

Data Path : Z:\HPCHEM1\BNA G\DATA\BG032718\
 Data File : BG033374.D
 Acq On : 27 Mar 2018 22:43
 Operator : SJ/JU
 Sample : J2076-01
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EFF-180326-05

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON Filtering: 5
 Sampling : 1 Min Area: 3 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG031918.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.479	25	32	43	rBV	145470	321938	7.50%	1.868%
2	3.573	43	48	57	rVB	100712	158005	3.68%	0.917%
3	6.287	503	510	523	rBV	220445	472165	11.00%	2.739%
4	7.967	789	796	808	rBV	132491	336856	7.85%	1.954%
5	8.373	857	865	888	rBV	417088	944254	22.01%	5.478%
6	8.866	941	949	958	rBV2	106804	241139	5.62%	1.399%
7	9.201	998	1006	1022	rBV	634705	1298280	30.26%	7.532%
8	10.065	1143	1153	1175	rBV	486240	1103613	25.72%	6.403%
9	11.751	1432	1440	1451	rBV	157328	350223	8.16%	2.032%
10	14.113	1835	1842	1863	rBV	1630145	2812515	65.55%	16.318%
11	15.482	2066	2075	2087	rBV2	334116	593698	13.84%	3.445%
12	16.957	2319	2326	2341	rBV	951592	1678726	39.12%	9.740%
13	18.214	2532	2540	2553	rBV	496294	825594	19.24%	4.790%
14	20.735	2962	2969	2979	rBV	3134493	4290783	100.00%	24.895%
15	22.556	3272	3279	3290	rBV2	466269	924114	21.54%	5.362%
16	26.328	3909	3921	3934	rBV	253963	883928	20.60%	5.128%

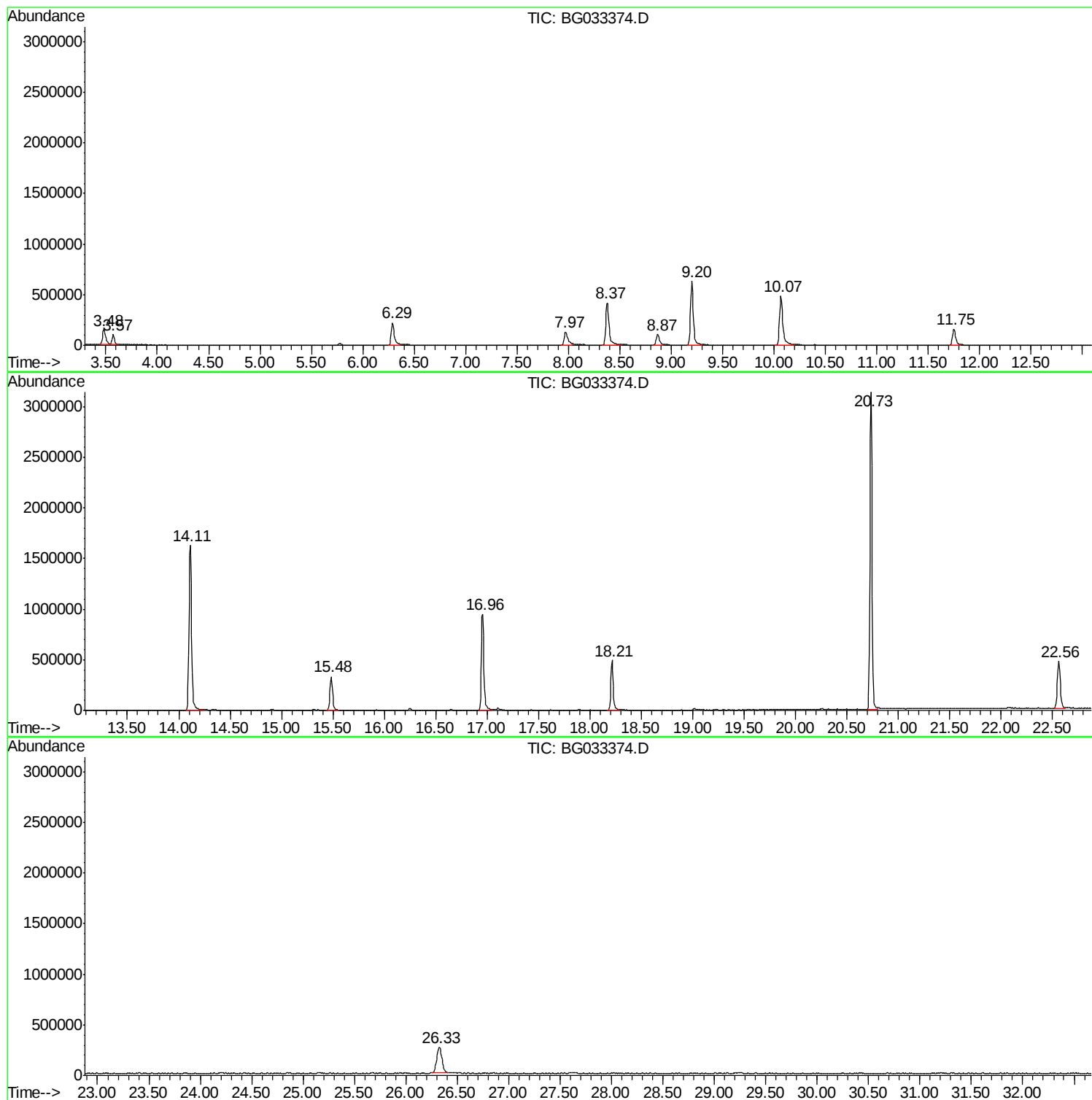
Sum of corrected areas: 17235831

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG031918.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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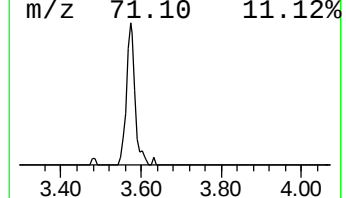
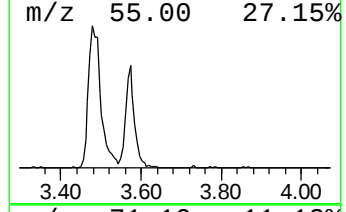
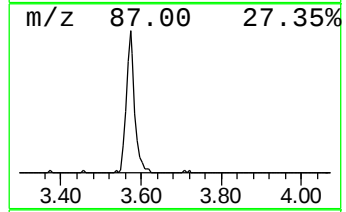
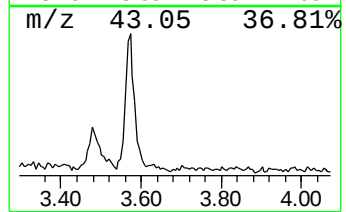
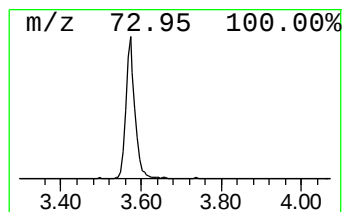
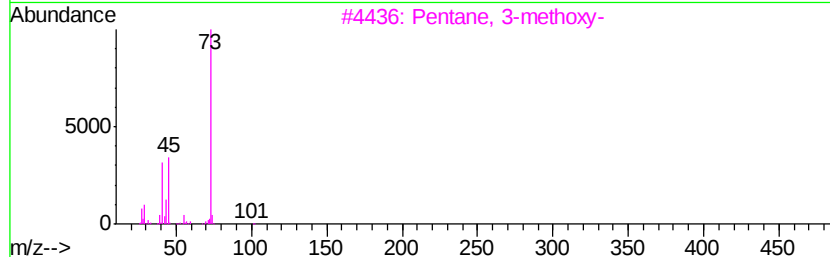
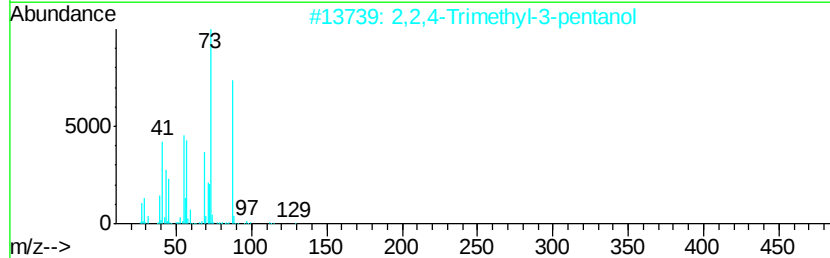
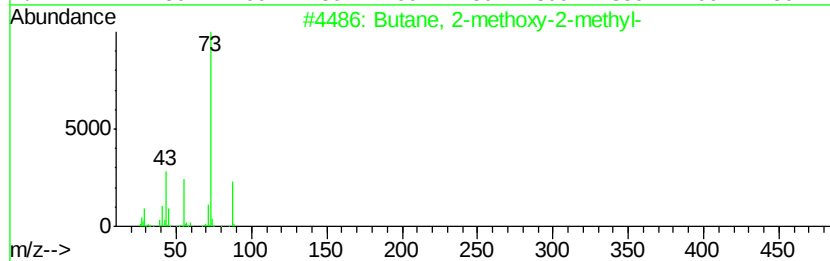
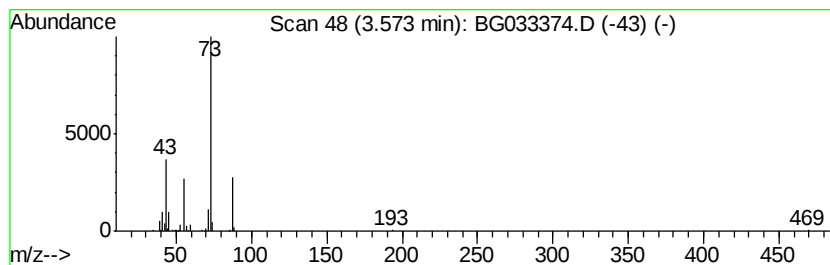
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Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.57	13.10 ng	158005	1,4-Dichlorobenzene-d4	8.87

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	10
4			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	10
5			N-Ethylformamide	73	C3H7NO	000627-45-2	9



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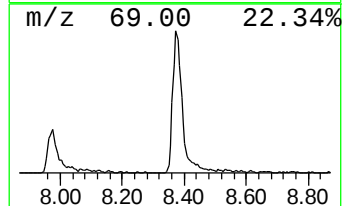
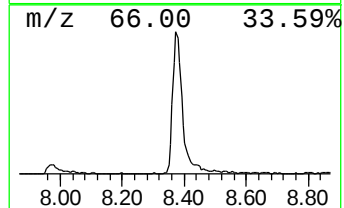
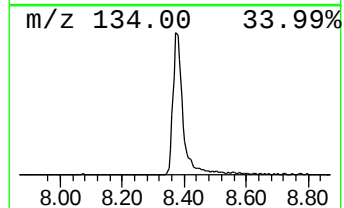
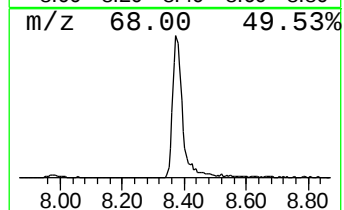
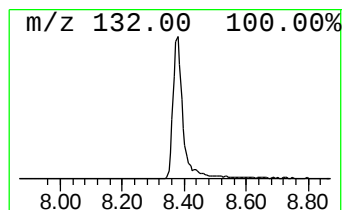
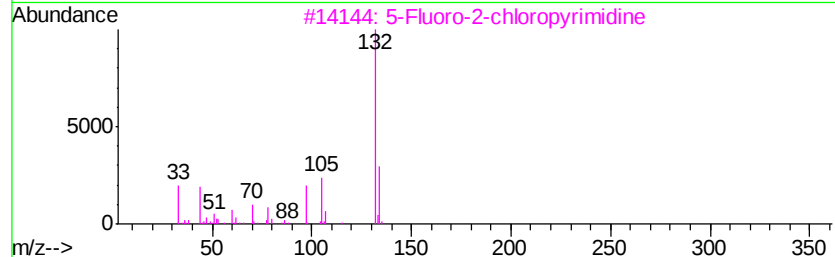
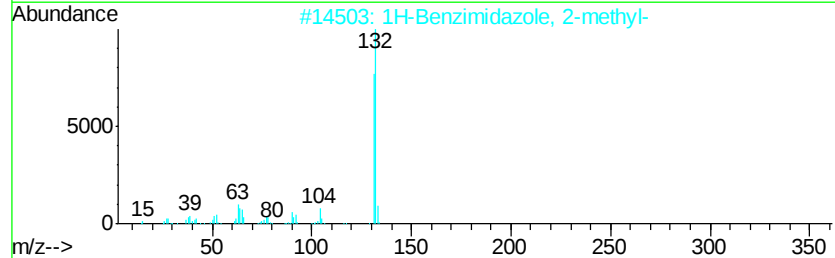
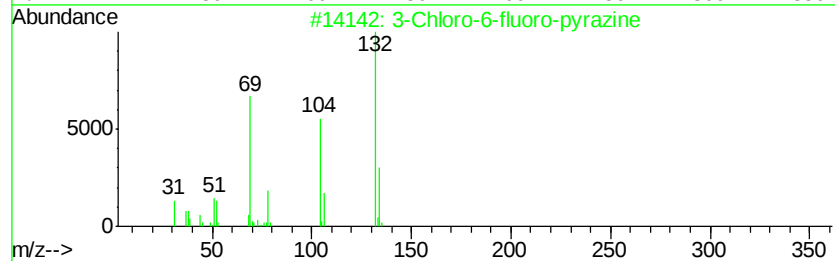
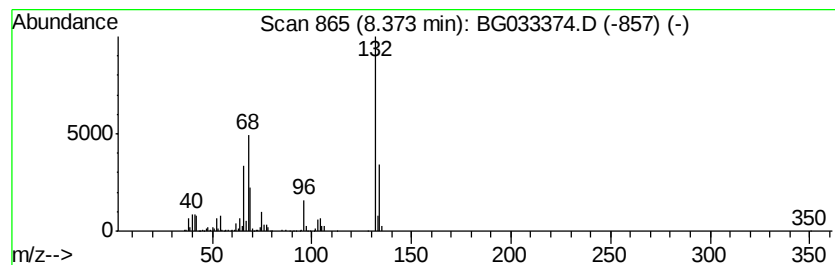
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Peak Number 3 unknown8.37 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.37	78.32 ng	944254	1,4-Dichlorobenzene-d4	8.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	3.57	13.1	ng	158005	1	8.87	241139	20.0
unknown8.37	8.37	78.3	ng	944254	1	8.87	241139	20.0