

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG110117\
 Data File : BG030643.D
 Acq On : 1 Nov 2017 20:46
 Operator : SJ/JU
 Sample : I6203-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 102717-TIDO-E-D-B

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-BG101617.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	5.241	326	332	339	rBV	69028	103085	1.69%	0.429%
2	5.717	406	413	425	rBV	414815	669041	10.94%	2.788%
3	7.315	674	685	699	rBV	298187	544637	8.90%	2.269%
4	7.691	741	749	759	rBV	738093	1256927	20.55%	5.237%
5	8.161	821	829	835	rBV	216024	375053	6.13%	1.563%
6	8.484	876	884	894	rBV	977671	1710601	27.96%	7.127%
7	9.324	1018	1027	1036	rBV	771240	1411531	23.07%	5.881%
8	10.975	1297	1308	1319	rBV	331237	611268	9.99%	2.547%
9	13.402	1713	1721	1730	rBV	2529262	3973357	64.95%	16.555%
10	14.777	1947	1955	1961	rBV2	571355	1003185	16.40%	4.180%
11	14.835	1961	1965	1972	rVB	72333	112221	1.83%	0.468%
12	16.257	2200	2207	2225	rBV	1260268	1995689	32.62%	8.315%
13	17.515	2410	2421	2428	rBV2	739978	1156246	18.90%	4.817%
14	19.553	2762	2768	2773	rBV	89686	128838	2.11%	0.537%
15	20.106	2855	2862	2868	rBV	4497636	6117615	100.00%	25.489%
16	21.786	3140	3148	3162	rVB	843833	1455424	23.79%	6.064%
17	25.094	3701	3711	3723	rBV2	472905	1376690	22.50%	5.736%

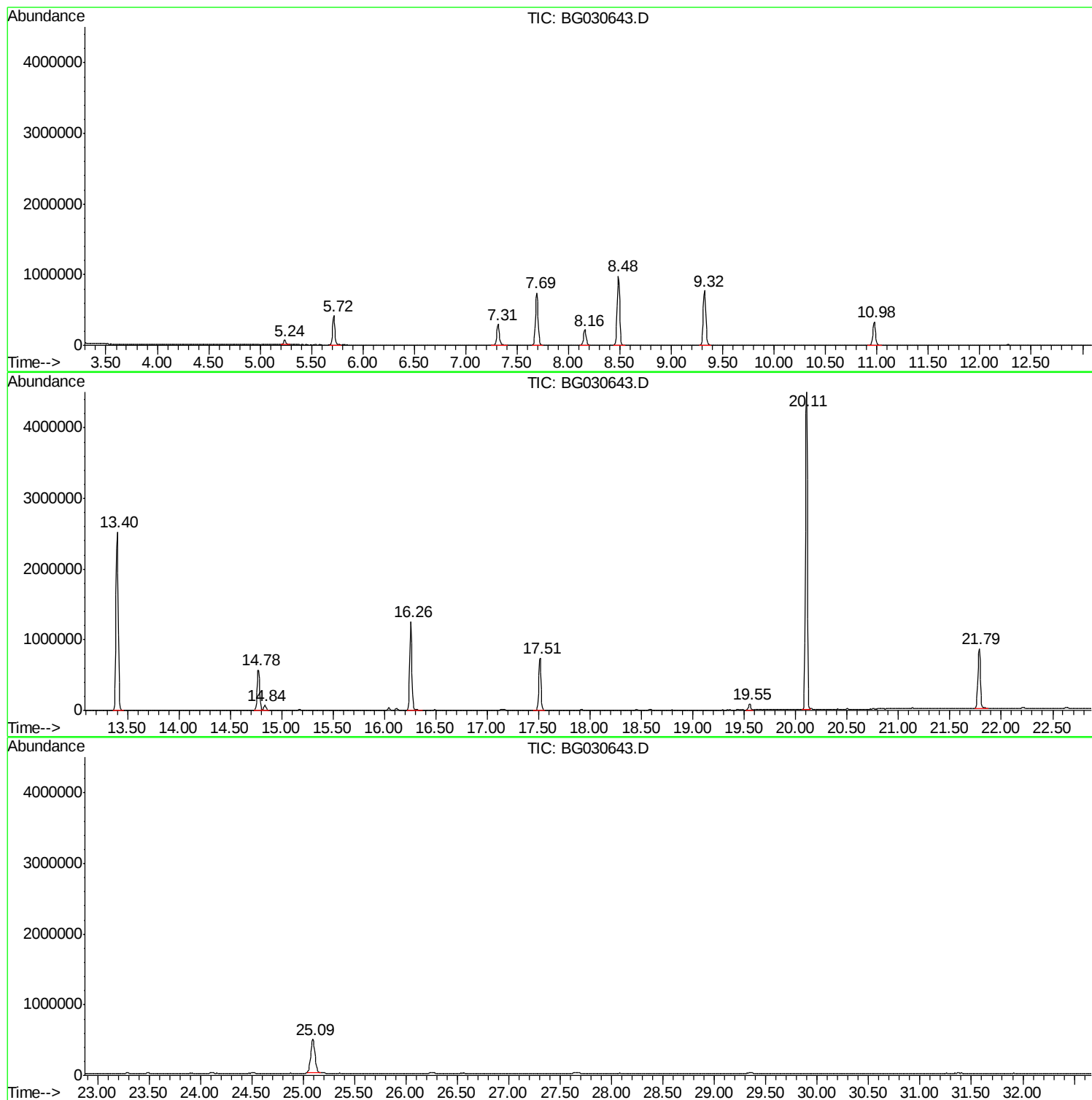
Sum of corrected areas: 24001408

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG101617.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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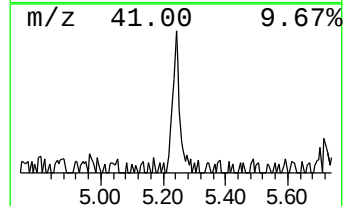
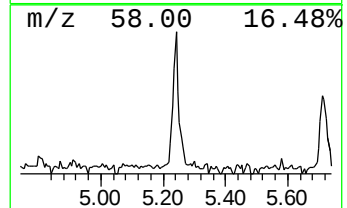
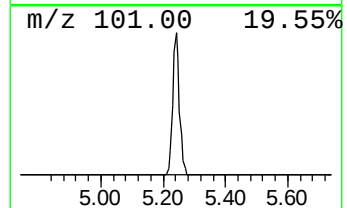
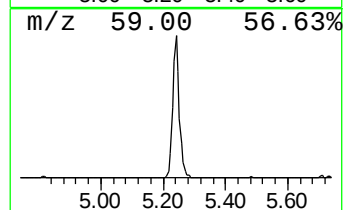
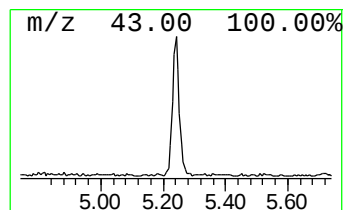
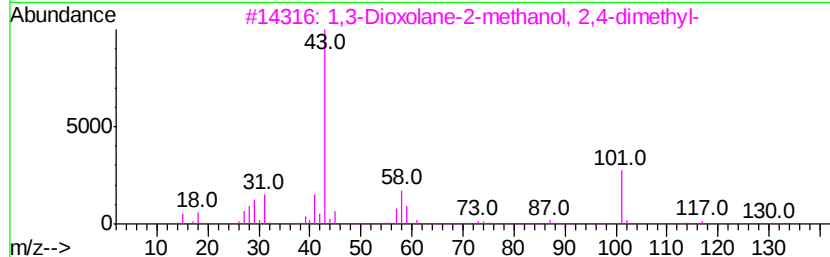
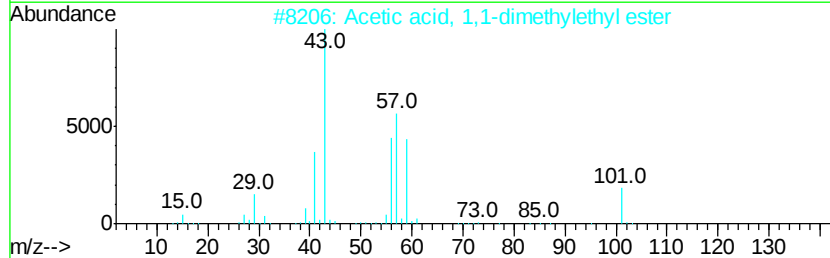
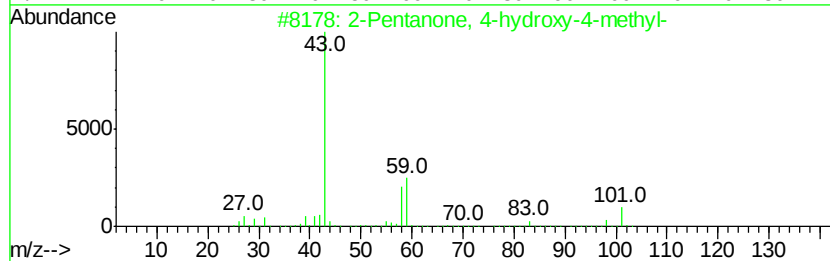
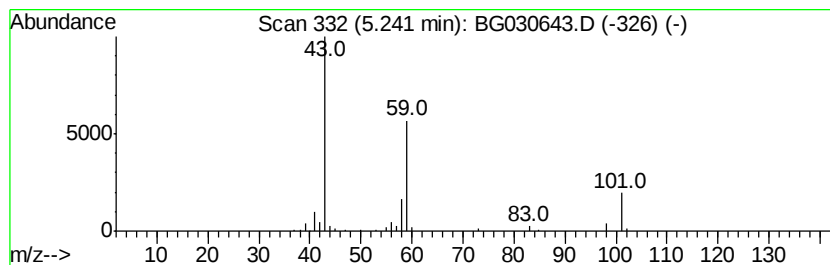
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.24	5.50 ng	103085	1,4-Dichlorobenzene-d4	8.16

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
3			1,3-Dioxolane-2-methanol, 2,4-di...	132	C6H12O3	053951-43-2	25
4			4-Hydroxy-3-hexanone	116	C6H12O2	004984-85-4	23
5			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17



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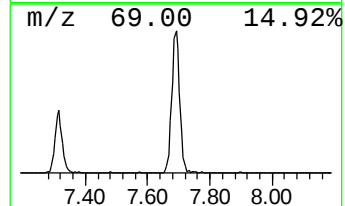
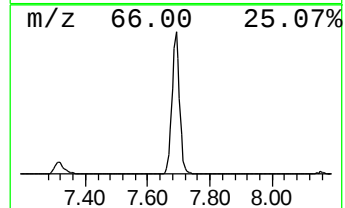
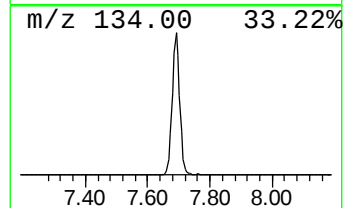
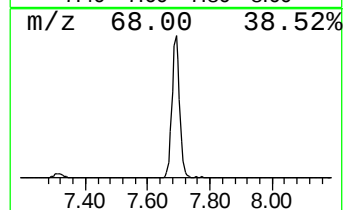
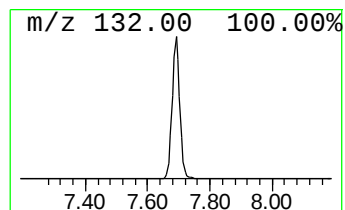
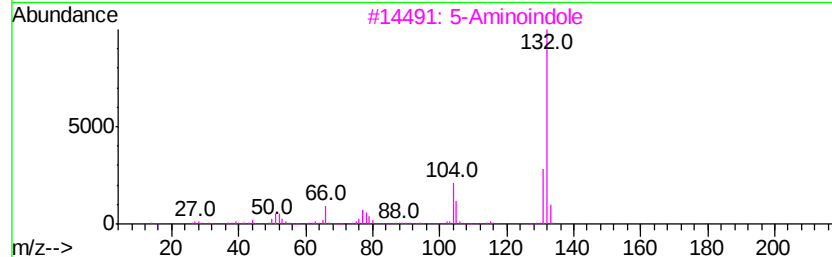
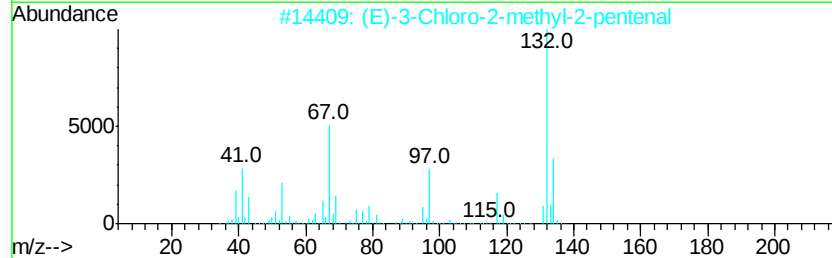
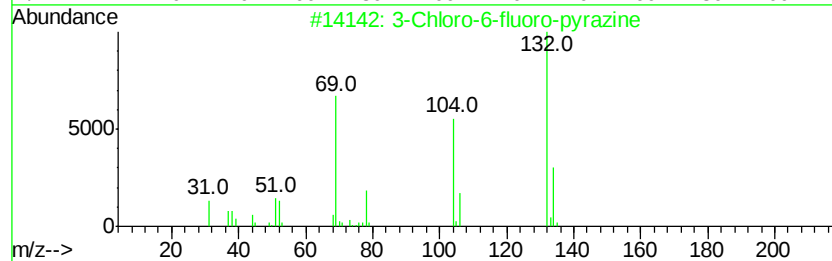
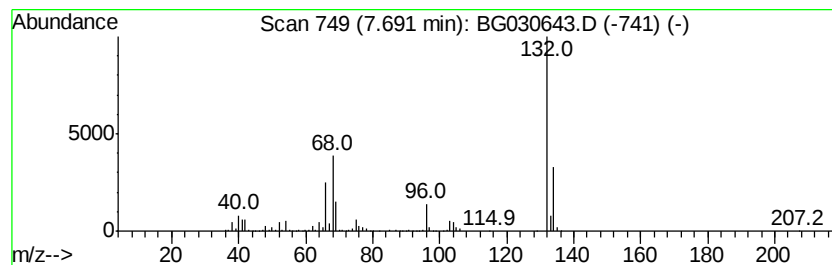
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Peak Number 2 unknown7.69 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.69	67.03 ng	1256930	1,4-Dichlorobenzene-d4	8.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.24	5.5	ng	103085	1	8.16	375053	20.0
unknown7.69	7.69	67.0	ng	1256930	1	8.16	375053	20.0