

Data Path : Z:\HPCHEM1\BNA G\DATA\BG122717\
 Data File : BG031807.D
 Acq On : 27 Dec 2017 21:44
 Operator : SJ/JU
 Sample : I7058-08
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 122117-SIDI-F-U-S

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-BG122117.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	3.432	17	24	35	rBV	77236	163606	4.12%	1.174%
2	3.520	35	39	51	rVB	37243	60965	1.54%	0.437%
3	6.223	491	499	507	rBV	149120	258880	6.53%	1.857%
4	7.898	776	784	798	rVB2	98081	183518	4.63%	1.317%
5	8.315	846	855	865	rBV	302183	560091	14.12%	4.019%
6	8.802	931	938	948	rBB	118756	213364	5.38%	1.531%
7	9.143	988	996	1008	rBV	461604	868381	21.89%	6.231%
8	10.001	1133	1142	1152	rBV	329914	599017	15.10%	4.298%
9	11.687	1420	1429	1440	rVB	166385	312723	7.88%	2.244%
10	14.061	1825	1833	1846	rBB	1345433	2157434	54.39%	15.480%
11	15.436	2061	2067	2077	rVB2	317145	539982	13.61%	3.874%
12	16.905	2310	2317	2329	rBV	785205	1237313	31.19%	8.878%
13	18.168	2524	2532	2541	rBV	480546	783202	19.74%	5.620%
14	20.694	2956	2962	2973	rBV	2715230	3966729	100.00%	28.462%
15	22.510	3263	3271	3283	rVB	536220	1009763	25.46%	7.245%
16	26.258	3896	3909	3925	rBV2	298627	1022093	25.77%	7.334%

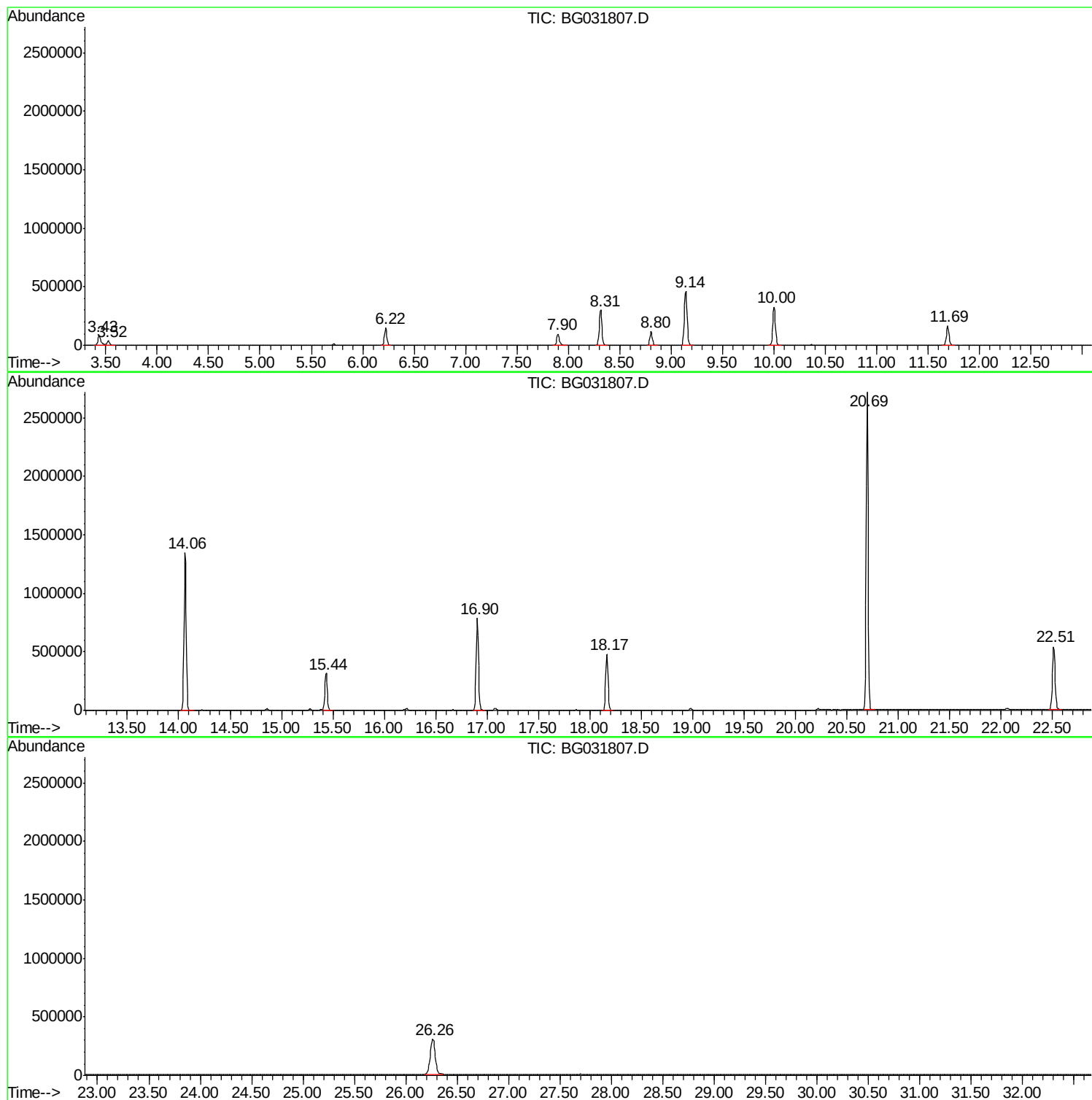
Sum of corrected areas: 13937061

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG122117.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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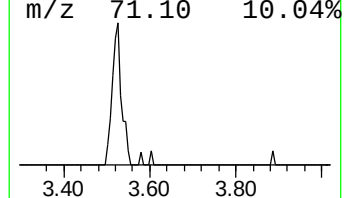
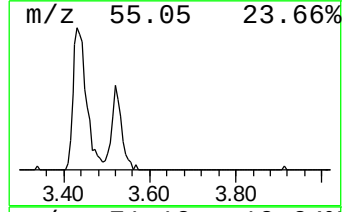
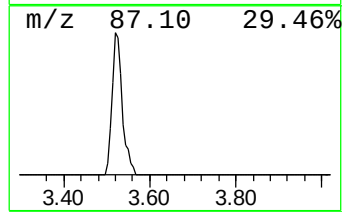
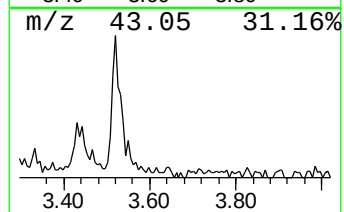
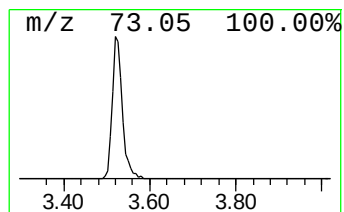
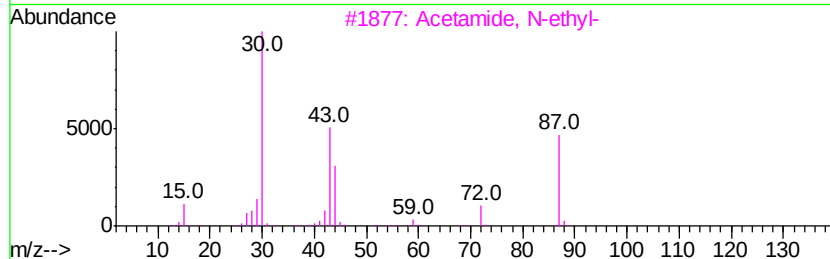
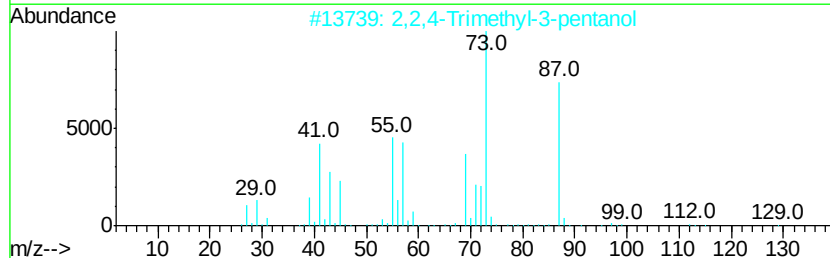
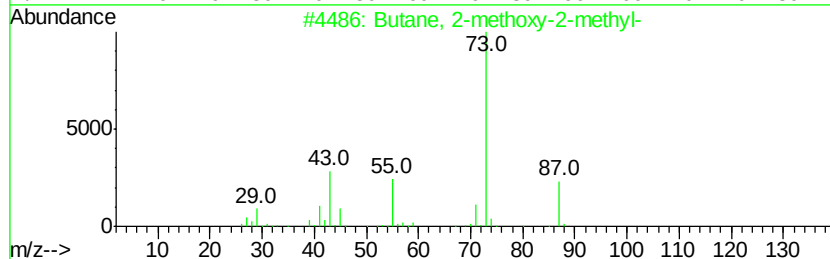
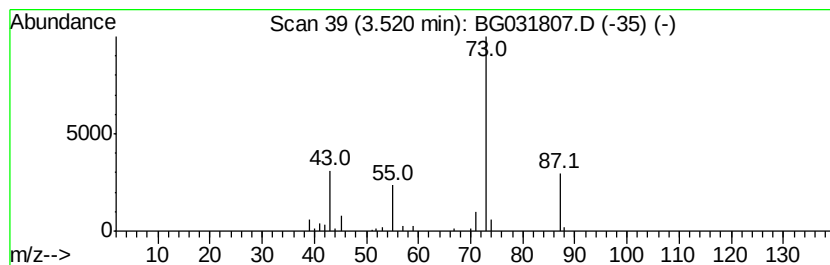
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Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.52	5.71 ng	60965	1,4-Dichlorobenzene-d4	8.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9
4		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9
5		1,4-Butanediamine	88	C4H12N2	000110-60-1	9



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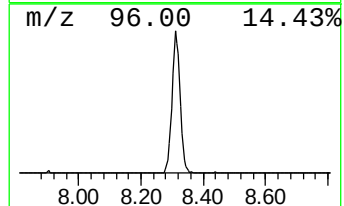
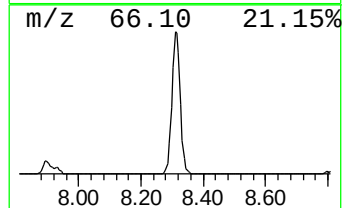
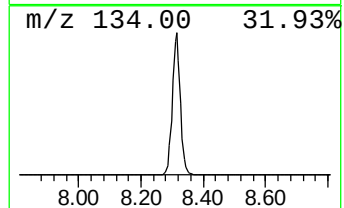
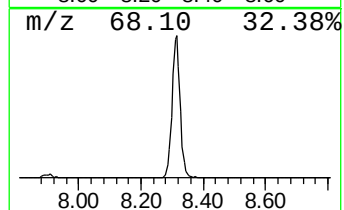
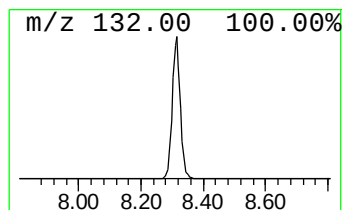
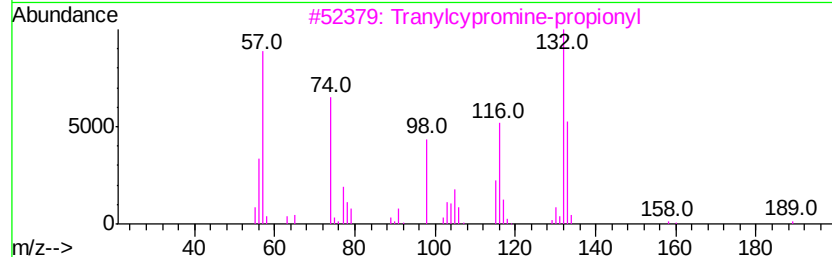
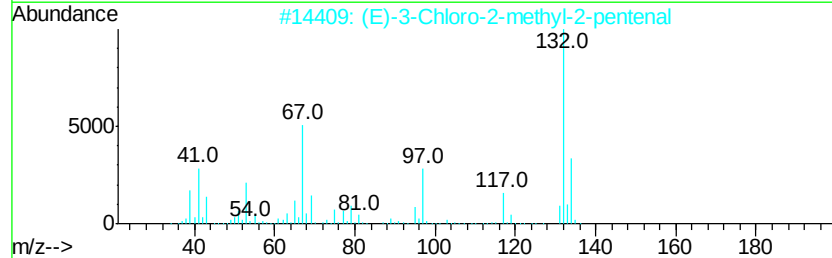
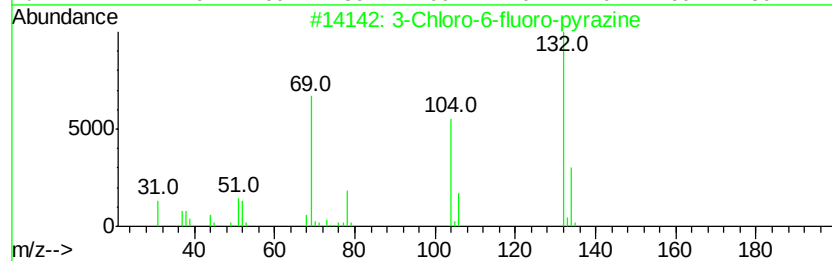
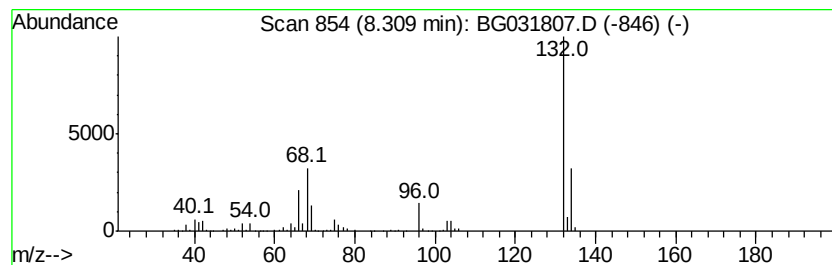
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Peak Number 3 unknown8.31 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.31	52.50 ng	560091	1,4-Dichlorobenzene-d4	8.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		1H-Indazole, 7-methyl-	132	C8H8N2	1000316-00-7	10
5		Xenon	132	Xe	007440-63-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
Butane, 2-methoxy...	3.52	5.7	ng	60965	1	8.80	213364	20.0
unknown8.31	8.31	52.5	ng	560091	1	8.80	213364	20.0