

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021920\
 Data File : BG044504.D
 Acq On : 20 Feb 2020 00:39
 Operator : CG/JU
 Sample : L1547-07
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MARCY-GREEN-BH-02-A

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:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG013020.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	5.467	399	405	427	rBV	794786	1336468	40.48%	5.324%
2	7.065	667	677	690	rBV	783733	1393452	42.20%	5.551%
3	7.888	809	817	826	rBV	197442	336701	10.20%	1.341%
4	8.211	863	872	882	rBV	664532	1140161	34.53%	4.542%
5	9.045	1004	1014	1026	rBV	464631	868643	26.31%	3.460%
6	10.690	1286	1294	1309	rBV	250429	479222	14.51%	1.909%
7	13.152	1706	1713	1730	rBV	1281299	2050064	62.09%	8.166%
8	13.986	1848	1855	1860	rBV	30599	51411	1.56%	0.205%
9	14.527	1939	1947	1954	rBV2	423911	688204	20.84%	2.741%
10	14.591	1954	1958	1966	rVB	20771	34396	1.04%	0.137%
11	15.579	2118	2126	2137	rBV2	21146	38052	1.15%	0.152%
12	16.019	2193	2201	2226	rBV	1110662	1821395	55.17%	7.255%
13	17.094	2380	2384	2393	rVB	20510	35708	1.08%	0.142%
14	17.271	2408	2414	2418	rBV	578921	900152	27.26%	3.586%
15	17.318	2418	2422	2428	rVV	313344	494472	14.98%	1.970%
16	17.406	2428	2437	2443	rVB	61905	109840	3.33%	0.438%
17	17.676	2477	2483	2488	rBV2	17394	34957	1.06%	0.139%
18	18.152	2557	2564	2568	rBV	61824	97292	2.95%	0.388%
19	18.199	2568	2572	2577	rVB	78201	111005	3.36%	0.442%
20	18.281	2582	2586	2590	rVV	24756	37237	1.13%	0.148%
21	18.340	2590	2596	2601	rVV3	85938	175826	5.33%	0.700%
22	18.381	2601	2603	2610	rVB	50773	60807	1.84%	0.242%
23	18.651	2644	2649	2652	rBV	40749	60961	1.85%	0.243%
24	18.687	2652	2655	2661	rVB3	31348	50066	1.52%	0.199%
25	19.069	2710	2720	2723	rBV2	54436	106698	3.23%	0.425%
26	19.204	2738	2743	2750	rBV3	37730	68265	2.07%	0.272%
27	19.327	2758	2764	2772	rVB	547184	765116	23.17%	3.048%
28	19.685	2816	2825	2834	rBV	518331	883574	26.76%	3.520%
29	19.762	2834	2838	2841	rBV2	24989	41921	1.27%	0.167%
30	19.797	2841	2844	2850	rVB	49720	64042	1.94%	0.255%
31	19.891	2852	2860	2865	rBV	2521483	3301640	100.00%	13.152%
32	20.014	2875	2881	2888	rBV4	39144	104738	3.17%	0.417%
33	20.150	2899	2904	2906	rBV3	34524	62139	1.88%	0.248%
34	20.179	2906	2909	2920	rVB3	72491	125673	3.81%	0.501%

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35	20.279	2923	2926	2930	rVV	27663	35871	1.09%	0.143%
36	20.338	2931	2936	2941	rVV2	50141	85793	2.60%	0.342%
37	20.479	2956	2960	2963	rBV	34815	46661	1.41%	0.186%
38	20.526	2964	2968	2972	rVB	40113	56979	1.73%	0.227%
39	20.737	3000	3004	3009	rBV	43008	64509	1.95%	0.257%
40	20.931	3033	3037	3045	rVB	49567	83942	2.54%	0.334%
41	21.113	3061	3068	3071	rBV3	47882	100674	3.05%	0.401%
42	21.148	3071	3074	3077	rVV2	45323	70520	2.14%	0.281%
43	21.190	3077	3081	3088	rVV3	283707	517974	15.69%	2.063%
44	21.254	3088	3092	3099	rVB2	75562	135148	4.09%	0.538%
45	21.419	3117	3120	3124	rVB	45515	53563	1.62%	0.213%
46	21.513	3128	3136	3141	rVV3	720948	1527594	46.27%	6.085%
47	21.560	3141	3144	3158	rVB	230380	374268	11.34%	1.491%
48	21.718	3166	3171	3177	rVB3	86461	149790	4.54%	0.597%
49	22.218	3254	3256	3263	rVB	42778	64201	1.94%	0.256%
50	22.306	3265	3271	3280	rBV4	113425	252195	7.64%	1.005%
51	22.970	3378	3384	3390	rVB	106903	206050	6.24%	0.821%
52	23.622	3488	3495	3503	rBV	157251	491931	14.90%	1.960%
53	23.734	3509	3514	3522	rVB2	116993	228498	6.92%	0.910%
54	24.309	3605	3612	3624	rVB2	102739	297538	9.01%	1.185%
55	24.445	3628	3635	3648	rVB2	144918	397348	12.03%	1.583%
56	24.597	3651	3661	3680	rBV	404515	1412393	42.78%	5.626%
57	25.714	3844	3851	3862	rVB3	79583	221776	6.72%	0.883%
58	28.070	4246	4252	4272	rVB3	67389	298344	9.04%	1.188%

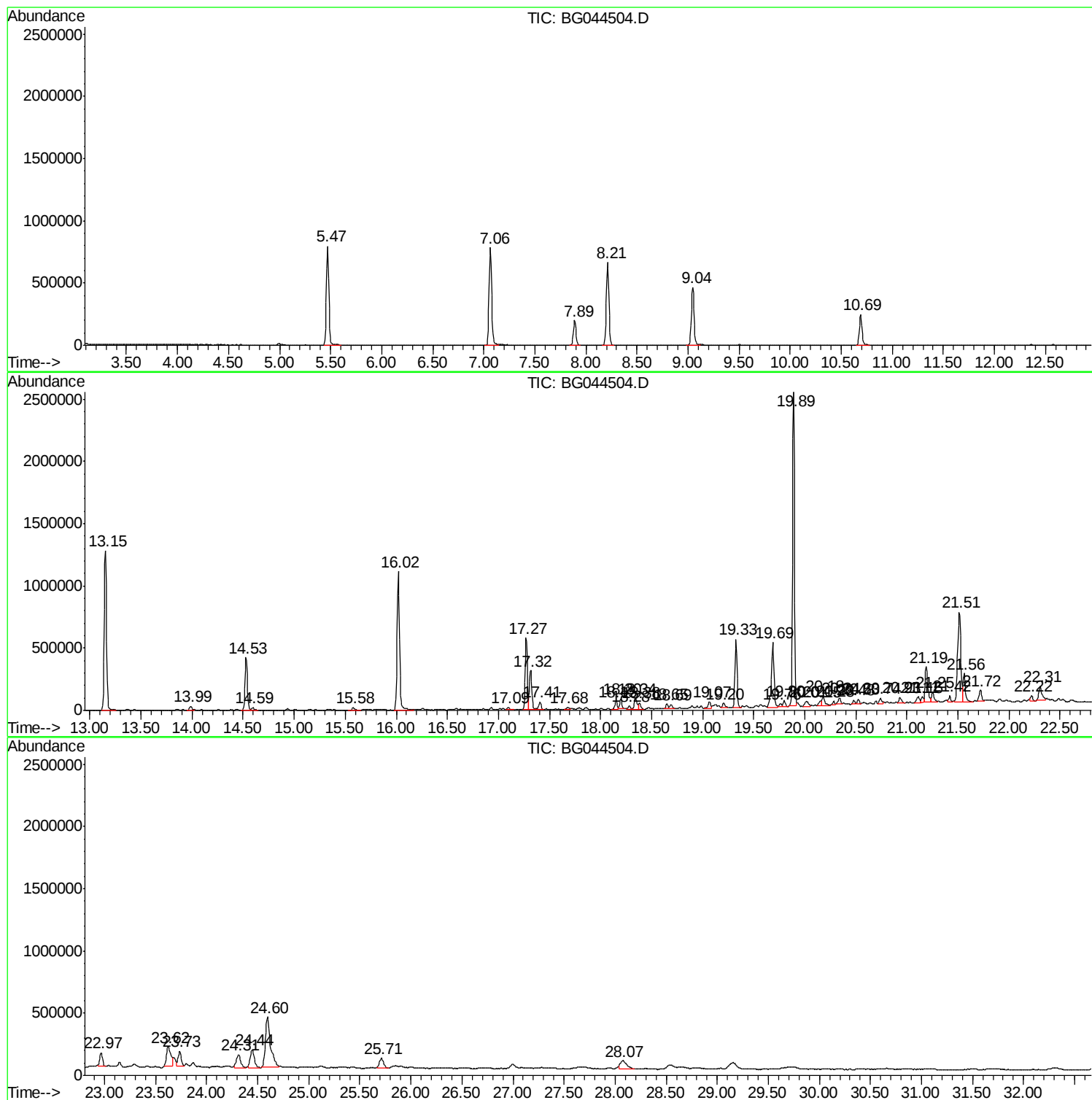
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG013020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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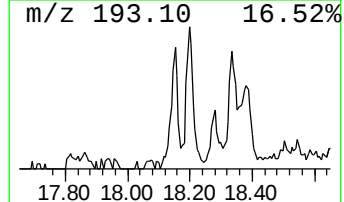
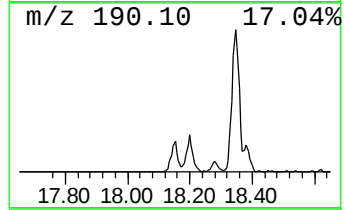
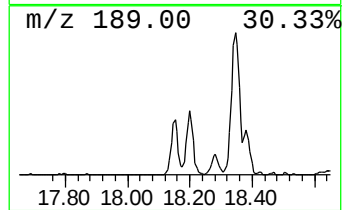
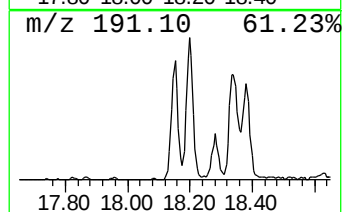
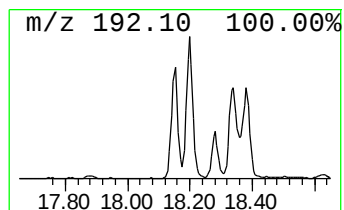
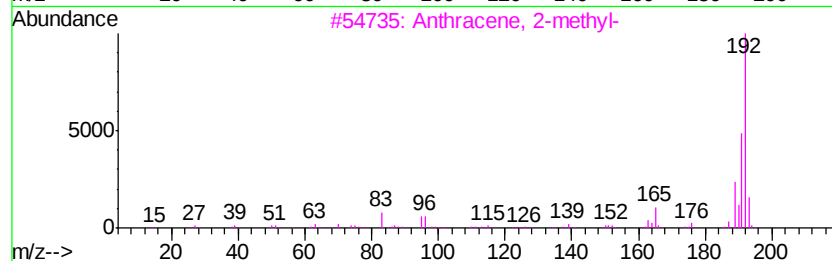
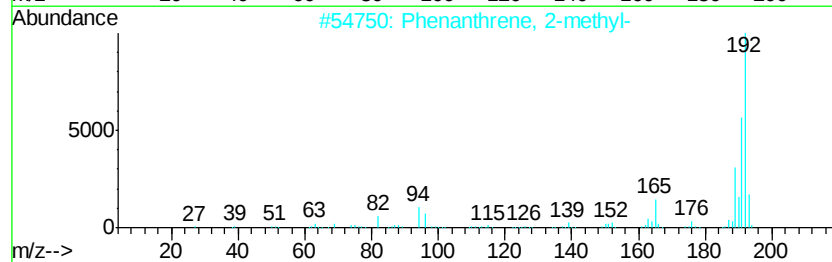
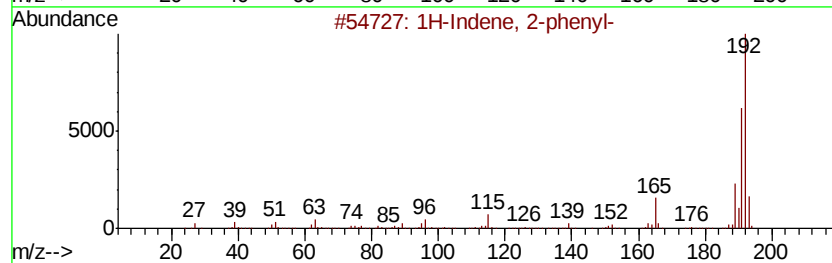
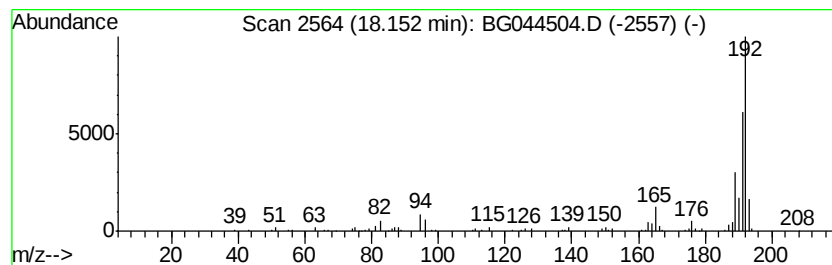
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 1H-Indene, 2-phenyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.15	2.16 ng	97292	Phenanthrene-d10	17.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	94
2		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
3		Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
4		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
5		Naphtho[2,3-b]norbornadiene	192	C15H12	107426-38-0	91



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Misc :
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Instrument :
BNA_G
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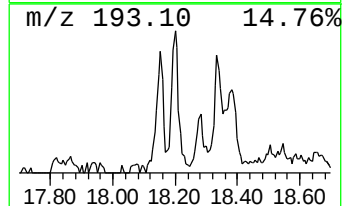
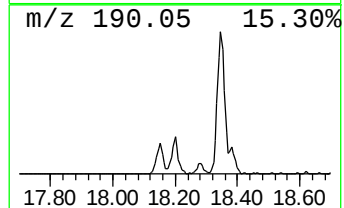
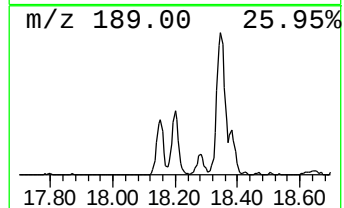
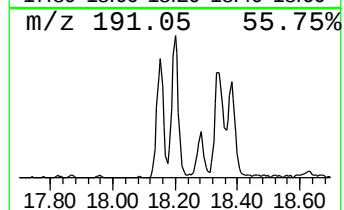
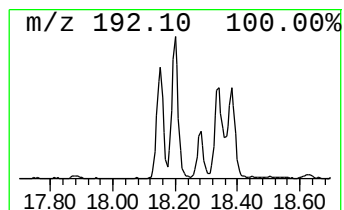
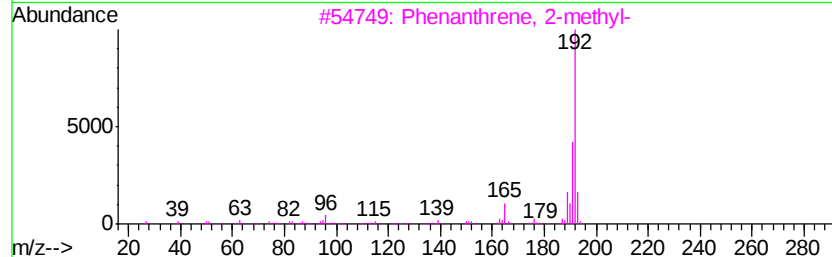
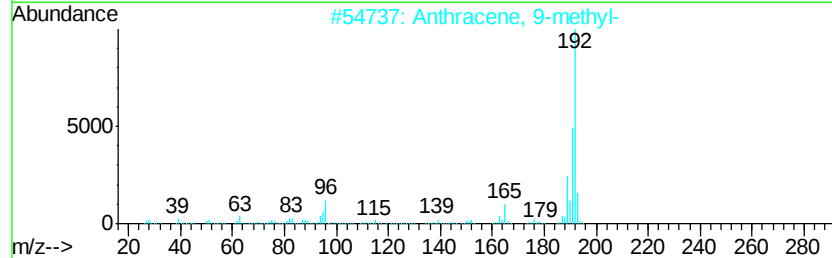
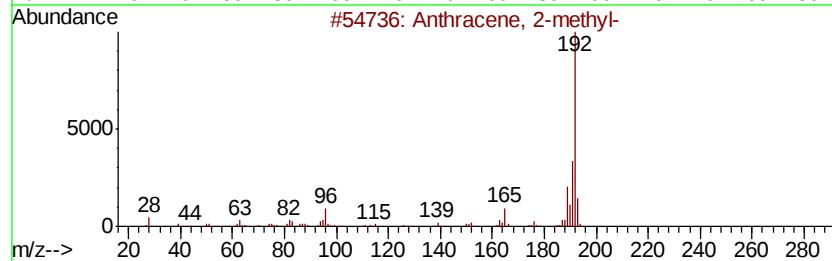
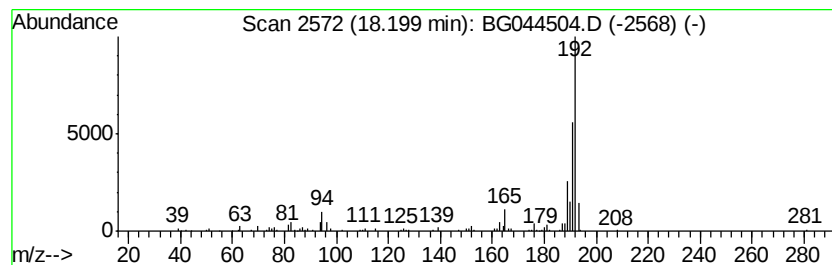
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TIC Library : C:\Database\NIST11.L
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Peak Number 3 Anthracene, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.20	2.47 ng	111005	Phenanthrene-d10	17.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Anthracene, 2-methyl-	192	C15H12	000613-12-7	94
2		Anthracene, 9-methyl-	192	C15H12	000779-02-2	91
3		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	90
4		Anthracene, 1-methyl-	192	C15H12	000610-48-0	90
5		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	90



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Misc :
ALS Vial : 18 Sample Multiplier: 1

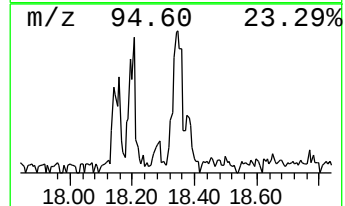
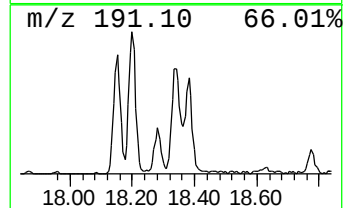
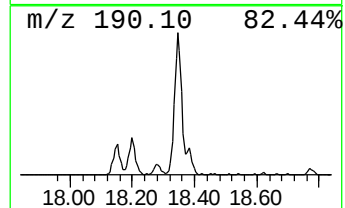
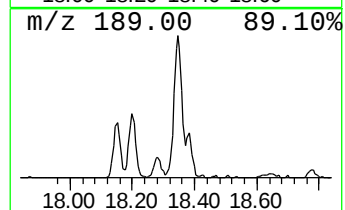
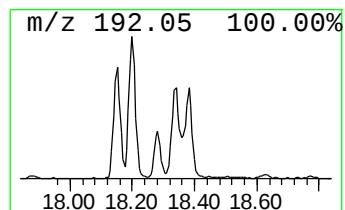
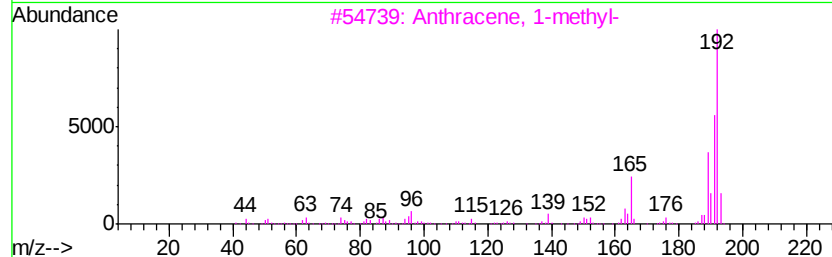
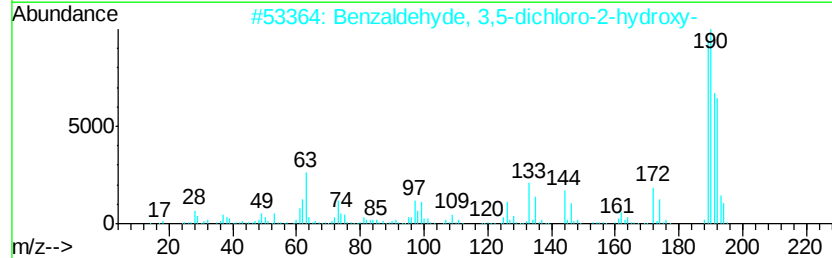
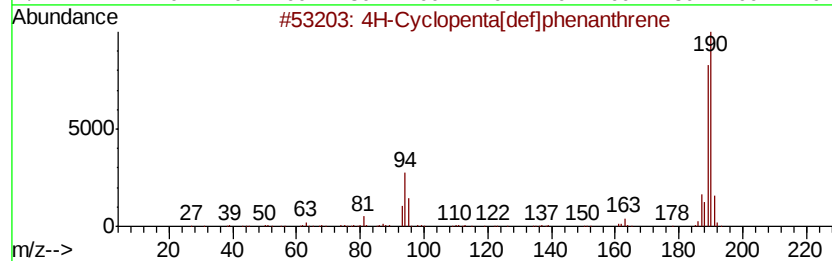
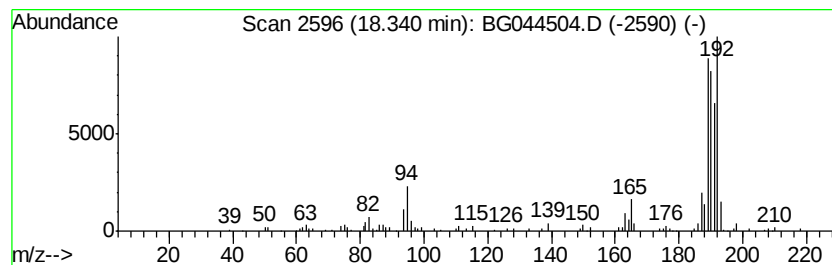
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ClientSampleId :
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.34	3.91 ng	175826	Phenanthrene-d10	17.27	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	91
2	Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	53
3	Anthracene, 1-methyl-	192	C15H12	000610-48-0	42
4	5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	30
5	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	30



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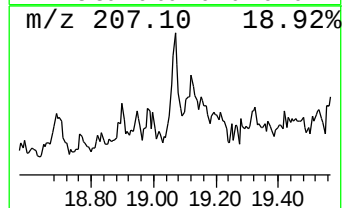
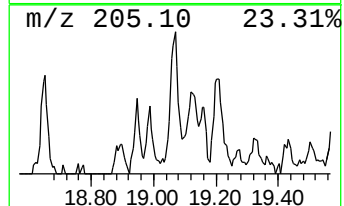
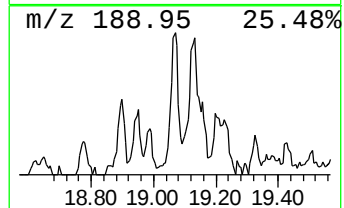
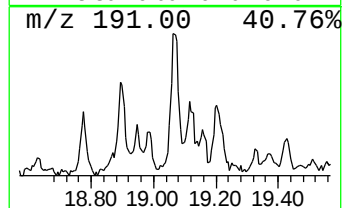
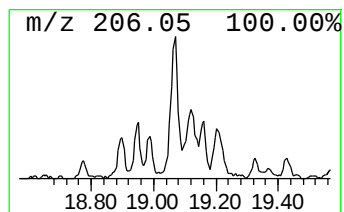
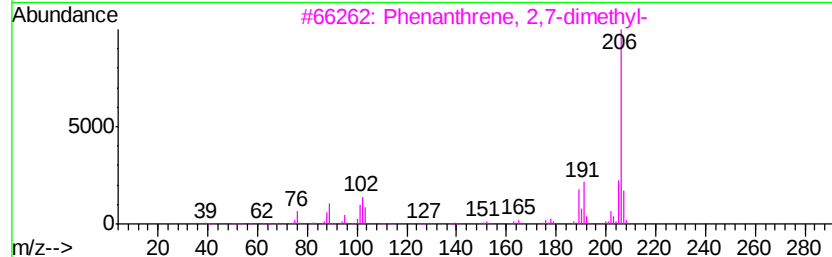
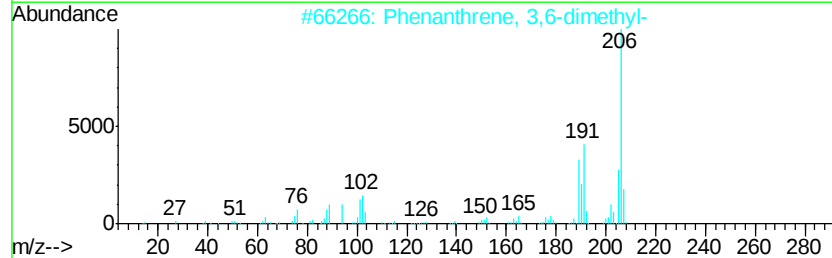
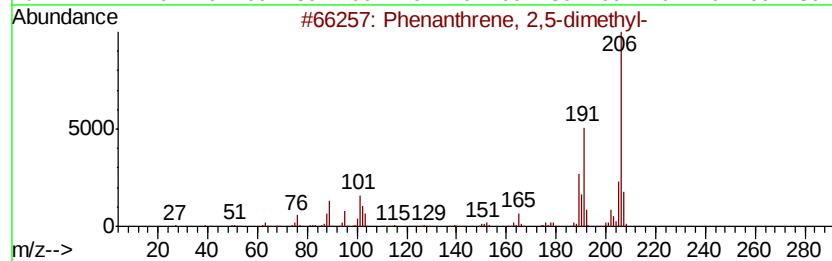
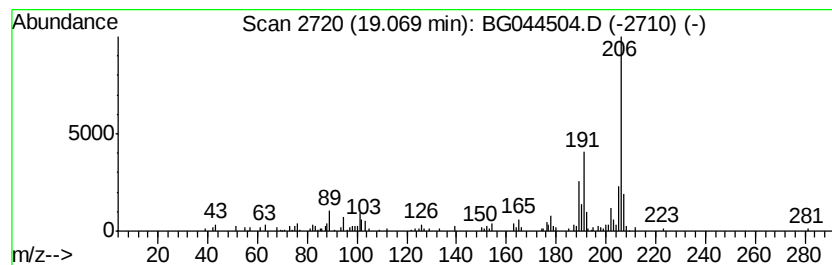
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
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Peak Number 5 Phenanthrene, 2,5-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.07	2.37 ng	106698	Phenanthrene-d10	17.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	97
2		Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	94
3		Phenanthrene, 2,7-dimethyl-	206	C16H14	001576-69-8	93
4		9,10-Dimethylantracene	206	C16H14	000781-43-1	91
5		Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	91



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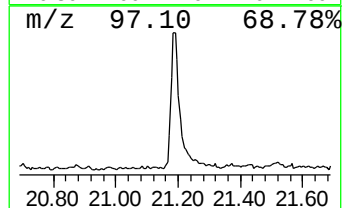
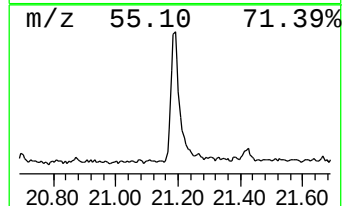
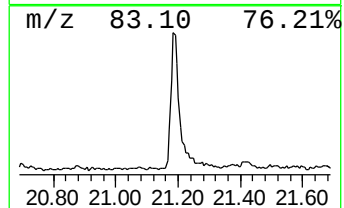
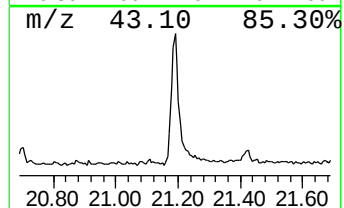
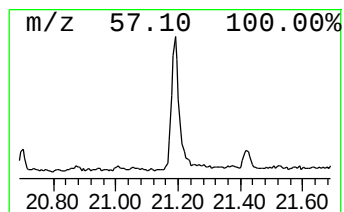
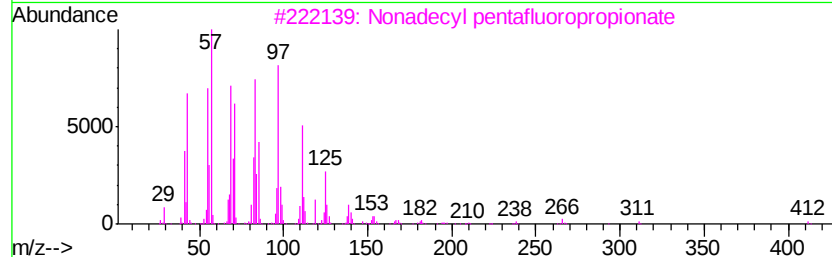
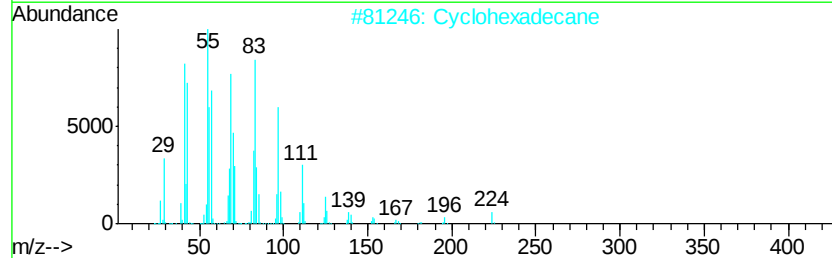
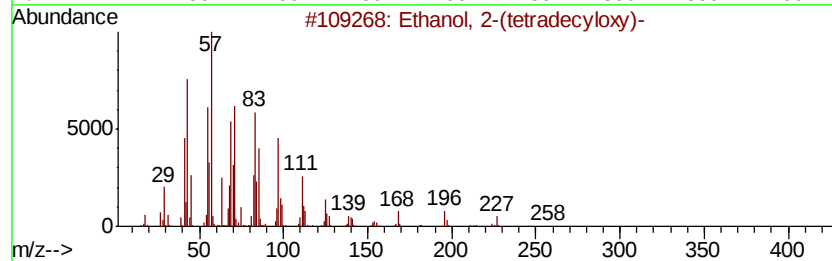
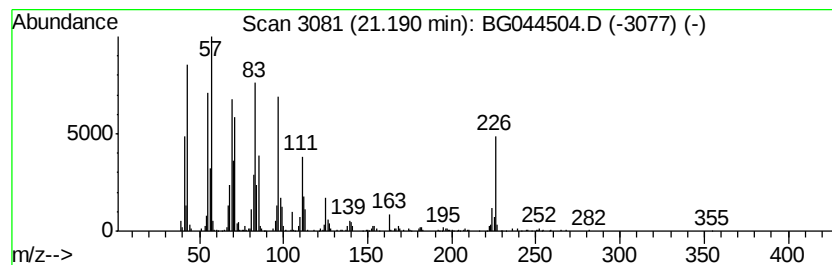
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MARCY-GREEN-BH-02-A

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Ethanol, 2-(tetradecyloxy)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.		
21.19	6.78 ng	517974	Chrysene-d12	21.52		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	94
2		Cyclohexadecane	224	C16H32	000295-65-8	93
3		Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	81
4		E-15-Heptadecenal	252	C17H32O	1000130-97-9	78
5		Trihexadecyl borate	735	C48H99B03	002665-11-4	74



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021920\
Data File : BG044504.D
Acq On : 20 Feb 2020 00:39
Operator : CG/JU
Sample : L1547-07
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
MARCY-GREEN-BH-02-A

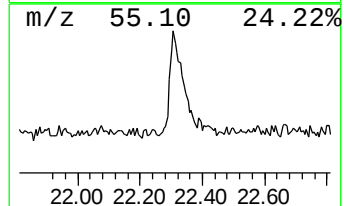
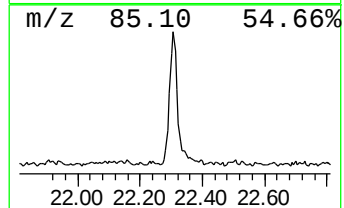
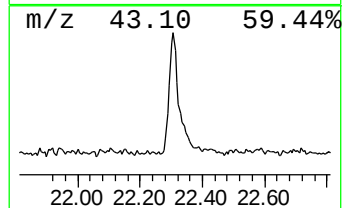
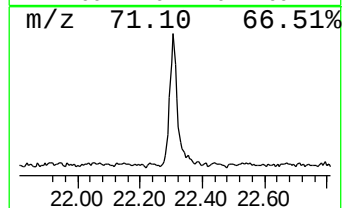
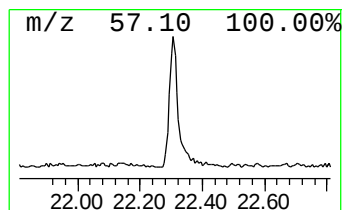
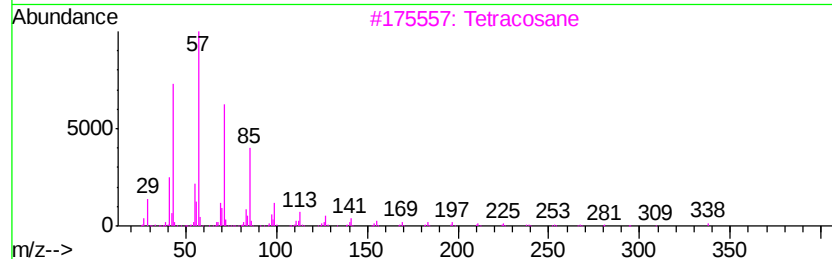
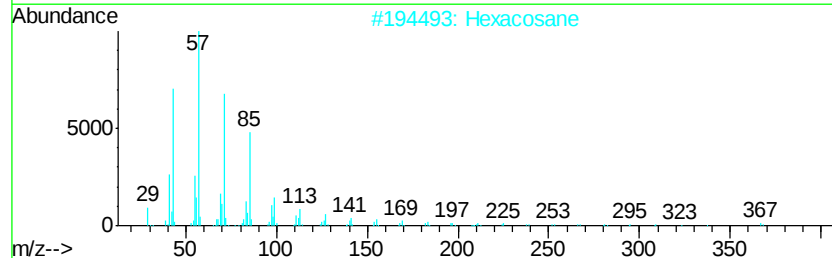
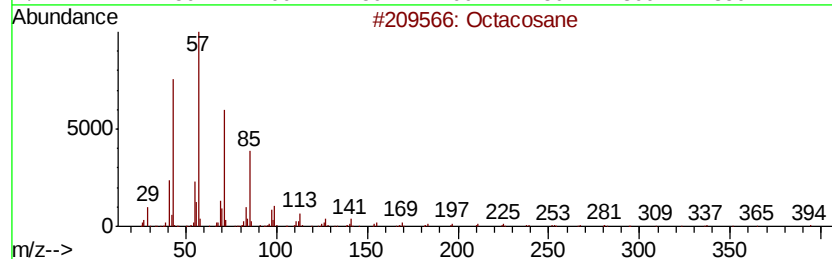
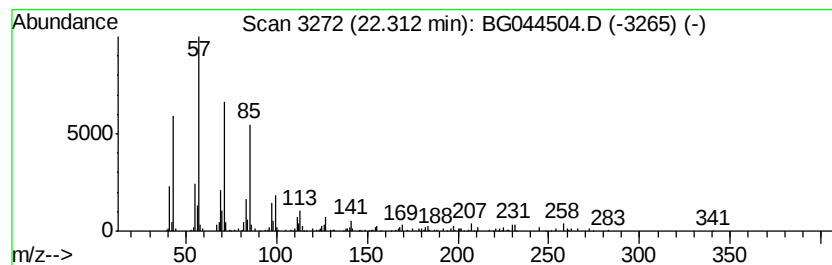
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Octacosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.31	3.30 ng	252195	Chrysene-d12	21.52

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	90
2		Hexacosane	366	C26H54	000630-01-3	90
3		Tetracosane	338	C24H50	000646-31-1	90
4		Heneicosane	296	C21H44	000629-94-7	90
5		Tetratetracontane	619	C44H90	007098-22-8	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021920\
Data File : BG044504.D
Acq On : 20 Feb 2020 00:39
Operator : CG/JU
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ALS Vial : 18 Sample Multiplier: 1

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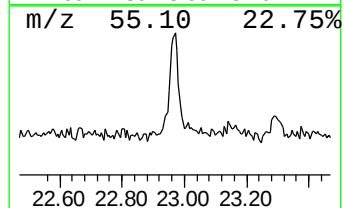
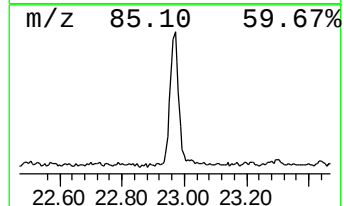
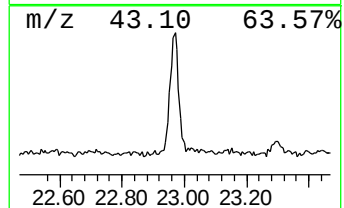
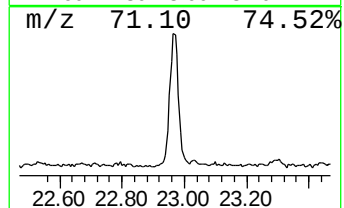
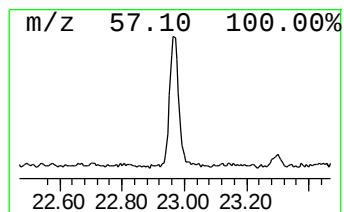
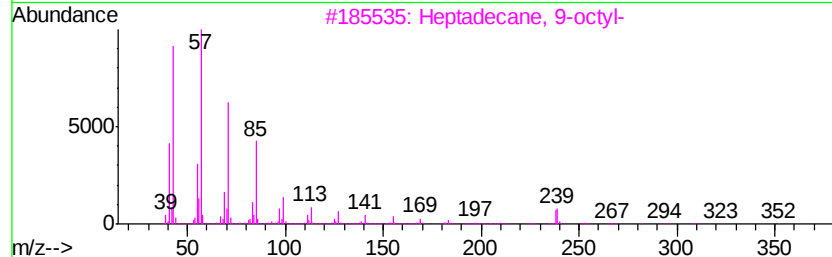
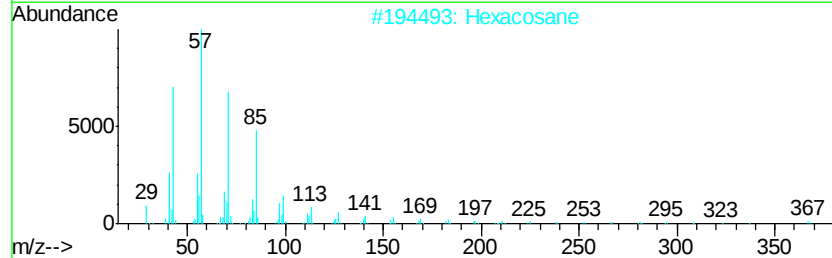
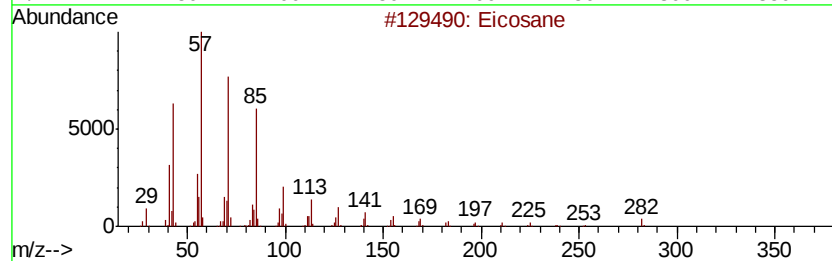
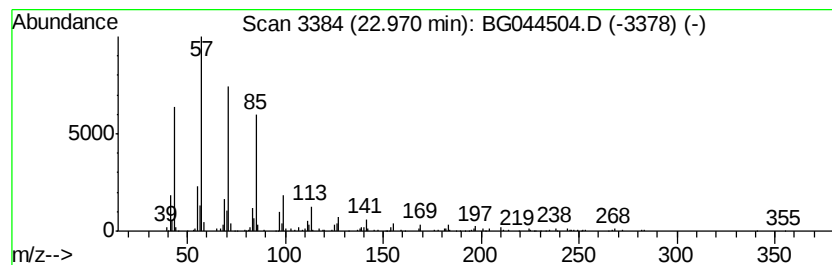
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Eicosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.97	2.70 ng	206050	Chrysene-d12	21.52

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	97
2	Hexacosane		366	C26H54	000630-01-3	91
3	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	91
4	Heneicosane		296	C21H44	000629-94-7	91
5	Tetracosane		338	C24H50	000646-31-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021920\
Data File : BG044504.D
Acq On : 20 Feb 2020 00:39
Operator : CG/JU
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Misc :
ALS Vial : 18 Sample Multiplier: 1

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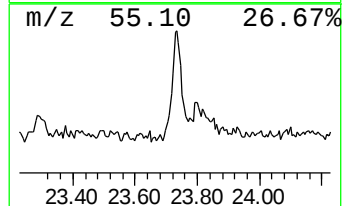
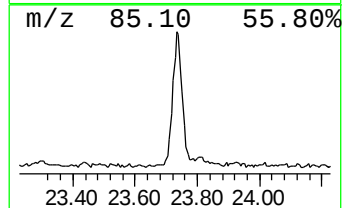
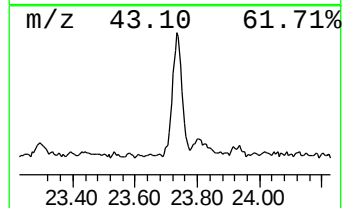
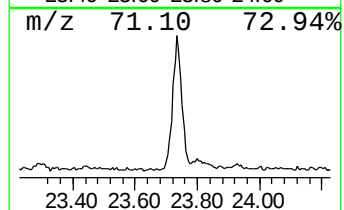
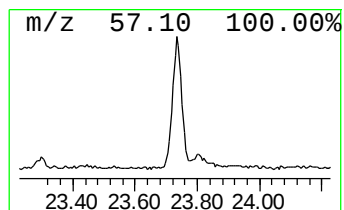
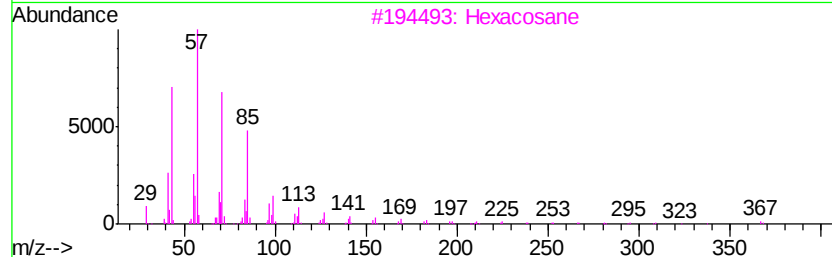
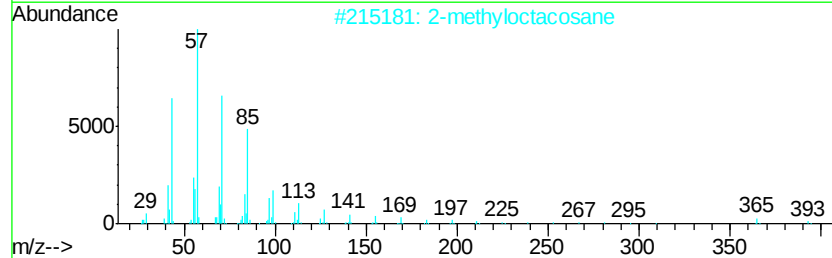
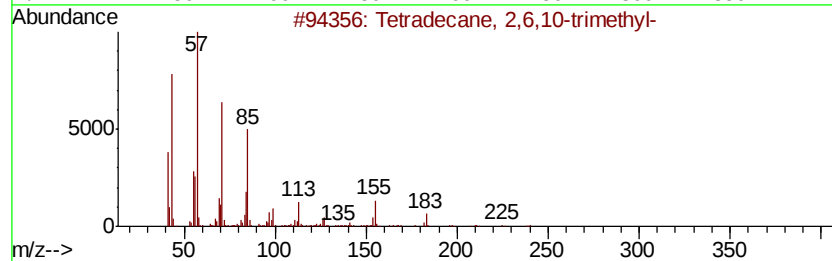
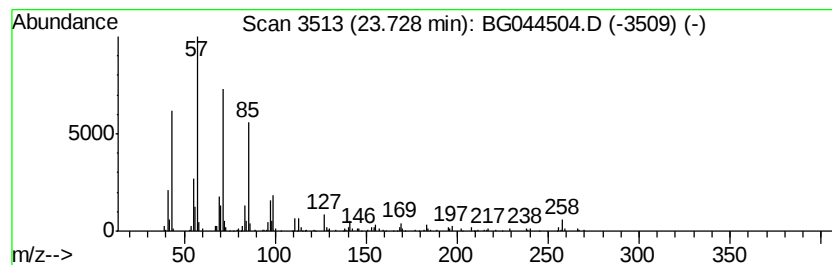
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Tetradecane, 2,6,10-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.73	3.24 ng	228498	Perylene-d12	24.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	91
2		2-methyloctacosane	408	C29H60	1000376-72-8	91
3		Hexacosane	366	C26H54	000630-01-3	90
4		Eicosane	282	C20H42	000112-95-8	87
5		Tetracosane	338	C24H50	000646-31-1	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021920\
Data File : BG044504.D
Acq On : 20 Feb 2020 00:39
Operator : CG/JU
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Misc :
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ClientSampleId :
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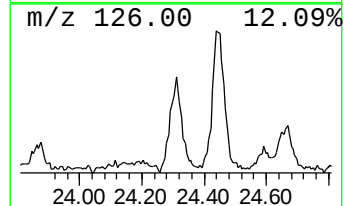
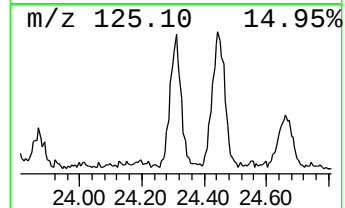
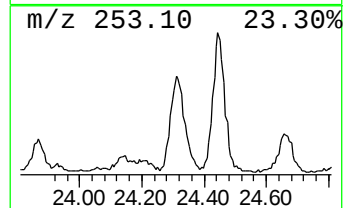
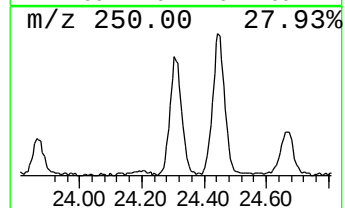
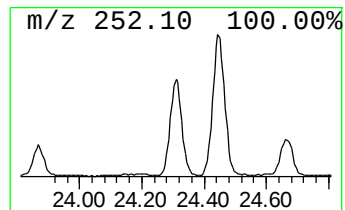
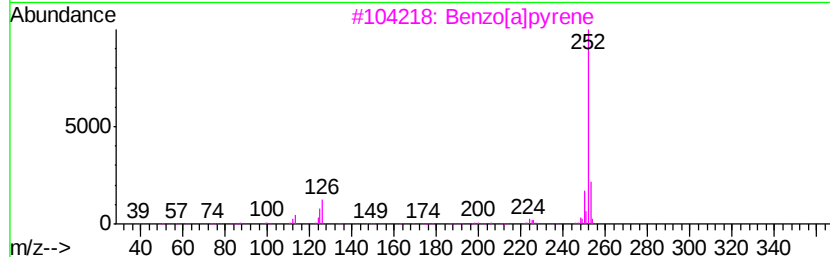
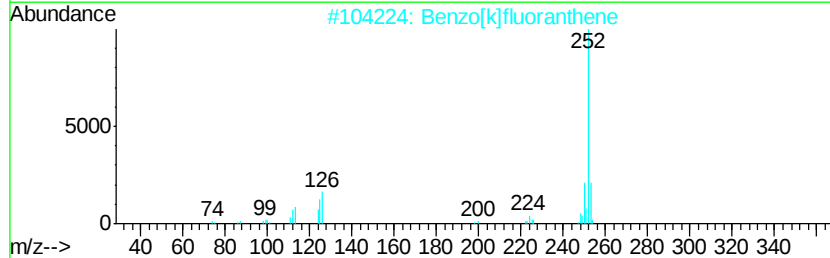
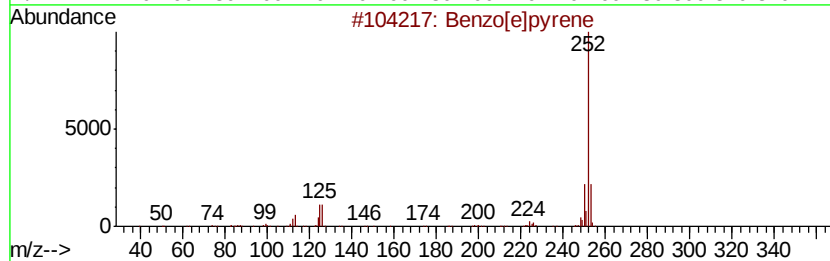
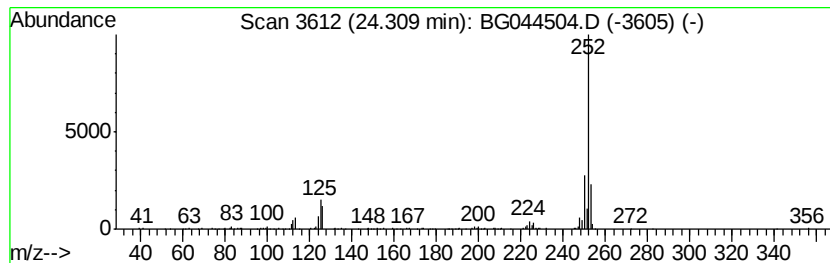
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Benzo[e]pyrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.31	4.21 ng	297538	Perylene-d12	24.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[e]pyrene	252	C20H12	000192-97-2	98
2		Benzo[k]fluoranthene	252	C20H12	000207-08-9	97
3		Benzo[a]pyrene	252	C20H12	000050-32-8	97
4		Perylene	252	C20H12	000198-55-0	97
5		Benz[e]acephenanthrylene	252	C20H12	000205-99-2	95



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021920\
Data File : BG044504.D
Acq On : 20 Feb 2020 00:39
Operator : CG/JU
Sample : L1547-07
Misc :
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Instrument :
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ClientSampleId :
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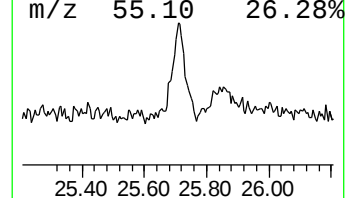
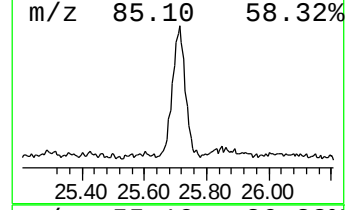
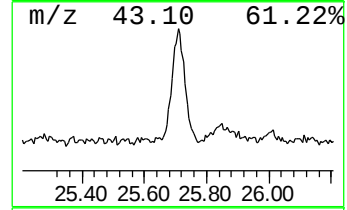
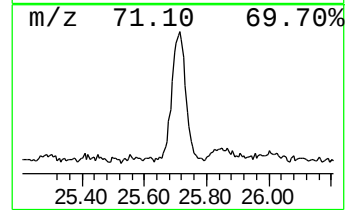
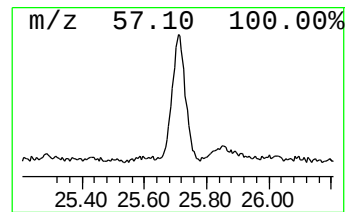
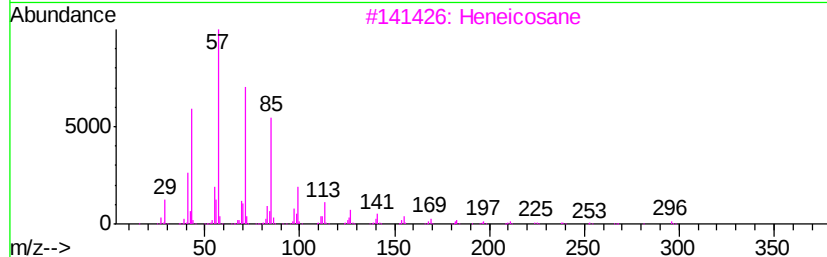
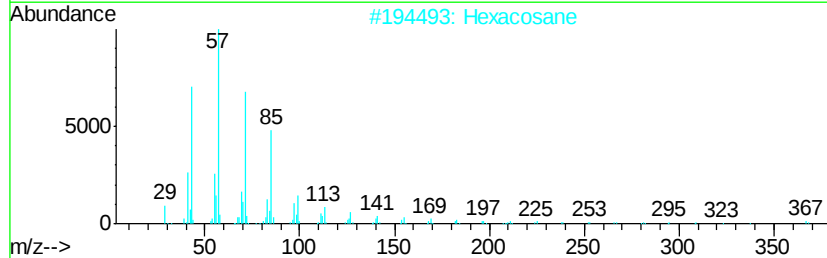
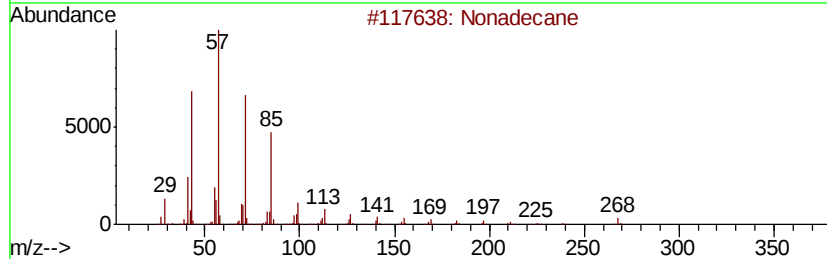
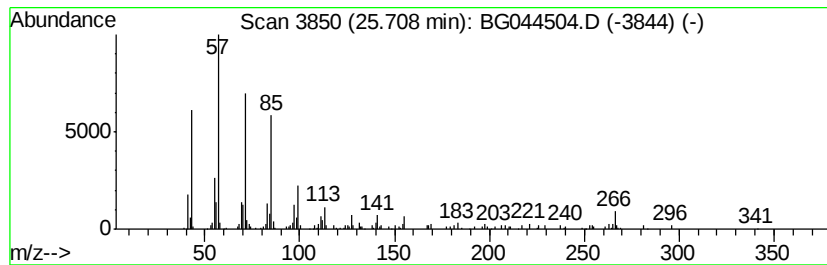
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Nonadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.71	3.14 ng	221776	Perylene-d12	24.60

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	93
2			Hexacosane	366	C26H54	000630-01-3	90
3			Heneicosane	296	C21H44	000629-94-7	90
4			Hexadecane	226	C16H34	000544-76-3	89
5			Heptacosane	380	C27H56	000593-49-7	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG021920\
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 Operator : CG/JU
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 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
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 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
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Anthracene, 2-met...	18.20	2.5	ng	111005	4	17.27	900152	20.0
4H-Cyclopenta[def...	18.34	3.9	ng	175826	4	17.27	900152	20.0
Phenanthrene, 2,5...	19.07	2.4	ng	106698	4	17.27	900152	20.0
Ethanol, 2-(tetra...	21.19	6.8	ng	517974	5	21.52	1527590	20.0
Octacosane	22.31	3.3	ng	252195	5	21.52	1527590	20.0
Eicosane	22.97	2.7	ng	206050	5	21.52	1527590	20.0
Tetradecane, 2,6,...	23.73	3.2	ng	228498	6	24.60	1412390	20.0
Benzo[e]pyrene	24.31	4.2	ng	297538	6	24.60	1412390	20.0
Nonadecane	25.71	3.1	ng	221776	6	24.60	1412390	20.0