

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG082718\
 Data File : BG036433.D
 Acq On : 27 Aug 2018 16:03
 Operator : JU/SJ
 Sample : J4522-09
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 FB081618

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-BG082318.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	5.167	315	320	330	rVB	40021	60247	1.08%	0.270%
2	5.631	391	399	409	rBV	500619	785558	14.13%	3.520%
3	7.218	660	669	679	rBV	306879	527992	9.49%	2.366%
4	7.594	726	733	742	rBV	926268	1579386	28.40%	7.077%
5	8.064	806	813	820	rBV	206784	350273	6.30%	1.570%
6	8.387	859	868	879	rVB	869717	1520813	27.35%	6.815%
7	9.227	1003	1011	1023	rBV	737975	1312489	23.60%	5.881%
8	10.872	1283	1291	1298	rVB	283006	499663	8.98%	2.239%
9	13.316	1699	1707	1717	rBV	2101670	3259764	58.62%	14.607%
10	14.691	1933	1941	1949	rVB2	443999	730653	13.14%	3.274%
11	16.172	2186	2193	2210	rBV	1631764	2446556	43.99%	10.963%
12	17.429	2400	2407	2415	rBV	658139	982589	17.67%	4.403%
13	20.044	2845	2852	2860	rBV	4072842	5561092	100.00%	24.920%
14	21.695	3126	3133	3142	rBV2	765926	1222240	21.98%	5.477%
15	21.912	3165	3170	3175	rBV	43846	61380	1.10%	0.275%
16	22.523	3270	3274	3286	rVB3	43690	86716	1.56%	0.389%
17	23.222	3383	3393	3399	rBV4	43248	102494	1.84%	0.459%
18	24.039	3525	3532	3539	rVB3	33877	70938	1.28%	0.318%
19	24.915	3670	3681	3690	rBV2	410661	1155154	20.77%	5.176%

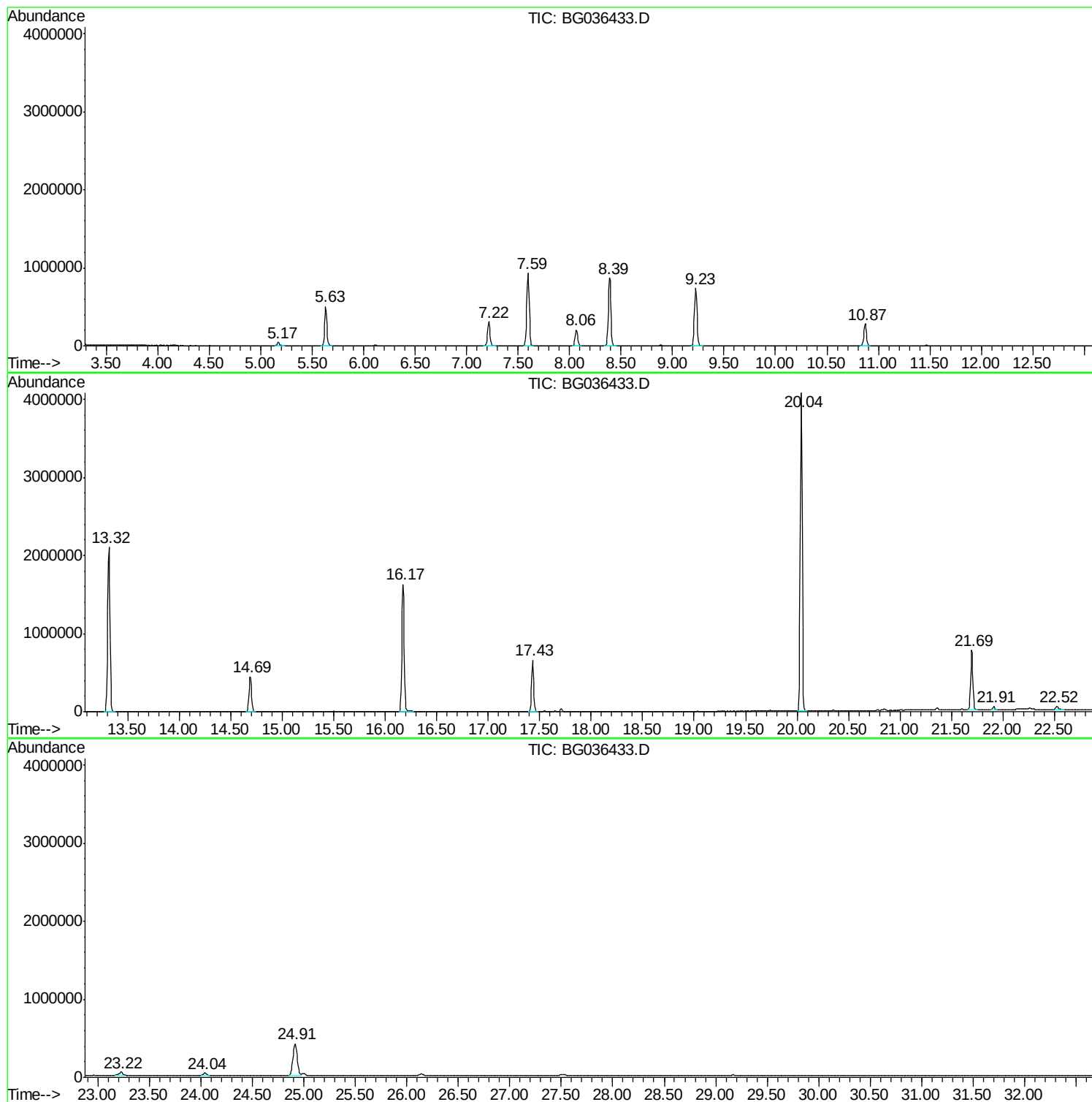
Sum of corrected areas: 22315997

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082318.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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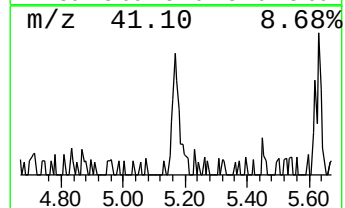
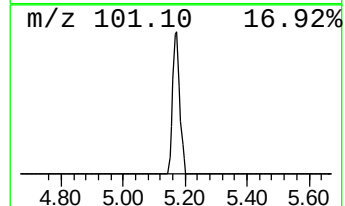
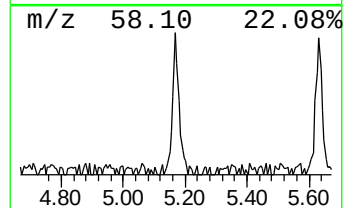
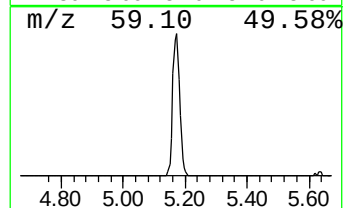
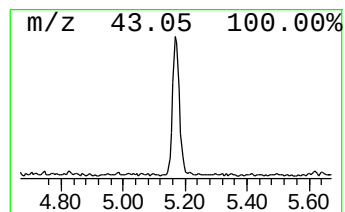
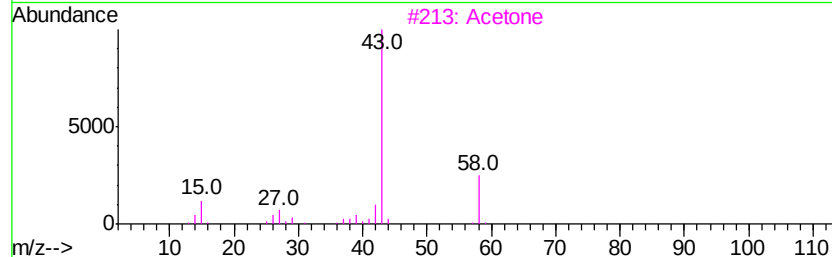
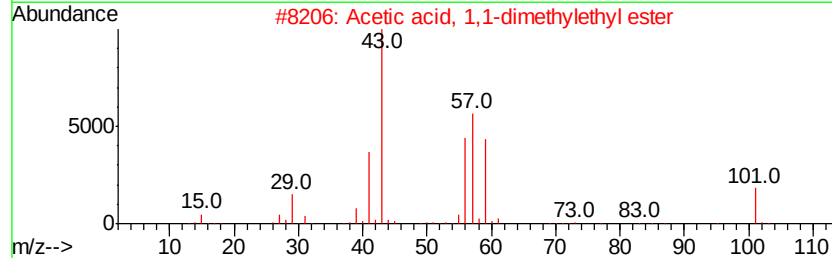
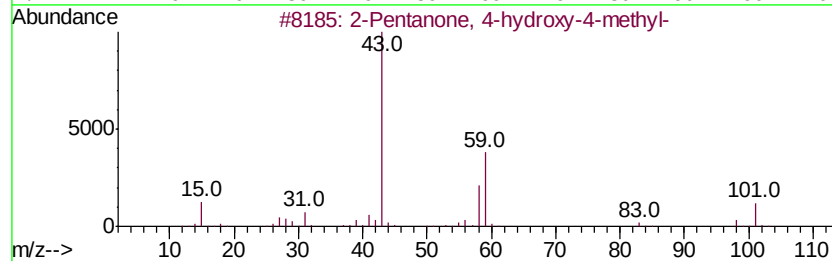
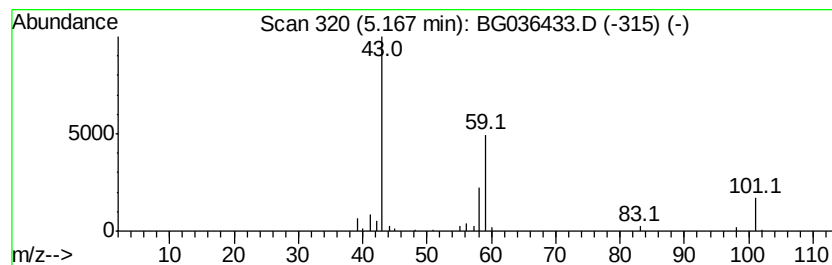
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.17	3.44 ng	60247	1,4-Dichlorobenzene-d4	8.06

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	78
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	36
3			Acetone	58	C3H6O	000067-64-1	9
4			4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	9
5			Formamide, N,N-diethyl-	101	C5H11NO	000617-84-5	9



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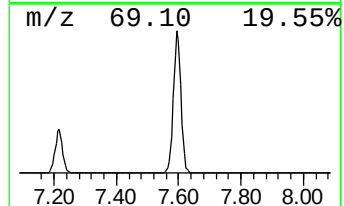
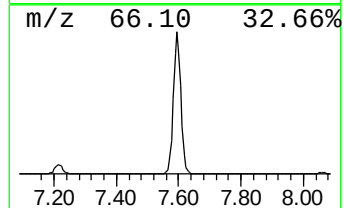
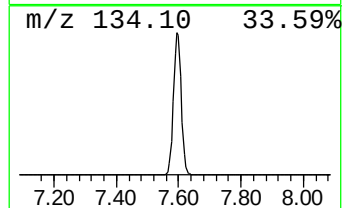
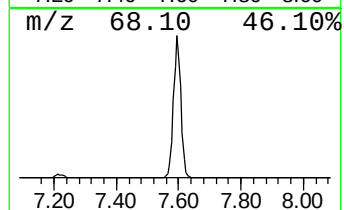
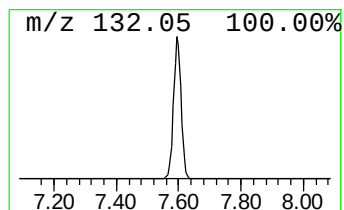
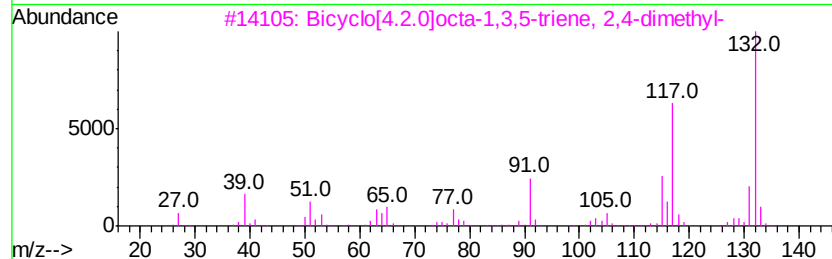
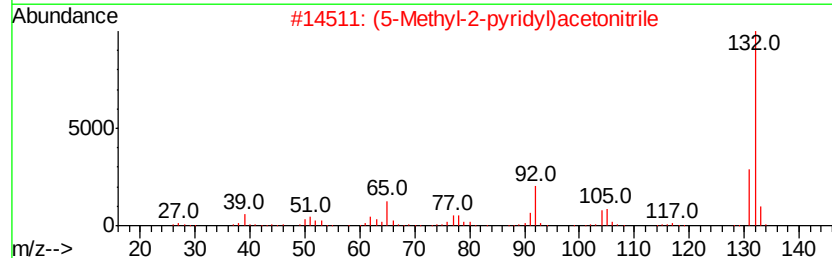
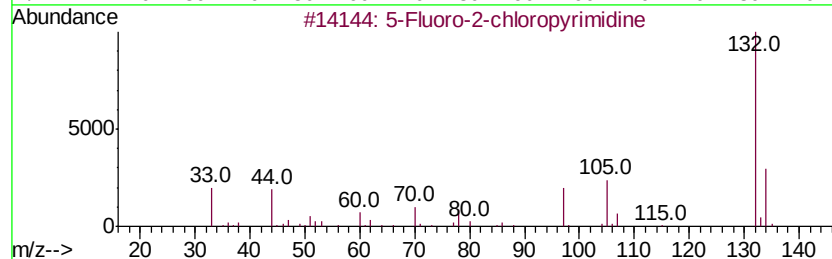
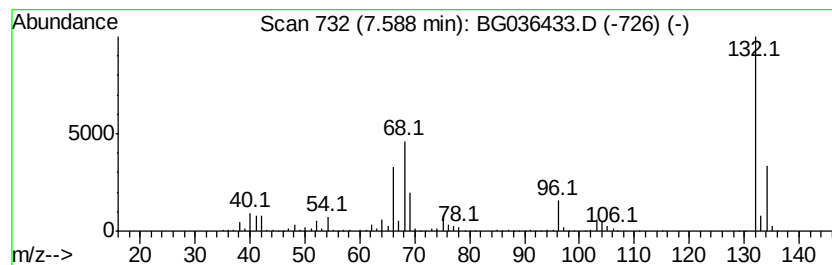
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Peak Number 2 unknown7.59 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.59	90.18 ng	1579390	1,4-Dichlorobenzene-d4	8.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
2		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		Xenon	132	Xe	007440-63-3	9
5		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.17	3.4	ng	60247	1	8.06	350273	20.0
unknown7.59	7.59	90.2	ng	1579390	1	8.06	350273	20.0