

Data Path : U:\HPCHEM1\BNA G\DATA\BG122617\
 Data File : BG031785.D
 Acq On : 27 Dec 2017 6:31
 Operator : SJ/JU
 Sample : I7009-03
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 121917-JETT-F-D-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	3.434	19	25	36	rBV2	73698	157476	3.70%	1.048%
2	3.523	36	40	52	rVB	28041	48015	1.13%	0.319%
3	6.225	493	500	509	rBV	166692	290323	6.83%	1.932%
4	7.900	778	785	797	rVB	113787	211453	4.97%	1.407%
5	8.311	847	855	869	rBV	322565	612288	14.40%	4.074%
6	8.805	932	939	950	rVB	132900	236457	5.56%	1.573%
7	9.140	987	996	1009	rVB	500828	964152	22.67%	6.415%
8	10.003	1134	1143	1155	rVB	351942	650206	15.29%	4.326%
9	11.689	1423	1430	1443	rBV	190938	347369	8.17%	2.311%
10	14.063	1825	1834	1844	rBV	1469904	2366075	55.64%	15.743%
11	15.432	2059	2067	2076	rVB2	338725	580752	13.66%	3.864%
12	16.907	2311	2318	2333	rBV	771207	1243156	29.23%	8.272%
13	18.164	2526	2532	2542	rBV2	539559	841563	19.79%	5.600%
14	20.697	2957	2963	2972	rBV	3039893	4252501	100.00%	28.295%
15	22.512	3263	3272	3283	rBV2	581495	1104793	25.98%	7.351%
16	26.267	3899	3911	3925	rBV2	336500	1122670	26.40%	7.470%

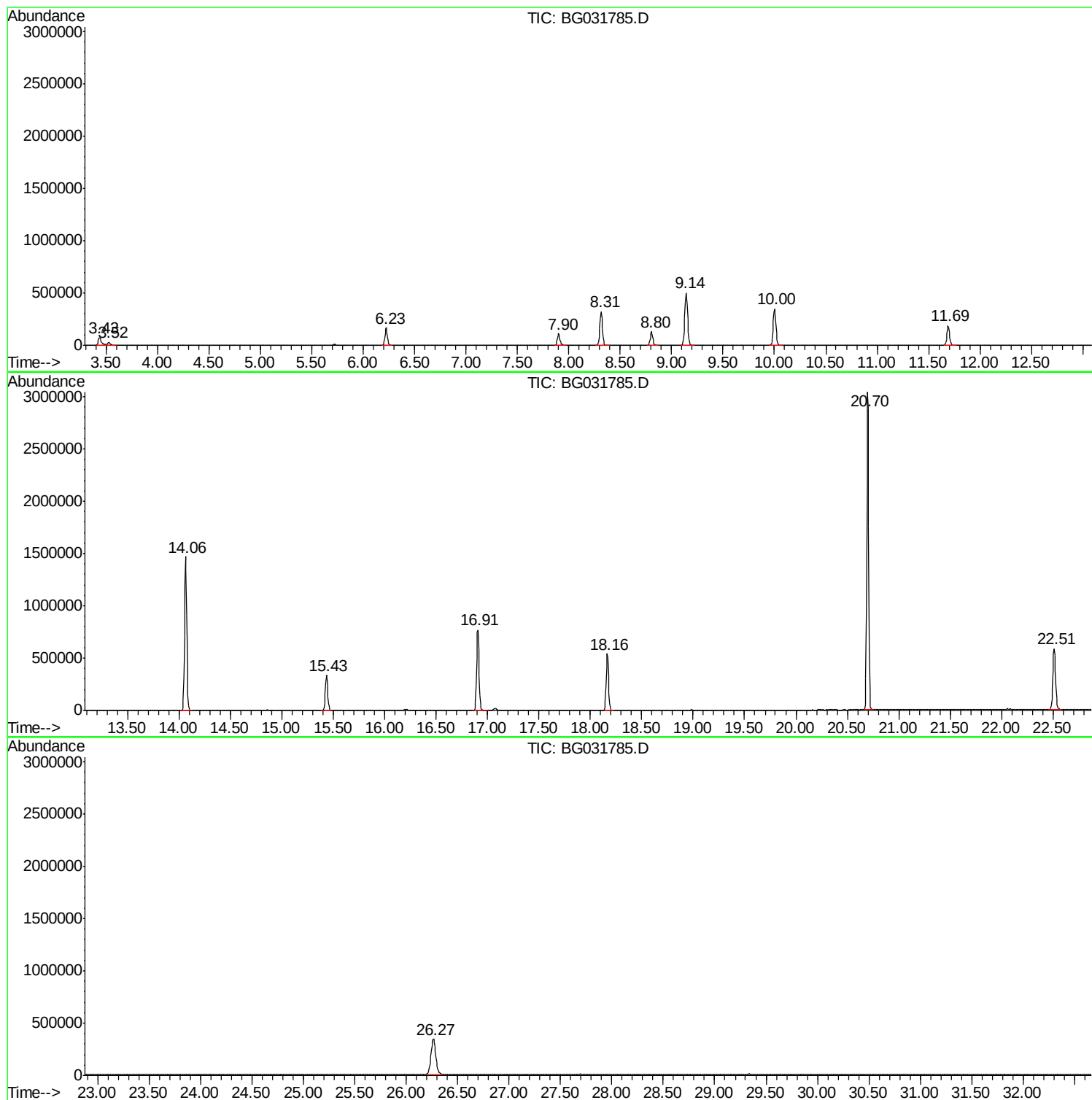
Sum of corrected areas: 15029249

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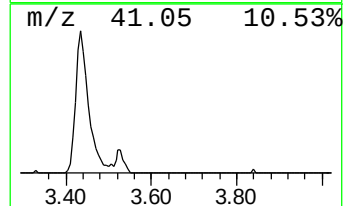
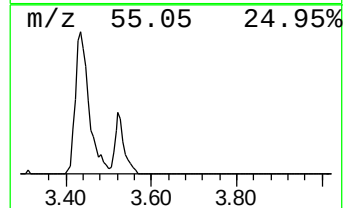
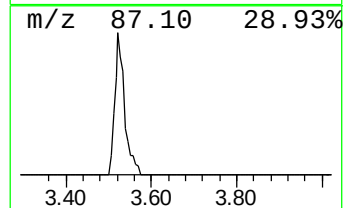
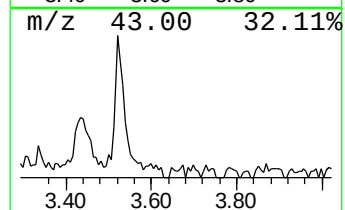
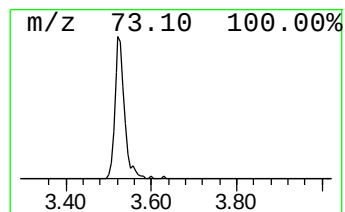
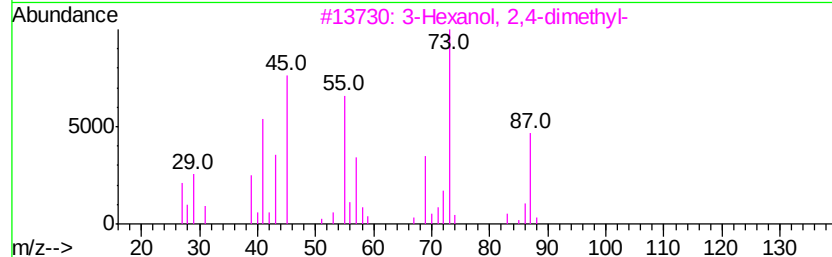
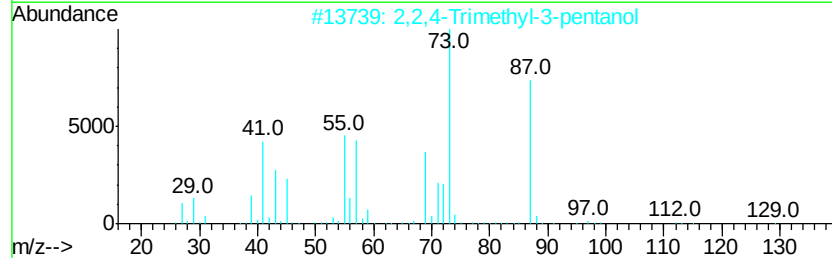
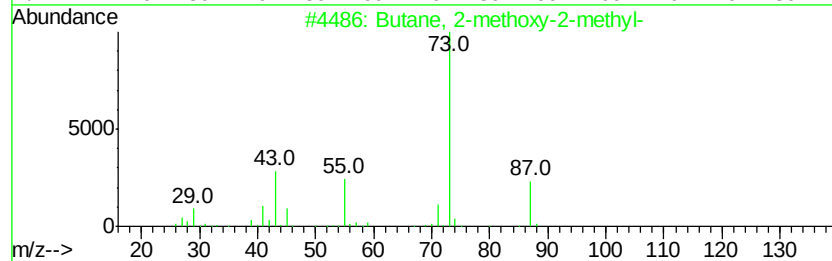
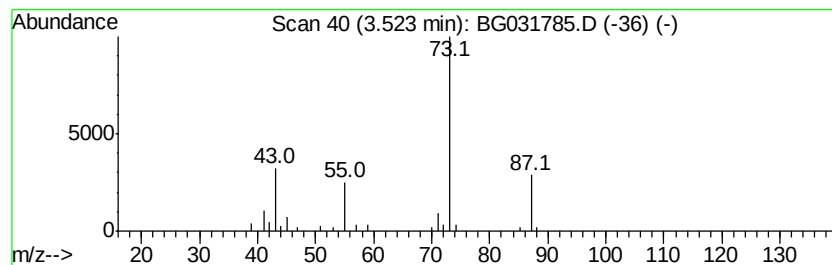
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Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.52	4.06 ng	48015	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	72
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3		3-Hexanol, 2,4-dimethyl-	130	C8H18O	013432-25-2	28
4		2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	12
5		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	10



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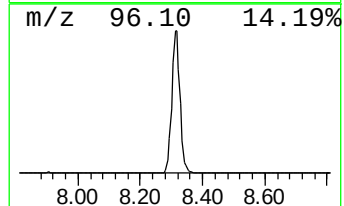
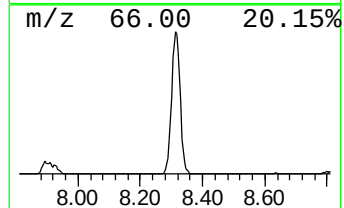
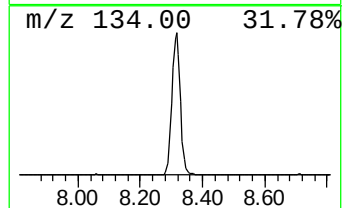
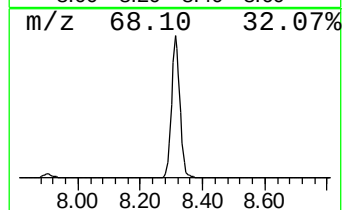
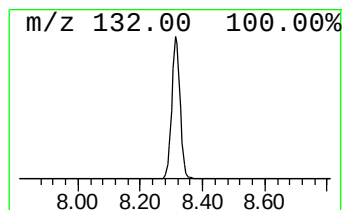
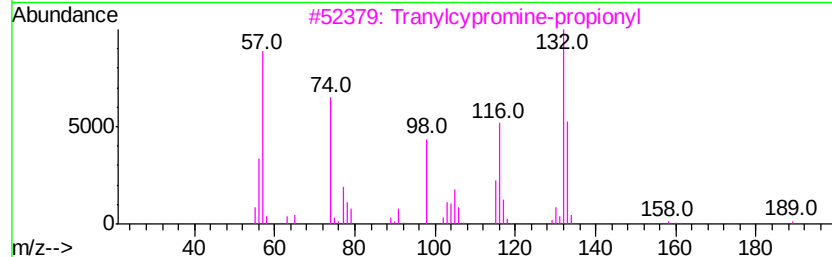
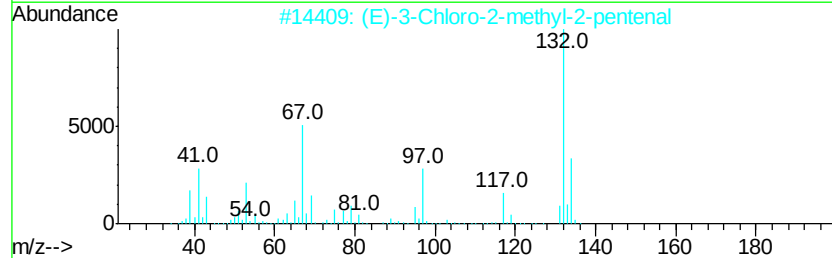
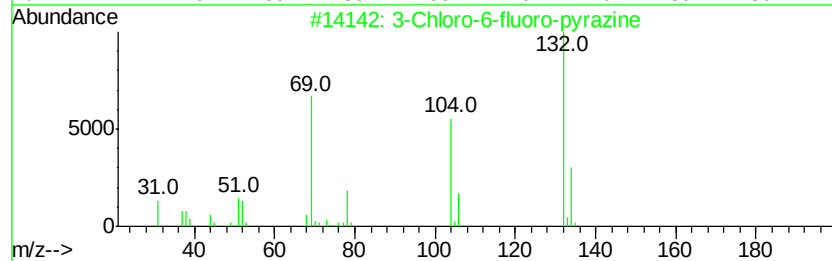
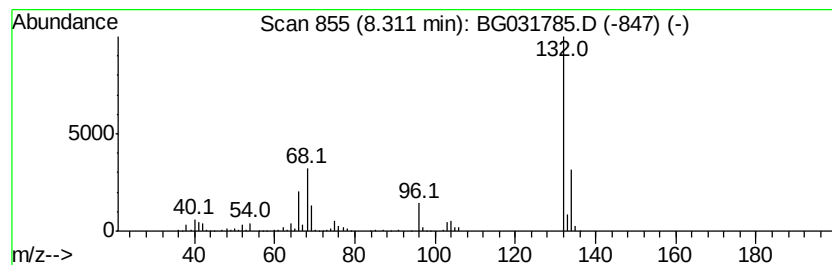
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Peak Number 3 unknown8.31 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.31	51.79 ng	612288	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	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	12
4		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9
5		5-Aminoindole	132	C8H8N2	005192-03-0	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	3.52	4.1	ng	48015	1	8.80	236457	20.0
unknown8.31	8.31	51.8	ng	612288	1	8.80	236457	20.0