

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG121217\
 Data File : BG031504.D
 Acq On : 12 Dec 2017 23:18
 Operator : SJ/JU
 Sample : I6763-13
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 120517-FOPI-F-U-B

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-BG121117.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	5.185	317	323	336	rBV	56806	94998	1.46%	0.431%
2	5.667	399	405	428	rVB	300318	573715	8.80%	2.601%
3	7.277	673	679	693	rBV	195821	389287	5.97%	1.765%
4	7.641	734	741	759	rBV	609003	1099882	16.87%	4.987%
5	8.111	814	821	835	rVB	194229	343974	5.28%	1.559%
6	8.434	868	876	888	rBV	879304	1570329	24.09%	7.119%
7	9.275	1011	1019	1032	rBV	610827	1212040	18.59%	5.495%
8	10.926	1293	1300	1319	rBV	223138	487447	7.48%	2.210%
9	13.352	1705	1713	1738	rBV	2090367	3507873	53.81%	15.904%
10	14.733	1938	1948	1963	rBV2	422493	786212	12.06%	3.564%
11	16.219	2194	2201	2218	rBV	1088414	1763766	27.06%	7.996%
12	17.477	2408	2415	2429	rBV2	638295	1055756	16.19%	4.787%
13	20.068	2849	2856	2870	rBV	4940899	6519149	100.00%	29.556%
14	21.748	3135	3142	3155	rBV2	751806	1378988	21.15%	6.252%
15	25.032	3689	3701	3715	rBV2	417135	1273335	19.53%	5.773%

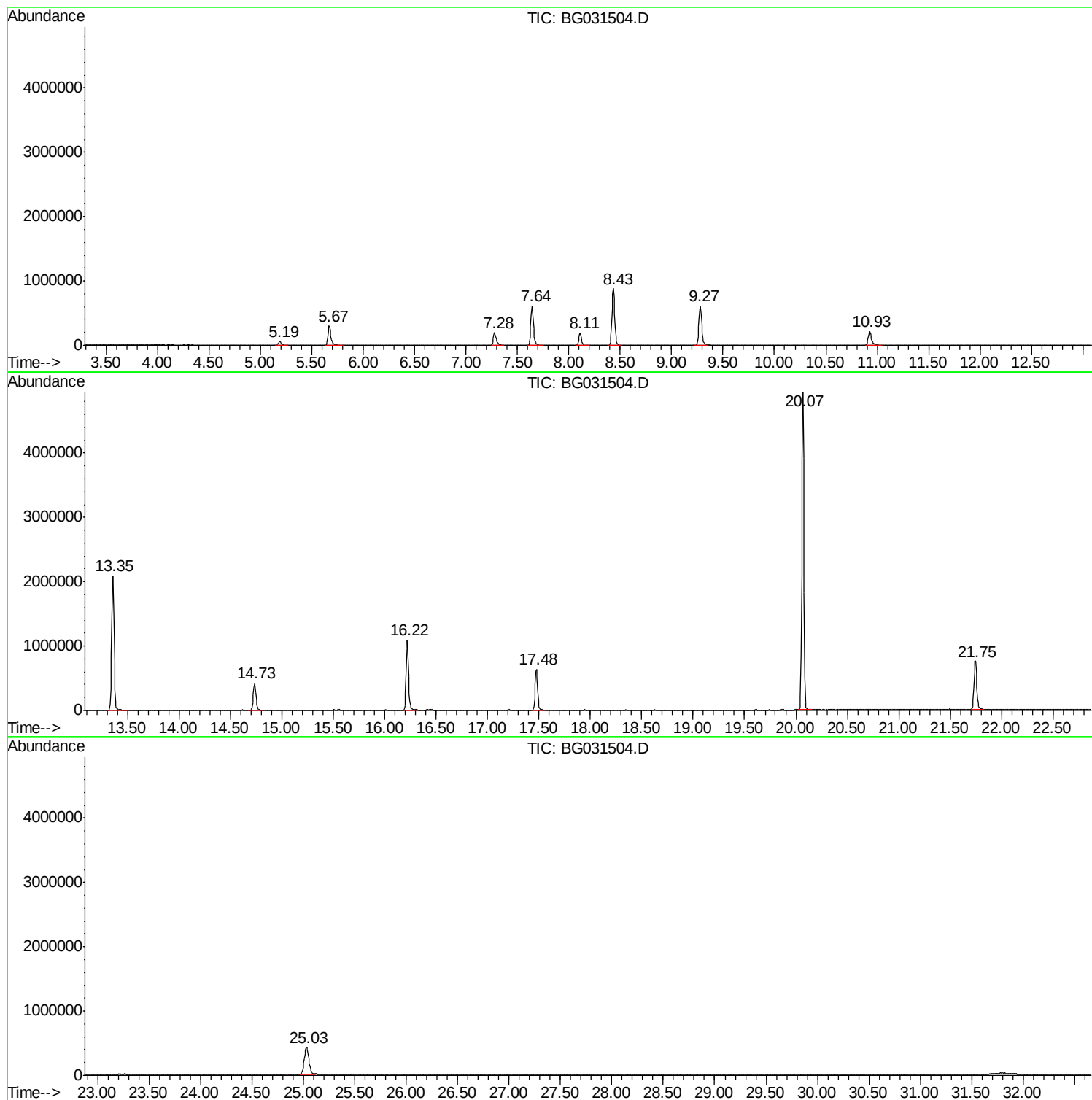
Sum of corrected areas: 22056751

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG121117.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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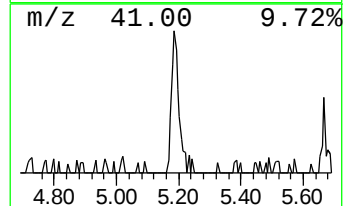
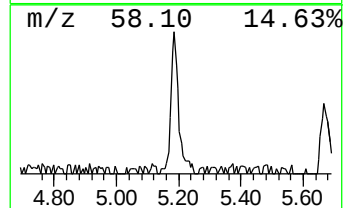
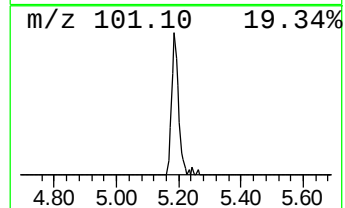
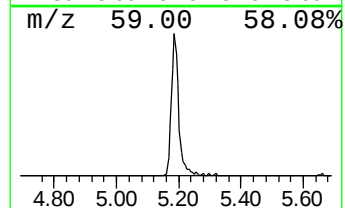
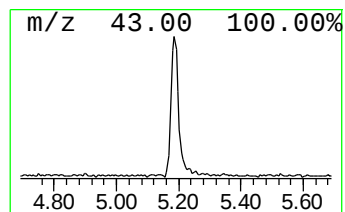
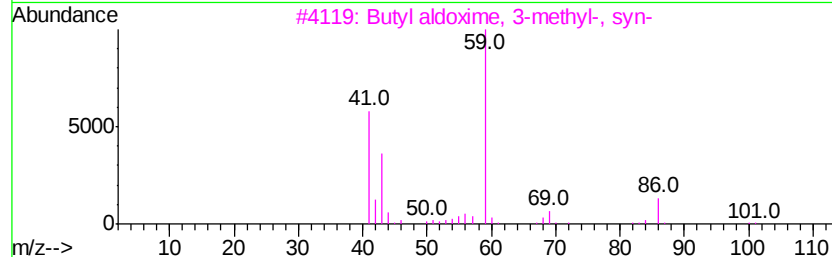
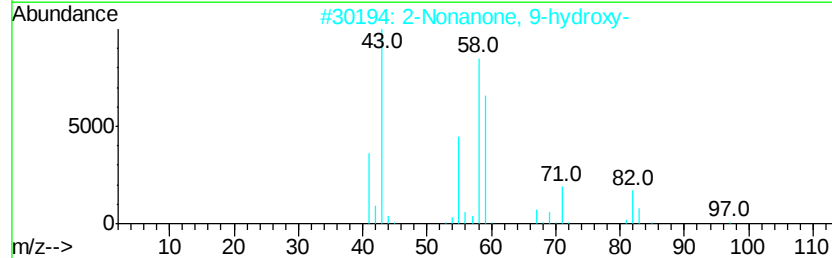
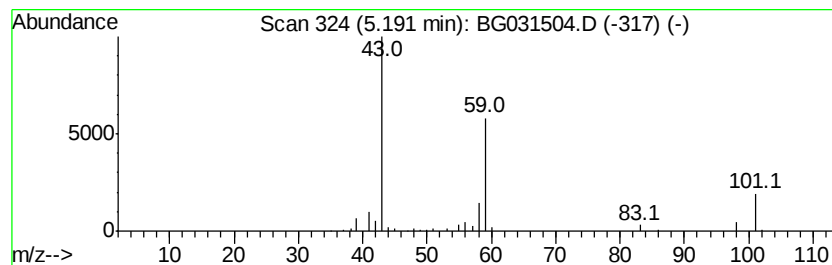
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.19	5.52 ng	94998	1,4-Dichlorobenzene-d4	8.11

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2			2-Nonanone, 9-hydroxy-	158	C9H18O2	025368-56-3	33
3			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	32
4			Acetone	58	C3H6O	000067-64-1	9
5			5-Hexen-2-one	98	C6H10O	000109-49-9	9



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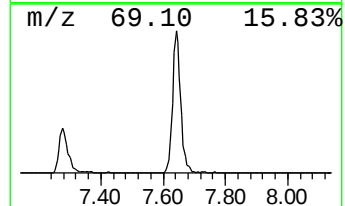
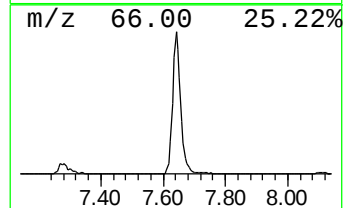
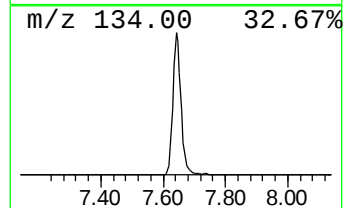
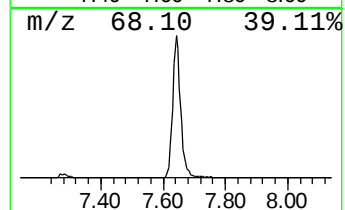
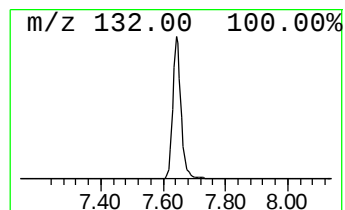
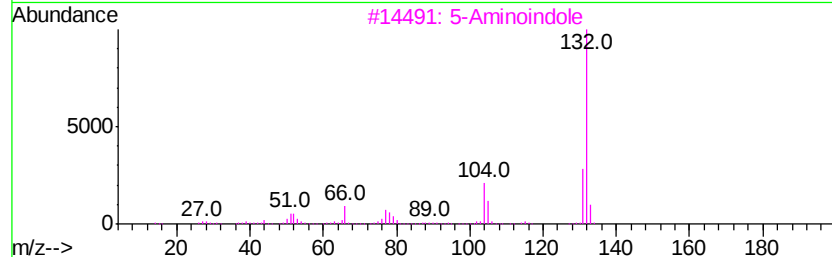
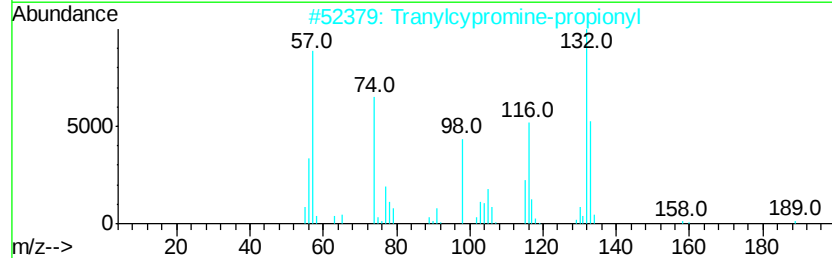
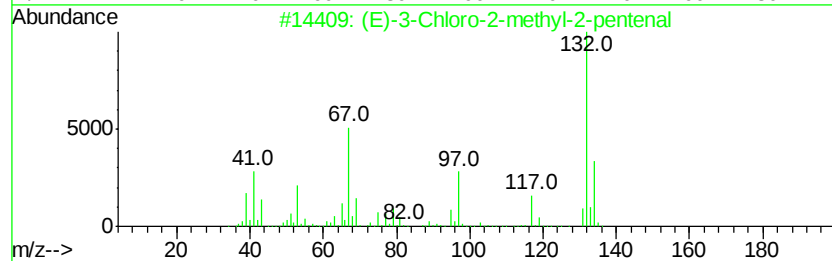
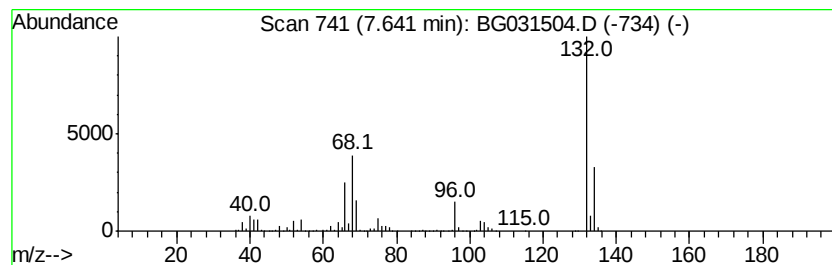
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Peak Number 2 unknown7.64 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.64	63.95 ng	1099880	1,4-Dichlorobenzene-d4	8.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		5-Aminoindole	132	C8H8N2	005192-03-0	10
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.19	5.5	ng	94998	1	8.11	343974	20.0
unknown7.64	7.64	64.0	ng	1099880	1	8.11	343974	20.0