

Data Path : Z:\HPCHEM1\BNA G\DATA\BG032818\
 Data File : BG033403.D
 Acq On : 29 Mar 2018 1:52
 Operator : SJ/JU
 Sample : PB107692BL
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB107692BL

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-BG031918.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	5.779	418	424	432	rBV	45966	76899	1.78%	0.379%
2	6.301	506	513	531	rBV	832779	1581085	36.56%	7.794%
3	7.988	793	800	823	rBV	824817	1780128	41.16%	8.775%
4	8.381	859	867	880	rBV	809765	1629934	37.69%	8.035%
5	8.863	943	949	960	rBV	118766	257096	5.94%	1.267%
6	9.198	998	1006	1020	rBV	611631	1216606	28.13%	5.997%
7	10.068	1146	1154	1171	rBV	459531	1018574	23.55%	5.021%
8	11.748	1431	1440	1455	rBV	187913	378306	8.75%	1.865%
9	14.110	1835	1842	1860	rBV	1526352	2583579	59.74%	12.736%
10	15.485	2069	2076	2088	rBV	349914	594638	13.75%	2.931%
11	16.959	2320	2327	2344	rBV	1230765	2146062	49.62%	10.579%
12	18.217	2534	2541	2551	rBV	474638	813990	18.82%	4.013%
13	20.737	2963	2970	2984	rBV	3013684	4324978	100.00%	21.320%
14	22.559	3273	3280	3291	rBV	481807	945418	21.86%	4.660%
15	26.325	3909	3921	3936	rBV2	276619	938948	21.71%	4.628%

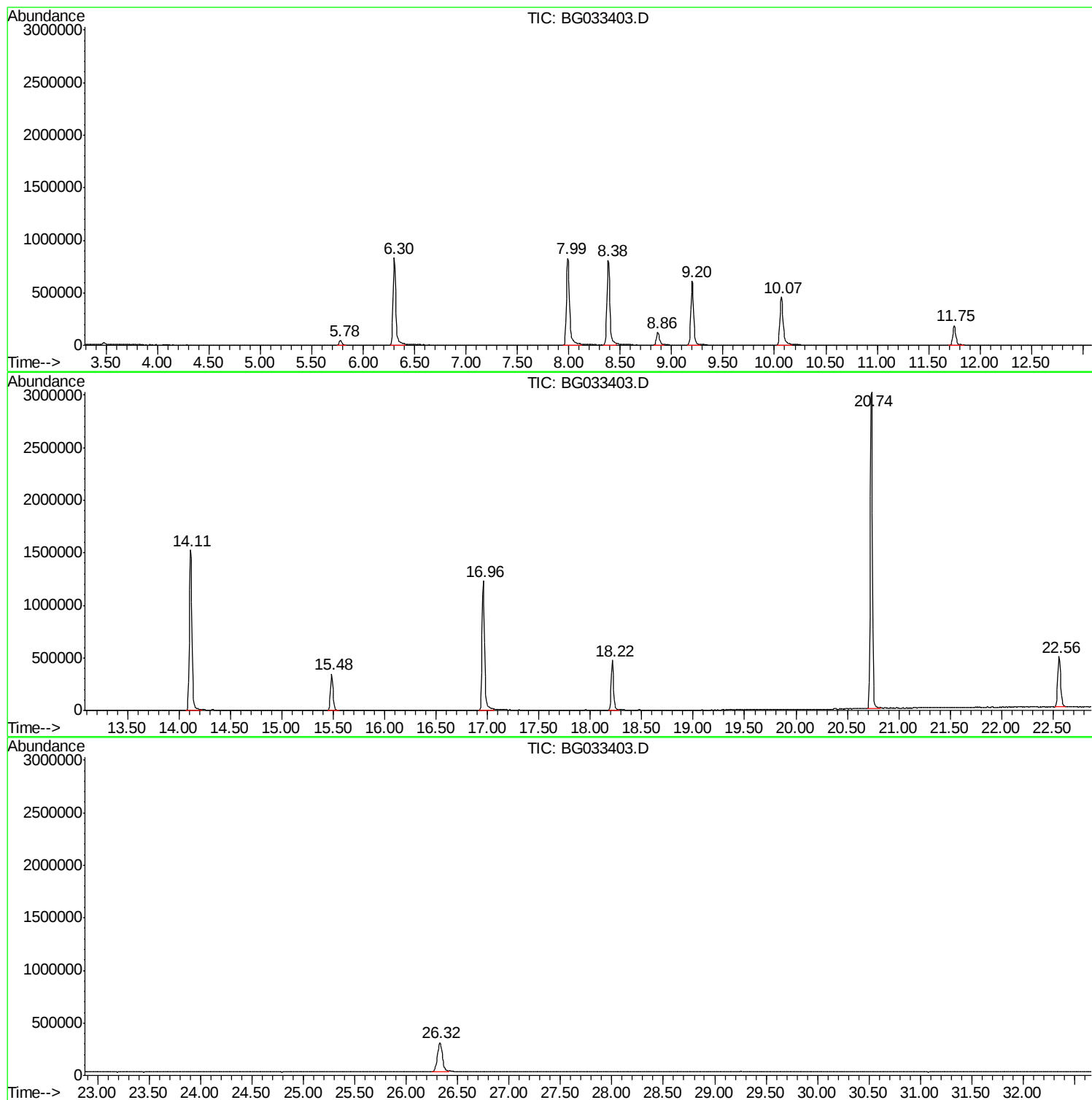
Sum of corrected areas: 20286241

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG031918.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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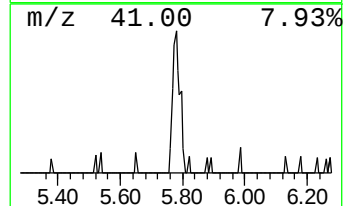
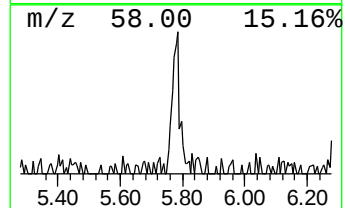
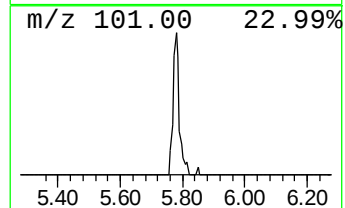
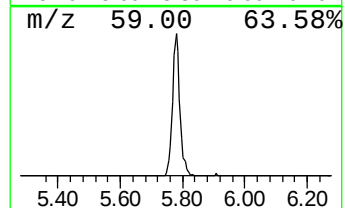
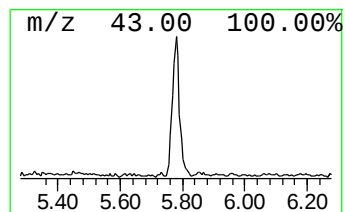
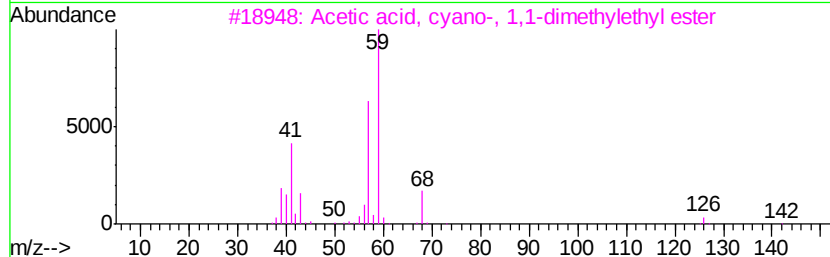
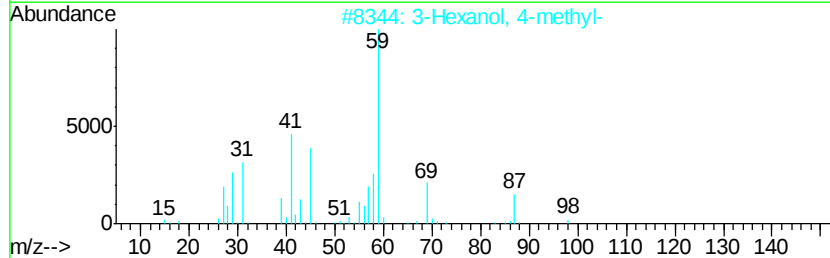
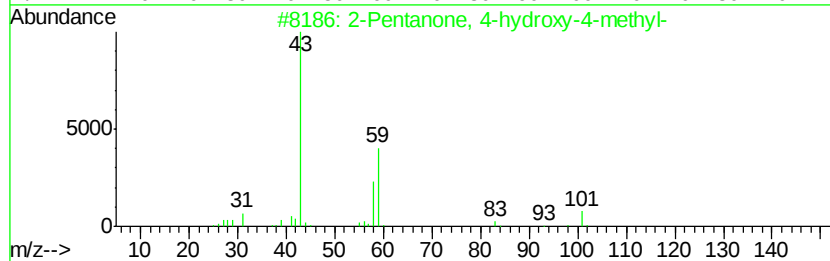
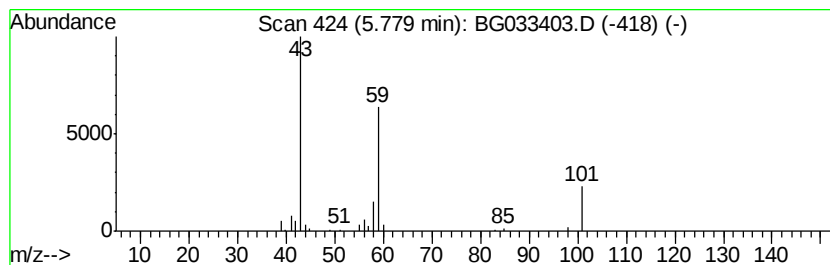
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.78	5.98 ng	76899	1,4-Dichlorobenzene-d4	8.86

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56		
2	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	25		
3	Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23		
4	2-Propyn-1-ol, acetate	98	C5H6O2	000627-09-8	9		
5	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9		



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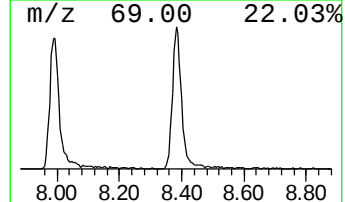
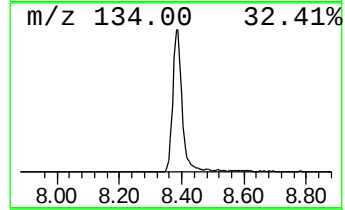
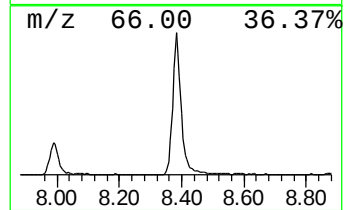
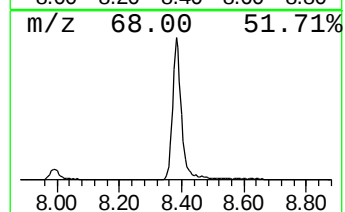
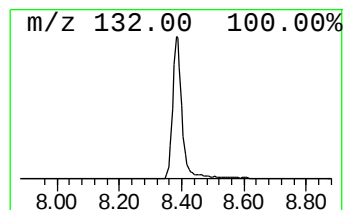
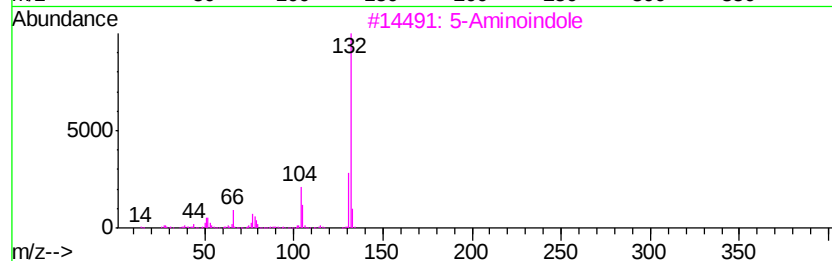
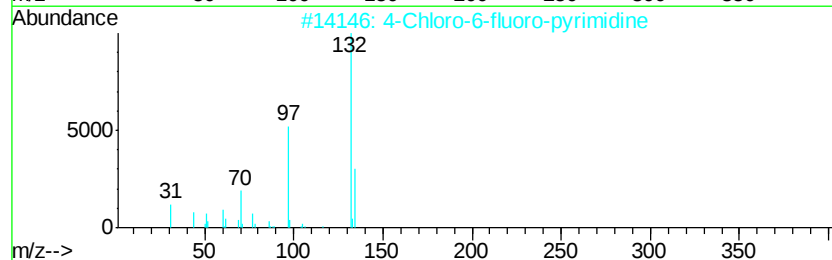
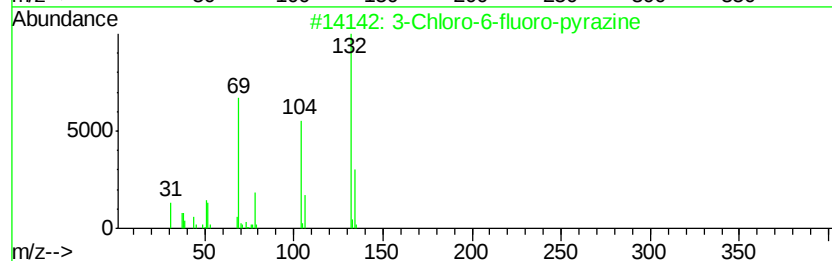
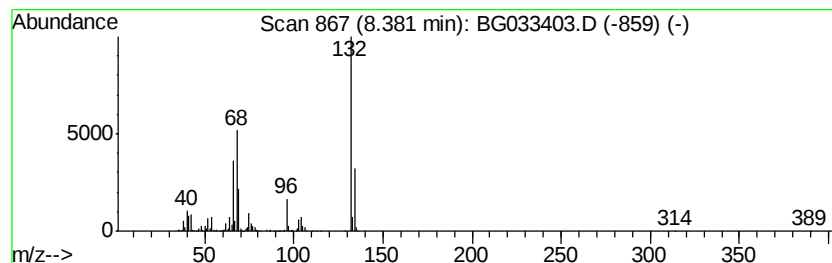
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Peak Number 2 unknown8.38 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.38	126.80 ng	1629930	1,4-Dichlorobenzene-d4	8.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
3		5-Aminoindole	132	C8H8N2	005192-03-0	22
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
5		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.78	6.0	ng	76899	1	8.86	257096	20.0
unknown8.38	8.38	126.8	ng	1629930	1	8.86	257096	20.0