

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF122618\
 Data File : BF111732.D
 Acq On : 26 Dec 2018 23:46
 Operator : JU/SJ
 Sample : J6505-01
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-57(0-12)-COMP

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-BF121918.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	2.251	24	28	34	rVB	45815	64655	1.43%	0.166%
2	5.134	514	518	524	rVB	250735	291138	6.43%	0.749%
3	5.545	583	588	591	rBV	2912701	2764044	61.04%	7.109%
4	6.545	753	758	764	rBV	3011443	3000489	66.26%	7.718%
5	6.681	776	781	784	rBV	3358881	2907889	64.22%	7.479%
6	6.904	815	819	822	rBV	1134857	875000	19.32%	2.251%
7	6.969	826	830	838	rBV	1009154	937062	20.69%	2.410%
8	7.063	842	846	849	rVB	2923259	2389100	52.76%	6.145%
9	7.463	909	914	917	rBV	1973329	1804477	39.85%	4.641%
10	7.669	945	949	952	rBV	1247538	1184670	26.16%	3.047%
11	7.804	968	972	975	rBV	85721	67385	1.49%	0.173%
12	8.186	1031	1037	1041	rBV	1205569	1005178	22.20%	2.585%
13	9.257	1215	1219	1223	rBV	3615853	3134407	69.22%	8.062%
14	9.533	1260	1266	1269	rBV	3154422	2485084	54.88%	6.392%
15	9.645	1282	1285	1289	rBV	248662	210280	4.64%	0.541%
16	9.939	1331	1335	1339	rBV	1080862	980463	21.65%	2.522%
17	10.104	1357	1363	1366	rBV2	35794	57497	1.27%	0.148%
18	10.727	1465	1469	1475	rBV	1929508	1661858	36.70%	4.275%
19	11.139	1535	1539	1549	rVB	333133	446811	9.87%	1.149%
20	11.427	1584	1588	1591	rBV	1080914	904608	19.98%	2.327%
21	11.951	1674	1677	1680	rBV	310853	267581	5.91%	0.688%
22	12.227	1721	1724	1727	rBV	115245	123974	2.74%	0.319%
23	12.280	1731	1733	1735	rBV2	70405	59026	1.30%	0.152%
24	12.468	1761	1765	1768	rBV	5223564	4528234	100.00%	11.647%
25	12.563	1778	1781	1784	rBV	406169	332595	7.34%	0.855%
26	12.733	1808	1810	1813	rBV2	115398	126270	2.79%	0.325%
27	12.951	1844	1847	1850	rBV	541433	491805	10.86%	1.265%
28	13.010	1853	1857	1861	rVB	3176903	2794387	61.71%	7.188%
29	13.321	1908	1910	1914	rVB	720099	578828	12.78%	1.489%
30	13.674	1968	1970	1974	rBV	399562	406174	8.97%	1.045%
31	14.015	2026	2028	2031	rBV	259959	193050	4.26%	0.497%
32	14.045	2031	2033	2034	rVV	349813	275929	6.09%	0.710%
33	14.062	2034	2036	2041	rVB	863778	825773	18.24%	2.124%
34	14.386	2088	2091	2096	rVB	560437	702685	15.52%	1.807%

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Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF121918.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

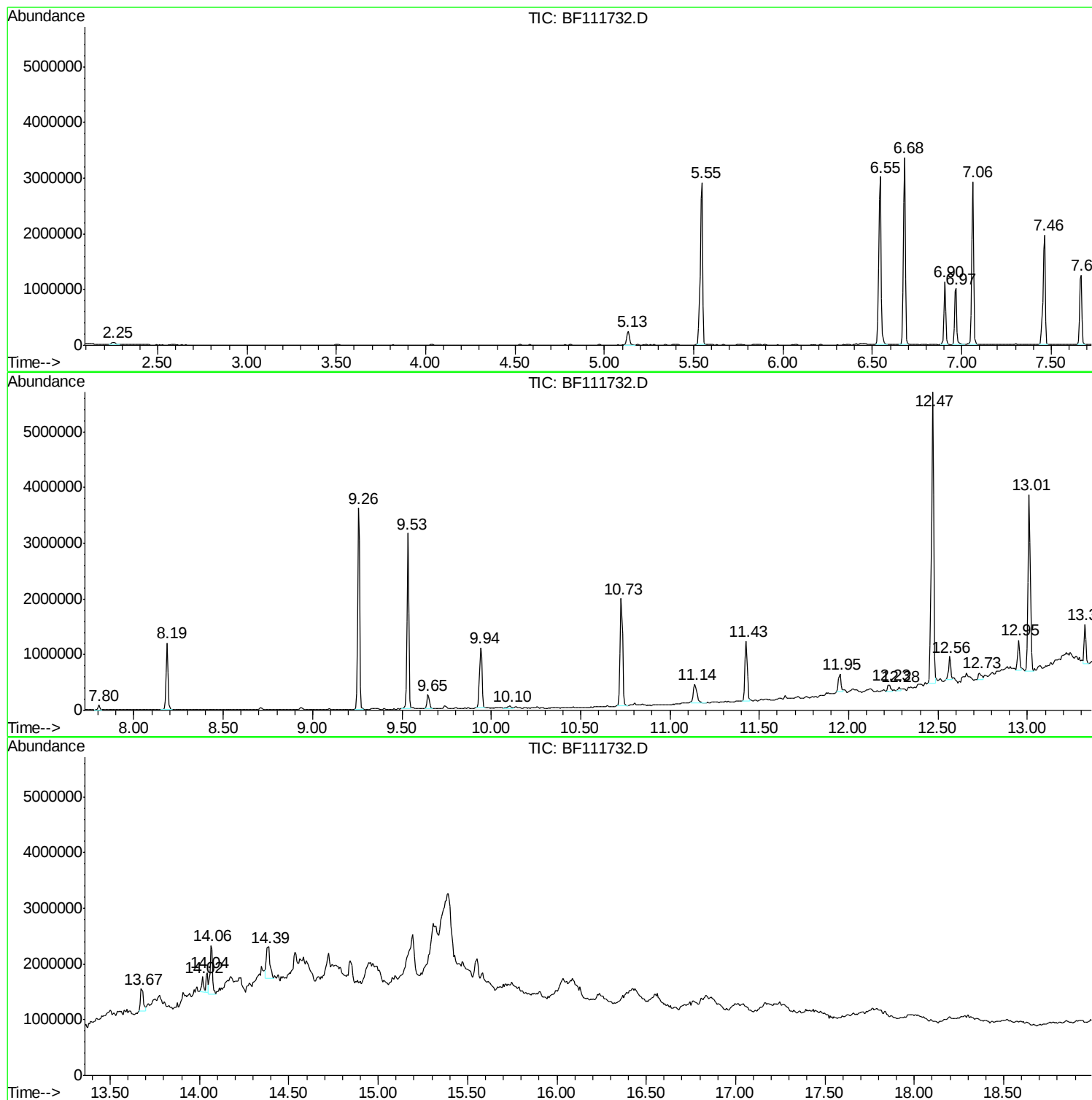
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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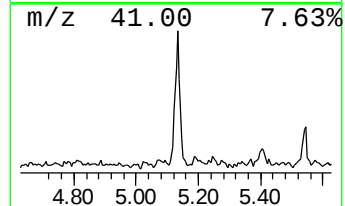
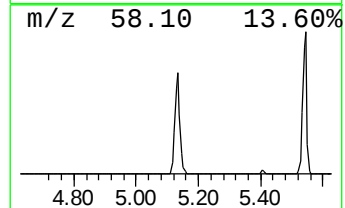
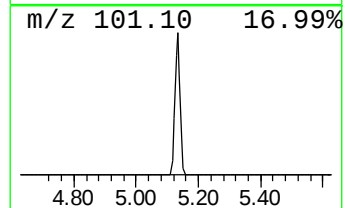
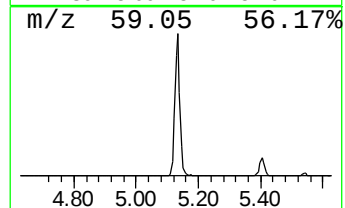
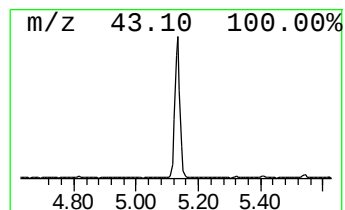
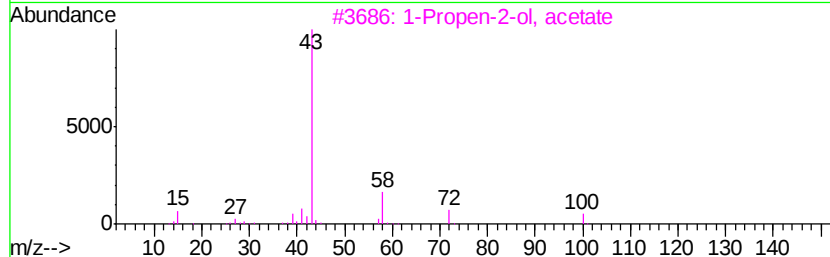
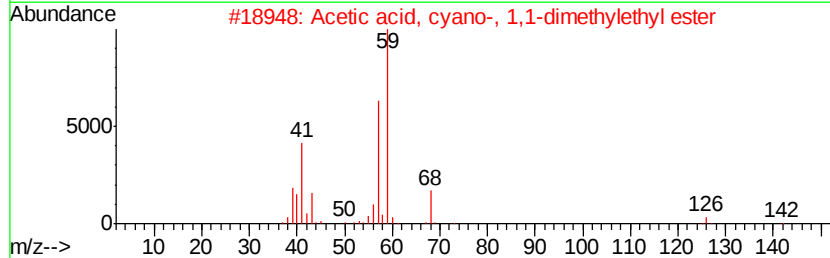
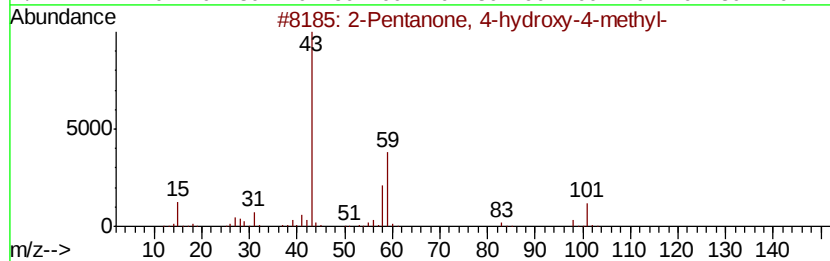
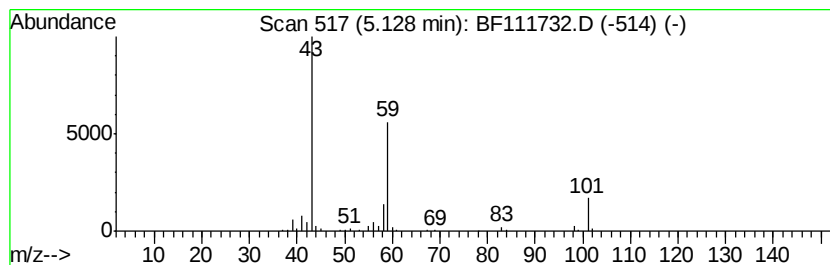
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.13	6.65 ng	291138	1,4-Dichlorobenzene-d4	6.90

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	32
3			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
4			Butane	58	C4H10	000106-97-8	9
5			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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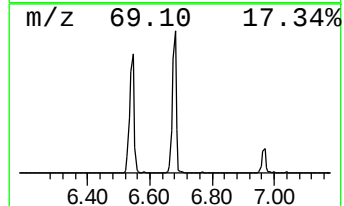
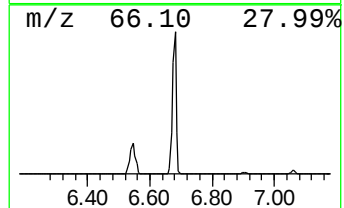
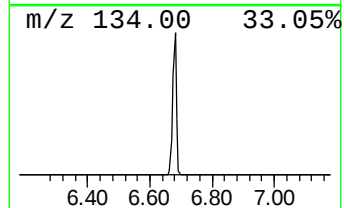
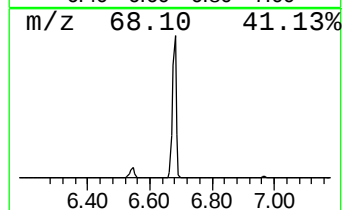
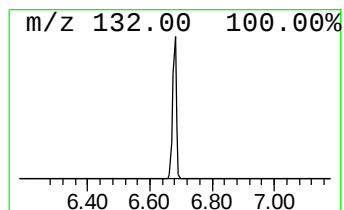
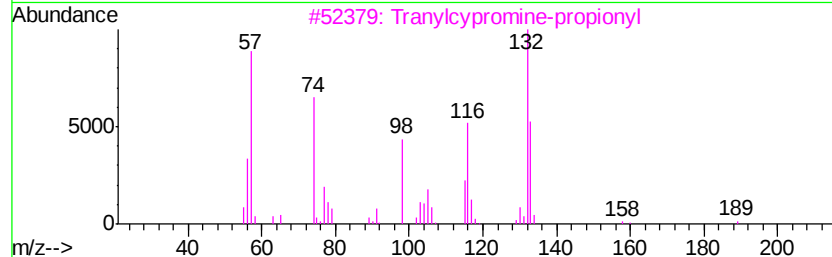
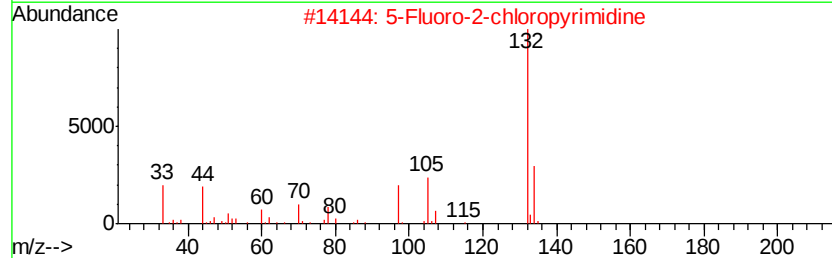
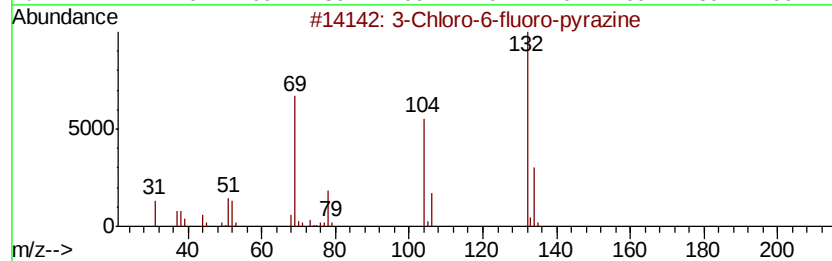
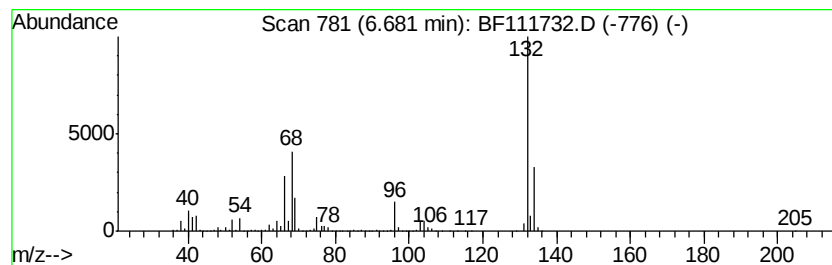
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.68 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.68	66.47 ng	2907890	1,4-Dichlorobenzene-d4	6.90

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		Tranvlcypramine-propionyl	189	C12H15NO	1000123-86-3	12
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



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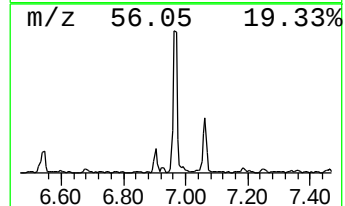
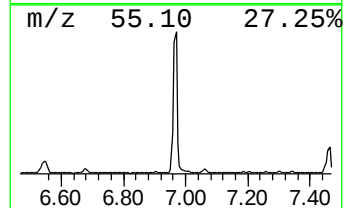
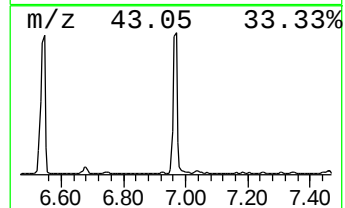
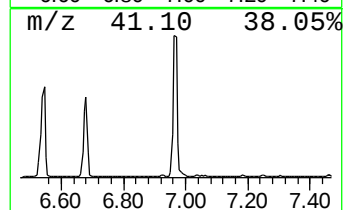
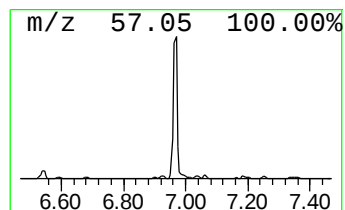
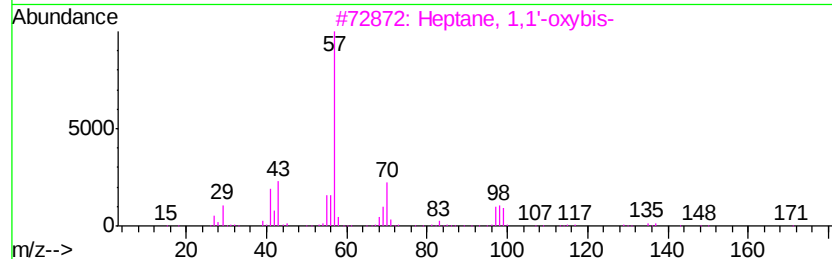
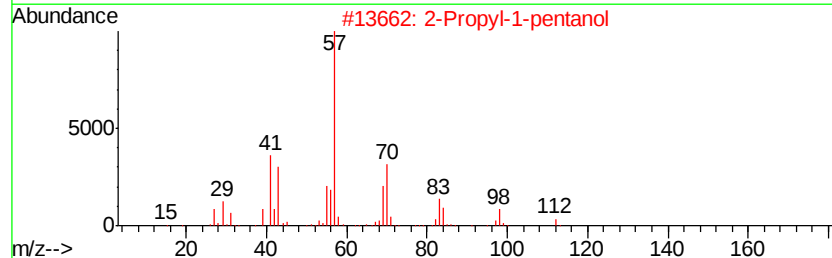
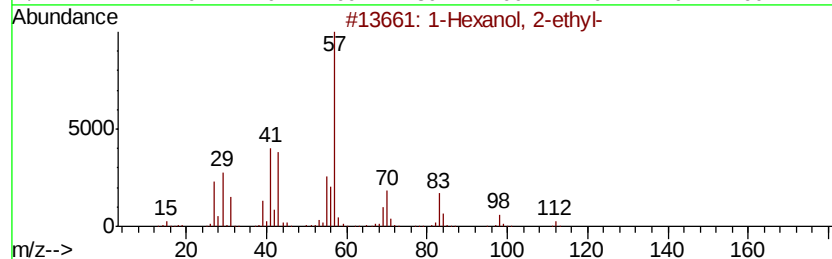
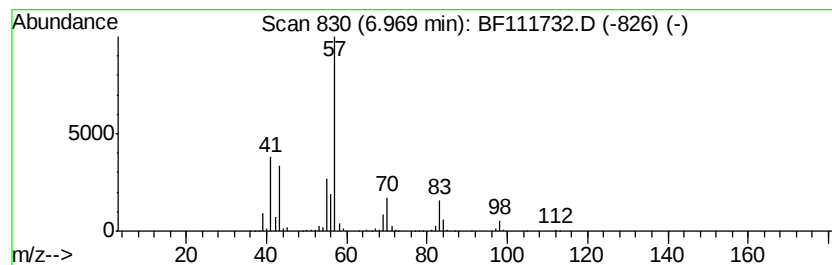
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 1-Hexanol, 2-ethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.97	21.42 ng	937062	1,4-Dichlorobenzene-d4	6.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	83
2		2-Propyl-1-pentanol	130	C8H18O	058175-57-8	64
3		Heptane, 1,1'-oxybis-	214	C14H30O	000629-64-1	59
4		1-Isobutoxy-2-ethylhexane	186	C12H26O	1000139-90-3	56
5		1-Pentene, 2,4,4-trimethyl-	112	C8H16	000107-39-1	43



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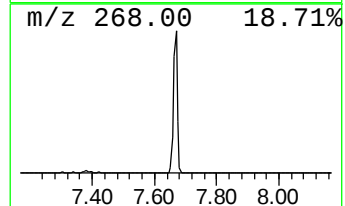
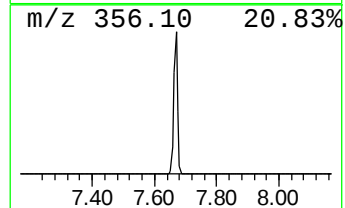
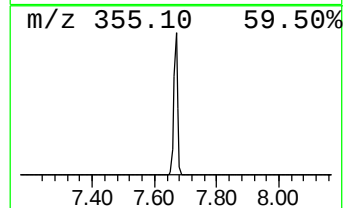
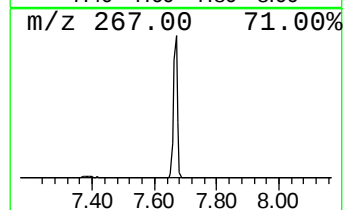
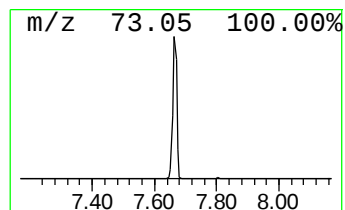
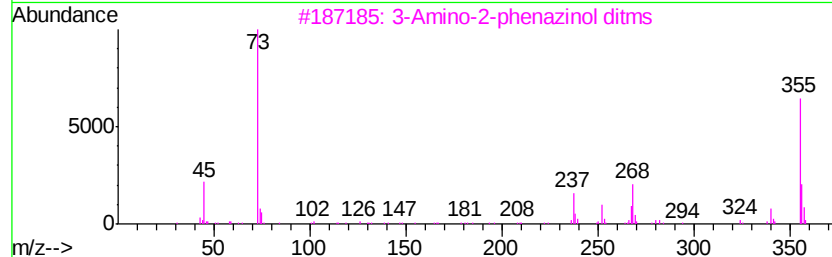
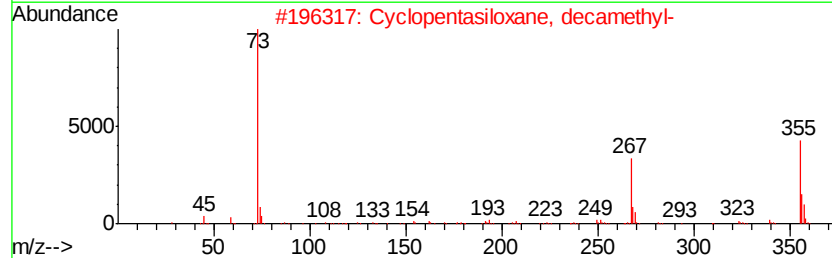
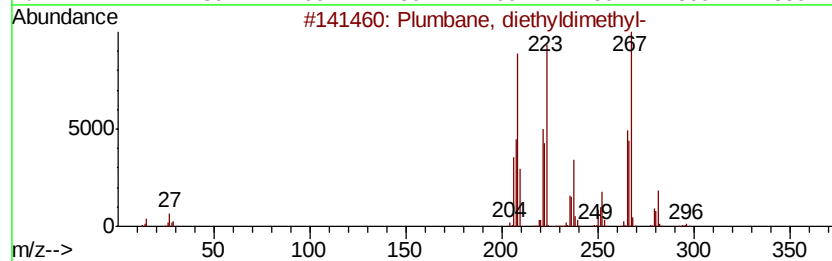
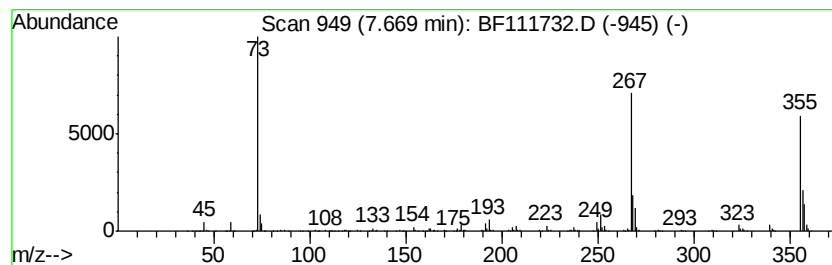
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Plumbane, diethyldimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.67	23.57 ng	1184670	Naphthalene-d8	8.19

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Plumbane, diethyldimethyl-	296	C6H16Pb	001762-27-2	92
2		Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	91
3		3-Amino-2-phenazinol ditms	355	C18H25N3O5Si2	1000332-15-5	52
4		Terbutaline, N-trifluoroacetyl-0...	537	C23H42F3N04Si3	325836-92-8	50
5		N-(Trifluoroacetyl)-0,0',0''-tris...	495	C20H36F3N04Si3	054135-51-2	47



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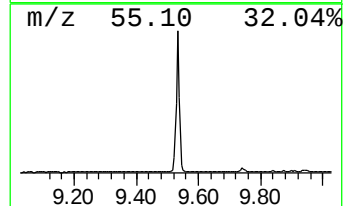
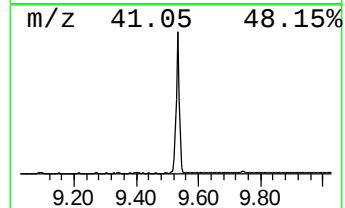
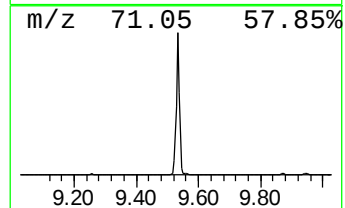
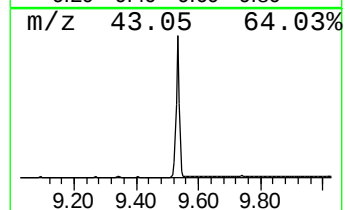
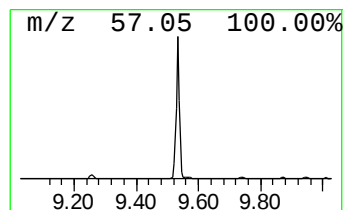
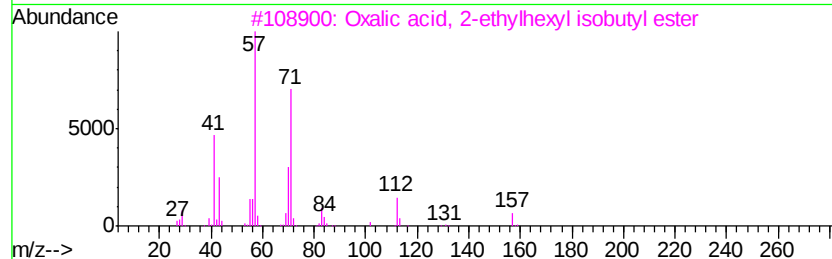
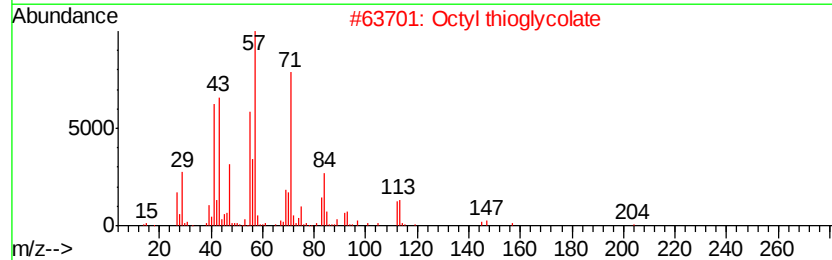
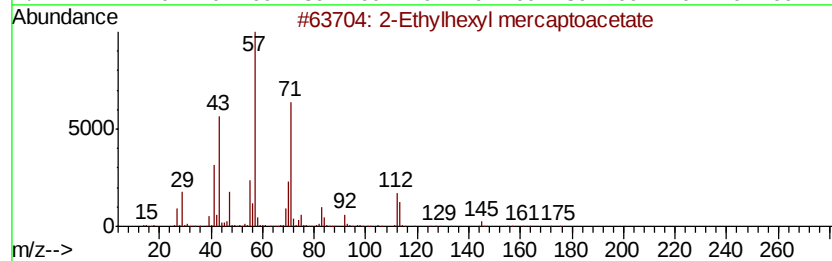
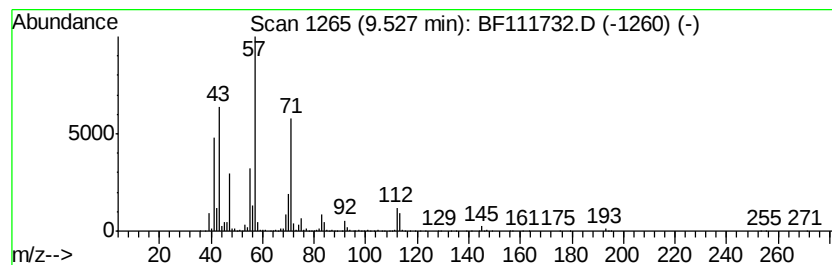
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TIC Library : C:\DATABASE\NIST11.L
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 Peak Number 6 2-Ethylhexyl mercaptoacetate Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.53	50.69 ng	2485080	Acenaphthene-d10	9.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Ethylhexyl mercaptoacetate	204	C10H20O2S	007659-86-1	80
2		Octyl thio glycolate	204	C10H20O2S	007664-80-4	64
3		Oxalic acid, 2-ethylhexyl isobut...	258	C14H26O4	1000309-38-5	50
4		Decane, 4-methyl-	156	C11H24	002847-72-5	50
5		Nonane, 4-methyl-5-propyl-	184	C13H28	062185-55-1	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122618\
Data File : BF111732.D
Acq On : 26 Dec 2018 23:46
Operator : JU/SJ
Sample : J6505-01
Misc :
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Instrument :
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ClientSampleId :
SB-57(0-12)-COMP

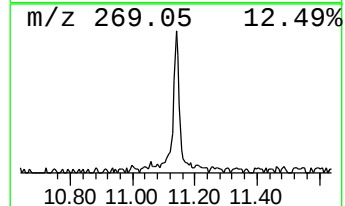
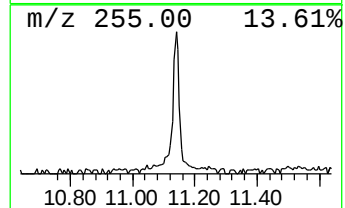
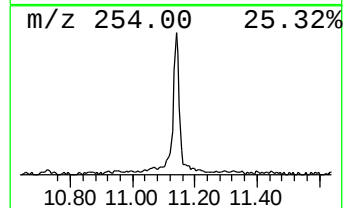
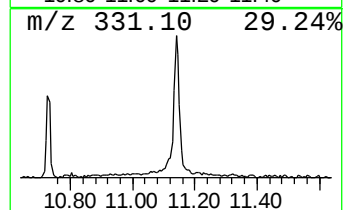
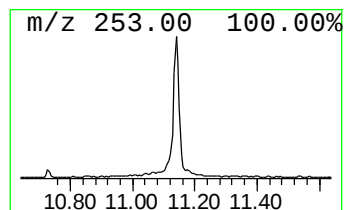
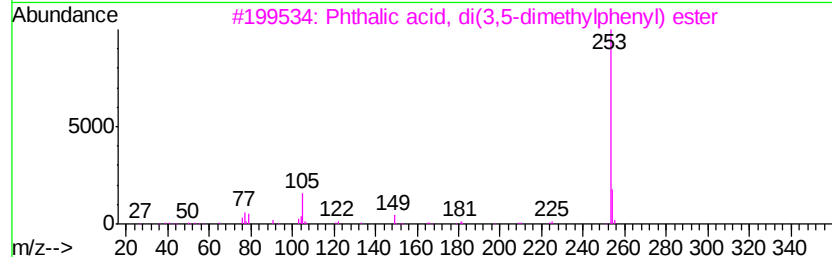
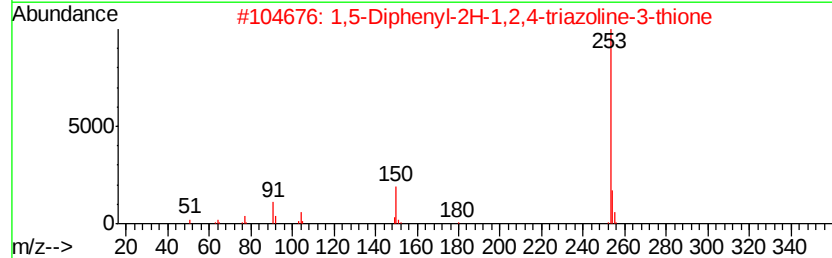
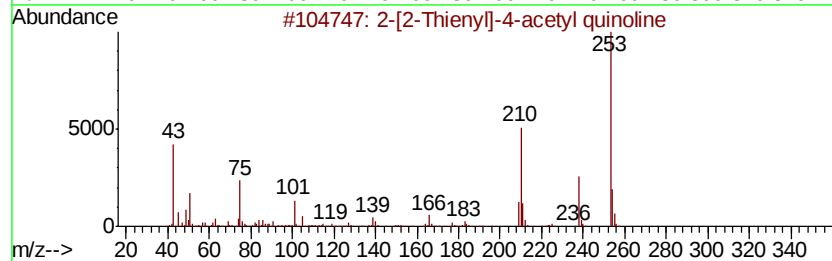
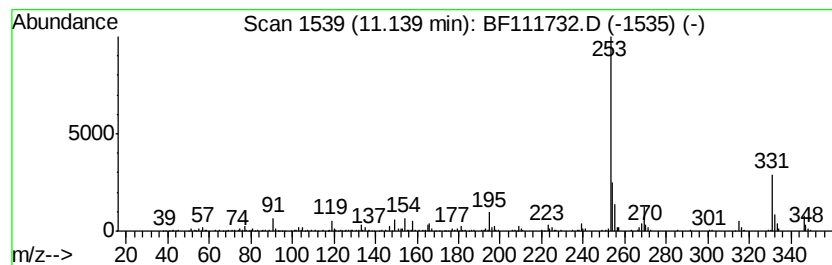
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 2-[2-Thienyl]-4-acetyl quin... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.14	9.88 ng	446811	Phenanthrene-d10	11.43

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-[2-Thienyl]-4-acetyl quinoline	253	C15H11NOS	027302-83-6	50
2		1,5-Diphenyl-2H-1,2,4-triazoline...	253	C14H11N3S	005055-74-3	47
3		Phthalic acid, di(3,5-dimethylph...	374	C24H22O4	1000309-79-1	38
4		Phthalic acid, 3,5-dimethylpheny...	374	C23H18O5	1000315-73-8	38
5		6-Methyl-2-(3-nitrophenyl)imidaz...	253	C14H11N3O2	1000224-30-4	37



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122618\
Data File : BF111732.D
Acq On : 26 Dec 2018 23:46
Operator : JU/SJ
Sample : J6505-01
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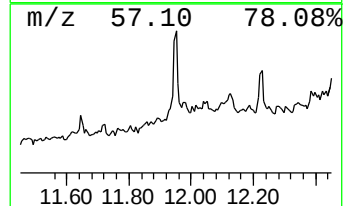
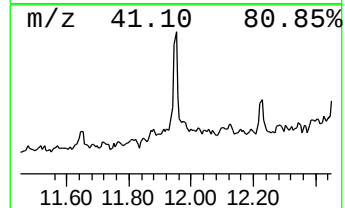
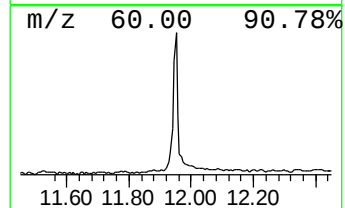
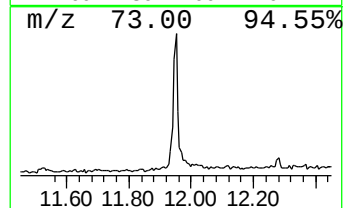
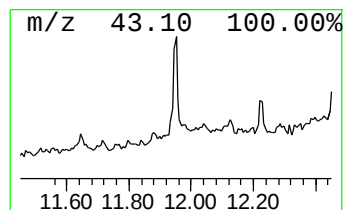
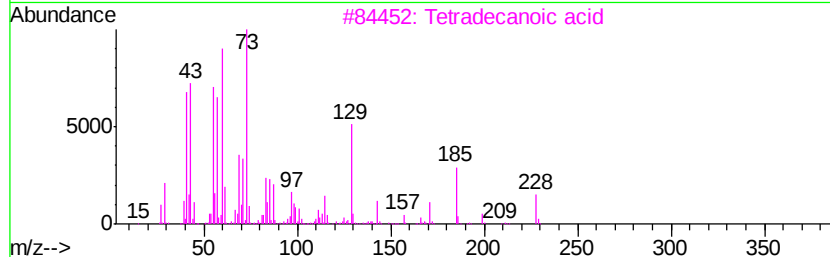
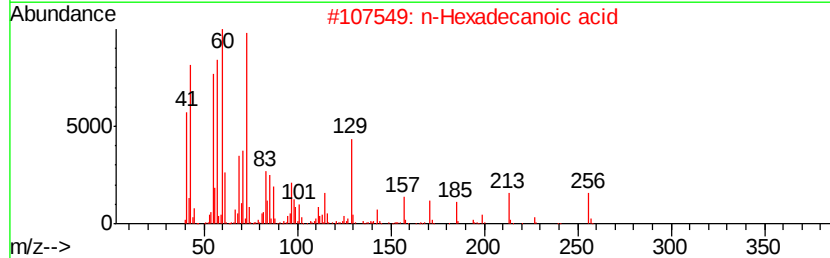
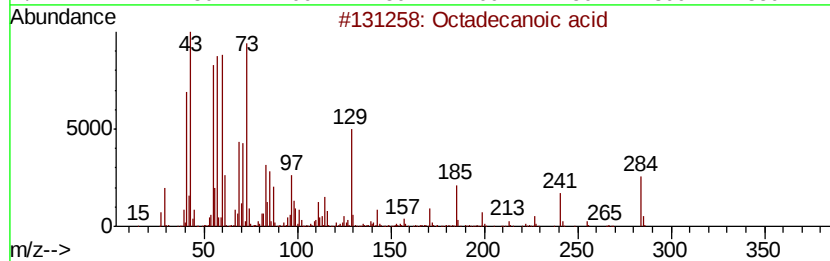
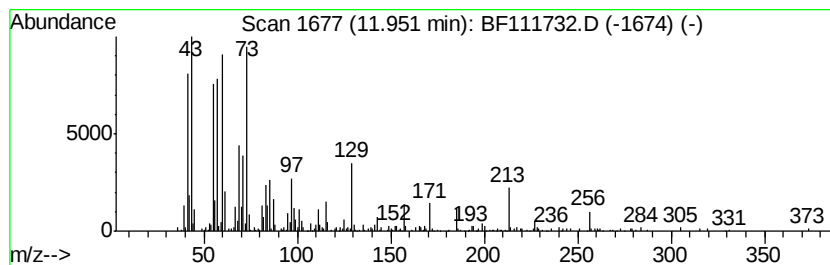
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ClientSampleId :
SB-57(0-12)-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Octadecanoic acid Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.95	5.92 ng	267581	Phenanthrene-d10	11.43	
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	98
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	96
4	Dodecanoic acid	200	C12H24O2	000143-07-7	93
5	Pentadecanoic acid	242	C15H30O2	001002-84-2	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122618\
Data File : BF111732.D
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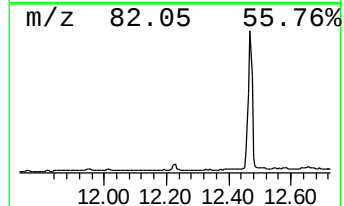
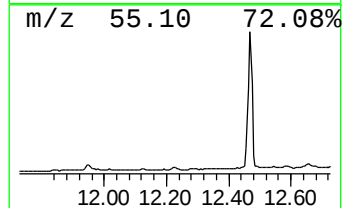
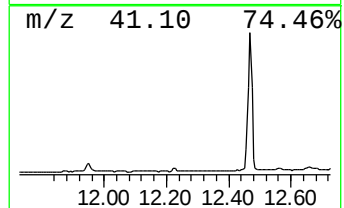
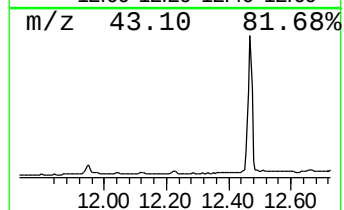
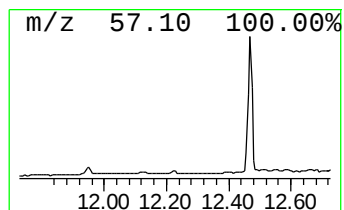
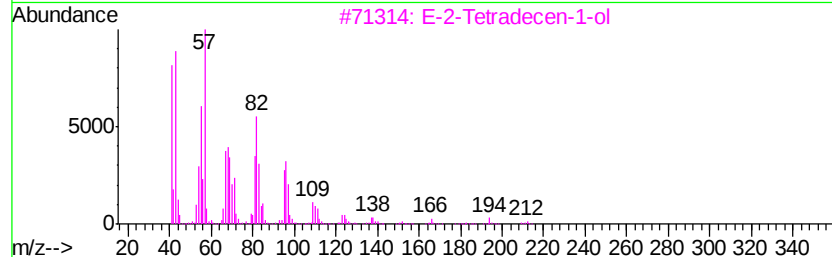
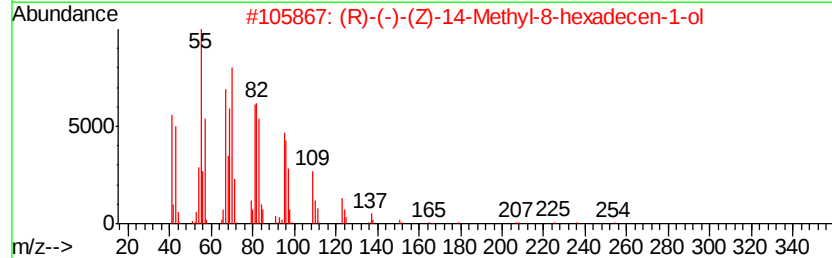
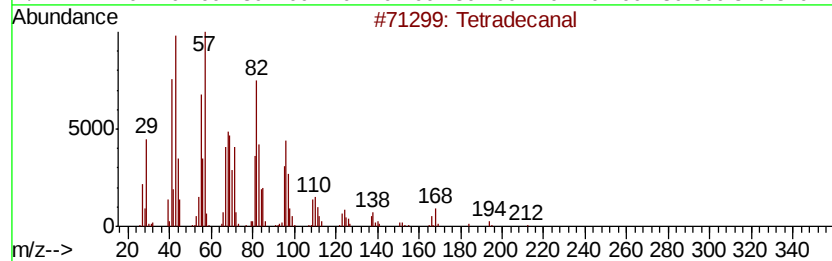
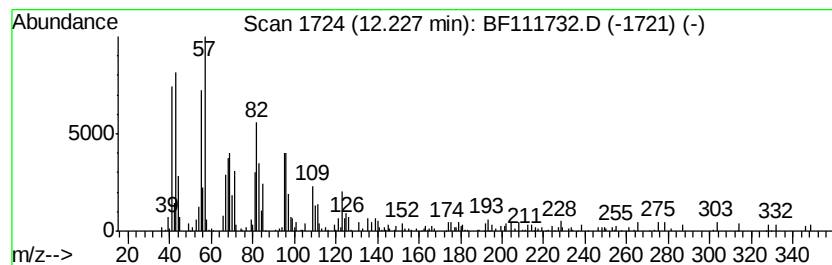
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Tetradecanal Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.23	2.74 ng	123974	Phenanthrene-d10	11.43	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecanal	212	C14H28O	000124-25-4	90
2	(R)-(-)-(Z)-14-Methyl-8-hexadec...	254	C17H34O	030689-78-2	86
3	E-2-Tetradecen-1-ol	212	C14H28O	1000130-83-7	70
4	1,12-Tridecadiene	180	C13H24	021964-48-7	58
5	Z,Z-2,15-Octadecediene-1-ol acetate	308	C20H36O2	1000130-95-1	58



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122618\
Data File : BF111732.D
Acq On : 26 Dec 2018 23:46
Operator : JU/SJ
Sample : J6505-01
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ALS Vial : 14 Sample Multiplier: 1

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ClientSampleId :
SB-57(0-12)-COMP

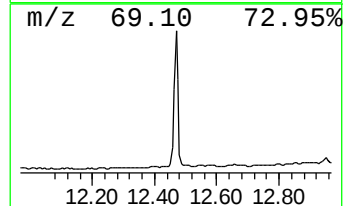
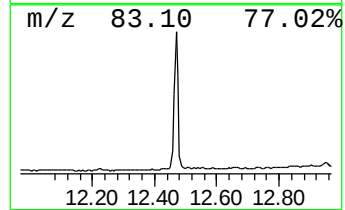
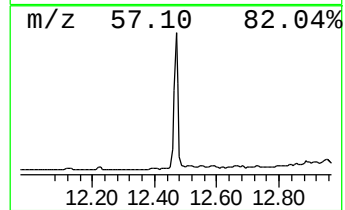
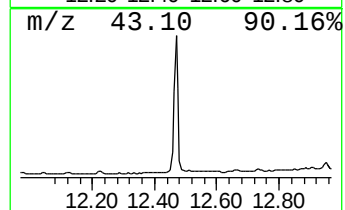
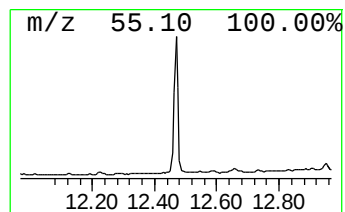
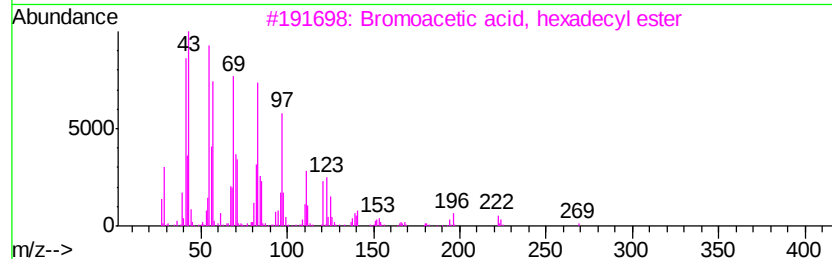
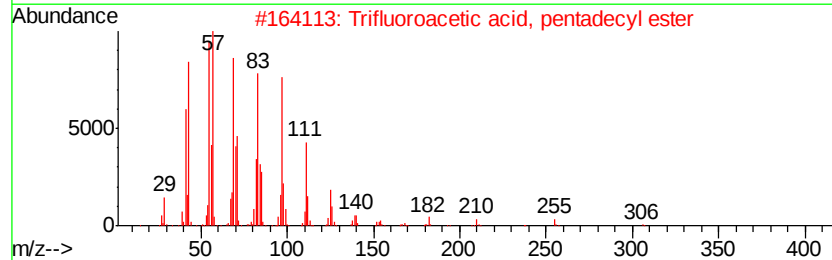
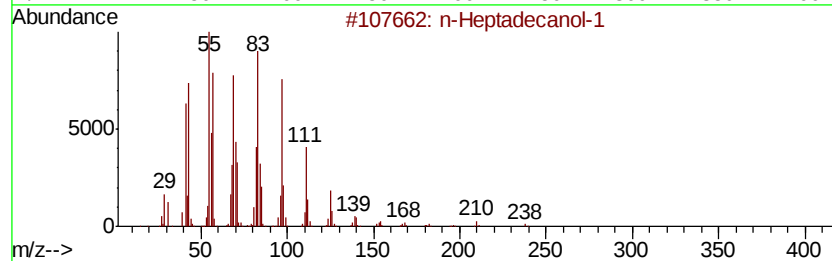
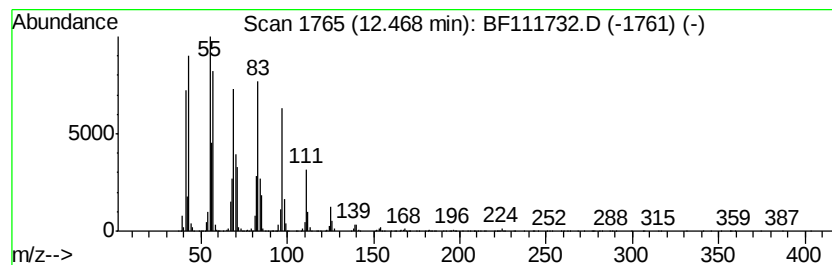
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 n-Heptadecanol-1 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.47	100.12 ng	4528230	Phenanthrene-d10	11.43

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Heptadecanol-1	256	C17H36O	001454-85-9	95
2		Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	94
3		Bromoacetic acid, hexadecyl ester	362	C18H35BrO2	005454-48-8	94
4		1-Hexadecanol	242	C16H34O	036653-82-4	94
5		Pentafluoropropionic acid, tride...	346	C16H27F5O2	959261-22-4	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122618\
Data File : BF111732.D
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Operator : JU/SJ
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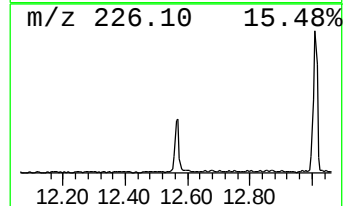
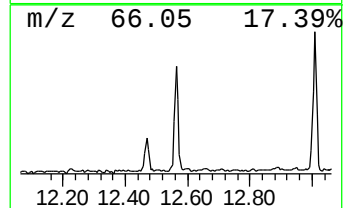
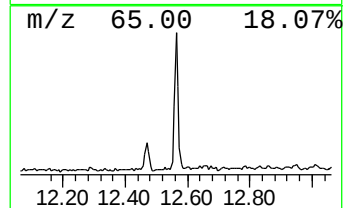
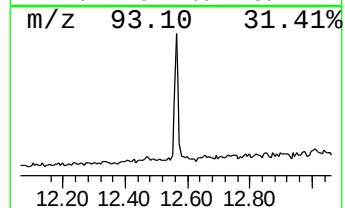
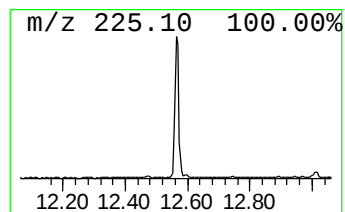
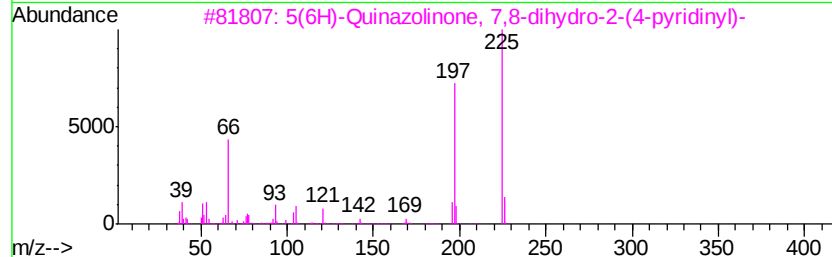
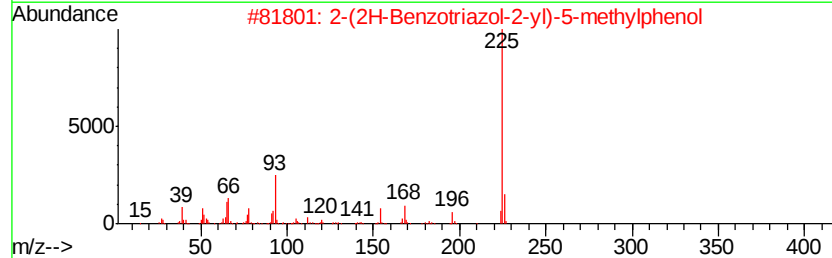
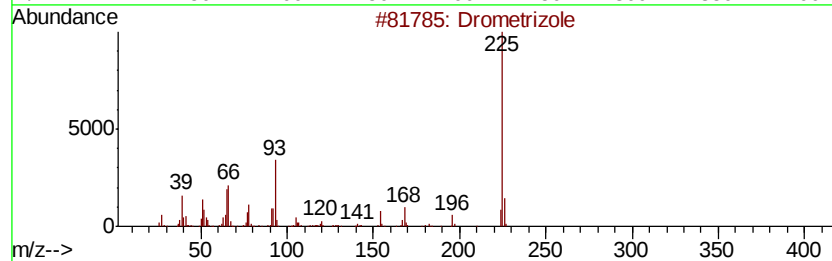
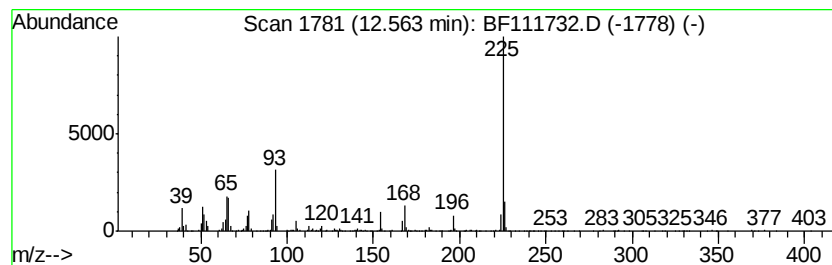
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Drometrizole Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.		
12.56	7.35 ng	332595	Phenanthrene-d10	11.43		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Drometrizole	225	C13H11N3O	002440-22-4	98
2		2-(2H-Benzotriazol-2-yl)-5-methy...	225	C13H11N3O	004998-48-5	98
3		5(6H)-Quinazolinone, 7,8-dihydro...	225	C13H11N3O	1000337-49-5	50
4		Phenazine, 2-amino-8-methoxy-	225	C13H11N3O	018450-21-0	50
5		2H-Benzotriazole, 2-(2-butoxy-5-...	281	C17H19N3O	1000262-87-5	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122618\
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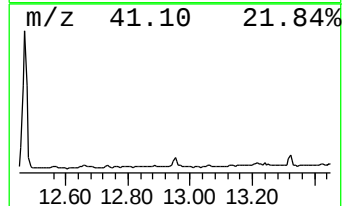
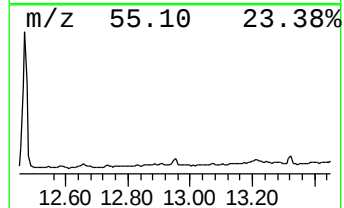
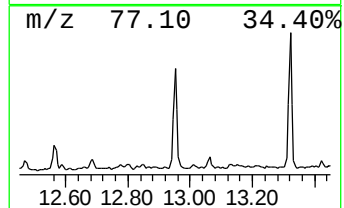
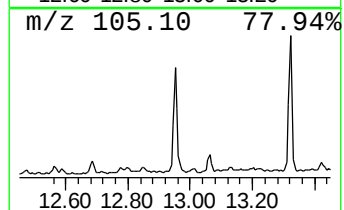
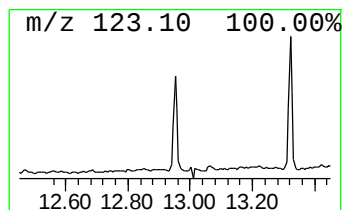
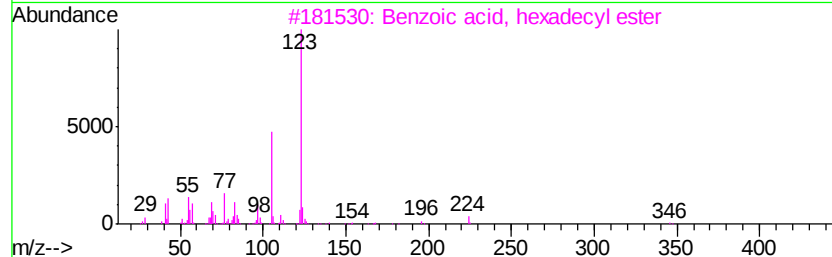
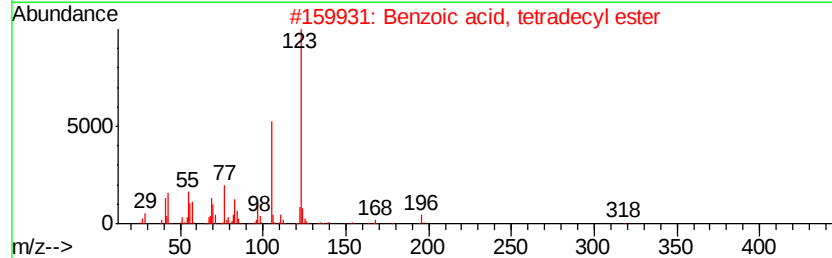
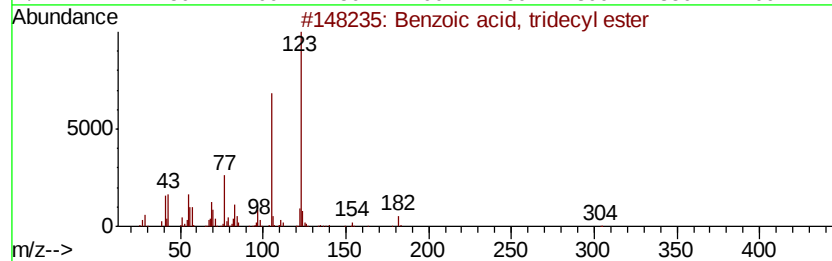
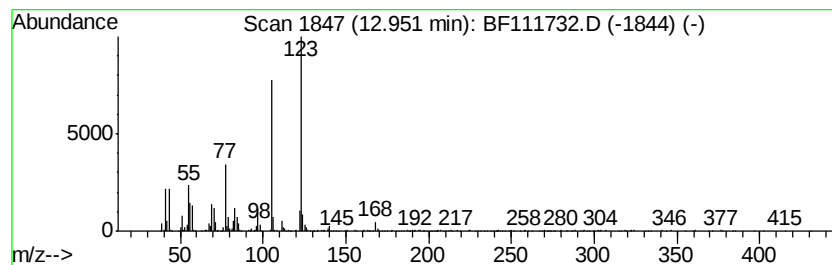
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Benzoic acid, tridecyl ester Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.95	11.91 ng	491805	Chrysene-d12	14.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzoic acid, tridecyl ester	304	C20H32O2	1000340-22-6	98
2		Benzoic acid, tetradecyl ester	318	C21H34O2	1000340-22-7	98
3		Benzoic acid, hexadecyl ester	346	C23H38O2	1000340-22-9	93
4		Benzoic acid, pentadecyl ester	332	C22H36O2	1000340-22-8	90
5		Benzoic acid, octadecyl ester	374	C25H42O2	1000340-23-1	74



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122618\
Data File : BF111732.D
Acq On : 26 Dec 2018 23:46
Operator : JU/SJ
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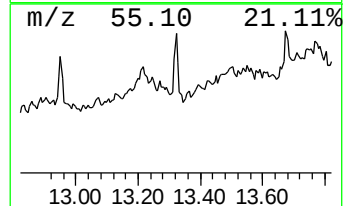
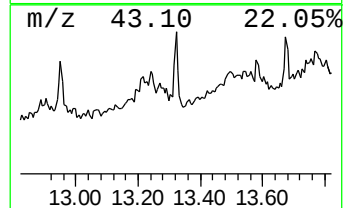
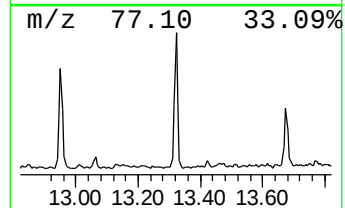
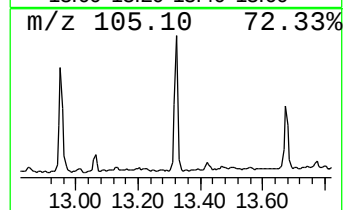
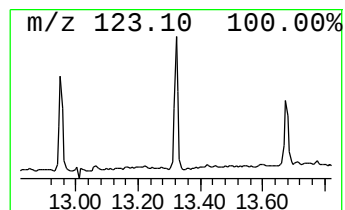
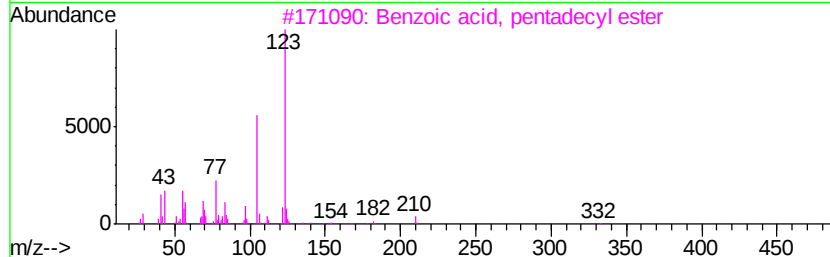
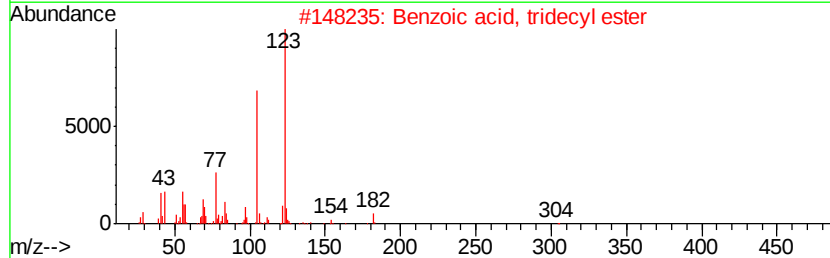
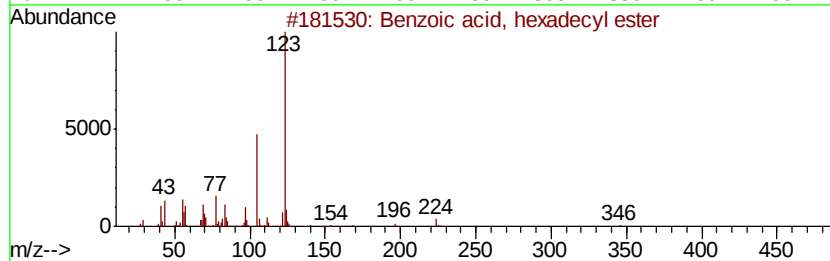
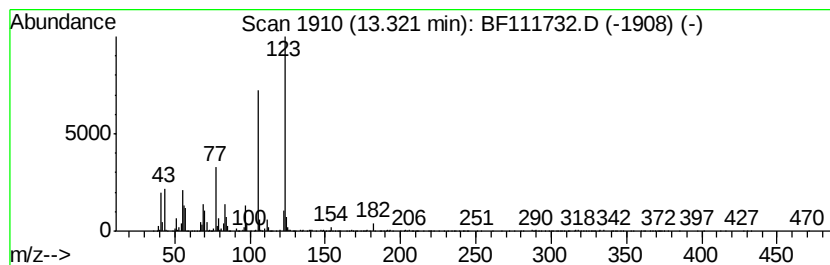
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Benzoic acid, hexadecyl ester Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.32	14.02 ng	578828	Chrysene-d12	14.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzoic acid, hexadecyl ester	346	C23H38O2	1000340-22-9	95
2		Benzoic acid, tridecyl ester	304	C20H32O2	1000340-22-6	90
3		Benzoic acid, pentadecyl ester	332	C22H36O2	1000340-22-8	83
4		Benzoic acid, octadecyl ester	374	C25H42O2	1000340-23-1	83
5		Benzoic acid, tetradecyl ester	318	C21H34O2	1000340-22-7	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122618\
Data File : BF111732.D
Acq On : 26 Dec 2018 23:46
Operator : JU/SJ
Sample : J6505-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-57(0-12)-COMP

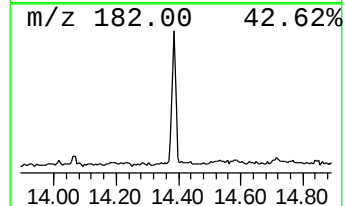
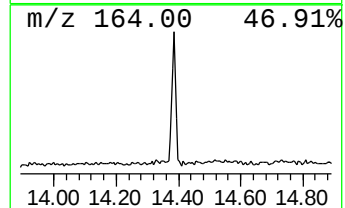
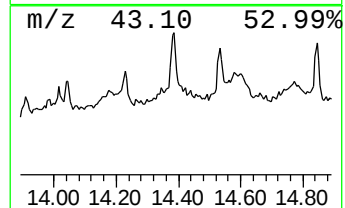
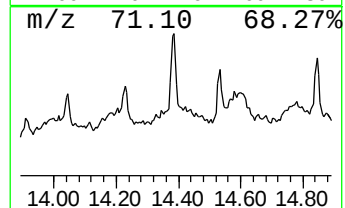
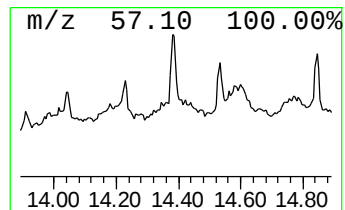
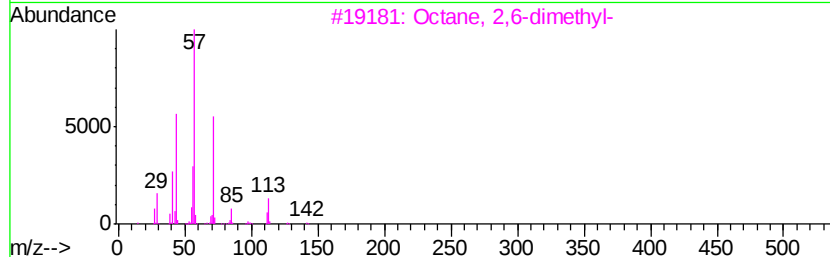
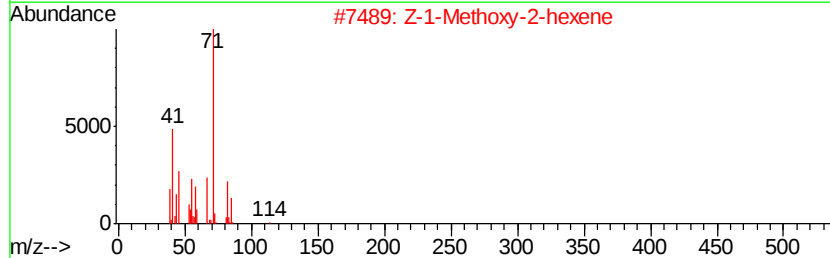
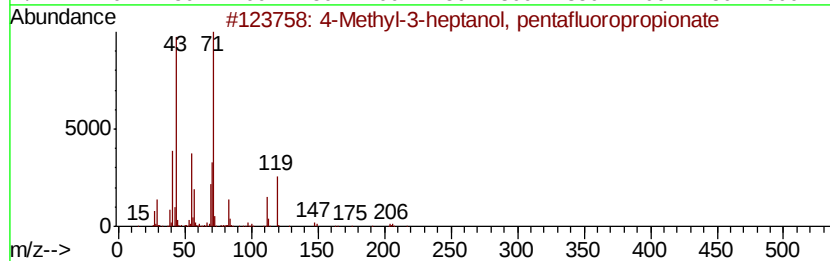
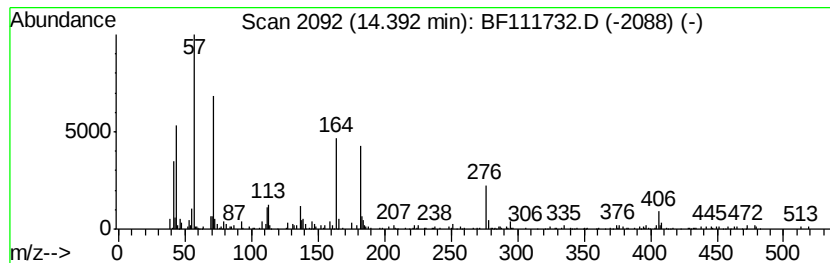
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 unknown14.39 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.39	17.02 ng	702685	Chrysene-d12	14.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Methyl-3-heptanol, pentafluoro...	276	C11H17F5O2	1000365-51-7	14
2		Z-1-Methoxy-2-hexene	114	C7H14O	1000279-60-9	14
3		Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	14
4		Nonane, 3-methyl-	142	C10H22	005911-04-6	14
5		Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	14



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF122618\
 Data File : BF111732.D
 Acq On : 26 Dec 2018 23:46
 Operator : JU/SJ
 Sample : J6505-01
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-57(0-12)-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF121918.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...	5.13	6.7	ng	291138	1	6.90	875000	20.0
unknown6.68	6.68	66.5	ng	2907890	1	6.90	875000	20.0
1-Hexanol, 2-ethyl-	6.97	21.4	ng	937062	1	6.90	875000	20.0
Plumbane, diethyl...	7.67	23.6	ng	1184670	2	8.19	1005180	20.0
2-Ethylhexyl merc...	9.53	50.7	ng	2485080	3	9.94	980463	20.0
2-[2-Thienyl]-4-a...	11.14	9.9	ng	446811	4	11.43	904608	20.0
Octadecanoic acid	11.95	5.9	ng	267581	4	11.43	904608	20.0
Tetradecanal	12.23	2.7	ng	123974	4	11.43	904608	20.0
n-Heptadecanol-1	12.47	100.1	ng	4528230	4	11.43	904608	20.0
Drometrizole	12.56	7.3	ng	332595	4	11.43	904608	20.0
Benzoic acid, tri...	12.95	11.9	ng	491805	5	14.06	825773	20.0
Benzoic acid, hex...	13.32	14.0	ng	578828	5	14.06	825773	20.0
unknown14.39	14.39	17.0	ng	702685	5	14.06	825773	20.0