

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060815\
 Data File : BG017558.D
 Acq On : 9 Jun 2015 4:18
 Operator : TP/IZ
 Sample : G2450-34
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 AVE-E-NBS13.2(3)

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:\HPCHEM1\BNA_G\METHODS\8270-BG060315.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	2.738	4	8	15	rVB	63976	76871	6.59%	0.635%
2	4.707	338	343	351	rBV	21743	32765	2.81%	0.270%
3	5.200	420	427	438	rBV	421398	606765	52.00%	5.009%
4	6.816	695	702	712	rBV	458978	693971	59.47%	5.729%
5	7.145	750	758	767	rBV	453964	706983	60.59%	5.837%
6	7.603	830	836	843	rVB	99975	152990	13.11%	1.263%
7	7.920	884	890	898	rVB	313492	498069	42.68%	4.112%
8	8.767	1027	1034	1041	rBV	277549	443303	37.99%	3.660%
9	10.394	1305	1311	1318	rBV	128258	225645	19.34%	1.863%
10	12.074	1593	1597	1603	rVB3	8847	14221	1.22%	0.117%
11	12.298	1628	1635	1641	rVB	8791	15687	1.34%	0.130%
12	12.885	1725	1735	1744	rBV	580717	810622	69.47%	6.692%
13	13.091	1761	1770	1776	rBV5	9480	16532	1.42%	0.136%
14	13.590	1850	1855	1859	rBV6	6941	12430	1.07%	0.103%
15	13.731	1873	1879	1889	rBV	36038	62853	5.39%	0.519%
16	14.172	1949	1954	1960	rVB2	11609	15786	1.35%	0.130%
17	14.272	1965	1971	1978	rVV2	198861	297406	25.49%	2.455%
18	14.671	2035	2039	2044	rVB3	9007	12768	1.09%	0.105%
19	15.130	2113	2117	2121	rVB2	15317	19658	1.68%	0.162%
20	15.318	2144	2149	2153	rVB6	10086	15998	1.37%	0.132%
21	15.770	2217	2226	2238	rBV	591043	826890	70.86%	6.826%
22	15.993	2259	2264	2266	rBV2	19254	27146	2.33%	0.224%
23	16.563	2353	2361	2365	rBV4	7084	15693	1.34%	0.130%
24	16.675	2377	2380	2387	rVB5	8260	13428	1.15%	0.111%
25	16.787	2393	2399	2402	rBV6	11712	18669	1.60%	0.154%
26	17.022	2433	2439	2443	rBV	255775	348126	29.83%	2.874%
27	17.063	2443	2446	2452	rVB	82716	111579	9.56%	0.921%
28	17.157	2458	2462	2466	rVB	21604	33781	2.89%	0.279%
29	17.897	2584	2588	2590	rBV2	16804	16918	1.45%	0.140%
30	17.932	2590	2594	2601	rVB4	60506	115680	9.91%	0.955%
31	18.032	2606	2611	2616	rVV6	9409	16558	1.42%	0.137%
32	18.091	2616	2621	2626	rVV4	33408	61997	5.31%	0.512%
33	18.132	2626	2628	2633	rVB	13865	13691	1.17%	0.113%
34	18.214	2638	2642	2647	rBV5	9949	16213	1.39%	0.134%

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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:\HPCHEM1\BNA_G\METHODS\8270-BG060315.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	18.408	2670	2675	2678	rBV2	15974	24586	2.11%	0.203%
36	18.449	2679	2682	2686	rVB5	14570	20189	1.73%	0.167%
37	18.825	2742	2746	2750	rBV4	15312	25512	2.19%	0.211%
38	18.966	2767	2770	2777	rVB7	16463	26210	2.25%	0.216%
39	19.090	2785	2791	2796	rBV	319139	419449	35.95%	3.463%
40	19.190	2805	2808	2811	rBV5	7626	12288	1.05%	0.101%
41	19.360	2833	2837	2841	rVB	76596	92702	7.94%	0.765%
42	19.454	2843	2853	2862	rVV	259394	476063	40.80%	3.930%
43	19.530	2862	2866	2872	rVB2	16267	26152	2.24%	0.216%
44	19.660	2879	2888	2893	rBV	918311	1166874	100.00%	9.633%
45	19.695	2893	2894	2898	rVB2	16213	15955	1.37%	0.132%
46	19.771	2904	2907	2915	rBV6	21526	54053	4.63%	0.446%
47	19.953	2928	2938	2944	rVB2	59428	109329	9.37%	0.903%
48	20.053	2951	2955	2958	rBV	49245	63367	5.43%	0.523%
49	20.147	2969	2971	2975	rVB5	18854	20242	1.73%	0.167%
50	20.194	2976	2979	2984	rBV7	13259	20204	1.73%	0.167%
51	20.247	2986	2988	2992	rVB5	11782	12468	1.07%	0.103%
52	20.294	2993	2996	2999	rVB2	14408	19606	1.68%	0.162%
53	20.406	3012	3015	3022	rVB3	69036	86671	7.43%	0.716%
54	20.470	3022	3026	3029	rBV3	30715	37522	3.22%	0.310%
55	20.864	3090	3093	3096	rBV2	24103	31184	2.67%	0.257%
56	20.899	3097	3099	3101	rVV3	28567	29421	2.52%	0.243%
57	20.929	3101	3104	3111	rVB3	106767	165685	14.20%	1.368%
58	21.134	3136	3139	3142	rBV	36028	36655	3.14%	0.303%
59	21.217	3146	3153	3157	rVV2	423161	789461	67.66%	6.517%
60	21.252	3157	3159	3168	rVB	201213	242896	20.82%	2.005%
61	21.369	3176	3179	3184	rBV2	81417	93813	8.04%	0.774%
62	21.798	3249	3252	3262	rVB4	41667	87078	7.46%	0.719%
63	22.774	3412	3418	3424	rBV	188110	485210	41.58%	4.006%
64	23.255	3495	3500	3507	rVB	113894	196517	16.84%	1.622%
65	23.349	3511	3516	3525	rVB	185454	329549	28.24%	2.721%
66	23.449	3527	3533	3538	rBV	194183	354578	30.39%	2.927%
67	25.711	3914	3918	3927	rVB4	69480	173899	14.90%	1.436%

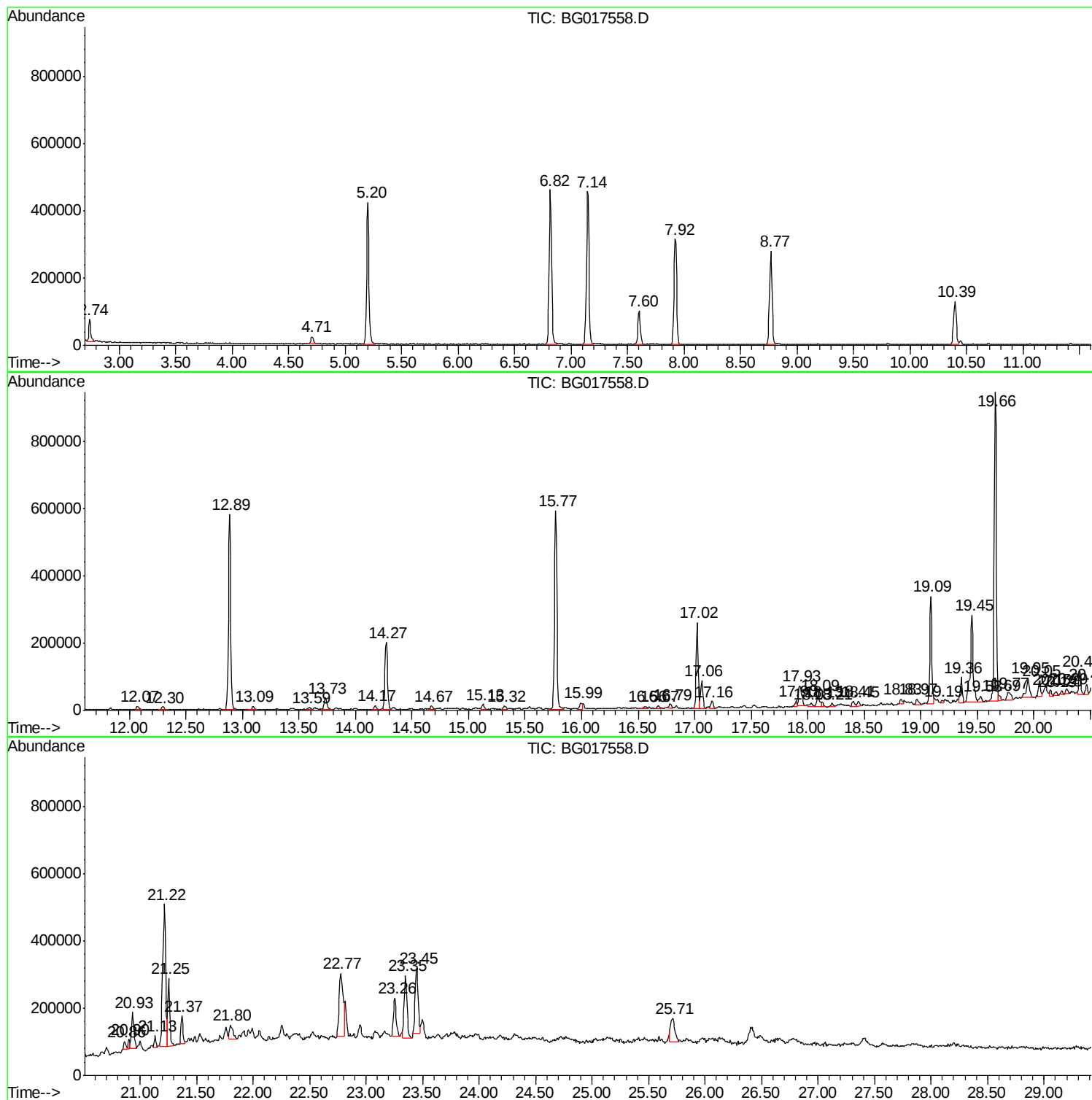
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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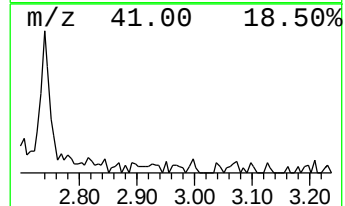
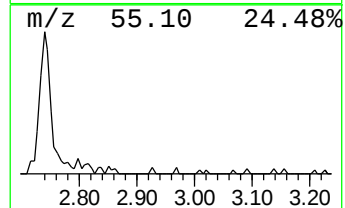
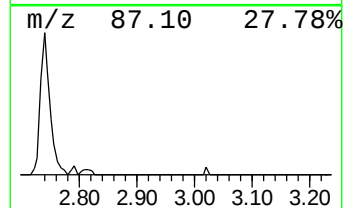
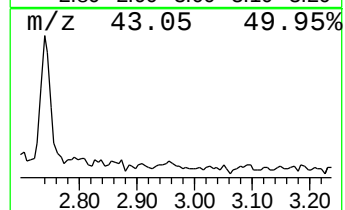
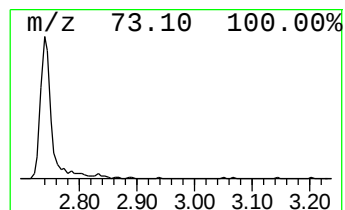
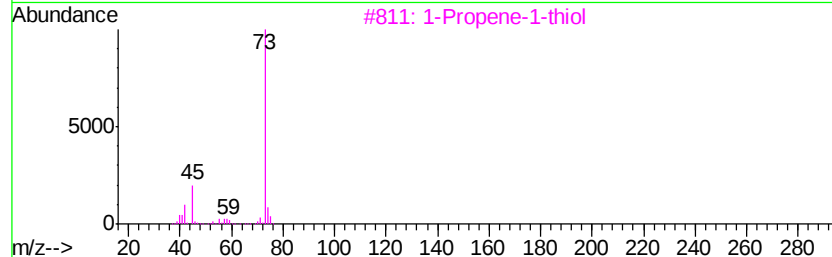
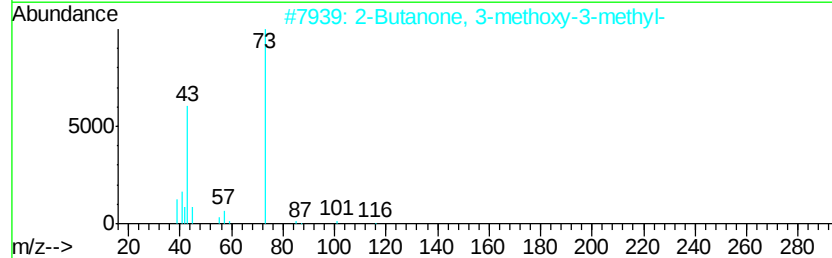
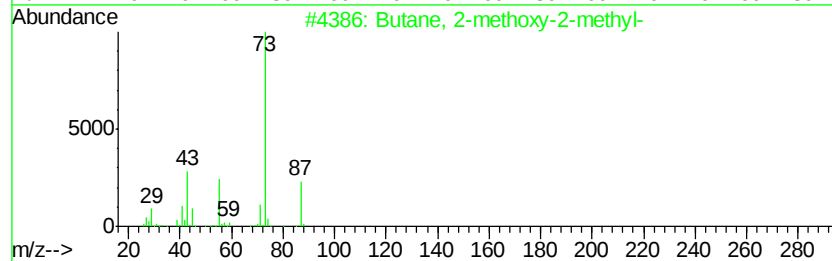
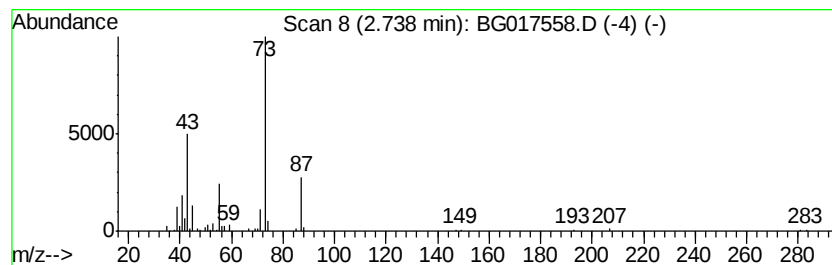
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.74	10.05 ng	76871	1,4-Dichlorobenzene-d4	7.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
2		2-Butanone, 3-methoxy-3-methyl-	116	C6H12O2	036687-98-6	16
3		1-Propene-1-thiol	74	C3H6S	000925-89-3	9
4		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9
5		Silane, trimethyl[2-methylene-1-...	194	C12H22Si	097778-15-9	9



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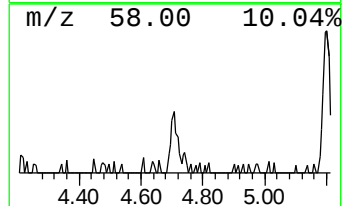
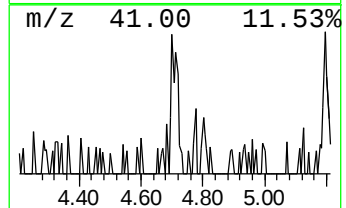
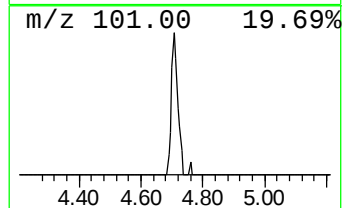
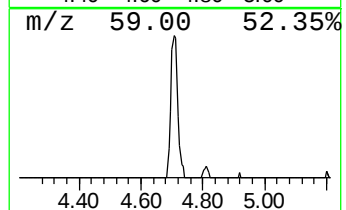
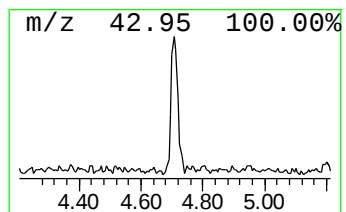
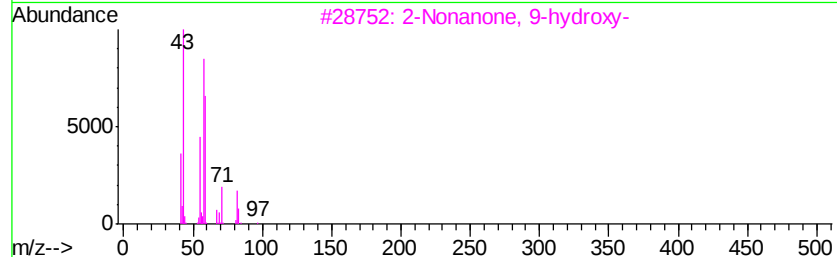
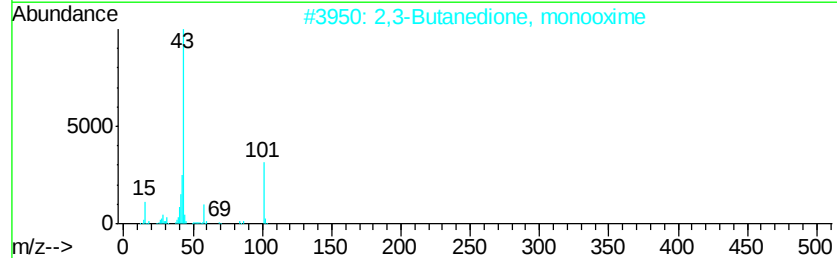
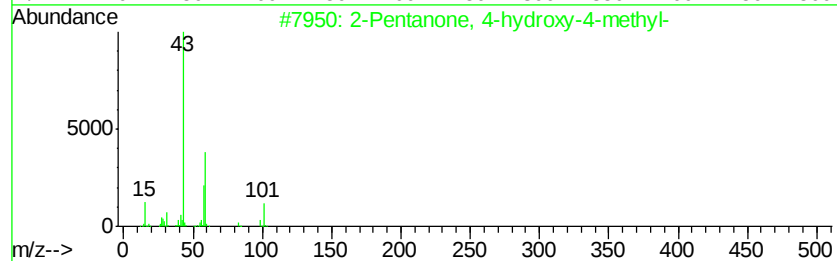
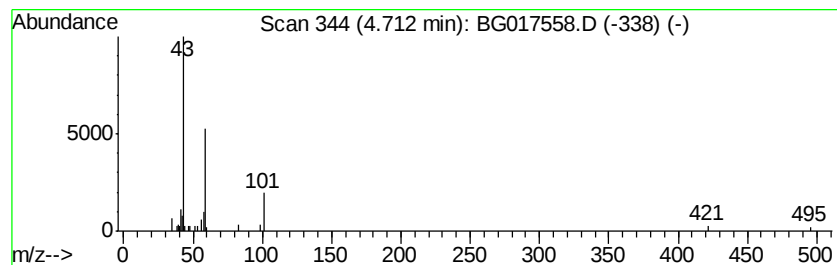
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.71	4.28 ng	32765	1,4-Dichlorobenzene-d4	7.60

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	35
3			2-Nonanone, 9-hydroxy-	158	C9H18O2	025368-56-3	25
4			5-Hexen-2-one	98	C6H10O	000109-49-9	10
5			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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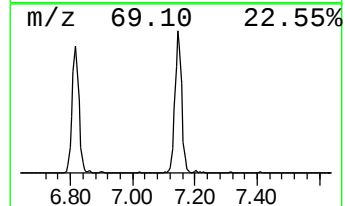
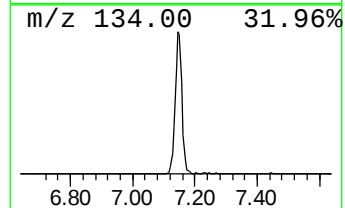
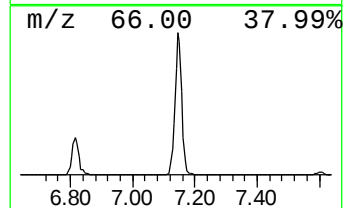
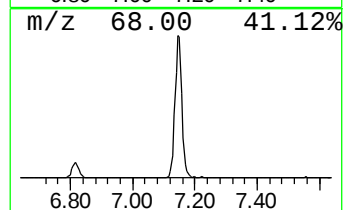
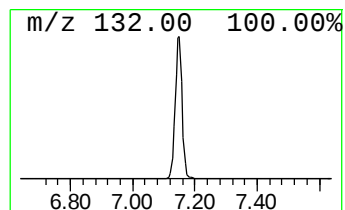
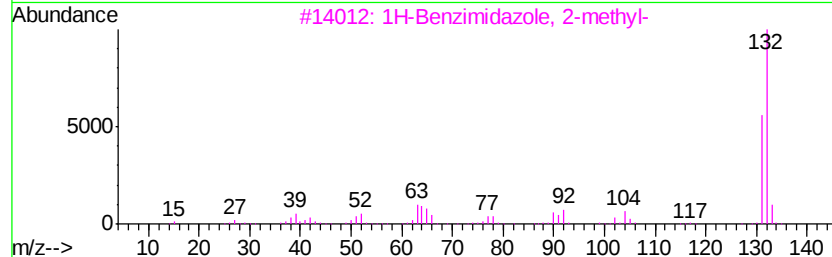
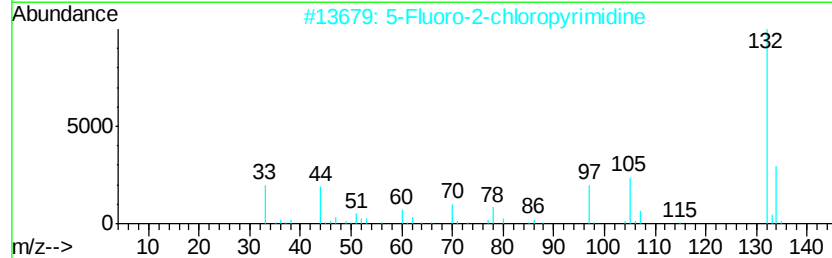
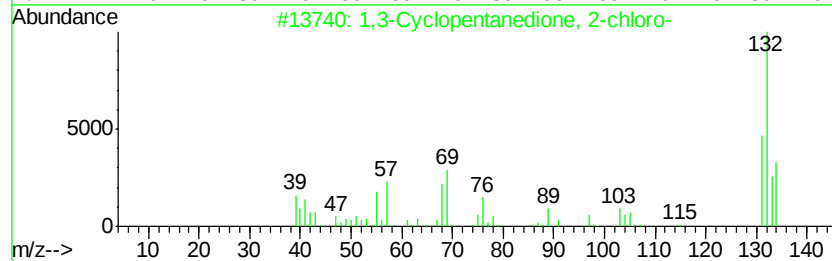
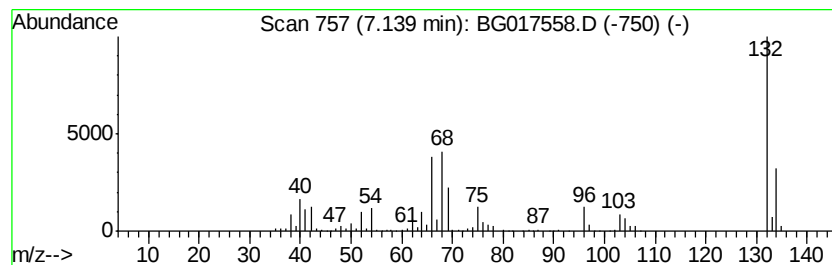
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown7.14 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.14	92.42 ng	706983	1,4-Dichlorobenzene-d4	7.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	35
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	27
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	27
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	22
5		Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	22



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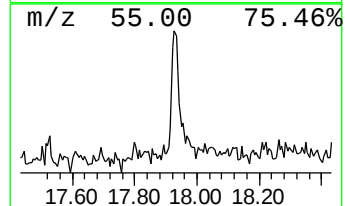
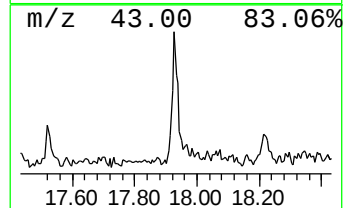
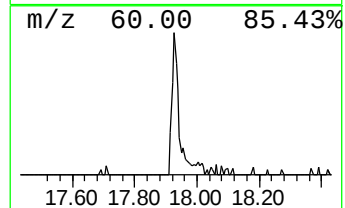
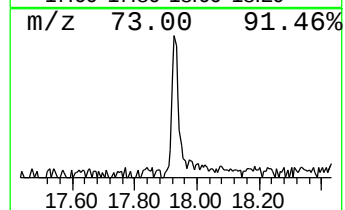
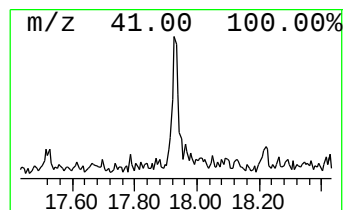
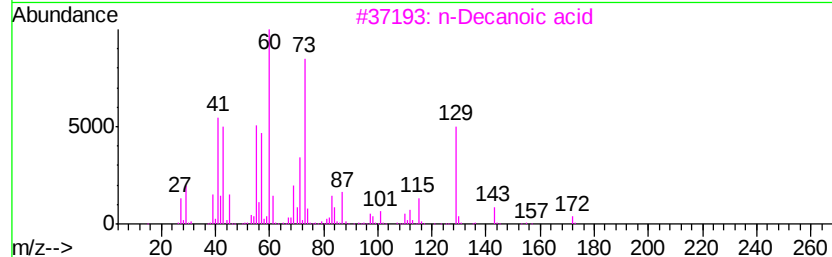
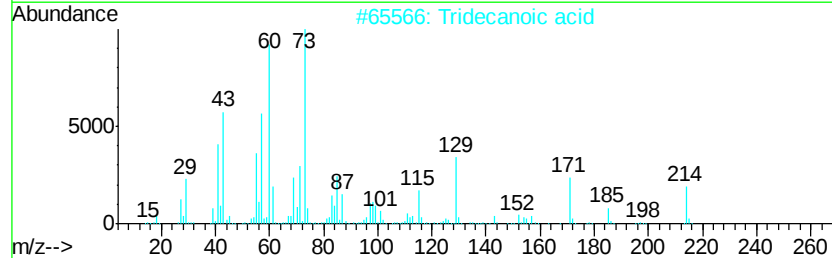
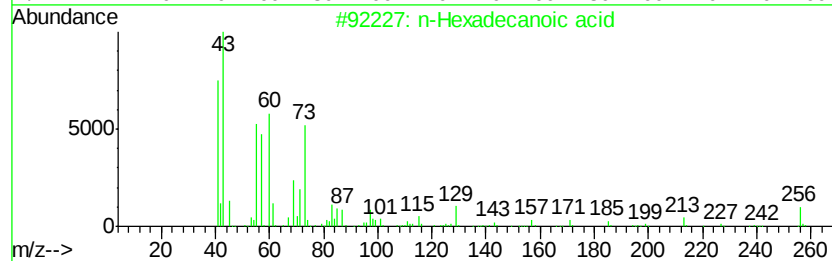
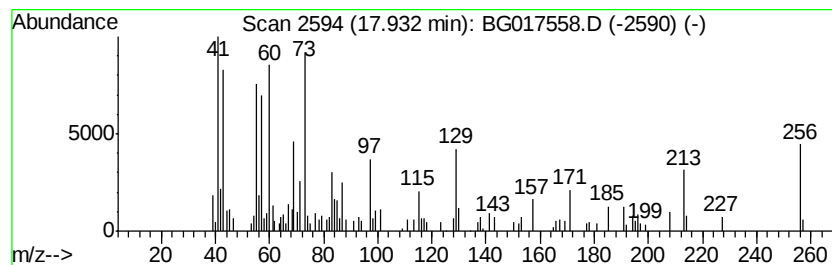
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.	
17.93	6.65 ng	115680	Phenanthrene-d10		17.02	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	80
2		Tridecanoic acid	214	C13H26O2	000638-53-9	58
3		n-Decanoic acid	172	C10H20O2	000334-48-5	43
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	35
5		Heptanoic acid	130	C7H14O2	000111-14-8	27



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Sample : G2450-34
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
AVE-E-NBS13.2(3)

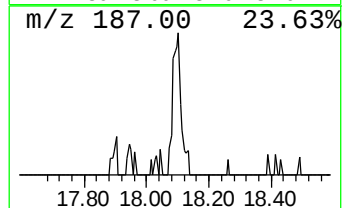
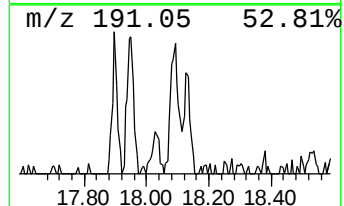
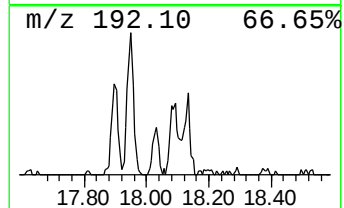
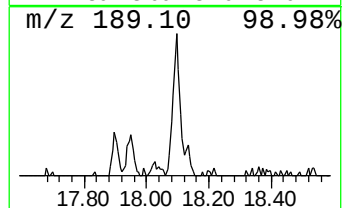
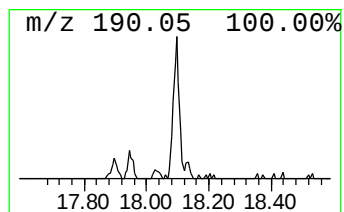
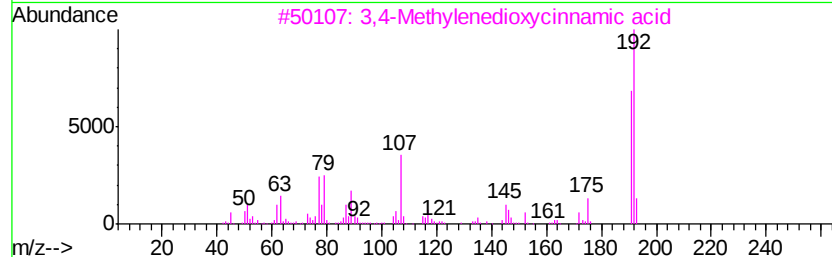
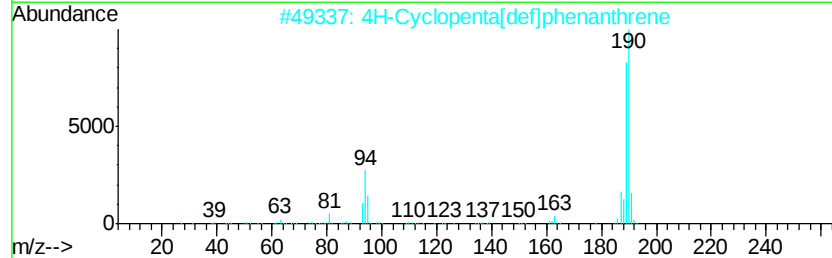
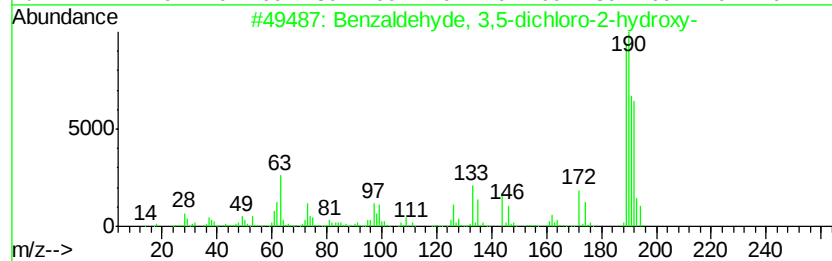
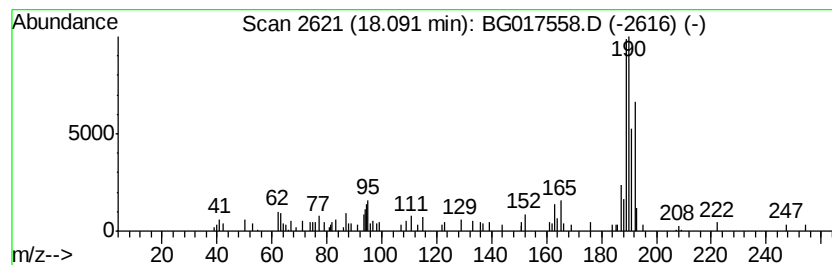
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzaldehyde, 3,5-dichloro-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.09	3.56 ng	61997	Phenanthrene-d10	17.02

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	74
2	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	49
3	3,4-Methylenedioxycinnamic acid	192	C10H8O4	002373-80-0	25
4	Naphtho[1,2-b]norbornadiene	192	C15H12	1000210-14-8	15
5	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	15



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060815\
Data File : BG017558.D
Acq On : 9 Jun 2015 4:18
Operator : TP/IZ
Sample : G2450-34
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
AVE-E-NBS13.2(3)

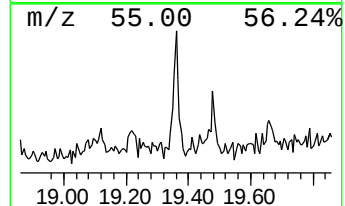
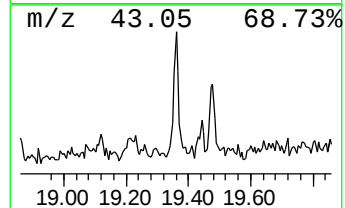
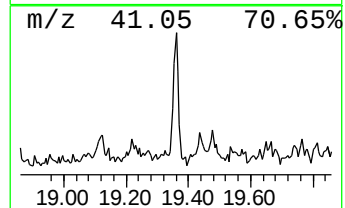
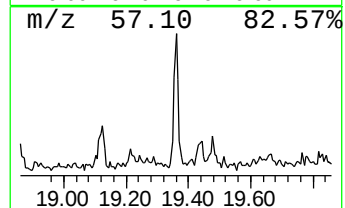
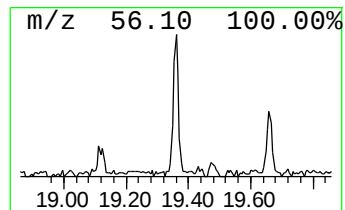
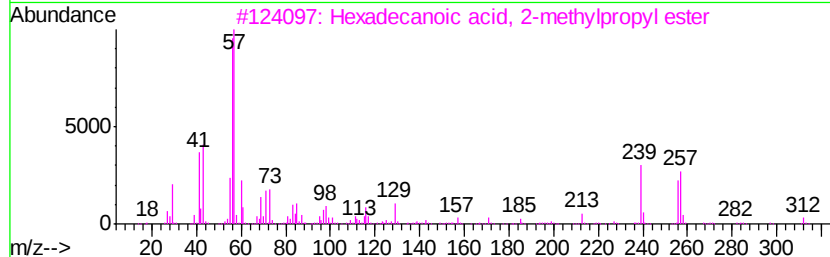
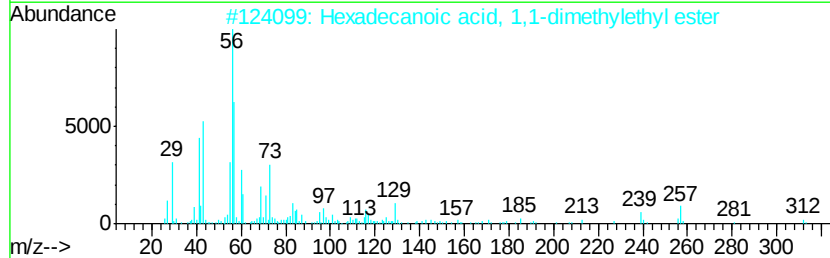
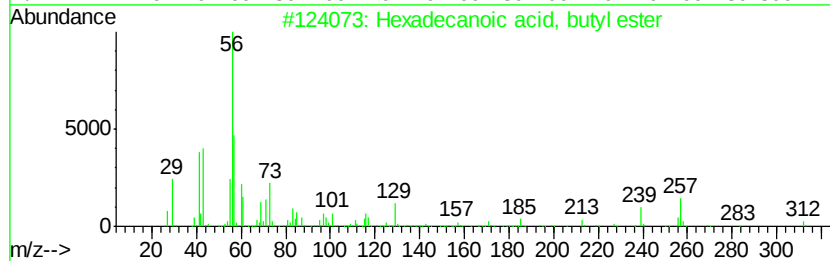
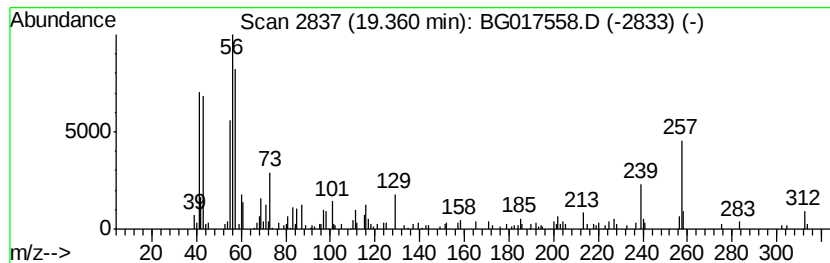
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Hexadecanoic acid, butyl ester Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.36	2.35 ng	92702	Chrysene-d12	21.22

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	94
2		Hexadecanoic acid, 1,1-dimethyle...	312	C20H40O2	031158-91-5	60
3		Hexadecanoic acid, 2-methylpropy...	312	C20H40O2	000110-34-9	42
4		Pentanoic acid, 4-oxo-	116	C5H8O3	000123-76-2	38
5		Cyclopentane, 1,1-dimethyl-	98	C7H14	001638-26-2	30



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060815\
Data File : BG017558.D
Acq On : 9 Jun 2015 4:18
Operator : TP/IZ
Sample : G2450-34
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
AVE-E-NBS13.2(3)

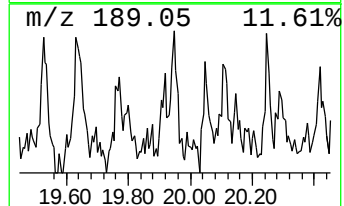
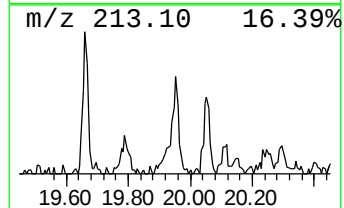
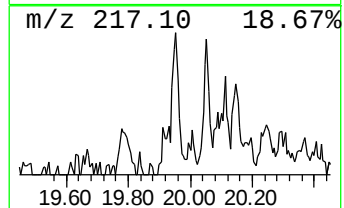
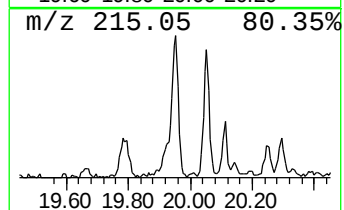
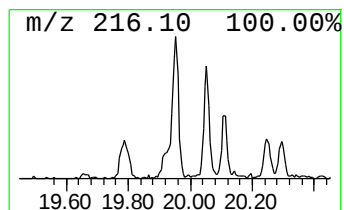
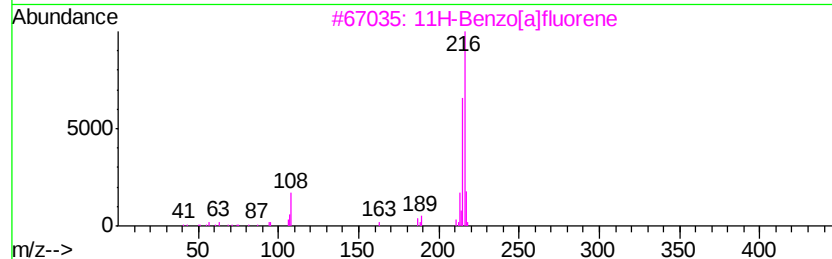
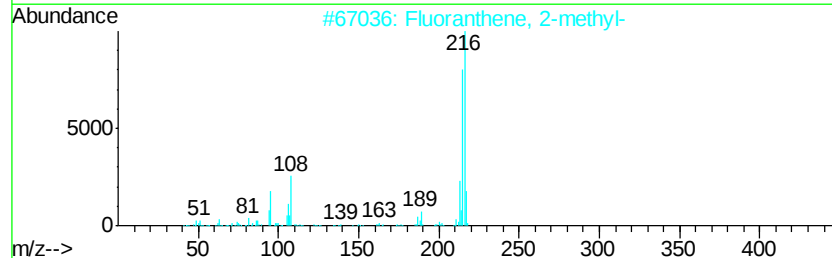
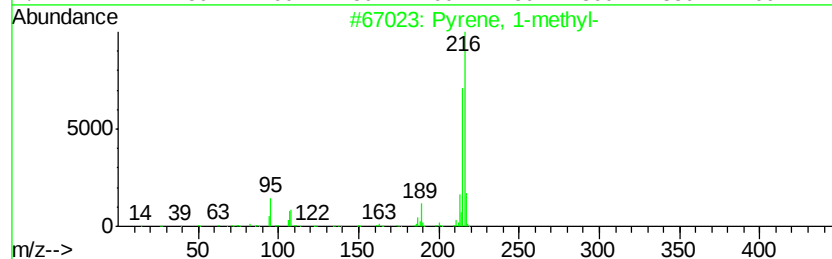
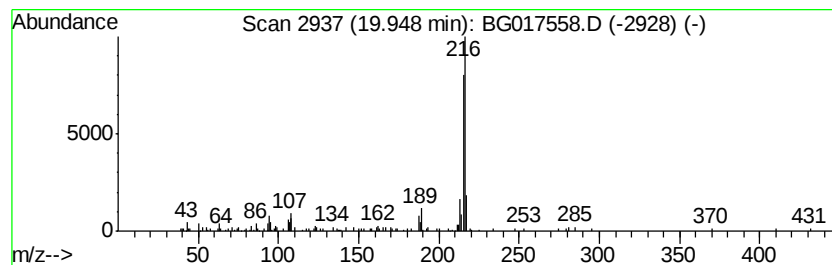
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Pyrene, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.95	2.77 ng	109329	Chrysene-d12	21.22

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pyrene, 1-methyl-	216	C17H12	002381-21-7	94
2		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	91
3		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	91
4		Pyrene, 2-methyl-	216	C17H12	003442-78-2	91
5		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	90



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060815\
Data File : BG017558.D
Acq On : 9 Jun 2015 4:18
Operator : TP/IZ
Sample : G2450-34
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
AVE-E-NBS13.2(3)

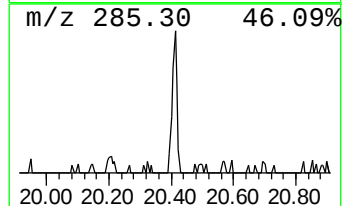
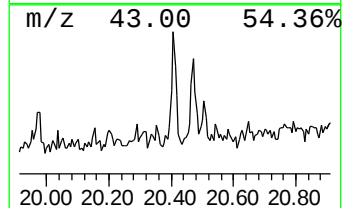
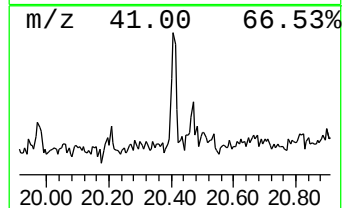
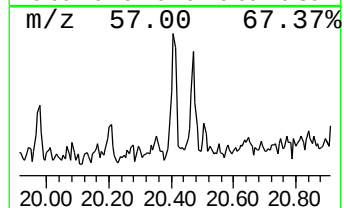
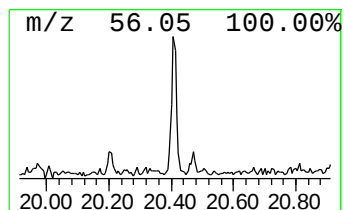
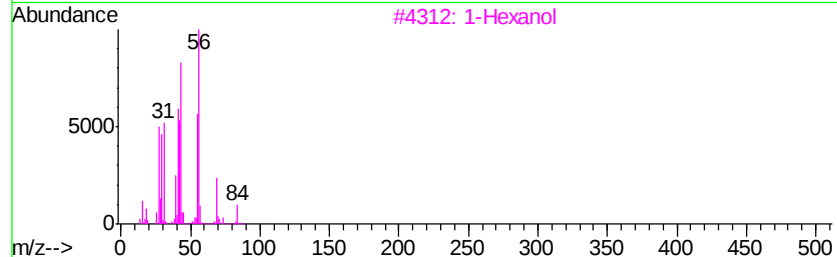
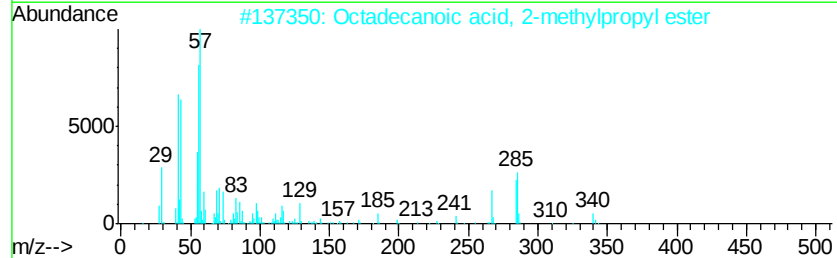
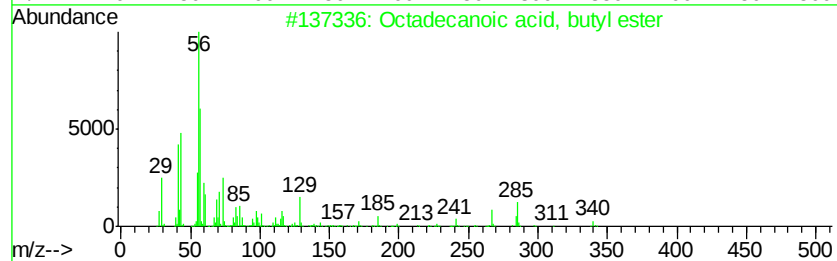
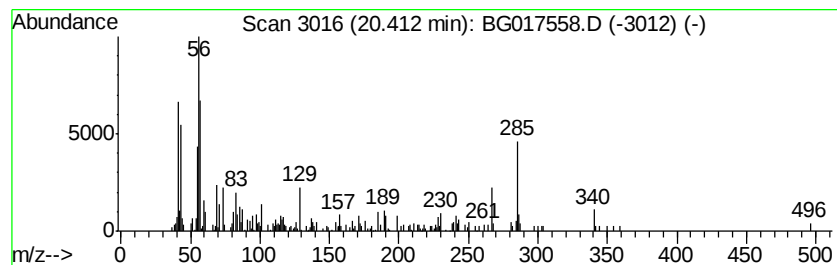
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Octadecanoic acid, butyl ester Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.41	2.20 ng	86671	Chrysene-d12	21.22

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid, butyl ester	340	C22H44O2	000123-95-5	94
2		Octadecanoic acid, 2-methylpropy...	340	C22H44O2	000646-13-9	93
3		1-Hexanol	102	C6H14O	000111-27-3	35
4		2-Methyl-n-1-tridecene	196	C14H28	018094-01-4	30
5		1-Hexene, 5-methyl-	98	C7H14	003524-73-0	30



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060815\
Data File : BG017558.D
Acq On : 9 Jun 2015 4:18
Operator : TP/IZ
Sample : G2450-34
Misc :
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Instrument :
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ClientSampleId :
AVE-E-NBS13.2(3)

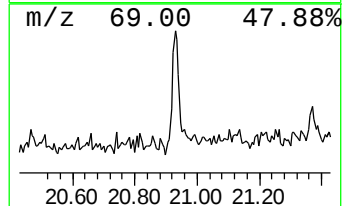
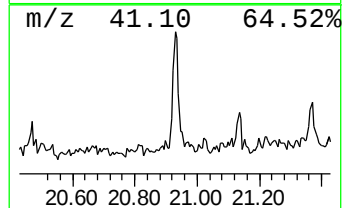
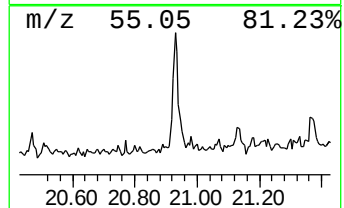
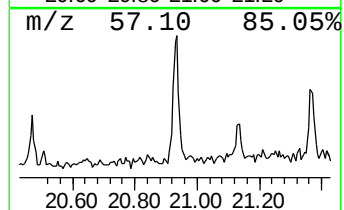
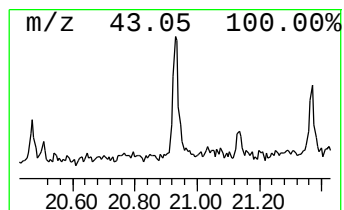
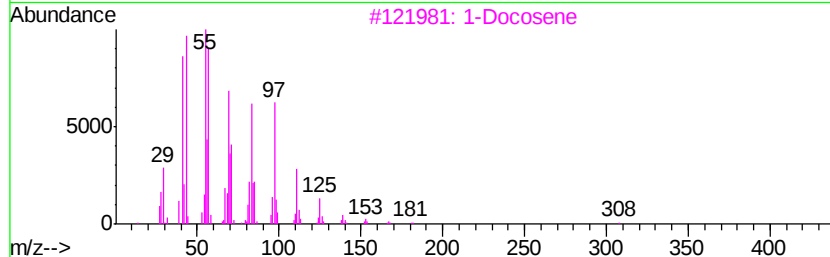
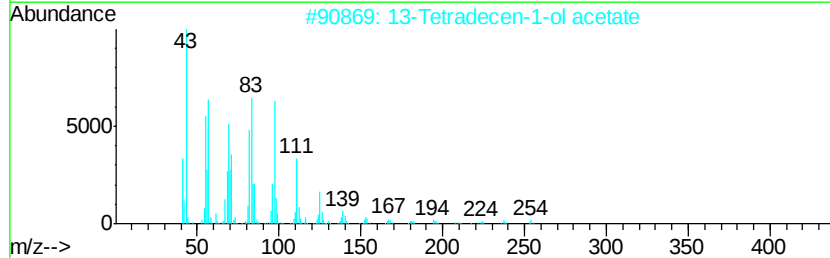
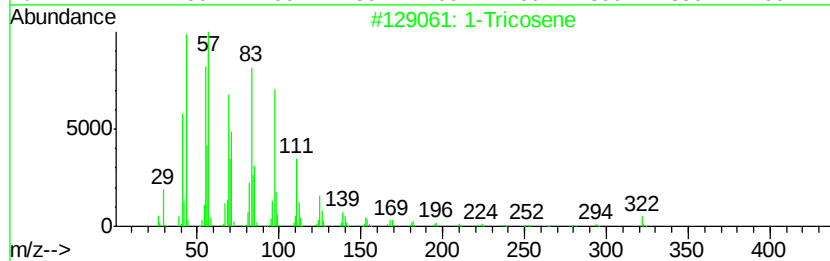
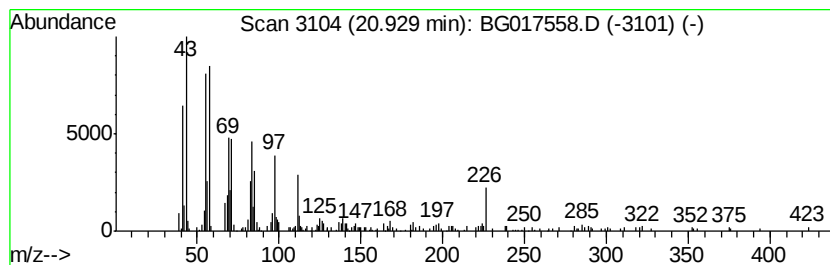
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 1-Tricosene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.93	4.20 ng	165685	Chrysene-d12	21.22

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Tricosene	322	C23H46	018835-32-0	64
2			13-Tetradecen-1-ol acetate	254	C16H30O2	056221-91-1	64
3			1-Docosene	308	C22H44	001599-67-3	58
4			1-Hexacosanol	382	C26H54O	000506-52-5	58
5			Cycloeicosane	280	C20H40	000296-56-0	58



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060815\
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Acq On : 9 Jun 2015 4:18
Operator : TP/IZ
Sample : G2450-34
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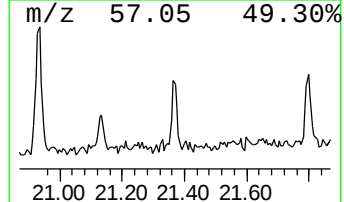
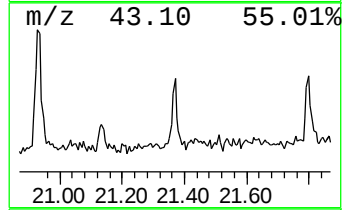
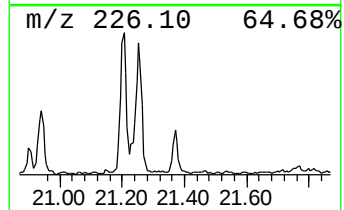
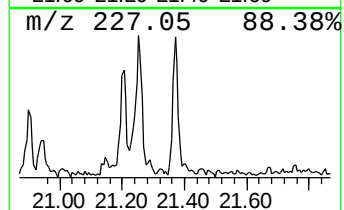
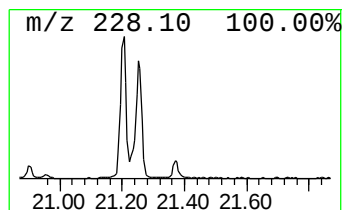
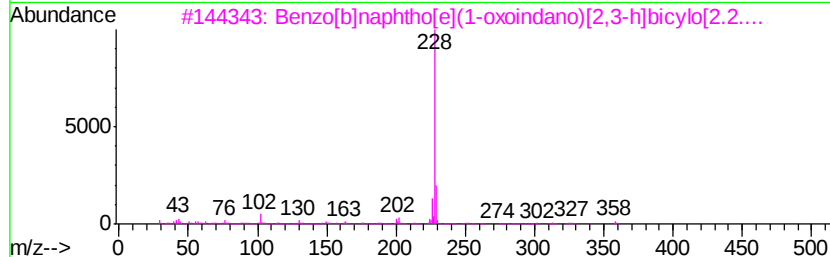
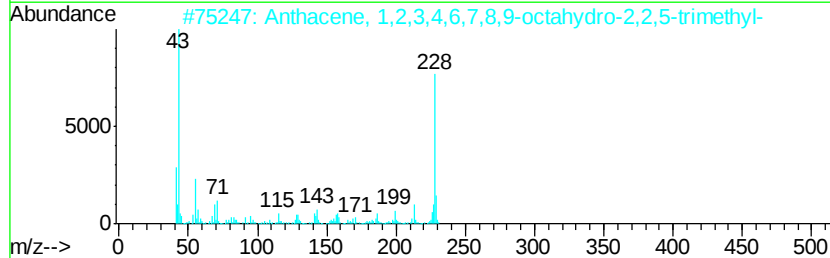
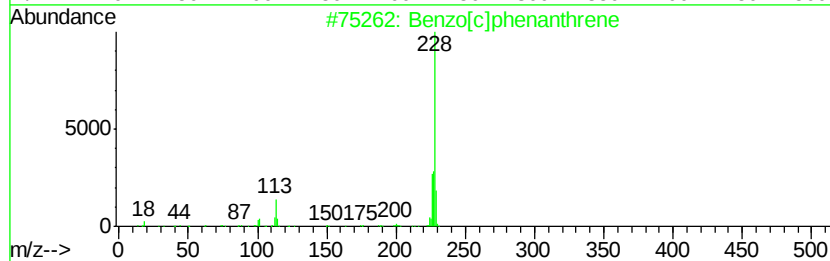
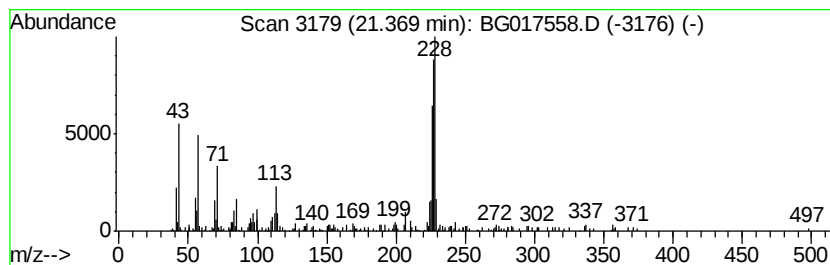
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ClientSampleId :
AVE-E-NBS13.2(3)

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Benzo[c]phenanthrene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.37	2.38 ng	93813	Chrysene-d12	21.22
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzo[c]phenanthrene	228 C18H12	000195-19-7 55
2		Anthracene, 1,2,3,4,6,7,8,9-octah...	228 C17H24	1000214-19-9 30
3		Benzo[b]naphtho[e](1-oxoindano)[...]	358 C27H18O	1000193-97-9 30
4		Triphenylene	228 C18H12	000217-59-4 30
5		Chrysene	228 C18H12	000218-01-9 27



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060815\
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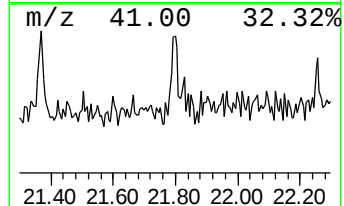
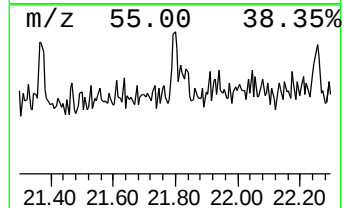
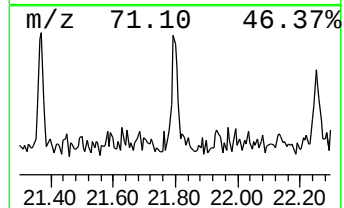
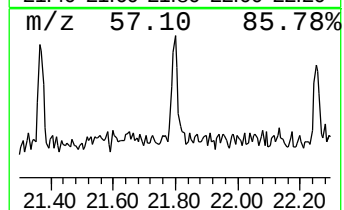
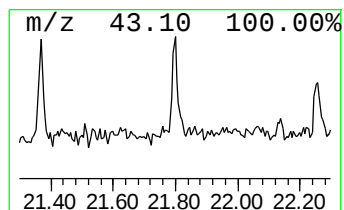
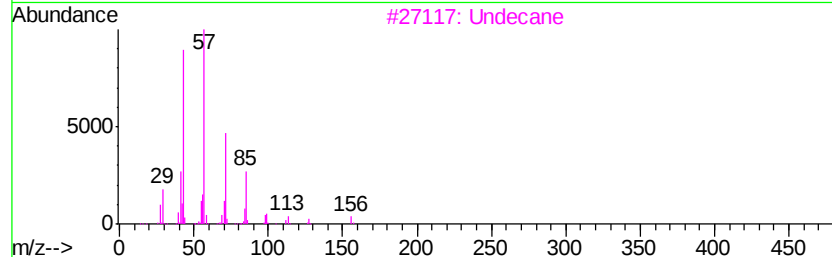
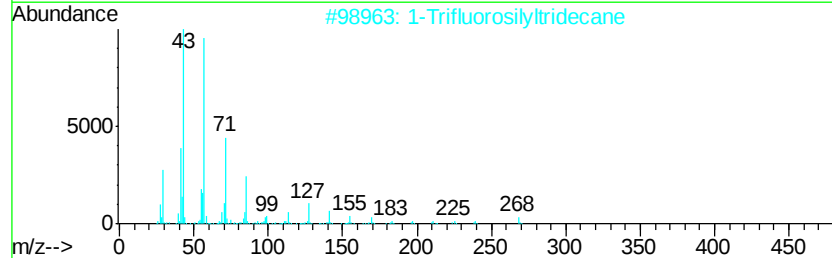
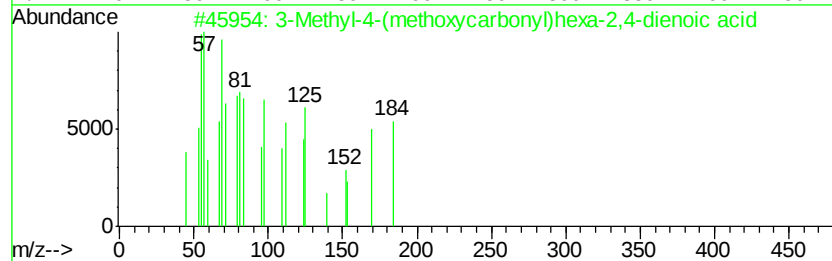
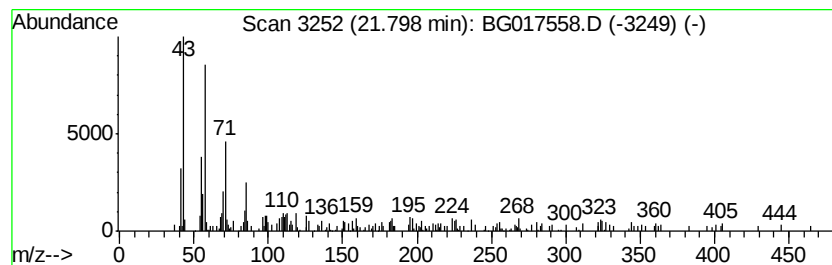
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AVE-E-NBS13.2(3)

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 3-Methyl-4-(methoxycarbonyl... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.		
21.80	2.21 ng	87078	Chrysene-d12	21.22		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Methyl-4-(methoxycarbonyl)hexa...	184	C9H12O4	1000104-10-8	86	
2	1-Trifluorosilyltridecane	268	C13H27F3Si	1000217-08-2	62	
3	Undecane	156	C11H24	001120-21-4	62	
4	Hexadecane, 1-chloro-	260	C16H33Cl	004860-03-1	62	
5	Tridecane, 6-methyl-	198	C14H30	013287-21-3	58	



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060815\
Data File : BG017558.D
Acq On : 9 Jun 2015 4:18
Operator : TP/IZ
Sample : G2450-34
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
AVE-E-NBS13.2(3)

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	2.74	10.1	ng	76871	1	7.60	152990	20.0
2-Pentanone, 4-hy...	4.71	4.3	ng	32765	1	7.60	152990	20.0
unknown7.14	7.14	92.4	ng	706983	1	7.60	152990	20.0
n-Hexadecanoic acid	17.93	6.7	ng	115680	4	17.02	348126	20.0
Benzaldehyde, 3,5...	18.09	3.6	ng	61997	4	17.02	348126	20.0
Hexadecanoic acid...	19.36	2.4	ng	92702	5	21.22	789461	20.0
Pyrene, 1-methyl-	19.95	2.8	ng	109329	5	21.22	789461	20.0
Octadecanoic acid...	20.41	2.2	ng	86671	5	21.22	789461	20.0
1-Tricosene	20.93	4.2	ng	165685	5	21.22	789461	20.0
Benzo[c]phenanthrene	21.37	2.4	ng	93813	5	21.22	789461	20.0
3-Methyl-4-(metho...	21.80	2.2	ng	87078	5	21.22	789461	20.0