

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120319\
 Data File : BF118066.D
 Acq On : 3 Dec 2019 17:05
 Operator : JU
 Sample : K6069-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-3

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-BF111319.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	2.169	6	14	25	rVB	795487	1074487	24.76%	2.862%
2	5.092	506	511	523	rVB	517850	664368	15.31%	1.770%
3	5.498	575	580	583	rBV	3353370	3209629	73.95%	8.550%
4	6.504	746	751	754	rBV	3230554	3456144	79.63%	9.207%
5	6.645	770	775	779	rBV	3808775	3641247	83.89%	9.700%
6	6.875	811	814	818	rBV	954083	774429	17.84%	2.063%
7	7.034	837	841	844	rBV	3435997	2913617	67.13%	7.761%
8	7.439	905	910	913	rBV	2298327	2166759	49.92%	5.772%
9	8.163	1029	1033	1040	rBV	1197770	1000789	23.06%	2.666%
10	9.245	1211	1217	1220	rBV	4770739	4340409	100.00%	11.562%
11	9.633	1280	1283	1287	rBV	432240	328813	7.58%	0.876%
12	9.839	1312	1318	1325	rVB3	24710	54465	1.25%	0.145%
13	9.927	1328	1333	1336	rBV	1258905	1154749	26.60%	3.076%
14	10.716	1463	1467	1471	rBV	2883002	2681812	61.79%	7.144%
15	11.416	1582	1586	1590	rBV	1285779	1182914	27.25%	3.151%
16	11.480	1595	1597	1600	rVB2	61118	51312	1.18%	0.137%
17	11.633	1619	1623	1630	rBV2	34798	63785	1.47%	0.170%
18	11.939	1672	1675	1680	rBV	296128	272680	6.28%	0.726%
19	12.727	1806	1809	1817	rBV2	133955	174156	4.01%	0.464%
20	12.886	1833	1836	1844	rVB2	106071	142166	3.28%	0.379%
21	13.010	1852	1857	1860	rBV	4525181	4332207	99.81%	11.540%
22	13.521	1941	1944	1947	rBV	69477	68187	1.57%	0.182%
23	13.580	1951	1954	1957	rVB	69305	61710	1.42%	0.164%
24	13.727	1976	1979	1984	rVB	116280	114489	2.64%	0.305%
25	13.904	2005	2009	2017	rVB2	288196	390442	9.00%	1.040%
26	14.068	2033	2037	2041	rVB	1018875	856658	19.74%	2.282%
27	14.227	2060	2064	2069	rBV	135856	131387	3.03%	0.350%
28	14.533	2113	2116	2121	rVB	182168	164871	3.80%	0.439%
29	14.845	2163	2169	2175	rBV	220905	256238	5.90%	0.683%
30	15.192	2223	2228	2235	rBV	228783	271761	6.26%	0.724%
31	15.562	2286	2291	2307	rVB2	682726	1136881	26.19%	3.028%
32	16.033	2366	2371	2377	rBV	133799	209094	4.82%	0.557%
33	16.556	2454	2460	2466	rBV2	79277	149373	3.44%	0.398%
34	17.174	2560	2565	2570	rVB6	26632	48197	1.11%	0.128%

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ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-3

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

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

Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF111319.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

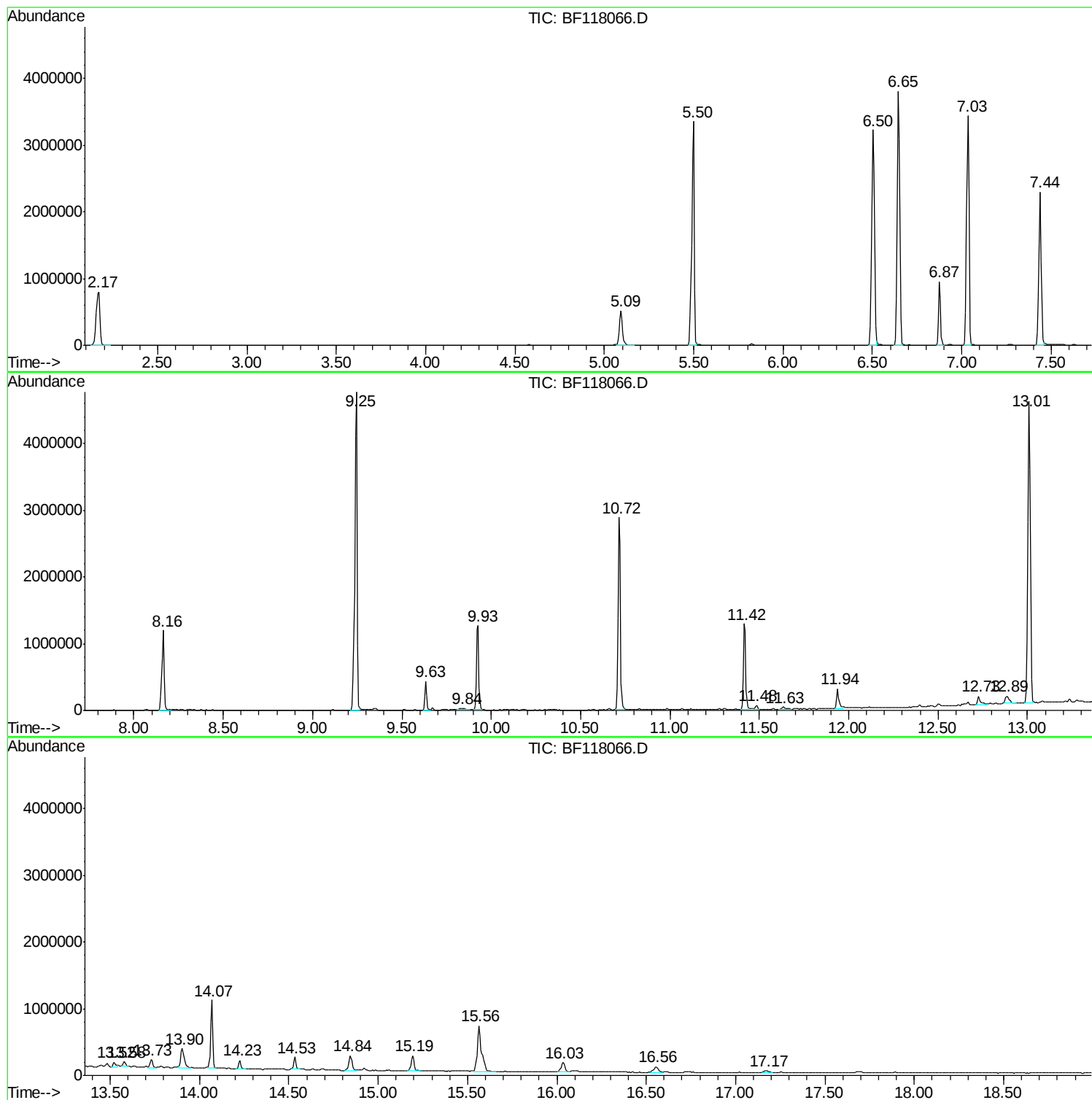
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ClientSampled :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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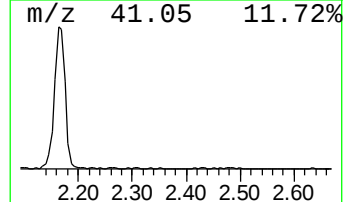
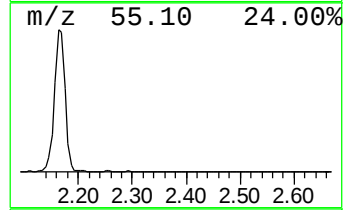
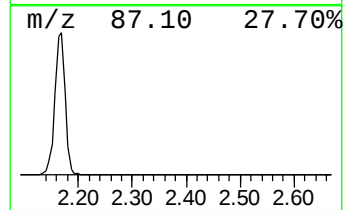
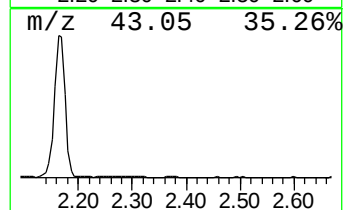
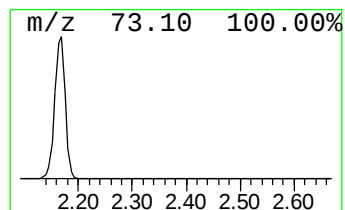
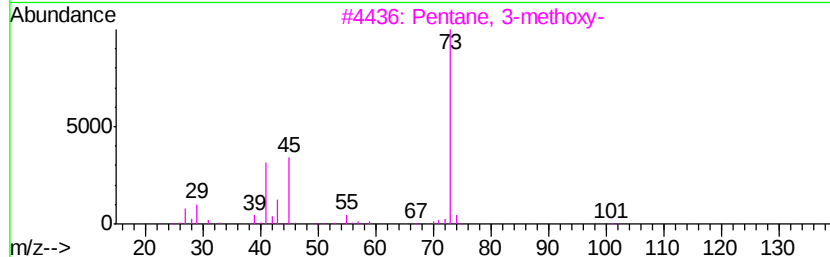
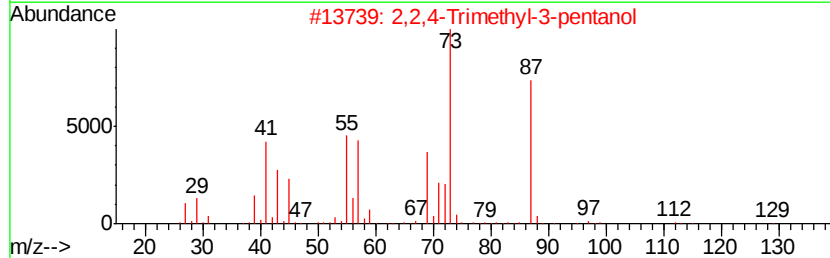
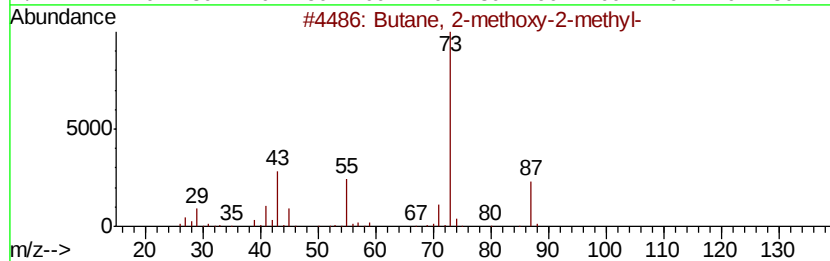
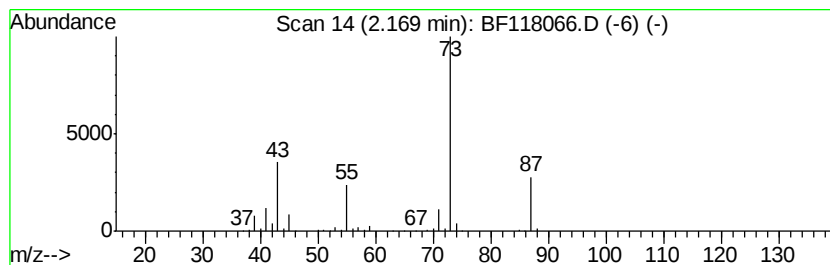
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.17	27.75 ng	1074490	1,4-Dichlorobenzene-d4	6.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
4		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	22
5		Silane, tetramethyl-	88	C4H12Si	000075-76-3	17



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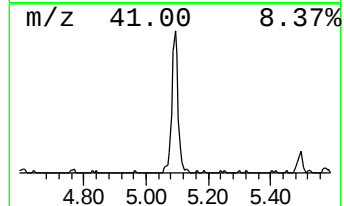
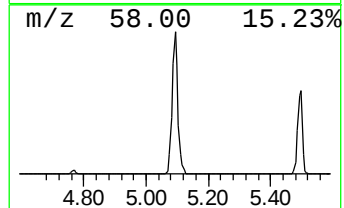
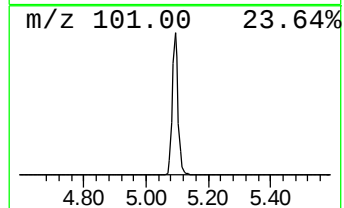
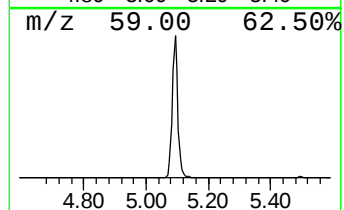
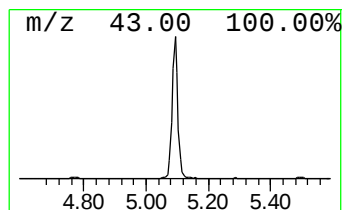
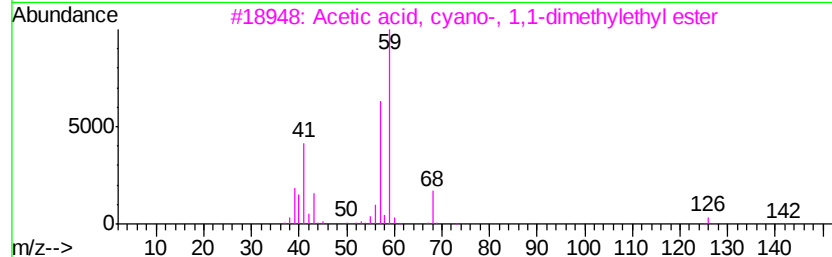
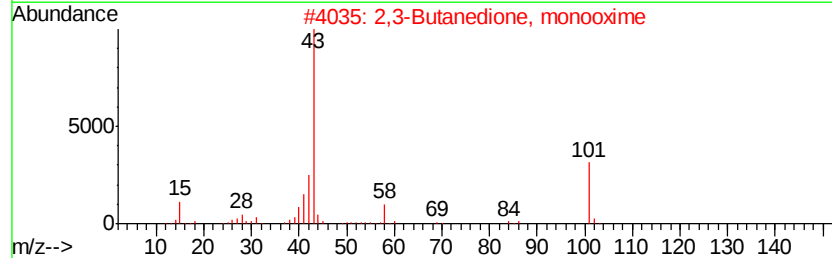
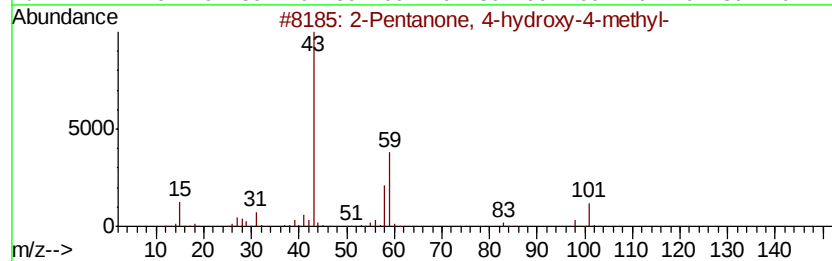
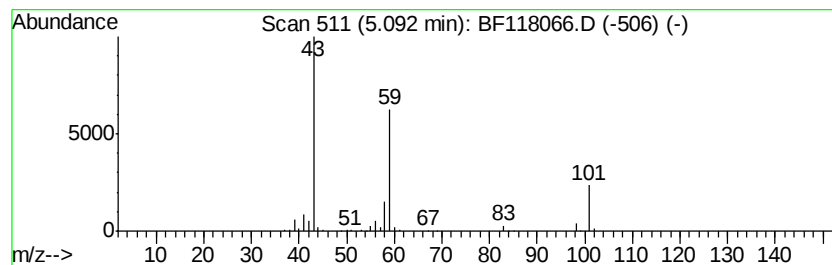
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.09	17.16 ng	664368	1,4-Dichlorobenzene-d4	6.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	17
3		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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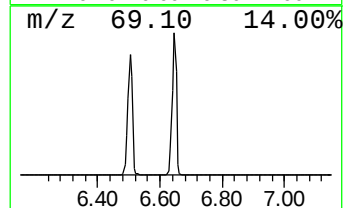
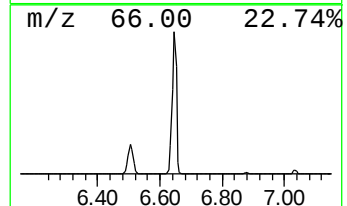
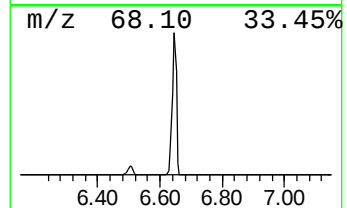
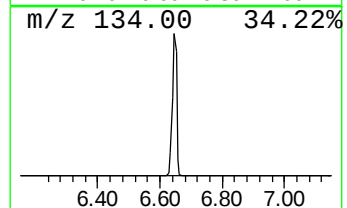
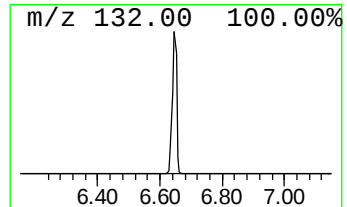
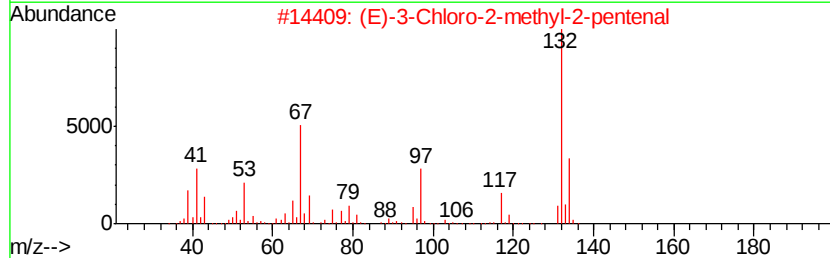
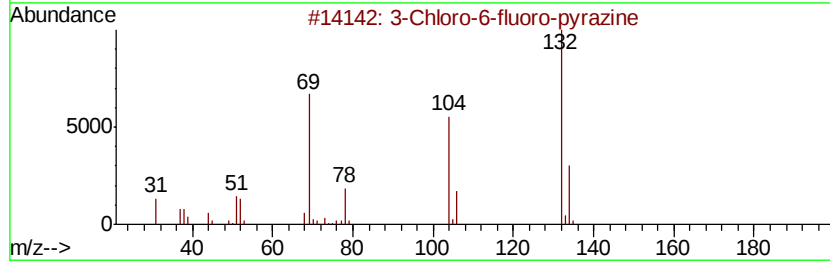
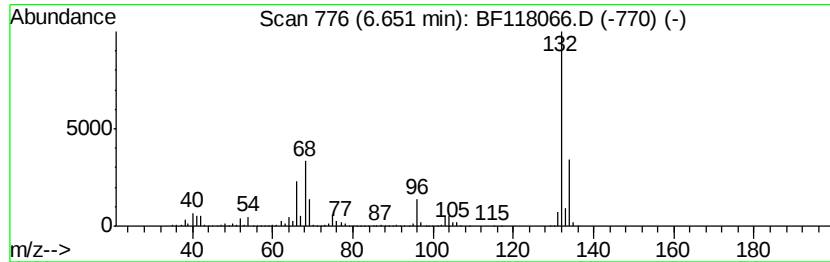
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.65 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.65	94.04 ng	3641250	1,4-Dichlorobenzene-d4	6.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



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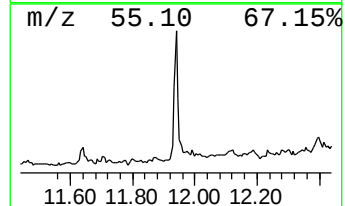
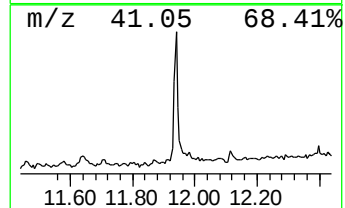
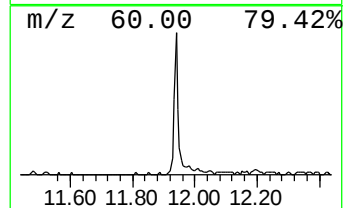
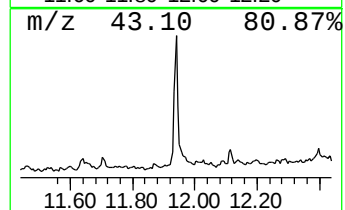
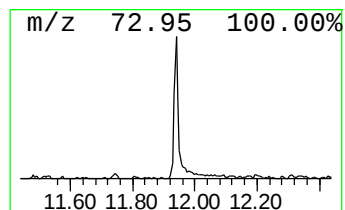
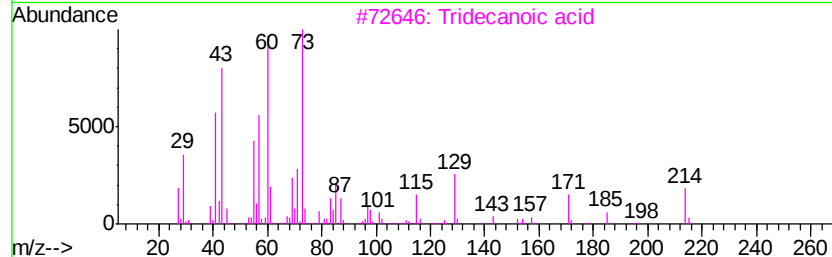
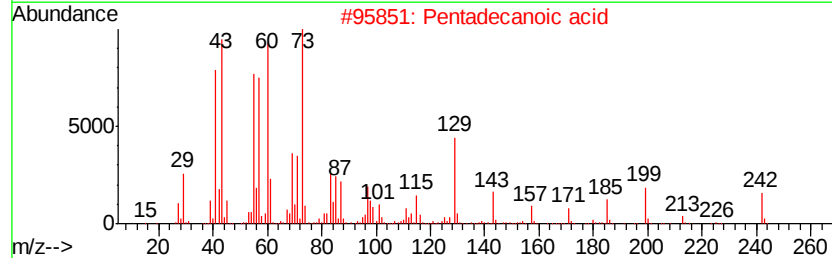
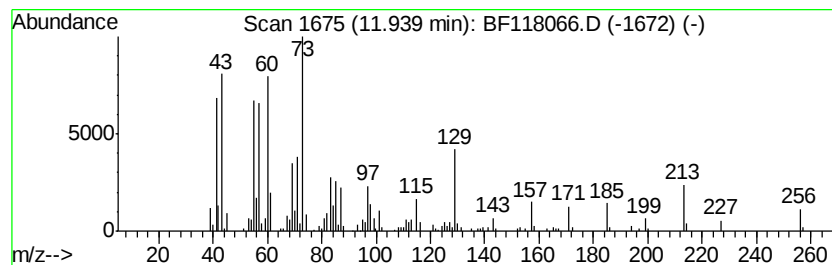
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.94	4.61 ng	272680	Phenanthrene-d10	11.42

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	93
3		Tridecanoic acid	214	C13H26O2	000638-53-9	91
4		Undecanoic acid	186	C11H22O2	000112-37-8	80
5		Tetradecanoic acid	228	C14H28O2	000544-63-8	58



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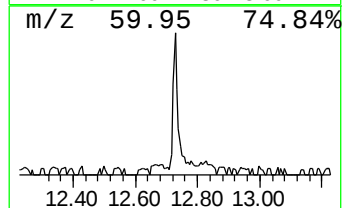
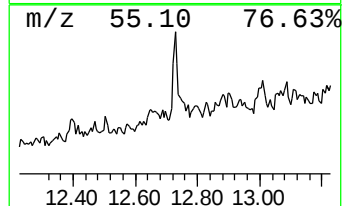
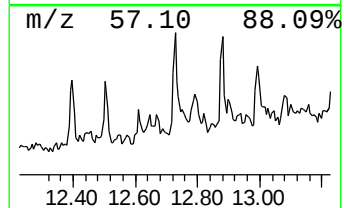
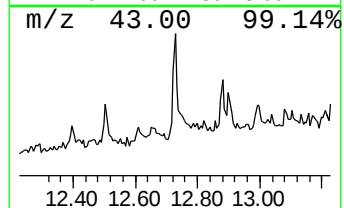
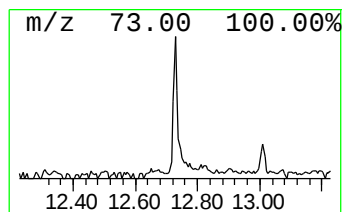
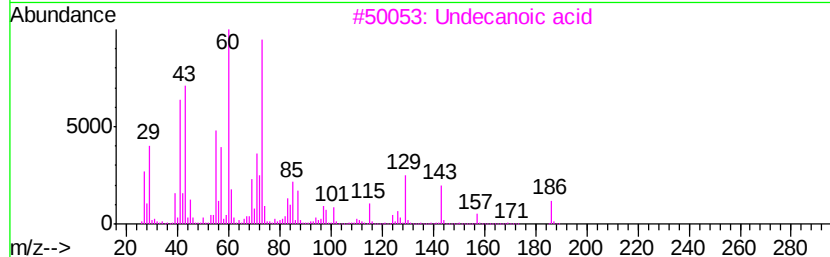
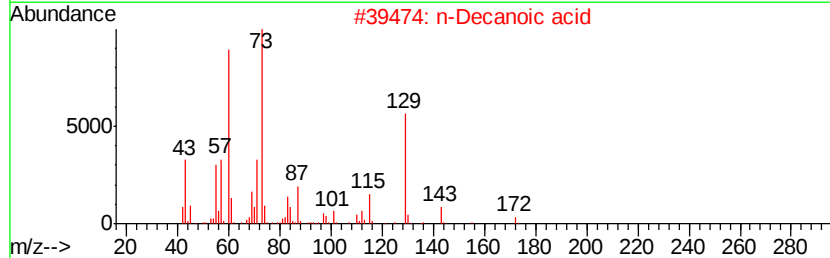
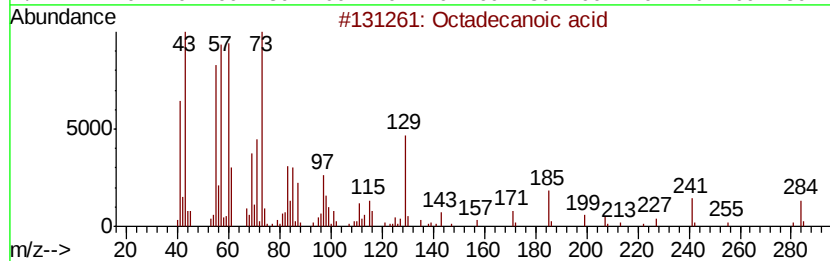
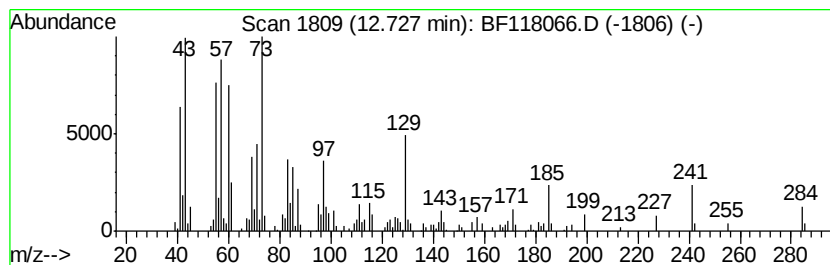
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Octadecanoic acid Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.73	2.94 ng	174156	Phenanthrene-d10	11.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		n-Decanoic acid	172	C10H20O2	000334-48-5	91
3		Undecanoic acid	186	C11H22O2	000112-37-8	74
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	74
5		Pentadecanoic acid	242	C15H30O2	001002-84-2	64



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SB-3

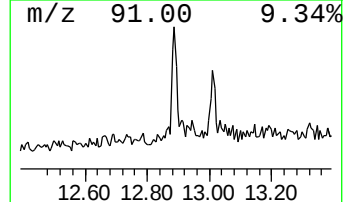
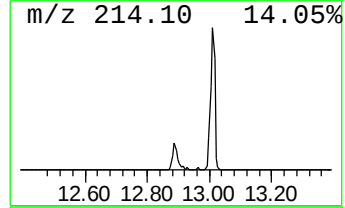
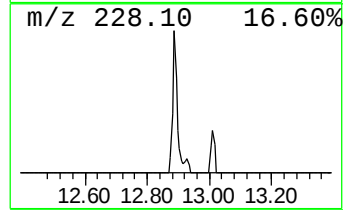
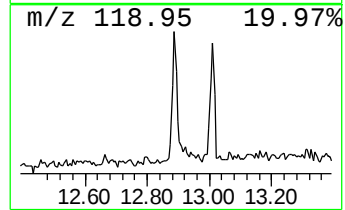
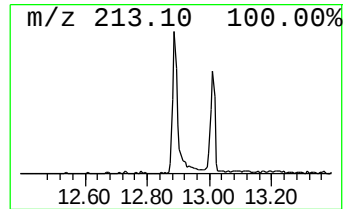
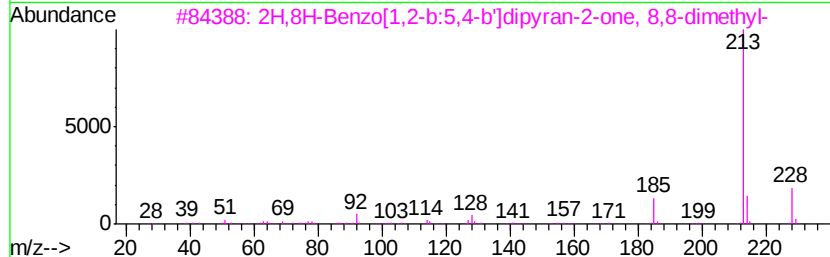
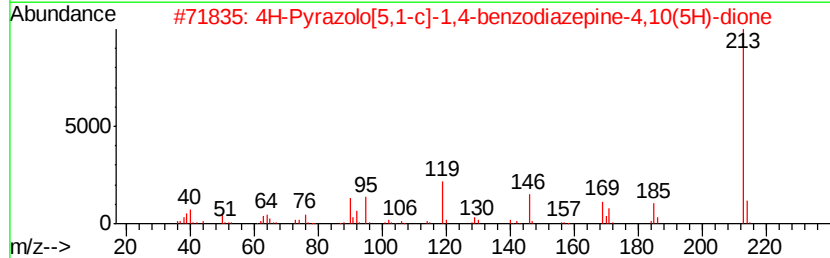
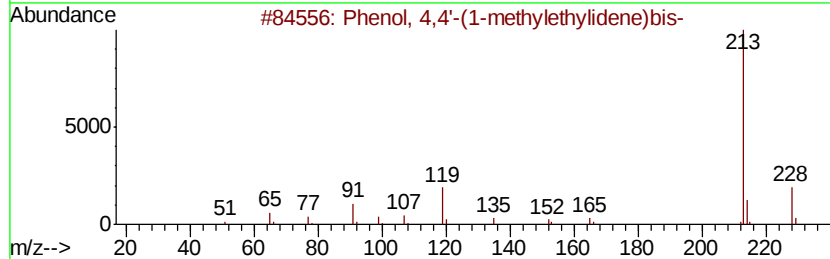
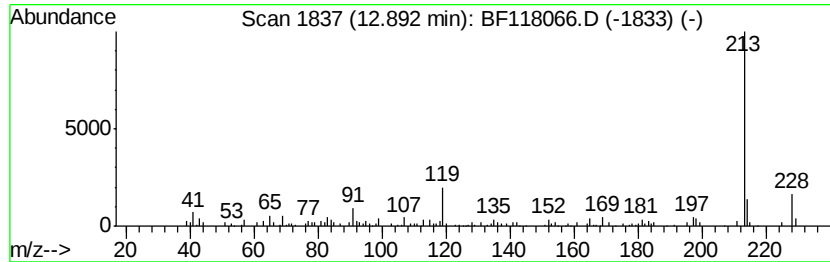
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Phenol, 4,4'-(1-methylethyl... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.89	3.32 ng	142166	Chrysene-d12	14.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenol, 4,4'-(1-methylethylidene)bis-	228	C15H16O2	000080-05-7	97
2		4H-Pyrazolo[5,1-c]-1,4-benzodiazepine-4,10(5H)-dione	213	C11H7N3O2	1000277-20-1	64
3		2H,8H-Benzo[1,2-b:5,4-b']dipyrans-2-one, 8,8-dimethyl-	228	C14H12O3	000553-19-5	59
4		2H,8H-Benzo[1,2-b:3,4-b']dipyrans-2-one, 8,8-dimethyl-	228	C14H12O3	000523-59-1	59
5		Carbamic acid, [1,1'-biphenyl]-3-ylidenebis-	213	C13H11NO2	055030-29-0	52



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120319\
Data File : BF118066.D
Acq On : 3 Dec 2019 17:05
Operator : JU
Sample : K6069-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

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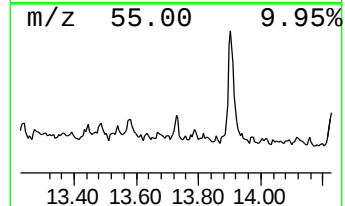
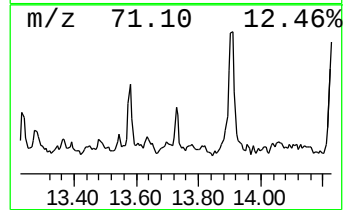
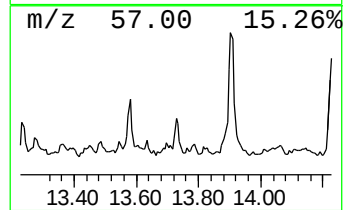
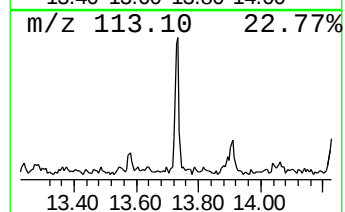
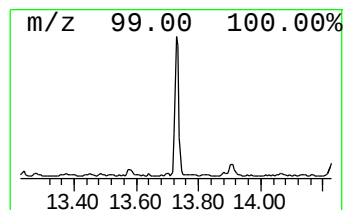
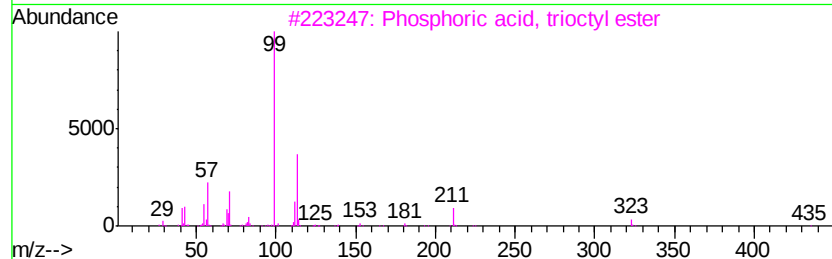
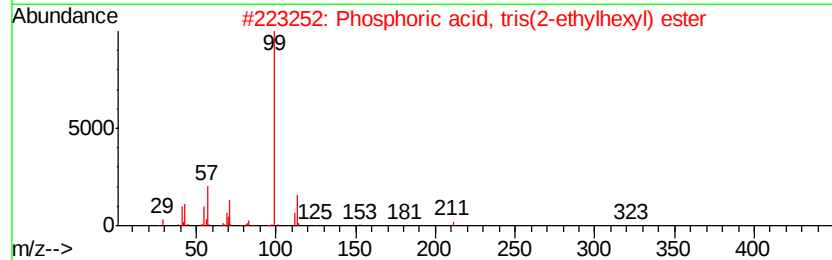
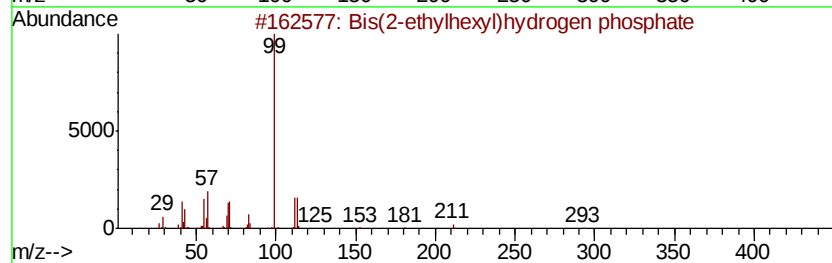
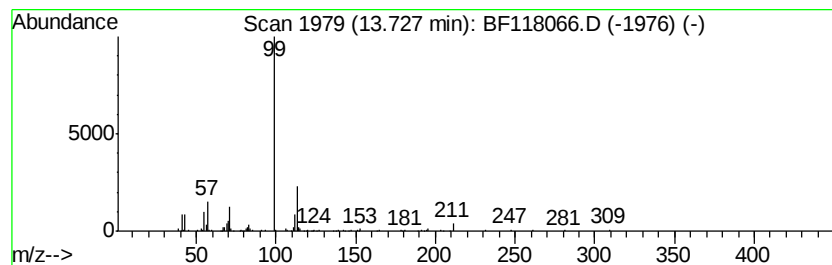
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Bis(2-ethylhexyl)hydrogen p... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.73	2.67 ng	114489	Chrysene-d12	14.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bis(2-ethylhexyl)hydrogen phosphate	322	C16H35O4P	000298-07-7	64
2		Phosphoric acid, tris(2-ethylhex...	434	C24H51O4P	000078-42-2	59
3		Phosphoric acid, trioctyl ester	434	C24H51O4P	001806-54-8	47
4		Phosphoric acid, dihexyl nonyl e...	392	C21H45O4P	1000308-89-2	38
5		Didodecyl phosphate	434	C24H51O4P	007057-92-3	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120319\
Data File : BF118066.D
Acq On : 3 Dec 2019 17:05
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Sample : K6069-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

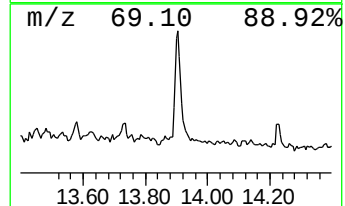
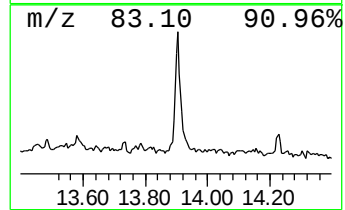
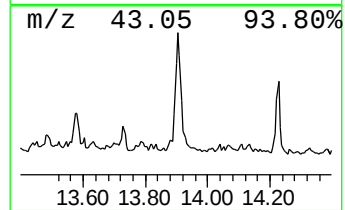
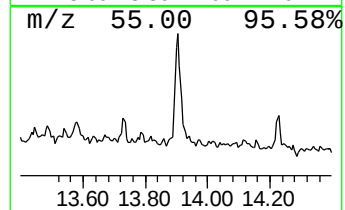
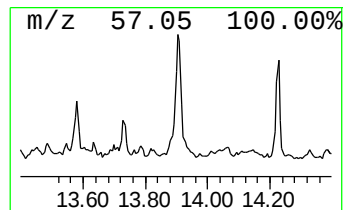
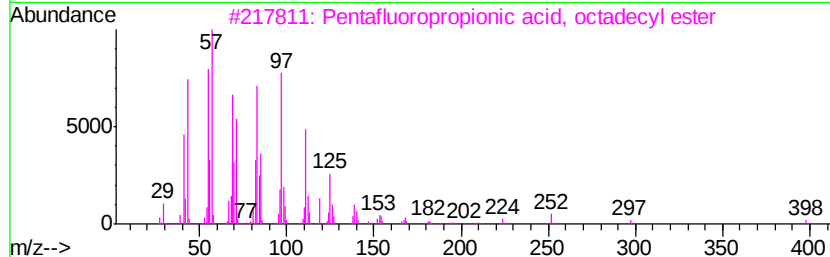
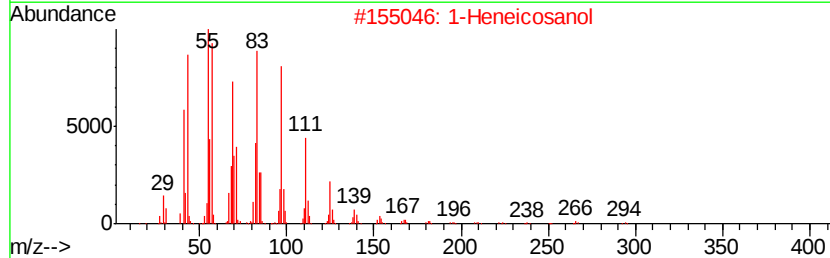
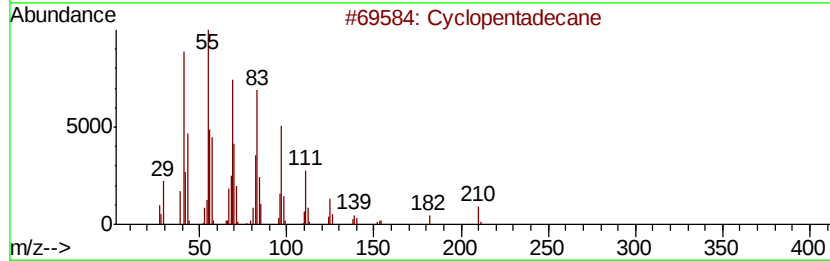
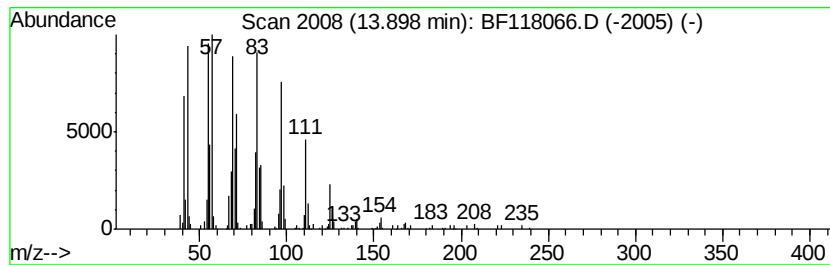
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Cyclopentadecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.90	9.12 ng	390442	Chrysene-d12	14.07	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cvclopentadecane	210	C15H30	000295-48-7	94
2	1-Heneicosanol	312	C21H44O	015594-90-8	93
3	Pentafluoropropionic acid, octad...	416	C21H37F5O2	959261-25-7	91
4	Heptafluorobutvric acid, n-octad...	466	C22H37F7O2	000400-57-7	91
5	Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120319\
 Data File : BF118066.D
 Acq On : 3 Dec 2019 17:05
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 Sample : K6069-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

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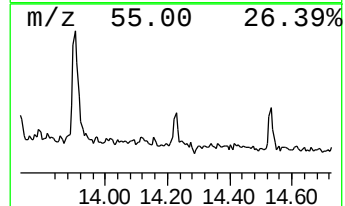
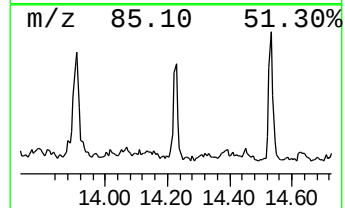
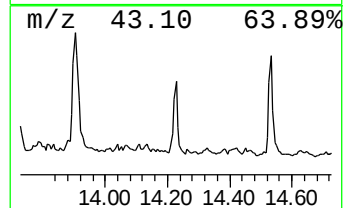
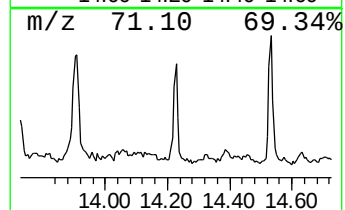
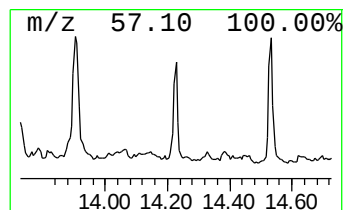
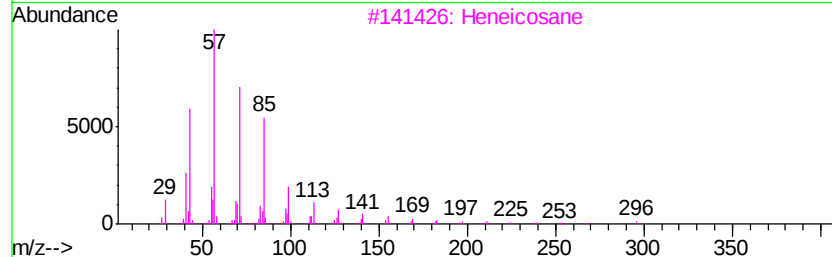
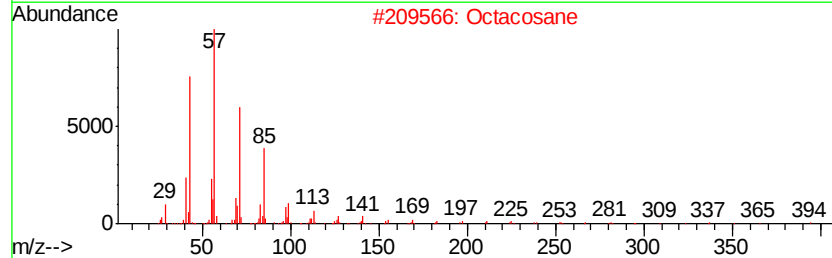
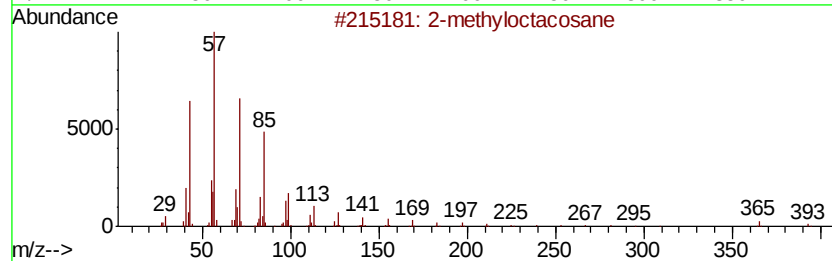
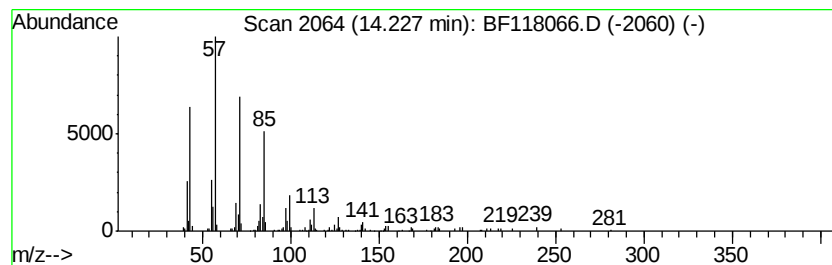
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 10 2-methyloctacosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.23	3.07 ng	131387	Chrysene-d12	14.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-methyloctacosane	408	C29H60	1000376-72-8	91
2		Octacosane	394	C28H58	000630-02-4	91
3		Heneicosane	296	C21H44	000629-94-7	91
4		Hexacosane	366	C26H54	000630-01-3	91
5		Hexatriacontane	507	C36H74	000630-06-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120319\
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Sample : K6069-01
Misc :
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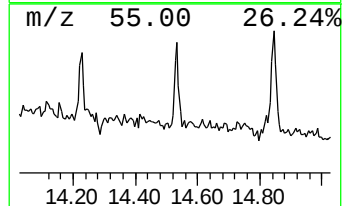
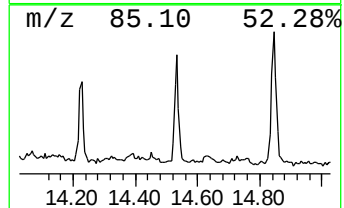
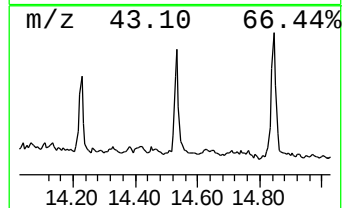
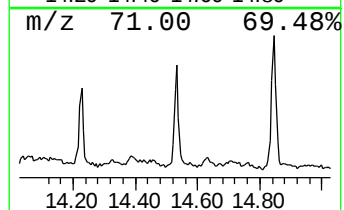
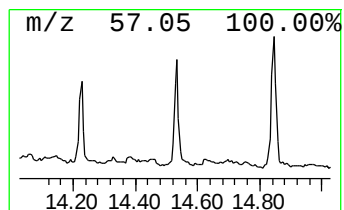
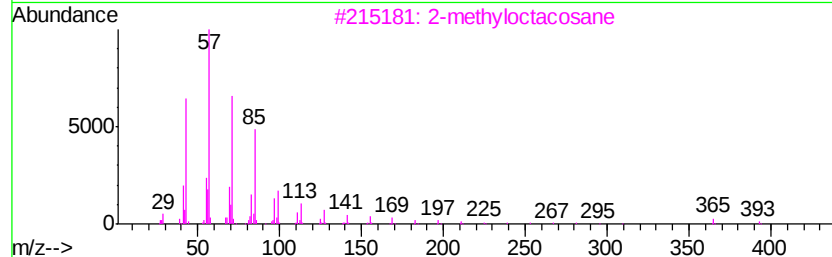
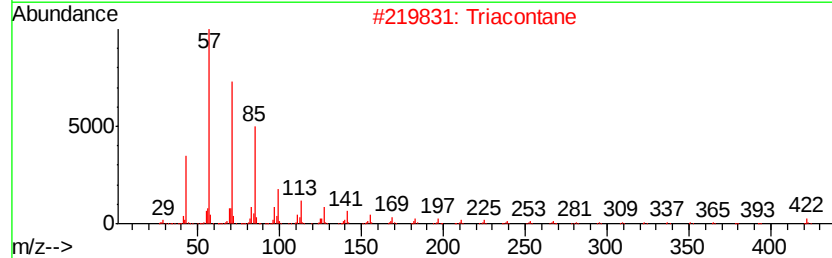
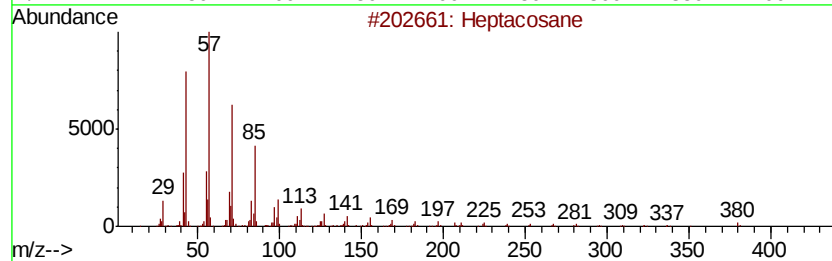
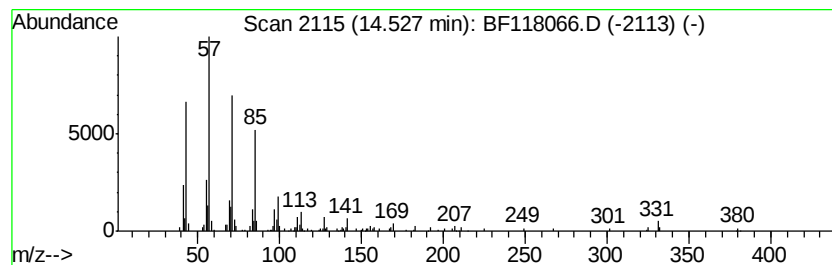
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Heptacosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.53	3.85 ng	164871	Chrysene-d12	14.07
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Heptacosane	380 C27H56	000593-49-7	91
2	triacontane	422 C30H62	000638-68-6	90
3	2-methyloctacosane	408 C29H60	1000376-72-8	90
4	Hexacosane	366 C26H54	000630-01-3	90
5	Heptacosane, 1-chloro-	414 C27H55Cl	062016-79-9	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120319\
Data File : BF118066.D
Acq On : 3 Dec 2019 17:05
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Sample : K6069-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

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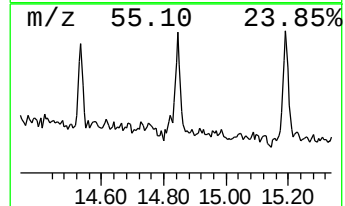
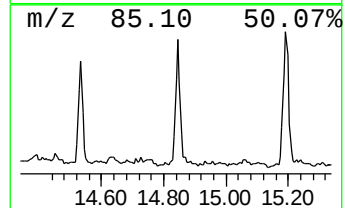
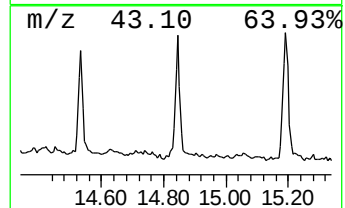
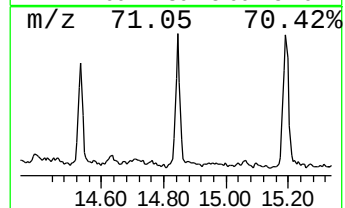
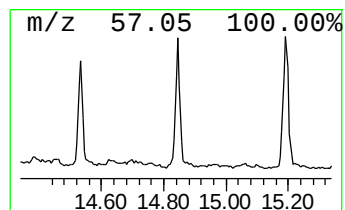
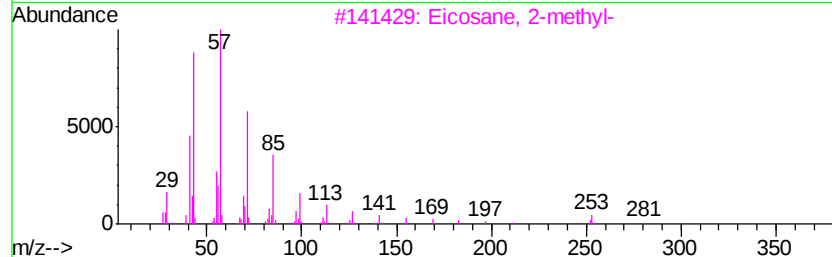
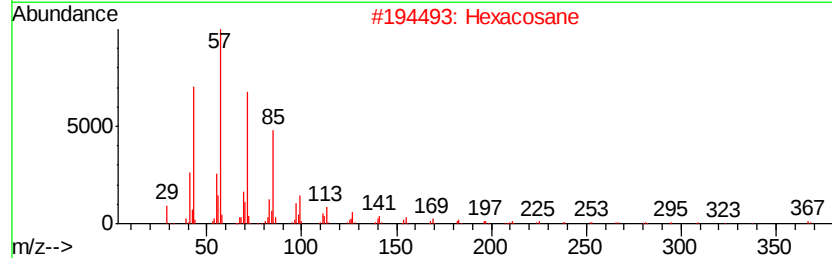
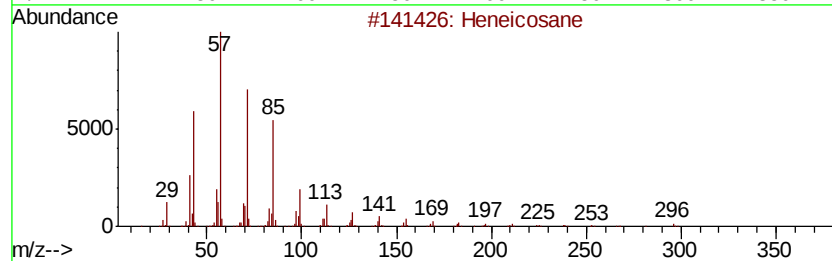
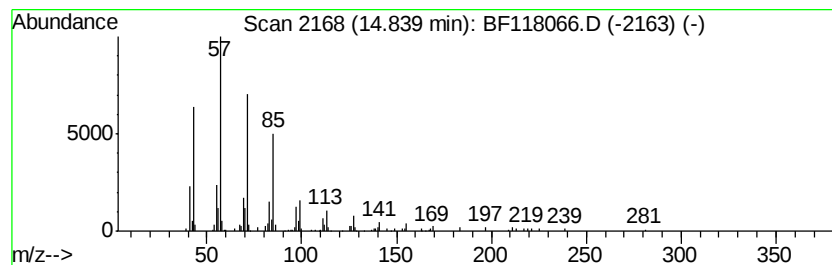
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Heneicosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.84	4.51 ng	256238	Perylene-d12	15.56

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	91
2			Hexacosane	366	C26H54	000630-01-3	91
3			Eicosane, 2-methyl-	296	C21H44	001560-84-5	91
4			Heptadecane	240	C17H36	000629-78-7	91
5			triacontane	422	C30H62	000638-68-6	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120319\
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Misc :
ALS Vial : 9 Sample Multiplier: 1

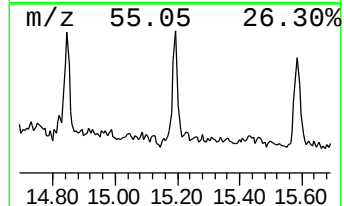
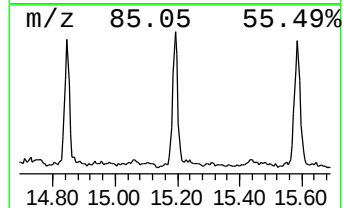
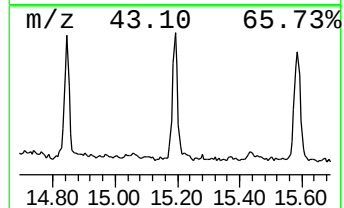
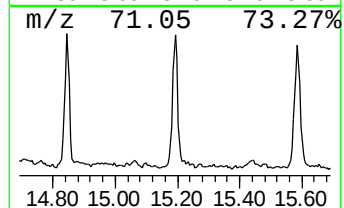
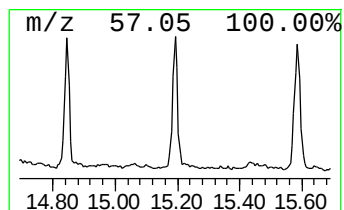
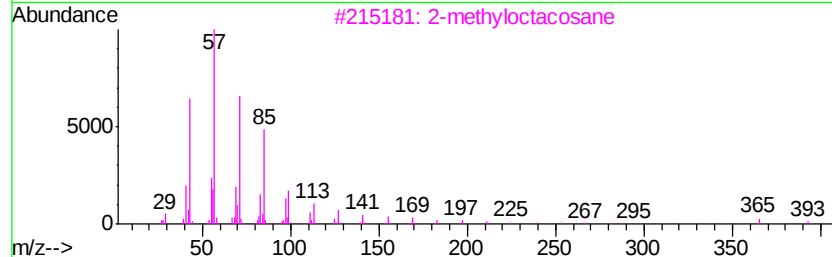
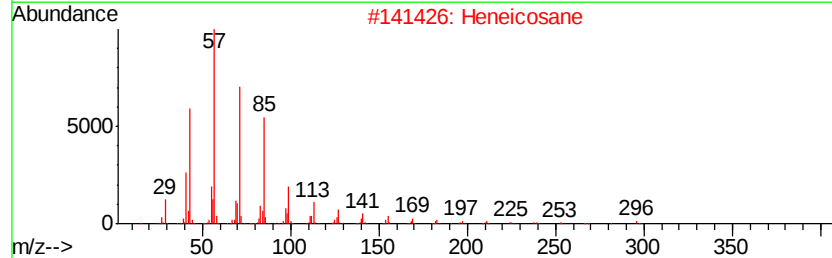
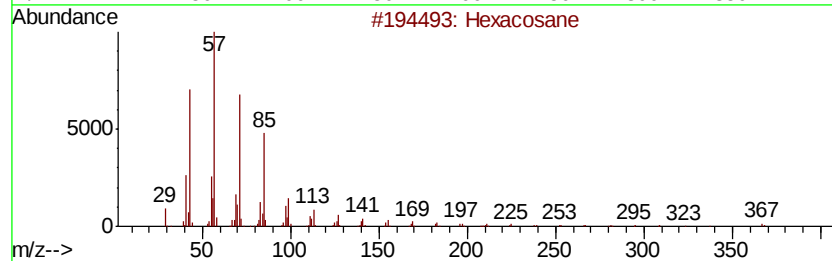
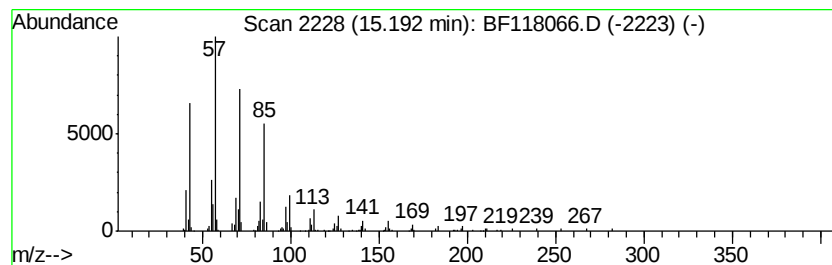
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Hexacosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.19	4.78 ng	271761	Perylene-d12	15.56
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Hexacosane	366 C26H54	000630-01-3	94
2	Heneicosane	296 C21H44	000629-94-7	91
3	2-methyloctacosane	408 C29H60	1000376-72-8	91
4	Heptadecane, 9-octyl-	352 C25H52	007225-64-1	91
5	Tetratetracontane	619 C44H90	007098-22-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120319\
Data File : BF118066.D
Acq On : 3 Dec 2019 17:05
Operator : JU
Sample : K6069-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-3

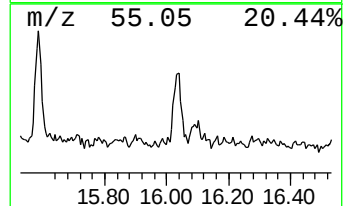
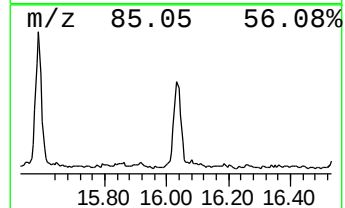
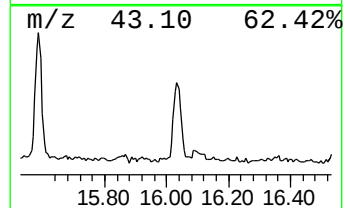
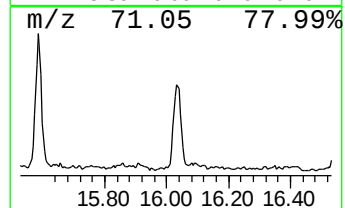
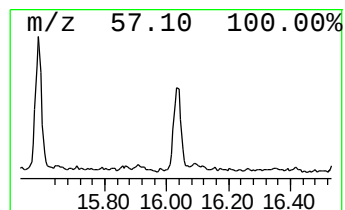
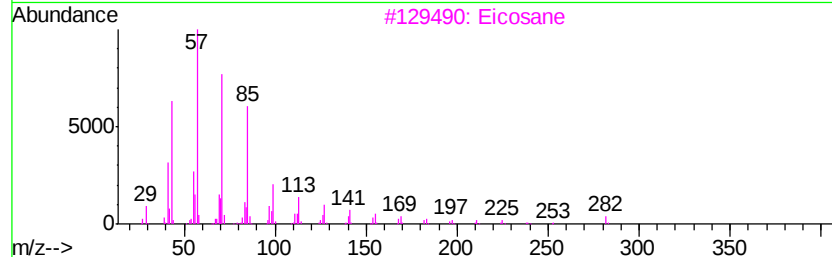
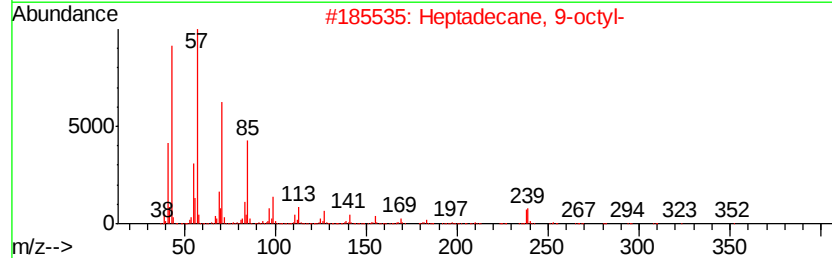
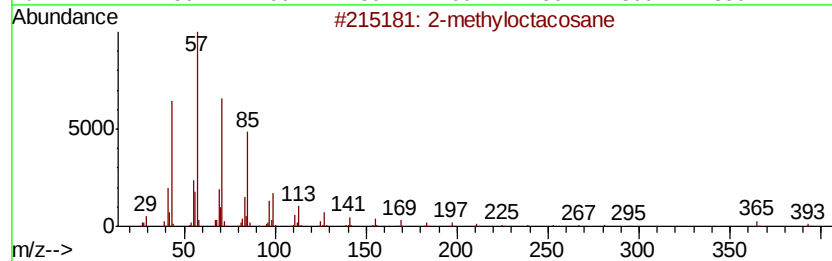
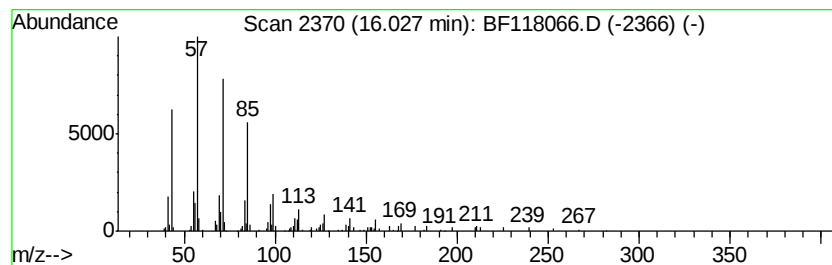
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Heptadecane, 9-octyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.03	3.68 ng	209094	Perylene-d12	15.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-methyloctacosane	408	C29H60	1000376-72-8	91
2		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
3		Eicosane	282	C20H42	000112-95-8	90
4		Octacosane	394	C28H58	000630-02-4	90
5		Octadecane, 1-iodo-	380	C18H37I	000629-93-6	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120319\
Data File : BF118066.D
Acq On : 3 Dec 2019 17:05
Operator : JU
Sample : K6069-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-3

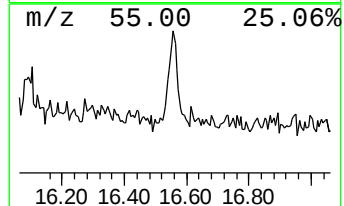
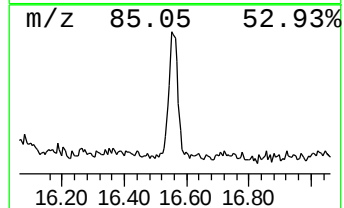
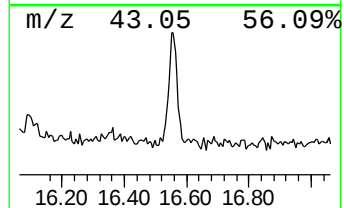
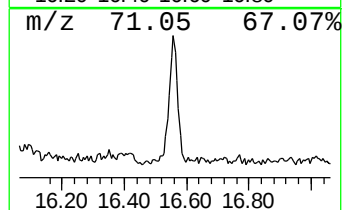
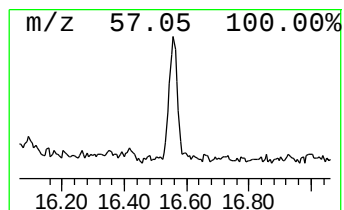
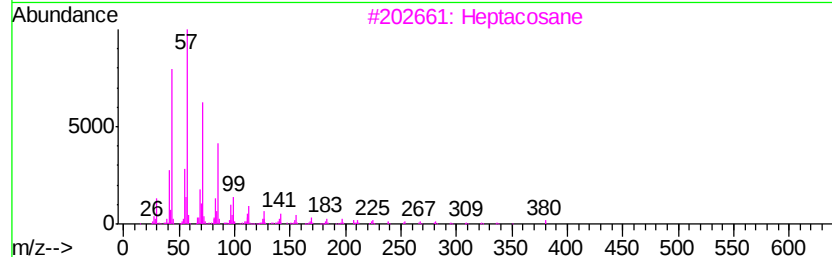
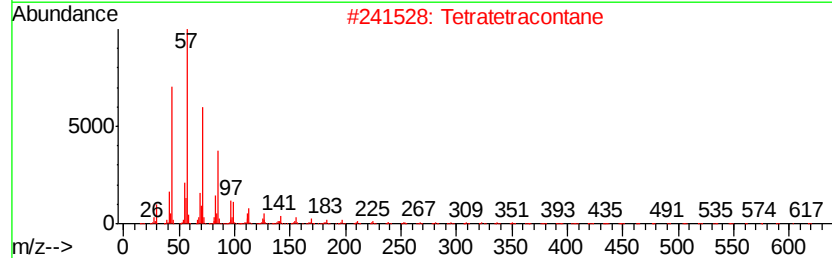
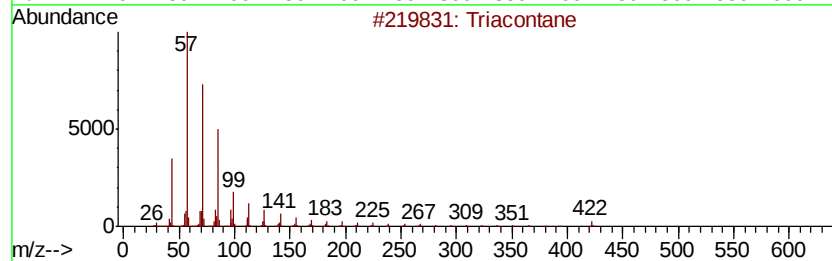
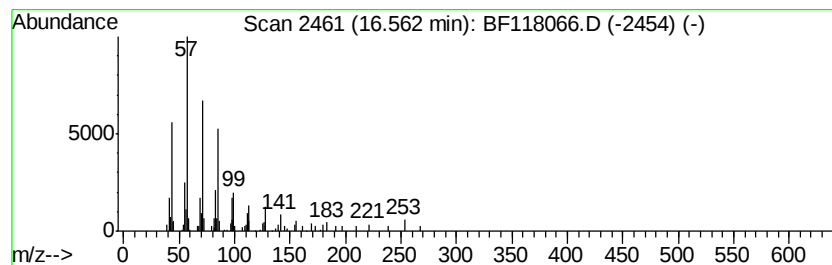
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Triacontane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.56	2.63 ng	149373	Perylene-d12	15.56

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triacontane	422	C30H62	000638-68-6	91
2			Tetratetracontane	619	C44H90	007098-22-8	91
3			Heptacosane	380	C27H56	000593-49-7	91
4			Octacosane	394	C28H58	000630-02-4	91
5			Hexacosane	366	C26H54	000630-01-3	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF120319\
Data File : BF118066.D
Acq On : 3 Dec 2019 17:05
Operator : JU
Sample : K6069-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF111319.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...	2.17	27.8	ng	1074490	1	6.87	774429	20.0
2-Pentanone, 4-hy...	5.09	17.2	ng	664368	1	6.87	774429	20.0
unknown6.65	6.65	94.0	ng	3641250	1	6.87	774429	20.0
n-Hexadecanoic acid	11.94	4.6	ng	272680	4	11.42	1182910	20.0
Octadecanoic acid	12.73	2.9	ng	174156	4	11.42	1182910	20.0
Phenol, 4,4'-(1-m...	12.89	3.3	ng	142166	5	14.07	856658	20.0
Bis(2-ethylhexyl)...	13.73	2.7	ng	114489	5	14.07	856658	20.0
Cyclopentadecane	13.90	9.1	ng	390442	5	14.07	856658	20.0
2-methyloctacosane	14.23	3.1	ng	131387	5	14.07	856658	20.0
Heptacosane	14.53	3.9	ng	164871	5	14.07	856658	20.0
Heneicosane	14.84	4.5	ng	256238	6	15.56	1136880	20.0
Hexacosane	15.19	4.8	ng	271761	6	15.56	1136880	20.0
Heptadecane, 9-oc...	16.03	3.7	ng	209094	6	15.56	1136880	20.0
Triacontane	16.56	2.6	ng	149373	6	15.56	1136880	20.0