

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF100918\
 Data File : BF109705.D
 Acq On : 9 Oct 2018 19:59
 Operator : JU/SJ
 Sample : J5324-01 10X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OILY-DEBRIS

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-BF100818.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	5.316	545	549	564	rBV	29604	75071	13.30%	1.432%
2	6.351	721	725	739	rBV	37243	120004	21.26%	2.288%
3	6.469	742	745	755	rVV	90845	132889	23.54%	2.534%
4	6.692	780	783	793	rBV	284974	400166	70.89%	7.631%
5	6.763	793	795	806	rVV3	22800	51211	9.07%	0.977%
6	6.851	806	810	830	rVB2	93116	157293	27.87%	2.999%
7	7.263	877	880	892	rBV2	26022	68334	12.11%	1.303%
8	7.522	922	924	935	rVB	10282	15198	2.69%	0.290%
9	7.786	965	969	980	rBV3	23285	57917	10.26%	1.104%
10	7.975	998	1001	1028	rBV	502106	564454	100.00%	10.764%
11	8.339	1059	1063	1073	rBV2	8318	17755	3.15%	0.339%
12	8.869	1150	1153	1160	rBV2	8582	13317	2.36%	0.254%
13	9.051	1179	1184	1193	rBV	74699	91319	16.18%	1.741%
14	9.210	1208	1211	1224	rBV	71332	131668	23.33%	2.511%
15	9.445	1247	1251	1256	rBV2	17642	30077	5.33%	0.574%
16	9.604	1273	1278	1279	rBV3	10179	14593	2.59%	0.278%
17	9.639	1281	1284	1287	rVB2	16840	15699	2.78%	0.299%
18	9.669	1287	1289	1291	rBV3	8291	10174	1.80%	0.194%
19	9.722	1293	1298	1304	rBV	558570	555849	98.48%	10.600%
20	9.886	1324	1326	1328	rBV2	9794	10658	1.89%	0.203%
21	9.986	1341	1343	1345	rBV2	17725	16623	2.94%	0.317%
22	10.063	1350	1356	1358	rBV6	22876	39942	7.08%	0.762%
23	10.098	1359	1362	1365	rBV4	18331	23396	4.14%	0.446%
24	10.127	1365	1367	1370	rBV2	41074	43100	7.64%	0.822%
25	10.186	1374	1377	1378	rBV3	17636	17148	3.04%	0.327%
26	10.280	1391	1393	1396	rVB3	26364	24745	4.38%	0.472%
27	10.516	1430	1433	1436	rBV3	38763	53326	9.45%	1.017%
28	10.592	1444	1446	1451	rBV3	99023	152926	27.09%	2.916%
29	10.651	1453	1456	1460	rVB	103969	114173	20.23%	2.177%
30	11.110	1532	1534	1538	rVB	134444	134349	23.80%	2.562%
31	11.204	1547	1550	1562	rVB	389136	457268	81.01%	8.720%
32	12.257	1727	1729	1734	rVB2	89661	83103	14.72%	1.585%
33	12.780	1816	1818	1823	rVB	95117	85350	15.12%	1.628%
34	13.827	1993	1996	2006	rVB	490135	549794	97.40%	10.484%

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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 >
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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	14.304	2075	2077	2080	rVB	26847	21734	3.85%	0.414%
36	14.598	2124	2127	2133	rVB	66422	67971	12.04%	1.296%
37	14.904	2176	2179	2183	rVB	90551	98953	17.53%	1.887%
38	15.227	2229	2234	2237	rBV	293538	429647	76.12%	8.193%
39	15.651	2302	2306	2314	rVB	109546	162745	28.83%	3.103%
40	16.109	2379	2384	2388	rVB	60634	101056	17.90%	1.927%
41	16.639	2471	2474	2478	rBV4	19585	33030	5.85%	0.630%

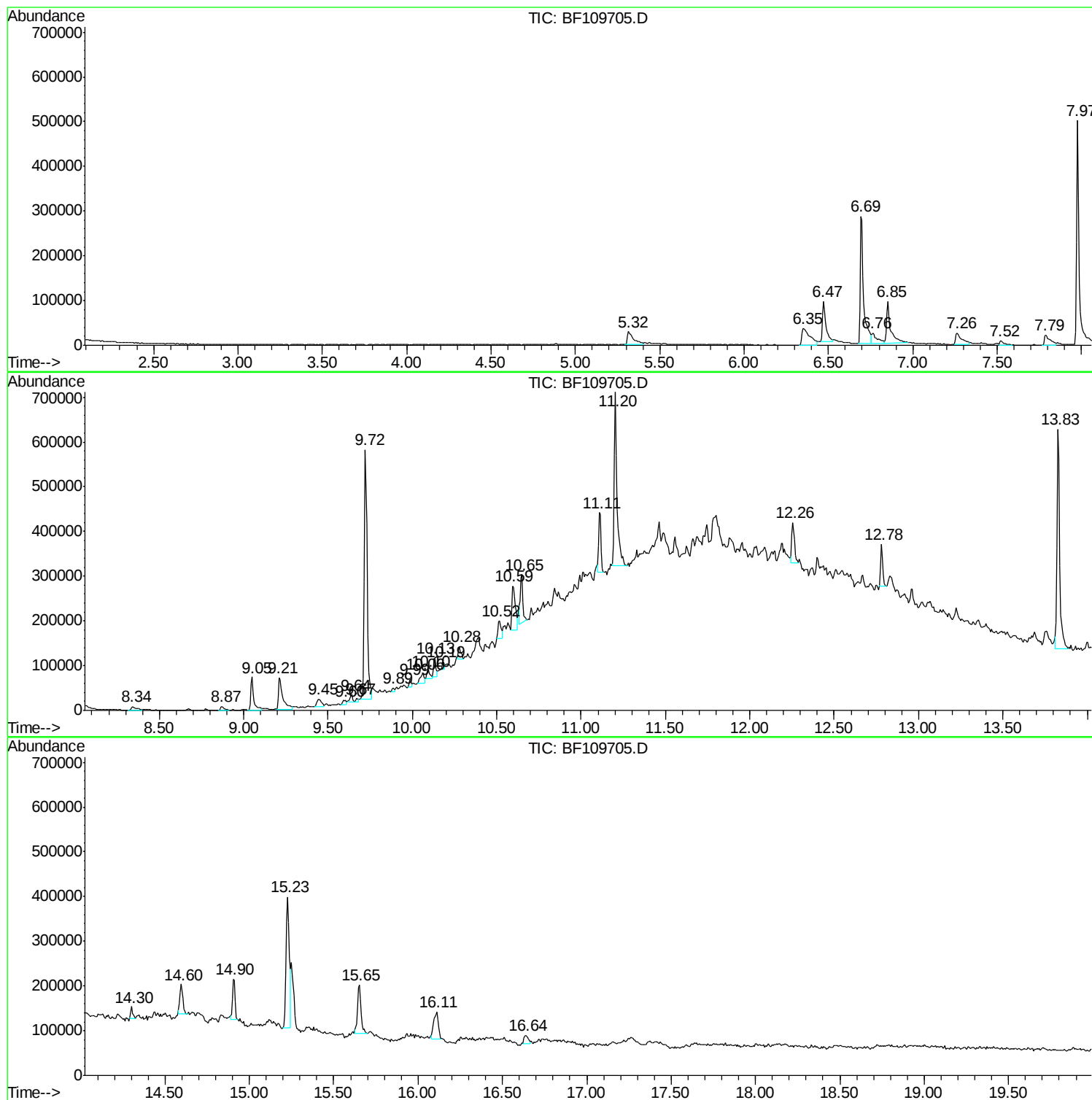
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF100818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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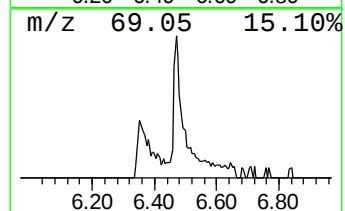
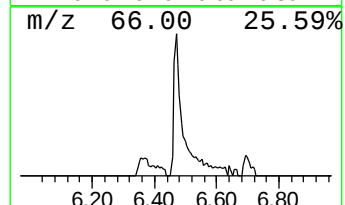
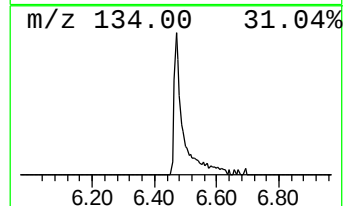
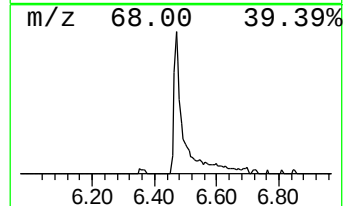
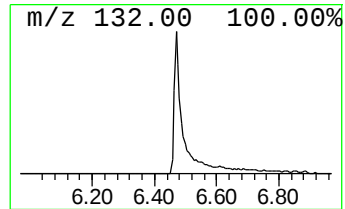
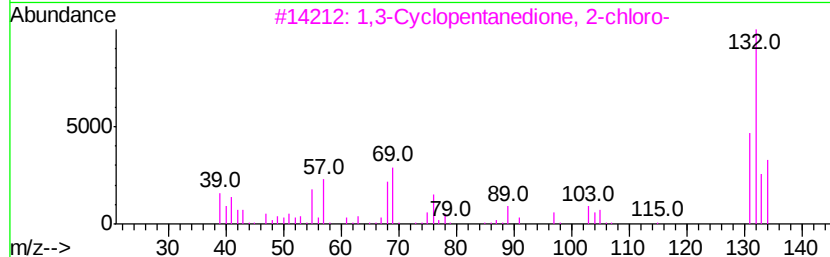
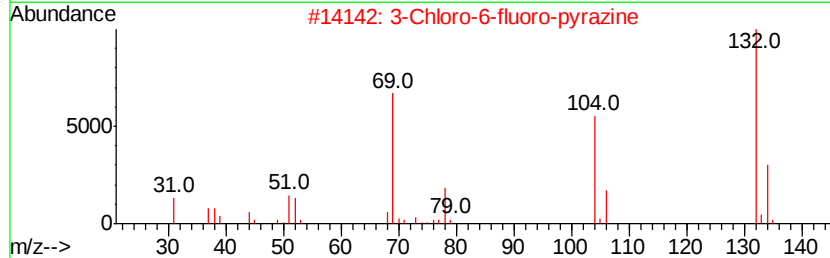
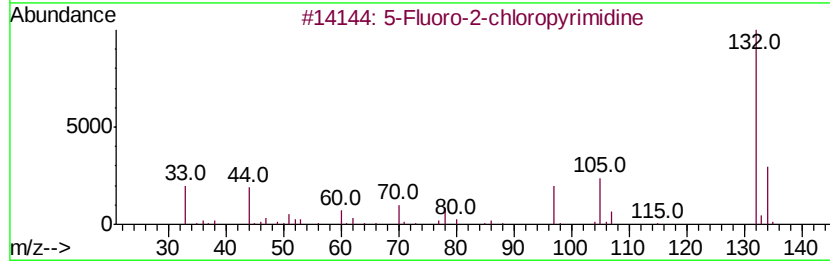
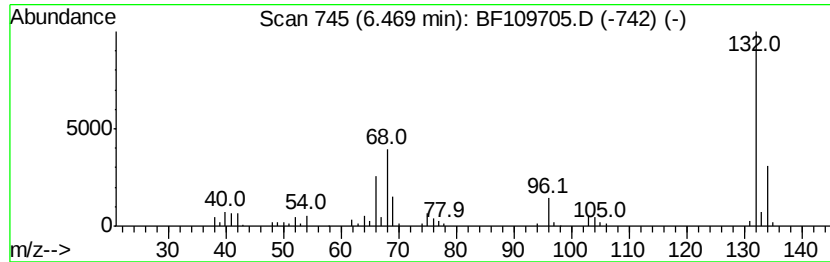
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF100818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown6.47 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.47	6.64 ng	132889	1,4-Dichlorobenzene-d4	6.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	17
4		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
5		5-Aminoindole	132	C8H8N2	005192-03-0	12



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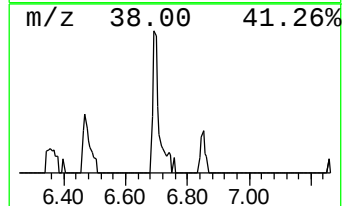
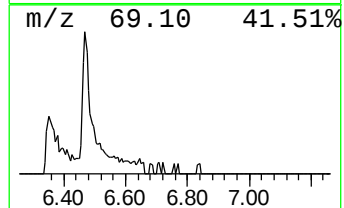
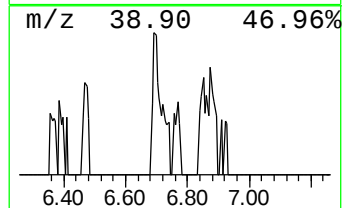
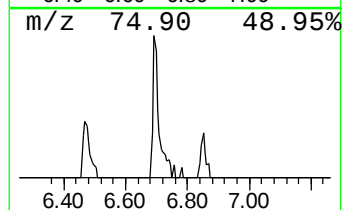
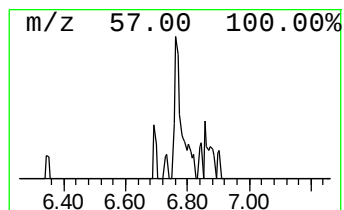
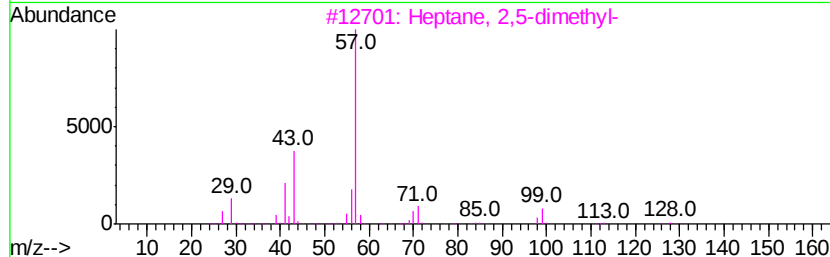
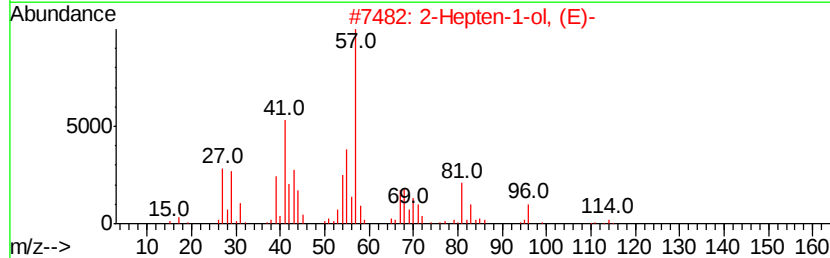
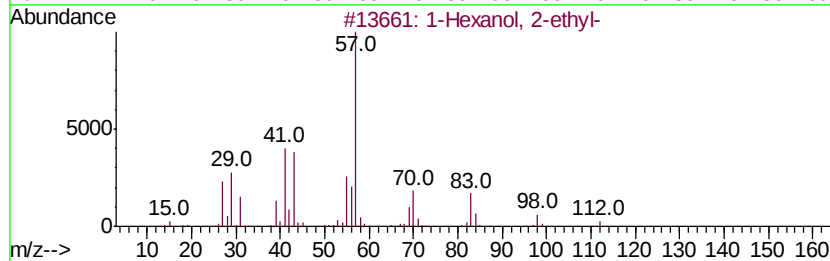
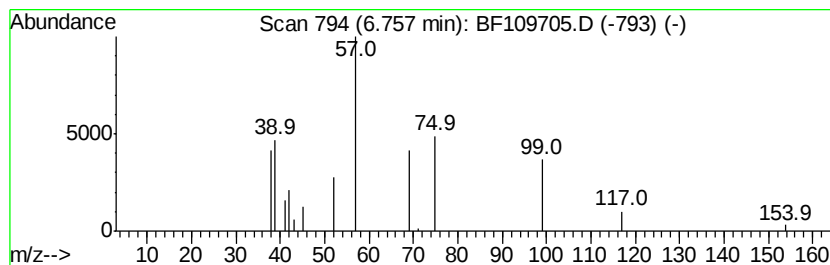
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 1-Hexanol, 2-ethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.76	2.56 ng	51211	1,4-Dichlorobenzene-d4	6.70

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	59
2			2-Hepten-1-ol, (E)-	114	C7H14O	033467-76-4	28
3			Heptane, 2,5-dimethyl-	128	C9H20	002216-30-0	9
4			Pentane, 3,3-diethyl-	128	C9H20	001067-20-5	9
5			2-Propyl-1-pentanol	130	C8H18O	058175-57-8	9



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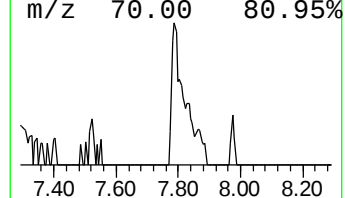
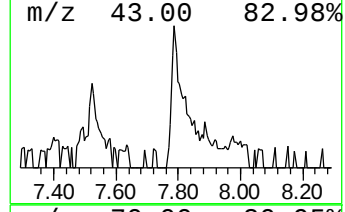
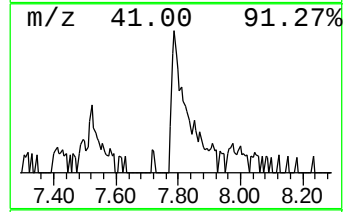
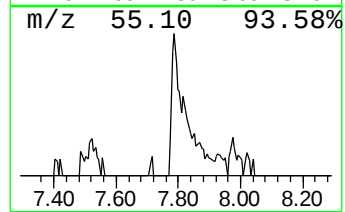
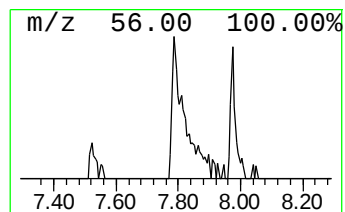
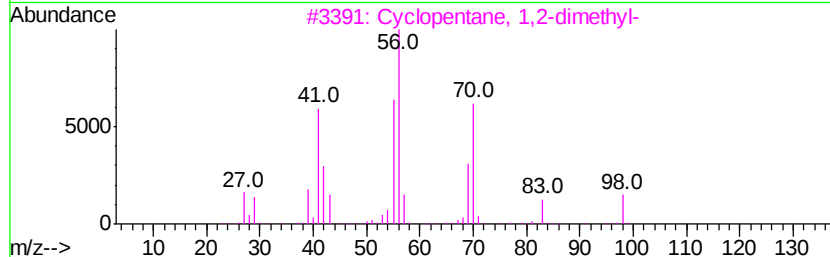
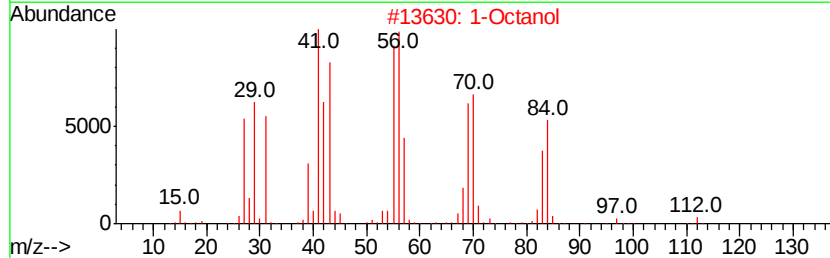
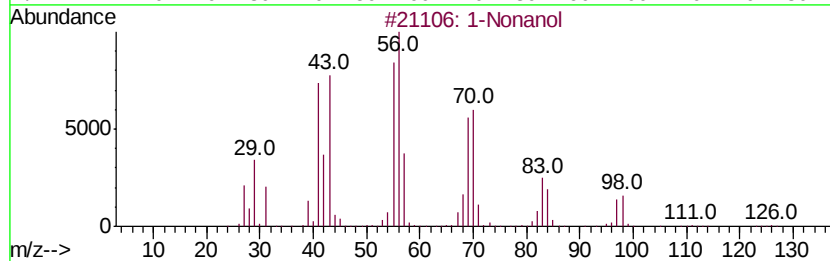
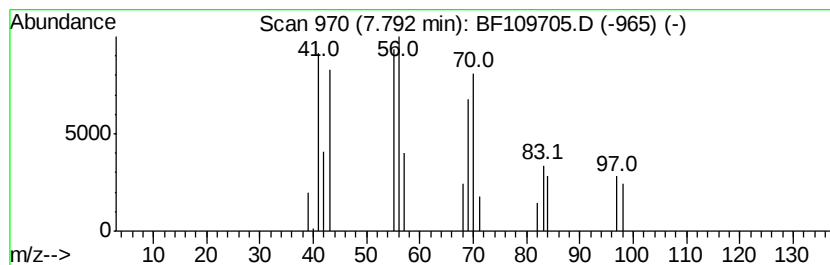
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 1-Nonanol Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.79	2.05 ng	57917	Napthalene-d8	7.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Nonanol	144	C9H20O	000143-08-8	90
2		1-Octanol	130	C8H18O	000111-87-5	80
3		Cyclopentane, 1,2-dimethyl-	98	C7H14	002452-99-5	64
4		1-Heptene	98	C7H14	000592-76-7	64
5		Formic acid, heptyl ester	144	C8H16O2	000112-23-2	53



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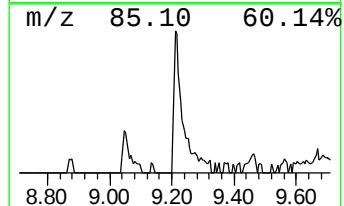
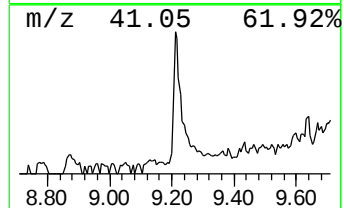
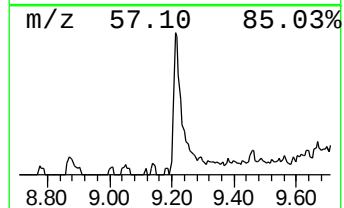
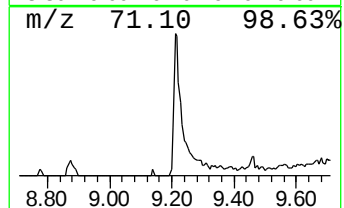
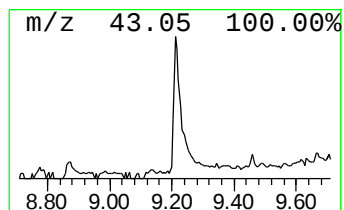
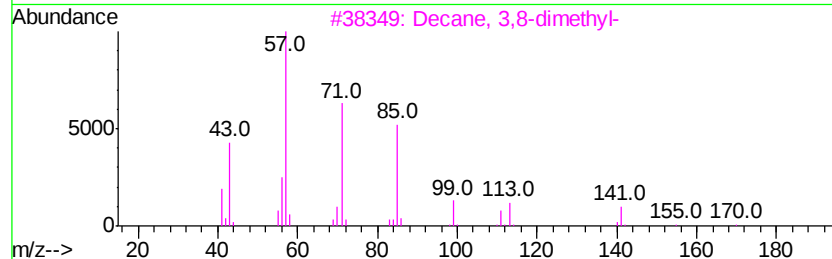
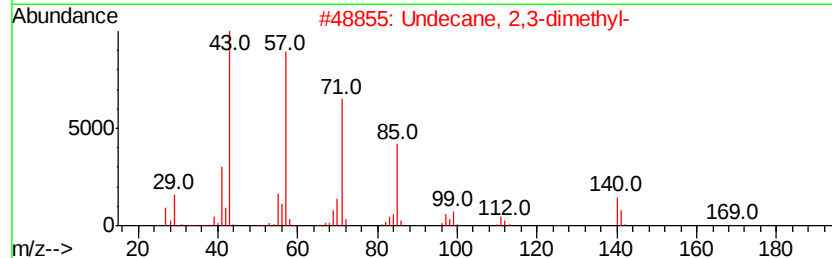
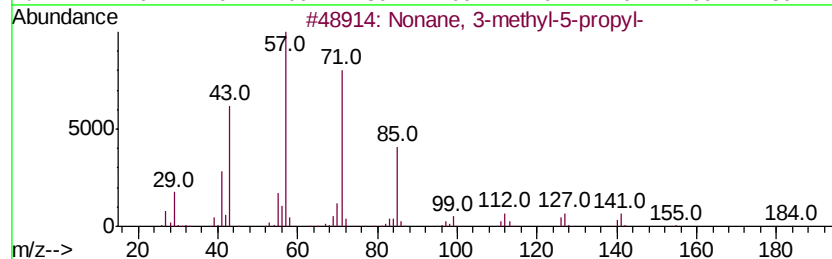
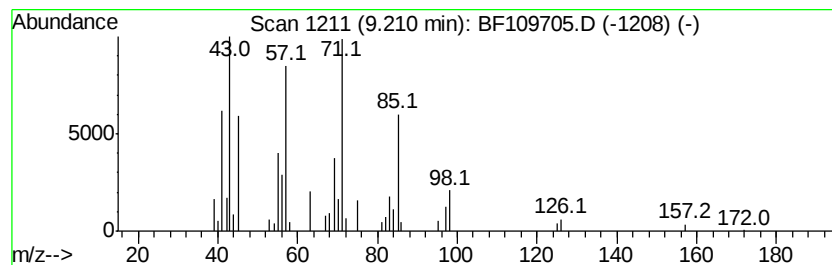
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 unknown9.21 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.21	4.74 ng	131668	Acenaphthene-d10	9.72

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonane, 3-methyl-5-propyl-	184	C13H28	031081-18-2	47
2		Undecane, 2,3-dimethyl-	184	C13H28	017312-77-5	47
3		Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	43
4		Nonane, 1-iodo-	254	C9H19I	004282-42-2	43
5		Sulfurous acid, hexyl pentyl ester	236	C11H24O3S	1000309-14-1	43



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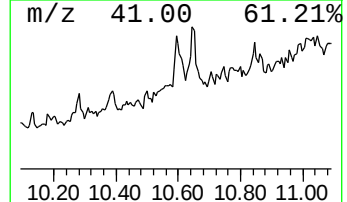
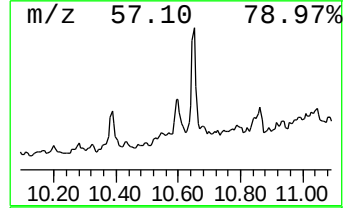
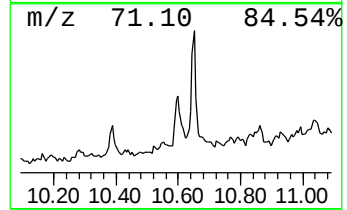
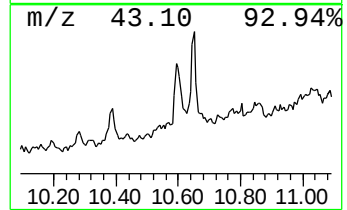
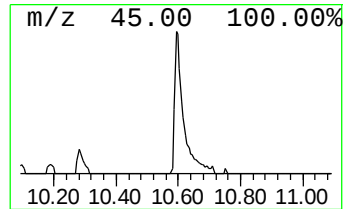
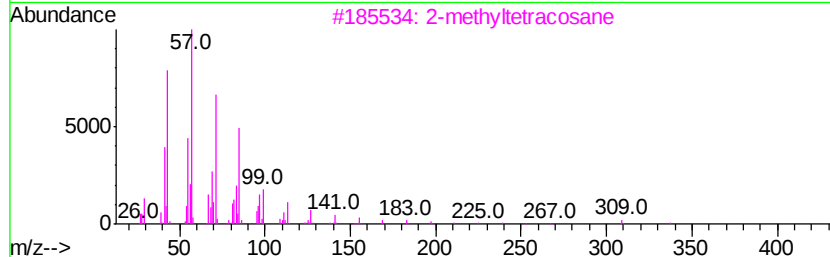
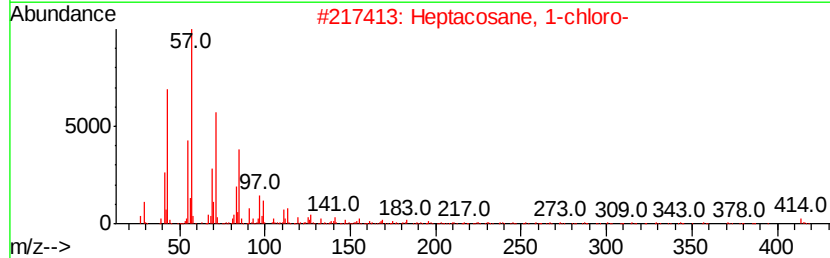
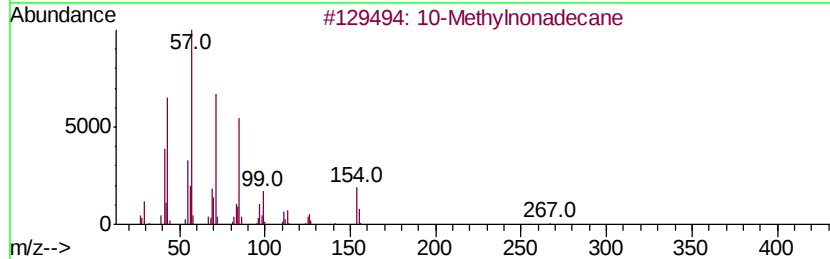
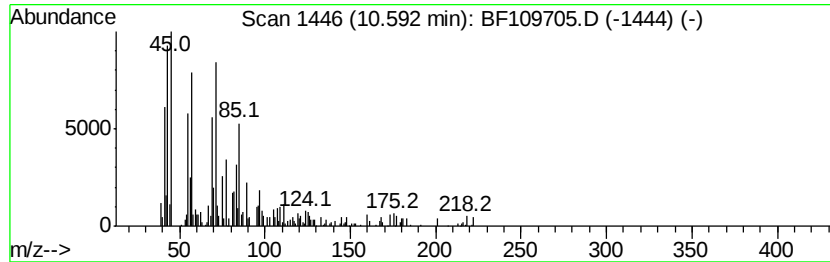
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 unknown10.59 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.59	6.69 ng	152926	Phenanthrene-d10	11.20

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	10-Methylnonadecane		282	C20H42	056862-62-5	43
2	Heptacosane, 1-chloro-		414	C27H55Cl	062016-79-9	43
3	2-methyltetracosane		352	C25H52	1000376-72-6	43
4	Methoxyacetic acid, pentadecyl e...		300	C18H36O3	1000281-82-2	43
5	Octadecane, 1-chloro-		288	C18H37Cl	003386-33-2	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100918\
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Operator : JU/SJ
Sample : J5324-01 10X
Misc :
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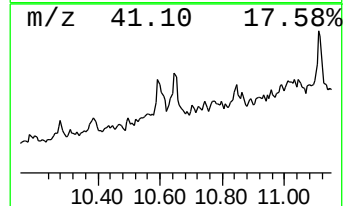
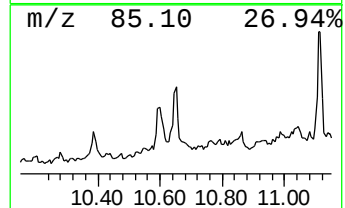
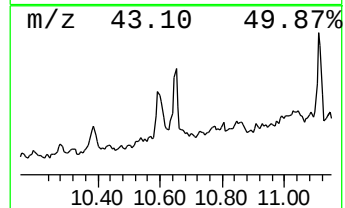
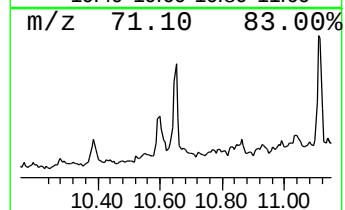
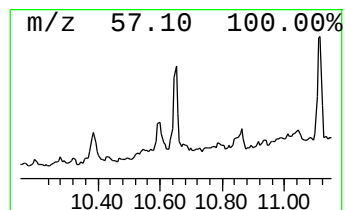
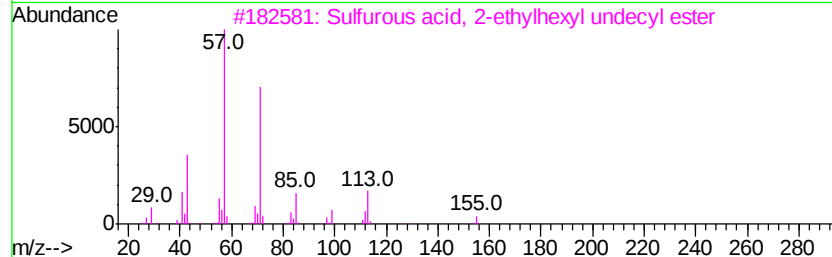
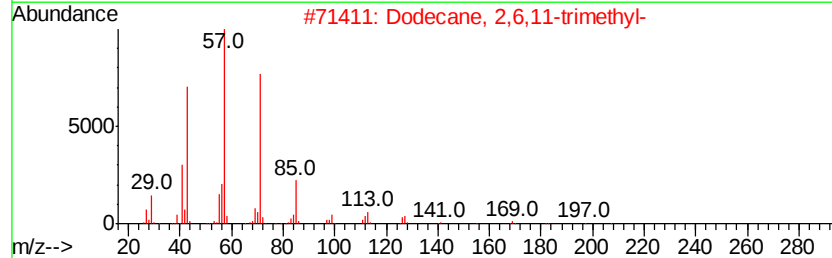
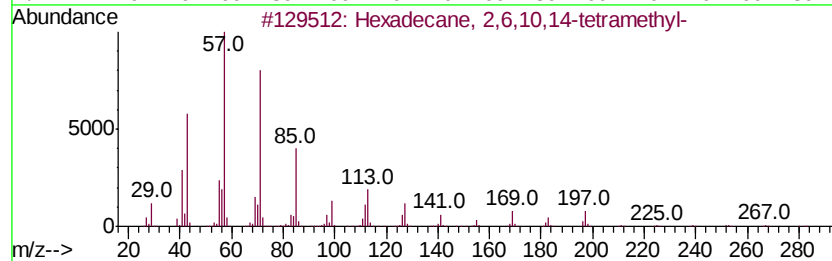
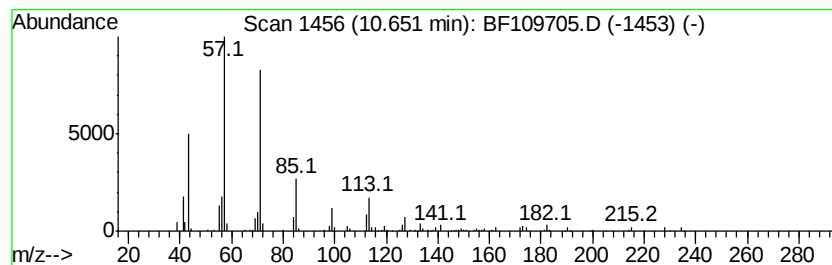
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Hexadecane, 2,6,10,14-tetra... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.65	4.99 ng	114173	Phenanthrene-d10	11.20	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	78
2	Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	72
3	Sulfurous acid, 2-ethylhexyl und...	348	C19H40O3S	1000309-19-4	72
4	Tridecane	184	C13H28	000629-50-5	59
5	Undecane, 6,6-dimethyl-	184	C13H28	017312-76-4	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100918\
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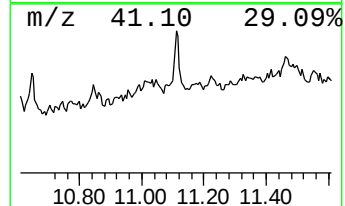
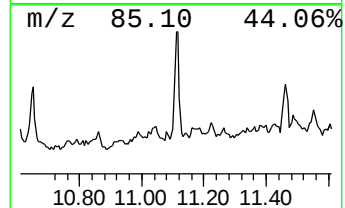
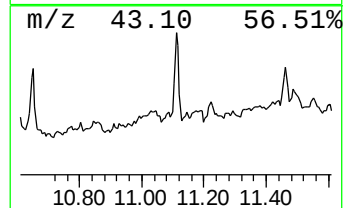
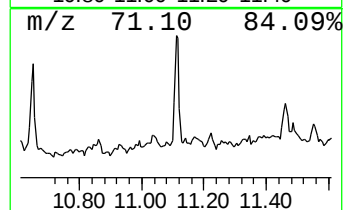
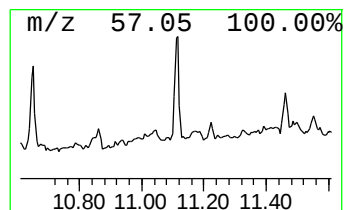
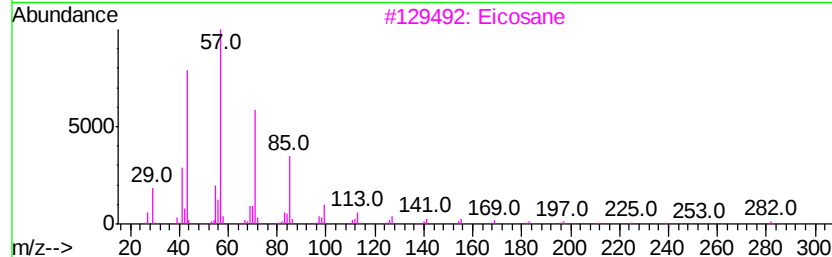
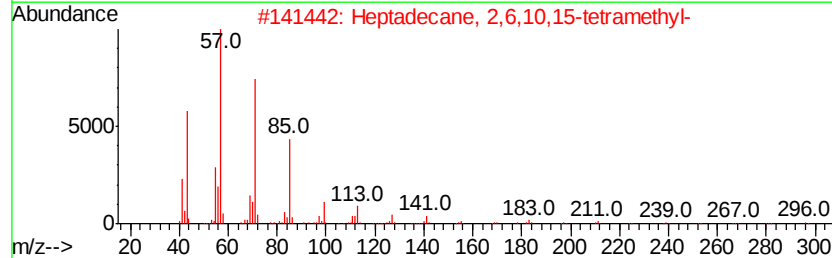
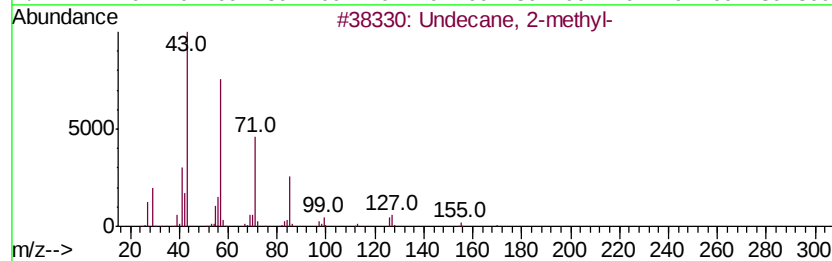
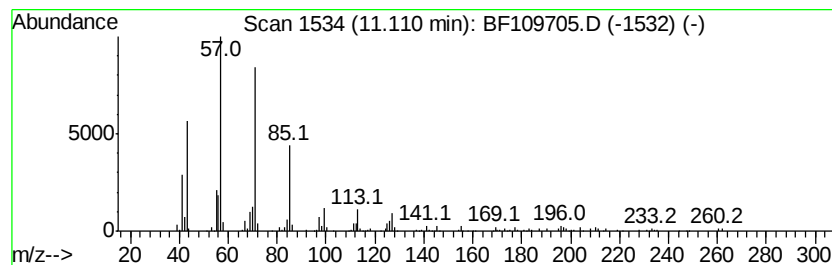
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Undecane, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.11	5.88 ng	134349	Phenanthrene-d10	11.20	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane, 2-methyl-	170	C12H26	007045-71-8	86
2	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	86
3	Eicosane	282	C20H42	000112-95-8	86
4	Heptacosane	380	C27H56	000593-49-7	86
5	Heptadecane, 8-methyl-	254	C18H38	013287-23-5	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100918\
Data File : BF109705.D
Acq On : 9 Oct 2018 19:59
Operator : JU/SJ
Sample : J5324-01 10X
Misc :
ALS Vial : 9 Sample Multiplier: 1

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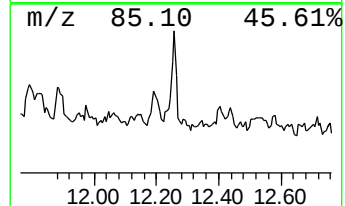
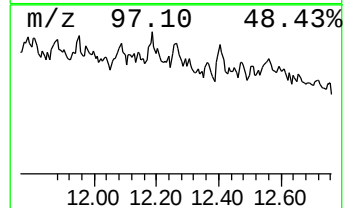
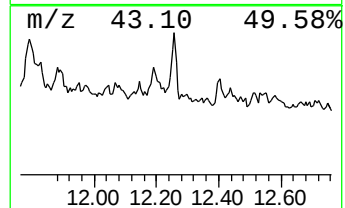
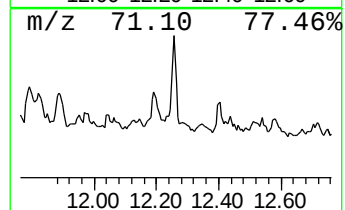
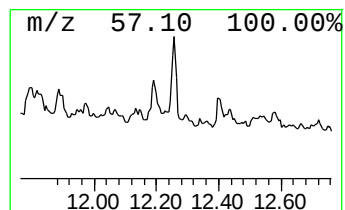
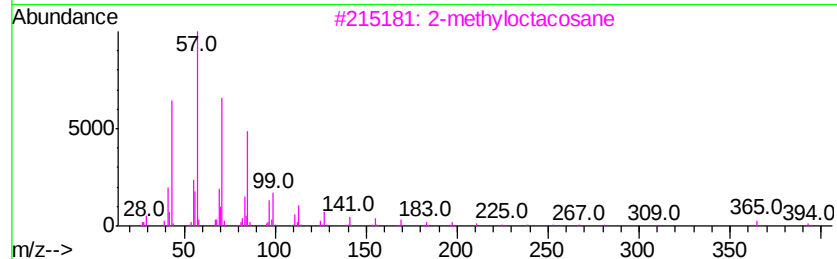
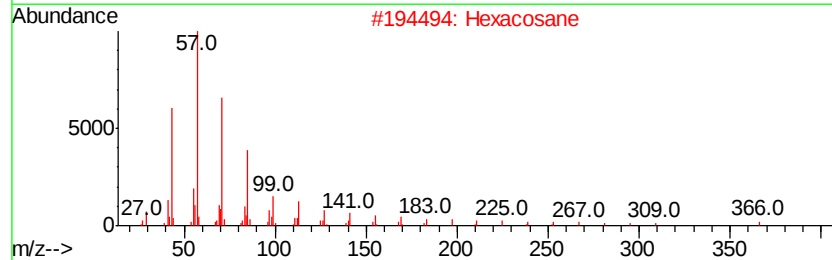
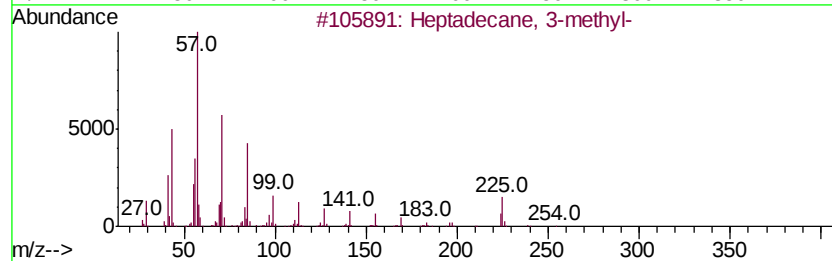
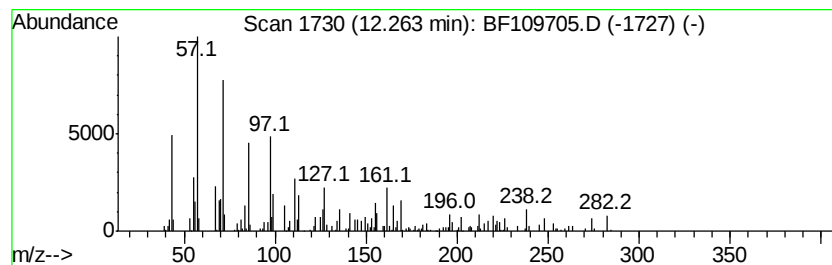
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Heptadecane, 3-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.26	3.63 ng	83103	Phenanthrene-d10	11.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 3-methyl-	254	C18H38	006418-44-6	91
2		Hexacosane	366	C26H54	000630-01-3	91
3		2-methyloctacosane	408	C29H60	1000376-72-8	90
4		Octadecane, 3-ethyl-5-(2-ethylbu...	366	C26H54	055282-12-7	90
5		Pentacosane	352	C25H52	000629-99-2	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100918\
Data File : BF109705.D
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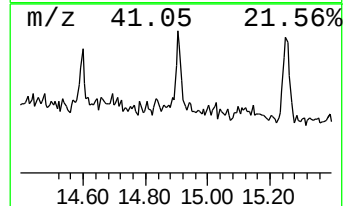
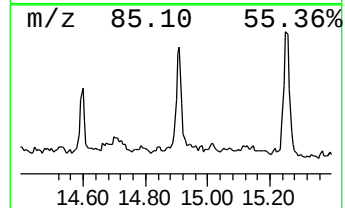
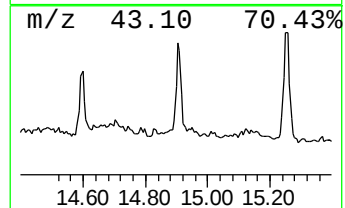
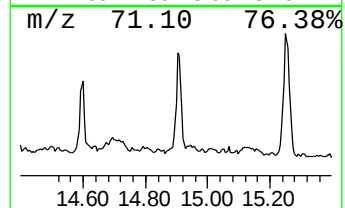
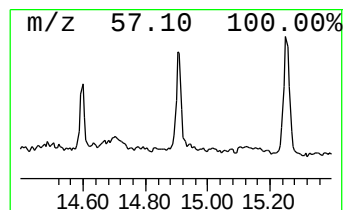
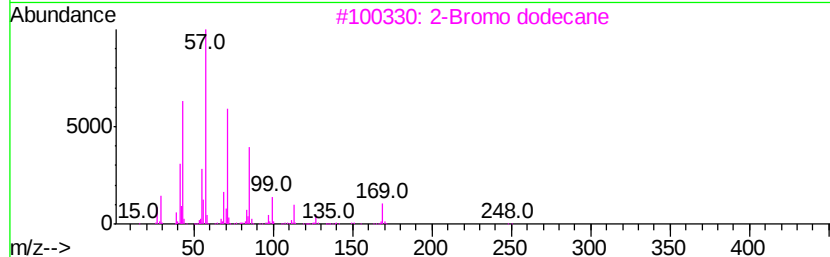
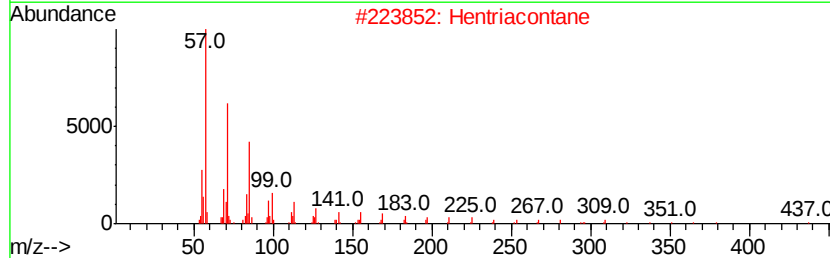
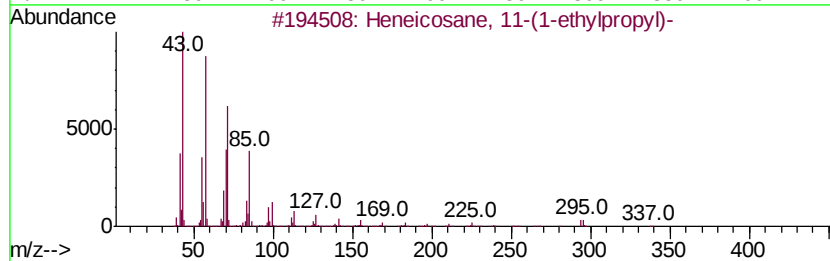
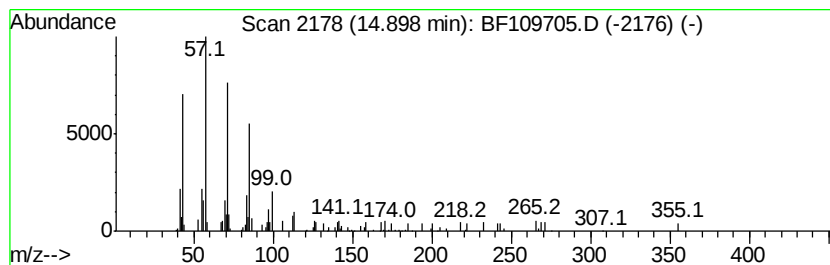
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Heneicosane, 11-(1-ethylpro... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.90	4.61 ng	98953	Perylene-d12	15.23

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	90
2			Hentriacontane	437	C31H64	000630-04-6	90
3			2-Bromo dodecane	248	C12H25Br	013187-99-0	87
4			Nonadecane	268	C19H40	000629-92-5	86
5			Octadecane, 1-iodo-	380	C18H37I	000629-93-6	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100918\
Data File : BF109705.D
Acq On : 9 Oct 2018 19:59
Operator : JU/SJ
Sample : J5324-01 10X
Misc :
ALS Vial : 9 Sample Multiplier: 1

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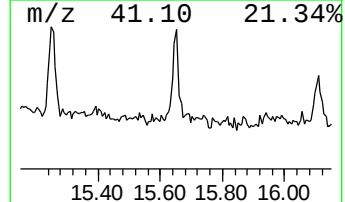
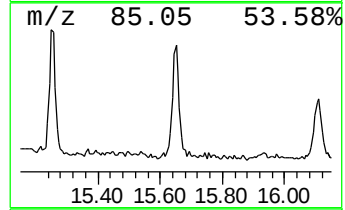
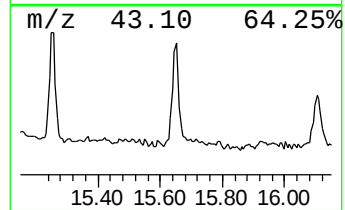
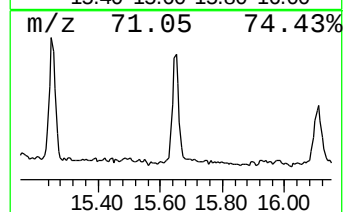
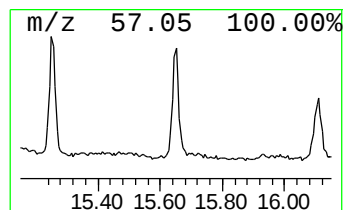
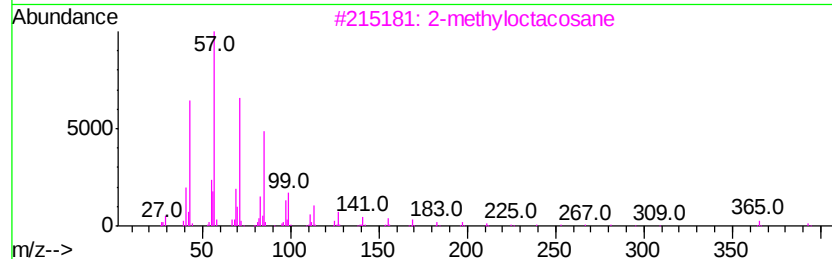
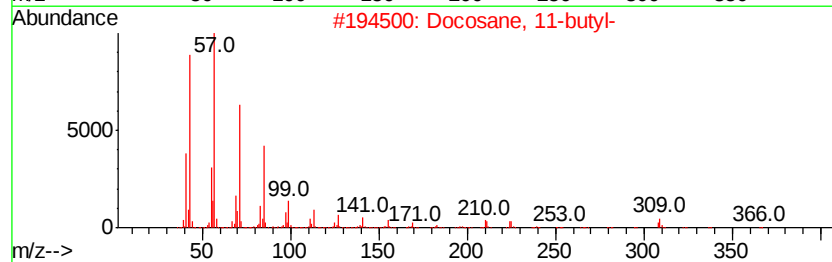
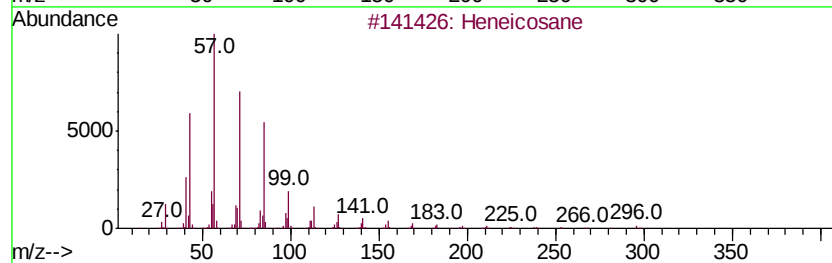
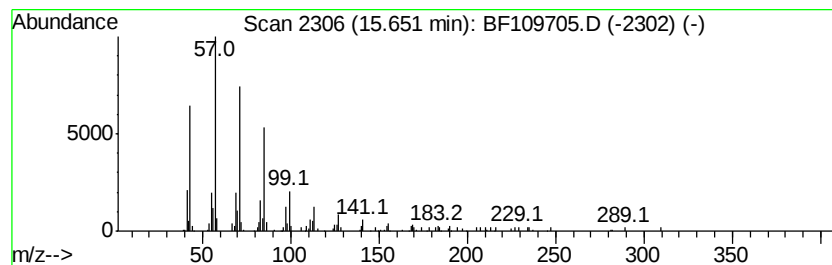
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Heneicosane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.65	7.58 ng	162745	Perylene-d12	15.23

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	91
2	Docosane, 11-butyl-		366	C26H54	013475-76-8	91
3	2-methyloctacosane		408	C29H60	1000376-72-8	91
4	Octacosane		394	C28H58	000630-02-4	91
5	Dodecane, 2-methyl-		184	C13H28	001560-97-0	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100918\
Data File : BF109705.D
Acq On : 9 Oct 2018 19:59
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ClientSampleId :
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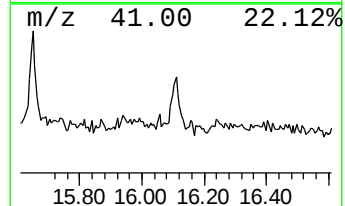
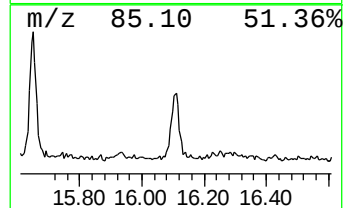
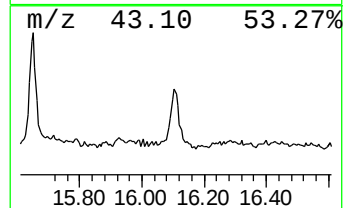
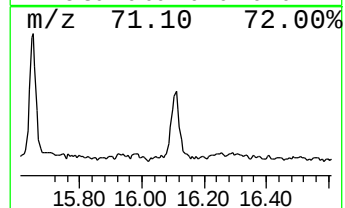
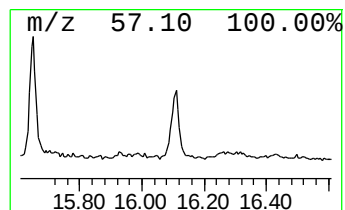
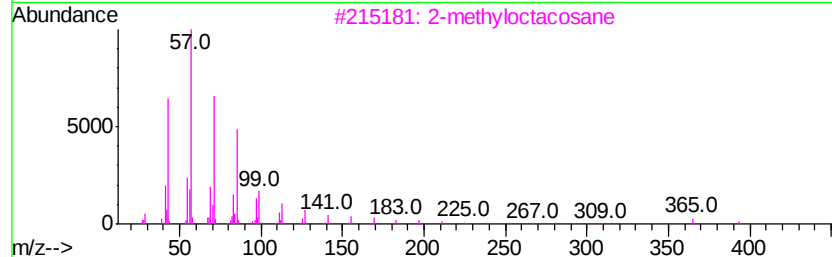
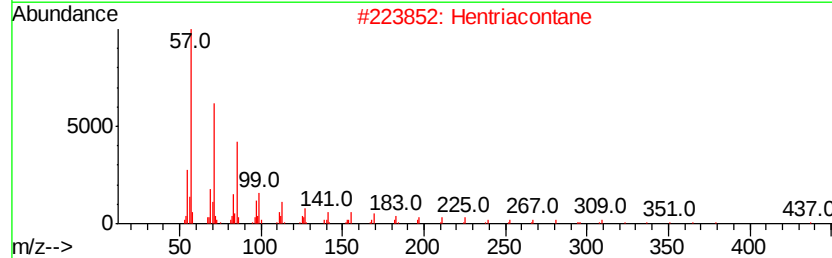
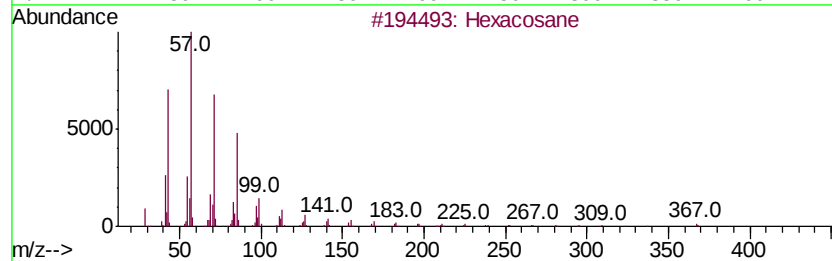
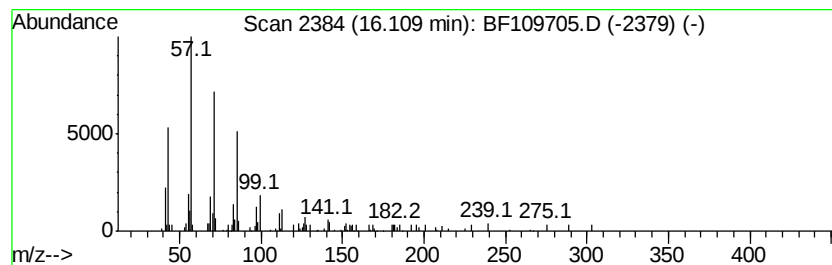
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Hexacosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.11	4.70 ng	101056	Perylene-d12	15.23

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosane	366	C26H54	000630-01-3	87
2			Hentriacontane	437	C31H64	000630-04-6	87
3			2-methyloctacosane	408	C29H60	1000376-72-8	87
4			Hexatriacontane	507	C36H74	000630-06-8	86
5			Octacosane	394	C28H58	000630-02-4	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF100918\
 Data File : BF109705.D
 Acq On : 9 Oct 2018 19:59
 Operator : JU/SJ
 Sample : J5324-01 10X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF100818.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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					#	RT	Resp	Conc
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1-Nonanol	7.79	2.0	ng	57917	2	7.97	564454	20.0
unknown9.21	9.21	4.7	ng	131668	3	9.72	555849	20.0
unknown10.59	10.59	6.7	ng	152926	4	11.20	457268	20.0
Hexadecane, 2,6,1...	10.65	5.0	ng	114173	4	11.20	457268	20.0
Undecane, 2-methyl-	11.11	5.9	ng	134349	4	11.20	457268	20.0
Heptadecane, 3-me...	12.26	3.6	ng	83103	4	11.20	457268	20.0
Heneicosane, 11-(...	14.90	4.6	ng	98953	6	15.23	429647	20.0
Heneicosane	15.65	7.6	ng	162745	6	15.23	429647	20.0
Hexacosane	16.11	4.7	ng	101056	6	15.23	429647	20.0