

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF062218\
 Data File : BF106746.D
 Acq On : 23 Jun 2018 8:39
 Operator : JU/SJ
 Sample : J3597-02
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 W-43

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	493	rBV	32502	36564	2.72%	0.375%
2	5.284	498	501	504	rVB	33586	30276	2.25%	0.311%
3	5.601	552	555	567	rBB	633411	566458	42.14%	5.813%
4	6.601	721	725	733	rBV	428970	375011	27.90%	3.848%
5	6.736	744	748	751	rBV	994833	916999	68.21%	9.409%
6	6.960	782	786	791	rBV	346001	296567	22.06%	3.043%
7	7.119	808	813	816	rBV	1015356	918761	68.34%	9.428%
8	7.525	877	882	885	rBV	752470	717812	53.40%	7.366%
9	7.636	899	901	906	rBV	27569	25213	1.88%	0.259%
10	8.242	1000	1004	1010	rBV	373264	311182	23.15%	3.193%
11	8.977	1125	1129	1132	rBV	31937	32097	2.39%	0.329%
12	9.083	1145	1147	1151	rVB3	21371	21867	1.63%	0.224%
13	9.183	1161	1164	1167	rVB4	12874	16759	1.25%	0.172%
14	9.219	1167	1170	1173	rBV4	17209	16560	1.23%	0.170%
15	9.319	1182	1187	1190	rBV	1239326	1183688	88.05%	12.146%
16	9.360	1190	1194	1196	rVB4	16197	22323	1.66%	0.229%
17	9.519	1218	1221	1224	rVB	29965	26026	1.94%	0.267%
18	9.660	1242	1245	1248	rBV2	35801	35783	2.66%	0.367%
19	9.707	1251	1253	1257	rVB	24547	21611	1.61%	0.222%
20	9.913	1282	1288	1292	rBV6	9162	14417	1.07%	0.148%
21	10.001	1299	1303	1307	rBV	358230	301016	22.39%	3.089%
22	10.654	1412	1414	1417	rVB	18426	14318	1.07%	0.147%
23	10.795	1434	1438	1446	rBV	760678	685405	50.99%	7.033%
24	10.883	1450	1453	1456	rBV	14566	13632	1.01%	0.140%
25	10.913	1456	1458	1463	rVB2	17538	15991	1.19%	0.164%
26	10.966	1463	1467	1469	rBV	33079	27423	2.04%	0.281%
27	10.995	1469	1472	1475	rBV2	29742	30743	2.29%	0.315%
28	11.030	1475	1478	1481	rBV2	39995	31629	2.35%	0.325%
29	11.083	1484	1487	1488	rBV2	18255	17233	1.28%	0.177%
30	11.142	1494	1497	1501	rVB2	30939	32907	2.45%	0.338%
31	11.183	1501	1504	1506	rVV	33398	31649	2.35%	0.325%
32	11.224	1509	1511	1514	rVB	20703	16154	1.20%	0.166%
33	11.377	1532	1537	1541	rVB4	15855	18478	1.37%	0.190%
34	11.495	1553	1557	1560	rBV	357975	306252	22.78%	3.142%

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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	12.960	1803	1806	1813	rBV	106959	95974	7.14%	0.985%
36	13.077	1822	1826	1830	rBV	1319334	1344300	100.00%	13.794%
37	13.624	1916	1919	1921	rVB	28881	23724	1.76%	0.243%
38	13.954	1972	1975	1982	rVB	55681	58569	4.36%	0.601%
39	14.142	2004	2007	2012	rBV	372281	372506	27.71%	3.822%
40	14.277	2028	2030	2035	rVB	29524	28952	2.15%	0.297%
41	14.595	2080	2084	2088	rBV2	53474	57165	4.25%	0.587%
42	14.912	2135	2138	2143	rVB	77963	75728	5.63%	0.777%
43	15.265	2195	2198	2203	rVB	84497	94533	7.03%	0.970%
44	15.683	2262	2269	2279	rVB2	218408	355712	26.46%	3.650%
45	16.130	2341	2345	2352	rVB	50882	73530	5.47%	0.755%
46	16.665	2433	2436	2442	rVB3	24141	36001	2.68%	0.369%

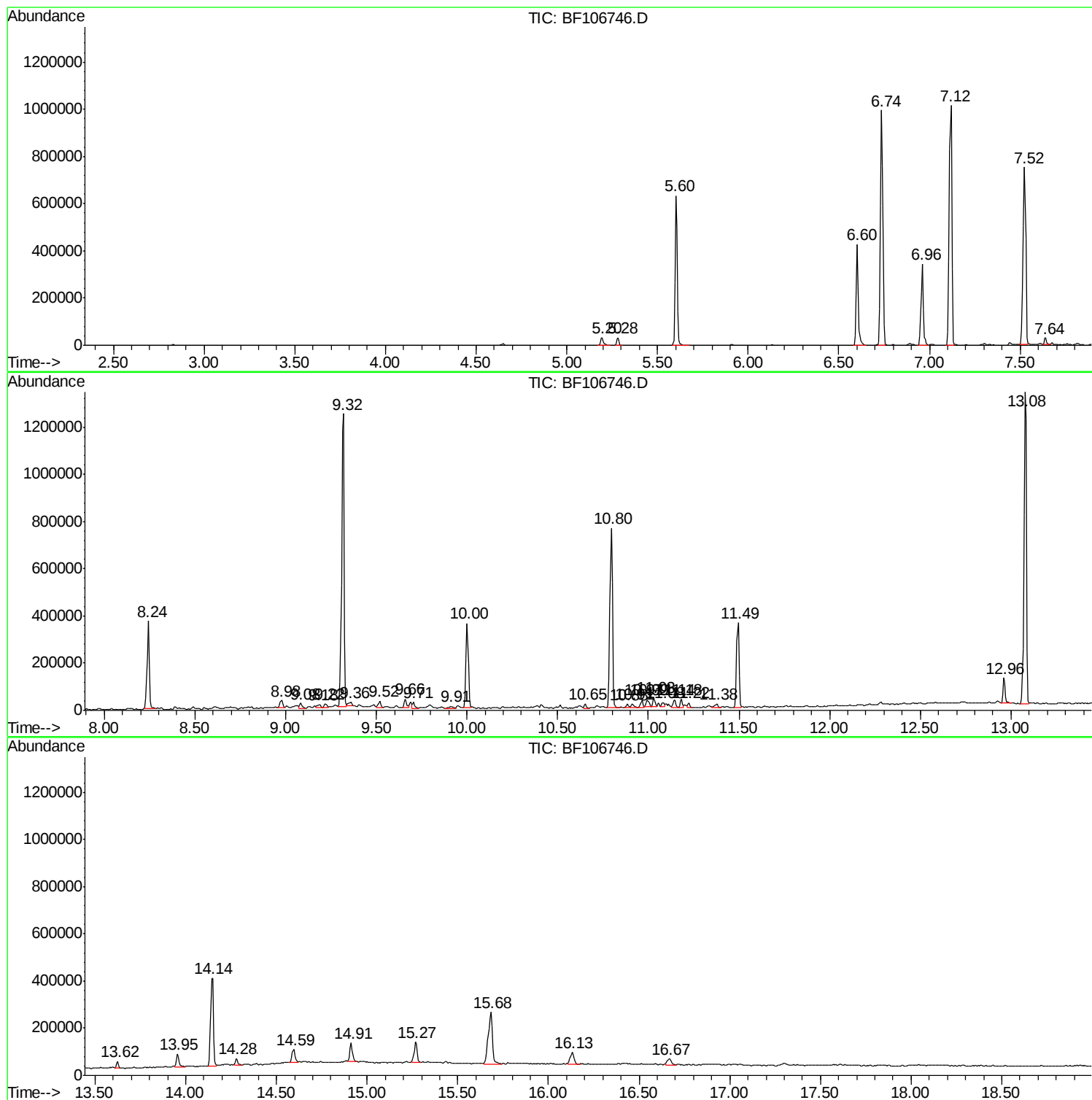
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ClientSampled :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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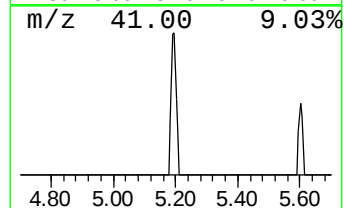
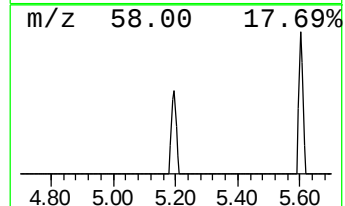
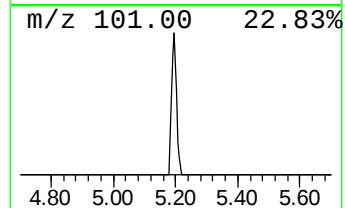
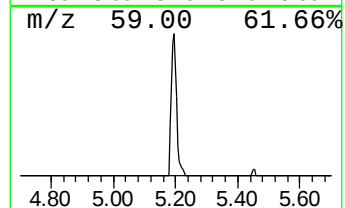
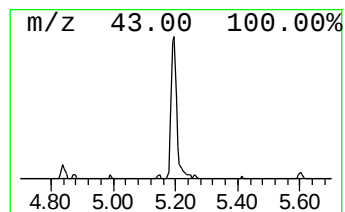
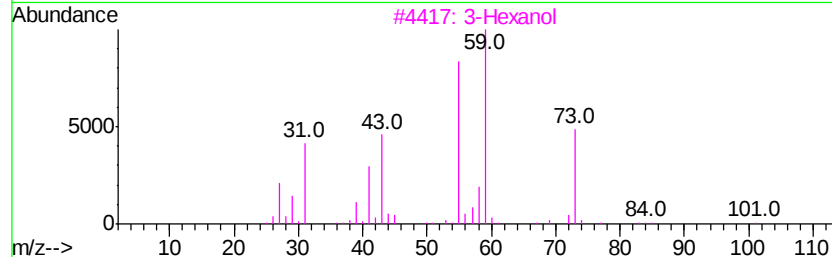
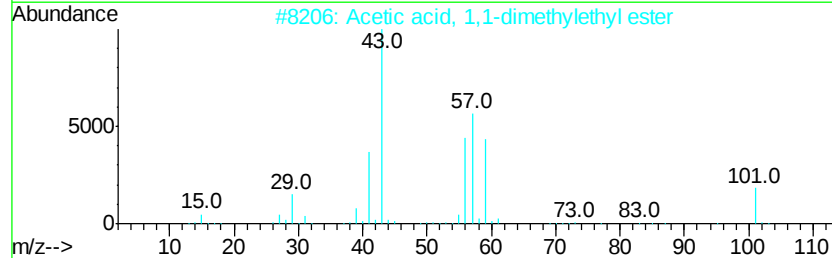
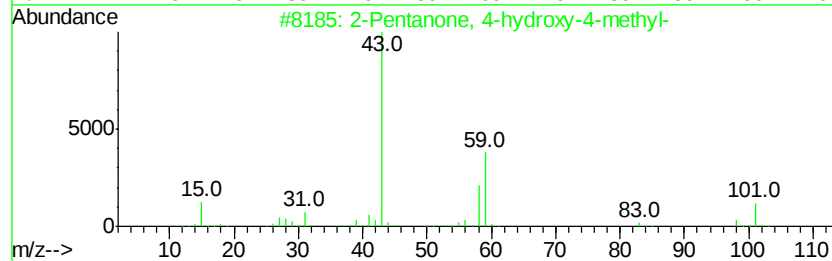
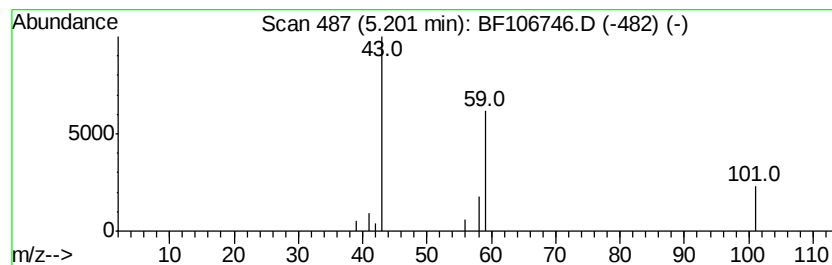
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.20	2.47 ng	36564	1,4-Dichlorobenzene-d4	6.96

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	40
3			3-Hexanol	102	C6H14O	000623-37-0	33
4			5-Hexen-2-one	98	C6H10O	000109-49-9	9
5			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	9



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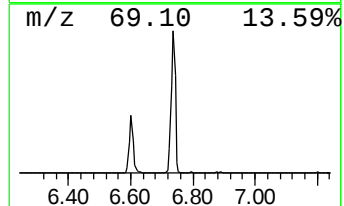
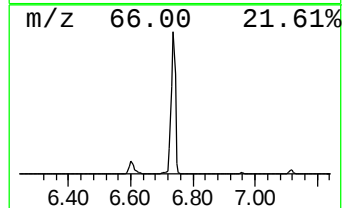
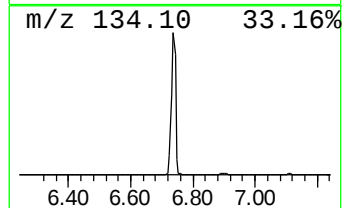
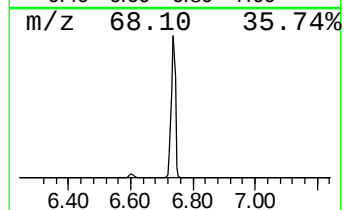
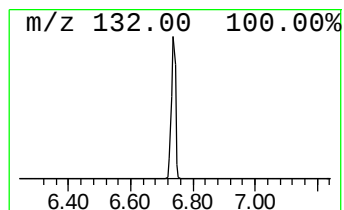
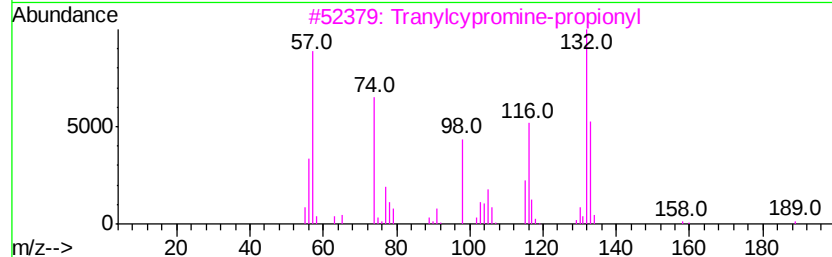
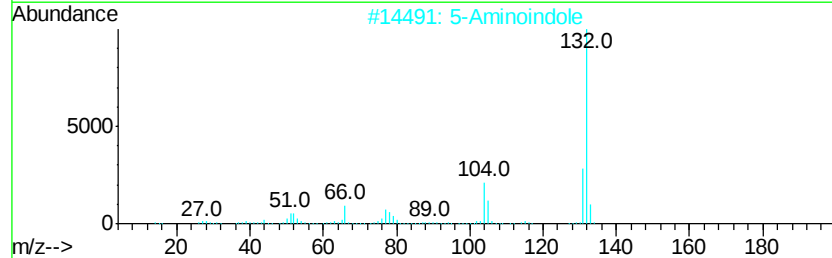
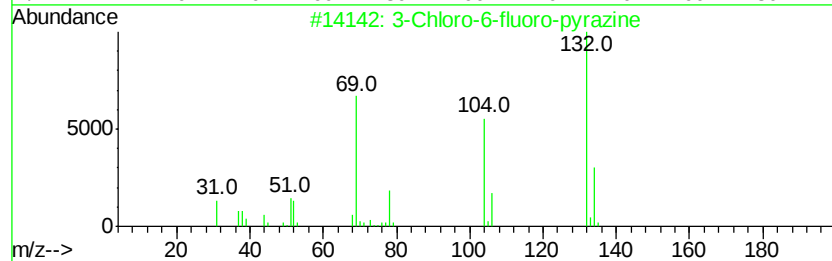
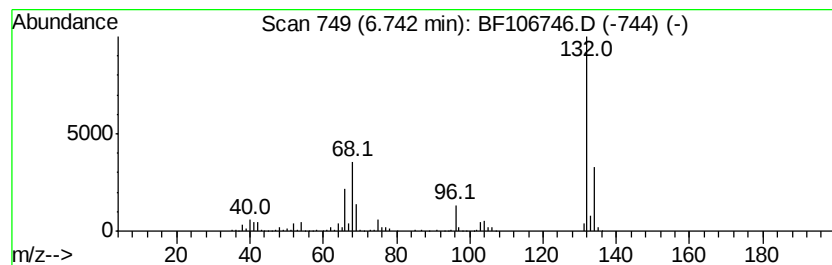
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.74 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.74	61.84 ng	916999	1,4-Dichlorobenzene-d4	6.96

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine			132	C4H2ClFN2	1000146-10-7	16
2	5-Aminoindole			132	C8H8N2	005192-03-0	12
3	Tranvlcvpromine-propionyl			189	C12H15NO	1000123-86-3	10
4	4-Chloro-2-fluoro-pyrimidine			132	C4H2ClFN2	051422-00-5	9
5	1H-Benzimidazole, 2-methyl-			132	C8H8N2	000615-15-6	9



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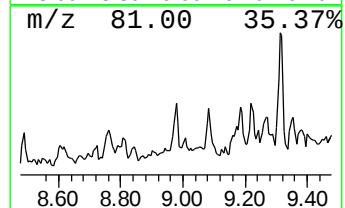
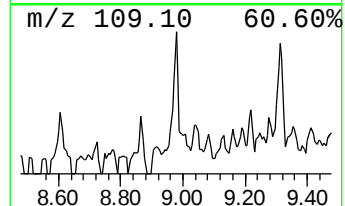
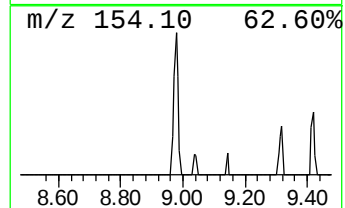
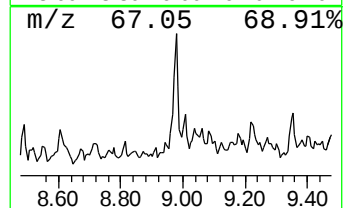
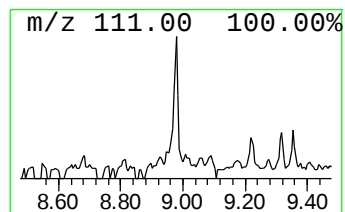
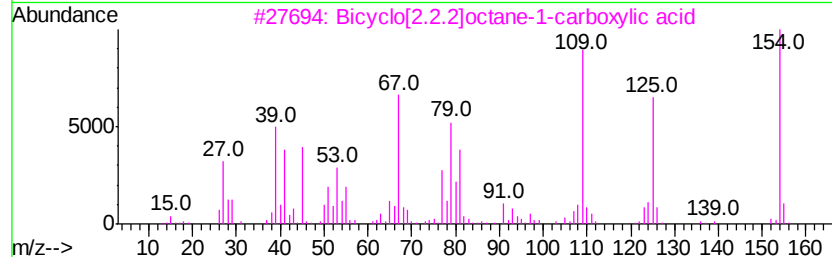
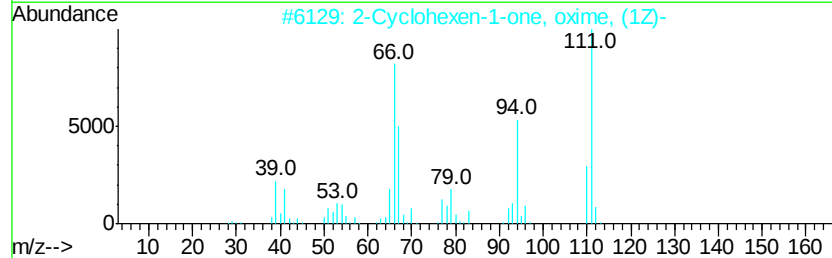
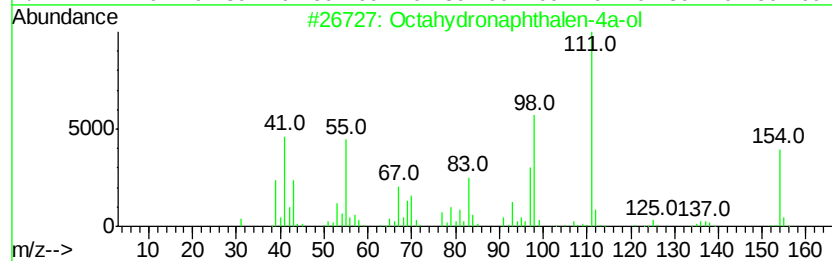
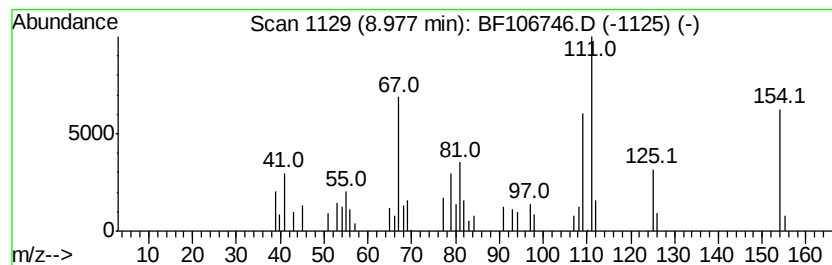
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 unknown8.98 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.98	2.06 ng	32097	Naphthalene-d8	8.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octahydronaphthalen-4a-ol	154	C10H18O	055693-34-0	38
2		2-Cyclohexen-1-one, oxime, (1Z)-	111	C6H9NO	1000362-09-7	35
3		Bicyclo[2.2.2]octane-1-carboxylic acid...	154	C9H14O2	000699-55-8	35
4		2-Cyclopentene-1-carboxylic acid...	154	C9H14O2	006894-69-5	35
5		4-Butyl-5-methylpyrazalone	154	C8H14N2O	1000289-28-0	25



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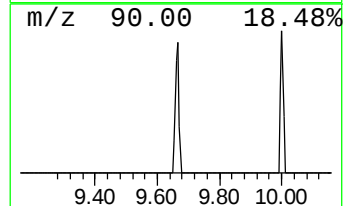
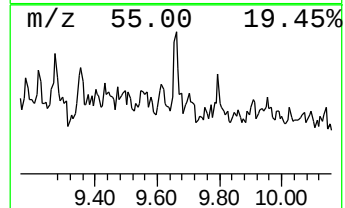
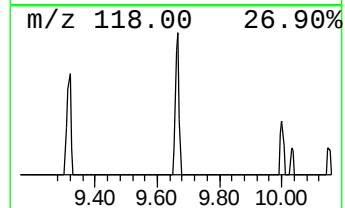
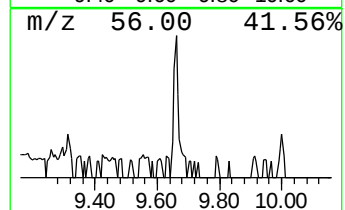
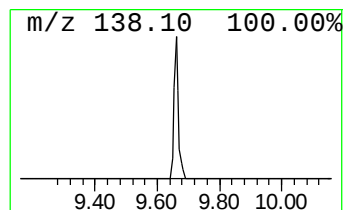
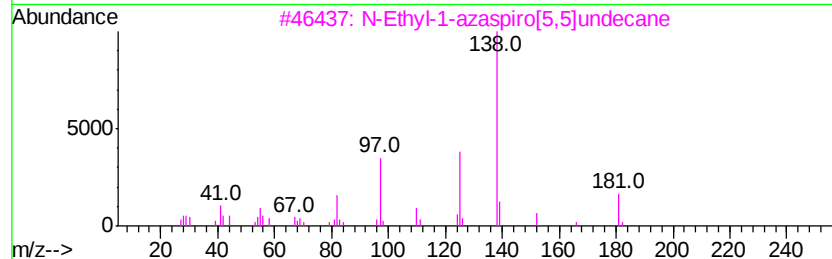
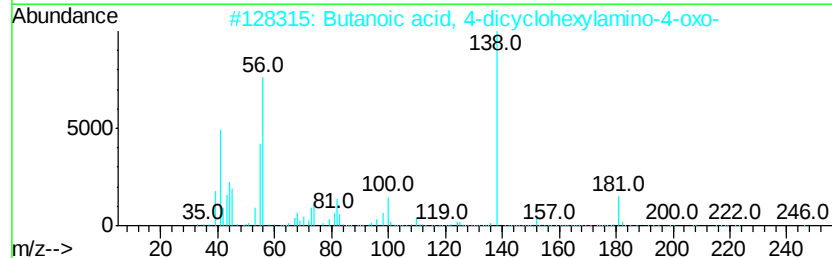
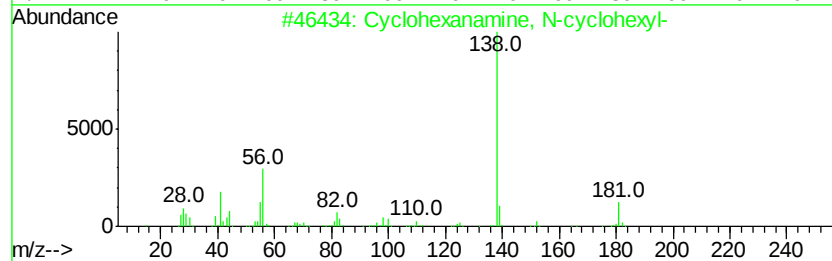
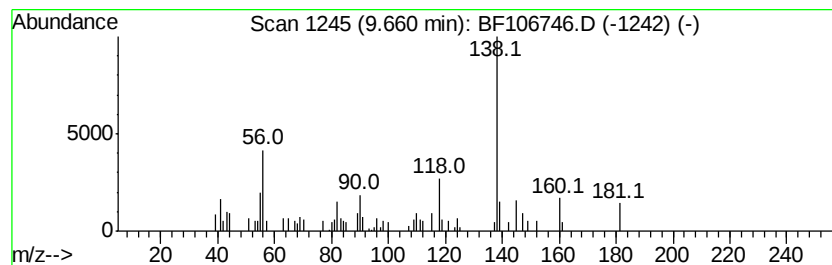
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Cyclohexanamine, N-cyclohexyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.66	2.38 ng	35783	Acenaphthene-d10	10.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	64
2		Butanoic acid, 4-dicyclohexylami...	281	C16H27NO3	328013-95-2	50
3		N-Ethyl-1-azaspiro[5.5]undecane	181	C12H23N	1000101-78-6	38
4		4H-Pyran-4-one, 2-ethyl-6-methyl-	138	C8H10O2	057276-03-6	35
5		2-Imidazolidinone, 1,3-diethenyl-	138	C7H10N2O	013811-50-2	32



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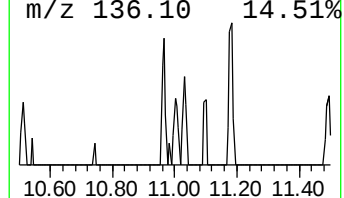
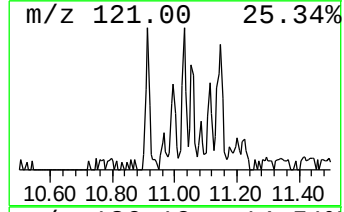
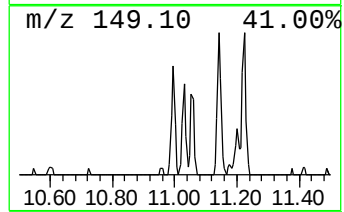
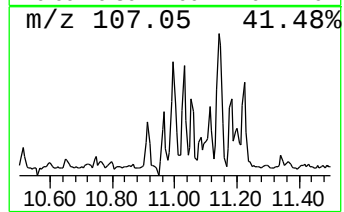
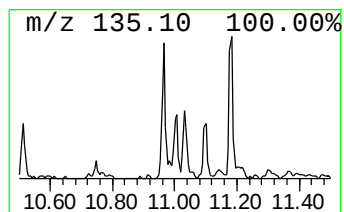
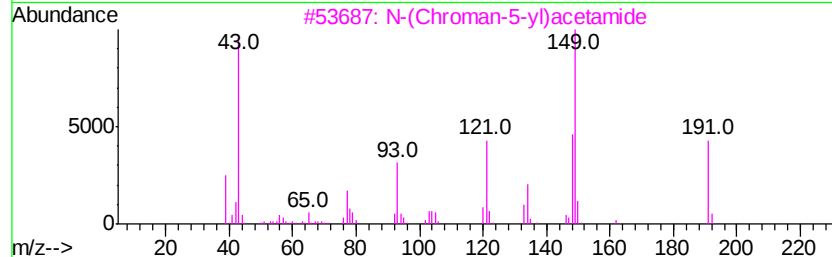
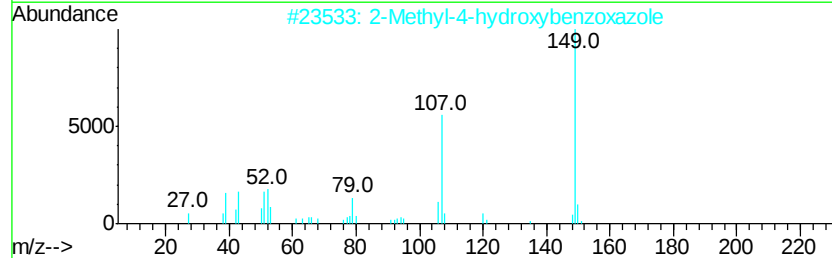
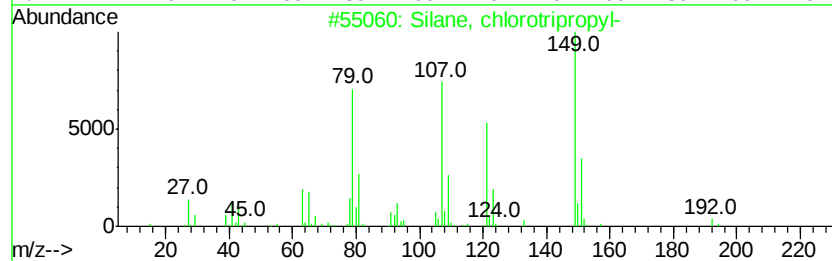
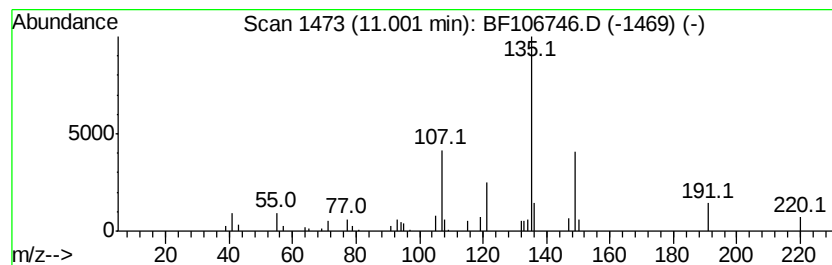
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Silane, chlorotripropyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.00	2.01 ng	30743	Phenanthrene-d10	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Silane, chlorotripropyl-	192	C9H21ClSi	000995-25-5	58
2		2-Methyl-4-hydroxybenzoxazole	149	C8H7NO2	051110-60-2	53
3		N-(Chroman-5-yl)acetamide	191	C11H13NO2	1000306-69-8	38
4		1,3-Cyclohexadiene-1-carboxaldeh...	150	C10H14O	000116-26-7	35
5		Pyridine-3-carbonitrile, 1,2-dih...	191	C9H9N3O2	220098-92-0	35



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\
Data File : BF106746.D
Acq On : 23 Jun 2018 8:39
Operator : JU/SJ
Sample : J3597-02
Misc :
ALS Vial : 36 Sample Multiplier: 1

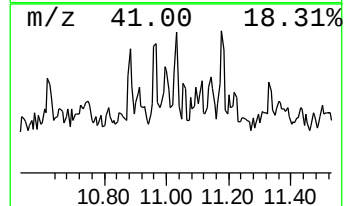
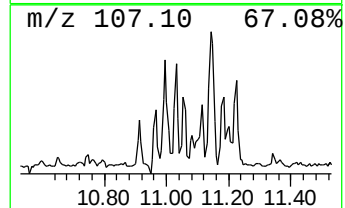
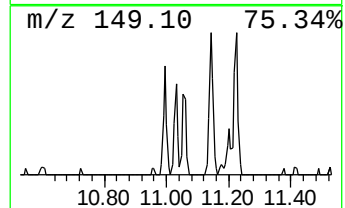
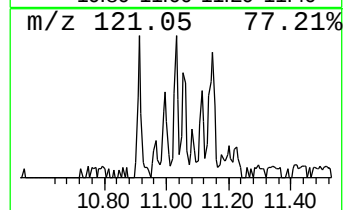
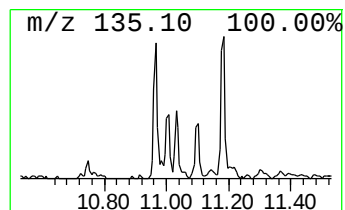
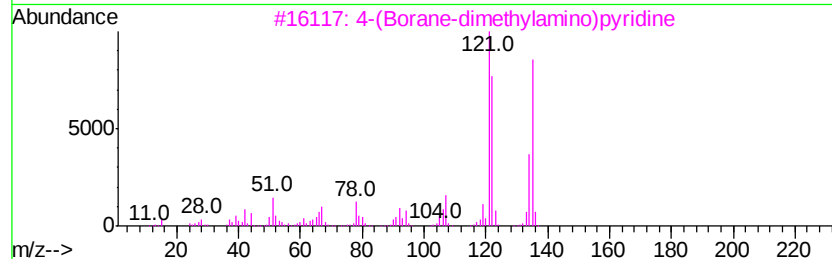
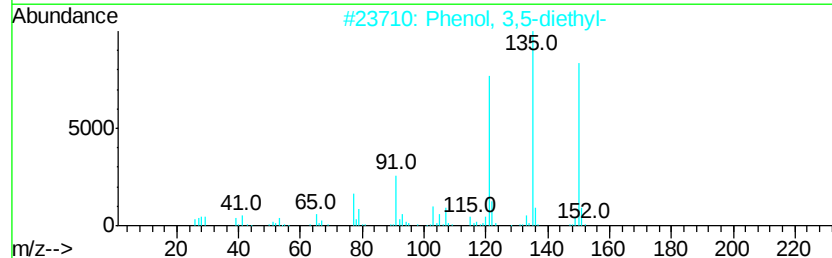
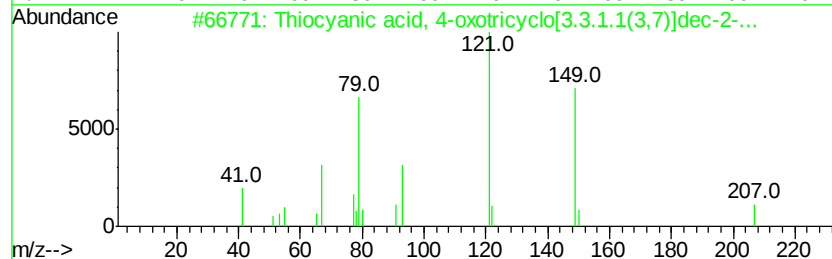
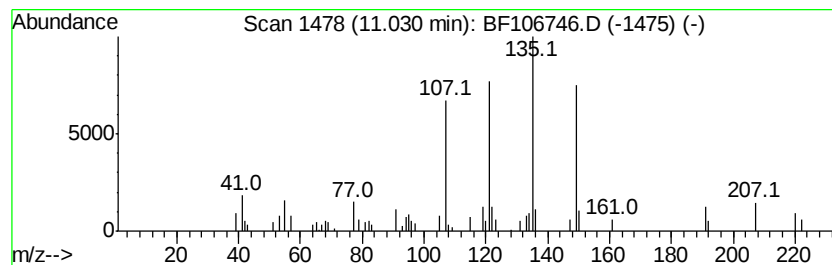
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Thiocyanic acid, 4-oxotricy... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD		R.T.		
11.03	2.07 ng	31629	Phenanthrene-d10		11.49		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Thiocyanic acid, 4-oxotricvclo[3...	207	C11H13NOS	056781-89-6	55
2			Phenol, 3,5-diethyl-	150	C10H14O	001197-34-8	30
3			4-(Borane-dimethylamino)pvridine	136	C7H13BN2	001769-74-0	27
4			Thiocyanic acid, 4-oxotricvclo[3...	207	C11H13NOS	056781-88-5	25
5			Methyl 2,3,4-tri-O-acetyl-1-(3-(...	483	C20H25N3O9S	1000242-58-9	22



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\
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ClientSampleId :
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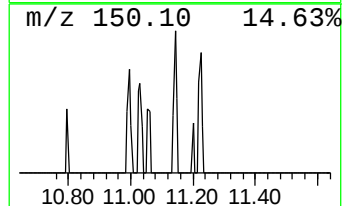
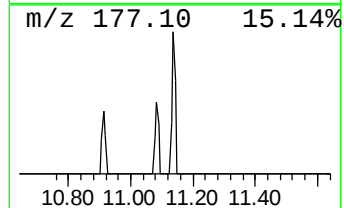
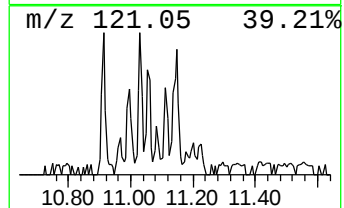
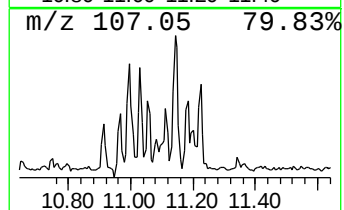
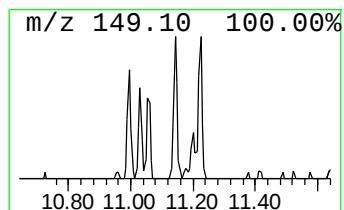
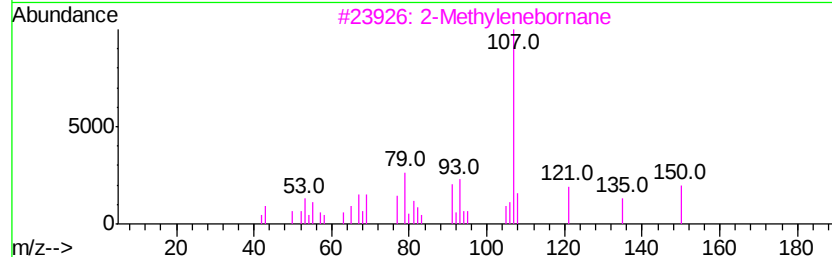
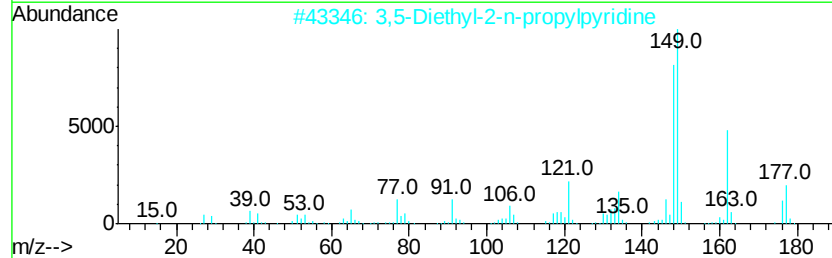
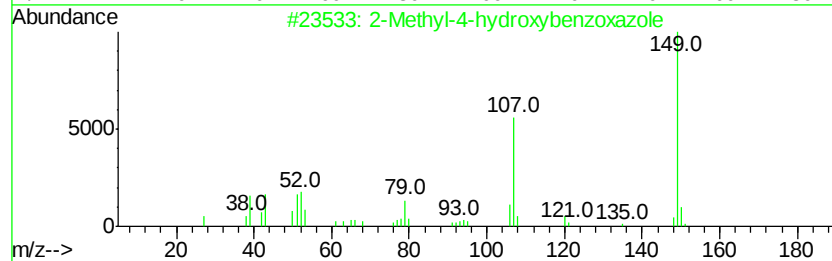
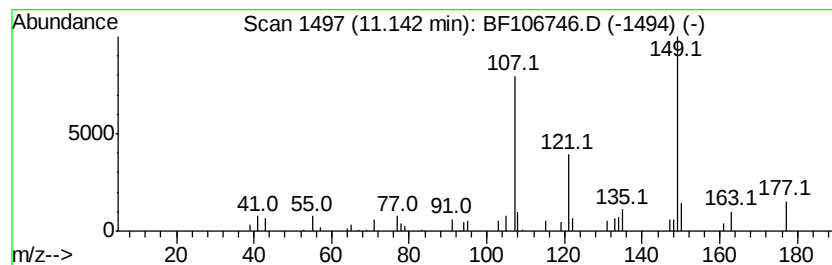
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 2-Methyl-4-hydroxybenzoxazole Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.14	2.15 ng	32907	Phenanthrene-d10	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Methyl-4-hydroxybenzoxazole	149	C8H7NO2	051110-60-2	64
2		3,5-Diethyl-2-n-propylpyridine	177	C12H19N	004808-75-7	43
3		2-Methylenebornane	150	C11H18	027538-47-2	38
4		Phenol, 2-butyl-	150	C10H14O	003180-09-4	38
5		Phthalic acid, ethyl 2-(2-nitrop...	343	C18H17NO6	1000377-99-1	35



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\
Data File : BF106746.D
Acq On : 23 Jun 2018 8:39
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Sample : J3597-02
Misc :
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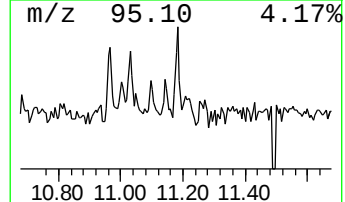
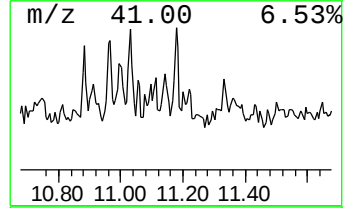
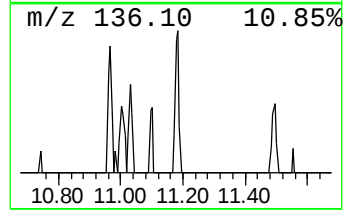
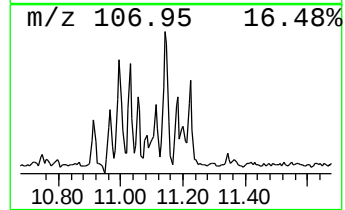
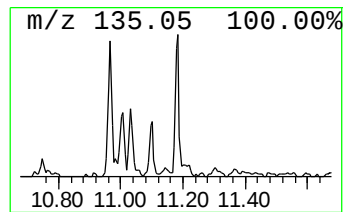
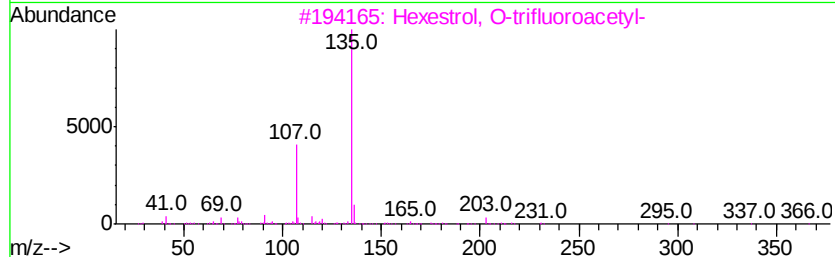
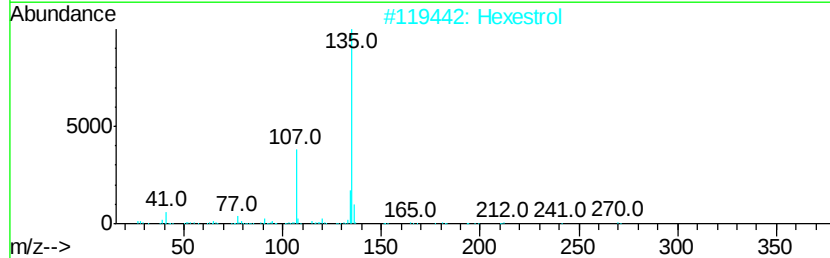
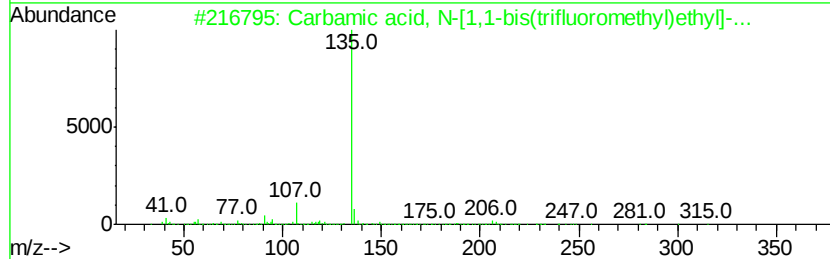
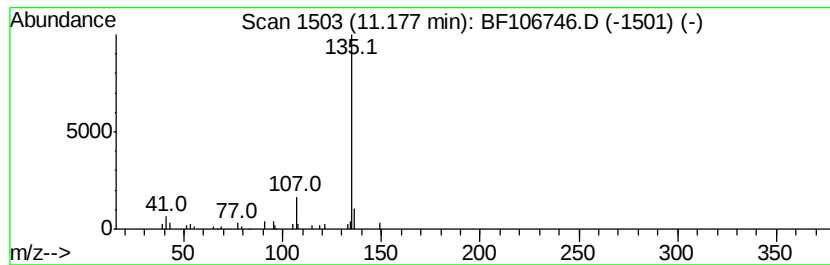
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Carbamic acid, N-[1,1-bis(t... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.18	2.07 ng	31649	Phenanthrene-d10	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Carbamic acid, N-[1,1-bis(triflu...	413	C19H25F6N02	296242-69-8	78
2		Hexestrol	270	C18H22O2	005635-50-7	45
3		Hexestrol, O-trifluoroacetyl-	366	C20H21F3O3	1000365-31-7	45
4		Hexestrol	270	C18H22O2	000084-16-2	45
5		1-Adamantaneacetic acid	194	C12H18O2	004942-47-6	40



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\
Data File : BF106746.D
Acq On : 23 Jun 2018 8:39
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Sample : J3597-02
Misc :
ALS Vial : 36 Sample Multiplier: 1

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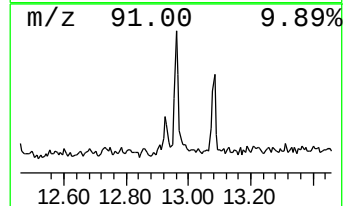
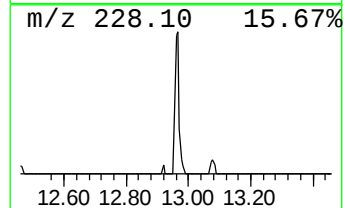
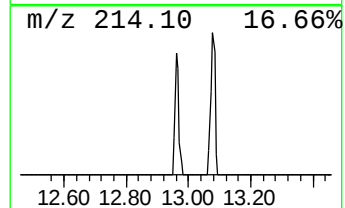
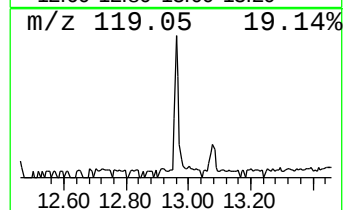
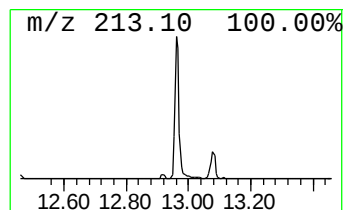
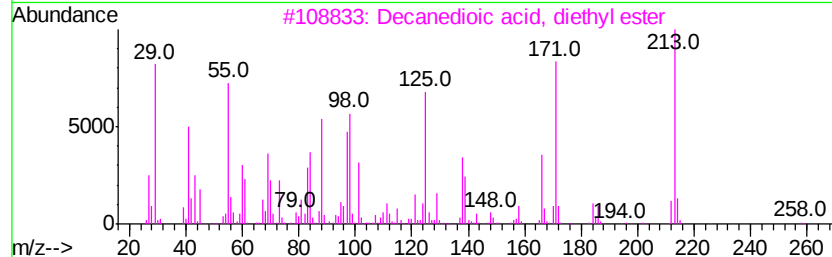
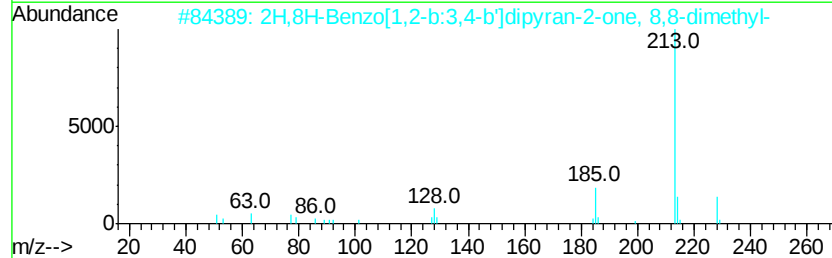
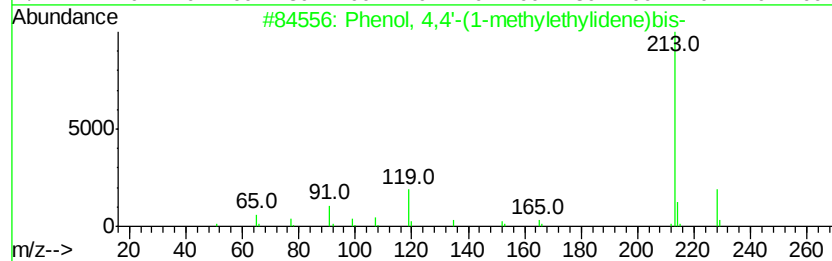
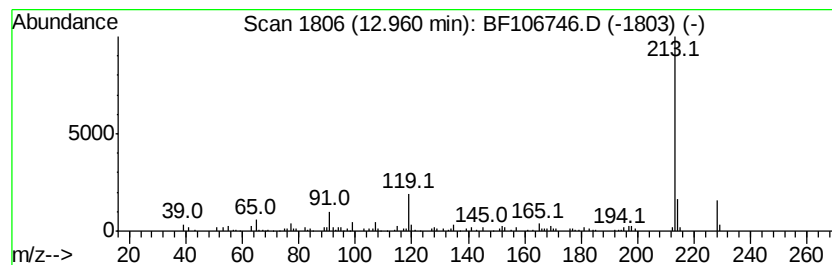
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Phenol, 4,4'-(1-methylethyl... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.96	5.15 ng	95974	Chrysene-d12	14.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenol, 4,4'-(1-methylethylidene)bis-	228	C15H16O2	000080-05-7	96
2		2H,8H-Benzo[1,2-b:3,4-b']dipyran-2-one, 8,8-dimethyl-	228	C14H12O3	000523-59-1	58
3		Decanedioic acid, diethyl ester	258	C14H26O4	000110-40-7	50
4		4-Adamantan-1-yl-2-(2-methylpropyl)-1H-imidazole	333	C19H27N02S	1000311-66-6	50
5		4-Bromo-2-(methylamino)tropone	213	C8H8BrNO	1000241-45-0	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\
Data File : BF106746.D
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ClientSampleId :
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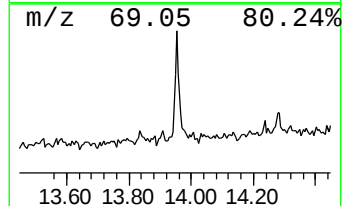
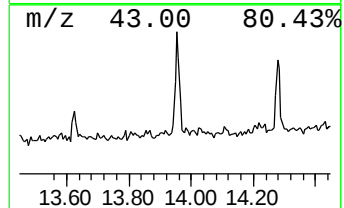
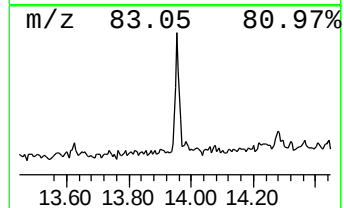
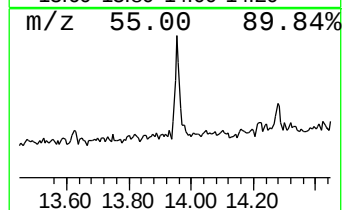
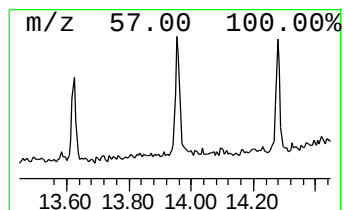
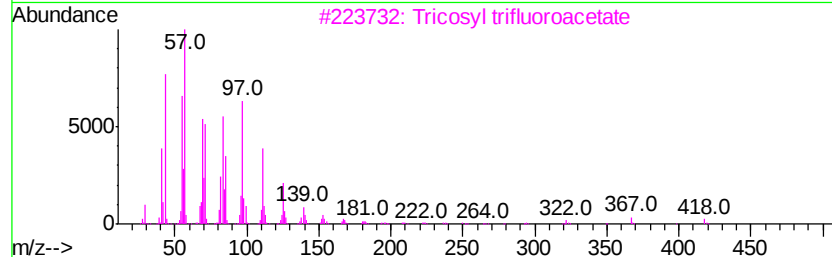
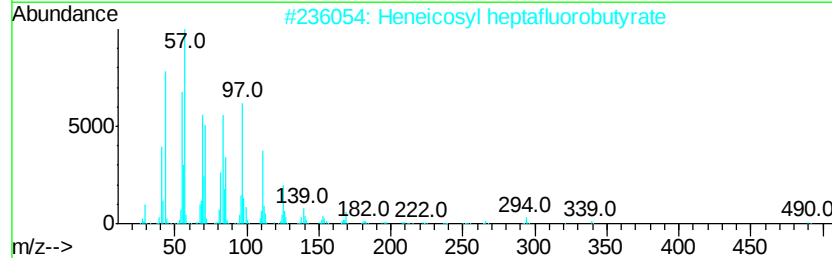
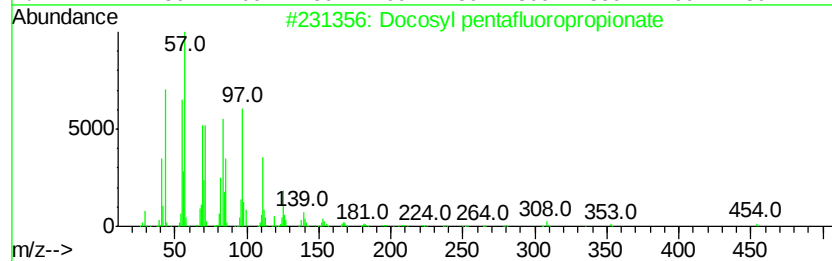
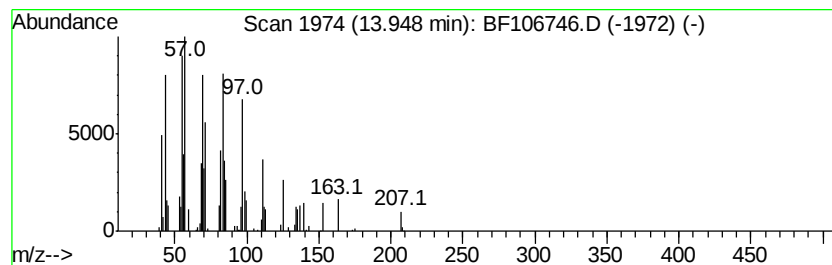
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Docosyl pentafluoropropionate Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.95	3.14 ng	58569	Chrysene-d12	14.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Docosyl pentafluoropropionate	472	C25H45F5O2	1000351-80-9	91
2		Heneicosyl heptafluorobutyrate	508	C25H43F7O2	1000351-83-8	91
3		Tricosyl trifluoroacetate	436	C25H47F3O2	1000351-75-1	91
4		Tricosyl heptafluorobutyrate	536	C27H47F7O2	1000351-83-4	90
5		Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\
Data File : BF106746.D
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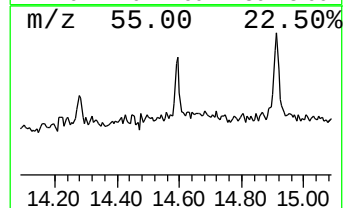
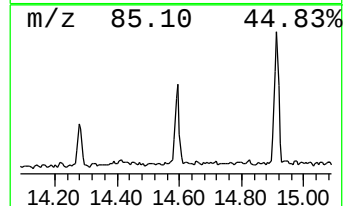
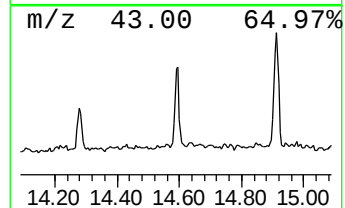
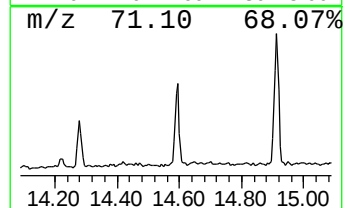
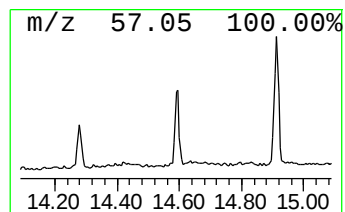
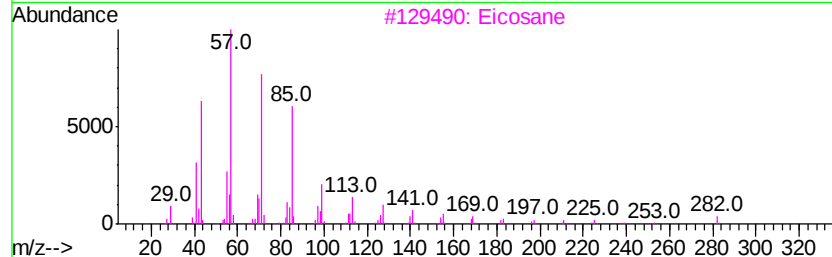
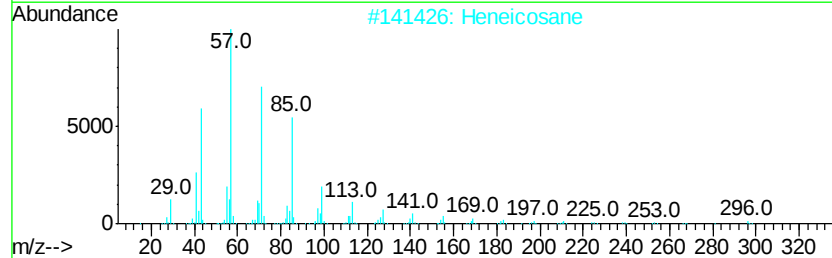
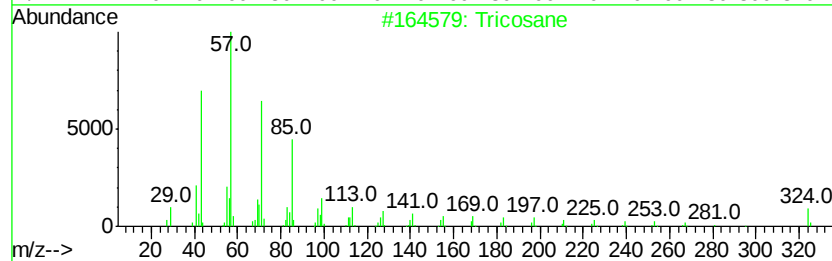
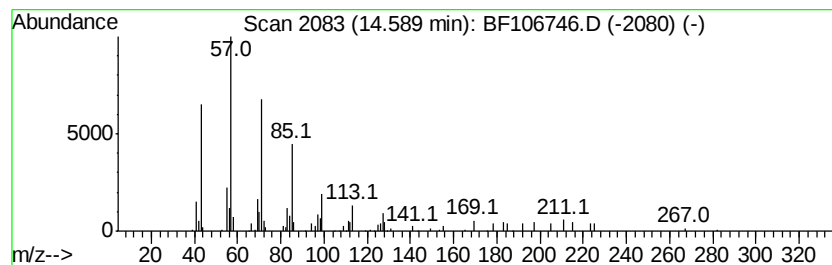
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Tricosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.59	3.07 ng	57165	Chrysene-d12	14.15

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tricosane	324	C23H48	000638-67-5	96
2			Heneicosane	296	C21H44	000629-94-7	90
3			Eicosane	282	C20H42	000112-95-8	90
4			Pentacosane	352	C25H52	000629-99-2	87
5			Hexacosane	366	C26H54	000630-01-3	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\
Data File : BF106746.D
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Sample : J3597-02
Misc :
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ClientSampleId :
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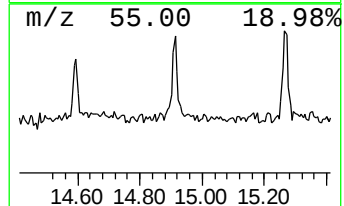
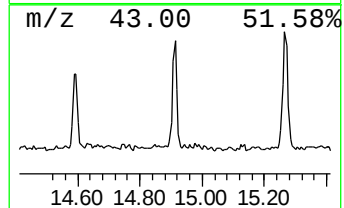
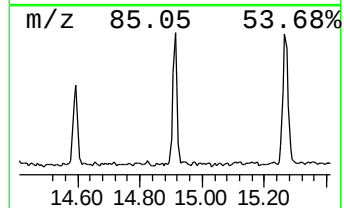
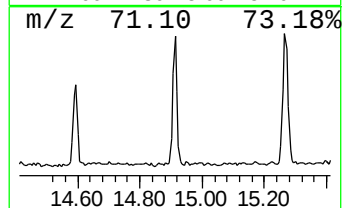
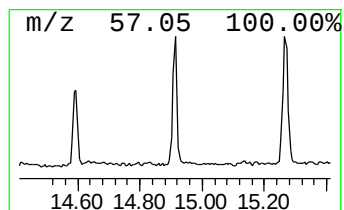
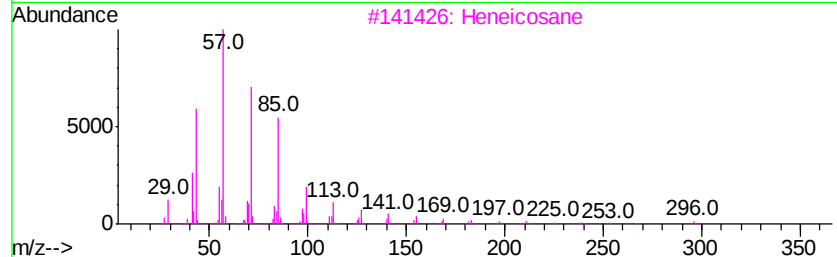
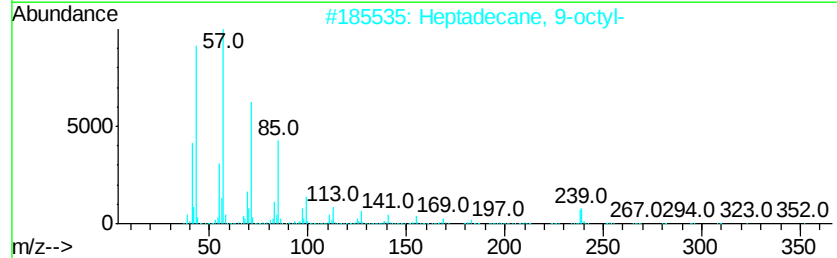
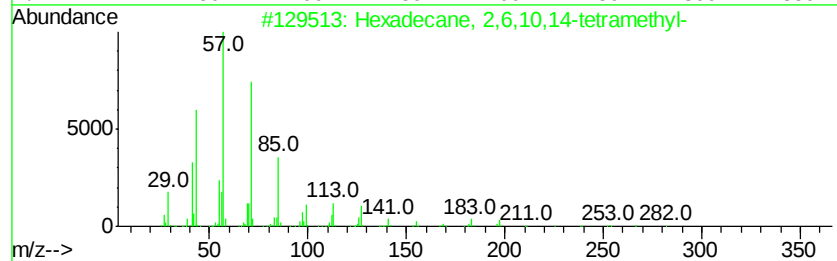
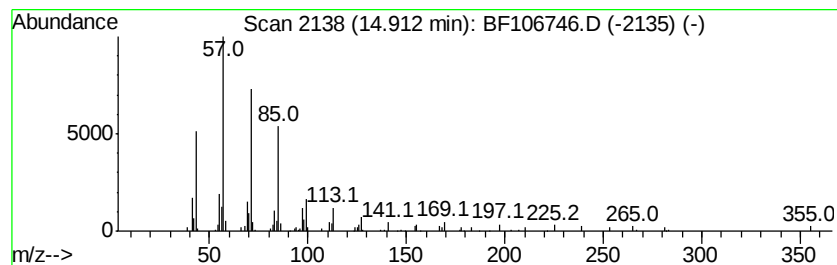
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.91	4.07 ng	75728	Chrysene-d12	14.15

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	93
2			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
3			Heneicosane	296	C21H44	000629-94-7	91
4			Hexacosane	366	C26H54	000630-01-3	91
5			Octacosane	394	C28H58	000630-02-4	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\
Data File : BF106746.D
Acq On : 23 Jun 2018 8:39
Operator : JU/SJ
Sample : J3597-02
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
W-43

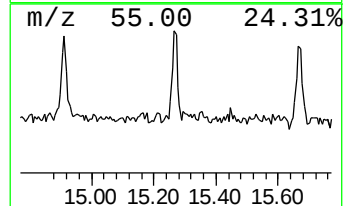
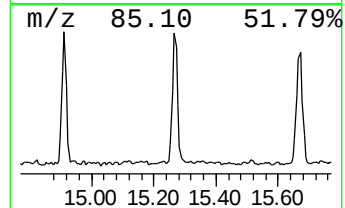
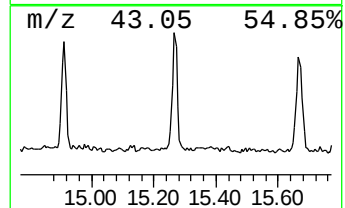
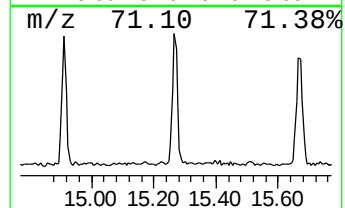
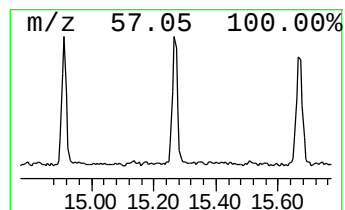
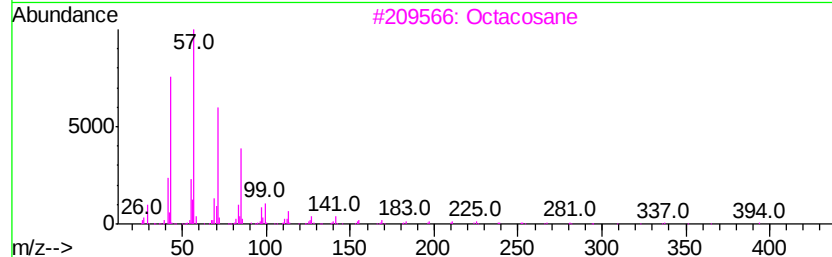
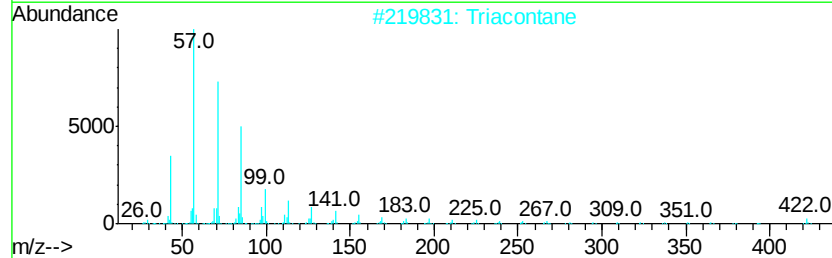
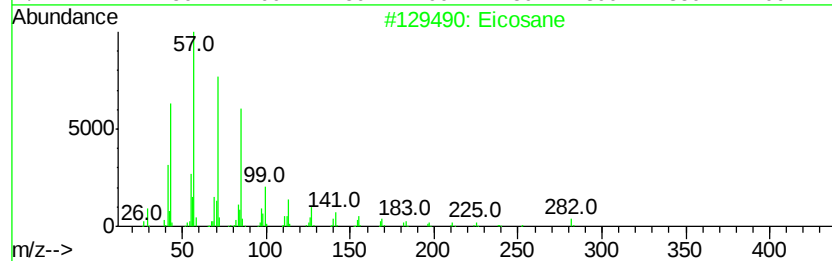
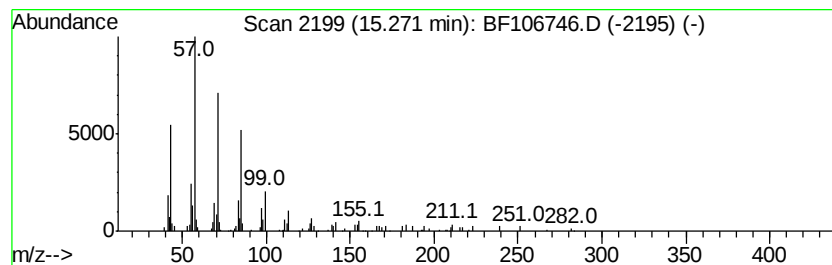
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Eicosane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.27	5.32 ng	94533	Perylene-d12	15.68

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane			282	C20H42	000112-95-8	96
2	Triacontane			422	C30H62	000638-68-6	91
3	Octacosane			394	C28H58	000630-02-4	91
4	Hexacosane			366	C26H54	000630-01-3	90
5	Heneicosane			296	C21H44	000629-94-7	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF062218\
Data File : BF106746.D
Acq On : 23 Jun 2018 8:39
Operator : JU/SJ
Sample : J3597-02
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
W-43

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	2.5	ng	36564	1	6.96	296567	20.0
unknown6.74	6.74	61.8	ng	916999	1	6.96	296567	20.0
unknown8.98	8.98	2.1	ng	32097	2	8.24	311182	20.0
Cyclohexanamine, ...	9.66	2.4	ng	35783	3	10.00	301016	20.0
Silane, chlorotri...	11.00	2.0	ng	30743	4	11.49	306252	20.0
Thiocyanic acid, ...	11.03	2.1	ng	31629	4	11.49	306252	20.0
2-Methyl-4-hydrox...	11.14	2.1	ng	32907	4	11.49	306252	20.0
Carbamic acid, N-...	11.18	2.1	ng	31649	4	11.49	306252	20.0
Phenol, 4,4'-(1-m...	12.96	5.2	ng	95974	5	14.15	372506	20.0
Docosyl pentaflu...	13.95	3.1	ng	58569	5	14.15	372506	20.0
Tricosane	14.59	3.1	ng	57165	5	14.15	372506	20.0
Hexadecane, 2,6,1...	14.91	4.1	ng	75728	5	14.15	372506	20.0
Eicosane	15.27	5.3	ng	94533	6	15.68	355712	20.0