

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050417\
 Data File : BF094884.D
 Acq On : 4 May 2017 20:38
 Operator : SJ/MA
 Sample : I2916-04
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-06-7.0-8.0

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-BF050217.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	5.025	520	525	541	rBV	272052	543591	21.53%	2.319%
2	5.666	629	634	662	rBV	1579124	2302312	91.19%	9.820%
3	6.242	725	732	740	rBV	20752	37491	1.49%	0.160%
4	7.266	900	906	921	rBV	1703690	2333966	92.45%	9.955%
5	7.383	921	926	943	rVB	1951369	2524628	100.00%	10.768%
6	7.719	979	983	992	rVB	392229	456220	18.07%	1.946%
7	7.960	1019	1024	1033	rBV	1450674	1941927	76.92%	8.283%
8	8.436	1101	1105	1110	rBV	28286	30258	1.20%	0.129%
9	8.619	1130	1136	1156	rBV	1175408	1472208	58.31%	6.279%
10	9.736	1322	1326	1340	rBV	465076	610666	24.19%	2.605%
11	11.513	1623	1628	1645	rBV	1745847	2053436	81.34%	8.758%
12	11.807	1673	1678	1683	rBV	18290	26983	1.07%	0.115%
13	12.207	1741	1746	1753	rBV	214499	264575	10.48%	1.128%
14	12.571	1803	1808	1821	rBV2	520801	722174	28.61%	3.080%
15	13.412	1947	1951	1957	rVB	28319	33584	1.33%	0.143%
16	13.865	2022	2028	2040	rBV	1023700	1423149	56.37%	6.070%
17	14.189	2076	2083	2085	rBV	32805	41102	1.63%	0.175%
18	14.918	2197	2207	2211	rBV3	27775	49886	1.98%	0.213%
19	14.965	2211	2215	2226	rVB	574597	784059	31.06%	3.344%
20	15.183	2247	2252	2264	rBV5	56966	113448	4.49%	0.484%
21	15.659	2326	2333	2351	rBV3	83550	199785	7.91%	0.852%
22	15.812	2355	2359	2367	rVV3	26552	37598	1.49%	0.160%
23	15.942	2376	2381	2391	rBV	208099	277860	11.01%	1.185%
24	17.024	2561	2565	2576	rBV3	39818	62933	2.49%	0.268%
25	17.295	2606	2611	2615	rBV	1596094	1758522	69.65%	7.500%
26	17.659	2668	2673	2676	rBV3	21255	38205	1.51%	0.163%
27	18.542	2818	2823	2834	rBV	235208	469828	18.61%	2.004%
28	18.636	2834	2839	2846	rVV	593422	739822	29.30%	3.155%
29	18.700	2846	2850	2856	rVB	36368	44256	1.75%	0.189%
30	18.953	2886	2893	2901	rBV4	23174	46689	1.85%	0.199%
31	19.342	2952	2959	2967	rBV3	217935	465192	18.43%	1.984%
32	19.512	2985	2988	2992	rVB2	30552	34262	1.36%	0.146%
33	19.659	3007	3013	3017	rBV4	38280	53569	2.12%	0.228%
34	19.771	3029	3032	3036	rVB	94397	93694	3.71%	0.400%

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

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
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35	20.053	3076	3080	3083	rBV	120784	139336	5.52%	0.594%
36	20.089	3083	3086	3090	rVV2	58347	87518	3.47%	0.373%
37	20.294	3117	3121	3132	rVB	424433	568099	22.50%	2.423%
38	20.400	3135	3139	3143	rBV5	25506	34440	1.36%	0.147%
39	20.777	3199	3203	3208	rBV	72313	91405	3.62%	0.390%
40	20.830	3208	3212	3217	rBV4	33327	72904	2.89%	0.311%
41	21.700	3355	3360	3363	rBV6	13538	25861	1.02%	0.110%
42	21.835	3373	3383	3396	rVB8	28641	102984	4.08%	0.439%
43	22.100	3423	3428	3433	rVB6	31866	56267	2.23%	0.240%
44	22.324	3460	3466	3480	rVB6	49261	137206	5.43%	0.585%
45	22.912	3560	3566	3573	rBV9	17269	41767	1.65%	0.178%

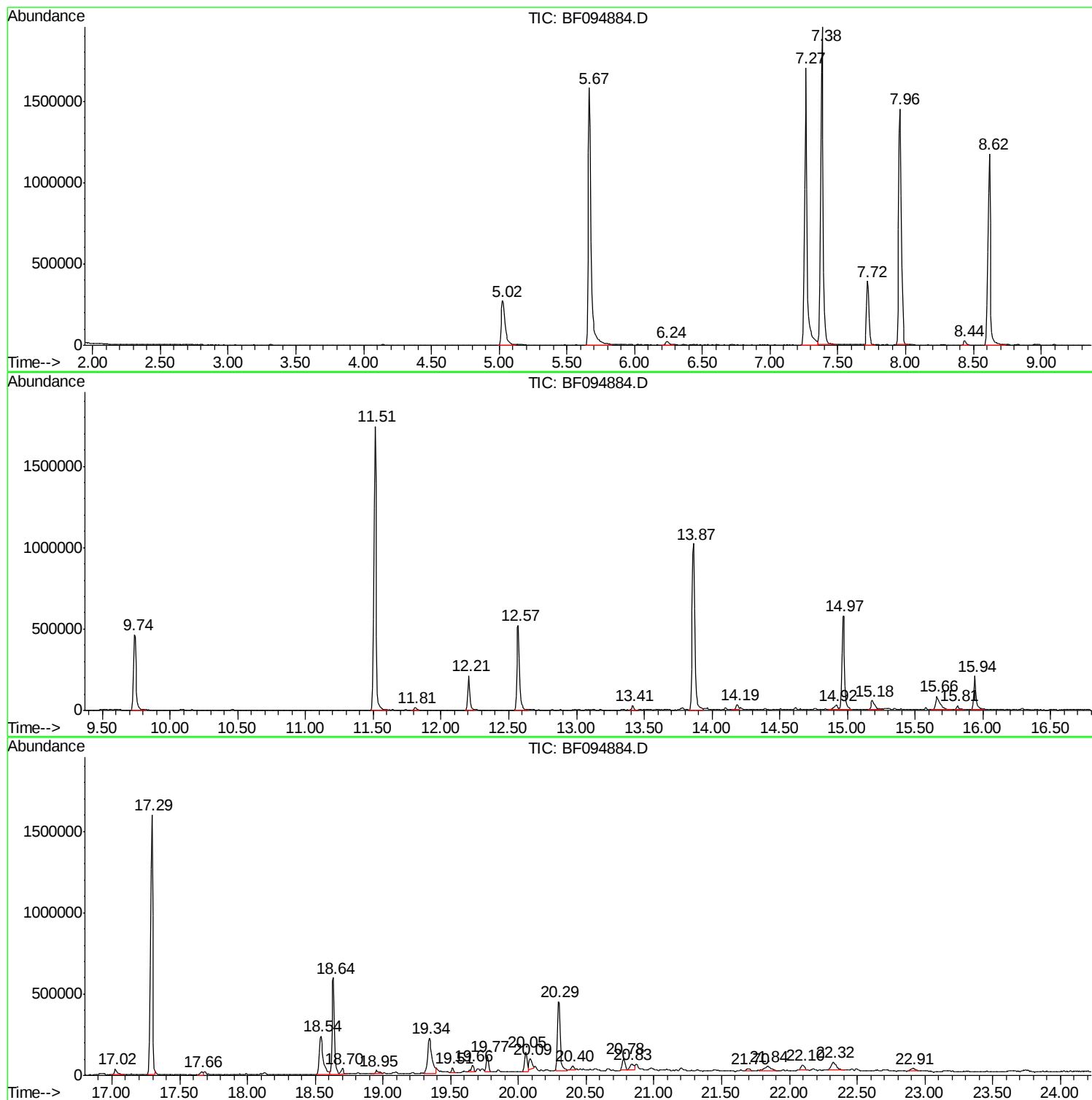
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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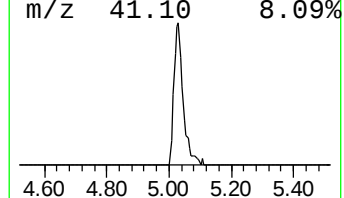
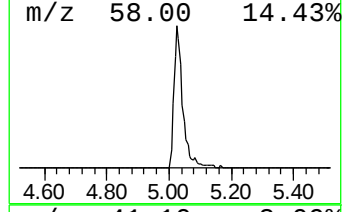
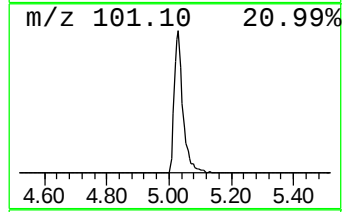
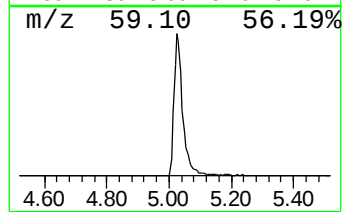
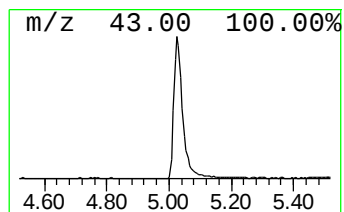
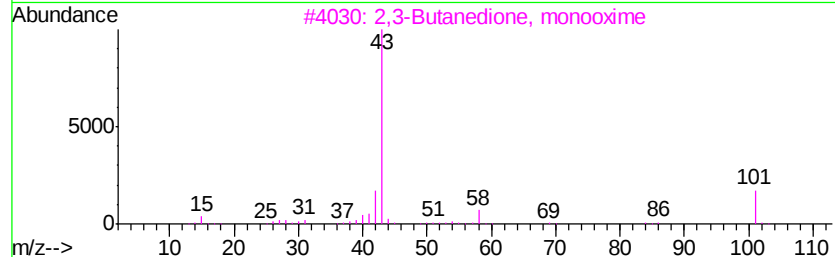
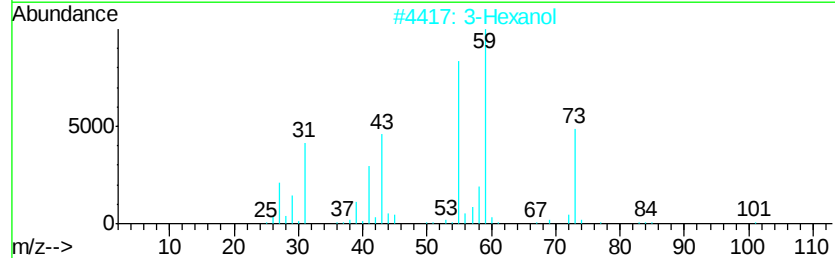
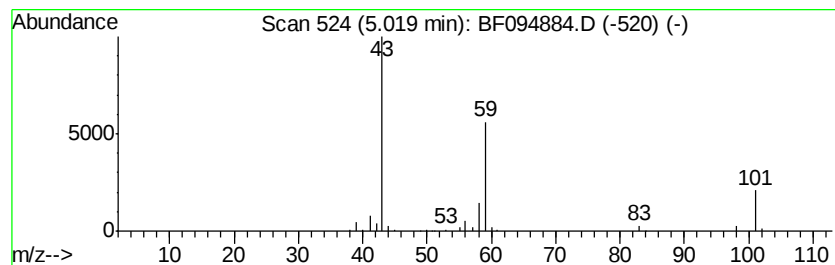
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.02	23.83 ng	543591	1,4-Dichlorobenzene-d4	7.72

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2		3-Hexanol	102	C6H14O	000623-37-0	28
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
4		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
5		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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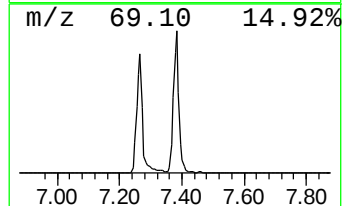
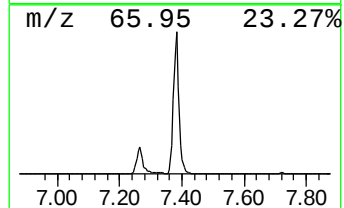
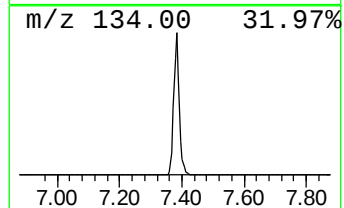
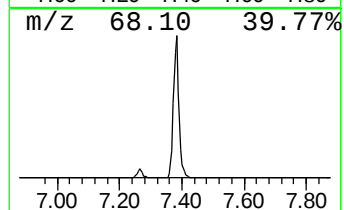
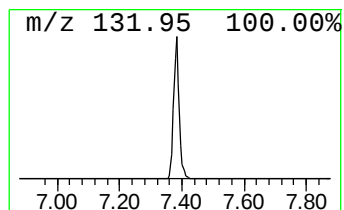
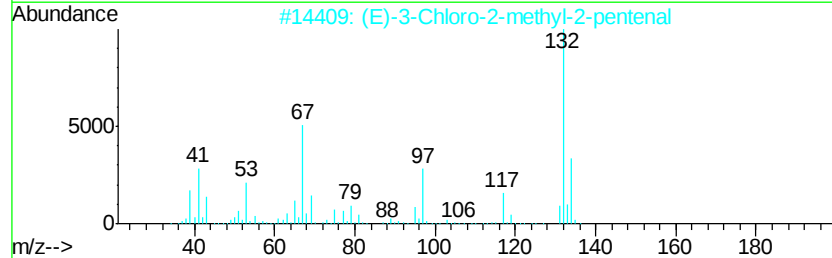
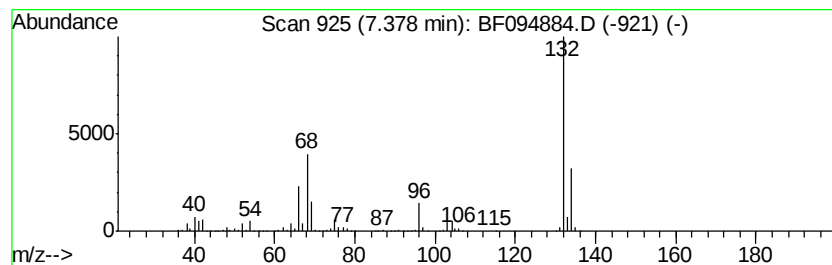
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown7.38 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.38	110.68 ng	2524630	1,4-Dichlorobenzene-d4	7.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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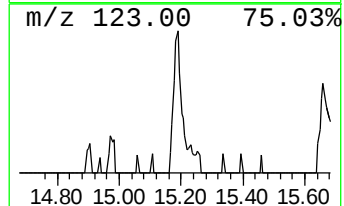
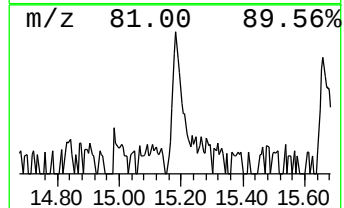
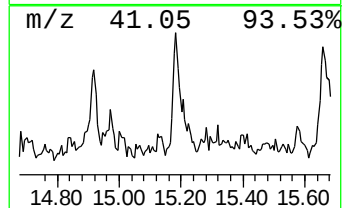
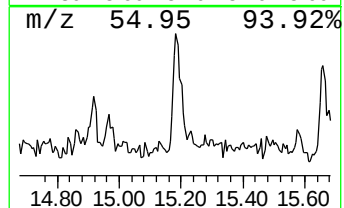
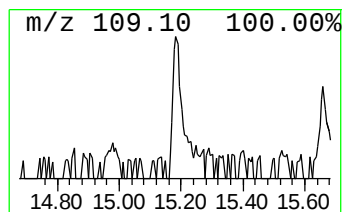
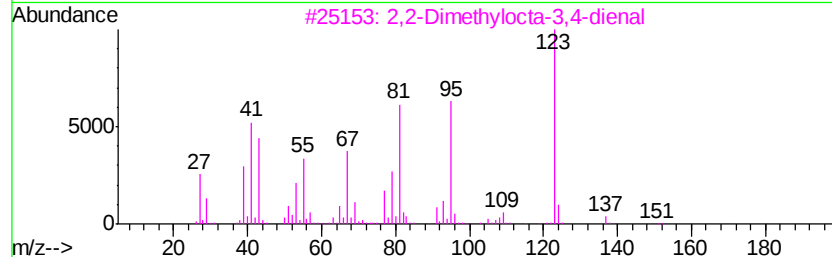
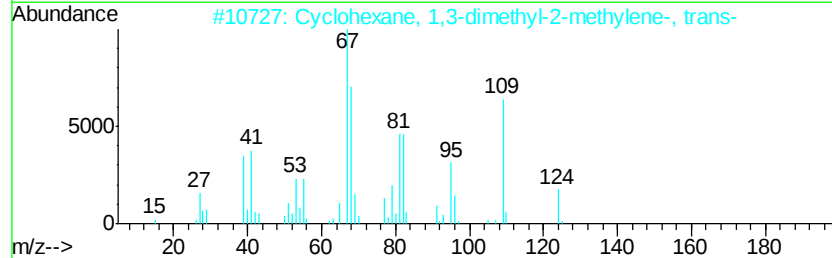
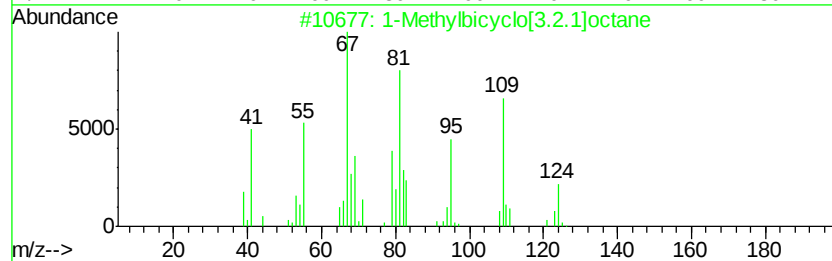
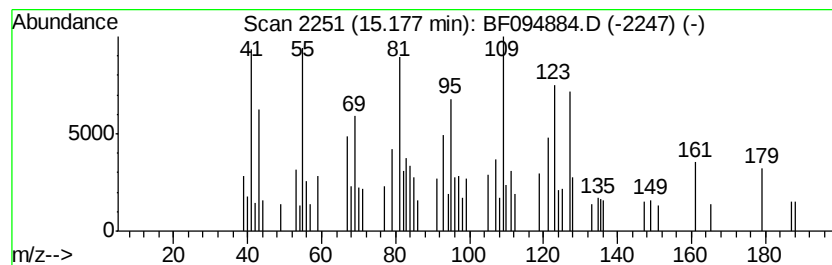
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 unknown15.18 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.18	2.89 ng	113448	Phenanthrene-d10	14.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Methylbicyclo[3.2.1]octane	124	C9H16	119972-41-7	38
2		Cyclohexane, 1,3-dimethyl-2-meth...	124	C9H16	020348-74-7	27
3		2,2-Dimethylocta-3,4-dienal	152	C10H16O	000590-71-6	27
4		4-Acetylocta-1,2-diene	152	C10H16O	1000250-44-5	27
5		Cyclopentane, 1-methyl-1-(2-meth...	138	C10H18	074764-47-9	27



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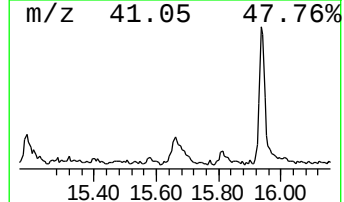
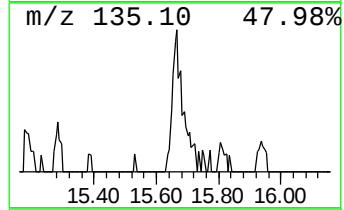
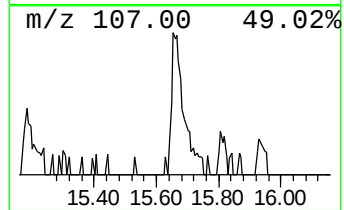
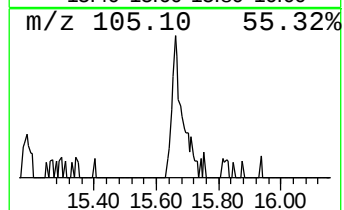
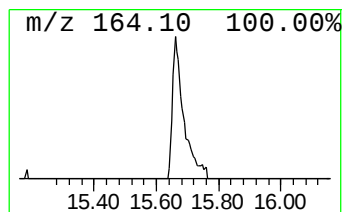
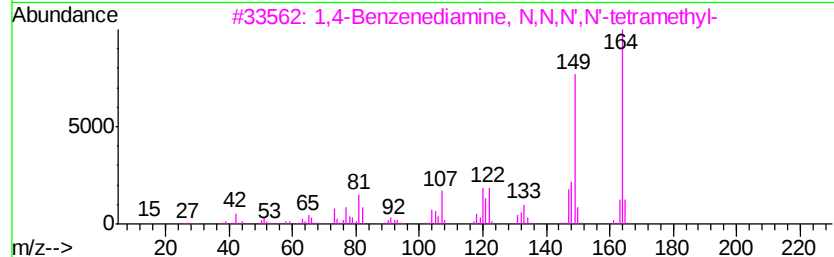
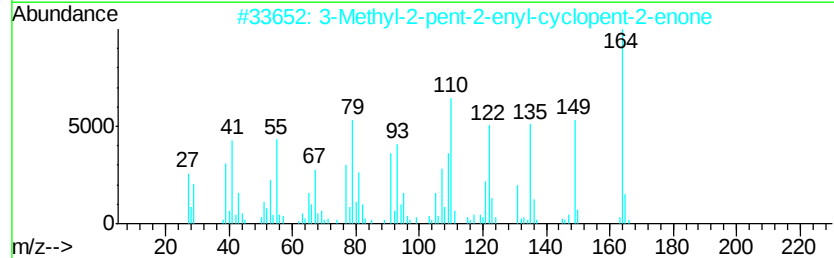
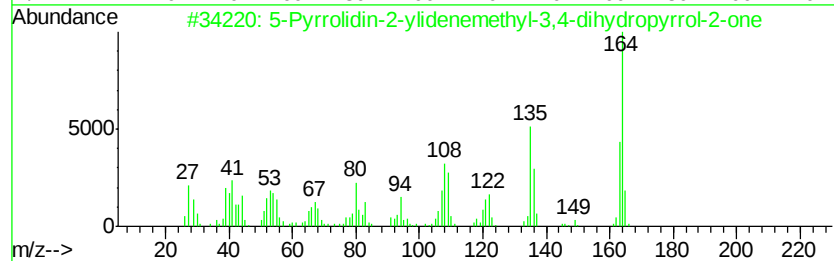
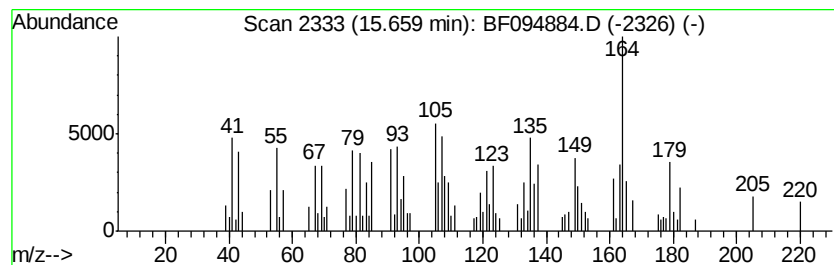
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TIC Integration Parameters: LSCINT.P

Peak Number 5 5-Pyrrolidin-2-ylidenemethy... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.66	5.10 ng	199785	Phenanthrene-d10	14.97

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		5-Pyrrolidin-2-ylidenemethyl-3,4...	164	C9H12N2O	1000192-09-5	55
2		3-Methyl-2-pent-2-enyl-cyclopent...	164	C11H16O	1000193-43-6	43
3		1,4-Benzenediamine, N,N,N',N'-te...	164	C10H16N2	000100-22-1	38
4		1,3-Benzenediamine, N,N,N',N'-te...	164	C10H16N2	022440-93-3	35
5		(1R,9aS,Z)-1-Methyl-3-(2-methylp...	207	C14H25N	1000367-15-3	27



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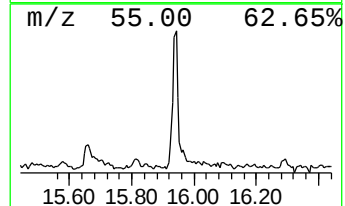
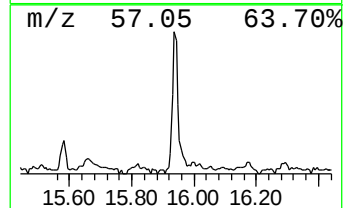
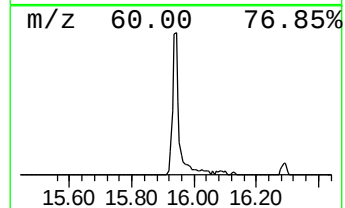
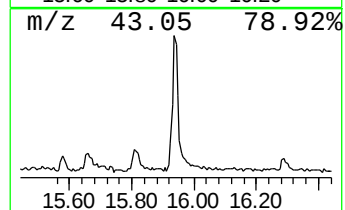
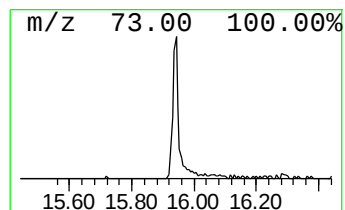
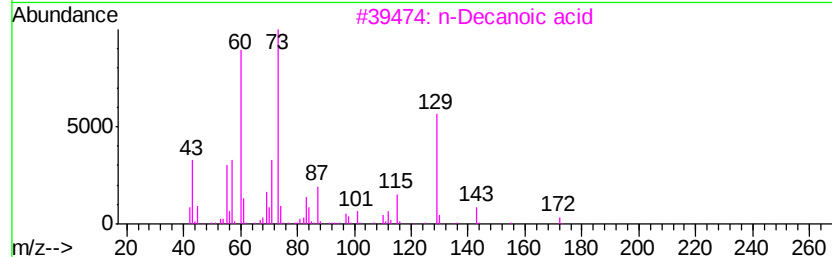
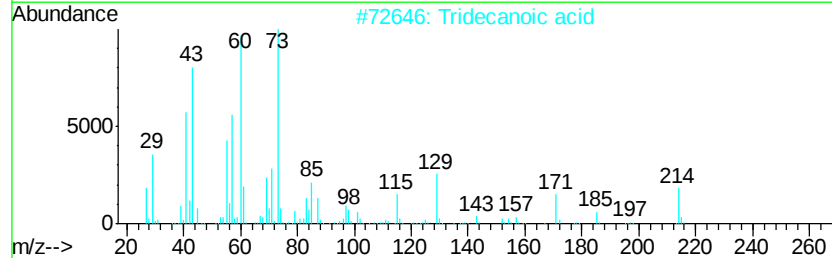
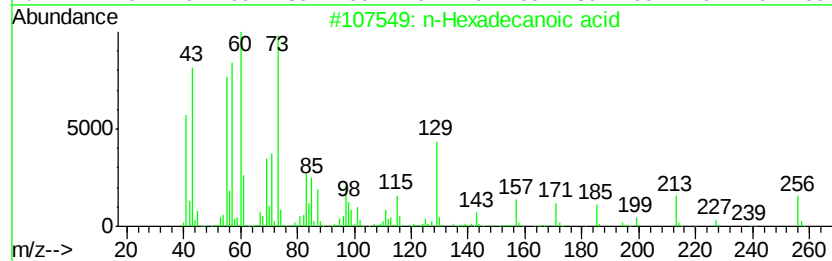
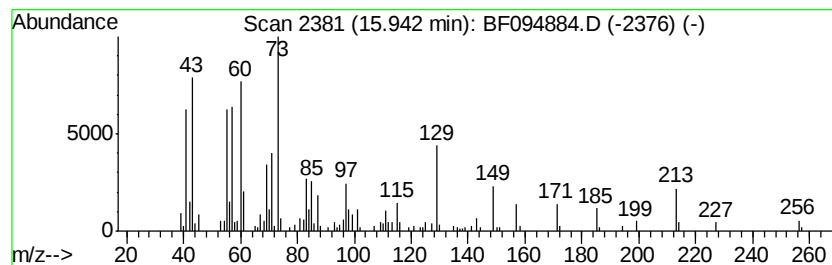
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Peak Number 6 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.94	7.09 ng	277860	Phenanthrene-d10	14.97

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tridecanoic acid	214	C13H26O2	000638-53-9	90
3		n-Decanoic acid	172	C10H20O2	000334-48-5	90
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	81
5		Oxalic acid, cyclobutyl hexadecy...	368	C22H40O4	1000309-70-6	18



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Data File : BF094884.D
Acq On : 4 May 2017 20:38
Operator : SJ/MA
Sample : I2916-04
Misc :
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ClientSampleId :
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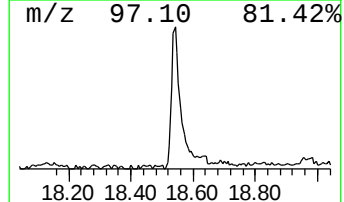
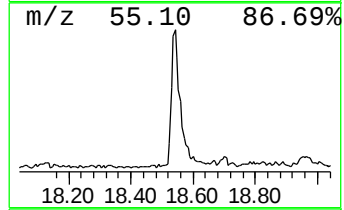
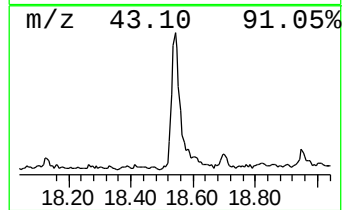
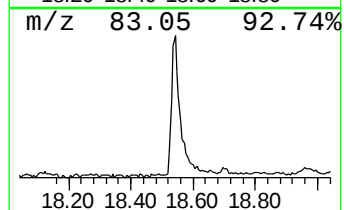
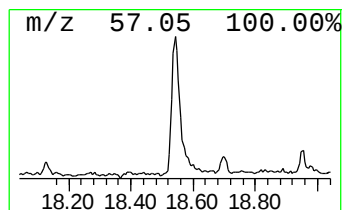
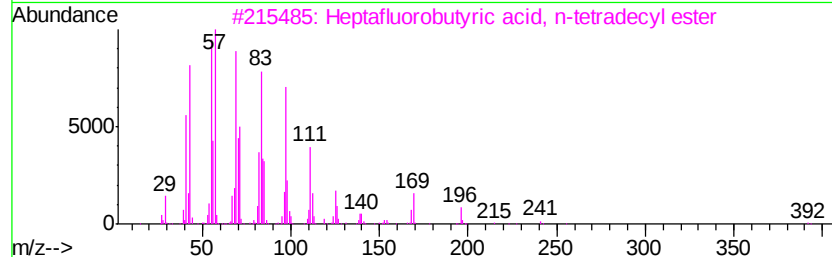
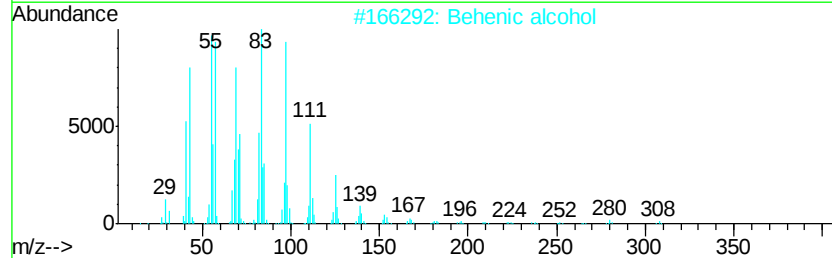
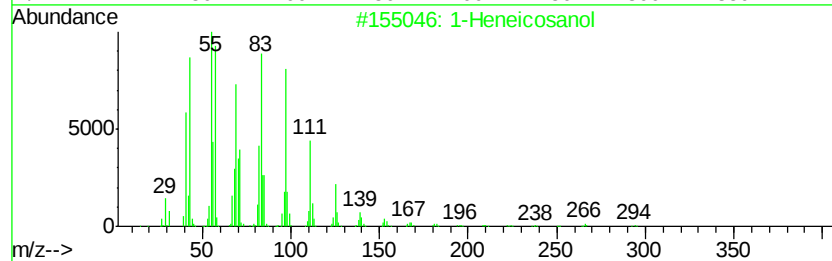
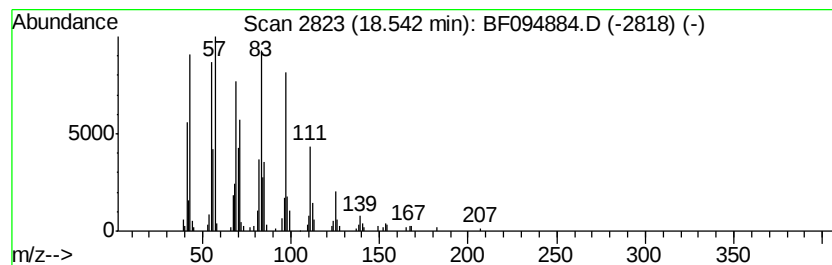
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 1-Heneicosanol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.54	12.70 ng	469828	Chrysene-d12	18.64

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosanol		312	C21H44O	015594-90-8	95
2	Behenic alcohol		326	C22H46O	000661-19-8	95
3	Heptafluorobutyric acid, n-tetra...		410	C18H29F7O2	007365-36-8	94
4	Trifluoroacetoxy hexadecane		338	C18H33F3O2	006222-03-3	94
5	n-Tetracosanol-1		354	C24H50O	000506-51-4	93



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050417\
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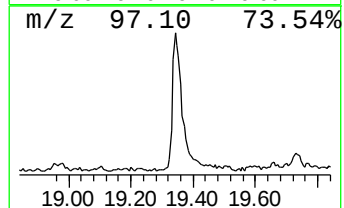
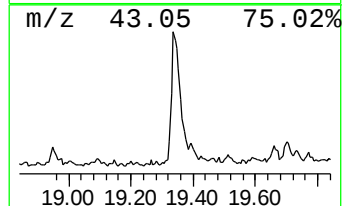
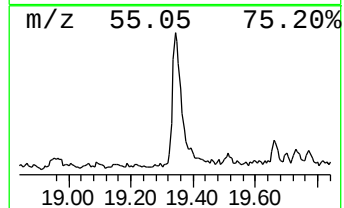
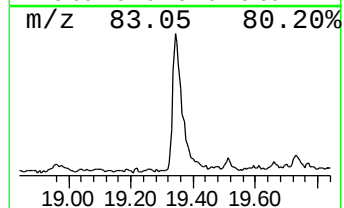
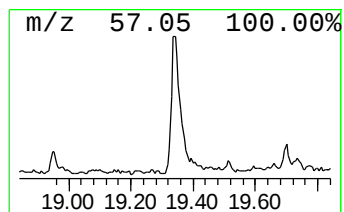
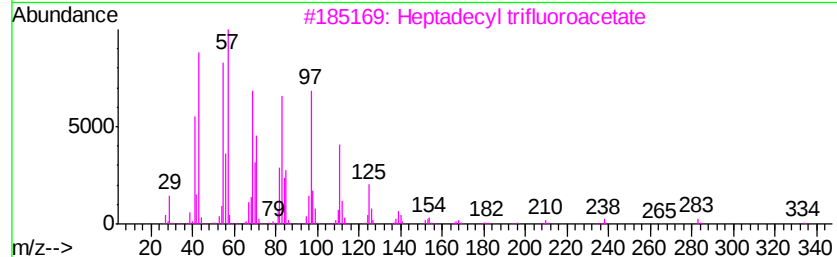
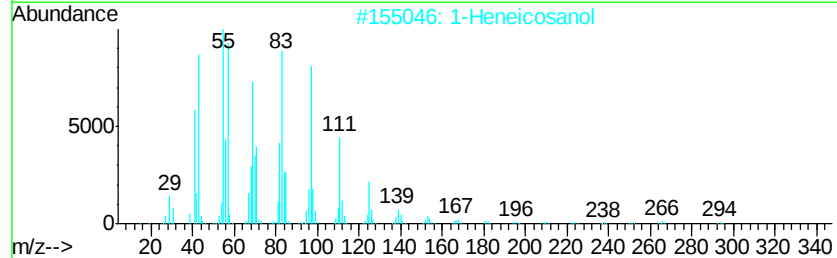
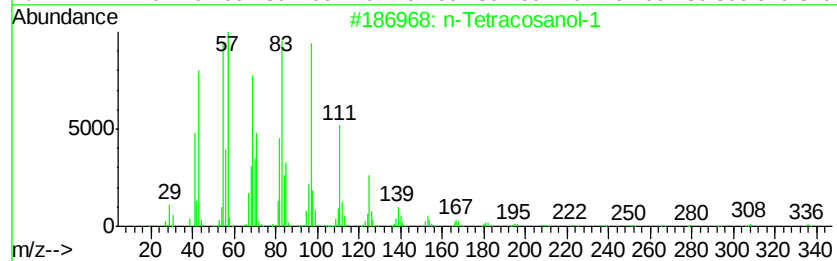
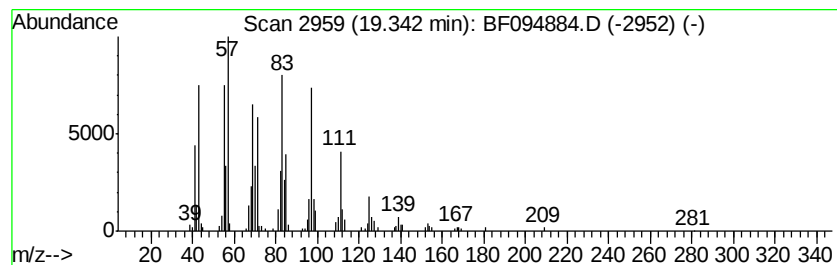
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 n-Tetracosanol-1 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.34	12.58 ng	465192	Chrysene-d12	18.64	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Tetracosanol-1	354	C24H50O	000506-51-4	93
2	1-Heneicosanol	312	C21H44O	015594-90-8	93
3	Heptadecyl trifluoroacetate	352	C19H35F3O2	1000351-87-0	93
4	Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	91
5	1-Docosene	308	C22H44	001599-67-3	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050417\
Data File : BF094884.D
Acq On : 4 May 2017 20:38
Operator : SJ/MA
Sample : I2916-04
Misc :
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Instrument :
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ClientSampleId :
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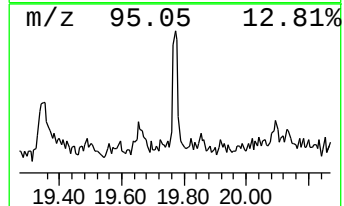
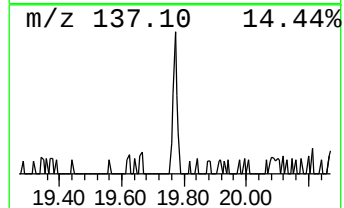
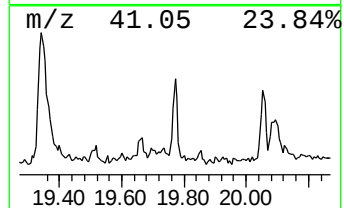
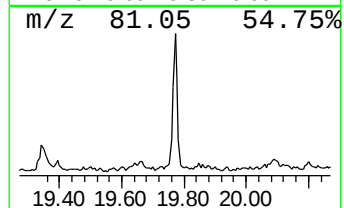
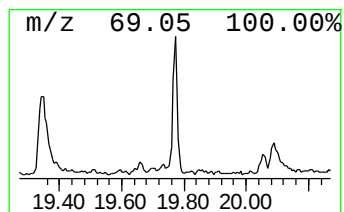
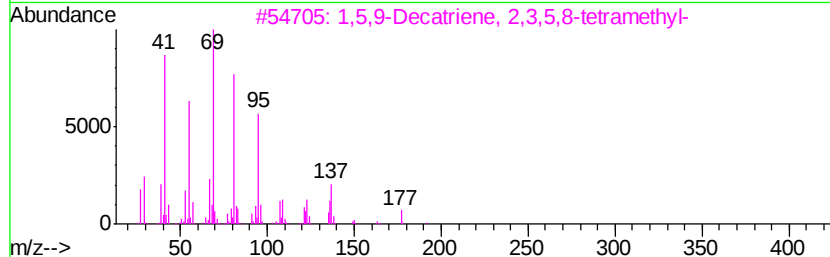
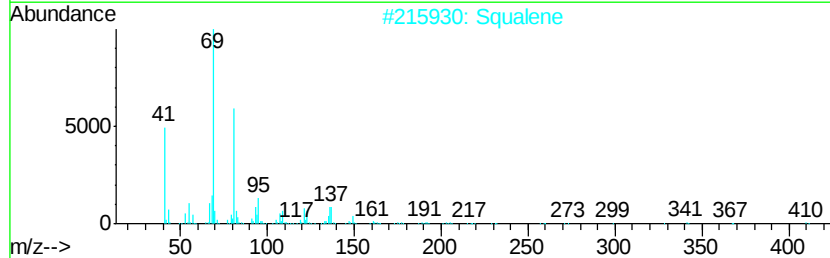
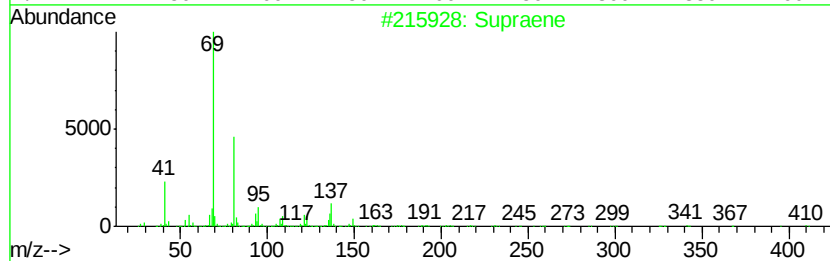
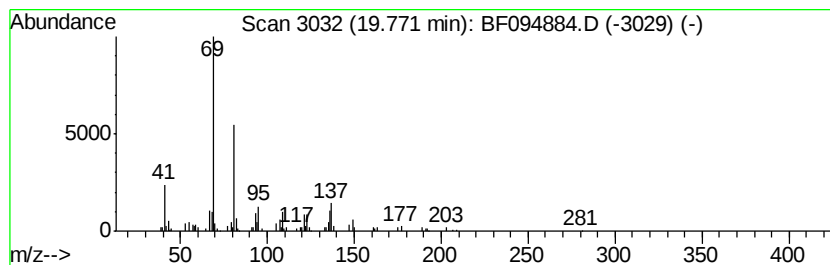
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Supraene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.77	3.30 ng	93694	Perylene-d12	20.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Supraene	410	C30H50	007683-64-9	91
2		Squalene	410	C30H50	000111-02-4	91
3		1,5,9-Decatriene, 2,3,5,8-tetram...	192	C14H24	230646-72-7	86
4		trans-Farnesol	222	C15H26O	000106-28-5	72
5		4,8,12-Tetradecatrienal, 5,9,13-...	248	C17H28O	066408-55-7	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050417\
Data File : BF094884.D
Acq On : 4 May 2017 20:38
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Misc :
ALS Vial : 12 Sample Multiplier: 1

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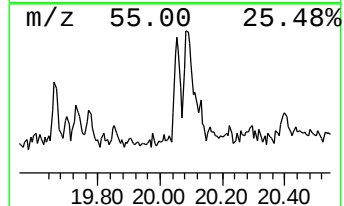
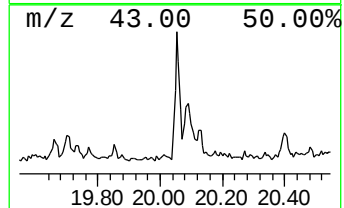
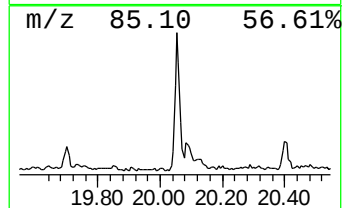
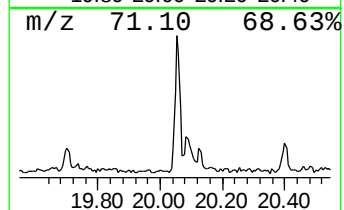
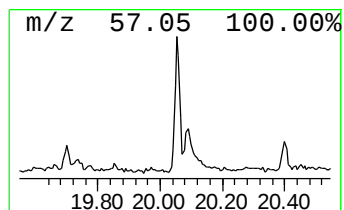
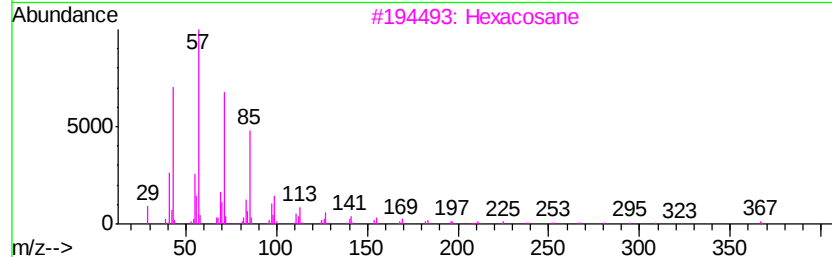
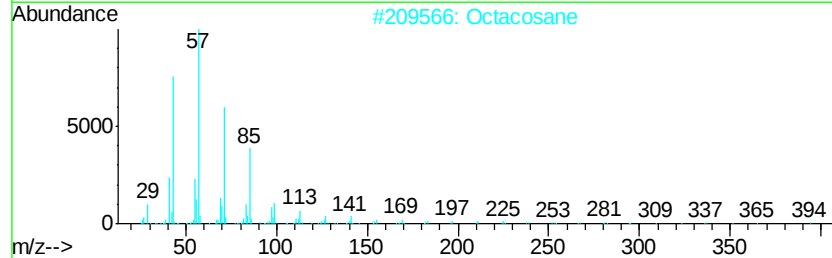
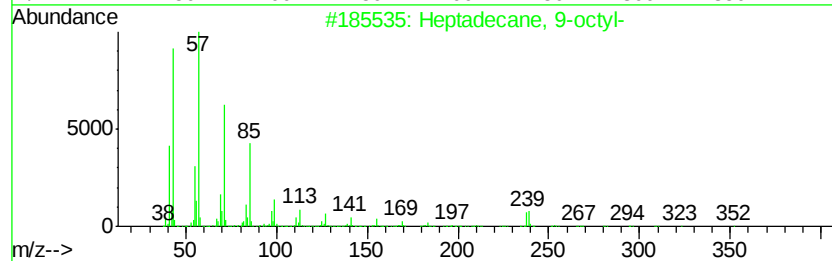
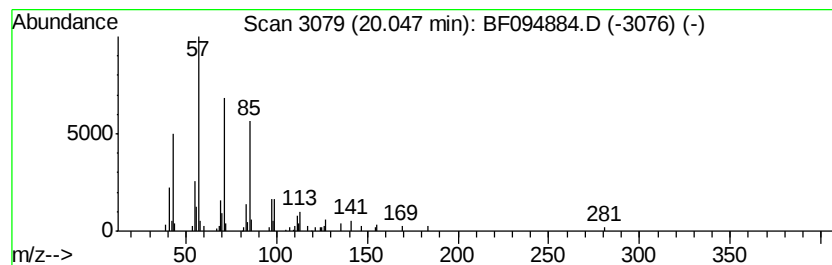
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Heptadecane, 9-octyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.05	4.91 ng	139336	Perylene-d12	20.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
2		Octacosane	394	C28H58	000630-02-4	91
3		Hexacosane	366	C26H54	000630-01-3	91
4		triacontane	422	C30H62	000638-68-6	91
5		2-methyloctacosane	408	C29H60	1000376-72-8	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050417\
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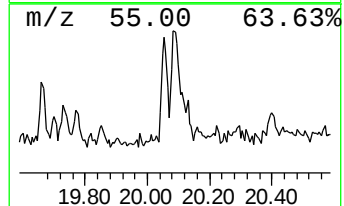
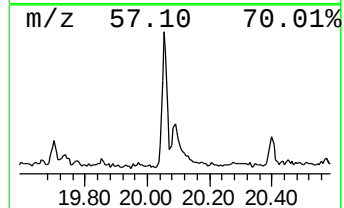
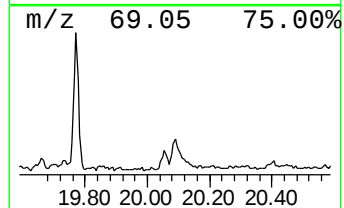
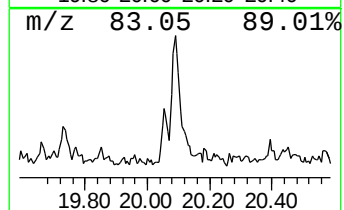
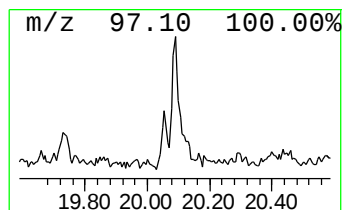
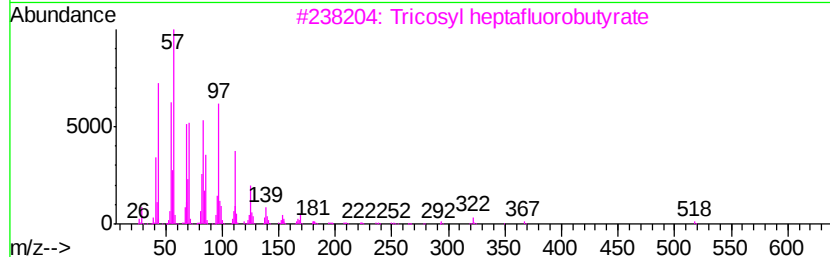
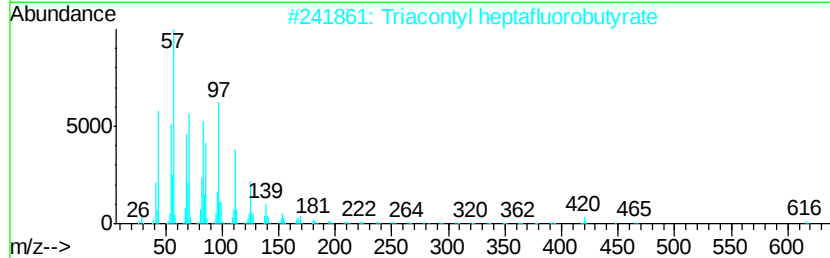
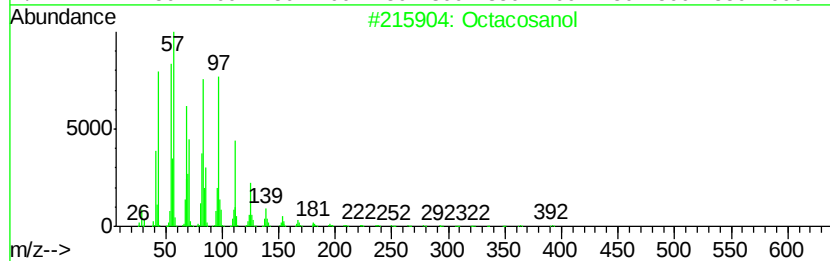
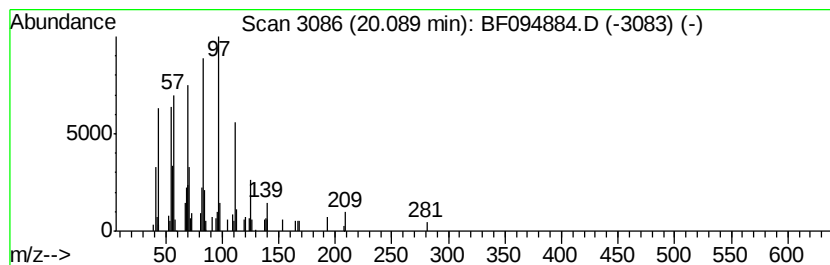
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Octacosanol Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.09	3.08 ng	87518	Perylene-d12	20.30
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Octacosanol	410 C28H58O	000557-61-9	62
2	Triacontyl heptafluorobutyrate	634 C34H61F7O2	1000351-83-2	58
3	Tricosyl heptafluorobutyrate	536 C27H47F7O2	1000351-83-4	58
4	Heptafluorobutyric acid, pentade...	424 C19H31F7O2	959261-23-5	53
5	Trifluoroacetic acid, pentadecyl...	324 C17H31F3O2	959010-23-2	53



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Operator : SJ/MA
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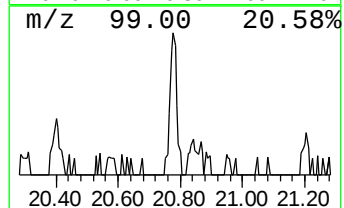
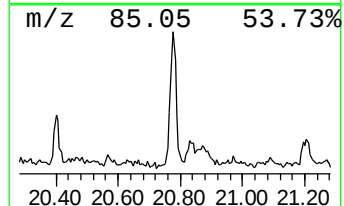
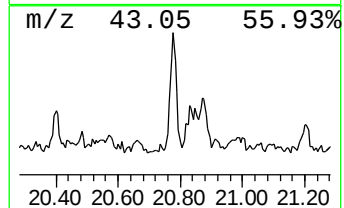
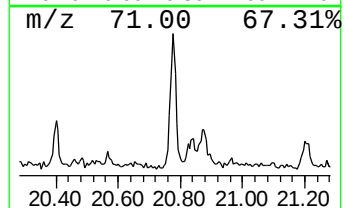
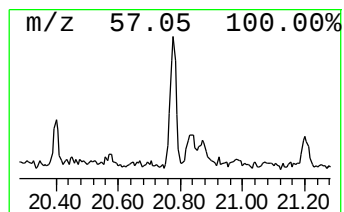
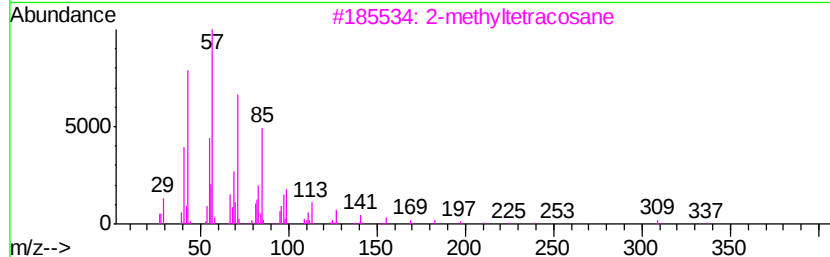
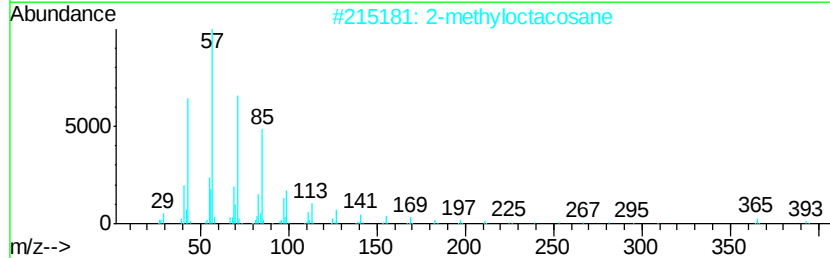
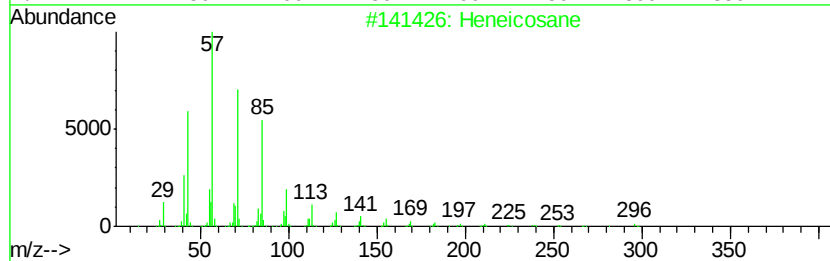
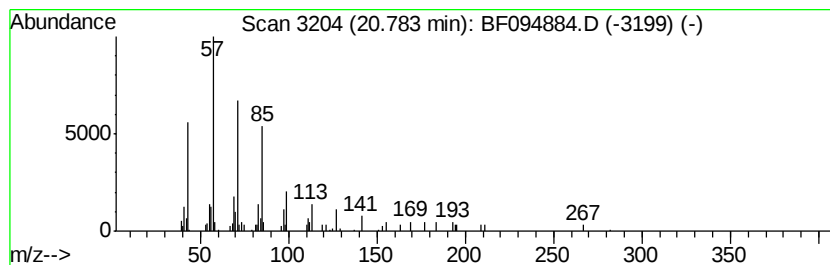
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Heneicosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.78	3.22 ng	91405	Perylene-d12	20.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		2-methyloctacosane	408	C29H60	1000376-72-8	91
3		2-methyltetracosane	352	C25H52	1000376-72-6	91
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	90
5		Heptadecane	240	C17H36	000629-78-7	90



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Acq On : 4 May 2017 20:38
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ALS Vial : 12 Sample Multiplier: 1

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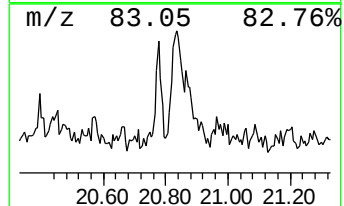
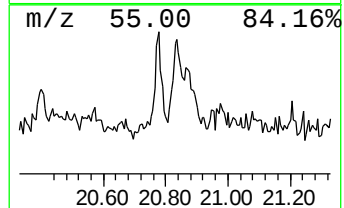
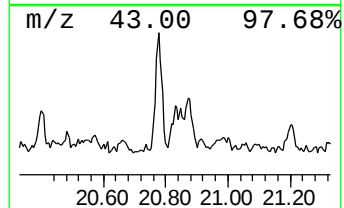
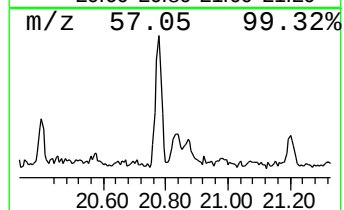
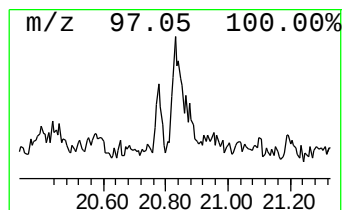
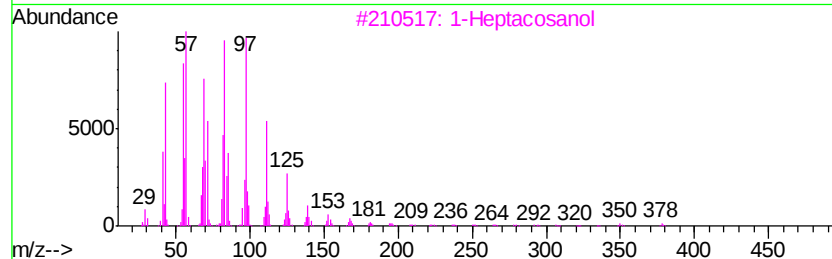
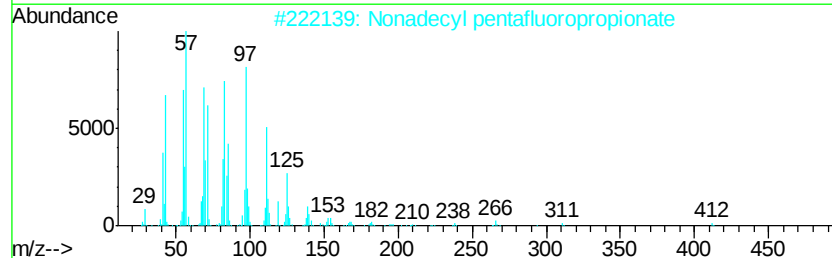
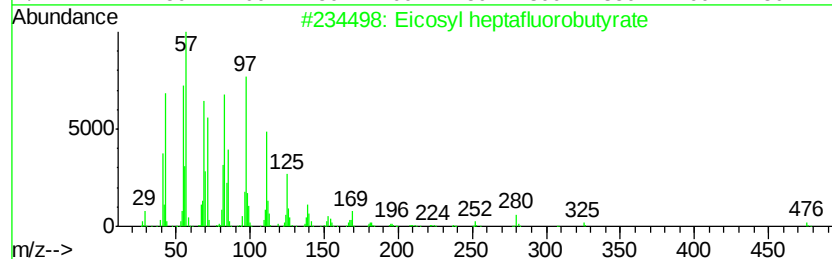
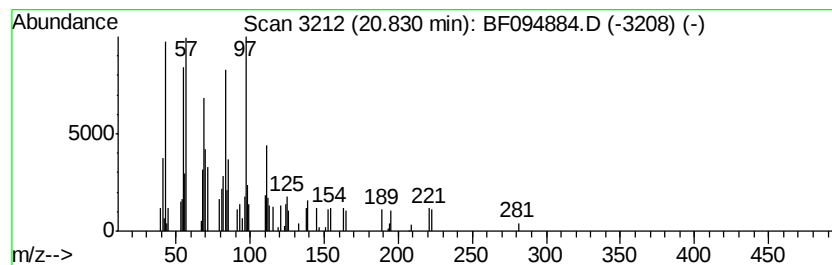
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Eicosyl heptafluorobutyrate Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.83	2.57 ng	72904	Perylene-d12	20.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosyl heptafluorobutyrate	494	C24H41F7O2	1000351-83-5	72
2		Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	72
3		1-Heptacosanol	396	C27H56O	002004-39-9	62
4		n-Tetracosanol-1	354	C24H50O	000506-51-4	62
5		Behenic alcohol	326	C22H46O	000661-19-8	58



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050417\
Data File : BF094884.D
Acq On : 4 May 2017 20:38
Operator : SJ/MA
Sample : I2916-04
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-06-7.0-8.0

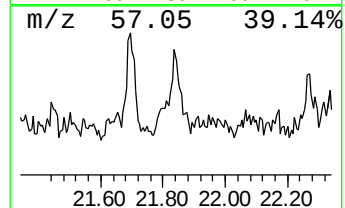
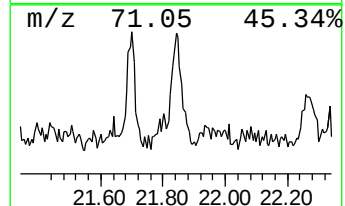
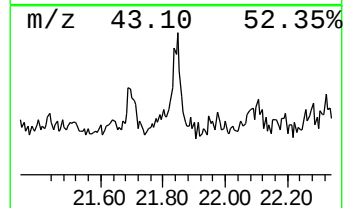
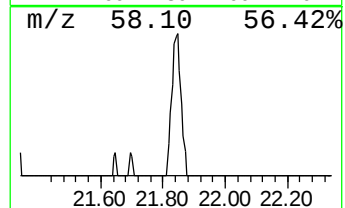
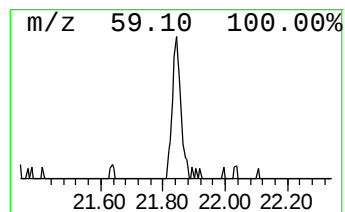
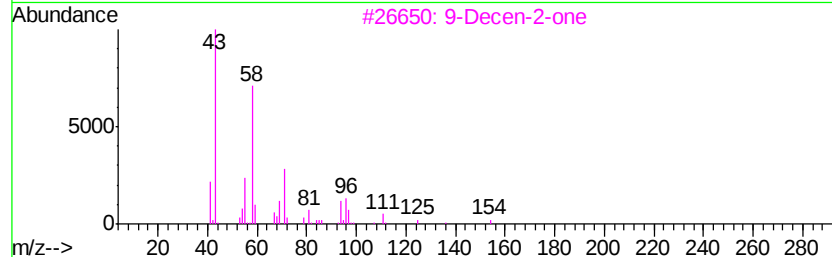
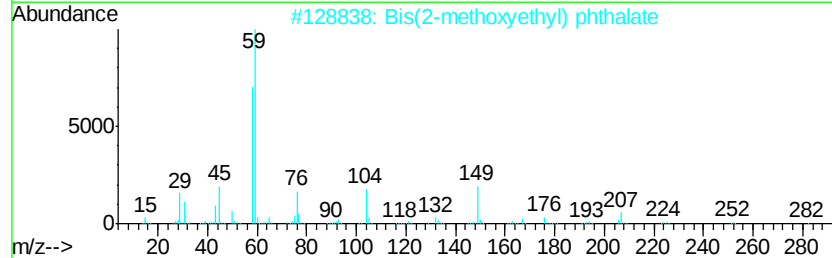
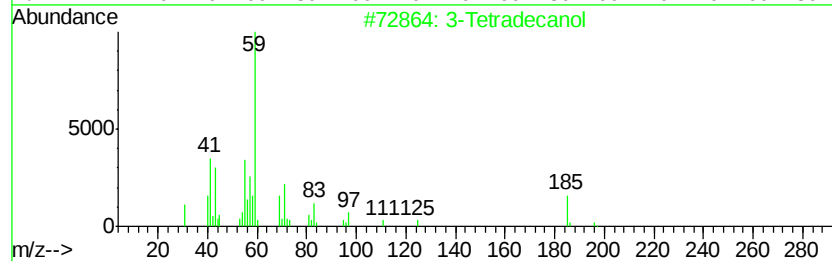
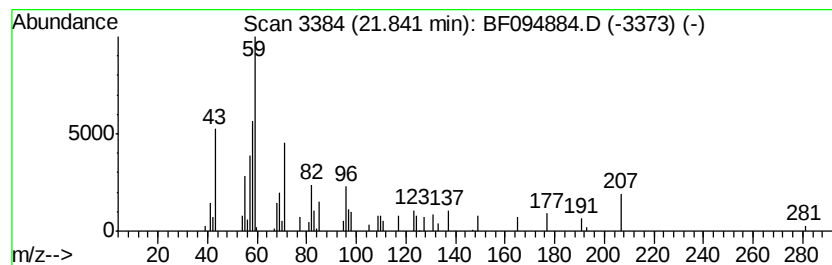
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown21.84 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.84	3.63 ng	102984	Perylene-d12	20.30

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Tetradecanol	214	C14H30O	001653-32-3	43
2			Bis(2-methoxyethyl) phthalate	282	C14H18O6	000117-82-8	37
3			9-Decen-2-one	154	C10H18O	035194-30-0	25
4			8-Nonen-2-one	140	C9H16O	005009-32-5	22
5			Menthane, 3-(2-methyl-2-hydroxy-...	212	C14H28O	1000160-06-0	22



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050417\
Data File : BF094884.D
Acq On : 4 May 2017 20:38
Operator : SJ/MA
Sample : I2916-04
Misc :
ALS Vial : 12 Sample Multiplier: 1

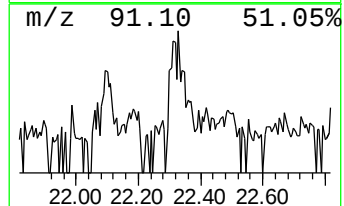
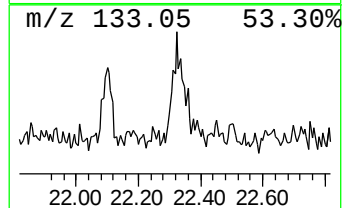
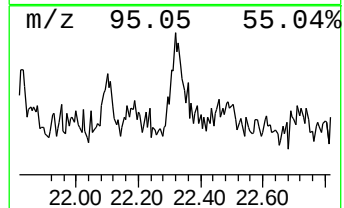
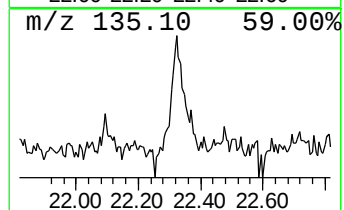
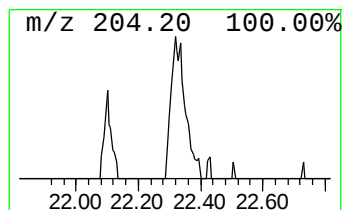
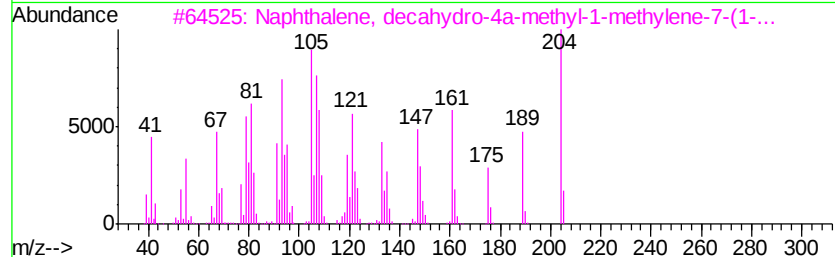
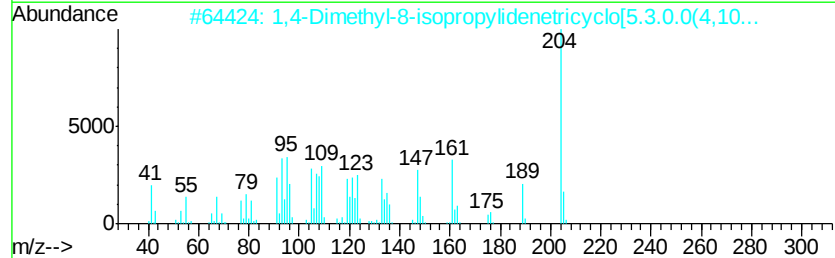
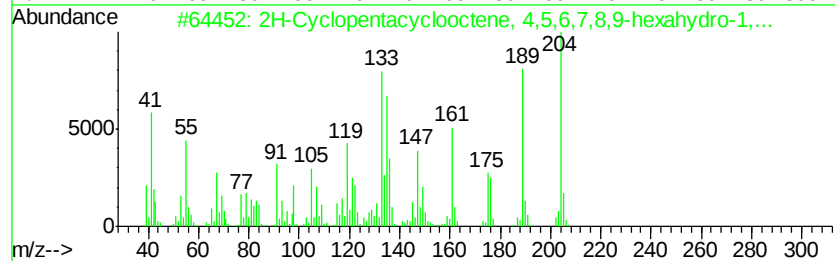
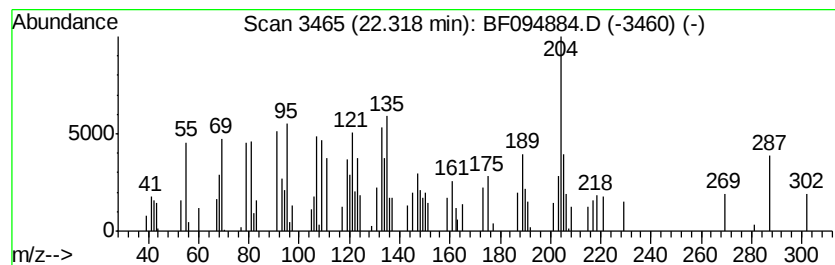
Instrument :
BNA_F
ClientSampleId :
SB-06-7.0-8.0

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 2H-Cyclopentacyclooctene, 4... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.		
22.32	4.83 ng	137206	Perylene-d12	20.30		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		2H-Cyclopentacyclooctene, 4,5,6,...	204	C15H24	1000221-85-8	50
2		1,4-Dimethyl-8-isopropylidenetri...	204	C15H24	1000140-07-7	50
3		Naphthalene, decahydro-4a-methyl...	204	C15H24	017066-67-0	46
4		Guaia-3,9-diene	204	C15H24	000489-83-8	43
5		1H-3a,7-Methanoazulene, octahydr...	204	C15H24	000508-55-4	42



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050417\
 Data File : BF094884.D
 Acq On : 4 May 2017 20:38
 Operator : SJ/MA
 Sample : I2916-04
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-06-7.0-8.0

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.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...	5.02	23.8	ng	543591	1	7.72	456220	20.0
unknown7.38	7.38	110.7	ng	2524630	1	7.72	456220	20.0
unknown15.18	15.18	2.9	ng	113448	4	14.97	784059	20.0
5-Pyrrolidin-2-yl...	15.66	5.1	ng	199785	4	14.97	784059	20.0
n-Hexadecanoic acid	15.94	7.1	ng	277860	4	14.97	784059	20.0
1-Heneicosanol	18.54	12.7	ng	469828	5	18.64	739822	20.0
n-Tetracosanol-1	19.34	12.6	ng	465192	5	18.64	739822	20.0
Supraene	19.77	3.3	ng	93694	6	20.30	568099	20.0
Heptadecane, 9-oc...	20.05	4.9	ng	139336	6	20.30	568099	20.0
Octacosanol	20.09	3.1	ng	87518	6	20.30	568099	20.0
Heneicosane	20.78	3.2	ng	91405	6	20.30	568099	20.0
Eicosyl heptaflu...	20.83	2.6	ng	72904	6	20.30	568099	20.0
unknown21.84	21.84	3.6	ng	102984	6	20.30	568099	20.0
2H-Cyclopentacycl...	22.32	4.8	ng	137206	6	20.30	568099	20.0