

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042718\
 Data File : BG034058.D
 Acq On : 28 Apr 2018 4:33
 Operator : SJ/JU
 Sample : J2571-11
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 042418-SIDI-E-U-M

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG041918.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	4.891	268	273	286	rBV	51243	107845	1.64%	0.516%
2	5.379	349	356	369	rBV2	57420	233306	3.54%	1.115%
3	6.989	621	630	638	rBV5	25306	110876	1.68%	0.530%
4	7.294	676	682	708	rBV	207542	775159	11.76%	3.706%
5	7.758	754	761	781	rBV2	78036	289502	4.39%	1.384%
6	8.064	807	813	842	rBV	569281	1445481	21.93%	6.910%
7	8.916	951	958	984	rBV2	218846	888376	13.48%	4.247%
8	10.543	1227	1235	1254	rBV	112887	401888	6.10%	1.921%
9	13.011	1649	1655	1689	rBV	1479564	3380010	51.27%	16.159%
10	14.392	1883	1890	1916	rBV2	388718	857201	13.00%	4.098%
11	15.884	2138	2144	2174	rBV3	689208	1770236	26.85%	8.463%
12	16.125	2181	2185	2198	rVB2	27626	66534	1.01%	0.318%
13	17.147	2352	2359	2384	rBV2	470373	1171202	17.77%	5.599%
14	19.780	2801	2807	2839	rBV	3905712	6592028	100.00%	31.514%
15	21.078	3024	3028	3039	rBV6	38029	100185	1.52%	0.479%
16	21.389	3075	3081	3098	rBV2	562705	1399485	21.23%	6.690%
17	24.392	3582	3592	3614	rBV	355185	1328433	20.15%	6.351%

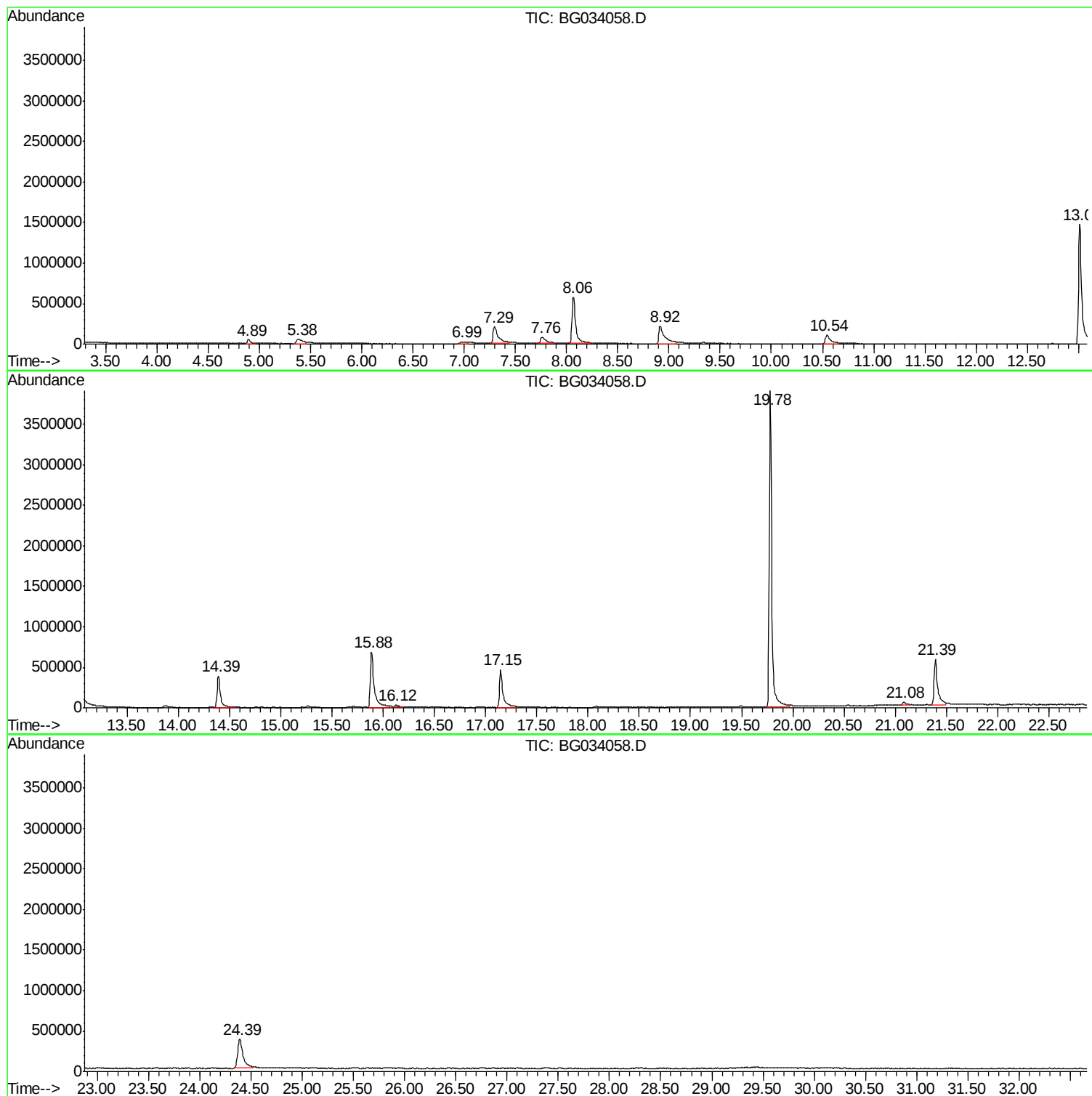
Sum of corrected areas: 20917747

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG041918.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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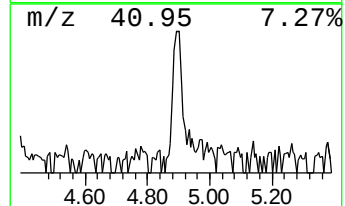
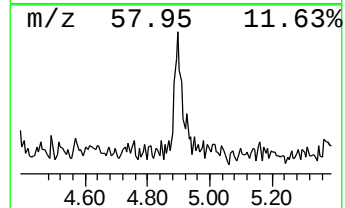
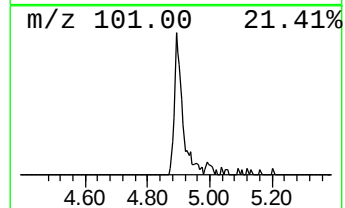
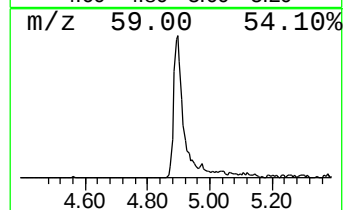
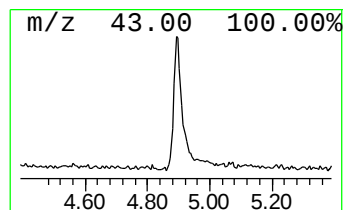
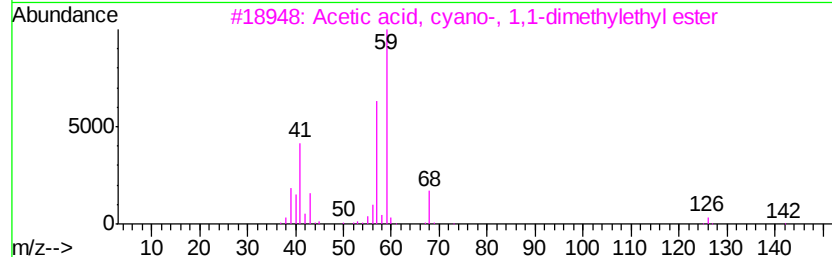
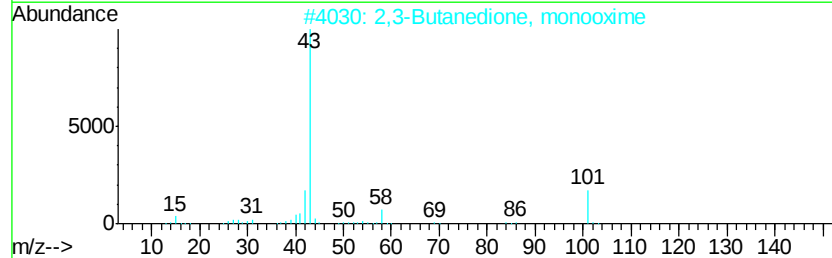
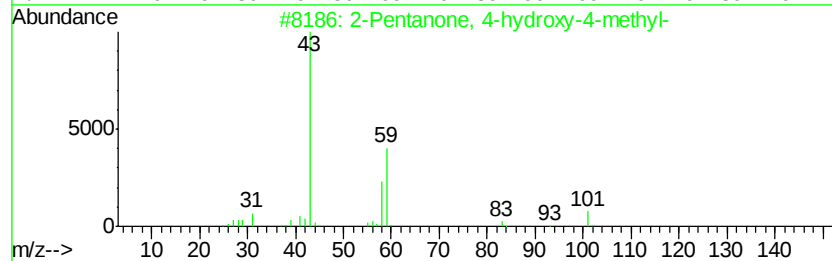
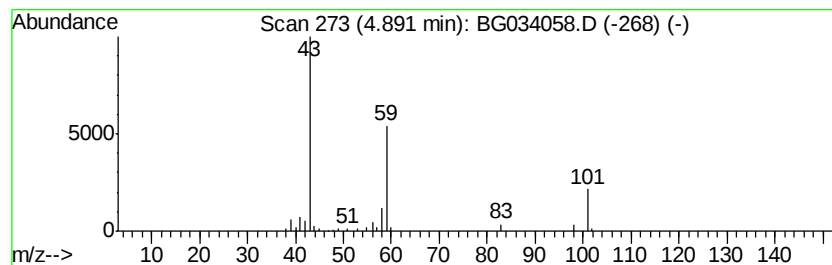
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.89	7.45 ng	107845	1,4-Dichlorobenzene-d4	7.76

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	35
3			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
4			Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
5			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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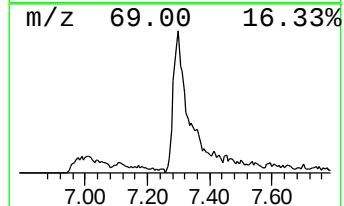
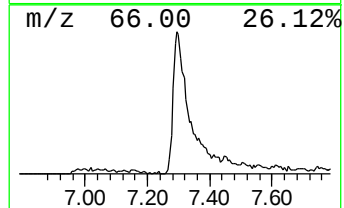
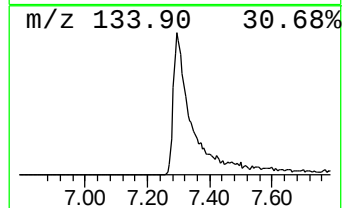
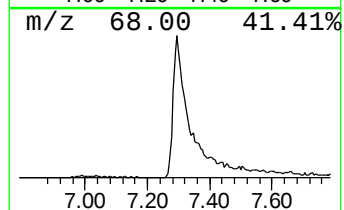
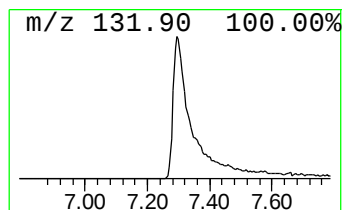
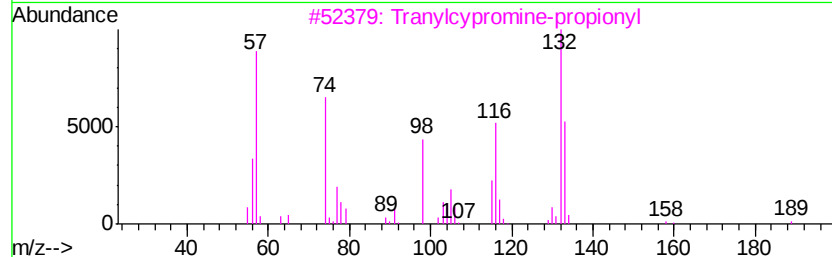
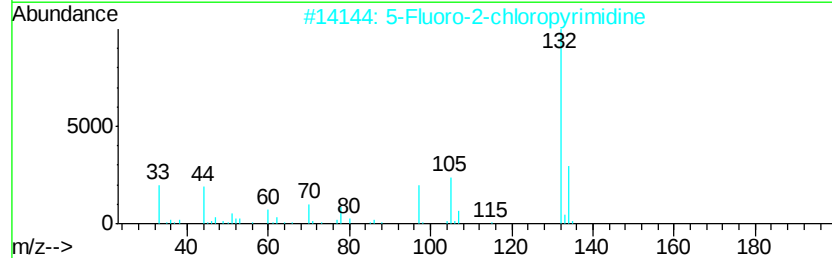
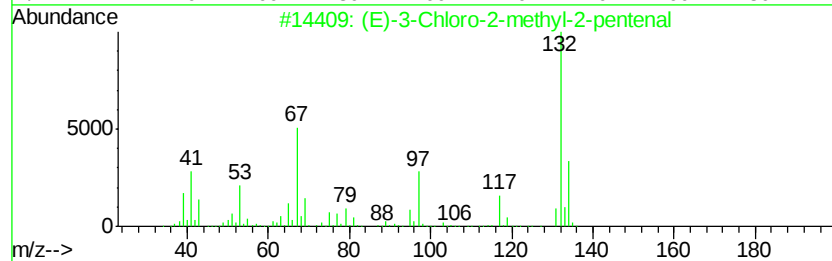
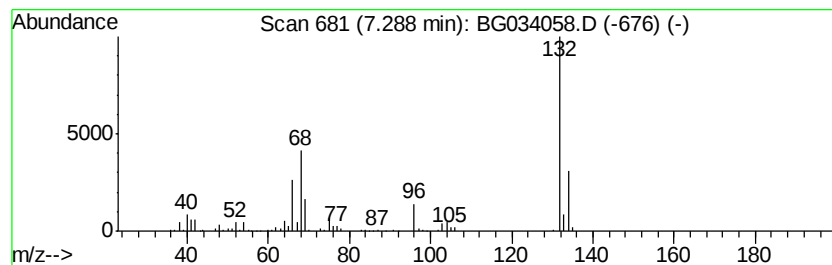
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Peak Number 2 unknown7.29 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.29	53.55 ng	775159	1,4-Dichlorobenzene-d4	7.76

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		Tranvlcvpromine-propionyl	189	C12H15NO	1000123-86-3	10
4		Xenon	132	Xe	007440-63-3	9
5		1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.89	7.5	ng	107845	1	7.76	289502	20.0
unknown7.29	7.29	53.5	ng	775159	1	7.76	289502	20.0