

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF072617\
 Data File : BF097097.D
 Acq On : 27 Jul 2017 1:24
 Operator : SJ/JU
 Sample : I4359-03
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CEMETERY-2-(0-2)

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-BF072517.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	4.646	465	469	487	rVB	173522	275637	10.93%	1.351%
2	5.098	541	546	560	rBV	2316833	2506866	99.37%	12.289%
3	6.163	721	727	741	rBV	2112462	2522786	100.00%	12.367%
4	6.269	741	745	748	rBV	2481530	2280313	90.39%	11.179%
5	6.493	780	783	793	rBV	718227	655779	25.99%	3.215%
6	6.651	806	810	819	rBV	2088385	1890809	74.95%	9.269%
7	7.063	875	880	883	rBV	1566210	1443377	57.21%	7.076%
8	7.781	998	1002	1005	rBV	1004895	872712	34.59%	4.278%
9	8.857	1180	1185	1188	rBV	2459599	2220978	88.04%	10.888%
10	9.257	1249	1253	1258	rBV	264310	218307	8.65%	1.070%
11	9.522	1294	1298	1302	rBV	832315	762096	30.21%	3.736%
12	9.981	1373	1376	1379	rVB	46624	35620	1.41%	0.175%
13	10.310	1428	1432	1440	rVB	1126894	1026244	40.68%	5.031%
14	10.451	1452	1456	1463	rBV	63398	69794	2.77%	0.342%
15	10.998	1545	1549	1553	rBV	700343	571575	22.66%	2.802%
16	11.551	1641	1643	1646	rBV	43702	43413	1.72%	0.213%
17	12.586	1814	1819	1822	rBV	1691240	1599972	63.42%	7.843%
18	13.498	1970	1974	1980	rVB	95783	117537	4.66%	0.576%
19	13.621	1991	1995	2003	rBV	594394	589703	23.38%	2.891%
20	14.421	2127	2131	2134	rBV6	31096	58806	2.33%	0.288%
21	14.974	2221	2225	2231	rVB	387914	439847	17.43%	2.156%
22	16.668	2511	2513	2517	rBV5	18478	29781	1.18%	0.146%
23	16.845	2540	2543	2548	rBV7	26269	53554	2.12%	0.263%
24	17.580	2665	2668	2671	rBV4	20783	32771	1.30%	0.161%
25	18.656	2846	2851	2855	rBV8	19213	41404	1.64%	0.203%
26	18.833	2877	2881	2882	rBV4	26089	39147	1.55%	0.192%

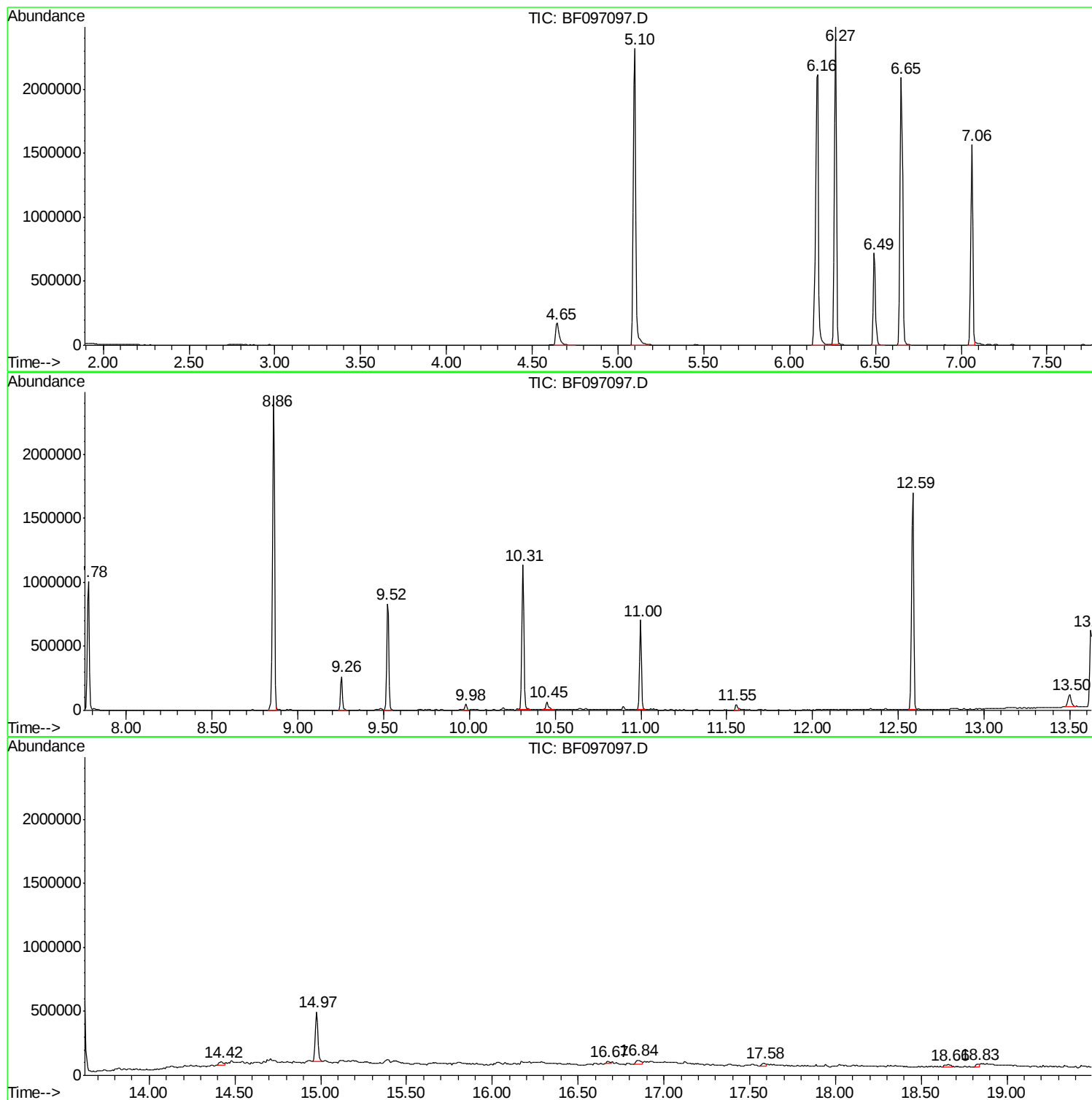
Sum of corrected areas: 20398828

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF072617\
Data File : BF097097.D
Acq On : 27 Jul 2017 1:24
Operator : SJ/JU
Sample : I4359-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CEMETERY-2-(0-2)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072517.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\BF072617\
Data File : BF097097.D
Acq On : 27 Jul 2017 1:24
Operator : SJ/JU
Sample : I4359-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CEMETERY-2-(0-2)

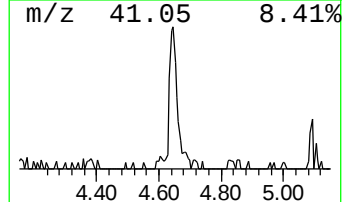
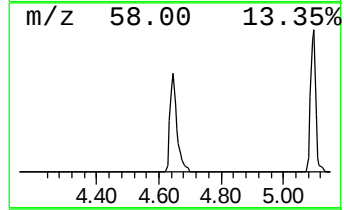
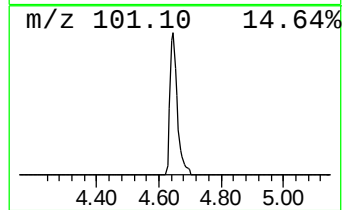
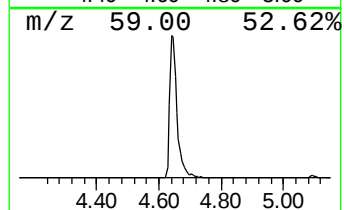
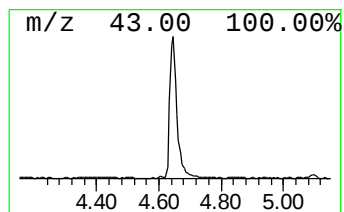
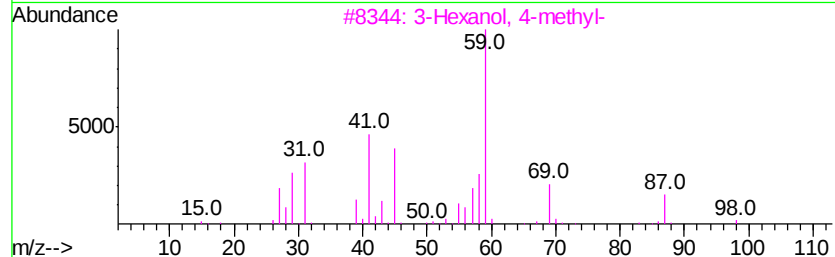
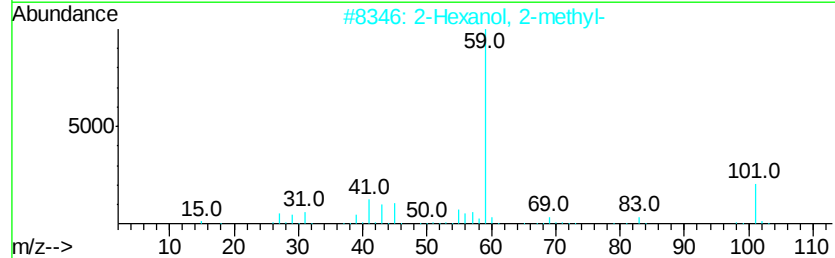
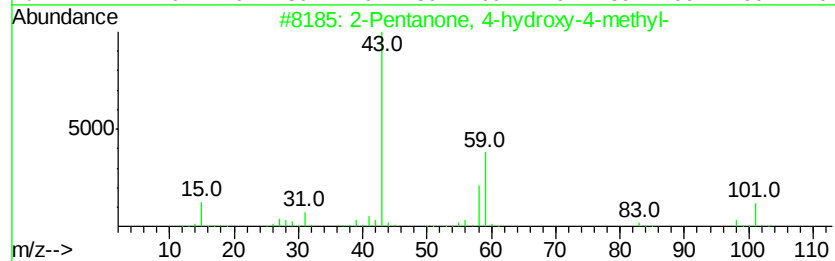
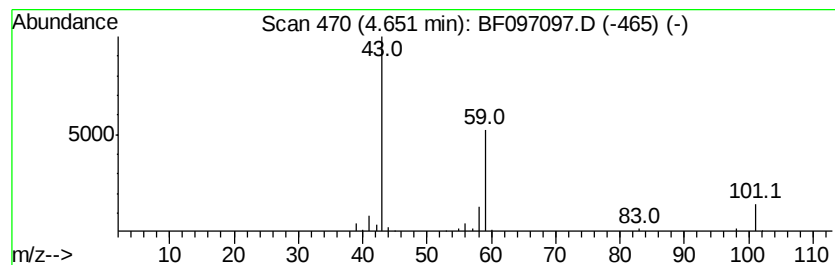
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
4.65	8.41 ng	275637	1,4-Dichlorobenzene-d4	6.49

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
3			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
4			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



Data Path : Z:\HPCHEM1\BNA F\DATA\BF072617\
Data File : BF097097.D
Acq On : 27 Jul 2017 1:24
Operator : SJ/JU
Sample : I4359-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CEMETERY-2-(0-2)

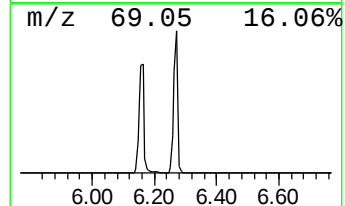
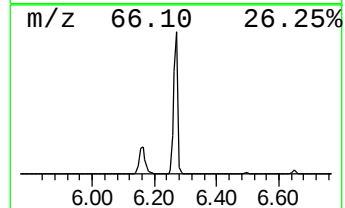
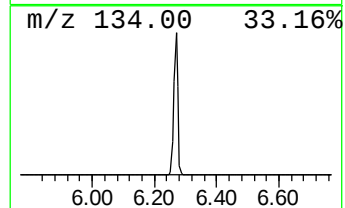
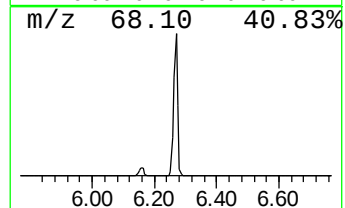
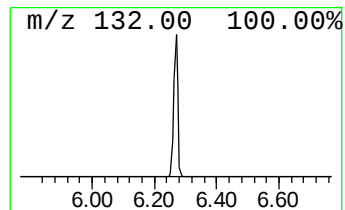
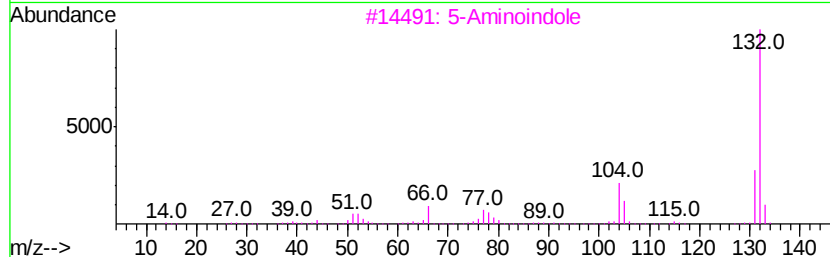
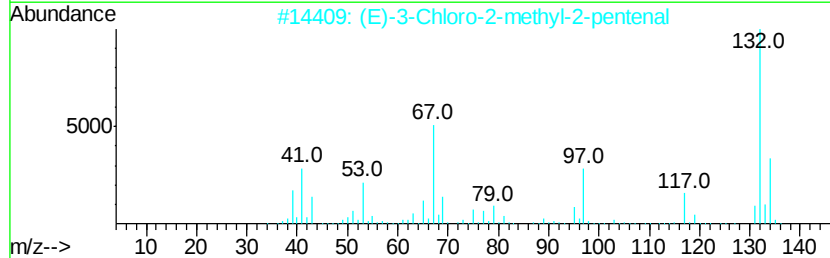
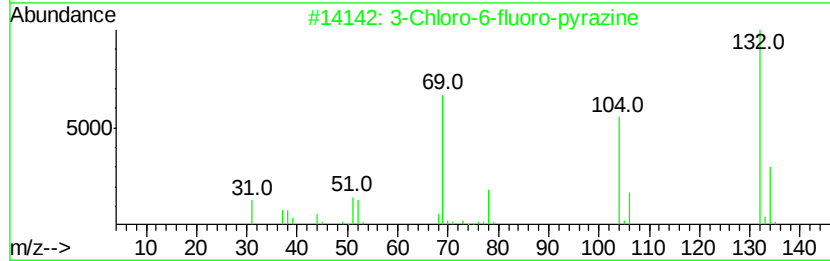
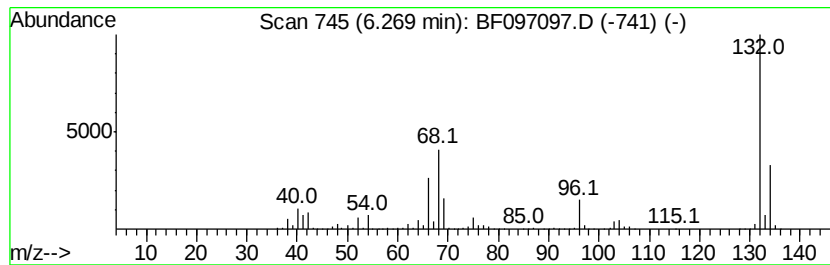
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 unknown6.27 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.27	69.55 ng	2280310	1,4-Dichlorobenzene-d4	6.49

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	27
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9



Data Path : Z:\HPCHEM1\BNA F\DATA\BF072617\
Data File : BF097097.D
Acq On : 27 Jul 2017 1:24
Operator : SJ/JU
Sample : I4359-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

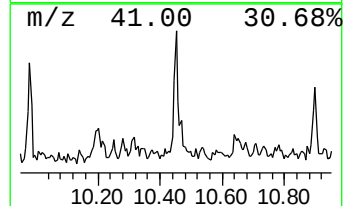
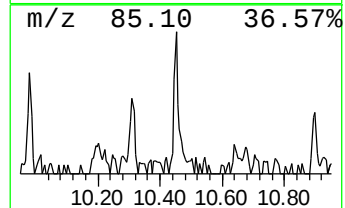
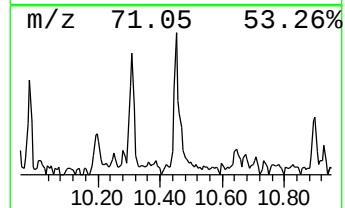
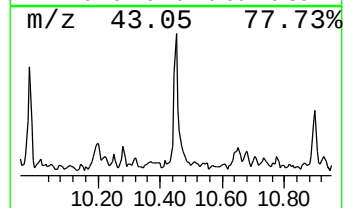
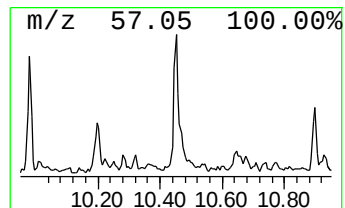
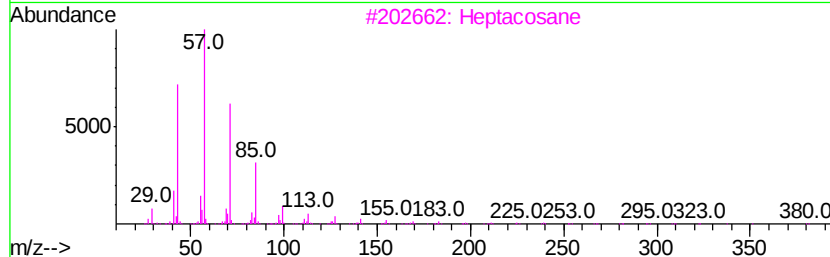
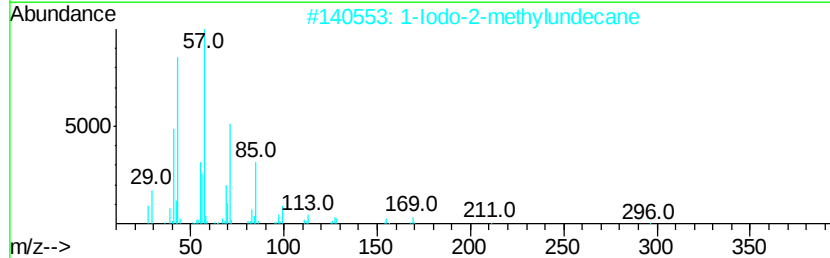
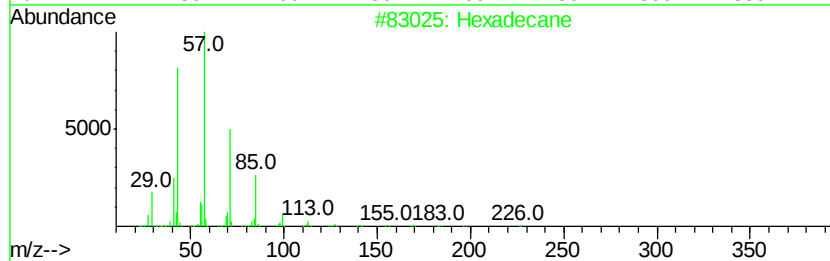
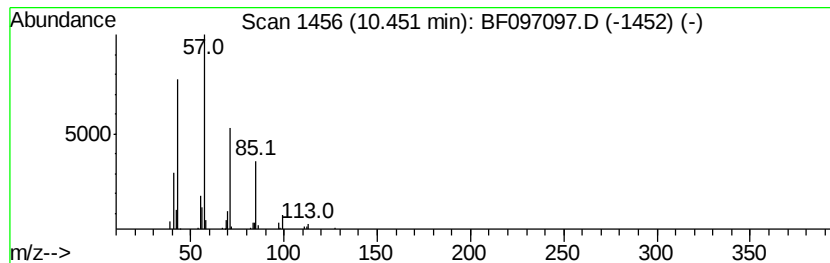
Instrument :
BNA_F
ClientSampleId :
CEMETERY-2-(0-2)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Hexadecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.45	2.44 ng	69794	Phenanthrene-d10	11.00	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	86
2	1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	86
3	Heptacosane	380	C27H56	000593-49-7	78
4	Tridecane	184	C13H28	000629-50-5	72
5	Tetradecane	198	C14H30	000629-59-4	72



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF072617\
Data File : BF097097.D
Acq On : 27 Jul 2017 1:24
Operator : SJ/JU
Sample : I4359-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CEMETERY-2-(0-2)

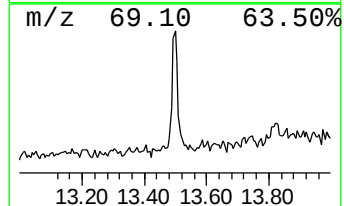
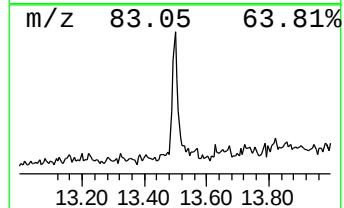
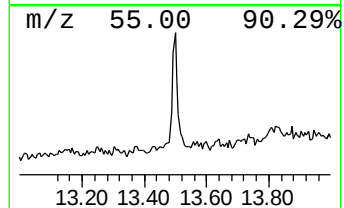
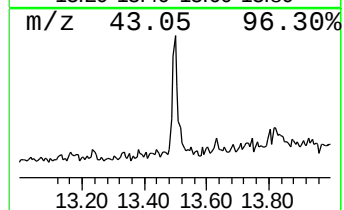
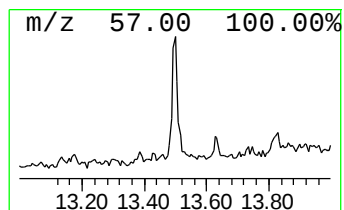
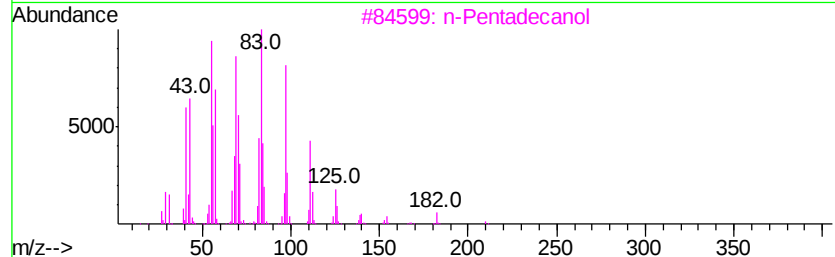
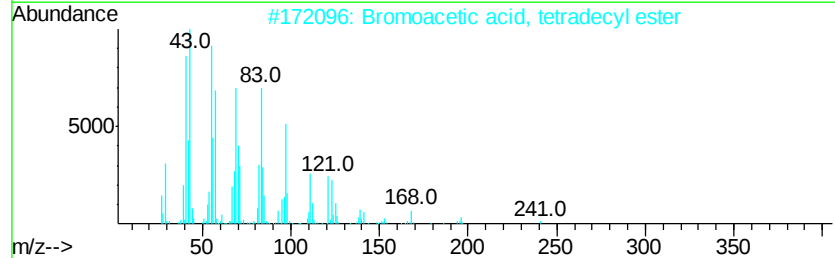
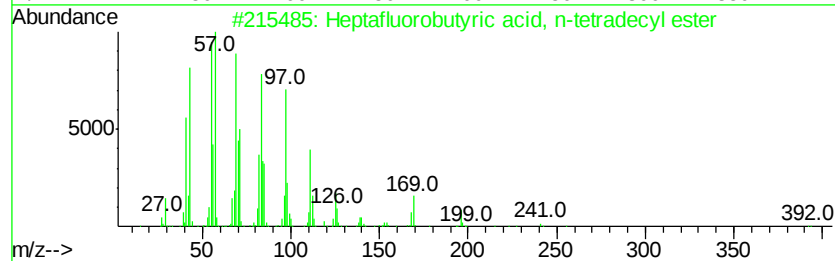
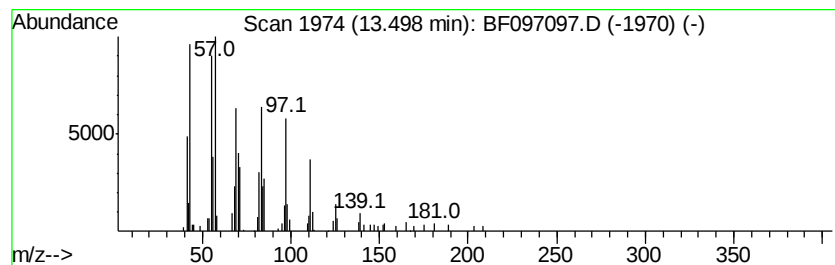
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Heptafluorobutyric acid, n-... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.50	3.99 ng	117537	Chrysene-d12	13.62

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptafluorobutyric acid, n-tetra...	410	C18H29F7O2	007365-36-8	91
2		Bromoacetic acid, tetradecyl ester	334	C16H31BrO2	018992-01-3	90
3		n-Pentadecanol	228	C15H32O	000629-76-5	90
4		1-Hexadecanol	242	C16H34O	036653-82-4	90
5		n-Tetracosanol-1	354	C24H50O	000506-51-4	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF072617\
Data File : BF097097.D
Acq On : 27 Jul 2017 1:24
Operator : SJ/JU
Sample : I4359-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

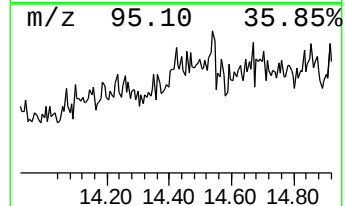
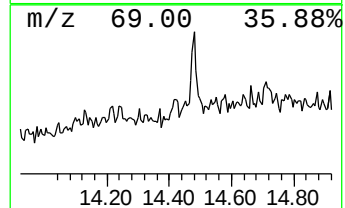
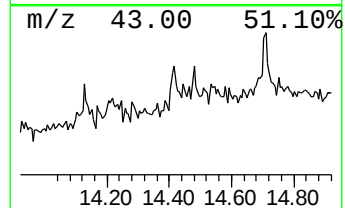
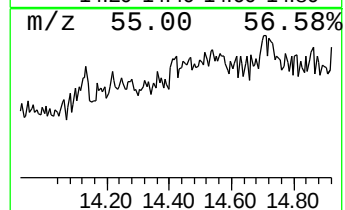
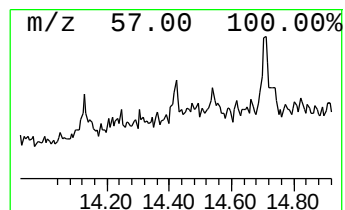
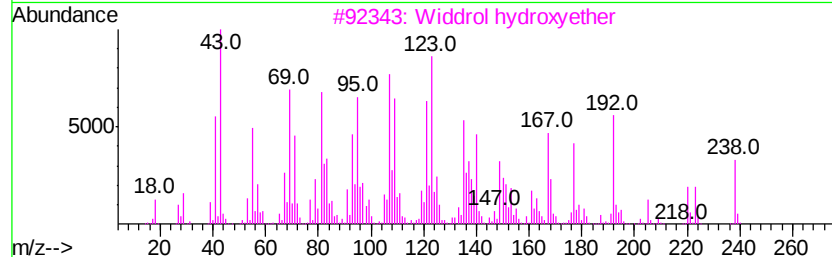
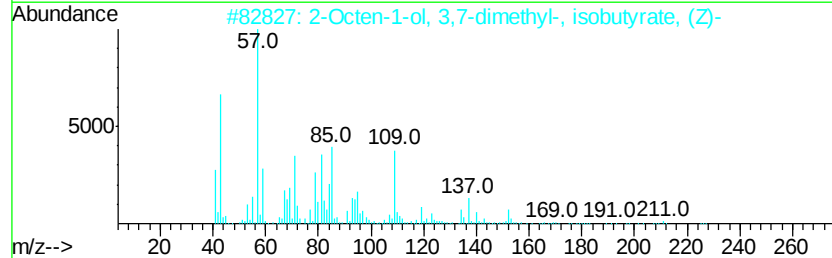
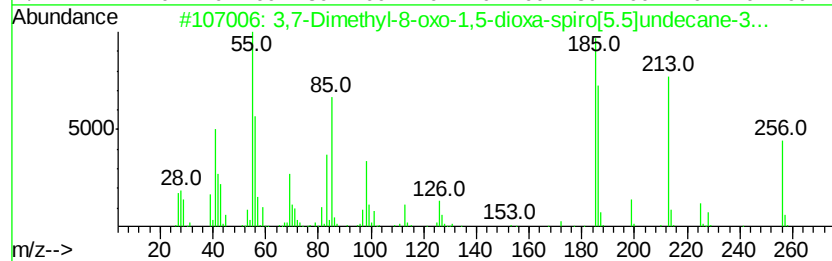
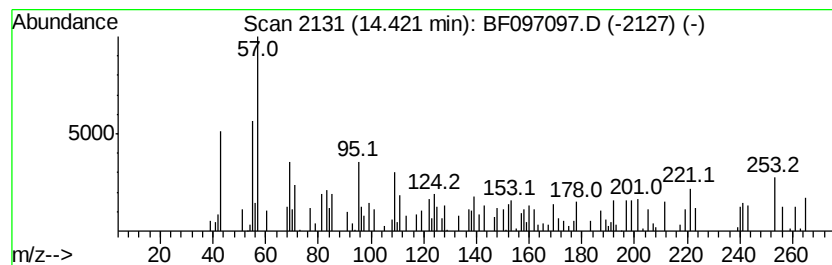
Instrument :
BNA_F
ClientSampleId :
CEMETERY-2-(0-2)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 unknown14.42 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.42	2.67 ng	58806	Perylene-d12	14.97	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,7-Dimethyl-8-oxo-1,5-dioxo-spi...	256	C13H20O5	1000193-79-8	10
2	2-Octen-1-ol, 3,7-dimethyl-, iso...	226	C14H26O2	1000132-45-6	10
3	Widdrol hvdroxyether	238	C15H26O2	029620-91-5	10
4	p-Nitrophenyl nonyl ether	265	C15H23NO3	086702-46-7	10
5	2,5,5,8a-Tetramethyl-4-methylene...	238	C14H22O3	1000192-73-0	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF072617\
Data File : BF097097.D
Acq On : 27 Jul 2017 1:24
Operator : SJ/JU
Sample : I4359-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CEMETERY-2-(0-2)

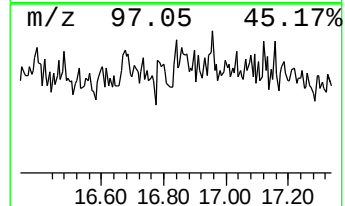
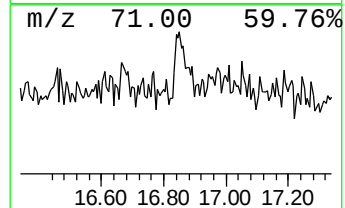
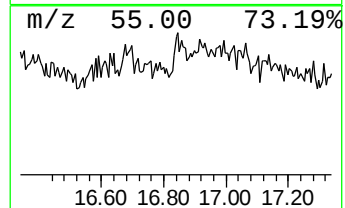
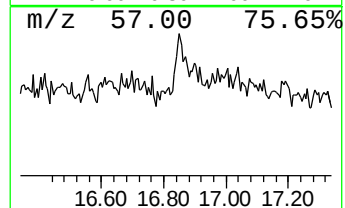
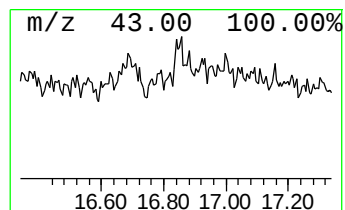
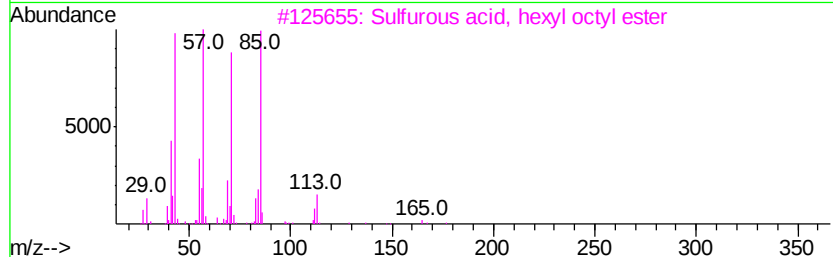
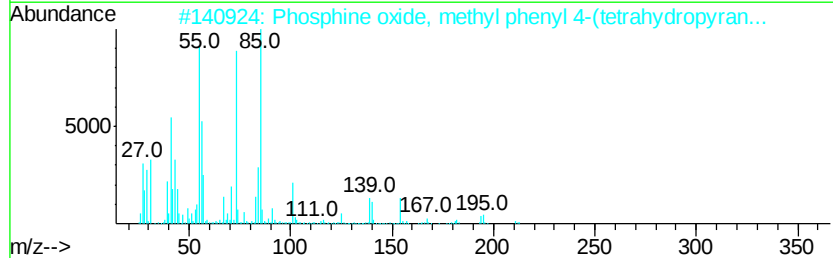
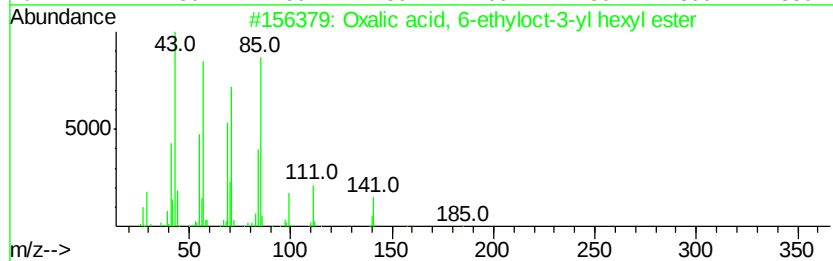
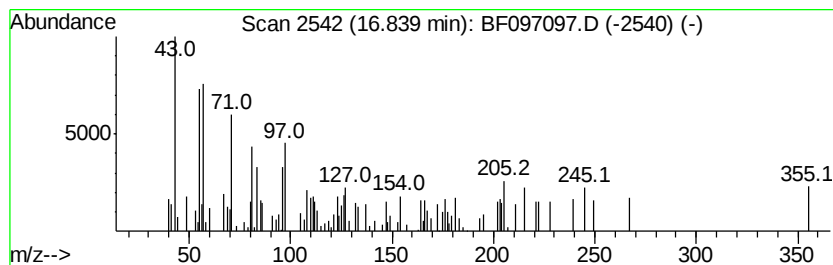
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown16.84 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.84	2.44 ng	53554	Perylene-d12	14.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Oxalic acid, 6-ethyloct-3-yl hex...	314	C18H34O4	1000309-34-4	10
2		Phosphine oxide, methyl phenyl 4...	296	C16H25O3P	1000195-93-9	10
3		Sulfurous acid, hexyl octyl ester	278	C14H30O3S	1000309-13-0	10
4		4,6-Dioxoheptanoic acid	158	C7H10O4	051568-18-4	10
5		Sulfurous acid, 2-ethylhexyl iso...	278	C14H30O3S	1000309-19-0	10



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF072617\
Data File : BF097097.D
Acq On : 27 Jul 2017 1:24
Operator : SJ/JU
Sample : I4359-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CEMETERY-2-(0-2)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072517.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...	4.65	8.4	ng	275637	1	6.49	655779	20.0
unknown6.27	6.27	69.5	ng	2280310	1	6.49	655779	20.0
Hexadecane	10.45	2.4	ng	69794	4	11.00	571575	20.0
Heptafluorobutyri...	13.50	4.0	ng	117537	5	13.62	589703	20.0
unknown14.42	14.42	2.7	ng	58806	6	14.97	439847	20.0
unknown16.84	16.84	2.4	ng	53554	6	14.97	439847	20.0