

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG110417\
 Data File : BG030713.D
 Acq On : 3 Nov 2017 23:24
 Operator : SJ/JU
 Sample : PB103952BL
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB103952BL

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-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	5.227	325	330	340	rVB	151101	233770	4.82%	0.991%
2	5.709	405	412	423	rBV	894097	1448470	29.85%	6.140%
3	7.307	677	684	699	rBV	938747	1612803	33.24%	6.836%
4	7.683	741	748	758	rBV	931186	1576143	32.49%	6.681%
5	8.153	820	828	835	rBV	216518	380134	7.83%	1.611%
6	8.476	876	883	891	rBV	691500	1212126	24.98%	5.138%
7	9.316	1018	1026	1033	rBV	532583	958039	19.75%	4.061%
8	10.967	1299	1307	1319	rBV	317994	592418	12.21%	2.511%
9	13.394	1713	1720	1741	rVB	1799924	2840778	58.55%	12.041%
10	14.769	1945	1954	1963	rBV2	601719	991583	20.44%	4.203%
11	16.250	2199	2206	2226	rBV	1440439	2333798	48.10%	9.892%
12	17.507	2412	2420	2432	rBV2	882142	1337446	27.57%	5.669%
13	20.098	2855	2861	2872	rBV	3691756	4851910	100.00%	20.565%
14	21.778	3140	3147	3160	rVB	960560	1609767	33.18%	6.823%
15	25.074	3698	3708	3723	rVB2	557389	1613380	33.25%	6.839%

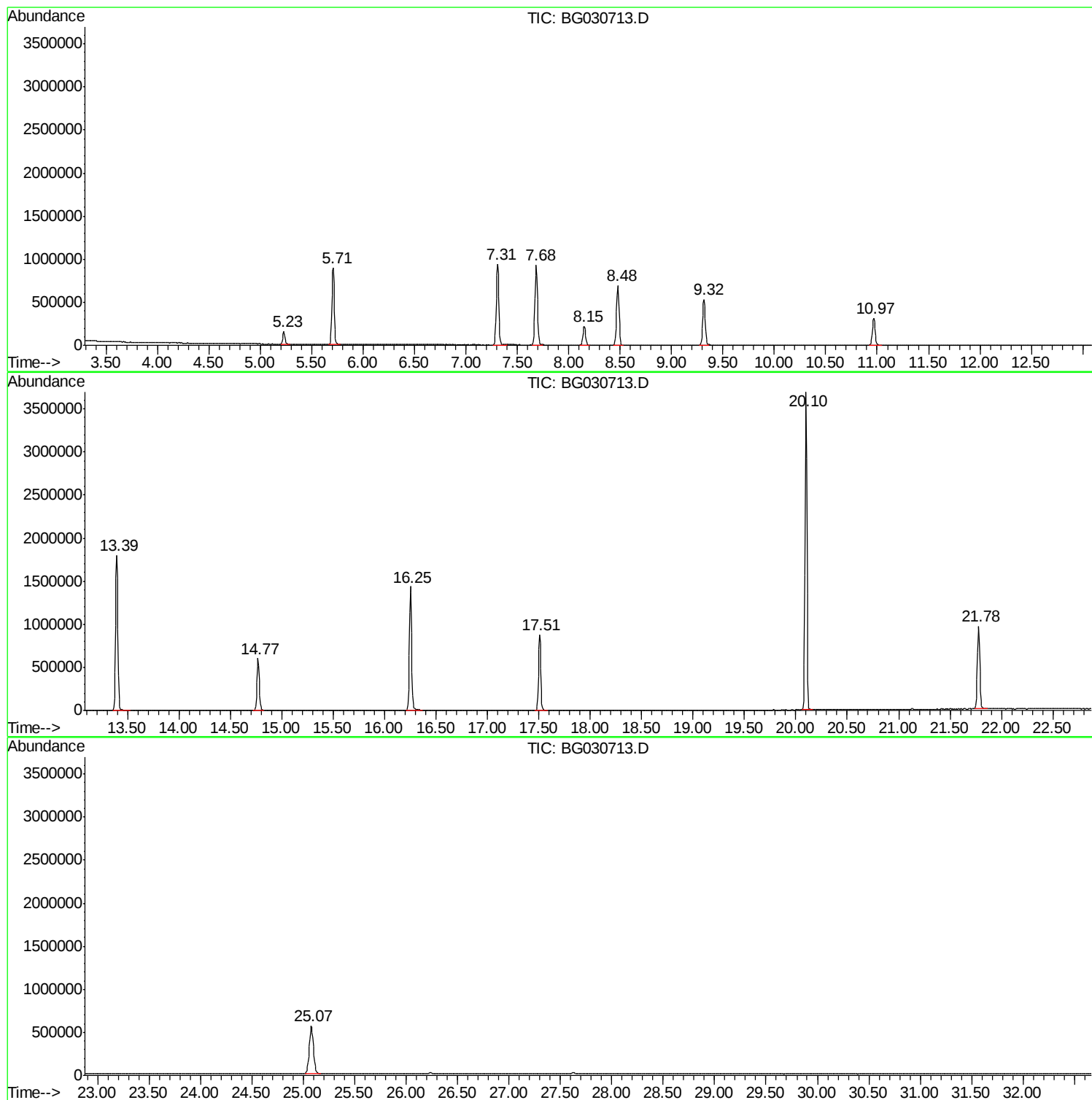
Sum of corrected areas: 23592565

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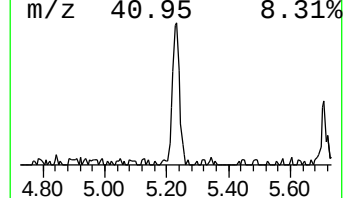
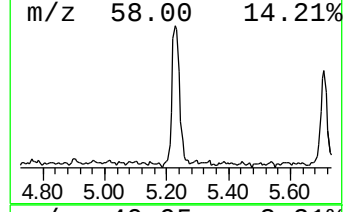
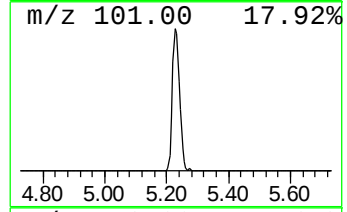
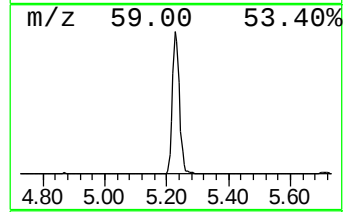
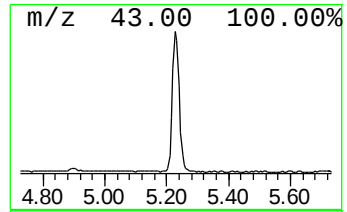
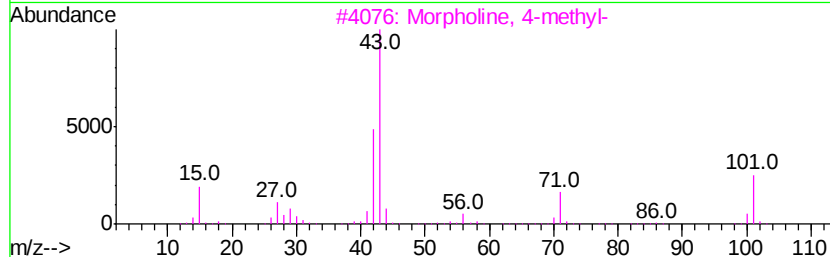
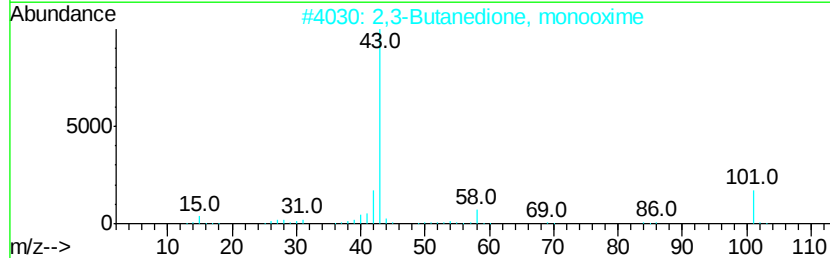
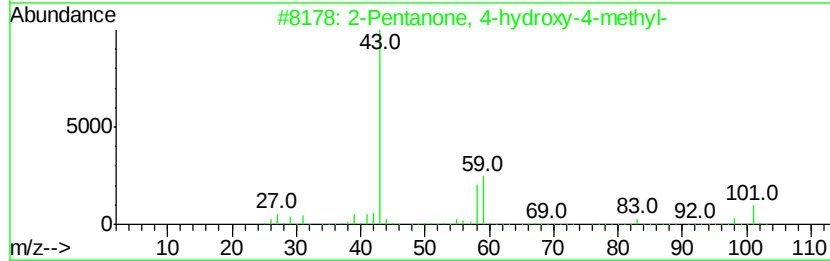
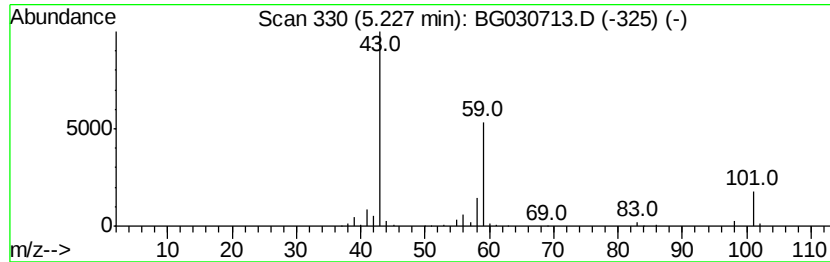
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.23	12.30 ng	233770	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	50
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
3			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4			Acetamide, N-(aminocarbonyl)-	102	C3H6N2O2	000591-07-1	9
5			5-Aminovaleric acid	117	C5H11NO2	000660-88-8	9



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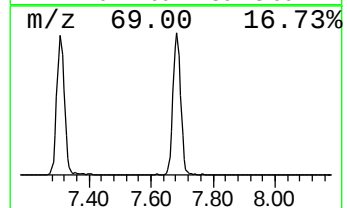
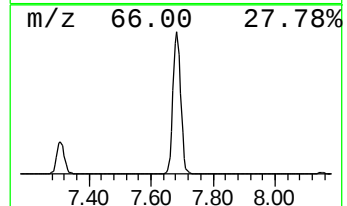
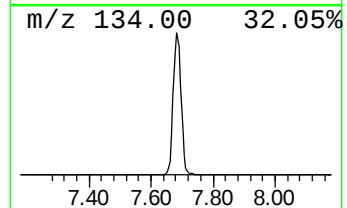
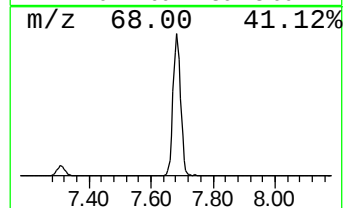
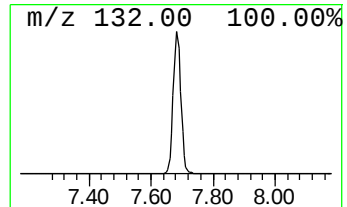
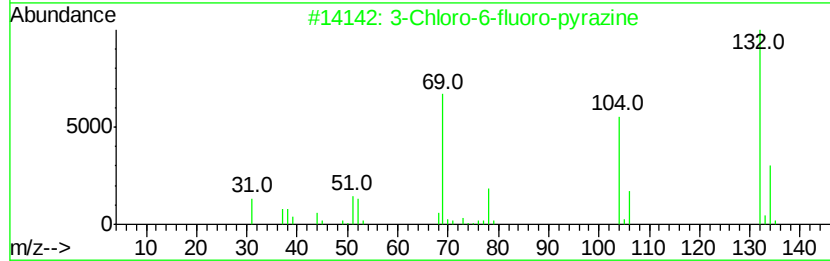
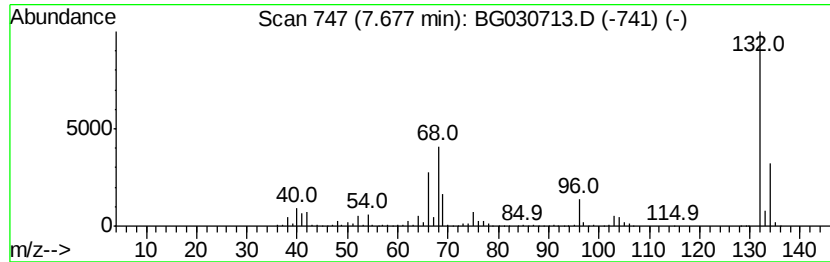
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Peak Number 2 unknown7.68 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.68	82.93 ng	1576140	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		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.23	12.3	ng	233770	1	8.15	380134	20.0
unknown7.68	7.68	82.9	ng	1576140	1	8.15	380134	20.0