

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP070720\
 Data File : BP002657.D
 Acq On : 07 Jul 2020 14:41
 Operator : CG/JU
 Sample : L3217-03
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E3-41-COMP-1-FLOOR

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_P\METHODS\8270-BP062920.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	5.187	385	391	410	rBV	669709	1063682	14.83%	2.463%
2	6.758	652	658	678	rBV	967558	1679755	23.42%	3.890%
3	7.569	787	796	806	rBV	595307	971628	13.54%	2.250%
4	7.887	842	850	870	rVB	1809406	2981624	41.56%	6.905%
5	8.710	981	990	1011	rBV	1324596	2340819	32.63%	5.421%
6	10.340	1260	1267	1283	rVB	777336	1347977	18.79%	3.122%
7	12.834	1684	1691	1719	rVB	3236424	5219194	72.76%	12.088%
8	13.681	1830	1835	1843	rBV	139730	235646	3.28%	0.546%
9	14.216	1919	1926	1947	rVB	1233627	1815955	25.31%	4.206%
10	15.722	2176	2182	2193	rBV	568133	931375	12.98%	2.157%
11	16.975	2389	2395	2408	rBV	1496322	2170436	30.26%	5.027%
12	18.010	2566	2571	2574	rBV	101982	160539	2.24%	0.372%
13	18.057	2574	2579	2582	rVV2	161736	326729	4.55%	0.757%
14	18.086	2582	2584	2588	rVV	138378	195695	2.73%	0.453%
15	18.139	2588	2593	2598	rVV2	135765	266113	3.71%	0.616%
16	18.198	2599	2603	2604	rVV2	94925	130050	1.81%	0.301%
17	18.216	2604	2606	2609	rVV2	122587	177932	2.48%	0.412%
18	18.251	2609	2612	2617	rVV2	153767	257335	3.59%	0.596%
19	18.298	2617	2620	2623	rVV3	74355	111582	1.56%	0.258%
20	18.339	2623	2627	2636	rVV2	143319	340552	4.75%	0.789%
21	18.410	2636	2639	2646	rVV	95279	165498	2.31%	0.383%
22	18.639	2674	2678	2683	rBV	177667	219761	3.06%	0.509%
23	18.745	2692	2696	2701	rVB4	52749	93511	1.30%	0.217%
24	18.810	2704	2707	2711	rVB	98668	114960	1.60%	0.266%
25	19.375	2800	2803	2807	rBV	114758	117208	1.63%	0.271%
26	19.563	2830	2835	2840	rBV	5598527	7173580	100.00%	16.614%
27	19.610	2840	2843	2851	rVB	210943	292433	4.08%	0.677%
28	19.710	2855	2860	2865	rBV2	468228	924691	12.89%	2.142%
29	19.769	2866	2870	2874	rVB3	318278	438666	6.12%	1.016%
30	19.810	2874	2877	2882	rVB	345056	404702	5.64%	0.937%
31	19.892	2882	2891	2894	rBV3	563498	1156646	16.12%	2.679%
32	19.963	2900	2903	2911	rVB	432756	594695	8.29%	1.377%
33	20.027	2911	2914	2920	rVB	276845	351269	4.90%	0.814%
34	20.380	2971	2974	2978	rVB	161197	161009	2.24%	0.373%

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ClientSampleId :
E3-41-COMP-1-FLOOR

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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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	20.457	2984	2987	2993	rVB	127712	148518	2.07%	0.344%
36	20.827	3046	3050	3061	rVB2	485353	786014	10.96%	1.820%
37	21.022	3079	3083	3088	rVB	1411307	1478134	20.61%	3.423%
38	21.086	3089	3094	3103	rVV	1800032	2372740	33.08%	5.495%
39	21.257	3121	3123	3131	rVB	194314	233721	3.26%	0.541%
40	21.686	3193	3196	3207	rVB	183364	257030	3.58%	0.595%
41	22.145	3271	3274	3280	rVB2	167686	211847	2.95%	0.491%
42	22.645	3356	3359	3363	rBV	133657	165110	2.30%	0.382%
43	23.274	3460	3466	3478	rVB2	1355062	2591505	36.13%	6.002%

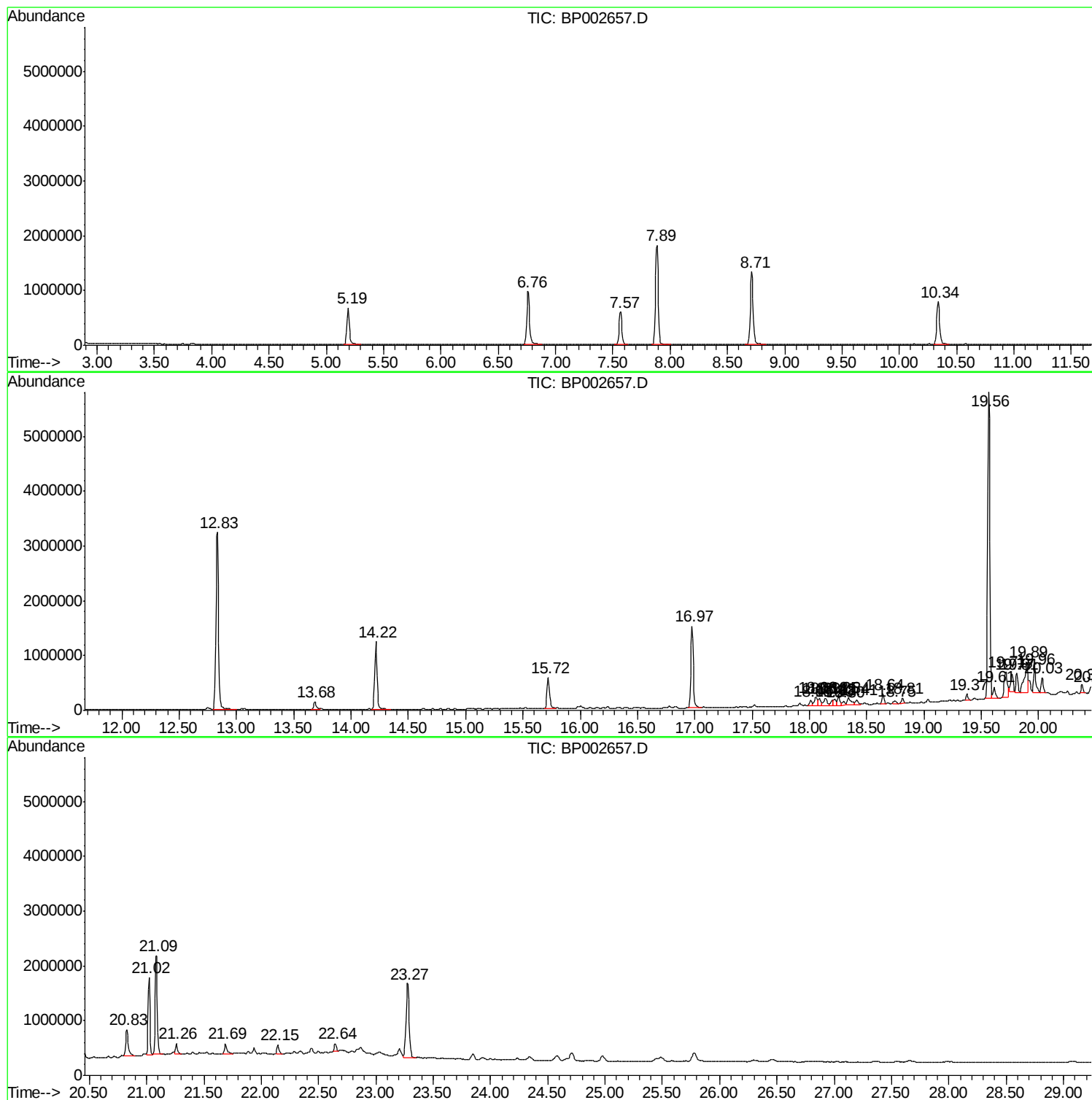
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP062920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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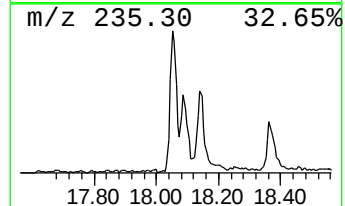
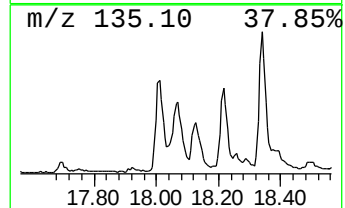
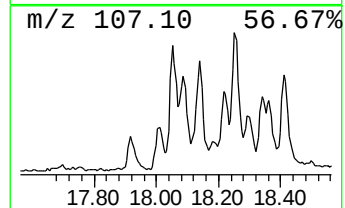
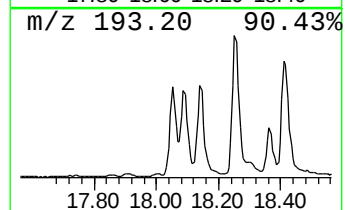
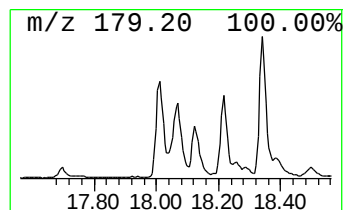
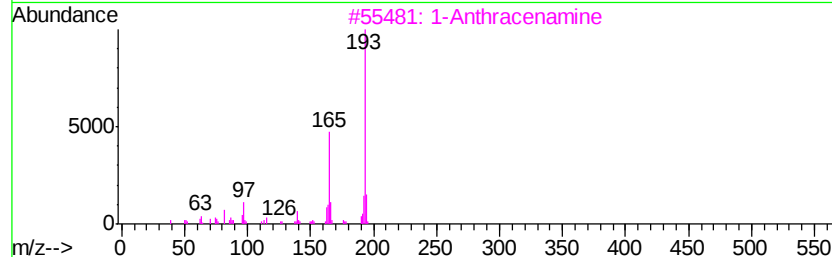
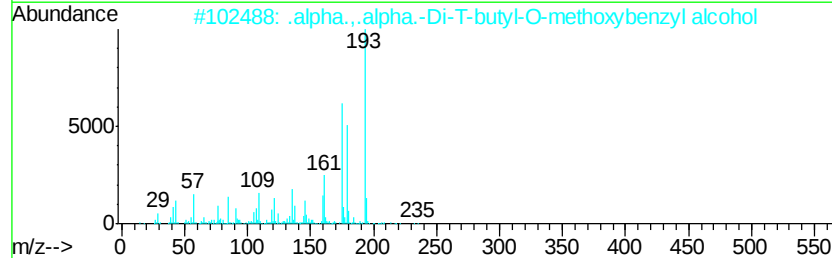
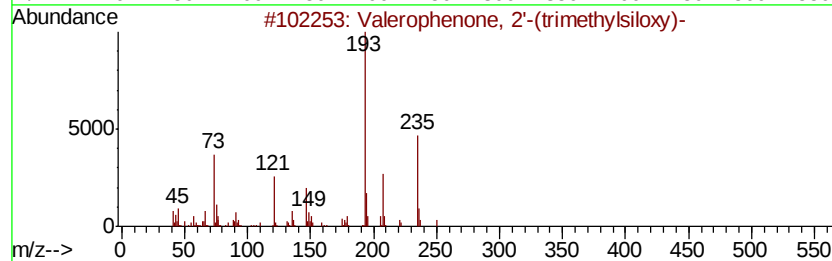
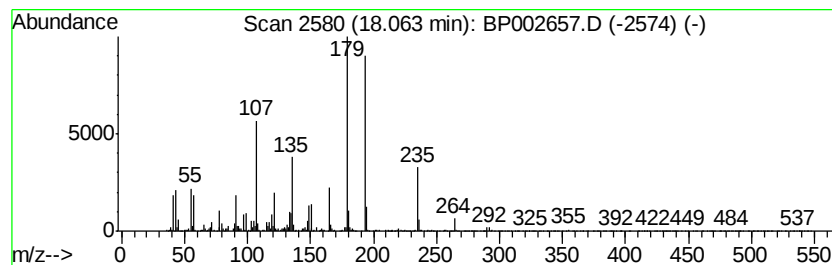
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP062920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 unknown18.06 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.06	3.01 ng	326729	Phenanthrene-d10	16.97	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Valerophenone, 2'-(trimethylsilo...	250	C14H22O2Si	033342-91-5	43
2	.alpha.,.alpha.-Di-T-butyl-O-met...	250	C16H26O2	016750-65-5	37
3	1-Anthracenamine	193	C14H11N	000610-49-1	35
4	Acetophenone, 3'-(trimethylsiloxy)-	208	C11H16O2Si	033342-86-8	32
5	Acetamide, N-9-phenanthrenyl-	235	C16H13NO	004235-09-0	32



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Instrument :
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ClientSampleId :
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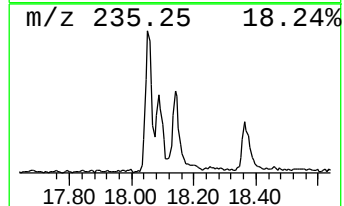
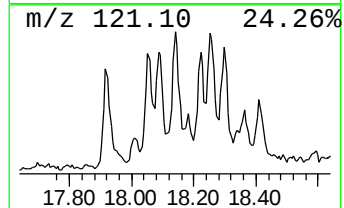
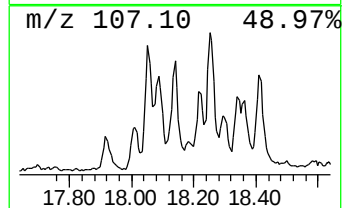
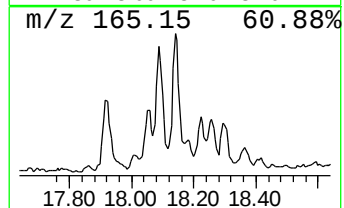
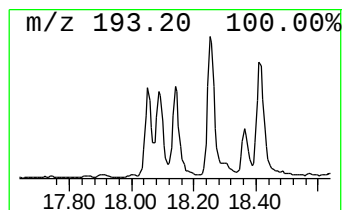
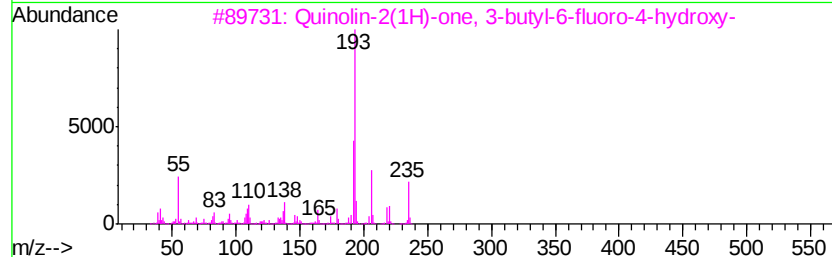
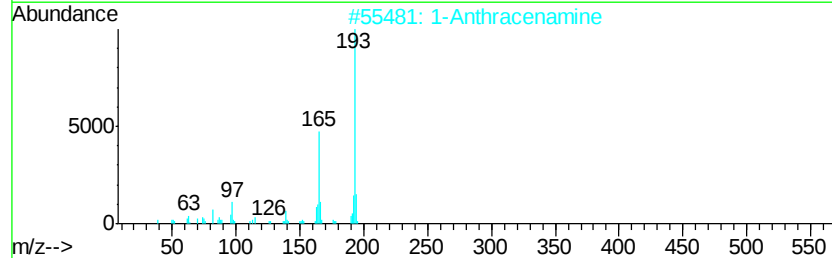
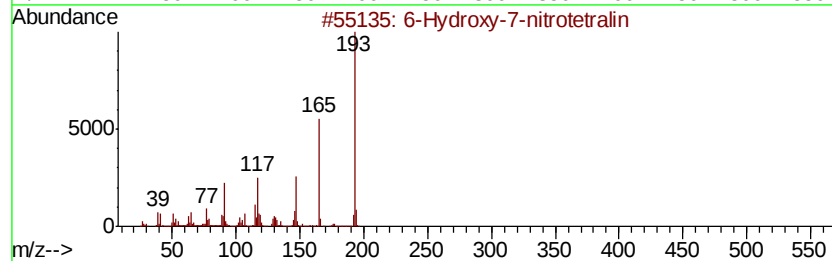
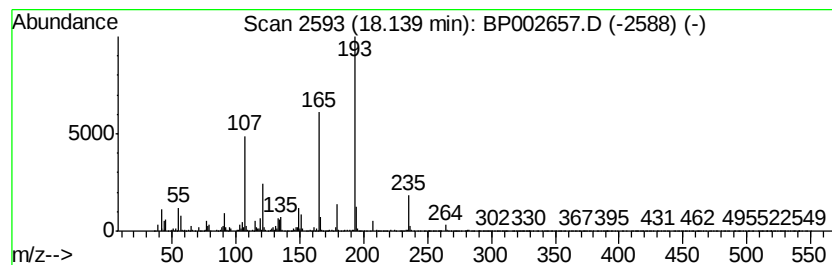
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 unknown18.14 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.14	2.45 ng	266113	Phenanthrene-d10	16.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	6-Hydroxy-7-nitrotetralin	193	C10H11NO3	006240-79-5	47
2		1-Anthracenamine	193	C14H11N	000610-49-1	47
3		Quinolin-2(1H)-one, 3-butyl-6-fluoro-4-hydroxy-	235	C13H14FN02	279691-75-7	43
4		[1,1'-Biphenyl]-4-acetonitrile	193	C14H11N	031603-77-7	38
5		1H-Indole, 2-phenyl-	193	C14H11N	000948-65-2	38



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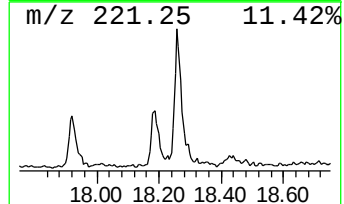
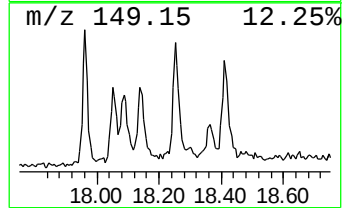
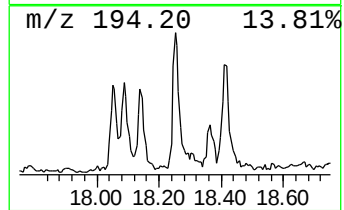
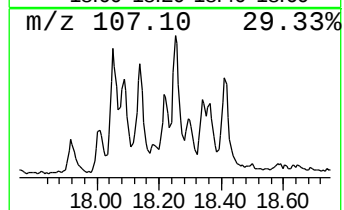
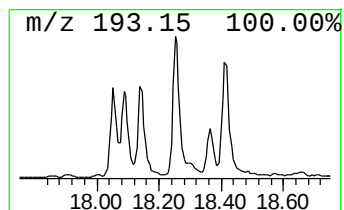
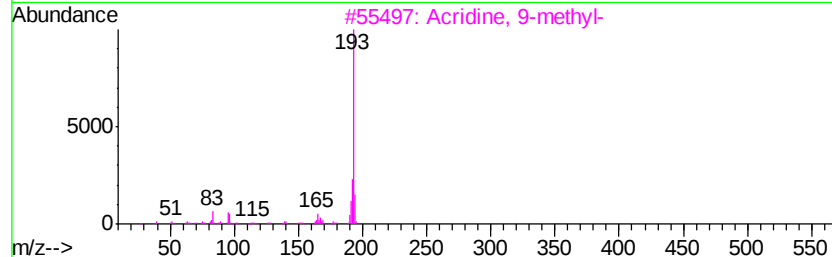
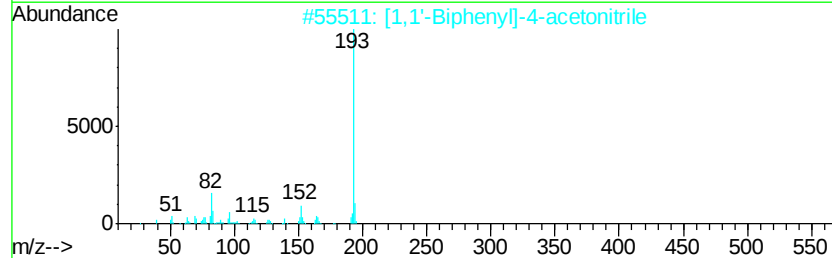
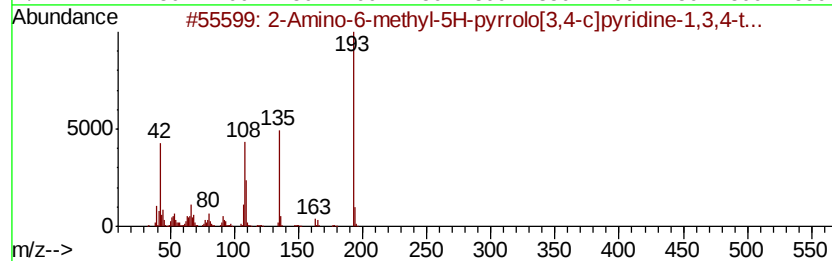
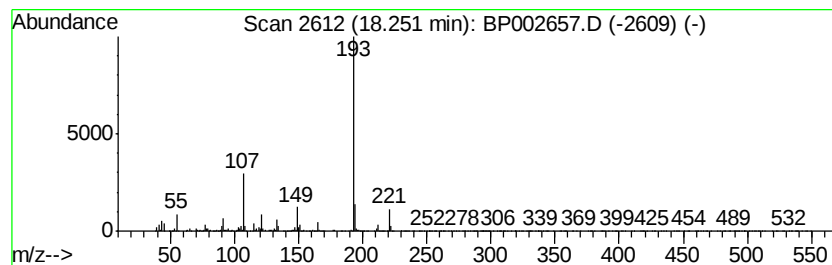
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP062920.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 2-Amino-6-methyl-5H-pyrrolo... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.	
18.25	2.37 ng	257335	Phenanthrene-d10		16.97	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Amino-6-methyl-5H-pyrrolo[3,4-...	193	C8H7N3O3	1000300-48-5	53	
2	[1,1'-Biphenyl]-4-acetonitrile	193	C14H11N	031603-77-7	50	
3	Acridine, 9-methyl-	193	C14H11N	000611-64-3	49	
4	1,5-Methano-8H-pyrido[1,2-a][1,5...	234	C14H22N2O	000550-43-6	47	
5	Anthracene, 9-ethyl-9,10-dihydro...	222	C17H18	036778-20-8	40	



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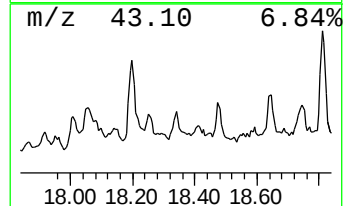
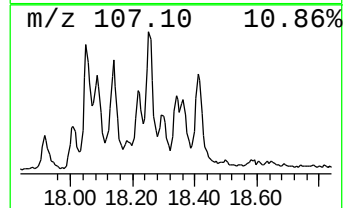
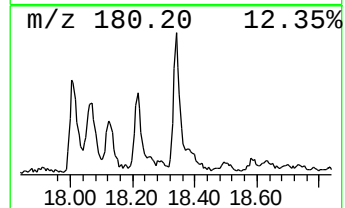
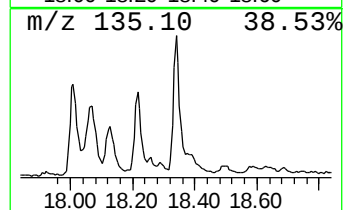
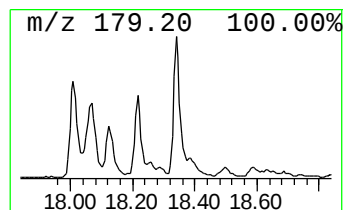
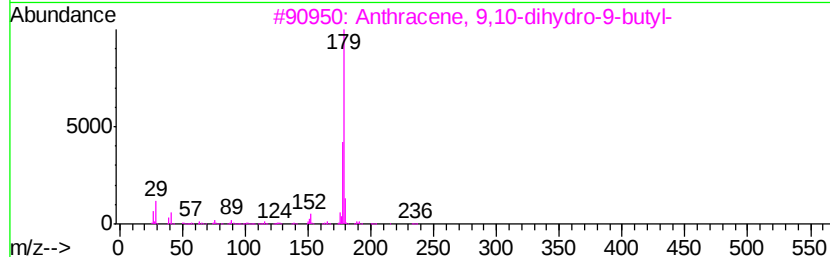
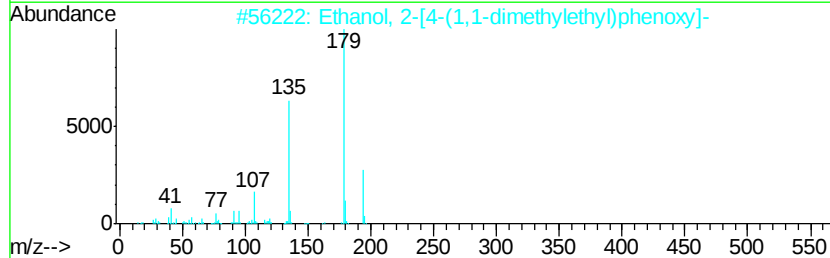
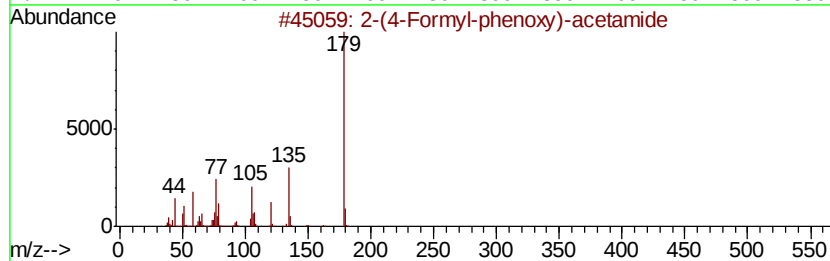
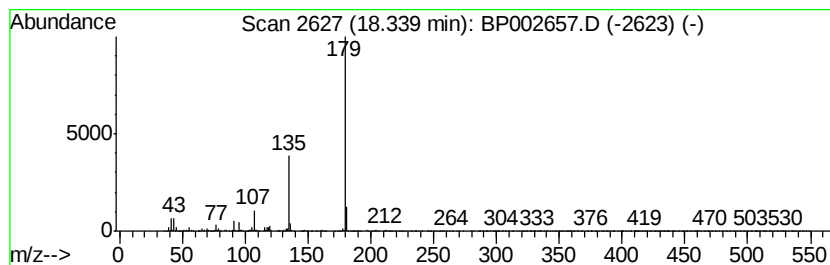
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 2-(4-Formyl-phenoxy)-acetamide Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.34	3.14 ng	340552	Phenanthrene-d10	16.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-(4-Formyl-phenoxy)-acetamide	179	C9H9NO3	135857-20-4	72
2		Ethanol, 2-[4-(1,1-dimethylethyl)phenoxy]-	194	C12H18O2	000713-46-2	52
3		Anthracene, 9,10-dihydro-9-butyl-	236	C18H20	010394-60-2	43
4		(p-(Allyloxy)carbonyl)phenoxy)ace...	236	C12H12O5	1000241-83-7	37
5		Acetic acid, [2-[(2-propenylamin...	235	C12H13NO4	000119-45-9	37



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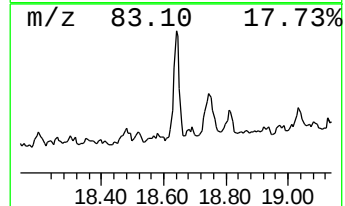
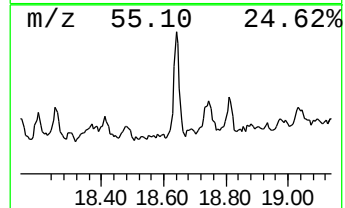
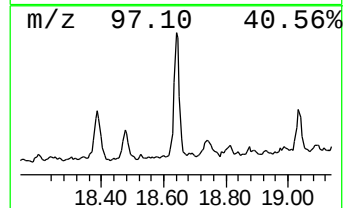
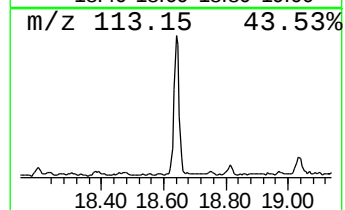
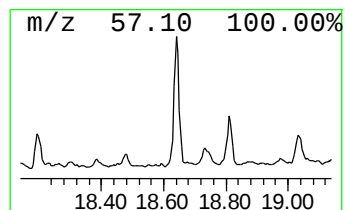
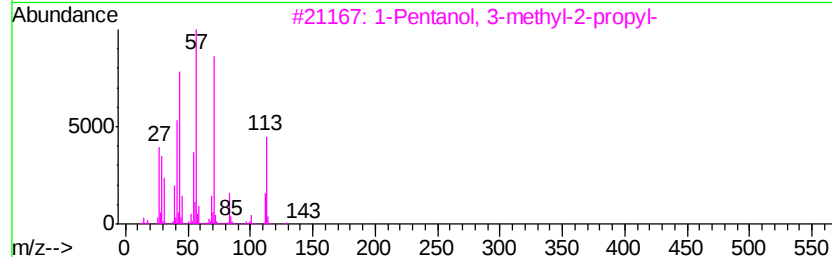
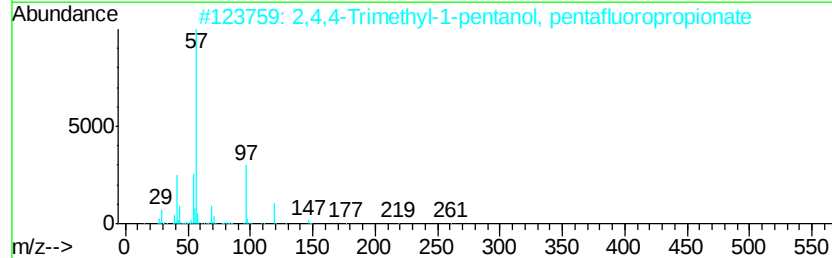
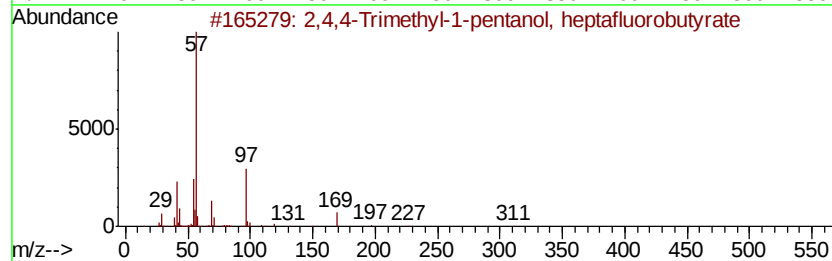
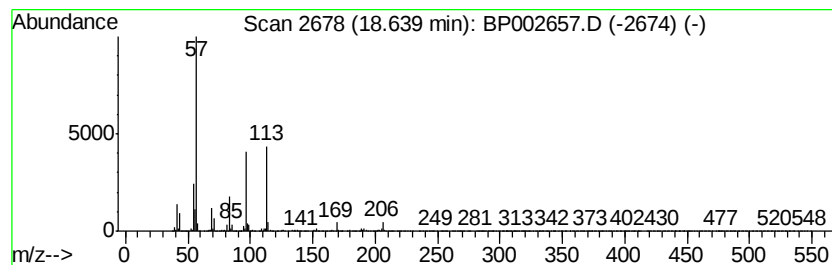
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 unknown18.64 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.64	2.03 ng	219761	Phenanthrene-d10	16.97

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,4-Trimethyl-1-pentanol, hept...	326	C12H17F7O2	1000365-01-4	43		
2	2,4,4-Trimethyl-1-pentanol, pent...	276	C11H17F5O2	1000365-51-0	38		
3	1-Pentanol, 3-methyl-2-propyl-	144	C9H20O	054004-40-9	37		
4	2,4,4,6,6,8,8-Heptamethyl-1-nonene	224	C16H32	015796-04-0	33		
5	Octane, 2-bromo-	192	C8H17Br	000557-35-7	32		



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP070720\
Data File : BP002657.D
Acq On : 07 Jul 2020 14:41
Operator : CG/JU
Sample : L3217-03
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E3-41-COMP-1-FLOOR

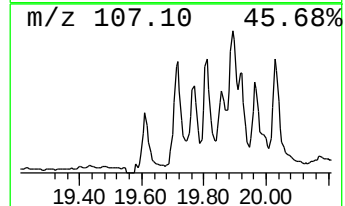
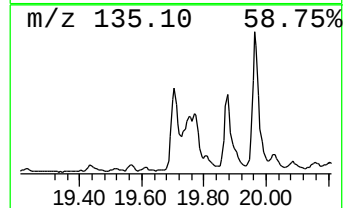
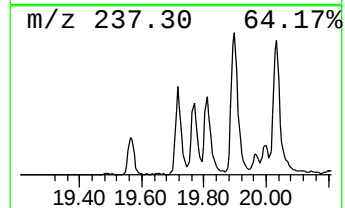
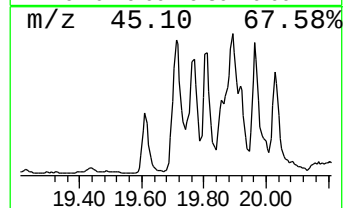
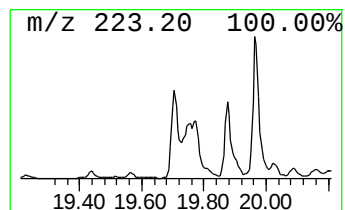
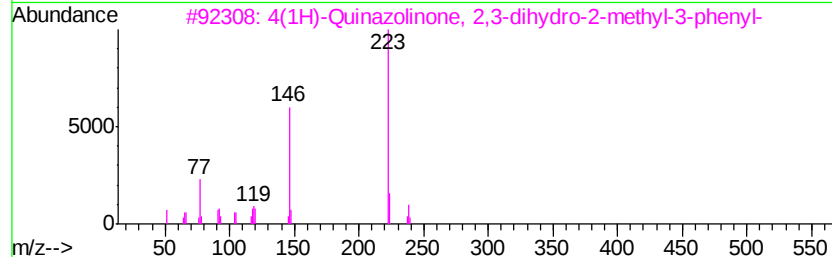
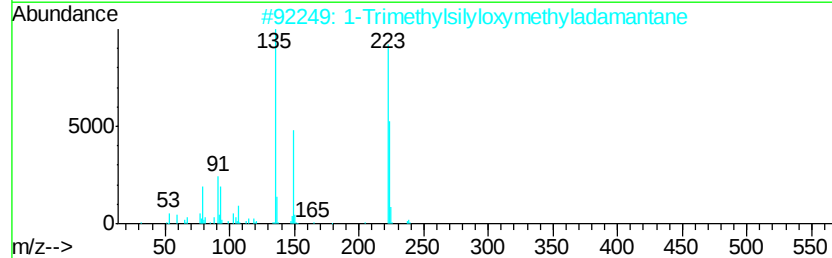
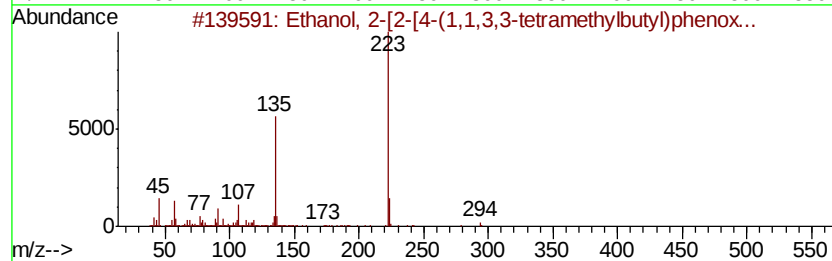
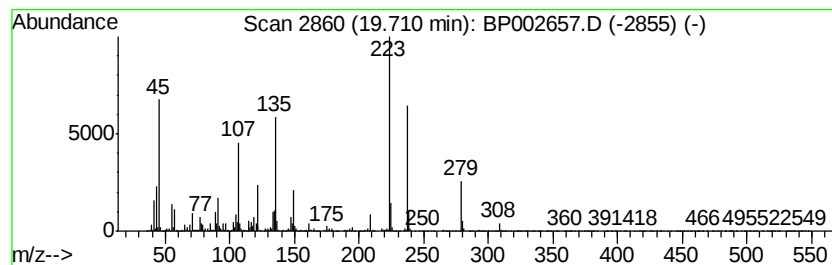
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown19.71 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.71	7.79 ng	924691	Chrysene-d12	21.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol, 2-[2-[4-(1,1,3,3-tetramethylbutyl)phenoxy]ethyl]phenyl	294	C18H30O3	002315-61-9	38
2		1-Trimethylsilyloxymethyladamantane	238	C14H26OSi	079671-68-4	35
3		4(1H)-Quinazolinone, 2,3-dihydro-2-methyl-3-phenyl-	238	C15H14N2O	017761-74-9	25
4		1-(2-Fluoro-4-nitro-phenyl)-2-methyl-2H-imidazole	238	C12H15FN2O2	1000278-20-9	22
5		1H-Benzimidazole, 2-(5-ethyl-2-pyridyl)-	223	C14H13N3	067273-43-2	18



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP070720\
Data File : BP002657.D
Acq On : 07 Jul 2020 14:41
Operator : CG/JU
Sample : L3217-03
Misc :
ALS Vial : 11 Sample Multiplier: 1

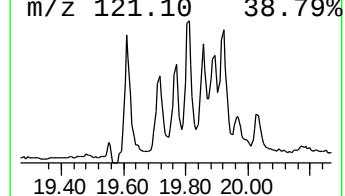
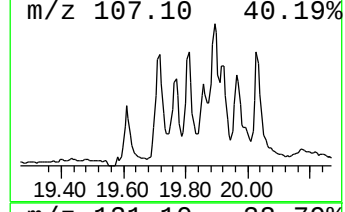
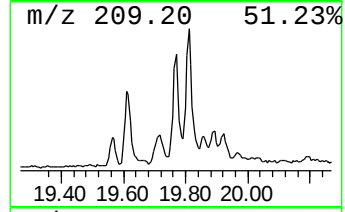
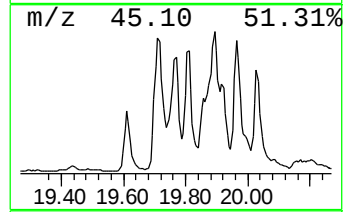
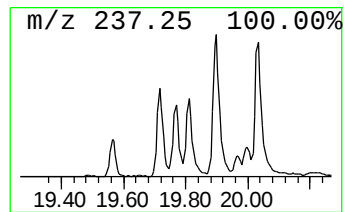
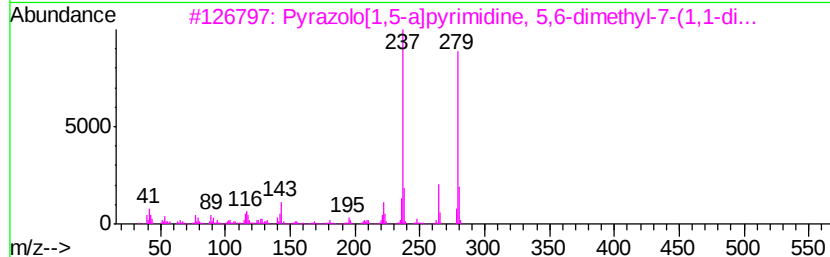
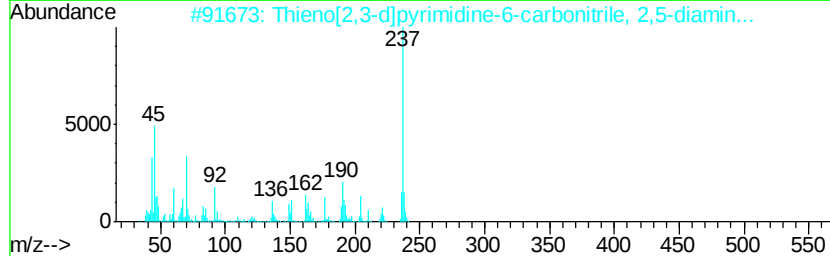
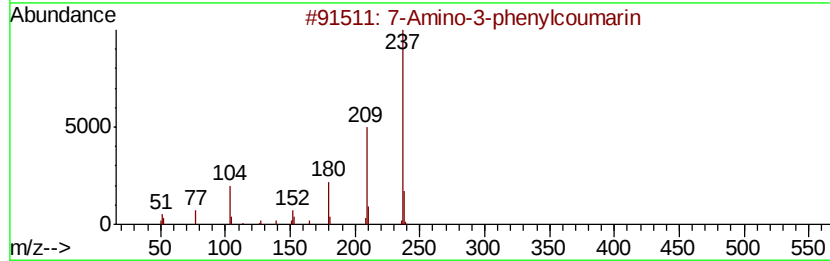
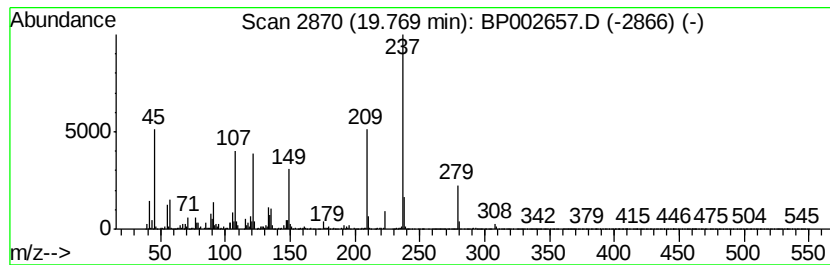
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ClientSampleId :
E3-41-COMP-1-FLOOR

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP062920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown19.77 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.		
19.77	3.70 ng	438666	Chrysene-d12	21.09		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	7-Amino-3-phenylcoumarin	237	C15H11N02	004108-61-6	43	
2	Thieno[2,3-d]pyrimidine-6-carbon...	237	C8H7N5S2	1000311-13-6	38	
3	Pvrazolo[1,5-alpyrimidine, 5,6-d...	279	C18H21N3	1000274-10-7	35	
4	Coumarin-6-ol, 3,4-dihydro-4,4-d...	237	C11H11N05	1000126-61-0	27	
5	4,6-Difluoro-2-(4-methoxyphenyla...	237	C11H9F2N3O	1000070-53-0	27	



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP070720\
Data File : BP002657.D
Acq On : 07 Jul 2020 14:41
Operator : CG/JU
Sample : L3217-03
Misc :
ALS Vial : 11 Sample Multiplier: 1

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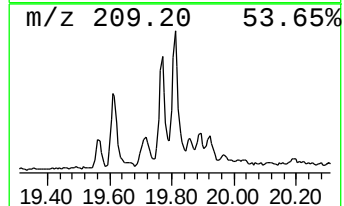
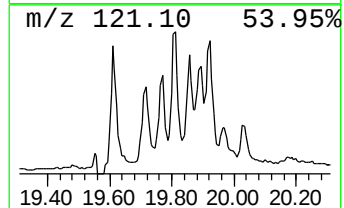
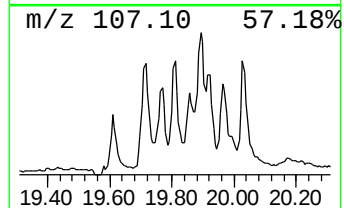
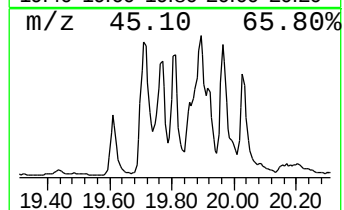
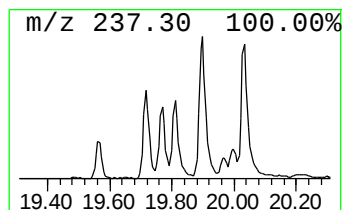
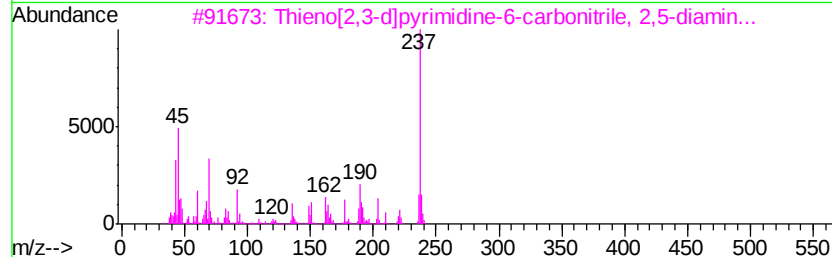
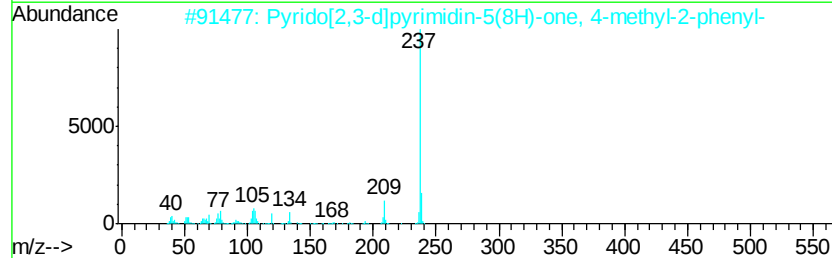
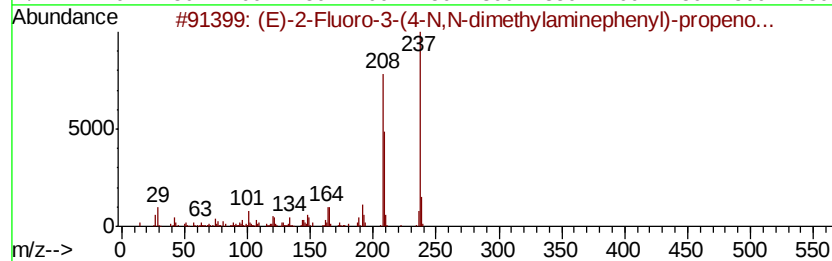
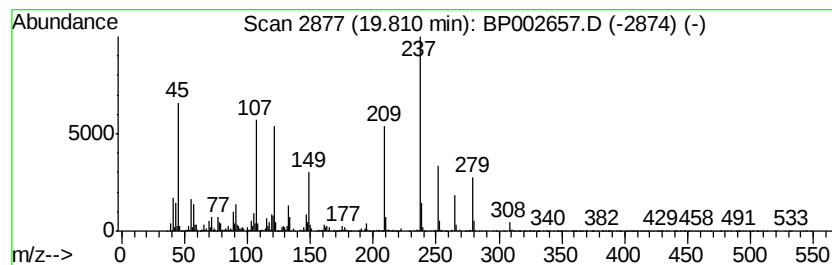
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 unknown19.81 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.81	3.41 ng	404702	Chrysene-d12	21.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-2-Fluoro-3-(4-N,N-dimethylam...	237	C13H16FN02	110915-65-6	38
2		Pyrido[2,3-d]pyrimidin-5(8H)-one...	237	C14H11N3O	161465-98-1	38
3		Thieno[2,3-d]pyrimidine-6-carbon...	237	C8H7N5S2	1000311-13-6	35
4		Coumarine, 6-amino-3-phenyl-	237	C15H11N02	021408-16-2	35
5		Pyrazolo[1,5-a]pyrimidine, 5,6-d...	279	C18H21N3	1000273-96-0	27



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP070720\
Data File : BP002657.D
Acq On : 07 Jul 2020 14:41
Operator : CG/JU
Sample : L3217-03
Misc :
ALS Vial : 11 Sample Multiplier: 1

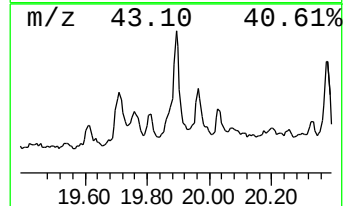
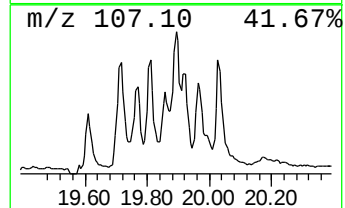
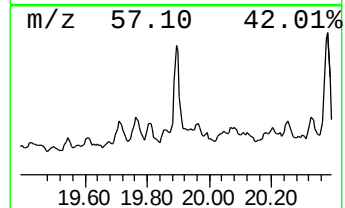
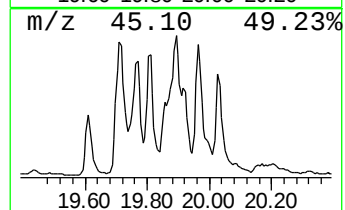
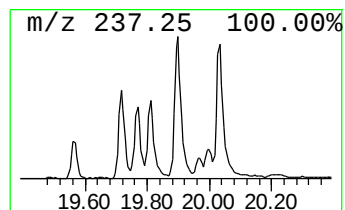
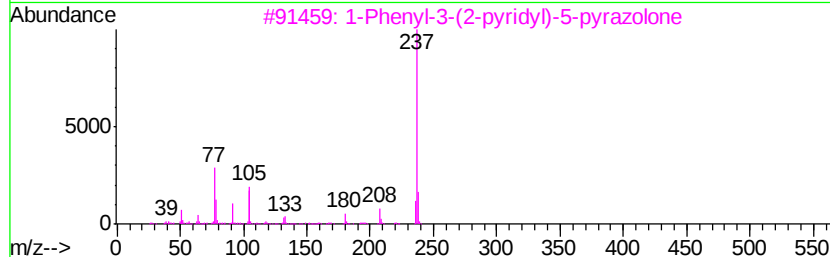
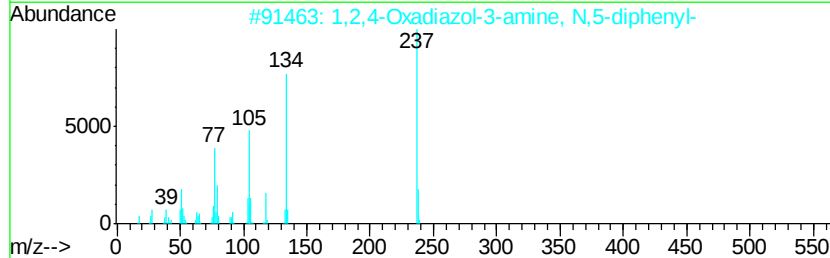
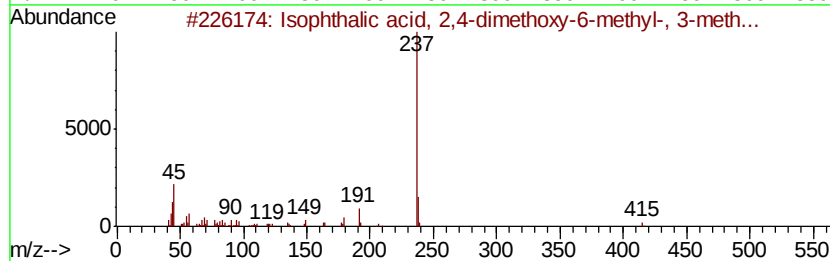
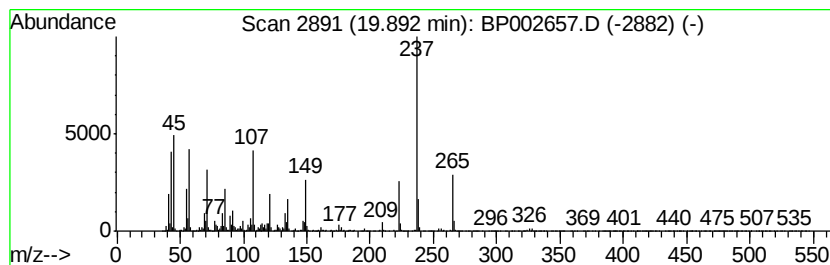
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ClientSampleId :
E3-41-COMP-1-FLOOR

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP062920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown19.89 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.		
19.89	9.75 ng	1156650	Chrysene-d12	21.09		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Isophthalic acid, 2,4-dimethoxyv...	446	C23H26O9	019314-77-3	32
2		1,2,4-Oxadiazol-3-amine, N,5-dip...	237	C14H11N3O	058599-06-7	22
3		1-Phenvl-3-(2-pyridvl)-5-pvrazolone	237	C14H11N3O	021683-60-3	22
4		Propanamide, N-(2,5-dimethoxyvphe...	237	C13H19NO3	1000307-19-9	22
5		(E)-2-Fluoro-3-(4-N,N-dimethylam...	237	C13H16FN02	110915-65-6	22



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP070720\
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Acq On : 07 Jul 2020 14:41
Operator : CG/JU
Sample : L3217-03
Misc :
ALS Vial : 11 Sample Multiplier: 1

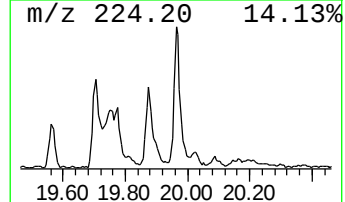
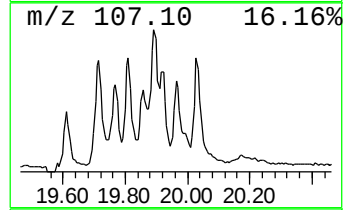
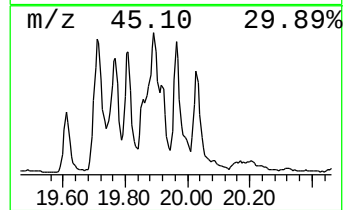
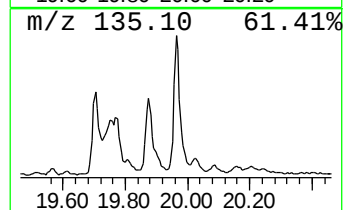
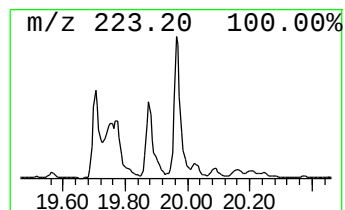
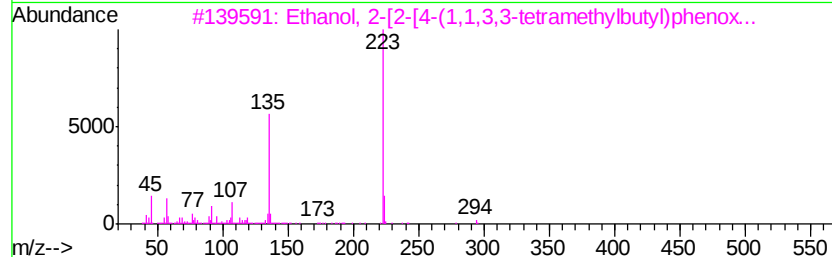
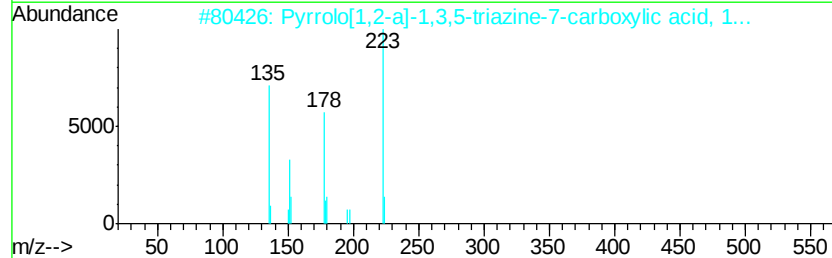
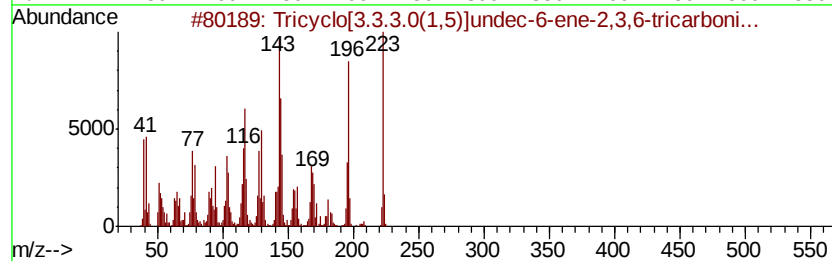
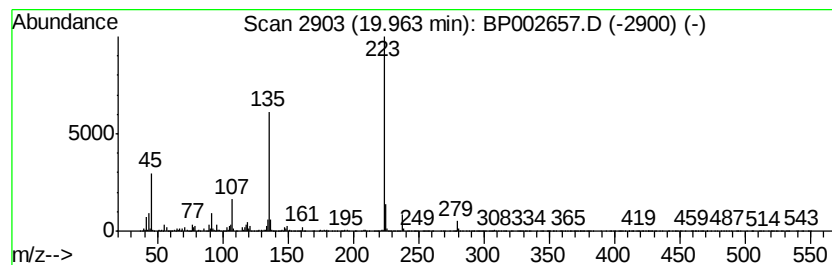
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Tricyclo[3.3.3.0(1,5)]undec... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.96	5.01 ng	594695	Chrysene-d12	21.09	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tricyclo[3.3.3.0(1,5)]undec-6-en...	223	C14H13N3	118894-71-6	86
2	Pyrrolo[1,2-a]-1,3,5-triazine-7-...	223	C9H9N3O4	054449-89-7	86
3	Ethanol, 2-[2-[4-(1,1,3,3-tetram...	294	C18H30O3	002315-61-9	80
4	Glycine, N-(m-anisoyl)-, methyl ...	223	C11H13NO4	1000299-70-7	59
5	Malonodinitrile, 2-(3-dimethylam...	223	C14H13N3	065996-13-6	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP070720\
Data File : BP002657.D
Acq On : 07 Jul 2020 14:41
Operator : CG/JU
Sample : L3217-03
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
E3-41-COMP-1-FLOOR

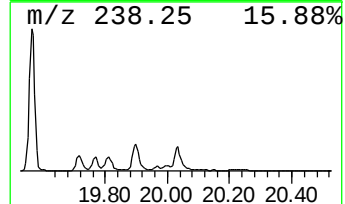
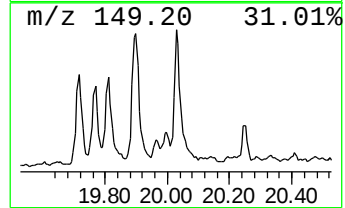
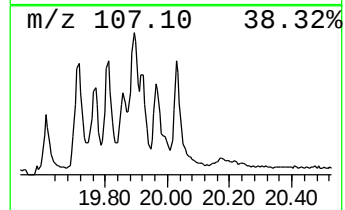
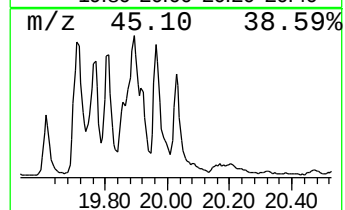
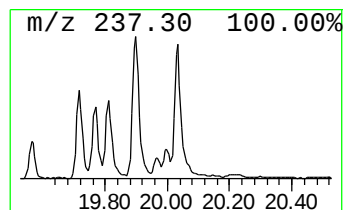
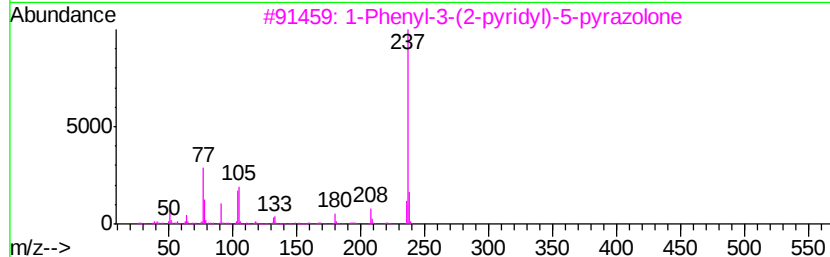
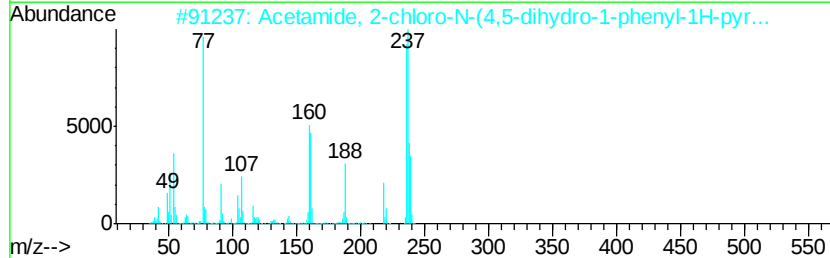
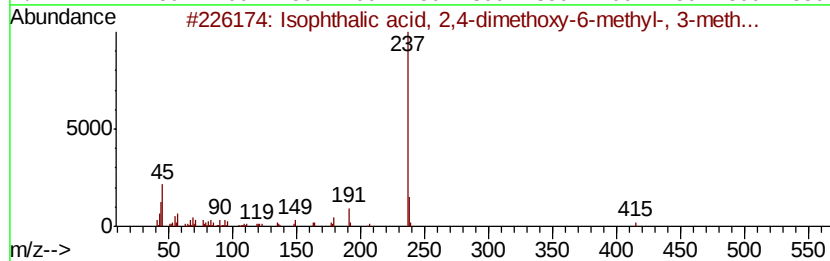
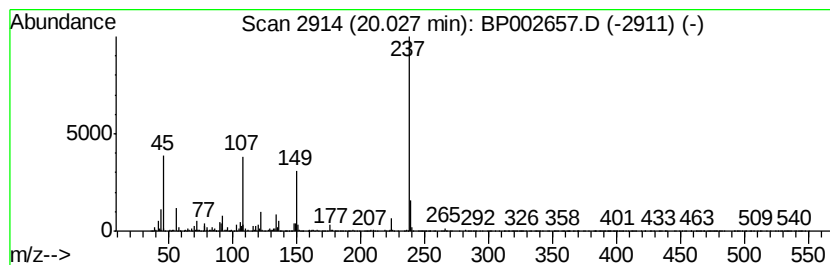
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown20.03 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.03	2.96 ng	351269	Chrysene-d12	21.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Isophthalic acid, 2,4-dimethoxyv-...	446	C23H26O9	019314-77-3	40
2		Acetamide, 2-chloro-N-(4,5-dihyd...	237	C11H12ClN3O	1000337-23-3	38
3		1-Phenvl-3-(2-pyridyl)-5-pyrazolone	237	C14H11N3O	021683-60-3	27
4		Indolo[2,3-b]quinoxaline, 1-fluoro-	237	C14H8FN3	296244-71-8	16
5		Pyrido[2,3-d]pyrimidin-5(8H)-one...	237	C14H11N3O	161465-98-1	16



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP070720\
 Data File : BP002657.D
 Acq On : 07 Jul 2020 14:41
 Operator : CG/JU
 Sample : L3217-03
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E3-41-COMP-1-FLOOR

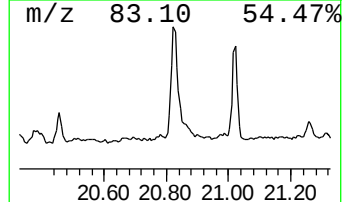
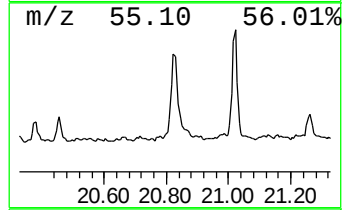
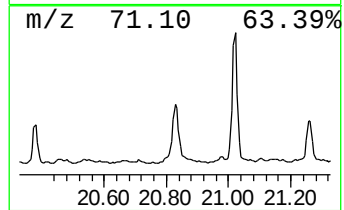
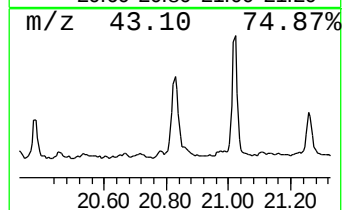
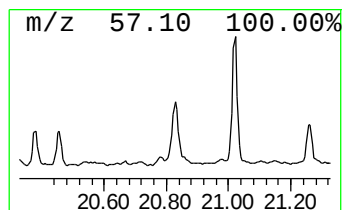
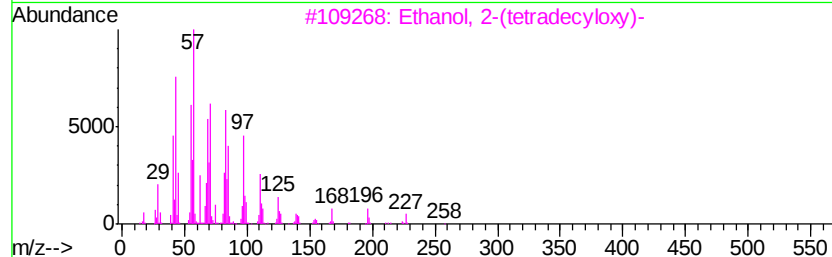
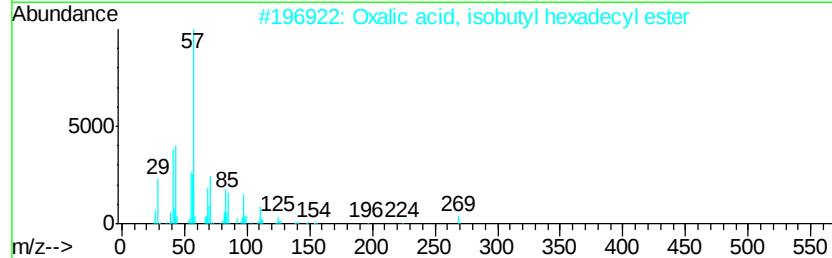
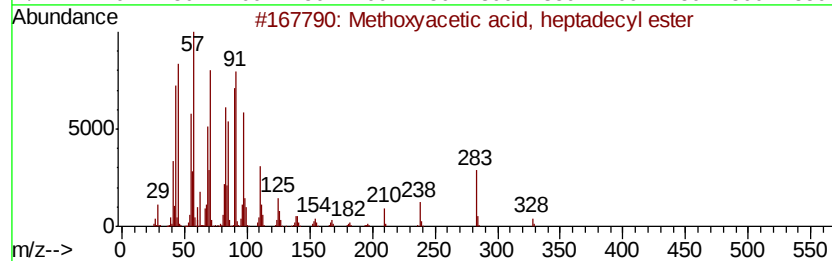
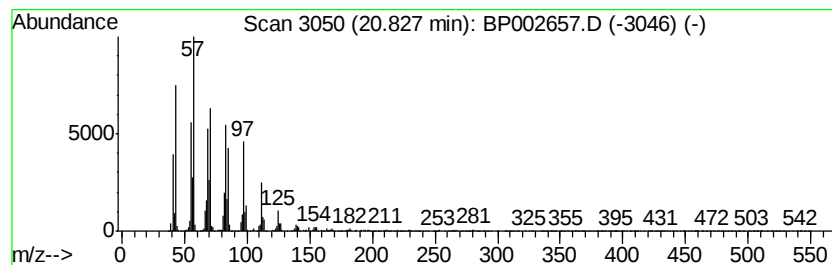
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 13 Methoxyacetic acid, heptade... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.83	6.63 ng	786014	Chrysene-d12	21.09

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methoxvacetic acid, heptadecyl e...	328	C20H40O3	1000282-99-1	91
2			Oxalic acid, isobutyl hexadecyl ...	370	C22H42O4	1000309-38-1	91
3			Ethanol, 2-(tetradecyl)-	258	C16H34O2	002136-70-1	90
4			1-Decanol, 2-hexyl-	242	C16H34O	002425-77-6	89
5			9-Eicosene, (E)-	280	C20H40	074685-29-3	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP070720\
Data File : BP002657.D
Acq On : 07 Jul 2020 14:41
Operator : CG/JU
Sample : L3217-03
Misc :
ALS Vial : 11 Sample Multiplier: 1

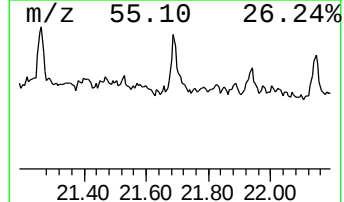
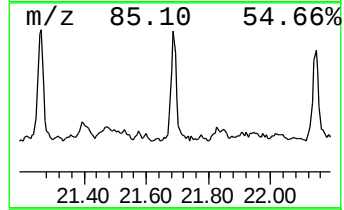
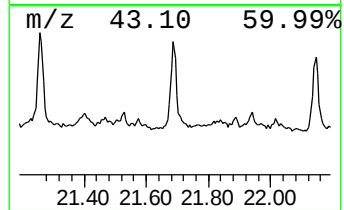
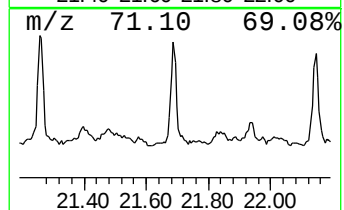
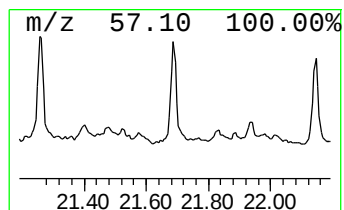
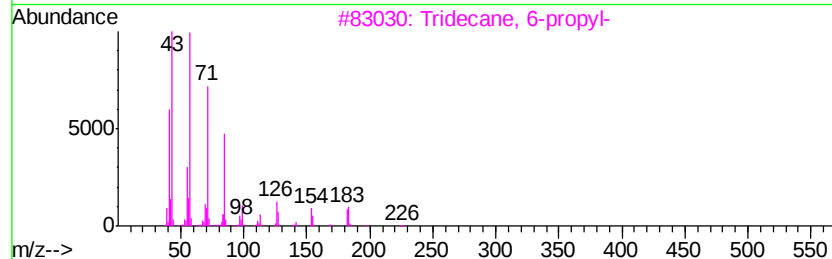
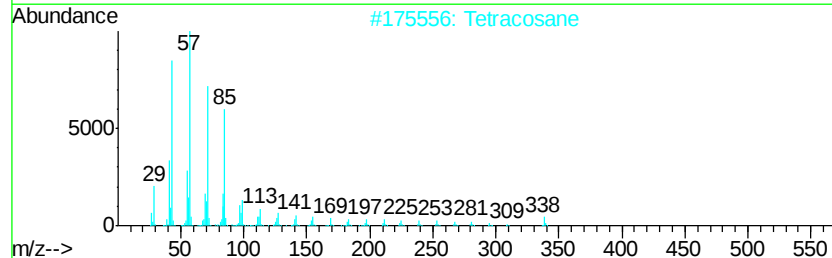
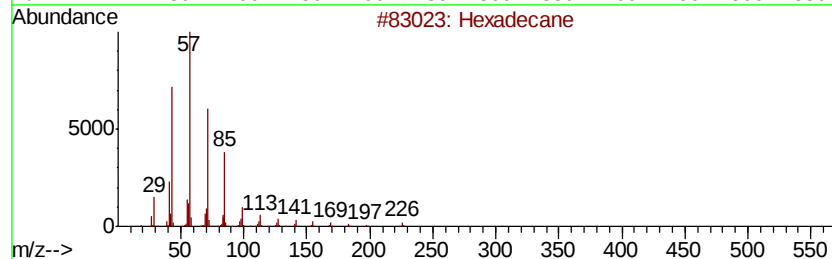
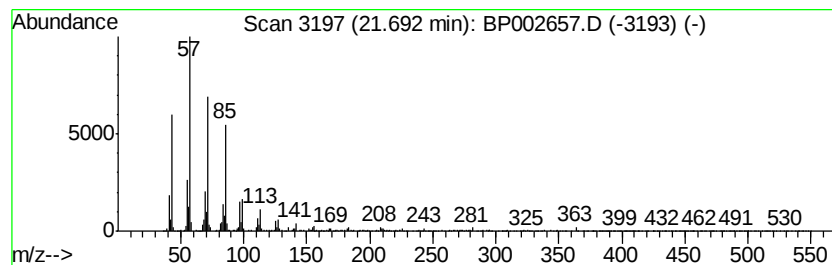
Instrument :
BNA_P
ClientSampleId :
E3-41-COMP-1-FLOOR

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP062920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Hexadecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.69	2.17 ng	257030	Chrysene-d12	21.09
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Hexadecane	226 C16H34	000544-76-3	93
2	Tetracosane	338 C24H50	000646-31-1	93
3	Tridecane, 6-propyl-	226 C16H34	055045-10-8	87
4	Heneicosane	296 C21H44	000629-94-7	87
5	Octadecane	254 C18H38	000593-45-3	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP070720\
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 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E3-41-COMP-1-FLOOR

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP062920.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
unknown18.06	18.06	3.0	ng	326729	4	16.97	2170440	20.0
unknown18.14	18.14	2.5	ng	266113	4	16.97	2170440	20.0
2-Amino-6-methyl-...	18.25	2.4	ng	257335	4	16.97	2170440	20.0
2-(4-Formyl-pheno...	18.34	3.1	ng	340552	4	16.97	2170440	20.0
unknown18.64	18.64	2.0	ng	219761	4	16.97	2170440	20.0
unknown19.71	19.71	7.8	ng	924691	5	21.09	2372740	20.0
unknown19.77	19.77	3.7	ng	438666	5	21.09	2372740	20.0
unknown19.81	19.81	3.4	ng	404702	5	21.09	2372740	20.0
unknown19.89	19.89	9.8	ng	1156650	5	21.09	2372740	20.0
Tricyclo[3.3.3.0(...	19.96	5.0	ng	594695	5	21.09	2372740	20.0
unknown20.03	20.03	3.0	ng	351269	5	21.09	2372740	20.0
Methoxyacetic aci...	20.83	6.6	ng	786014	5	21.09	2372740	20.0
Hexadecane	21.69	2.2	ng	257030	5	21.09	2372740	20.0