

Data Path : U:\HPCHEM1\BNA G\DATA\BG122617\
 Data File : BG031784.D
 Acq On : 27 Dec 2017 5:53
 Operator : SJ/JU
 Sample : I7009-02
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 121917-JETT-F-D-M

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	3.438	18	25	35	rBV	77179	160635	3.55%	0.984%
2	3.526	35	40	51	rVB	32421	54396	1.20%	0.333%
3	6.223	492	499	511	rBV	218019	378317	8.37%	2.317%
4	7.897	777	784	797	rBB	141842	279129	6.18%	1.710%
5	8.314	846	855	864	rBV	399128	726852	16.08%	4.452%
6	8.802	931	938	948	rVB	126115	235780	5.22%	1.444%
7	9.143	988	996	1008	rVB	578205	1070144	23.68%	6.555%
8	10.006	1133	1143	1155	rBV	375576	723624	16.01%	4.432%
9	11.687	1422	1429	1442	rVB	178590	350766	7.76%	2.149%
10	14.066	1825	1834	1844	rBV	1583254	2573499	56.94%	15.763%
11	15.435	2061	2067	2076	rVB2	366340	606332	13.42%	3.714%
12	16.904	2310	2317	2330	rBV	896236	1440148	31.87%	8.821%
13	18.167	2525	2532	2540	rBV2	584239	891811	19.73%	5.463%
14	20.700	2957	2963	2973	rBV	3257239	4519375	100.00%	27.682%
15	22.515	3264	3272	3279	rBV	620868	1121900	24.82%	6.872%
16	26.270	3897	3911	3931	rBV2	341511	1193192	26.40%	7.309%

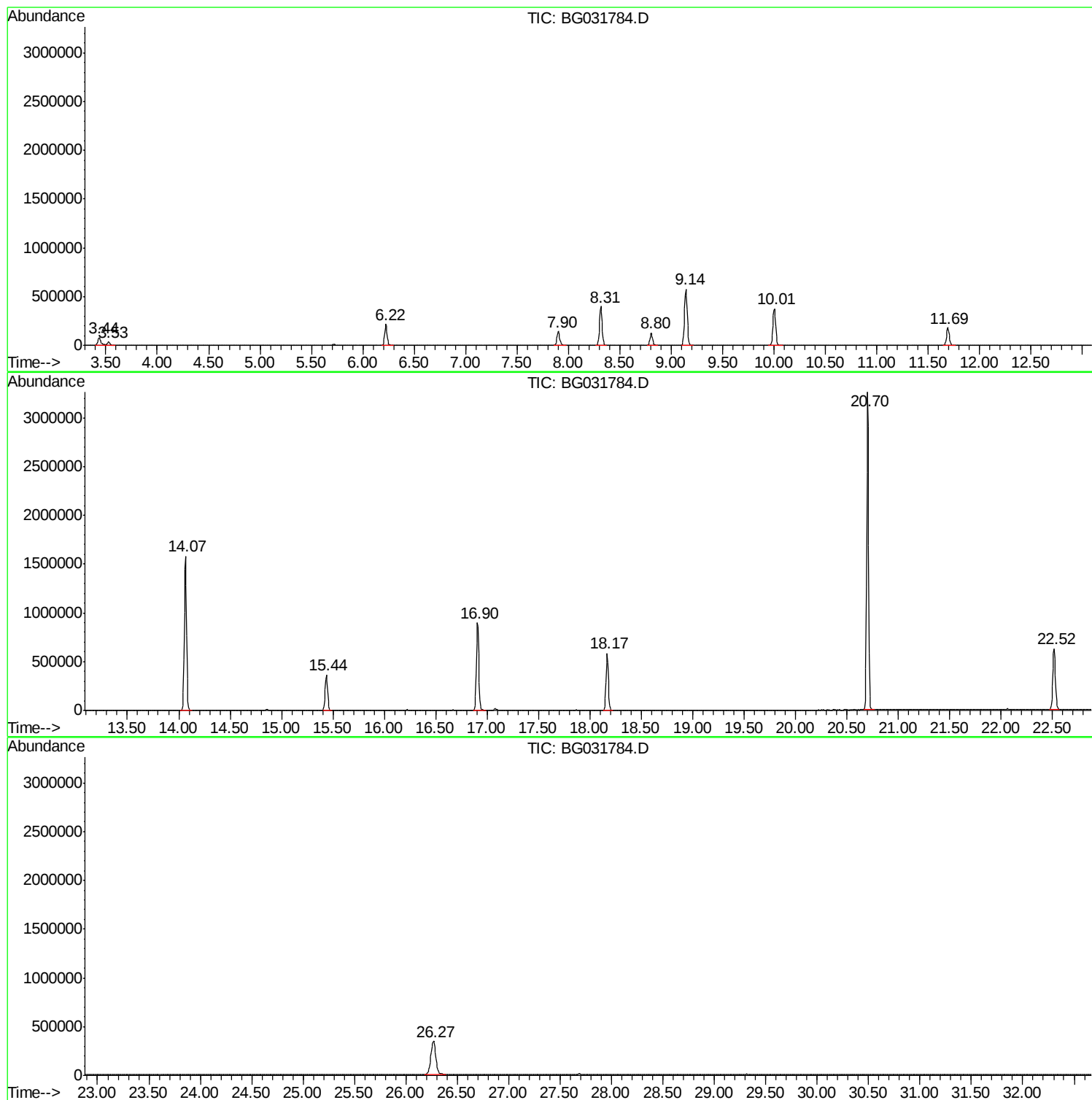
Sum of corrected areas: 16325900

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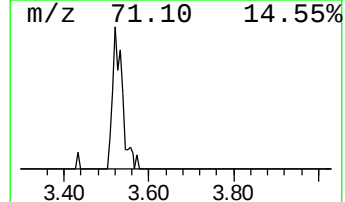
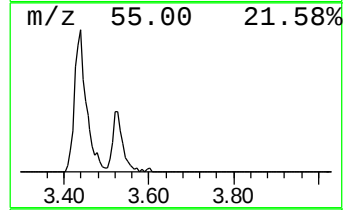
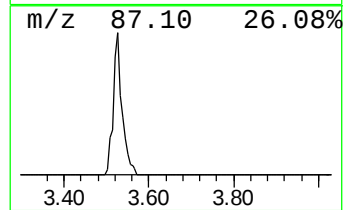
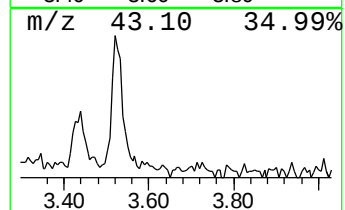
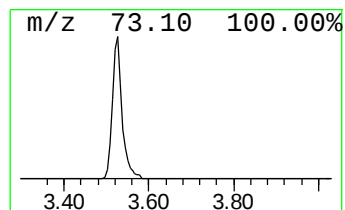
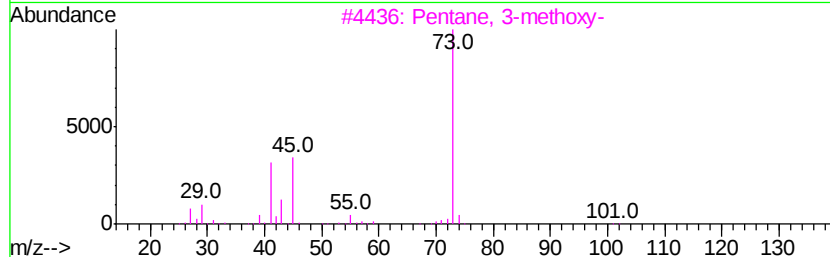
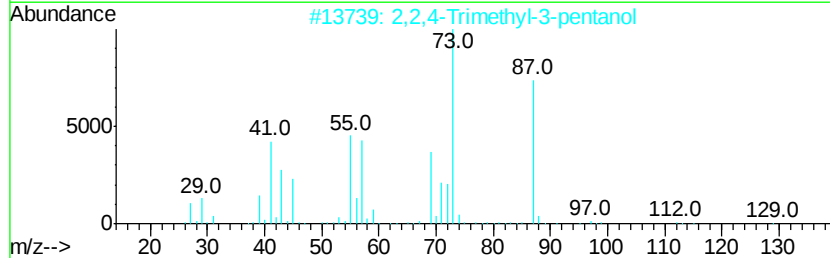
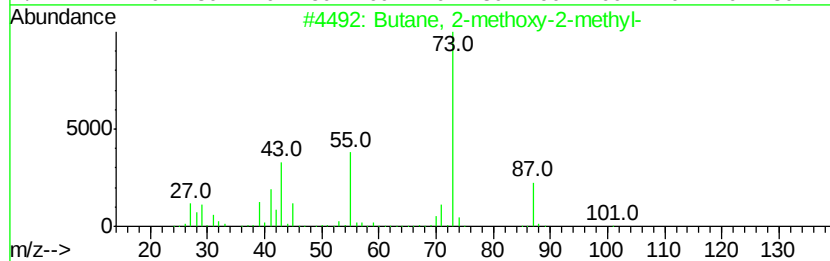
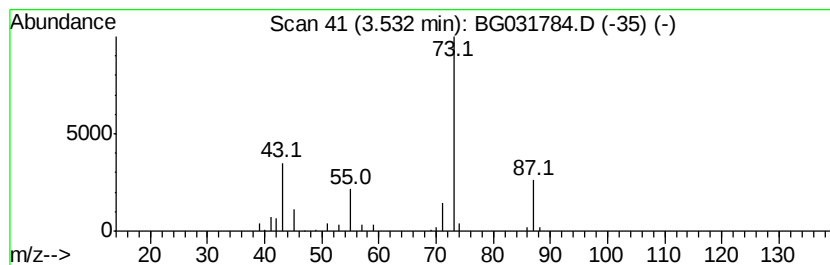
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Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.53	4.61 ng	54396	1,4-Dichlorobenzene-d4	8.80

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	56
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
4		1-Propene-1-thiol	74	C3H6S	000925-89-3	9
5		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9



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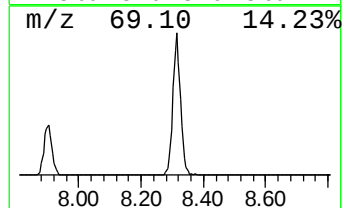
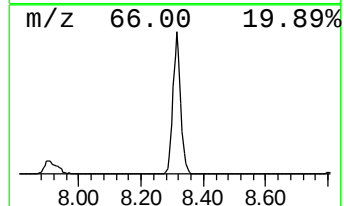
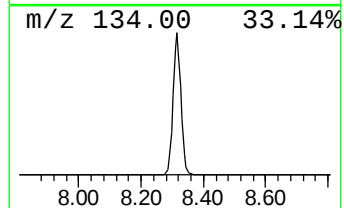
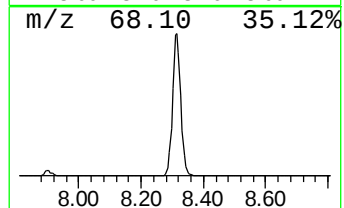
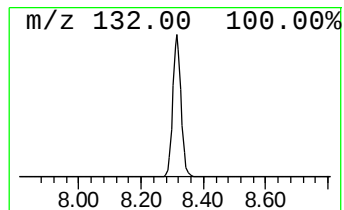
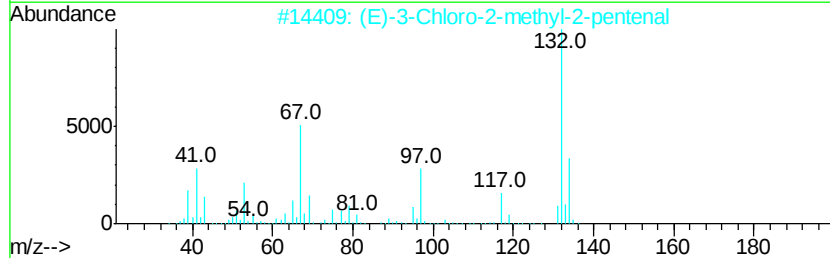
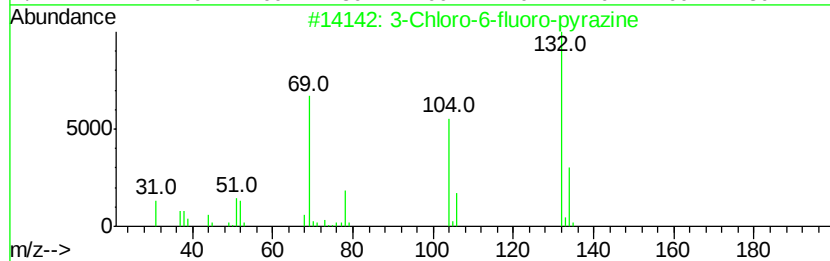
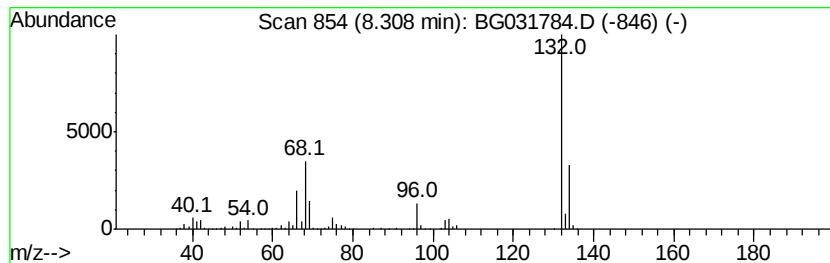
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Peak Number 3 unknown8.31 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.31	61.66 ng	726852	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	25
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Tranvlcypromine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	3.53	4.6	ng	54396	1	8.80	235780	20.0
unknown8.31	8.31	61.7	ng	726852	1	8.80	235780	20.0