

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG110117\
 Data File : BG030647.D
 Acq On : 1 Nov 2017 23:16
 Operator : SJ/JU
 Sample : I6204-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 102617-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	5.239	326	332	340	rVB	79840	125348	2.11%	0.537%
2	5.715	406	413	427	rVB	412603	663355	11.15%	2.844%
3	7.313	678	685	696	rBV	265517	501282	8.42%	2.149%
4	7.689	742	749	760	rBV	716427	1243152	20.89%	5.329%
5	8.159	822	829	836	rBV	216410	369439	6.21%	1.584%
6	8.482	877	884	896	rVB	925619	1659155	27.88%	7.113%
7	9.323	1019	1027	1037	rBV	757964	1383507	23.25%	5.931%
8	10.974	1300	1308	1321	rBV	324977	599137	10.07%	2.568%
9	13.400	1713	1721	1732	rBV	2477510	3866117	64.97%	16.574%
10	14.775	1947	1955	1965	rVB2	569304	985923	16.57%	4.227%
11	16.256	2197	2207	2221	rBV	1279072	1988453	33.41%	8.524%
12	17.513	2414	2421	2432	rBV2	756121	1173872	19.73%	5.032%
13	20.104	2856	2862	2868	rBV	4535431	5950866	100.00%	25.511%
14	21.785	3141	3148	3161	rVB	823396	1415891	23.79%	6.070%
15	25.092	3701	3711	3730	rVB2	473980	1400873	23.54%	6.006%

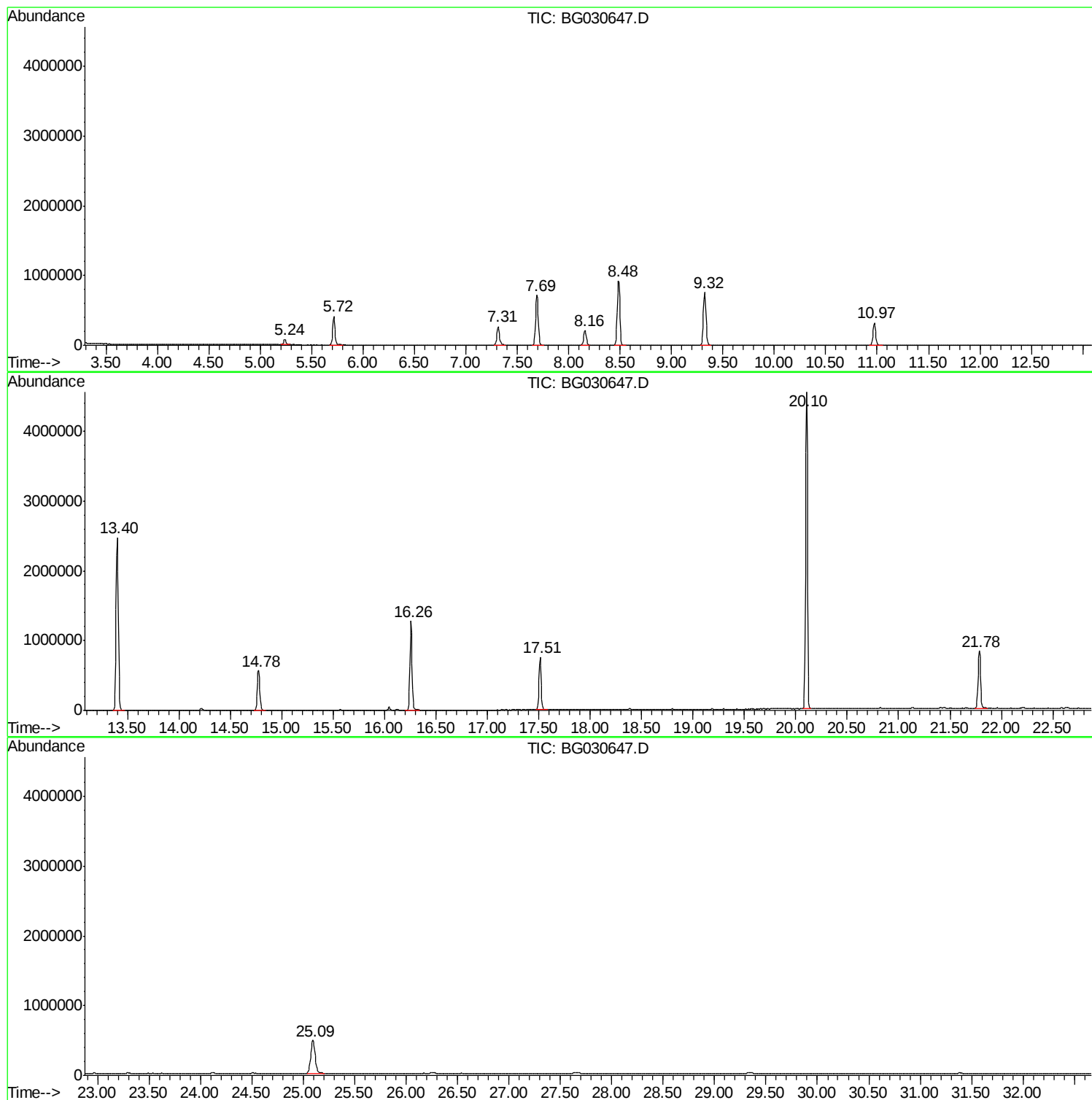
Sum of corrected areas: 23326370

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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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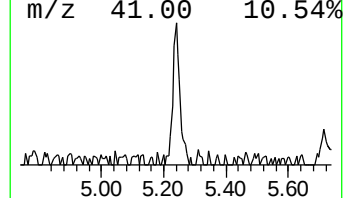
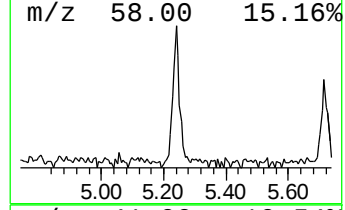
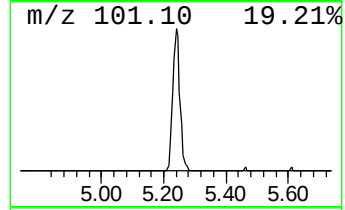
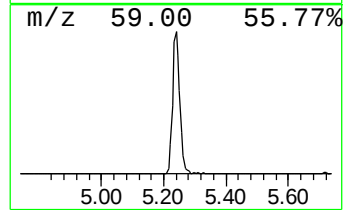
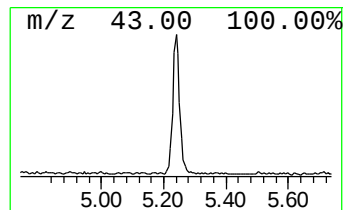
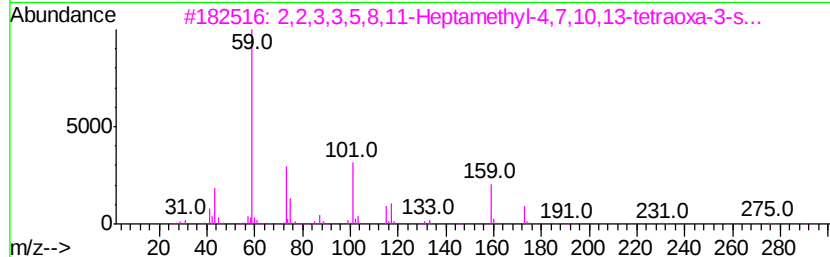
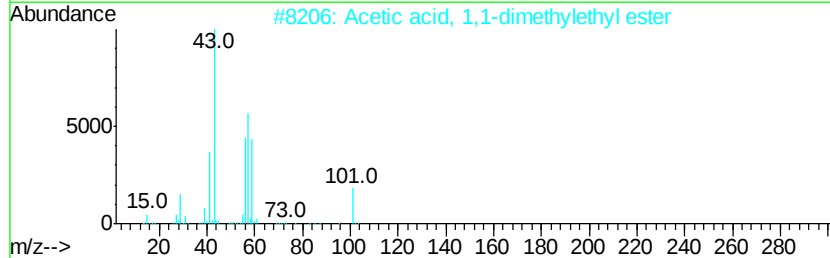
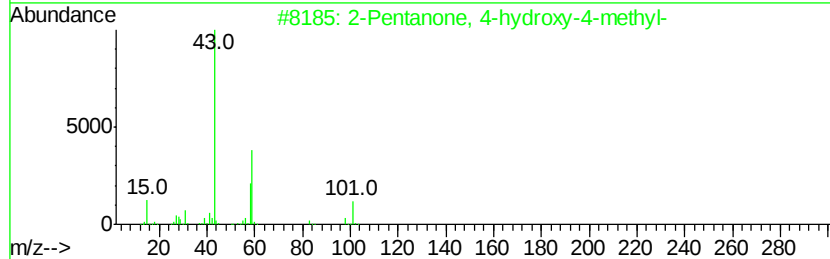
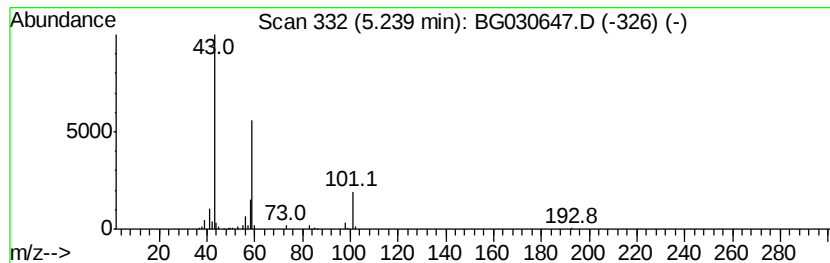
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.24	6.79 ng	125348	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	50
3			2,2,3,3,5,8,11-Heptamethyl-4,7,1...	348	C18H40O4Si	1000367-07-2	28
4			4-Ethyl-3-octanol	158	C10H22O	019781-26-1	25
5			Tetraglyme	222	C10H22O5	000143-24-8	25



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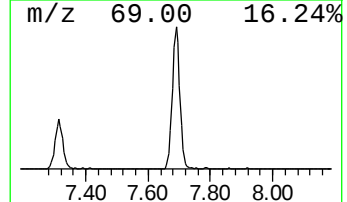
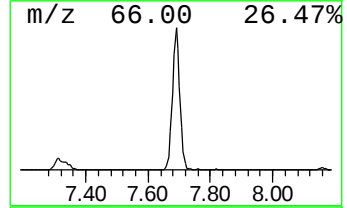
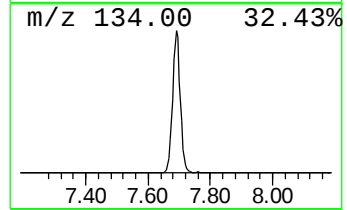
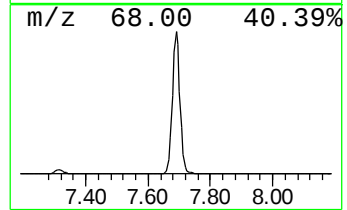
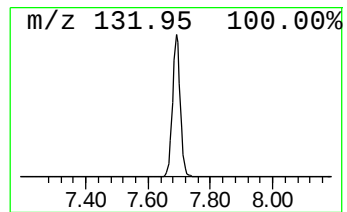
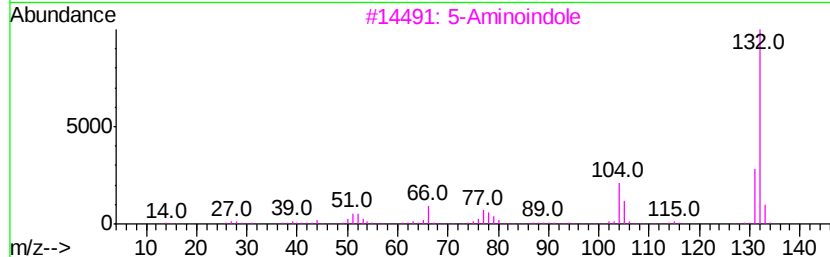
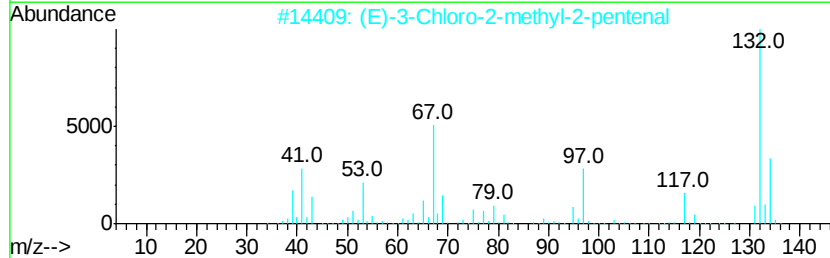
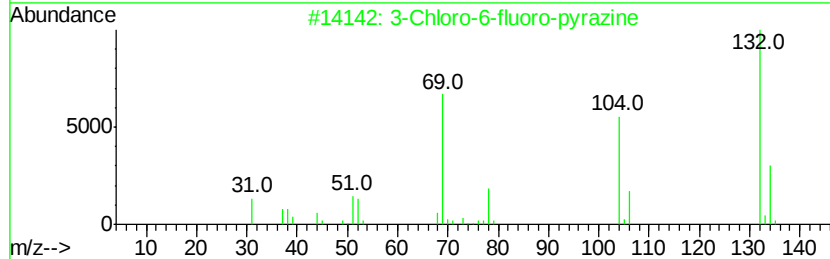
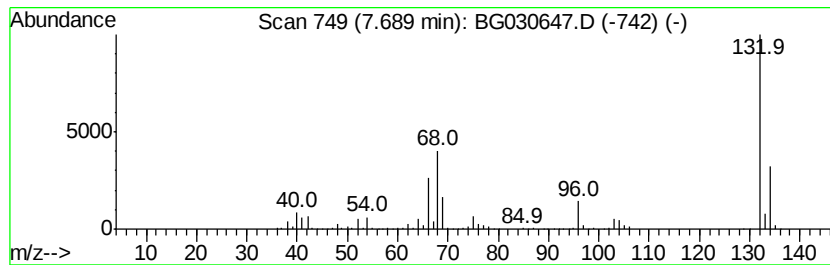
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Peak Number 2 unknown7.69 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.69	67.30 ng	1243150	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	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
5		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.24	6.8	ng	125348	1	8.16	369439	20.0
unknown7.69	7.69	67.3	ng	1243150	1	8.16	369439	20.0