

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042018\
 Data File : BG033913.D
 Acq On : 21 Apr 2018 9:28
 Operator : SJ/JU
 Sample : J2473-16
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 041818-SIDW-F-U-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-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	4.991	285	290	298	rVB	56222	78765	1.27%	0.360%
2	5.437	359	366	376	rBV	396867	618298	9.93%	2.829%
3	6.994	623	631	640	rBV	249134	427611	6.87%	1.957%
4	7.365	686	694	705	rBV	720483	1194305	19.19%	5.465%
5	7.829	763	773	780	rBV	218490	369438	5.94%	1.691%
6	8.146	819	827	838	rVB	858328	1501502	24.12%	6.871%
7	8.980	960	969	977	rBV	762171	1319805	21.21%	6.039%
8	10.614	1239	1247	1256	rVB	301956	525251	8.44%	2.404%
9	13.087	1660	1668	1675	rBV	2019035	3141692	50.48%	14.376%
10	14.462	1894	1902	1911	rBV2	481092	766831	12.32%	3.509%
11	15.949	2149	2155	2166	rBV	1183287	1780248	28.60%	8.146%
12	17.212	2360	2370	2380	rBV	680420	1048175	16.84%	4.796%
13	18.128	2521	2526	2534	rBV3	38275	63218	1.02%	0.289%
14	19.856	2813	2820	2830	rBV	4564533	6223950	100.00%	28.481%
15	21.472	3087	3095	3103	rBV	907929	1418619	22.79%	6.492%
16	24.527	3603	3615	3626	rBV2	500072	1375419	22.10%	6.294%

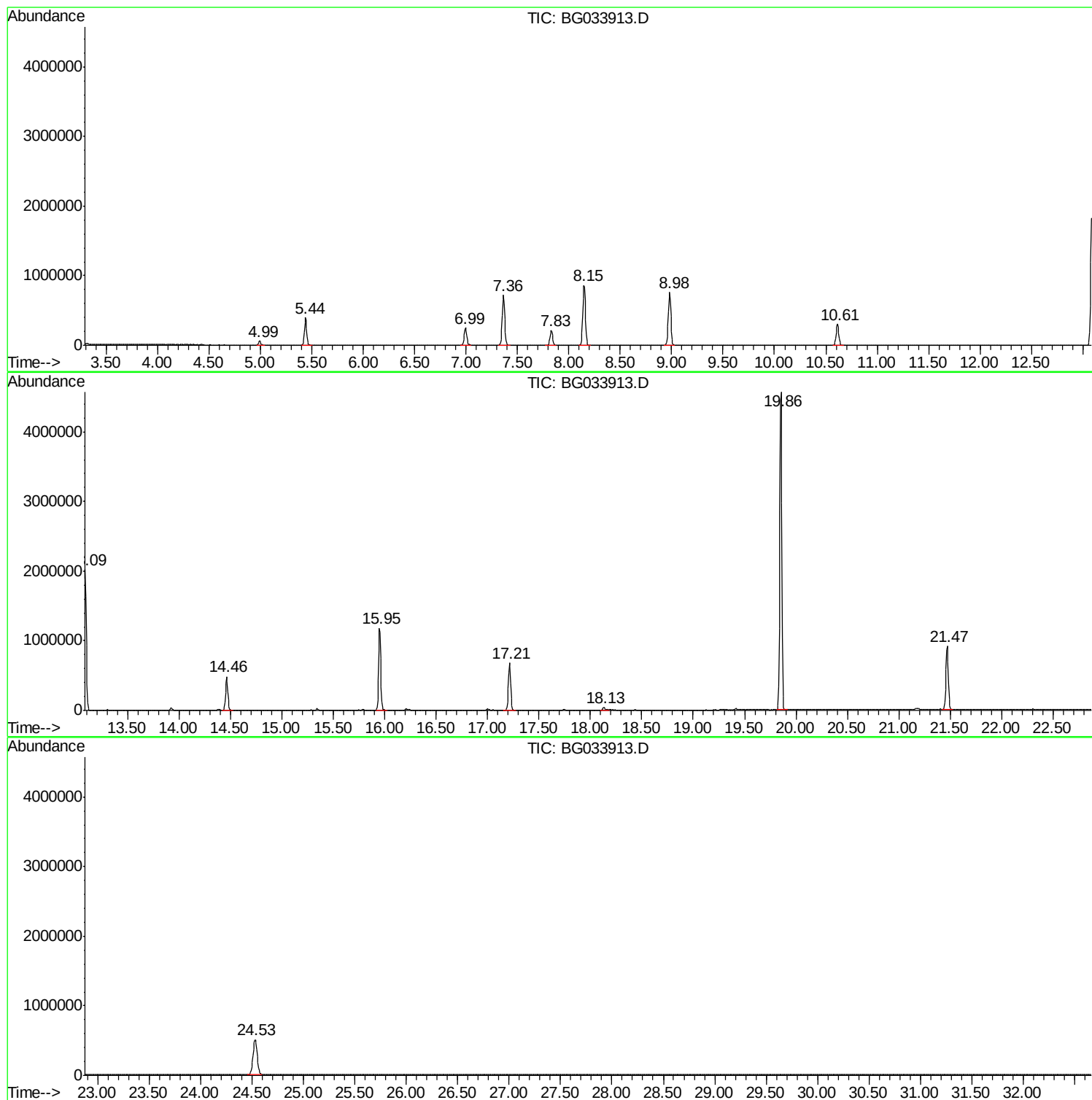
Sum of corrected areas: 21853127

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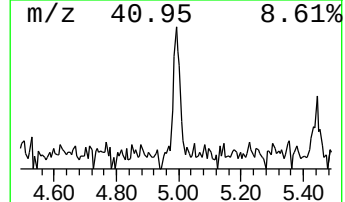
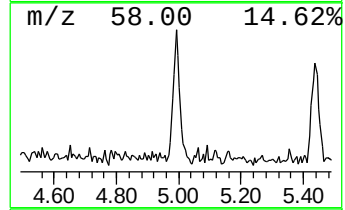
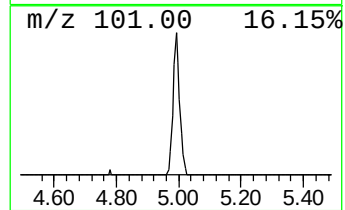
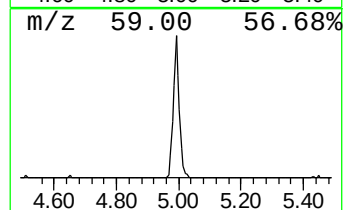
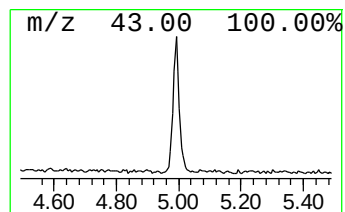
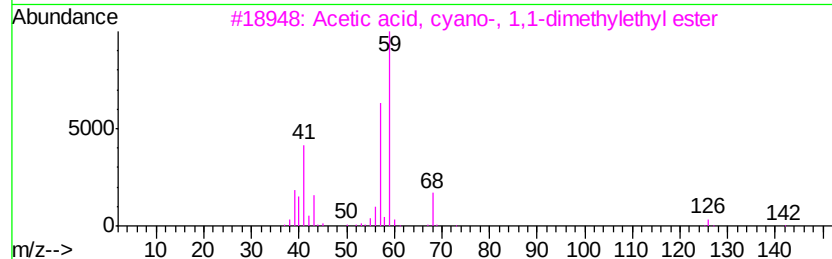
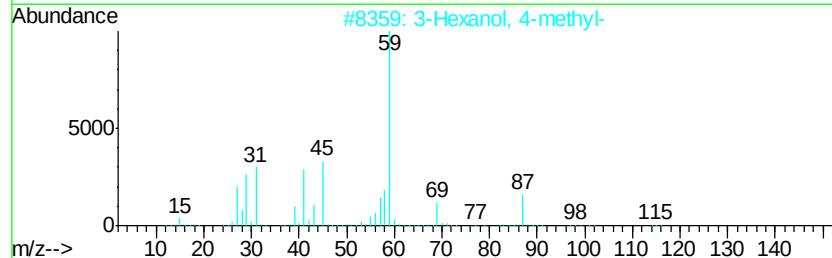
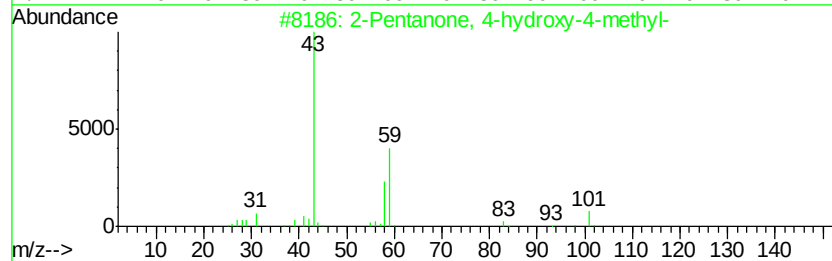
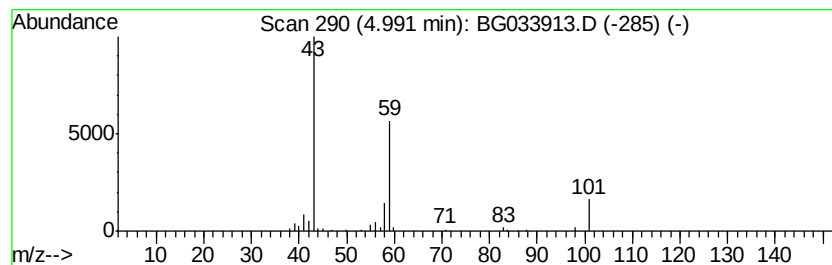
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.99	4.26 ng	78765	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	50
2			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	25
3			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
4			Propane, 2-ethoxy-2-methyl-	102	C6H14O	000637-92-3	23
5			Acetamide	59	C2H5NO	000060-35-5	9



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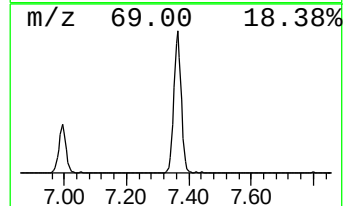
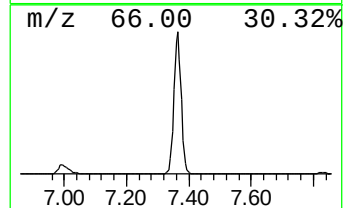
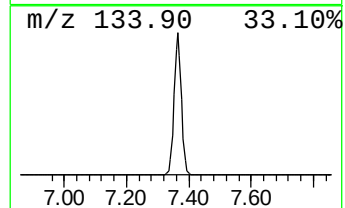
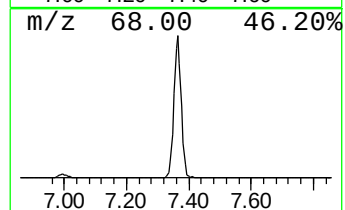
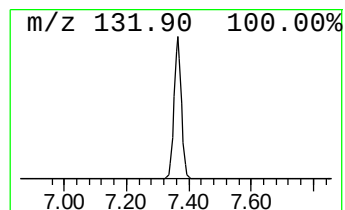
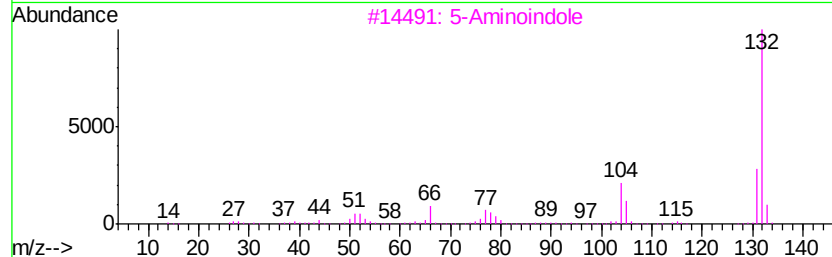
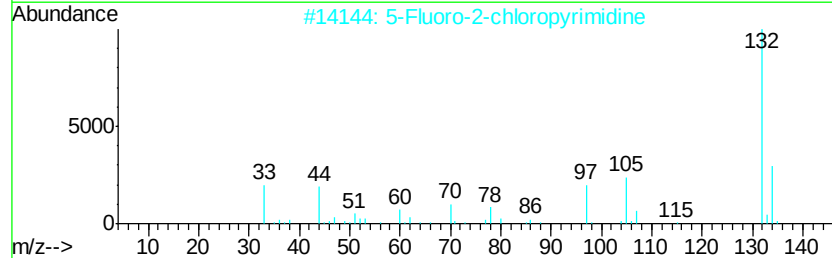
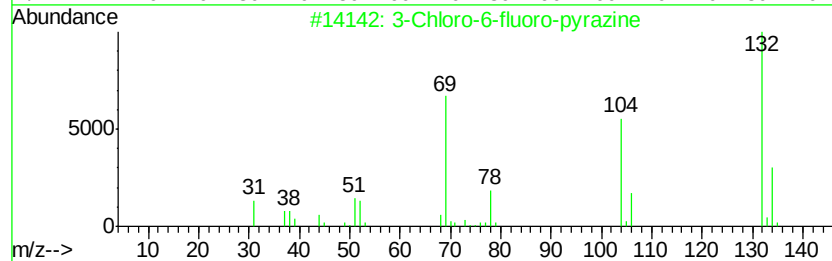
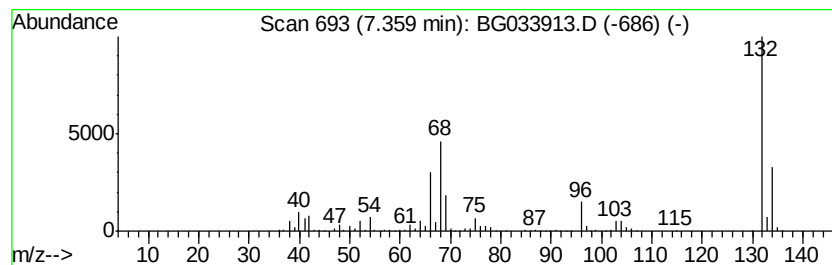
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Peak Number 2 unknown7.36 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.36	64.66 ng	1194310	1,4-Dichlorobenzene-d4	7.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	10
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	9
5		Xenon	132	Xe	007440-63-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.99	4.3	ng	78765	1	7.83	369438	20.0
unknown7.36	7.36	64.7	ng	1194310	1	7.83	369438	20.0