

Data Path : Z:\HPCHEM1\BNA G\DATA\BG122117\
 Data File : BG031688.D
 Acq On : 22 Dec 2017 00:21
 Operator : SJ/JU
 Sample : I7002-02
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EPE-106-3DS(20-30)

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-BG122117.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	3.447	21	27	39	rBV	40191	84755	2.30%	0.478%
2	5.738	412	417	425	rVB	34395	58916	1.60%	0.332%
3	6.243	496	503	517	rVB	625057	1072565	29.15%	6.049%
4	7.924	780	789	802	rVB	635270	1194808	32.47%	6.738%
5	8.335	851	859	868	rBV	653859	1213307	32.98%	6.843%
6	8.829	935	943	950	rBV2	138014	254521	6.92%	1.435%
7	9.163	992	1000	1011	rBV	504776	938405	25.50%	5.292%
8	10.027	1138	1147	1160	rBV	356485	660878	17.96%	3.727%
9	11.719	1426	1435	1444	rBV	203490	382254	10.39%	2.156%
10	14.093	1831	1839	1848	rVB	1307598	2118094	57.57%	11.945%
11	14.886	1968	1974	1981	rBV	93484	142415	3.87%	0.803%
12	15.462	2065	2072	2081	rVB2	383024	631821	17.17%	3.563%
13	16.937	2316	2323	2335	rVB	1236202	1947196	52.92%	10.981%
14	18.194	2531	2537	2547	rVB2	557256	907017	24.65%	5.115%
15	19.017	2673	2677	2686	rBV3	40326	64135	1.74%	0.362%
16	20.256	2884	2888	2895	rBV4	29206	39579	1.08%	0.223%
17	20.732	2962	2969	2976	rBV	2734816	3679441	100.00%	20.751%
18	22.554	3272	3279	3289	rVB	600839	1166243	31.70%	6.577%
19	26.343	3912	3924	3938	rVB	358699	1175457	31.95%	6.629%

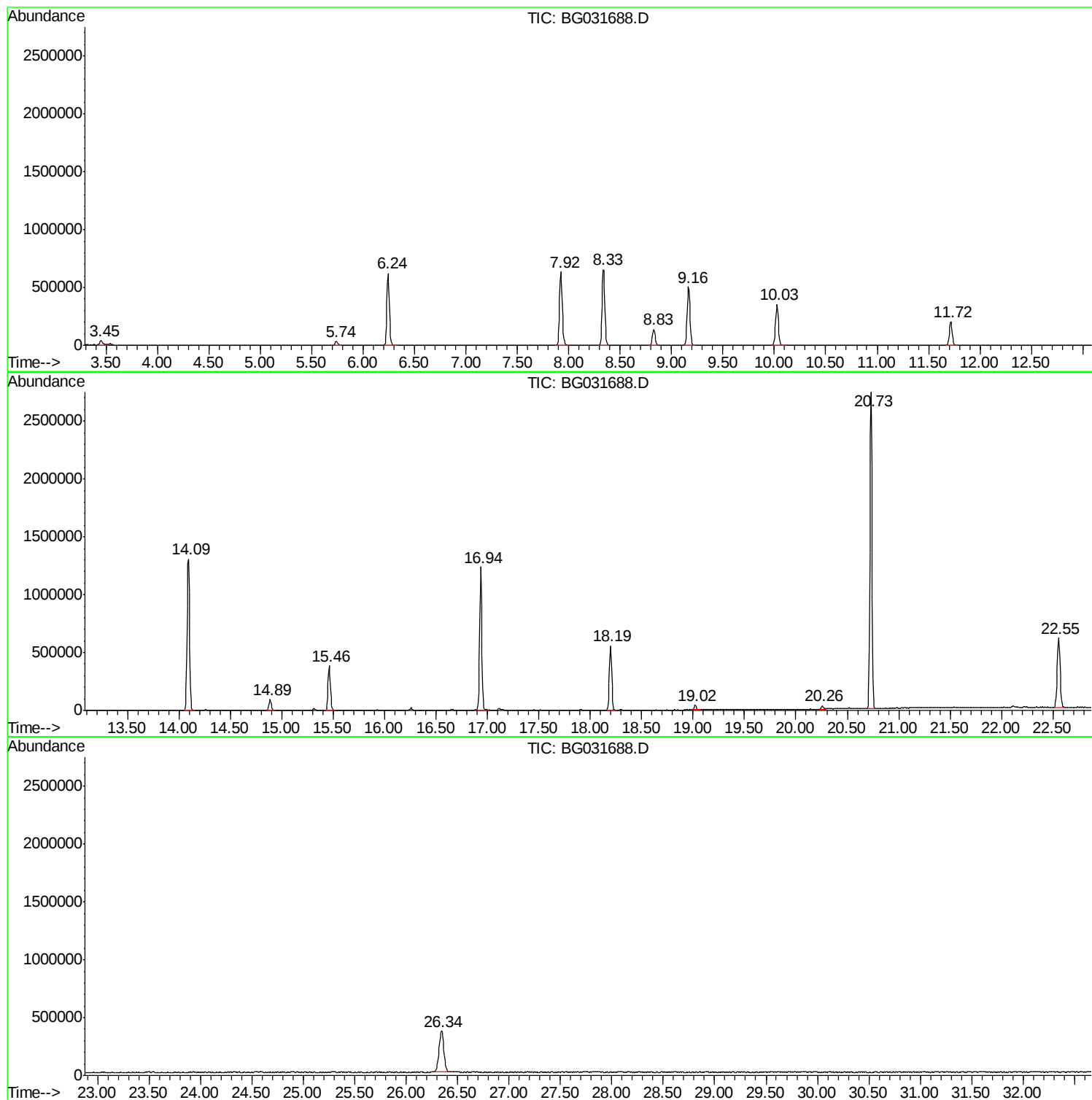
Sum of corrected areas: 17731807

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG122117.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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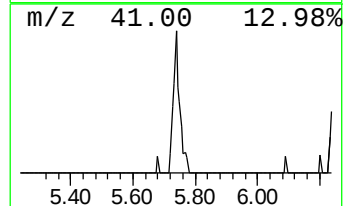
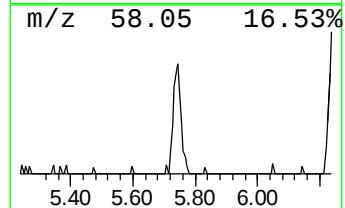
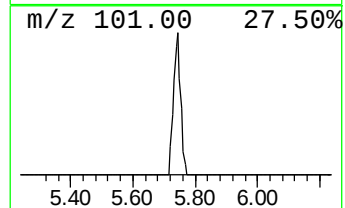
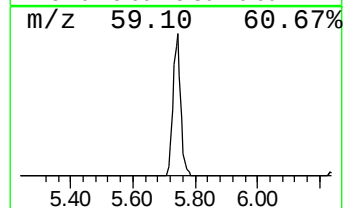
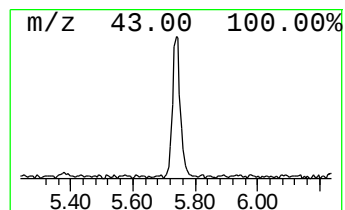
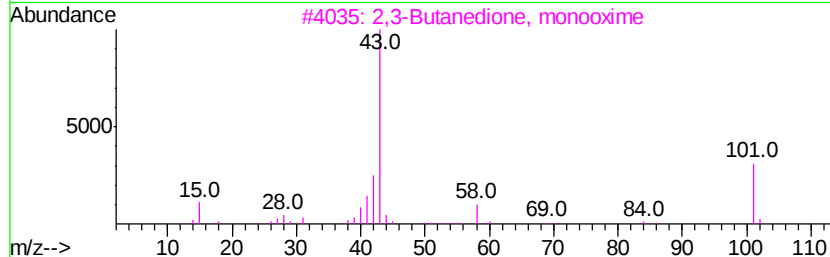
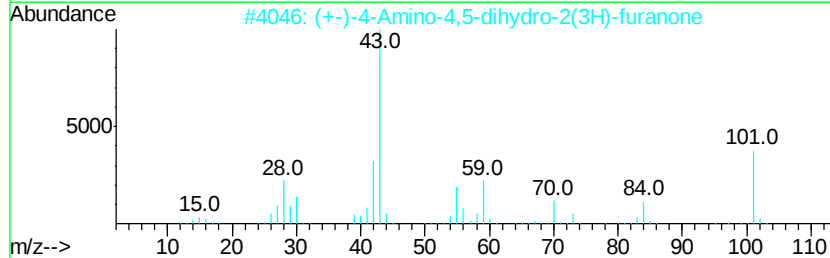
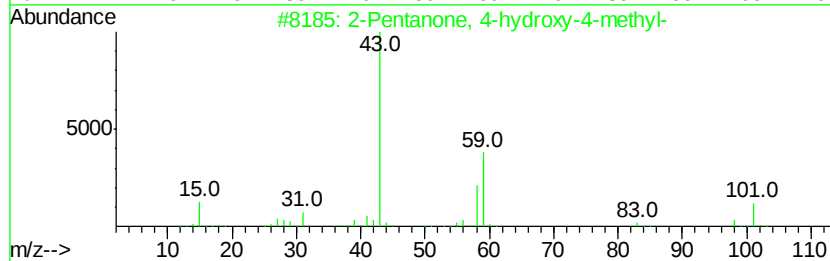
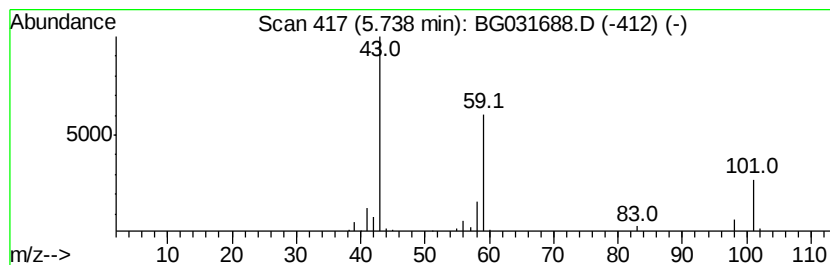
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.74	4.63 ng	58916	1,4-Dichlorobenzene-d4	8.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		(+)-4-Amino-4,5-dihydro-2(3H)-furanone	101	C4H7NO2	016504-58-8	47
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
4		3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	25
5		Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	16



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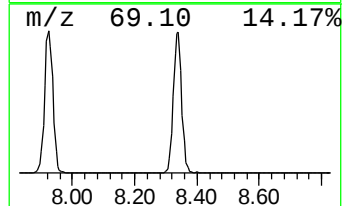
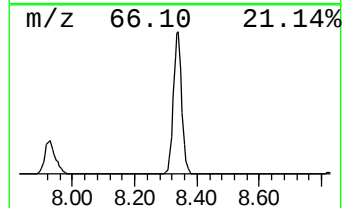
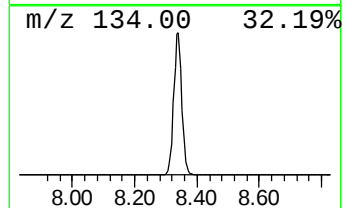
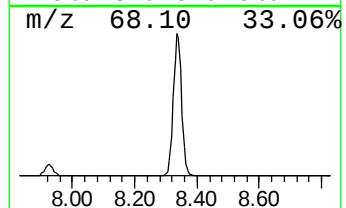
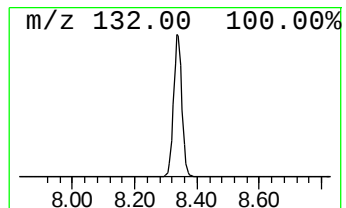
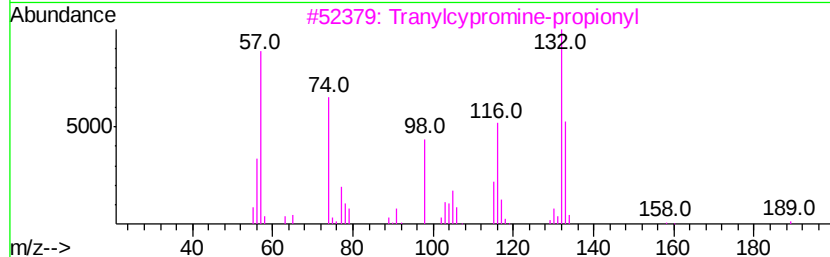
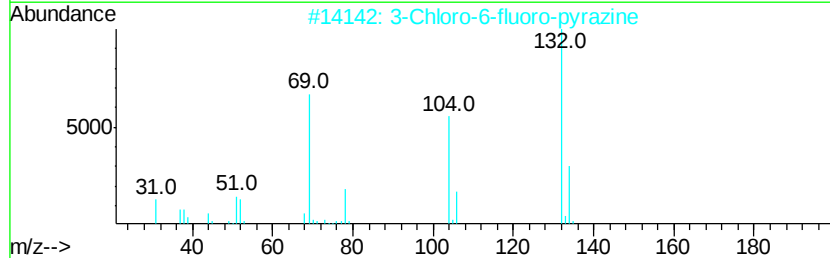
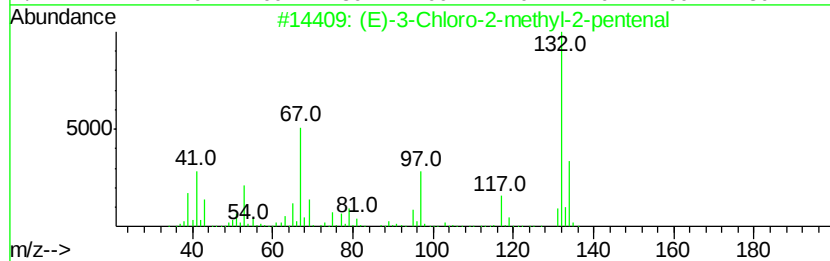
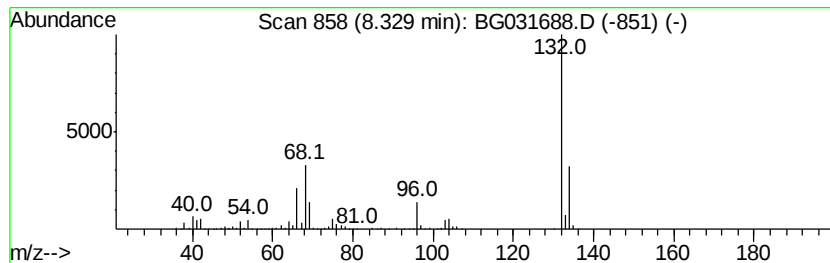
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Peak Number 3 unknown8.33 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.33	95.34 ng	1213310	1,4-Dichlorobenzene-d4	8.83

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	35	
2	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25	
3	Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12	
4	5-Aminoindole	132	C8H8N2	005192-03-0	12	
5	1H-Indazole, 7-methyl-	132	C8H8N2	1000316-00-7	10	



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.74	4.6	ng	58916	1	8.83	254521	20.0
unknown8.33	8.33	95.3	ng	1213310	1	8.83	254521	20.0