

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF062918\
 Data File : BF106970.D
 Acq On : 29 Jun 2018 13:59
 Operator : JU/SJ
 Sample : J3709-05
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EME-FW-TS005-01

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.195	482	486	500	rBB	129235	158505	10.47%	0.969%
2	5.595	550	554	565	rBV	1453840	1362269	89.98%	8.328%
3	6.225	659	661	664	rBV	21865	17535	1.16%	0.107%
4	6.595	720	724	731	rBV	1288508	1414218	93.41%	8.646%
5	6.731	743	747	750	rBV	1612295	1432404	94.61%	8.757%
6	6.954	781	785	788	rBV	473565	411322	27.17%	2.515%
7	7.107	807	811	815	rVB	1307554	1139810	75.29%	6.968%
8	7.519	876	881	884	rBV	975001	904701	59.76%	5.531%
9	8.236	999	1003	1015	rBV	575918	520502	34.38%	3.182%
10	9.307	1181	1185	1188	rBV	1703586	1513971	100.00%	9.256%
11	9.701	1248	1252	1257	rBV	183275	155927	10.30%	0.953%
12	9.901	1281	1286	1288	rBV2	31061	32378	2.14%	0.198%
13	9.995	1298	1302	1308	rBV	645876	585894	38.70%	3.582%
14	10.101	1317	1320	1323	rVB	38207	27417	1.81%	0.168%
15	10.401	1369	1371	1374	rVB	49777	33883	2.24%	0.207%
16	10.454	1378	1380	1384	rVB	21301	17551	1.16%	0.107%
17	10.619	1403	1408	1411	rBV	13588	17729	1.17%	0.108%
18	10.742	1426	1429	1433	rBV4	12673	18597	1.23%	0.114%
19	10.789	1433	1437	1440	rBV	1086712	967748	63.92%	5.916%
20	10.872	1448	1451	1460	rVB3	38930	53157	3.51%	0.325%
21	11.001	1468	1473	1475	rBV2	20957	24358	1.61%	0.149%
22	11.324	1525	1528	1530	rBV	21877	18170	1.20%	0.111%
23	11.489	1552	1556	1558	rBV	577653	548457	36.23%	3.353%
24	11.513	1558	1560	1563	rVB2	65795	60581	4.00%	0.370%
25	11.671	1584	1587	1593	rBV3	14158	17690	1.17%	0.108%
26	12.495	1724	1727	1731	rBV	21088	21181	1.40%	0.129%
27	12.548	1732	1736	1738	rVB4	28359	29639	1.96%	0.181%
28	12.707	1759	1763	1766	rBV	120559	114871	7.59%	0.702%
29	12.936	1799	1802	1807	rVB	111088	110024	7.27%	0.673%
30	13.071	1820	1825	1828	rBV	1358334	1292828	85.39%	7.904%
31	13.260	1855	1857	1863	rVB4	22064	32772	2.16%	0.200%
32	13.495	1895	1897	1901	rVB3	60306	55109	3.64%	0.337%
33	13.565	1906	1909	1912	rVB	51166	48679	3.22%	0.298%
34	13.618	1915	1918	1919	rBV2	20513	20457	1.35%	0.125%

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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-BF061918.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	13.948	1971	1974	1978	rBV3	106838	135465	8.95%	0.828%
36	14.142	2002	2007	2010	rBV	597288	616080	40.69%	3.766%
37	14.165	2010	2011	2014	rVV	63643	51977	3.43%	0.318%
38	14.242	2018	2024	2032	rVB5	118861	333134	22.00%	2.037%
39	14.407	2050	2052	2054	rVB2	30624	25993	1.72%	0.159%
40	14.601	2080	2085	2091	rBV2	116158	163281	10.78%	0.998%
41	14.901	2134	2136	2140	rVB2	45717	50720	3.35%	0.310%
42	15.065	2162	2164	2167	rVB3	37154	34961	2.31%	0.214%
43	15.230	2189	2192	2194	rBV	58494	65412	4.32%	0.400%
44	15.265	2195	2198	2202	rVB2	147290	168025	11.10%	1.027%
45	15.554	2244	2247	2252	rVB	69001	82782	5.47%	0.506%
46	15.689	2266	2270	2274	rVB	307091	349233	23.07%	2.135%
47	16.189	2351	2355	2361	rBV2	71802	133252	8.80%	0.815%
48	16.430	2393	2396	2408	rVB	103842	202050	13.35%	1.235%
49	17.565	2585	2589	2602	rVB5	107731	271161	17.91%	1.658%
50	17.842	2632	2636	2646	rVB5	59028	140480	9.28%	0.859%
51	18.453	2736	2740	2753	rVB10	91267	285382	18.85%	1.745%
52	18.942	2821	2823	2828	rVB4	40734	67674	4.47%	0.414%

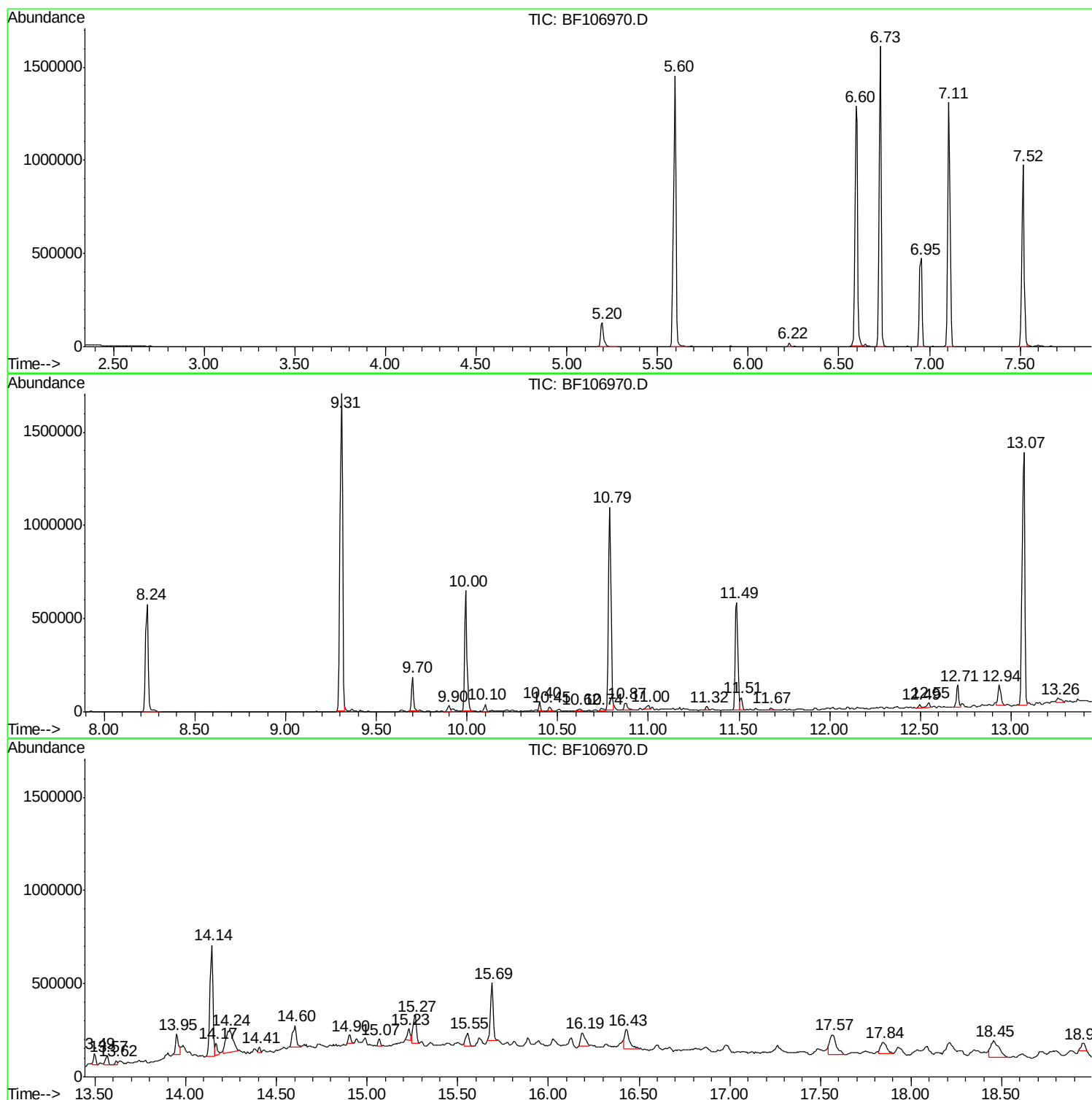
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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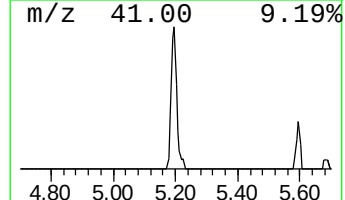
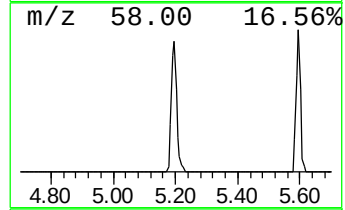
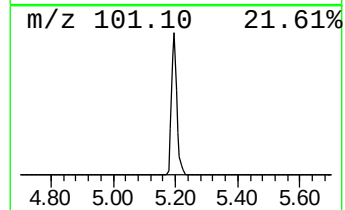
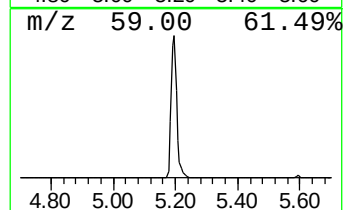
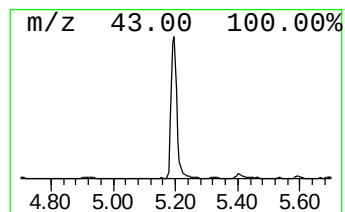
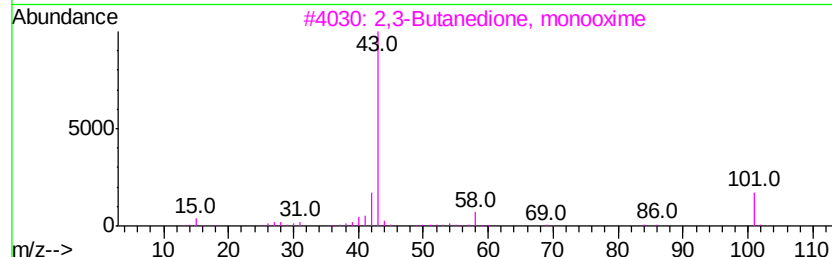
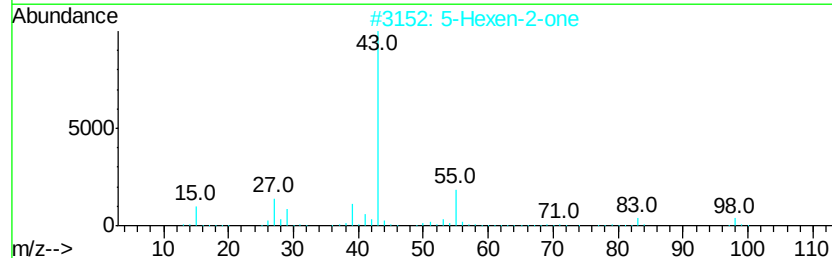
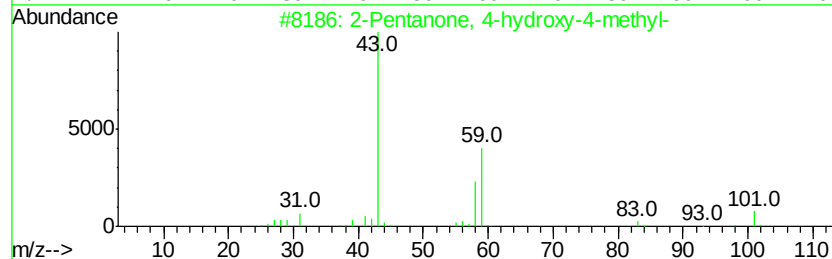
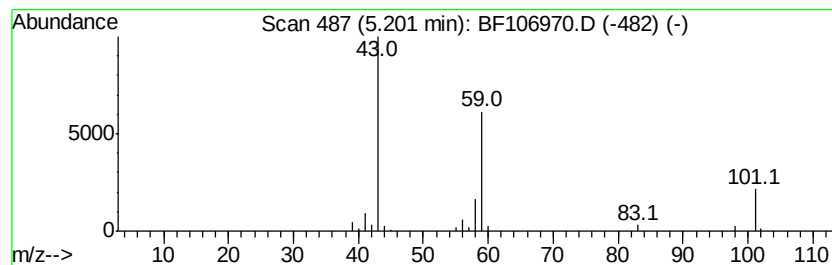
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.20	7.71 ng	158505	1,4-Dichlorobenzene-d4	6.95

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56		
2	5-Hexen-2-one	98	C6H10O	000109-49-9	9		
3	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9		
4	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9		
5	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	9		



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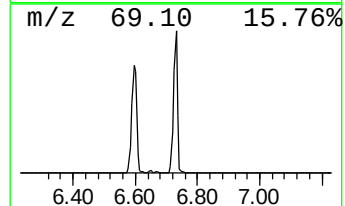
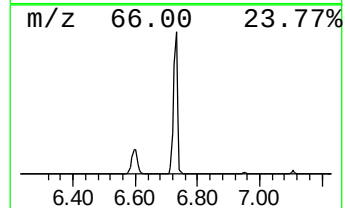
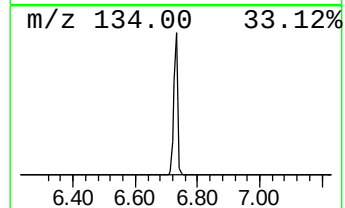
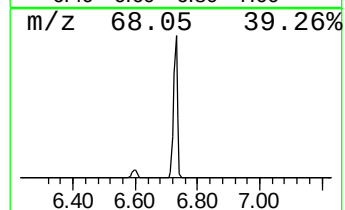
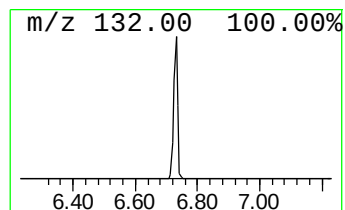
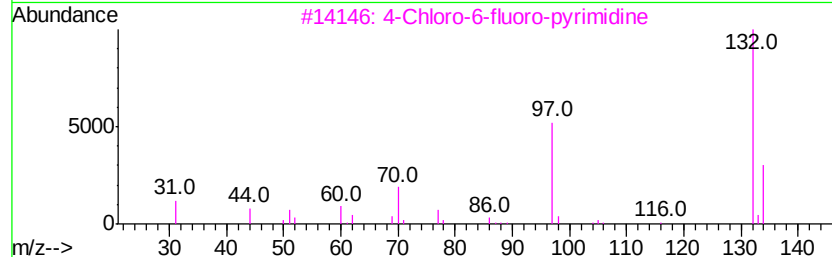
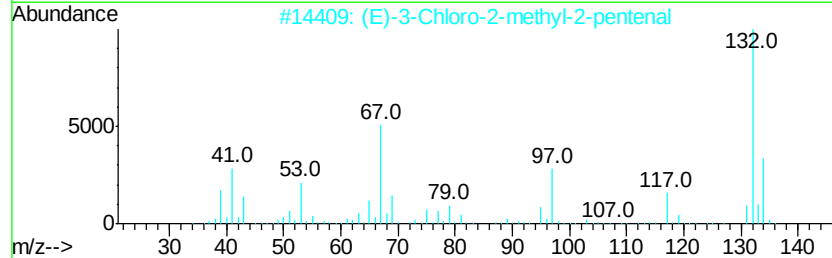
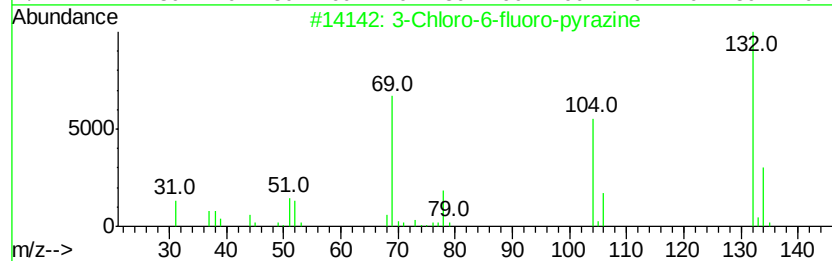
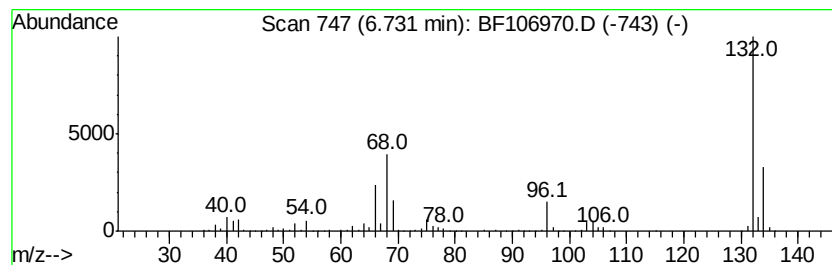
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.73 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.73	69.65 ng	1432400	1,4-Dichlorobenzene-d4	6.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
3		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9
4		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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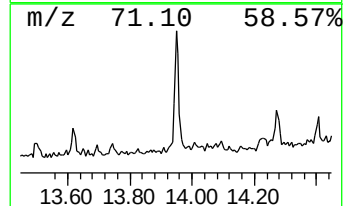
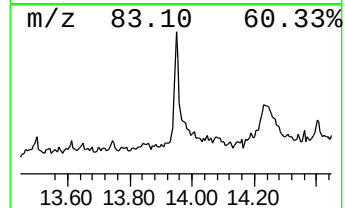
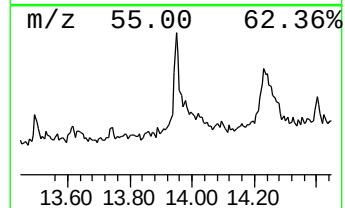
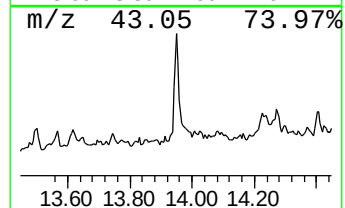
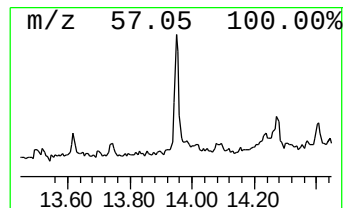
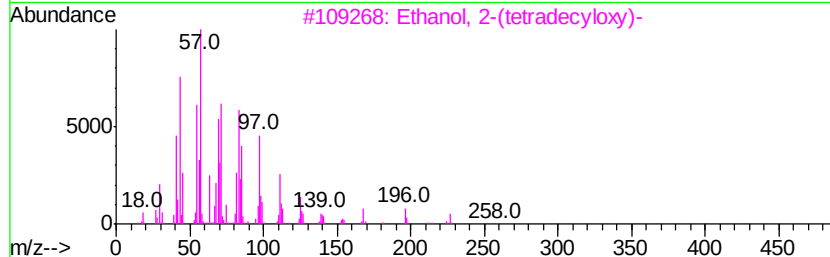
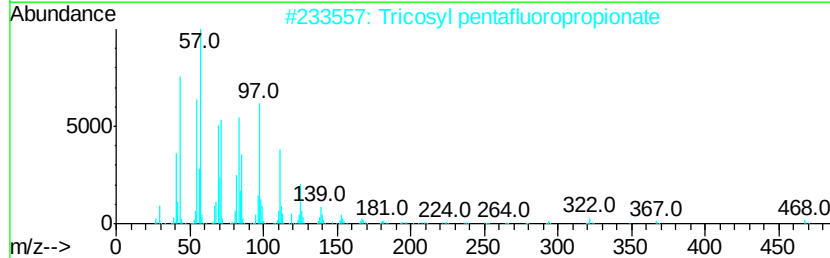
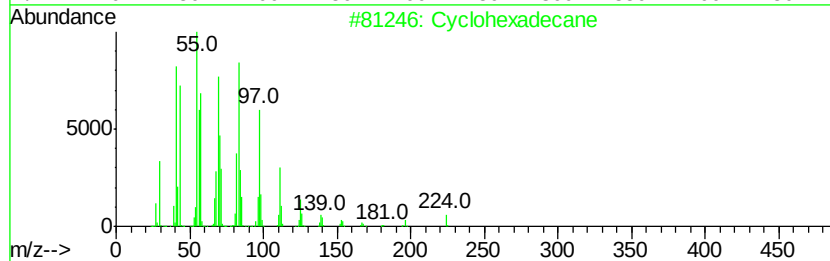
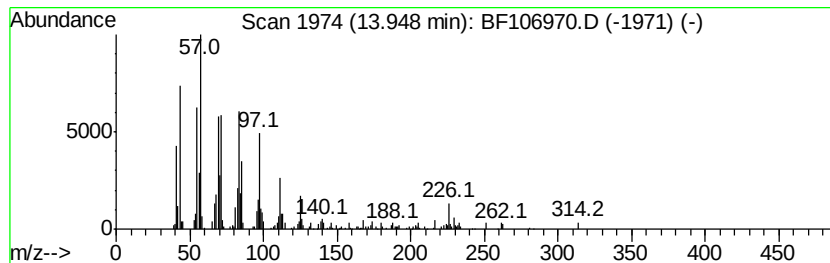
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Cyclohexadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.95	4.40 ng	135465	Chrysene-d12	14.14	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexadecane	224	C16H32	000295-65-8	96
2	Tricosyl pentafluoropropionate	486	C26H47F5O2	1000351-81-0	90
3	Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	90
4	Tricosyl trifluoroacetate	436	C25H47F3O2	1000351-75-1	90
5	Docosyl pentafluoropropionate	472	C25H45F5O2	1000351-80-9	90



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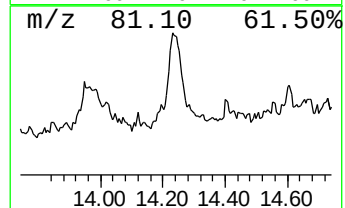
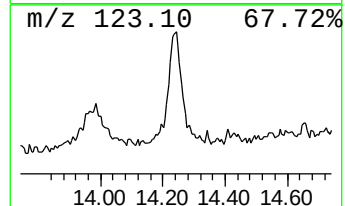
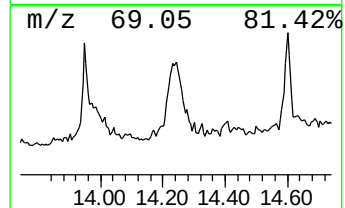
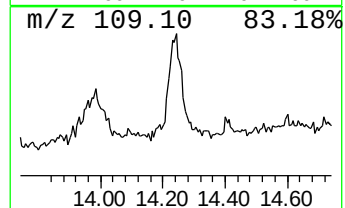
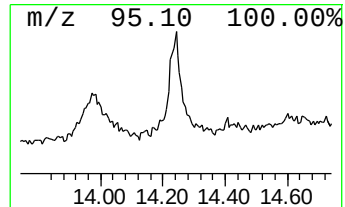
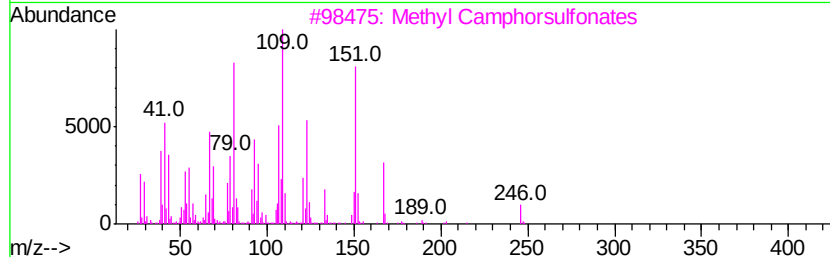
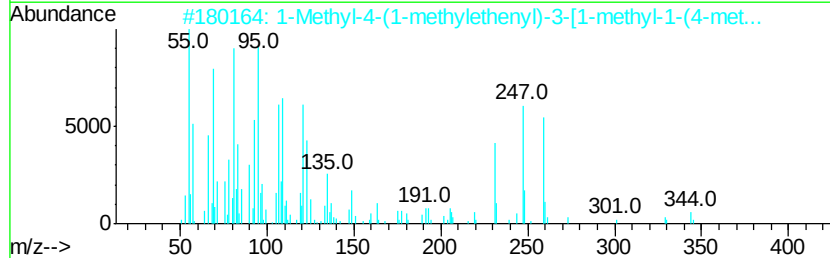
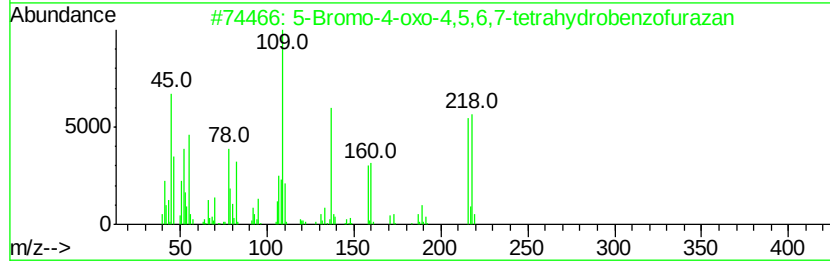
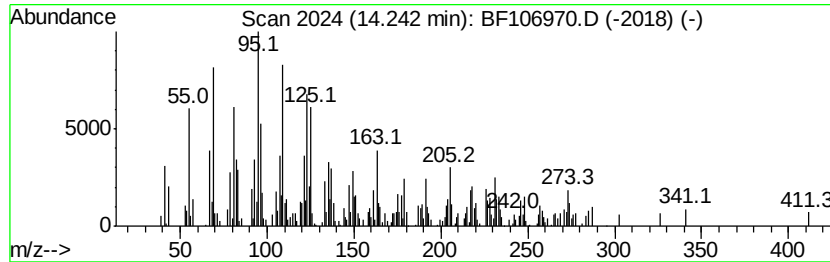
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 5-Bromo-4-oxo-4,5,6,7-tetra... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.24	10.81 ng	333134	Chrysene-d12	14.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Bromo-4-oxo-4,5,6,7-tetrahydro...	216	C6H5BrN2O2	300574-36-1	90
2		1-Methyl-4-(1-methylethenyl)-3-[...	344	C25H44	1000370-41-3	56
3		Methyl Camphorsulfonates	246	C11H18O4S	046471-67-4	55
4		Bicyclo[3.1.1]heptan-3-one, 2-(b...	192	C13H20O	1000163-96-1	49
5		1,1,6-trimethyl-3-methylene-2-(3...	452	C33H56	1000373-94-5	45



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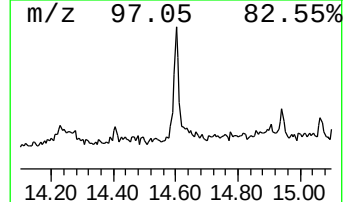
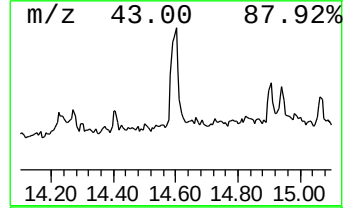
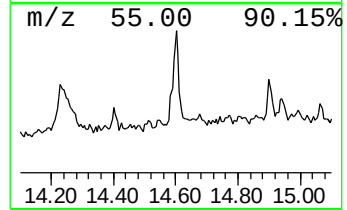
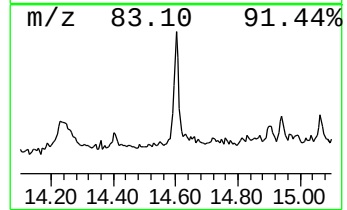
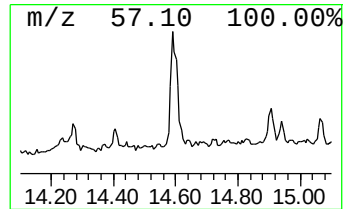
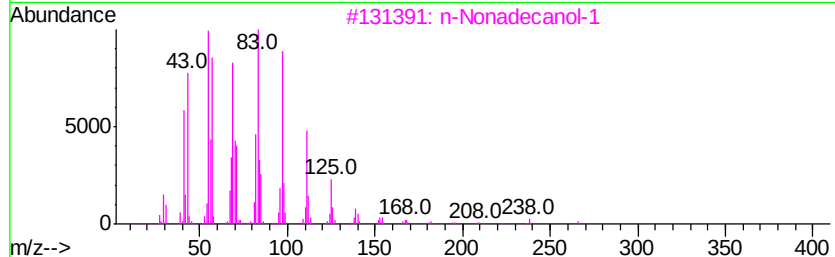
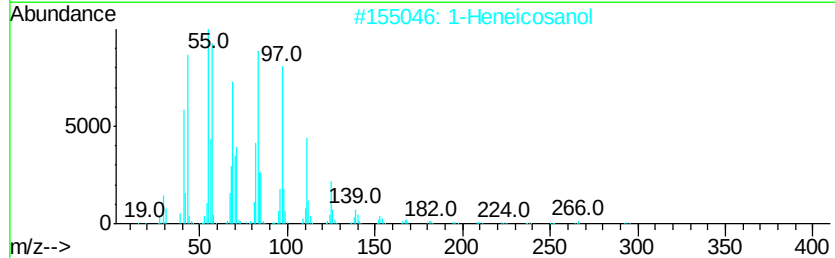
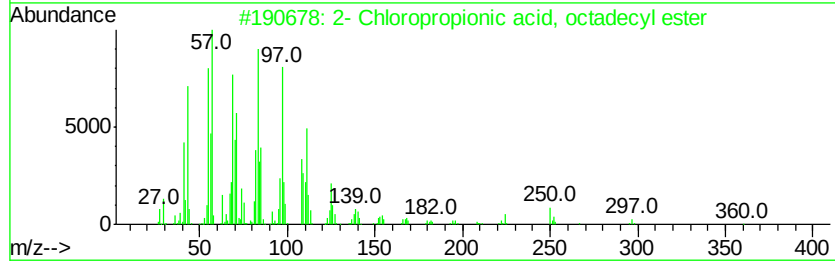
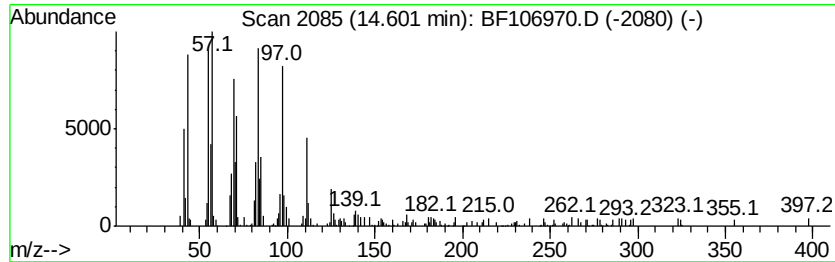
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 2- Chloropropionic acid, oc... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.60	5.30 ng	163281	Chrysene-d12	14.14

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	94
2			1-Heneicosanol	312	C21H44O	015594-90-8	93
3			n-Nonadecanol-1	284	C19H40O	001454-84-8	93
4			Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	93
5			n-Tetracosanol-1	354	C24H50O	000506-51-4	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062918\
Data File : BF106970.D
Acq On : 29 Jun 2018 13:59
Operator : JU/SJ
Sample : J3709-05
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
EME-FW-TS005-01

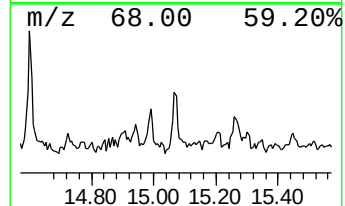
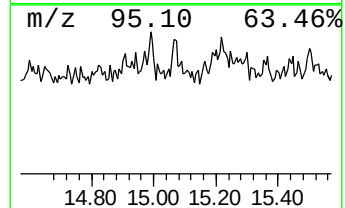
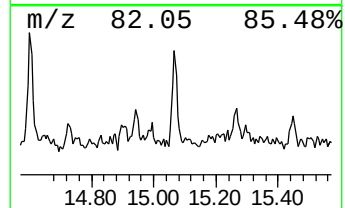
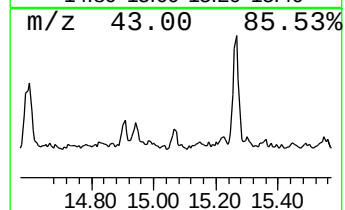
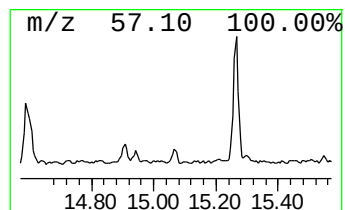
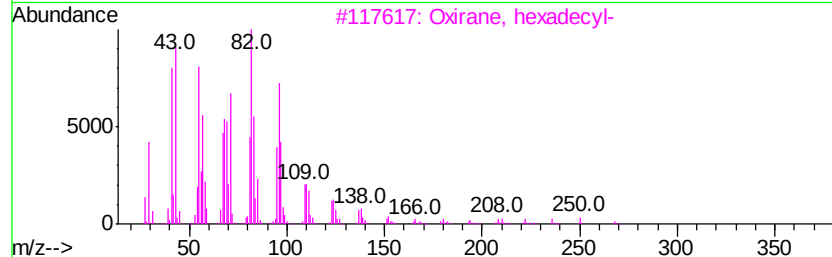
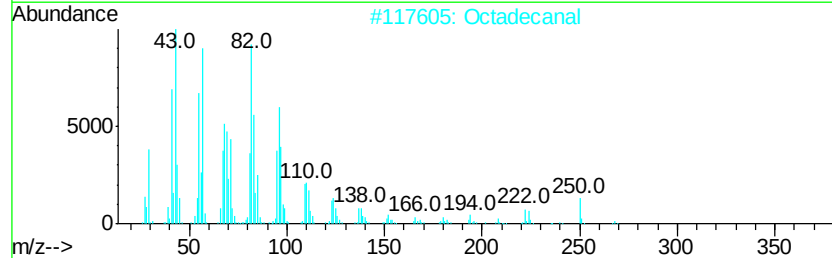
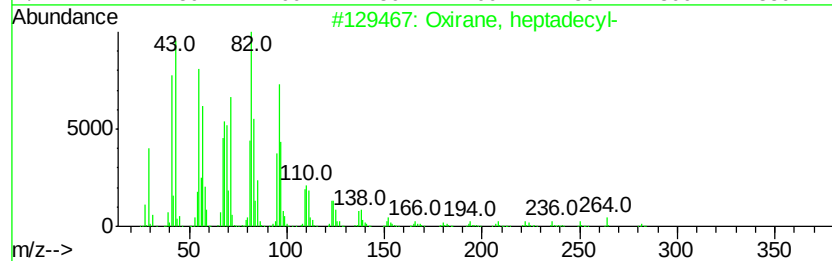
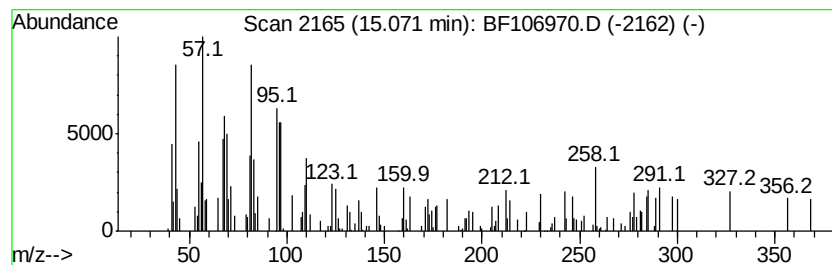
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Oxirane, heptadecyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.07	2.00 ng	34961	Perylene-d12	15.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Oxirane, heptadecyl-	282	C19H38O	067860-04-2	74
2		Octadecanal	268	C18H36O	000638-66-4	72
3		Oxirane, hexadecyl-	268	C18H36O	007390-81-0	72
4		16-Heptadecenal	252	C17H32O	1000144-57-9	62
5		cis-11,12-Epoxytetradecen-1-ol	270	C16H30O3	1000130-82-5	60



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062918\
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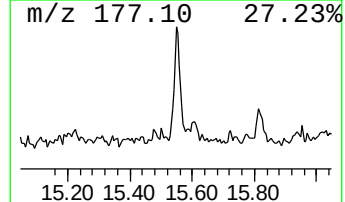
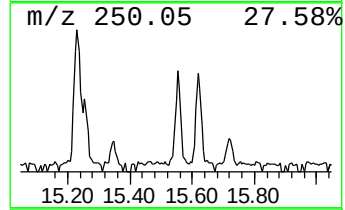
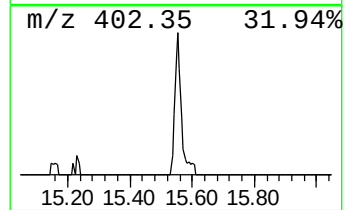
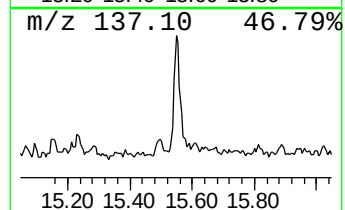
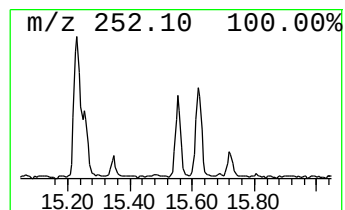
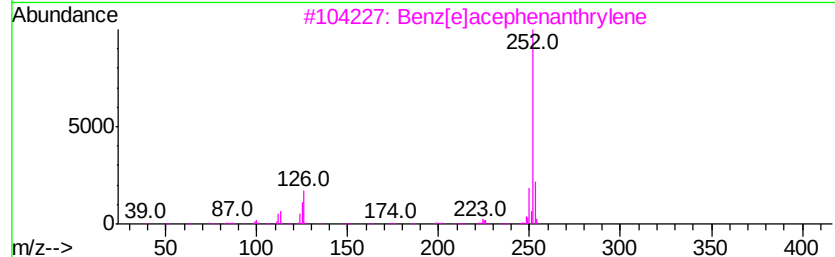
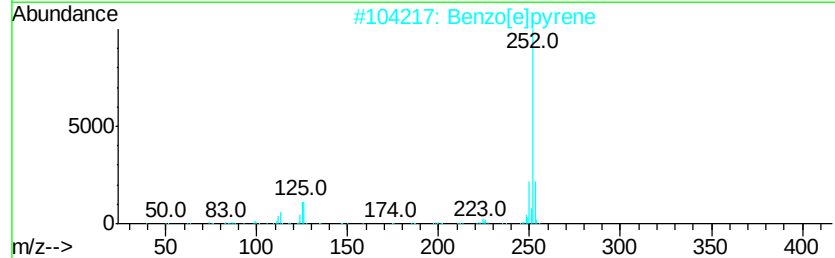
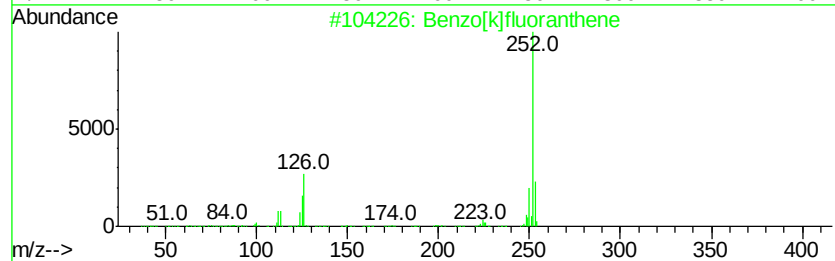
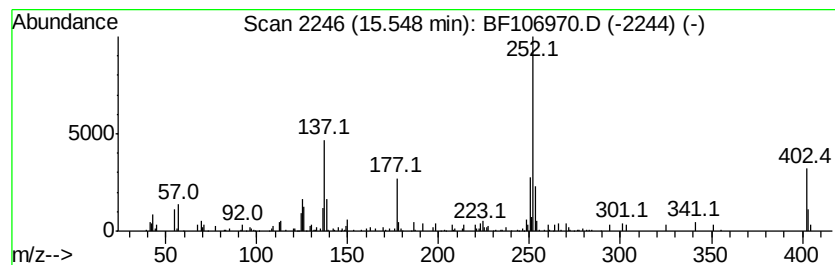
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Benzo[k]fluoranthene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.55	4.74 ng	82782	Perylene-d12	15.69

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzo[k]fluoranthene	252	C20H12	000207-08-9	78
2		Benzo[e]pyrene	252	C20H12	000192-97-2	78
3		Benzo[e]acephenanthrylene	252	C20H12	000205-99-2	78
4		Perylene	252	C20H12	000198-55-0	60
5		Benzo[a]pyrene	252	C20H12	000050-32-8	60



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Data File : BF106970.D
Acq On : 29 Jun 2018 13:59
Operator : JU/SJ
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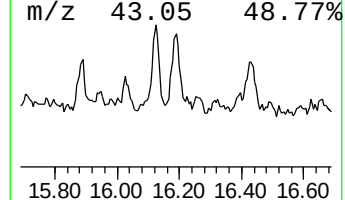
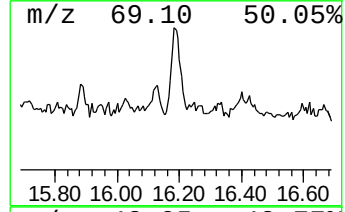
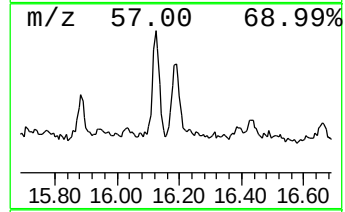
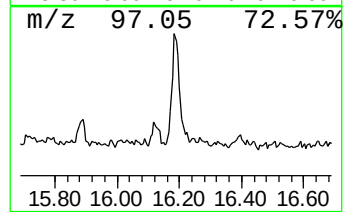
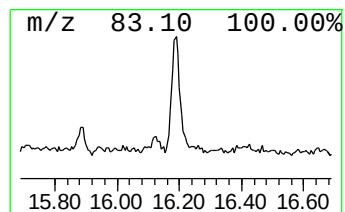
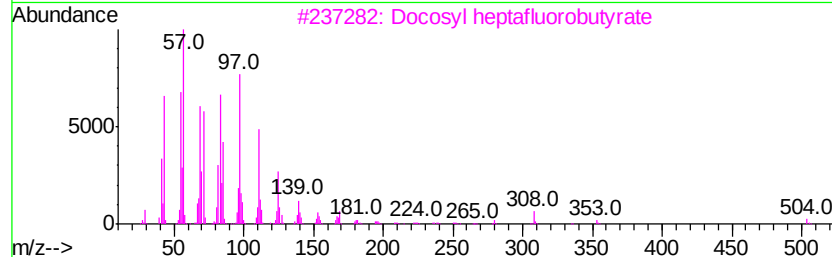
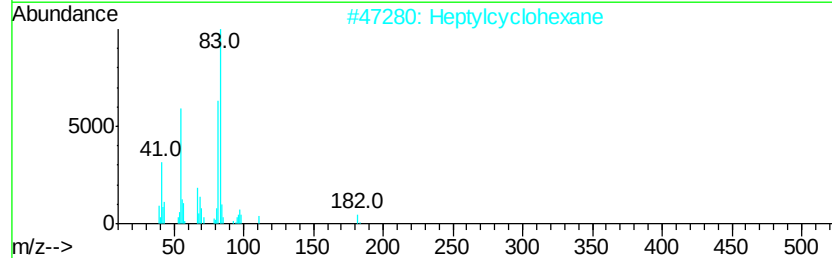
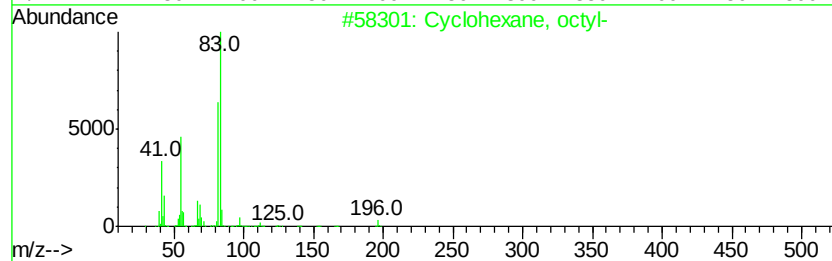
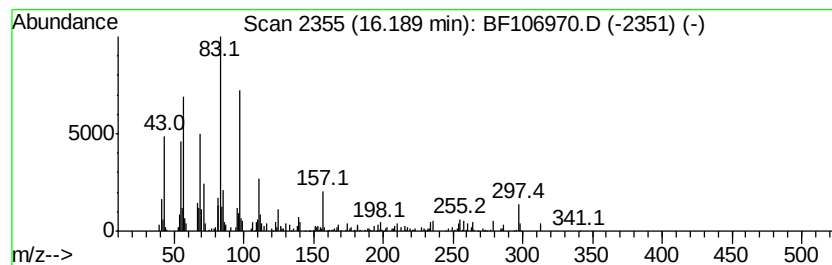
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 unknown16.19 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.19	7.63 ng	133252	Perylene-d12	15.69

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexane, octyl-	196	C14H28	001795-15-9	30
2		Heptylcyclohexane	182	C13H26	005617-41-4	30
3		Docosyl heptafluorobutvrate	522	C26H45F7O2	1000351-83-1	25
4		Glutamic acid, N-tialyl-, dimeth...	257	C12H19N05	1000153-21-0	25
5		Hexatriacontyl trifluoroacetate	619	C38H73F3O2	1000351-88-6	18



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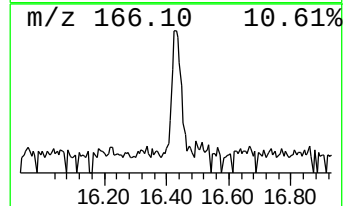
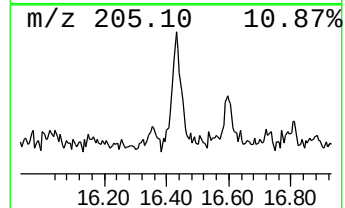
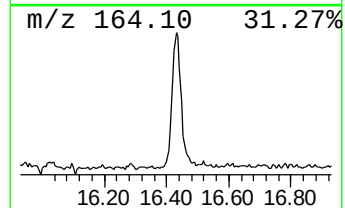
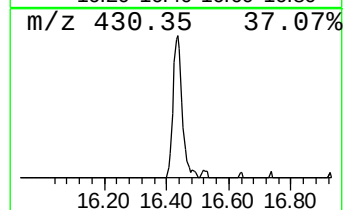
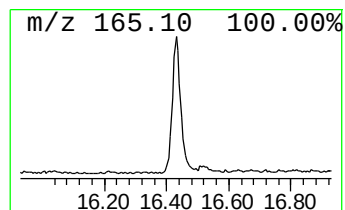
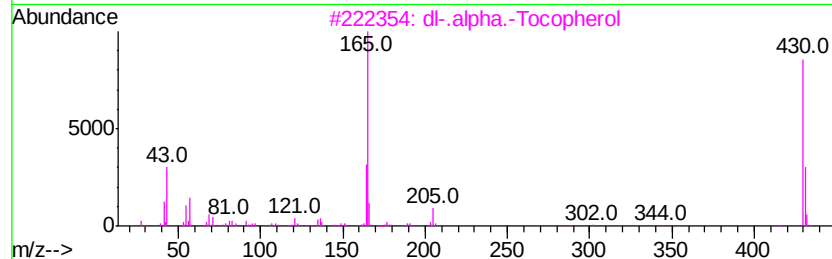
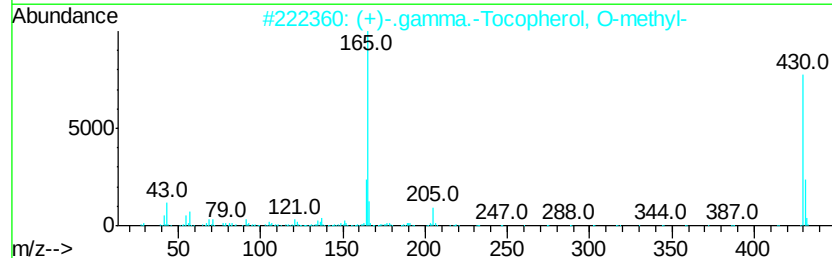
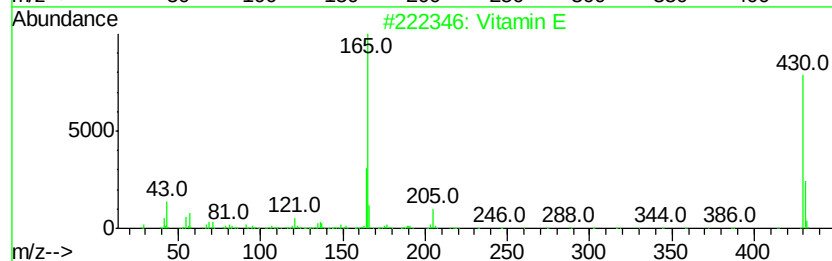
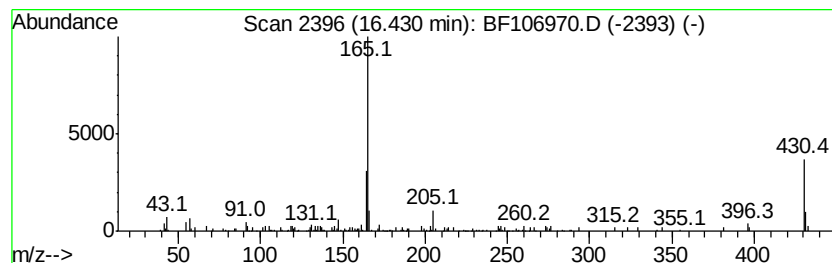
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Vitamin E Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.43	11.57 ng	202050	Perylene-d12	15.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Vitamin E	430	C29H50O2	000059-02-9	87
2		(+)-.gamma.-Tocopherol, O-methyl-	430	C29H50O2	1000374-72-2	83
3		dl-.alpha.-Tocopherol	430	C29H50O2	010191-41-0	80
4		.alpha.-Tocopherol-.beta.-D-mann...	592	C35H60O7	1000156-68-2	64
5		9H-Fluorene, 9-bromo-	244	C13H9Br	001940-57-4	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062918\
Data File : BF106970.D
Acq On : 29 Jun 2018 13:59
Operator : JU/SJ
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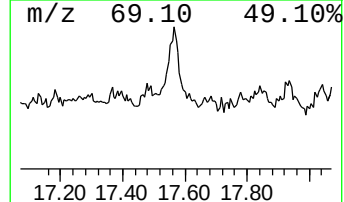
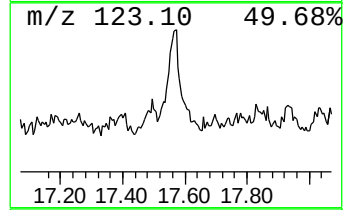
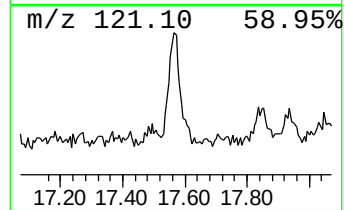
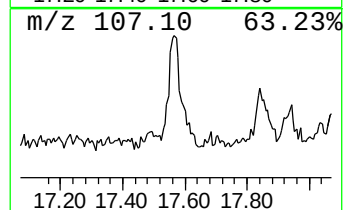
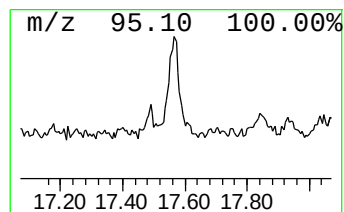
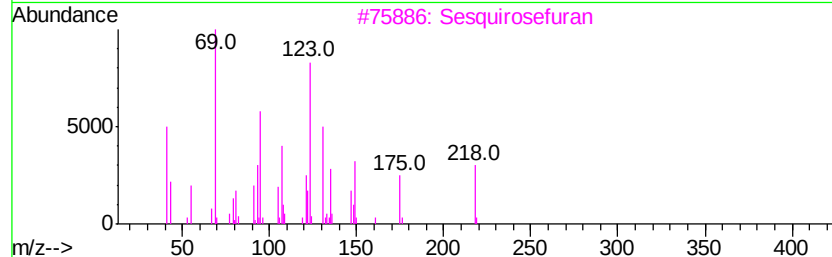
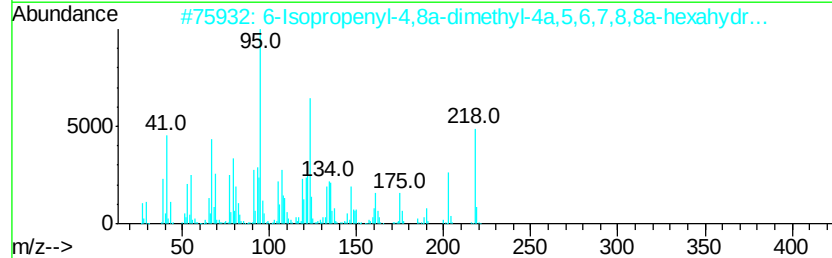
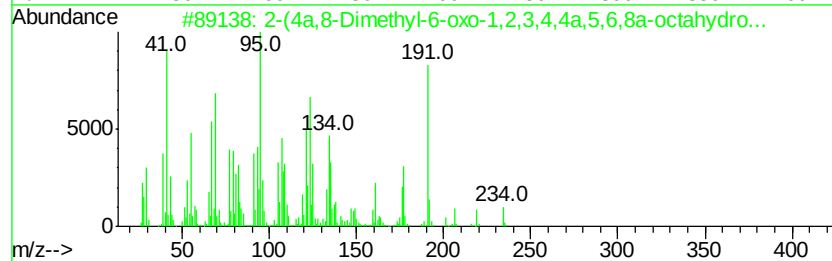
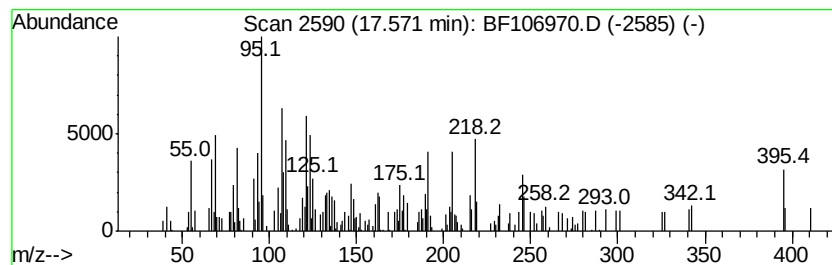
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown17.57 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.57	15.53 ng	271161	Perylene-d12	15.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-(4a,8-Dimethyl-6-oxo-1,2,3,4,4...	234	C15H22O2	1000190-21-8	47
2		6-Isopropenyl-4,8a-dimethyl-4a,5...	218	C15H22O	086917-79-5	46
3		Sesquirosefuran	218	C15H22O	039007-93-7	38
4		Cycloprop[7,8]ergost-22-en-3-one...	410	C29H46O	053755-18-3	37
5		Anthracene, 9-cyclohexyltetradec...	274	C20H34	055255-70-4	25



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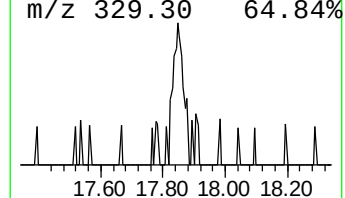
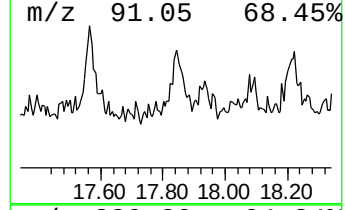
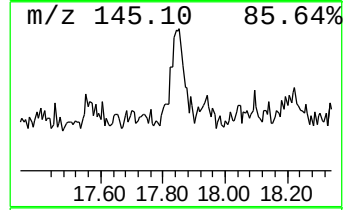
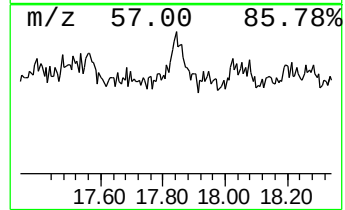
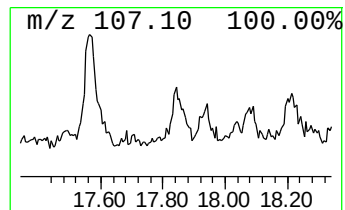
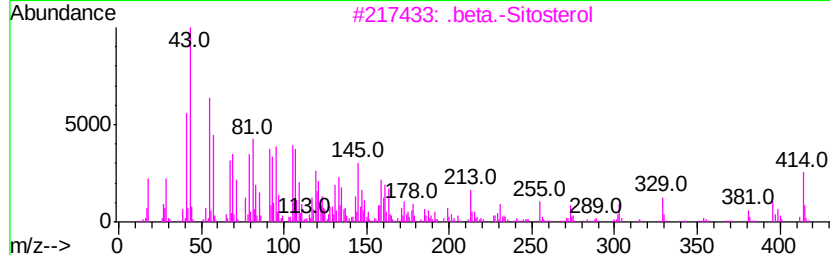
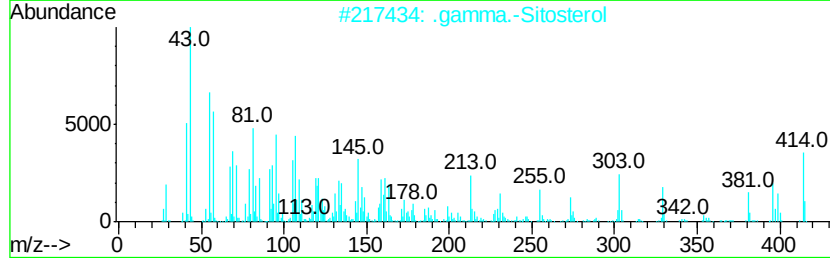
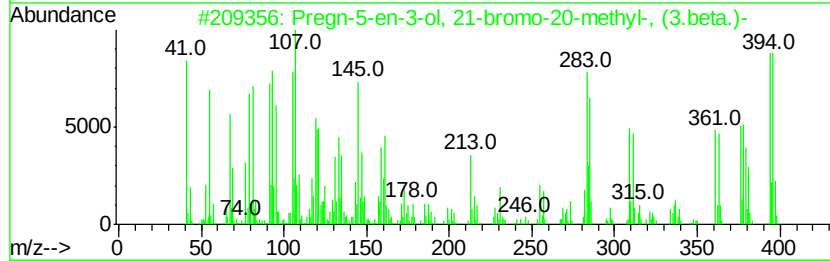
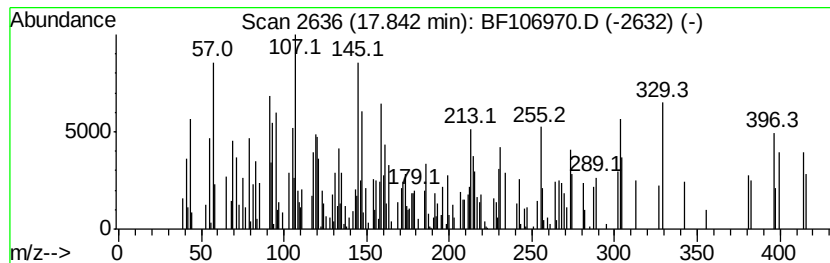
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Pregn-5-en-3-ol, 21-bromo-2... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.
17.84	8.05 ng	140480	Perylene-d12		15.69
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pregn-5-en-3-ol, 21-bromo-20-met...	394	C22H35BrO	055103-80-5	92
2	.gamma.-Sitosterol	414	C29H50O	000083-47-6	70
3	.beta.-Sitosterol	414	C29H50O	000083-46-5	35
4	5-O-Tolylcarbamoyl-3H-imidazole-...	273	C14H15N3O3	313251-31-9	11
5	Norflurazon	303	C12H9ClF3N3O	027314-13-2	10



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ClientSampleId :
EME-FW-TS005-01

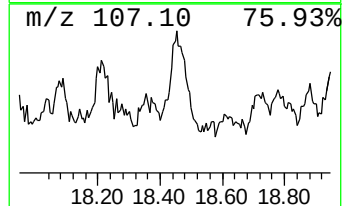
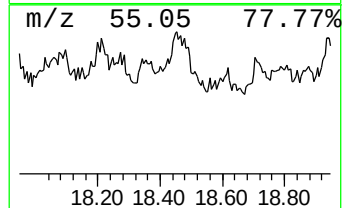
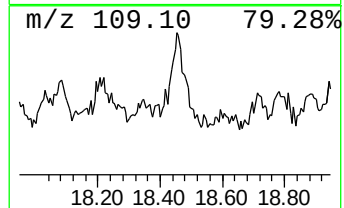
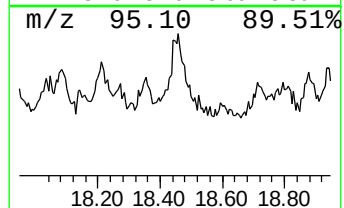
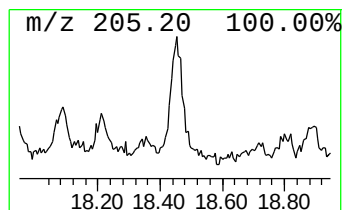
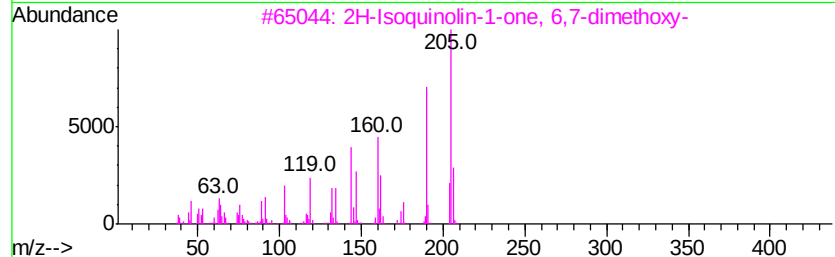
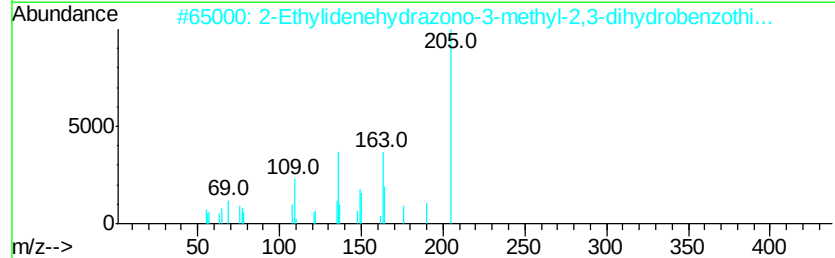
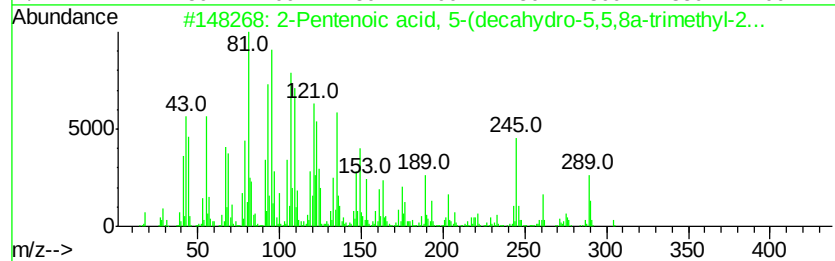
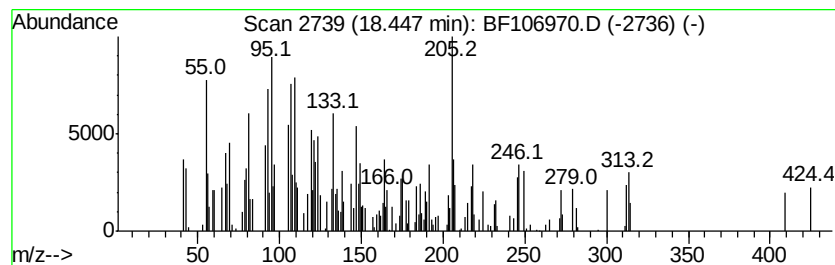
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown18.45 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.45	16.34 ng	285382	Perylene-d12	15.69

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentenoic acid, 5-(decahydro-5...	304	C20H32O2	024470-48-2	27		
2	2-Ethylidenehydrazono-3-methyl-2...	205	C10H11N3S	1000147-60-5	20		
3	2H-Isoquinolin-1-one, 6,7-dimeth...	205	C11H11NO3	1000305-07-9	15		
4	6.beta.Bicyclo[4.3.0]nonane, 5.b...	332	C15H25I	1000195-85-9	14		
5	9-Borabicyclo[3.3.1]nonane, 9-pi...	205	C13H24BN	022516-42-3	11		



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062918\
Data File : BF106970.D
Acq On : 29 Jun 2018 13:59
Operator : JU/SJ
Sample : J3709-05
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
EME-FW-TS005-01

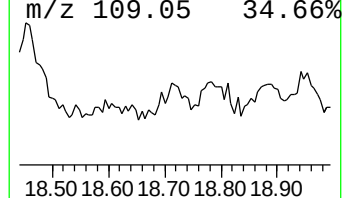
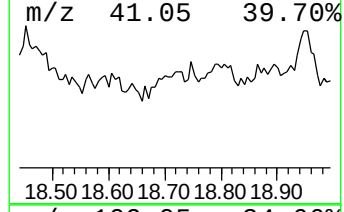
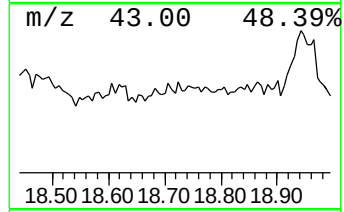
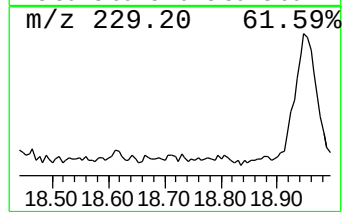
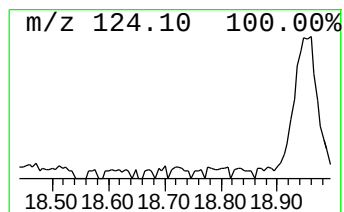
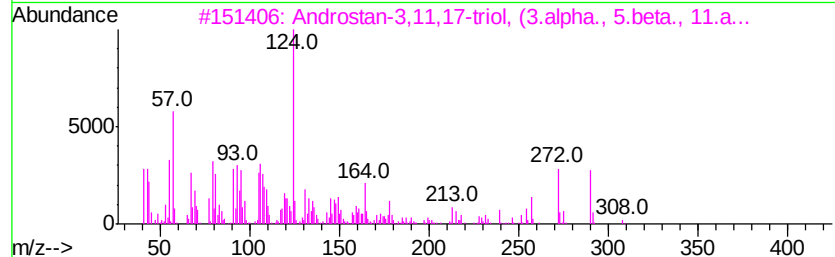
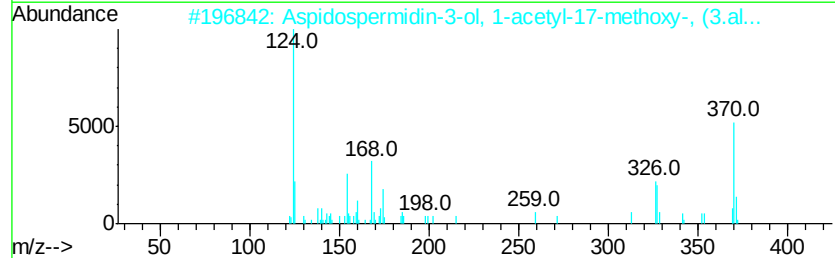
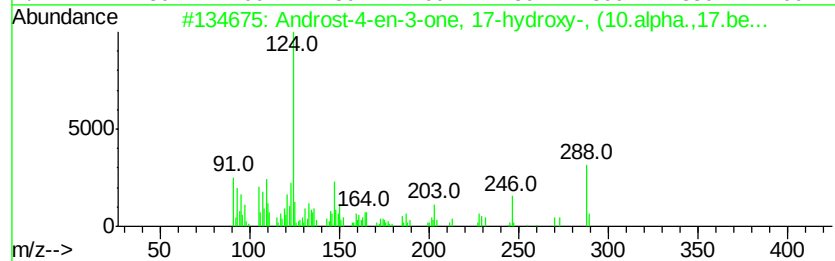
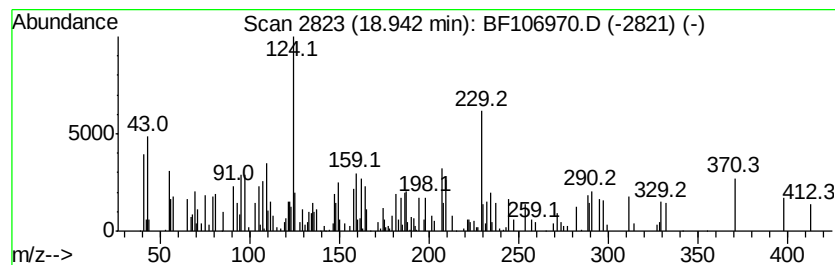
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown18.94 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.94	3.88 ng	67674	Perylene-d12	15.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Androst-4-en-3-one, 17-hydroxy-,...	288	C19H28O2	000604-39-7	35
2		Aspidospermidin-3-ol, 1-acetyl-1...	370	C22H30N2O3	054725-60-9	27
3		Androstan-3,11,17-triol, (3.alpha...	308	C19H32O3	004484-95-1	27
4		1,3-Benzenediol, 5-hexyl-	194	C12H18O2	005465-20-3	14
5		1,5-Dioxaspiro[5.5]undec-3-en-2-...	278	C17H26O3	1000155-15-1	14



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF062918\
 Data File : BF106970.D
 Acq On : 29 Jun 2018 13:59
 Operator : JU/SJ
 Sample : J3709-05
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EME-FW-TS005-01

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

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.20	7.7	ng	158505	1	6.95	411322	20.0
unknown6.73	6.73	69.7	ng	1432400	1	6.95	411322	20.0
Cyclohexadecane	13.95	4.4	ng	135465	5	14.14	616080	20.0
5-Bromo-4-oxo-4,5...	14.24	10.8	ng	333134	5	14.14	616080	20.0
2-Chloropropioni...	14.60	5.3	ng	163281	5	14.14	616080	20.0
Oxirane, heptadecyl-	15.07	2.0	ng	34961	6	15.69	349233	20.0
Benzo[k]fluoranthene	15.55	4.7	ng	82782	6	15.69	349233	20.0
unknown16.19	16.19	7.6	ng	133252	6	15.69	349233	20.0
Vitamin E	16.43	11.6	ng	202050	6	15.69	349233	20.0
unknown17.57	17.57	15.5	ng	271161	6	15.69	349233	20.0
Pregn-5-en-3-ol, ...	17.84	8.1	ng	140480	6	15.69	349233	20.0
unknown18.45	18.45	16.3	ng	285382	6	15.69	349233	20.0
unknown18.94	18.94	3.9	ng	67674	6	15.69	349233	20.0