

Data Path : Z:\HPCHEM1\BNA F\DATA\BF052217\
 Data File : BF095318.D
 Acq On : 22 May 2017 17:22
 Operator : SJ/MA
 Sample : I3157-03
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FLOOR-B

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:\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	5.131	539	543	560	rBV	143469	356686	11.82%	1.541%
2	5.725	640	644	682	rBV	1044224	2373340	78.62%	10.255%
3	7.301	908	912	928	rBV	1343199	2209361	73.19%	9.547%
4	7.425	928	933	943	rVB	1592134	2544086	84.27%	10.993%
5	7.760	986	990	1001	rVB	371909	446629	14.79%	1.930%
6	7.995	1025	1030	1045	rBV	1598748	1959285	64.90%	8.466%
7	8.660	1139	1143	1165	rBV	809651	1408664	46.66%	6.087%
8	9.783	1330	1334	1363	rBV	335116	644146	21.34%	2.783%
9	11.548	1628	1634	1660	rBV	2006701	3018866	100.00%	13.045%
10	12.248	1749	1753	1768	rBV	183581	330045	10.93%	1.426%
11	12.613	1810	1815	1834	rBV2	467269	769103	25.48%	3.323%
12	13.442	1948	1956	1961	rBV	30468	38186	1.26%	0.165%
13	13.907	2029	2035	2056	rBV	920054	1685050	55.82%	7.281%
14	14.213	2083	2087	2090	rBV2	26540	33730	1.12%	0.146%
15	15.013	2218	2223	2228	rBV	464367	702271	23.26%	3.035%
16	15.160	2244	2248	2258	rVB2	17977	32086	1.06%	0.139%
17	15.960	2375	2384	2388	rBV3	32297	69889	2.32%	0.302%
18	16.801	2523	2527	2537	rBV	118600	176572	5.85%	0.763%
19	17.107	2574	2579	2588	rBV	132534	187825	6.22%	0.812%
20	17.330	2612	2617	2630	rBV2	2255637	2767239	91.66%	11.957%
21	18.583	2821	2830	2837	rBV6	19176	59315	1.96%	0.256%
22	18.683	2842	2847	2867	rVB	441057	761549	25.23%	3.291%
23	19.971	3062	3066	3069	rBV	31205	46317	1.53%	0.200%
24	20.312	3121	3124	3128	rVV	26854	30466	1.01%	0.132%
25	20.359	3128	3132	3141	rVV	247232	430221	14.25%	1.859%
26	20.424	3141	3143	3149	rVB5	24245	31496	1.04%	0.136%
27	21.741	3360	3367	3373	rBV6	14292	30200	1.00%	0.130%

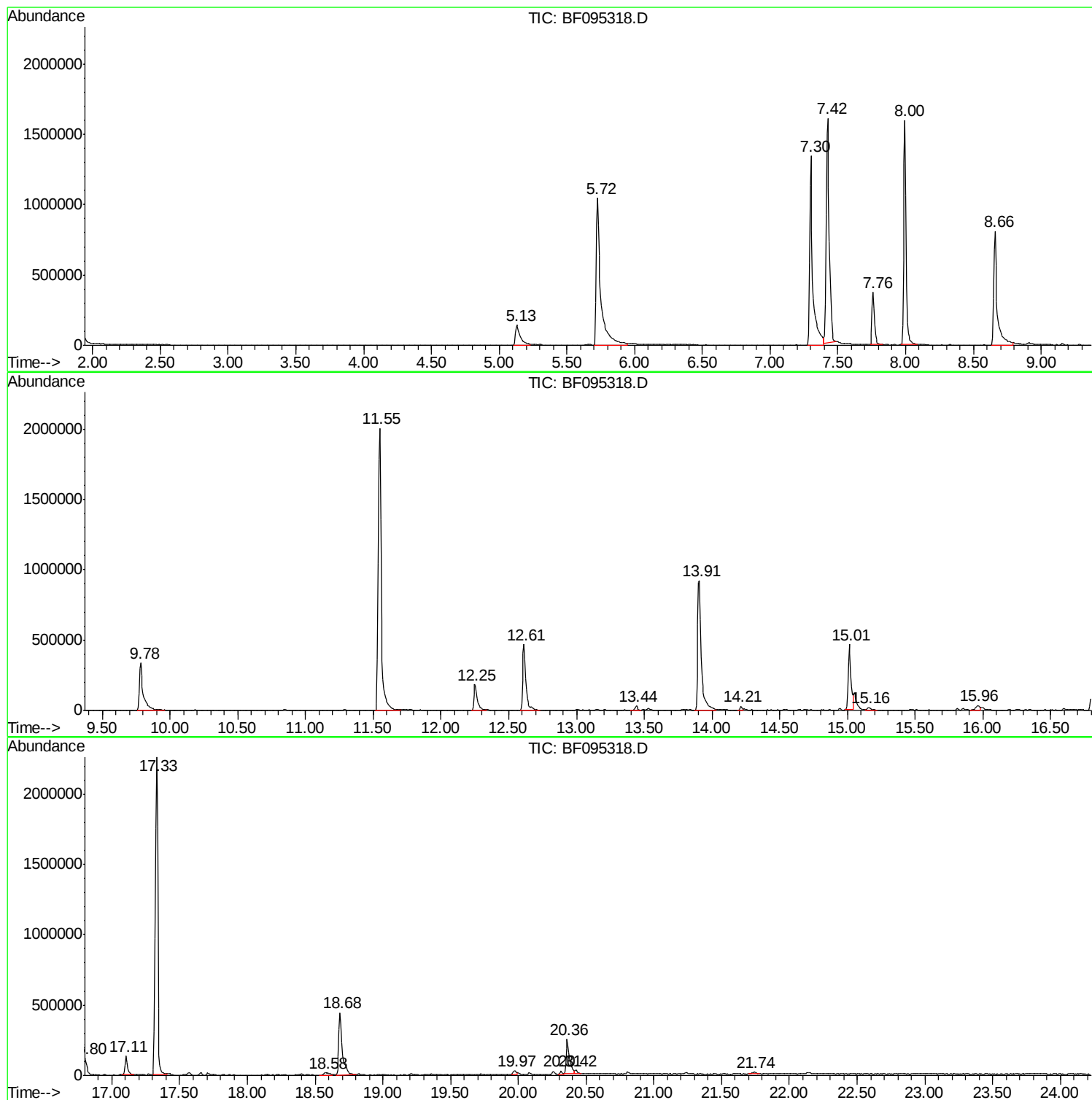
Sum of corrected areas: 23142623

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052217\
Data File : BF095318.D
Acq On : 22 May 2017 17:22
Operator : SJ/MA
Sample : I3157-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
FLOOR-B

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



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052217\
Data File : BF095318.D
Acq On : 22 May 2017 17:22
Operator : SJ/MA
Sample : I3157-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
FLOOR-B

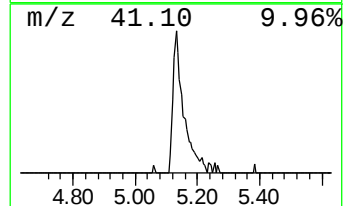
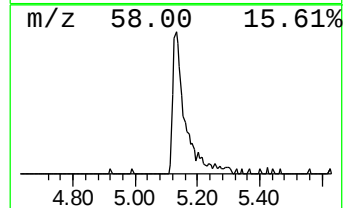
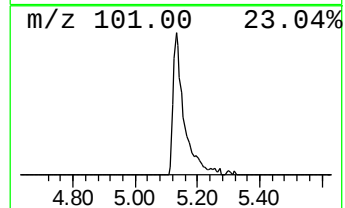
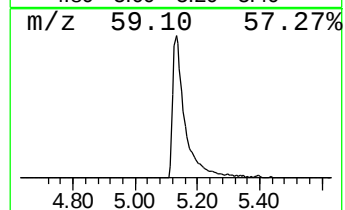
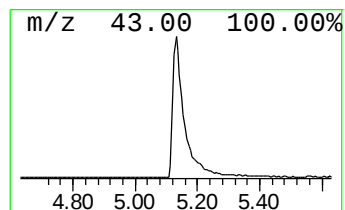
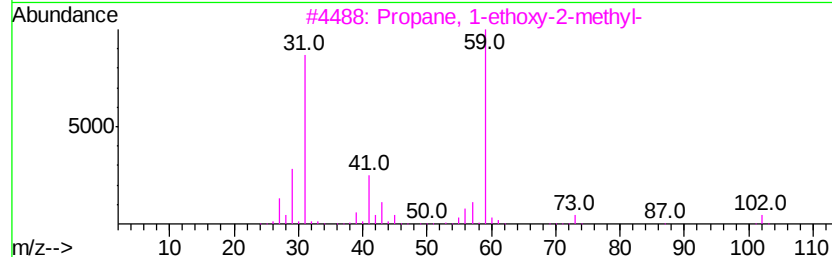
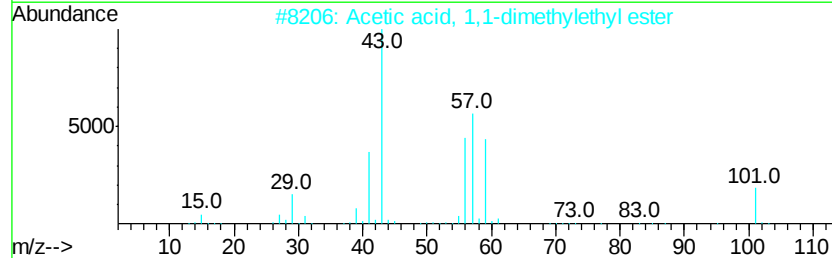
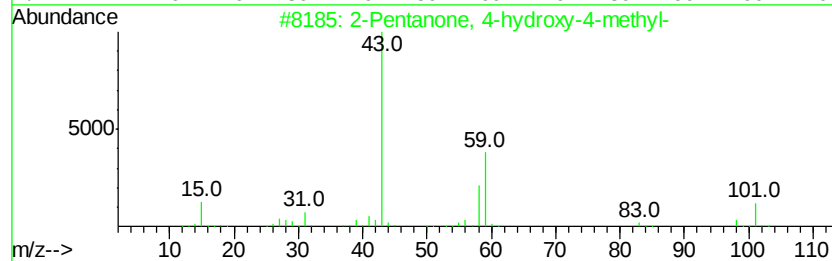
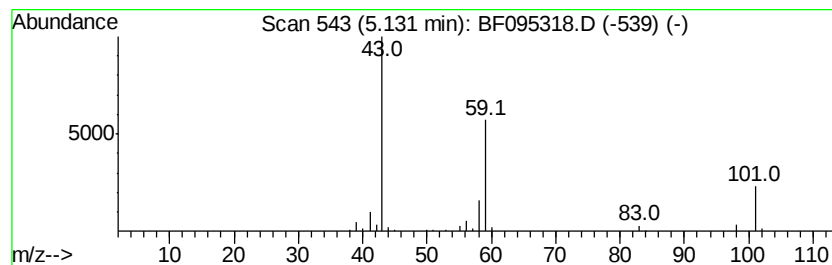
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.13	15.97 ng	356686	1,4-Dichlorobenzene-d4	7.76

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	25
4			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
5			Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052217\
Data File : BF095318.D
Acq On : 22 May 2017 17:22
Operator : SJ/MA
Sample : I3157-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
FLOOR-B

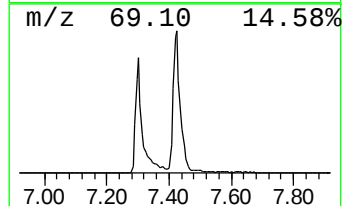
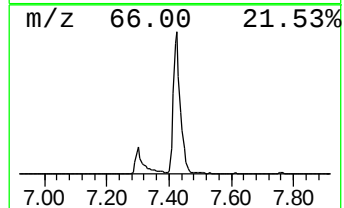
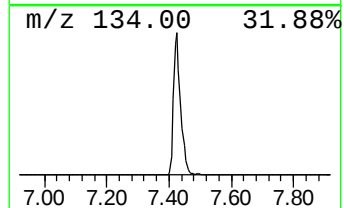
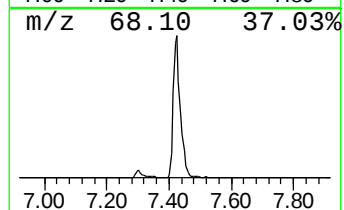
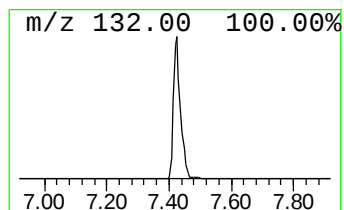
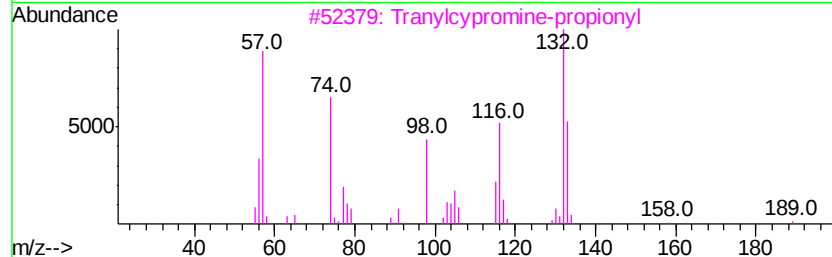
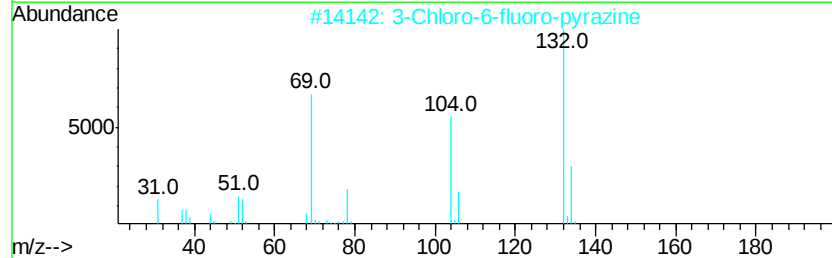
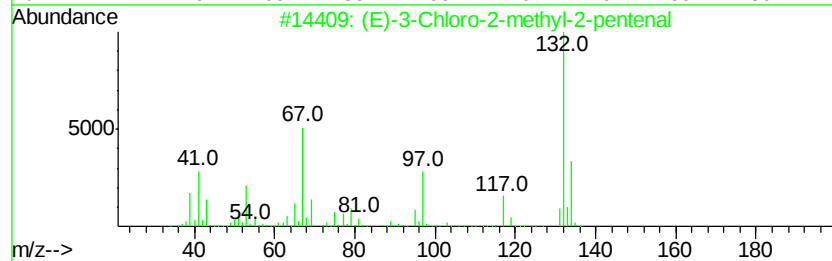
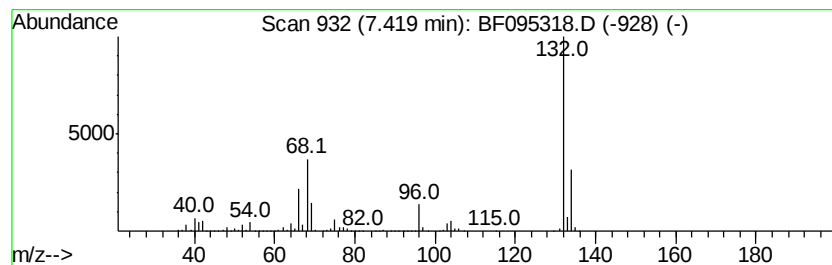
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 unknown7.42 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.42	113.92 ng	2544090	1,4-Dichlorobenzene-d4	7.76

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	35
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		Xenon	132	Xe	007440-63-3	9



Data Path : Z:\HPCHEM1\BNA F\DATA\BF052217\
Data File : BF095318.D
Acq On : 22 May 2017 17:22
Operator : SJ/MA
Sample : I3157-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
FLOOR-B

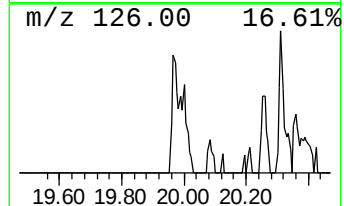
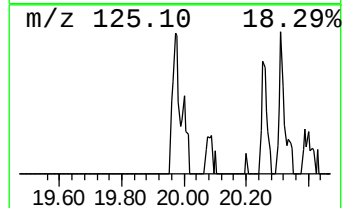
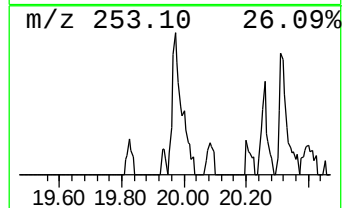
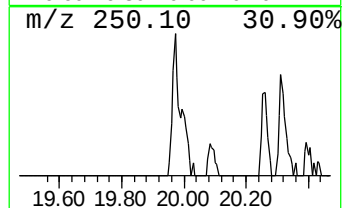
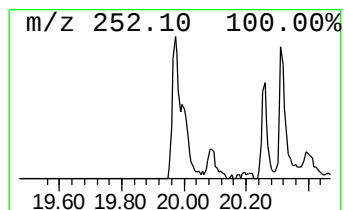
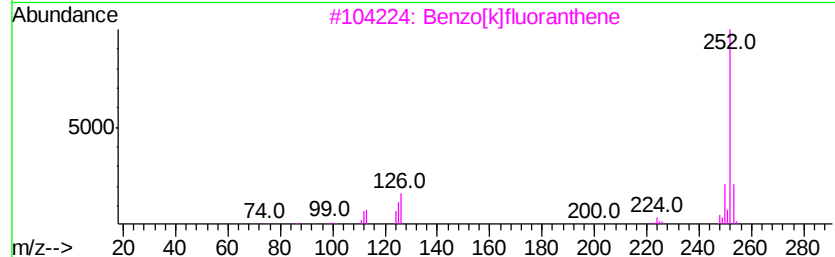
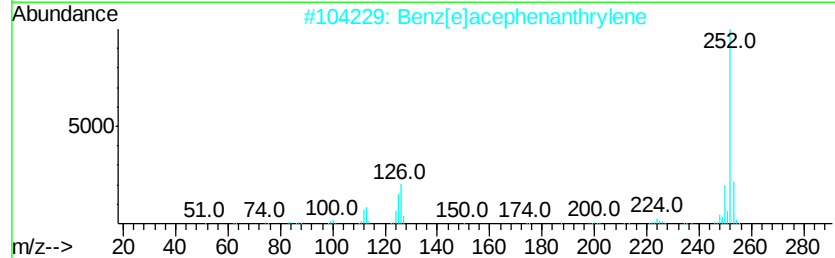
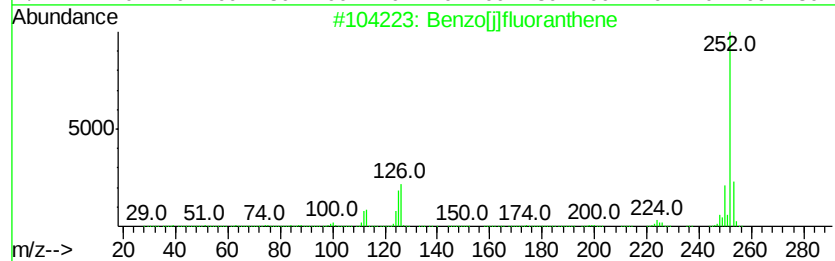
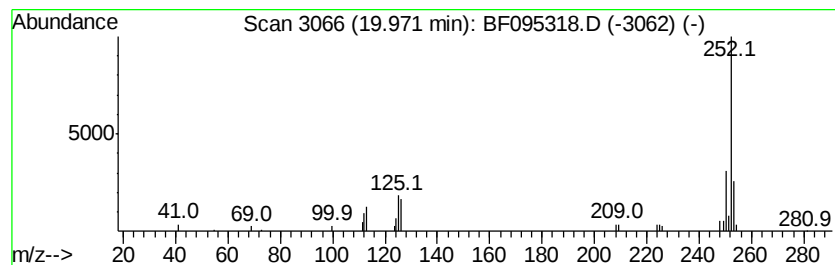
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Benzo[j]fluoranthene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.97	2.15 ng	46317	Perylene-d12	20.36

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[i]fluoranthene	252	C20H12	000205-82-3	95
2			Benzo[e]acephenanthrylene	252	C20H12	000205-99-2	95
3			Benzo[k]fluoranthene	252	C20H12	000207-08-9	95
4			Benzo[e]pyrene	252	C20H12	000192-97-2	94
5			Benzo[a]pyrene	252	C20H12	000050-32-8	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052217\
Data File : BF095318.D
Acq On : 22 May 2017 17:22
Operator : SJ/MA
Sample : I3157-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
FLOOR-B

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...	5.13	16.0	ng	356686	1	7.76	446629	20.0
unknown7.42	7.42	113.9	ng	2544090	1	7.76	446629	20.0
Benzo[j]fluoranthene	19.97	2.1	ng	46317	6	20.36	430221	20.0