

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040820\
 Data File : BF119862.D
 Acq On : 9 Apr 2020 9:00
 Operator : MA/SJ
 Sample : L2151-11
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CBH-140

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-BF040820.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	4.963	485	489	497	rVB	109740	128713	1.66%	0.257%
2	5.375	554	559	562	rBV	2910835	2497974	32.20%	4.988%
3	6.393	728	732	743	rBV	1773048	1537883	19.83%	3.071%
4	6.763	791	795	799	rBV	1431666	1180972	15.22%	2.358%
5	6.922	817	822	826	rBV	4888352	4957861	63.91%	9.900%
6	7.334	886	892	895	rBV	3894575	4242361	54.69%	8.472%
7	8.051	1009	1014	1017	rBV	1735154	1567737	20.21%	3.131%
8	9.134	1192	1198	1201	rBV	7322648	7757276	100.00%	15.491%
9	9.522	1260	1264	1267	rBV	554790	435270	5.61%	0.869%
10	9.810	1306	1313	1316	rBV	1960663	1825624	23.53%	3.646%
11	10.604	1441	1448	1451	rBV	5064072	5419352	69.86%	10.822%
12	11.292	1561	1565	1569	rBV	2147311	1889383	24.36%	3.773%
13	11.833	1652	1657	1667	rBV	1536927	1496099	19.29%	2.988%
14	12.522	1771	1774	1775	rBV2	114032	116765	1.51%	0.233%
15	12.557	1778	1780	1786	rVV	131340	122364	1.58%	0.244%
16	12.616	1787	1790	1802	rVV	867960	902144	11.63%	1.801%
17	12.892	1831	1837	1840	rBV	6692605	6955877	89.67%	13.890%
18	13.122	1872	1876	1879	rBV	110203	115482	1.49%	0.231%
19	13.274	1899	1902	1905	rVB	133131	105088	1.35%	0.210%
20	13.339	1910	1913	1915	rBV4	56641	81199	1.05%	0.162%
21	13.369	1915	1918	1921	rVB	286822	268330	3.46%	0.536%
22	13.463	1931	1934	1936	rBV	172675	122943	1.58%	0.246%
23	13.792	1987	1990	1999	rVB	721653	822705	10.61%	1.643%
24	13.933	2010	2014	2019	rBV	1568946	1396552	18.00%	2.789%
25	14.110	2040	2044	2045	rBV	240865	227902	2.94%	0.455%
26	14.133	2045	2048	2055	rVB	495104	647815	8.35%	1.294%
27	14.416	2092	2096	2098	rBV	285721	277611	3.58%	0.554%
28	14.521	2111	2114	2118	rBV	429162	371906	4.79%	0.743%
29	14.716	2143	2147	2150	rBV	310382	304694	3.93%	0.608%
30	14.786	2157	2159	2162	rVB	97923	81190	1.05%	0.162%
31	15.039	2199	2202	2208	rVB	267579	323643	4.17%	0.646%
32	15.374	2254	2259	2262	rBV	901120	1089484	14.04%	2.176%
33	15.410	2262	2265	2270	rVB	196745	245504	3.16%	0.490%
34	15.715	2314	2317	2325	rVB	94936	143887	1.85%	0.287%

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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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35	15.833	2333	2337	2342	rVB	131497	186904	2.41%	0.373%
36	16.057	2372	2375	2384	rVB6	64393	123473	1.59%	0.247%
37	16.321	2415	2420	2424	rBV2	60214	107624	1.39%	0.215%

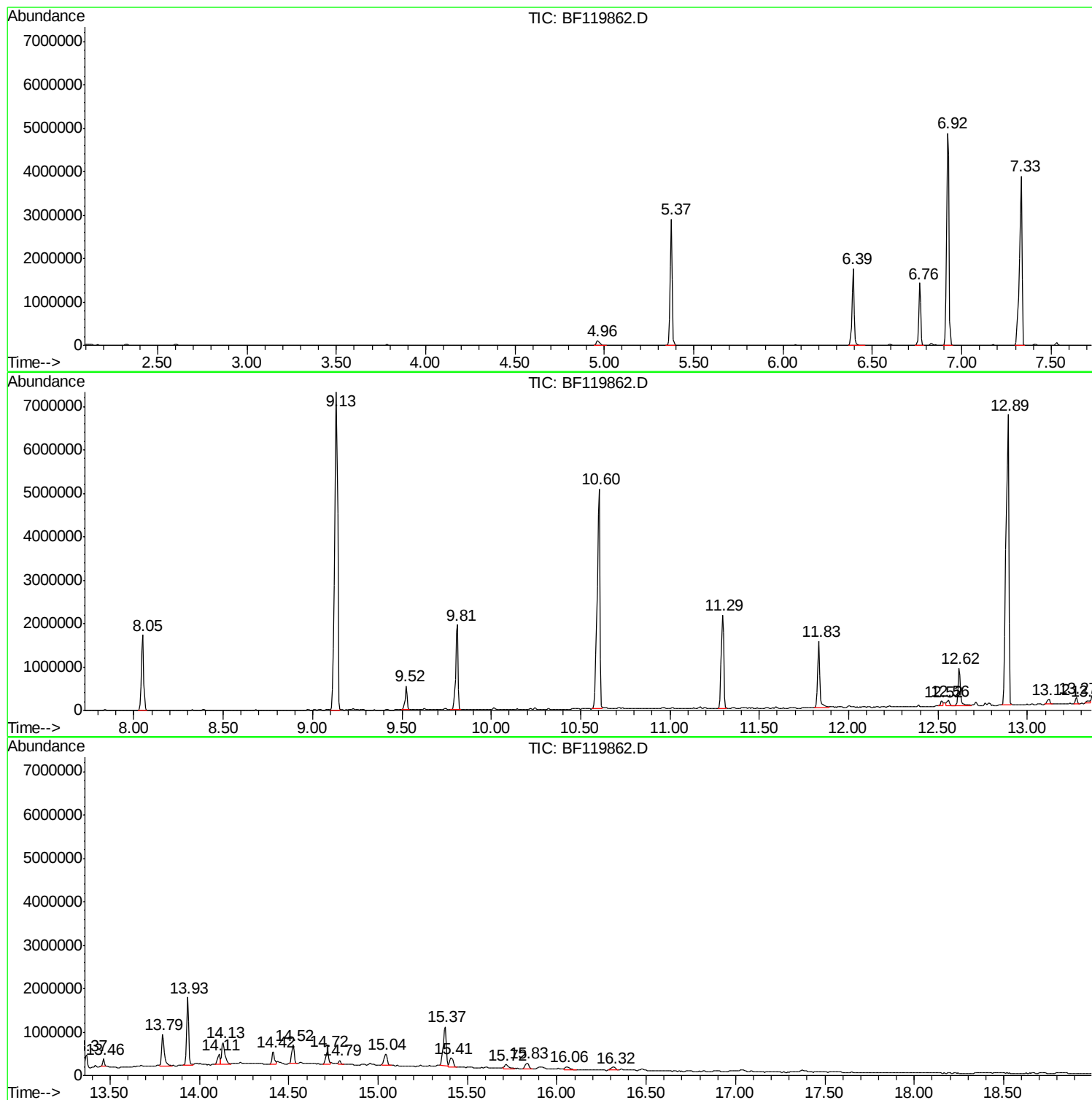
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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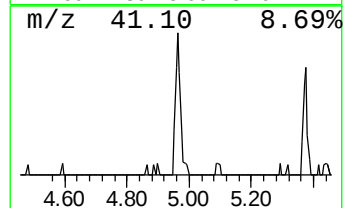
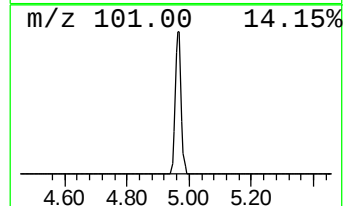
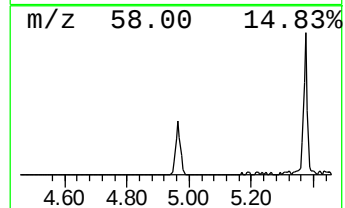
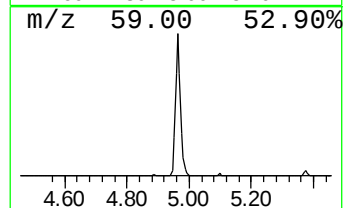
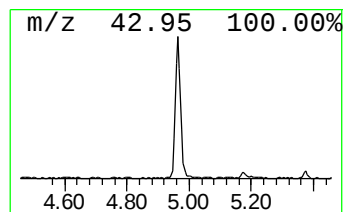
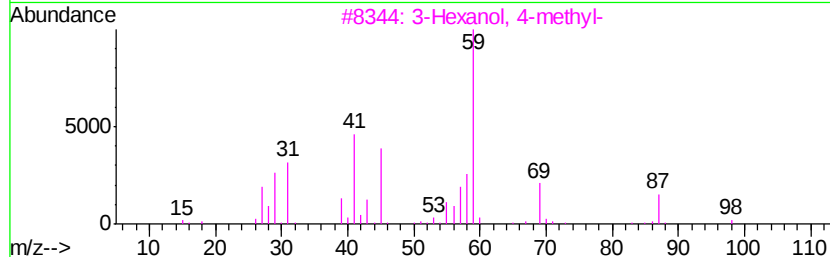
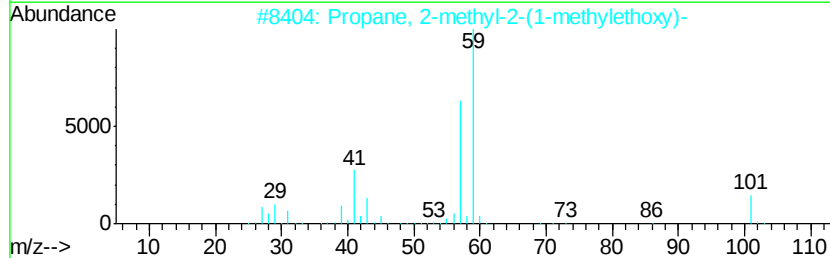
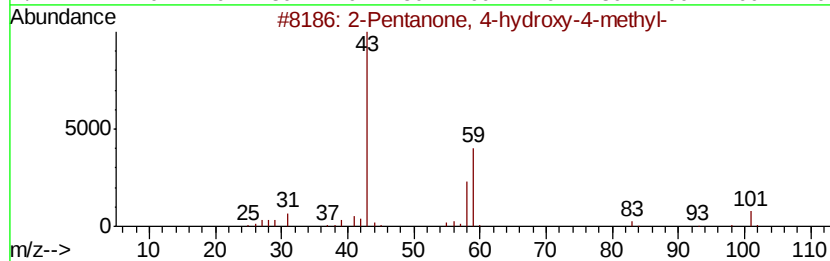
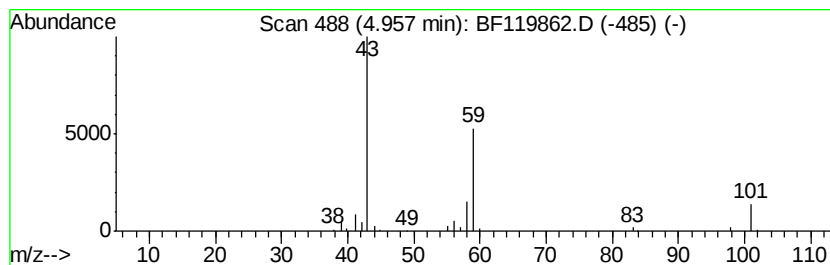
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.	
4.96	2.18 ng	128713	1,4-Dichlorobenzene-d4	6.76	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2	Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	42
3	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4	3-Octanol	130	C8H18O	000589-98-0	33
5	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28



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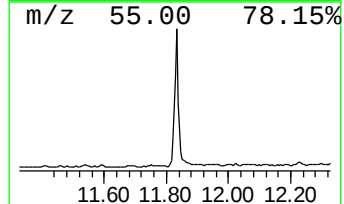
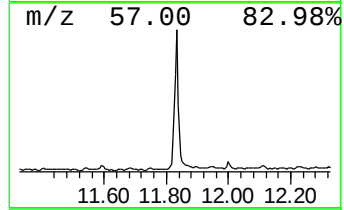
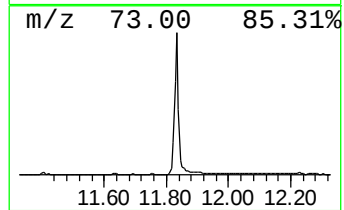
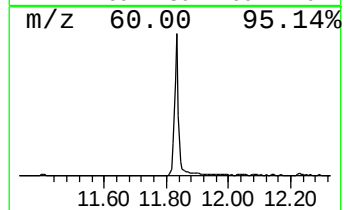
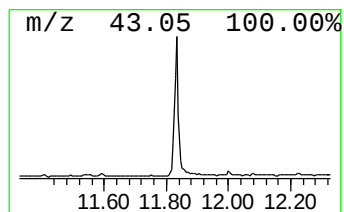
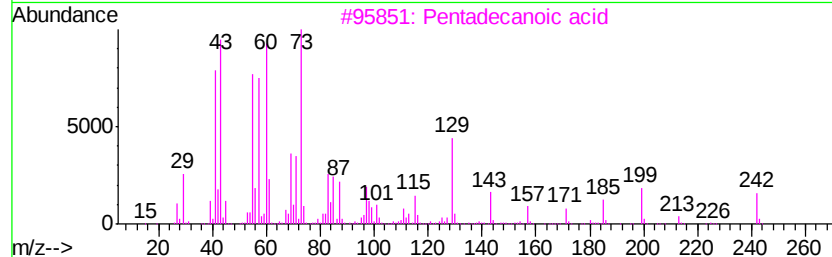
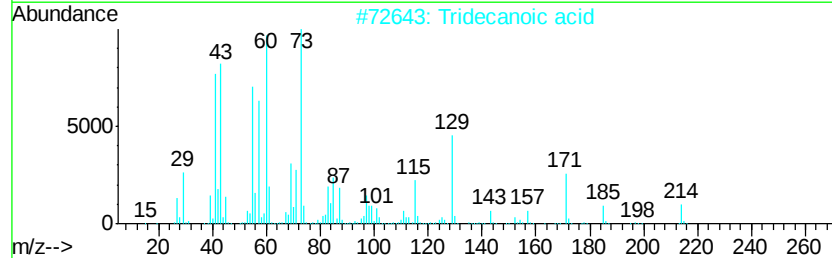
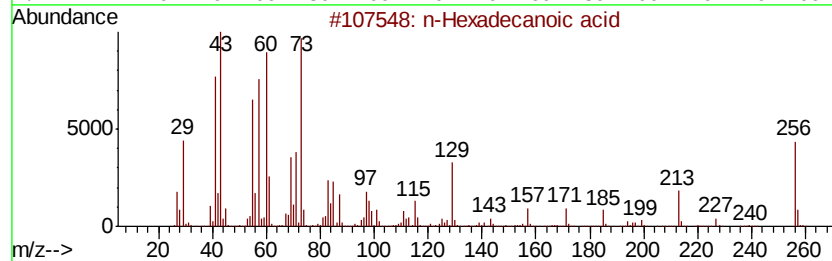
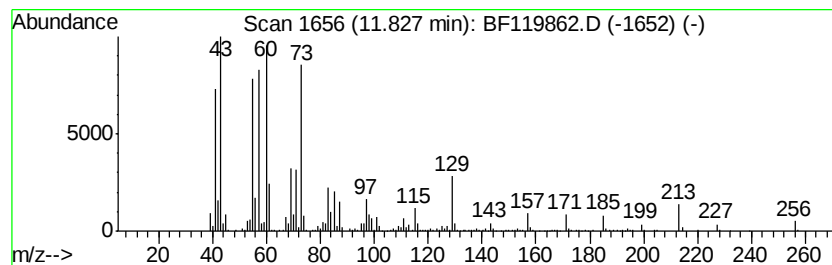
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.83	15.84 ng	1496100	Phenanthrene-d10	11.29

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94
2	Tridecanoic acid	214	C13H26O2	000638-53-9	90
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	90
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	86
5	Dodecanoic acid	200	C12H24O2	000143-07-7	50



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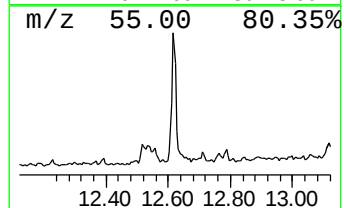
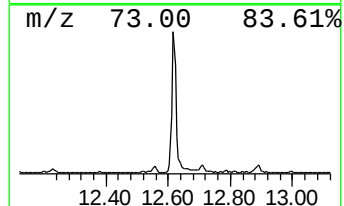
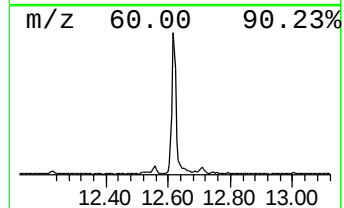
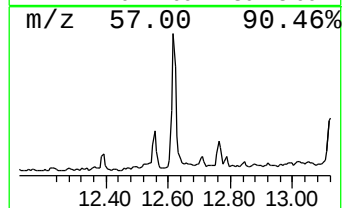
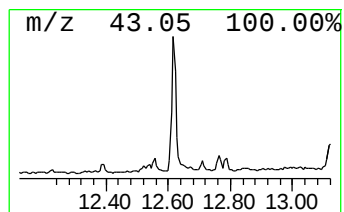
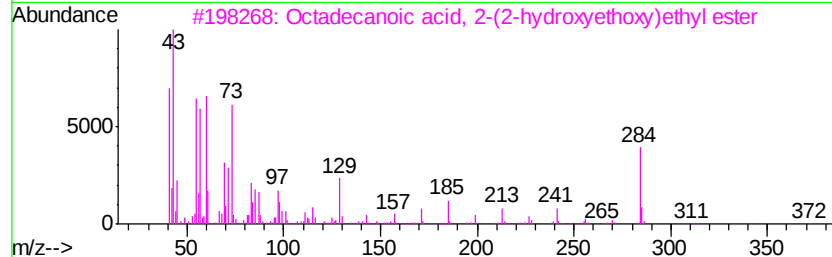
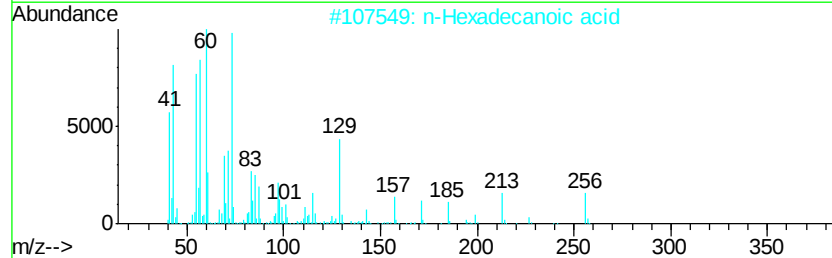
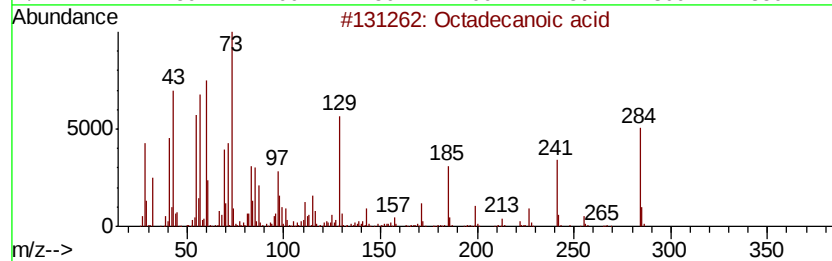
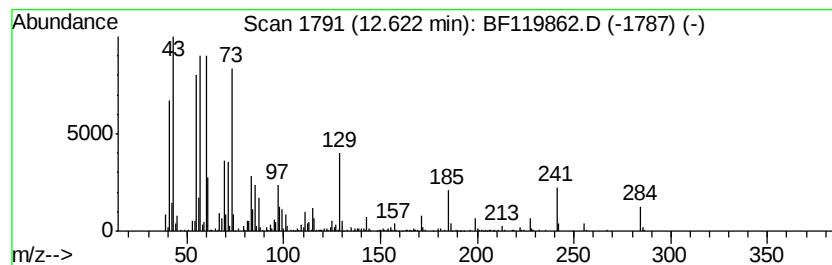
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Octadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.62	12.92 ng	902144	Chrysene-d12	13.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91
3		Octadecanoic acid, 2-(2-hydroxyve...	372	C22H44O4	000106-11-6	87
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	64
5		n-Decanoic acid	172	C10H20O2	000334-48-5	53



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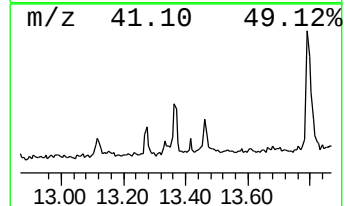
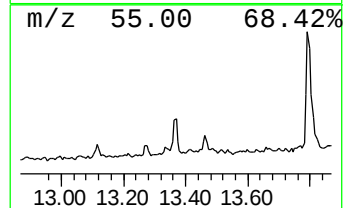
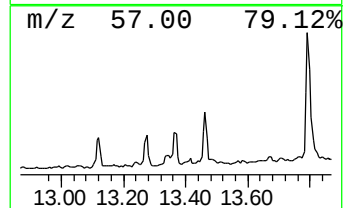
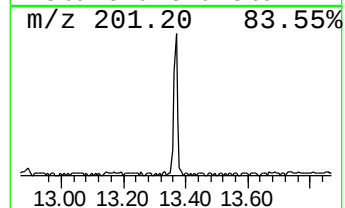
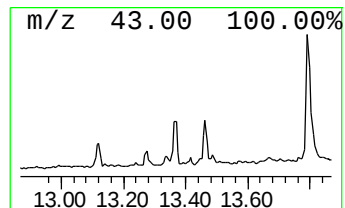
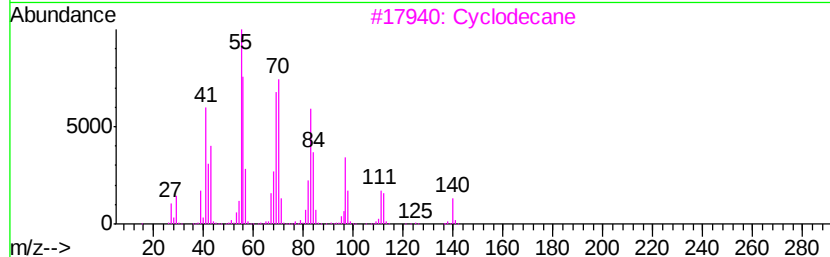
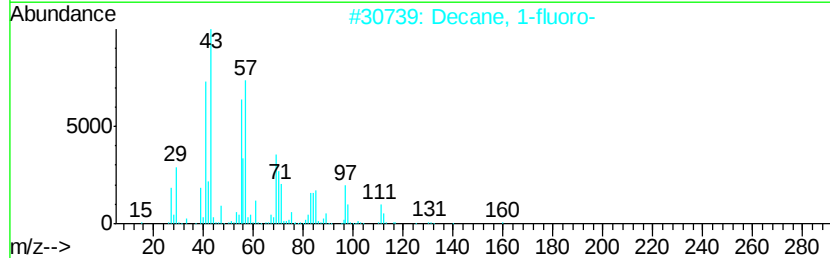
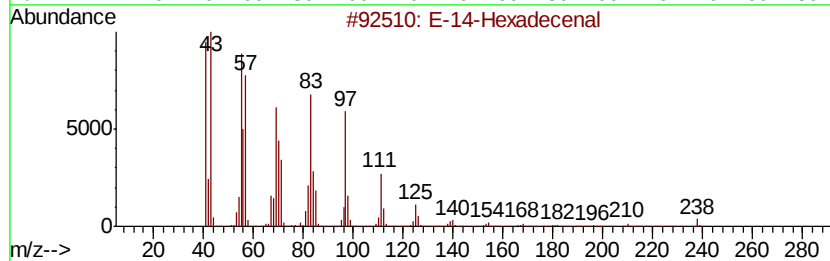
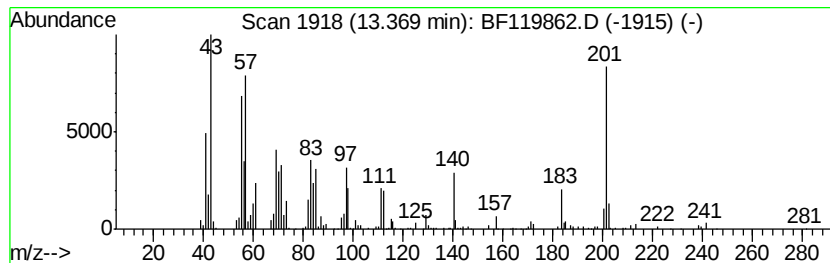
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 unknown13.37 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.37	3.84 ng	268330	Chrysene-d12	13.93

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		E-14-Hexadecenal	238	C16H30O	330207-53-9	44
2		Decane, 1-fluoro-	160	C10H21F	000334-56-5	35
3		Cyclodecane	140	C10H20	000293-96-9	27
4		1-Hexene-3,5-dione	112	C6H8O2	052204-69-0	18
5		1-Heptadecene	238	C17H34	006765-39-5	15



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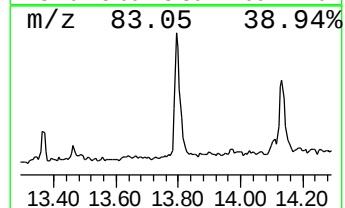
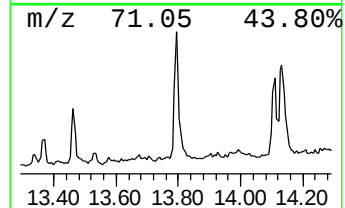
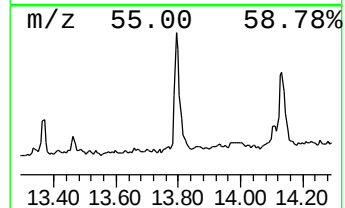
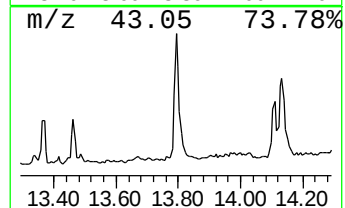
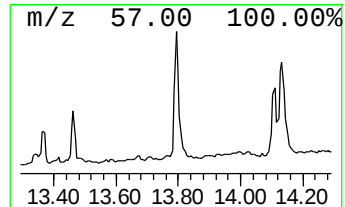
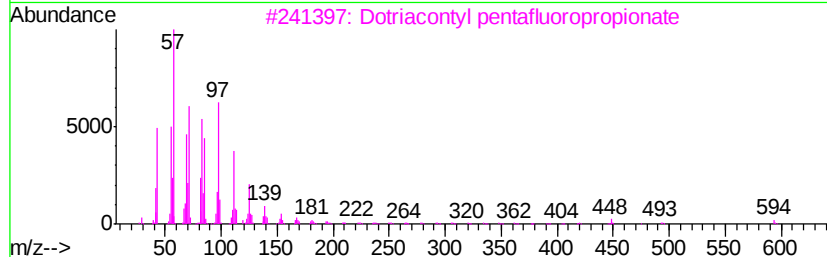
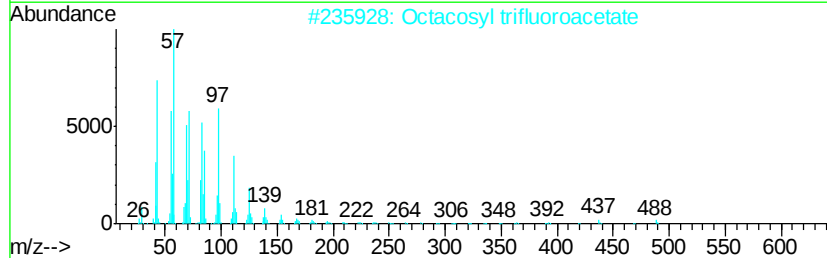
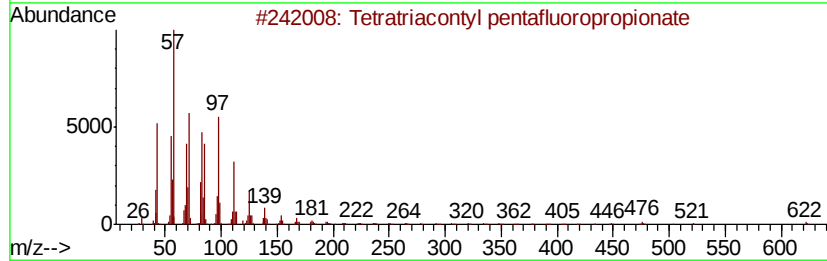
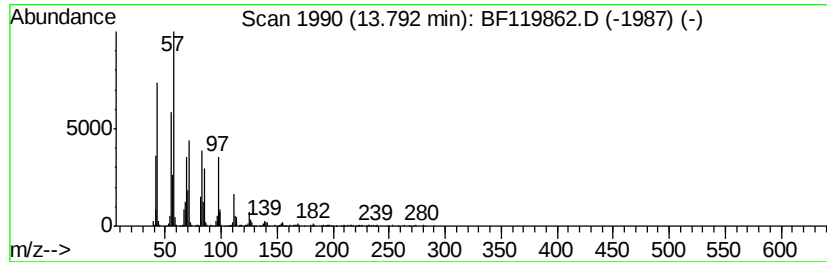
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Tetratriacontyl pentafluoro... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.79	11.78 ng	822705	Chrysene-d12	13.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetratriacontyl pentafluoropropi...	641	C37H69F5O2	1000351-81-5	91
2		Octacosyl trifluoroacetate	506	C30H57F3O2	1000351-74-9	91
3		Dotriacontyl pentafluoropropionate	612	C35H65F5O2	1000351-81-4	91
4		Tetratriacontyl heptafluorobutyrate	691	C38H69F7O2	1000351-84-1	91
5		Triacetyl pentafluoropropionate	584	C33H61F5O2	1000351-80-0	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040820\
Data File : BF119862.D
Acq On : 9 Apr 2020 9:00
Operator : MA/SJ
Sample : L2151-11
Misc :
ALS Vial : 30 Sample Multiplier: 1

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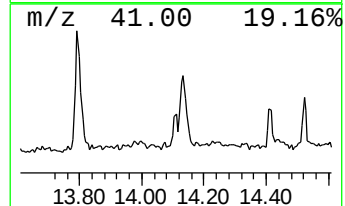
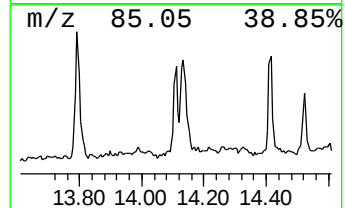
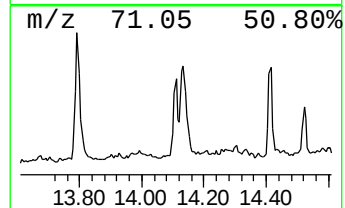
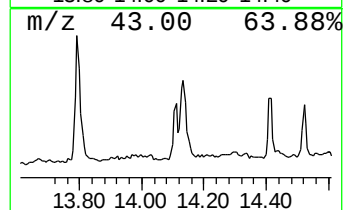
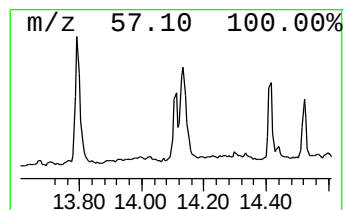
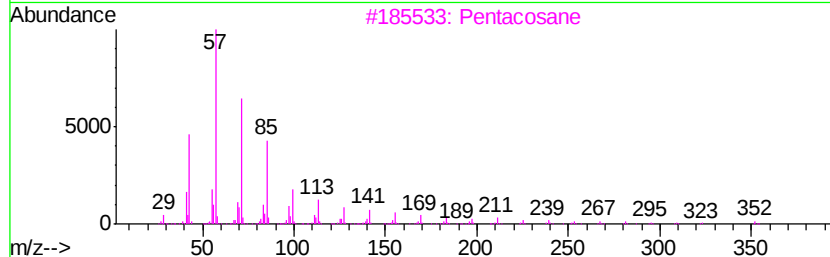
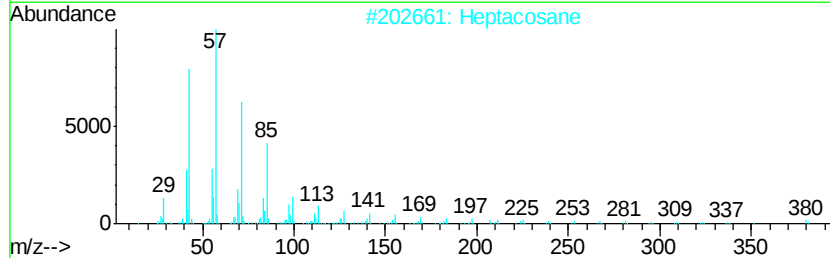
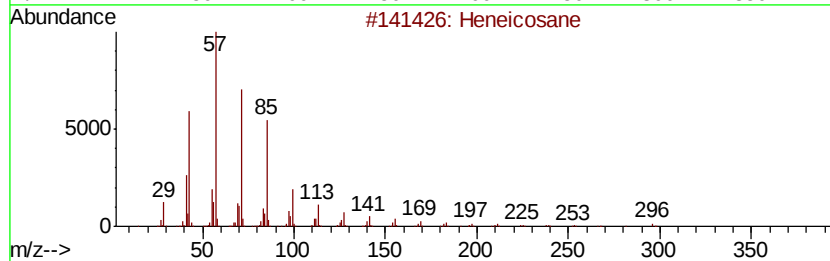
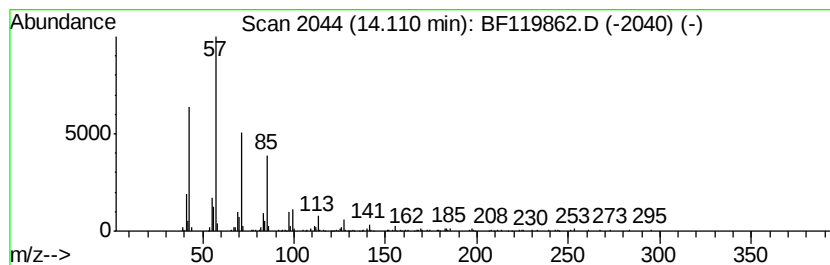
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Heneicosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.11	3.26 ng	227902	Chrysene-d12	13.93

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	87
2	Heptacosane		380	C27H56	000593-49-7	81
3	Pentacosane		352	C25H52	000629-99-2	80
4	Hexatriacontane		507	C36H74	000630-06-8	74
5	Hexacosane		366	C26H54	000630-01-3	74



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040820\
Data File : BF119862.D
Acq On : 9 Apr 2020 9:00
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Sample : L2151-11
Misc :
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ClientSampleId :
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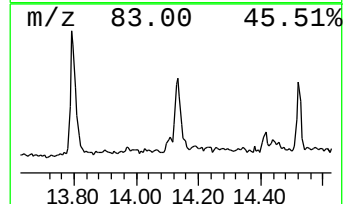
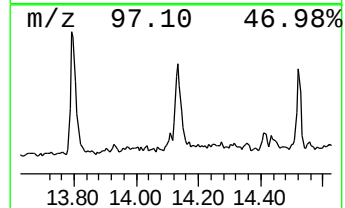
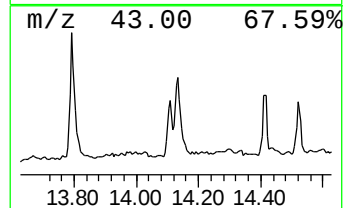
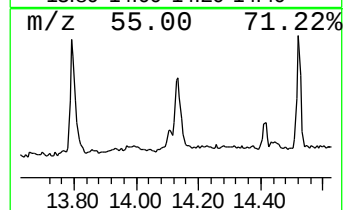
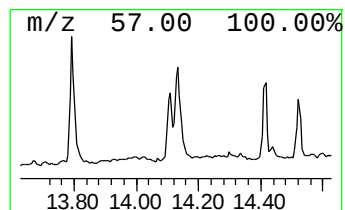
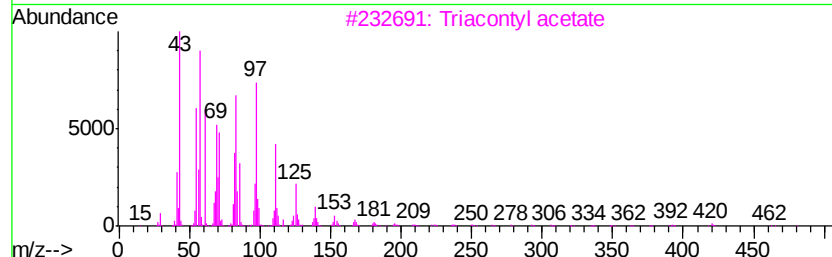
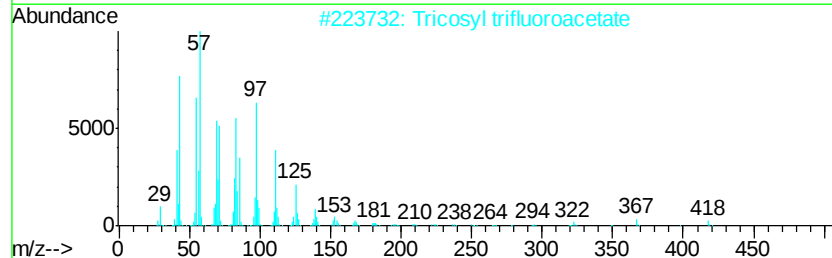
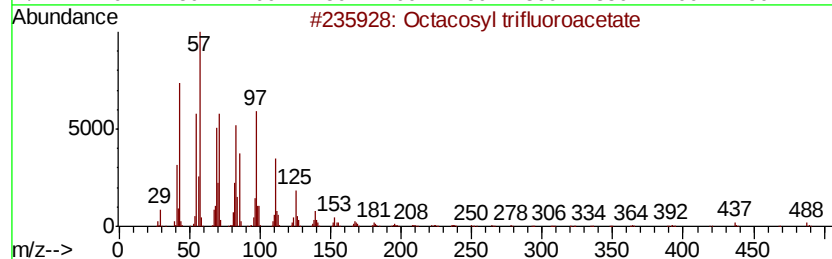
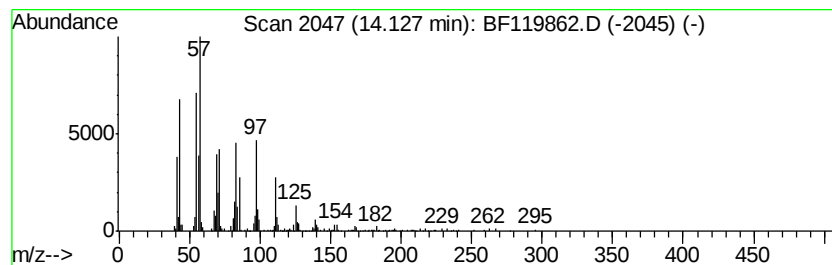
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Octacosyl trifluoroacetate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.13	9.28 ng	647815	Chrysene-d12	13.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosyl trifluoroacetate	506	C30H57F3O2	1000351-74-9	95
2		Tricosyl trifluoroacetate	436	C25H47F3O2	1000351-75-1	94
3		triacontyl acetate	480	C32H64O2	041755-58-2	91
4		Oxalic acid, isobutyl heptadecyl...	384	C23H44O4	1000309-38-2	91
5		Carbonic acid, hexadecyl isobuty...	342	C21H42O3	1000314-61-3	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040820\
Data File : BF119862.D
Acq On : 9 Apr 2020 9:00
Operator : MA/SJ
Sample : L2151-11
Misc :
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ClientSampleId :
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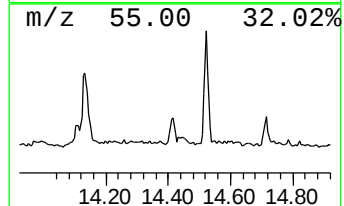
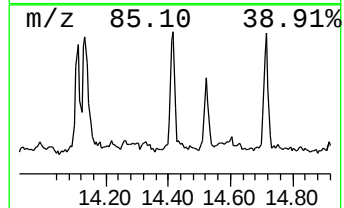
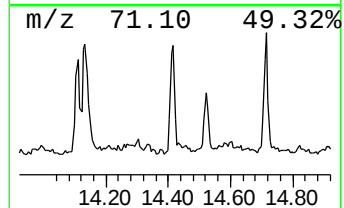
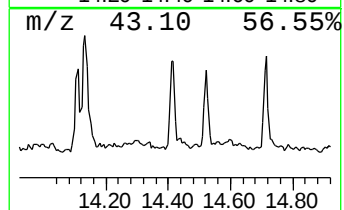
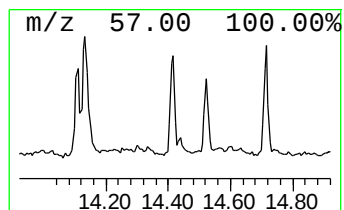
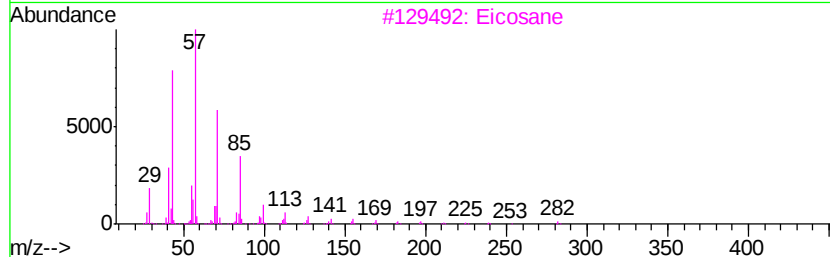
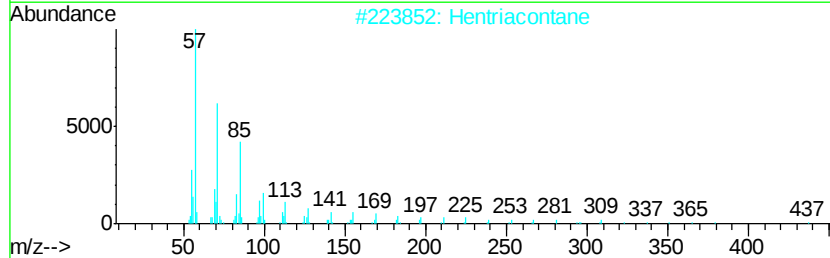
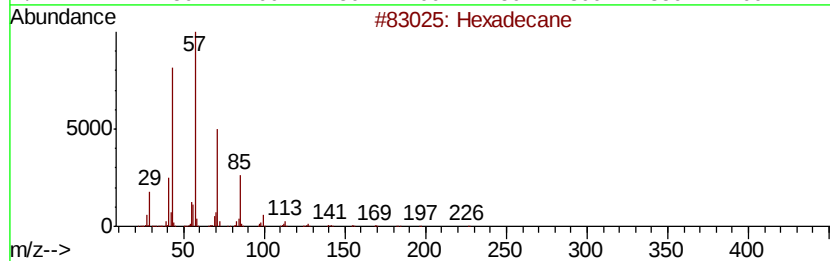
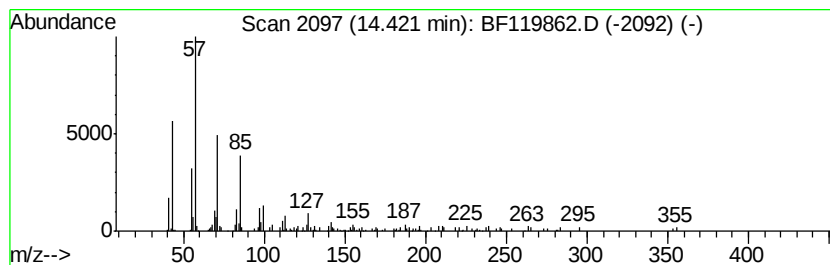
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Hexadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.42	3.98 ng	277611	Chrysene-d12	13.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	94
2		Hentriacontane	437	C31H64	000630-04-6	90
3		Eicosane	282	C20H42	000112-95-8	87
4		Pentacosane	352	C25H52	000629-99-2	83
5		Nonadecane	268	C19H40	000629-92-5	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040820\
Data File : BF119862.D
Acq On : 9 Apr 2020 9:00
Operator : MA/SJ
Sample : L2151-11
Misc :
ALS Vial : 30 Sample Multiplier: 1

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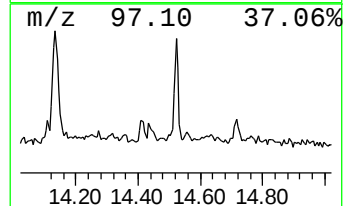
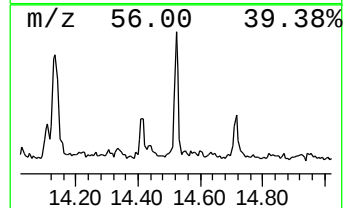
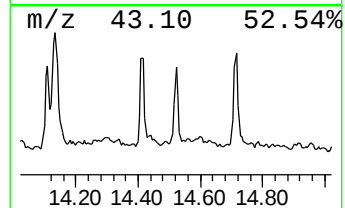
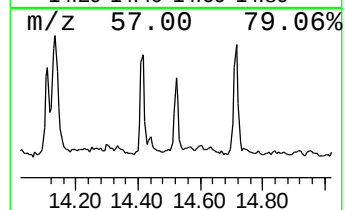
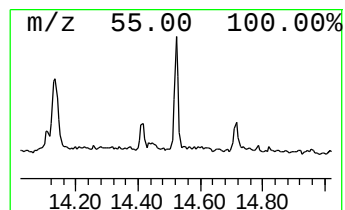
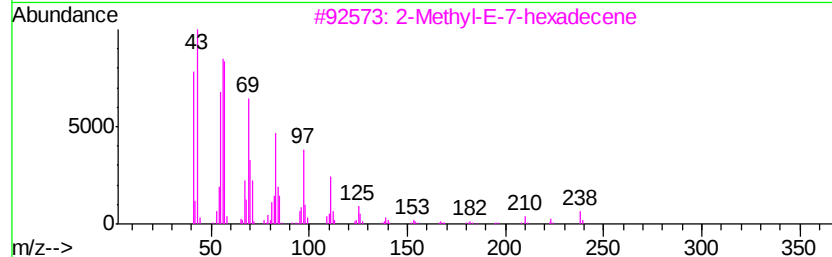
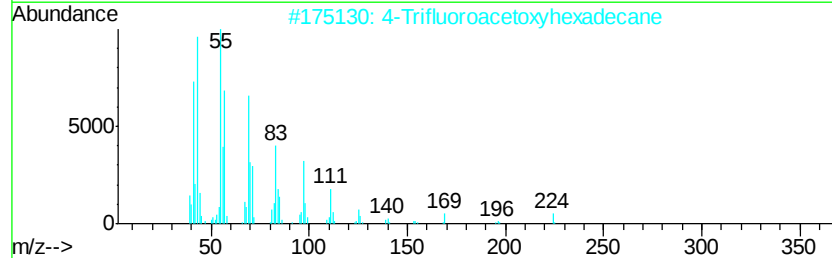
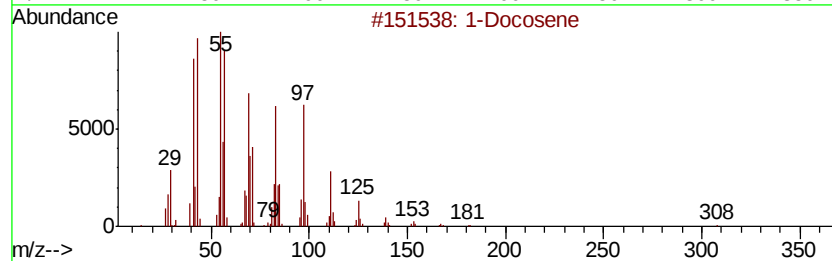
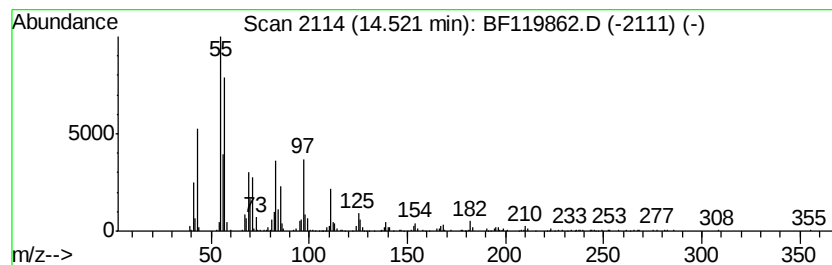
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 1-Docosene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.52	5.33 ng	371906	Chrysene-d12	13.93

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Docosene	308	C22H44	001599-67-3	76
2			4-Trifluoroacetoxyhexadecane	338	C18H33F3O2	1000215-97-5	72
3			2-Methyl-E-7-hexadecene	238	C17H34	064183-52-4	70
4			Cyclotetracosane	336	C24H48	000297-03-0	58
5			Hexacosyl trifluoroacetate	478	C28H53F3O2	1000351-75-0	52



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040820\
Data File : BF119862.D
Acq On : 9 Apr 2020 9:00
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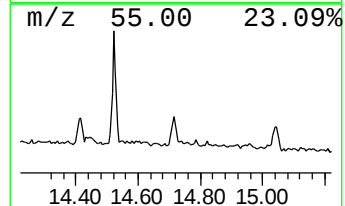
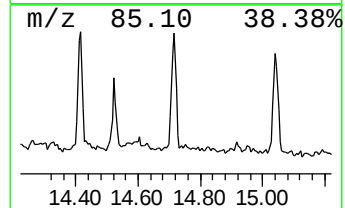
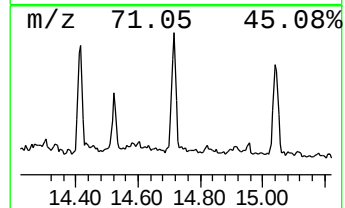
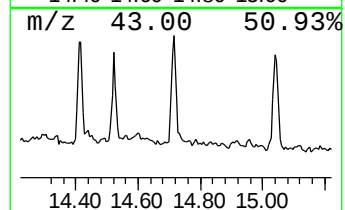
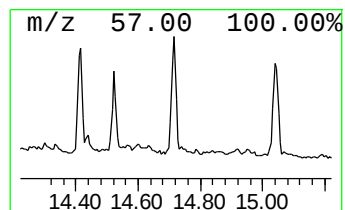
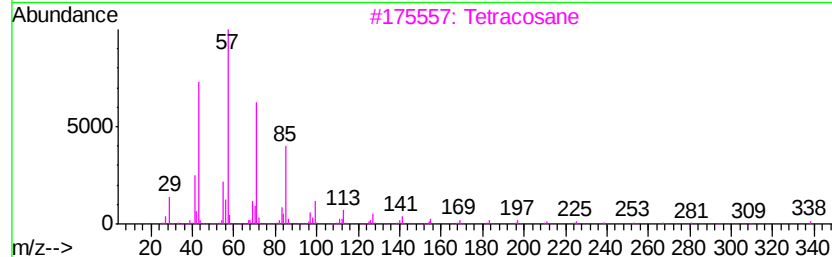
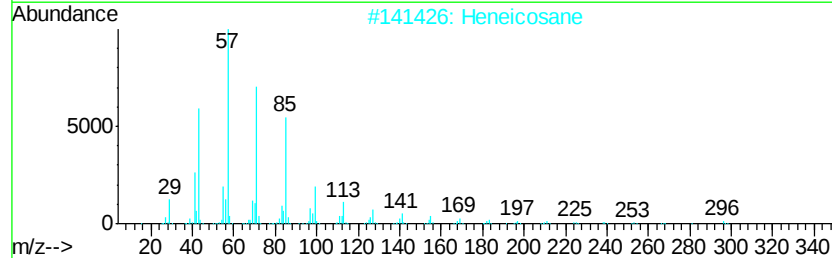
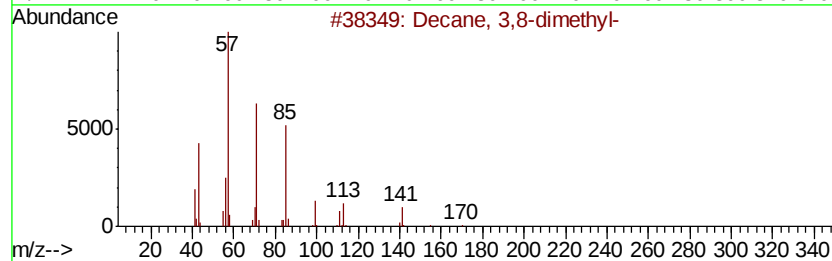
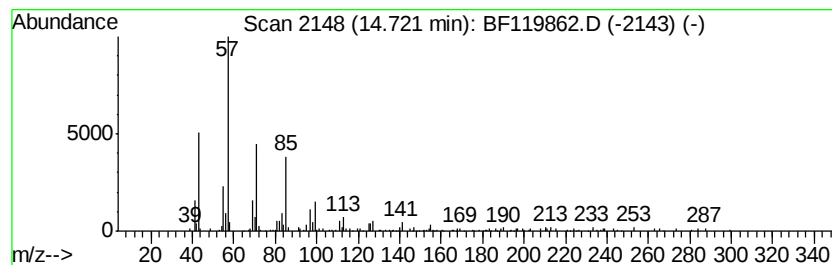
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Decane, 3,8-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.72	5.59 ng	304694	Perylene-d12	15.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	90
2		Heneicosane	296	C21H44	000629-94-7	90
3		Tetracosane	338	C24H50	000646-31-1	90
4		Pentacosane	352	C25H52	000629-99-2	90
5		Hexacosane	366	C26H54	000630-01-3	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040820\
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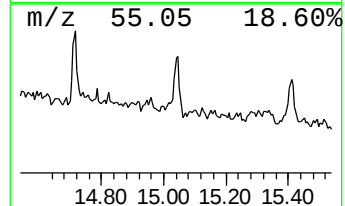
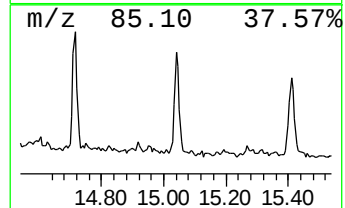
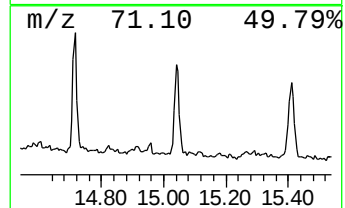
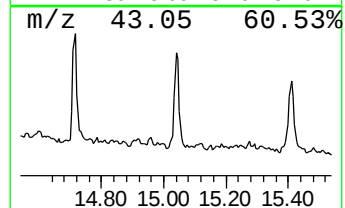
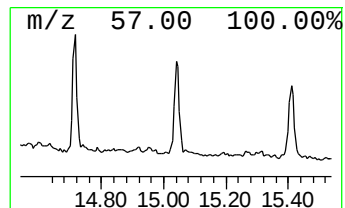
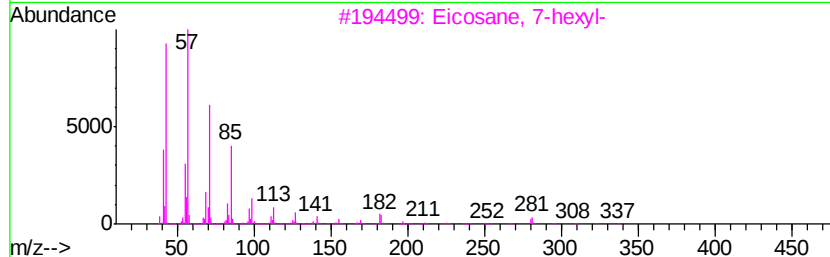
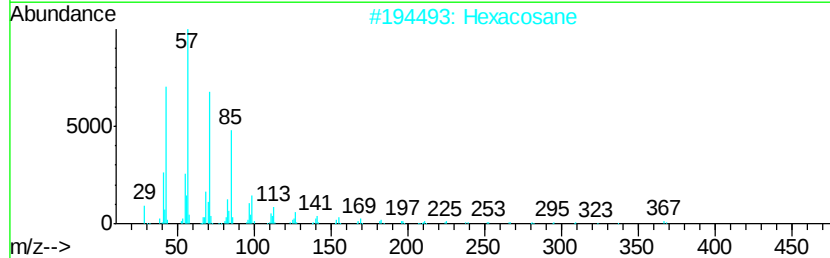
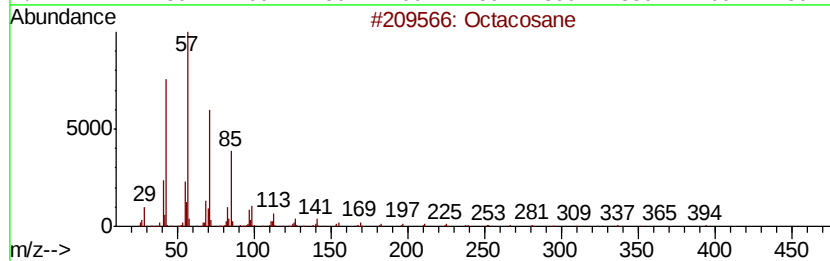
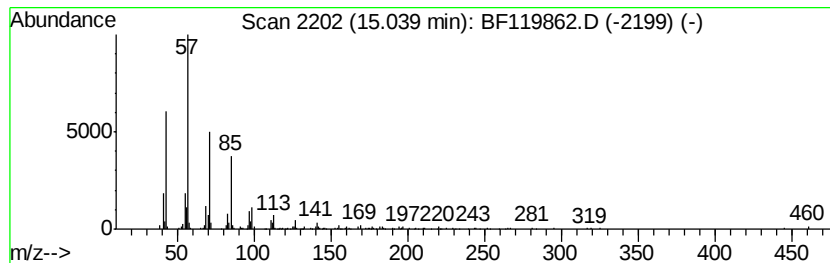
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Octacosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.04	5.94 ng	323643	Perylene-d12	15.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	87
2		Hexacosane	366	C26H54	000630-01-3	81
3		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	74
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	74
5		Eicosane	282	C20H42	000112-95-8	74



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040820\
Data File : BF119862.D
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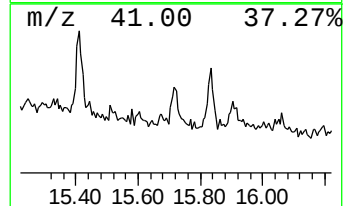
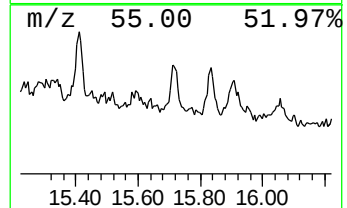
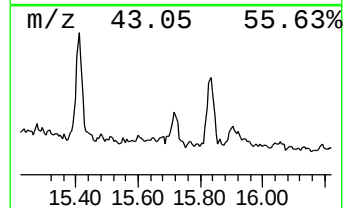
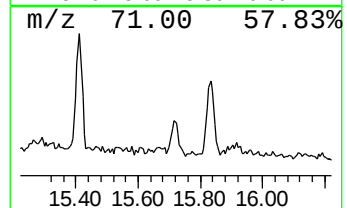
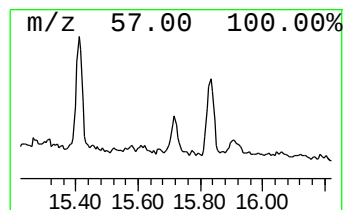
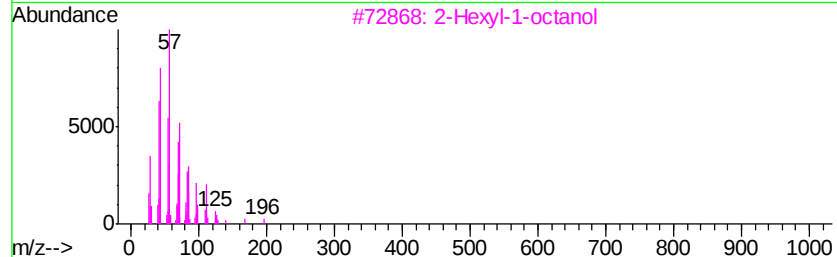
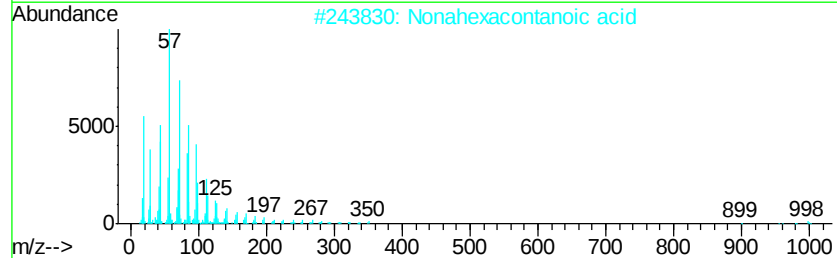
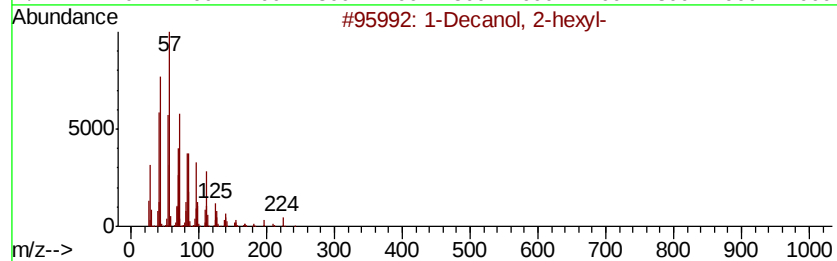
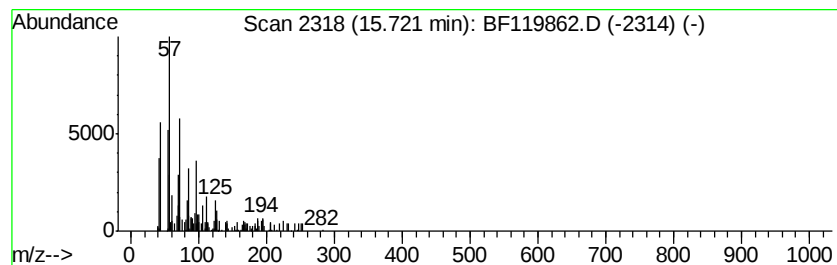
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 1-Decanol, 2-hexyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.72	2.64 ng	143887	Perylene-d12	15.37

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Decanol, 2-hexyl-	242	C16H34O	002425-77-6	72
2			Nonahexacontanoic acid	999	C69H138O2	040710-32-5	64
3			2-Hexyl-1-octanol	214	C14H30O	019780-79-1	64
4			Sulfurous acid, butyl tetradecyl...	334	C18H38O3S	1000309-18-1	52
5			Sulfurous acid, 2-propyl tetradecyl...	320	C17H36O3S	1000309-12-5	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040820\
Data File : BF119862.D
Acq On : 9 Apr 2020 9:00
Operator : MA/SJ
Sample : L2151-11
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CBH-140

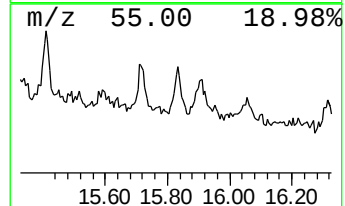
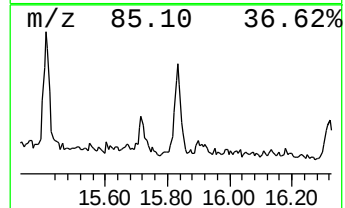
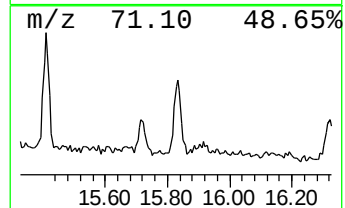
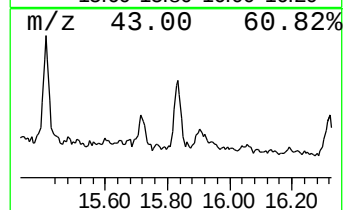
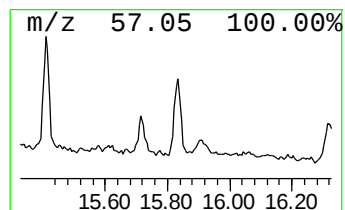
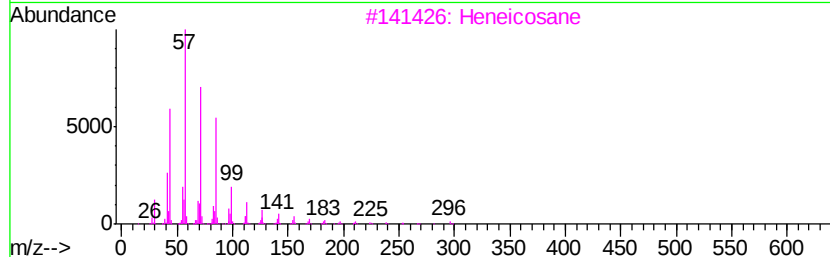
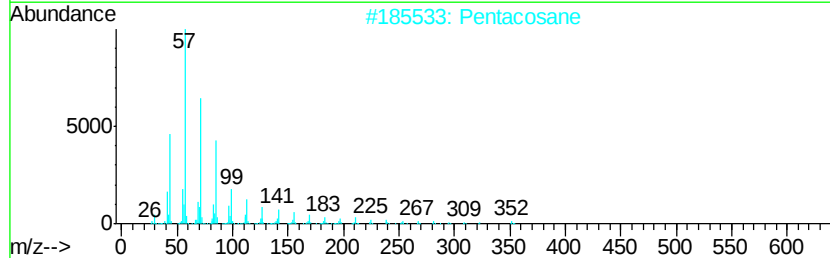
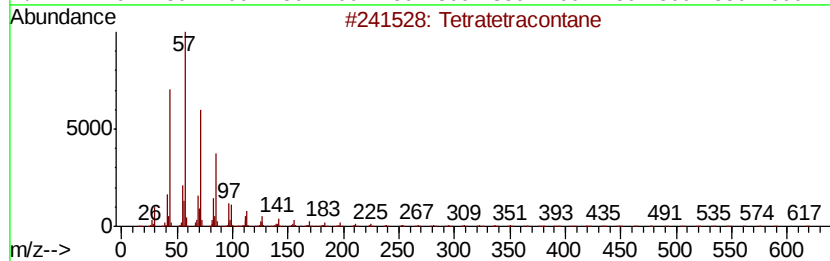
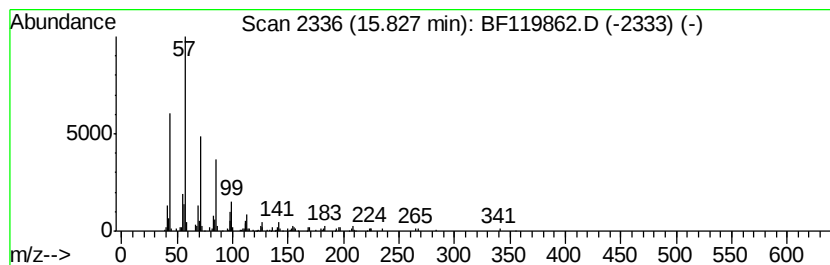
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Tetratetracontane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.83	3.43 ng	186904	Perylene-d12	15.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetratetracontane	619	C44H90	007098-22-8	86
2		Pentacosane	352	C25H52	000629-99-2	86
3		Heneicosane	296	C21H44	000629-94-7	81
4		Heptadecane	240	C17H36	000629-78-7	80
5		Docosane	310	C22H46	000629-97-0	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040820\
Data File : BF119862.D
Acq On : 9 Apr 2020 9:00
Operator : MA/SJ
Sample : L2151-11
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CBH-140

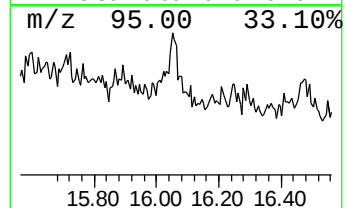
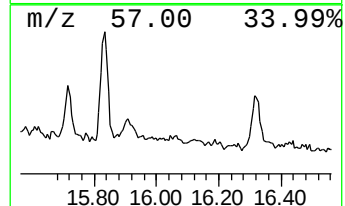
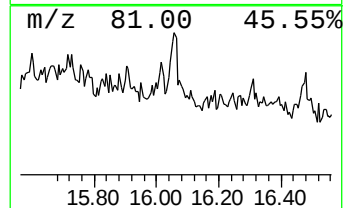
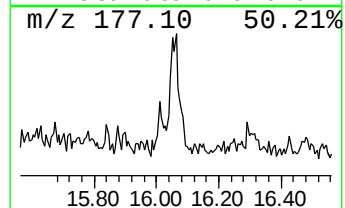
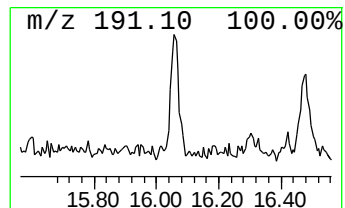
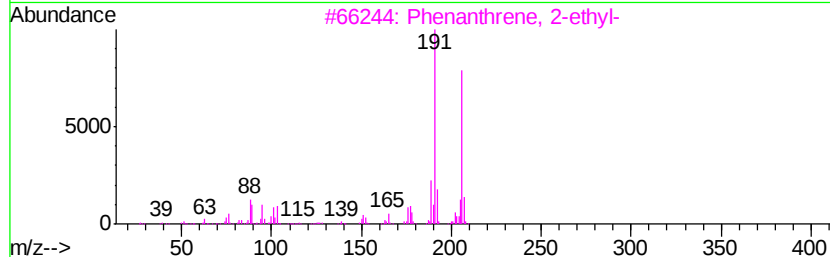
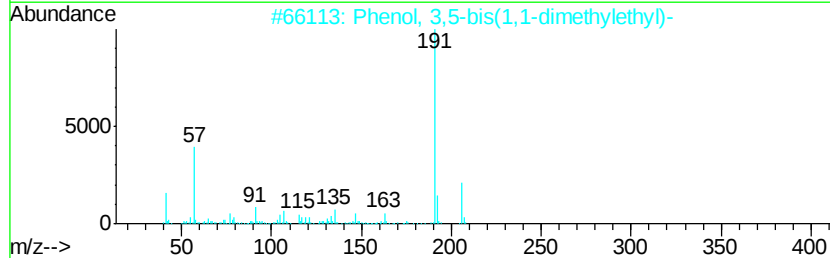
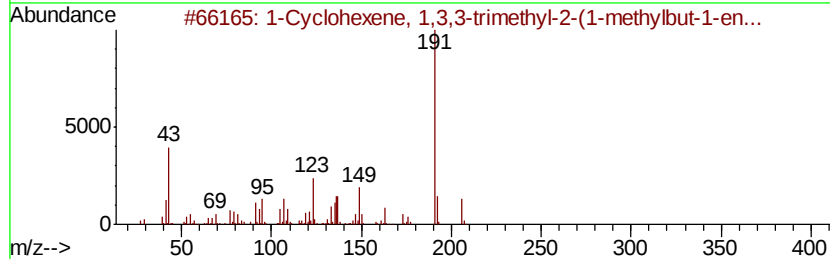
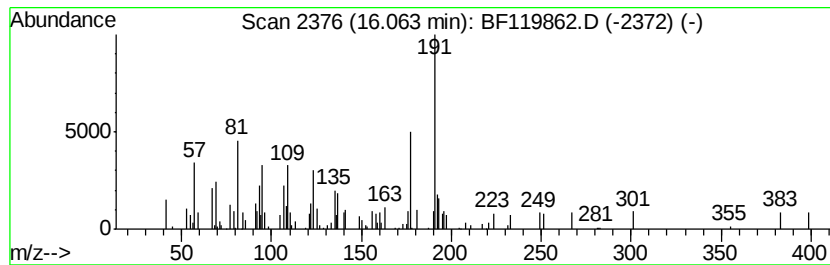
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 unknown16.06 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.06	2.27 ng	123473	Perylene-d12	15.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Cyclohexene, 1,3,3-trimethyl-2...	206	C14H22O	1000197-08-4	45
2		Phenol, 3,5-bis(1,1-dimethylethyl)-	206	C14H22O	001138-52-9	38
3		Phenanthrene, 2-ethyl-	206	C16H14	003674-74-6	35
4		Propanamide, 2,2-dimethyl-N-(3-m...	191	C12H17NO	032597-29-8	30
5		Phenol, 2,5-bis(1,1-dimethylethyl)-	206	C14H22O	005875-45-6	30



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040820\
 Data File : BF119862.D
 Acq On : 9 Apr 2020 9:00
 Operator : MA/SJ
 Sample : L2151-11
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CBH-140

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF040820.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...	4.96	2.2	ng	128713	1	6.76	1180970	20.0
n-Hexadecanoic acid	11.83	15.8	ng	1496100	4	11.29	1889380	20.0
Octadecanoic acid	12.62	12.9	ng	902144	5	13.93	1396550	20.0
unknown13.37	13.37	3.8	ng	268330	5	13.93	1396550	20.0
Tetratriacontyl p...	13.79	11.8	ng	822705	5	13.93	1396550	20.0
Heneicosane	14.11	3.3	ng	227902	5	13.93	1396550	20.0
Octacosyl trifluo...	14.13	9.3	ng	647815	5	13.93	1396550	20.0
Hexadecane	14.42	4.0	ng	277611	5	13.93	1396550	20.0
1-Docosene	14.52	5.3	ng	371906	5	13.93	1396550	20.0
Decane, 3,8-dimet...	14.72	5.6	ng	304694	6	15.37	1089480	20.0
Octacosane	15.04	5.9	ng	323643	6	15.37	1089480	20.0
1-Decanol, 2-hexyl-	15.72	2.6	ng	143887	6	15.37	1089480	20.0
Tetratetracontane	15.83	3.4	ng	186904	6	15.37	1089480	20.0
unknown16.06	16.06	2.3	ng	123473	6	15.37	1089480	20.0