

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG121517\
 Data File : BG031596.D
 Acq On : 15 Dec 2017 20:13
 Operator : SJ/JU
 Sample : PB105057BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB105057BL

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-BG121117.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	5.170	315	320	330	rBV	47172	77812	1.58%	0.405%
2	5.658	396	403	418	rBV	559712	981120	19.95%	5.107%
3	7.268	670	677	697	rBV	537373	1028522	20.92%	5.354%
4	7.632	732	739	750	rBV	580554	1038675	21.12%	5.407%
5	8.096	811	818	827	rBV	188961	330995	6.73%	1.723%
6	8.420	866	873	885	rBV	673500	1220105	24.81%	6.351%
7	9.266	1010	1017	1034	rBV	432651	886762	18.03%	4.616%
8	10.917	1290	1298	1318	rBV	205451	470042	9.56%	2.447%
9	13.343	1704	1711	1724	rBV	1477680	2537977	51.61%	13.212%
10	14.724	1940	1946	1963	rBV2	382934	713469	14.51%	3.714%
11	16.216	2193	2200	2226	rBV	685165	1304840	26.54%	6.793%
12	17.468	2406	2413	2425	rBV2	600251	1044401	21.24%	5.437%
13	20.059	2848	2854	2863	rBV	3584183	4917221	100.00%	25.597%
14	21.739	3133	3140	3157	rBV2	801570	1456882	29.63%	7.584%
15	25.018	3687	3698	3715	rBV2	377842	1200990	24.42%	6.252%

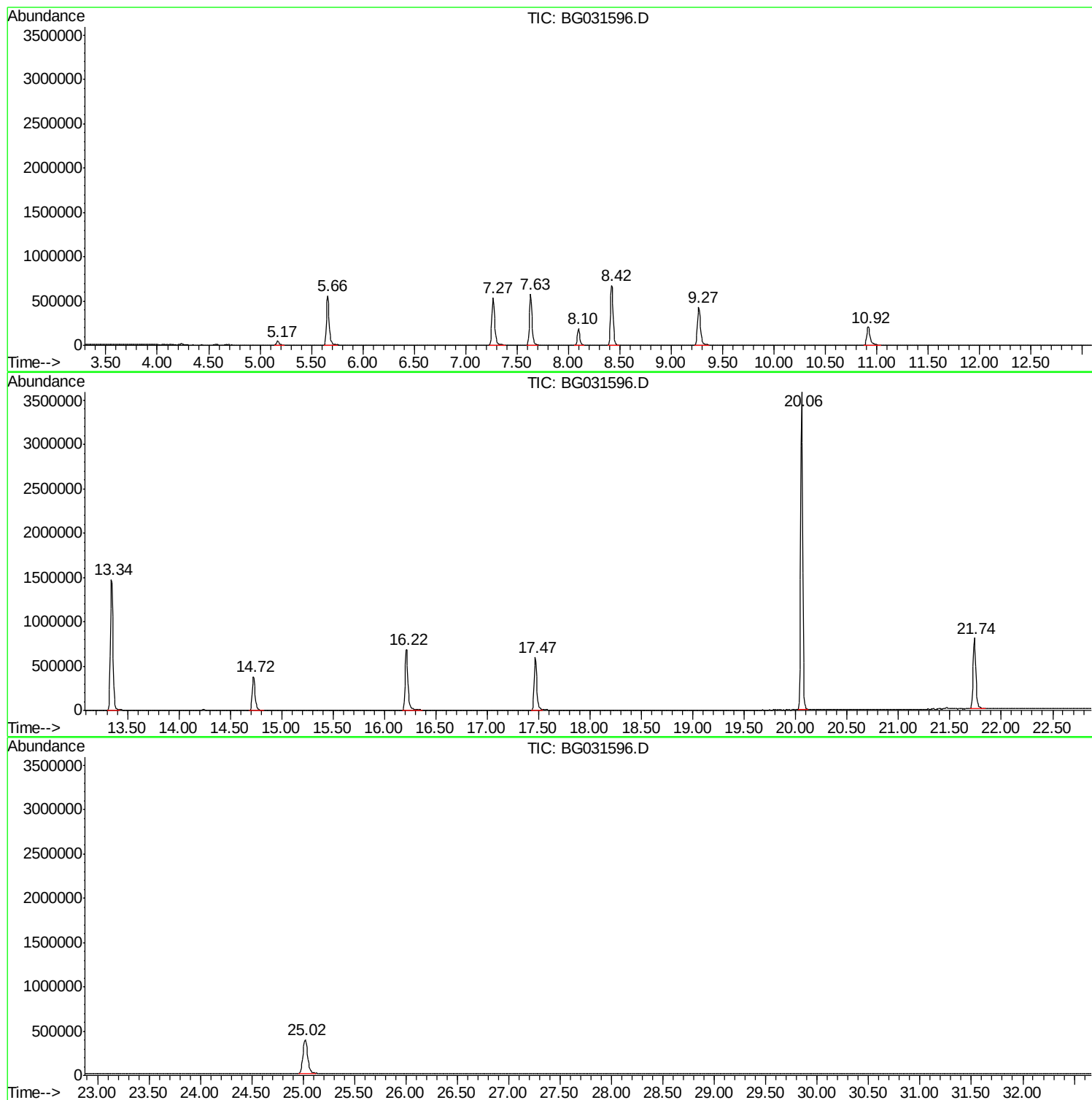
Sum of corrected areas: 19209813

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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG121117.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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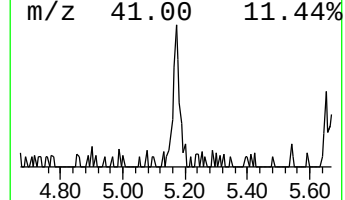
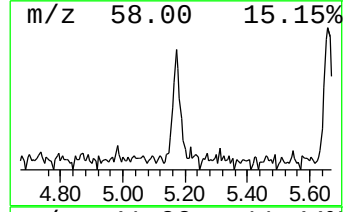
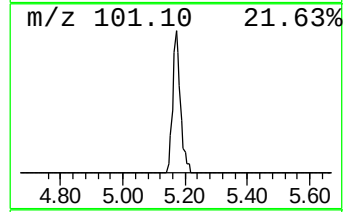
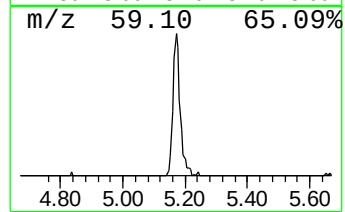
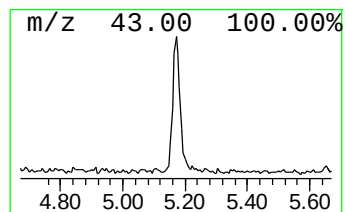
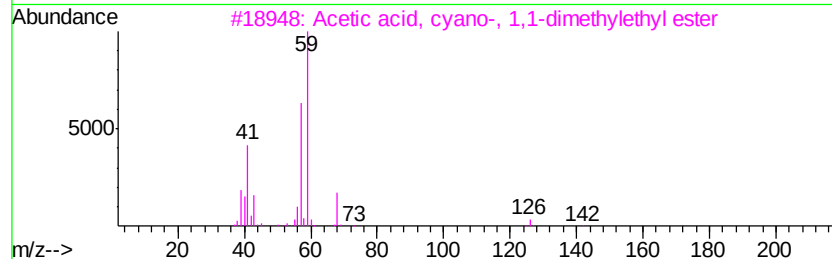
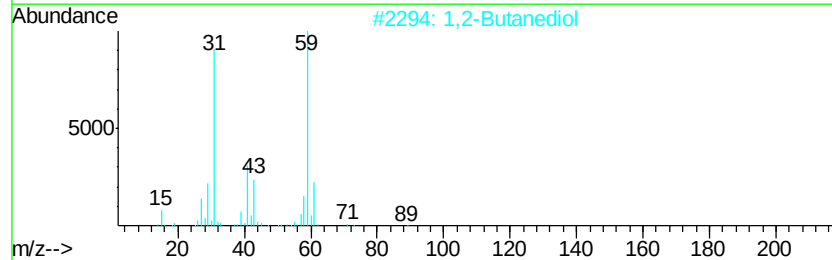
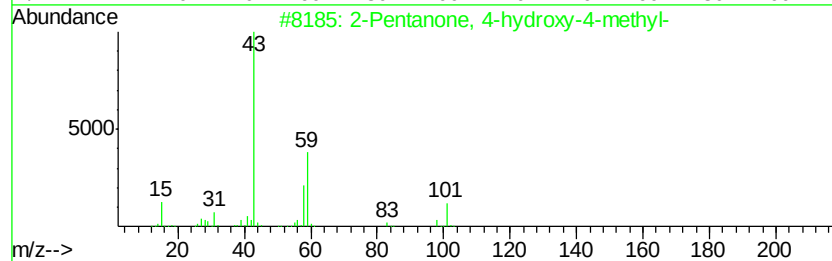
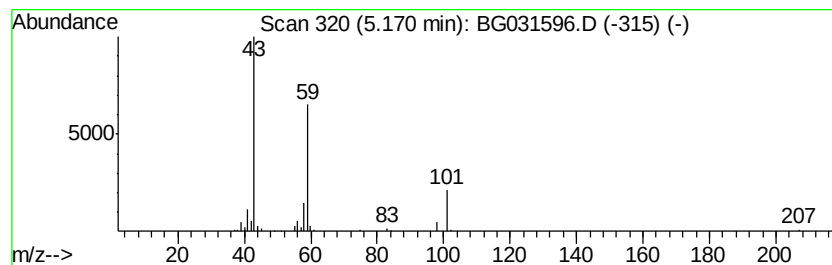
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.17	4.70 ng	77812	1,4-Dichlorobenzene-d4	8.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40
2		1,2-Butanediol	90	C4H10O2	000584-03-2	23
3		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17
4		1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	17
5		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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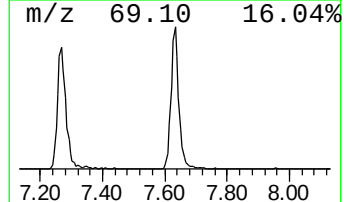
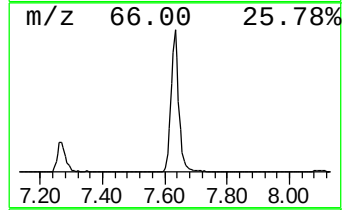
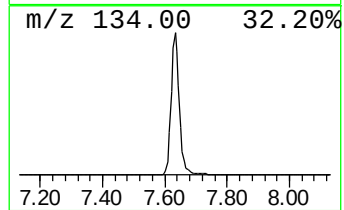
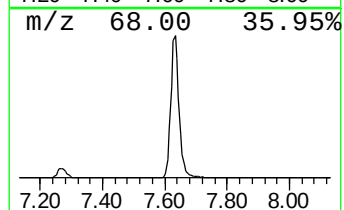
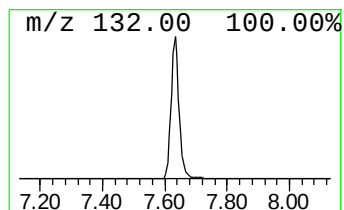
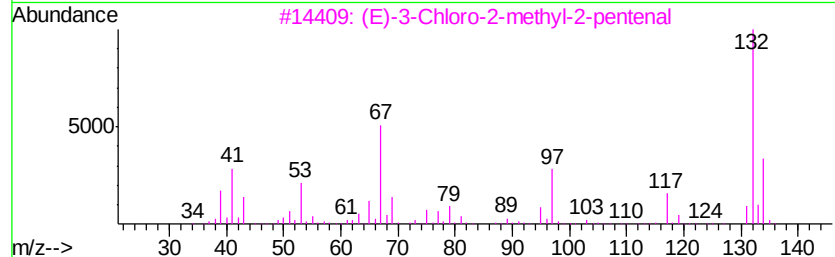
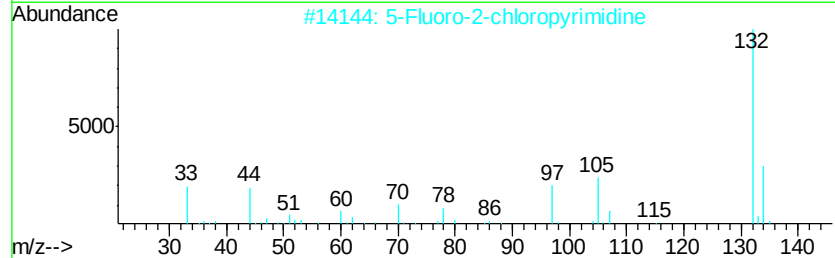
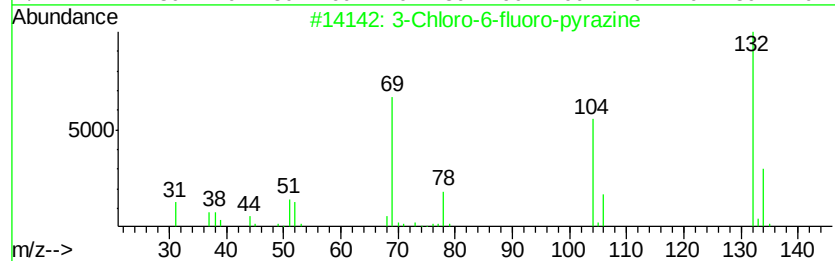
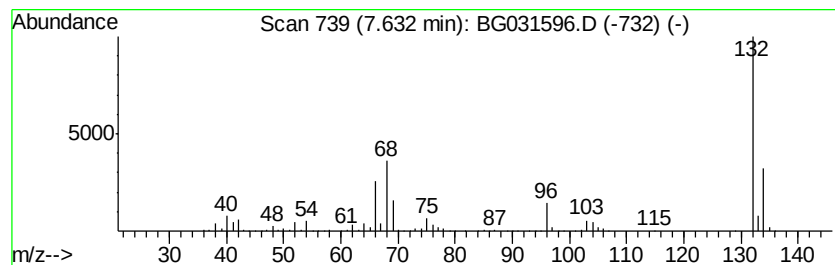
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Peak Number 2 unknown7.63 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.63	62.76 ng	1038680	1,4-Dichlorobenzene-d4	8.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	17
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
4		Tranlylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
5		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.17	4.7	ng	77812	1	8.10	330995	20.0
unknown7.63	7.63	62.8	ng	1038680	1	8.10	330995	20.0