

Data Path : Z:\HPCHEM1\BNA G\DATA\BG060916\
 Data File : BG022272.D
 Acq On : 9 Jun 2016 17:48
 Operator : UM/SJ
 Sample : H3399-14
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SS-33

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_G\METHODS\8270-BG060616.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	2.984	19	25	45	rVB	1200913	2020169	100.00%	10.554%
2	5.188	392	400	409	rBV2	38325	73801	3.65%	0.386%
3	5.669	474	482	499	rBV	396064	752707	37.26%	3.932%
4	7.279	747	756	771	rBV	433963	863555	42.75%	4.512%
5	7.649	808	819	835	rBV	489592	939274	46.49%	4.907%
6	8.119	892	899	909	rVB	128280	245413	12.15%	1.282%
7	8.443	946	954	968	rBV	377422	735111	36.39%	3.840%
8	9.289	1091	1098	1109	rBV	254100	529205	26.20%	2.765%
9	10.940	1369	1379	1386	rBV	209905	430005	21.29%	2.247%
10	13.372	1783	1793	1803	rBV	861158	1490777	73.79%	7.788%
11	14.065	1902	1911	1918	rBV4	9976	21939	1.09%	0.115%
12	14.195	1924	1933	1938	rBV	72242	121377	6.01%	0.634%
13	14.753	2018	2028	2034	rBV2	313956	567663	28.10%	2.966%
14	14.812	2034	2038	2045	rVB	51770	87707	4.34%	0.458%
15	15.147	2086	2095	2103	rBV2	26033	48885	2.42%	0.255%
16	15.793	2198	2205	2213	rBV2	42851	80852	4.00%	0.422%
17	16.081	2251	2254	2264	rVB	11662	23089	1.14%	0.121%
18	16.239	2274	2281	2288	rBV	510282	853857	42.27%	4.461%
19	16.421	2308	2312	2315	rBV3	16195	25966	1.29%	0.136%
20	17.144	2431	2435	2442	rVB3	21490	38749	1.92%	0.202%
21	17.215	2443	2447	2453	rBV5	27286	47727	2.36%	0.249%
22	17.315	2459	2464	2470	rVB2	27519	50382	2.49%	0.263%
23	17.497	2488	2495	2498	rBV	326832	575639	28.49%	3.007%
24	17.538	2498	2502	2509	rVV	514813	846636	41.91%	4.423%
25	17.626	2512	2517	2523	rVB	102033	168459	8.34%	0.880%
26	17.896	2556	2563	2569	rBV2	43007	95047	4.70%	0.497%
27	17.955	2571	2573	2577	rVB2	18715	22054	1.09%	0.115%
28	18.002	2578	2581	2588	rVV4	12268	20236	1.00%	0.106%
29	18.078	2589	2594	2601	rVB6	12589	28678	1.42%	0.150%
30	18.302	2627	2632	2635	rBV4	10770	21196	1.05%	0.111%
31	18.366	2638	2643	2647	rVV	62123	100901	4.99%	0.527%
32	18.419	2647	2652	2658	rVB2	76784	134509	6.66%	0.703%
33	18.496	2661	2665	2670	rVB2	29048	45351	2.24%	0.237%
34	18.566	2670	2677	2687	rBV3	97748	264934	13.11%	1.384%

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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

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35	18.860	2722	2727	2731	rBV2	44914	69341	3.43%	0.362%
36	18.907	2731	2735	2742	rVB3	40724	65394	3.24%	0.342%
37	19.271	2793	2797	2803	rBV2	35804	70339	3.48%	0.367%
38	19.418	2818	2822	2830	rVB3	32286	59192	2.93%	0.309%
39	19.541	2836	2843	2849	rBV	743351	1165275	57.68%	6.088%
40	19.870	2894	2899	2900	rBV2	104823	145069	7.18%	0.758%
41	19.900	2900	2904	2912	rVV	631774	1000959	49.55%	5.229%
42	19.964	2912	2915	2922	rVB2	30717	44477	2.20%	0.232%
43	20.088	2930	2936	2942	rBV	1064734	1509340	74.71%	7.885%
44	20.487	3001	3004	3008	rBV2	45714	51674	2.56%	0.270%
45	21.157	3114	3118	3124	rVB	44046	78112	3.87%	0.408%
46	21.380	3153	3156	3161	rVB3	66407	103852	5.14%	0.543%
47	21.762	3214	3221	3227	rBV2	405498	1079437	53.43%	5.639%
48	21.821	3227	3231	3241	rVB	232682	462253	22.88%	2.415%
49	24.065	3601	3613	3626	rVB3	191986	864539	42.80%	4.517%

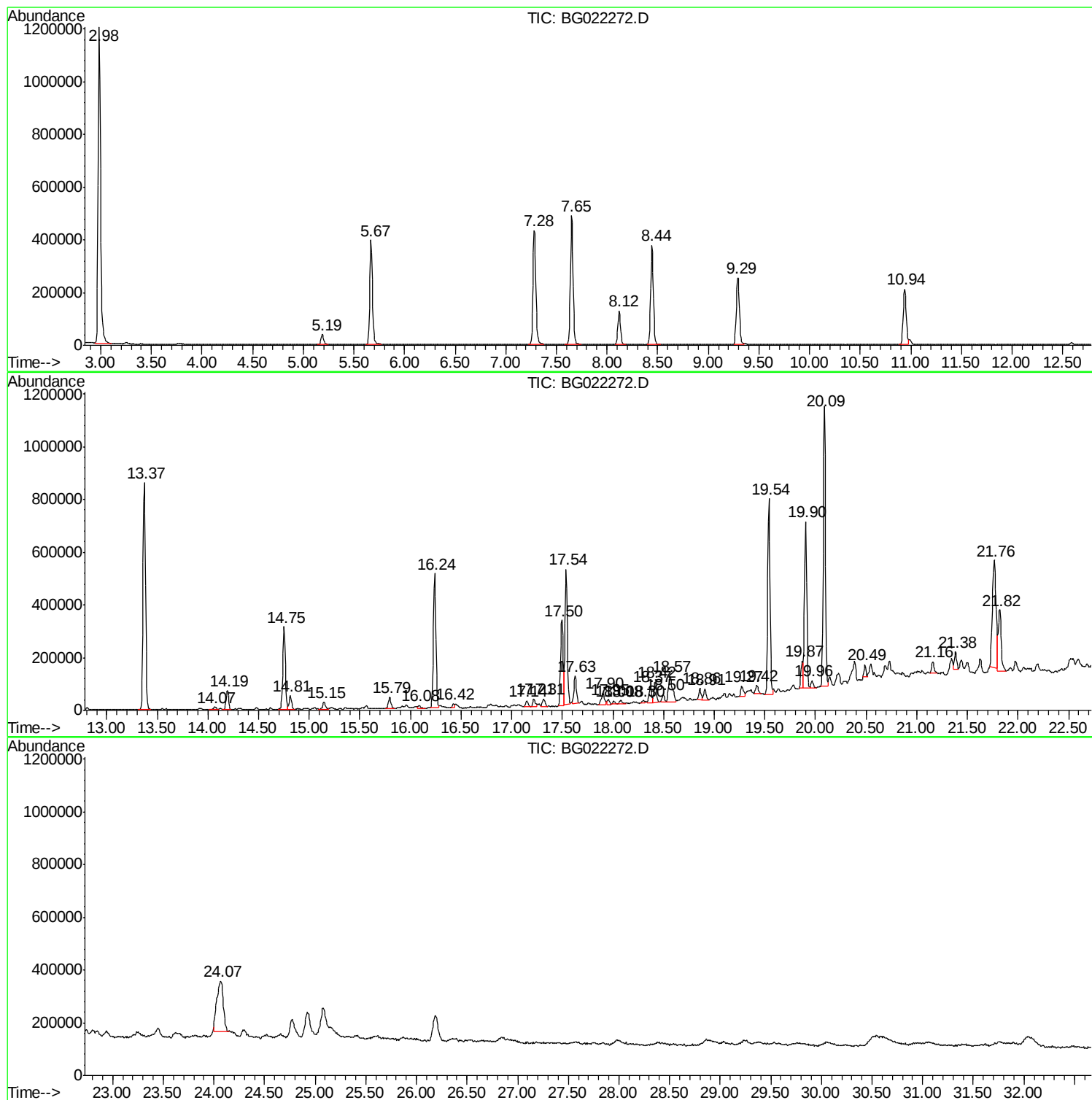
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ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG060616.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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Operator : UM/SJ
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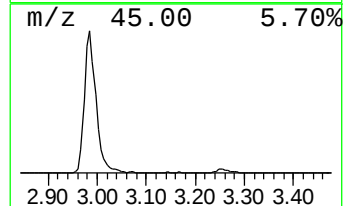
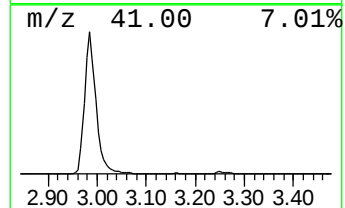
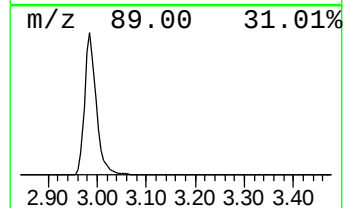
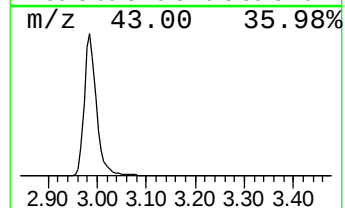
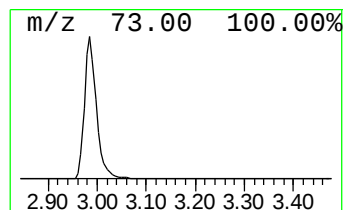
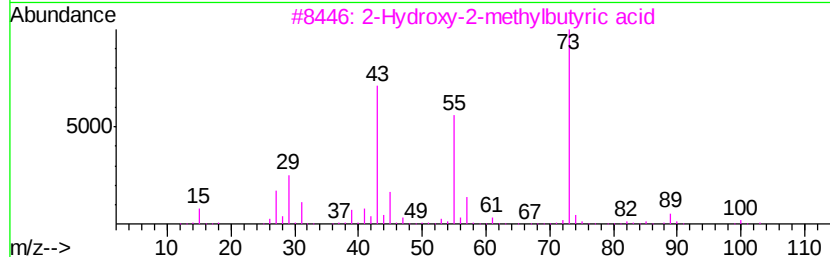
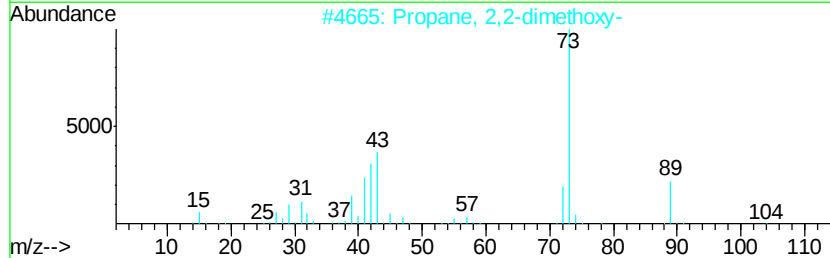
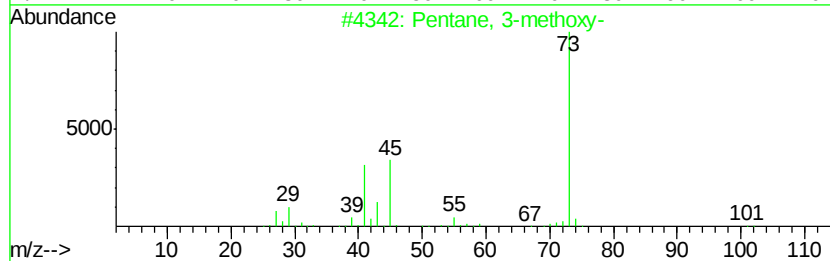
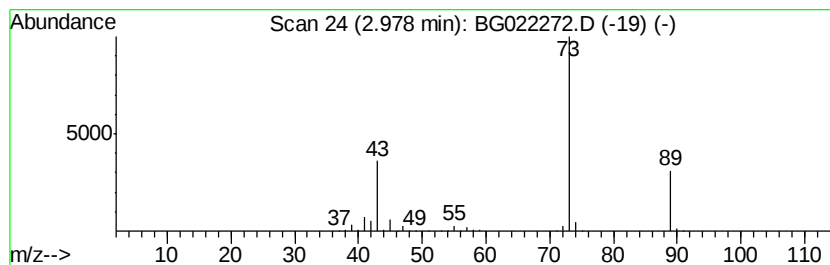
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG060616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown2.98 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.98	164.63 ng	2020170	1,4-Dichlorobenzene-d4	8.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
2		Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	9
3		2-Hydroxy-2-methylbutyric acid	118	C5H10O3	003739-30-8	9
4		Methane, isothiocyanato-	73	C2H3NS	000556-61-6	7
5		N-Ethylformamide	73	C3H7NO	000627-45-2	7



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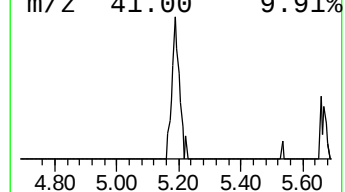
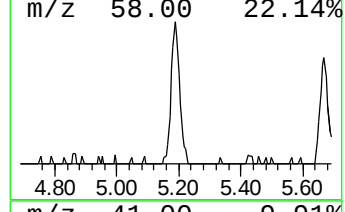
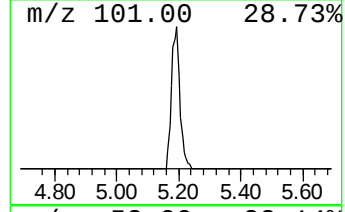
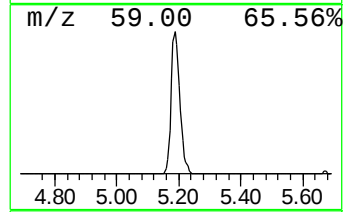
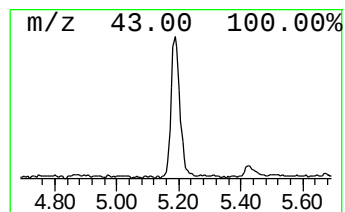
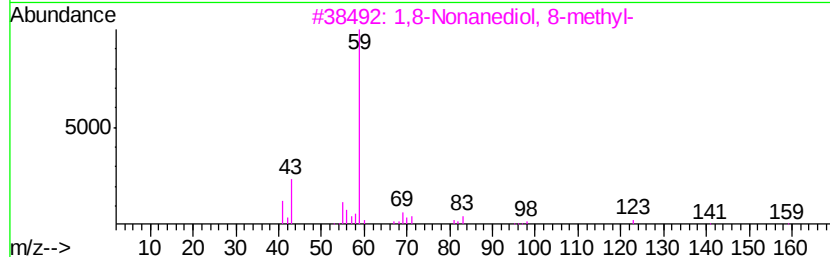
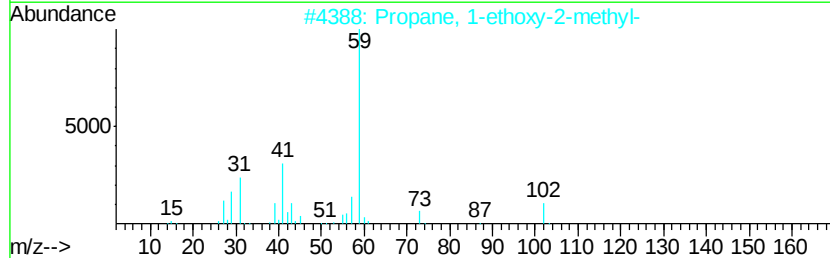
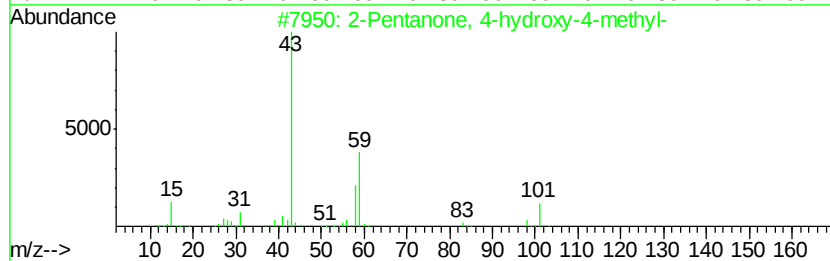
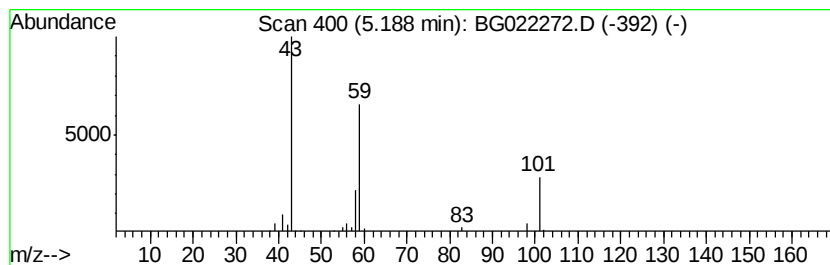
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.19	6.01 ng	73801	1,4-Dichlorobenzene-d4	8.12

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	42
2			Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	35
3			1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	17
4			3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	12
5			Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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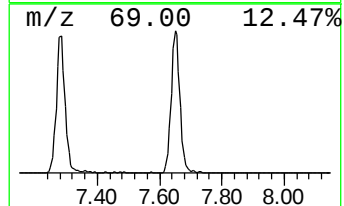
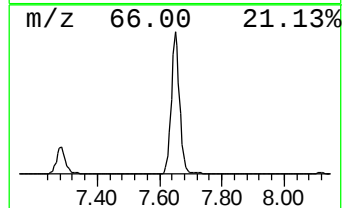
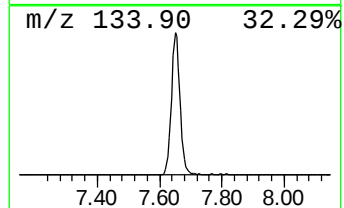
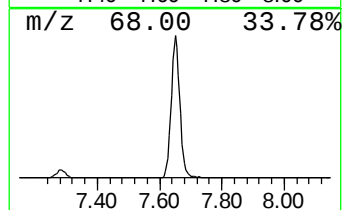
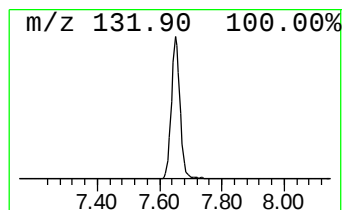
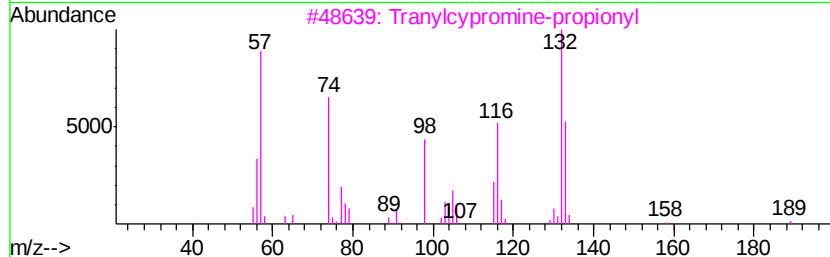
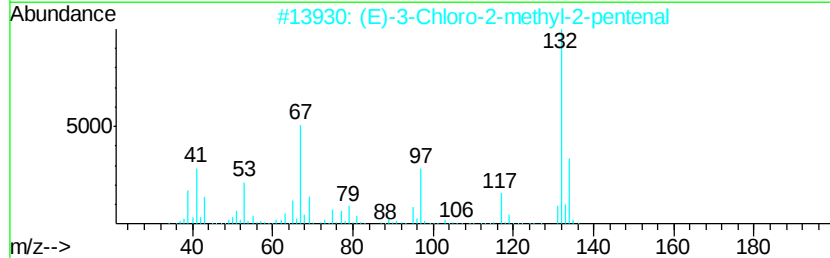
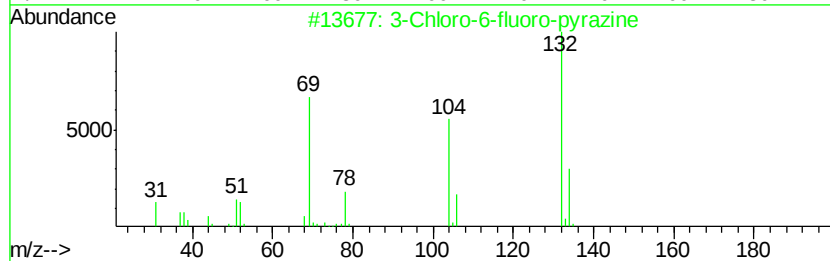
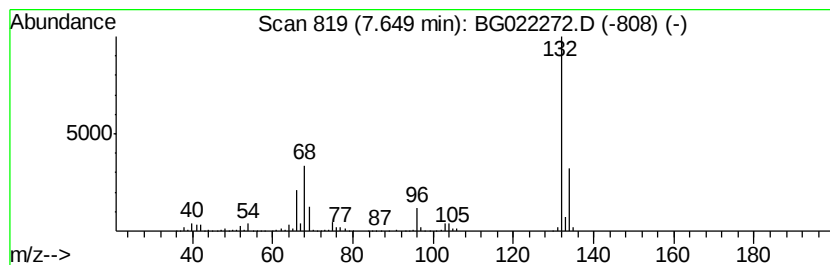
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown7.65 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.65	76.55 ng	939274	1,4-Dichlorobenzene-d4	8.12

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	35
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		Amphetaminil	250	C17H18N2	017590-01-1	12
5		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9



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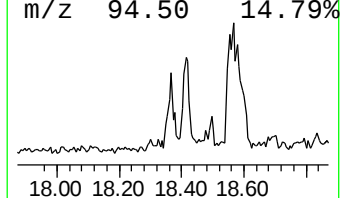
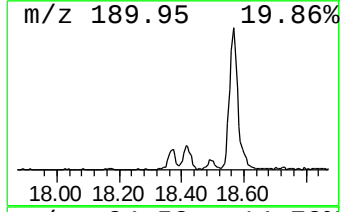
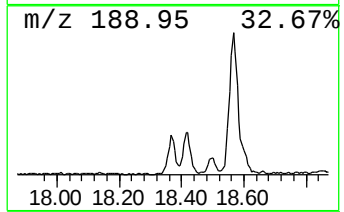
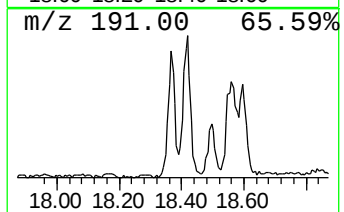
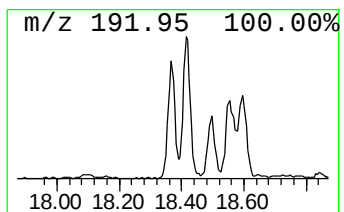
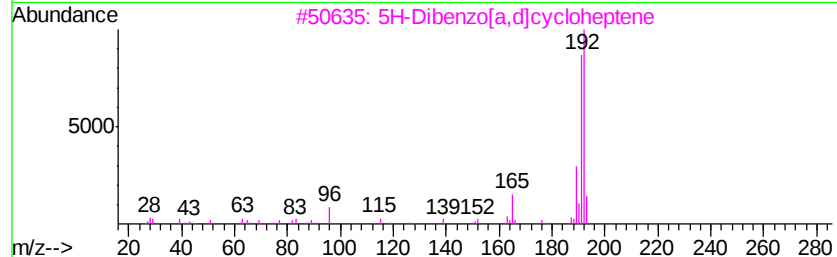
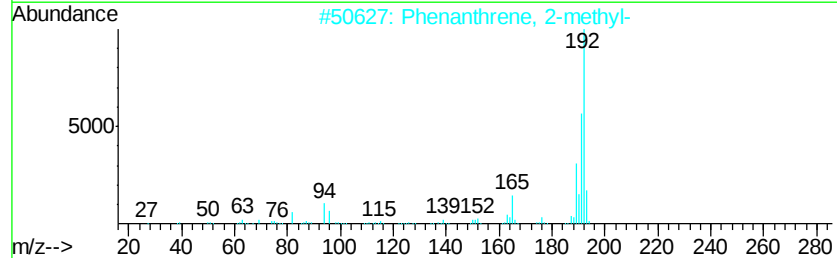
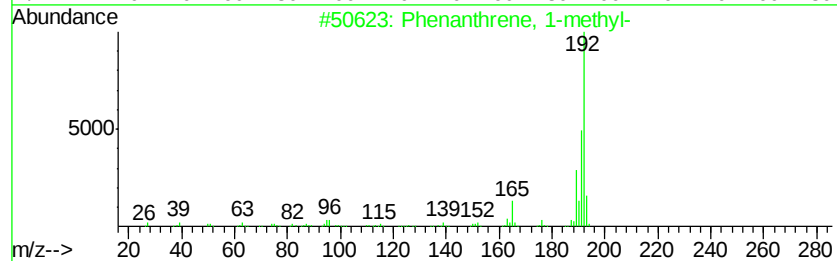
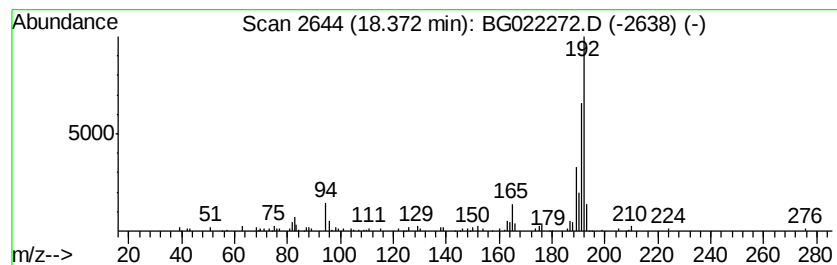
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 5 Phenanthrene, 1-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.37	3.51 ng	100901	Phenanthrene-d10	17.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
2		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	91
3		5H-Dibenzof[a,d]cycloheptene	192	C15H12	000256-81-5	90
4		Anthracene, 2-methyl-	192	C15H12	000613-12-7	90
5		1H-Cyclopropa[1]phenanthrene, 1a,...	192	C15H12	000949-41-7	89



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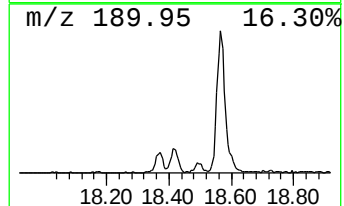
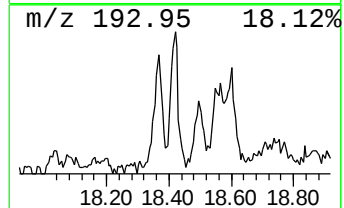
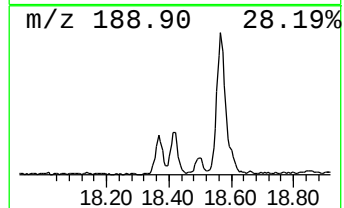
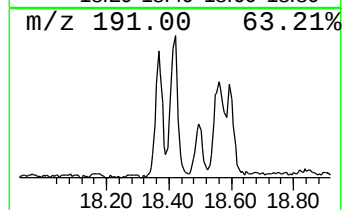
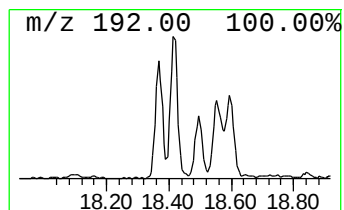
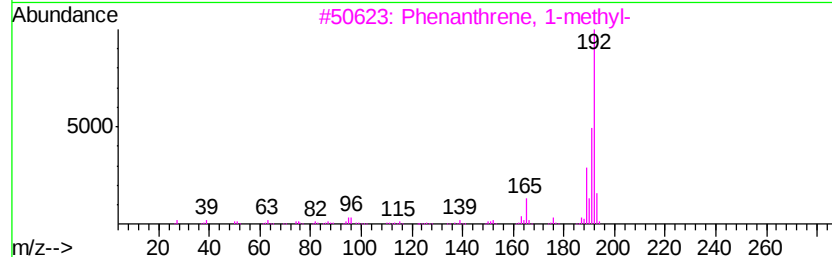
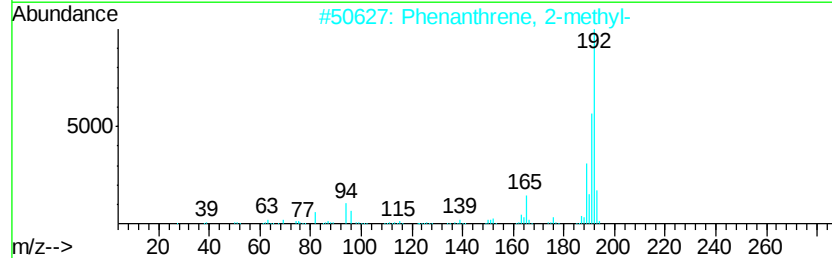
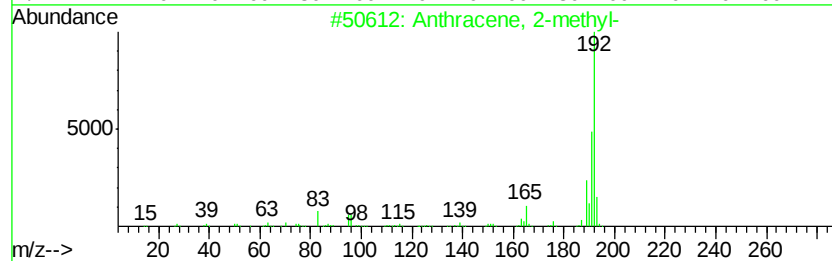
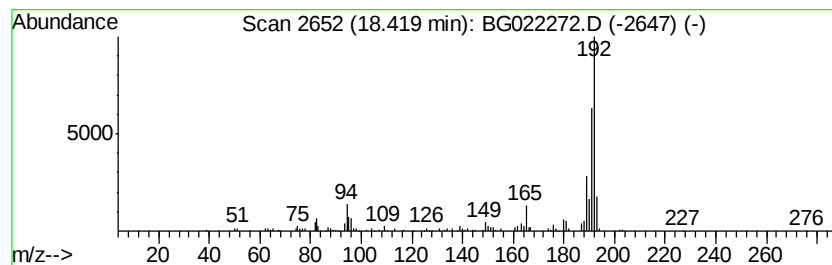
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG060616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Anthracene, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.42	4.67 ng	134509	Phenanthrene-d10	17.50	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 2-methyl-	192	C15H12	000613-12-7	97
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
3	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	94
4	1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	94
5	Anthracene, 1-methyl-	192	C15H12	000610-48-0	93



Data Path : Z:\HPCHEM1\BNA G\DATA\BG060916\
Data File : BG02272.D
Acq On : 9 Jun 2016 17:48
Operator : UM/SJ
Sample : H3399-14
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SS-33

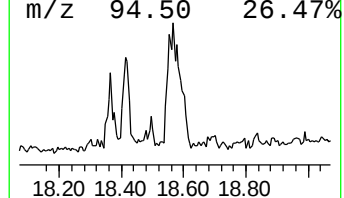
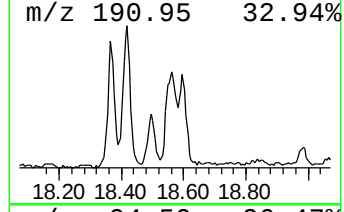
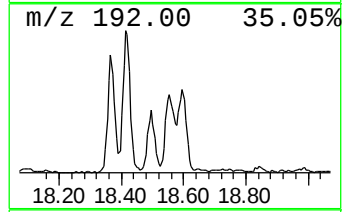
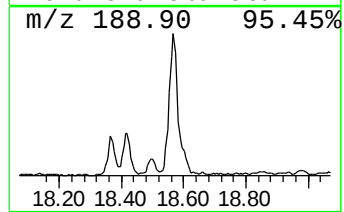
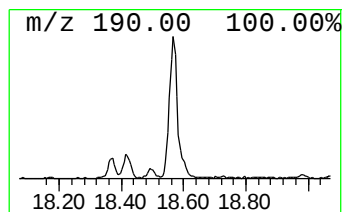
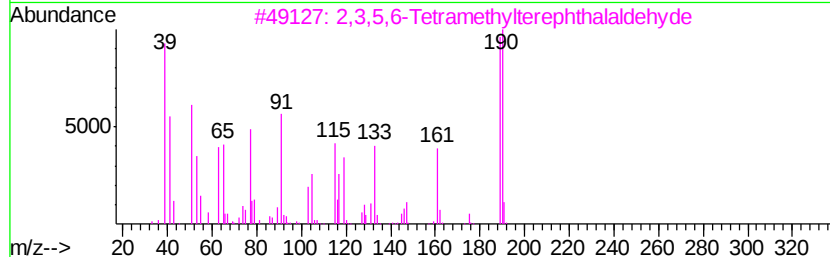
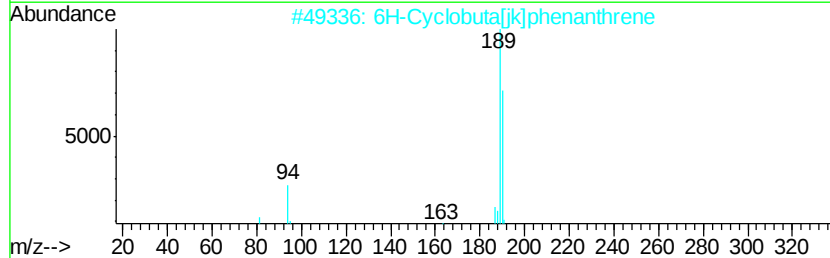
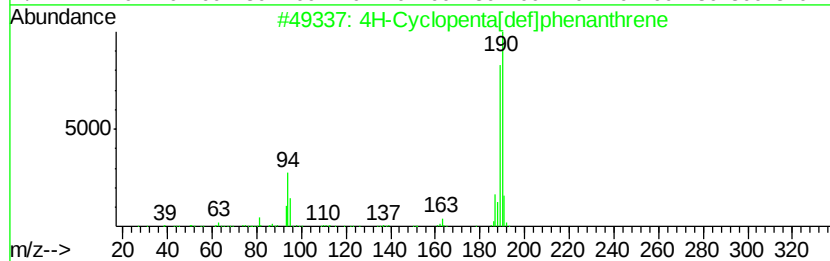
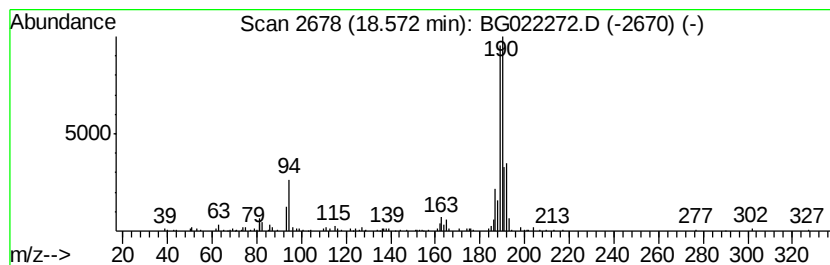
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 4H-Cyclopenta[def]phenanthrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.57	9.20 ng	264934	Phenanthrene-d10	17.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	94
2		6H-Cyclobuta[ik]phenanthrene	190	C15H10	083469-43-6	59
3		2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	47
4		2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	42
5		Hydrazinecarboxaldehyde, 2-[3-(m...	190	C4H6N4O2S	038362-25-3	14



Data Path : Z:\HPCHEM1\BNA G\DATA\BG060916\
Data File : BG022272.D
Acq On : 9 Jun 2016 17:48
Operator : UM/SJ
Sample : H3399-14
Misc :
ALS Vial : 17 Sample Multiplier: 1

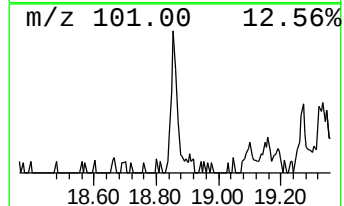
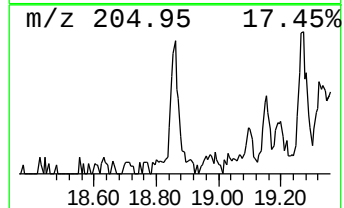
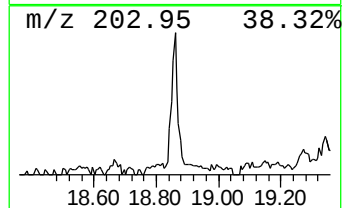
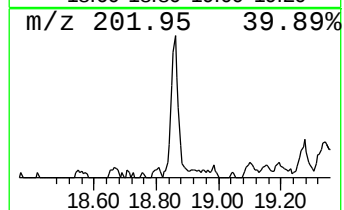
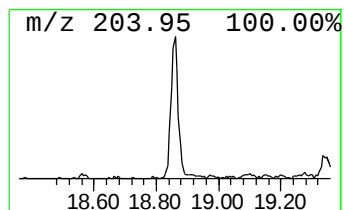
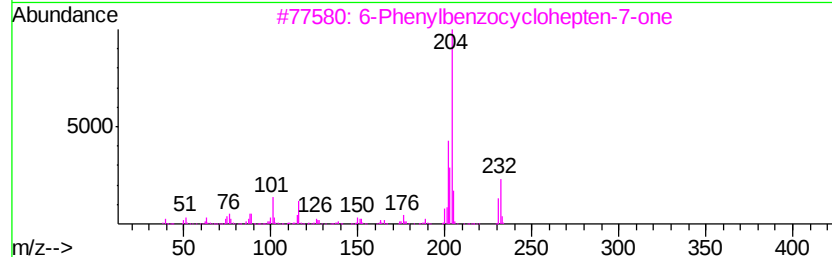
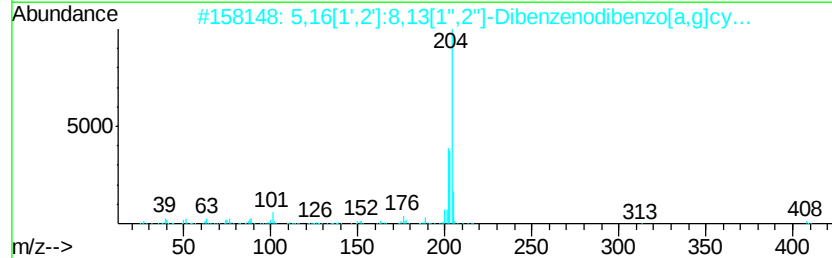
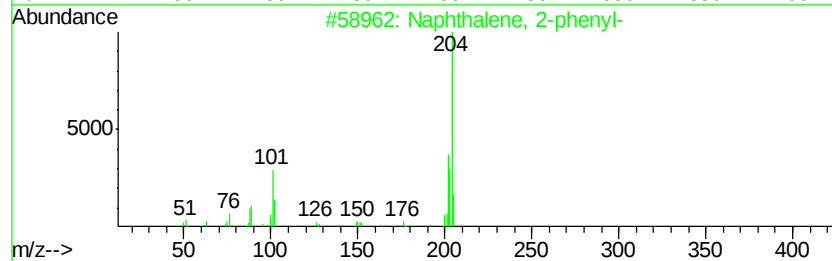
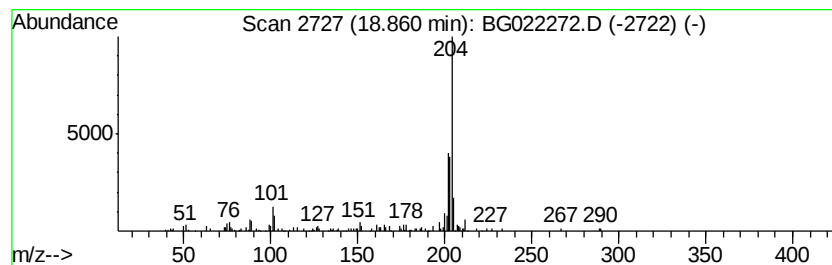
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Naphthalene, 2-phenyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.		
18.86	2.41 ng	69341	Phenanthrene-d10	17.50		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	86
2		5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	78
3		6-Phenylbenzocyclohepten-7-one	232	C17H12O	093327-56-1	64
4		9,10-Bis(bromomethyl)anthracene	362	C16H12Br2	034373-96-1	64
5		Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	58



Data Path : Z:\HPCHEM1\BNA G\DATA\BG060916\
Data File : BG022272.D
Acq On : 9 Jun 2016 17:48
Operator : UM/SJ
Sample : H3399-14
Misc :
ALS Vial : 17 Sample Multiplier: 1

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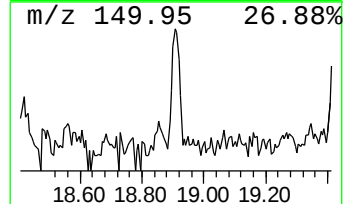
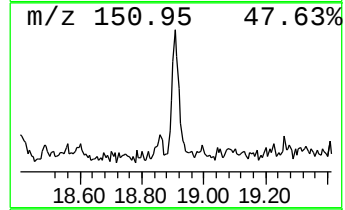
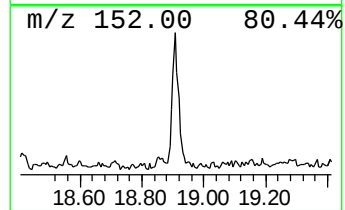
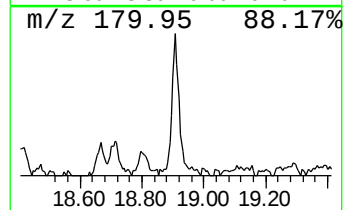
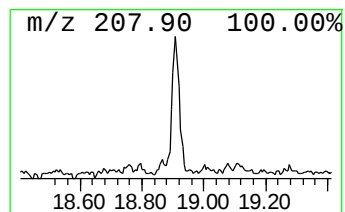
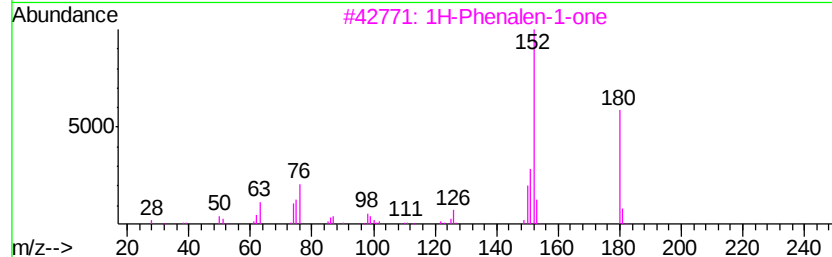
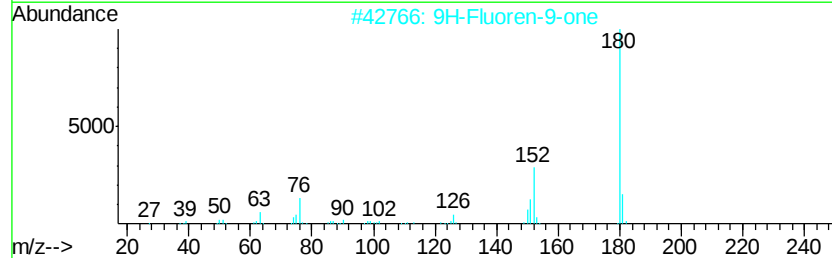
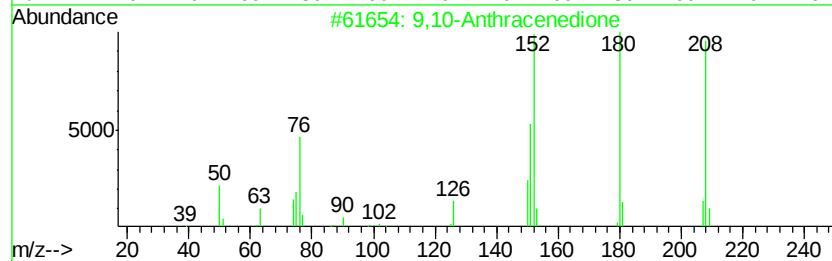
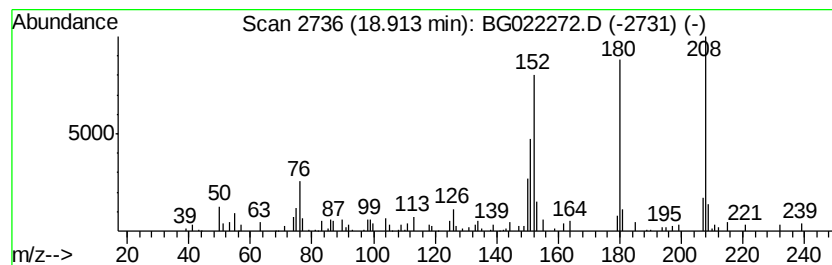
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 9,10-Anthracenedione Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.91	2.27 ng	65394	Phenanthrene-d10	17.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9,10-Anthracenedione	208	C14H8O2	000084-65-1	94
2		9H-Fluoren-9-one	180	C13H8O	000486-25-9	91
3		1H-Phenalen-1-one	180	C13H8O	000548-39-0	64
4		Benzo[h]cinnoline	180	C12H8N2	000230-31-9	64
5		Benzo[c]cinnoline	180	C12H8N2	000230-17-1	62



Data Path : Z:\HPCHEM1\BNA G\DATA\BG060916\
Data File : BG02272.D
Acq On : 9 Jun 2016 17:48
Operator : UM/SJ
Sample : H3399-14
Misc :
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Instrument :
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ClientSampleId :
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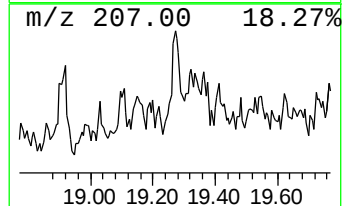
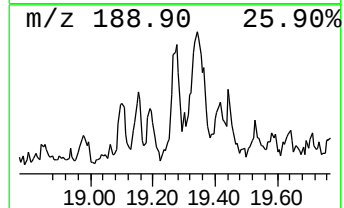
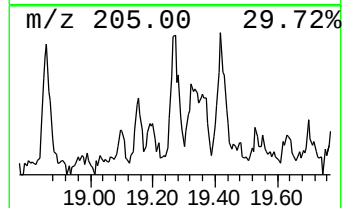
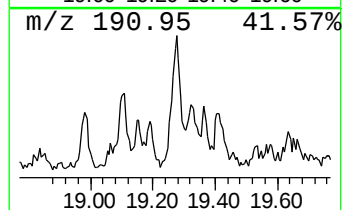
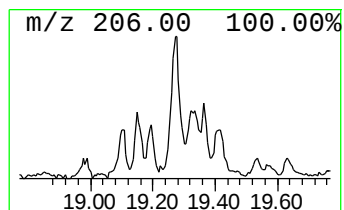
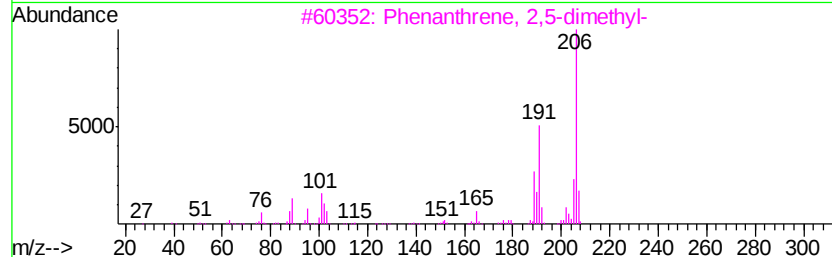
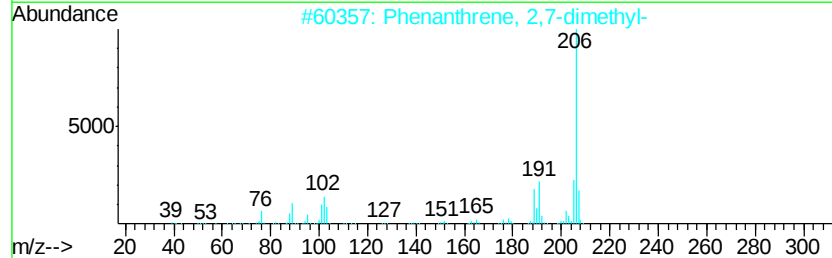
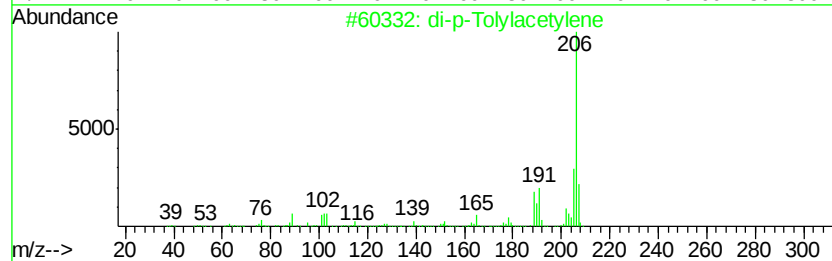
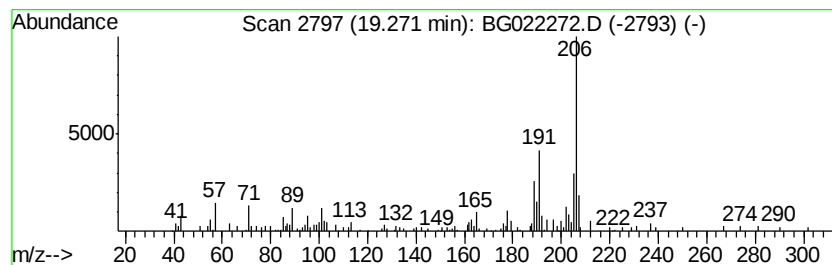
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 di-p-Tolylacetylene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.27	2.44 ng	70339	Phenanthrene-d10	17.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	di-p-Tolylacetylene	206	C16H14	002789-88-0	93
2		Phenanthrene, 2,7-dimethyl-	206	C16H14	001576-69-8	91
3		Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	91
4		Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	91
5		Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	91



Data Path : Z:\HPCHEM1\BNA G\DATA\BG060916\
Data File : BG02272.D
Acq On : 9 Jun 2016 17:48
Operator : UM/SJ
Sample : H3399-14
Misc :
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Instrument :
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ClientSampled :
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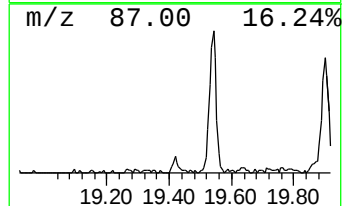
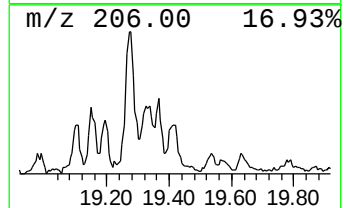
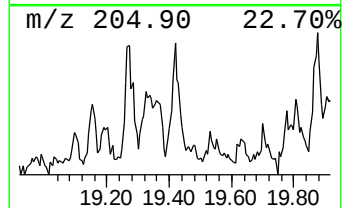
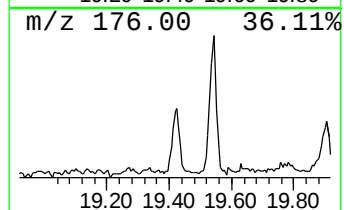
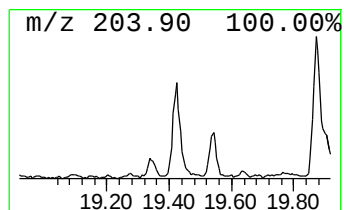
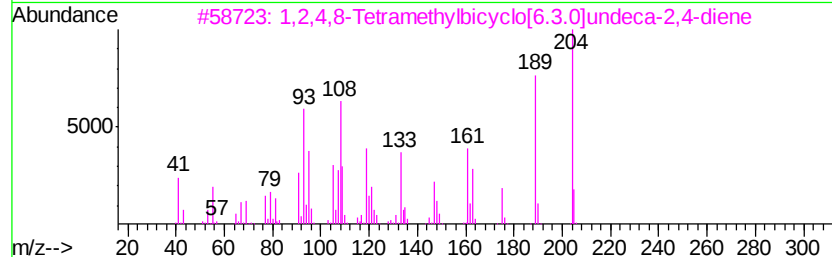
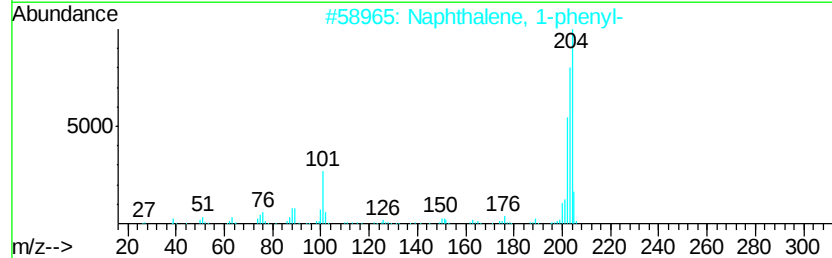
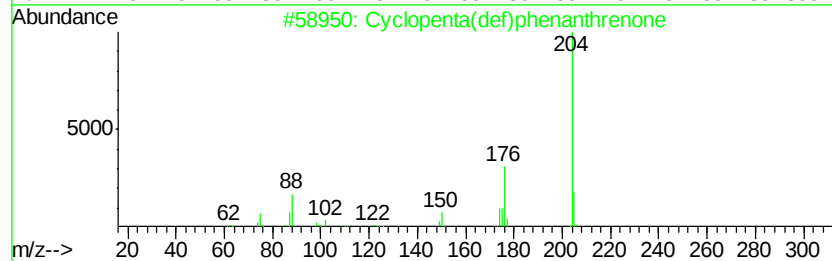
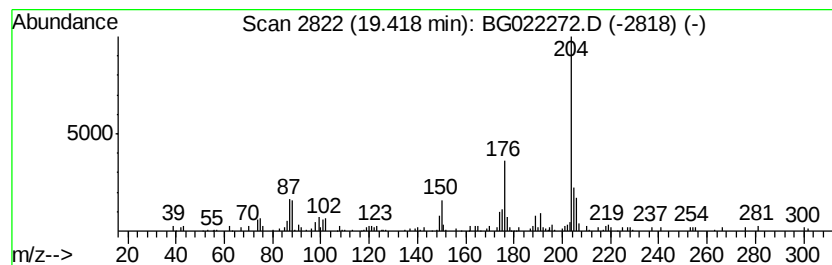
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Cyclopenta(def)phenanthrenone Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.42	2.06 ng	59192	Phenanthrene-d10	17.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	87
2		Naphthalene, 1-phenyl-	204	C16H12	000605-02-7	50
3		1,2,4,8-Tetramethylbicyclo[6.3.0...	204	C15H24	137235-51-9	49
4		.alpha.-l-Mannopyranoside, methy...	394	C16H38O5Si3	056271-60-4	47
5		[1,1'-Biphenyl]-2,2'-dicarbonitrile	204	C14H8N2	004341-02-0	47



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060916\
Data File : BG022272.D
Acq On : 9 Jun 2016 17:48
Operator : UM/SJ
Sample : H3399-14
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060616.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
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2-Pentanone, 4-hy...	5.19	6.0	ng	73801	1	8.12	245413	20.0
unknown7.65	7.65	76.5	ng	939274	1	8.12	245413	20.0
Phenanthrene, 1-m...	18.37	3.5	ng	100901	4	17.50	575639	20.0
Anthracene, 2-met...	18.42	4.7	ng	134509	4	17.50	575639	20.0
4H-Cyclopenta[def...	18.57	9.2	ng	264934	4	17.50	575639	20.0
Naphthalene, 2-ph...	18.86	2.4	ng	69341	4	17.50	575639	20.0
9,10-Anthracenedione	18.91	2.3	ng	65394	4	17.50	575639	20.0
di-p-Tolylacetylene	19.27	2.4	ng	70339	4	17.50	575639	20.0
Cyclopenta(def)ph...	19.42	2.1	ng	59192	4	17.50	575639	20.0