

Data Path : Z:\HPCHEM1\BNA G\DATA\BG022318\
 Data File : BG032806.D
 Acq On : 24 Feb 2018 1:12
 Operator : SJ/JU
 Sample : PB106787BL
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB106787BL

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	427	432	440	rVB2	44665	74136	1.85%	0.399%
2	6.333	510	518	531	rBV	686947	1192355	29.76%	6.417%
3	8.014	796	804	818	rBV	670372	1242504	31.01%	6.687%
4	8.431	867	875	886	rBV	676749	1255151	31.32%	6.755%
5	8.924	951	959	966	rBV	199053	355781	8.88%	1.915%
6	9.265	1008	1017	1030	rVB	753249	1396712	34.86%	7.517%
7	10.122	1152	1163	1176	rBV	561780	1053897	26.30%	5.672%
8	11.814	1441	1451	1460	rBV	282004	514758	12.85%	2.770%
9	14.170	1845	1852	1863	rBV	1552406	2472946	61.72%	13.308%
10	15.539	2078	2085	2093	rVB2	405097	667064	16.65%	3.590%
11	17.007	2328	2335	2349	rBV	814483	1233279	30.78%	6.637%
12	18.270	2542	2550	2561	rBV	560697	874920	21.83%	4.709%
13	20.414	2910	2915	2920	rBV	41620	57444	1.43%	0.309%
14	20.790	2972	2979	2991	rBV	2885172	4007027	100.00%	21.564%
15	22.629	3283	3292	3300	rBV	588261	1078773	26.92%	5.806%
16	26.442	3928	3941	3958	rBV	326245	1104957	27.58%	5.946%

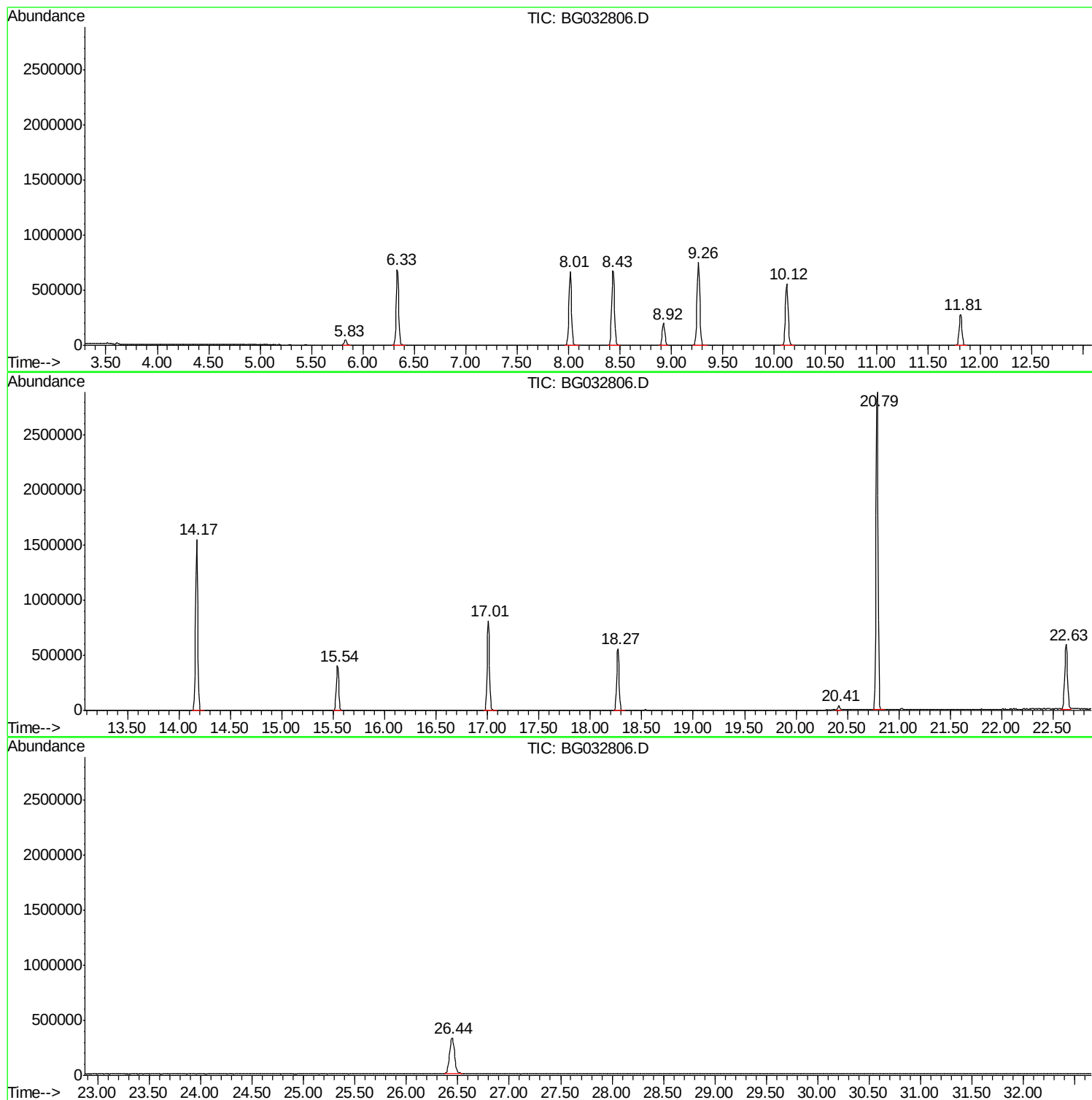
Sum of corrected areas: 18581704

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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 unknown5.83 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
5.83	4.17 ng	74136	1,4-Dichlorobenzene-d4	8.92	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	47
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
3	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	36
4	Silane, ethyldimethyl-	88	C4H12Si	000758-21-4	25
5	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10

Peak Number 2 unknown8.43 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
8.43	70.56 ng	1255150	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	16
3	5-Aminoindole	132	C8H8N2	005192-03-0	12
4	2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	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
unknown5.83	5.83	4.2	ng	74136	1	8.92	355781	20.0
unknown8.43	8.43	70.6	ng	1255150	1	8.92	355781	20.0