

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120618\
 Data File : BF111137.D
 Acq On : 6 Dec 2018 14:41
 Operator : JU/SJ
 Sample : J6204-08
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HLSP-SB-2B-(2-9.75)

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-BF120518.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	3.246	194	197	205	rBV	155353	284625	1.76%	0.194%
2	5.028	497	500	509	rVB	662312	879928	5.45%	0.601%
3	5.428	562	568	571	rBV	13732588	14631215	90.59%	9.998%
4	6.440	733	740	744	rBV	13598231	16150971	100.00%	11.037%
5	6.575	758	763	767	rBV	13227135	14846442	91.92%	10.145%
6	6.804	799	802	806	rVB	3758511	3057383	18.93%	2.089%
7	6.963	825	829	832	rBV	11802250	11165724	69.13%	7.630%
8	7.369	892	898	901	rBV	9291930	9996812	61.90%	6.831%
9	8.087	1016	1020	1023	rBV	4790461	3978535	24.63%	2.719%
10	9.169	1198	1204	1207	rBV	16208891	15600534	96.59%	10.661%
11	9.557	1266	1270	1273	rBV	2508666	1970231	12.20%	1.346%
12	9.839	1314	1318	1322	rVB2	5037996	4354640	26.96%	2.976%
13	10.628	1446	1452	1455	rBV	10152336	10004550	61.94%	6.837%
14	11.322	1566	1570	1573	rBV	5309369	4361874	27.01%	2.981%
15	11.863	1656	1662	1680	rBV	4380012	4571475	28.30%	3.124%
16	12.645	1791	1795	1806	rBV	2505975	2443218	15.13%	1.670%
17	12.916	1836	1841	1844	rBV	13868180	14469758	89.59%	9.888%
18	13.392	1914	1922	1927	rBV4	106085	201120	1.25%	0.137%
19	13.810	1990	1993	2010	rBV	2884654	3156324	19.54%	2.157%
20	13.957	2014	2018	2021	rVV	4326758	3669755	22.72%	2.508%
21	14.139	2046	2049	2053	rBV	192340	182302	1.13%	0.125%
22	14.445	2098	2101	2111	rVB	374434	464054	2.87%	0.317%
23	14.745	2149	2152	2156	rBV	289218	287385	1.78%	0.196%
24	14.821	2161	2165	2169	rVB	683955	658836	4.08%	0.450%
25	15.080	2205	2209	2212	rBV	323130	381243	2.36%	0.261%
26	15.392	2257	2262	2267	rBV2	2234980	2694687	16.68%	1.841%
27	15.451	2268	2272	2277	rVB	319206	371081	2.30%	0.254%
28	15.874	2340	2344	2348	rBV	253754	358705	2.22%	0.245%
29	15.921	2348	2352	2360	rVV	357692	666067	4.12%	0.455%
30	16.368	2424	2428	2434	rVB2	137040	237821	1.47%	0.163%
31	17.039	2537	2542	2549	rVB4	100871	240534	1.49%	0.164%

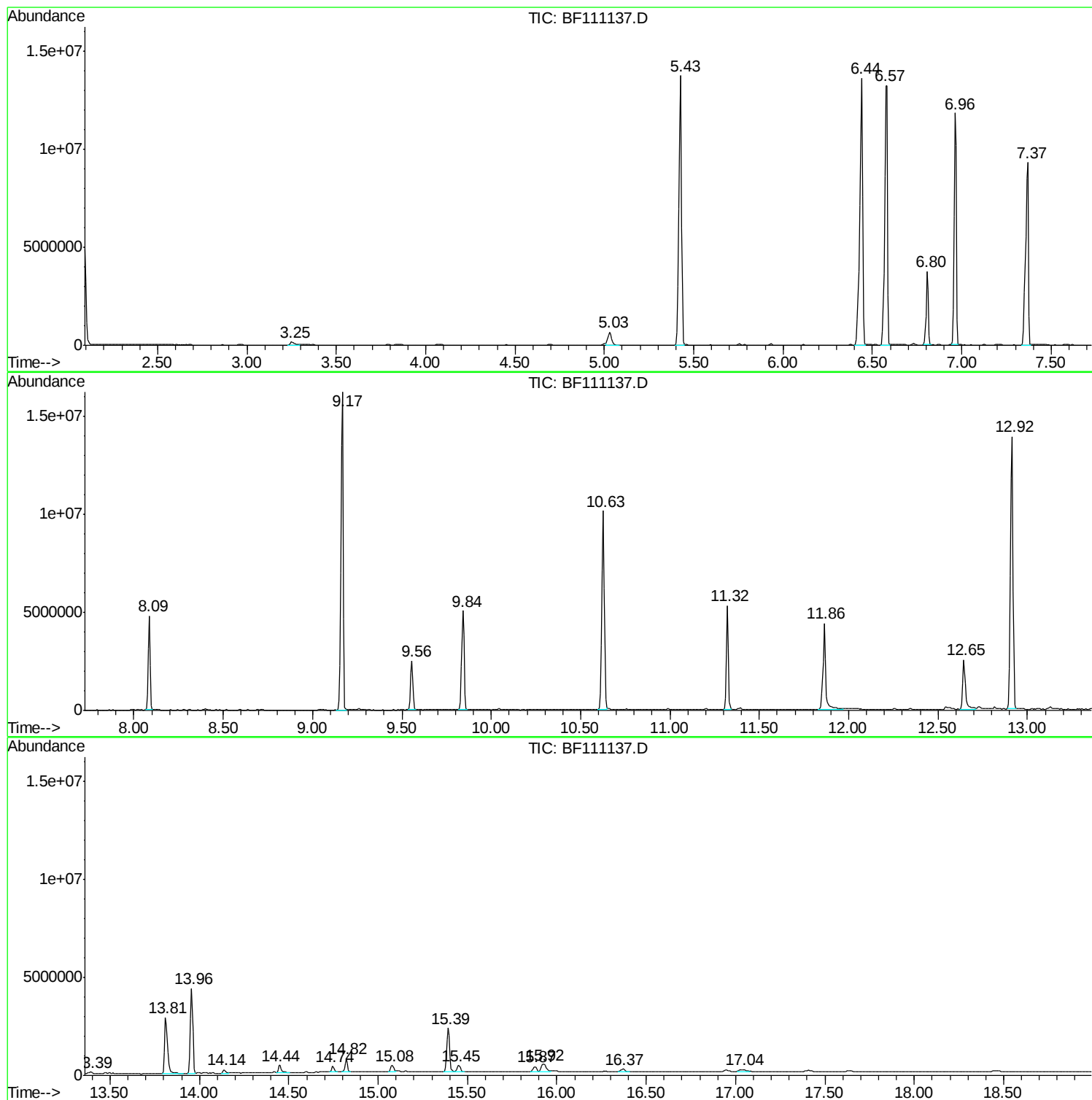
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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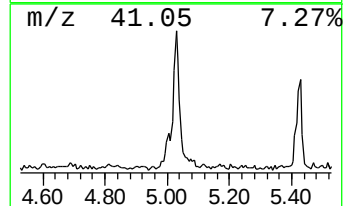
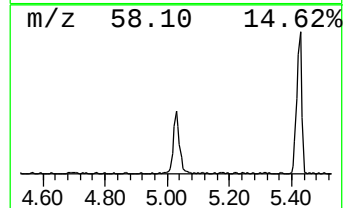
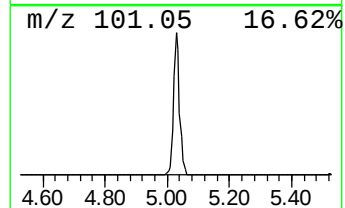
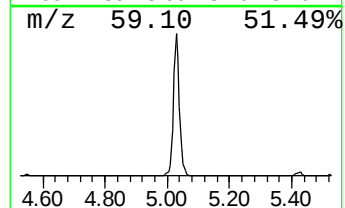
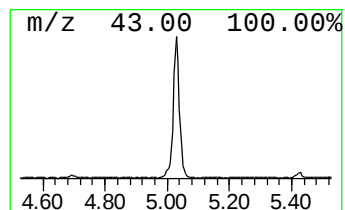
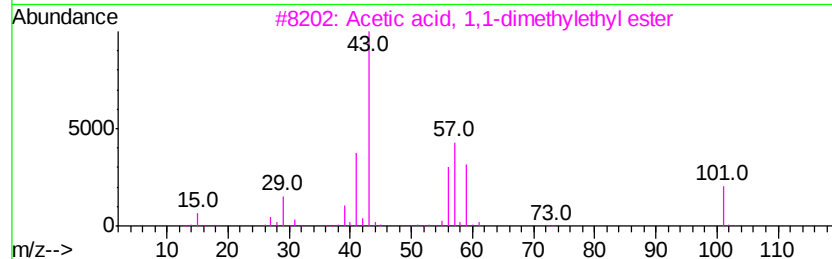
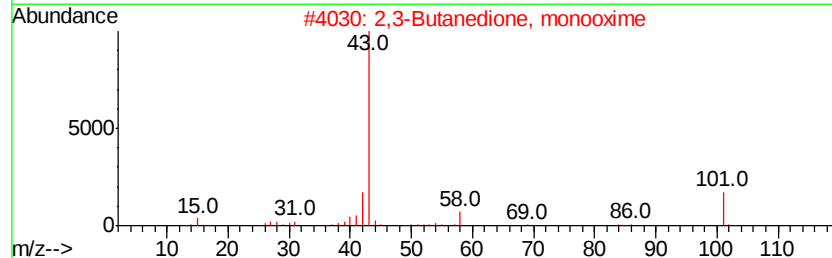
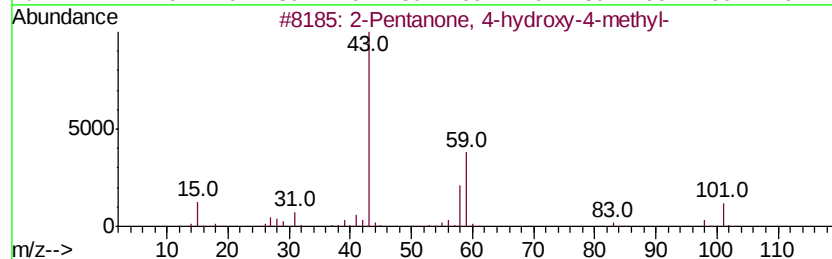
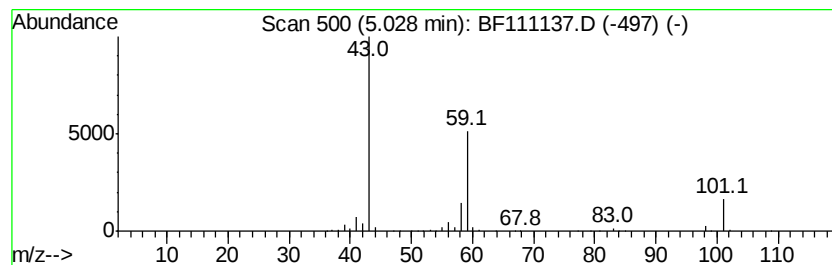
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
5.03	5.76 ng	879928	1,4-Dichlorobenzene-d4	6.80	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
3	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	9
4	Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	9
5	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	9



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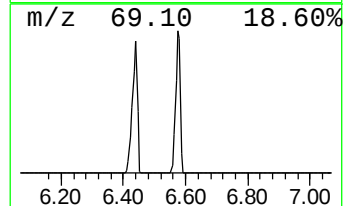
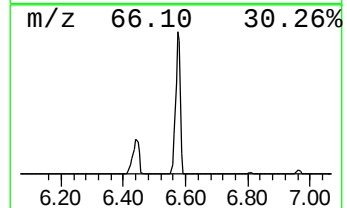
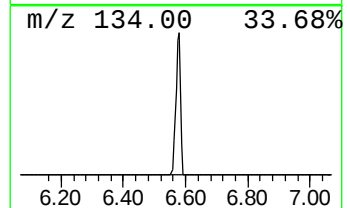
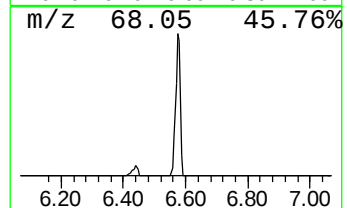
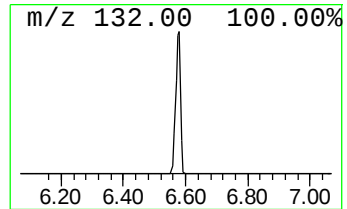
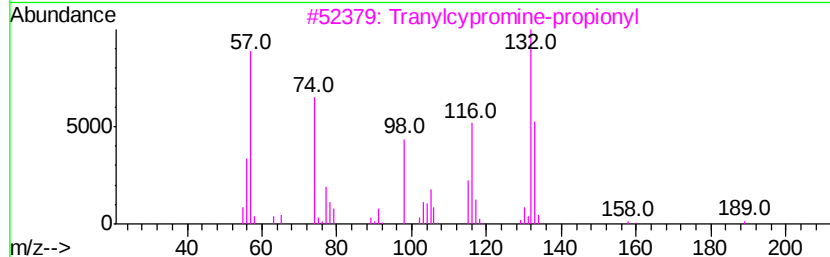
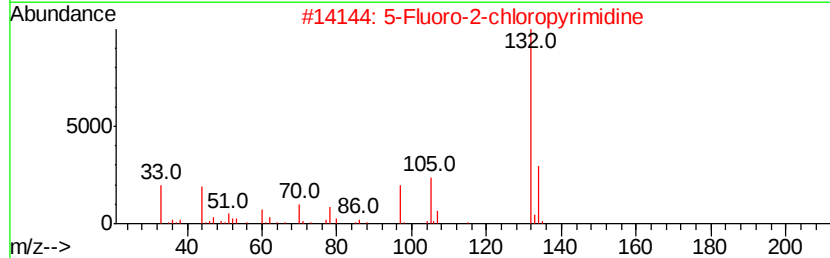
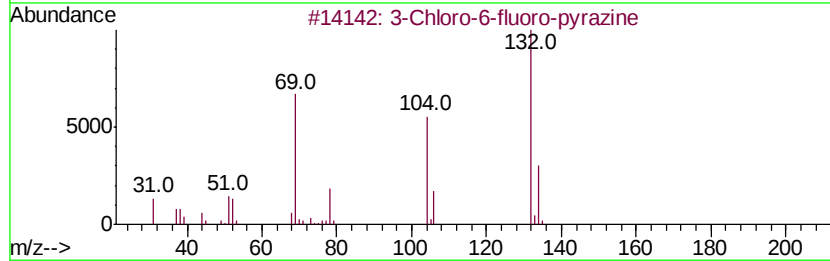
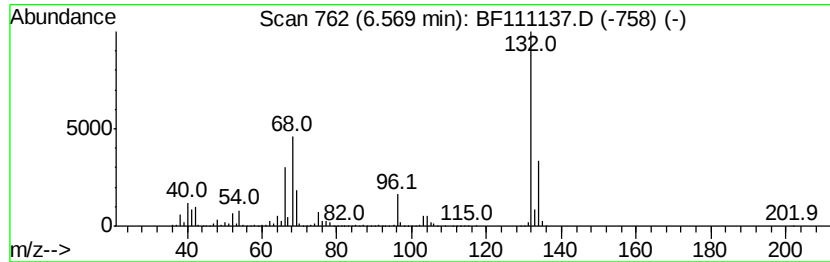
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Peak Number 2 unknown6.57 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.57	97.12 ng	14846400	1,4-Dichlorobenzene-d4	6.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
5		1,2,4-Trifluorobenzene	132	C6H3F3	000367-23-7	9



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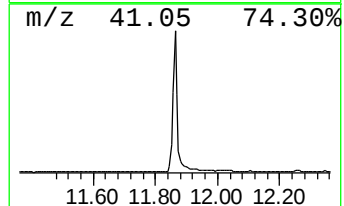
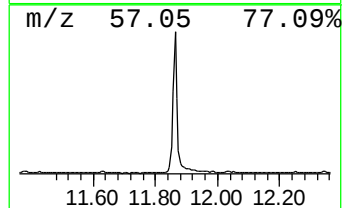
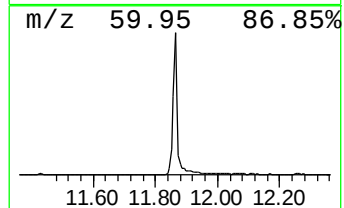
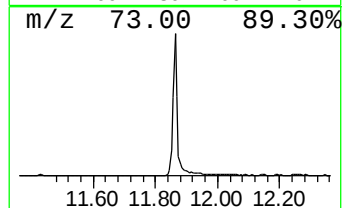
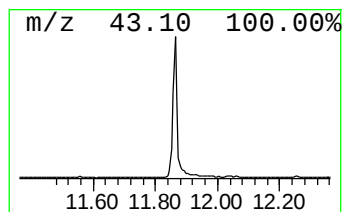
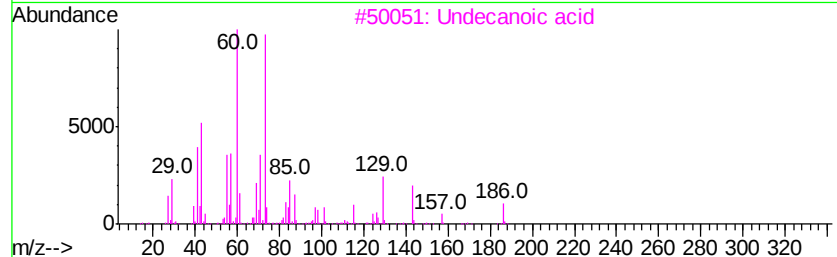
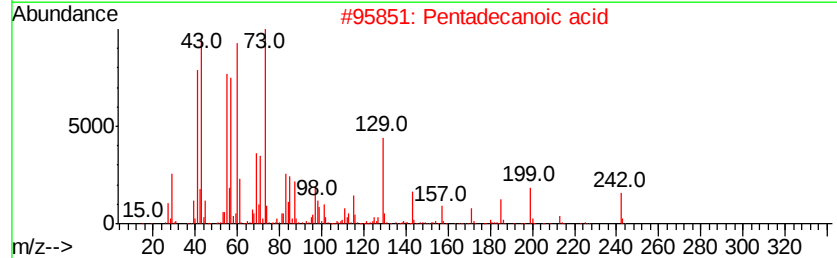
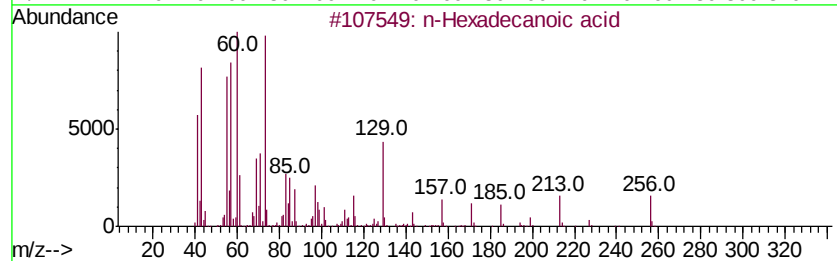
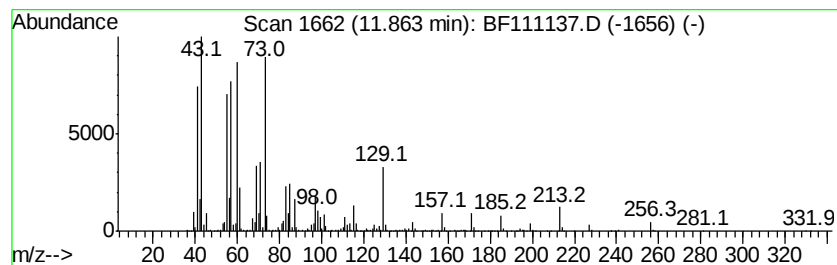
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.86	20.96 ng	4571480	Phenanthrene-d10	11.32

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	90
3		Undecanoic acid	186	C11H22O2	000112-37-8	72
4		Tridecanoic acid	214	C13H26O2	000638-53-9	64
5		n-Decanoic acid	172	C10H20O2	000334-48-5	64



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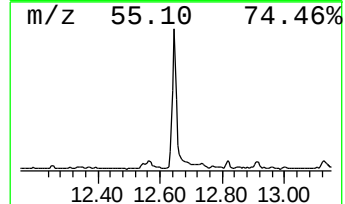
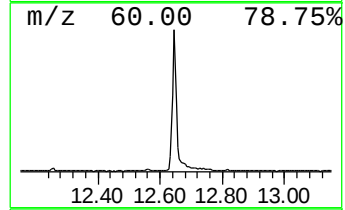
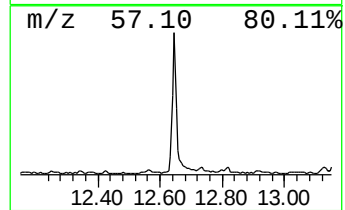
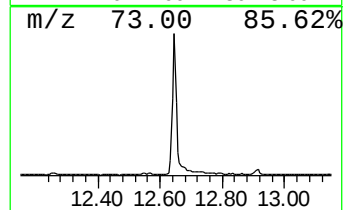
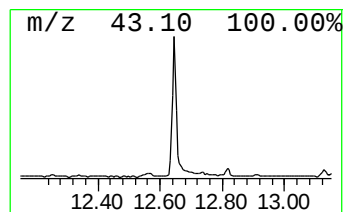
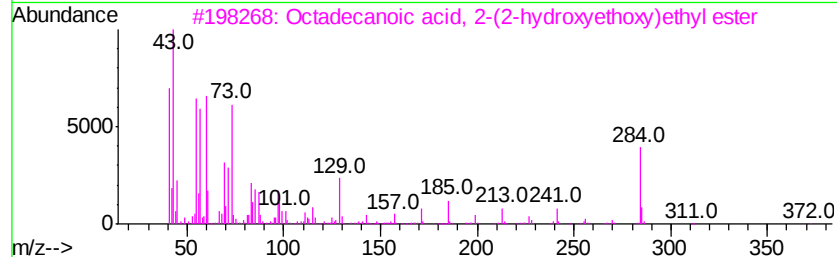
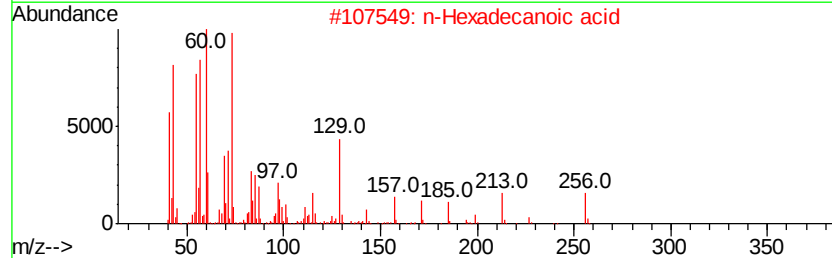
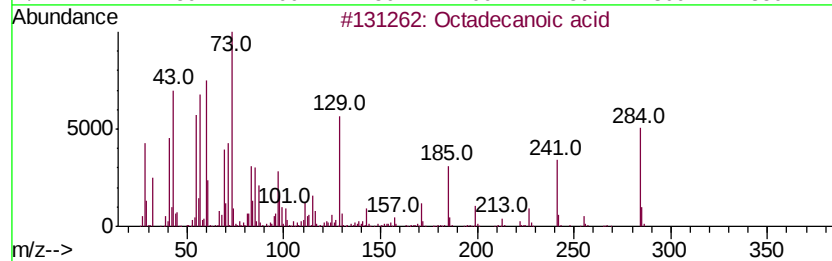
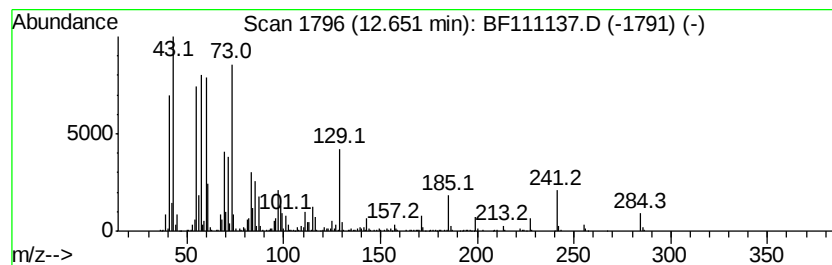
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Peak Number 5 Octadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.65	13.32 ng	2443220	Chrysene-d12	13.96
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Octadecanoic acid	284 C18H36O2	000057-11-4 99
2		n-Hexadecanoic acid	256 C16H32O2	000057-10-3 87
3		Octadecanoic acid, 2-(2-hydroxyve...	372 C22H44O4	000106-11-6 87
4		Tridecanoic acid	214 C13H26O2	000638-53-9 76
5		Tetradecanoic acid	228 C14H28O2	000544-63-8 59



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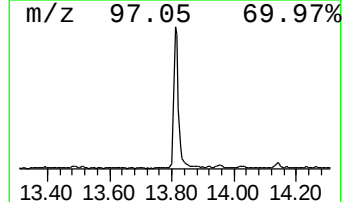
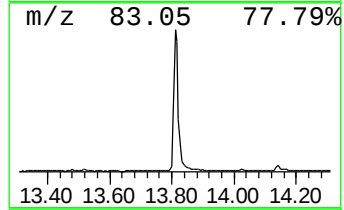
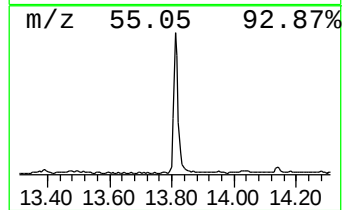
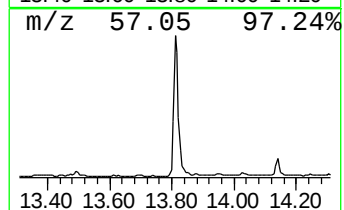
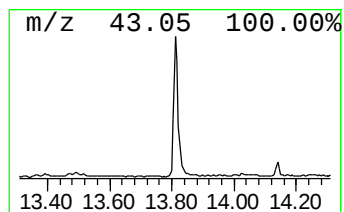
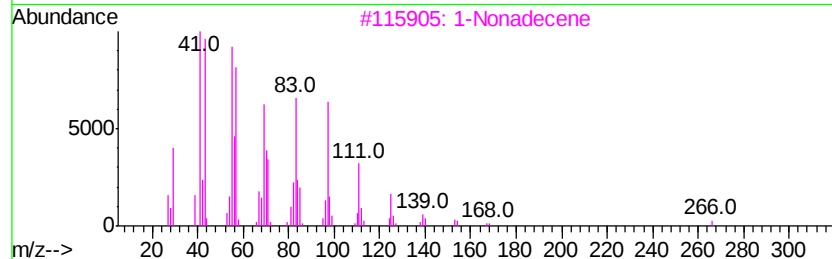
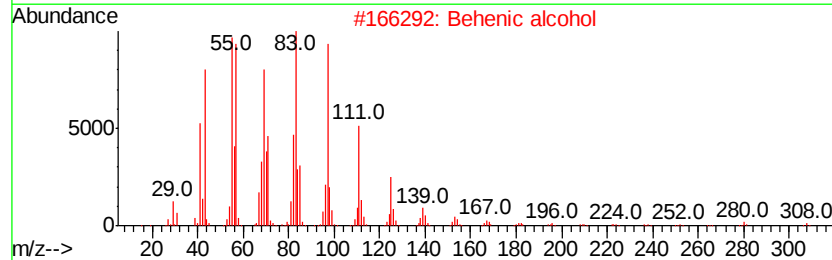
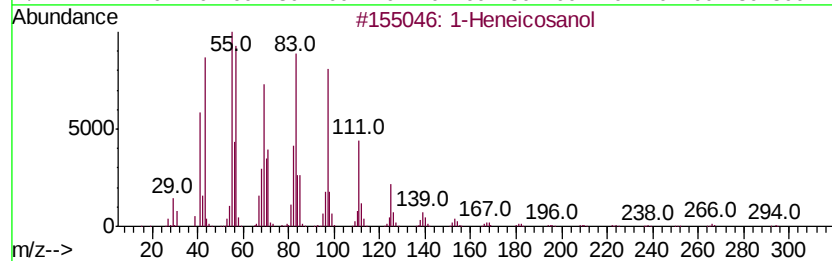
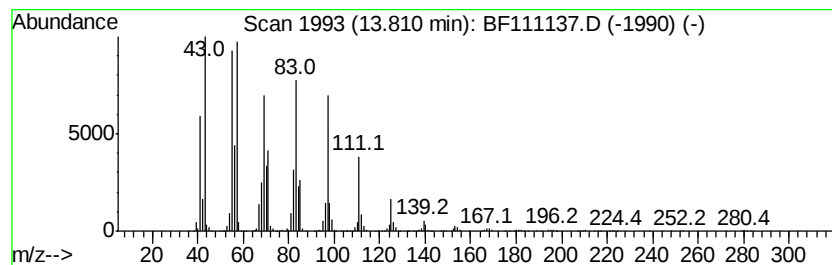
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 6 1-Heneicosanol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.81	17.20 ng	3156320	Chrysene-d12	13.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Heneicosanol	312	C21H44O	015594-90-8	93
2		Behenic alcohol	326	C22H46O	000661-19-8	91
3		1-Nonadecene	266	C19H38	018435-45-5	91
4		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	90
5		Bromoacetic acid, hexadecyl ester	362	C18H35BrO2	005454-48-8	90



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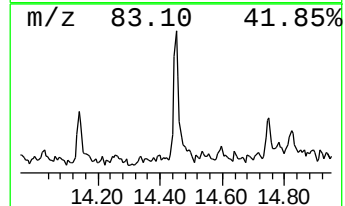
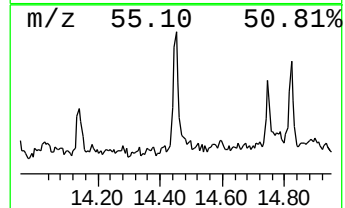
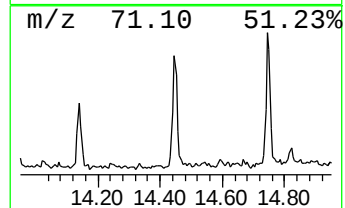
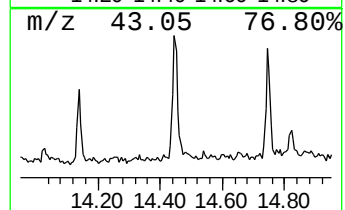
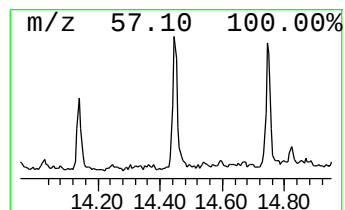
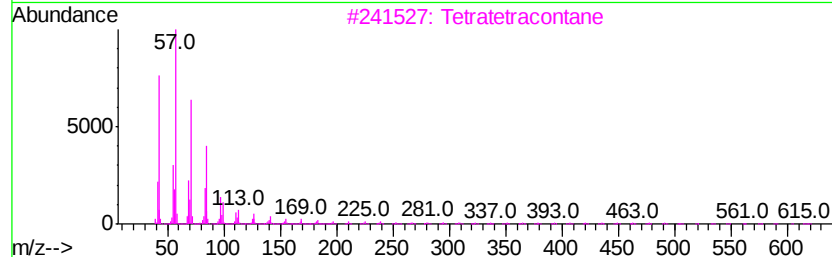
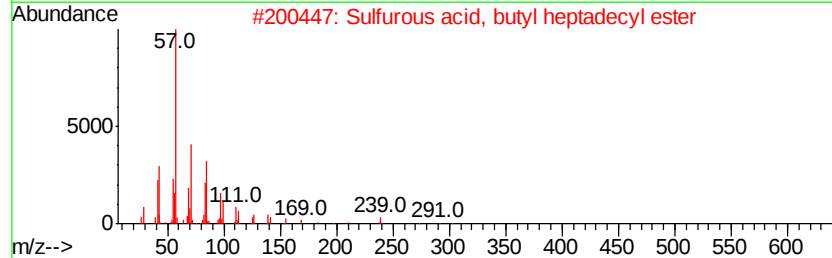
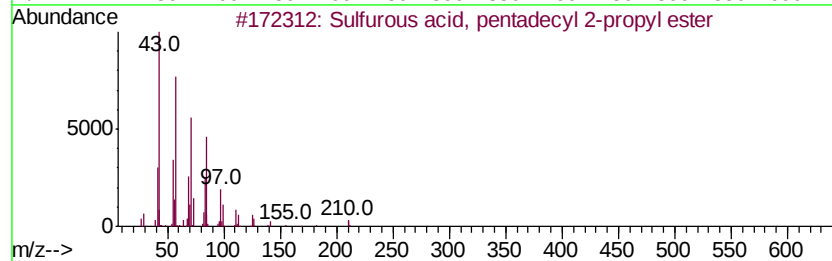
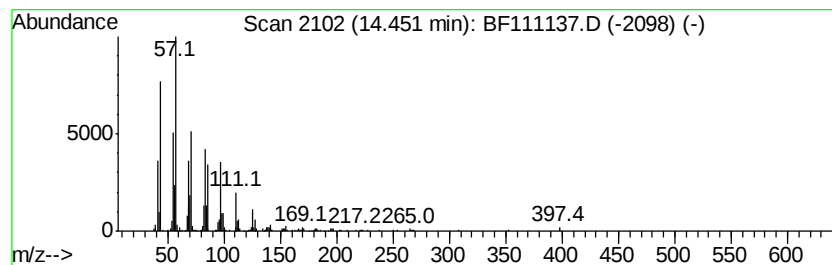
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Sulfurous acid, pentadecyl ... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.45	2.53 ng	464054	Chrysene-d12	13.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Sulfurous acid, pentadecyl 2-pro...	334	C18H38O3S	1000309-12-6	90
2		Sulfurous acid, butyl heptadecyl...	376	C21H44O3S	1000309-18-4	83
3		Tetratetracontane	619	C44H90	007098-22-8	72
4		Nonadecane	268	C19H40	000629-92-5	64
5		Heneicosane	296	C21H44	000629-94-7	62



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120618\
Data File : BF111137.D
Acq On : 6 Dec 2018 14:41
Operator : JU/SJ
Sample : J6204-08
Misc :
ALS Vial : 9 Sample Multiplier: 1

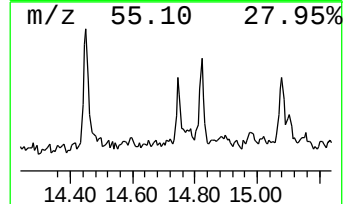
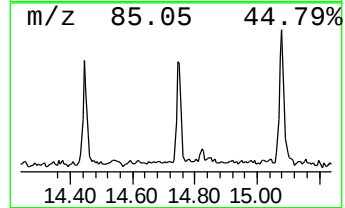
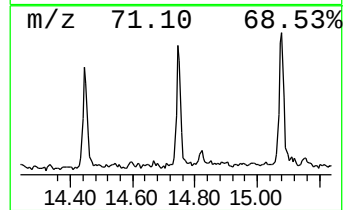
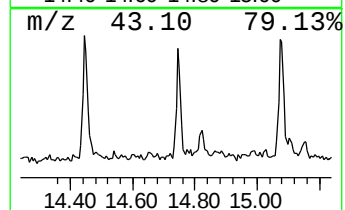
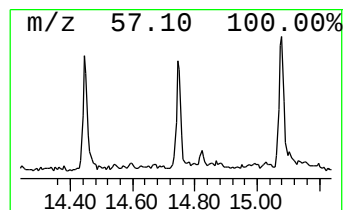
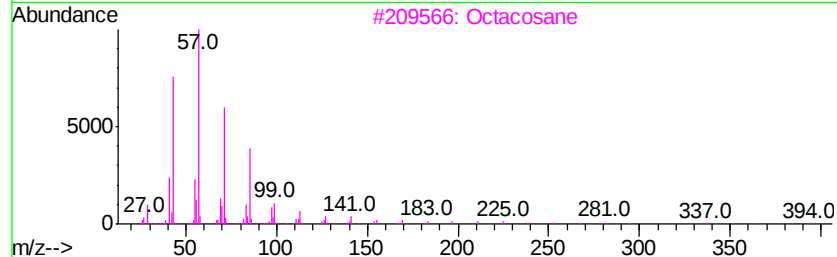
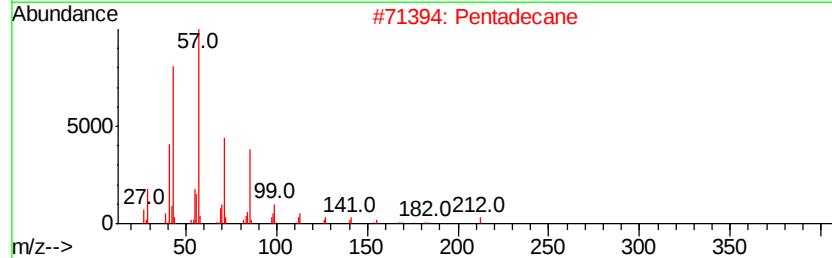
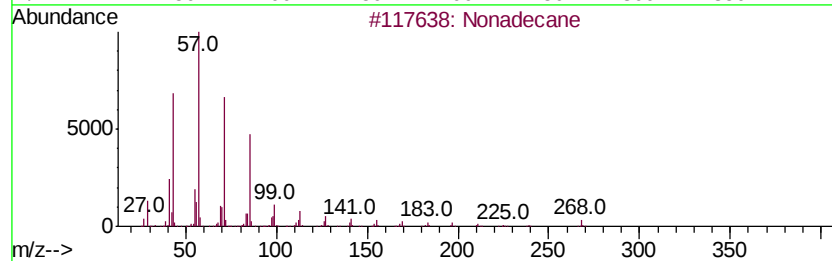
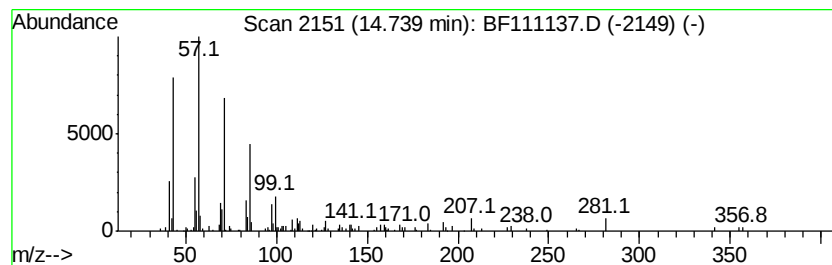
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HLSP-SB-2B-(2-9.75)

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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 Nonadecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.74	2.13 ng	287385	Perylene-d12	15.39
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Nonadecane	268 C19H40	000629-92-5	95
2	Pentadecane	212 C15H32	000629-62-9	92
3	Octacosane	394 C28H58	000630-02-4	90
4	Hexacosane	366 C26H54	000630-01-3	90
5	2-methyloctacosane	408 C29H60	1000376-72-8	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120618\
Data File : BF111137.D
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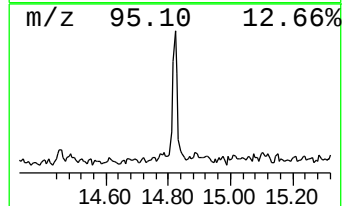
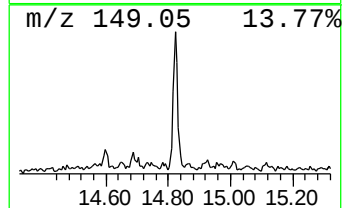
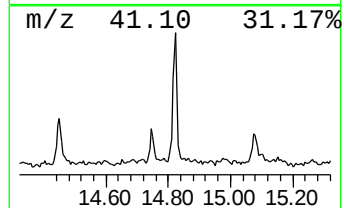
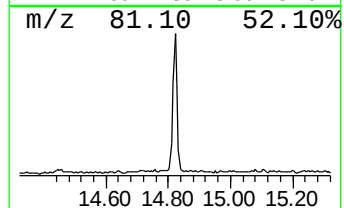
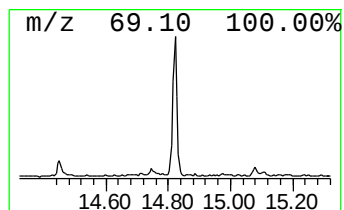
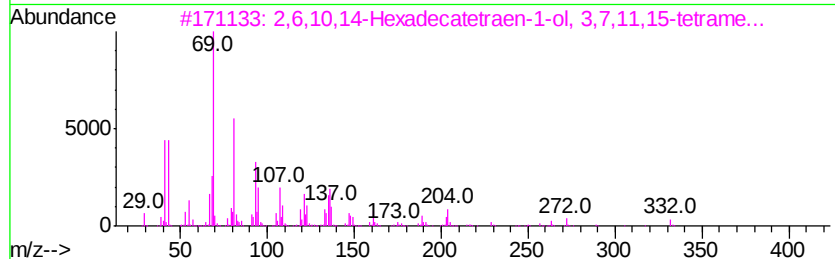
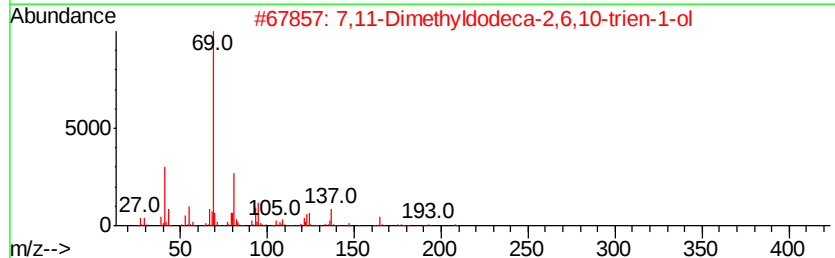
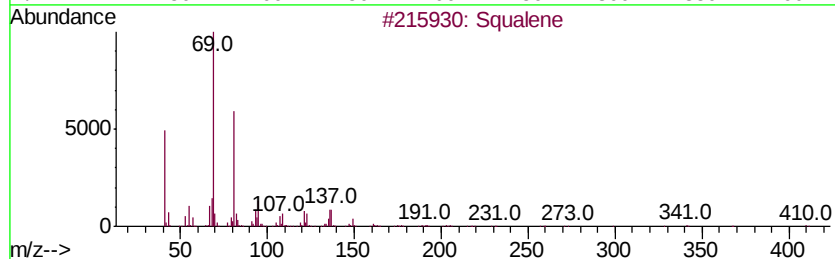
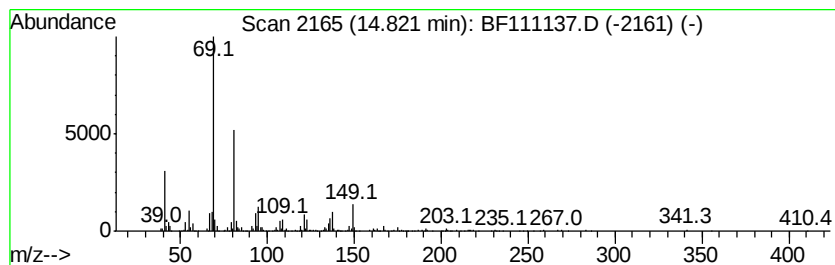
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Squalene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.82	4.89 ng	658836	Perylene-d12	15.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Squalene	410	C30H50	000111-02-4	94
2		7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24O	125529-10-4	72
3		2,6,10,14-Hexadecatetraen-1-ol, ...	332	C22H36O2	061691-98-3	64
4		2,6,10,14-Hexadecatetraenoic aci...	332	C22H36O2	024035-35-6	64
5		4-Oxo-.beta.-damascone	206	C13H18O2	053398-08-6	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120618\
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Sample : J6204-08
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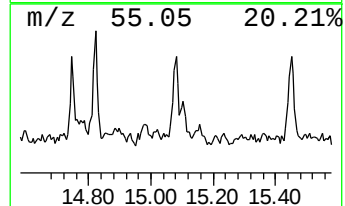
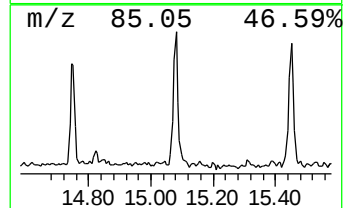
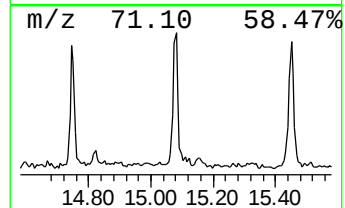
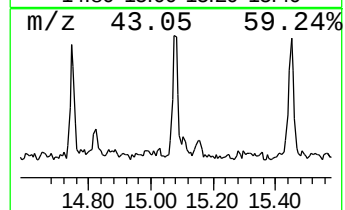
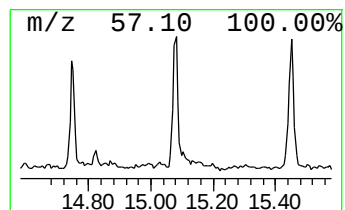
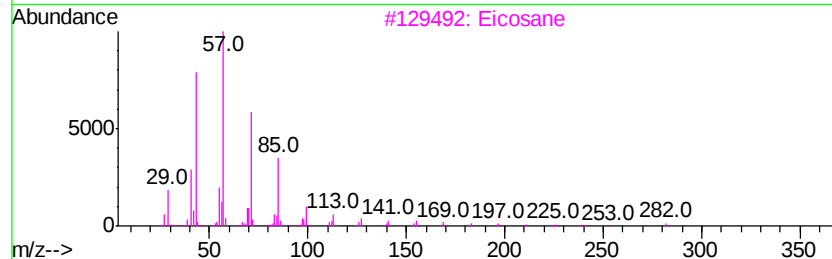
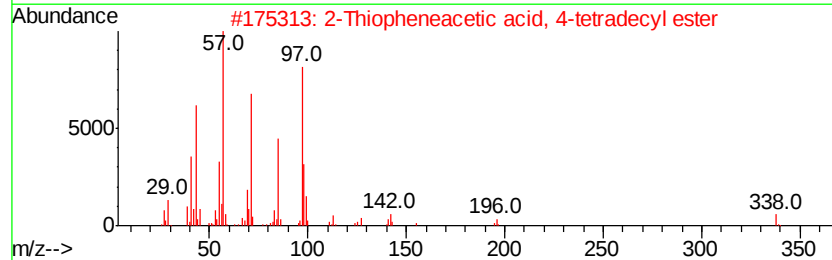
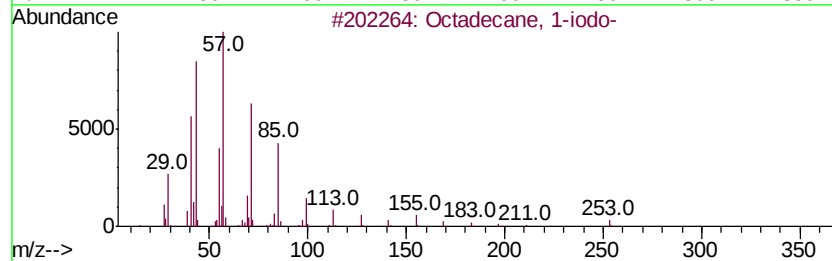
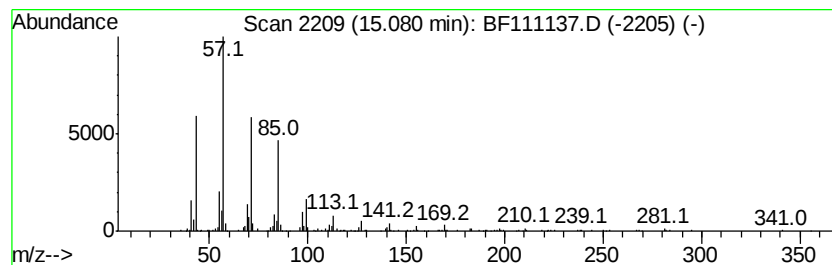
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Octadecane, 1-iodo- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.		
15.08	2.83 ng	381243	Perylene-d12	15.39		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecane, 1-iodo-	380	C18H37I	000629-93-6	87
2		2-Thiopheneacetic acid, 4-tetrad...	338	C20H34O2S	1000280-67-0	86
3		Eicosane	282	C20H42	000112-95-8	86
4		Decane, 1-iodo-	268	C10H21I	002050-77-3	81
5		1-Iodoundecane	282	C11H23I	004282-44-4	81



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120618\
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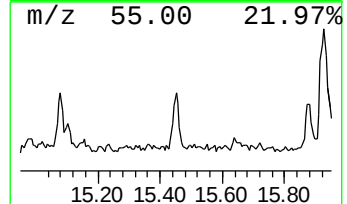
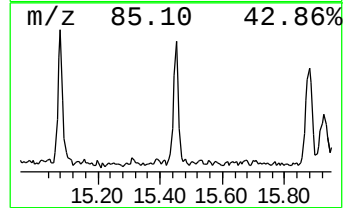
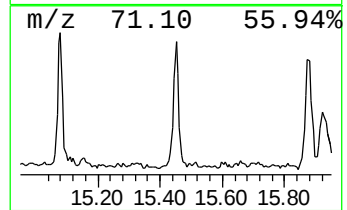
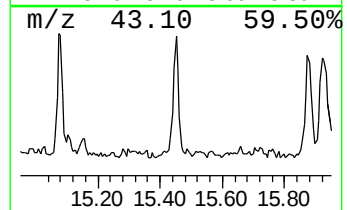
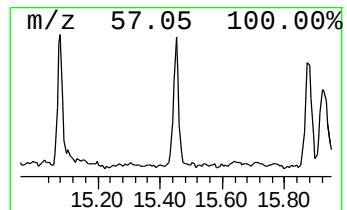
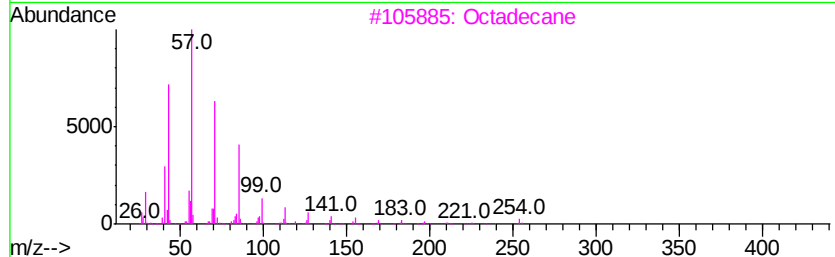
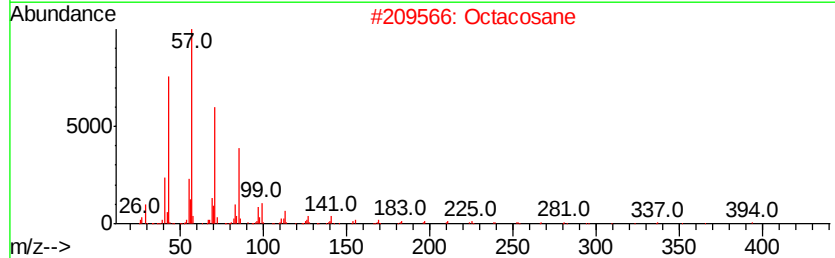
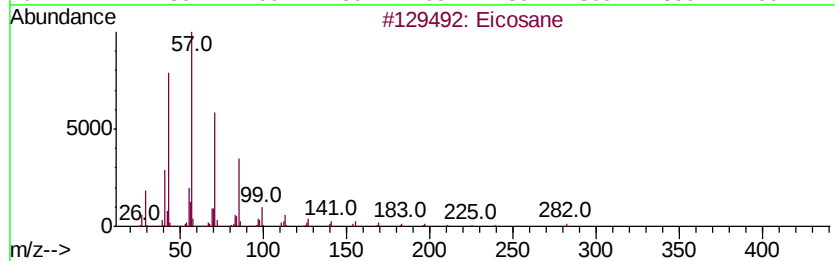
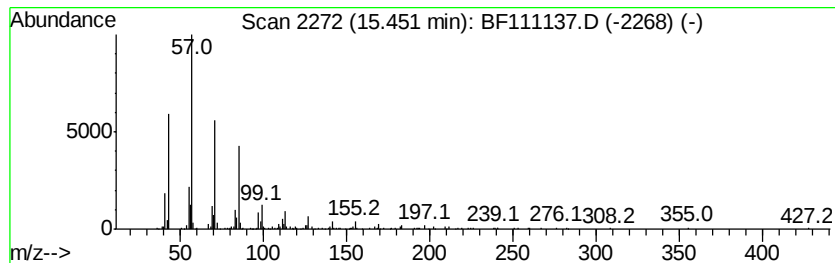
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Eicosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.45	2.75 ng	371081	Perylene-d12	15.39
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Eicosane	282 C20H42	000112-95-8	96
2	Octacosane	394 C28H58	000630-02-4	91
3	Octadecane	254 C18H38	000593-45-3	91
4	2-methyloctacosane	408 C29H60	1000376-72-8	91
5	Pentacosane	352 C25H52	000629-99-2	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120618\
Data File : BF111137.D
Acq On : 6 Dec 2018 14:41
Operator : JU/SJ
Sample : J6204-08
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
HLSP-SB-2B-(2-9.75)

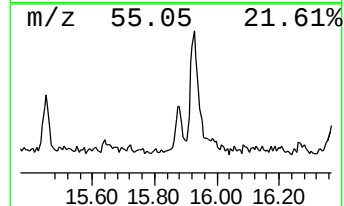
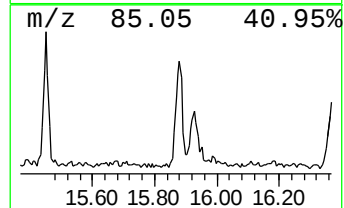
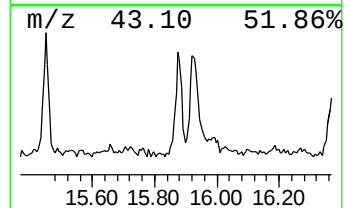
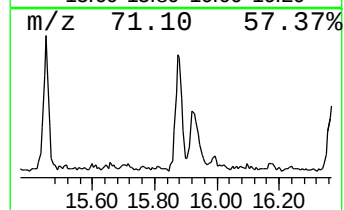
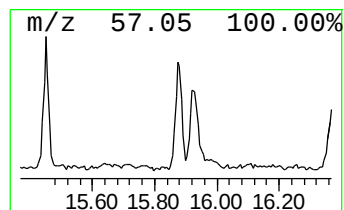
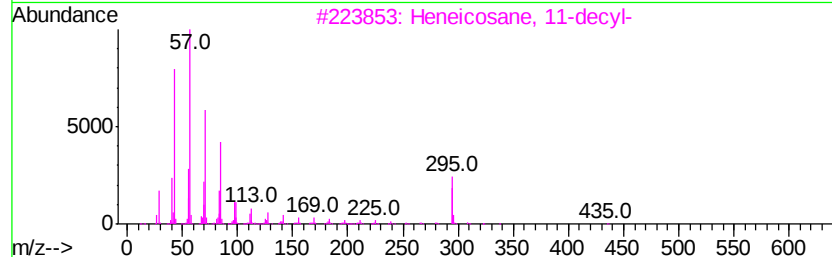
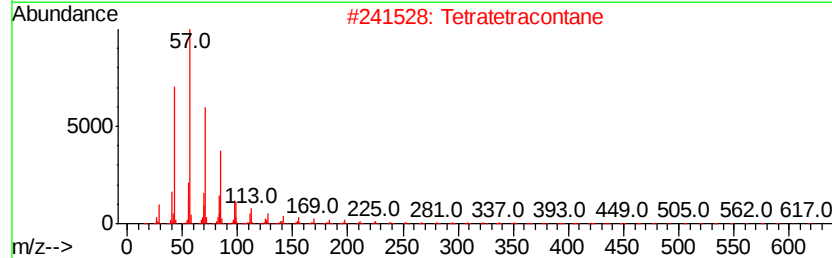
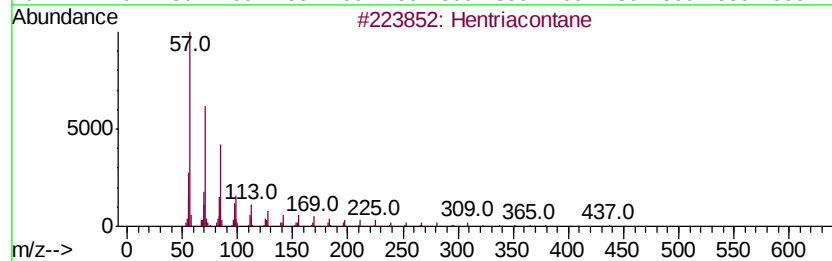
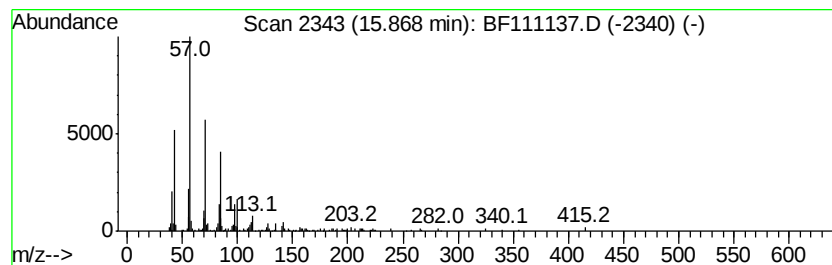
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Hentriacontane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.87	2.66 ng	358705	Perylene-d12	15.39

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hentriacontane	437	C31H64	000630-04-6	91
2			Tetratetracontane	619	C44H90	007098-22-8	91
3			Heneicosane, 11-decyl-	437	C31H64	055320-06-4	90
4			Octacosane	394	C28H58	000630-02-4	90
5			Hexacosane	366	C26H54	000630-01-3	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120618\
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ALS Vial : 9 Sample Multiplier: 1

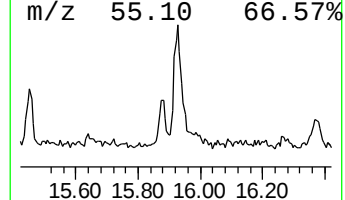
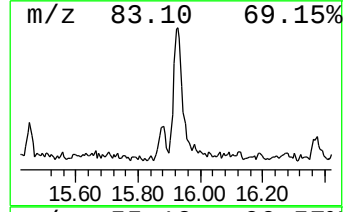
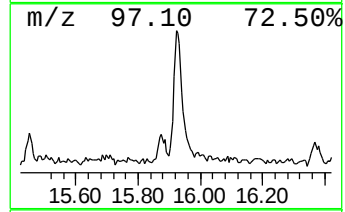
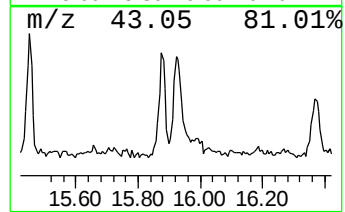
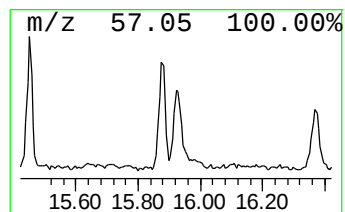
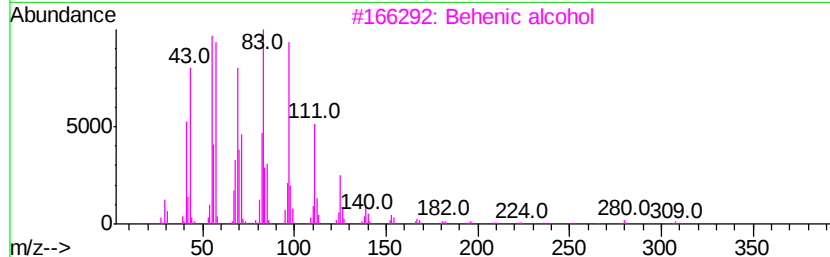
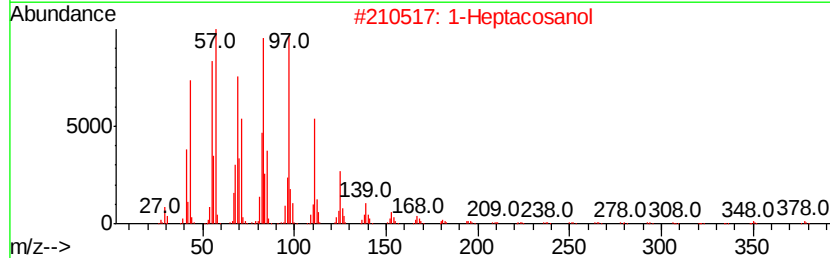
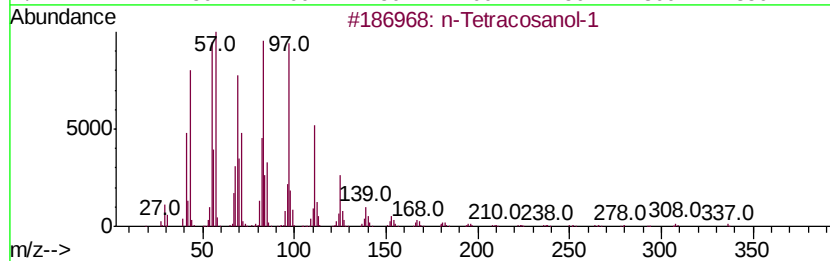
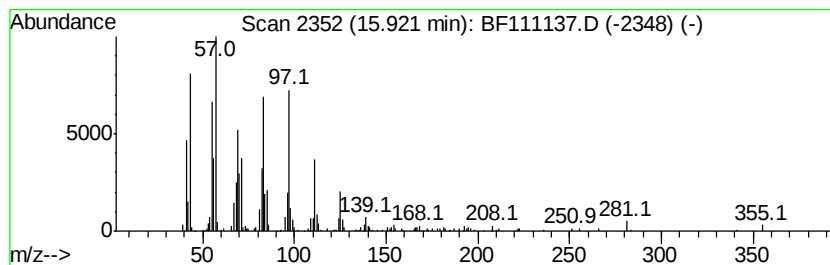
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 n-Tetracosanol-1 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.92	4.94 ng	666067	Perylene-d12	15.39	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Tetracosanol-1	354	C24H50O	000506-51-4	95
2	1-Heptacosanol	396	C27H56O	002004-39-9	95
3	Behenic alcohol	326	C22H46O	000661-19-8	93
4	Octacosyl acetate	452	C30H60O2	018206-97-8	91
5	Octacosanol	410	C28H58O	000557-61-9	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF120618\
 Data File : BF111137.D
 Acq On : 6 Dec 2018 14:41
 Operator : JU/SJ
 Sample : J6204-08
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HLSP-SB-2B-(2-9.75)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF120518.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...	5.03	5.8	ng	879928	1	6.80	3057380	20.0
unknown6.57	6.57	97.1	ng	14846400	1	6.80	3057380	20.0
n-Hexadecanoic acid	11.86	21.0	ng	4571480	4	11.32	4361870	20.0
Octadecanoic acid	12.65	13.3	ng	2443220	5	13.96	3669760	20.0
1-Heneicosanol	13.81	17.2	ng	3156320	5	13.96	3669760	20.0
Sulfurous acid, p...	14.45	2.5	ng	464054	5	13.96	3669760	20.0
Nonadecane	14.74	2.1	ng	287385	6	15.39	2694690	20.0
Squalene	14.82	4.9	ng	658836	6	15.39	2694690	20.0
Octadecane, 1-iodo-	15.08	2.8	ng	381243	6	15.39	2694690	20.0
Eicosane	15.45	2.8	ng	371081	6	15.39	2694690	20.0
Hentriacontane	15.87	2.7	ng	358705	6	15.39	2694690	20.0
n-Tetracosanol-1	15.92	4.9	ng	666067	6	15.39	2694690	20.0