

Data Path : Z:\HPCHEM1\BNA G\DATA\BG120117\
 Data File : BG031325.D
 Acq On : 1 Dec 2017 19:39
 Operator : SJ/JU
 Sample : I6671-06
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 ENV-106-1(40-50)

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	5.201	320	325	339	rBV	59044	103852	2.19%	0.459%
2	5.688	401	408	424	rBV	914195	1586278	33.44%	7.016%
3	7.292	674	681	705	rBV	923856	1729040	36.45%	7.647%
4	7.662	736	744	755	rBV	970871	1697193	35.78%	7.506%
5	8.127	816	823	831	rBV	189817	334644	7.06%	1.480%
6	8.450	869	878	893	rBV	676796	1254331	26.45%	5.548%
7	9.296	1015	1022	1039	rBV	475378	972468	20.50%	4.301%
8	10.947	1296	1303	1312	rBV	221929	467573	9.86%	2.068%
9	13.373	1708	1716	1728	rBV	1638214	2626979	55.39%	11.618%
10	14.202	1850	1857	1870	rBV	82621	187957	3.96%	0.831%
11	14.754	1940	1951	1965	rBV2	424442	758944	16.00%	3.357%
12	16.235	2197	2203	2223	rBV	1358714	2323054	48.98%	10.274%
13	17.492	2410	2417	2434	rBV2	664423	1110719	23.42%	4.912%
14	20.083	2852	2858	2870	rBV	3610538	4743020	100.00%	20.977%
15	21.763	3137	3144	3159	rBV2	795368	1405343	29.63%	6.215%
16	25.060	3693	3705	3719	rBV2	427738	1309000	27.60%	5.789%

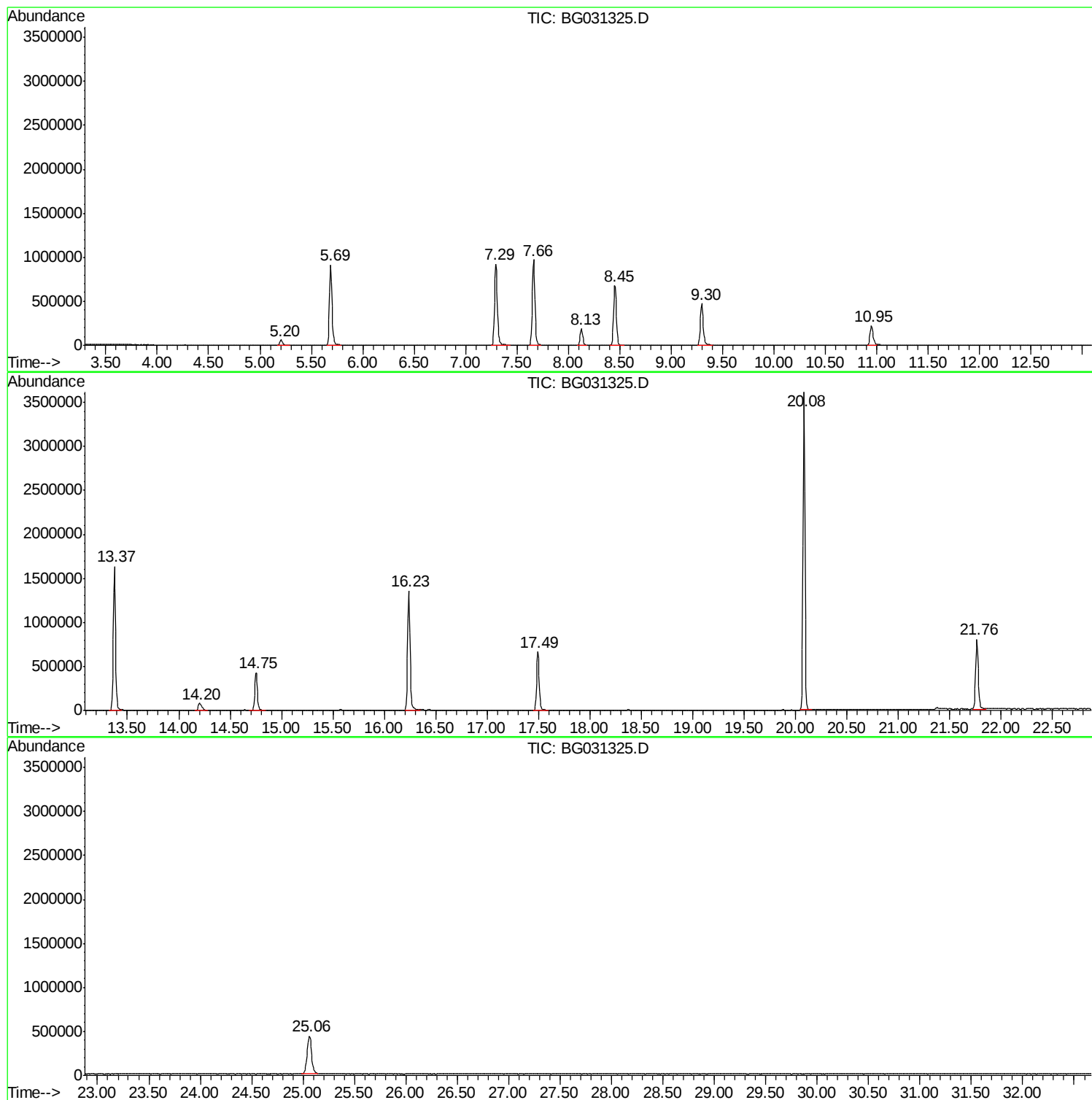
Sum of corrected areas: 22610395

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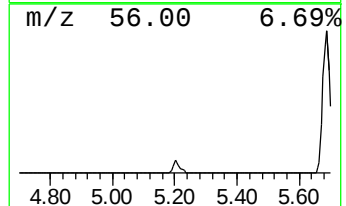
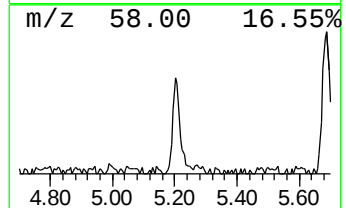
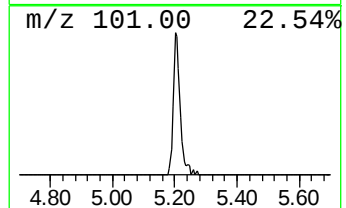
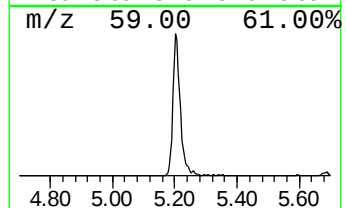
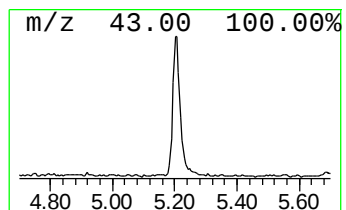
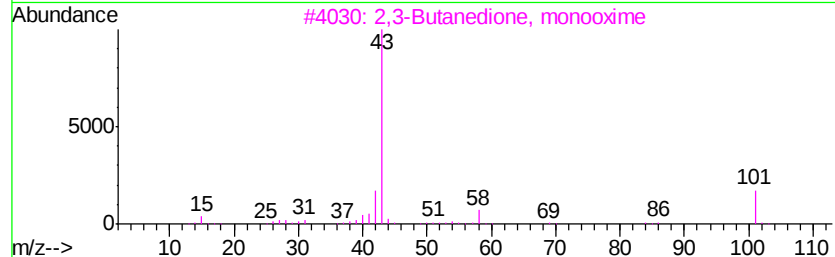
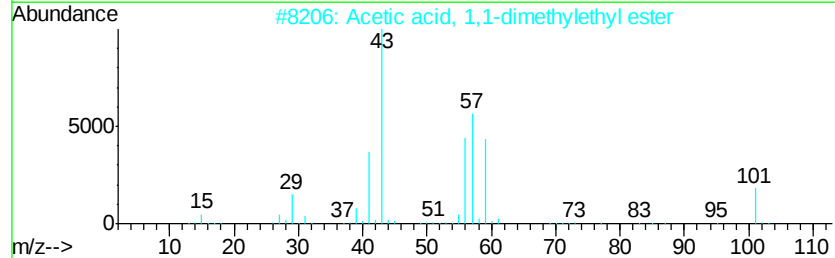
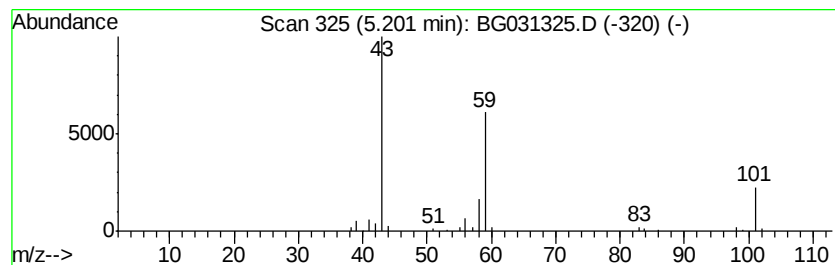
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.20	6.21 ng	103852	1,4-Dichlorobenzene-d4	8.13

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, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	38
4			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
5			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23



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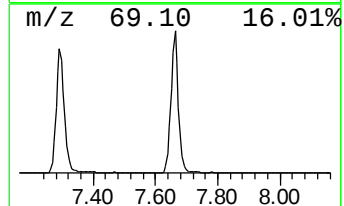
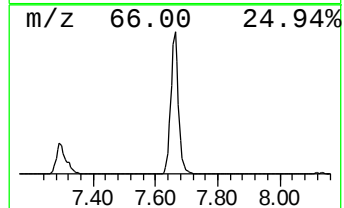
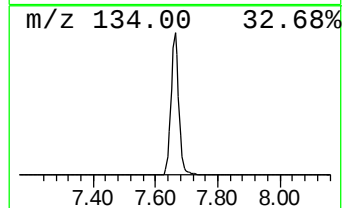
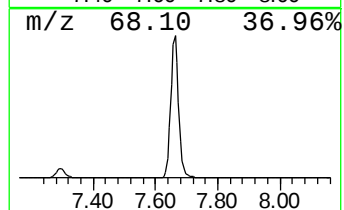
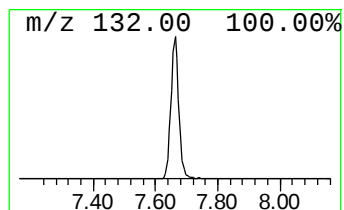
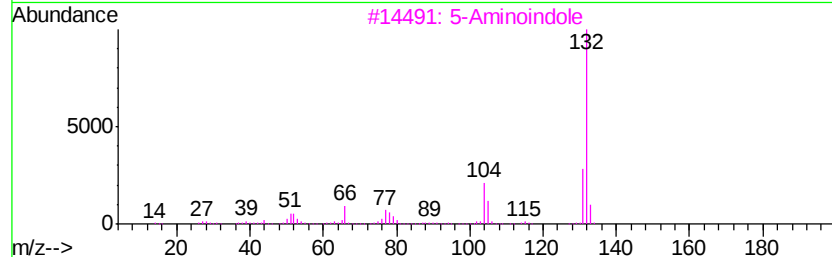
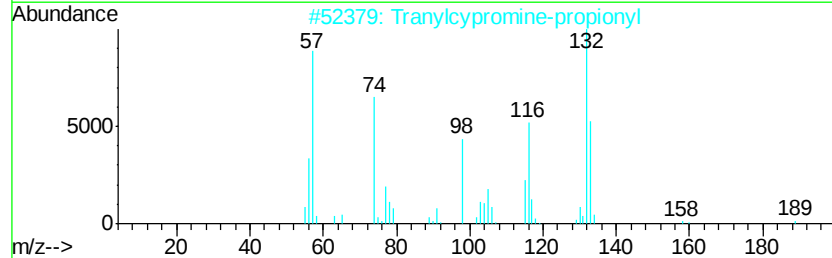
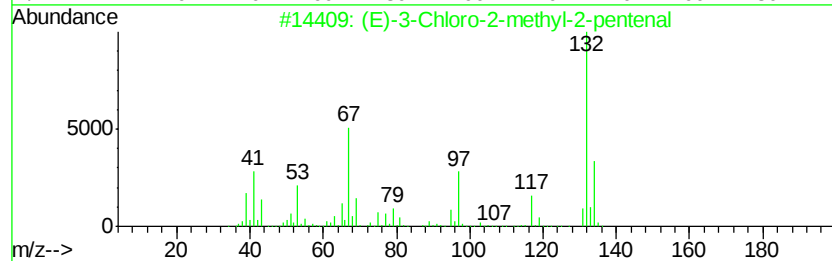
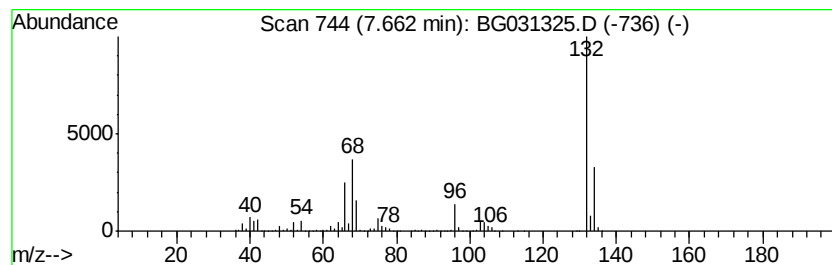
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Peak Number 2 unknown7.66 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.66	101.43 ng	1697190	1,4-Dichlorobenzene-d4	8.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
2	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
3	5-Aminoindole	132	C8H8N2	005192-03-0	10
4	Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.20	6.2	ng	103852	1	8.13	334644	20.0
unknown7.66	7.66	101.4	ng	1697190	1	8.13	334644	20.0