

Data Path : Z:\HPCHEM1\BNA G\DATA\BG022318\
 Data File : BG032807.D
 Acq On : 24 Feb 2018 1:50
 Operator : SJ/JU
 Sample : PB106794BL
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB106794BL

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-BG022118.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	5.828	425	432	440	rVB	38375	63250	1.93%	0.350%
2	6.339	511	519	531	rBV	878466	1493987	45.62%	8.259%
3	8.019	797	805	818	rBV	850488	1539913	47.03%	8.513%
4	8.436	867	876	887	rBV	846815	1546071	47.22%	8.547%
5	8.924	951	959	967	rBV	210456	375929	11.48%	2.078%
6	9.264	1007	1017	1026	rVB	634305	1189489	36.33%	6.576%
7	10.122	1154	1163	1174	rBV	473944	879374	26.86%	4.862%
8	11.814	1442	1451	1459	rVB	290634	540601	16.51%	2.989%
9	14.169	1845	1852	1860	rBV	1334650	2092264	63.90%	11.567%
10	15.538	2078	2085	2093	rVB2	413859	688827	21.04%	3.808%
11	17.007	2328	2335	2354	rBV	978482	1500182	45.81%	8.294%
12	18.270	2543	2550	2559	rBV	553041	866878	26.47%	4.792%
13	20.784	2972	2978	2986	rBV	2378190	3274509	100.00%	18.103%
14	22.623	3283	3291	3304	rBV2	531079	1017687	31.08%	5.626%
15	26.441	3928	3941	3956	rBV2	306472	1019480	31.13%	5.636%

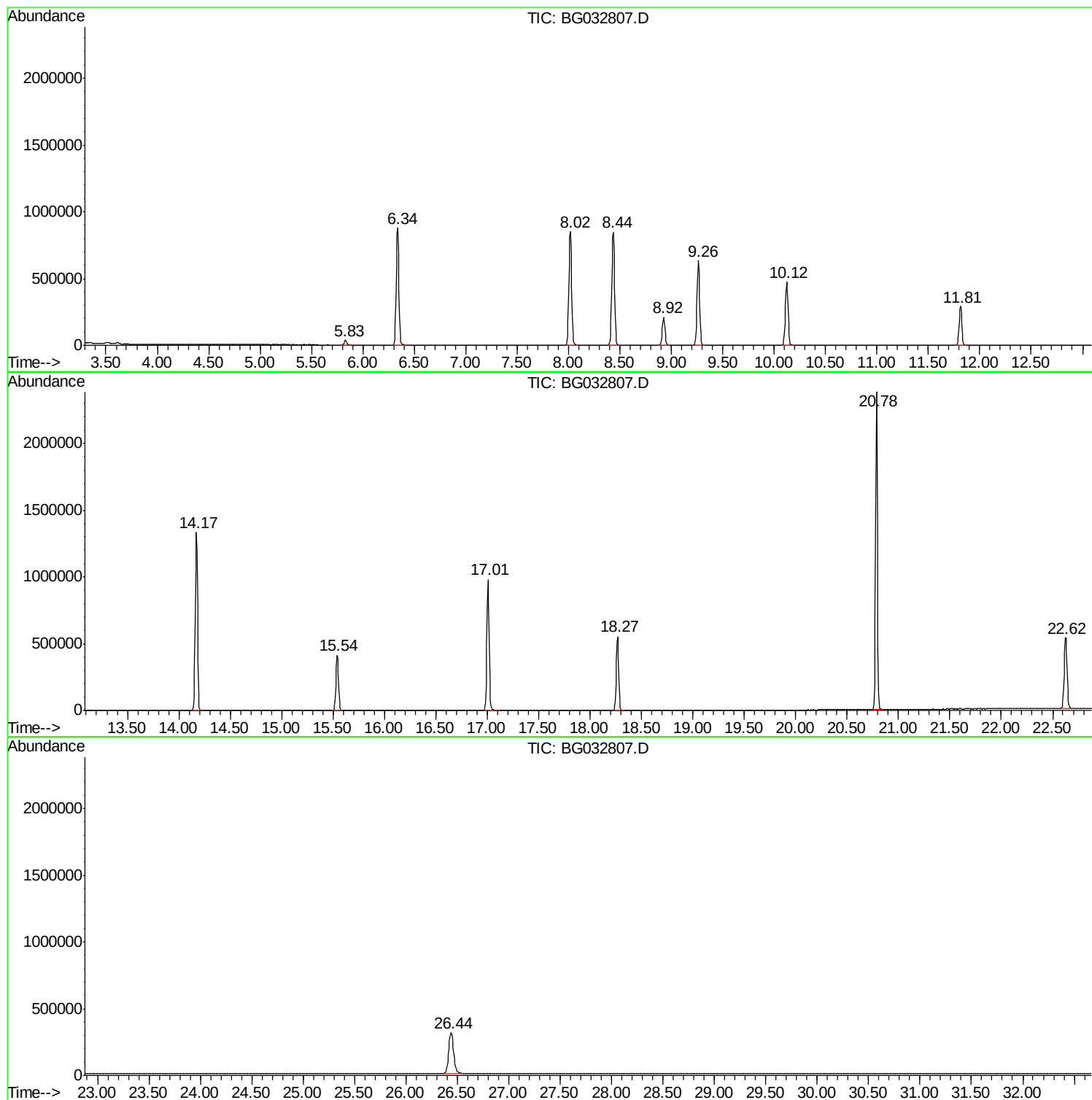
Sum of corrected areas: 18088441

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022118.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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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
5.83	3.37 ng	63250	1,4-Dichlorobenzene-d4	8.92	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
3	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
4	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	25
5	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23

Peak Number 2 unknown8.44 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
8.44	82.25 ng	1546070	1,4-Dichlorobenzene-d4	8.92	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3	1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
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...	5.83	3.4	ng	63250	1	8.92	375929	20.0
unknown8.44	8.44	82.3	ng	1546070	1	8.92	375929	20.0