

Data Path : Z:\HPCHEM1\BNA F\DATA\BF031918\
 Data File : BF103776.D
 Acq On : 20 Mar 2018 00:32
 Operator : JU/SJ
 Sample : J1971-07
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WEST-GARAGE-6-ST-LATERAL-NW

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:\HPCHEM1\BNA_F\METHODS\8270-BF031518.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	2.334	34	42	52	rVB	149219	208712	5.54%	0.650%
2	5.228	528	534	544	rVB	157416	210051	5.58%	0.654%
3	5.604	592	598	601	rBV	3433931	3731097	99.10%	11.621%
4	5.922	648	652	661	rVB	64278	59093	1.57%	0.184%
5	6.598	761	767	770	rBV	3278079	3765126	100.00%	11.727%
6	6.745	787	792	795	rBV	3715040	3762787	99.94%	11.720%
7	6.975	827	831	834	rBV	994312	830138	22.05%	2.586%
8	7.128	853	857	861	rBV	2910251	2894525	76.88%	9.016%
9	7.534	921	926	929	rBV	2201753	2408225	63.96%	7.501%
10	8.251	1044	1048	1052	rBV	1132319	1053938	27.99%	3.283%
11	9.328	1226	1231	1234	rBV	3568985	3138389	83.35%	9.775%
12	9.716	1292	1297	1300	rBV	310428	237577	6.31%	0.740%
13	9.927	1329	1333	1336	rBV	107004	80380	2.13%	0.250%
14	10.010	1343	1347	1351	rBV	1111746	1036089	27.52%	3.227%
15	10.427	1416	1418	1421	rVB	128625	90985	2.42%	0.283%
16	10.804	1478	1482	1485	rBV	2104608	1956527	51.96%	6.094%
17	10.904	1496	1499	1504	rVB	99917	98307	2.61%	0.306%
18	11.351	1571	1575	1578	rBV	56793	48192	1.28%	0.150%
19	11.504	1597	1601	1604	rBV	1047599	875342	23.25%	2.726%
20	12.016	1684	1688	1697	rBV2	147928	154414	4.10%	0.481%
21	13.092	1867	1871	1874	rBV	2750949	2520783	66.95%	7.851%
22	13.974	2018	2021	2030	rBV	438241	408341	10.85%	1.272%
23	14.157	2048	2052	2055	rBV	987765	852934	22.65%	2.657%
24	14.621	2127	2131	2135	rBV2	85863	94572	2.51%	0.295%
25	14.921	2178	2182	2192	rBV	466596	535580	14.22%	1.668%
26	15.015	2196	2198	2202	rVB	37623	40210	1.07%	0.125%
27	15.698	2308	2314	2321	rBV	590004	754373	20.04%	2.350%
28	16.221	2399	2403	2411	rBV5	57220	99873	2.65%	0.311%
29	17.451	2608	2612	2621	rVB5	37801	80164	2.13%	0.250%
30	17.892	2680	2687	2695	rBV5	29021	79076	2.10%	0.246%

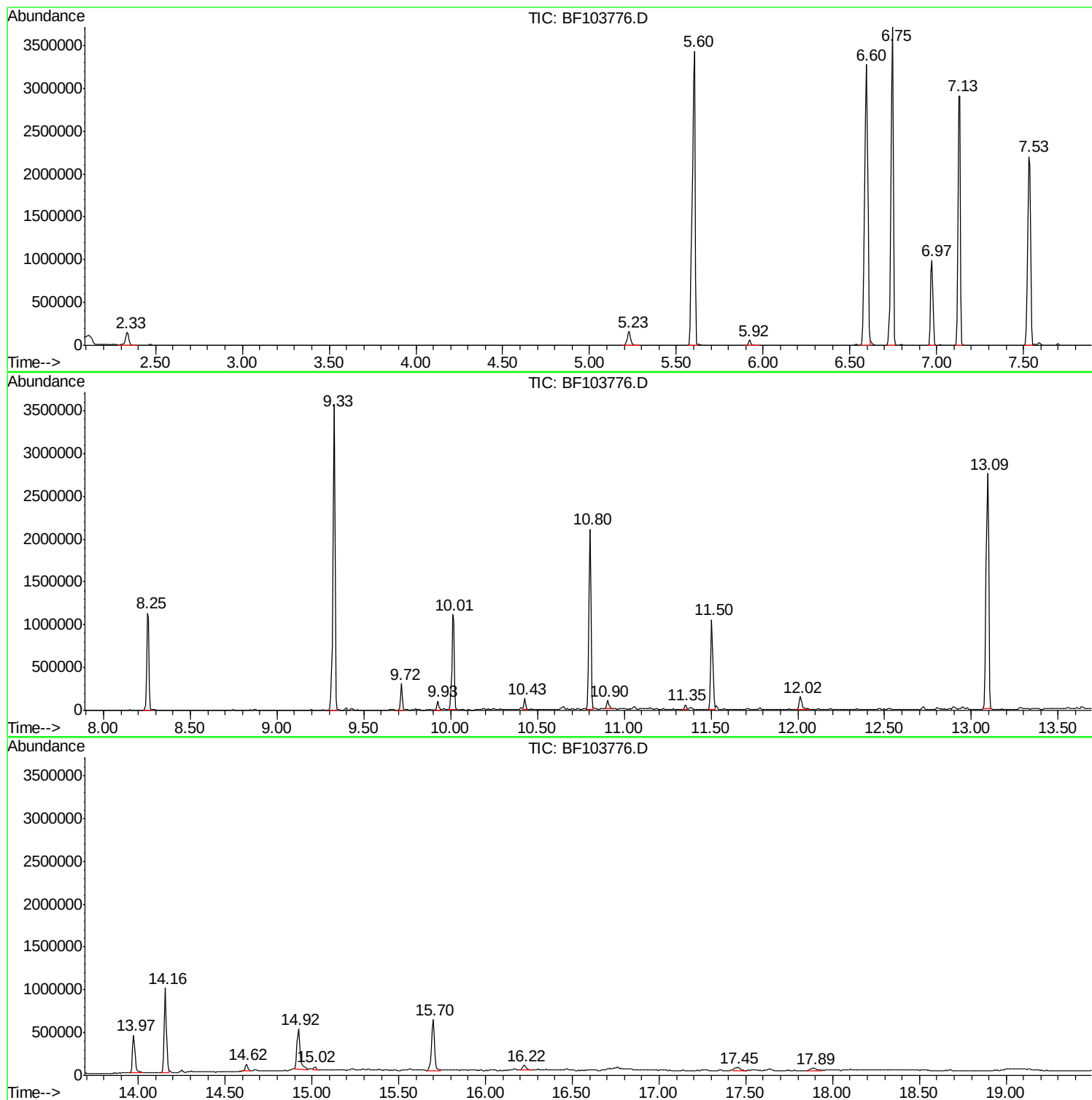
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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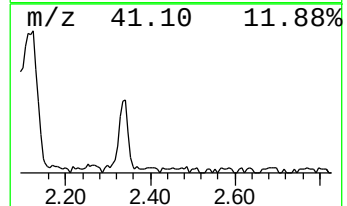
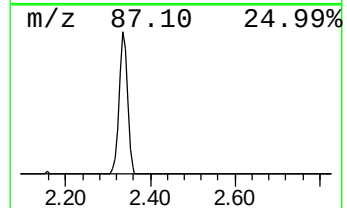
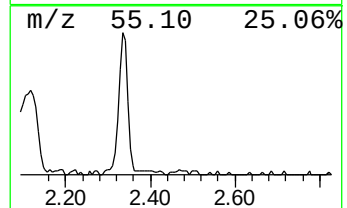
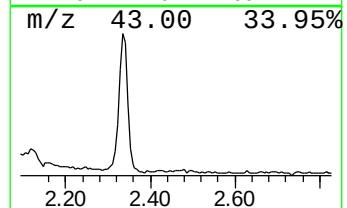
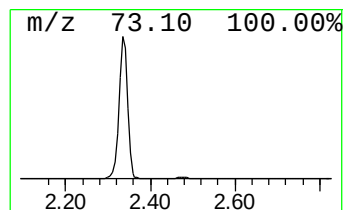
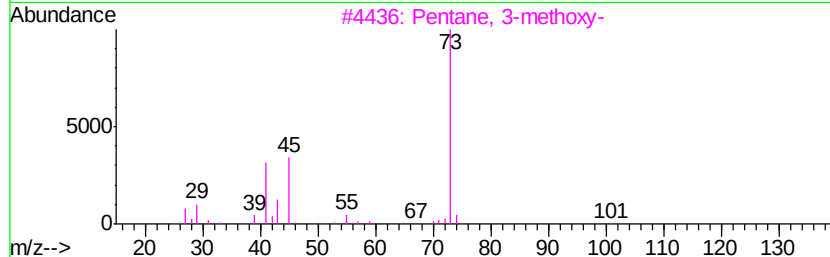
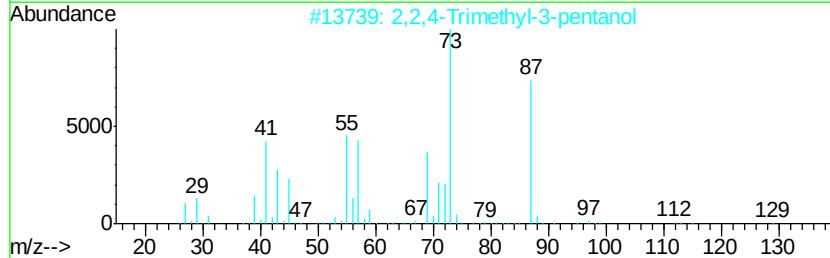
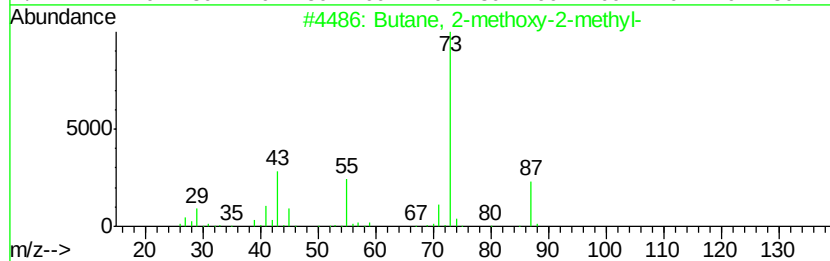
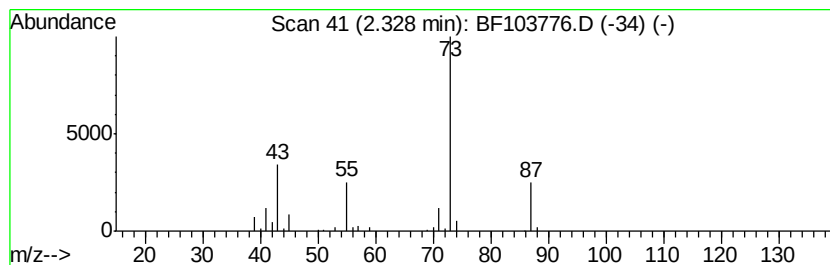
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.33	5.03 ng	208712	1,4-Dichlorobenzene-d4	6.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	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		3-Hexanol, 2-methyl-	116	C7H16O	000617-29-8	17
5		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	12



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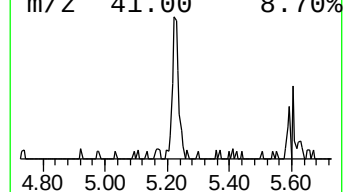
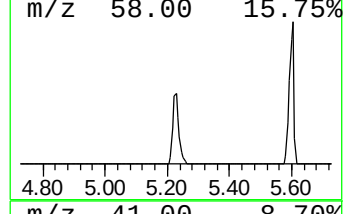
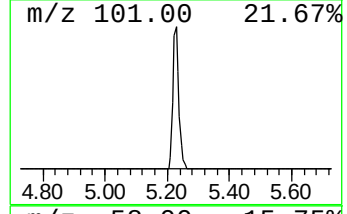
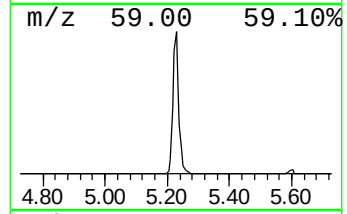
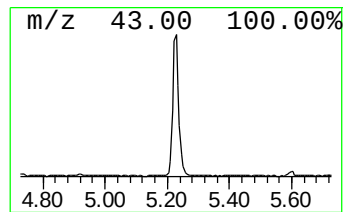
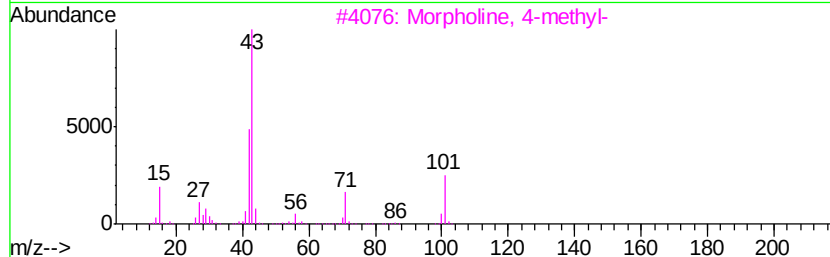
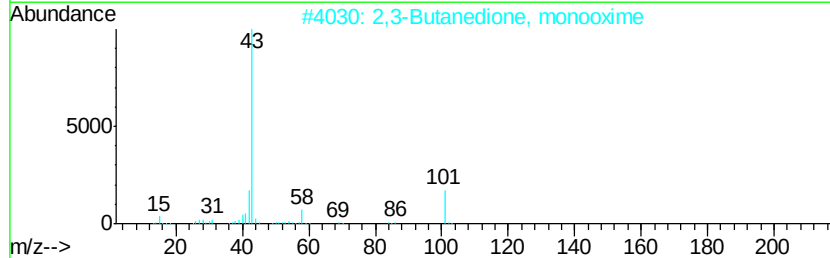
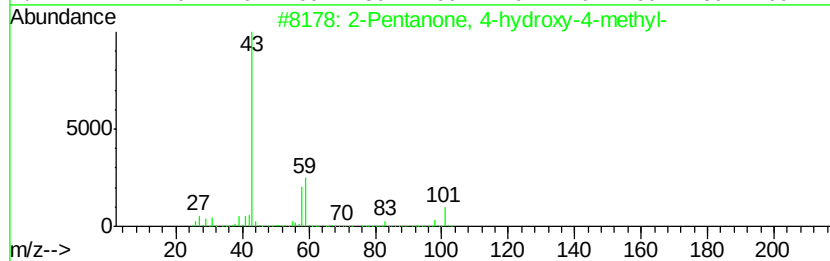
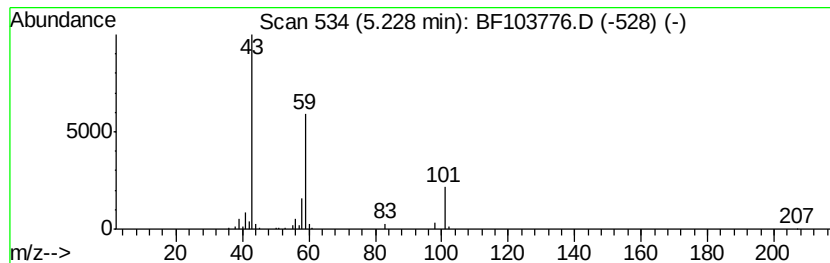
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
5.23	5.06 ng	210051	1,4-Dichlorobenzene-d4	6.97	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
3	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4	5-Hexen-2-one	98	C6H10O	000109-49-9	9
5	Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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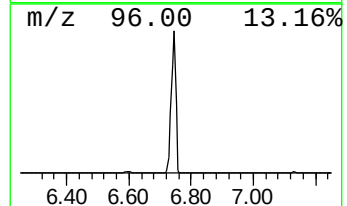
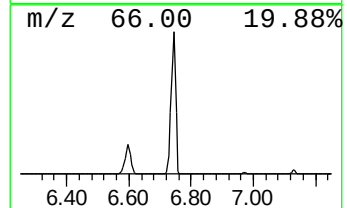
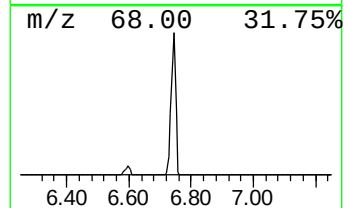
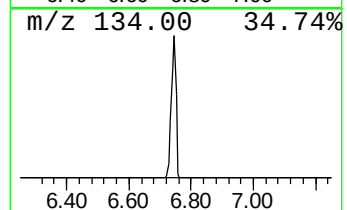
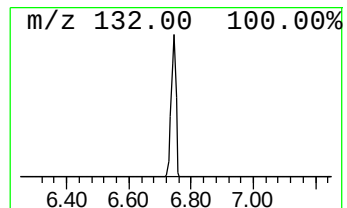
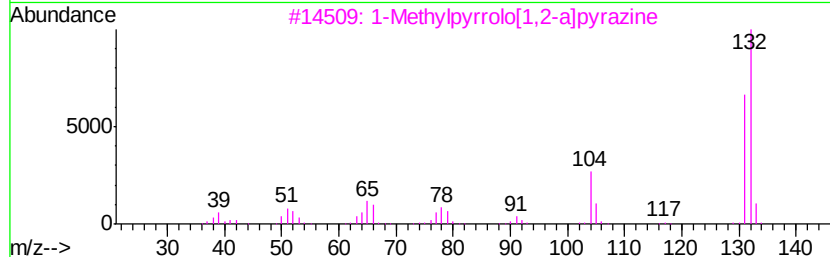
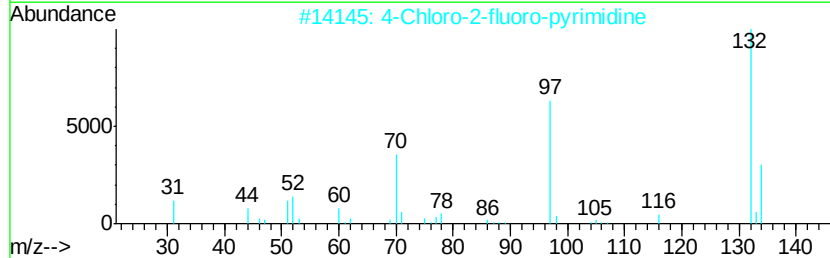
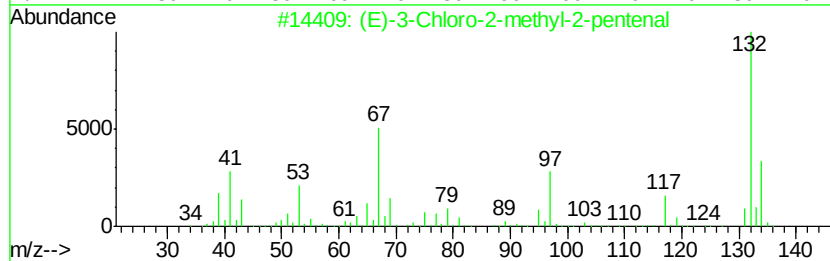
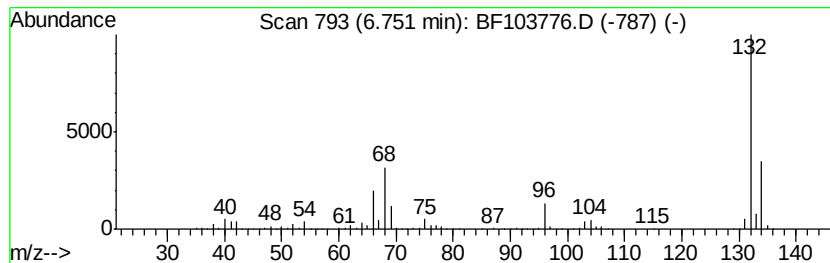
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Peak Number 3 unknown6.75 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.75	90.65 ng	3762790	1,4-Dichlorobenzene-d4	6.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
2		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
3		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9
4		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
5		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9



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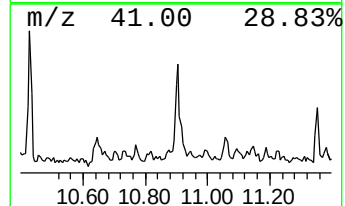
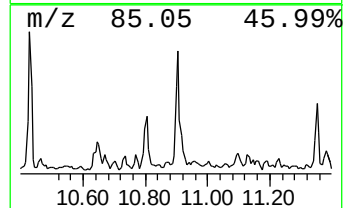
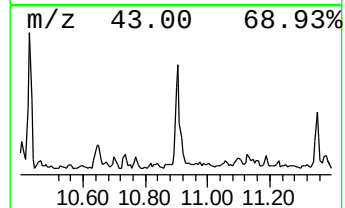
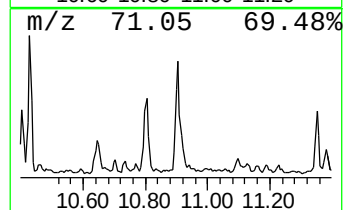
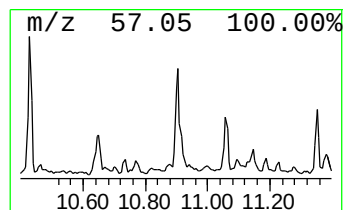
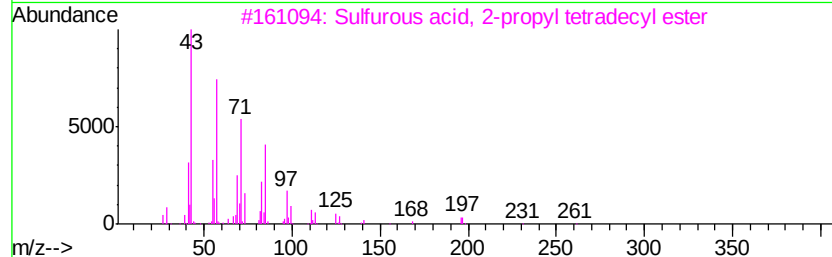
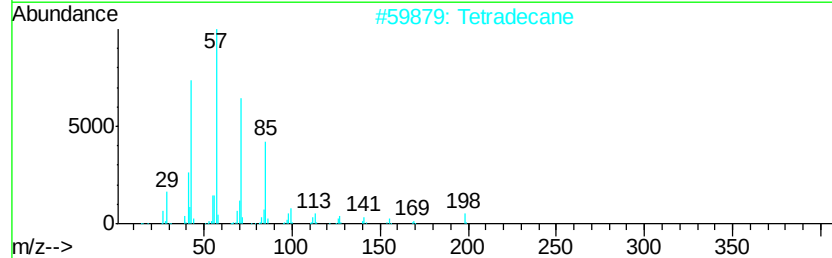
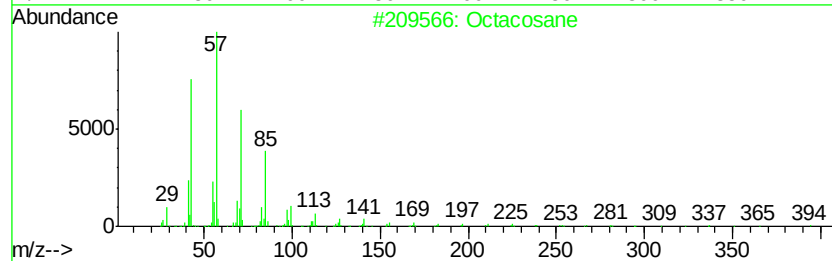
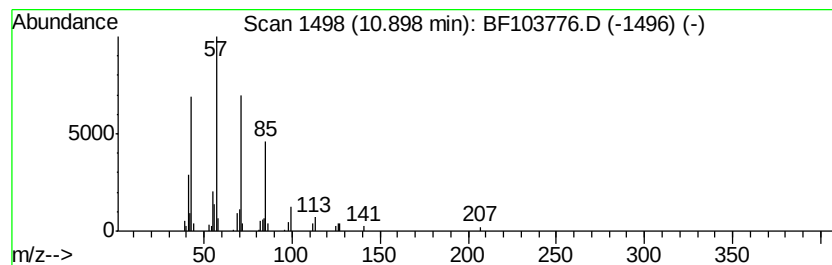
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Peak Number 5 Octacosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.90	2.25 ng	98307	Phenanthrene-d10		11.50
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octacosane	394	C28H58	000630-02-4	91
2	Tetradecane	198	C14H30	000629-59-4	90
3	Sulfurous acid, 2-propyl tetrade...	320	C17H36O3S	1000309-12-5	90
4	Pentadecane	212	C15H32	000629-62-9	90
5	Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	90



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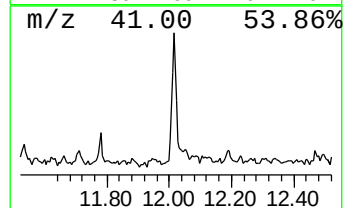
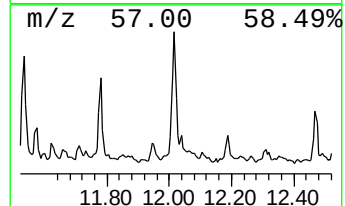
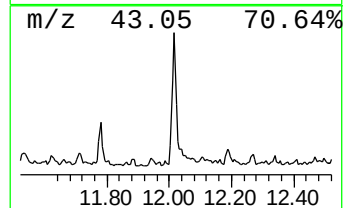
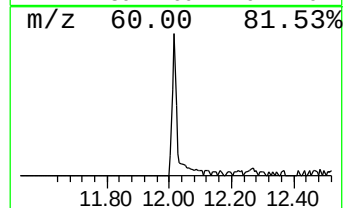
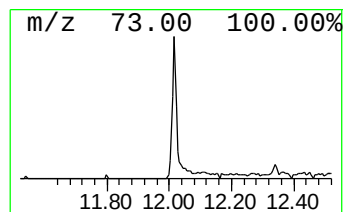
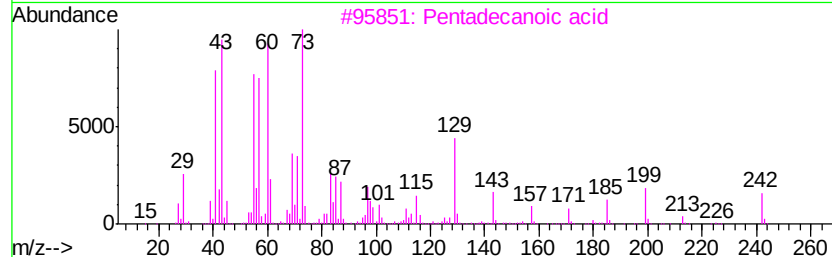
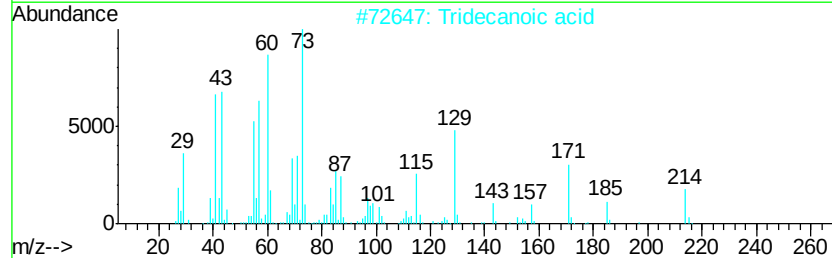
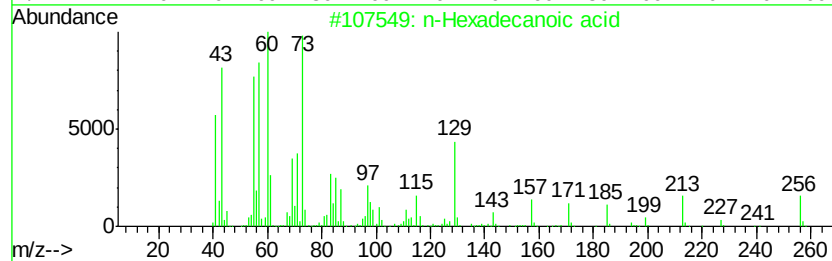
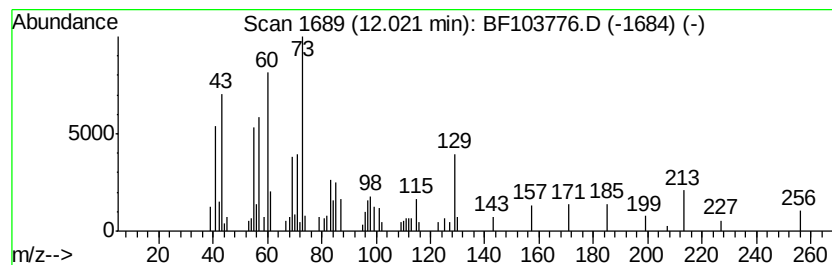
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 6 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.02	3.53 ng	154414	Phenanthrene-d10	11.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tridecanoic acid	214	C13H26O2	000638-53-9	90
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	87
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	86
5		Ethanone, 1-(4,5-dihydro-2-thiaz...	129	C5H7NOS	029926-41-8	43



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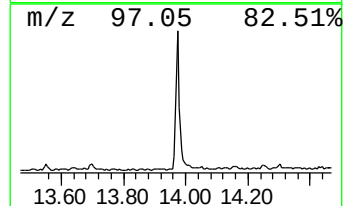
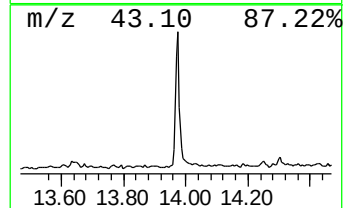
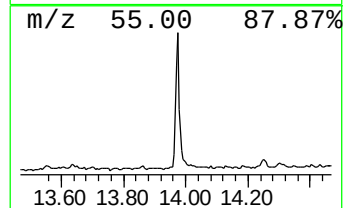
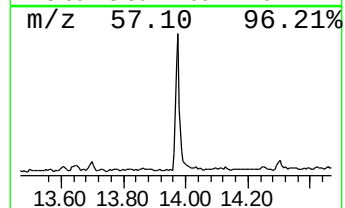
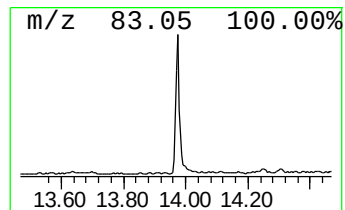
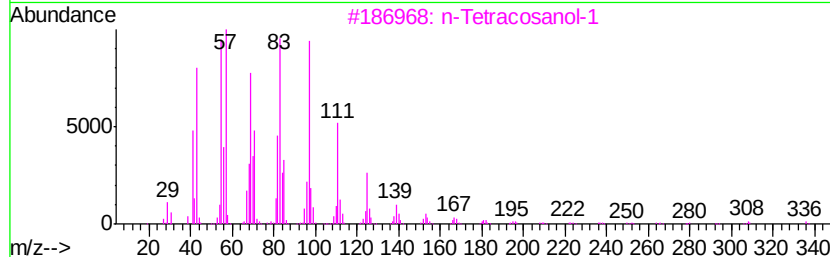
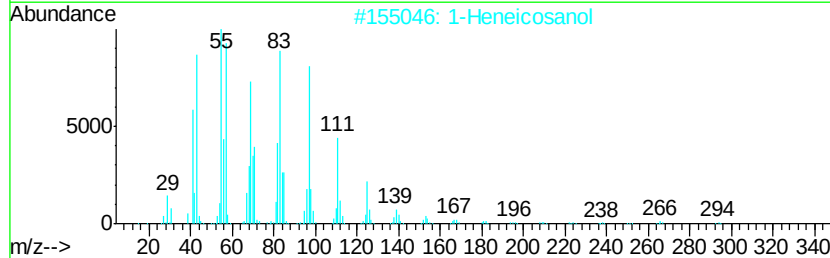
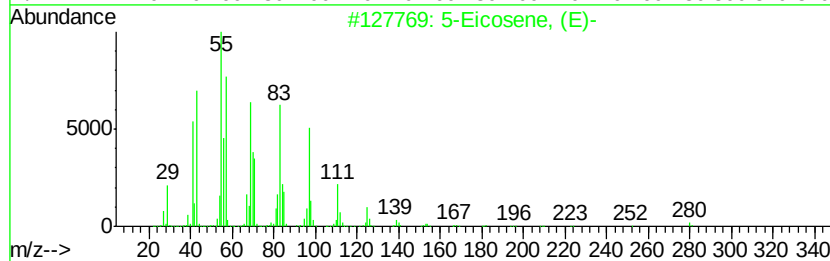
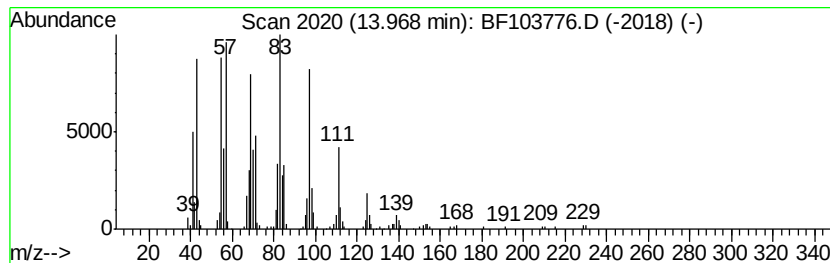
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BNA_F
ClientSampleId :
WEST-GARAGE-6-ST-LATERAL-NW

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 5-Eicosene, (E)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.97	9.57 ng	408341	Chrysene-d12	14.16		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Eicosene, (E)-		280	C20H40	074685-30-6	99
2	1-Heneicosanol		312	C21H44O	015594-90-8	95
3	n-Tetracosanol-1		354	C24H50O	000506-51-4	95
4	n-Nonadecanol-1		284	C19H40O	001454-84-8	95
5	2- Chloropropionic acid, octadec...		360	C21H41ClO2	088104-31-8	94



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031918\
Data File : BF103776.D
Acq On : 20 Mar 2018 00:32
Operator : JU/SJ
Sample : J1971-07
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WEST-GARAGE-6-ST-LATERAL-NW

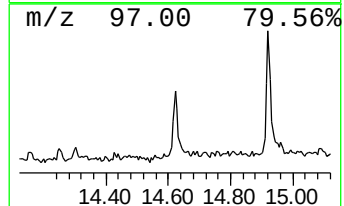
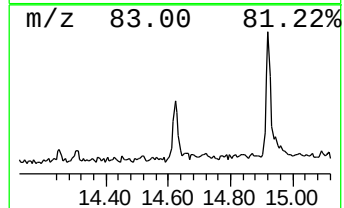
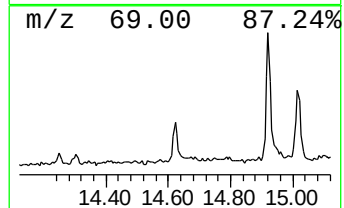
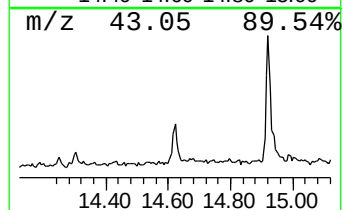
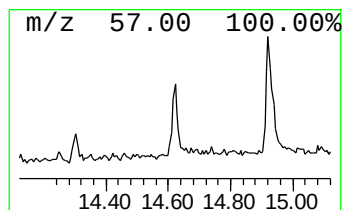
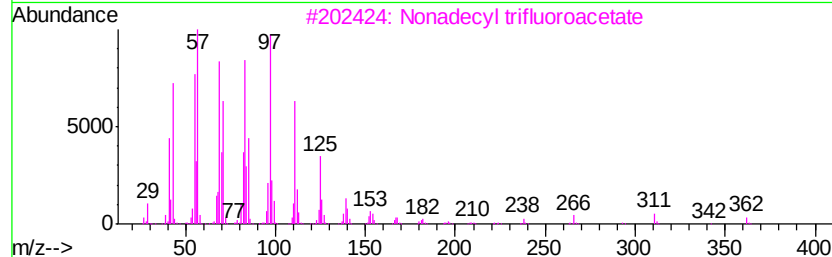
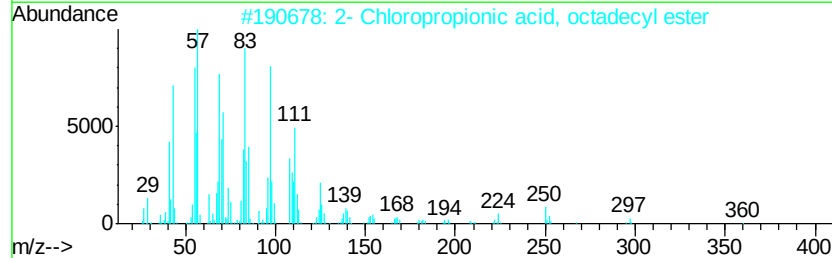
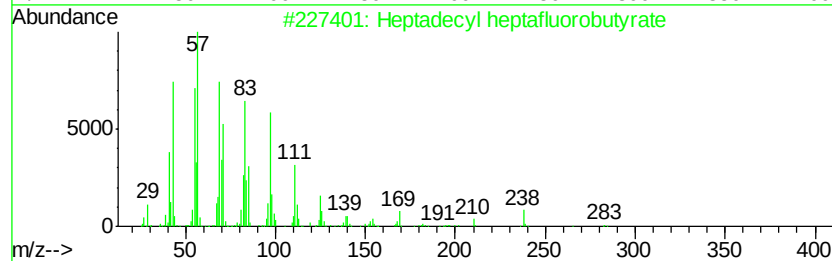
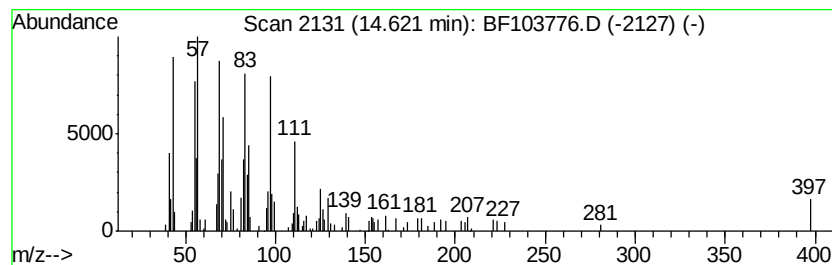
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Heptadecyl heptafluorobutyrate Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.62	2.22 ng	94572	Chrysene-d12	14.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	95
2		2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	94
3		Nonadecyl trifluoroacetate	380	C21H39F3O2	1000351-76-3	94
4		Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	93
5		1-Heptacosanol	396	C27H56O	002004-39-9	93



Data Path : Z:\HPCHEM1\BNA F\DATA\BF031918\
Data File : BF103776.D
Acq On : 20 Mar 2018 00:32
Operator : JU/SJ
Sample : J1971-07
Misc :
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Instrument :
BNA_F
ClientSampleId :
WEST-GARAGE-6-ST-LATERAL-NW

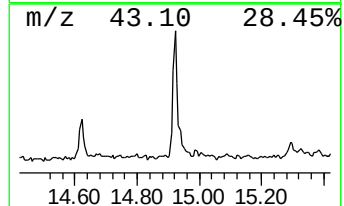
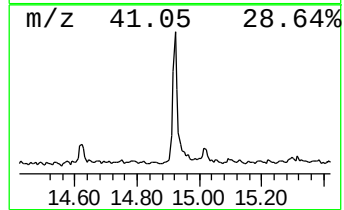
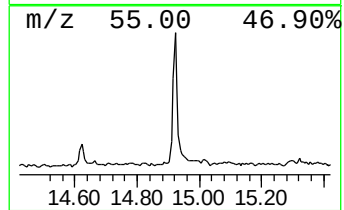
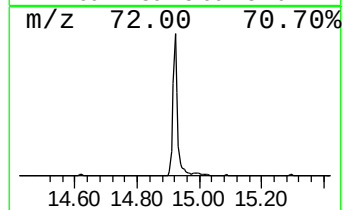
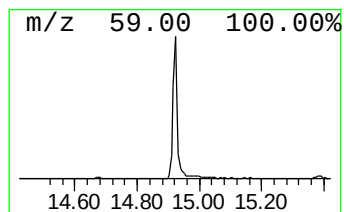
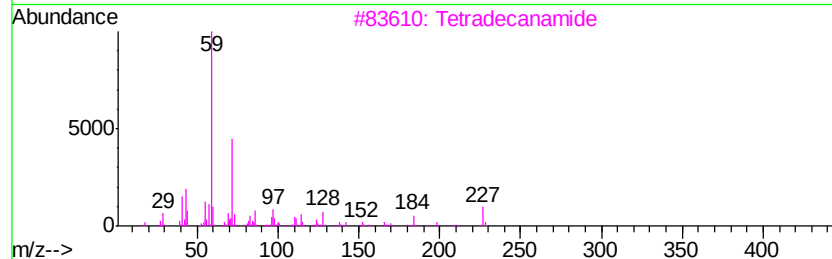
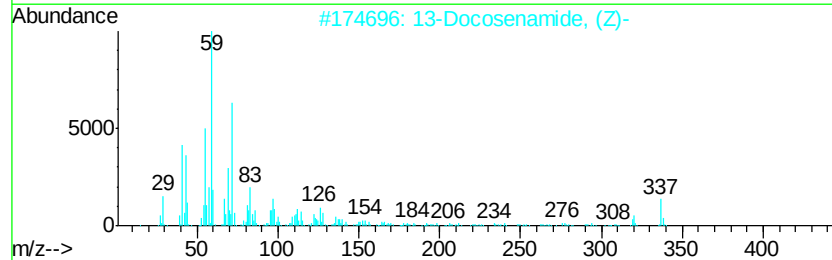
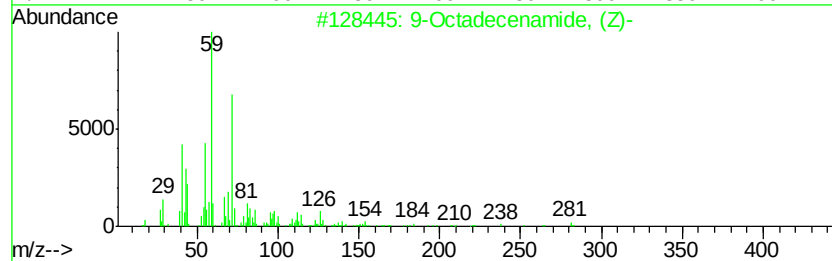
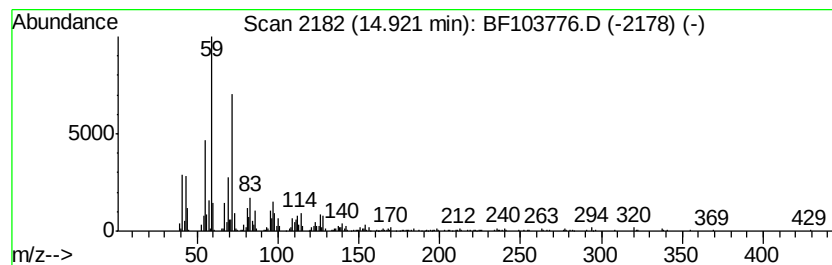
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 9-Octadecenamide, (Z)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.92	12.56 ng	535580	Chrysene-d12	14.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	91
2		13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	90
3		Tetradecanamide	227	C14H29NO	000638-58-4	86
4		Decanamide-	171	C10H21NO	002319-29-1	59
5		Dodecanamide	199	C12H25NO	001120-16-7	59



Data Path : Z:\HPCHEM1\BNA F\DATA\BF031918\
Data File : BF103776.D
Acq On : 20 Mar 2018 00:32
Operator : JU/SJ
Sample : J1971-07
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WEST-GARAGE-6-ST-LATERAL-NW

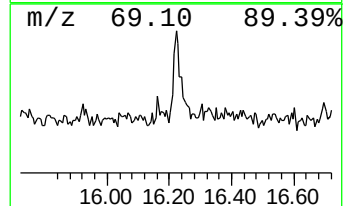
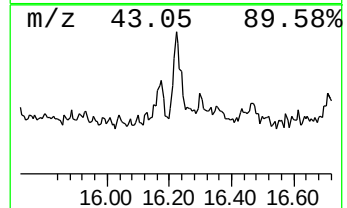
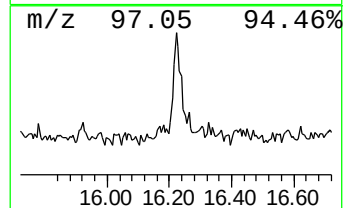
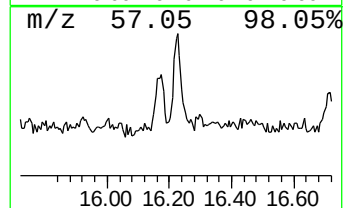
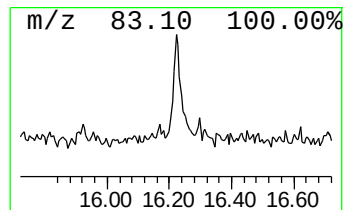
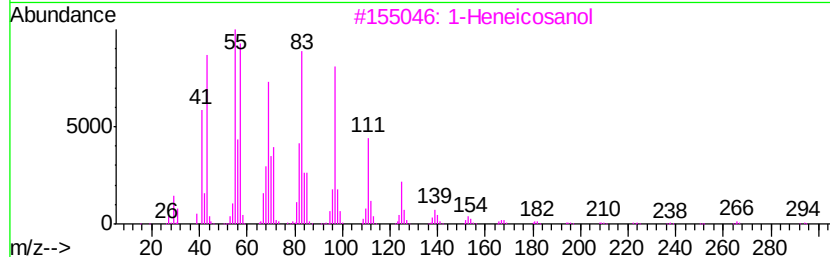
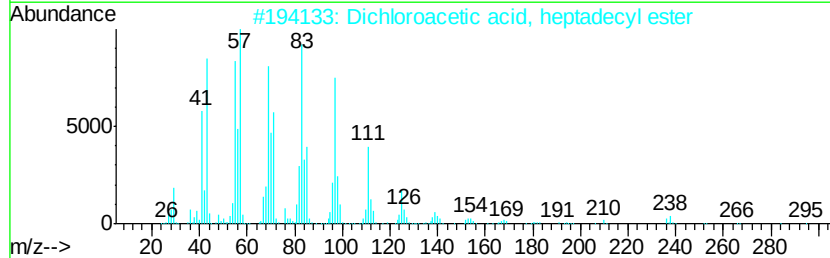
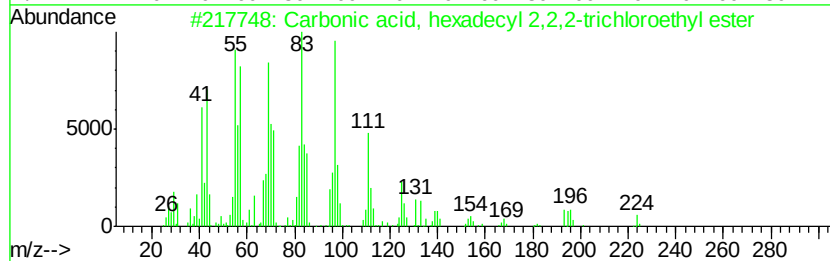
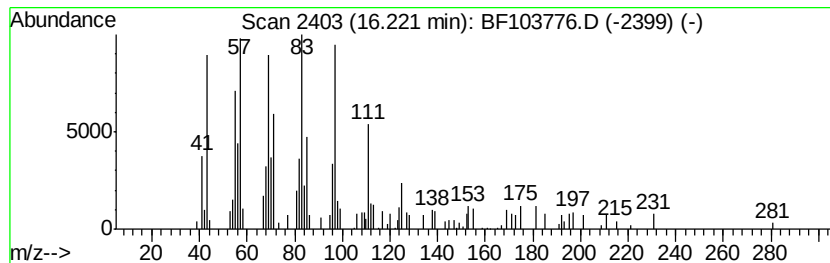
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Carbonic acid, hexadecyl 2,... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.22	2.65 ng	99873	Perylene-d12	15.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Carbonic acid, hexadecyl 2,2,2-t...	416	C19H35Cl3O3	1000314-56-2	91
2		Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	90
3		1-Heneicosanol	312	C21H44O	015594-90-8	87
4		Behenic alcohol	326	C22H46O	000661-19-8	70
5		Octatriacontyl pentafluoropropio...	697	C41H77F5O2	1000351-89-1	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031918\
Data File : BF103776.D
Acq On : 20 Mar 2018 00:32
Operator : JU/SJ
Sample : J1971-07
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WEST-GARAGE-6-ST-LATERAL-NW

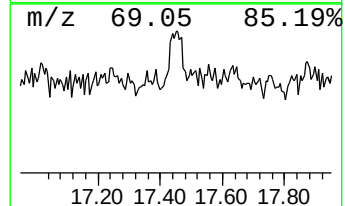
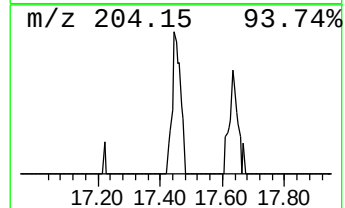
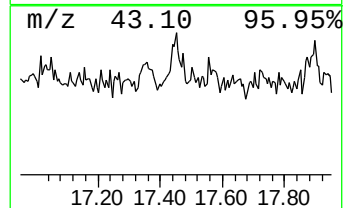
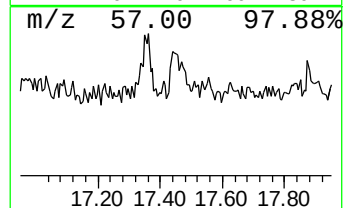
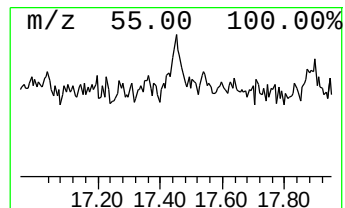
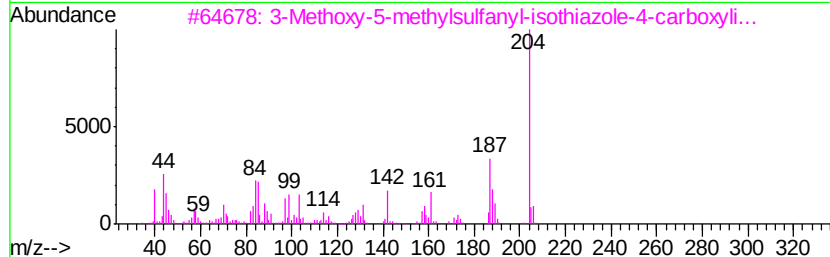
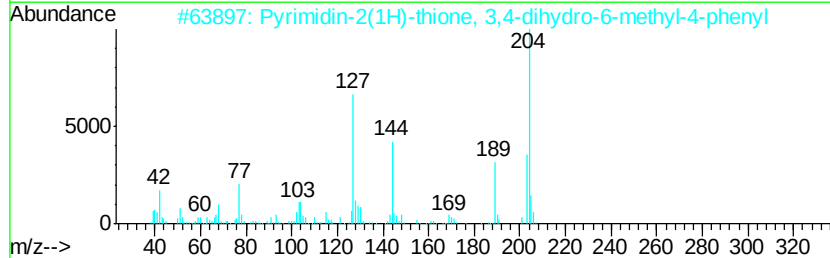
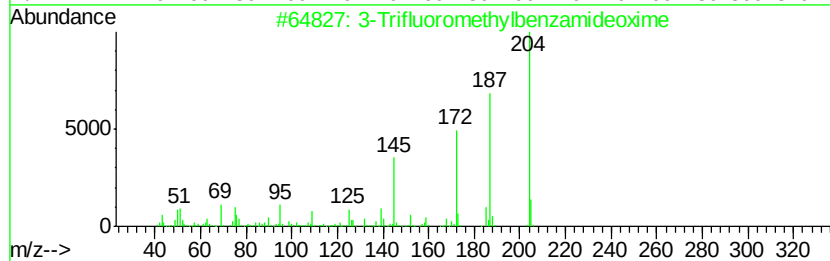
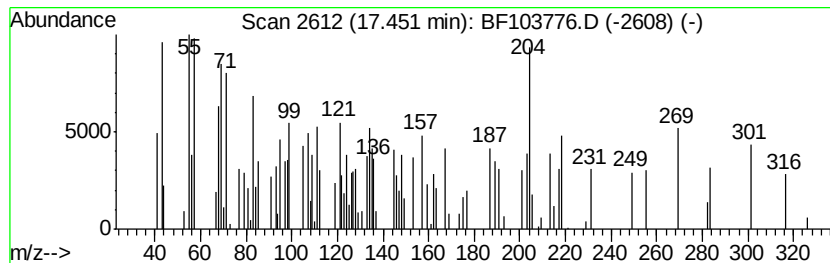
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown17.45 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.45	2.13 ng	80164	Perylene-d12	15.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Trifluoromethylbenzamidoxime	204	C8H7F3N2O	040067-80-9	18
2		Pyrimidin-2(1H)-thione, 3,4-dihy...	204	C11H12N2S	030570-19-5	14
3		3-Methoxy-5-methylsulfanyl-isoth...	204	C6H8N2O2S2	1000274-69-9	14
4		4-Trifluoromethylbenzamidoxime	204	C8H7F3N2O	022179-86-8	14
5		Naphthalene, 1,2,3,5,6,7,8,8a-oc...	204	C15H24	004630-07-3	11



Data Path : Z:\HPCHEM1\BNA F\DATA\BF031918\
Data File : BF103776.D
Acq On : 20 Mar 2018 00:32
Operator : JU/SJ
Sample : J1971-07
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WEST-GARAGE-6-ST-LATERAL-NW

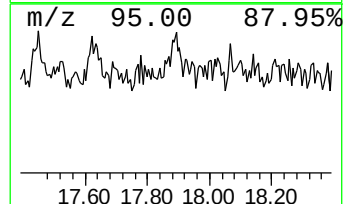
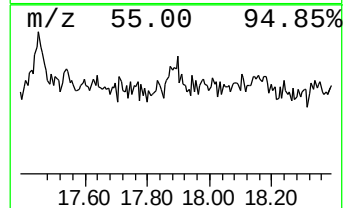
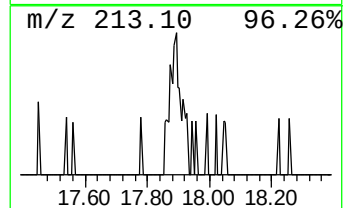
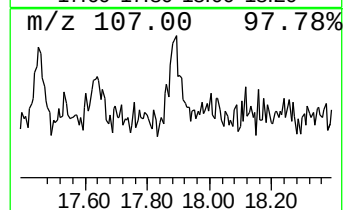
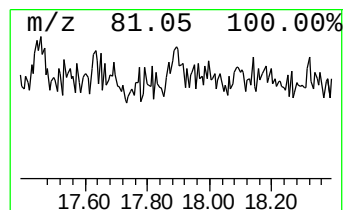
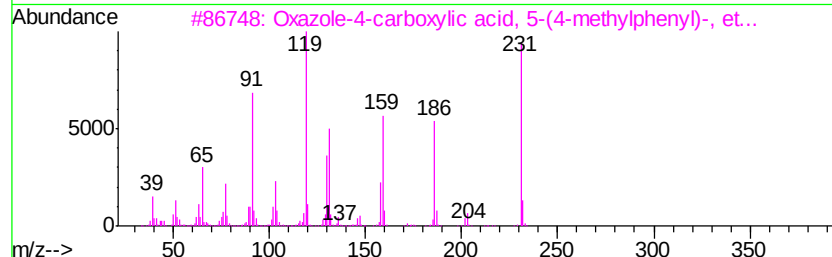
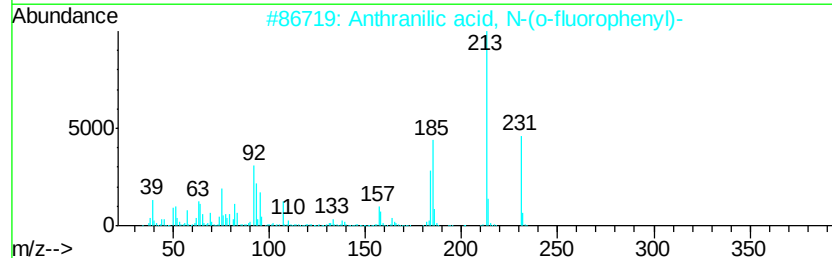
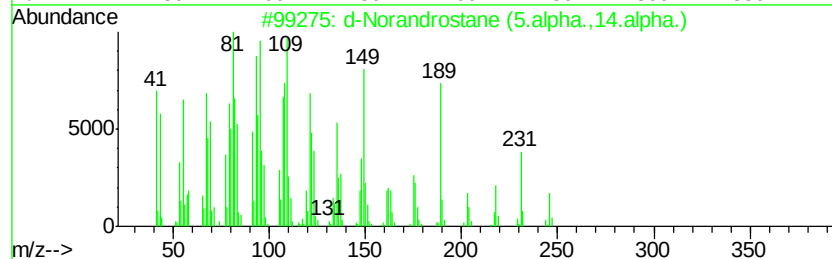
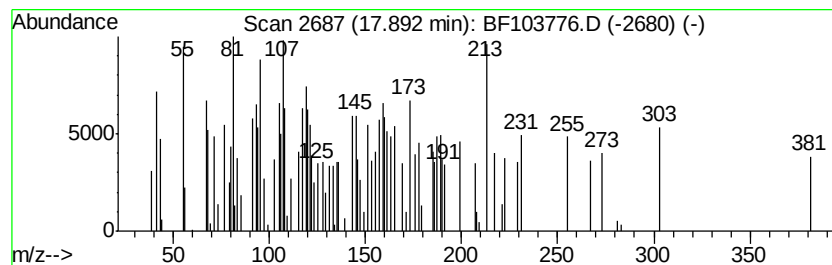
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown17.89 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.89	2.10 ng	79076	Perylene-d12	15.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	d-Norandrostande (5.alpha.,14.alp...	246	C18H30	1000281-13-8	11
2		Anthranilic acid, N-(o-fluorophe...	231	C13H10FN02	000054-58-0	10
3		Oxazole-4-carboxylic acid, 5-(4-...	231	C13H13N03	1000261-06-2	10
4		4,4,6-Trimethyl-2-(N-methylpheny...	232	C14H20N2O	053004-36-7	10
5		Andrographolide	350	C20H30O5	005508-58-7	10



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031918\
Data File : BF103776.D
Acq On : 20 Mar 2018 00:32
Operator : JU/SJ
Sample : J1971-07
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.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...	2.33	5.0	ng	208712	1	6.97	830138	20.0
2-Pentanone, 4-hy...	5.23	5.1	ng	210051	1	6.97	830138	20.0
unknown6.75	6.75	90.7	ng	3762790	1	6.97	830138	20.0
Octacosane	10.90	2.3	ng	98307	4	11.50	875342	20.0
n-Hexadecanoic acid	12.02	3.5	ng	154414	4	11.50	875342	20.0
5-Eicosene, (E)-	13.97	9.6	ng	408341	5	14.16	852934	20.0
Heptadecyl heptaf...	14.62	2.2	ng	94572	5	14.16	852934	20.0
9-Octadecenamide,...	14.92	12.6	ng	535580	5	14.16	852934	20.0
Carbonic acid, he...	16.22	2.6	ng	99873	6	15.70	754373	20.0
unknown17.45	17.45	2.1	ng	80164	6	15.70	754373	20.0
unknown17.89	17.89	2.1	ng	79076	6	15.70	754373	20.0