

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042018\
 Data File : BG033892.D
 Acq On : 20 Apr 2018 18:52
 Operator : SJ/JU
 Sample : PB108385BL
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB108385BL

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	4.992	285	290	299	rBV	67052	105173	1.99%	0.398%
2	5.445	360	367	375	rBV	1389807	2126121	40.30%	8.045%
3	7.002	624	632	641	rBV	1327425	2249452	42.64%	8.511%
4	7.372	686	695	705	rBV	1297238	2205679	41.81%	8.346%
5	7.836	767	774	781	rVB	256318	437539	8.29%	1.656%
6	8.153	820	828	836	rVB	995193	1660225	31.47%	6.282%
7	8.982	959	969	978	rBV	762029	1316306	24.95%	4.981%
8	10.615	1239	1247	1258	rVB	345630	616557	11.69%	2.333%
9	13.089	1661	1668	1675	rBV	2041447	3164821	59.99%	11.975%
10	14.464	1893	1902	1912	rBV2	543680	871094	16.51%	3.296%
11	15.956	2148	2156	2169	rBV	1729575	2511986	47.62%	9.505%
12	17.219	2362	2371	2381	rVB2	773789	1162219	22.03%	4.398%
13	19.857	2814	2820	2828	rVB	3944245	5275569	100.00%	19.961%
14	21.473	3089	3095	3104	rVB2	900357	1387642	26.30%	5.250%
15	24.534	3605	3616	3628	rVB2	496188	1338479	25.37%	5.064%

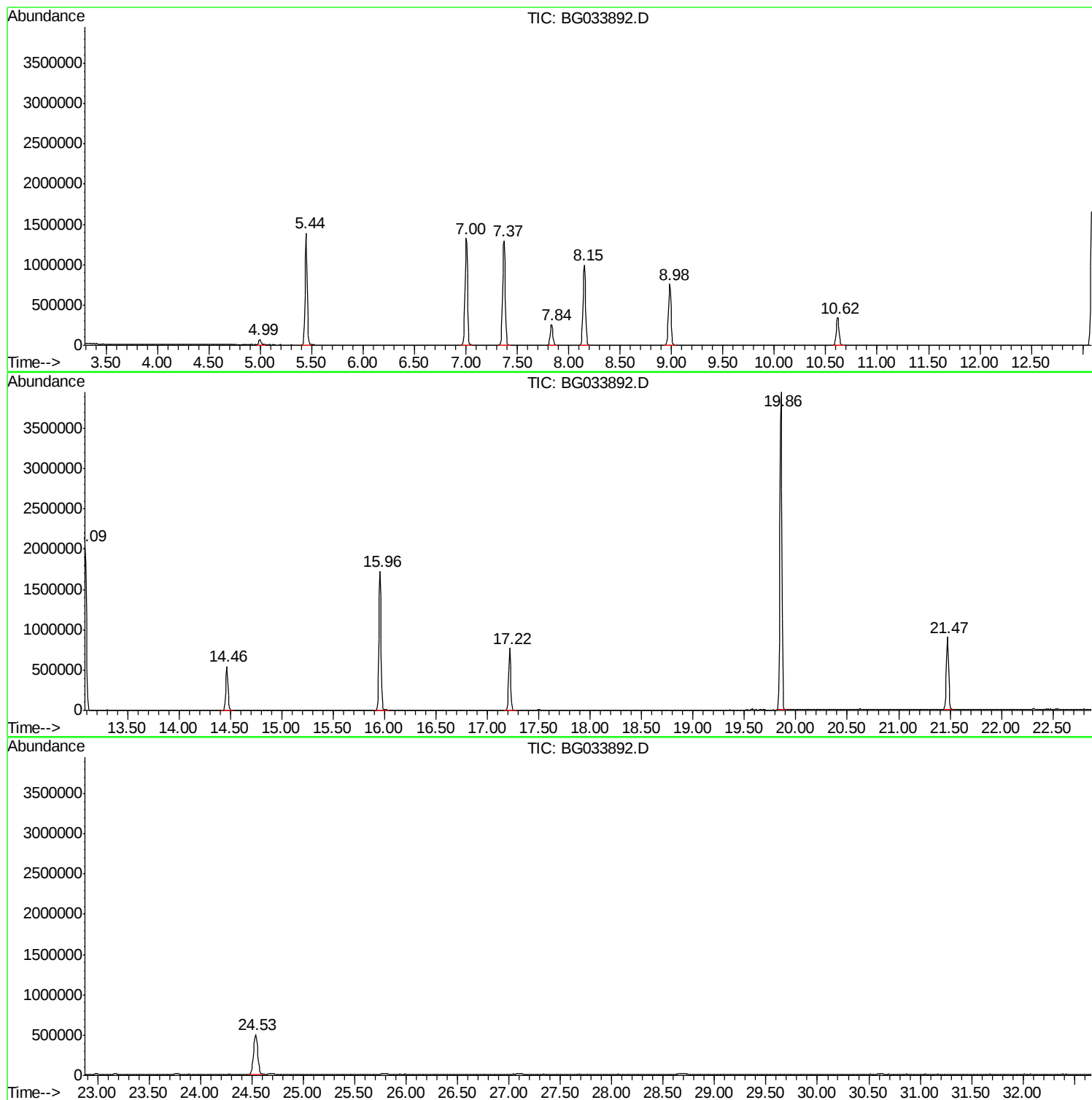
Sum of corrected areas: 26428862

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG041918.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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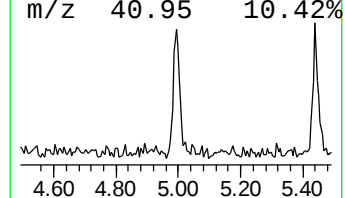
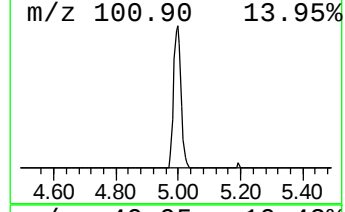
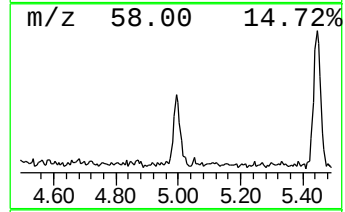
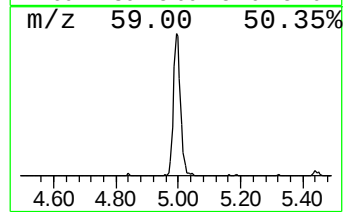
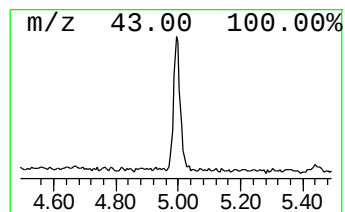
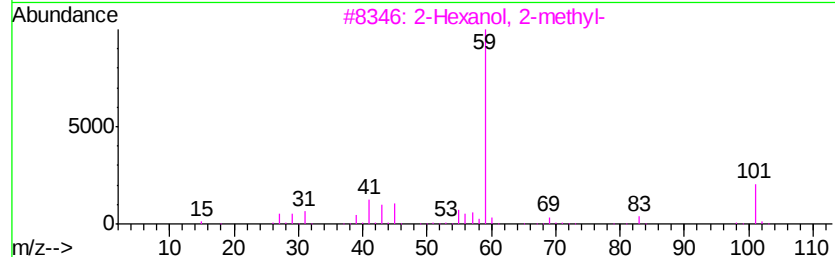
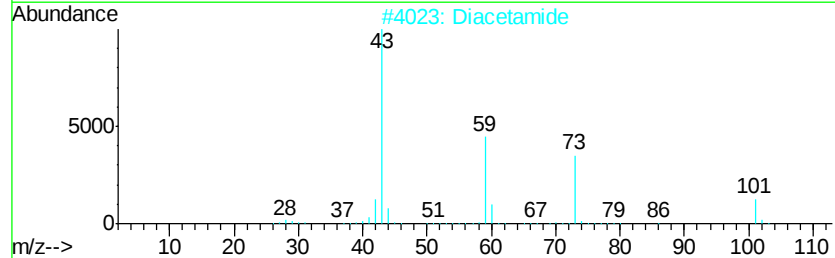
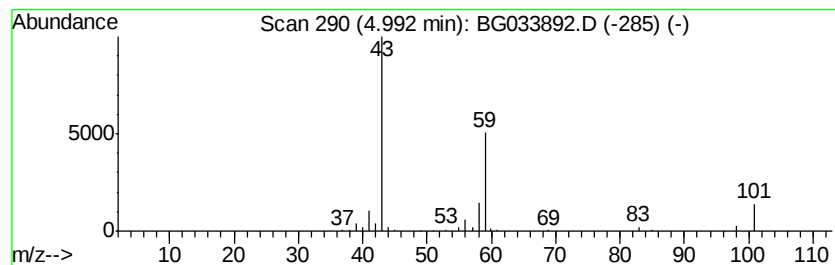
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.99	4.81 ng	105173	1,4-Dichlorobenzene-d4	7.84

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2	Diacetamide	101	C4H7NO2	000625-77-4	9
3	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	9
4	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9
5	Propane, 2-ethoxy-2-methyl-	102	C6H14O	000637-92-3	9



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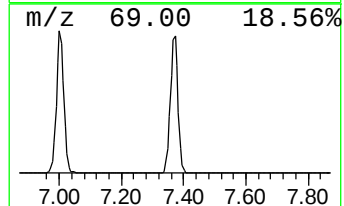
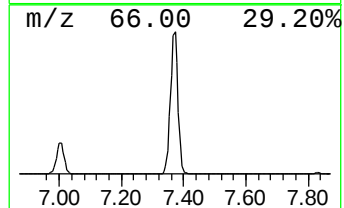
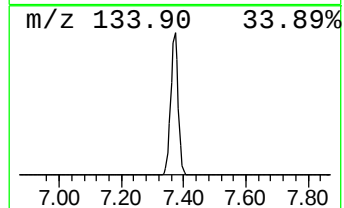
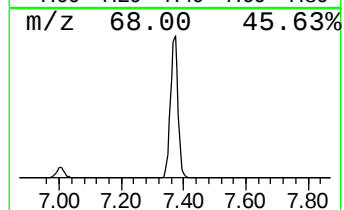
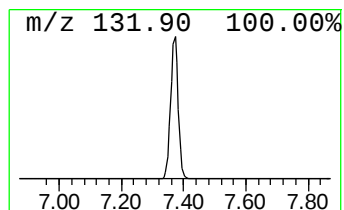
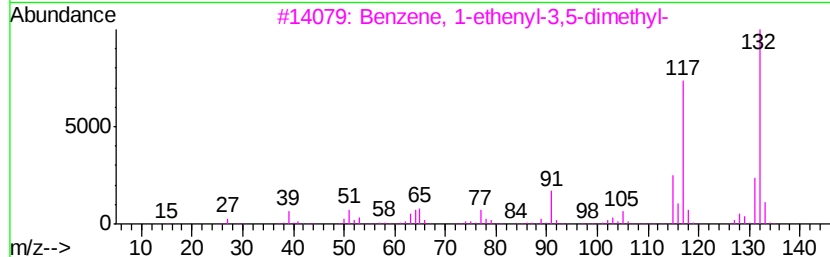
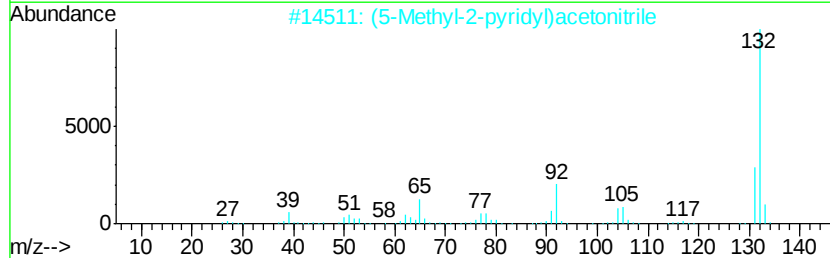
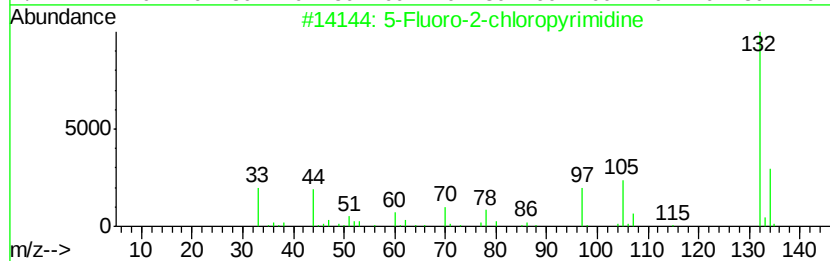
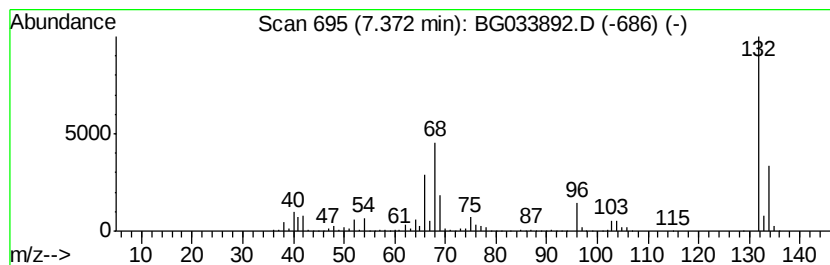
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Peak Number 2 unknown7.37 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.37	100.82 ng	2205680	1,4-Dichlorobenzene-d4	7.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
2		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
3		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.99	4.8	ng	105173	1	7.84	437539	20.0
unknown7.37	7.37	100.8	ng	2205680	1	7.84	437539	20.0