

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011819\
 Data File : BG039216.D
 Acq On : 18 Jan 2019 15:16
 Operator : JU/SJ
 Sample : K1157-09
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 WC-05-011519

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-BG010919.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	3.134	3	8	15	rVB3	4892	10436	1.26%	0.112%
2	4.486	233	238	245	rVB2	7110	11874	1.44%	0.128%
3	5.097	337	342	350	rVB	27957	42277	5.12%	0.456%
4	5.567	415	422	432	rBV	221049	375937	45.54%	4.051%
5	7.171	688	695	706	rBV	240778	459948	55.72%	4.957%
6	7.541	751	758	768	rBV	227328	397688	48.18%	4.286%
7	8.011	831	838	844	rBV	72990	123091	14.91%	1.327%
8	8.328	884	892	901	rBB	203946	341749	41.40%	3.683%
9	9.186	1030	1038	1050	rBV	154925	285894	34.63%	3.081%
10	10.825	1309	1317	1325	rBV	95189	176930	21.43%	1.907%
11	13.269	1725	1733	1744	rBB	375922	593772	71.93%	6.399%
12	14.092	1868	1873	1879	rBV	23794	35587	4.31%	0.384%
13	14.650	1959	1968	1975	rBB2	140552	226534	27.44%	2.441%
14	15.050	2031	2036	2043	rVB	7002	10931	1.32%	0.118%
15	16.143	2215	2222	2231	rBV	312926	500728	60.66%	5.396%
16	17.218	2401	2405	2410	rVB	5816	9118	1.10%	0.098%
17	17.400	2429	2436	2439	rBV	187690	300753	36.43%	3.241%
18	17.441	2439	2443	2449	rVB	142295	215605	26.12%	2.324%
19	17.529	2449	2458	2463	rBV	24612	40131	4.86%	0.432%
20	17.805	2498	2505	2511	rVB3	15699	25868	3.13%	0.279%
21	18.269	2576	2584	2588	rBV3	26137	52196	6.32%	0.563%
22	18.322	2588	2593	2599	rVV2	21126	37784	4.58%	0.407%
23	18.405	2602	2607	2611	rVB4	5911	10166	1.23%	0.110%
24	18.469	2611	2618	2621	rBV3	20283	40655	4.92%	0.438%
25	18.504	2621	2624	2629	rVB	9924	13867	1.68%	0.149%
26	18.769	2664	2669	2673	rVV	13065	17964	2.18%	0.194%
27	18.816	2673	2677	2683	rVB2	12997	19720	2.39%	0.213%
28	19.174	2733	2738	2743	rBV3	8229	15151	1.84%	0.163%
29	19.333	2759	2765	2771	rBV5	7412	14199	1.72%	0.153%
30	19.450	2779	2785	2791	rVB	248529	355374	43.05%	3.830%
31	19.779	2836	2841	2842	rBV	38117	52507	6.36%	0.566%
32	19.815	2842	2847	2854	rVV	206454	302784	36.68%	3.263%
33	19.879	2854	2858	2863	rVB	11421	15331	1.86%	0.165%
34	20.003	2873	2879	2884	rBV	615947	825506	100.00%	8.896%

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35	20.056	2885	2888	2894	rVB	11079	16415	1.99%	0.177%
36	20.126	2894	2900	2901	rBV2	14278	20282	2.46%	0.219%
37	20.202	2908	2913	2918	rBV3	6524	12446	1.51%	0.134%
38	20.273	2918	2925	2926	rVV6	18253	34019	4.12%	0.367%
39	20.302	2926	2930	2936	rVB	38200	56671	6.87%	0.611%
40	20.402	2942	2947	2950	rBV	27798	37929	4.59%	0.409%
41	20.461	2950	2957	2960	rVV2	19075	36381	4.41%	0.392%
42	20.490	2960	2962	2967	rVB5	6686	8484	1.03%	0.091%
43	20.596	2975	2980	2984	rBV	11151	19615	2.38%	0.211%
44	20.649	2984	2989	2992	rBV2	10567	16294	1.97%	0.176%
45	20.725	2999	3002	3005	rBV4	6360	8902	1.08%	0.096%
46	20.761	3005	3008	3019	rVB6	10132	15976	1.94%	0.172%
47	21.019	3048	3052	3056	rBV7	5497	10333	1.25%	0.111%
48	21.066	3056	3060	3067	rVB	19958	31041	3.76%	0.335%
49	21.266	3083	3094	3102	rBV5	61673	167903	20.34%	1.809%
50	21.342	3102	3107	3111	rVV2	24032	48224	5.84%	0.520%
51	21.395	3113	3116	3124	rVB2	17255	32457	3.93%	0.350%
52	21.524	3133	3138	3145	rVB3	8355	18745	2.27%	0.202%
53	21.671	3149	3163	3168	rVV2	271137	691119	83.72%	7.448%
54	21.718	3168	3171	3182	rVB	121731	202583	24.54%	2.183%
55	21.806	3182	3186	3192	rBV3	20371	35463	4.30%	0.382%
56	21.871	3192	3197	3204	rVB	14659	27293	3.31%	0.294%
57	22.089	3228	3234	3240	rVB4	14227	34615	4.19%	0.373%
58	22.324	3268	3274	3279	rBV8	5871	14537	1.76%	0.157%
59	22.406	3279	3288	3291	rVV4	39540	82107	9.95%	0.885%
60	22.441	3291	3294	3308	rVB4	36613	104672	12.68%	1.128%
61	22.617	3320	3324	3330	rVB6	7203	15212	1.84%	0.164%
62	22.682	3330	3335	3336	rBV5	8162	13059	1.58%	0.141%
63	22.729	3340	3343	3352	rVB8	11022	20758	2.51%	0.224%
64	23.099	3400	3406	3413	rVV5	28083	60251	7.30%	0.649%
65	23.463	3462	3468	3471	rBV6	12800	23095	2.80%	0.249%
66	23.516	3475	3477	3483	rVB6	9188	11076	1.34%	0.119%
67	23.657	3499	3501	3505	rBV4	6806	9111	1.10%	0.098%
68	23.892	3532	3541	3549	rBV2	123209	405787	49.16%	4.373%
69	24.145	3578	3584	3594	rVB	15812	41027	4.97%	0.442%
70	24.356	3614	3620	3624	rBV8	5132	12314	1.49%	0.133%
71	24.621	3658	3665	3677	rVB2	42930	122905	14.89%	1.325%

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72	24.774	3682	3691	3699	rBV	59468	181197	21.95%	1.953%
73	24.932	3709	3718	3728	rBV	121326	340718	41.27%	3.672%
74	25.009	3728	3731	3744	rVB4	20096	52556	6.37%	0.566%
75	25.373	3788	3793	3797	rBV8	8375	19243	2.33%	0.207%
76	25.520	3816	3818	3826	rVB8	5951	10687	1.29%	0.115%
77	25.966	3884	3894	3902	rBV2	28486	77503	9.39%	0.835%
78	27.300	4119	4121	4128	rBV7	5488	10064	1.22%	0.108%
79	28.128	4254	4262	4264	rBV8	7194	14896	1.80%	0.161%
80	28.663	4342	4353	4362	rBV3	22890	86494	10.48%	0.932%
81	29.809	4544	4548	4549	rBV4	7544	8671	1.05%	0.093%

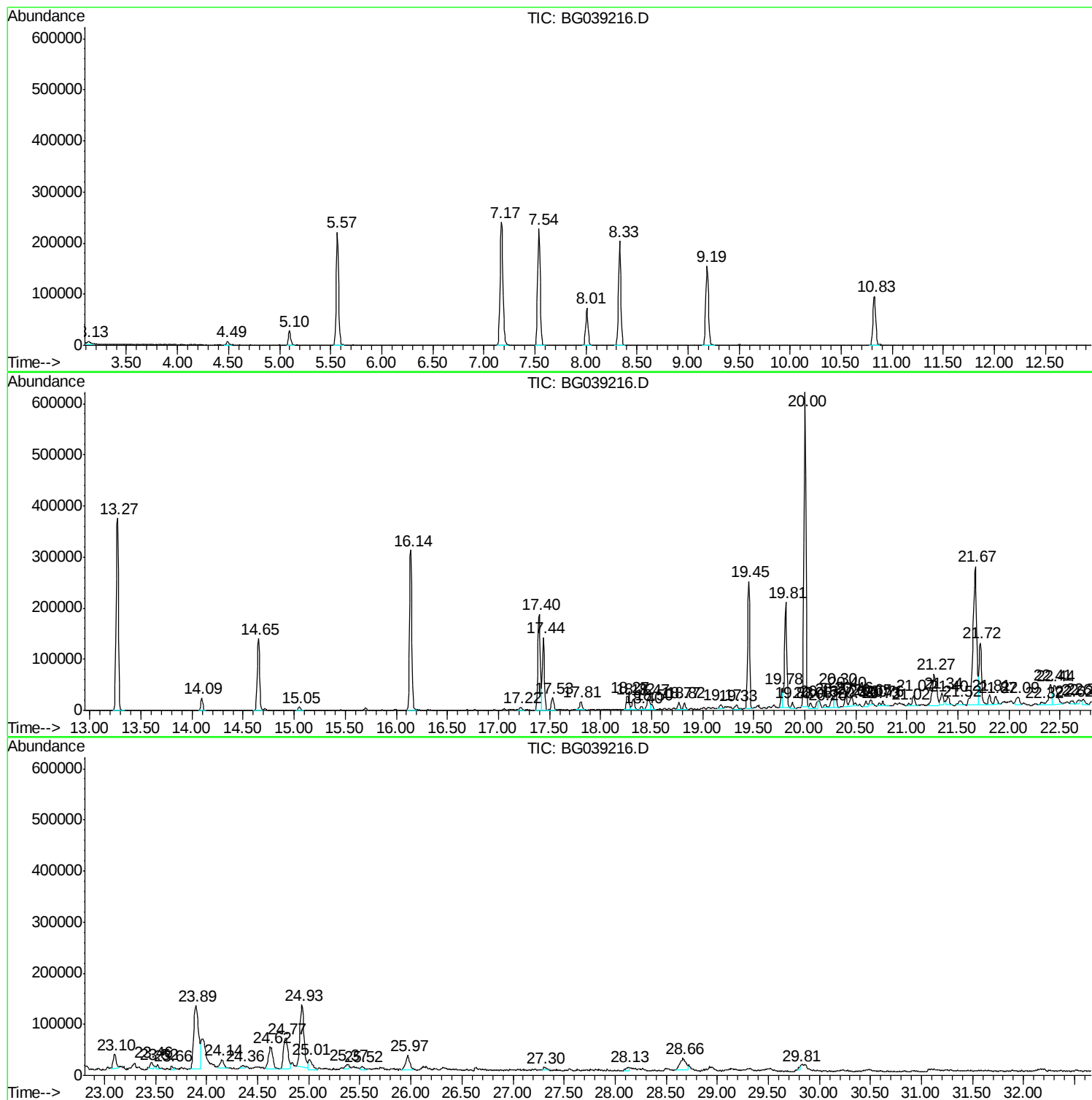
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG010919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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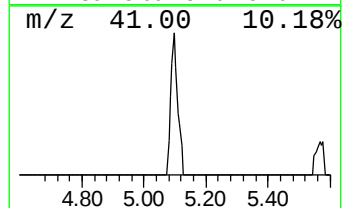
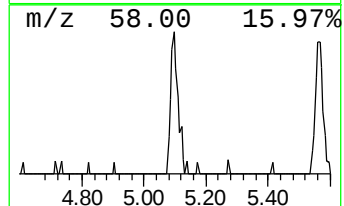
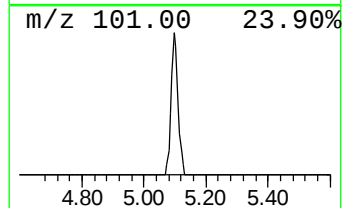
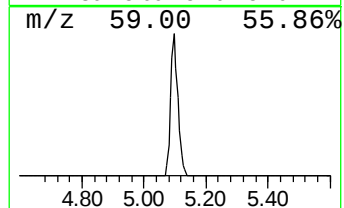
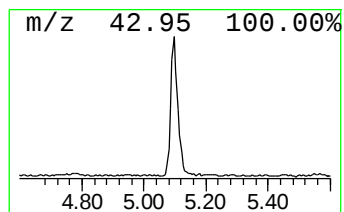
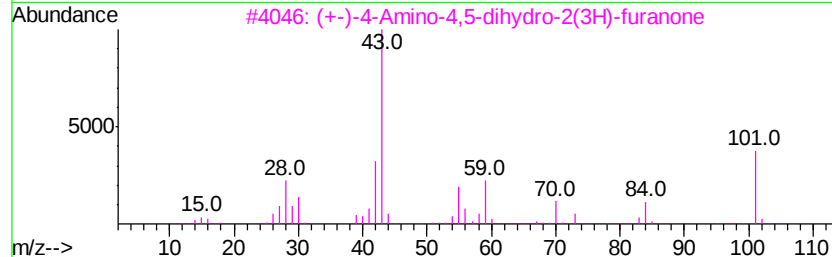
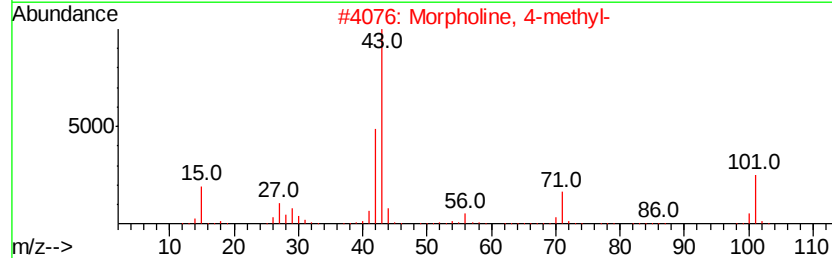
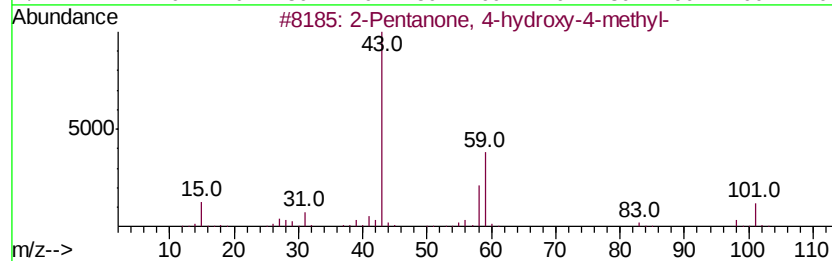
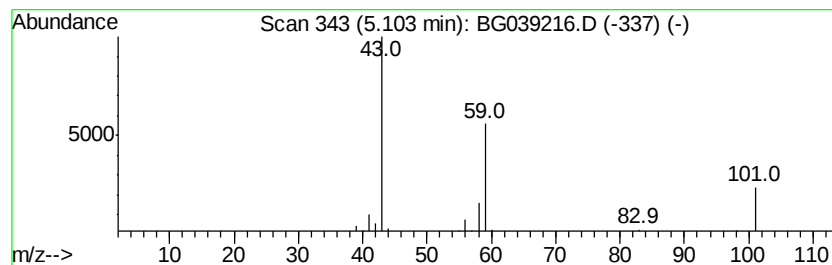
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG010919.M
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.10	6.87 ng	42277	1,4-Dichlorobenzene-d4	8.01

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	25
3			(+)-4-Amino-4,5-dihydro-2(3H)-furanone	101	C4H7NO2	016504-58-8	9
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9



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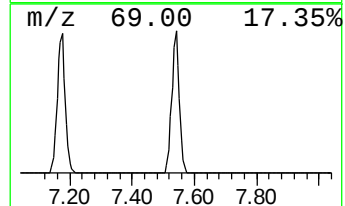
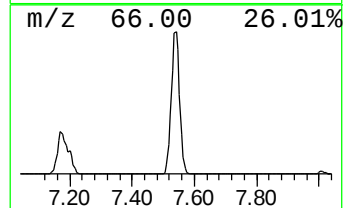
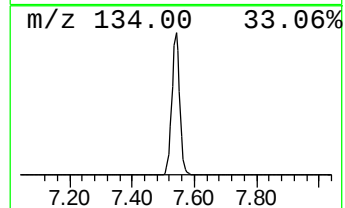
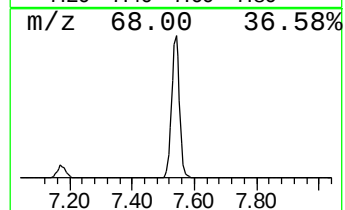
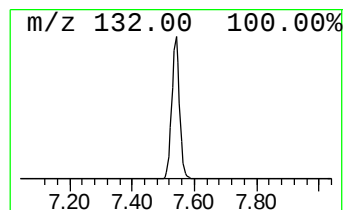
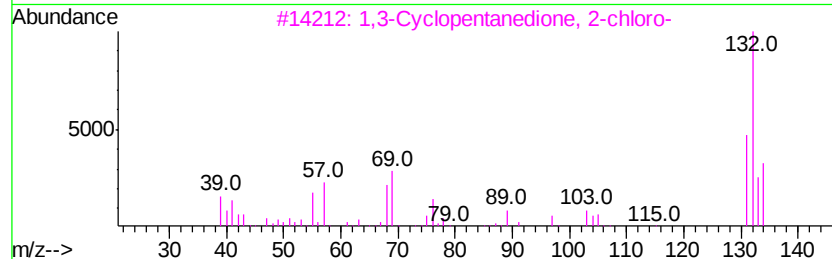
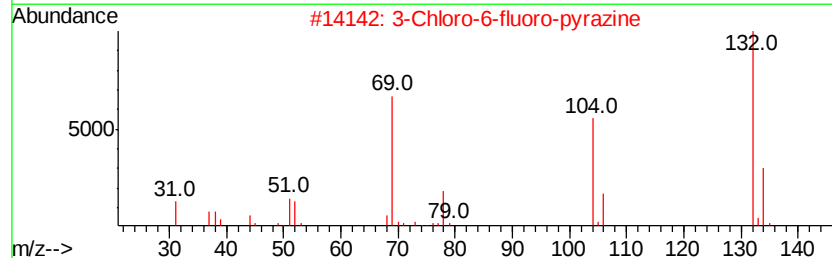
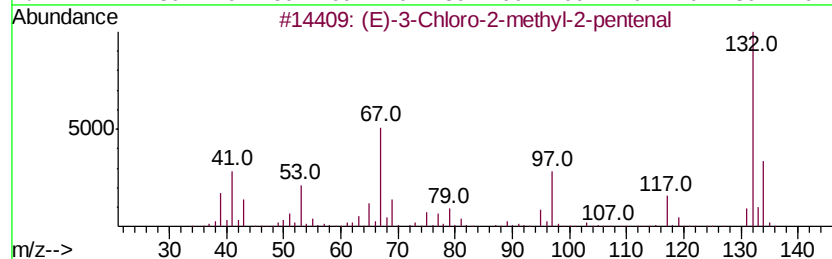
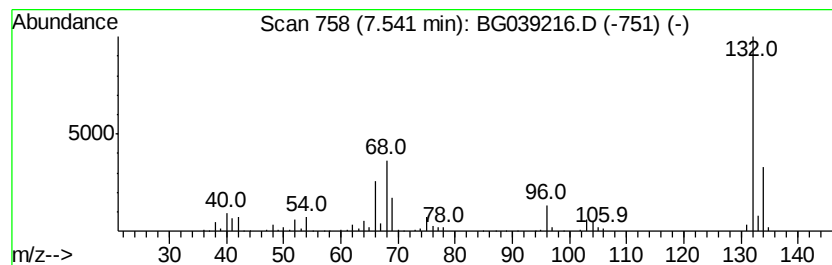
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TIC Library : C:\Database\NIST11.L
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Peak Number 2 unknown7.54 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.54	64.62 ng	397688	1,4-Dichlorobenzene-d4	8.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
3		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	23
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	17
5		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	16



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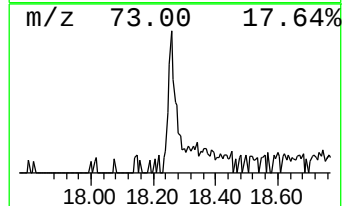
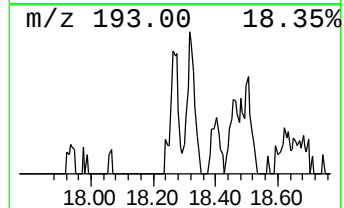
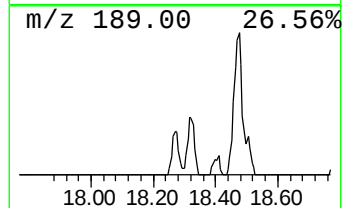
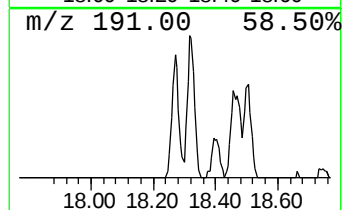
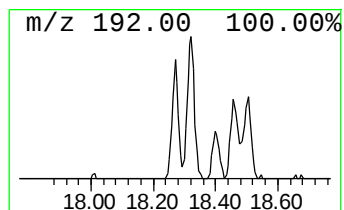
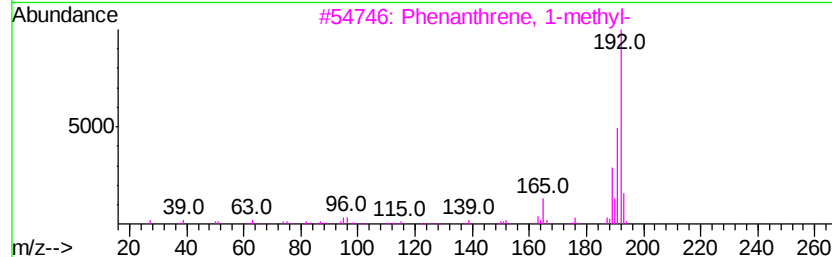
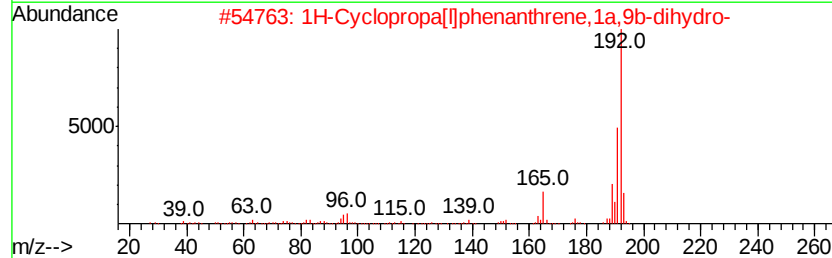
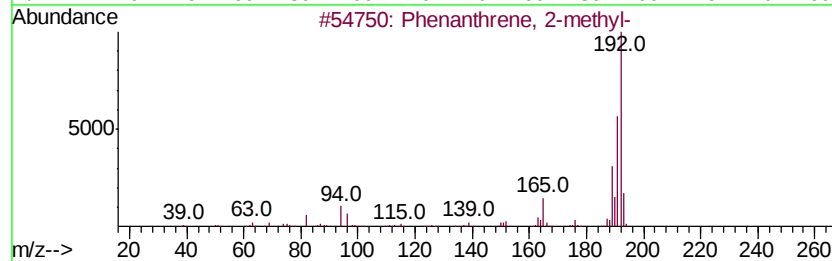
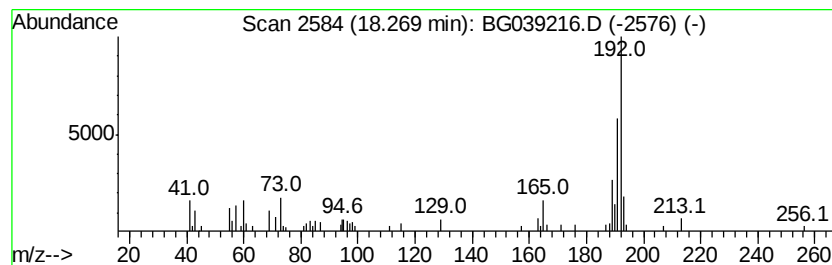
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TIC Library : C:\Database\NIST11.L
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Peak Number 4 Phenanthrene, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.27	3.47 ng	52196	Phenanthrene-d10	17.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	94
2		1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	94
3		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
4		Anthracene, 1-methyl-	192	C15H12	000610-48-0	93
5		Anthracene, 9-methyl-	192	C15H12	000779-02-2	81



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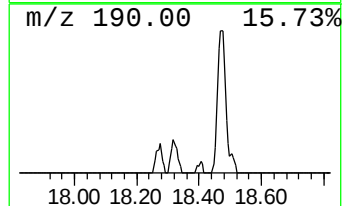
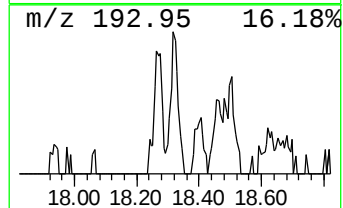
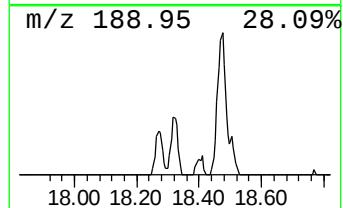
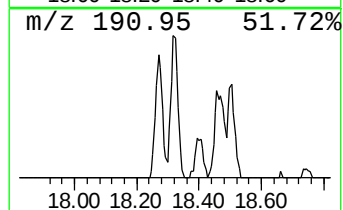
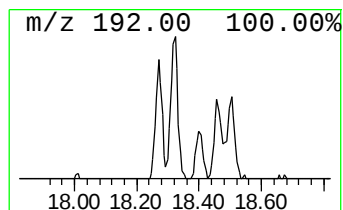
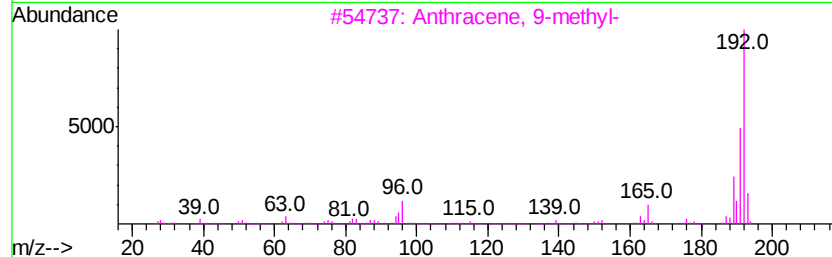
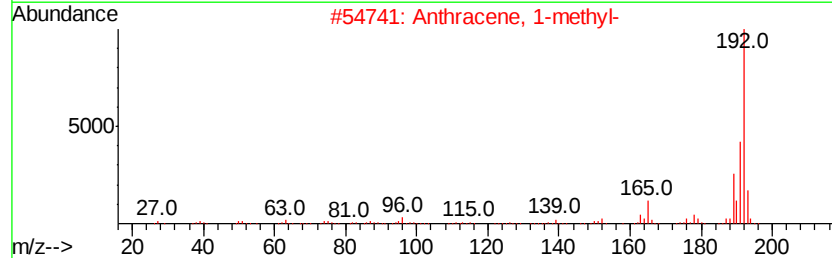
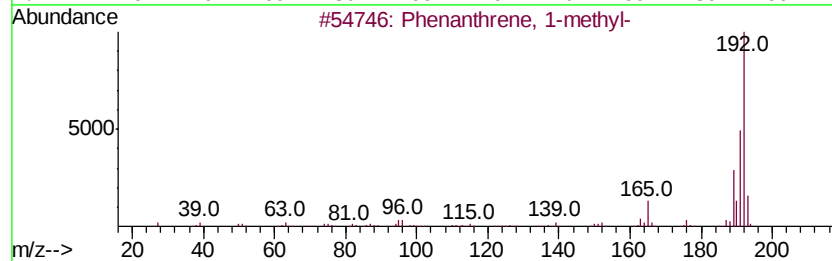
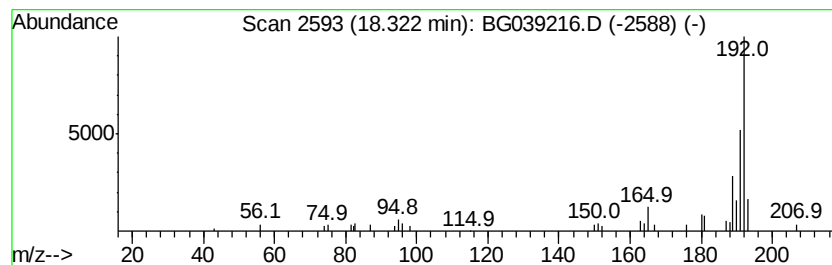
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Phenanthrene, 1-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.32	2.51 ng	37784	Phenanthrene-d10	17.40

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	94
2		Anthracene, 1-methyl-	192	C15H12	000610-48-0	93
3		Anthracene, 9-methyl-	192	C15H12	000779-02-2	91
4		1H-Cyclopropa[1,1]phenanthrene, 1a,...	192	C15H12	000949-41-7	81
5		Anthracene, 2-methyl-	192	C15H12	000613-12-7	81



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011819\
Data File : BG039216.D
Acq On : 18 Jan 2019 15:16
Operator : JU/SJ
Sample : K1157-09
Misc :
ALS Vial : 5 Sample Multiplier: 1

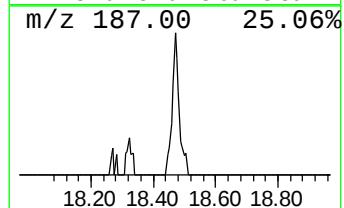
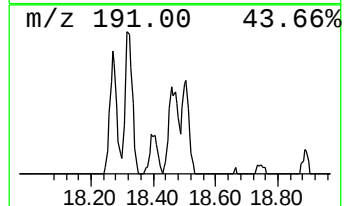
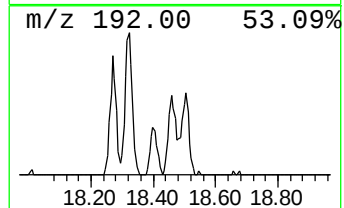
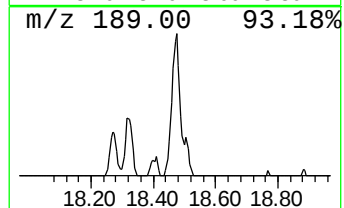
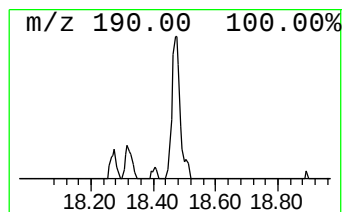
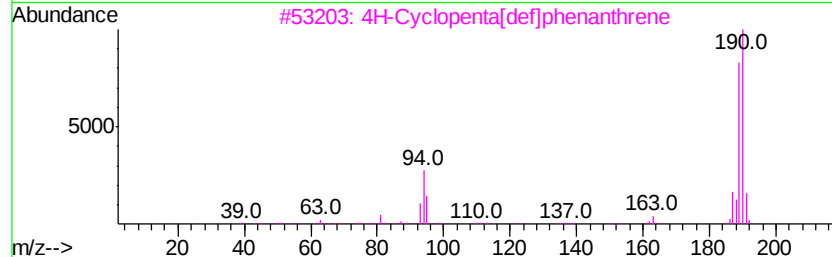
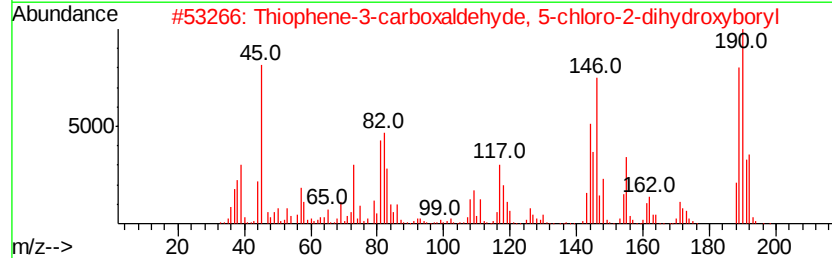
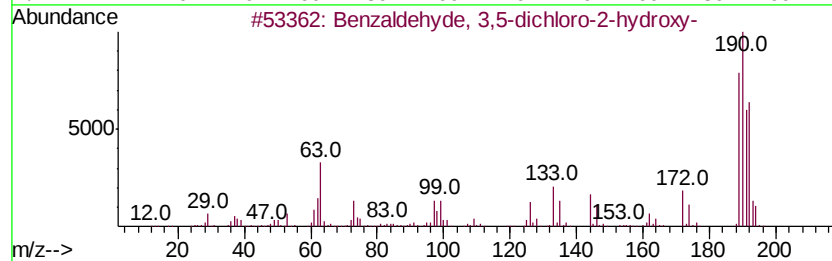
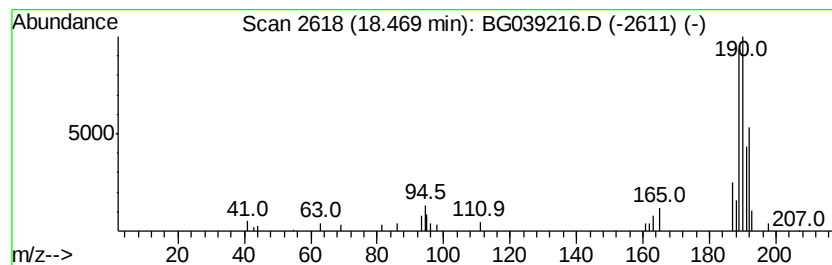
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Benzaldehyde, 3,5-dichloro-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.		
18.47	2.70 ng	40655	Phenanthrene-d10	17.40		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	52
2		Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	50
3		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	50
4		2,6-Dichlorobenzaldoxime	189	C7H5Cl2NO	025185-95-9	30
5		5-Thiazolecarboxamide, 4-amino-2...	189	C5H7N3OS2	1000338-08-9	27



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011819\
Data File : BG039216.D
Acq On : 18 Jan 2019 15:16
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Instrument :
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ClientSampleId :
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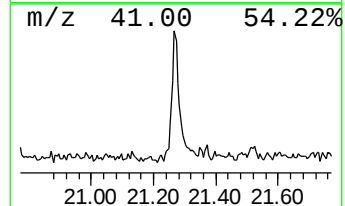
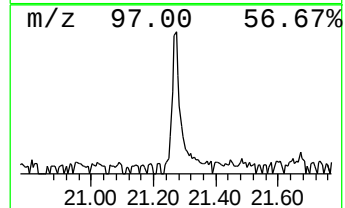
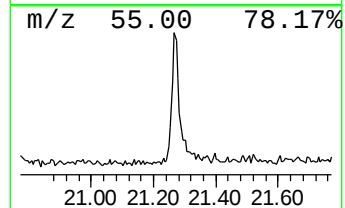
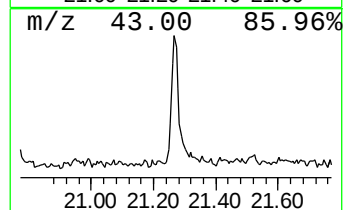
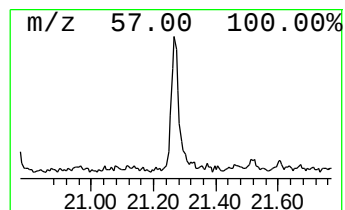
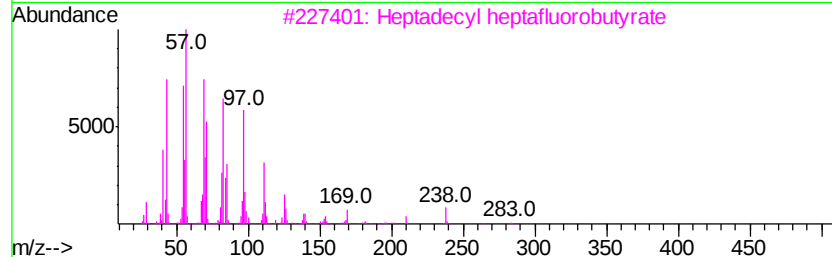
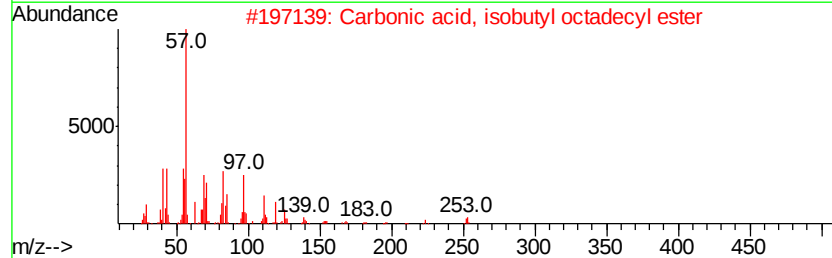
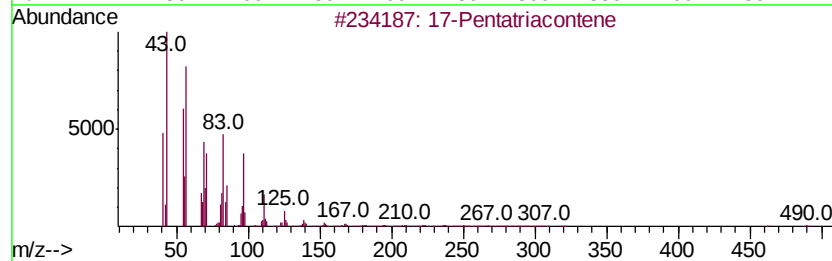
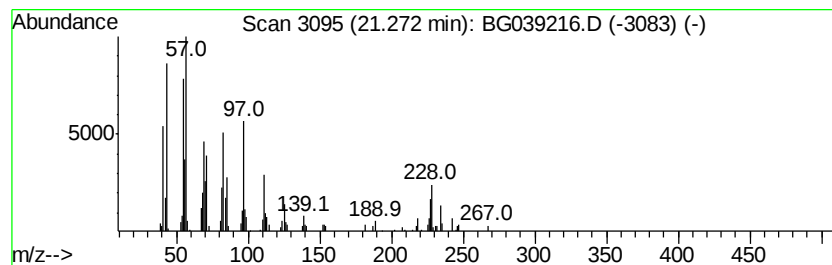
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 17-Pentatriacontene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.27	4.86 ng	167903	Chrysene-d12	21.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	17-Pentatriacontene	491	C35H70	006971-40-0	64
2		Carbonic acid, isobutyl octadecy...	370	C23H46O3	1000314-61-5	64
3		Heptadecyl heptafluorobutvrate	452	C21H35F7O2	959085-66-6	60
4		1-Docosene	308	C22H44	001599-67-3	60
5		Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	60



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011819\
Data File : BG039216.D
Acq On : 18 Jan 2019 15:16
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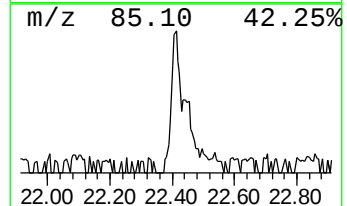
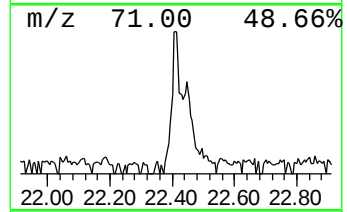
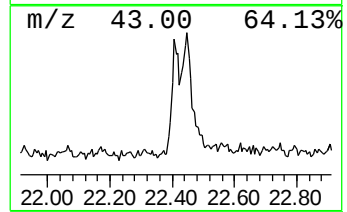
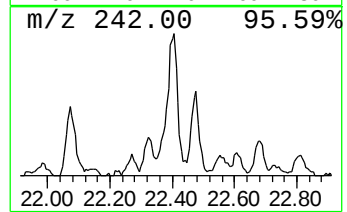
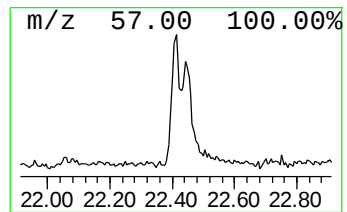
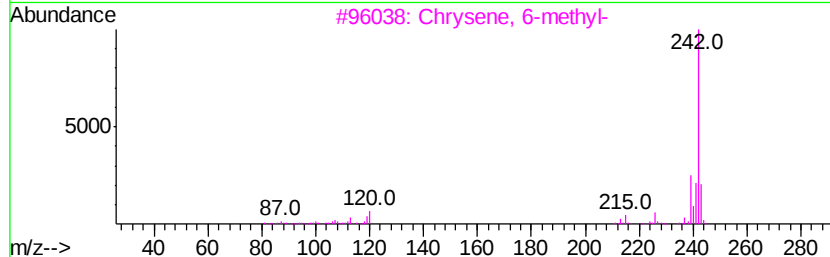
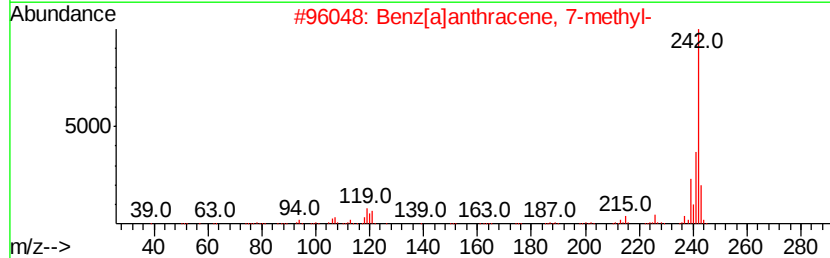
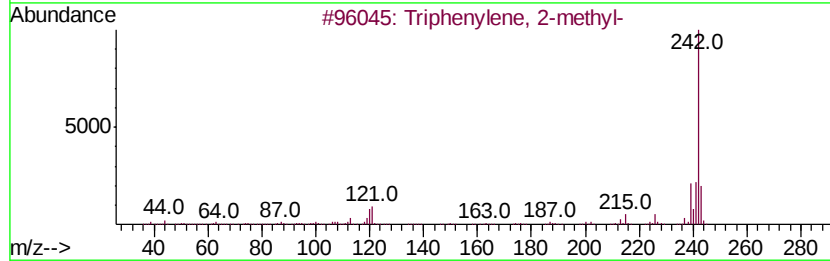
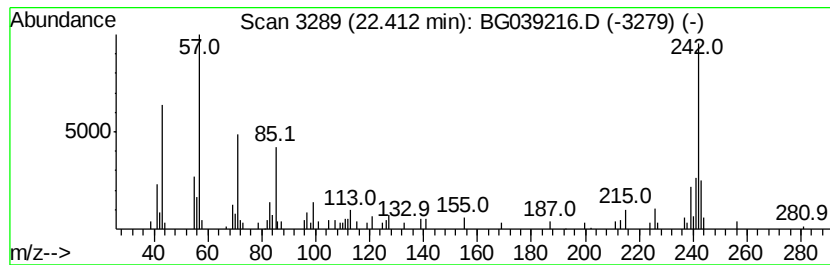
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Triphenylene, 2-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.41	2.38 ng	82107	Chrysene-d12	21.67	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Triphenylene, 2-methyl-	242	C19H14	001705-84-6	86
2	Benz[alanthracene, 7-methyl-	242	C19H14	002541-69-7	70
3	Chrysene, 6-methyl-	242	C19H14	001705-85-7	70
4	Benz[alanthracene, 1-methyl-	242	C19H14	002498-77-3	64
5	Benzo[c]phenanthrene, 2-methyl-	242	C19H14	002606-85-1	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011819\
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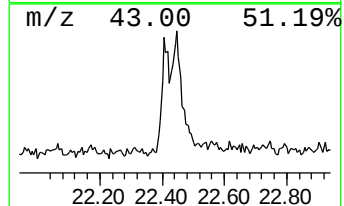
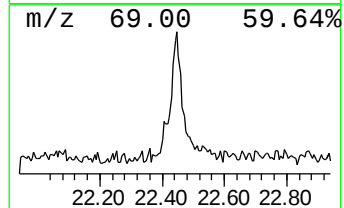
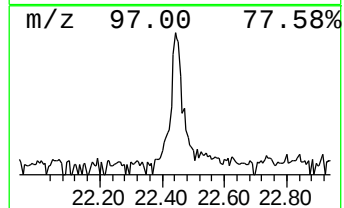
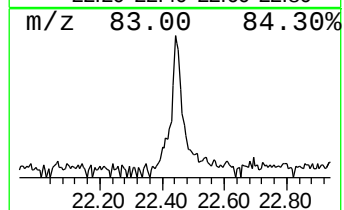
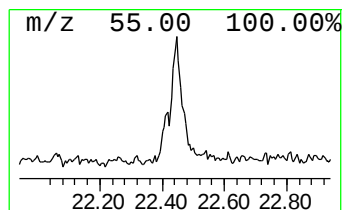
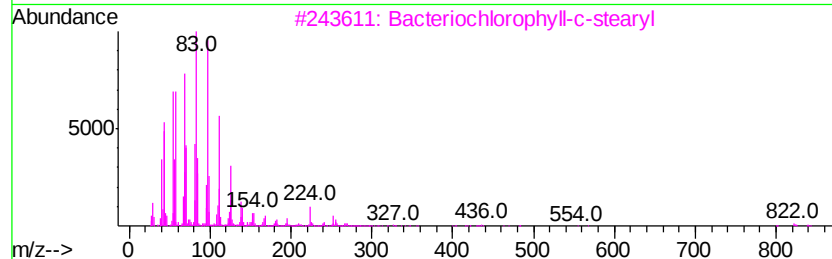
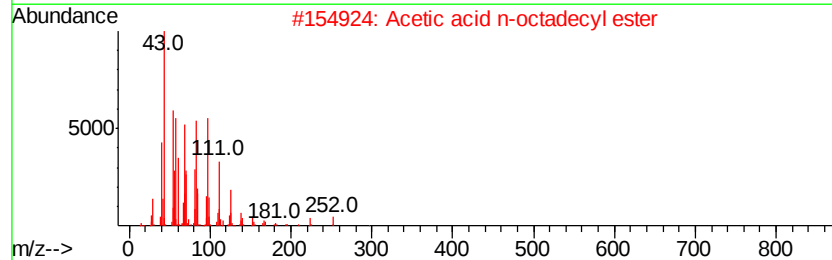
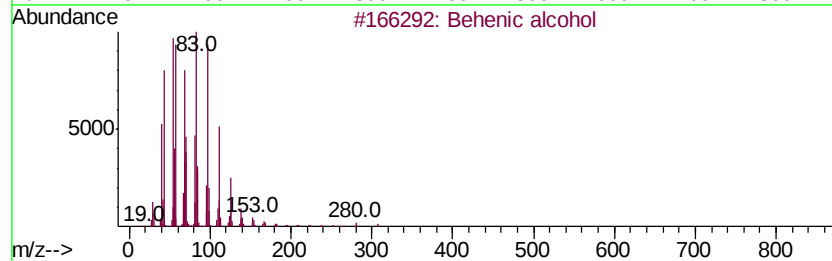
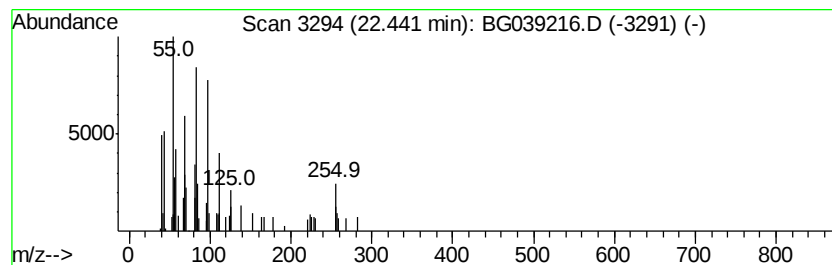
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG010919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Behenic alcohol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.44	3.03 ng	104672	Chrysene-d12	21.67
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Behenic alcohol	326 C22H46O	000661-19-8 72
2		Acetic acid n-octadecyl ester	312 C20H40O2	000822-23-1 52
3		Bacteriochlorophyll-c-stearyl	841 C52H72MgN4O4	1000164-49-7 50
4		Cyclopentane, 1,1,3-trimethyl-	112 C8H16	004516-69-2 46
5		Cyclohexadecane	224 C16H32	000295-65-8 46



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011819\
Data File : BG039216.D
Acq On : 18 Jan 2019 15:16
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Sample : K1157-09
Misc :
ALS Vial : 5 Sample Multiplier: 1

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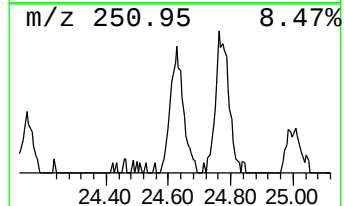
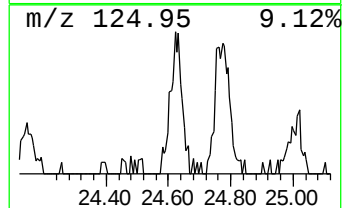
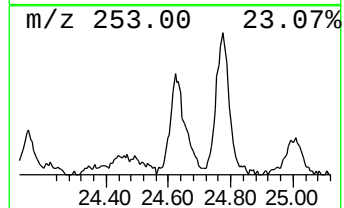
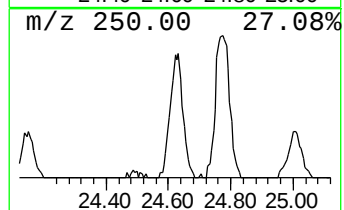
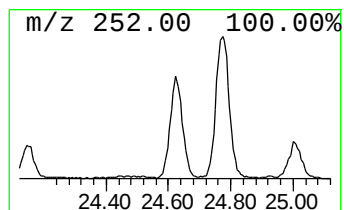
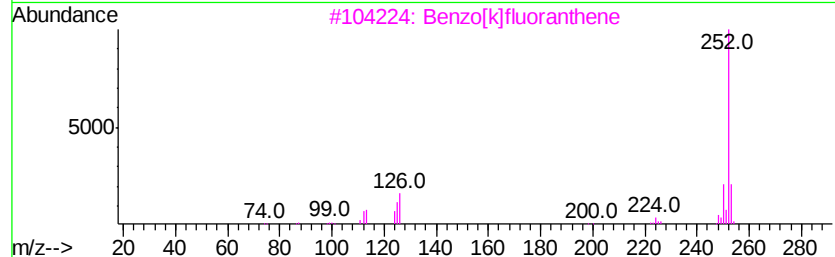
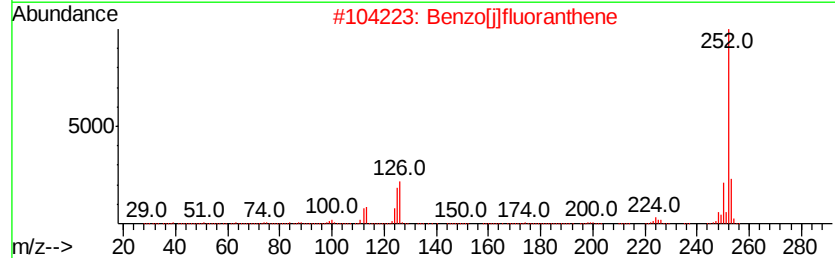
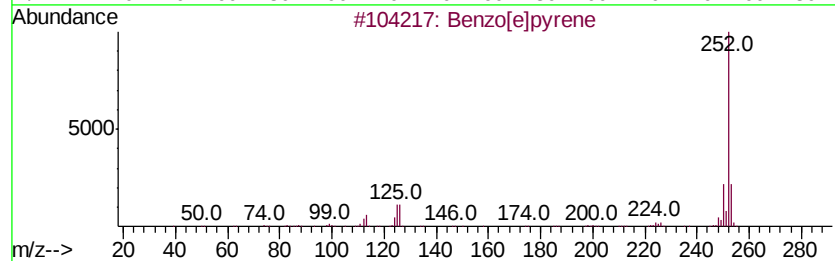
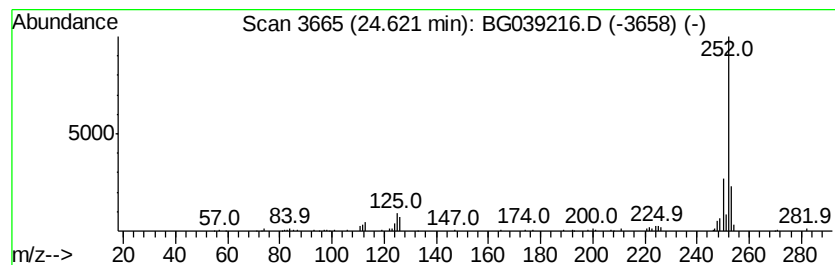
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Benzo[e]pyrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.62	7.21 ng	122905	Perylene-d12	24.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[e]pyrene	252	C20H12	000192-97-2	93
2		Benzo[i]fluoranthene	252	C20H12	000205-82-3	90
3		Benzo[k]fluoranthene	252	C20H12	000207-08-9	90
4		Perylene	252	C20H12	000198-55-0	90
5		Benzo[a]pyrene	252	C20H12	000050-32-8	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011819\
Data File : BG039216.D
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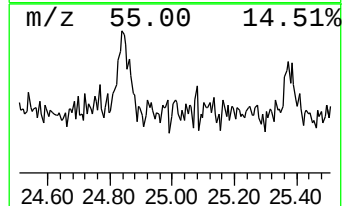
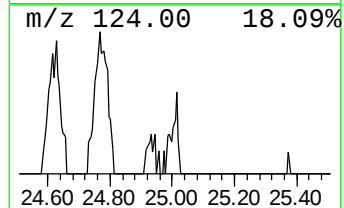
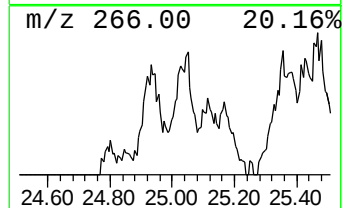
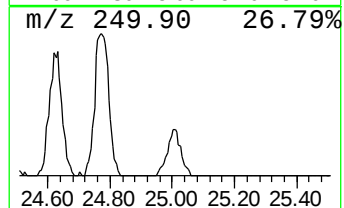
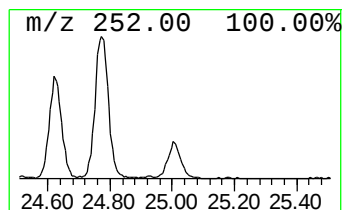
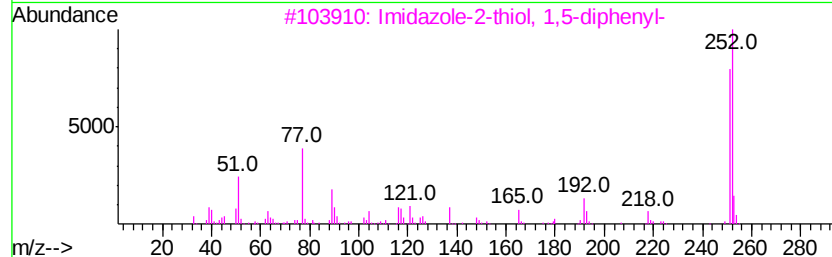
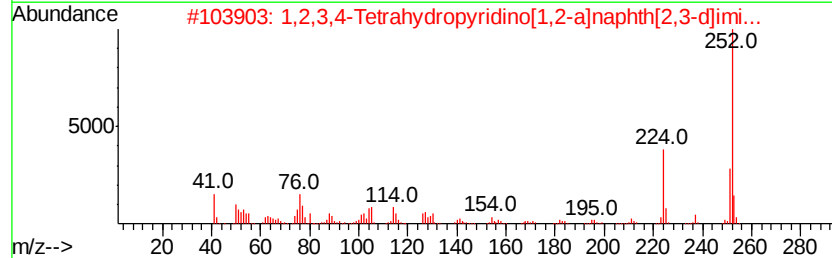
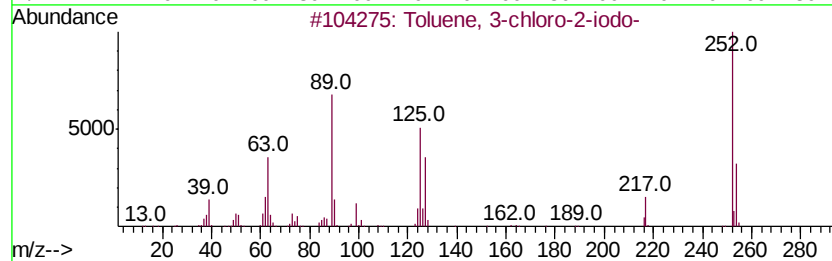
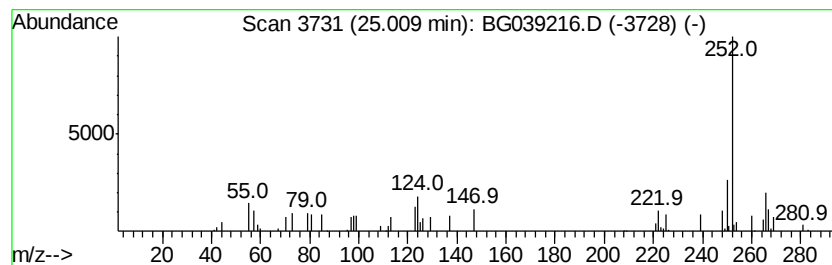
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown25.01 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.01	3.09 ng	52556	Perylene-d12	24.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Toluene, 3-chloro-2-iodo-	252	C7H6ClI	005100-98-1	43
2		1,2,3,4-Tetrahydropyridino[1,2-a...	252	C15H12N2O2	025219-83-4	38
3		Imidazole-2-thiol, 1,5-diphenyl-	252	C15H12N2S	136802-77-2	38
4		Phenol, 2,4-dibromo-	250	C6H4Br2O	000615-58-7	37
5		(1-Cyclopentenyl)ferrocene	252	C15H16Fe	012260-67-2	28



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011819\
Data File : BG039216.D
Acq On : 18 Jan 2019 15:16
Operator : JU/SJ
Sample : K1157-09
Misc :
ALS Vial : 5 Sample Multiplier: 1

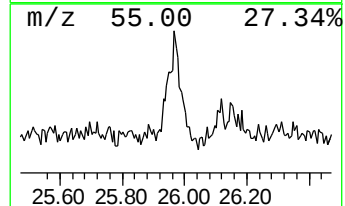
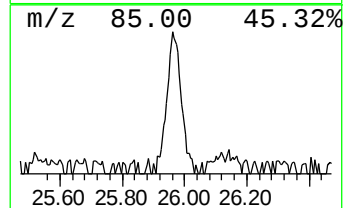
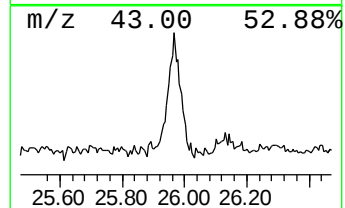
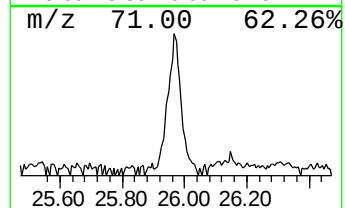
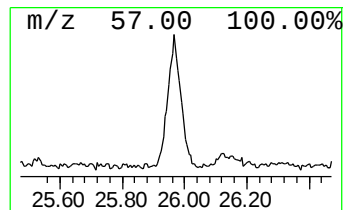
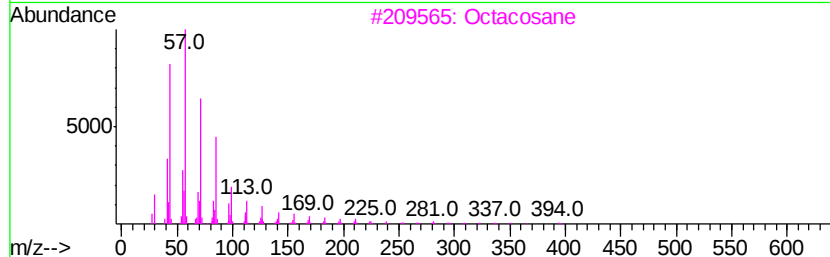
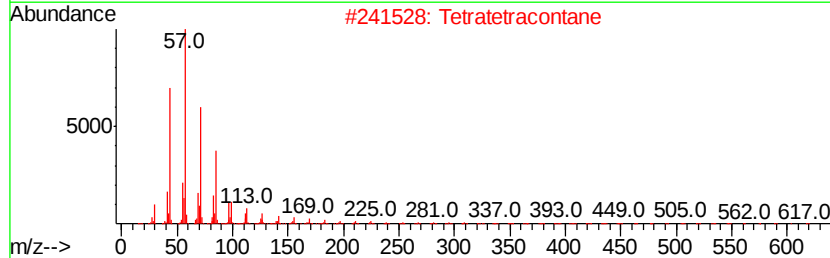
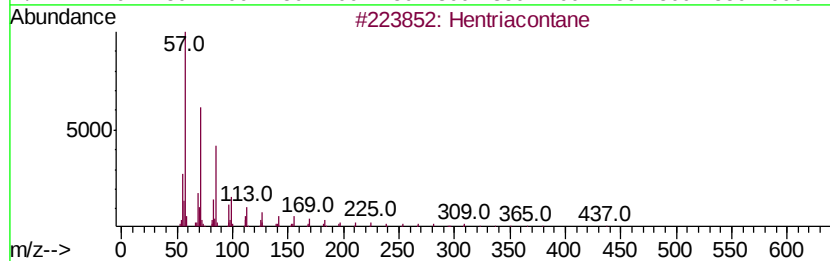
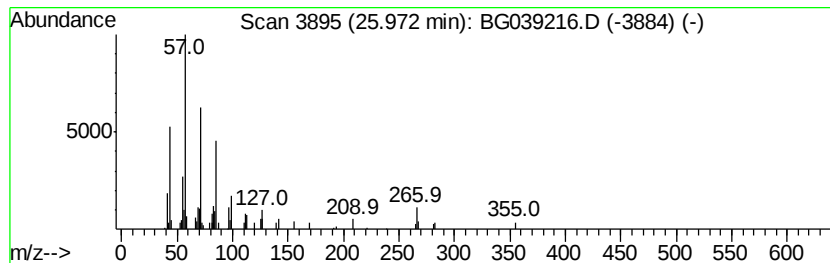
Instrument :
BNA_G
ClientSampleId :
WC-05-011519

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG010919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Hentriacontane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.
25.97	4.55 ng	77503	Perylene-d12		24.93
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hentriacontane	437	C31H64	000630-04-6	90
2	Tetratetracontane	619	C44H90	007098-22-8	90
3	Octacosane	394	C28H58	000630-02-4	87
4	Heptacosane	380	C27H56	000593-49-7	87
5	Hexacosane	366	C26H54	000630-01-3	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG011819\
 Data File : BG039216.D
 Acq On : 18 Jan 2019 15:16
 Operator : JU/SJ
 Sample : K1157-09
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 WC-05-011519

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG010919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.10	6.9	ng	42277	1	8.01	123091	20.0
unknown7.54	7.54	64.6	ng	397688	1	8.01	123091	20.0
Phenanthrene, 2-m...	18.27	3.5	ng	52196	4	17.40	300753	20.0
Phenanthrene, 1-m...	18.32	2.5	ng	37784	4	17.40	300753	20.0
Benzaldehyde, 3,5...	18.47	2.7	ng	40655	4	17.40	300753	20.0
17-Pentatriacontene	21.27	4.9	ng	167903	5	21.67	691119	20.0
Triphenylene, 2-m...	22.41	2.4	ng	82107	5	21.67	691119	20.0
Behenic alcohol	22.44	3.0	ng	104672	5	21.67	691119	20.0
Benzo[e]pyrene	24.62	7.2	ng	122905	6	24.93	340718	20.0
unknown25.01	25.01	3.1	ng	52556	6	24.93	340718	20.0
Hentriacontane	25.97	4.5	ng	77503	6	24.93	340718	20.0