

Data Path : Z:\HPCHEM1\BNA F\DATA\BF061115\
 Data File : BF079930.D
 Acq On : 12 Jun 2015 9:40
 Operator : TP/IZ
 Sample : G2449-10
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FM7

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-BF061115.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	2.170	85	86	92	rVV	137931	177806	3.61%	0.612%
2	2.273	92	95	97	rVV	3235188	4928220	100.00%	16.952%
3	2.353	100	102	109	rBV	573566	945981	19.20%	3.254%
4	2.695	130	132	136	rVB	122070	205790	4.18%	0.708%
5	5.610	385	387	390	rBV	131384	178506	3.62%	0.614%
6	6.010	419	422	424	rBV	1135589	1317508	26.73%	4.532%
7	6.993	505	508	510	rBV	1825491	1517061	30.78%	5.218%
8	7.165	520	523	525	rVB	1171430	1434137	29.10%	4.933%
9	7.393	541	543	545	rVB	459471	367511	7.46%	1.264%
10	7.553	555	557	559	rBV	1403390	1116134	22.65%	3.839%
11	7.953	589	592	594	rBV	716149	807959	16.39%	2.779%
12	8.685	654	656	658	rBV	670966	542444	11.01%	1.866%
13	9.759	748	750	752	rVB	2050371	1655902	33.60%	5.696%
14	10.148	782	784	786	rBV	149663	129822	2.63%	0.447%
15	10.456	809	811	813	rVB	658383	635194	12.89%	2.185%
16	11.256	878	881	883	rBV	1051645	1098782	22.30%	3.780%
17	11.976	942	944	946	rBV2	622952	935664	18.99%	3.218%
18	12.011	946	947	948	rVB	411557	285438	5.79%	0.982%
19	12.548	992	994	996	rVB	140825	140639	2.85%	0.484%
20	12.639	1000	1002	1006	rBV2	120094	234787	4.76%	0.808%
21	13.314	1059	1061	1063	rBV	747769	691058	14.02%	2.377%
22	13.588	1080	1085	1087	rBV	504076	875257	17.76%	3.011%
23	13.737	1096	1098	1101	rBV	1921123	1842091	37.38%	6.336%
24	13.828	1104	1106	1109	rBV2	79808	192285	3.90%	0.661%
25	14.320	1146	1149	1150	rBV	331590	466223	9.46%	1.604%
26	14.857	1194	1196	1201	rVB2	223358	359993	7.30%	1.238%
27	15.085	1212	1216	1218	rBV2	739336	1296287	26.30%	4.459%
28	15.120	1218	1219	1221	rVB	270390	189436	3.84%	0.652%
29	16.514	1338	1341	1345	rVB3	350351	1028543	20.87%	3.538%
30	16.903	1373	1375	1379	rVB	291291	427706	8.68%	1.471%
31	16.983	1380	1382	1384	rBV	173373	218410	4.43%	0.751%
32	17.063	1387	1389	1392	rVB	397996	614405	12.47%	2.113%
33	18.034	1471	1474	1478	rVB	297383	697249	14.15%	2.398%
34	18.606	1521	1524	1534	rVB	335590	889756	18.05%	3.061%

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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 >
Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061115.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	19.383	1589	1592	1597	rVB	161471	410089	8.32%	1.411%
36	19.474	1598	1600	1605	rVB	98460	217725	4.42%	0.749%

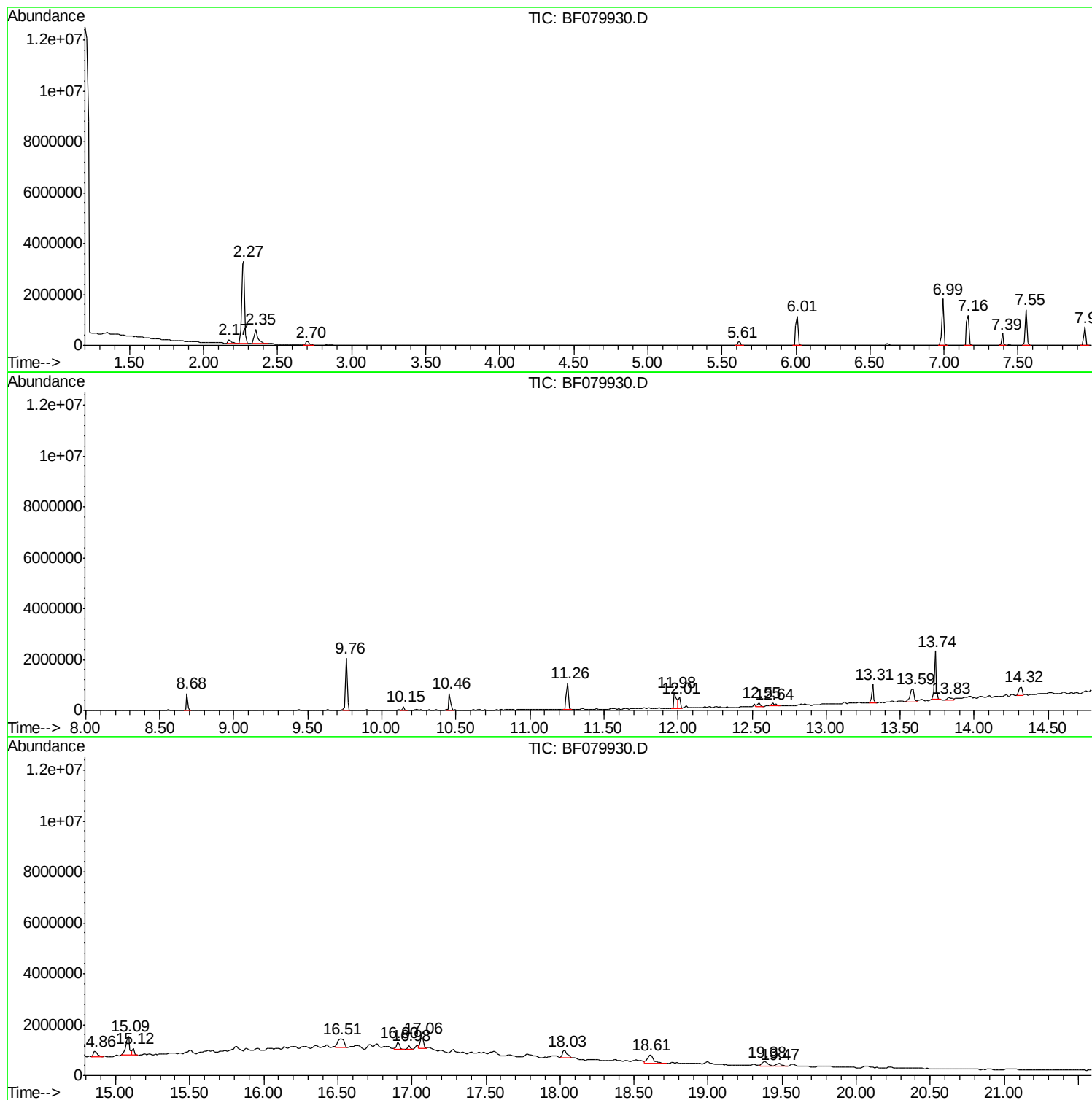
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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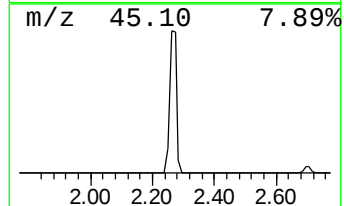
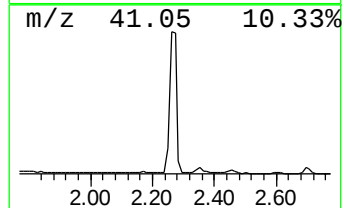
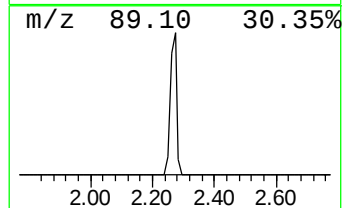
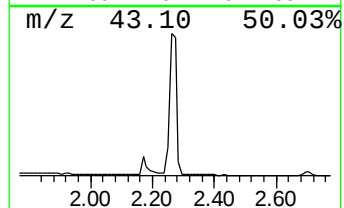
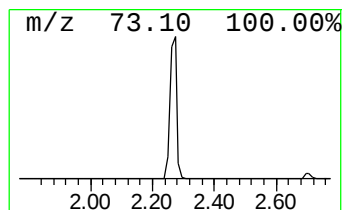
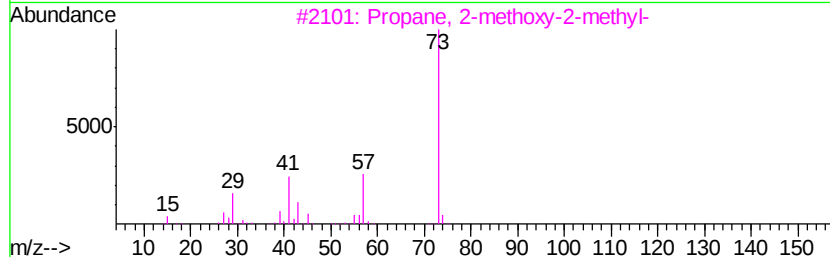
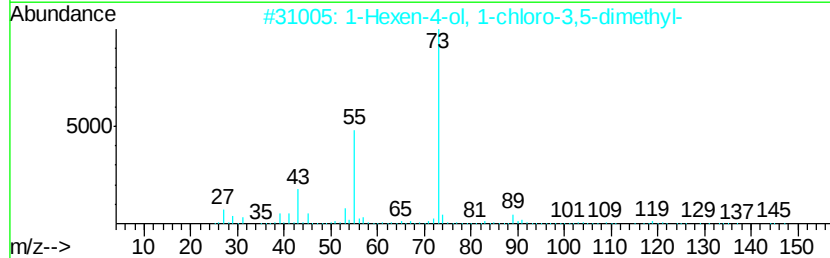
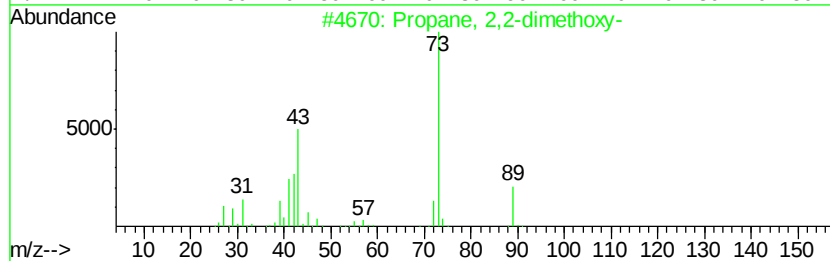
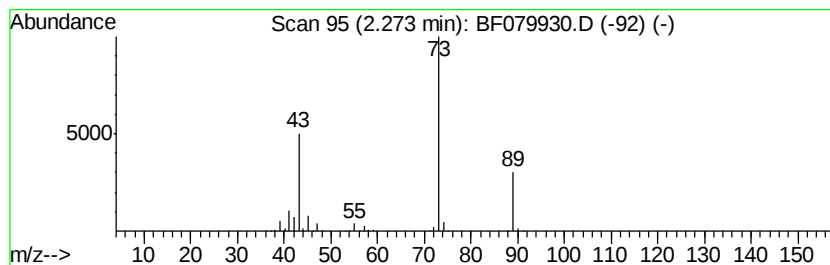
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown2.27 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.27	268.19 ng	4928220	1,4-Dichlorobenzene-d4	7.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	38
2		1-Hexen-4-ol, 1-chloro-3,5-dimet...	162	C8H15ClO	100244-09-5	9
3		Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	9
4		2-Hydroxy-2-methylbutyric acid	118	C5H10O3	003739-30-8	9
5		N-Ethylformamide	73	C3H7NO	000627-45-2	5



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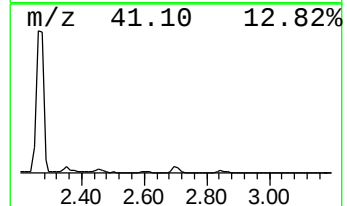
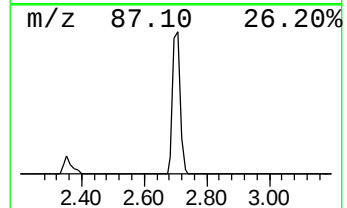
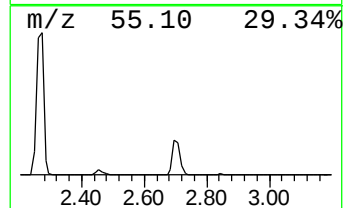
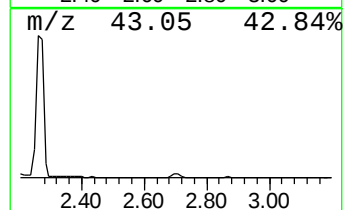
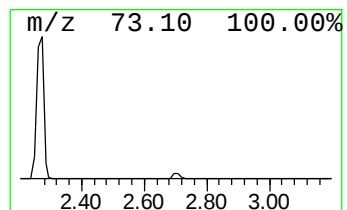
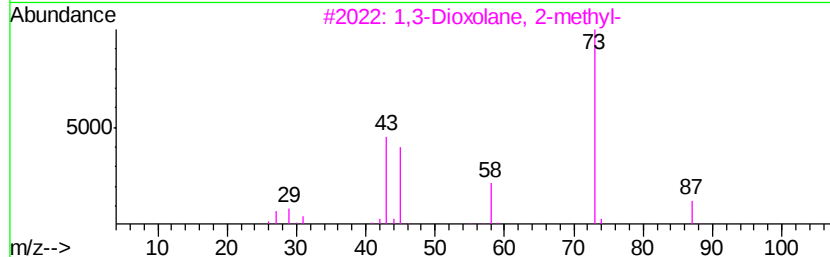
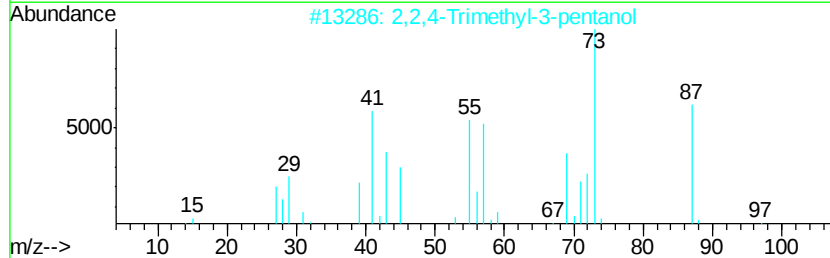
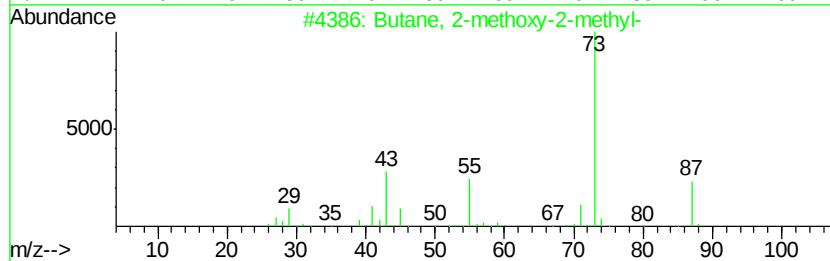
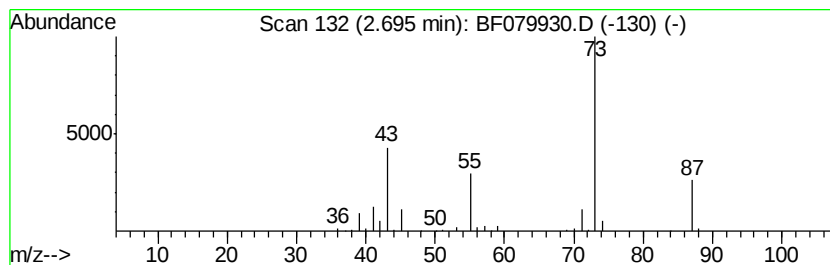
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Butane, 2-methoxy-2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.70	11.20 ng	205790	1,4-Dichlorobenzene-d4	7.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12
5		1-Propene-1-thiol	74	C3H6S	000925-89-3	12



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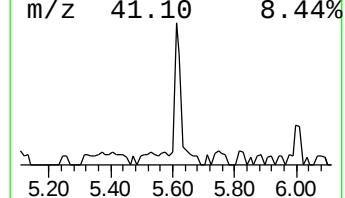
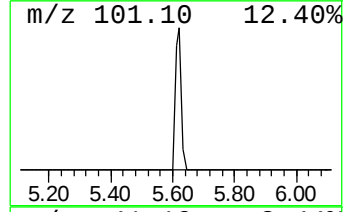
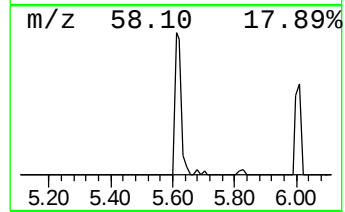
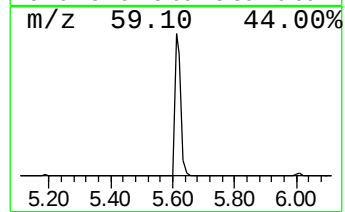
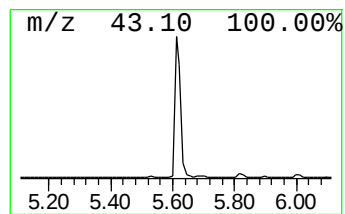
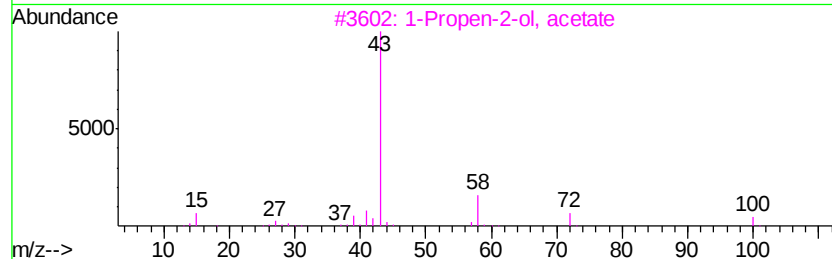
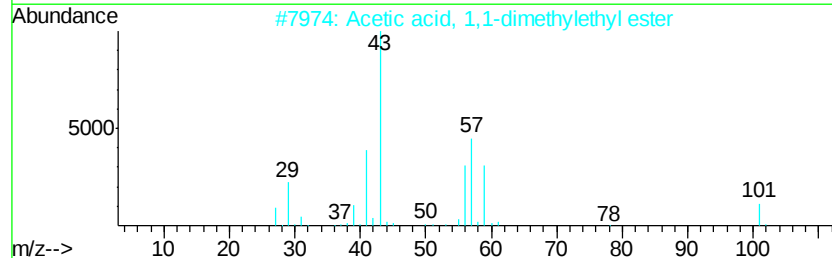
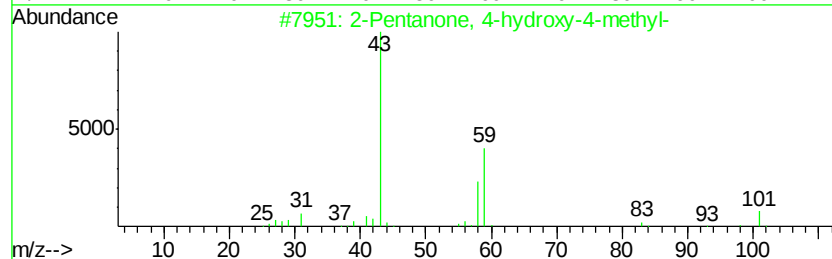
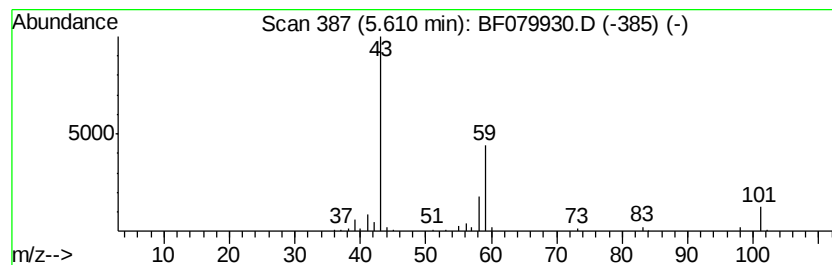
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.61	9.71 ng	178506	1,4-Dichlorobenzene-d4	7.39

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	78	
2	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	28	
3	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10	
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9	
5	Diacetamide	101	C4H7NO2	000625-77-4	9	



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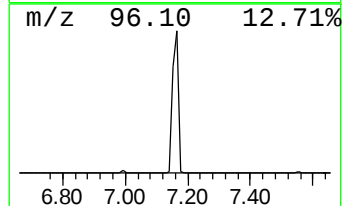
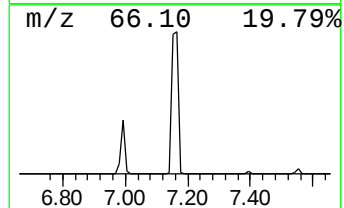
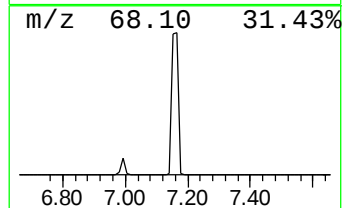
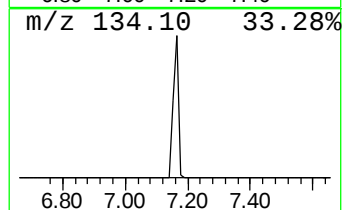
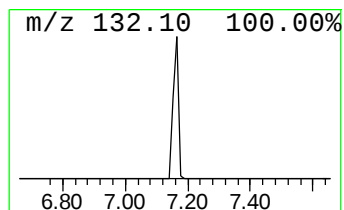
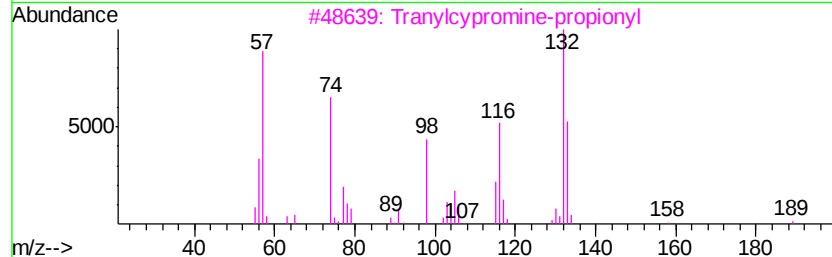
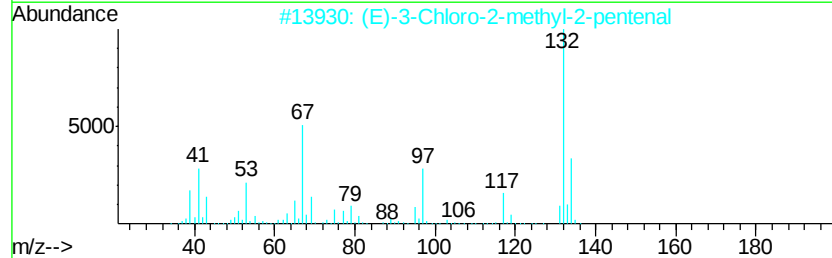
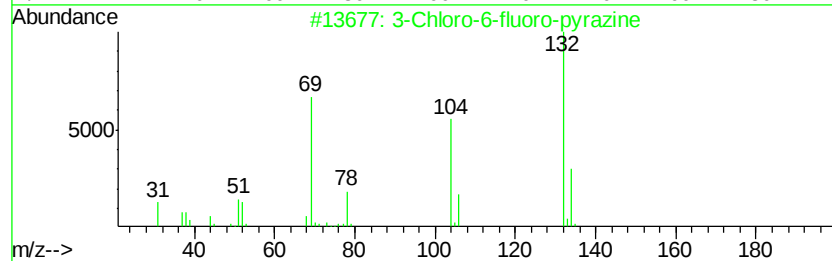
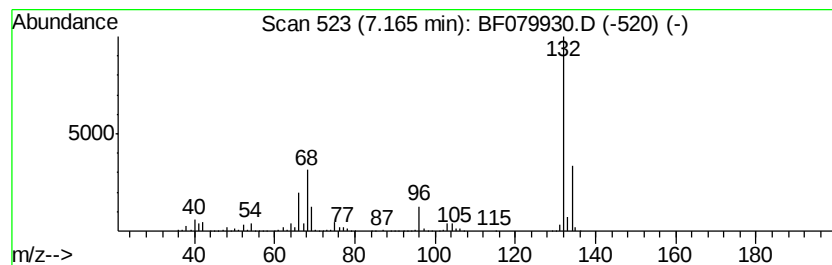
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 unknown7.16 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.16	78.05 ng	1434140	1,4-Dichlorobenzene-d4	7.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Tranvlcypromine-propionyl	189	C12H15NO	1000123-86-3	12
4		2H-Cyclopentadlpvridazine, 2-me...	132	C8H8N2	022291-85-6	9
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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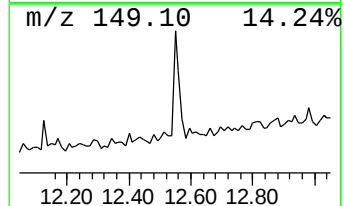
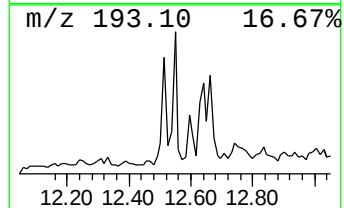
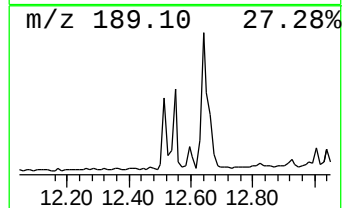
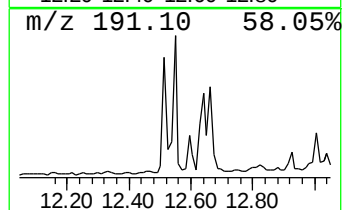
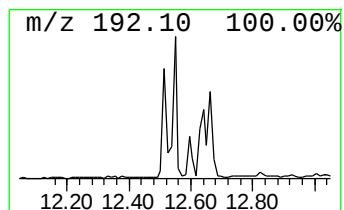
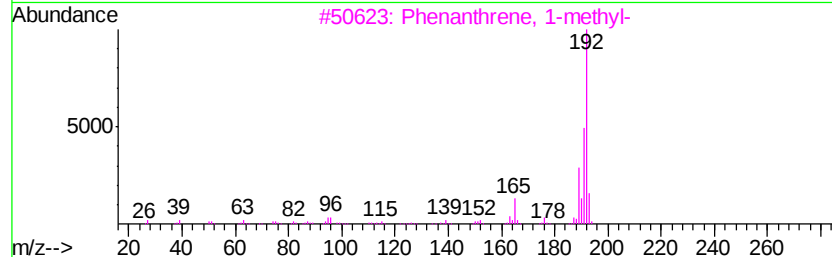
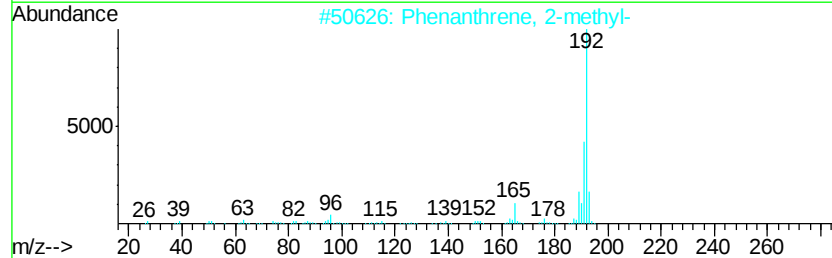
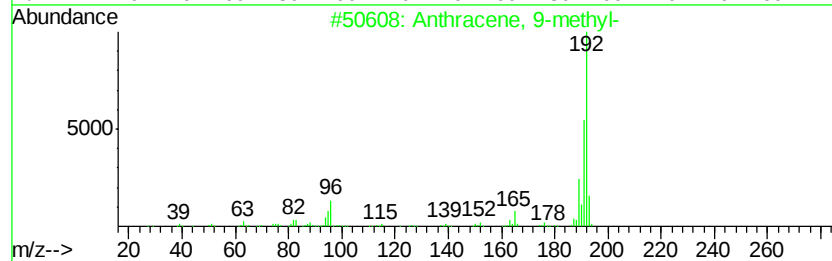
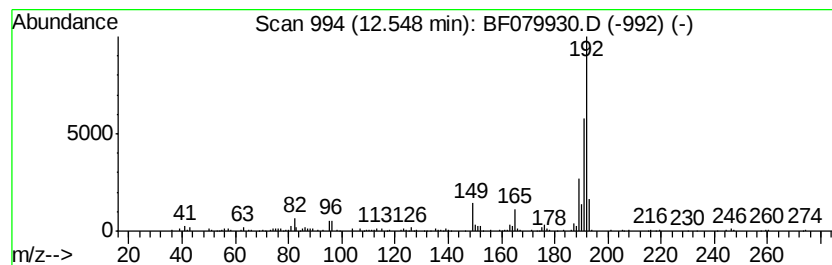
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Anthracene, 9-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.55	3.01 ng	140639	Phenanthrene-d10	11.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Anthracene, 9-methyl-	192	C15H12	000779-02-2	95
2		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	90
3		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	90
4		Anthracene, 1-methyl-	192	C15H12	000610-48-0	90
5		1H-Cyclopropa[1]phenanthrene, 1a, ...	192	C15H12	000949-41-7	89



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Sample : G2449-10
Misc :
ALS Vial : 41 Sample Multiplier: 1

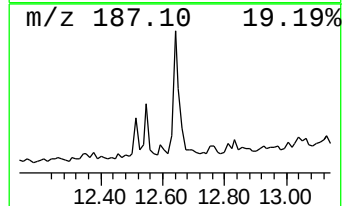
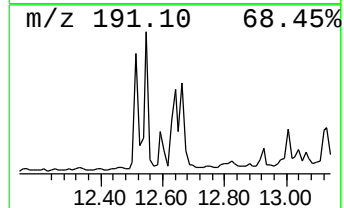
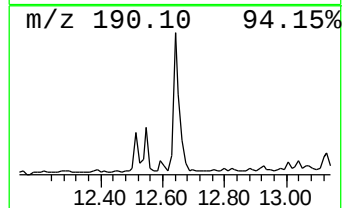
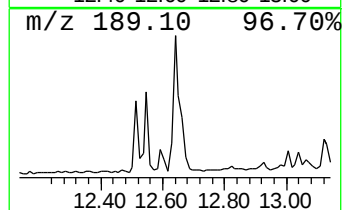
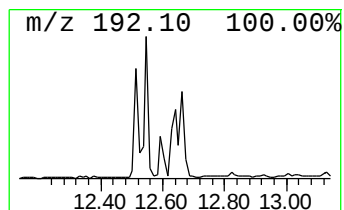
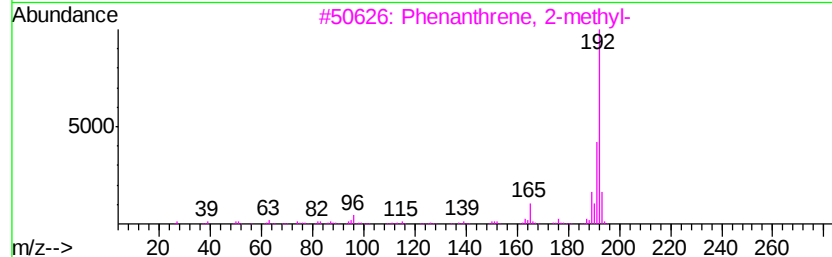
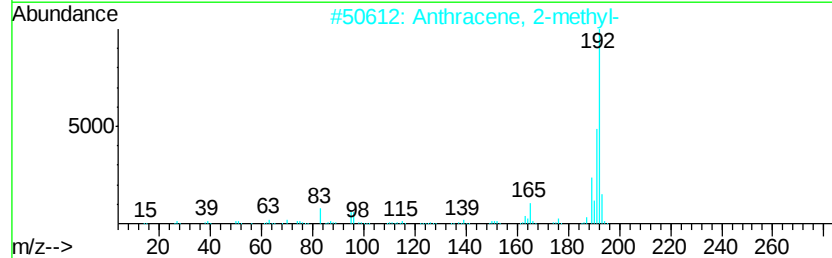
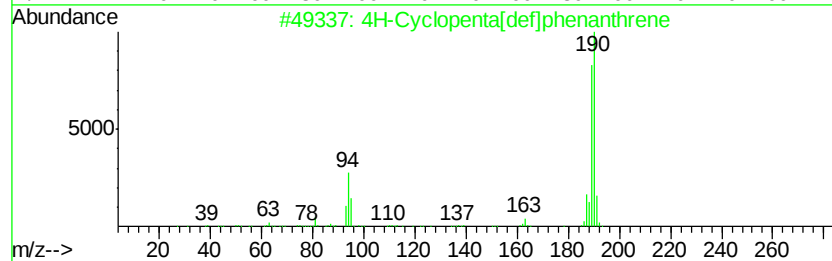
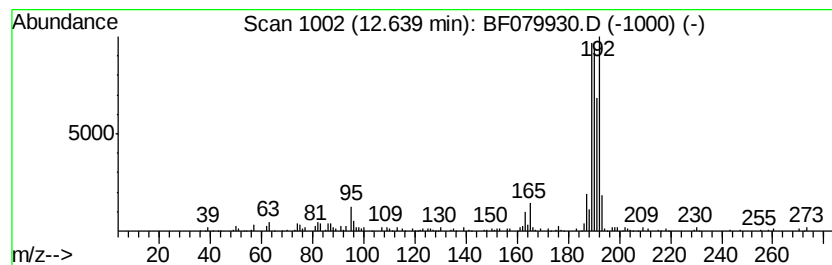
Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 unknown12.64 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.64	5.02 ng	234787	Phenanthrene-d10	11.98	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	49
2	Anthracene, 2-methyl-	192	C15H12	000613-12-7	46
3	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	46
4	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	42
5	Anthracene, 9-methyl-	192	C15H12	000779-02-2	41



Data Path : Z:\HPCHEM1\BNA F\DATA\BF061115\
Data File : BF079930.D
Acq On : 12 Jun 2015 9:40
Operator : TP/IZ
Sample : G2449-10
Misc :
ALS Vial : 41 Sample Multiplier: 1

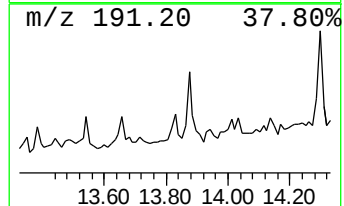
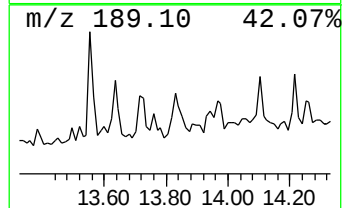
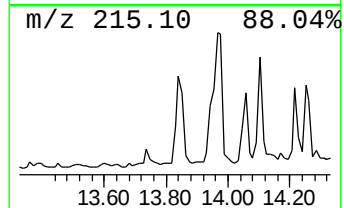
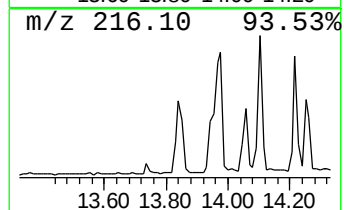
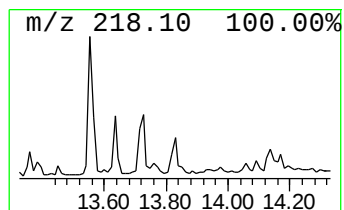
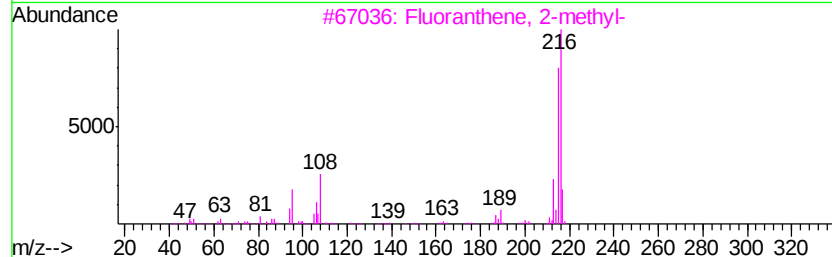
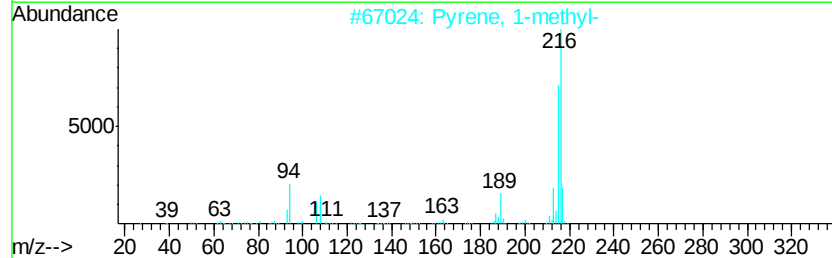
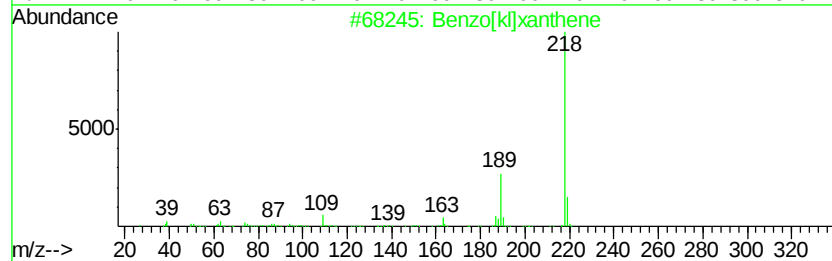
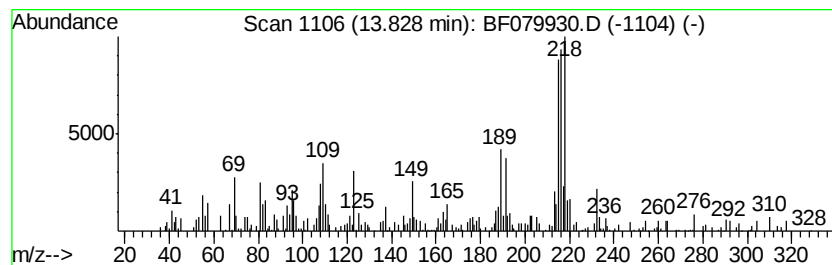
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF061115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Benzo[k]xanthene Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.83	2.97 ng	192285	Chrysene-d12	15.09
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzo[k]xanthene	218 C16H10O	000200-23-7 60
2		Pyrene, 1-methyl-	216 C17H12	002381-21-7 55
3		Fluoranthene, 2-methyl-	216 C17H12	033543-31-6 50
4		Pyrene, 4-methyl-	216 C17H12	003353-12-6 46
5		11H-Benzo[a]fluorene	216 C17H12	000238-84-6 46



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061115\
Data File : BF079930.D
Acq On : 12 Jun 2015 9:40
Operator : TP/IZ
Sample : G2449-10
Misc :
ALS Vial : 41 Sample Multiplier: 1

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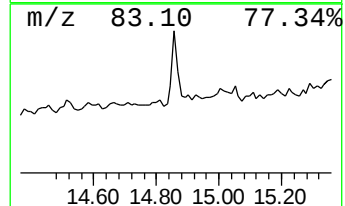
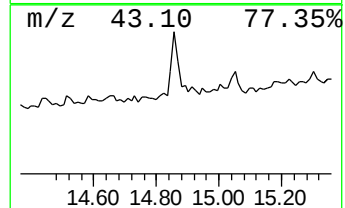
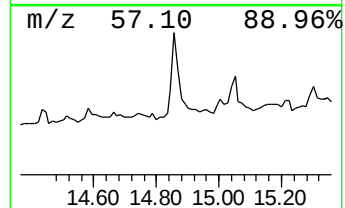
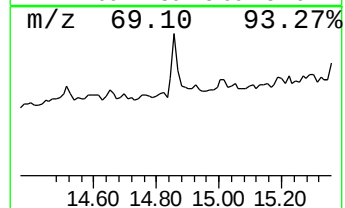
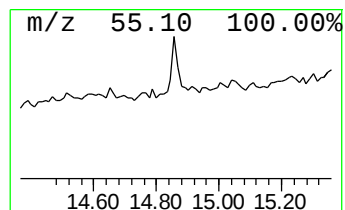
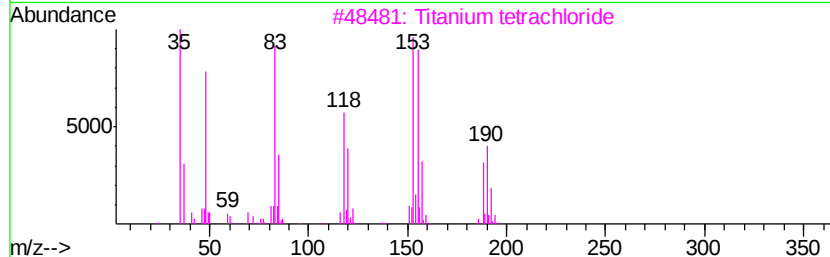
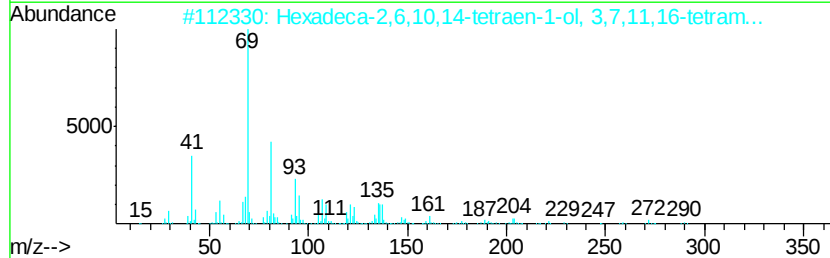
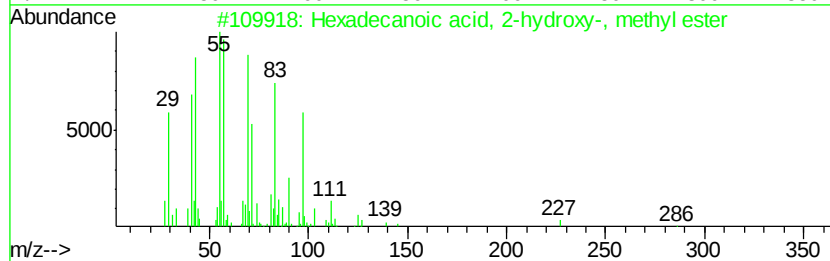
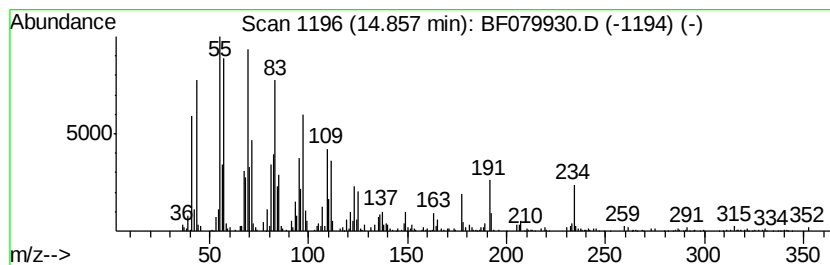
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Hexadecanoic acid, 2-hydrox... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.86	5.55 ng	359993	Chrysene-d12	15.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, 2-hydroxy-, m...	286	C17H34O3	016742-51-1	60
2		Hexadeca-2,6,10,14-tetraen-1-ol,...	290	C20H34O	007614-21-3	56
3		Titanium tetrachloride	188	Cl4Ti	007550-45-0	53
4		Dodecane, 1-cyclopentyl-4-(3-cyc...	348	C25H48	007225-68-5	52
5		Bromoacetic acid, octadecyl ester	390	C20H39BrO2	018992-03-5	52



Data Path : Z:\HPCHEM1\BNA F\DATA\BF061115\
Data File : BF079930.D
Acq On : 12 Jun 2015 9:40
Operator : TP/IZ
Sample : G2449-10
Misc :
ALS Vial : 41 Sample Multiplier: 1

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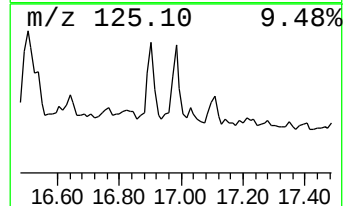
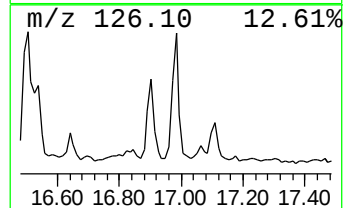
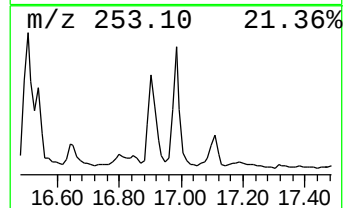
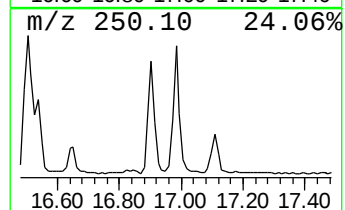
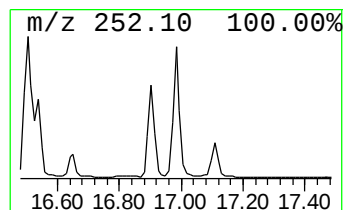
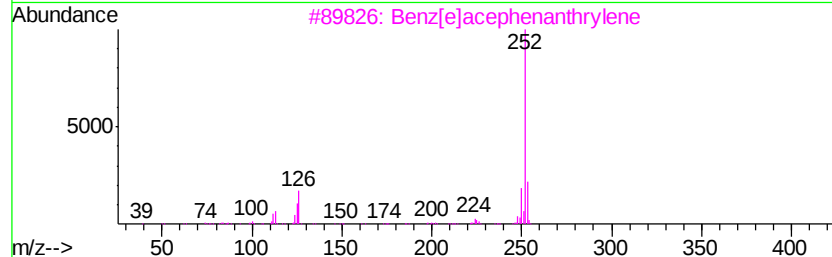
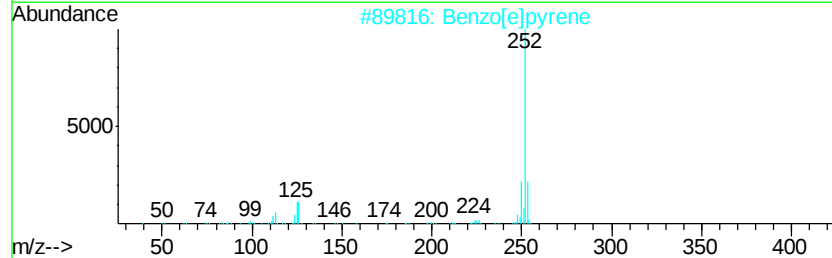
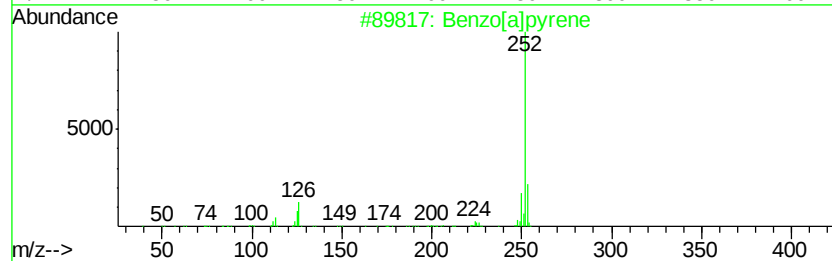
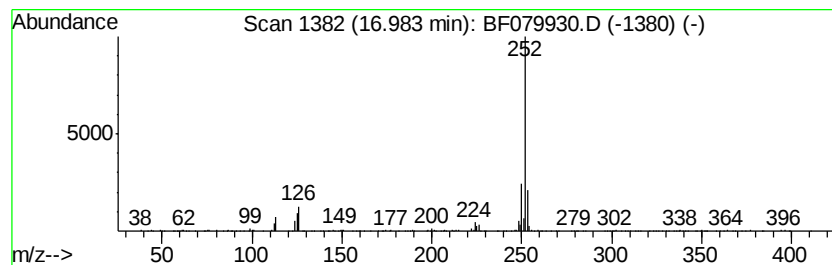
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Benzo[e]pyrene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.98	7.11 ng	218410	Perylene-d12	17.07

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[a]pyrene	252	C20H12	000050-32-8	98
2			Benzo[e]pyrene	252	C20H12	000192-97-2	98
3			Benz[el]acephenanthrylene	252	C20H12	000205-99-2	97
4			Perylene	252	C20H12	000198-55-0	95
5			Benzo[j]fluoranthene	252	C20H12	000205-82-3	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061115\
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Sample : G2449-10
Misc :
ALS Vial : 41 Sample Multiplier: 1

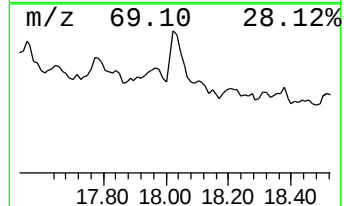
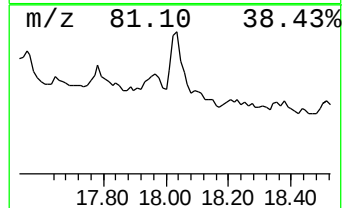
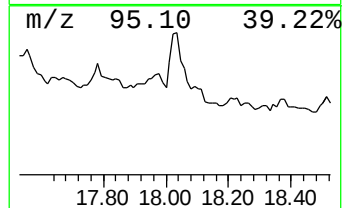
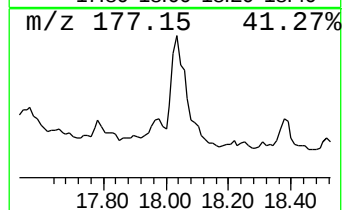
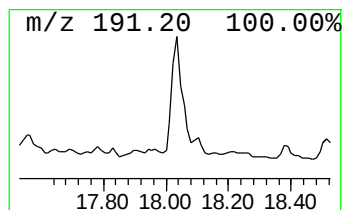
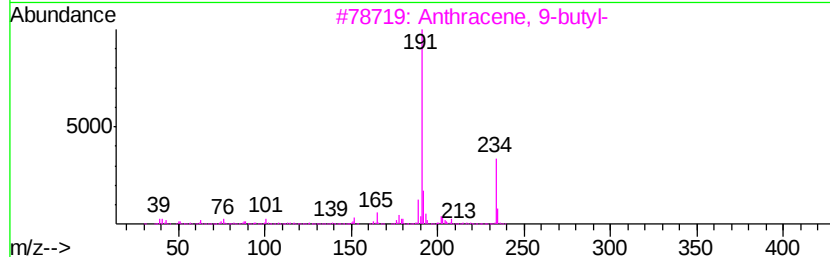
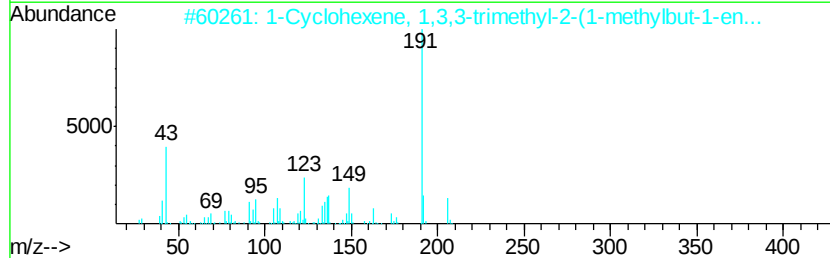
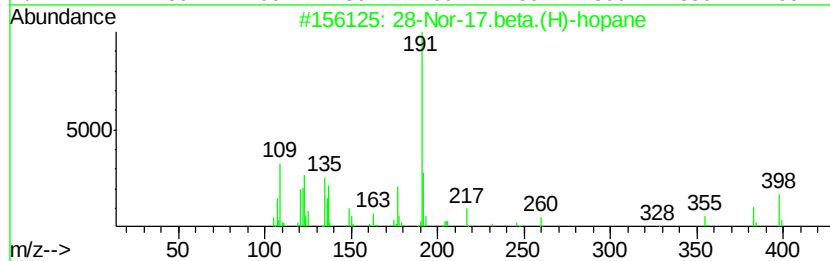
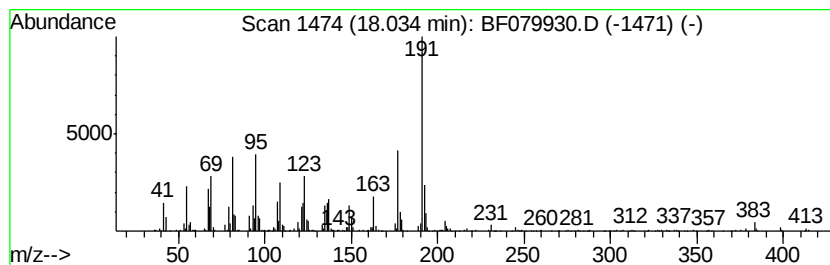
Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 28-Nor-17.beta.(H)-hopane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.03	22.70 ng	697249	Perylene-d12	17.07
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	28-Nor-17.beta.(H)-hopane	398 C29H50	036728-72-0	62
2	1-Cyclohexene, 1,3,3-trimethyl-2...	206 C14H22O	1000197-08-4	37
3	Anthracene, 9-butyl-	234 C18H18	001498-69-7	35
4	2-Amino-4-methyl-6-(2-thienyl)py...	191 C9H9N3S	026963-43-9	27
5	5-(p-Aminophenyl)-2-thiazolamine	191 C9H9N3S	090349-87-4	27



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061115\
Data File : BF079930.D
Acq On : 12 Jun 2015 9:40
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Sample : G2449-10
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ClientSampleId :
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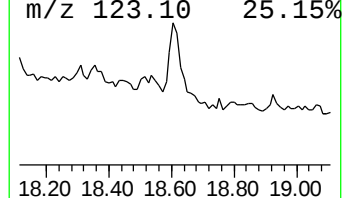
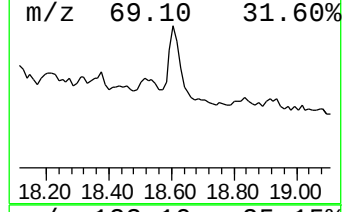
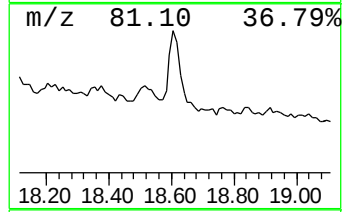
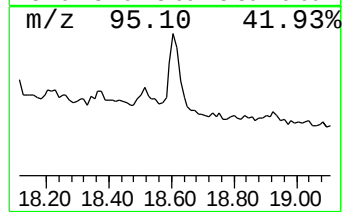
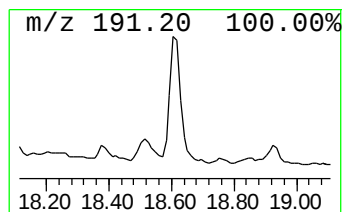
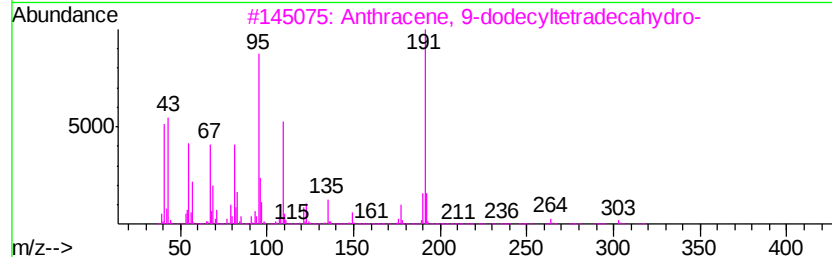
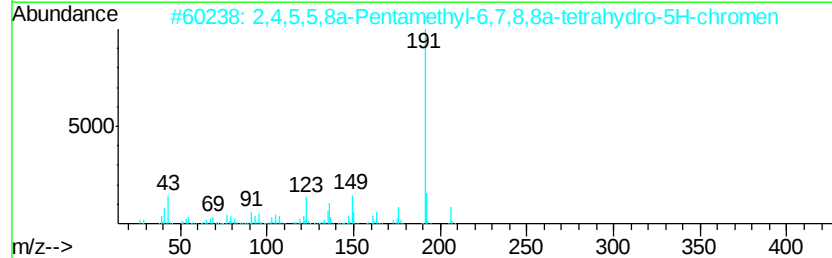
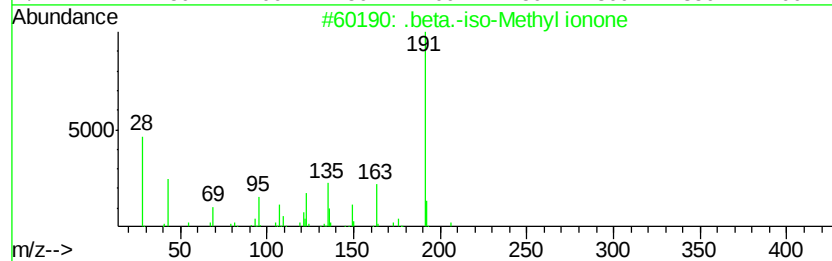
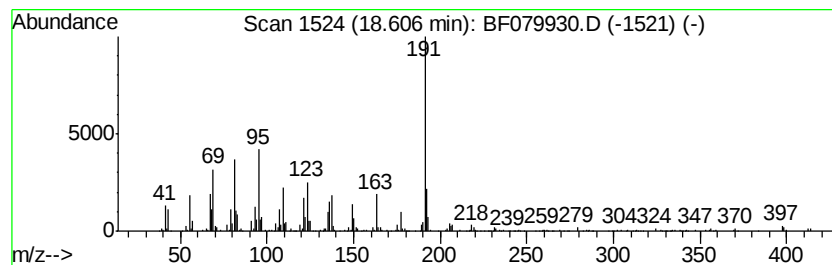
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 .beta.-iso-Methyl ionone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.61	28.96 ng	889756	Perylene-d12	17.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	.beta.-iso-Methyl ionone	206	C14H22O	1000285-40-2	60
2		2,4,5,5,8a-Pentamethyl-6,7,8,8a-...	206	C14H22O	1000195-40-9	47
3		Anthracene, 9-dodecyltetradecahv...	360	C26H48	055401-75-7	42
4		1-Penten-3-one, 1-(2,6,6-trimeth...	206	C14H22O	000127-43-5	38
5		28-Nor-17.beta.(H)-hopane	398	C29H50	036728-72-0	37



Data Path : Z:\HPCHEM1\BNA F\DATA\BF061115\
Data File : BF079930.D
Acq On : 12 Jun 2015 9:40
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Sample : G2449-10
Misc :
ALS Vial : 41 Sample Multiplier: 1

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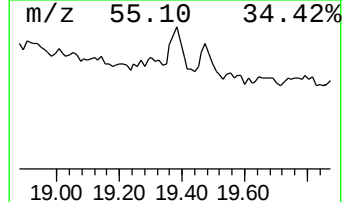
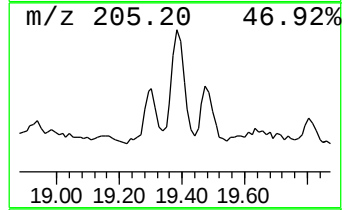
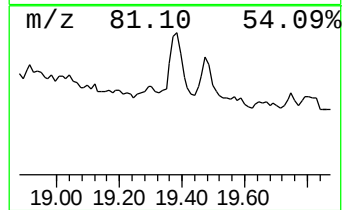
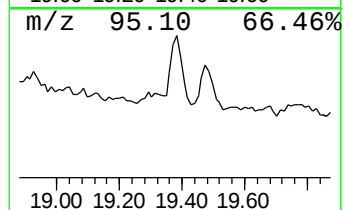
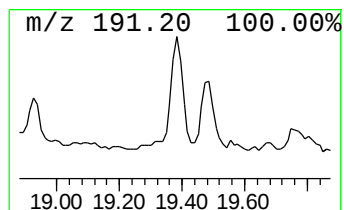
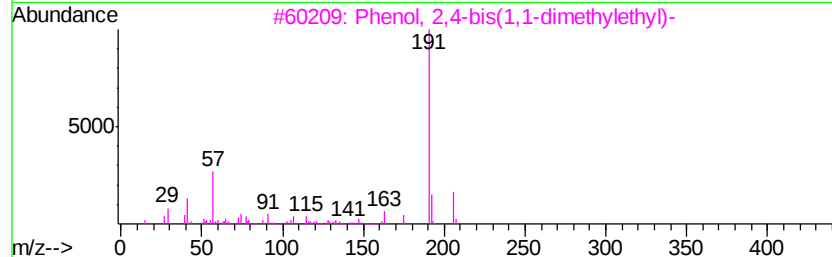
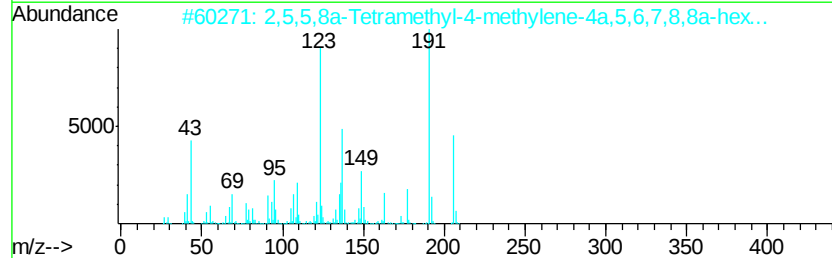
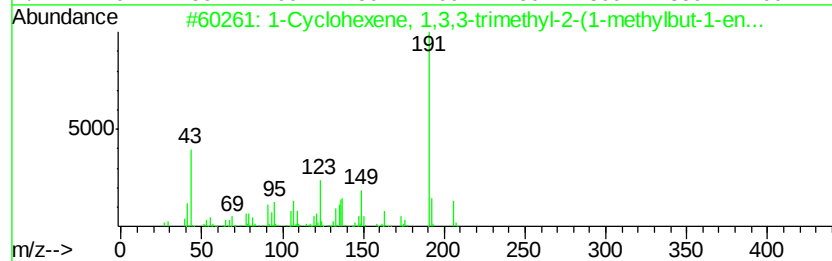
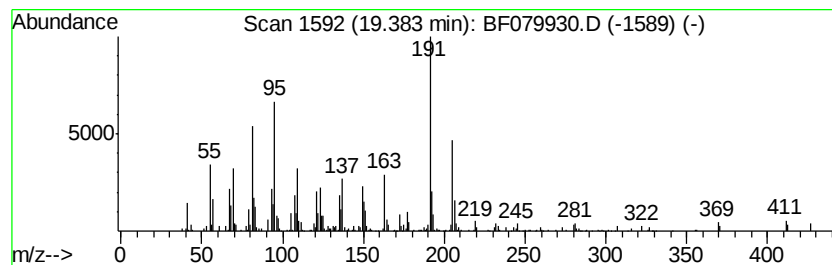
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 unknown19.38 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.38	13.35 ng	410089	Perylene-d12	17.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Cyclohexene, 1,3,3-trimethyl-2...	206	C14H22O	1000197-08-4	43
2		2,5,5,8a-Tetramethyl-4-methylene...	206	C14H22O	093175-76-9	43
3		Phenol, 2,4-bis(1,1-dimethylethyl)-	206	C14H22O	000096-76-4	41
4		Phenol, 3,5-bis(1,1-dimethylethyl)-	206	C14H22O	001138-52-9	35
5		1-Penten-3-one, 1-(2,6,6-trimeth...	206	C14H22O	000127-43-5	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061115\
Data File : BF079930.D
Acq On : 12 Jun 2015 9:40
Operator : TP/IZ
Sample : G2449-10
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
FM7

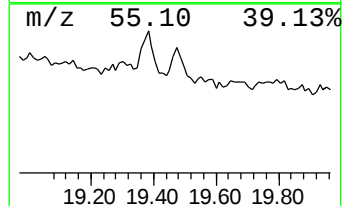
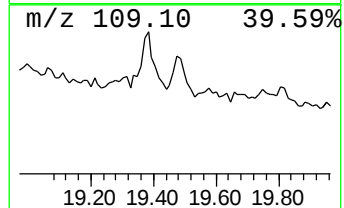
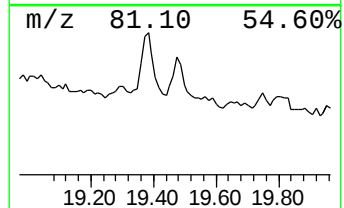
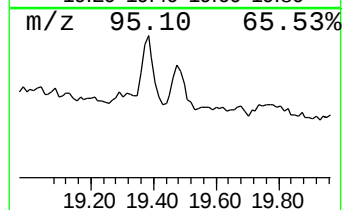
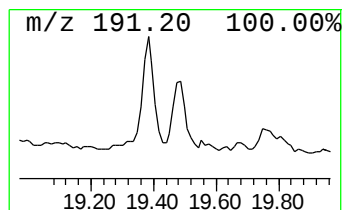
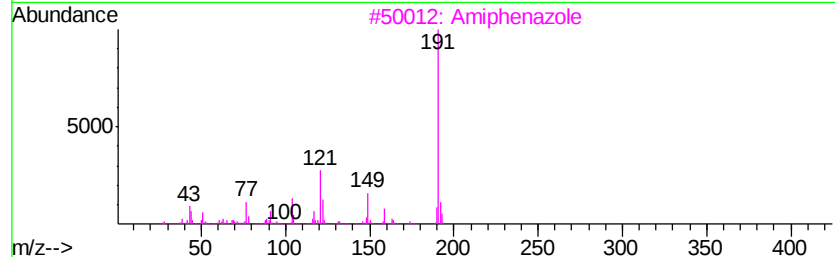
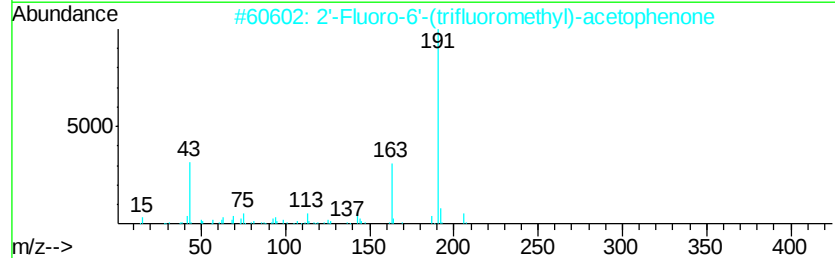
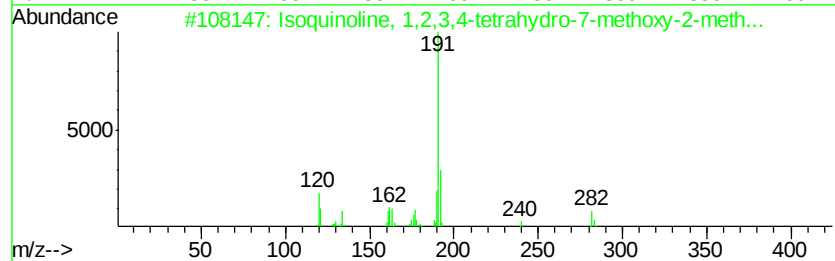
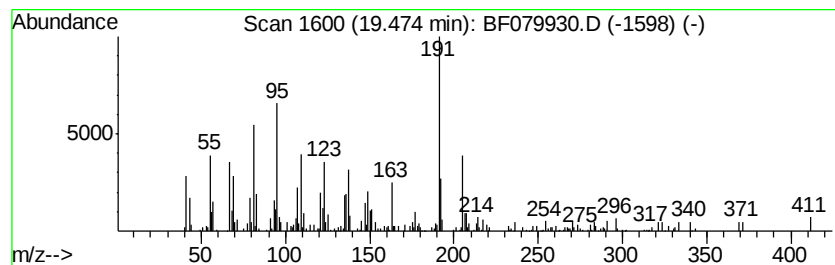
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 unknown19.47 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.47	7.09 ng	217725	Perylene-d12	17.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Isoquinoline, 1,2,3,4-tetrahydro...	283	C18H21N02	036646-87-4	32
2		2'-Fluoro-6'-(trifluoromethyl)-a...	206	C9H6F4O	174013-29-7	22
3		Amiphenazole	191	C9H9N3S	000490-55-1	22
4		5-(p-Aminophenyl)-2-thiazolamine	191	C9H9N3S	090349-87-4	22
5		1-Penten-3-one, 1-(2,6,6-trimeth...	206	C14H22O	000127-43-5	15



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061115\
Data File : BF079930.D
Acq On : 12 Jun 2015 9:40
Operator : TP/IZ
Sample : G2449-10
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
FM7

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown2.27	2.27	268.2	ng	4928220	1	7.39	367511	20.0
Butane, 2-methoxy...	2.70	11.2	ng	205790	1	7.39	367511	20.0
2-Pentanone, 4-hy...	5.61	9.7	ng	178506	1	7.39	367511	20.0
unknown7.16	7.16	78.0	ng	1434140	1	7.39	367511	20.0
Anthracene, 9-met...	12.55	3.0	ng	140639	4	11.98	935664	20.0
unknown12.64	12.64	5.0	ng	234787	4	11.98	935664	20.0
Benzo[kl]xanthene	13.83	3.0	ng	192285	5	15.09	1296290	20.0
Hexadecanoic acid...	14.86	5.5	ng	359993	5	15.09	1296290	20.0
Benzo[e]pyrene	16.98	7.1	ng	218410	6	17.07	614405	20.0
28-Nor-17.beta.(H...	18.03	22.7	ng	697249	6	17.07	614405	20.0
.beta.-iso-Methyl...	18.61	29.0	ng	889756	6	17.07	614405	20.0
unknown19.38	19.38	13.4	ng	410089	6	17.07	614405	20.0
unknown19.47	19.47	7.1	ng	217725	6	17.07	614405	20.0