

Data Path : U:\HPCHEM1\BNA G\DATA\BG012718\
 Data File : BG032362.D
 Acq On : 27 Jan 2018 18:39
 Operator : SJ/JU
 Sample : PB106052BL
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB106052BL

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-BG011218.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	3.470	24	31	38	rBV	30777	56707	1.82%	0.434%
2	6.155	481	488	501	rBV	472296	814188	26.18%	6.230%
3	7.836	765	774	783	rBV	451681	830101	26.69%	6.351%
4	8.235	835	842	855	rVB	501174	910221	29.27%	6.964%
5	8.723	916	925	933	rBV	112475	207796	6.68%	1.590%
6	9.058	974	982	992	rBV	430652	784224	25.21%	6.000%
7	9.916	1120	1128	1138	rBV	271388	515627	16.58%	3.945%
8	11.602	1407	1415	1427	rBB	145783	273606	8.80%	2.093%
9	13.982	1813	1820	1830	rBV	1036774	1625169	52.25%	12.435%
10	15.356	2047	2054	2064	rBV2	248787	412111	13.25%	3.153%
11	16.831	2298	2305	2317	rBV	857443	1331430	42.81%	10.187%
12	18.088	2513	2519	2529	rVB2	364207	572539	18.41%	4.381%
13	20.627	2945	2951	2959	rBV	2378785	3110240	100.00%	23.797%
14	22.413	3247	3255	3267	rBV	433820	812962	26.14%	6.220%
15	26.091	3870	3881	3895	rBV2	248143	812904	26.14%	6.220%

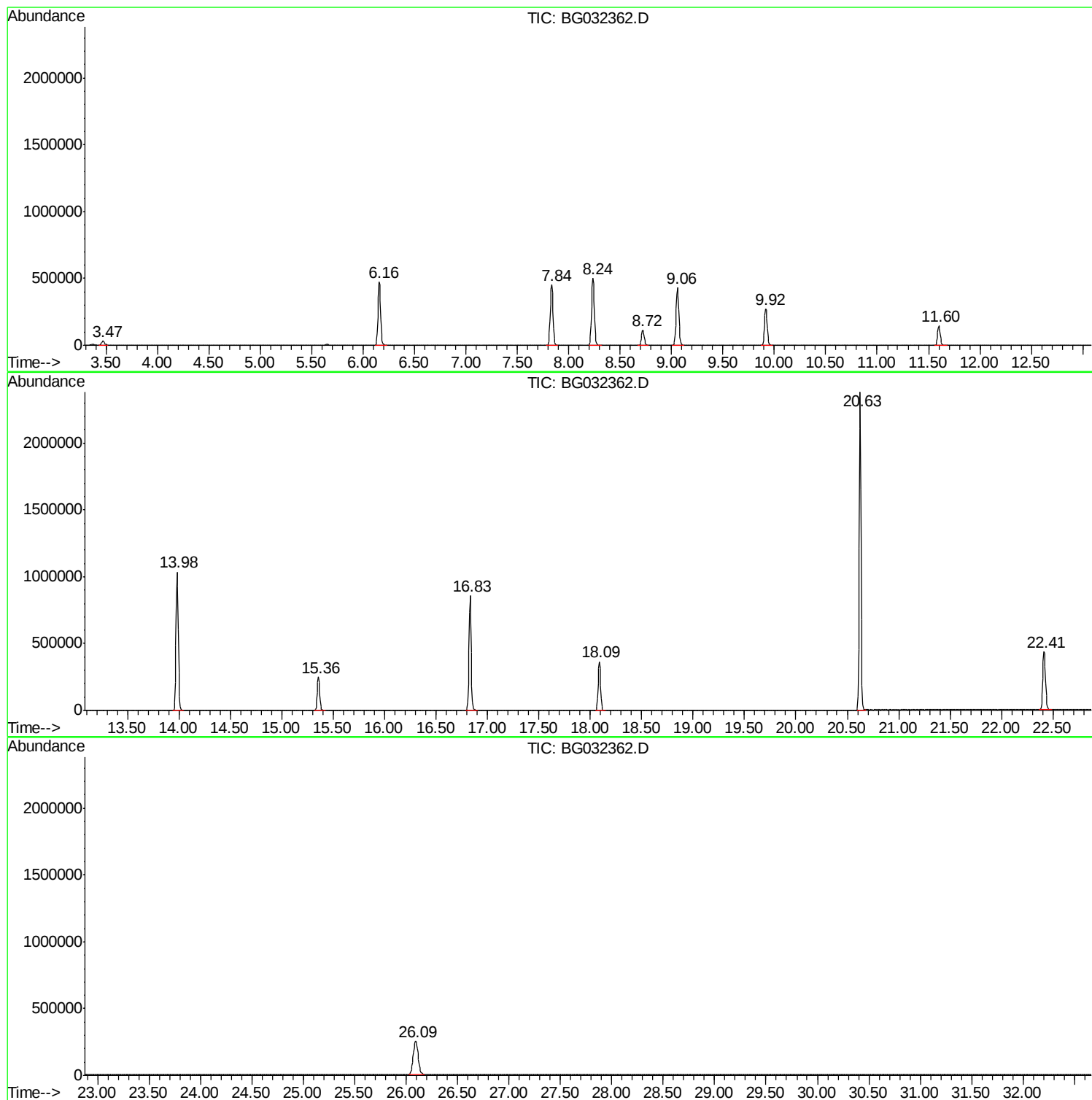
Sum of corrected areas: 13069825

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG011218.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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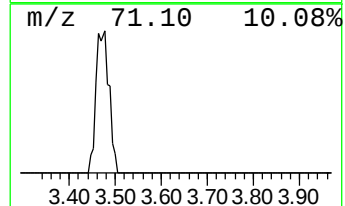
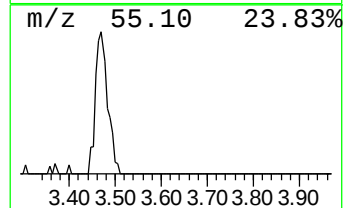
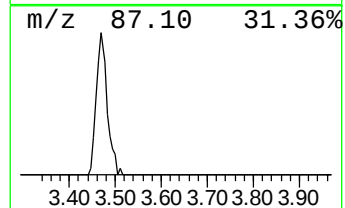
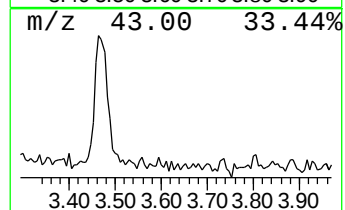
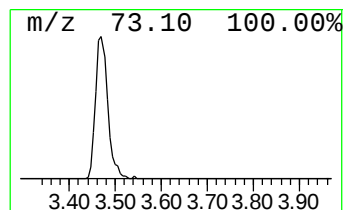
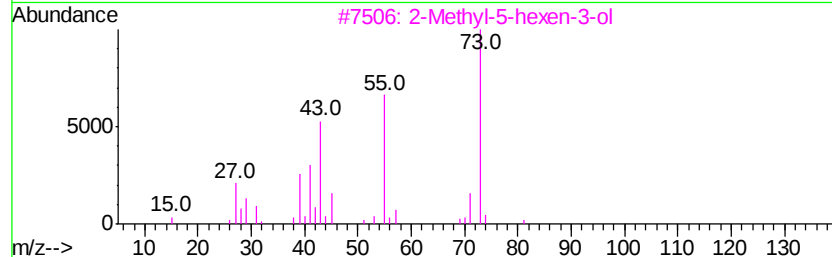
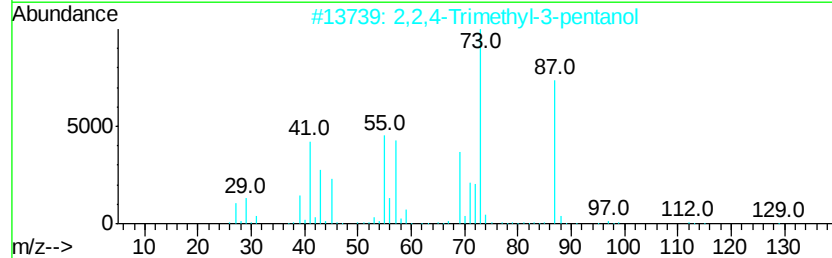
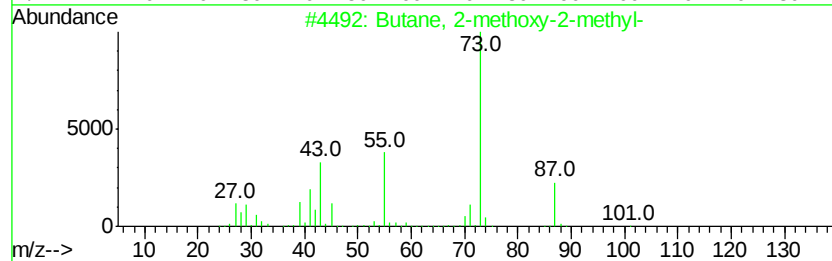
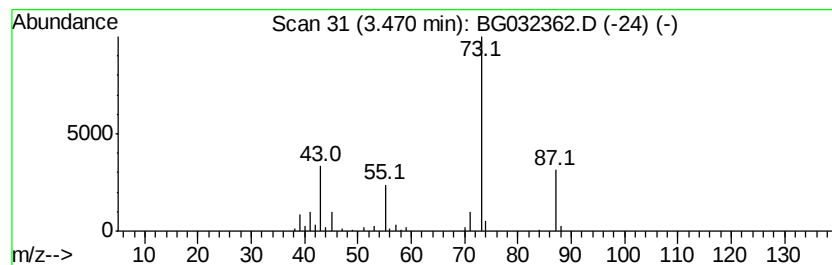
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Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.47	5.46 ng	56707	1,4-Dichlorobenzene-d4	8.72

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			2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	37
4			N-Ethylformamide	73	C3H7NO	000627-45-2	9
5			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9



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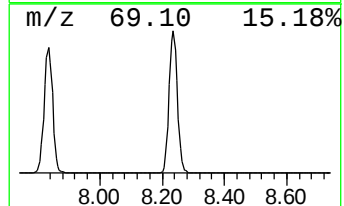
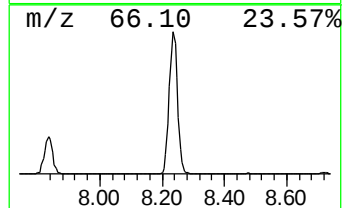
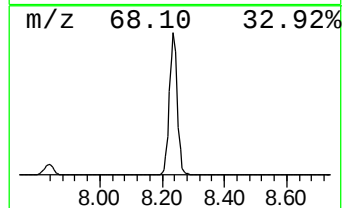
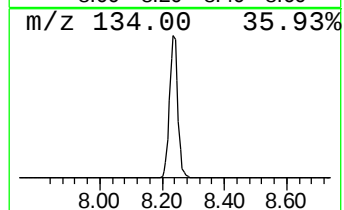
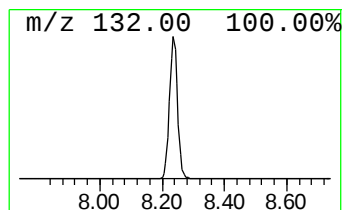
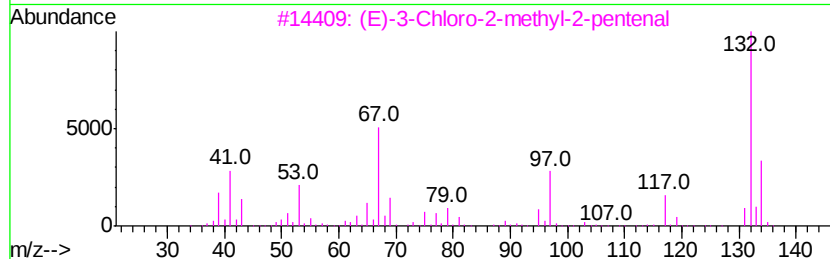
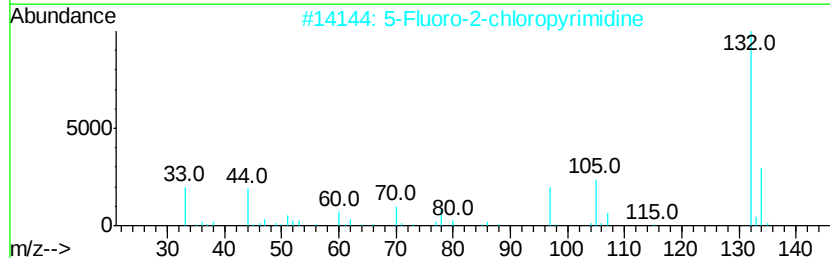
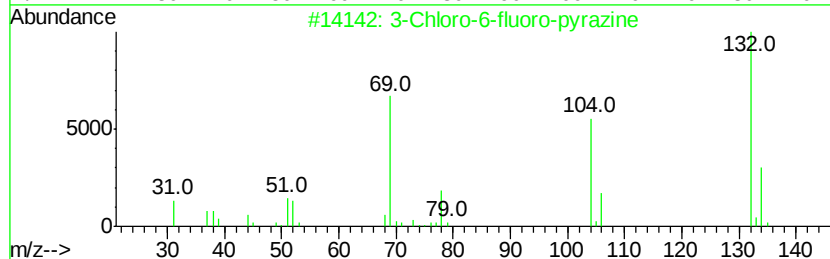
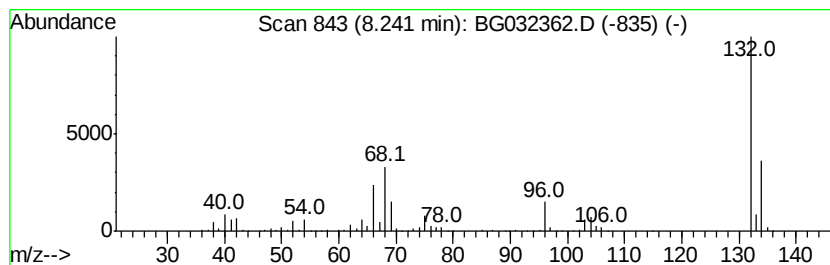
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Peak Number 2 unknown8.24 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.24	87.61 ng	910221	1,4-Dichlorobenzene-d4	8.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	17
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
5		5-Aminoindole	132	C8H8N2	005192-03-0	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	3.47	5.5	ng	56707	1	8.72	207796	20.0
unknown8.24	8.24	87.6	ng	910221	1	8.72	207796	20.0