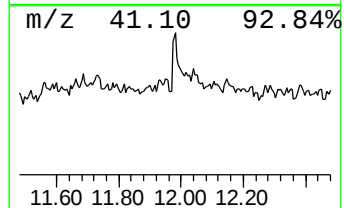
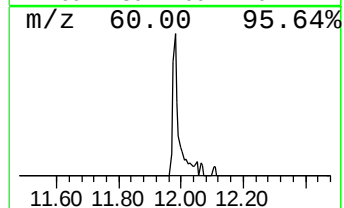
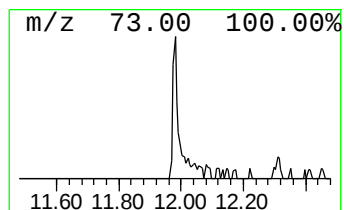
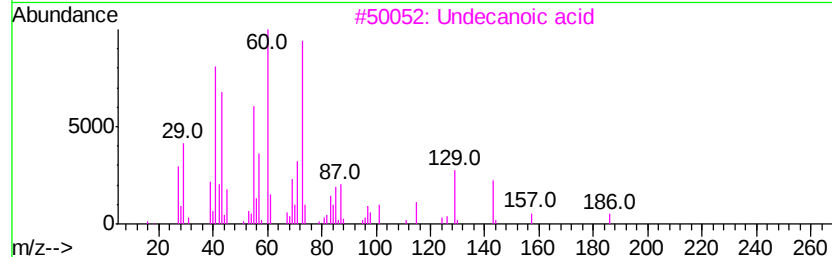
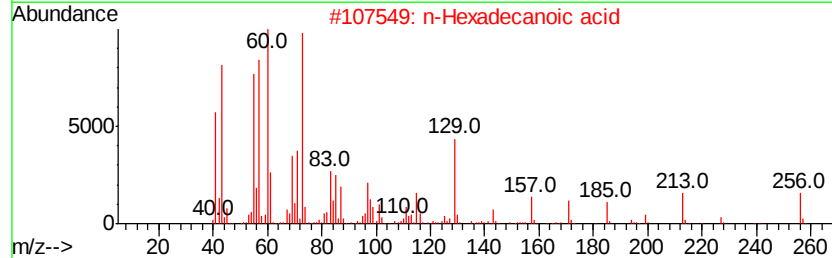
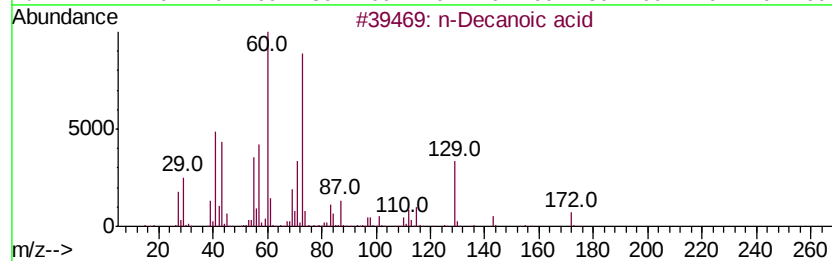
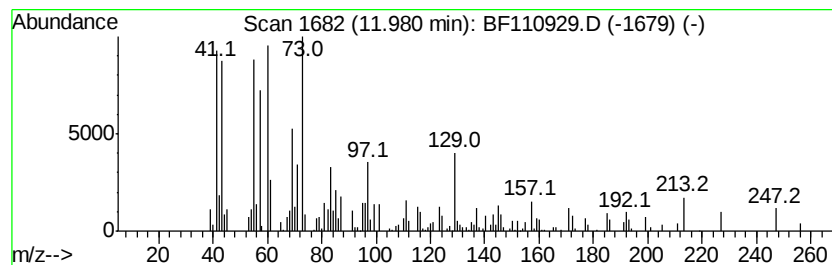
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 n-Decanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.98	3.46 ng	91088	Phenanthrene-d10	11.48

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Decanoic acid	172	C10H20O2	000334-48-5	83
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	81
3		Undecanoic acid	186	C11H22O2	000112-37-8	64
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	64
5		Tetradecanoic acid	228	C14H28O2	000544-63-8	58



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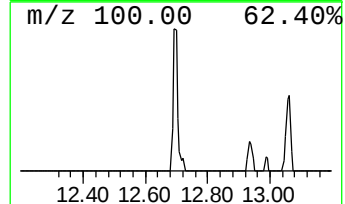
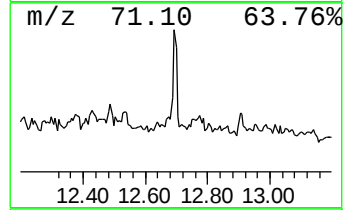
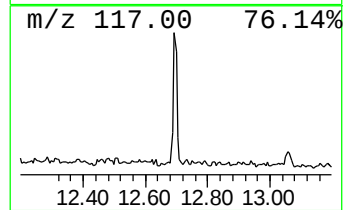
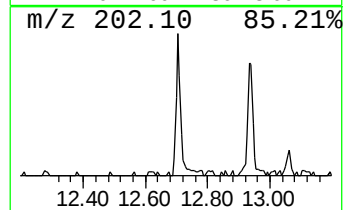
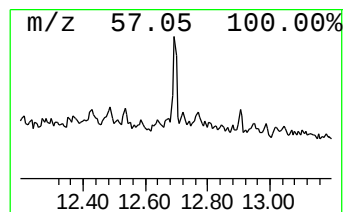
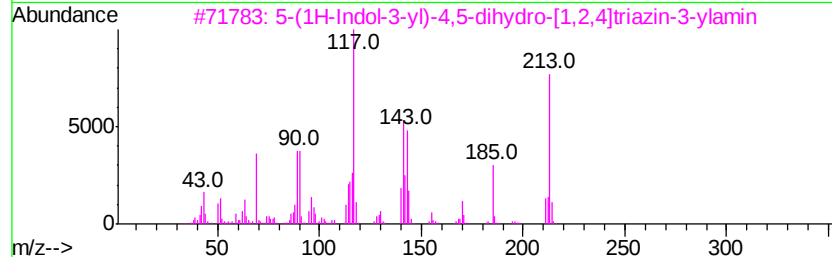
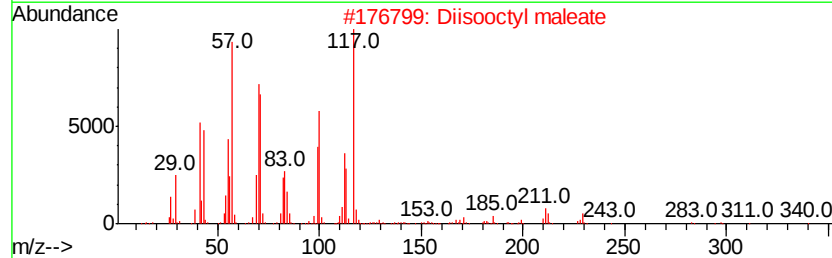
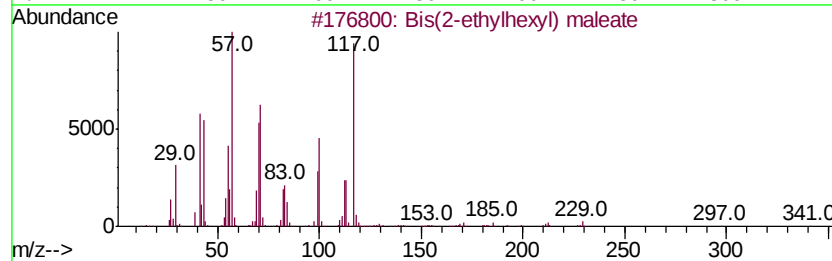
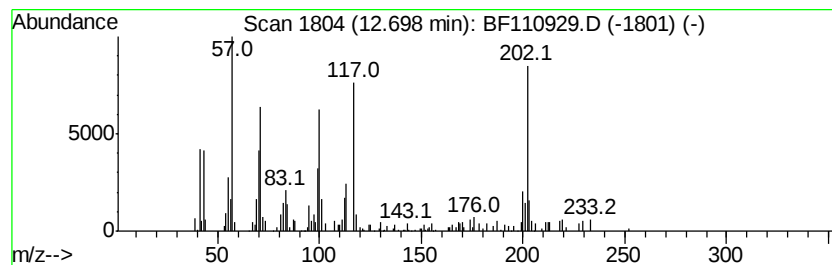
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Peak Number 6 unknown12.70 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.70	3.95 ng	103796	Phenanthrene-d10	11.48

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bis(2-ethylhexyl) maleate	340	C20H36O4	000142-16-5	49
2			Diisooctyl maleate	340	C20H36O4	001330-76-3	27
3			5-(1H-Indol-3-yl)-4,5-dihydro-[1...	213	C11H11N5	1000300-20-6	15
4			Pvrene	202	C16H10	000129-00-0	11
5			Oxalic acid, 2-ethylhexyl tetrad...	398	C24H46O4	1000309-39-7	11



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112118\
Data File : BF110929.D
Acq On : 21 Nov 2018 15:37
Operator : JU/SJ
Sample : J6000-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
FS-12072-11929

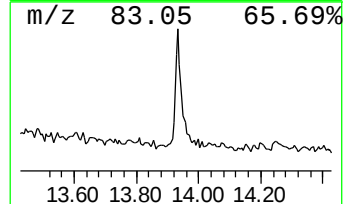
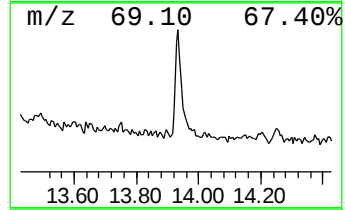
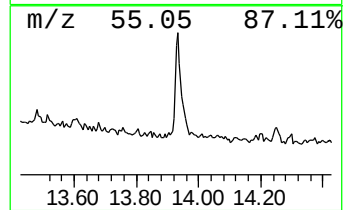
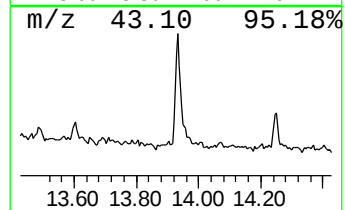
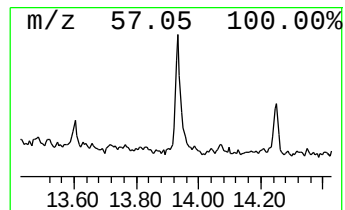
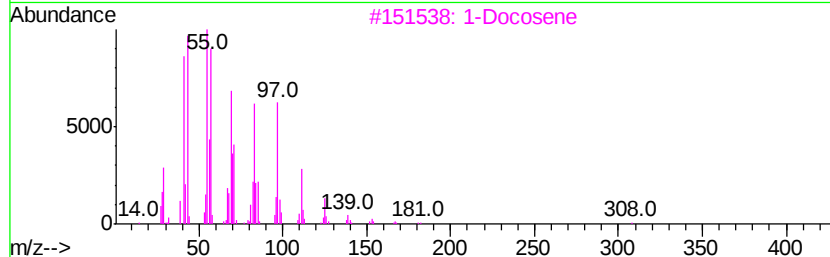
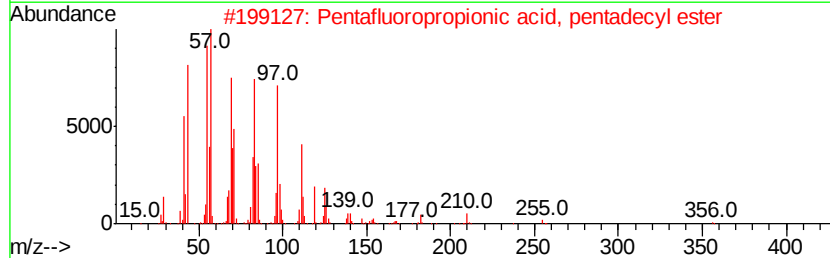
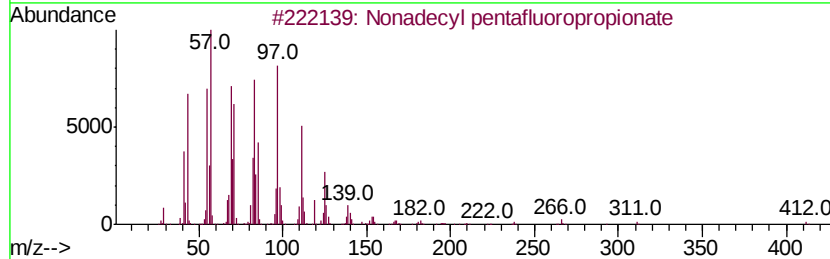
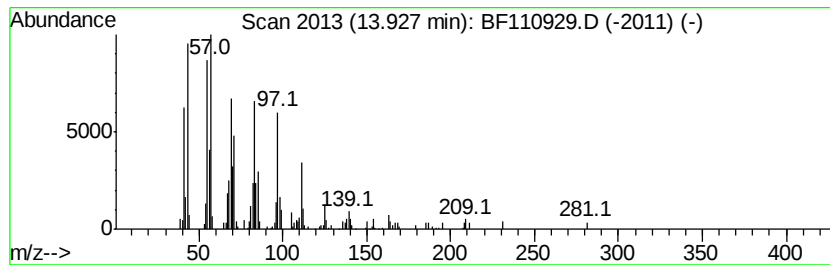
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Nonadecyl pentafluoropropio... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.93	7.48 ng	163486	Chrysene-d12	14.13

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	90
2			Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	90
3			1-Docosene	308	C22H44	001599-67-3	90
4			Heptadecyl trifluoroacetate	352	C19H35F3O2	1000351-87-0	87
5			Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112118\
Data File : BF110929.D
Acq On : 21 Nov 2018 15:37
Operator : JU/SJ
Sample : J6000-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

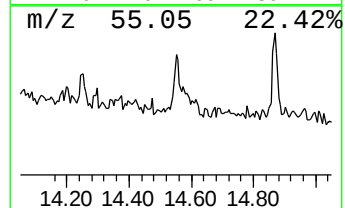
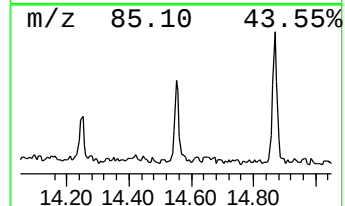
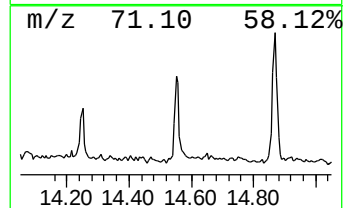
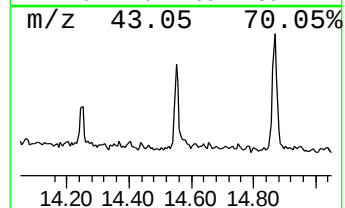
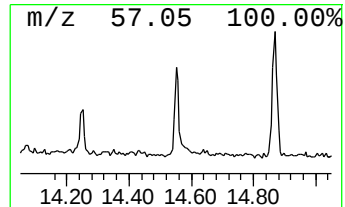
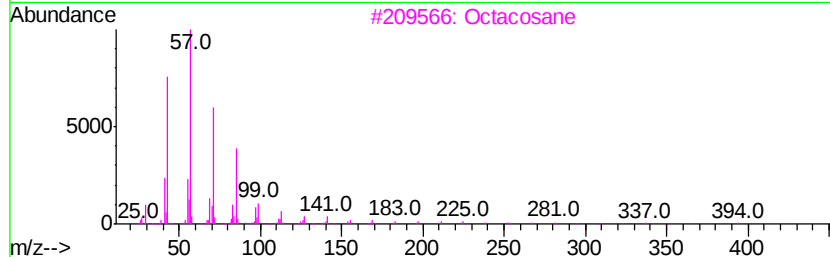
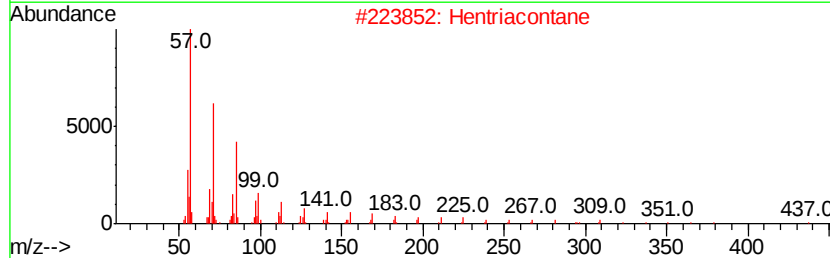
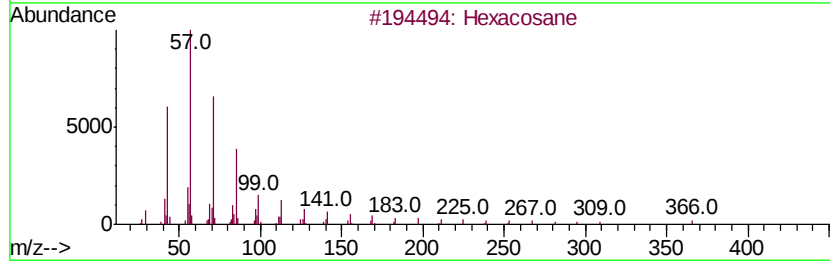
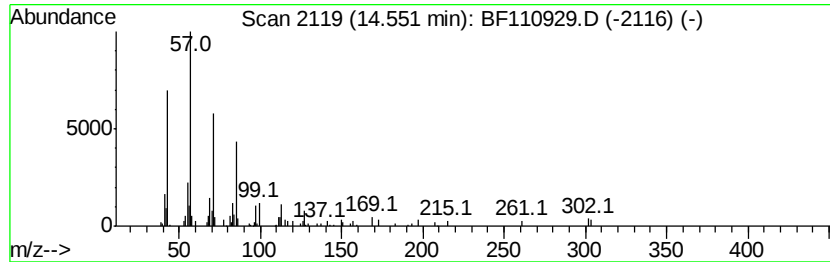
Instrument :
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ClientSampleId :
FS-12072-11929

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Hexacosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.
14.55	2.85 ng	62425	Chrysene-d12		14.13
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexacosane	366	C26H54	000630-01-3	87
2	Hentriacontane	437	C31H64	000630-04-6	87
3	Octacosane	394	C28H58	000630-02-4	87
4	Dodecane, 1-iodo-	296	C12H25I	004292-19-7	83
5	Heptacosane	380	C27H56	000593-49-7	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112118\
 Data File : BF110929.D
 Acq On : 21 Nov 2018 15:37
 Operator : JU/SJ
 Sample : J6000-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

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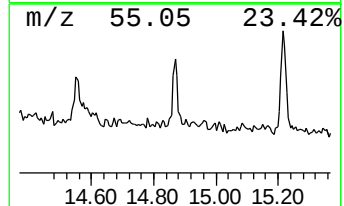
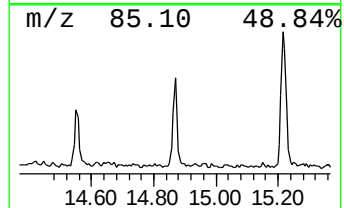
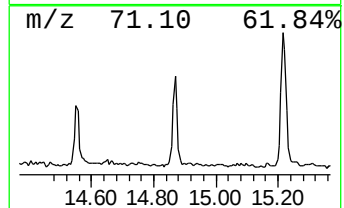
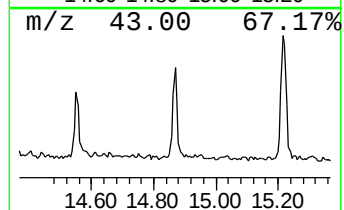
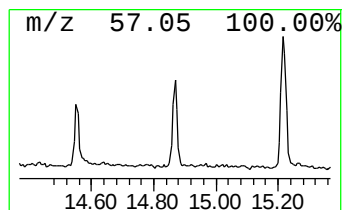
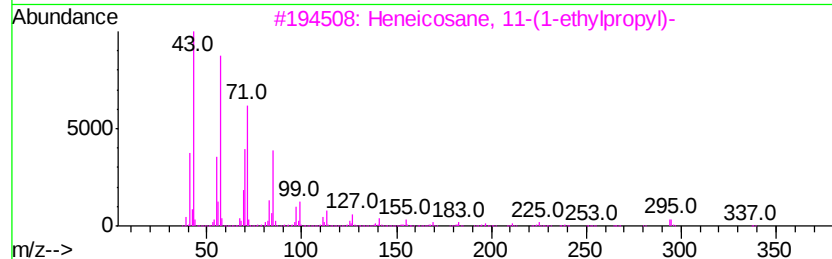
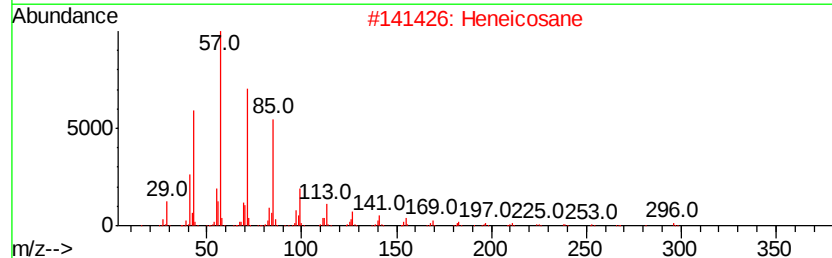
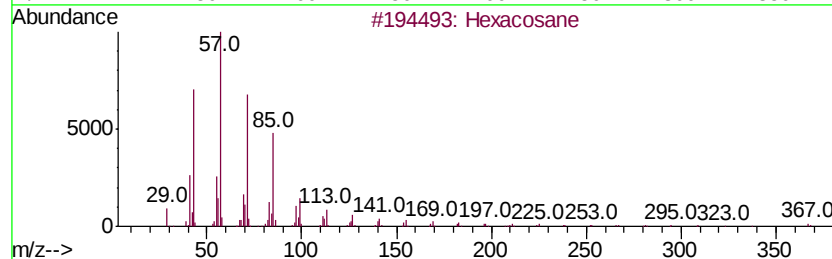
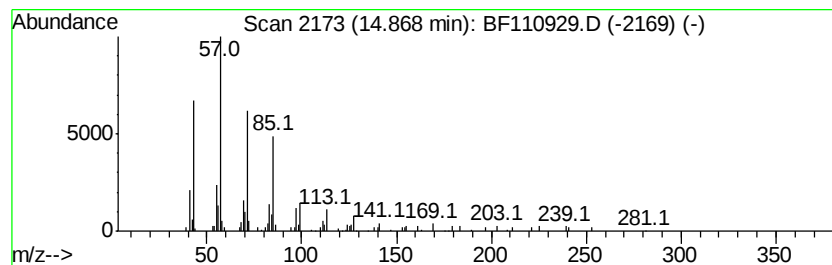
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 9 Heneicosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.87	4.00 ng	87542	Chrysene-d12	14.13

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexacosane		366	C26H54	000630-01-3	91
2	Heneicosane		296	C21H44	000629-94-7	91
3	Heneicosane, 11-(1-ethylpropyl)-		366	C26H54	055282-11-6	91
4	2-methyloctacosane		408	C29H60	1000376-72-8	91
5	Tetratetracontane		619	C44H90	007098-22-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112118\
Data File : BF110929.D
Acq On : 21 Nov 2018 15:37
Operator : JU/SJ
Sample : J6000-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
FS-12072-11929

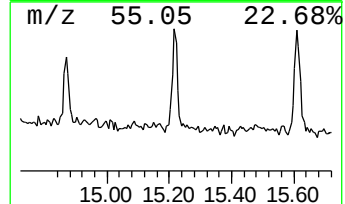
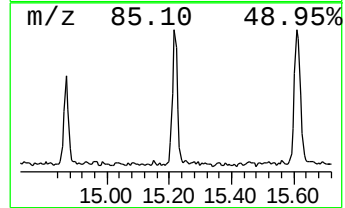
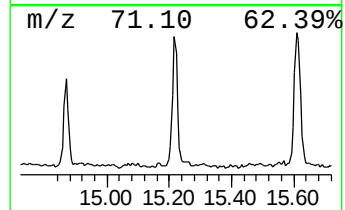
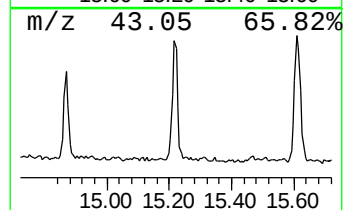
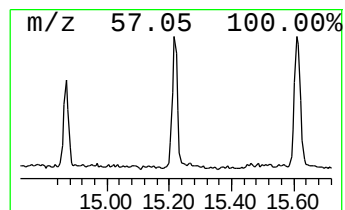
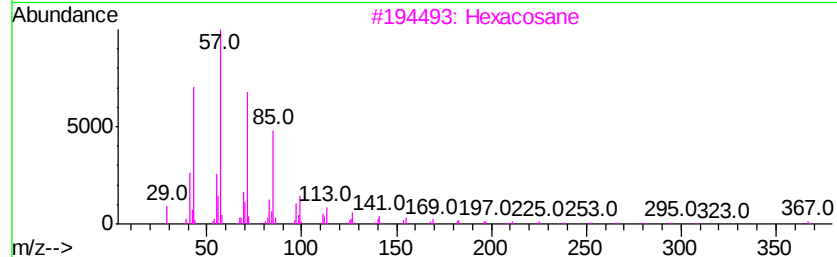
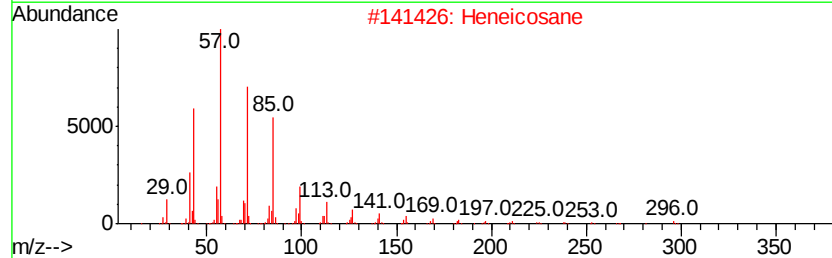
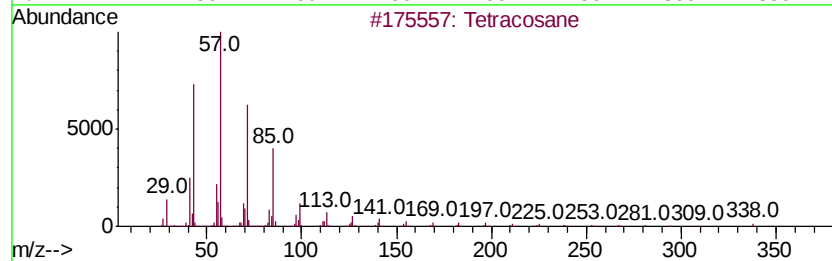
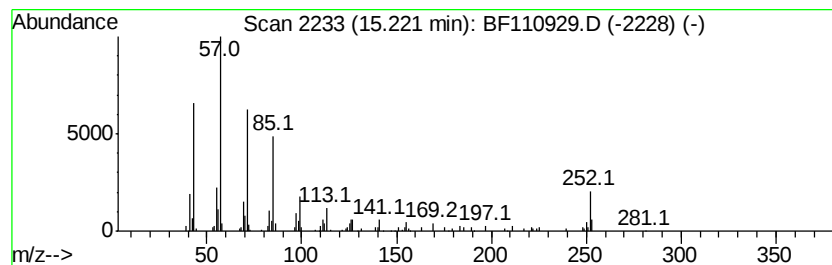
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Tetracosane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.22	10.73 ng	186313	Perylene-d12	15.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane	338	C24H50	000646-31-1	81
2		Heneicosane	296	C21H44	000629-94-7	81
3		Hexacosane	366	C26H54	000630-01-3	81
4		Docosane	310	C22H46	000629-97-0	81
5		Triacontane	422	C30H62	000638-68-6	81



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112118\
Data File : BF110929.D
Acq On : 21 Nov 2018 15:37
Operator : JU/SJ
Sample : J6000-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

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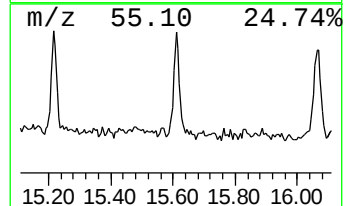
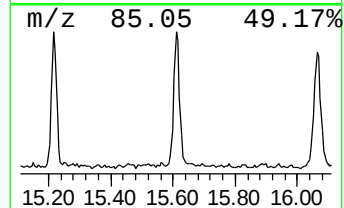
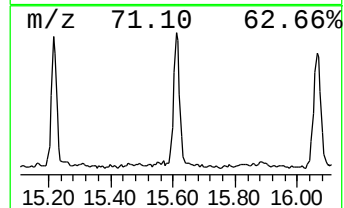
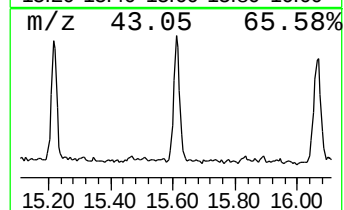
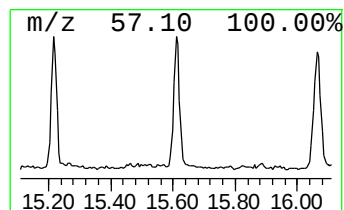
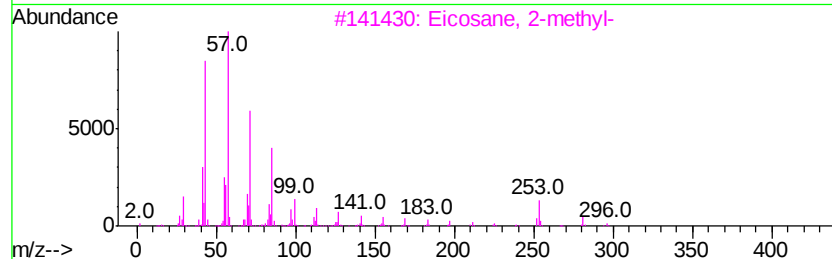
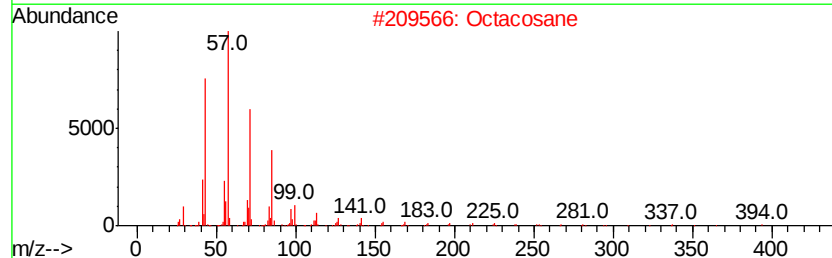
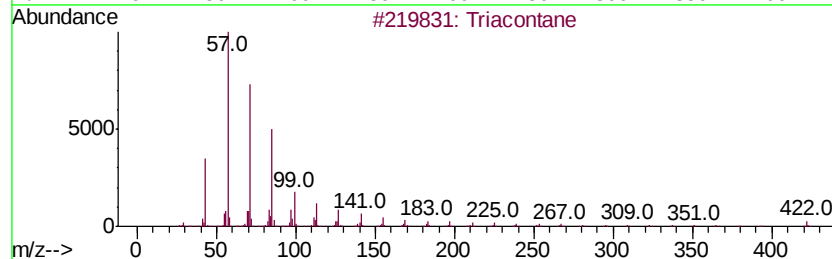
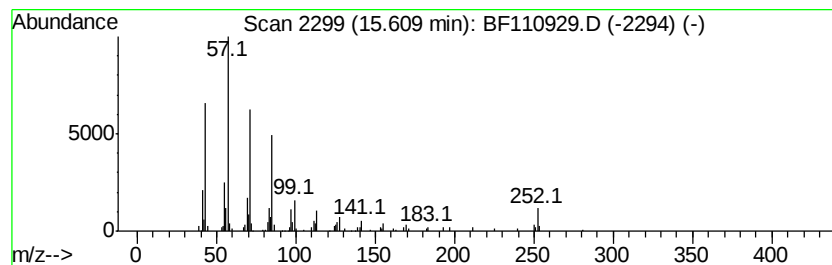
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Triacontane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.61	10.56 ng	183353	Perylene-d12	15.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Triacontane	422	C30H62	000638-68-6	90
2		Octacosane	394	C28H58	000630-02-4	90
3		Eicosane, 2-methyl-	296	C21H44	001560-84-5	87
4		Heneicosane	296	C21H44	000629-94-7	87
5		Docosane	310	C22H46	000629-97-0	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112118\
Data File : BF110929.D
Acq On : 21 Nov 2018 15:37
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Misc :
ALS Vial : 11 Sample Multiplier: 1

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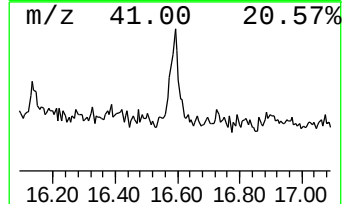
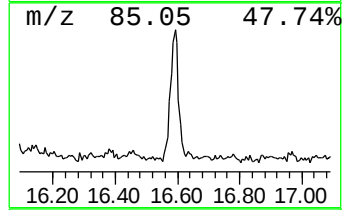
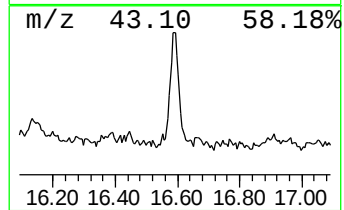
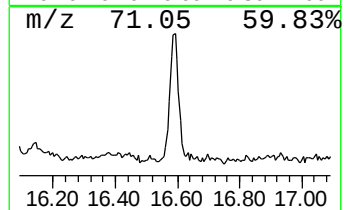
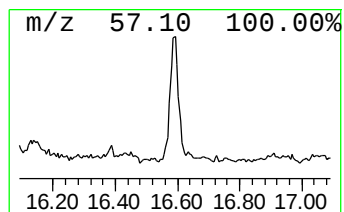
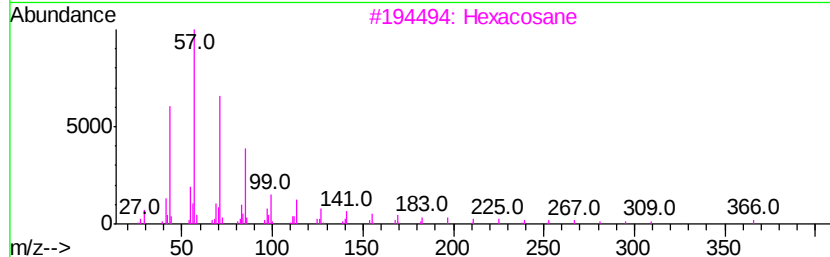
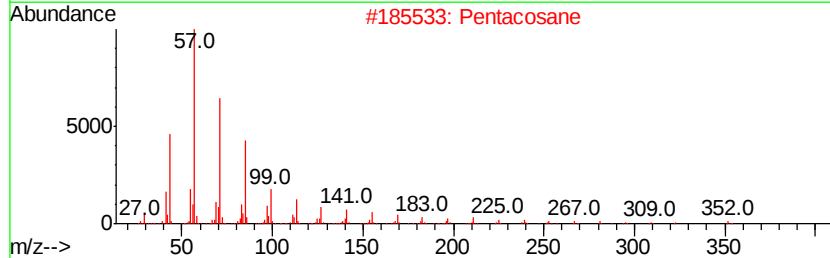
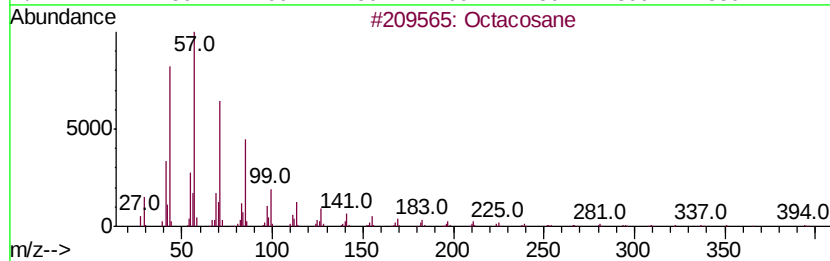
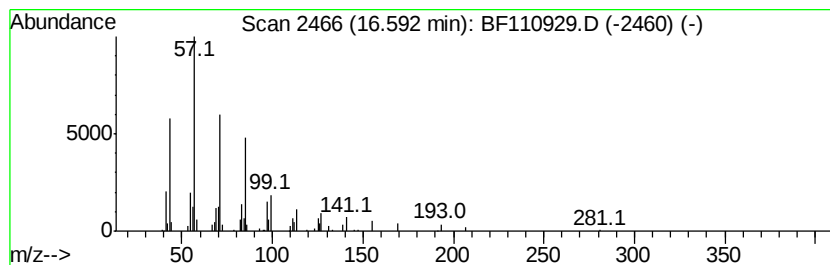
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Octacosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.59	5.79 ng	100543	Perylene-d12	15.67

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosane	394	C28H58	000630-02-4	91
2			Pentacosane	352	C25H52	000629-99-2	91
3			Hexacosane	366	C26H54	000630-01-3	91
4			Tetratetracontane	619	C44H90	007098-22-8	91
5			2-methyloctacosane	408	C29H60	1000376-72-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF112118\
Data File : BF110929.D
Acq On : 21 Nov 2018 15:37
Operator : JU/SJ
Sample : J6000-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110818.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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2-Pentanone, 4-hy...	5.18	2.8	ng	54674	1	6.94	396636	20.0
unknown6.71	6.71	76.3	ng	1513830	1	6.94	396636	20.0
n-Decanoic acid	11.98	3.5	ng	91088	4	11.48	526046	20.0
unknown12.70	12.70	4.0	ng	103796	4	11.48	526046	20.0
Nonadecyl pentafl...	13.93	7.5	ng	163486	5	14.13	437393	20.0
Hexacosane	14.55	2.9	ng	62425	5	14.13	437393	20.0
Heneicosane	14.87	4.0	ng	87542	5	14.13	437393	20.0
Tetracosane	15.22	10.7	ng	186313	6	15.67	347411	20.0
triacontane	15.61	10.6	ng	183353	6	15.67	347411	20.0
Octacosane	16.59	5.8	ng	100543	6	15.67	347411	20.0