

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG072018\
 Data File : BG035797.D
 Acq On : 20 Jul 2018 23:19
 Operator : JU/SJ
 Sample : J4083-01
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MH-6

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-BG071218.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	5.145	309	316	323	rBV	44382	69804	1.60%	0.302%
2	5.615	389	396	406	rBV	416590	712572	16.35%	3.082%
3	7.195	659	665	678	rBV	248134	495495	11.37%	2.143%
4	7.565	720	728	741	rBV	828640	1437329	32.98%	6.217%
5	8.029	799	807	817	rBV	189359	330799	7.59%	1.431%
6	8.352	852	862	870	rBV	709864	1233625	28.31%	5.336%
7	9.192	996	1005	1020	rBV	608838	1138706	26.13%	4.926%
8	10.831	1277	1284	1298	rBV	226032	409908	9.41%	1.773%
9	13.275	1692	1700	1709	rBV	1655424	2608128	59.85%	11.282%
10	14.097	1835	1840	1847	rBV	63456	102880	2.36%	0.445%
11	14.649	1928	1934	1946	rBV3	369487	618407	14.19%	2.675%
12	16.141	2181	2188	2201	rBV	1452166	2292989	52.62%	9.919%
13	16.658	2271	2276	2285	rBV2	22121	44246	1.02%	0.191%
14	17.393	2391	2401	2410	rBV	530662	805738	18.49%	3.485%
15	18.280	2547	2552	2559	rBV2	76769	121137	2.78%	0.524%
16	19.431	2740	2748	2753	rBV2	92310	155182	3.56%	0.671%
17	19.766	2802	2805	2809	rBV	49682	58911	1.35%	0.255%
18	19.807	2809	2812	2816	rVB2	54734	61827	1.42%	0.267%
19	20.001	2838	2845	2857	rBV	3123791	4357656	100.00%	18.850%
20	20.301	2891	2896	2900	rBV	95593	111368	2.56%	0.482%
21	20.659	2954	2957	2968	rVB10	47745	85721	1.97%	0.371%
22	20.794	2973	2980	2983	rVV2	167998	237702	5.45%	1.028%
23	21.299	3061	3066	3075	rVB	216881	289944	6.65%	1.254%
24	21.564	3097	3111	3115	rBV	67736	247677	5.68%	1.071%
25	21.657	3121	3127	3139	rVB	589985	1073829	24.64%	4.645%
26	21.840	3153	3158	3165	rVB	227773	312738	7.18%	1.353%
27	22.445	3255	3261	3266	rBV	222717	348342	7.99%	1.507%
28	23.132	3371	3378	3386	rVB	230224	418567	9.61%	1.811%
29	23.314	3406	3409	3418	rVB4	29531	56217	1.29%	0.243%
30	23.925	3506	3513	3520	rBV2	209936	433519	9.95%	1.875%
31	24.859	3661	3672	3681	rBV3	478101	1311722	30.10%	5.674%
32	25.981	3855	3863	3873	rBV5	128310	350613	8.05%	1.517%
33	27.250	4071	4079	4082	rBV6	27854	75471	1.73%	0.326%
34	27.320	4084	4091	4102	rVB5	71572	249049	5.72%	1.077%

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35	28.807	4324	4344	4345	rBV6	106114	402619	9.24%	1.742%
36	30.105	4558	4565	4567	rBV8	25626	57330	1.32%	0.248%

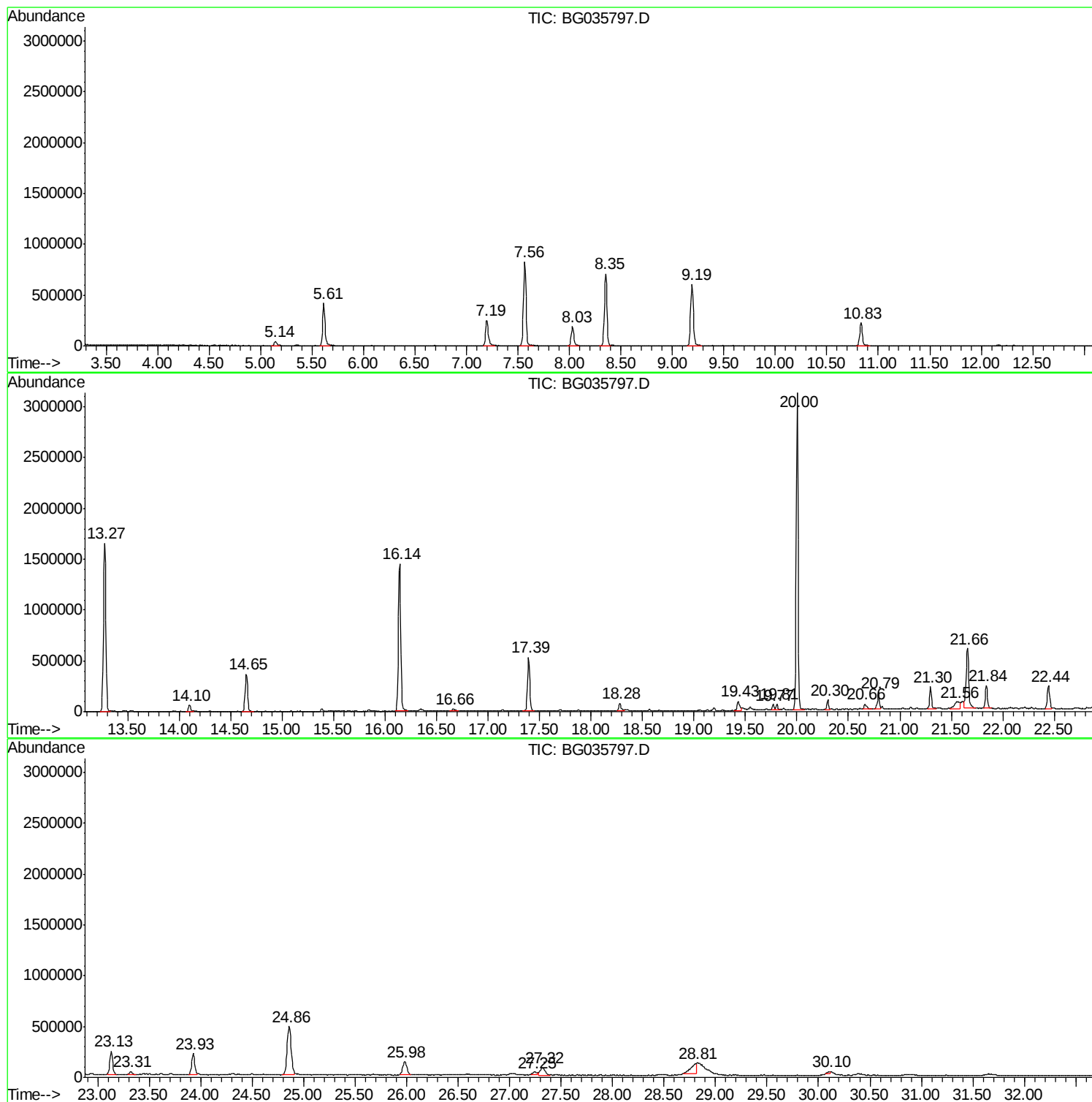
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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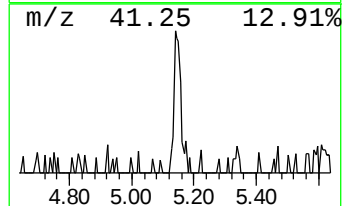
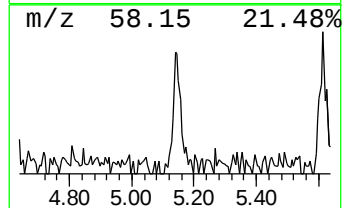
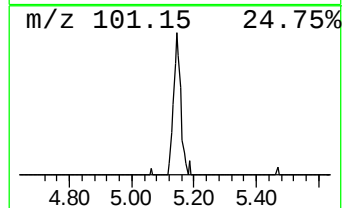
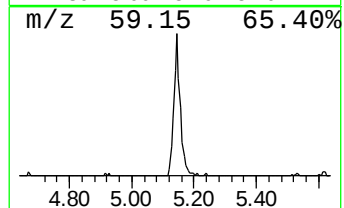
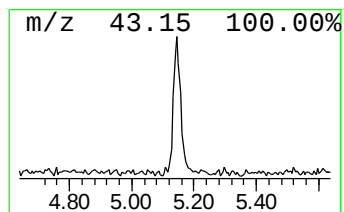
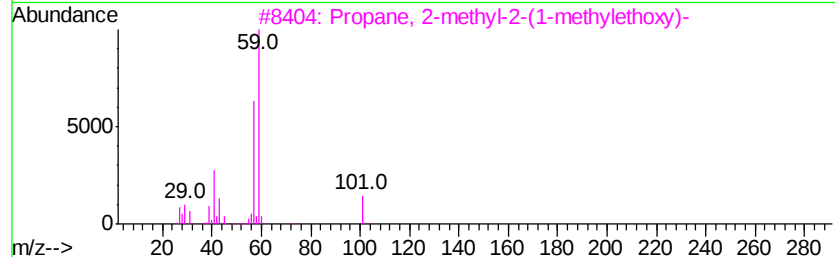
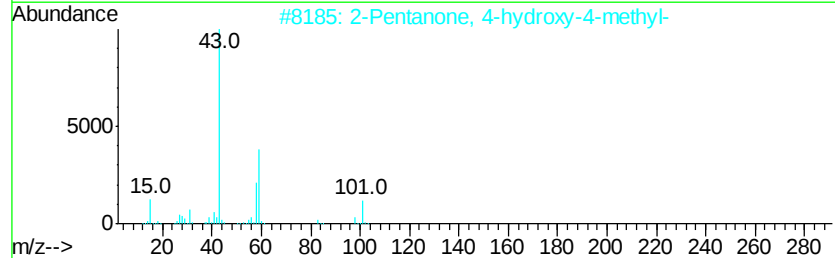
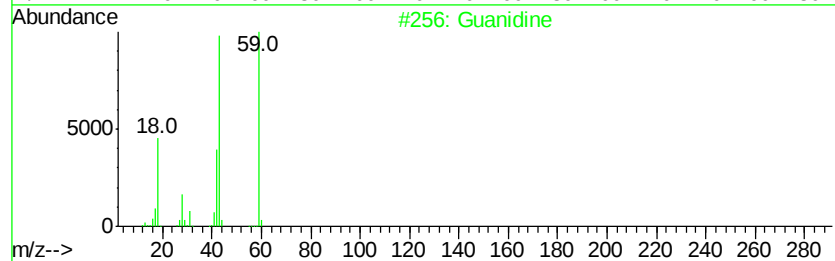
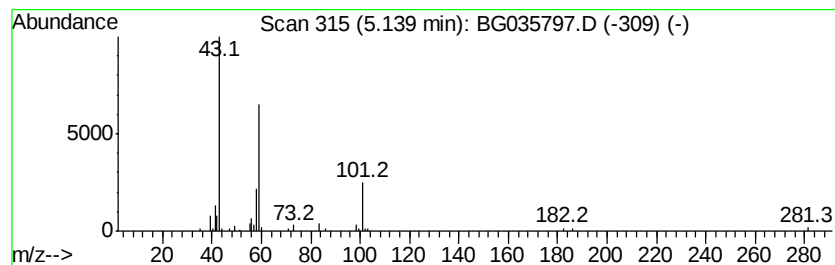
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Peak Number 1 Guanidine Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.14	4.22 ng	69804	1,4-Dichlorobenzene-d4	8.03

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Guanidine	59	CH5N3	000113-00-8	53
2		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
3		Propane, 2-methyl-2-(1-methylethyl)-	116	C7H16O	017348-59-3	45
4		Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	43
5		Dimethadione	129	C5H7NO3	000695-53-4	38



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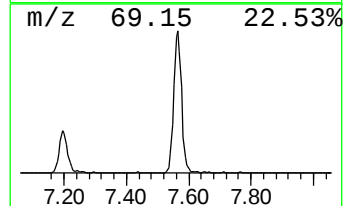
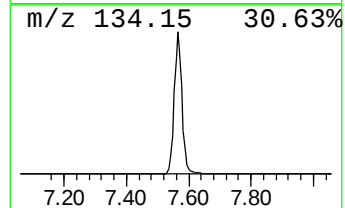
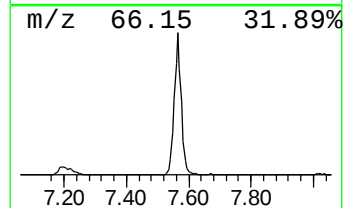
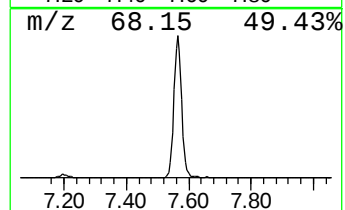
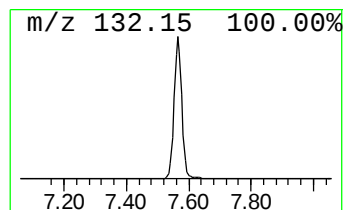
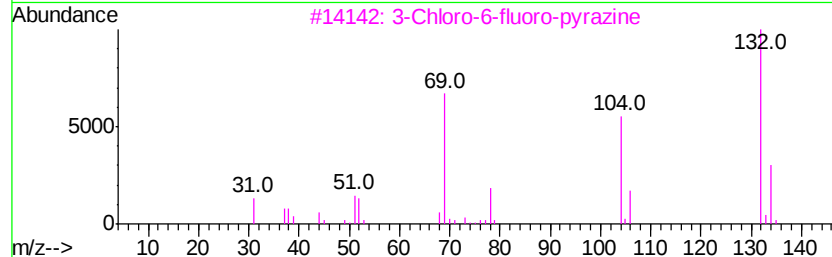
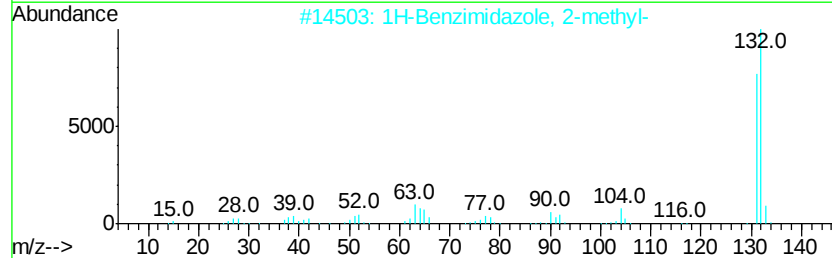
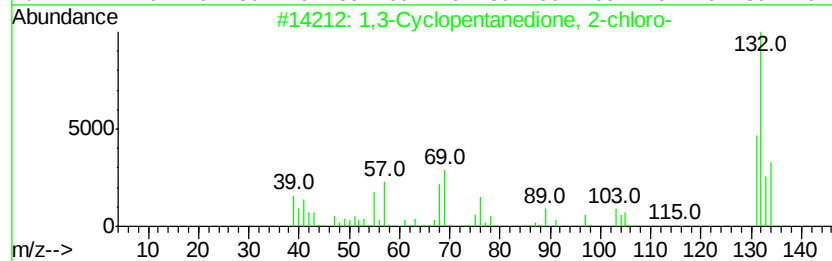
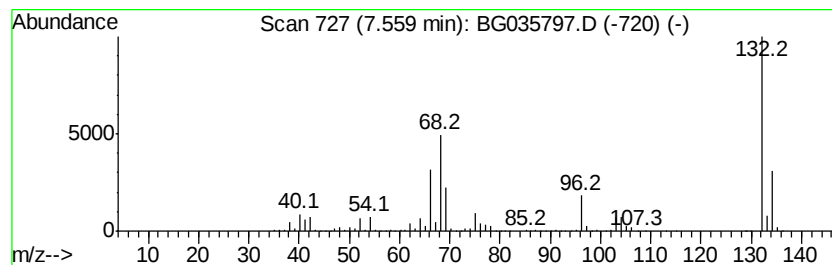
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Peak Number 2 unknown7.56 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.56	86.90 ng	1437330	1,4-Dichlorobenzene-d4	8.03

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	25
2		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
3		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
5		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10



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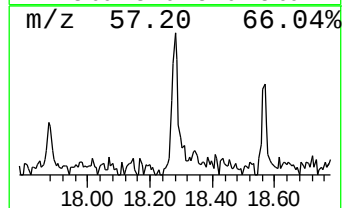
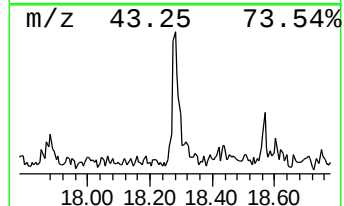
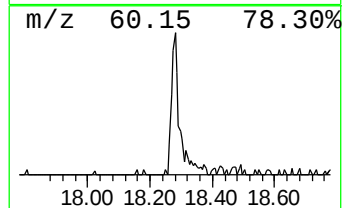
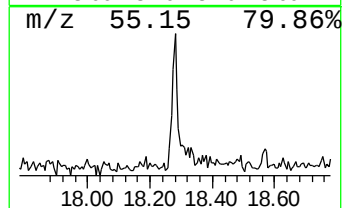
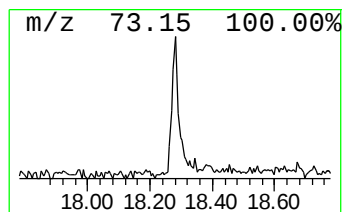
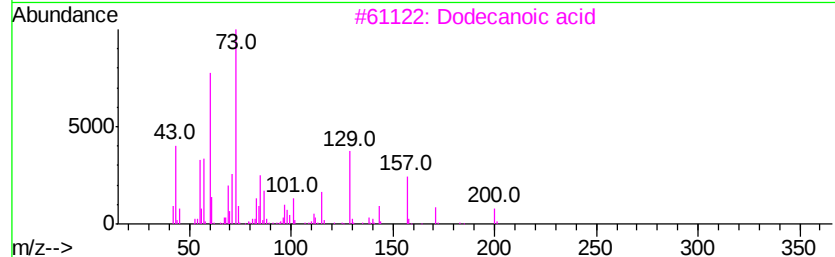
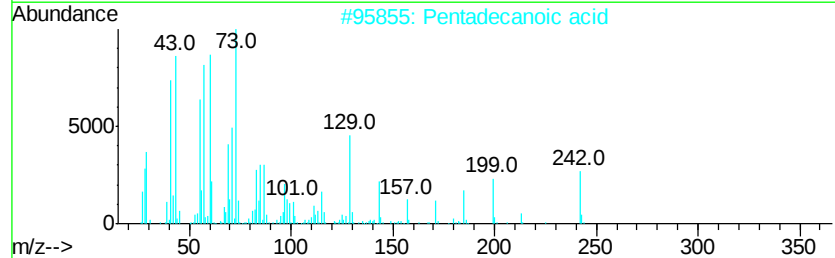
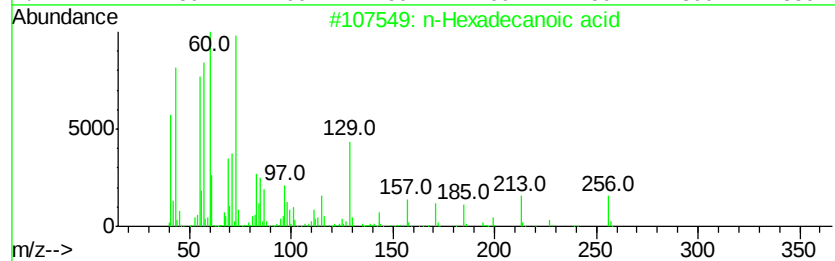
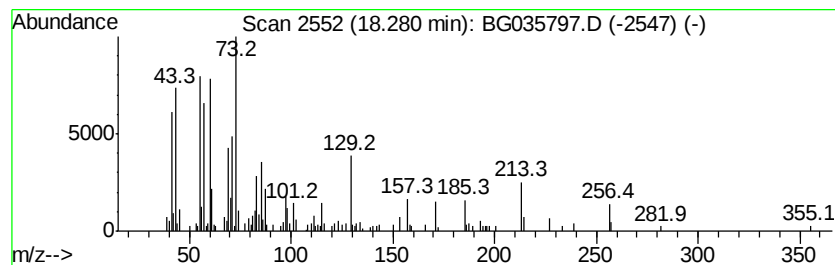
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Peak Number 4 n-Hexadecanoic acid Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.28	3.01 ng	121137	Phenanthrene-d10	17.39

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	87
3		Dodecanoic acid	200	C12H24O2	000143-07-7	76
4		Tridecanoic acid	214	C13H26O2	000638-53-9	74
5		Octadecanoic acid	284	C18H36O2	000057-11-4	74



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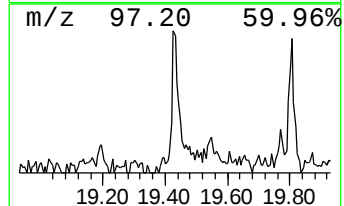
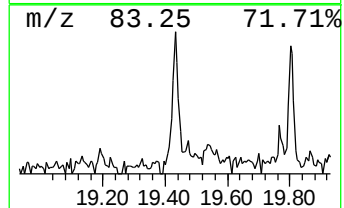
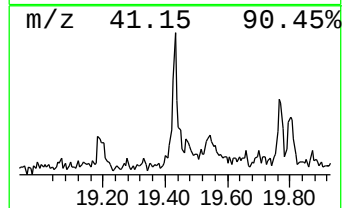
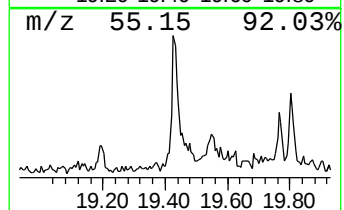
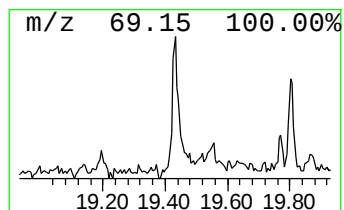
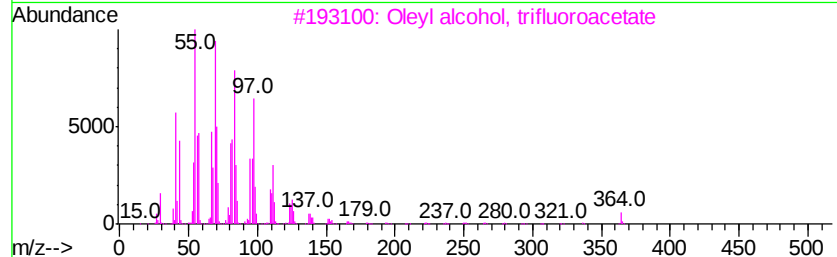
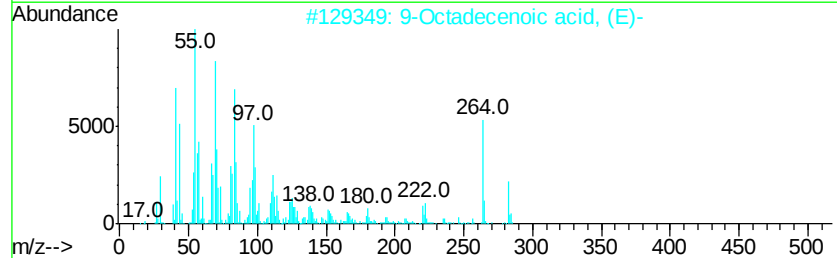
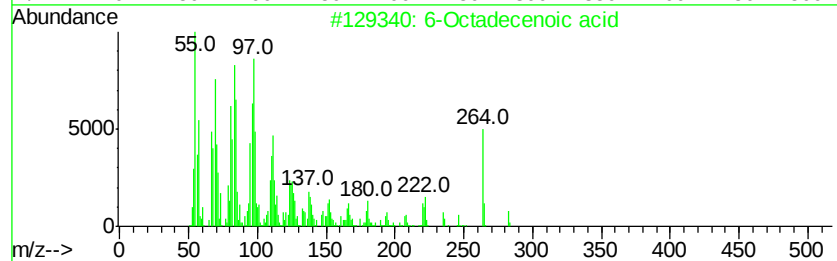
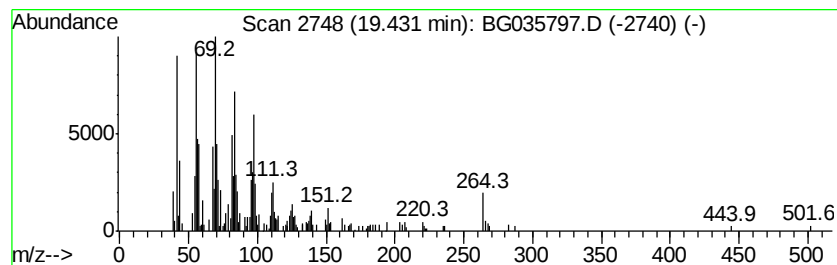
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Peak Number 5 6-Octadecenoic acid Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.43	3.85 ng	155182	Phenanthrene-d10	17.39

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	6-Octadecenoic acid	282	C18H34O2	1000336-66-8	83
2	9-Octadecenoic acid, (E)-	282	C18H34O2	000112-79-8	83
3	Olel alcohol, trifluoroacetate	364	C20H35F3O2	1000352-68-4	74
4	Oleic Acid	282	C18H34O2	000112-80-1	70
5	cis-13-Octadecenoic acid	282	C18H34O2	013126-39-1	64



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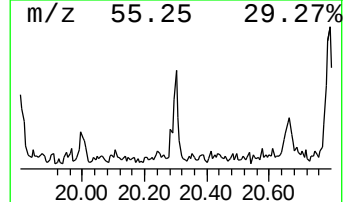
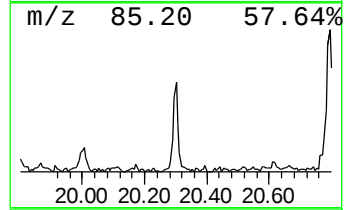
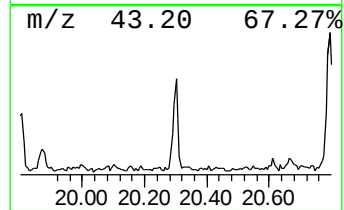
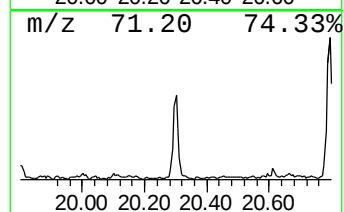
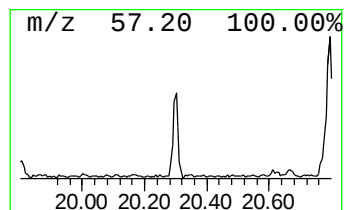
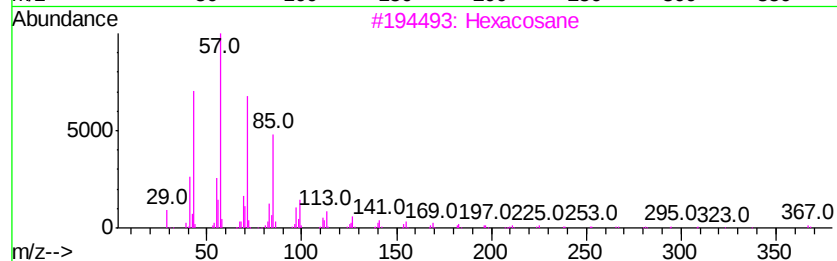
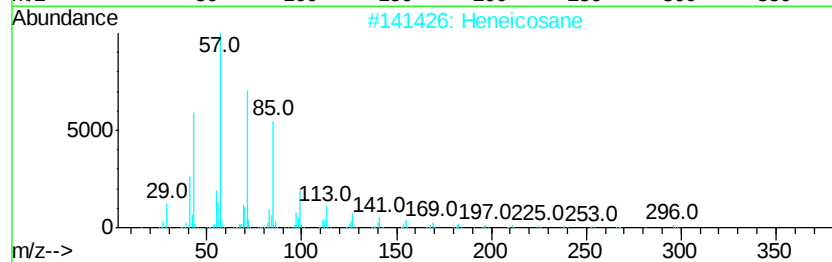
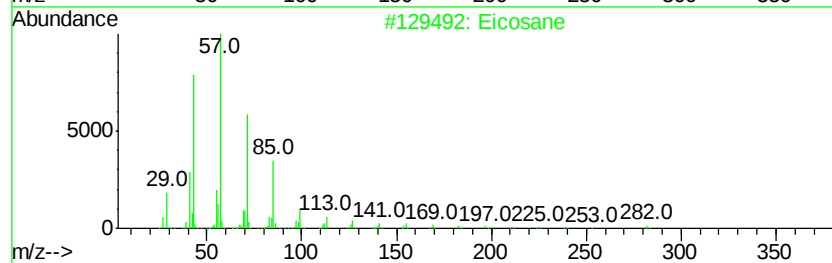
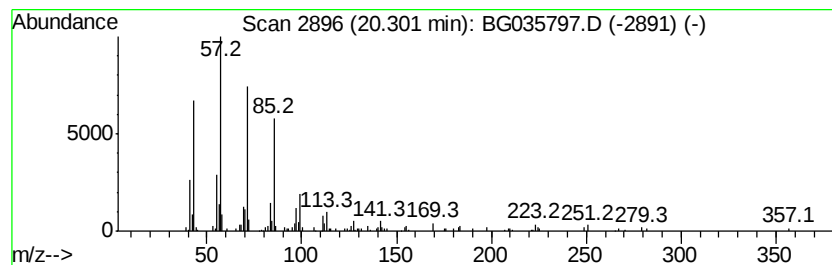
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Peak Number 6 Eicosane Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.30	2.07 ng	111368	Chrysene-d12	21.66

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	95
2	Heneicosane		296	C21H44	000629-94-7	91
3	Hexacosane		366	C26H54	000630-01-3	91
4	Triacontane		422	C30H62	000638-68-6	90
5	Dodecane, 1-iodo-		296	C12H25I	004292-19-7	86



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Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
MH-6

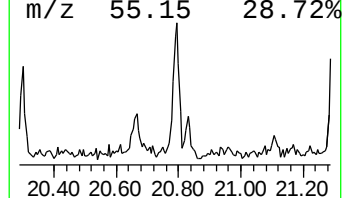
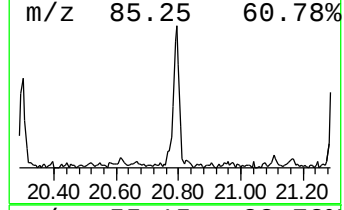
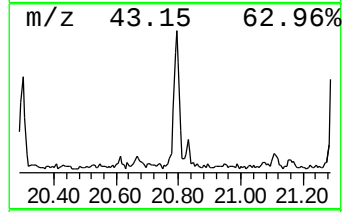
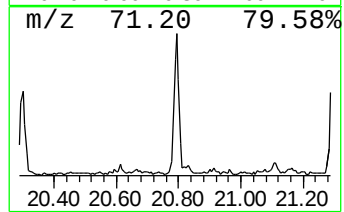
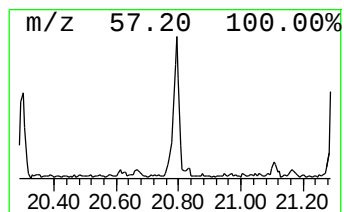
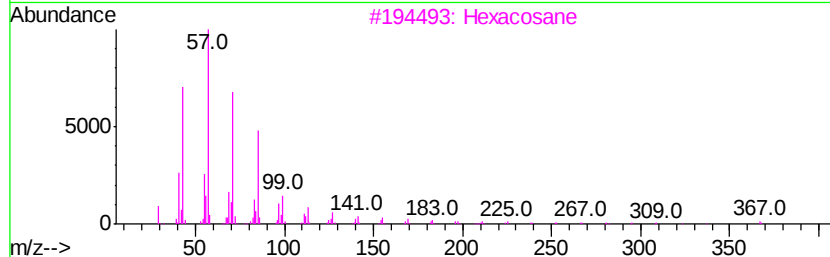
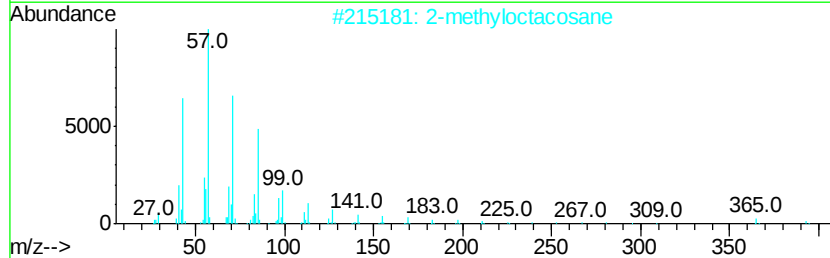
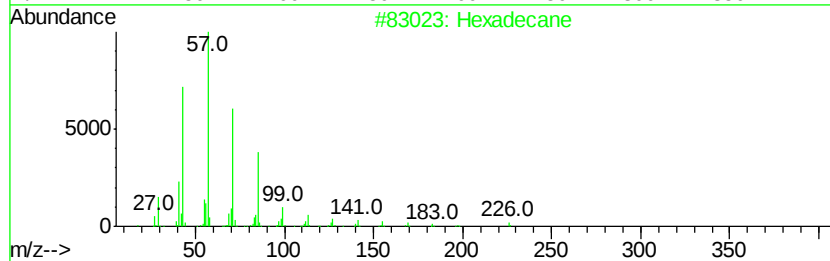
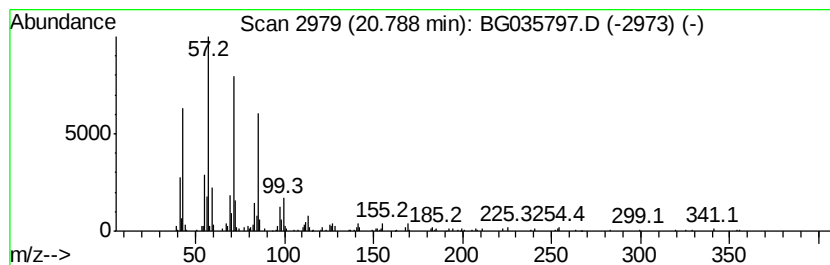
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Hexadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.79	4.43 ng	237702	Chrysene-d12	21.66

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	93
2			2-methyloctacosane	408	C29H60	1000376-72-8	91
3			Hexacosane	366	C26H54	000630-01-3	91
4			Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91
5			Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG072018\
Data File : BG035797.D
Acq On : 20 Jul 2018 23:19
Operator : JU/SJ
Sample : J4083-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

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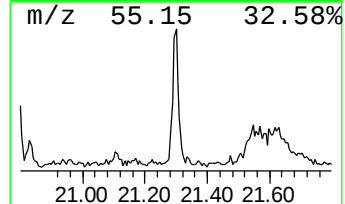
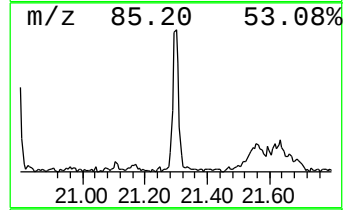
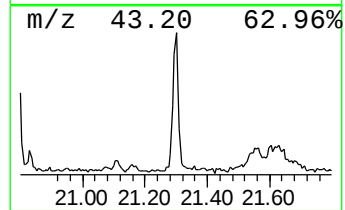
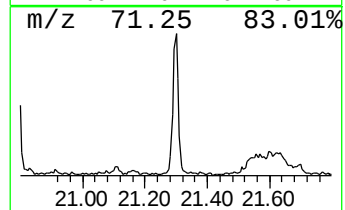
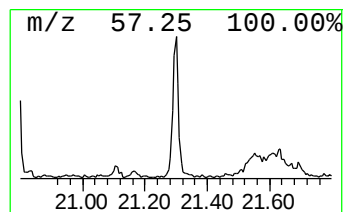
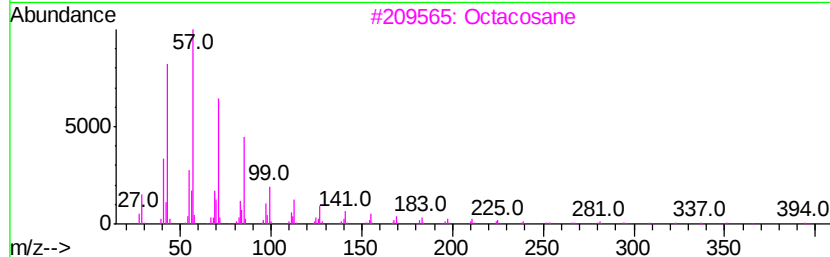
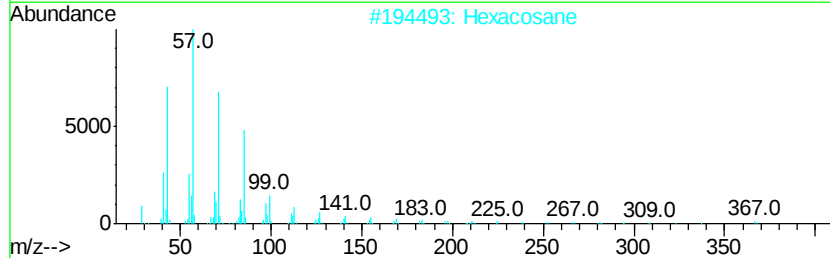
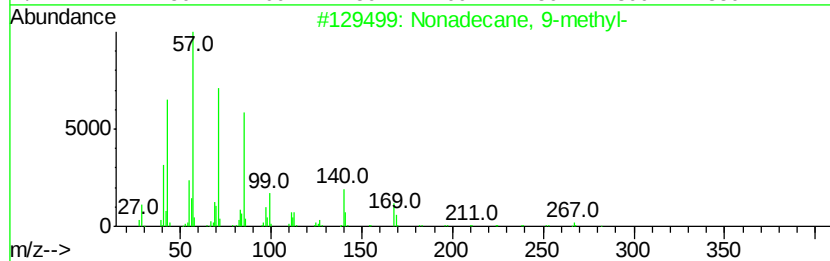
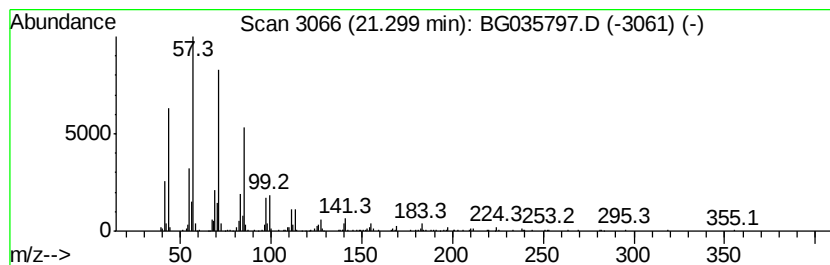
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Nonadecane, 9-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.30	5.40 ng	289944	Chrysene-d12	21.66

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonadecane, 9-methyl-	282	C20H42	013287-24-6	93
2		Hexacosane	366	C26H54	000630-01-3	90
3		Octacosane	394	C28H58	000630-02-4	90
4		Heneicosane, 11-decyl-	437	C31H64	055320-06-4	90
5		Heptadecane	240	C17H36	000629-78-7	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG072018\
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Sample : J4083-01
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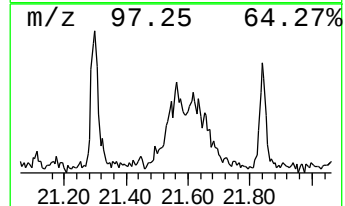
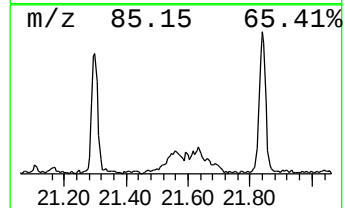
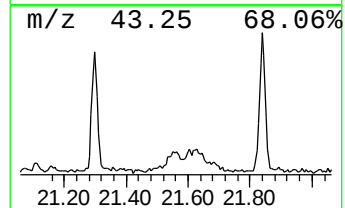
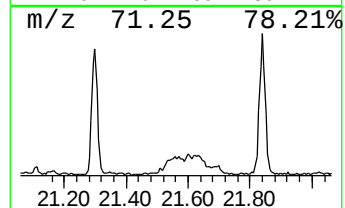
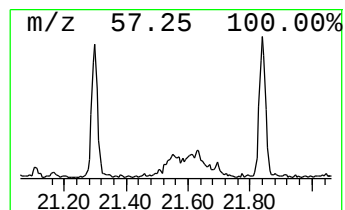
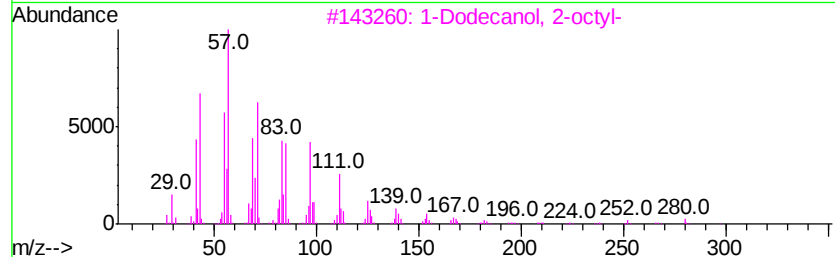
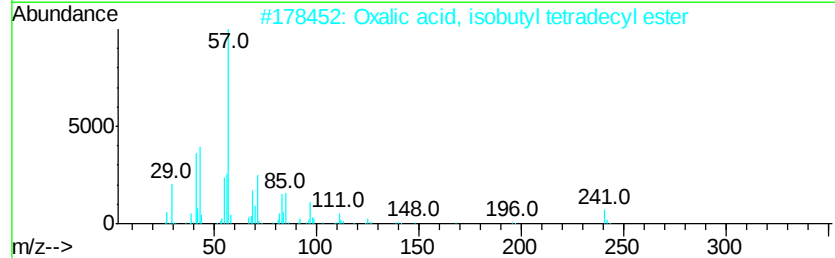
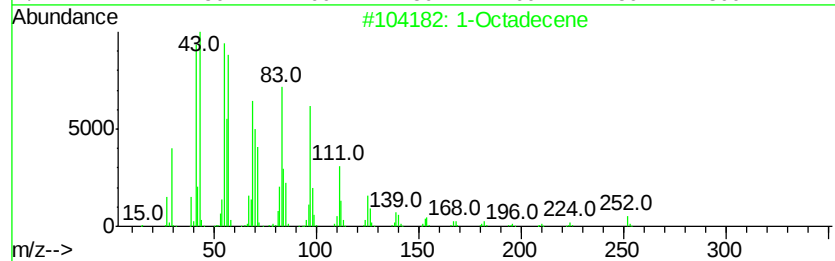
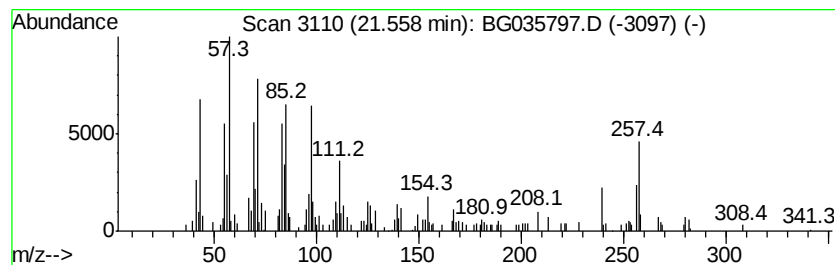
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 1-Octadecene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.56	4.61 ng	247677	Chrysene-d12	21.66
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1-Octadecene		252 C18H36	000112-88-9 55
2	Oxalic acid, isobutyl tetradecyl...		342 C20H38O4	1000309-37-9 53
3	1-Dodecanol, 2-octyl-		298 C20H42O	005333-42-6 49
4	2-Octyldodecyl isobutylate		368 C24H48O2	1000368-55-5 49
5	1-Decanol, 2-hexyl-		242 C16H34O	002425-77-6 46



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG072018\
Data File : BG035797.D
Acq On : 20 Jul 2018 23:19
Operator : JU/SJ
Sample : J4083-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

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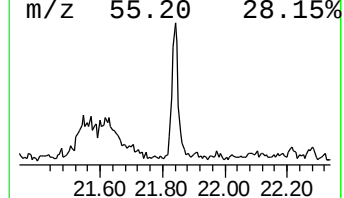
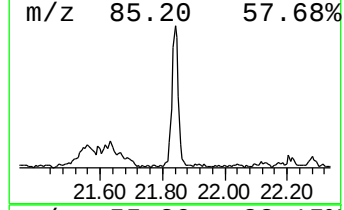
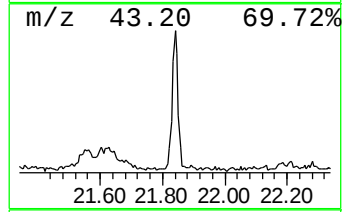
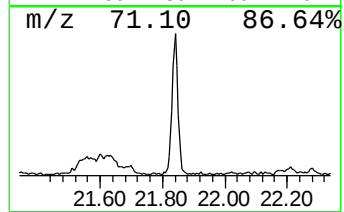
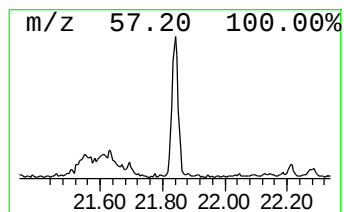
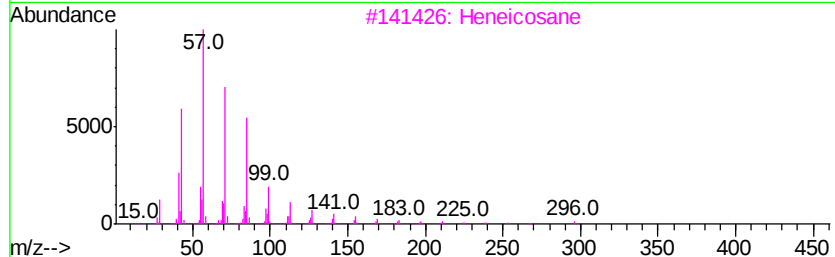
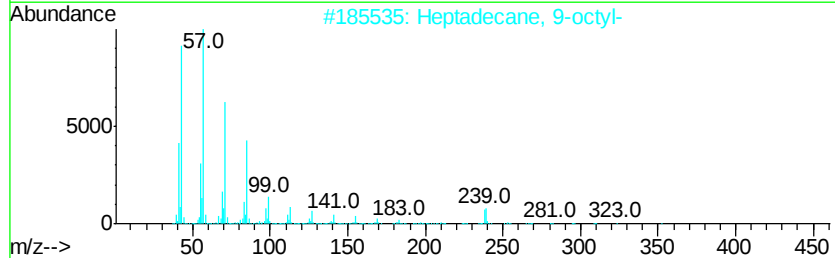
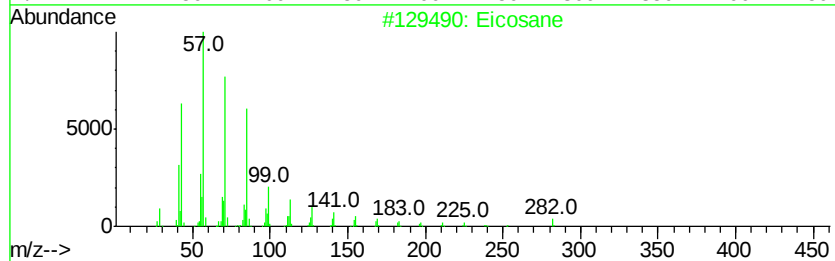
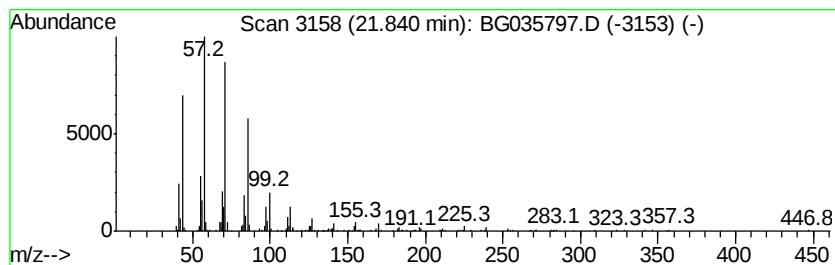
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Heneicosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.84	5.82 ng	312738	Chrysene-d12	21.66

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	94
2			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
3			Heneicosane	296	C21H44	000629-94-7	91
4			Hexadecane, 5-butyl-	282	C20H42	006912-07-8	90
5			Hexacosane	366	C26H54	000630-01-3	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG072018\
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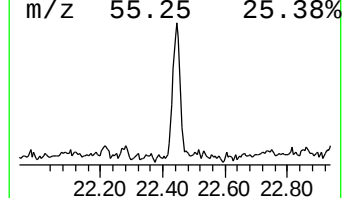
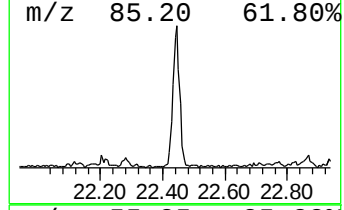
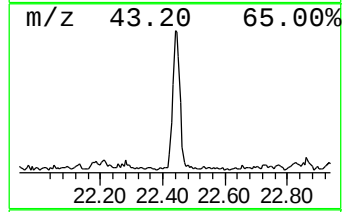
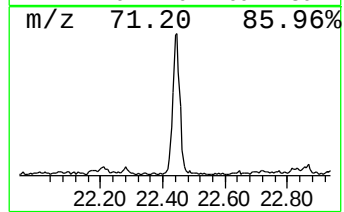
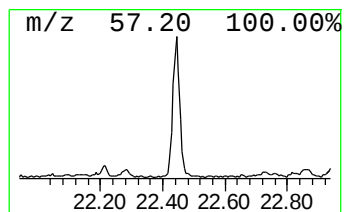
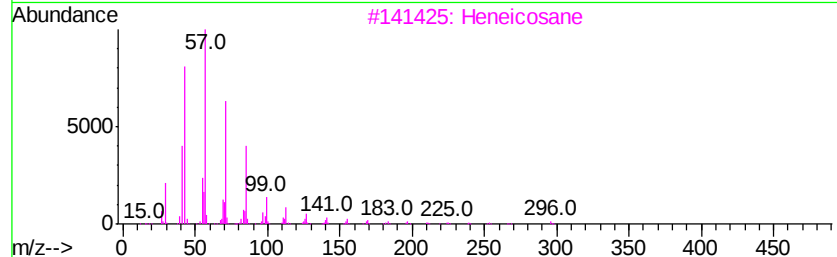
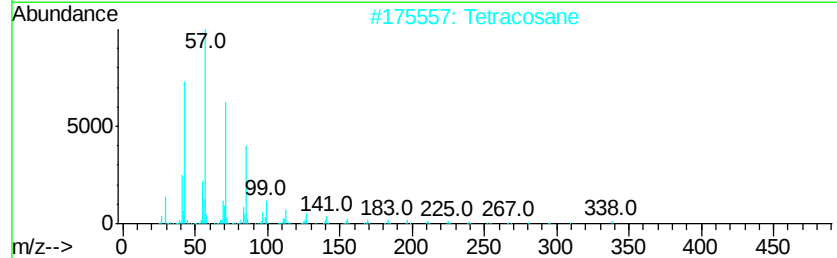
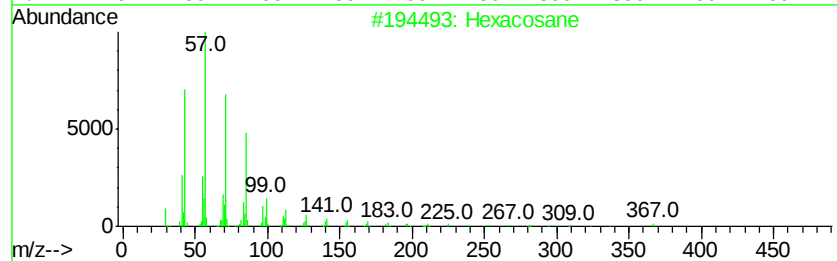
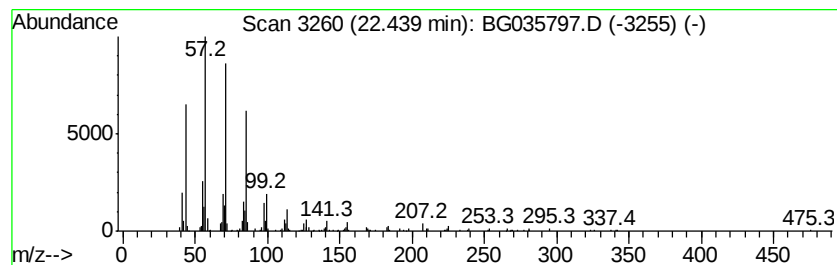
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Tetracosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.44	6.49 ng	348342	Chrysene-d12	21.66

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosane	366	C26H54	000630-01-3	91
2			Tetracosane	338	C24H50	000646-31-1	91
3			Heneicosane	296	C21H44	000629-94-7	91
4			2-methyloctacosane	408	C29H60	1000376-72-8	91
5			Heptadecane	240	C17H36	000629-78-7	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG072018\
Data File : BG035797.D
Acq On : 20 Jul 2018 23:19
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Sample : J4083-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

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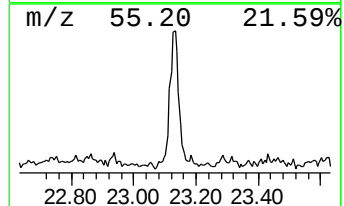
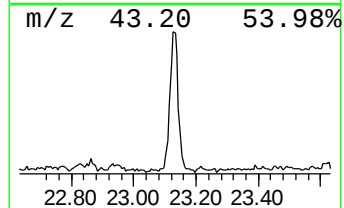
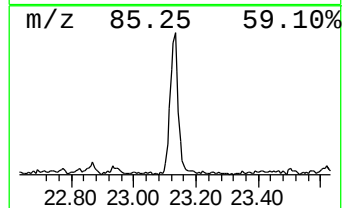
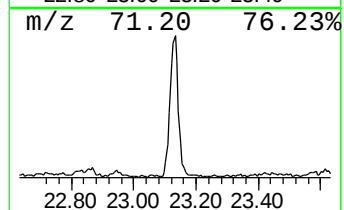
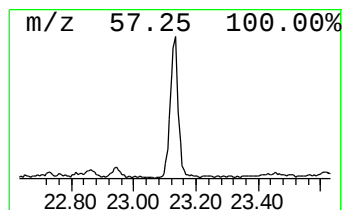
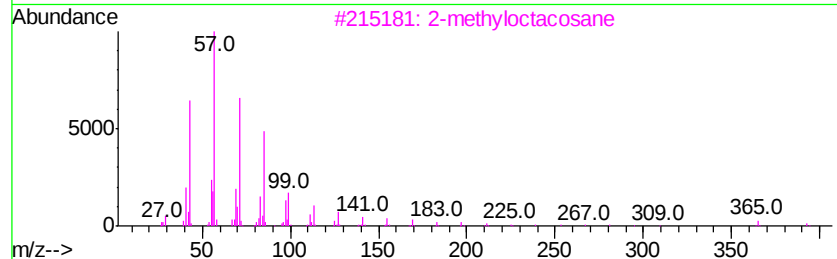
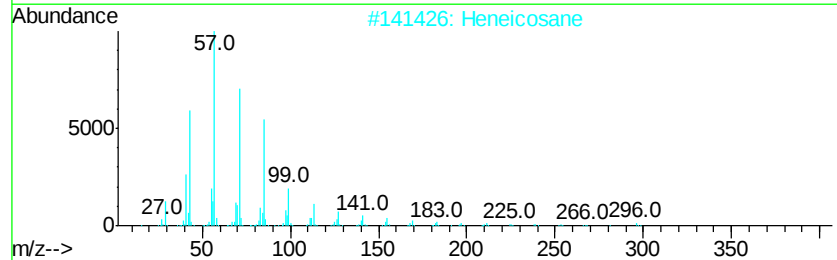
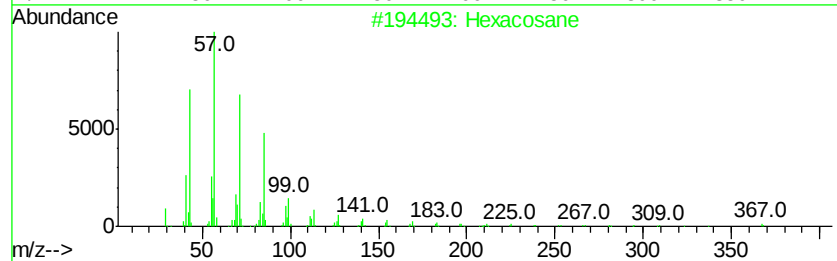
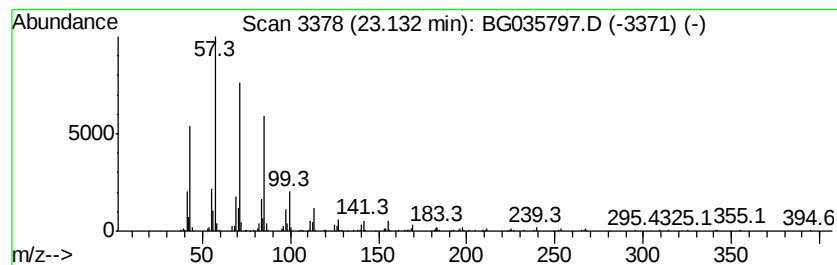
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Hexacosane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.13	7.80 ng	418567	Chrysene-d12	21.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	91
2		Heneicosane	296	C21H44	000629-94-7	91
3		2-methyloctacosane	408	C29H60	1000376-72-8	91
4		Triacontane	422	C30H62	000638-68-6	90
5		Octacosane	394	C28H58	000630-02-4	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG072018\
Data File : BG035797.D
Acq On : 20 Jul 2018 23:19
Operator : JU/SJ
Sample : J4083-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

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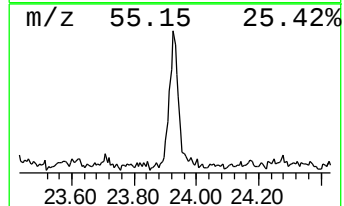
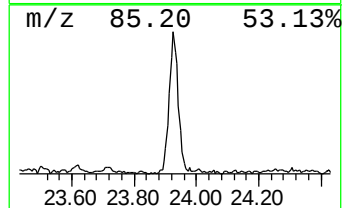
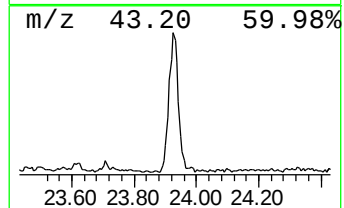
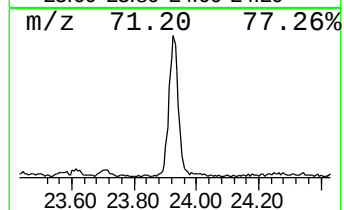
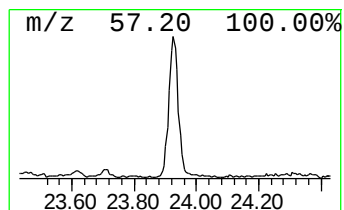
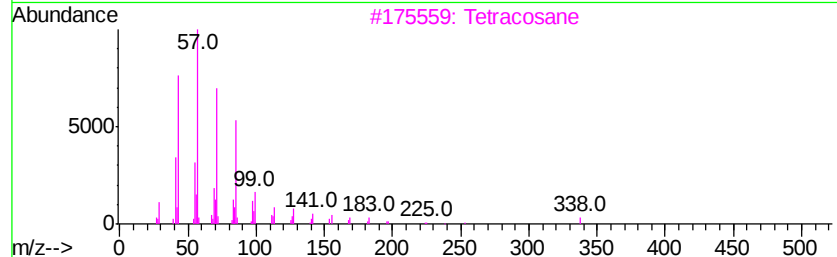
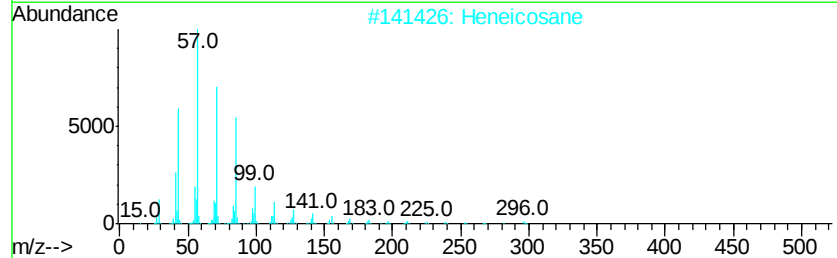
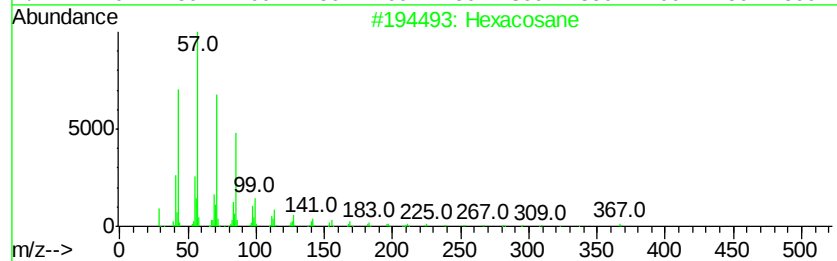
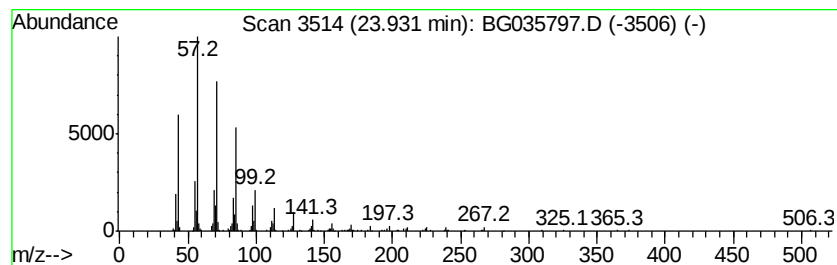
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Octacosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.93	6.61 ng	433519	Perylene-d12	24.85

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosane	366	C26H54	000630-01-3	91
2			Heneicosane	296	C21H44	000629-94-7	91
3			Tetracosane	338	C24H50	000646-31-1	91
4			Octacosane	394	C28H58	000630-02-4	91
5			triacontane	422	C30H62	000638-68-6	91



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Data File : BG035797.D
Acq On : 20 Jul 2018 23:19
Operator : JU/SJ
Sample : J4083-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
MH-6

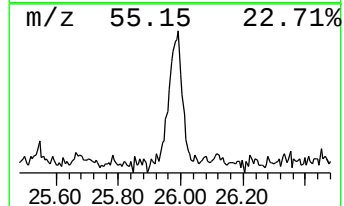
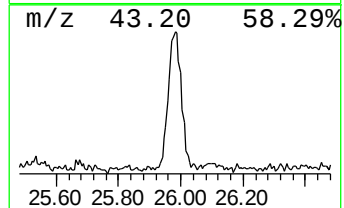
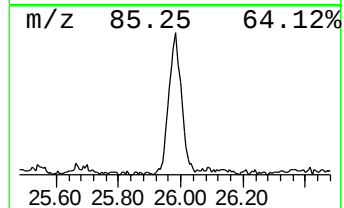
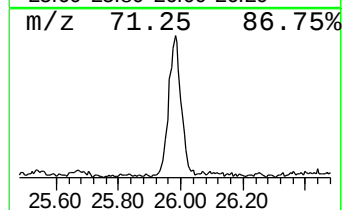
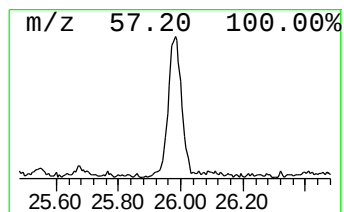
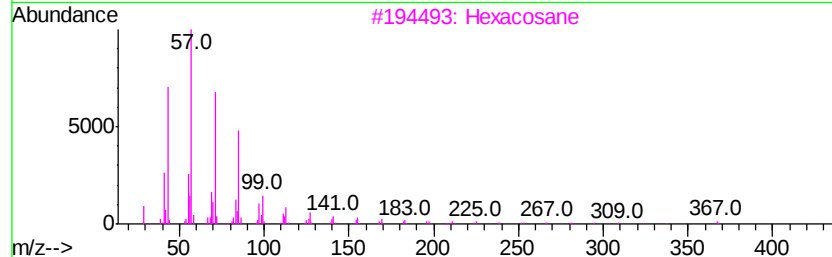
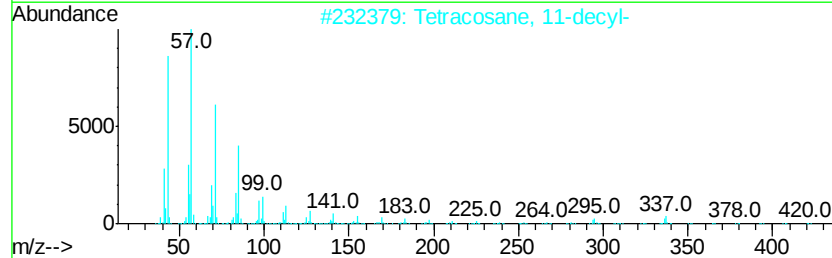
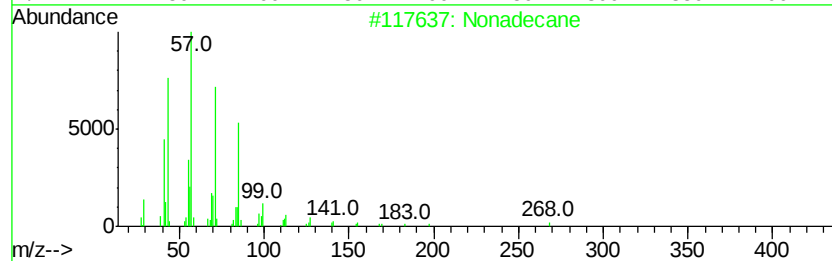
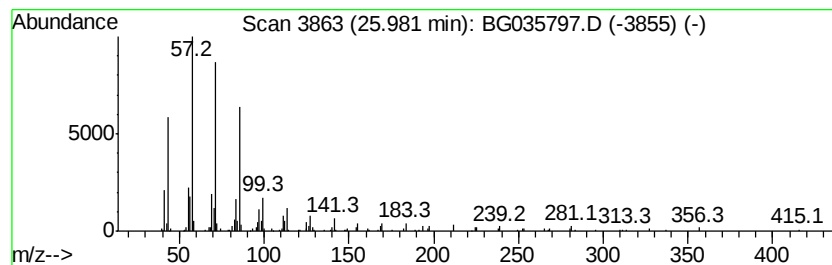
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Nonadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.98	5.35 ng	350613	Perylene-d12	24.85

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	94
2			Tetracosane, 11-decyl-	479	C34H70	055429-84-0	91
3			Hexacosane	366	C26H54	000630-01-3	91
4			triacontane	422	C30H62	000638-68-6	91
5			Heptacosane	380	C27H56	000593-49-7	91



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Operator : JU/SJ
Sample : J4083-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
MH-6

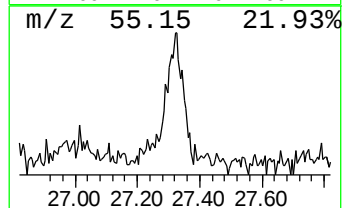
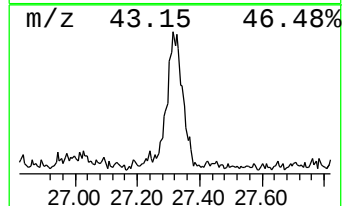
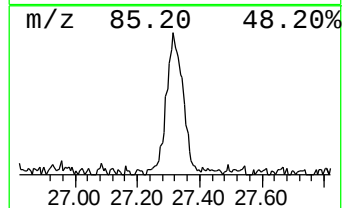
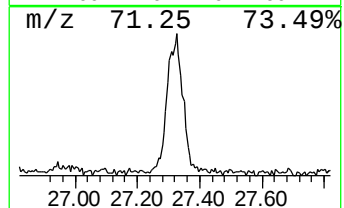
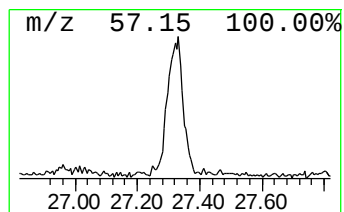
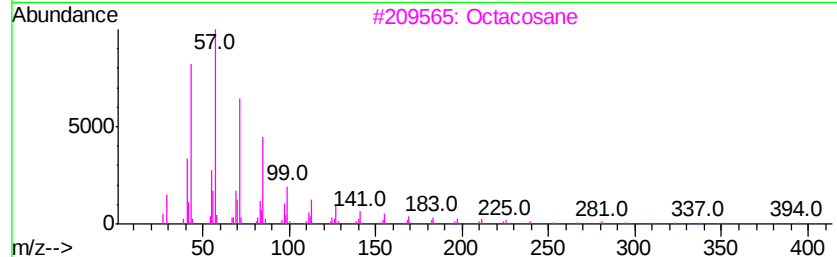
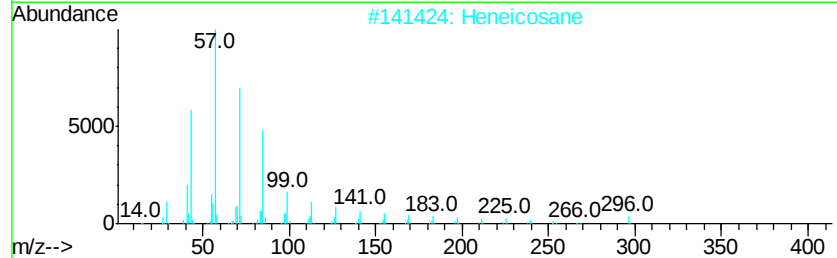
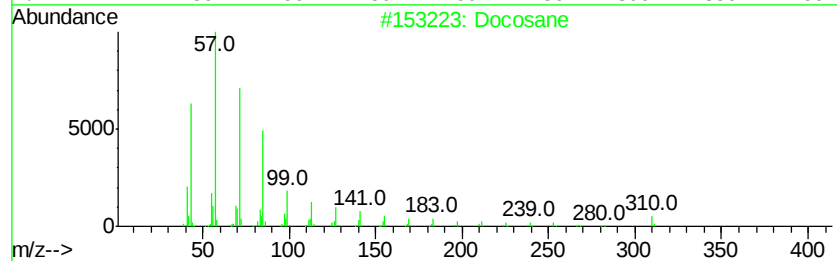
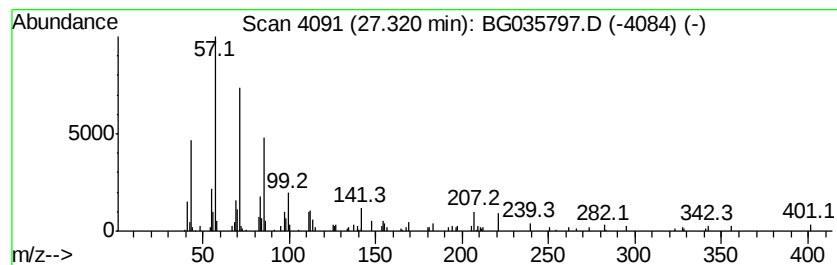
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Docosane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.32	3.80 ng	249049	Perylene-d12	24.85

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Docosane	310	C22H46	000629-97-0	72
2			Heneicosane	296	C21H44	000629-94-7	72
3			Octacosane	394	C28H58	000630-02-4	72
4			Hentriacontane	437	C31H64	000630-04-6	72
5			Tetratriacontane	479	C34H70	014167-59-0	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG072018\
Data File : BG035797.D
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Operator : JU/SJ
Sample : J4083-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
MH-6

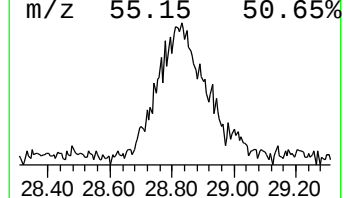
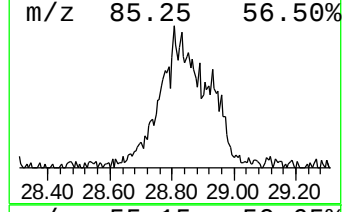
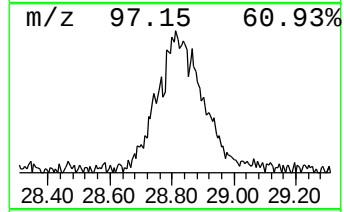
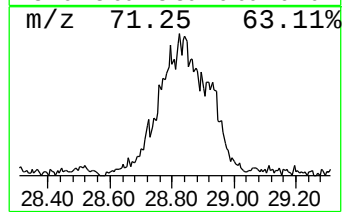
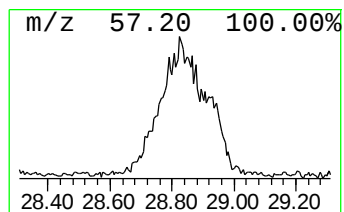
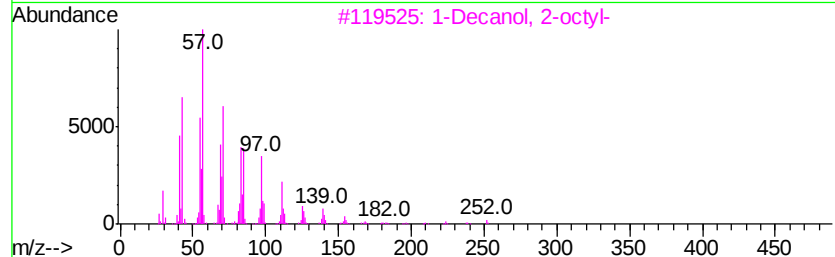
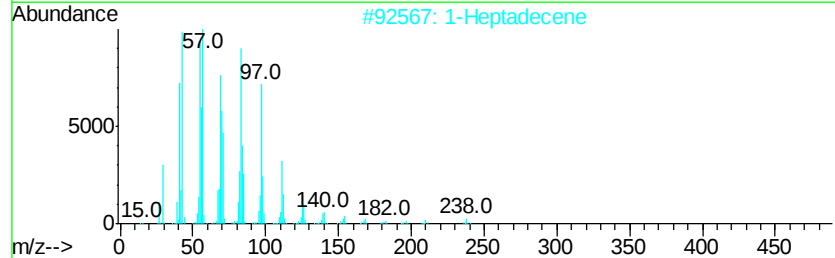
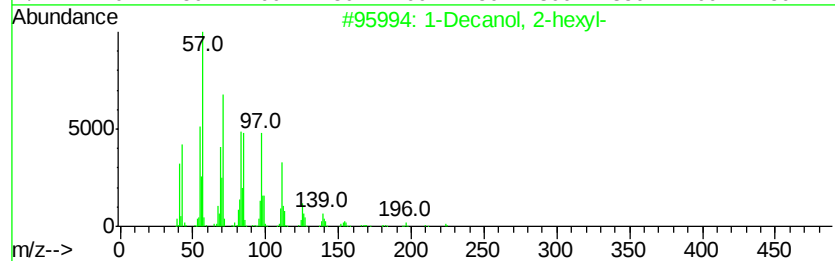
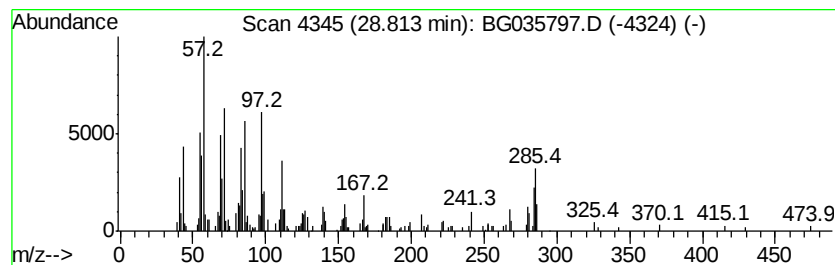
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 1-Decanol, 2-hexyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
28.81	6.14 ng	402619	Perylene-d12	24.85

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Decanol, 2-hexyl-	242	C16H34O	002425-77-6	68
2			1-Heptadecene	238	C17H34	006765-39-5	64
3			1-Decanol, 2-octyl-	270	C18H38O	045235-48-1	64
4			2-Octyldodecyl isobutylate	368	C24H48O2	1000368-55-5	58
5			1-Octadecene	252	C18H36	000112-88-9	55



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG072018\
 Data File : BG035797.D
 Acq On : 20 Jul 2018 23:19
 Operator : JU/SJ
 Sample : J4083-01
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MH-6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG071218.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
Guanidine	5.14	4.2	ng	69804	1	8.03	330799	20.0
unknown7.56	7.56	86.9	ng	1437330	1	8.03	330799	20.0
n-Hexadecanoic acid	18.28	3.0	ng	121137	4	17.39	805738	20.0
6-Octadecenoic acid	19.43	3.9	ng	155182	4	17.39	805738	20.0
Eicosane	20.30	2.1	ng	111368	5	21.66	1073830	20.0
Hexadecane	20.79	4.4	ng	237702	5	21.66	1073830	20.0
Nonadecane, 9-met...	21.30	5.4	ng	289944	5	21.66	1073830	20.0
1-Octadecene	21.56	4.6	ng	247677	5	21.66	1073830	20.0
Heneicosane	21.84	5.8	ng	312738	5	21.66	1073830	20.0
Tetracosane	22.44	6.5	ng	348342	5	21.66	1073830	20.0
Hexacosane	23.13	7.8	ng	418567	5	21.66	1073830	20.0
Octacosane	23.93	6.6	ng	433519	6	24.85	1311720	20.0
Nonadecane	25.98	5.3	ng	350613	6	24.85	1311720	20.0
Docosane	27.32	3.8	ng	249049	6	24.85	1311720	20.0
1-Decanol, 2-hexyl-	28.81	6.1	ng	402619	6	24.85	1311720	20.0