

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG111017\
 Data File : BG030926.D
 Acq On : 11 Nov 2017 3:22
 Operator : SJ/JU
 Sample : PB104155BL
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB104155BL

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-BG111017.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	5.214	322	328	337	rBV	76425	122628	4.10%	0.949%
2	5.702	404	411	426	rBV	470538	777073	25.96%	6.015%
3	7.306	677	684	701	rBV	444034	801223	26.76%	6.202%
4	7.676	738	747	759	rBV	479220	822249	27.47%	6.365%
5	8.140	816	826	834	rBV	125539	213630	7.14%	1.654%
6	8.469	875	882	893	rBV	365322	662476	22.13%	5.128%
7	9.310	1017	1025	1037	rBV	247596	475809	15.89%	3.683%
8	10.961	1298	1306	1317	rBV	144641	289157	9.66%	2.238%
9	13.387	1712	1719	1727	rBV	840194	1352010	45.16%	10.466%
10	14.762	1946	1953	1970	rBV2	252785	445040	14.87%	3.445%
11	16.249	2199	2206	2226	rBV	705468	1193170	39.86%	9.236%
12	17.506	2413	2420	2435	rBV	397003	655768	21.91%	5.076%
13	20.091	2854	2860	2869	rBV	2229004	2993658	100.00%	23.174%
14	21.777	3139	3147	3157	rBV	589833	1052888	35.17%	8.150%
15	25.079	3698	3709	3722	rBV2	342134	1061570	35.46%	8.218%

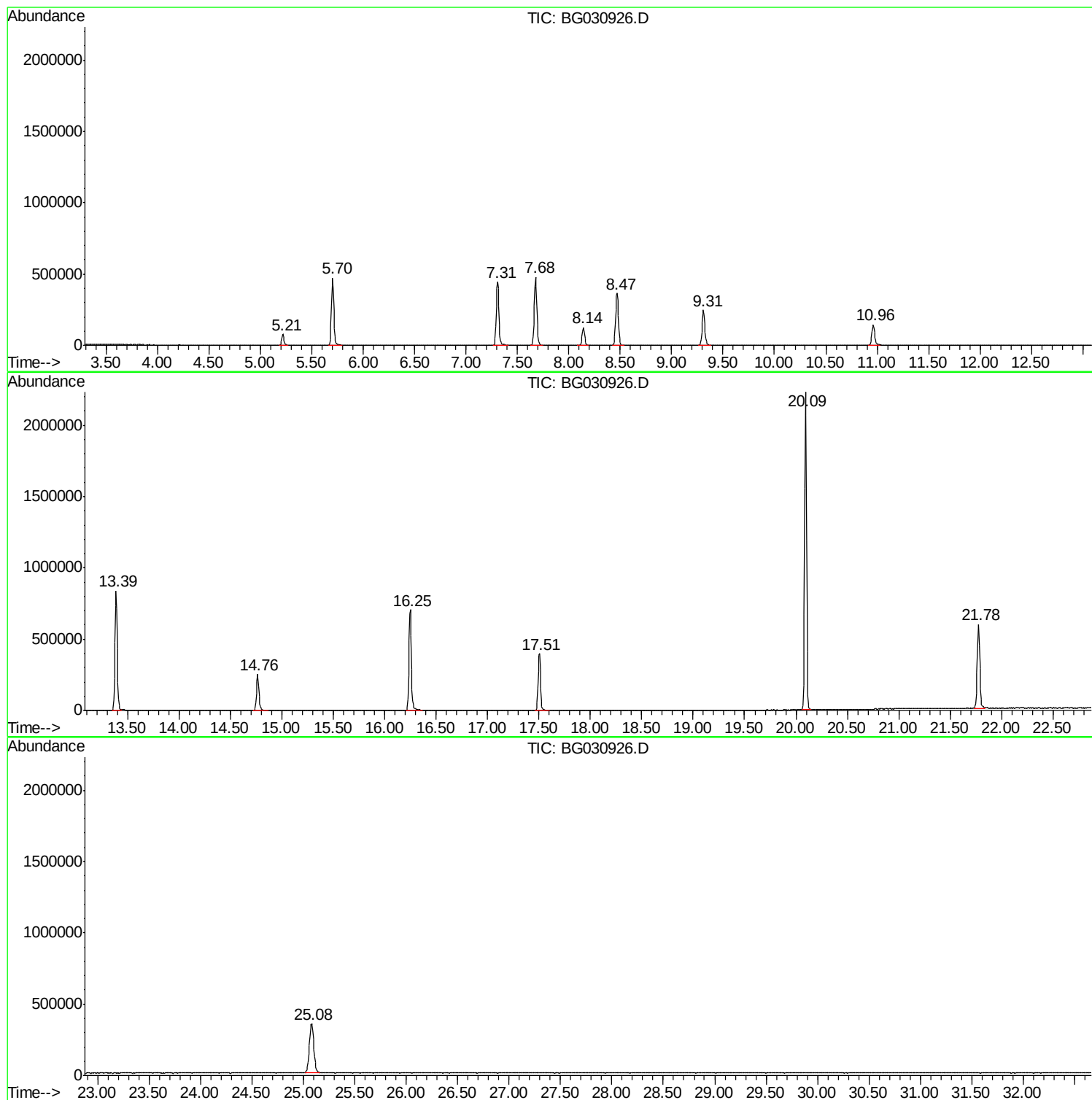
Sum of corrected areas: 12918349

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG111017.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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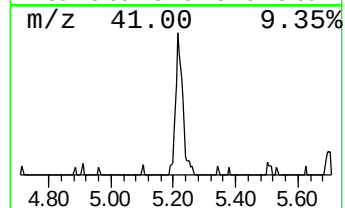
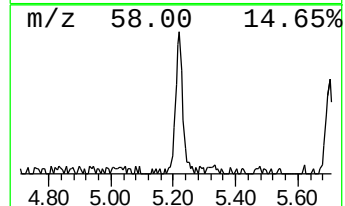
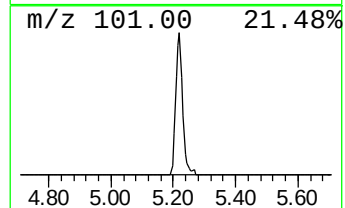
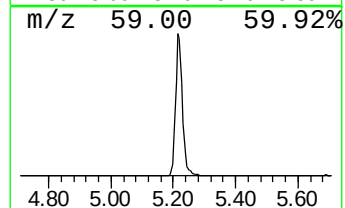
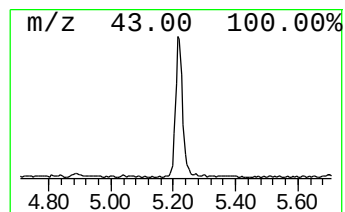
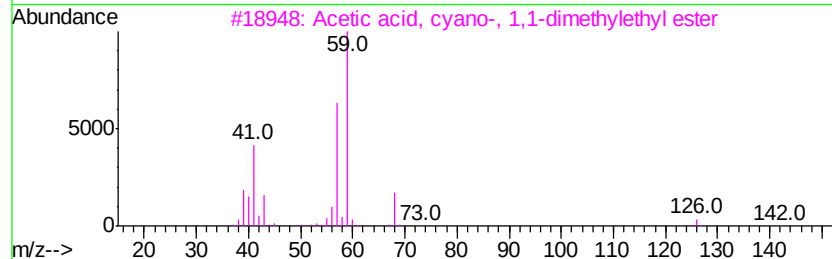
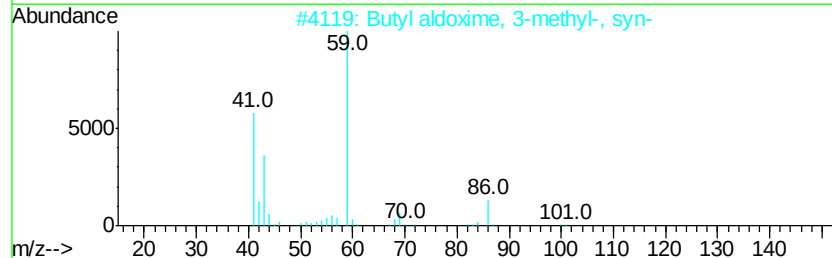
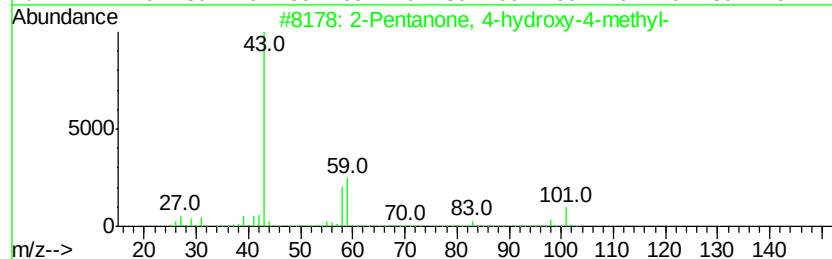
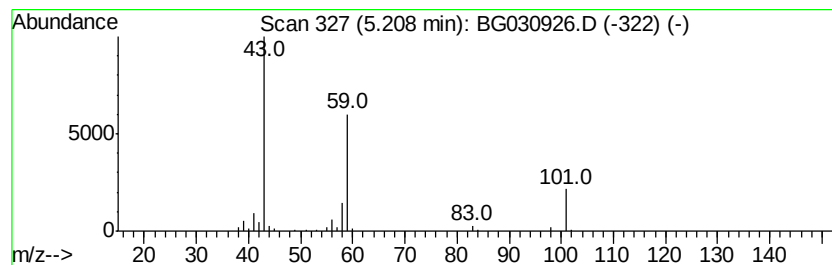
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.21	11.48 ng	122628	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	45	
2	Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	32	
3	Acetic acid, cyano-, 1,1-dimethv...	141	C7H11NO2	001116-98-9	23	
4	Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9	
5	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9	



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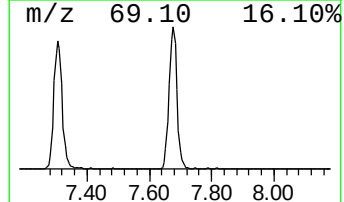
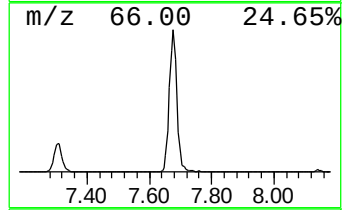
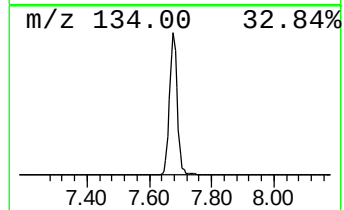
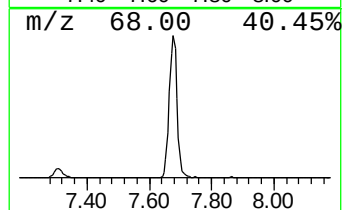
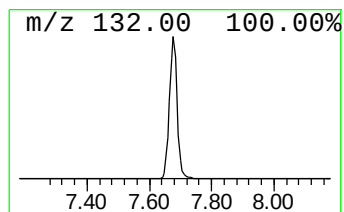
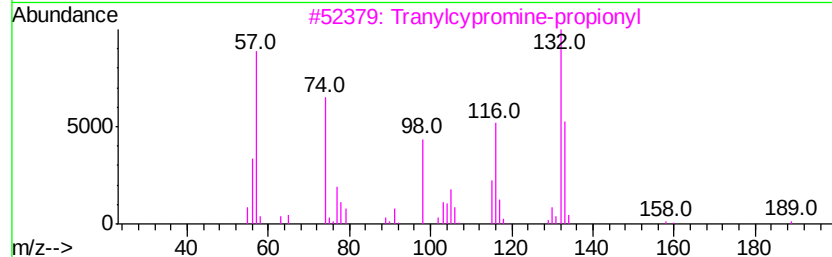
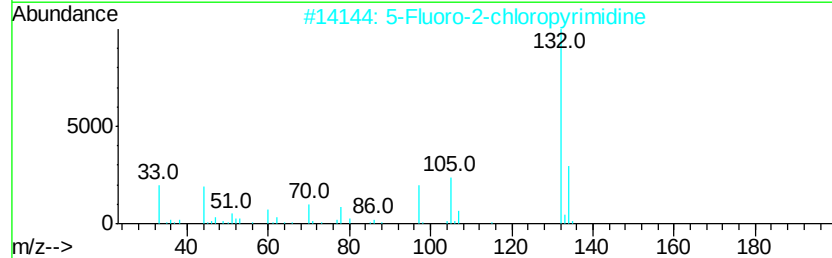
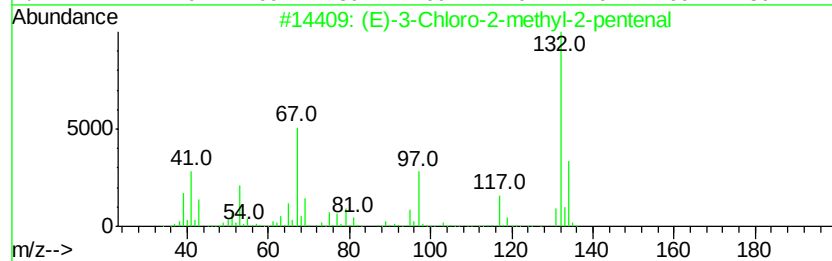
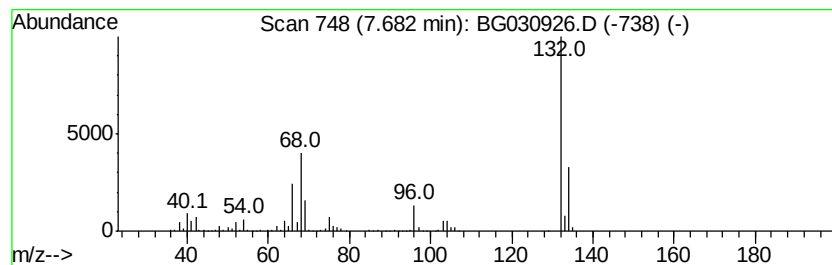
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Peak Number 2 unknown7.68 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.68	76.98 ng	822249	1,4-Dichlorobenzene-d4	8.15

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	12
3		Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	10
4		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.21	11.5	ng	122628	1	8.15	213630	20.0
unknown7.68	7.68	77.0	ng	822249	1	8.15	213630	20.0