

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081919\
 Data File : BG042344.D
 Acq On : 20 Aug 2019 2:17
 Operator : HP/JU
 Sample : K4290-06
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 WP-1

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-BG081919.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	3.124	3	6	21	rVB	1324087	1938674	47.61%	8.568%
2	5.092	333	341	354	rBV	1562991	2497880	61.34%	11.039%
3	5.568	414	422	442	rBV	546632	984166	24.17%	4.349%
4	7.172	686	695	713	rBV	607821	1130888	27.77%	4.998%
5	7.542	749	758	779	rBV	710611	1332432	32.72%	5.889%
6	8.006	829	837	844	rBV	166775	281408	6.91%	1.244%
7	8.329	884	892	903	rBV	648539	1137743	27.94%	5.028%
8	9.170	1028	1035	1050	rBV	471310	888616	21.82%	3.927%
9	10.827	1308	1317	1328	rBV	205747	398389	9.78%	1.761%
10	11.978	1505	1513	1525	rBV	36598	70053	1.72%	0.310%
11	13.147	1705	1712	1728	rBV5	13671	68397	1.68%	0.302%
12	13.283	1728	1735	1747	rVV	1602579	2522723	61.95%	11.149%
13	14.663	1964	1970	1982	rVB	395592	642116	15.77%	2.838%
14	15.821	2161	2167	2181	rBV2	37614	71386	1.75%	0.315%
15	16.167	2220	2226	2242	rBV	141542	396890	9.75%	1.754%
16	17.331	2418	2424	2429	rBV2	115684	175126	4.30%	0.774%
17	17.413	2429	2438	2447	rVB2	522407	838753	20.60%	3.707%
18	17.695	2481	2486	2493	rBV3	39503	57486	1.41%	0.254%
19	18.706	2653	2658	2665	rBV	62951	92529	2.27%	0.409%
20	20.028	2877	2883	2896	rBV	3168169	4072139	100.00%	17.996%
21	21.015	3047	3051	3056	rBV	44378	55982	1.37%	0.247%
22	21.338	3101	3106	3121	rBV	105116	228798	5.62%	1.011%
23	21.690	3160	3166	3178	rVB	598479	970469	23.83%	4.289%
24	21.890	3195	3200	3205	rBV	53082	77098	1.89%	0.341%
25	22.207	3250	3254	3261	rVB	38004	64098	1.57%	0.283%
26	22.501	3299	3304	3309	rBV	61373	102009	2.51%	0.451%
27	23.206	3417	3424	3432	rVB	66026	129916	3.19%	0.574%
28	24.017	3556	3562	3572	rVB2	62676	133150	3.27%	0.588%
29	24.934	3707	3718	3740	rBV2	351636	1168169	28.69%	5.163%
30	26.126	3915	3921	3931	rVB2	35418	100207	2.46%	0.443%

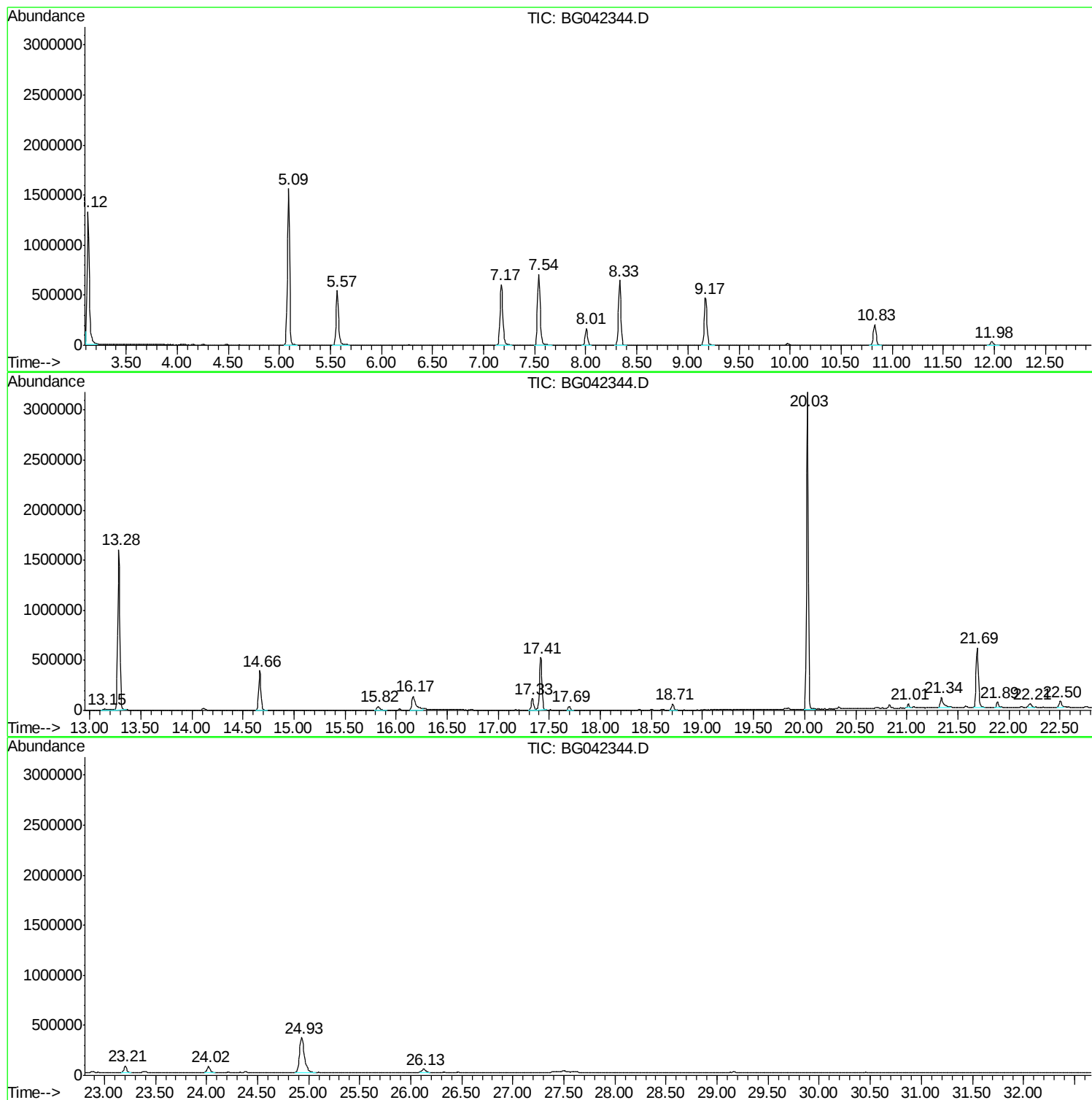
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Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG081919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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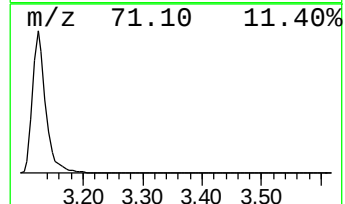
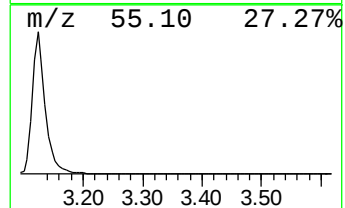
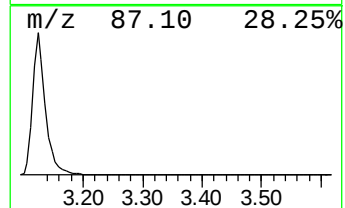
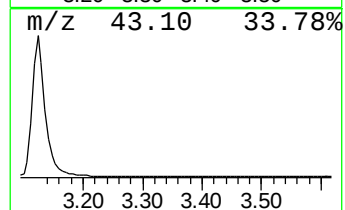
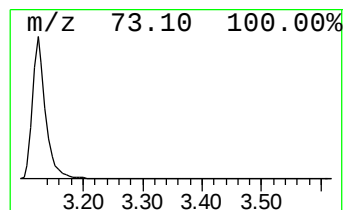
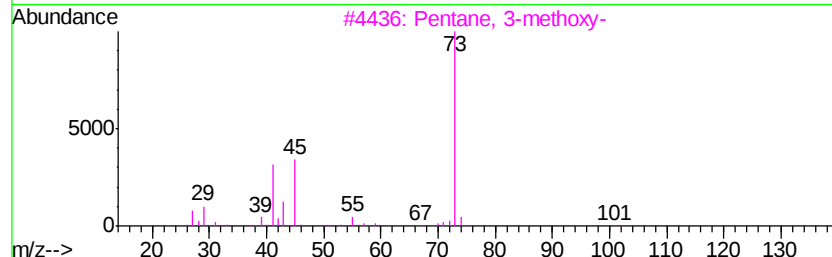
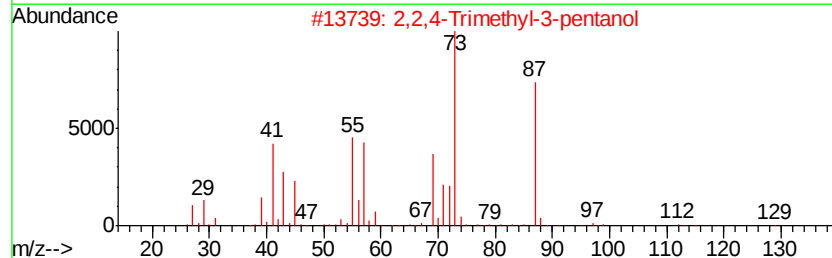
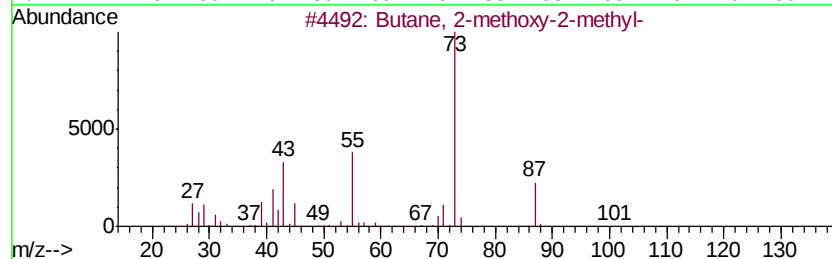
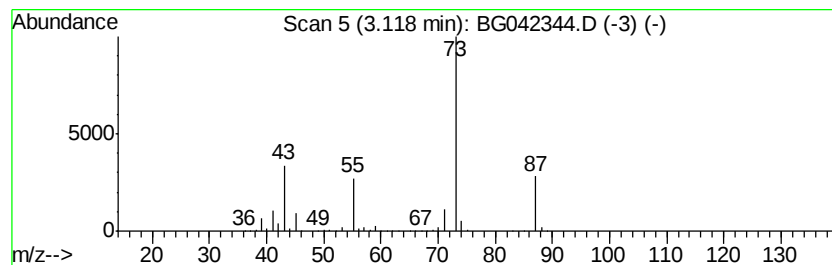
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.12	137.78 ng	1938670	1,4-Dichlorobenzene-d4	8.01

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	25
4		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9
5		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9



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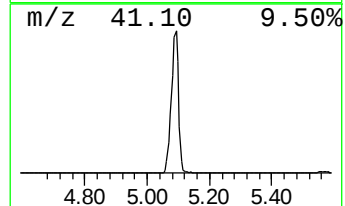
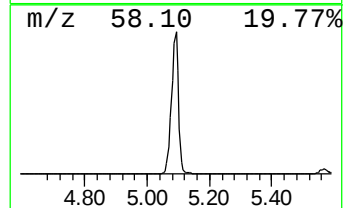
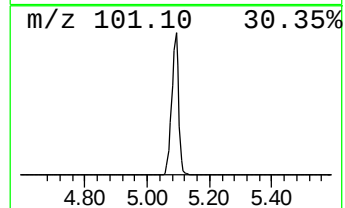
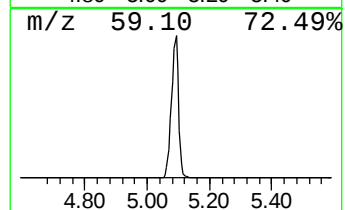
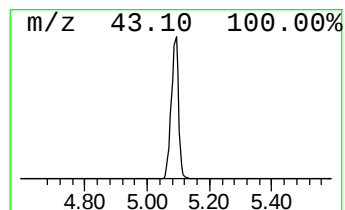
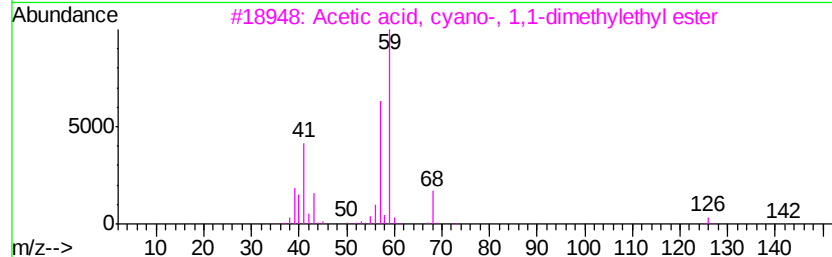
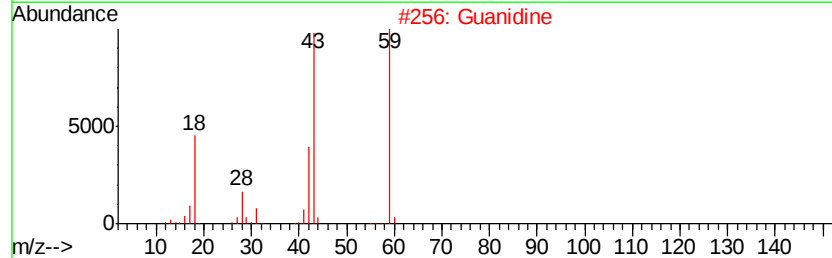
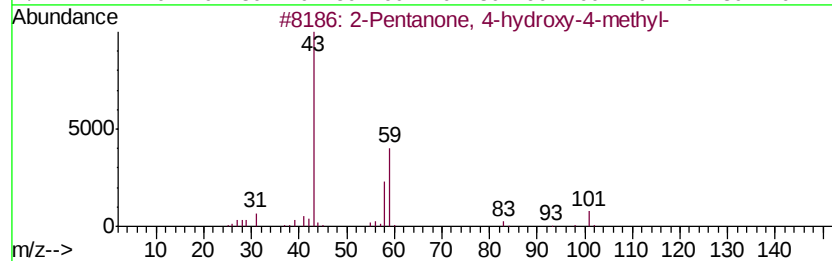
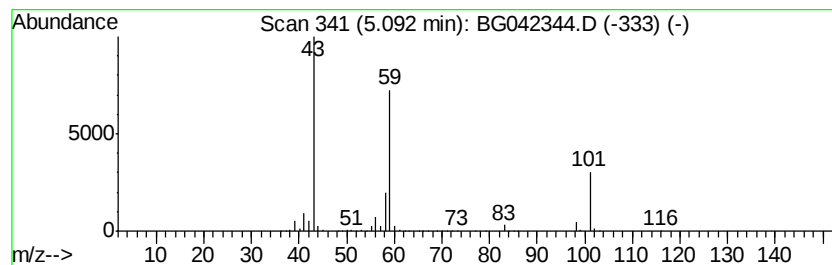
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TIC Library : C:\Database\NIST11.L
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Peak Number 2 unknown5.09 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.09	177.53 ng	2497880	1,4-Dichlorobenzene-d4	8.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2		Guanidine	59	CH5N3	000113-00-8	38
3		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	12
4		Acetone	58	C3H6O	000067-64-1	12
5		(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9



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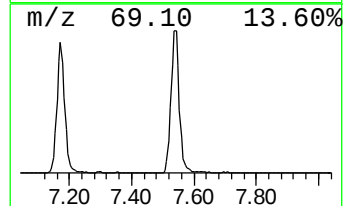
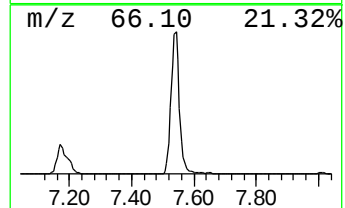
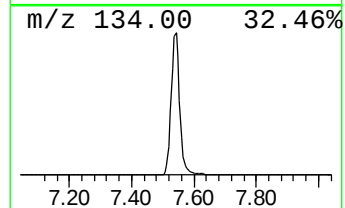
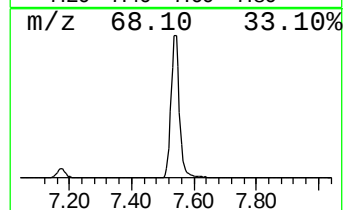
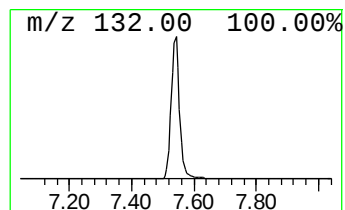
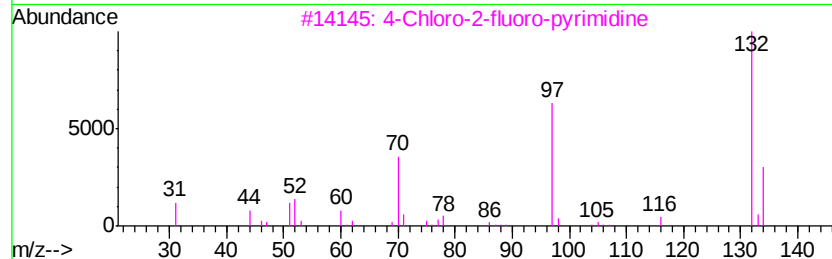
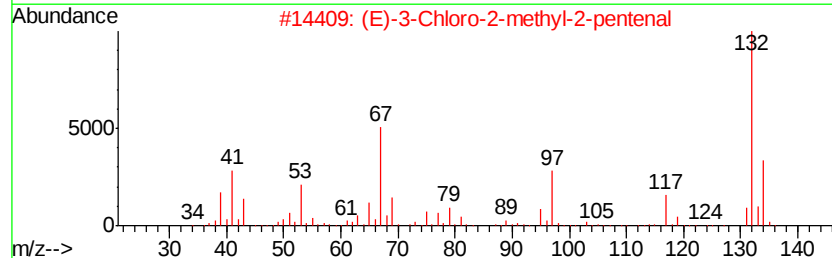
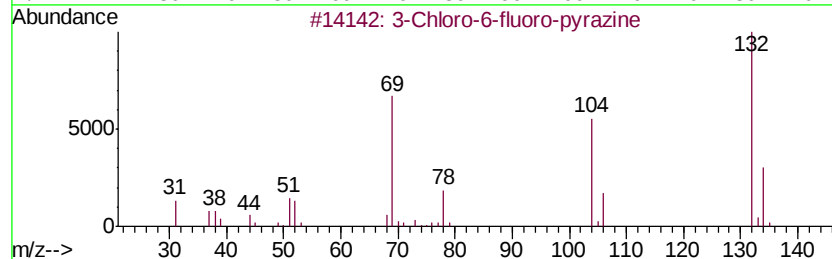
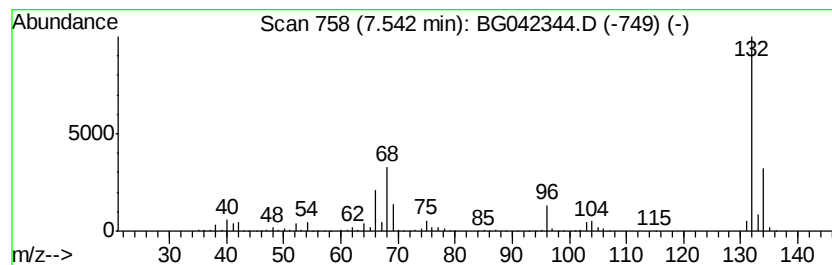
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TIC Library : C:\Database\NIST11.L
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Peak Number 3 unknown7.54 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.54	94.70 ng	1332430	1,4-Dichlorobenzene-d4	8.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
3		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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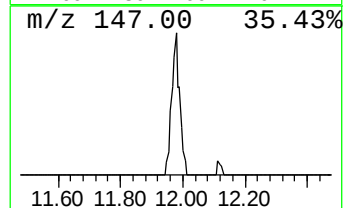
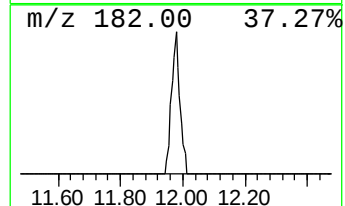
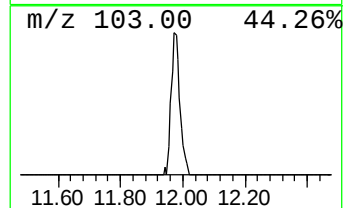
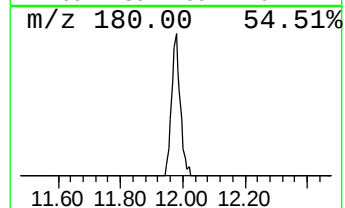
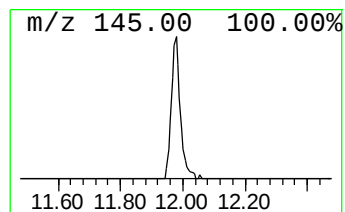
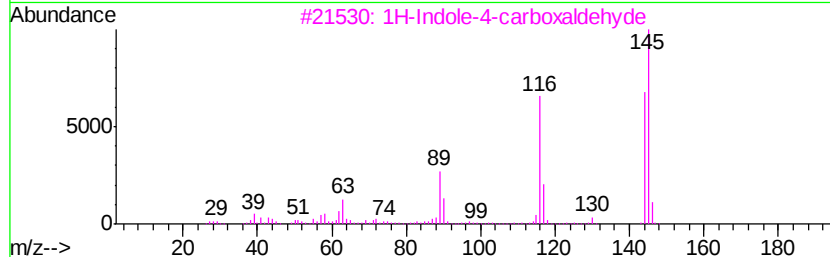
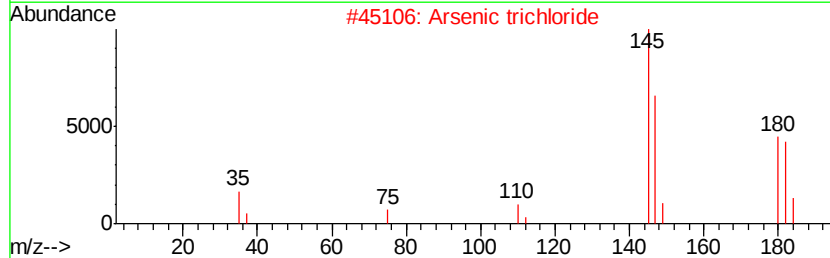
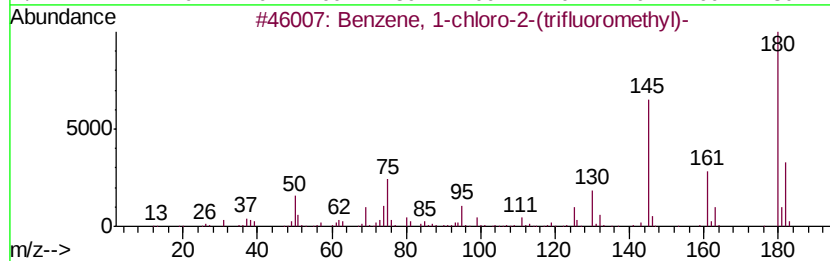
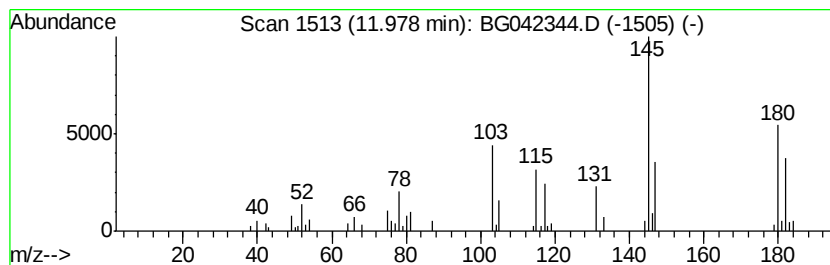
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Peak Number 5 unknown11.98 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.98	3.52 ng	70053	Naphthalene-d8	10.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-chloro-2-(trifluorome...	180	C7H4ClF3	000088-16-4	32
2		Arsenic trichloride	180	AsCl3	007784-34-1	32
3		1H-Indole-4-carboxaldehyd	145	C9H7NO	001074-86-8	27
4		5-Quinolinol	145	C9H7NO	000578-67-6	27
5		2(1H)-Pyridinone, 5-chloro-3-hyd...	145	C5H4ClN02	053233-89-9	27



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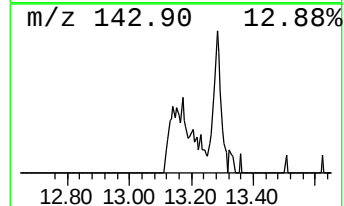
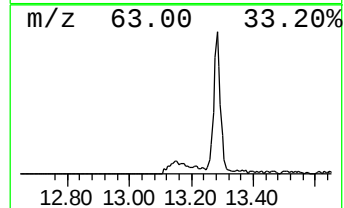
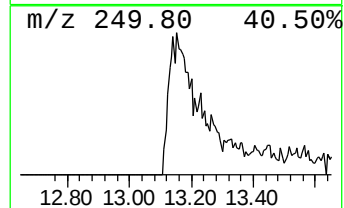
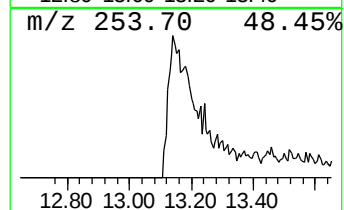
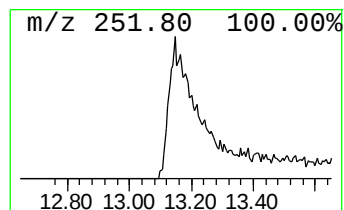
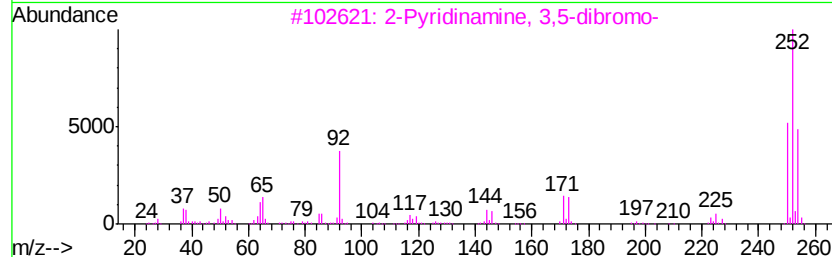
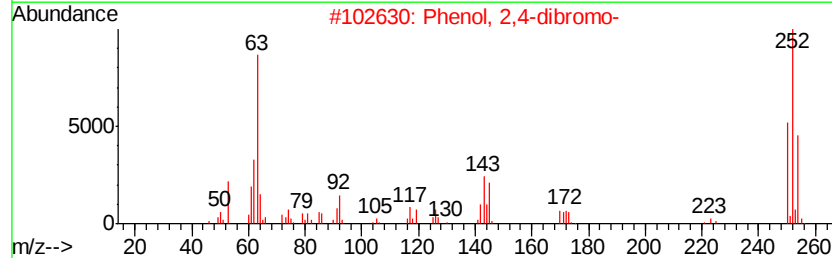
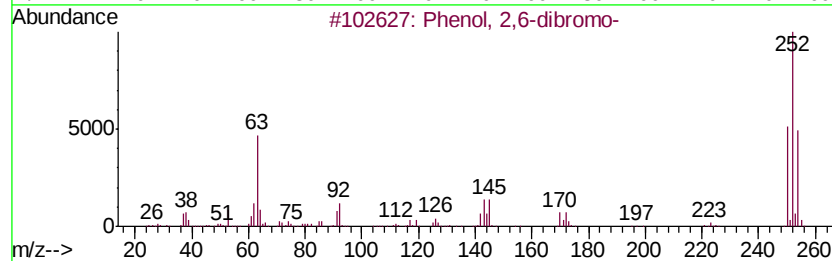
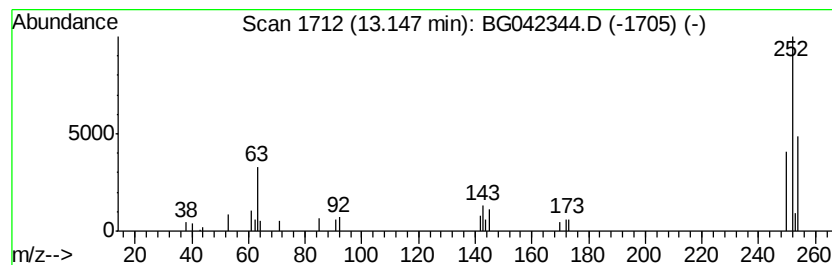
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Peak Number 6 Phenol, 2,6-dibromo- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.15	2.13 ng	68397	Acenaphthene-d10	14.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenol, 2,6-dibromo-	250	C6H4Br2O	000608-33-3	97
2		Phenol, 2,4-dibromo-	250	C6H4Br2O	000615-58-7	94
3		2-Pyridinamine, 3,5-dibromo-	250	C5H4Br2N2	035486-42-1	64
4		4-(2-Hydroxy-5-bromophenyl)pvrin...	250	C10H7BrN2O	097630-76-7	12
5		Thieno[3,2-d]isothiazole, 3-cyan...	252	C8H4N4S3	077729-68-1	9



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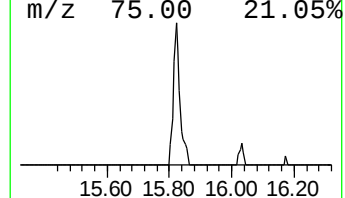
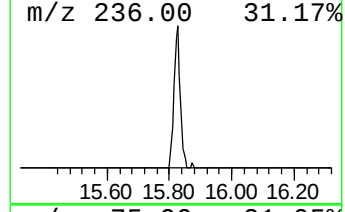
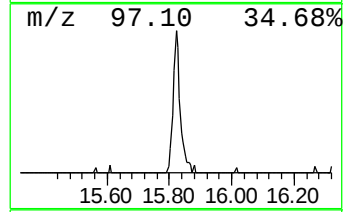
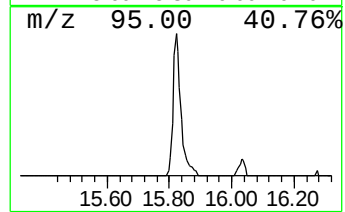
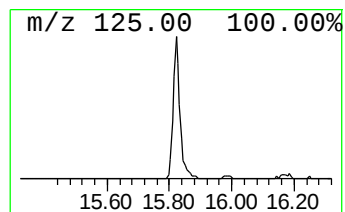
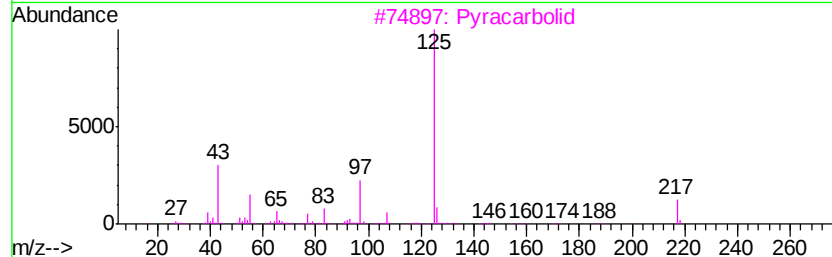
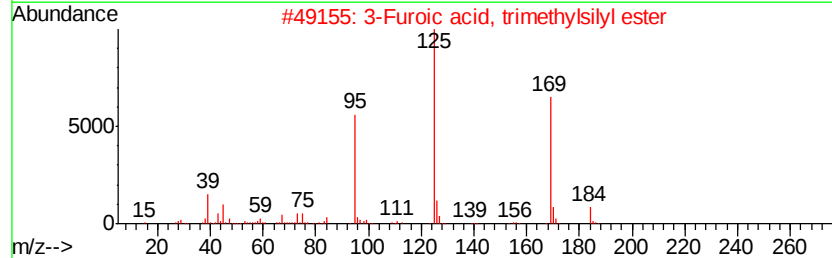
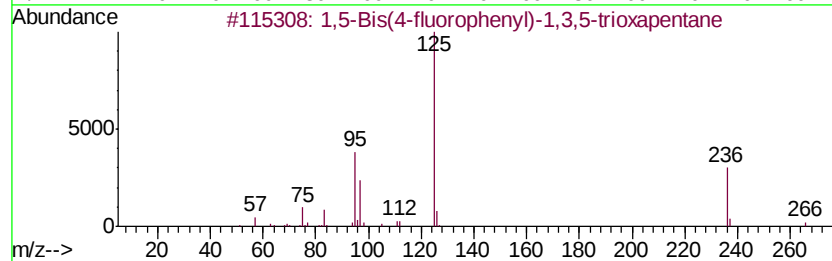
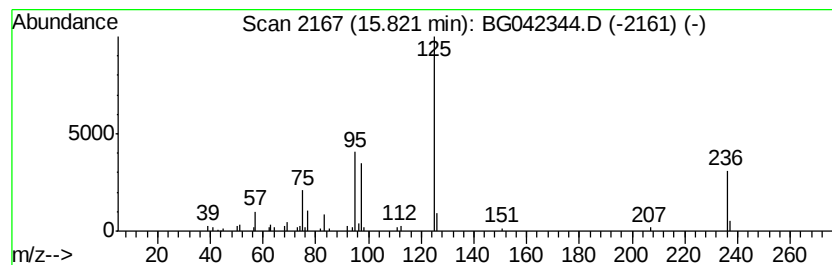
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 1,5-Bis(4-fluorophenyl)-1,3... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.82	2.22 ng	71386	Acenaphthene-d10	14.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,5-Bis(4-fluorophenyl)-1,3,5-tr...	266	C14H12F2O3	1000280-28-3	64
2		3-Furoic acid, trimethylsilyl ester	184	C8H12O3Si	1000332-78-1	47
3		Pyracarbolid	217	C13H15NO2	024691-76-7	43
4		2-Amino-4-hydroxy-6-methylpyrimi...	125	C5H7N3O	003977-29-5	43
5		Thiazole, 5-ethenyl-4-methyl-	125	C6H7NS	001759-28-0	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081919\
Data File : BG042344.D
Acq On : 20 Aug 2019 2:17
Operator : HP/JU
Sample : K4290-06
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
WP-1

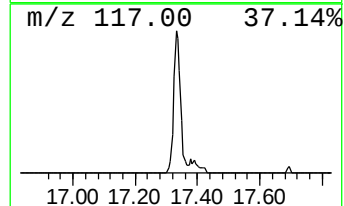
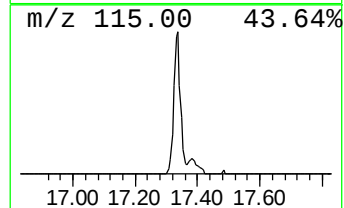
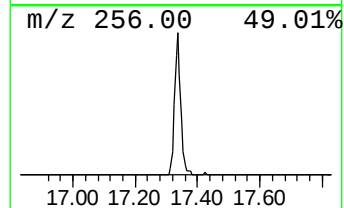
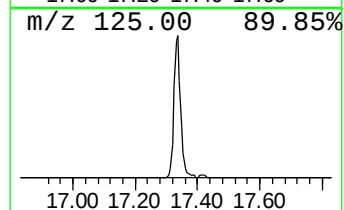
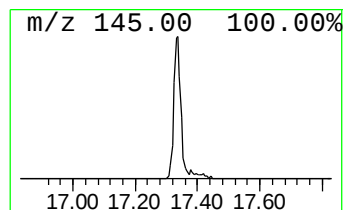
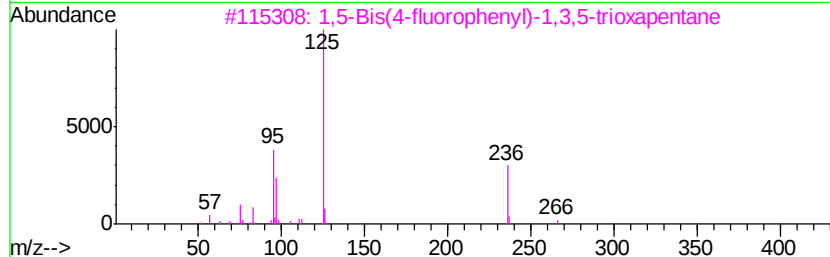
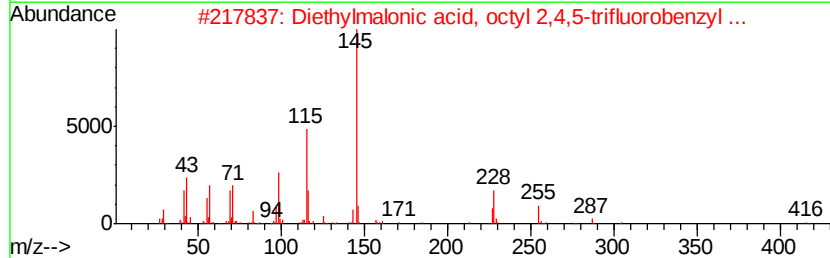
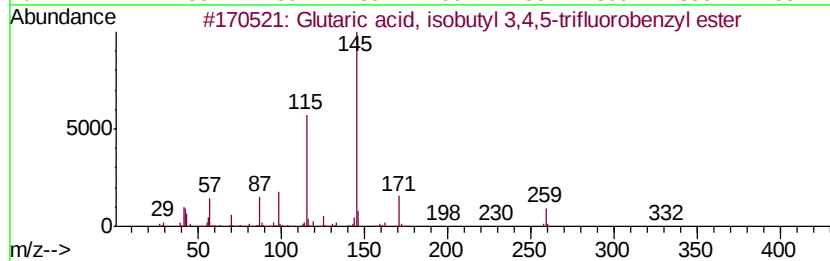
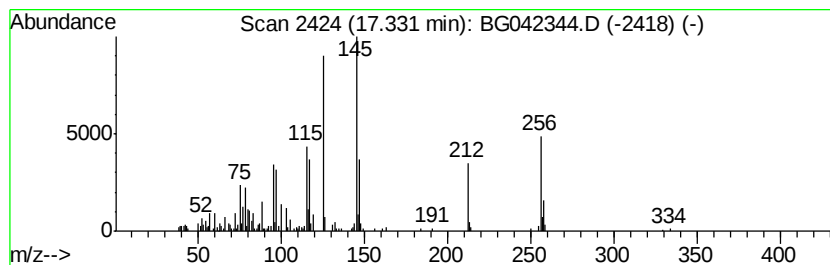
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown17.33 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.33	4.18 ng	175126	Phenanthrene-d10	17.41

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Glutaric acid, isobutyl 3,4,5-tr...	332	C16H19F3O4	1000377-56-7	22
2		Diethylmalonic acid, octyl 2,4,5...	416	C22H31F3O4	1000369-26-0	22
3		1,5-Bis(4-fluorophenyl)-1,3,5-tr...	266	C14H12F2O3	1000280-28-3	22
4		2(1H)-Pyridinone, 5-chloro-3-hyd...	145	C5H4ClN02	053233-89-9	18
5		2,4,5-Trifluorobenzyl alcohol, t...	258	C9H4F6O2	1000364-21-7	18



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081919\
Data File : BG042344.D
Acq On : 20 Aug 2019 2:17
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Sample : K4290-06
Misc :
ALS Vial : 17 Sample Multiplier: 1

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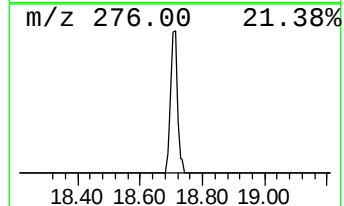
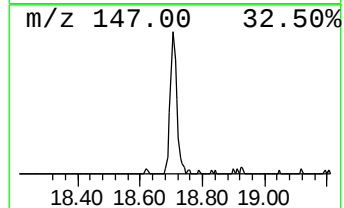
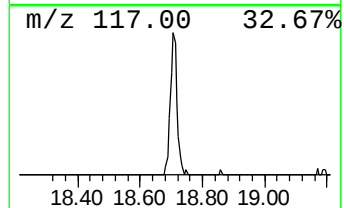
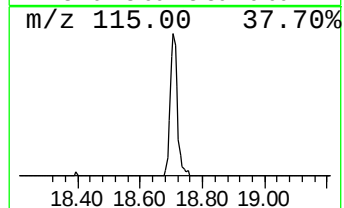
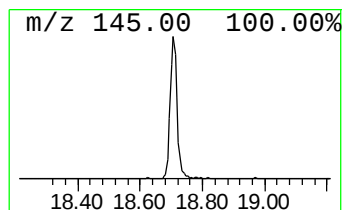
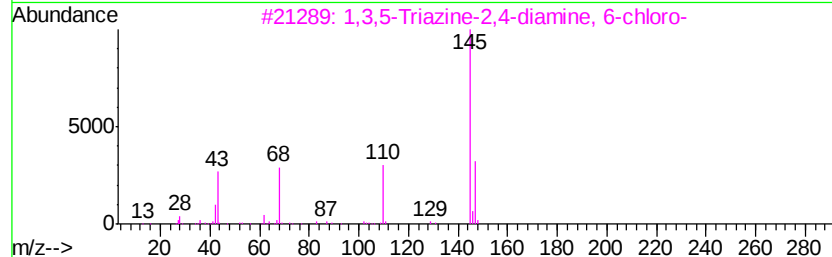
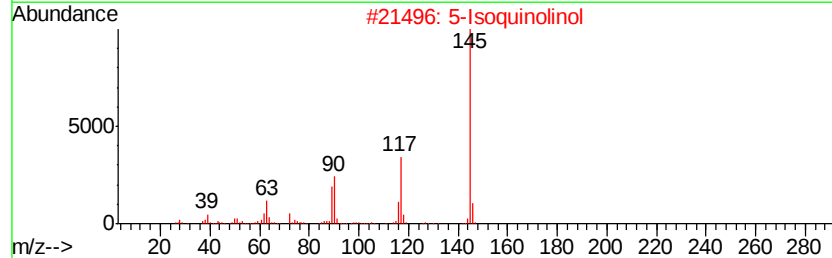
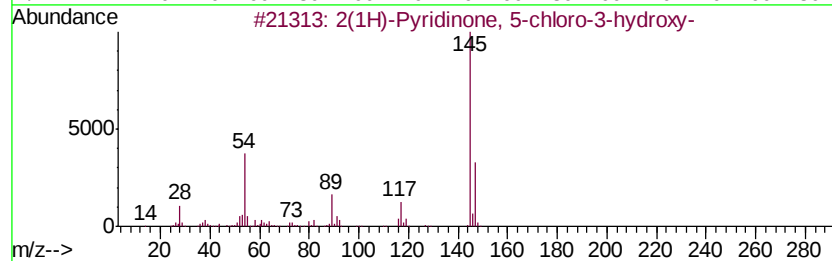
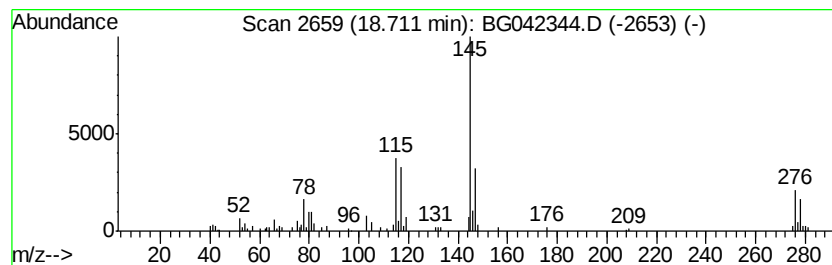
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 unknown18.71 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.71	2.21 ng	92529	Phenanthrene-d10	17.41

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2(1H)-Pyridinone, 5-chloro-3-hyd...	145	C5H4ClN02		053233-89-9	47
2	5-Isoquinolinol	145	C9H7NO		002439-04-5	43
3	1,3,5-Triazine-2,4-diamine, 6-ch...	145	C3H4ClN5		003397-62-4	43
4	Ethanone, 1-(2,3-dihydro-1H-inde...	160	C11H12O		004228-10-8	40
5	2-Quinazolinamine	145	C8H7N3		001687-51-0	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081919\
Data File : BG042344.D
Acq On : 20 Aug 2019 2:17
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Sample : K4290-06
Misc :
ALS Vial : 17 Sample Multiplier: 1

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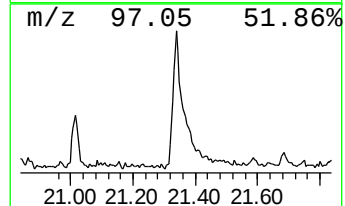
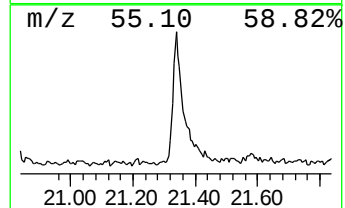
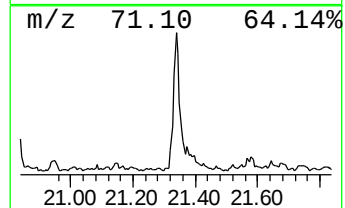
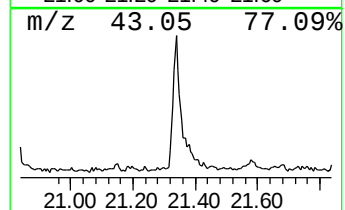
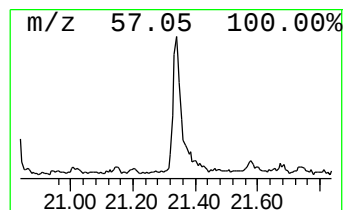
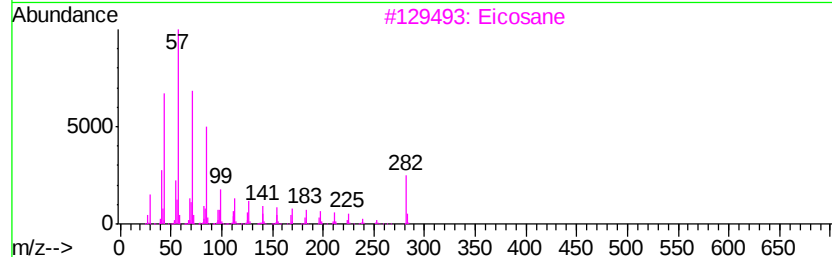
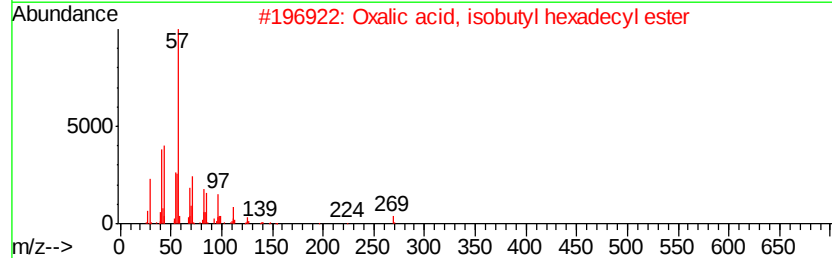
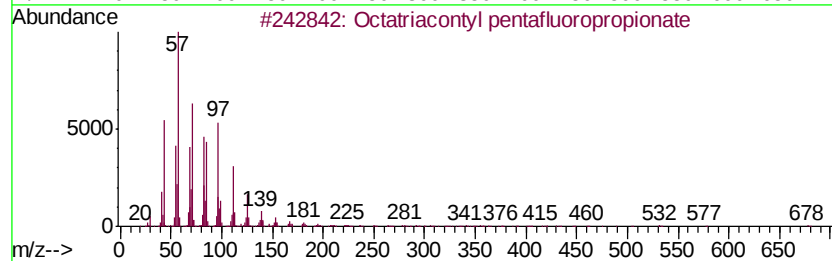
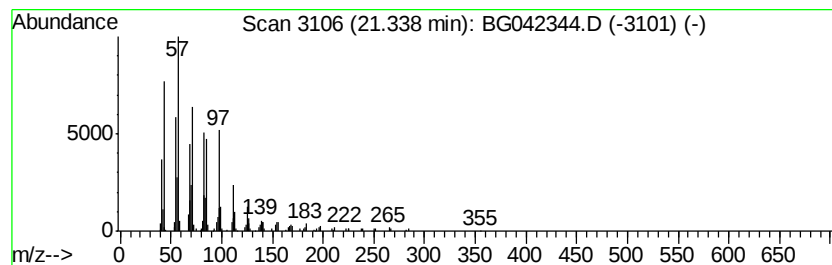
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Octatriacontyl pentafluorop... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.34	4.72 ng	228798	Chrysene-d12	21.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octatriacontyl pentafluoropropio...	697	C41H77F5O2	1000351-89-1	91
2		Oxalic acid, isobutyl hexadecyl ...	370	C22H42O4	1000309-38-1	91
3		Eicosane	282	C20H42	000112-95-8	83
4		Heptadecyl trifluoroacetate	352	C19H35F3O2	1000351-87-0	70
5		Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	70



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081919\
Data File : BG042344.D
Acq On : 20 Aug 2019 2:17
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Sample : K4290-06
Misc :
ALS Vial : 17 Sample Multiplier: 1

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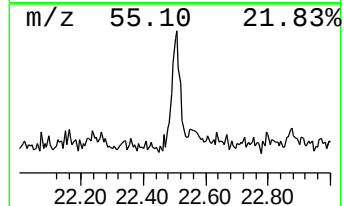
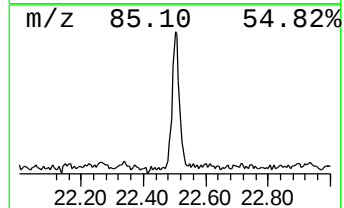
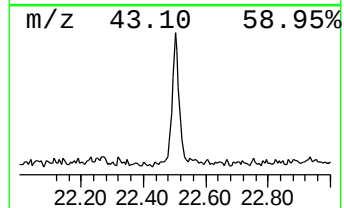
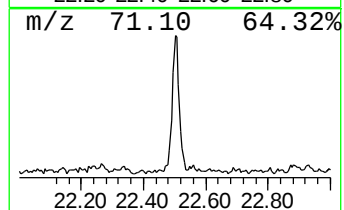
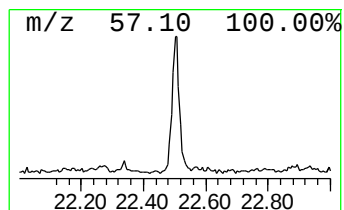
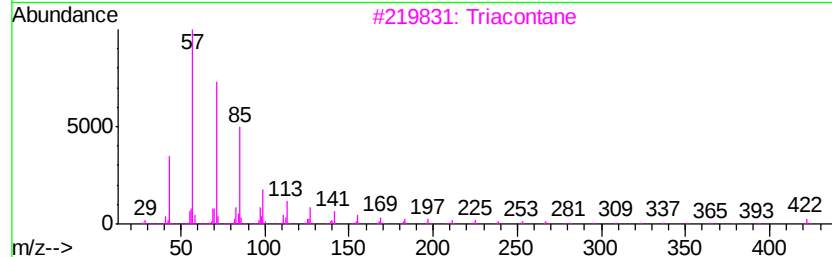
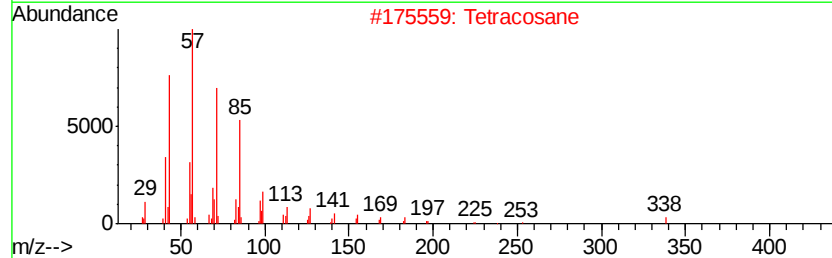
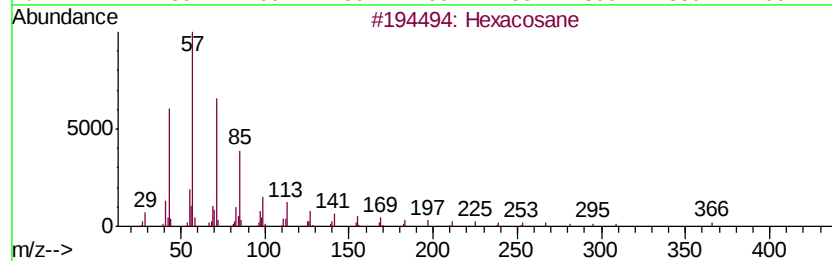
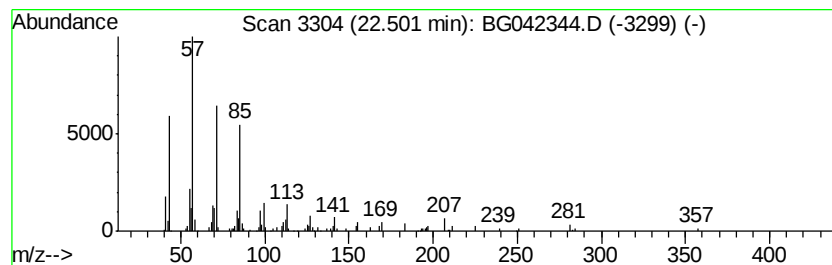
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Hexacosane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.50	2.10 ng	102009	Chrysene-d12	21.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	90
2		Tetracosane	338	C24H50	000646-31-1	87
3		Triacontane	422	C30H62	000638-68-6	87
4		Heptadecane	240	C17H36	000629-78-7	87
5		Docosane	310	C22H46	000629-97-0	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081919\
Data File : BG042344.D
Acq On : 20 Aug 2019 2:17
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Sample : K4290-06
Misc :
ALS Vial : 17 Sample Multiplier: 1

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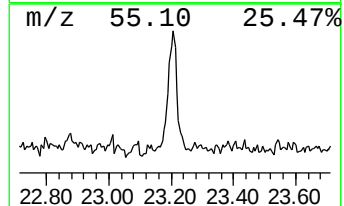
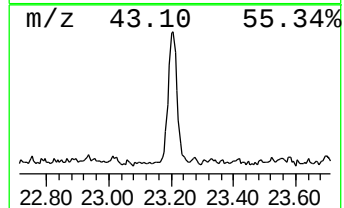
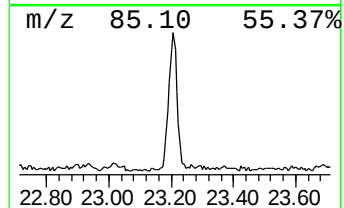
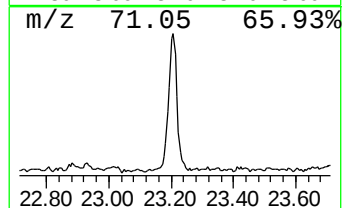
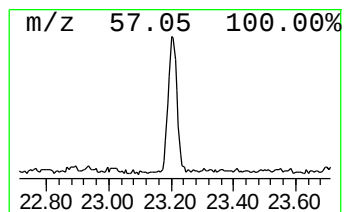
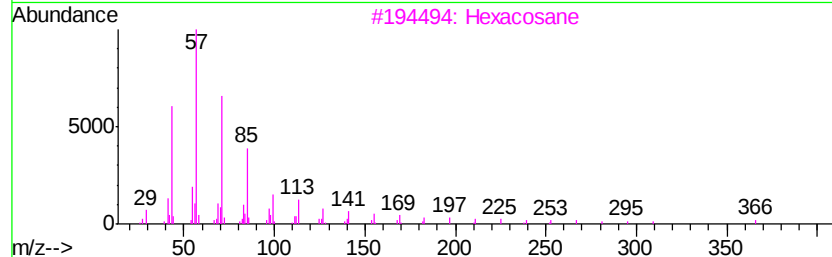
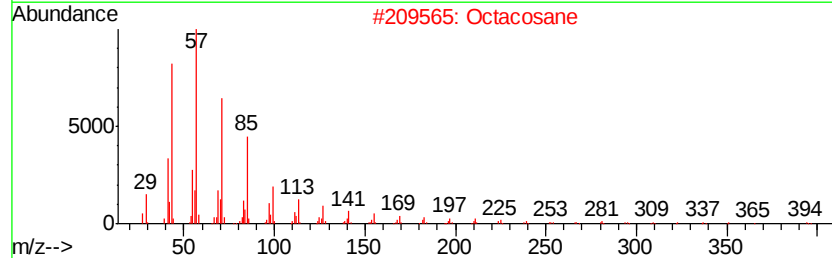
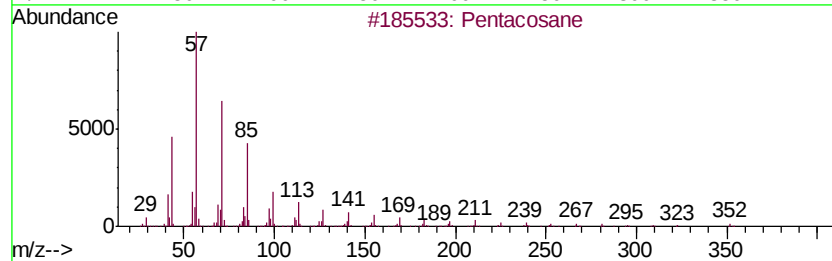
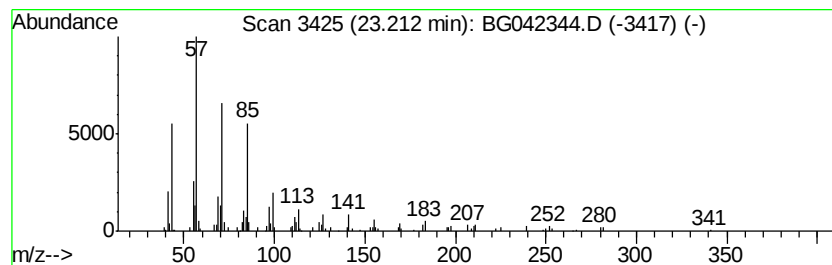
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Pentacosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.21	2.68 ng	129916	Chrysene-d12	21.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentacosane	352	C25H52	000629-99-2	91
2		Octacosane	394	C28H58	000630-02-4	91
3		Hexacosane	366	C26H54	000630-01-3	91
4		Docosane	310	C22H46	000629-97-0	91
5		Eicosane, 2-methyl-	296	C21H44	001560-84-5	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081919\
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Acq On : 20 Aug 2019 2:17
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Sample : K4290-06
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ALS Vial : 17 Sample Multiplier: 1

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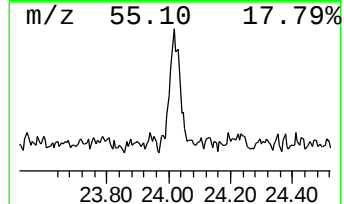
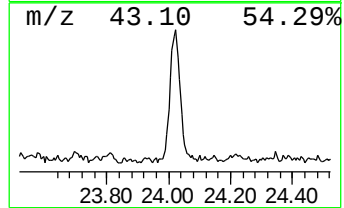
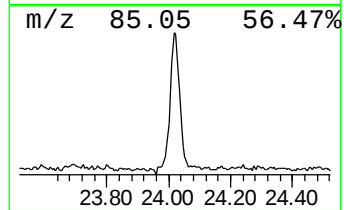
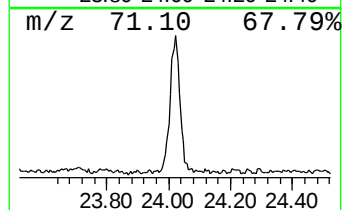
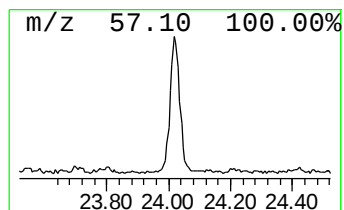
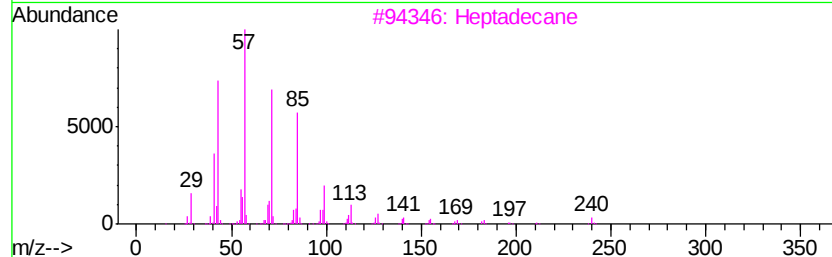
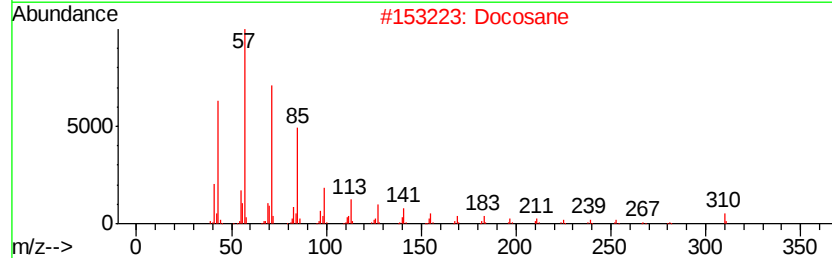
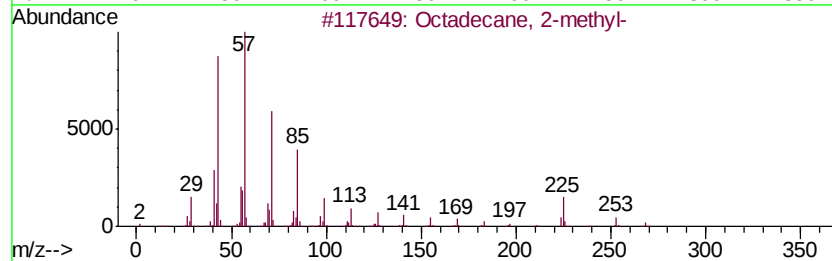
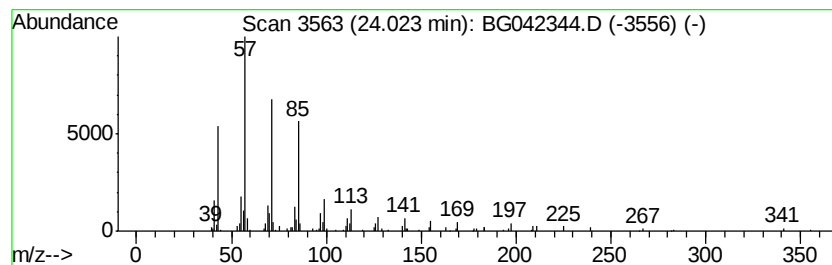
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Octadecane, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.02	2.28 ng	133150	Perylene-d12	24.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane, 2-methyl-	268	C19H40	001560-88-9	94
2		Docosane	310	C22H46	000629-97-0	90
3		Heptadecane	240	C17H36	000629-78-7	87
4		Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	87
5		Tetracosane	338	C24H50	000646-31-1	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG081919\
 Data File : BG042344.D
 Acq On : 20 Aug 2019 2:17
 Operator : HP/JU
 Sample : K4290-06
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 WP-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG081919.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
Butane, 2-methoxy...	3.12	137.8	ng	1938670	1	8.01	281408	20.0
unknown5.09	5.09	177.5	ng	2497880	1	8.01	281408	20.0
unknown7.54	7.54	94.7	ng	1332430	1	8.01	281408	20.0
unknown11.98	11.98	3.5	ng	70053	2	10.83	398389	20.0
Phenol, 2,6-dibromo-	13.15	2.1	ng	68397	3	14.66	642116	20.0
1,5-Bis(4-fluorop...	15.82	2.2	ng	71386	3	14.66	642116	20.0
unknown17.33	17.33	4.2	ng	175126	4	17.41	838753	20.0
unknown18.71	18.71	2.2	ng	92529	4	17.41	838753	20.0
Octatriacontyl pe...	21.34	4.7	ng	228798	5	21.69	970469	20.0
Hexacosane	22.50	2.1	ng	102009	5	21.69	970469	20.0
Pentacosane	23.21	2.7	ng	129916	5	21.69	970469	20.0
Octadecane, 2-met...	24.02	2.3	ng	133150	6	24.93	1168170	20.0