

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042518\
Data File : BG033999.D
Acq On : 25 Apr 2018 22:56
Operator : SJ/JU
Sample : PB108595BL
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
PB108595BL

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-BG041918.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	4.983	283	288	297	rVB	48359	74125	1.36%	0.308%
2	5.436	358	365	379	rVB	1034718	1615935	29.71%	6.712%
3	6.998	623	631	643	rBV	1079657	1781799	32.76%	7.401%
4	7.363	685	693	702	rVB	1048560	1720855	31.64%	7.148%
5	7.827	764	772	779	rBV	203673	337035	6.20%	1.400%
6	8.144	815	826	835	rBV	716852	1222225	22.47%	5.077%
7	8.973	959	967	977	rBV	602713	1051907	19.34%	4.369%
8	10.612	1238	1246	1256	rVB	290994	533346	9.81%	2.215%
9	13.080	1657	1666	1676	rBV	1859466	2842574	52.26%	11.807%
10	14.460	1892	1901	1913	rBV2	505788	829330	15.25%	3.445%
11	15.947	2147	2154	2168	rBV2	1566516	2424699	44.58%	10.071%
12	17.210	2362	2369	2382	rVB	810494	1187254	21.83%	4.931%
13	19.842	2811	2817	2823	rBV	4018838	5439387	100.00%	22.593%
14	21.464	3086	3093	3100	rBV	1028299	1549294	28.48%	6.435%
15	24.525	3602	3614	3625	rBV2	549294	1466270	26.96%	6.090%

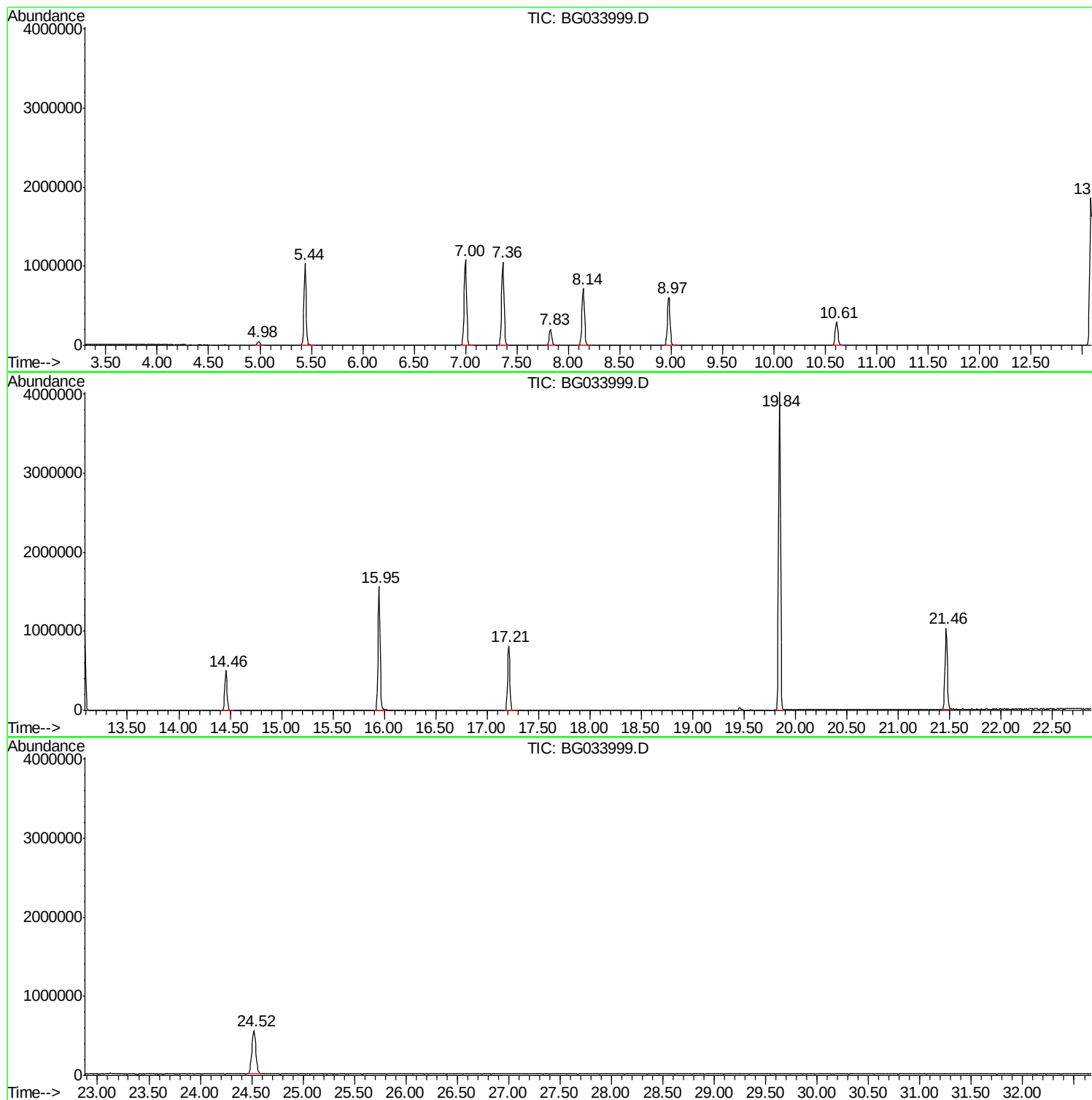
Sum of corrected areas: 24076035

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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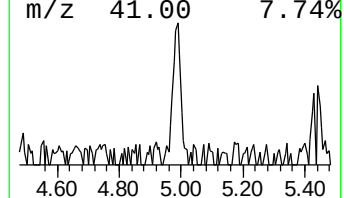
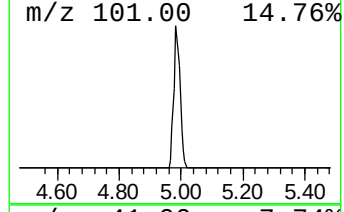
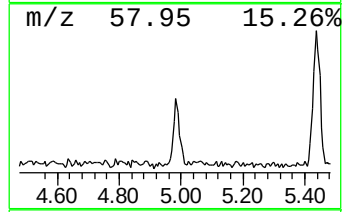
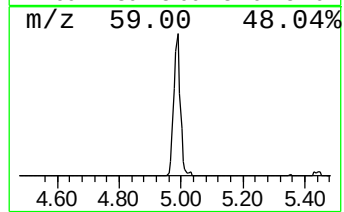
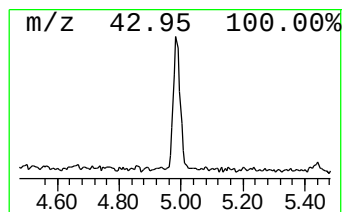
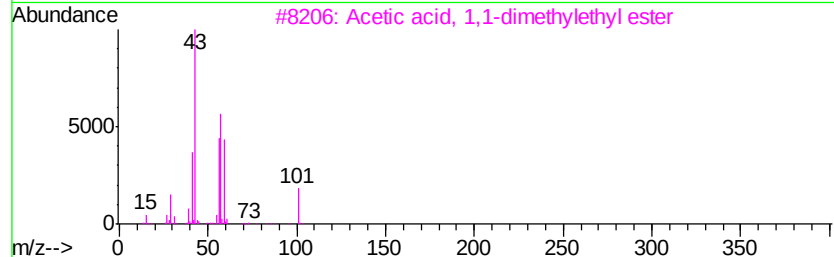
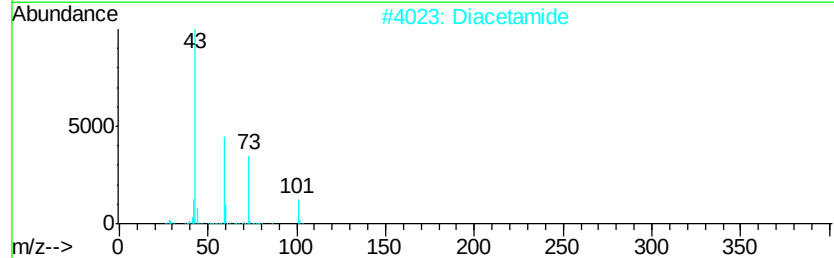
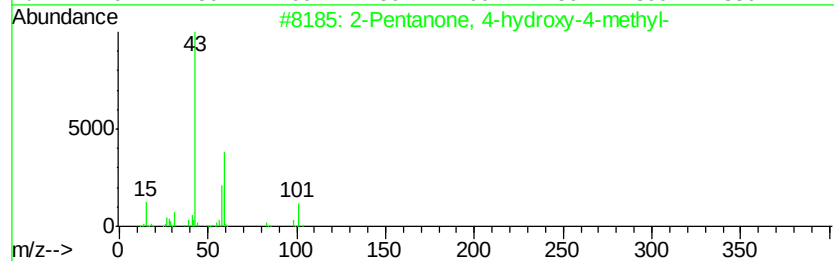
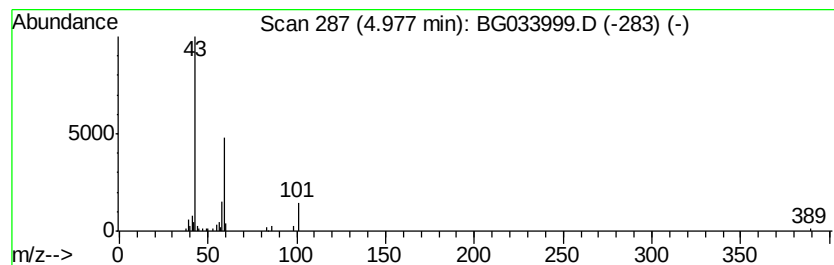
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.98	4.40 ng	74125	1,4-Dichlorobenzene-d4	7.83

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2			Diacetamide	101	C4H7NO2	000625-77-4	38
3			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	27
5			Acetamide, N-(aminocarbonyl)-	102	C3H6N2O2	000591-07-1	9



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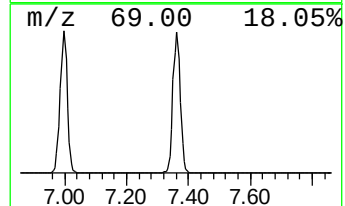
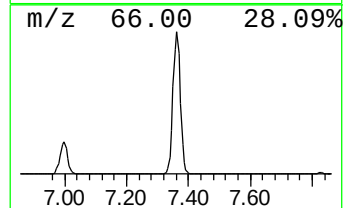
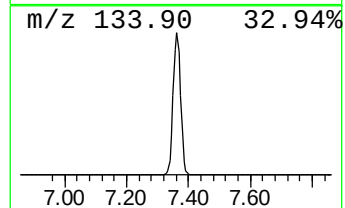
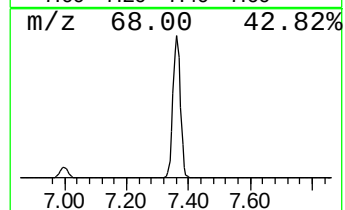
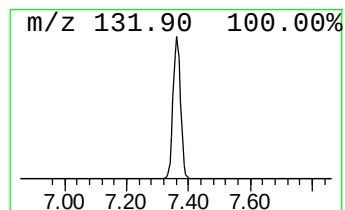
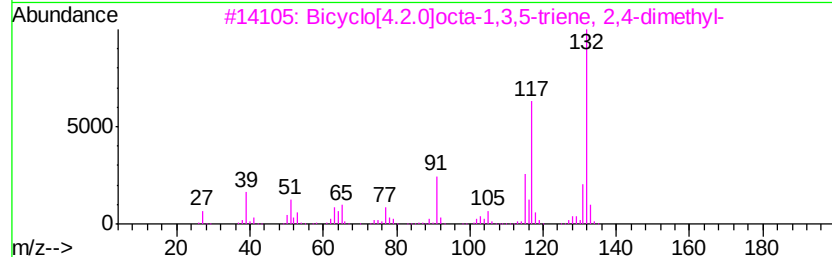
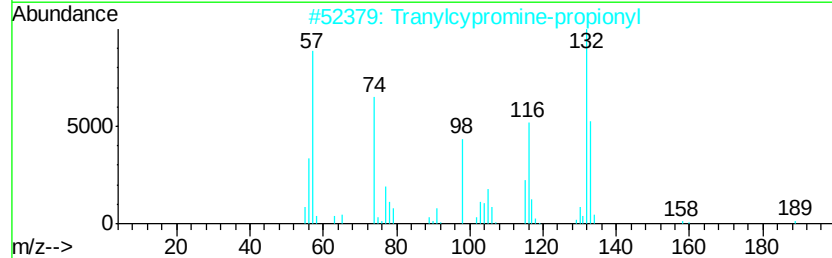
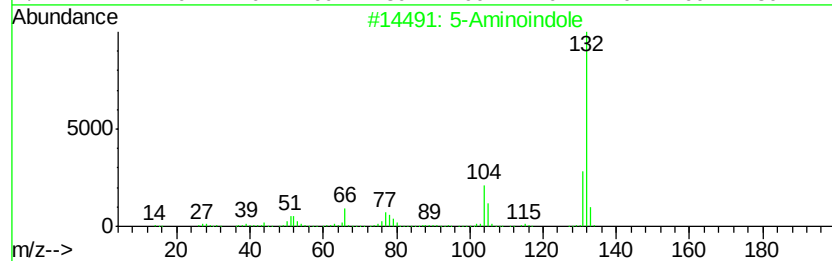
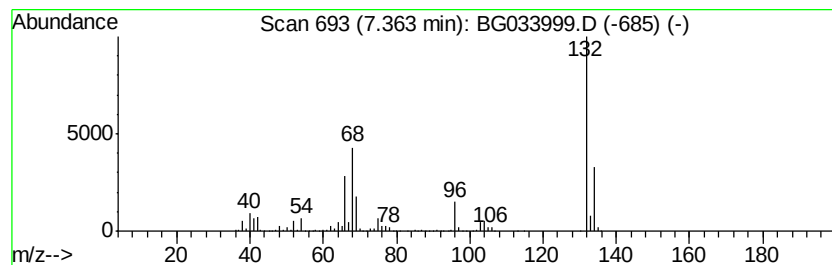
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Peak Number 2 unknown7.36 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.36	102.12 ng	1720860	1,4-Dichlorobenzene-d4	7.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Aminoindole	132	C8H8N2	005192-03-0	12
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	9
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.98	4.4	ng	74125	1	7.83	337035	20.0
unknown7.36	7.36	102.1	ng	1720860	1	7.83	337035	20.0