

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101118\
 Data File : BG037283.D
 Acq On : 12 Oct 2018 9:04
 Operator : JU/SJ
 Sample : J5371-04
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 100818-JETT-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:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG101118.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	5.194	319	324	332	rBV	35109	55576	1.07%	0.285%
2	5.652	395	402	411	rBV	296053	484847	9.36%	2.486%
3	7.250	668	674	686	rBV	193617	374204	7.22%	1.919%
4	7.644	732	741	749	rBV	531008	968533	18.70%	4.967%
5	8.120	814	822	833	rBV	199006	348565	6.73%	1.787%
6	8.443	869	877	886	rBV	793613	1403434	27.09%	7.197%
7	9.295	1014	1022	1037	rBV	665692	1218704	23.52%	6.250%
8	10.946	1295	1303	1311	rVB	282538	514019	9.92%	2.636%
9	13.378	1709	1717	1726	rBV	1950242	3084875	59.55%	15.820%
10	14.201	1850	1857	1863	rVB	96770	146809	2.83%	0.753%
11	14.753	1944	1951	1960	rVB2	476167	787697	15.20%	4.039%
12	16.239	2197	2204	2213	rBV	925585	1457869	28.14%	7.476%
13	17.497	2411	2418	2426	rBV	669457	1061019	20.48%	5.441%
14	20.105	2855	2862	2871	rBV	3611326	5180694	100.00%	26.567%
15	21.404	3078	3083	3088	rBV4	40895	63297	1.22%	0.325%
16	21.786	3141	3148	3158	rBV	689308	1180728	22.79%	6.055%
17	25.135	3707	3718	3734	rVB2	376696	1169360	22.57%	5.997%

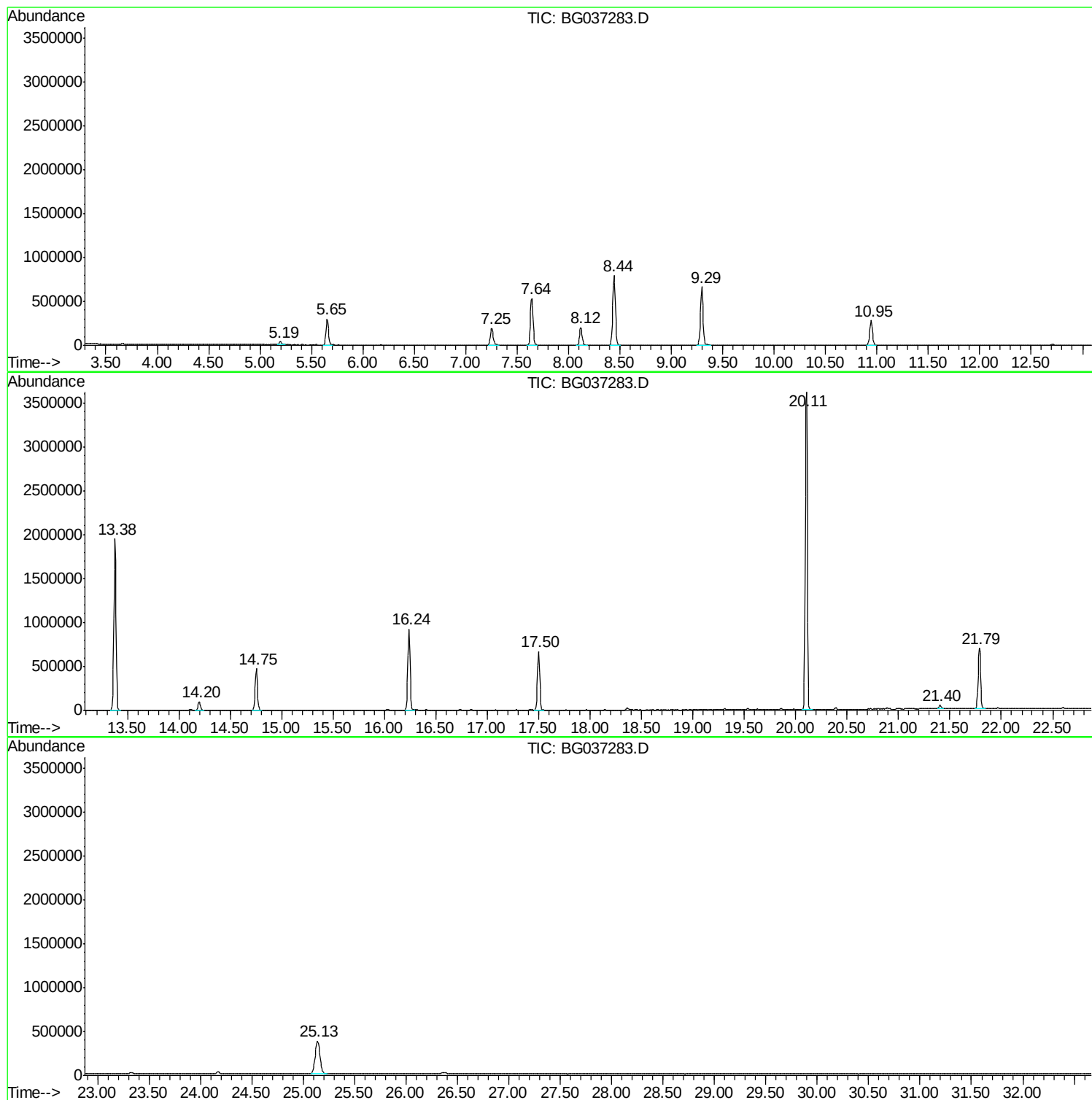
Sum of corrected areas: 19500230

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG101118.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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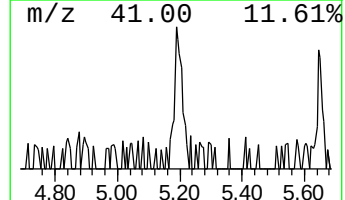
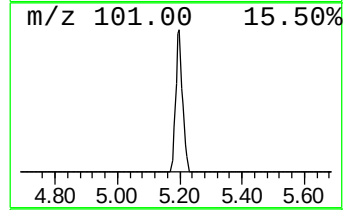
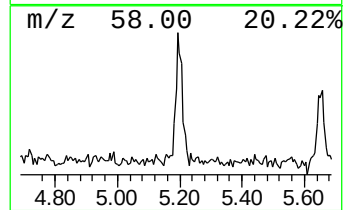
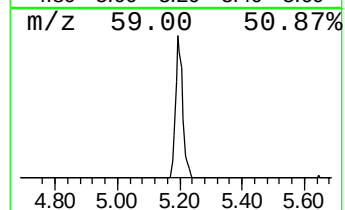
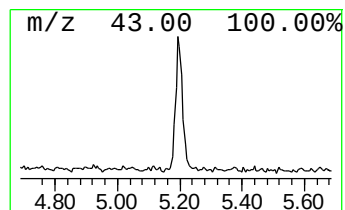
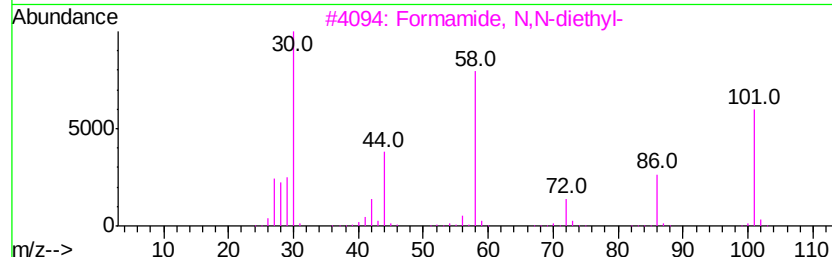
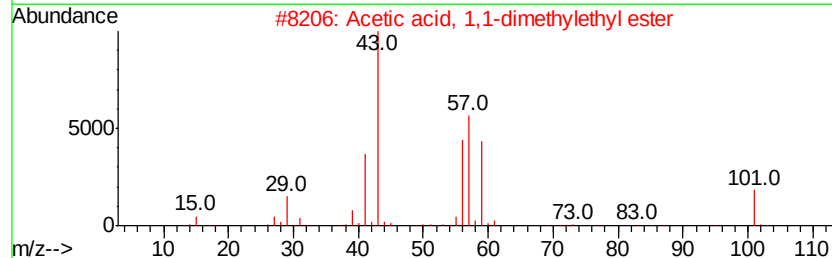
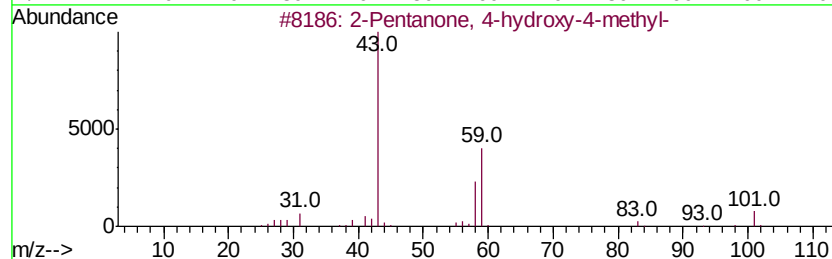
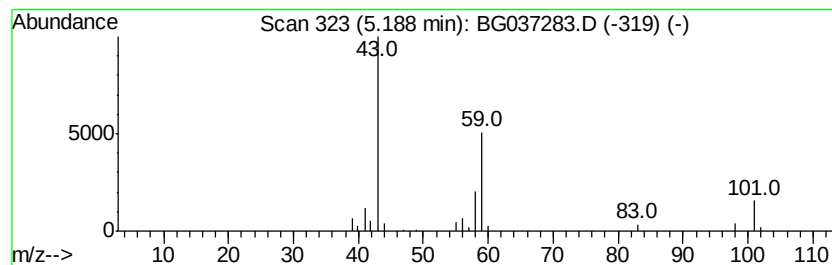
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.19	3.19 ng	55576	1,4-Dichlorobenzene-d4	8.12

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	36
3			Formamide, N,N-diethyl-	101	C5H11NO	000617-84-5	9
4			4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	9
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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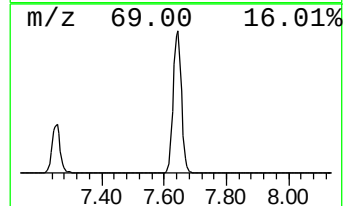
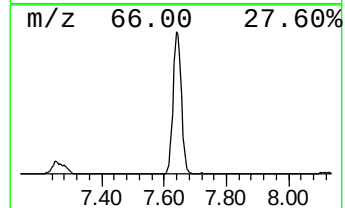
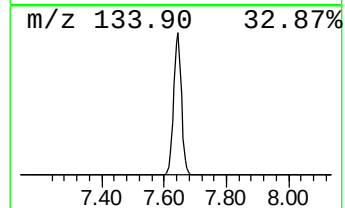
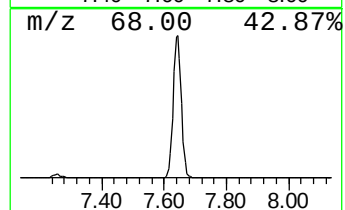
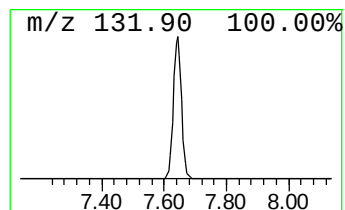
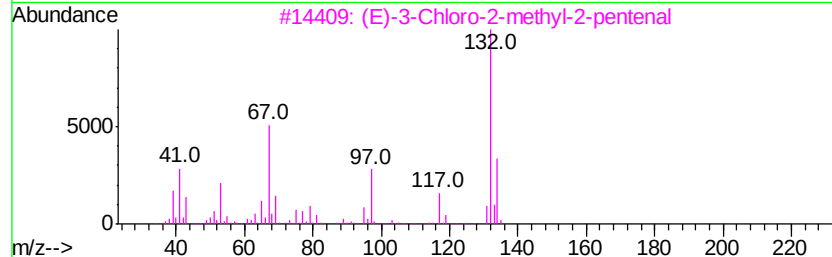
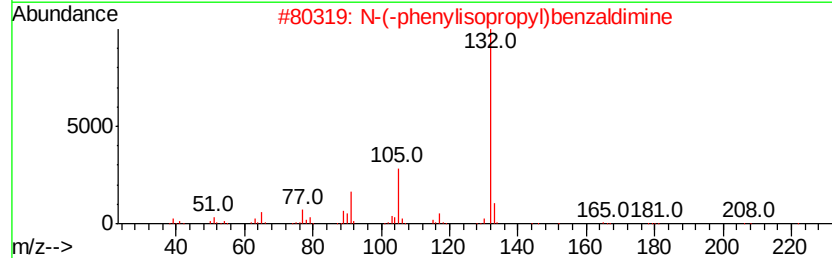
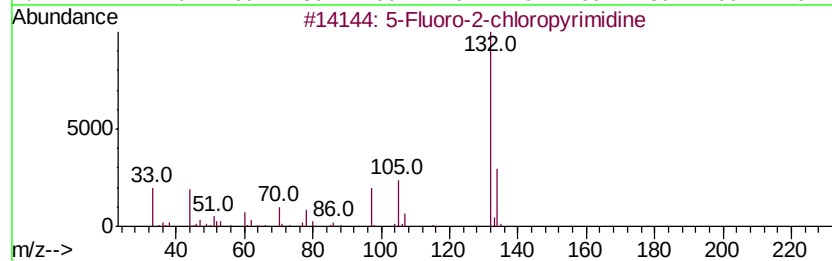
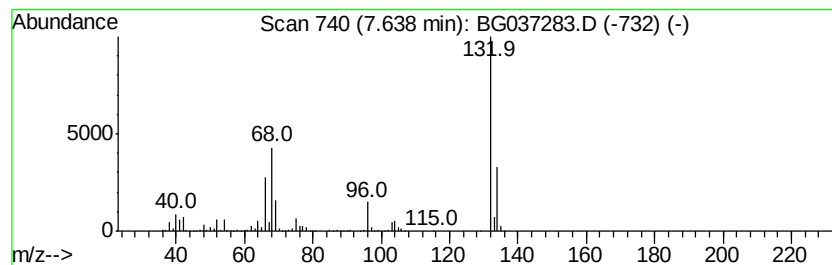
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Peak Number 2 unknown7.64 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.64	55.57 ng	968533	1,4-Dichlorobenzene-d4	8.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
2		N-(-phenylisopropyl)benzalimine	223	C16H17N	1000379-01-3	10
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	10
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	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	3.2	ng	55576	1	8.12	348565	20.0
unknown7.64	7.64	55.6	ng	968533	1	8.12	348565	20.0