

Data Path : Z:\HPCHEM1\BNA G\DATA\BG113017\
 Data File : BG031304.D
 Acq On : 1 Dec 2017 1:23
 Operator : SJ/JU
 Sample : I6635-04
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 112817-JETT-E-U-S

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	5.205	321	326	333	rBV	61143	97857	1.58%	0.430%
2	5.693	402	409	423	rBV	300343	554884	8.97%	2.440%
3	7.297	675	682	698	rBV	177375	363997	5.88%	1.601%
4	7.661	737	744	760	rBV	610440	1127259	18.22%	4.957%
5	8.131	816	824	835	rBV	236974	408982	6.61%	1.798%
6	8.454	871	879	890	rBV	897399	1649034	26.66%	7.251%
7	9.300	1014	1023	1037	rBV	624588	1242555	20.09%	5.464%
8	10.945	1296	1303	1325	rBV	272154	589628	9.53%	2.593%
9	13.378	1709	1717	1730	rBV	2134201	3560837	57.56%	15.658%
10	14.753	1940	1951	1964	rBV2	511392	946798	15.31%	4.163%
11	16.239	2197	2204	2222	rBV	1092360	1827542	29.54%	8.036%
12	17.491	2410	2417	2432	rBV2	739916	1247623	20.17%	5.486%
13	20.087	2852	2859	2869	rBV	4443328	6185878	100.00%	27.201%
14	21.768	3137	3145	3161	rBV	852454	1507076	24.36%	6.627%
15	25.064	3693	3706	3717	rBV3	473486	1431707	23.14%	6.296%

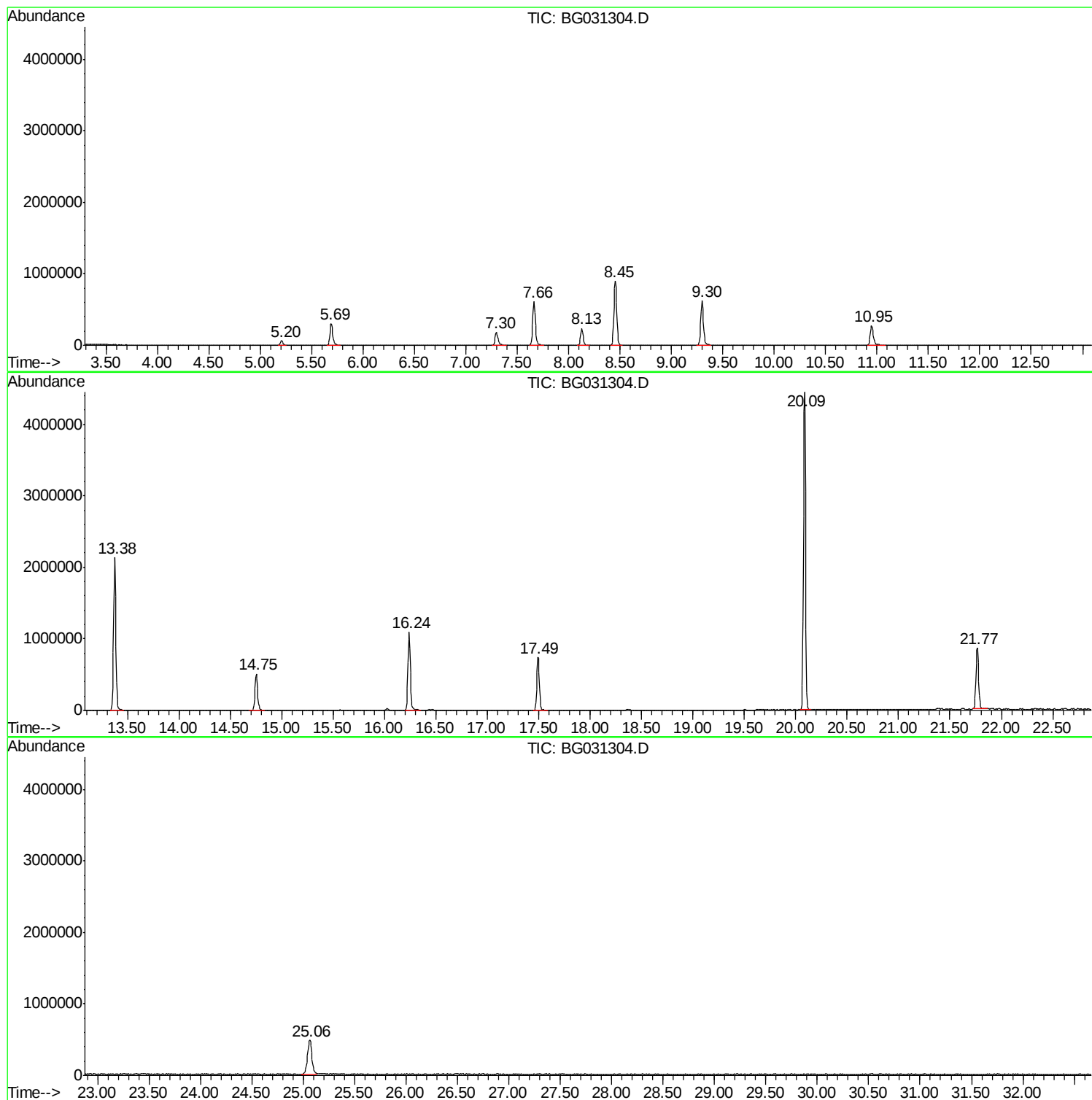
Sum of corrected areas: 22741657

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG111017.M
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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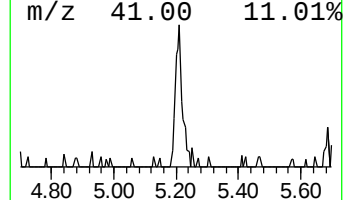
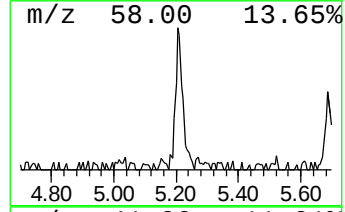
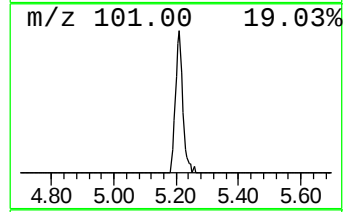
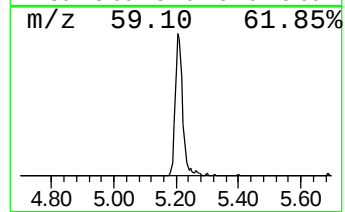
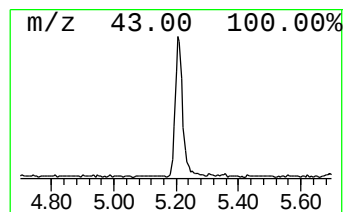
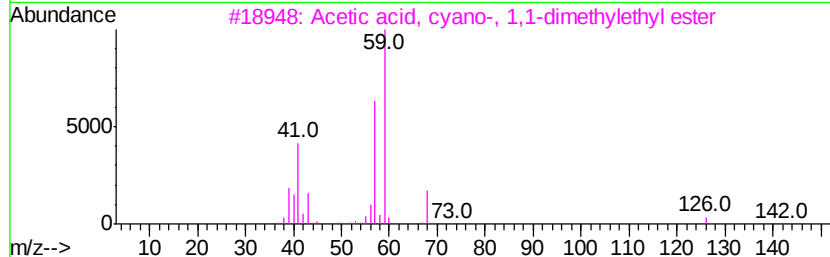
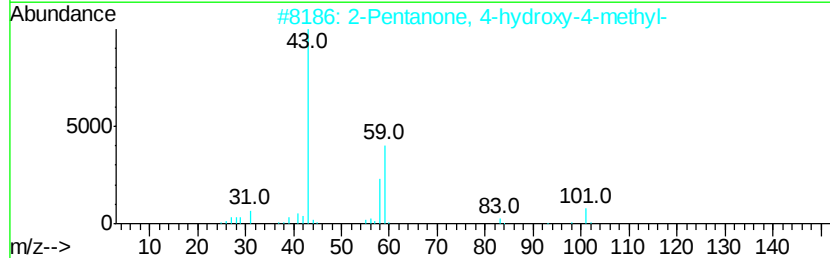
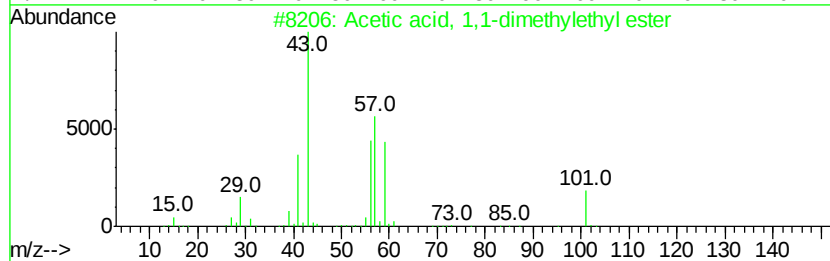
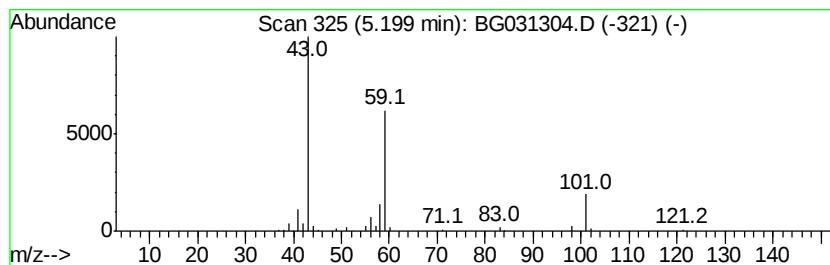
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Peak Number 1 unknown5.20 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.20	4.79 ng	97857	1,4-Dichlorobenzene-d4	8.13

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
2			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38
3			Acetic acid, cyano-, 1,1-dimethv...	141	C7H11NO2	001116-98-9	17
4			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5			Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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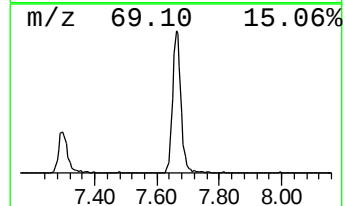
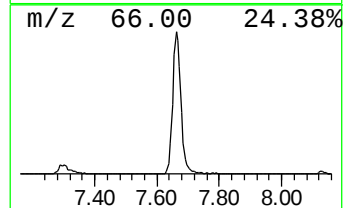
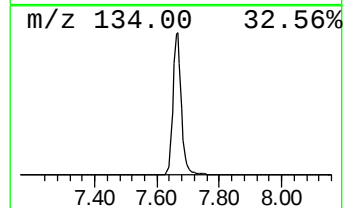
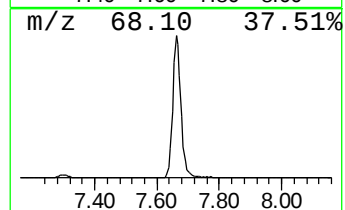
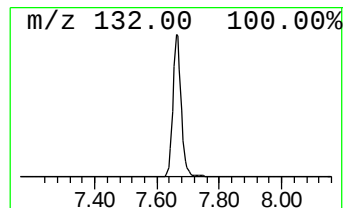
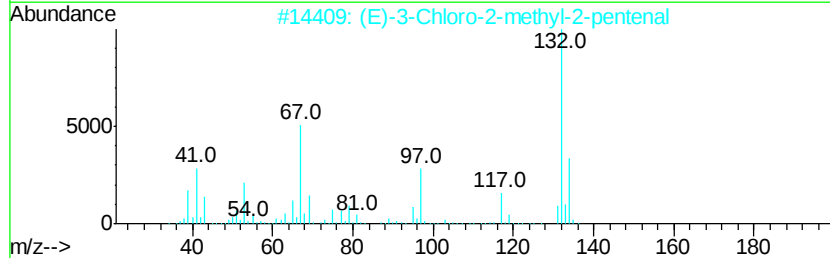
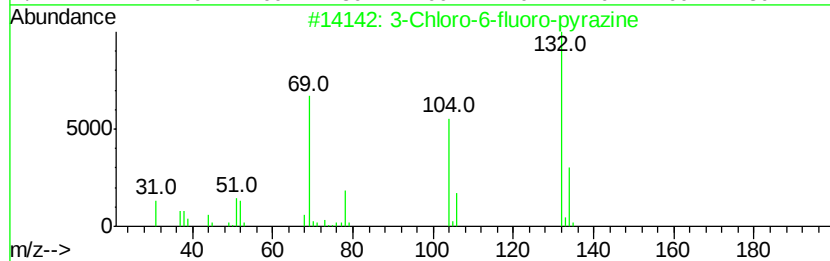
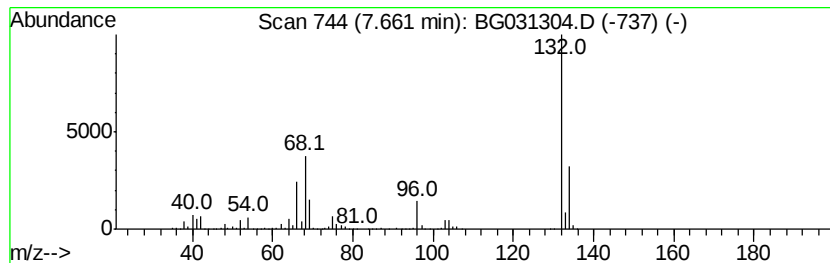
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Peak Number 2 unknown7.66 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.66	55.13 ng	1127260	1,4-Dichlorobenzene-d4	8.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Tranvlcypropane-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
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
unknown5.20	5.20	4.8	ng	97857	1	8.13	408982	20.0
unknown7.66	7.66	55.1	ng	1127260	1	8.13	408982	20.0