

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102518\
 Data File : BG037570.D
 Acq On : 26 Oct 2018 8:02
 Operator : JU/SJ
 Sample : J5655-05
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 S37-0003

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:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG101818.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	5.180	317	322	330	rVB2	22409	40896	1.11%	0.232%
2	5.645	394	401	412	rBV	390640	655251	17.82%	3.713%
3	7.243	665	673	687	rBV	256265	458865	12.48%	2.600%
4	7.625	730	738	756	rVB	707795	1308405	35.58%	7.414%
5	8.106	812	820	829	rVB	177944	328502	8.93%	1.861%
6	8.424	866	874	883	rBV	632031	1146205	31.17%	6.495%
7	9.281	1011	1020	1028	rBV	590695	1085922	29.53%	6.153%
8	10.921	1290	1299	1311	rVB	266699	510253	13.88%	2.891%
9	13.359	1706	1714	1726	rBV	1618549	2636150	71.69%	14.938%
10	14.734	1939	1948	1960	rBV2	465076	794423	21.61%	4.502%
11	16.220	2194	2201	2220	rBV	1257007	1976709	53.76%	11.201%
12	17.484	2409	2416	2427	rVB2	626091	975305	26.52%	5.527%
13	20.081	2852	2858	2864	rBV	2834354	3676963	100.00%	20.835%
14	21.767	3137	3145	3152	rBV	568667	1015053	27.61%	5.752%
15	24.088	3535	3540	3548	rVB6	19167	44956	1.22%	0.255%
16	25.098	3699	3712	3730	rBV3	307047	993865	27.03%	5.632%

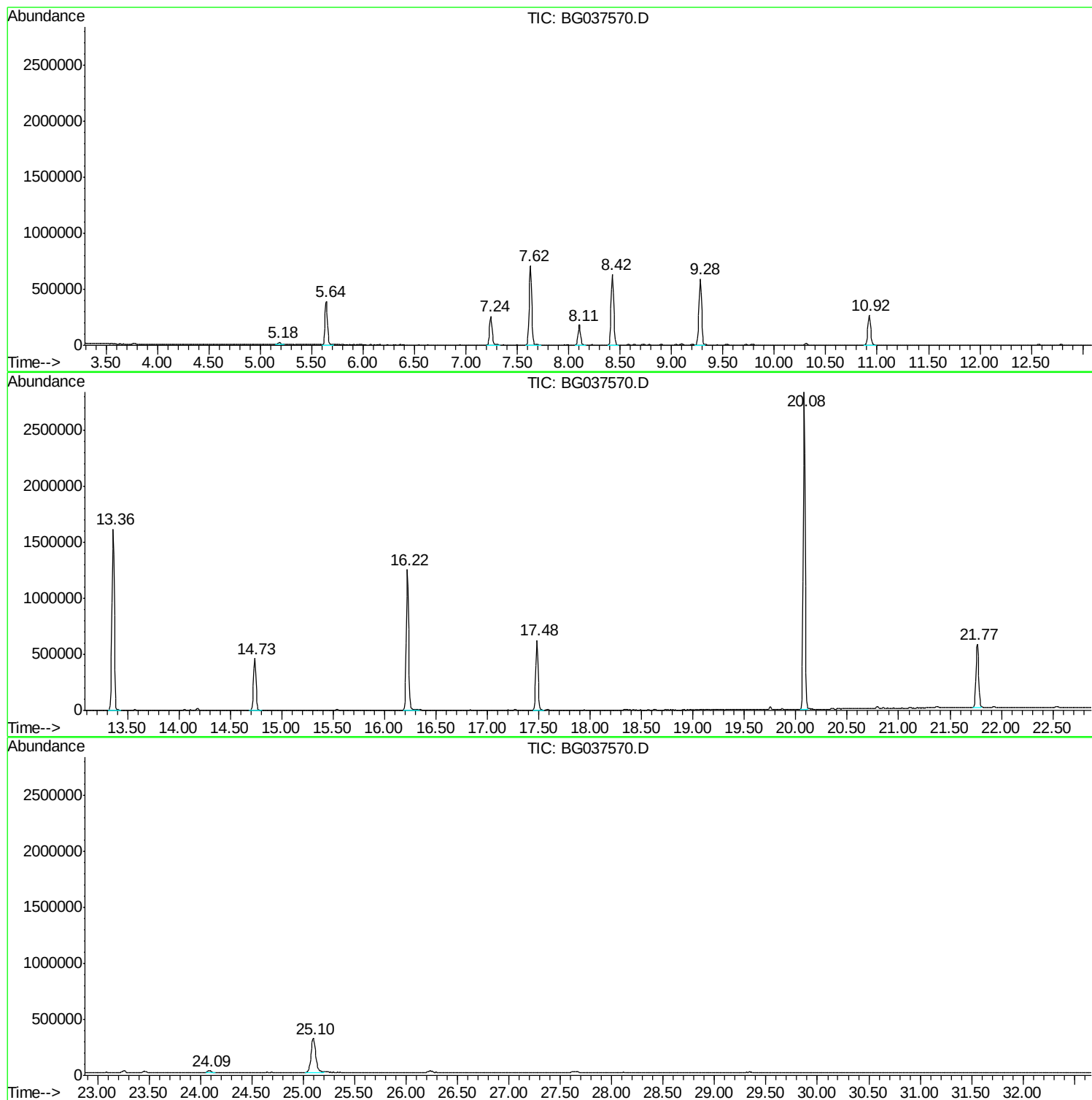
Sum of corrected areas: 17647723

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG101818.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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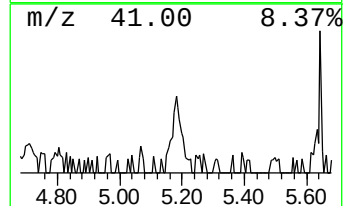
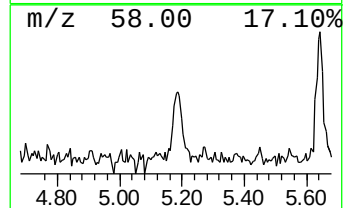
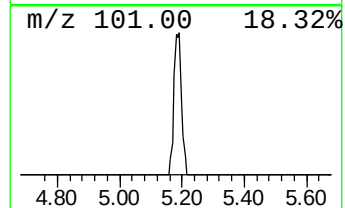
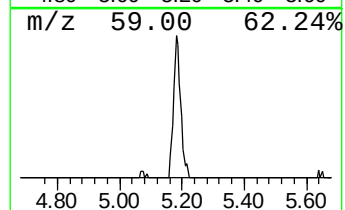
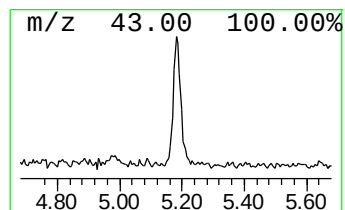
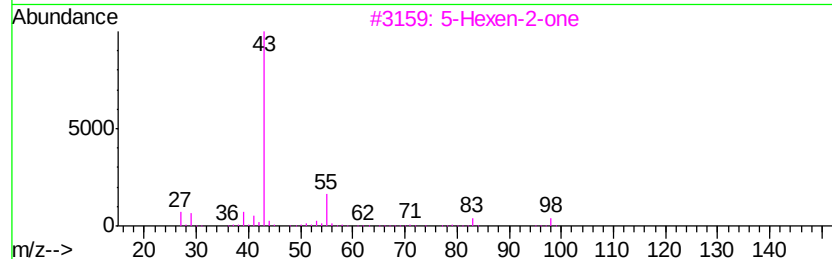
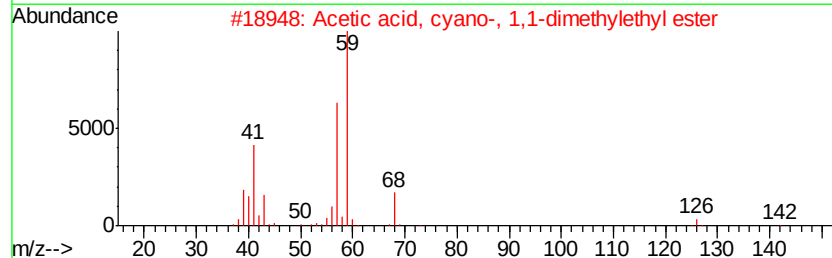
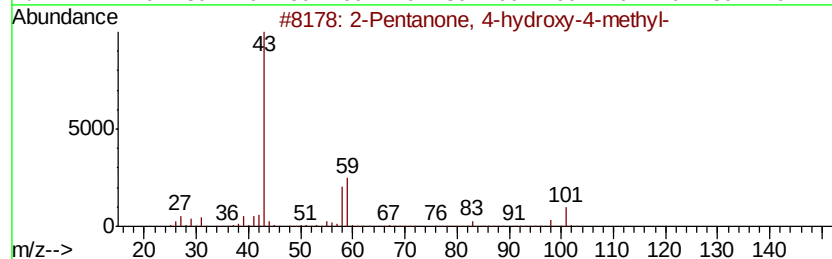
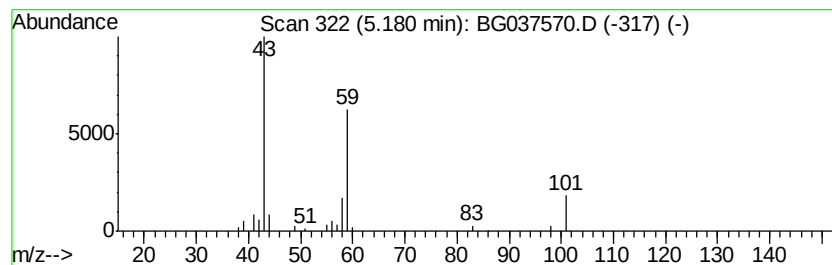
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.18	2.49 ng	40896	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	64
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
3		5-Hexen-2-one	98	C6H10O	000109-49-9	9
4		Diacetamide	101	C4H7NO2	000625-77-4	9
5		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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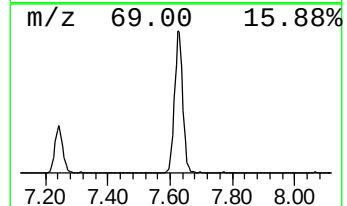
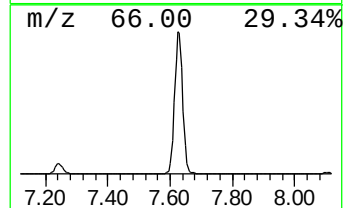
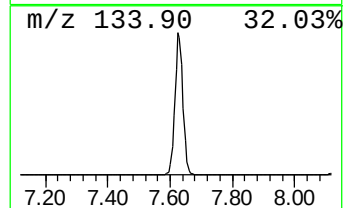
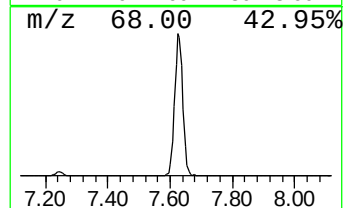
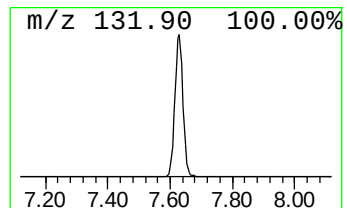
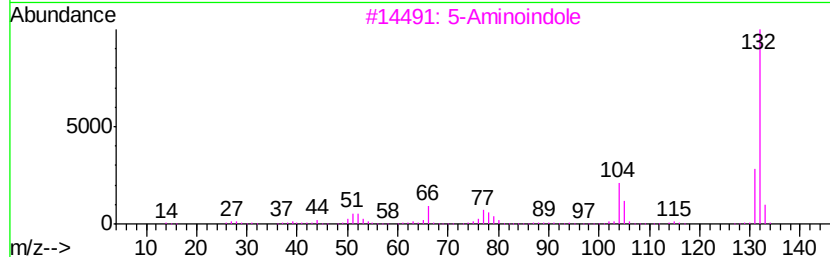
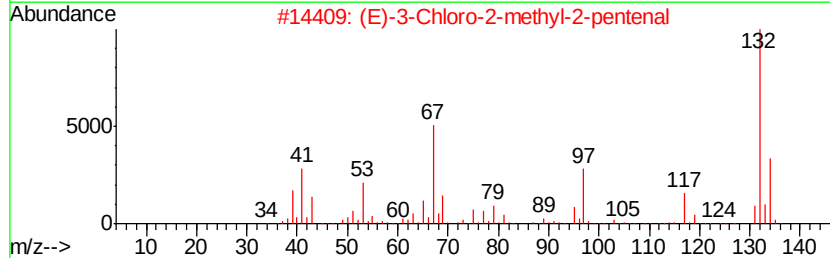
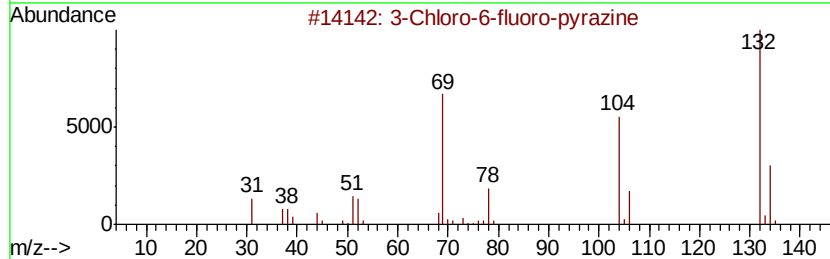
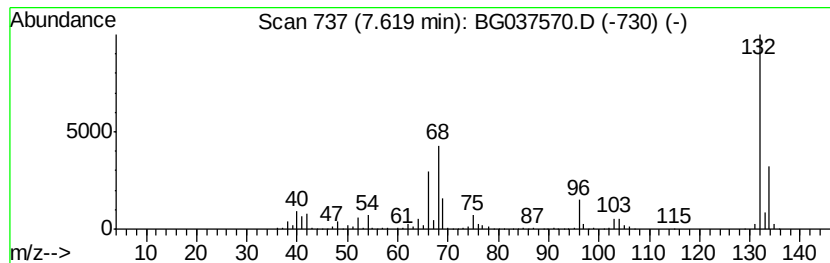
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Peak Number 2 unknown7.62 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.62	79.66 ng	1308410	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	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	27
3		5-Aminoindole	132	C8H8N2	005192-03-0	10
4		Xenon	132	Xe	007440-63-3	9
5		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.18	2.5	ng	40896	1	8.10	328502	20.0
unknown7.62	7.62	79.7	ng	1308410	1	8.10	328502	20.0