

Data Path : Z:\HPCHEM1\BNA G\DATA\BG102817\
 Data File : BG030544.D
 Acq On : 29 Oct 2017 2:41
 Operator : SJ/JU
 Sample : PB103762TB
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB103762TB

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	5.241	327	332	342	rVB	125307	189083	5.39%	1.039%
2	5.717	406	413	426	rVB	747713	1170000	33.33%	6.429%
3	7.315	678	685	698	rBV	712504	1210679	34.49%	6.652%
4	7.697	742	750	757	rBV	707790	1219210	34.73%	6.699%
5	8.161	823	829	839	rVB	218563	375058	10.68%	2.061%
6	8.490	878	885	895	rBV	547936	949644	27.05%	5.218%
7	9.330	1020	1028	1037	rBV	402180	715526	20.38%	3.932%
8	10.981	1299	1309	1316	rBV	290312	527002	15.01%	2.896%
9	13.408	1715	1722	1732	rBV	1288157	1953906	55.66%	10.736%
10	14.783	1949	1956	1969	rBV2	504980	811969	23.13%	4.462%
11	16.263	2201	2208	2225	rBV	1061092	1679996	47.86%	9.231%
12	17.520	2415	2422	2431	rBV	778092	1149548	32.75%	6.316%
13	20.106	2856	2862	2872	rBV	2712372	3510389	100.00%	19.288%
14	21.792	3142	3149	3159	rVB2	813241	1375896	39.19%	7.560%
15	25.100	3701	3712	3732	rVB	467957	1361530	38.79%	7.481%

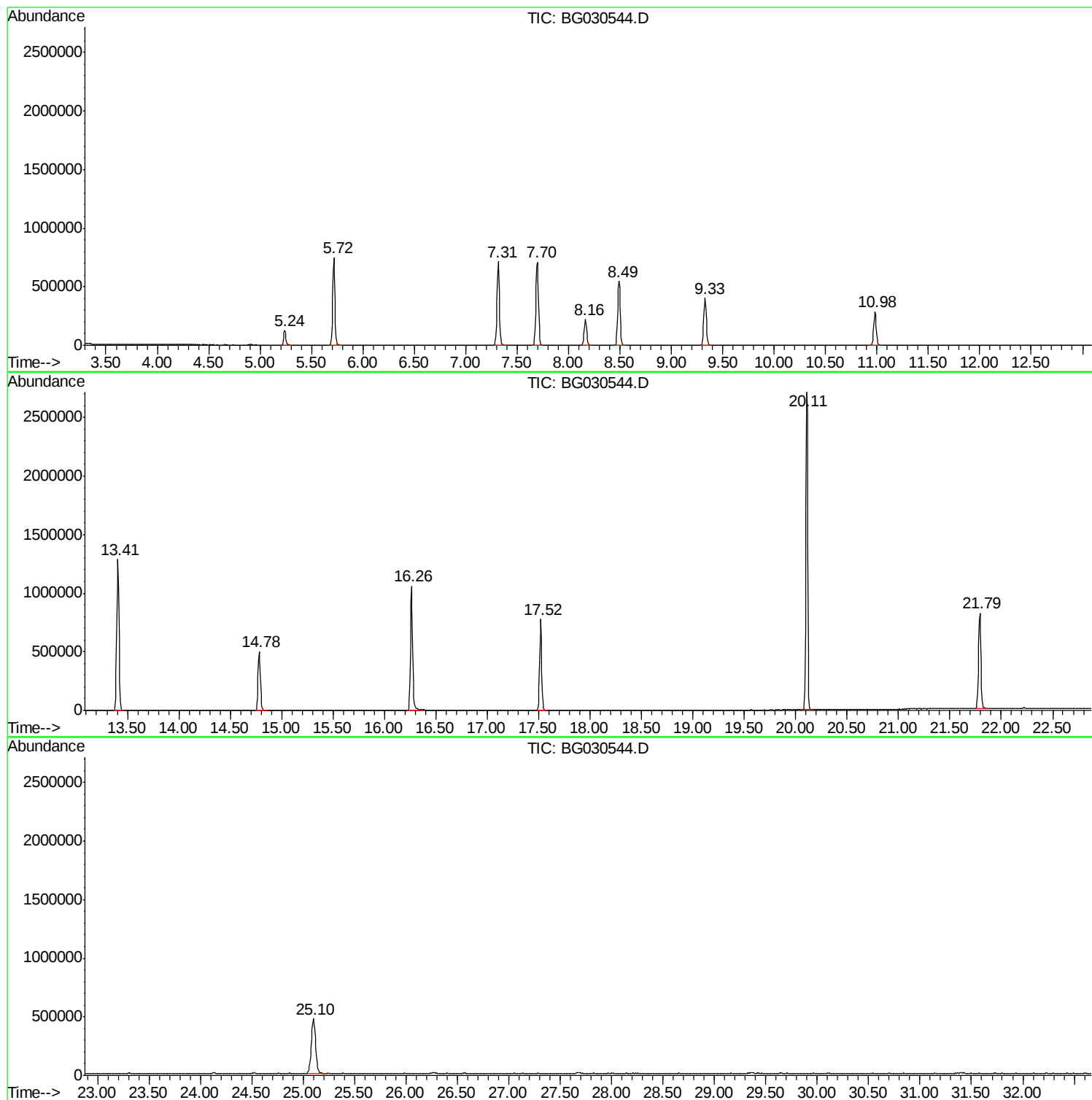
Sum of corrected areas: 18199436

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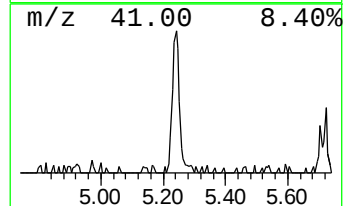
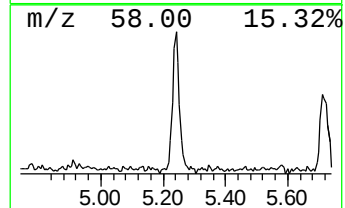
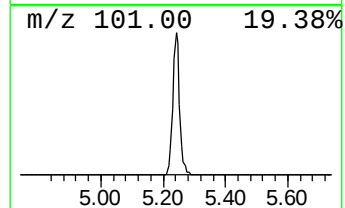
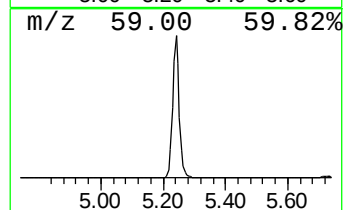
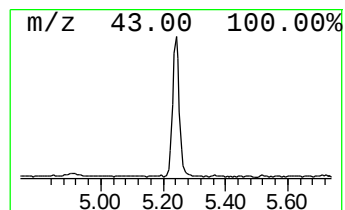
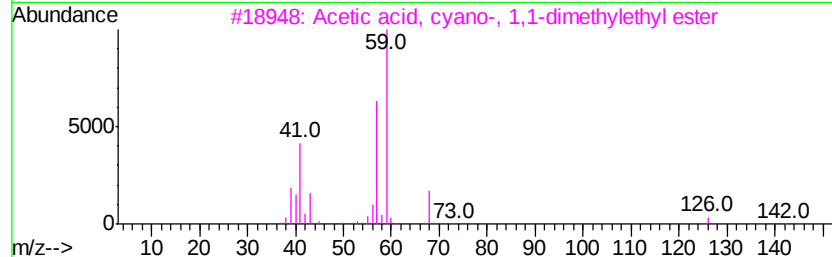
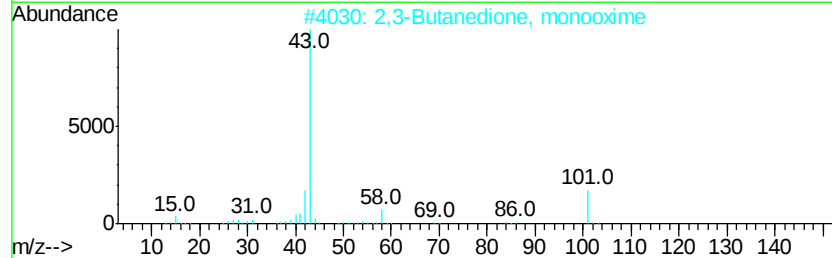
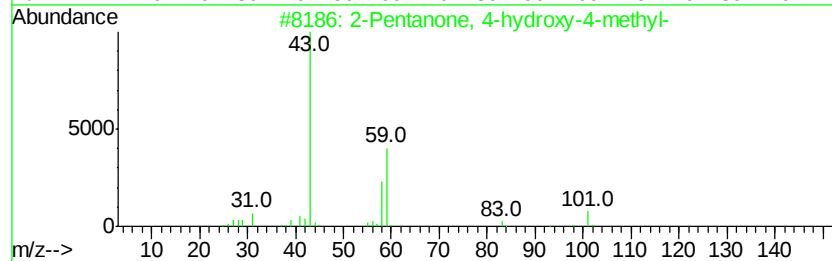
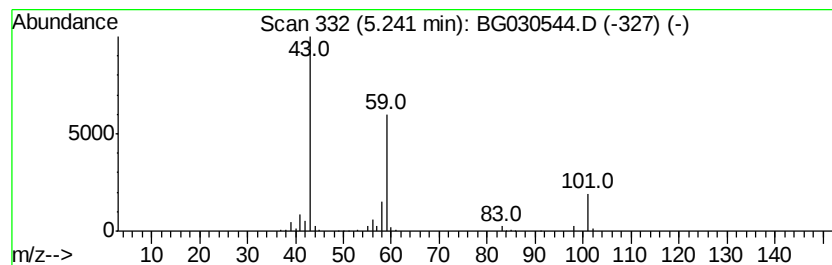
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.24	10.08 ng	189083	1,4-Dichlorobenzene-d4	8.17

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
3			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
4			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5			Acetone	58	C3H6O	000067-64-1	9



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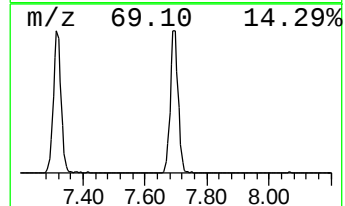
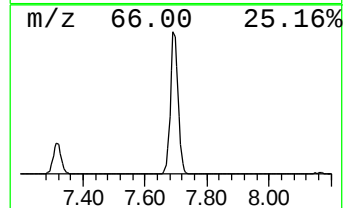
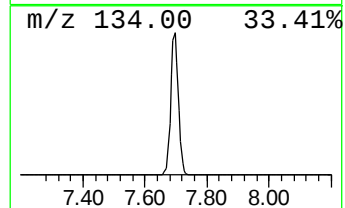
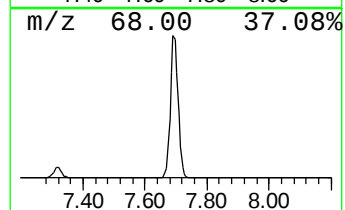
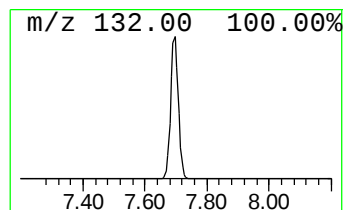
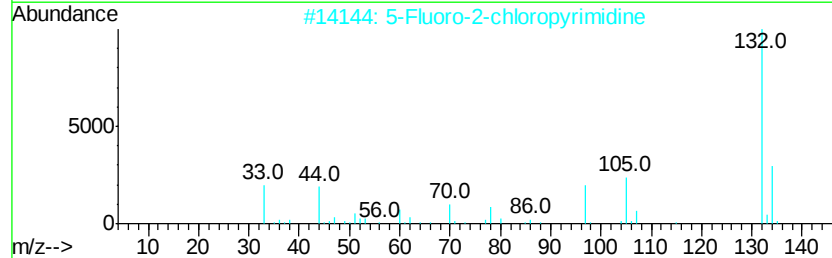
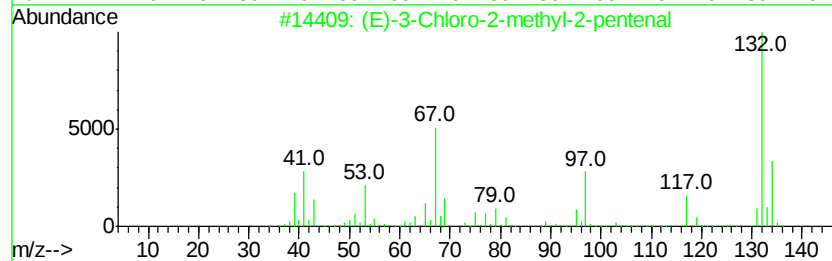
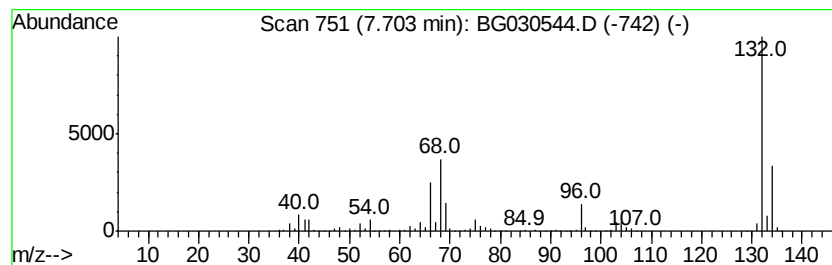
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Peak Number 2 unknown7.70 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.70	65.01 ng	1219210	1,4-Dichlorobenzene-d4	8.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		N-(-phenylisopropyl)benzalimine	223	C16H17N	1000379-01-3	10
5		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.24	10.1	ng	189083	1	8.17	375058	20.0
unknown7.70	7.70	65.0	ng	1219210	1	8.17	375058	20.0