

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060415\
 Data File : BG017451.D
 Acq On : 4 Jun 2015 20:43
 Operator : TP/IZ
 Sample : G2501-04
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MW-4-6-2-15

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	2.776	10	15	28	rVB	245871	315853	16.97%	1.113%
2	5.126	411	415	421	rVB	42503	58636	3.15%	0.207%
3	5.220	426	431	434	rVV	163044	238615	12.82%	0.841%
4	5.255	434	437	446	rVB	161701	243217	13.07%	0.857%
5	6.712	679	685	690	rBV	120471	170258	9.15%	0.600%
6	6.830	699	705	716	rVB2	118128	264937	14.24%	0.934%
7	7.012	729	736	742	rBV	64162	92936	4.99%	0.328%
8	7.165	756	762	771	rBV	316127	480665	25.83%	1.694%
9	7.271	773	780	789	rVB	274087	402976	21.66%	1.420%
10	7.623	832	840	845	rBV	66936	108555	5.83%	0.383%
11	7.694	845	852	856	rVV	158063	247438	13.30%	0.872%
12	7.741	856	860	869	rVB	150161	234069	12.58%	0.825%
13	7.946	883	895	900	rBV	281711	461436	24.80%	1.626%
14	7.993	900	903	909	rVB	57800	79877	4.29%	0.282%
15	8.169	929	933	938	rVB	26161	38372	2.06%	0.135%
16	8.263	944	949	954	rBV	47510	68432	3.68%	0.241%
17	8.428	970	977	985	rVB	23849	37560	2.02%	0.132%
18	8.575	997	1002	1006	rBV	30416	44018	2.37%	0.155%
19	8.628	1006	1011	1018	rVB	33522	49852	2.68%	0.176%
20	8.728	1023	1028	1032	rVV2	51451	78734	4.23%	0.277%
21	8.786	1032	1038	1050	rVB2	239848	438290	23.55%	1.545%
22	9.063	1079	1085	1091	rVV2	24848	37899	2.04%	0.134%
23	9.251	1112	1117	1122	rVV	25400	36766	1.98%	0.130%
24	9.309	1122	1127	1134	rVB	48488	71366	3.84%	0.252%
25	9.668	1184	1188	1193	rVB	26320	38931	2.09%	0.137%
26	9.820	1202	1214	1222	rBV3	103585	203274	10.92%	0.716%
27	10.050	1248	1253	1259	rVB2	52169	83064	4.46%	0.293%
28	10.414	1309	1315	1320	rVV2	91385	174543	9.38%	0.615%
29	10.467	1320	1324	1331	rVV	70529	120455	6.47%	0.425%
30	10.978	1404	1411	1420	rBV6	20254	52815	2.84%	0.186%
31	11.612	1514	1519	1524	rBV2	37241	59367	3.19%	0.209%
32	11.977	1576	1581	1586	rVB2	30942	47579	2.56%	0.168%
33	12.088	1591	1600	1608	rVB	128204	213485	11.47%	0.752%
34	12.312	1632	1638	1643	rBV	98179	155960	8.38%	0.550%

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

35	12.905	1733	1739	1746	rVB	620914	880119	47.30%	3.102%
36	13.440	1824	1830	1843	rVV5	31825	92383	4.96%	0.326%
37	13.604	1851	1858	1862	rVV3	42268	110091	5.92%	0.388%
38	13.657	1865	1867	1877	rVB2	39012	54895	2.95%	0.193%
39	14.004	1920	1926	1933	rVB2	45320	70756	3.80%	0.249%
40	14.198	1954	1959	1964	rBV5	18337	34056	1.83%	0.120%
41	14.292	1970	1975	1979	rVB2	123764	183276	9.85%	0.646%
42	14.703	2037	2045	2049	rBV9	14901	35488	1.91%	0.125%
43	15.238	2134	2136	2149	rVB7	20171	46162	2.48%	0.163%
44	15.338	2149	2153	2158	rBV6	17553	37418	2.01%	0.132%
45	15.467	2166	2175	2184	rVB4	25397	57440	3.09%	0.202%
46	15.743	2217	2222	2224	rBV	138936	176652	9.49%	0.623%
47	15.784	2224	2229	2236	rVV2	650744	961513	51.67%	3.389%
48	15.855	2236	2241	2248	rVB4	38465	72923	3.92%	0.257%
49	16.043	2264	2273	2278	rBV	214493	293648	15.78%	1.035%
50	16.137	2281	2289	2292	rVV	986100	1349164	72.50%	4.755%
51	16.195	2292	2299	2305	rVV2	883012	1860819	100.00%	6.558%
52	16.254	2305	2309	2313	rVV2	940544	1345414	72.30%	4.742%
53	16.301	2313	2317	2321	rVV	559147	746199	40.10%	2.630%
54	16.372	2321	2329	2333	rVV	437156	802456	43.12%	2.828%
55	16.454	2337	2343	2349	rVV4	607559	1076365	57.84%	3.794%
56	16.524	2349	2355	2358	rVV	984448	1341096	72.07%	4.727%
57	16.554	2358	2360	2363	rVV	215537	288114	15.48%	1.015%
58	16.595	2363	2367	2373	rVV	434035	598718	32.17%	2.110%
59	16.801	2399	2402	2407	rBV	24401	35611	1.91%	0.126%
60	17.041	2436	2443	2449	rBV	203991	289740	15.57%	1.021%
61	17.229	2470	2475	2483	rVB	46660	69254	3.72%	0.244%
62	17.741	2553	2562	2568	rBV2	24663	50908	2.74%	0.179%
63	17.964	2593	2600	2606	rBV	169513	270533	14.54%	0.953%
64	18.058	2611	2616	2620	rVV	291625	423496	22.76%	1.493%
65	18.105	2620	2624	2628	rVV2	339329	608240	32.69%	2.144%
66	18.140	2628	2630	2634	rVV	264476	366625	19.70%	1.292%
67	18.193	2634	2639	2644	rVV2	263553	547791	29.44%	1.931%
68	18.240	2644	2647	2650	rVV	66153	100793	5.42%	0.355%
69	18.281	2650	2654	2657	rVV2	316401	471218	25.32%	1.661%
70	18.316	2657	2660	2664	rVV2	388674	539362	28.99%	1.901%
71	18.358	2664	2667	2671	rVV	225112	294623	15.83%	1.038%

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72	18.405	2671	2675	2683	rVV2	366613	646270	34.73%	2.278%
73	18.481	2683	2688	2695	rVB	233221	365646	19.65%	1.289%
74	18.640	2708	2715	2724	rBV4	61146	124749	6.70%	0.440%
75	18.969	2767	2771	2777	rBV2	47823	66949	3.60%	0.236%
76	19.086	2786	2791	2792	rBV3	27940	34693	1.86%	0.122%
77	19.145	2797	2801	2807	rVV3	120839	214039	11.50%	0.754%
78	19.204	2807	2811	2813	rVV2	45923	63640	3.42%	0.224%
79	19.233	2813	2816	2821	rVV6	34096	61276	3.29%	0.216%
80	19.298	2821	2827	2833	rVB4	43191	76090	4.09%	0.268%
81	19.679	2886	2892	2896	rBV	1436600	1784497	95.90%	6.289%
82	19.715	2896	2898	2904	rVB	41083	47399	2.55%	0.167%
83	19.820	2909	2916	2922	rBV2	138368	344505	18.51%	1.214%
84	19.879	2922	2926	2929	rVV	130774	196183	10.54%	0.691%
85	19.915	2929	2932	2937	rVV2	77710	116555	6.26%	0.411%
86	19.967	2937	2941	2943	rVV	76723	106390	5.72%	0.375%
87	20.003	2943	2947	2950	rVV5	119057	206863	11.12%	0.729%
88	20.032	2950	2952	2956	rVV2	61148	81083	4.36%	0.286%
89	20.079	2956	2960	2968	rVV	177895	267258	14.36%	0.942%
90	20.144	2968	2971	2982	rVB	60325	107497	5.78%	0.379%
91	20.943	3104	3107	3114	rBV7	22612	35183	1.89%	0.124%
92	21.095	3129	3133	3139	rBV2	48155	61741	3.32%	0.218%
93	21.189	3145	3149	3151	rBV3	31085	39368	2.12%	0.139%
94	21.231	3151	3156	3162	rVV	309375	434073	23.33%	1.530%
95	21.283	3162	3165	3167	rVV4	41382	53369	2.87%	0.188%
96	21.313	3167	3170	3173	rVV	80645	117079	6.29%	0.413%
97	21.366	3173	3179	3184	rVV3	114002	258333	13.88%	0.910%
98	21.419	3184	3188	3191	rVV	29342	45093	2.42%	0.159%
99	21.477	3192	3198	3203	rVB2	31885	51707	2.78%	0.182%
100	23.463	3529	3536	3542	rBV	175763	332056	17.84%	1.170%

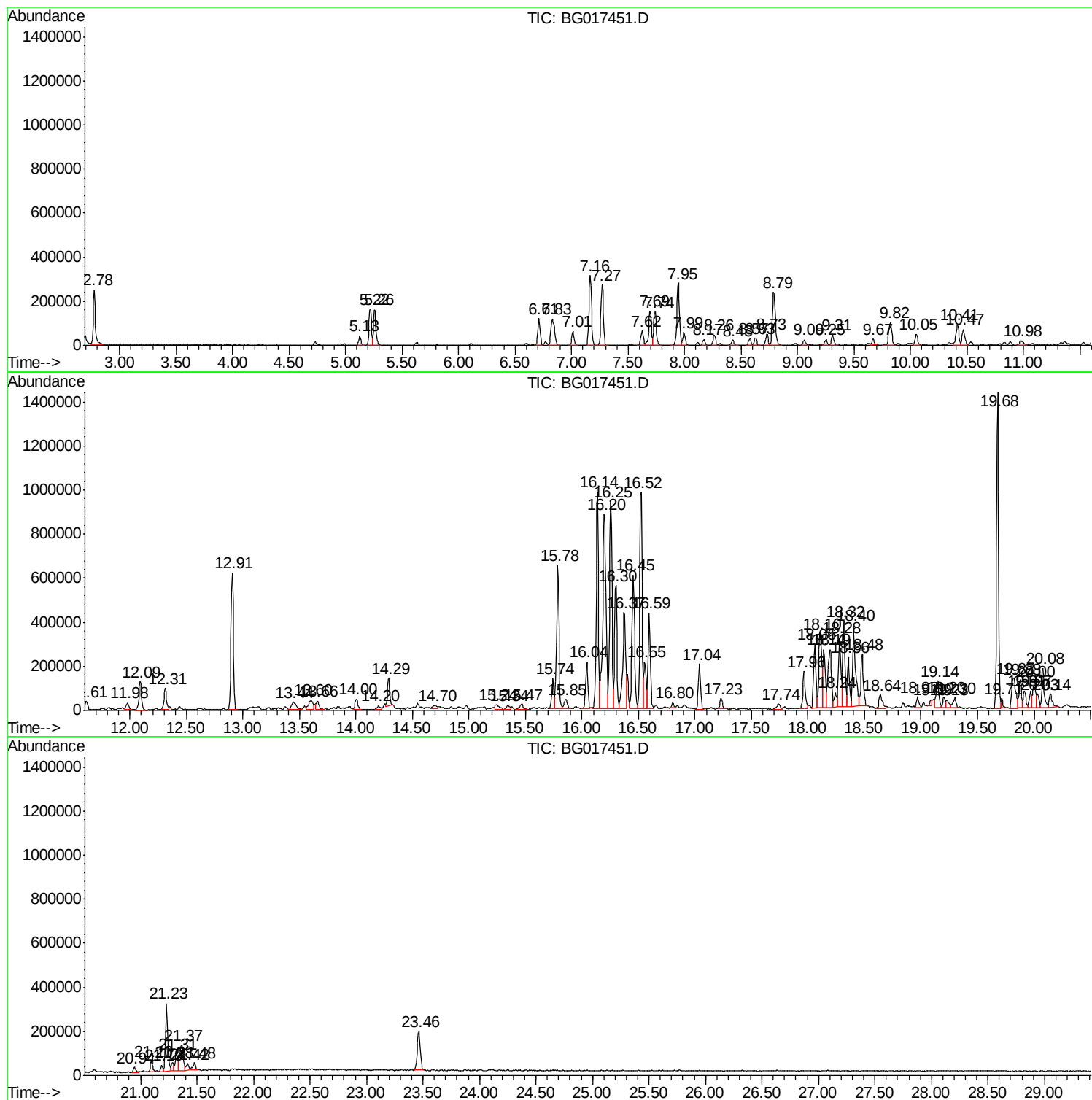
Sum of corrected areas: 28373173

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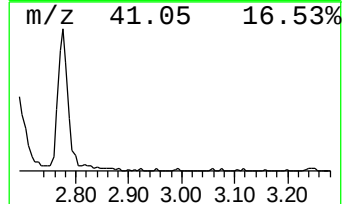
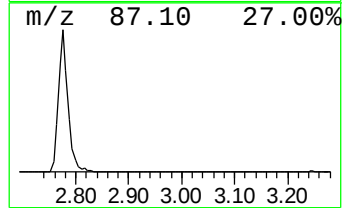
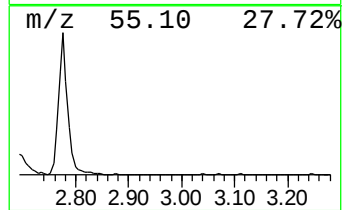
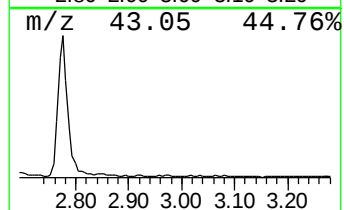
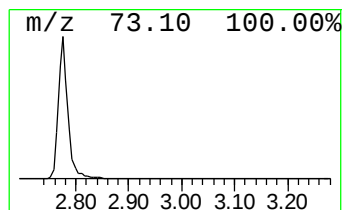
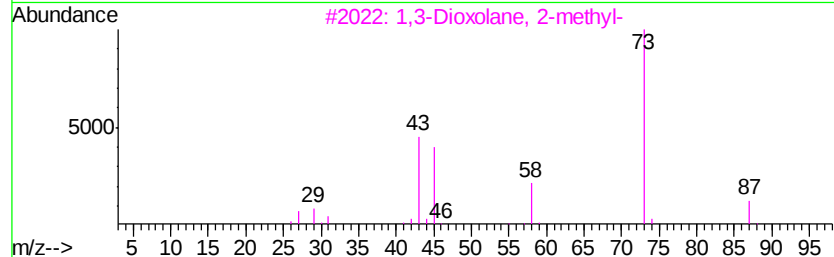
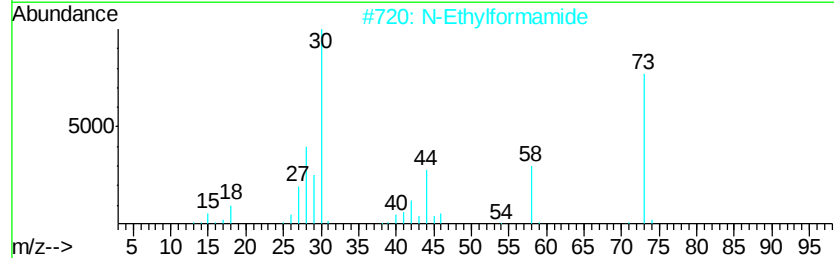
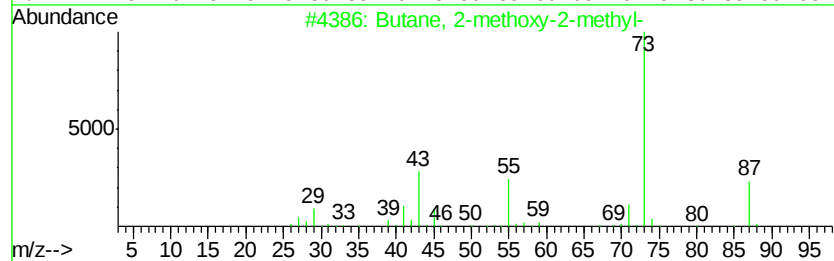
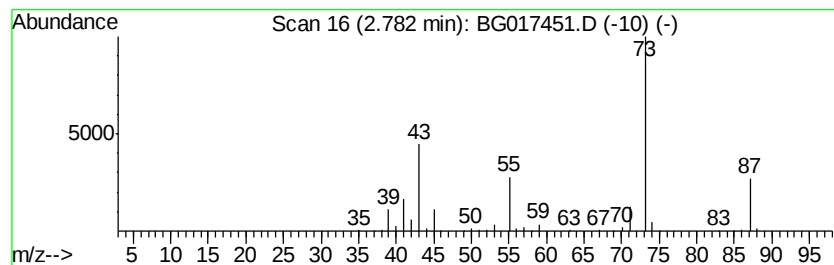
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.78	58.19 ng	315853	1,4-Dichlorobenzene-d4	7.62

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			N-Ethylformamide	73	C3H7NO	000627-45-2	9
3			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9
4			Silane, tetramethyl-	88	C4H12Si	000075-76-3	9
5			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	9



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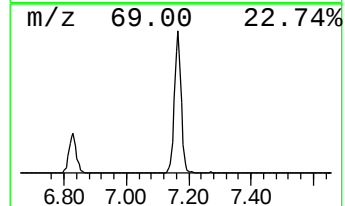
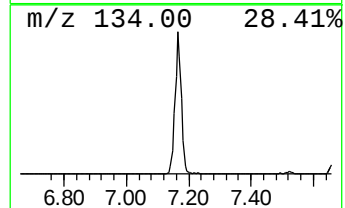
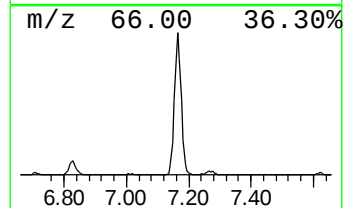
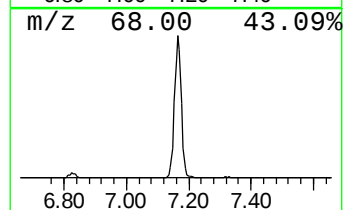
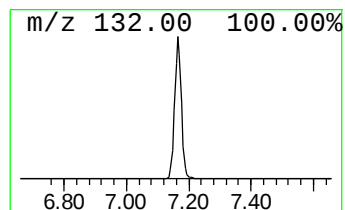
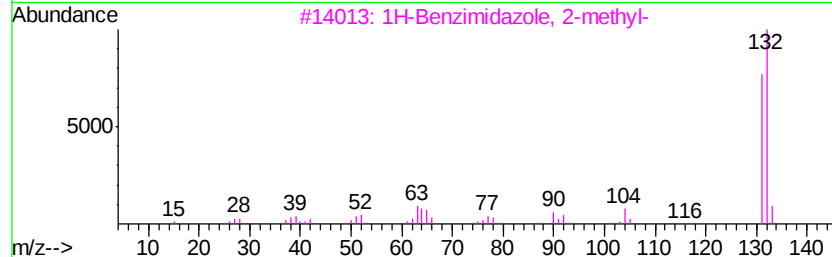
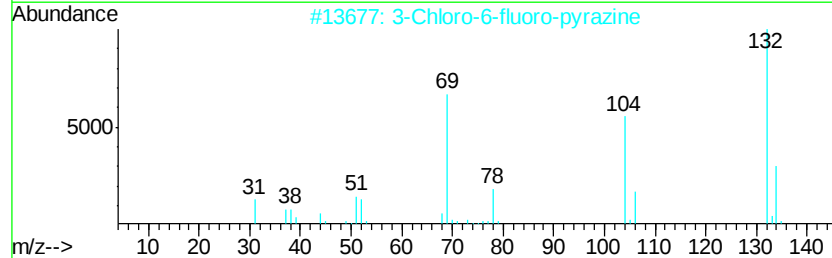
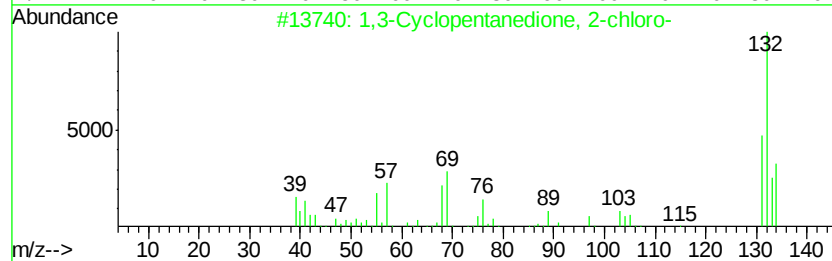
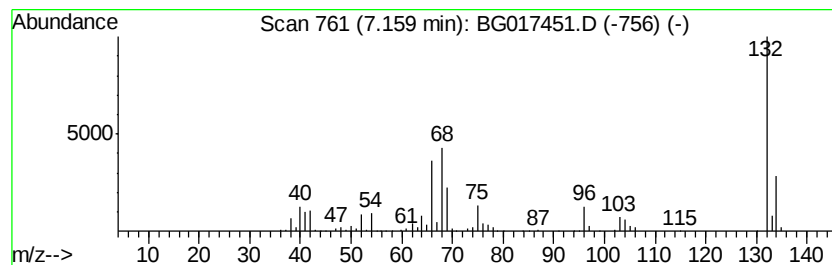
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 unknown7.16 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.16	88.56 ng	480665	1,4-Dichlorobenzene-d4	7.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	38
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4		Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	22
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16



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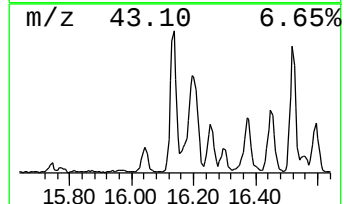
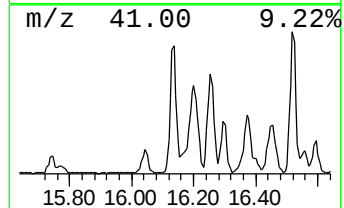
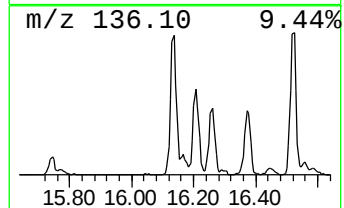
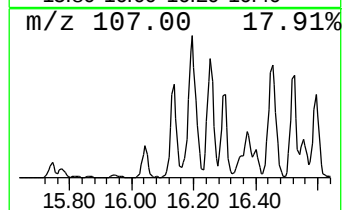
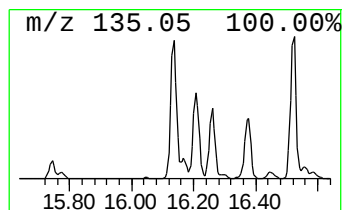
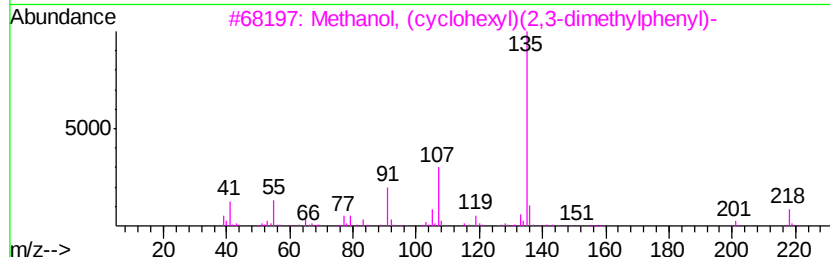
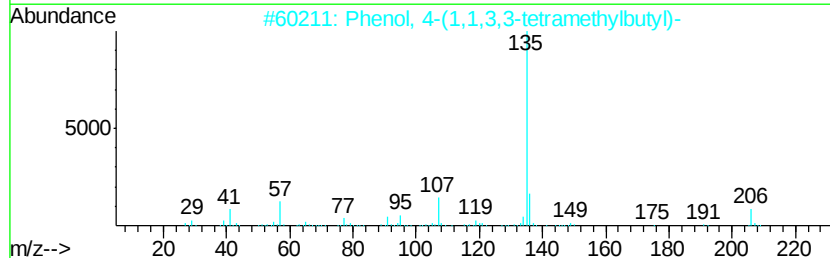
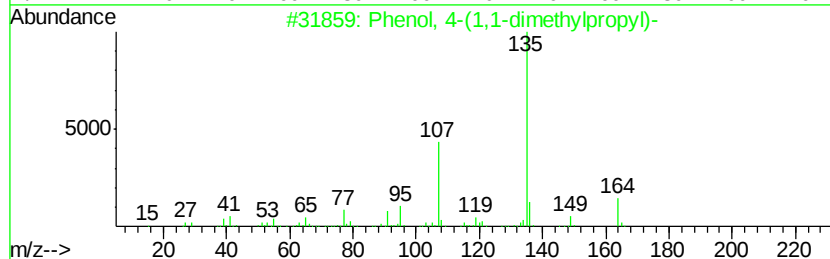
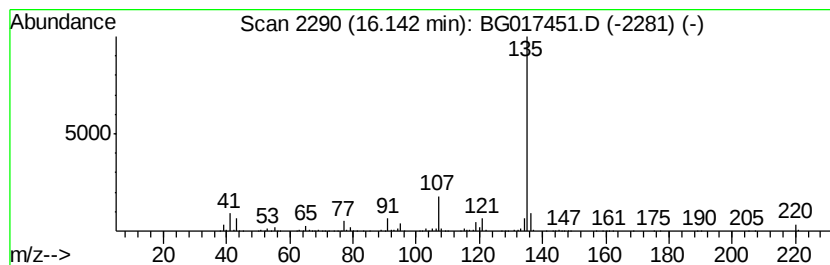
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TIC Integration Parameters: LSCINT.P

Peak Number 5 Phenol, 4-(1,1-dimethylprop... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.14	93.13 ng	1349160	Phenanthrene-d10	17.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	78
2		Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	72
3		Methanol, (cyclohexyl)(2,3-dimet...	218	C15H22O	349498-47-1	72
4		Hexestrol	270	C18H22O2	005635-50-7	64
5		((1,2-Diethylethylene)bis(p-phen...	354	C22H26O4	004547-76-6	64



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060415\
Data File : BG017451.D
Acq On : 4 Jun 2015 20:43
Operator : TP/IZ
Sample : G2501-04
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
MW-4-6-2-15

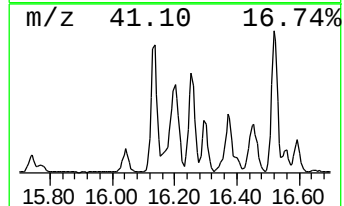
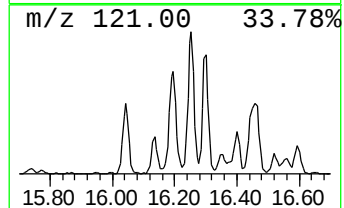
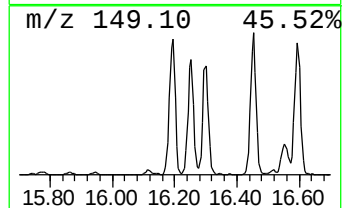
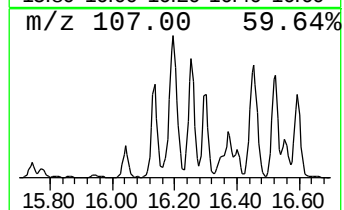
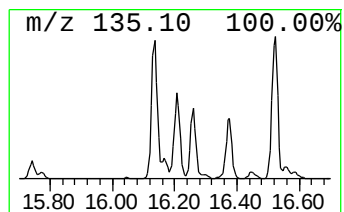
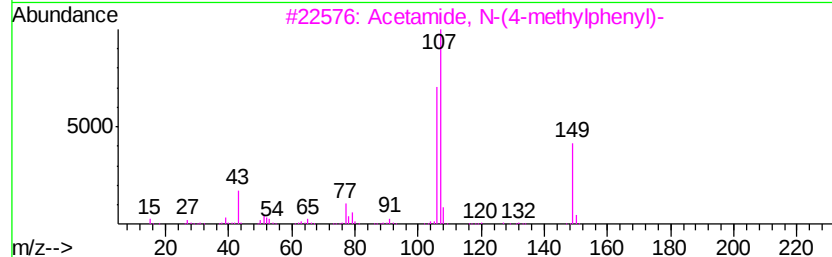
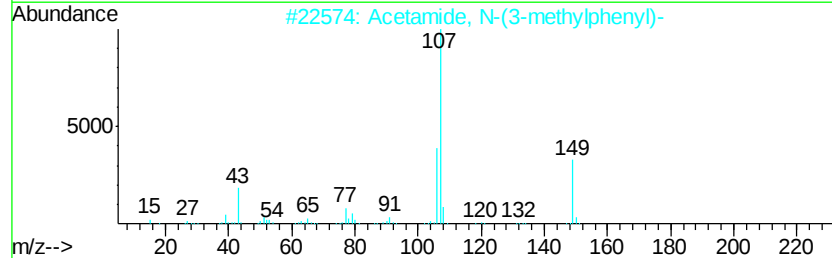
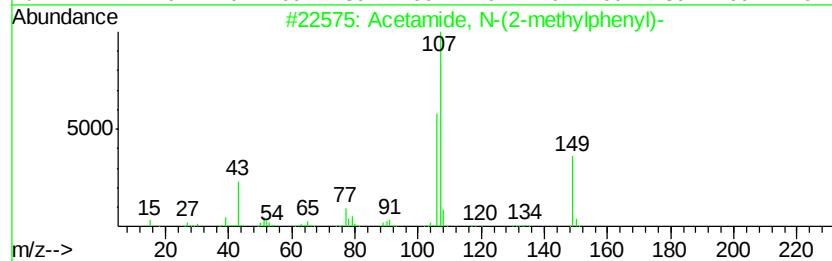
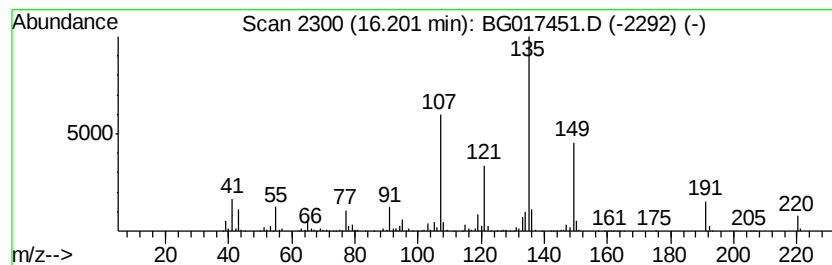
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 unknown16.20 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.20	128.45 ng	1860820	Phenanthrene-d10	17.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Acetamide, N-(2-methylphenyl)-	149	C9H11NO	000120-66-1	47
2		Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	35
3		Acetamide, N-(4-methylphenyl)-	149	C9H11NO	000103-89-9	35
4		4-Methyl-2-tert-octylphenol	220	C15H24O	004979-46-8	30
5		2-Methyl-2-(4'-methoxyphenyl)pen...	206	C13H18O2	185700-49-6	27



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060415\
Data File : BG017451.D
Acq On : 4 Jun 2015 20:43
Operator : TP/IZ
Sample : G2501-04
Misc :
ALS Vial : 13 Sample Multiplier: 1

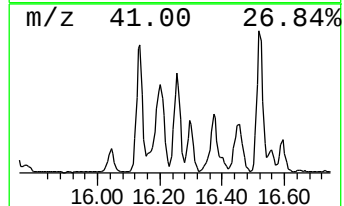
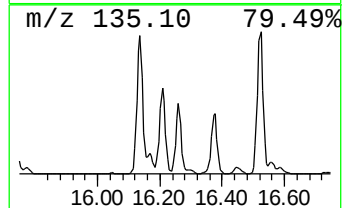
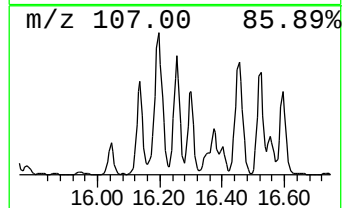
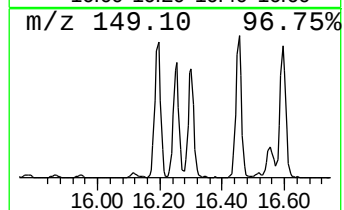
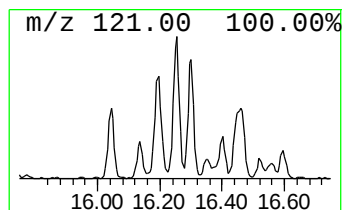
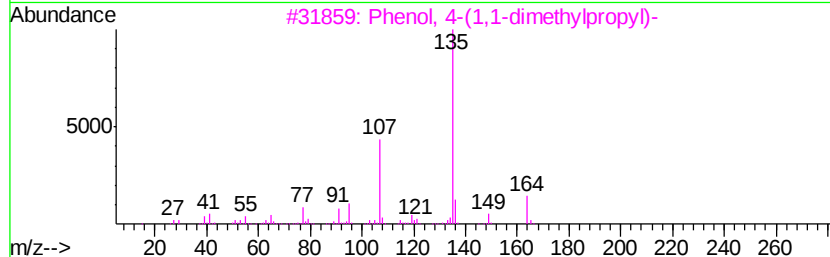
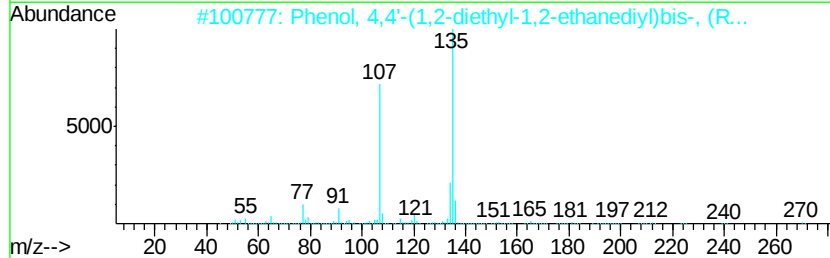
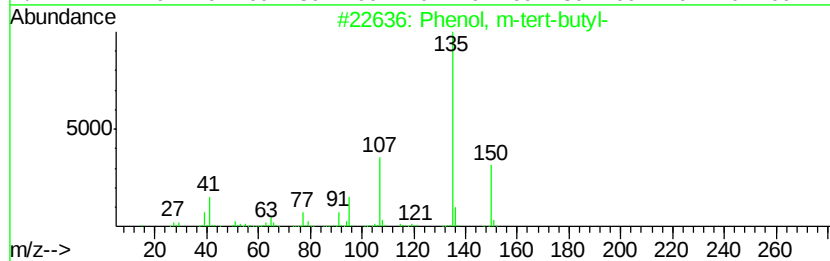
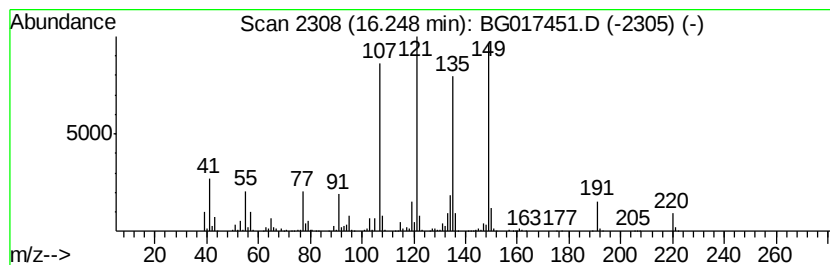
Instrument :
BNA_G
ClientSampleId :
MW-4-6-2-15

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 7 Phenol, m-tert-butyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.25	92.87 ng	1345410	Phenanthrene-d10	17.04
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Phenol, m-tert-butyl-	150 C10H14O	000585-34-2	60
2	Phenol, 4,4'-(1,2-diethyl-1,2-et...	270 C18H22O2	000084-16-2	50
3	Phenol, 4-(1,1-dimethylpropyl)-	164 C11H16O	000080-46-6	50
4	Hexestrol	270 C18H22O2	005635-50-7	50
5	1,5,6,7-Tetrahydro-4-indolone	135 C8H9NO	013754-86-4	47



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060415\
Data File : BG017451.D
Acq On : 4 Jun 2015 20:43
Operator : TP/IZ
Sample : G2501-04
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
MW-4-6-2-15

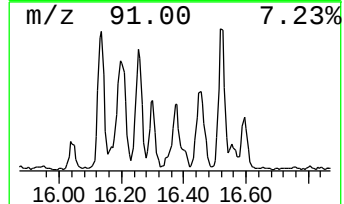
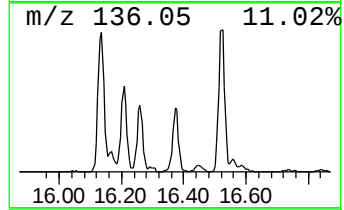
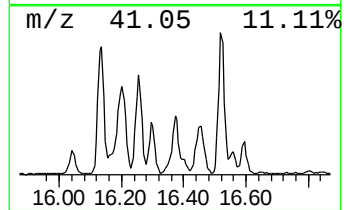
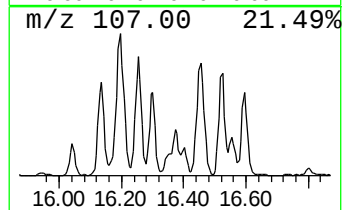
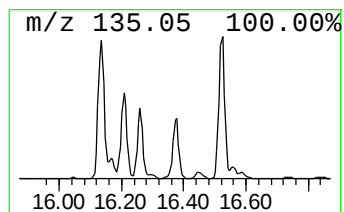
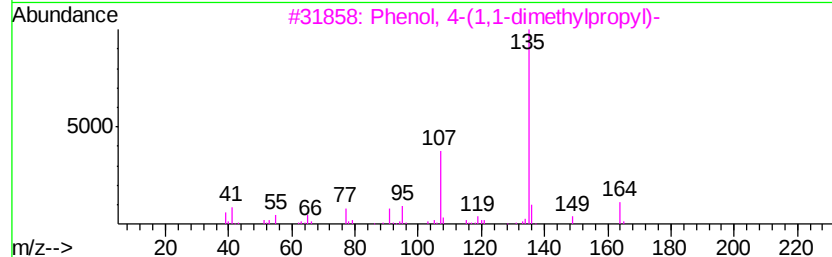
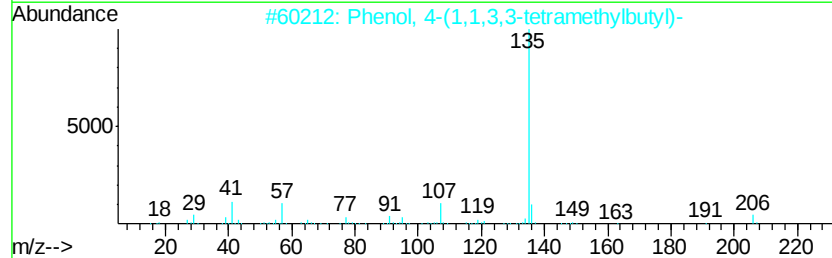
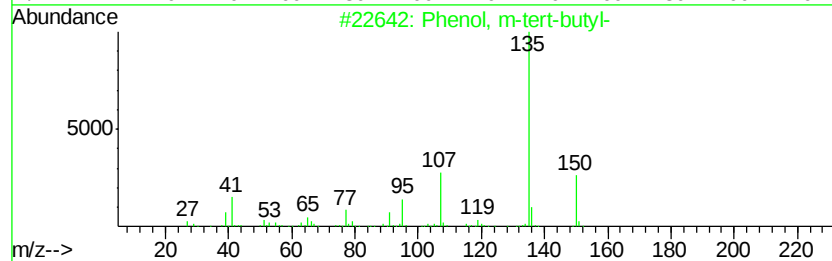
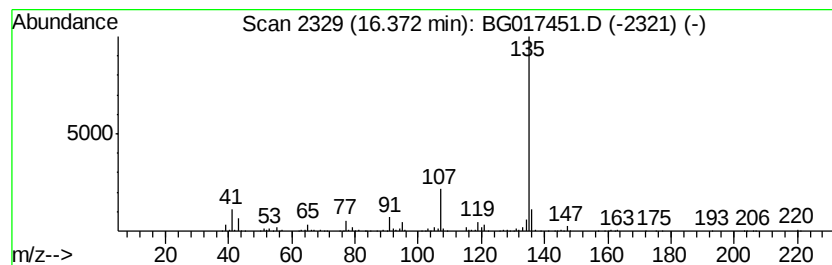
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Phenol, 4-(1,1,3,3-tetramet... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.37	55.39 ng	802456	Phenanthrene-d10	17.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	78
2		Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	72
3		Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	72
4		Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	64
5		2-Fluoro-4-cyanotoluene	135	C8H6FN	170572-49-3	64



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060415\
Data File : BG017451.D
Acq On : 4 Jun 2015 20:43
Operator : TP/IZ
Sample : G2501-04
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
MW-4-6-2-15

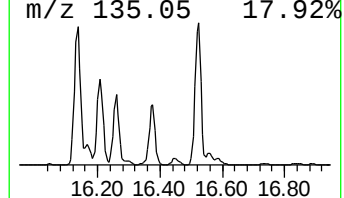
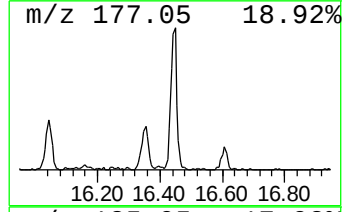
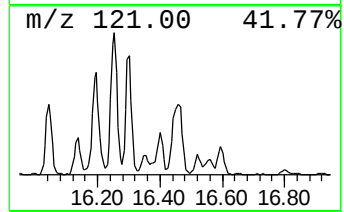
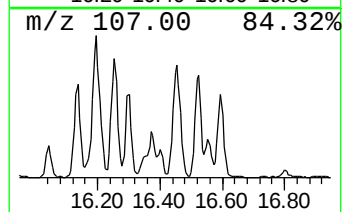
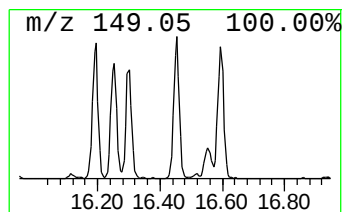
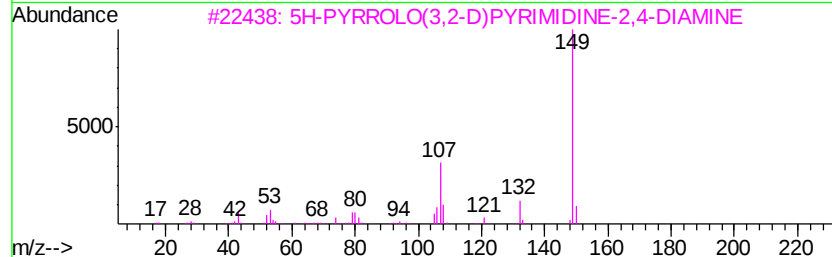
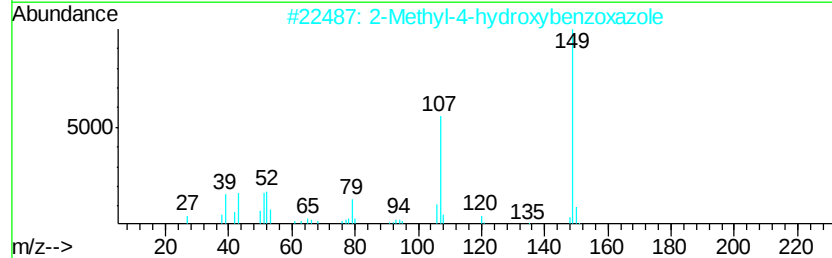
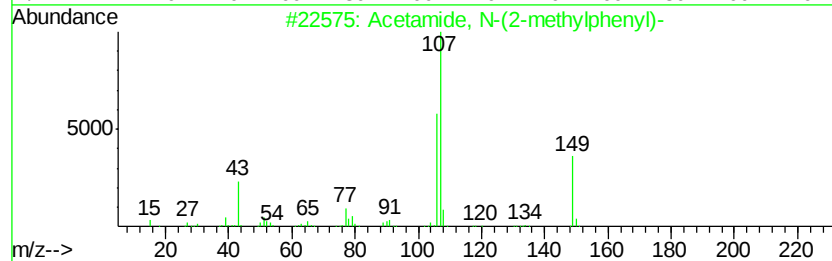
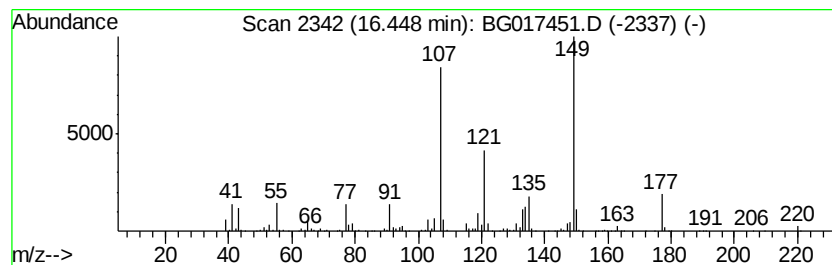
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Acetamide, N-(2-methylphenyl)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.45	74.30 ng	1076370	Phenanthrene-d10	17.04

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Acetamide, N-(2-methylphenyl)-	149	C9H11NO	000120-66-1	64
2		2-Methyl-4-hydroxybenzoxazole	149	C8H7NO2	051110-60-2	64
3		5H-PYRROLO(3,2-D)PYRIMIDINE-2,4-...	149	C6H7N5	1000244-21-4	64
4		Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	64
5		Acetamide, N-(4-methylphenyl)-	149	C9H11NO	000103-89-9	46



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060415\
Data File : BG017451.D
Acq On : 4 Jun 2015 20:43
Operator : TP/IZ
Sample : G2501-04
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
MW-4-6-2-15

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...	2.78	58.2	ng	315853	1	7.62	108555	20.0
unknown7.16	7.16	88.6	ng	480665	1	7.62	108555	20.0
Phenol, 4-(1,1-di...	16.14	93.1	ng	1349160	4	17.04	289740	20.0
unknown16.20	16.20	128.4	ng	1860820	4	17.04	289740	20.0
Phenol, m-tert-bu...	16.25	92.9	ng	1345410	4	17.04	289740	20.0
Phenol, 4-(1,1,3,...	16.37	55.4	ng	802456	4	17.04	289740	20.0
Acetamide, N-(2-m...	16.45	74.3	ng	1076370	4	17.04	289740	20.0