

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF101818\
 Data File : BF109980.D
 Acq On : 19 Oct 2018 1:02
 Operator : JU/SJ
 Sample : J5570-02
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TS-07

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-BF101518.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	2.357	40	46	52	rVB	666225	939178	13.53%	1.556%
2	3.622	259	261	271	rBV	101974	159082	2.29%	0.263%
3	5.239	532	536	546	rVB	208476	307876	4.44%	0.510%
4	5.610	592	599	602	rBV	5444461	6116733	88.13%	10.131%
5	6.604	761	768	772	rBV2	5150325	6940410	100.00%	11.495%
6	6.751	787	793	796	rBV	5708455	6285822	90.57%	10.411%
7	6.975	827	831	834	rBV	1719267	1362189	19.63%	2.256%
8	7.133	853	858	861	rBV	5393688	4863838	70.08%	8.056%
9	7.539	921	927	930	rBV	3754480	4158578	59.92%	6.888%
10	8.257	1045	1049	1052	rBV	2064046	1717985	24.75%	2.845%
11	9.333	1226	1232	1235	rBV	6253209	6253504	90.10%	10.357%
12	9.722	1294	1298	1301	rBV	1009752	777559	11.20%	1.288%
13	10.016	1344	1348	1352	rBV2	2035598	1791904	25.82%	2.968%
14	10.810	1478	1483	1486	rBV	3837714	3646230	52.54%	6.039%
15	11.510	1598	1602	1604	rBV	1831326	1556816	22.43%	2.578%
16	11.533	1604	1606	1610	rVB	212142	173743	2.50%	0.288%
17	12.015	1684	1688	1694	rBV2	344095	421132	6.07%	0.698%
18	12.721	1804	1808	1813	rBV	474601	398971	5.75%	0.661%
19	12.951	1841	1847	1851	rBV	339971	370183	5.33%	0.613%
20	13.098	1867	1872	1875	rBV	4139105	4205572	60.60%	6.966%
21	13.521	1939	1944	1947	rBV	204058	191572	2.76%	0.317%
22	13.562	1947	1951	1957	rVB2	412537	447614	6.45%	0.741%
23	13.974	2015	2021	2027	rBV4	367514	457105	6.59%	0.757%
24	14.151	2047	2051	2054	rBV	1518143	1381076	19.90%	2.287%
25	14.174	2054	2055	2061	rVB	181615	131607	1.90%	0.218%
26	14.298	2072	2076	2078	rBV	103508	108144	1.56%	0.179%
27	14.604	2125	2128	2135	rVB	236256	273480	3.94%	0.453%
28	14.921	2179	2182	2189	rBV	203245	255118	3.68%	0.423%
29	15.080	2206	2209	2214	rVB	96059	101335	1.46%	0.168%
30	15.221	2229	2233	2236	rBV	173350	245848	3.54%	0.407%
31	15.280	2240	2243	2246	rVV	358056	428504	6.17%	0.710%
32	15.304	2246	2247	2256	rVB3	131650	167607	2.41%	0.278%
33	15.545	2284	2288	2294	rVB	94572	133179	1.92%	0.221%
34	15.609	2295	2299	2304	rVV	116436	139859	2.02%	0.232%

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

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

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	15.680	2305	2311	2326	rVB2	1210484	1695610	24.43%	2.808%
36	16.151	2385	2391	2396	rBV	364092	576878	8.31%	0.955%
37	16.203	2396	2400	2407	rVB4	71391	135630	1.95%	0.225%
38	16.686	2477	2482	2488	rBV	176161	313310	4.51%	0.519%
39	17.233	2571	2575	2584	rVB3	66691	153648	2.21%	0.254%
40	17.327	2585	2591	2598	rVV2	94590	198879	2.87%	0.329%
41	17.415	2601	2606	2613	rVB9	58617	130855	1.89%	0.217%
42	17.703	2652	2655	2660	rVB2	40676	72553	1.05%	0.120%
43	17.850	2676	2680	2690	rVB7	78678	190182	2.74%	0.315%

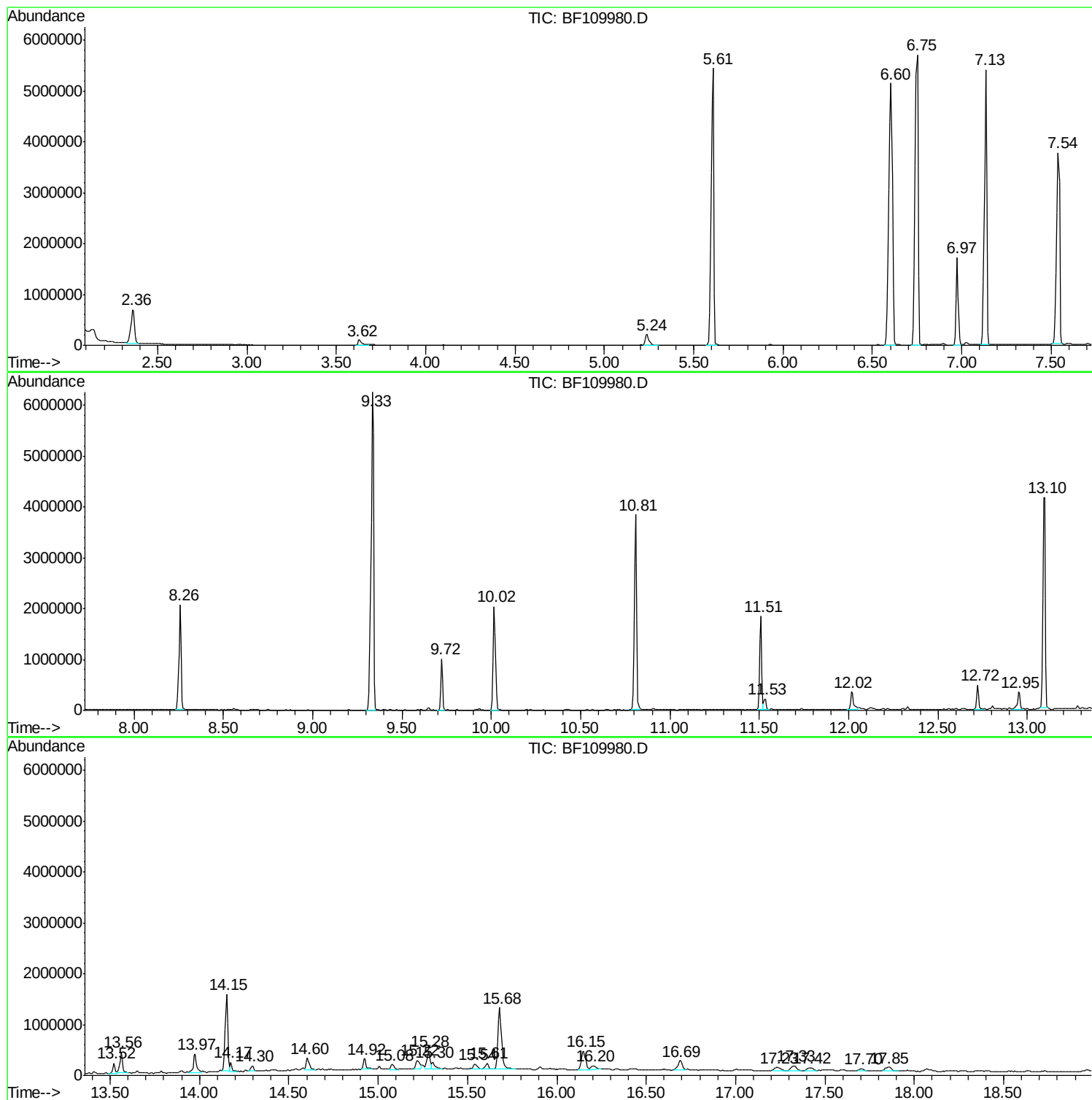
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ClientSampled :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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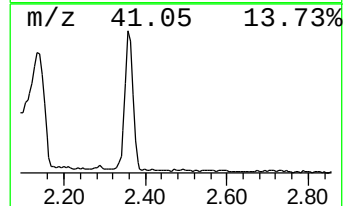
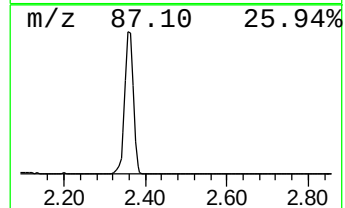
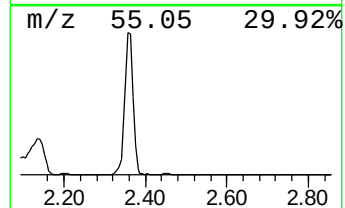
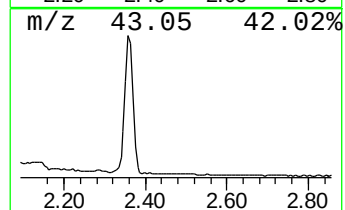
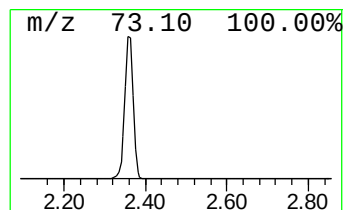
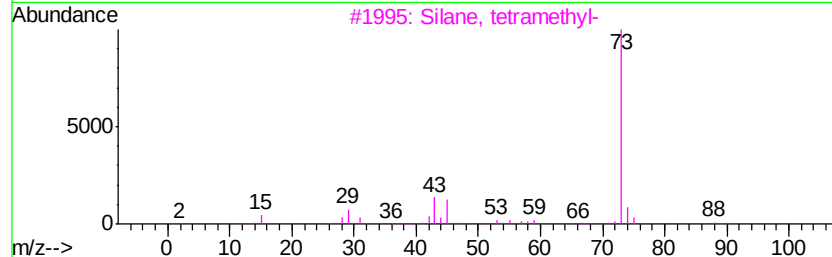
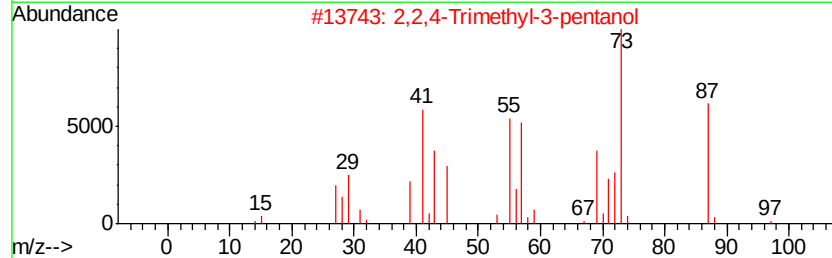
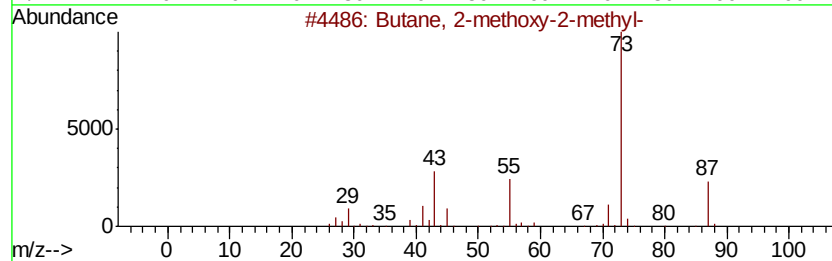
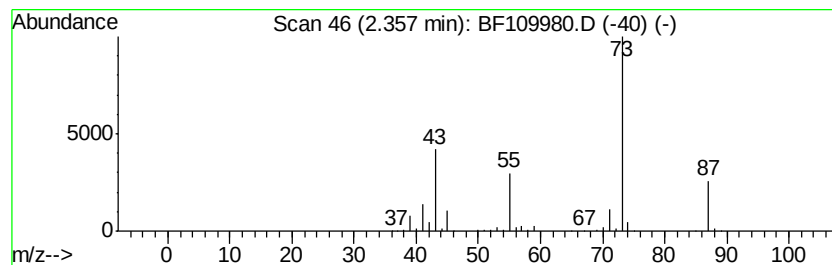
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.36	13.79 ng	939178	1,4-Dichlorobenzene-d4	6.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3		Silane, tetramethyl-	88	C4H12Si	000075-76-3	38
4		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	23
5		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12



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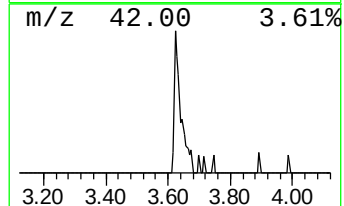
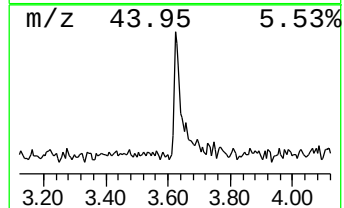
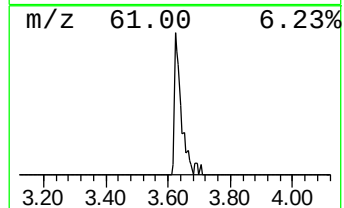
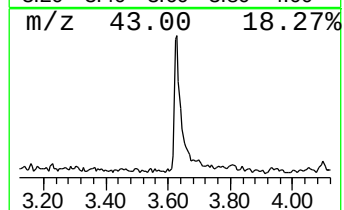
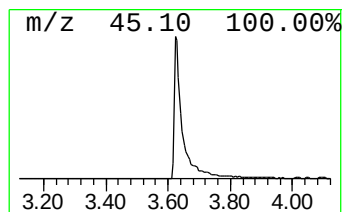
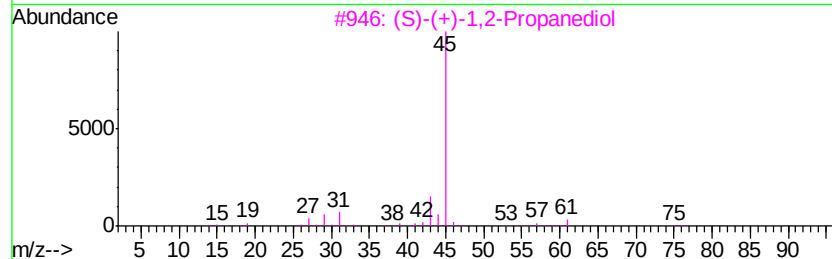
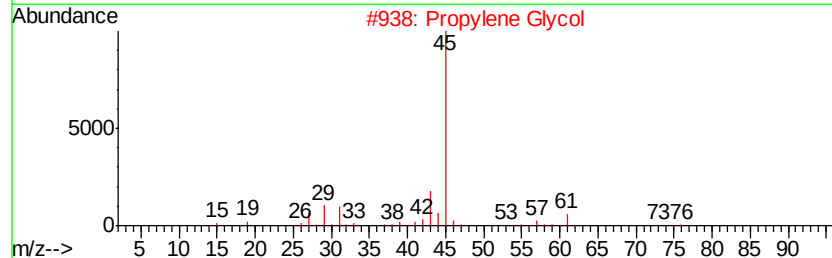
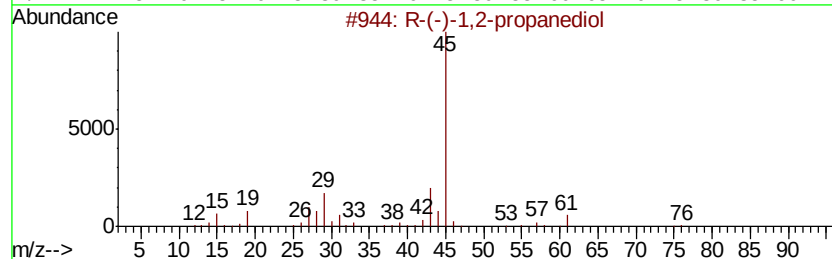
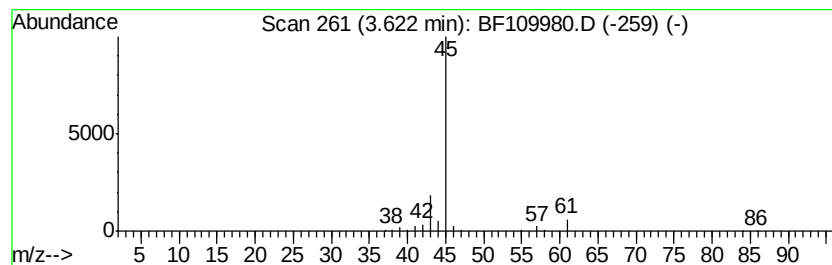
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Peak Number 2 unknown3.62 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.62	2.34 ng	159082	1,4-Dichlorobenzene-d4	6.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	43
2		Propylene Glycol	76	C3H8O2	000057-55-6	43
3		(S)-(+)-1,2-Propanediol	76	C3H8O2	004254-15-3	40
4		Isopropyl Alcohol	60	C3H8O	000067-63-0	9
5		L-Lactic acid	90	C3H6O3	000079-33-4	9



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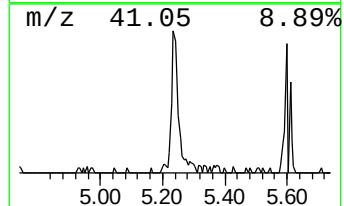
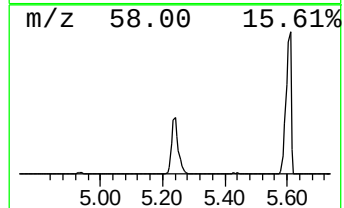
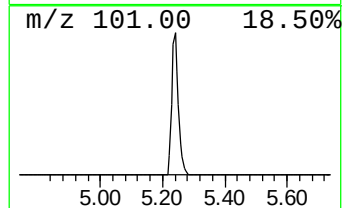
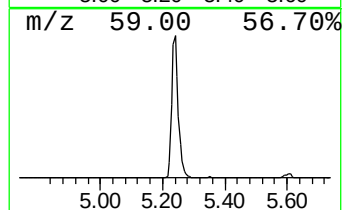
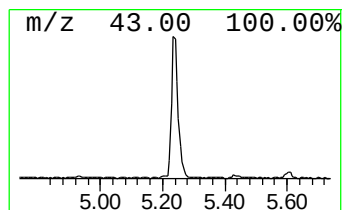
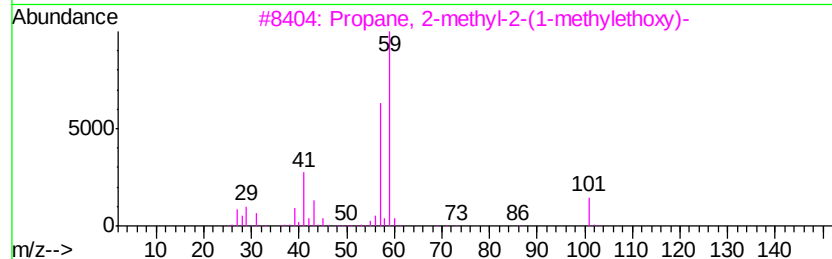
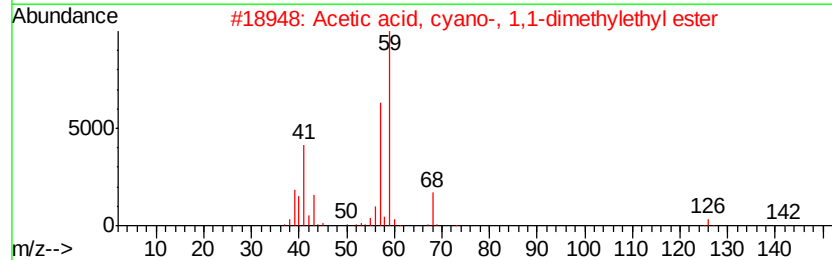
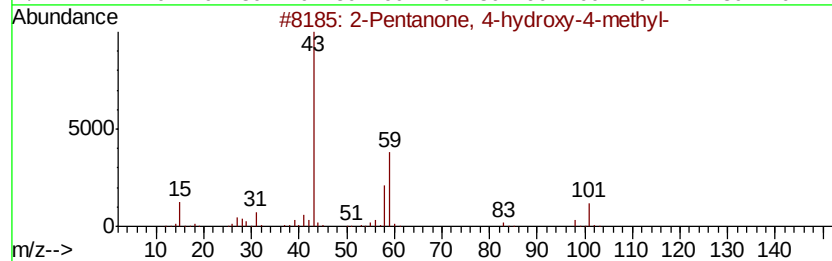
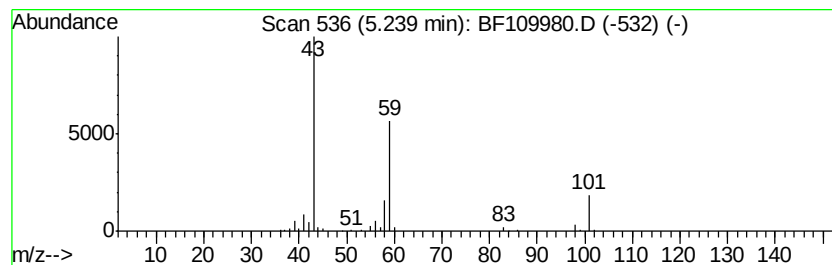
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.24	4.52 ng	307876	1,4-Dichlorobenzene-d4	6.97

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	37
3			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	33
4			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
5			5-Hexen-2-one	98	C6H10O	000109-49-9	9



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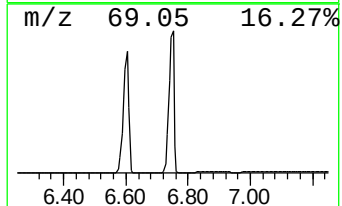
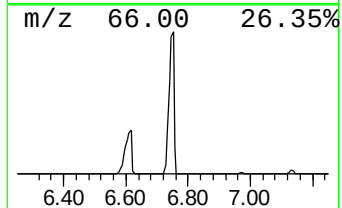
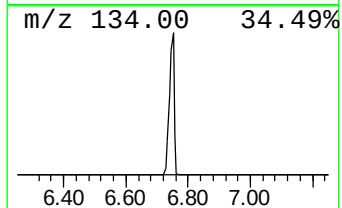
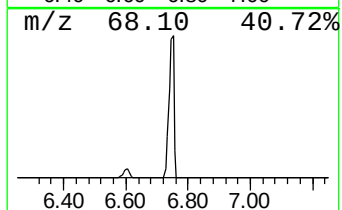
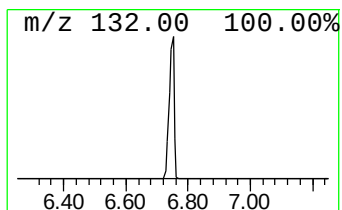
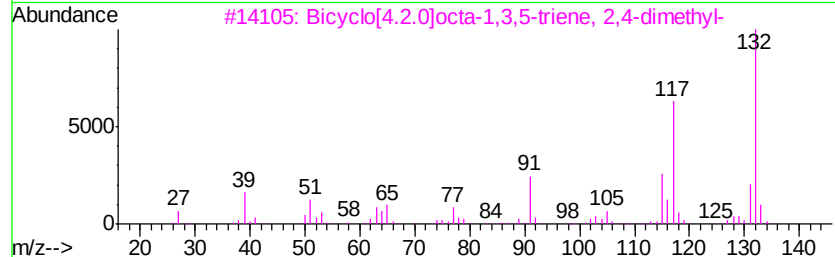
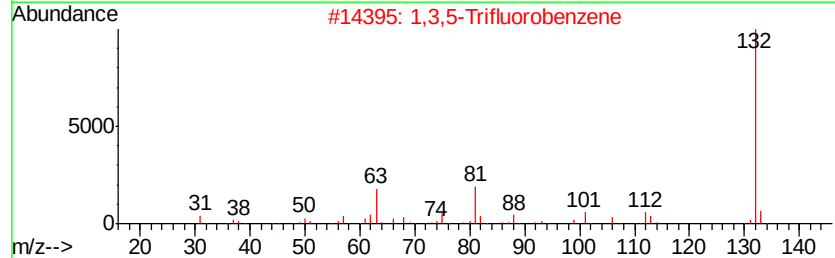
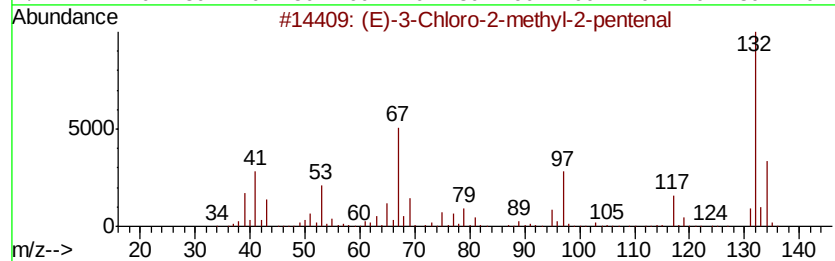
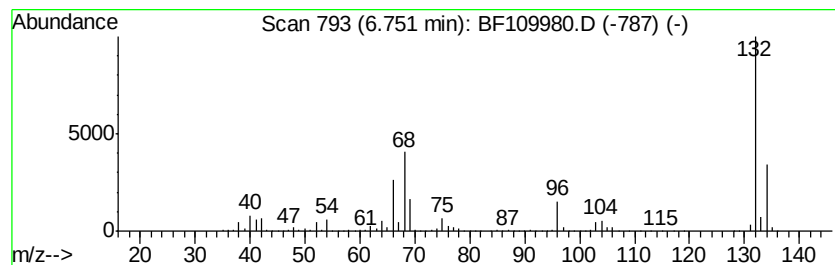
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown6.75 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.75	92.29 ng	6285820	1,4-Dichlorobenzene-d4	6.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
2		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		Xenon	132	Xe	007440-63-3	9
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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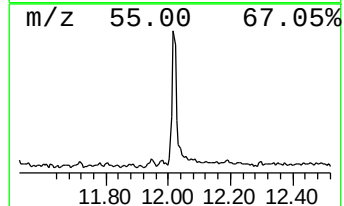
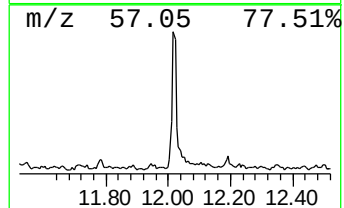
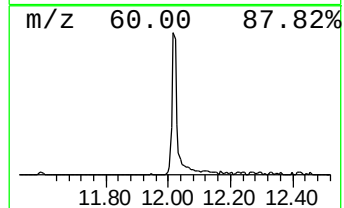
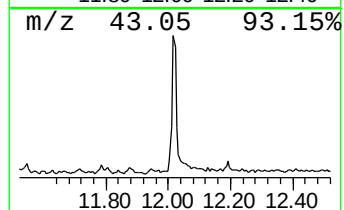
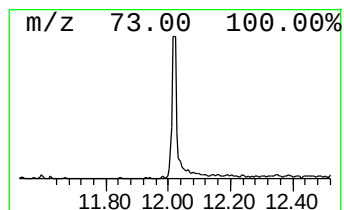
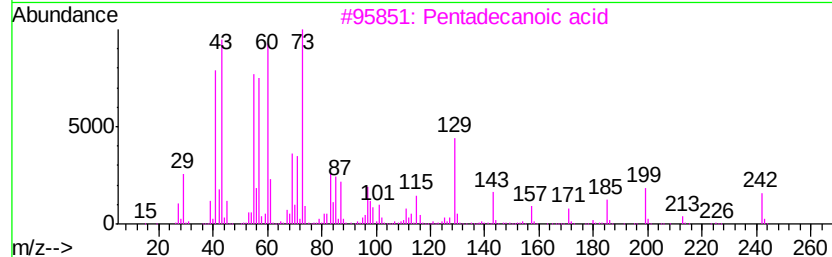
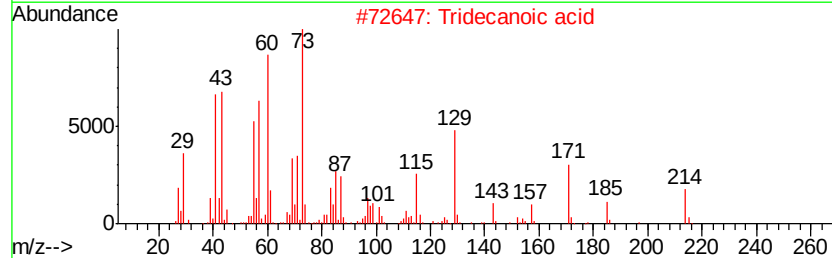
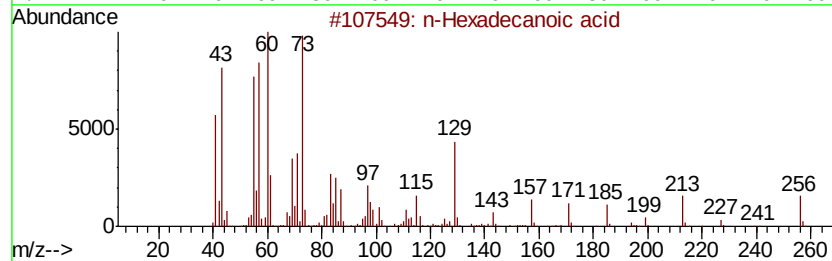
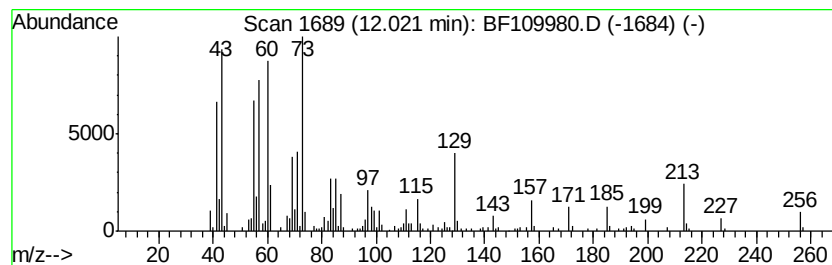
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 6 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.02	5.41 ng	421132	Phenanthrene-d10	11.51

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tridecanoic acid	214	C13H26O2	000638-53-9	96
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	87
4		Undecanoic acid	186	C11H22O2	000112-37-8	74
5		n-Decanoic acid	172	C10H20O2	000334-48-5	64



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Acq On : 19 Oct 2018 1:02
Operator : JU/SJ
Sample : J5570-02
Misc :
ALS Vial : 17 Sample Multiplier: 1

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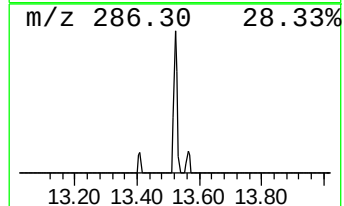
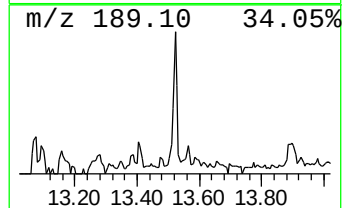
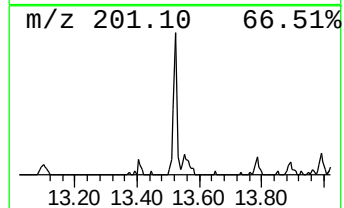
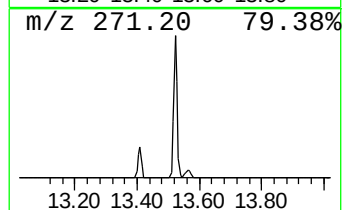
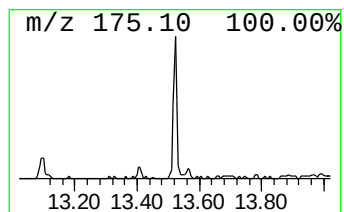
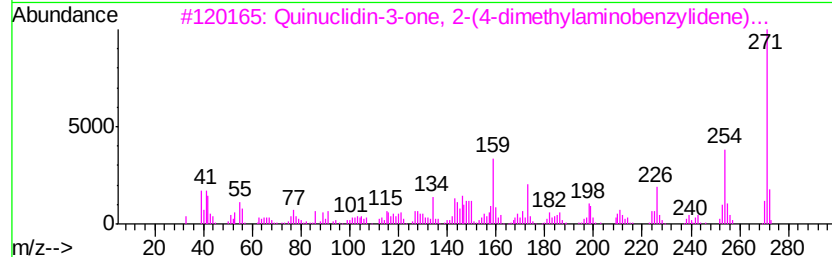
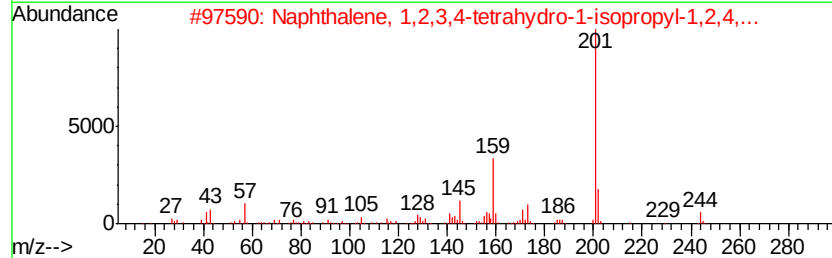
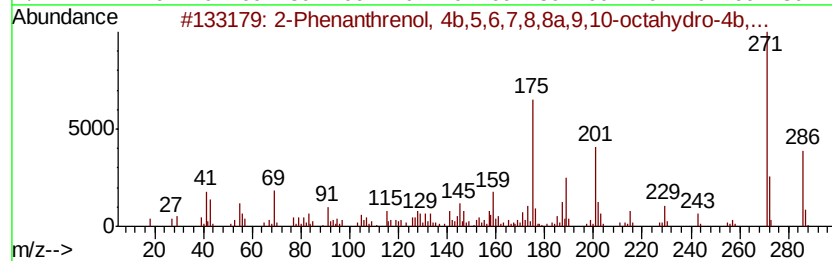
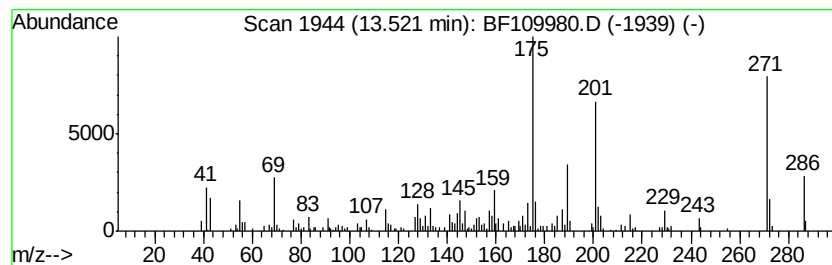
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 2-Phenanthrenol, 4b,5,6,7,8... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.52	2.77 ng	191572	Chrysene-d12	14.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Phenanthrenol, 4b,5,6,7,8,8a,9...	286	C20H30O	000511-15-9	93
2		Naphthalene, 1,2,3,4-tetrahydro-...	244	C18H28	029577-17-1	25
3		Quinuclidin-3-one, 2-(4-dimethyl...	271	C16H21N3O	1000266-14-6	22
4		2,2-Bis(4'-methoxyphenyl)-2-etho...	286	C18H22O3	1000283-53-7	22
5		Benz(a)acridine, 9,10,12-trimethyl-	271	C20H17N	063040-02-8	18



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101818\
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Sample : J5570-02
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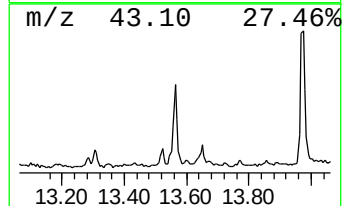
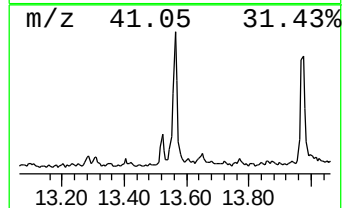
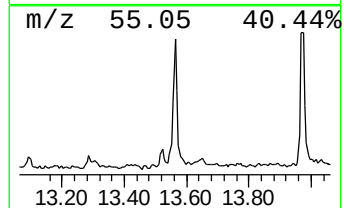
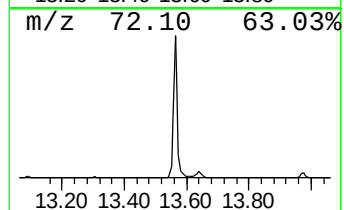
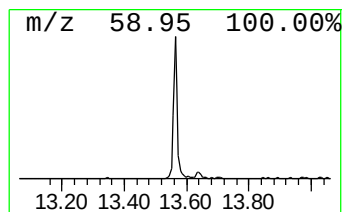
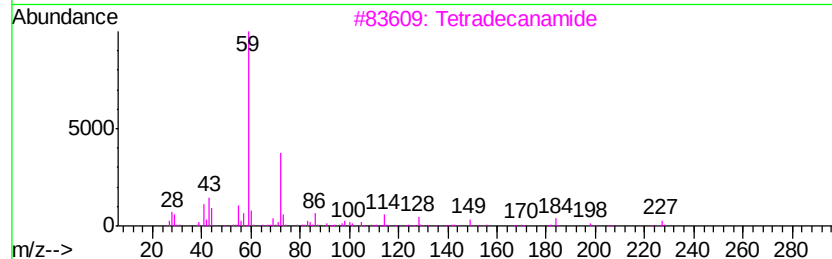
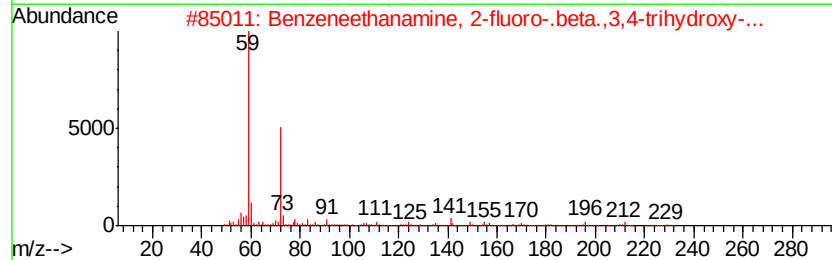
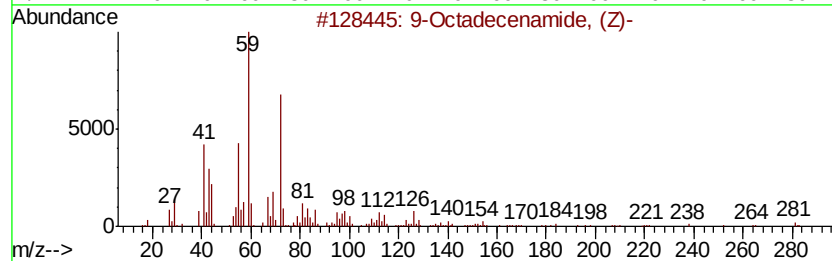
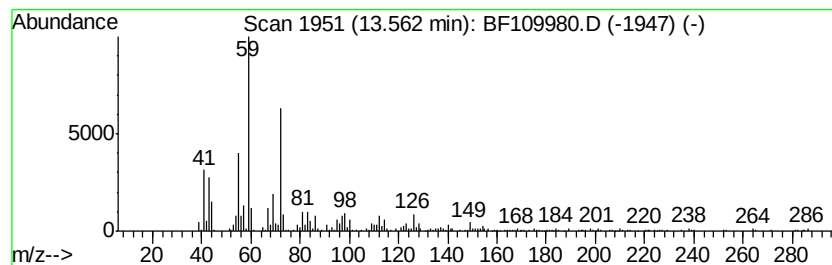
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 9-Octadecenamide, (Z)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.56	6.48 ng	447614	Chrysene-d12	14.15		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	97
2		Benzeneethanamine, 2-fluoro-.bet...	229	C11H16FN03	061338-98-5	72
3		Tetradecanamide	227	C14H29NO	000638-58-4	72
4		Octanamide	143	C8H17NO	000629-01-6	56
5		Nonanamide	157	C9H19NO	001120-07-6	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101818\
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Instrument :
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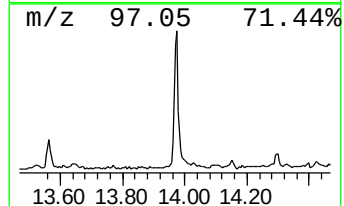
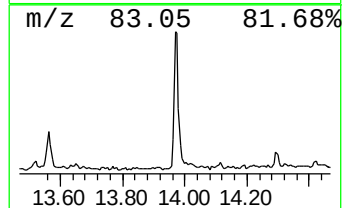
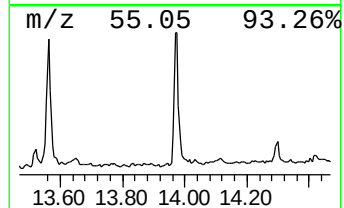
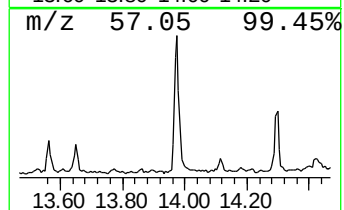
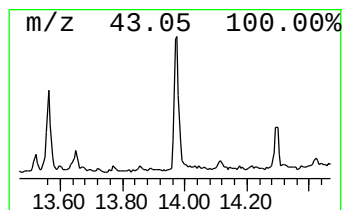
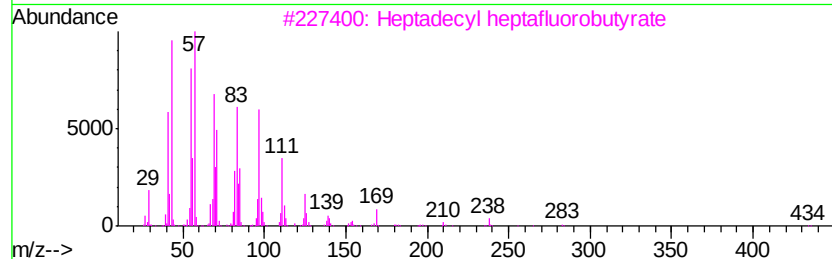
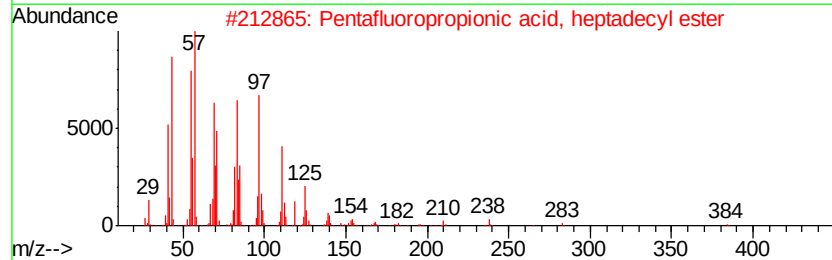
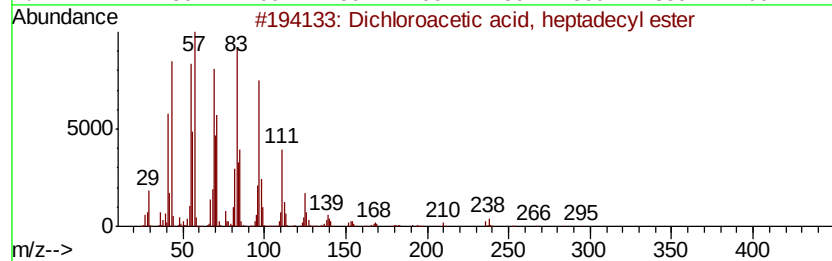
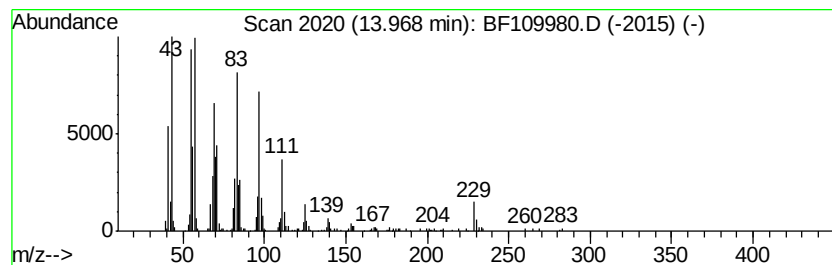
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Dichloroacetic acid, heptad... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.97	6.62 ng	457105	Chrysene-d12	14.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	95
2		Pentafluoropropionic acid, hepta...	402	C20H35F5O2	959218-78-1	94
3		Heptadecyl heptafluorobutyrat	452	C21H35F7O2	959085-66-6	94
4		Trifluoroacetoxv hexadecane	338	C18H33F3O2	006222-03-3	94
5		Heptafluorobutyric acid, hexadec...	438	C20H33F7O2	006385-15-5	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101818\
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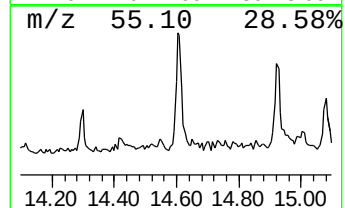
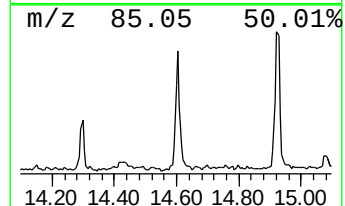
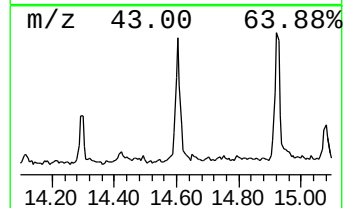
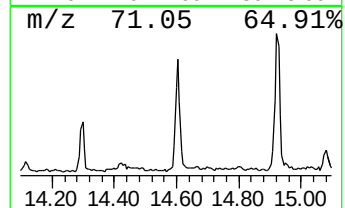
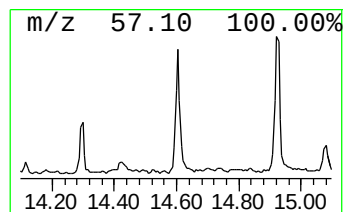
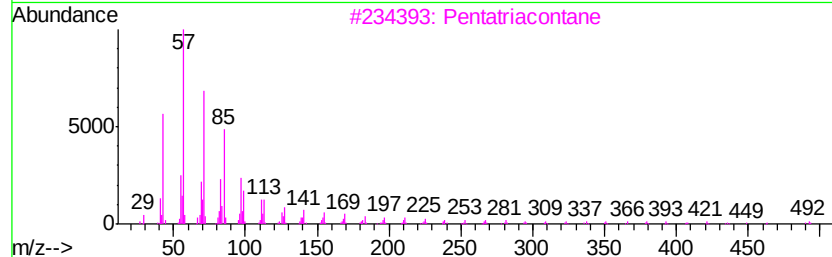
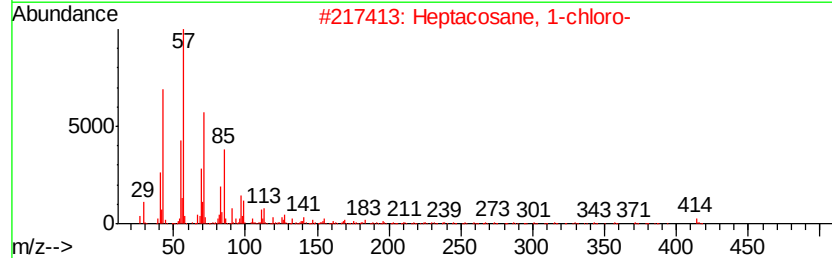
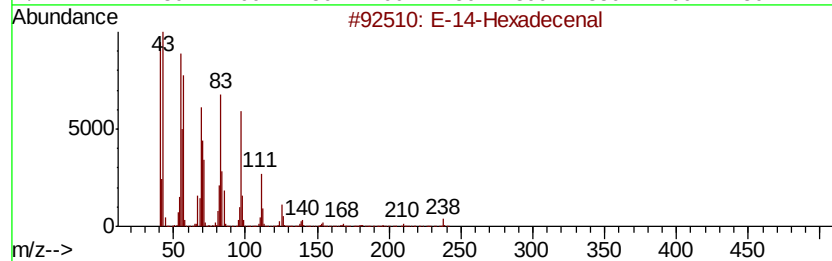
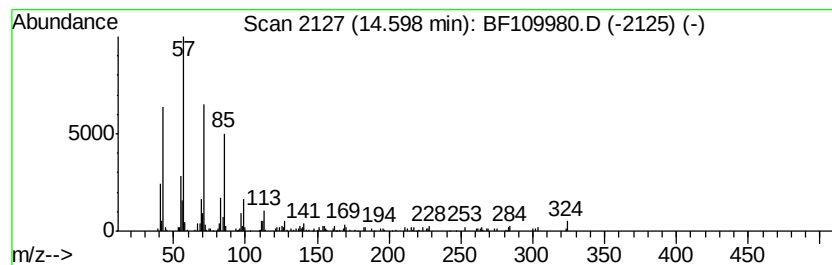
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 E-14-Hexadecenal Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.60	3.96 ng	273480	Chrysene-d12	14.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	E-14-Hexadecenal	238	C16H30O	330207-53-9	92
2		Heptacosane, 1-chloro-	414	C27H55Cl	062016-79-9	91
3		Pentatriacontane	493	C35H72	000630-07-9	91
4		Sulfurous acid, dodecyl 2-propyl...	292	C15H32O3S	1000309-12-3	91
5		Hexadecane, 1-chloro-	260	C16H33Cl	004860-03-1	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101818\
Data File : BF109980.D
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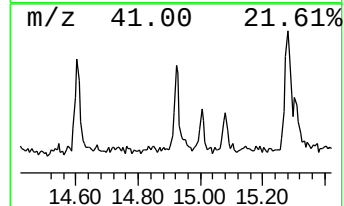
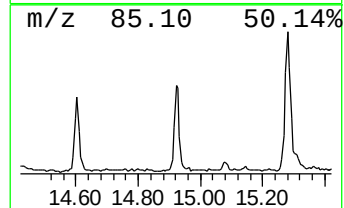
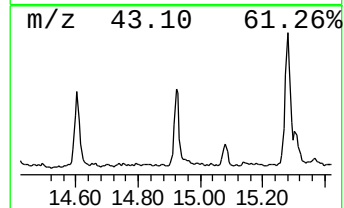
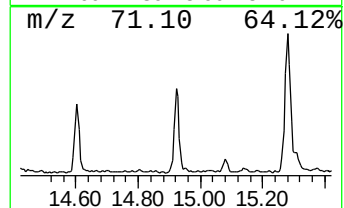
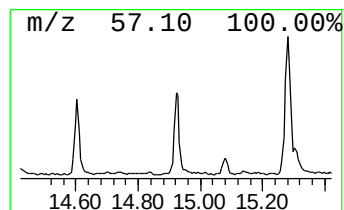
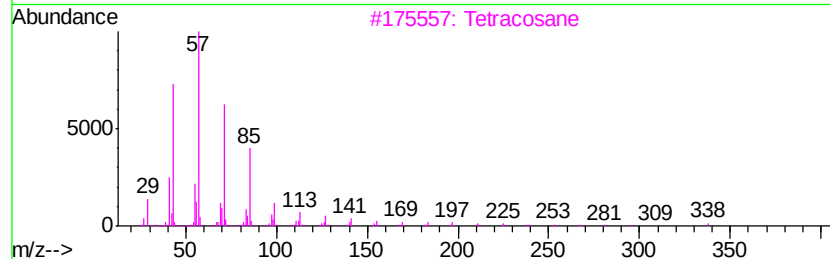
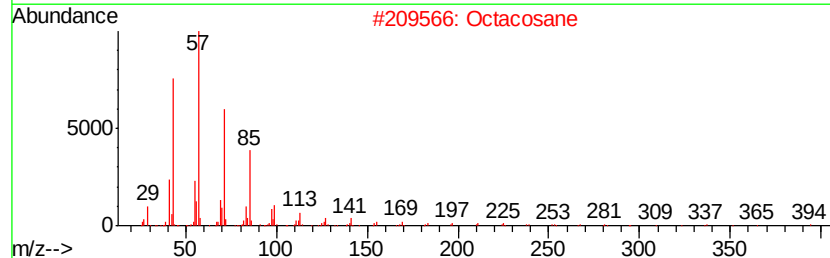
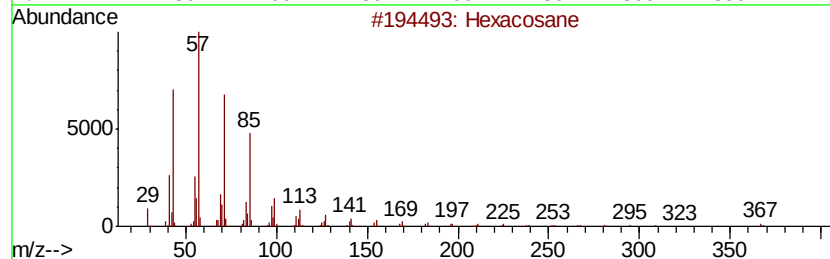
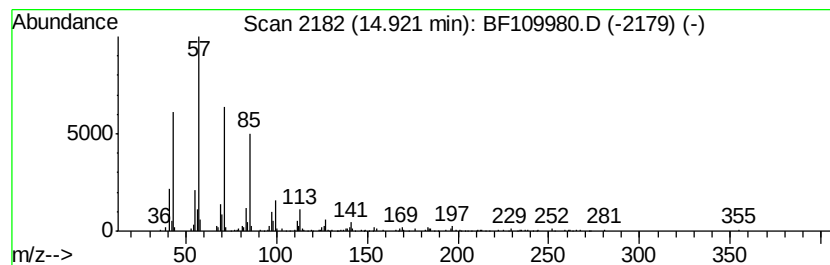
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Tetracosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.92	3.01 ng	255118	Perylene-d12	15.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	91
2		Octacosane	394	C28H58	000630-02-4	91
3		Tetracosane	338	C24H50	000646-31-1	90
4		Dodecane, 2-methyl-	184	C13H28	001560-97-0	86
5		Octadecane, 1-iodo-	380	C18H37I	000629-93-6	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101818\
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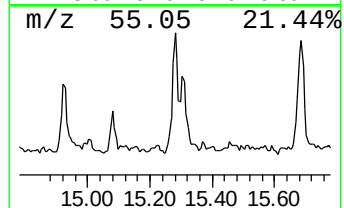
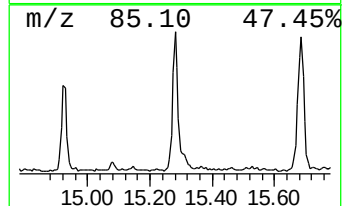
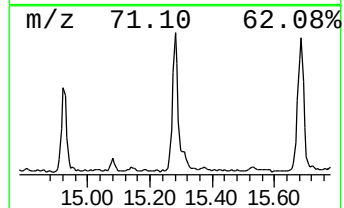
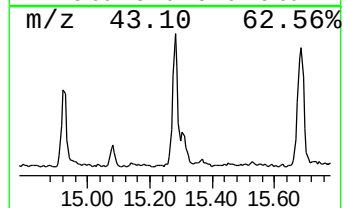
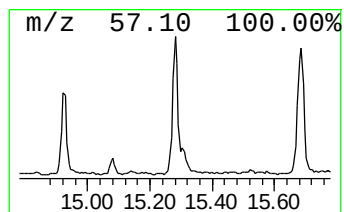
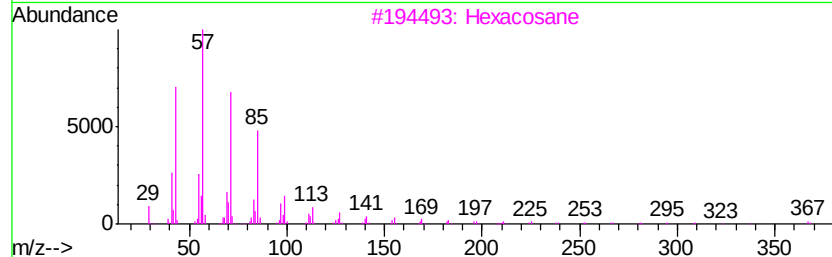
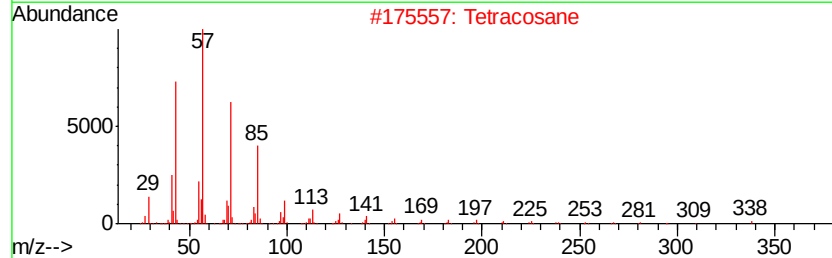
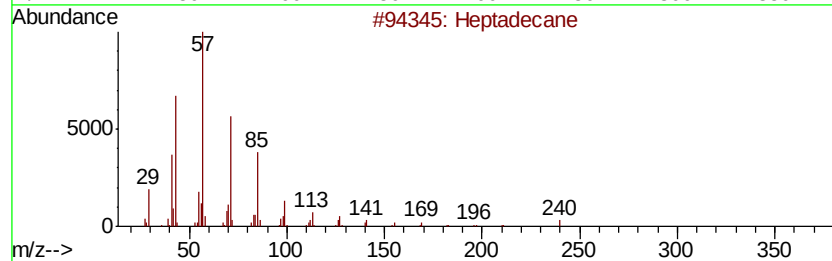
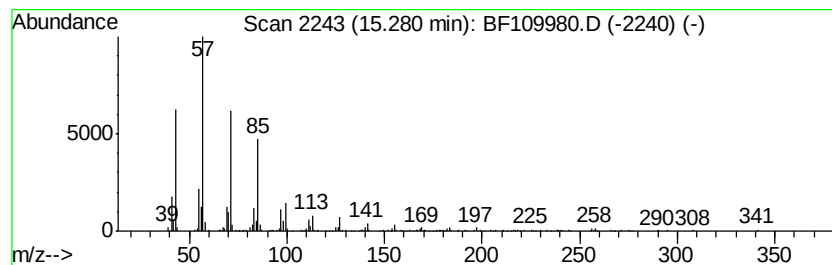
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Heptadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.28	5.05 ng	428504	Perylene-d12	15.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane	240	C17H36	000629-78-7	93
2		Tetracosane	338	C24H50	000646-31-1	91
3		Hexacosane	366	C26H54	000630-01-3	91
4		2-methyloctacosane	408	C29H60	1000376-72-8	91
5		Octacosane	394	C28H58	000630-02-4	91



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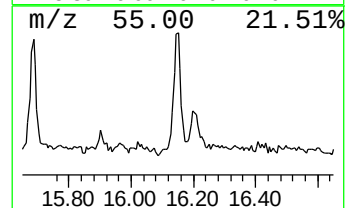
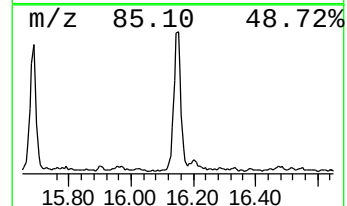
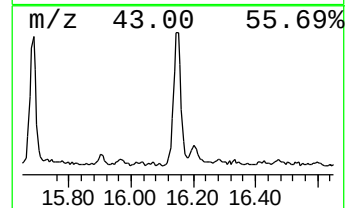
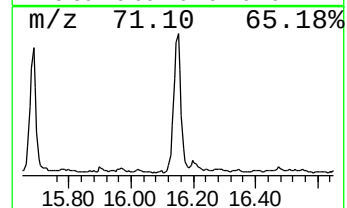
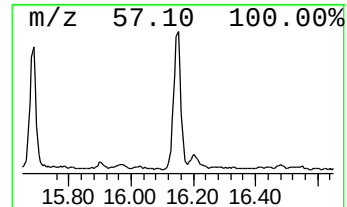
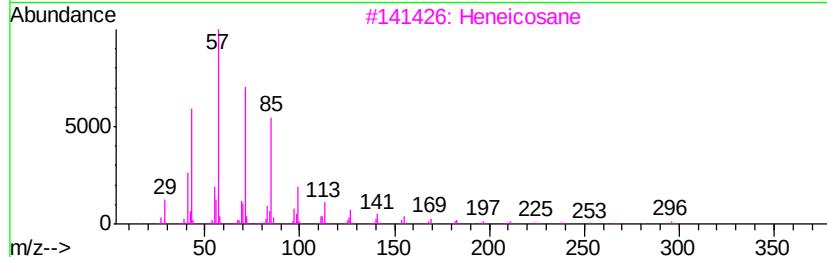
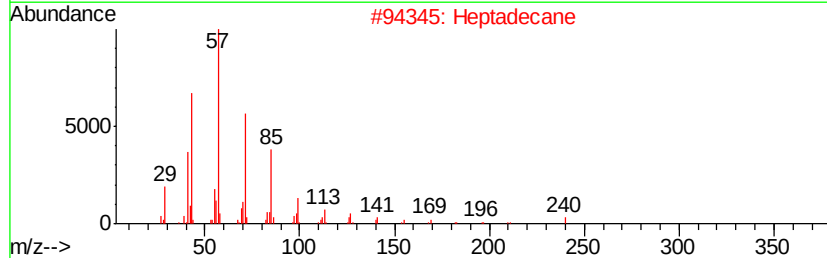
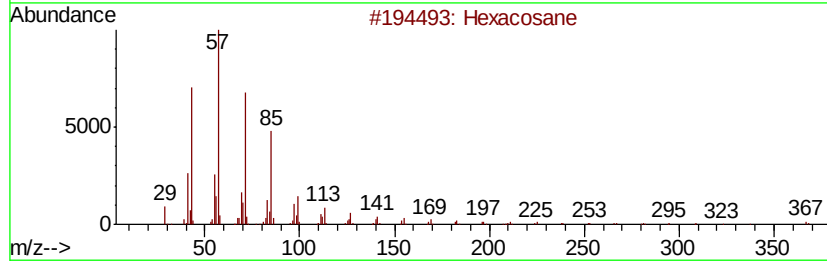
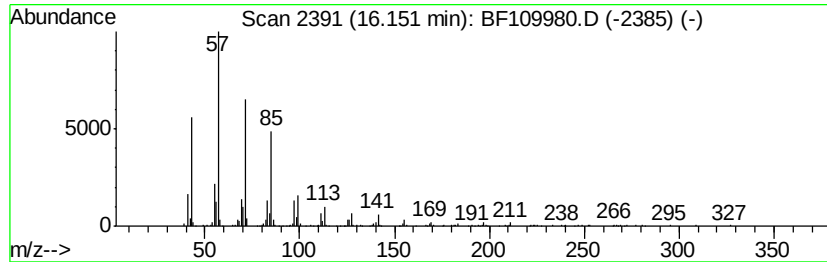
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Hexacosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.15	6.80 ng	576878	Perylene-d12	15.68

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosane	366	C26H54	000630-01-3	94
2			Heptadecane	240	C17H36	000629-78-7	93
3			Heneicosane	296	C21H44	000629-94-7	91
4			2-methyloctacosane	408	C29H60	1000376-72-8	91
5			Octacosane	394	C28H58	000630-02-4	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101818\
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 Acq On : 19 Oct 2018 1:02
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 Sample : J5570-02
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

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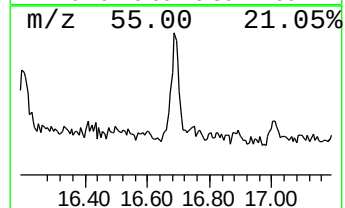
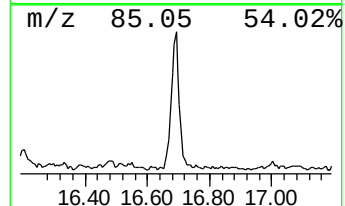
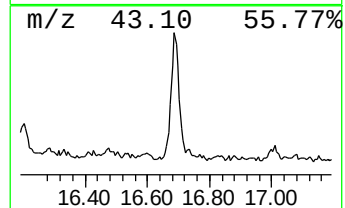
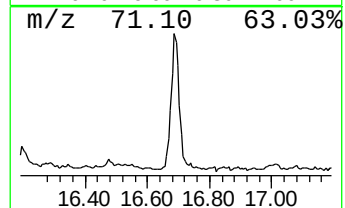
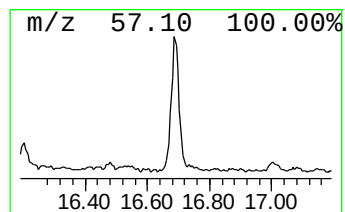
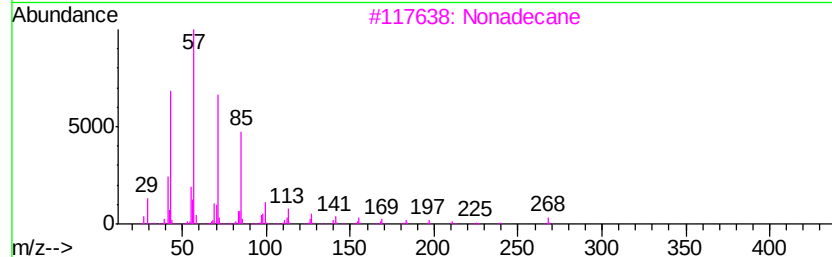
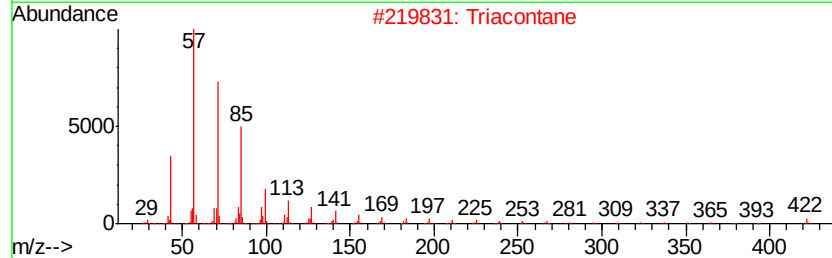
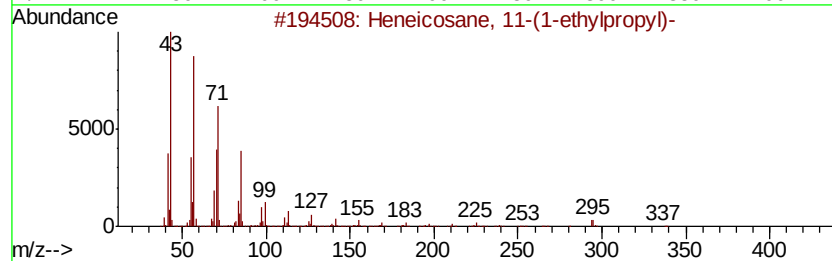
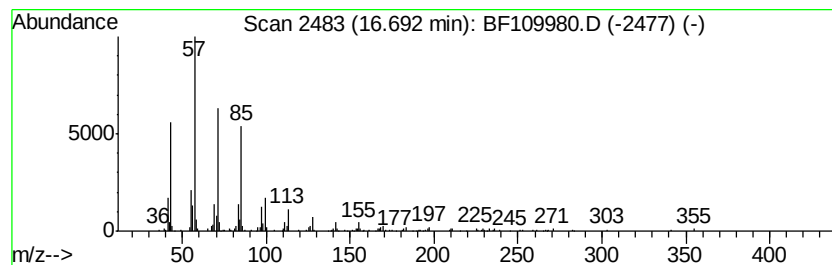
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 14 Heneicosane, 11-(1-ethylpro... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.69	3.70 ng	313310	Perylene-d12	15.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91
2		Triacontane	422	C30H62	000638-68-6	90
3		Nonadecane	268	C19H40	000629-92-5	87
4		Docosane	310	C22H46	000629-97-0	87
5		Hexacosane	366	C26H54	000630-01-3	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101818\
Data File : BF109980.D
Acq On : 19 Oct 2018 1:02
Operator : JU/SJ
Sample : J5570-02
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TS-07

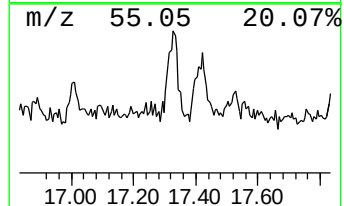
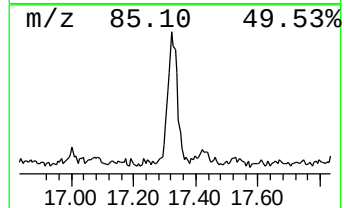
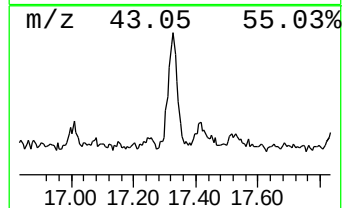
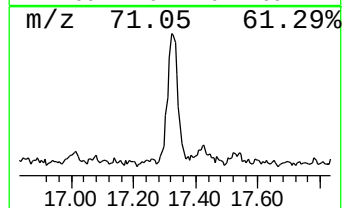
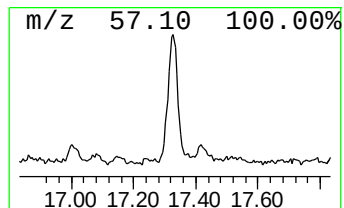
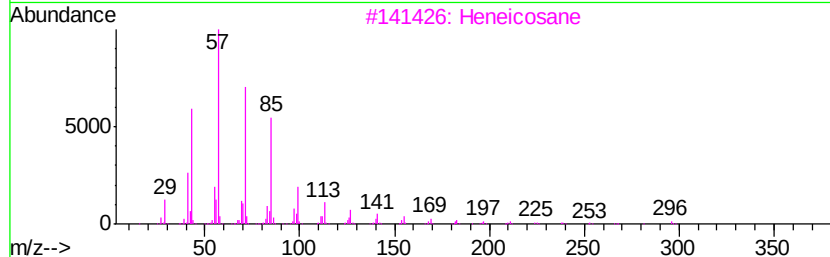
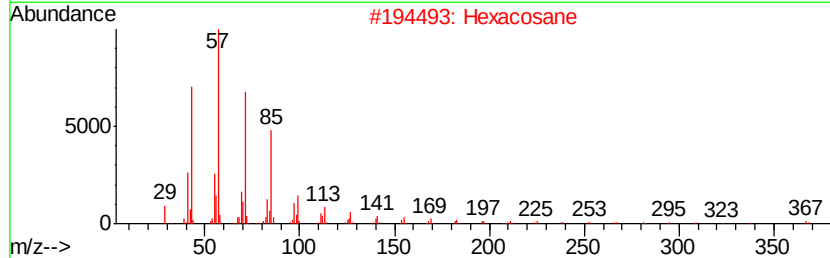
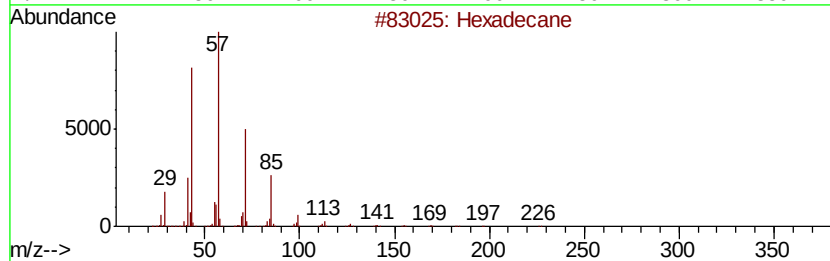
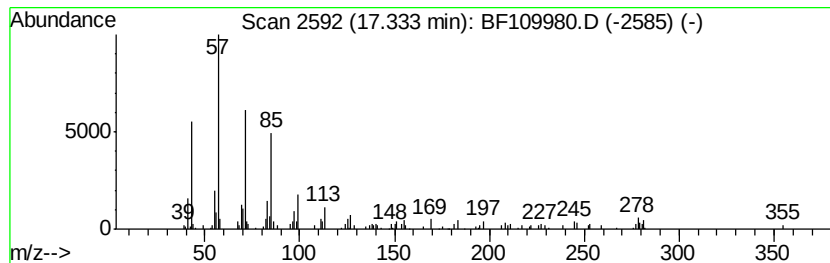
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Heptadecane, 9-octyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.33	2.35 ng	198879	Perylene-d12	15.68

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	91
2			Hexacosane	366	C26H54	000630-01-3	87
3			Heneicosane	296	C21H44	000629-94-7	87
4			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87
5			Eicosane	282	C20H42	000112-95-8	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101818\
Data File : BF109980.D
Acq On : 19 Oct 2018 1:02
Operator : JU/SJ
Sample : J5570-02
Misc :
ALS Vial : 17 Sample Multiplier: 1

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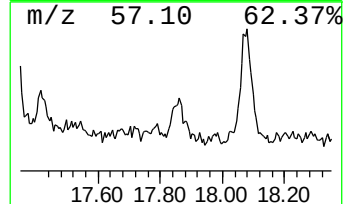
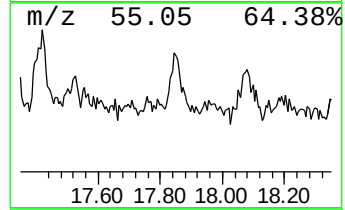
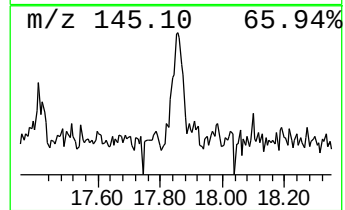
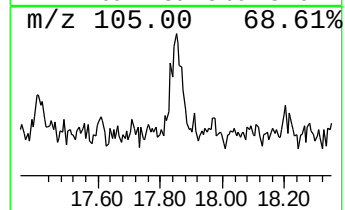
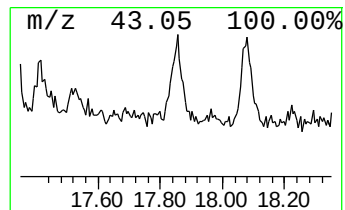
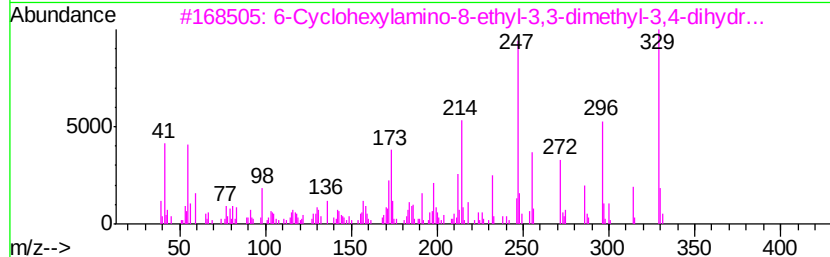
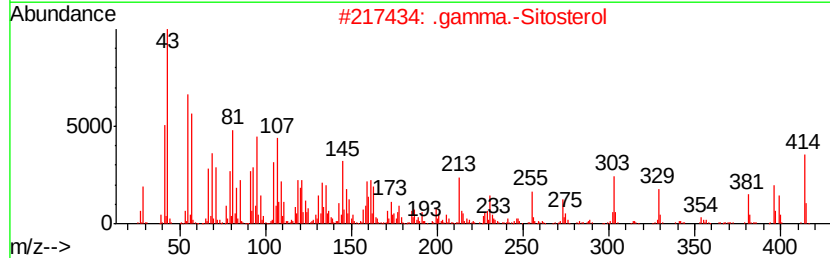
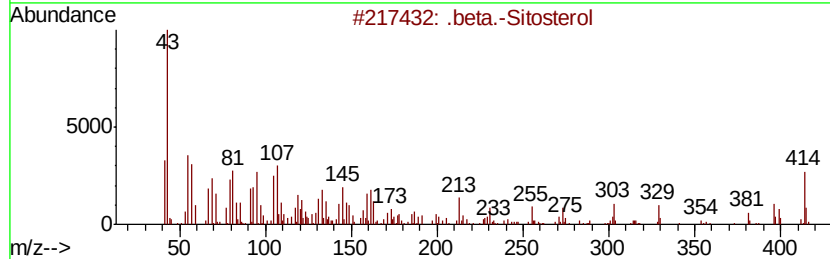
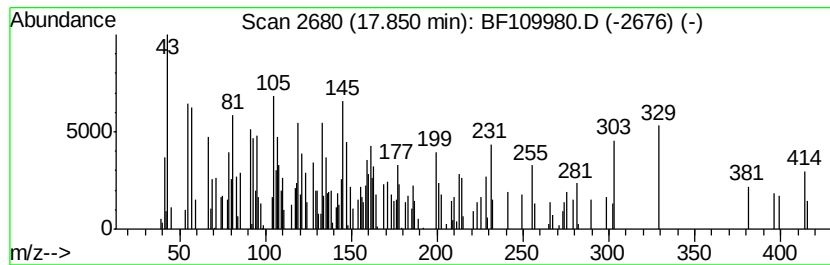
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 unknown17.85 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.85	2.24 ng	190182	Perylene-d12	15.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	.beta.-Sitosterol	414	C29H50O	000083-46-5	43
2		.gamma.-Sitosterol	414	C29H50O	000083-47-6	25
3		6-Cyclohexylamino-8-ethyl-3,3-di...	329	C19H27N3S	1000274-39-5	10
4		Ethanone, 1-(2,5-dimethyl-4-phen...	214	C14H14O2	088484-88-2	9
5		Propanoic acid, 2-(benzoylamino)...	402	C18H15F5N2O3	1000351-47-2	9



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF101818\
 Data File : BF109980.D
 Acq On : 19 Oct 2018 1:02
 Operator : JU/SJ
 Sample : J5570-02
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101518.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
Butane, 2-methoxy...	2.36	13.8	ng	939178	1	6.97	1362190	20.0
unknown3.62	3.62	2.3	ng	159082	1	6.97	1362190	20.0
2-Pentanone, 4-hy...	5.24	4.5	ng	307876	1	6.97	1362190	20.0
unknown6.75	6.75	92.3	ng	6285820	1	6.97	1362190	20.0
n-Hexadecanoic acid	12.02	5.4	ng	421132	4	11.51	1556820	20.0
2-Phenanthrenol, ...	13.52	2.8	ng	191572	5	14.15	1381080	20.0
9-Octadecenamide, ...	13.56	6.5	ng	447614	5	14.15	1381080	20.0
Dichloroacetic ac...	13.97	6.6	ng	457105	5	14.15	1381080	20.0
E-14-Hexadecenal	14.60	4.0	ng	273480	5	14.15	1381080	20.0
Tetracosane	14.92	3.0	ng	255118	6	15.68	1695610	20.0
Heptadecane	15.28	5.0	ng	428504	6	15.68	1695610	20.0
Hexacosane	16.15	6.8	ng	576878	6	15.68	1695610	20.0
Heneicosane, 11-(...	16.69	3.7	ng	313310	6	15.68	1695610	20.0
Heptadecane, 9-oc...	17.33	2.4	ng	198879	6	15.68	1695610	20.0
unknown17.85	17.85	2.2	ng	190182	6	15.68	1695610	20.0