

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG110217\
 Data File : BG030668.D
 Acq On : 2 Nov 2017 18:45
 Operator : SJ/JU
 Sample : PB103908BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB103908BL

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG101617.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	5.233	325	331	339	rVB	138786	214263	4.58%	1.001%
2	5.709	405	412	425	rBV	861679	1379776	29.49%	6.449%
3	7.307	677	684	703	rBV	866946	1499679	32.06%	7.009%
4	7.683	741	748	758	rBV	848073	1476350	31.56%	6.900%
5	8.153	821	828	836	rBV	188891	311933	6.67%	1.458%
6	8.482	873	884	893	rBV	652326	1150674	24.60%	5.378%
7	9.316	1018	1026	1038	rBV	486699	893427	19.10%	4.176%
8	10.967	1300	1307	1320	rBV	254049	477765	10.21%	2.233%
9	13.394	1713	1720	1734	rBV	1670843	2619987	56.00%	12.246%
10	14.769	1947	1954	1964	rBV2	485010	779604	16.66%	3.644%
11	16.249	2199	2206	2220	rBV	1406438	2173533	46.46%	10.159%
12	17.507	2413	2420	2433	rBV2	707602	1066814	22.80%	4.986%
13	20.098	2855	2861	2875	rBV	3692302	4678142	100.00%	21.865%
14	21.778	3140	3147	3155	rBV2	784488	1346557	28.78%	6.294%
15	25.080	3698	3709	3720	rVB	473680	1326667	28.36%	6.201%

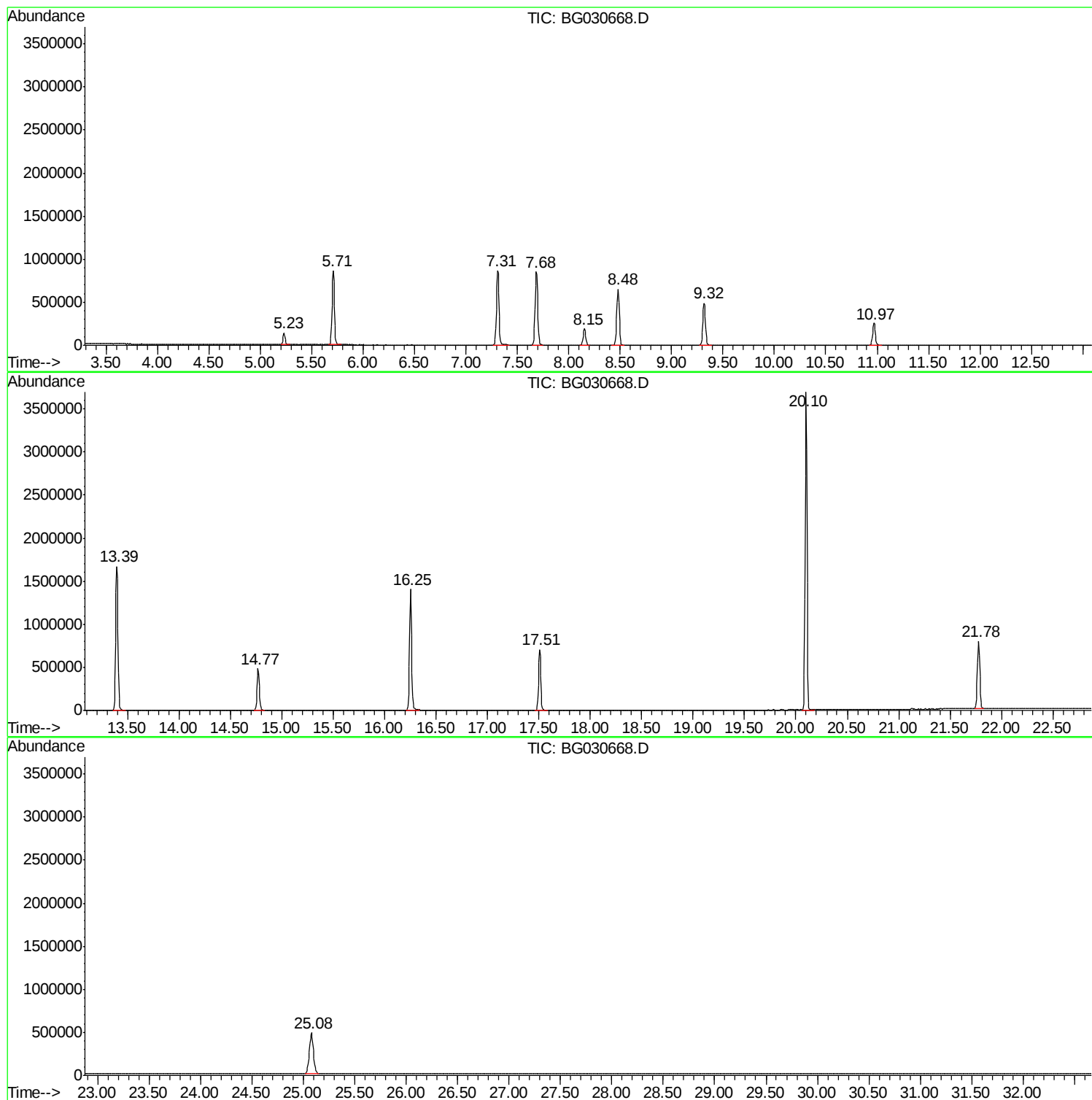
Sum of corrected areas: 21395171

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG101617.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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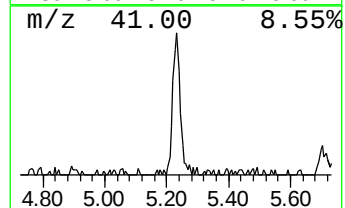
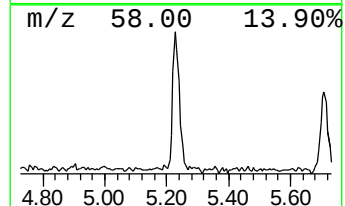
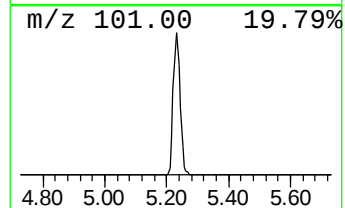
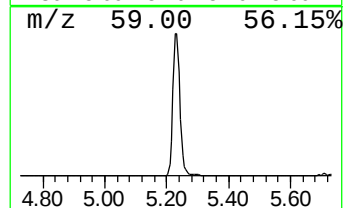
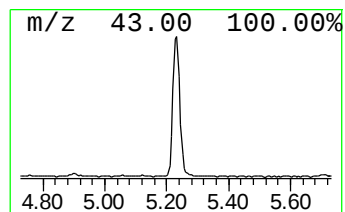
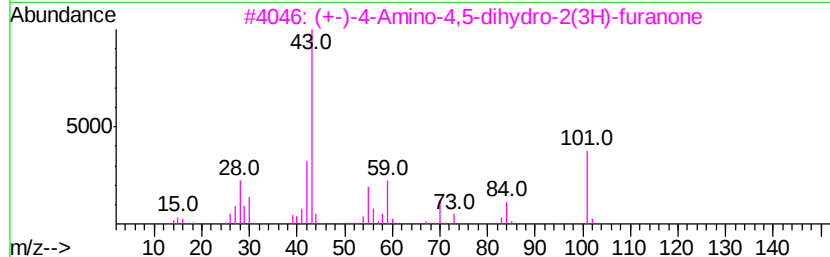
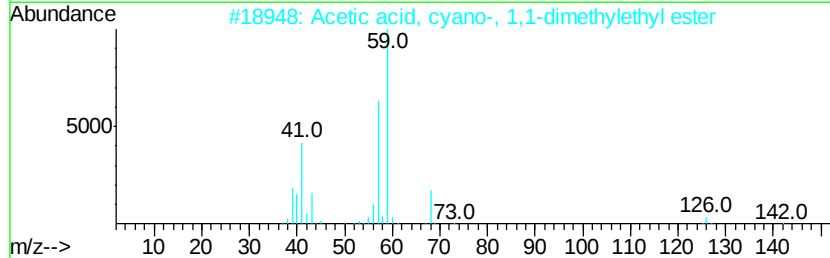
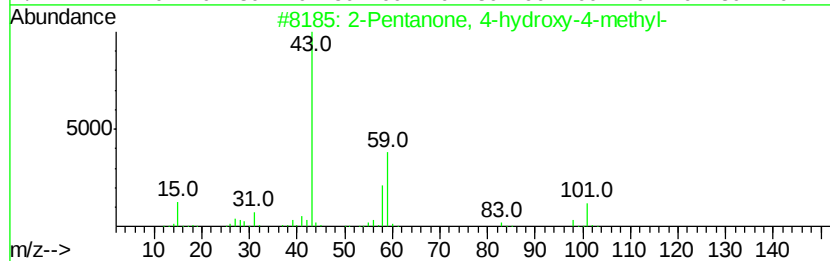
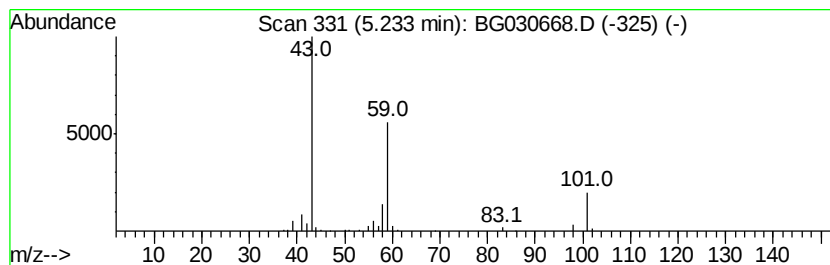
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.23	13.74 ng	214263	1,4-Dichlorobenzene-d4	8.15

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
3			(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9
4			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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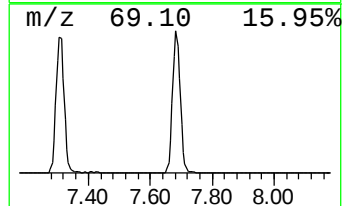
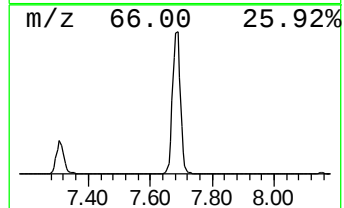
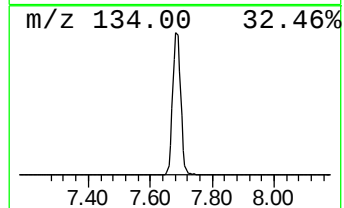
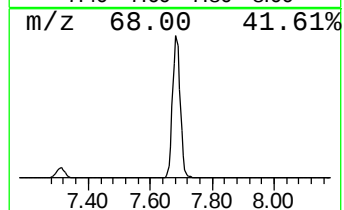
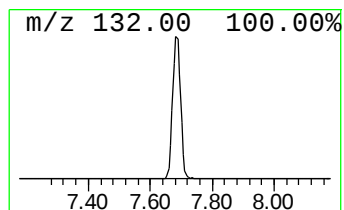
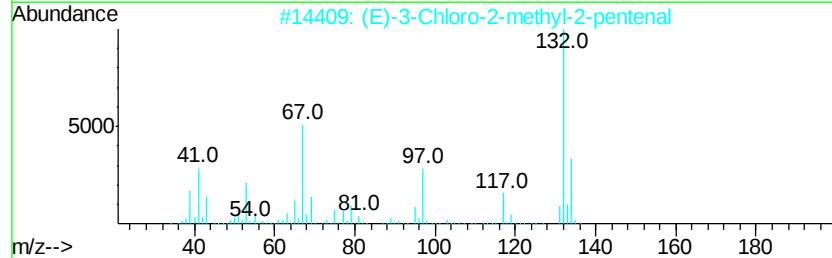
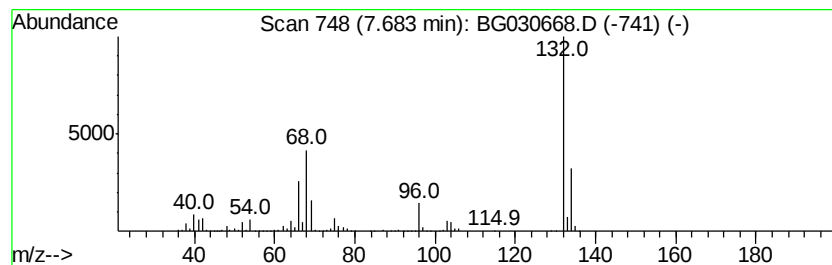
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Peak Number 2 unknown7.68 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.68	94.66 ng	1476350	1,4-Dichlorobenzene-d4	8.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.23	13.7	ng	214263	1	8.15	311933	20.0
unknown7.68	7.68	94.7	ng	1476350	1	8.15	311933	20.0