

Data Path : U:\HPCHEM1\BNA G\DATA\BG011118\
 Data File : BG031993.D
 Acq On : 11 Jan 2018 20:47
 Operator : SJ/JU
 Sample : J1070-01
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TW-1

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	3.419	16	22	32	rBV2	90333	212308	4.42%	1.168%
2	3.513	32	38	45	rVB	66252	106900	2.22%	0.588%
3	6.204	489	496	506	rBV	249952	434155	9.03%	2.389%
4	7.879	771	781	794	rBV2	155771	297660	6.19%	1.638%
5	8.290	842	851	863	rBV	547902	989104	20.58%	5.443%
6	8.778	926	934	944	rBV	141462	258776	5.38%	1.424%
7	9.118	984	992	1005	rVB	557375	1046763	21.78%	5.760%
8	9.976	1129	1138	1147	rBV	403279	757414	15.76%	4.168%
9	11.662	1416	1425	1434	rBV	207030	388390	8.08%	2.137%
10	14.036	1818	1829	1841	rBV	1691771	2726256	56.71%	15.002%
11	15.405	2056	2062	2073	rVB3	384829	679264	14.13%	3.738%
12	16.880	2306	2313	2323	rBV	1394651	2295210	47.75%	12.630%
13	18.137	2520	2527	2537	rBV2	564420	929742	19.34%	5.116%
14	20.675	2952	2959	2966	rBV	3392264	4807077	100.00%	26.452%
15	22.479	3259	3266	3278	rBV	601173	1141552	23.75%	6.282%
16	26.204	3887	3900	3917	rBV2	329768	1102459	22.93%	6.066%

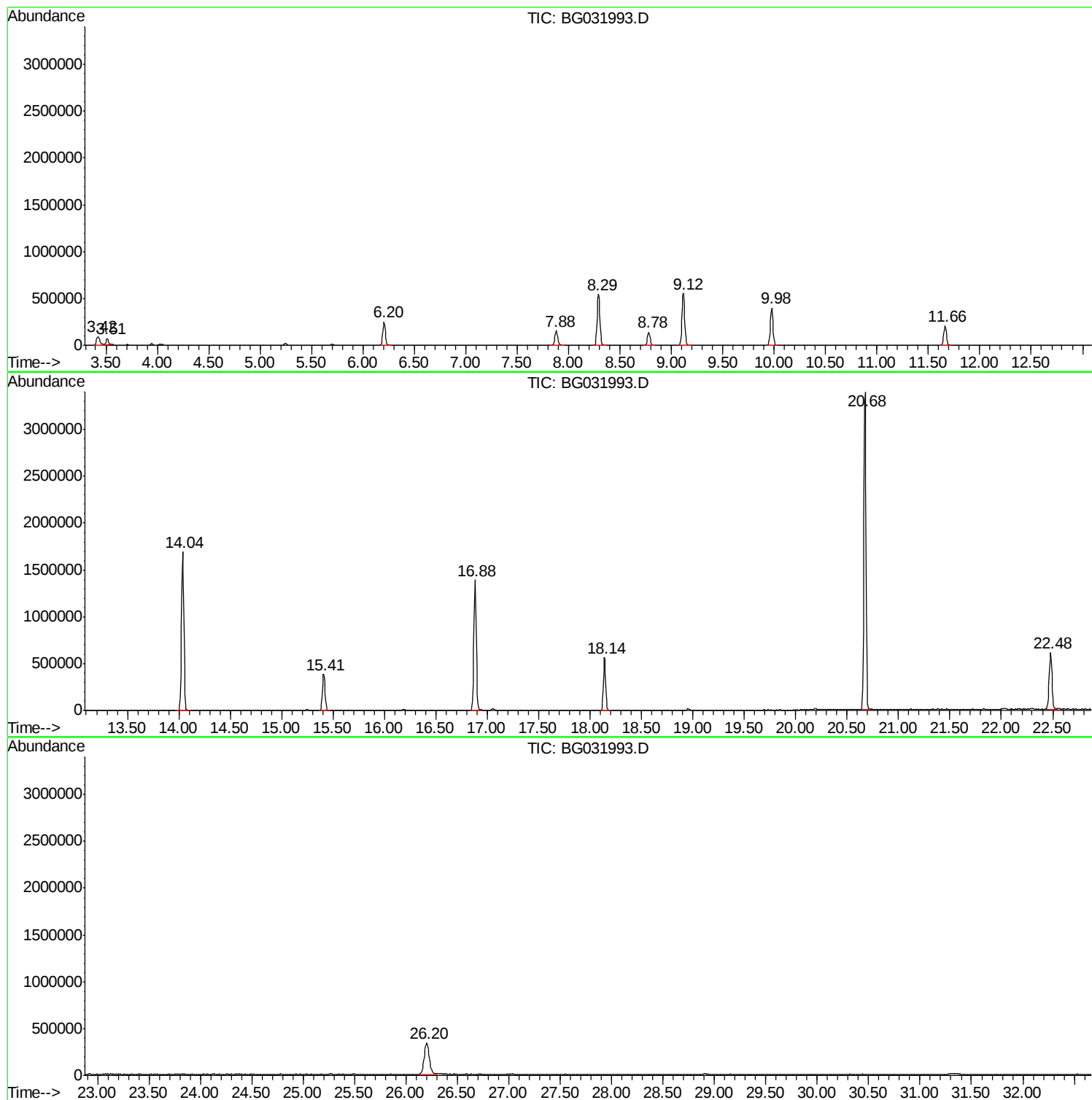
Sum of corrected areas: 18173030

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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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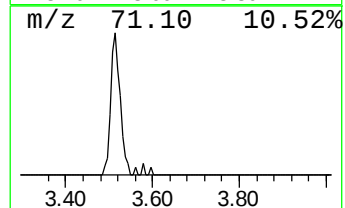
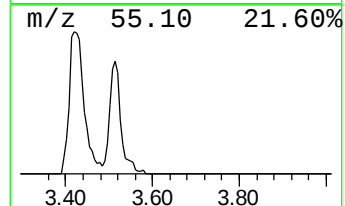
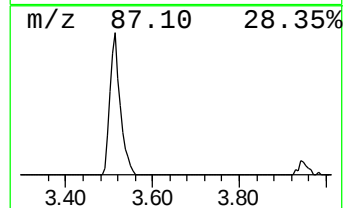
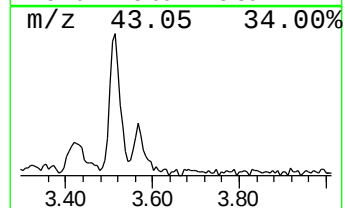
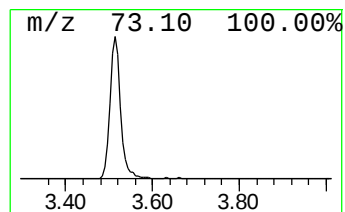
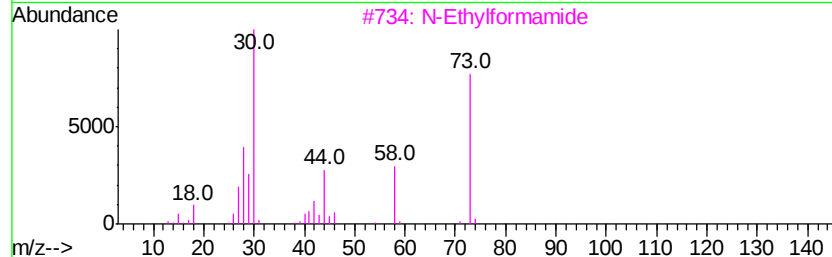
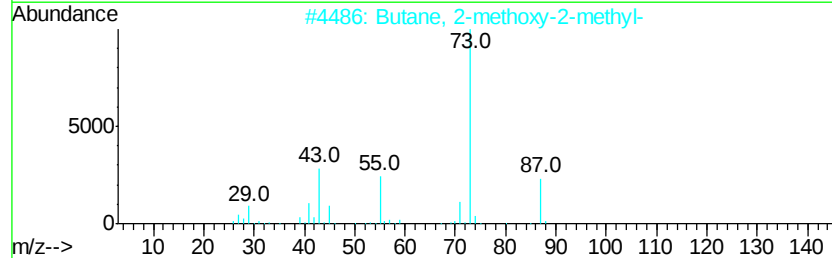
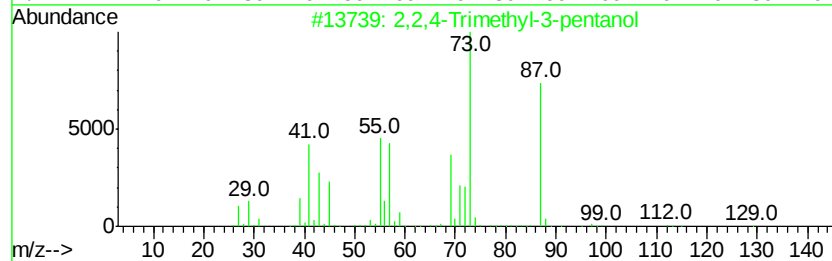
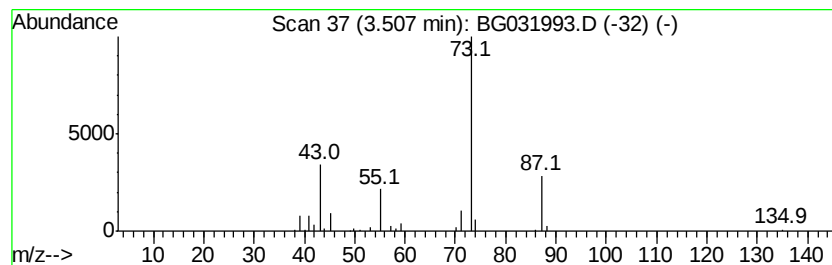
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Peak Number 2 2,2,4-Trimethyl-3-pentanol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.51	8.26 ng	106900	1,4-Dichlorobenzene-d4	8.78

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



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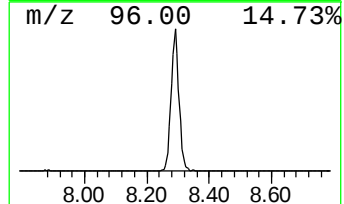
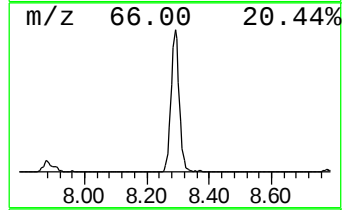
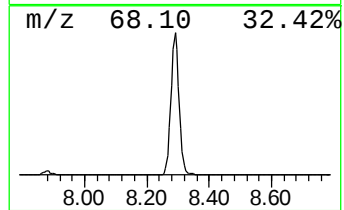
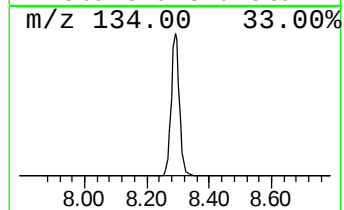
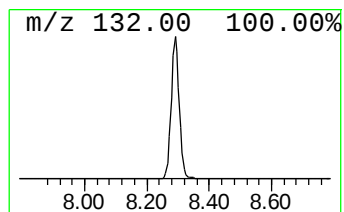
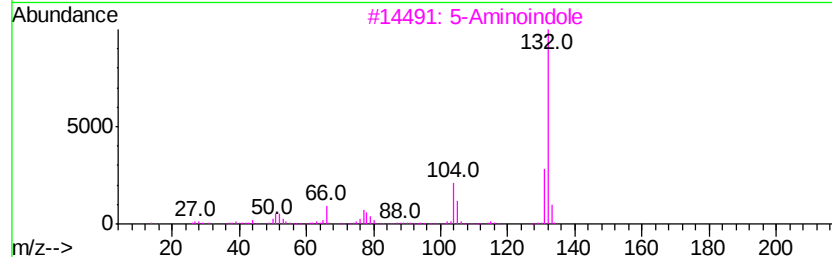
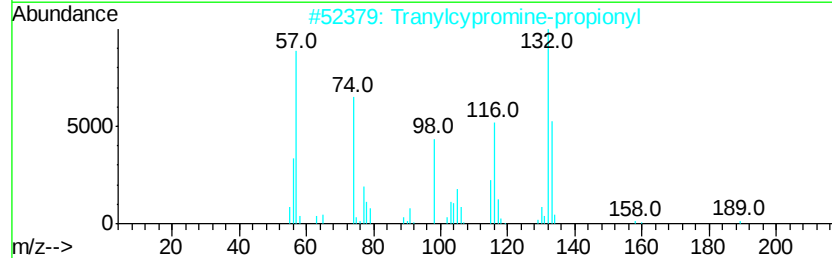
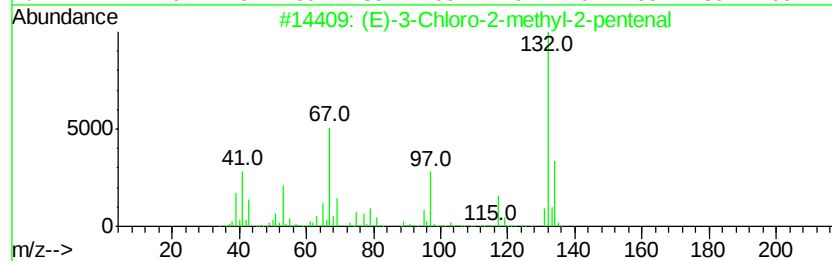
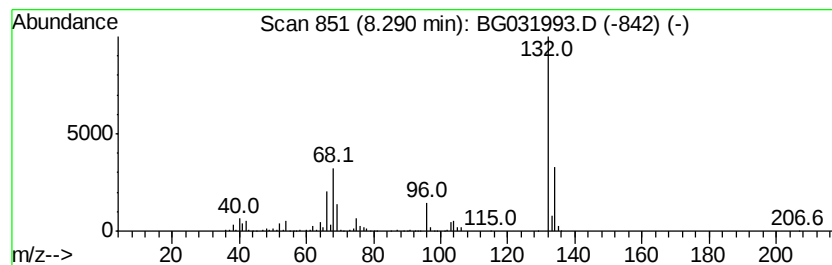
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Peak Number 3 unknown8.29 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.29	76.44 ng	989104	1,4-Dichlorobenzene-d4	8.78

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
3		5-Aminoindole	132	C8H8N2	005192-03-0	10
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	10
5		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2,2,4-Trimethyl-3...	3.51	8.3	ng	106900	1	8.78	258776	20.0
unknown8.29	8.29	76.4	ng	989104	1	8.78	258776	20.0