

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG042919\
 Data File : BG040666.D
 Acq On : 29 Apr 2019 17:56
 Operator : JU/SJ
 Sample : PB119264BL
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB119264BL

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-BG042319.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	3.153	6	11	17	rBV2	55988	99344	2.48%	0.553%
2	5.221	355	363	369	rBV	71061	112278	2.80%	0.625%
3	5.680	433	441	449	rBV	621661	965462	24.07%	5.377%
4	7.284	706	714	723	rBV	647636	1089191	27.16%	6.066%
5	7.671	772	780	788	rBV	591946	1020586	25.45%	5.684%
6	8.153	855	862	869	rVB	191872	330800	8.25%	1.842%
7	8.476	909	917	925	rBV	650317	1102957	27.50%	6.143%
8	9.328	1053	1062	1071	rBV	568072	1000229	24.94%	5.571%
9	10.973	1335	1342	1350	rBV	241667	443330	11.05%	2.469%
10	13.406	1747	1756	1764	rBV	1548734	2318766	57.82%	12.914%
11	14.781	1984	1990	1997	rBV2	444673	691454	17.24%	3.851%
12	16.261	2236	2242	2252	rBV	967340	1462343	36.46%	8.145%
13	17.525	2449	2457	2465	rBV2	680666	1026079	25.59%	5.715%
14	20.121	2892	2899	2904	rBV	3044732	4010443	100.00%	22.336%
15	21.808	3179	3186	3195	rBV2	691743	1162008	28.97%	6.472%
16	25.174	3749	3759	3774	rVB2	379514	1119622	27.92%	6.236%

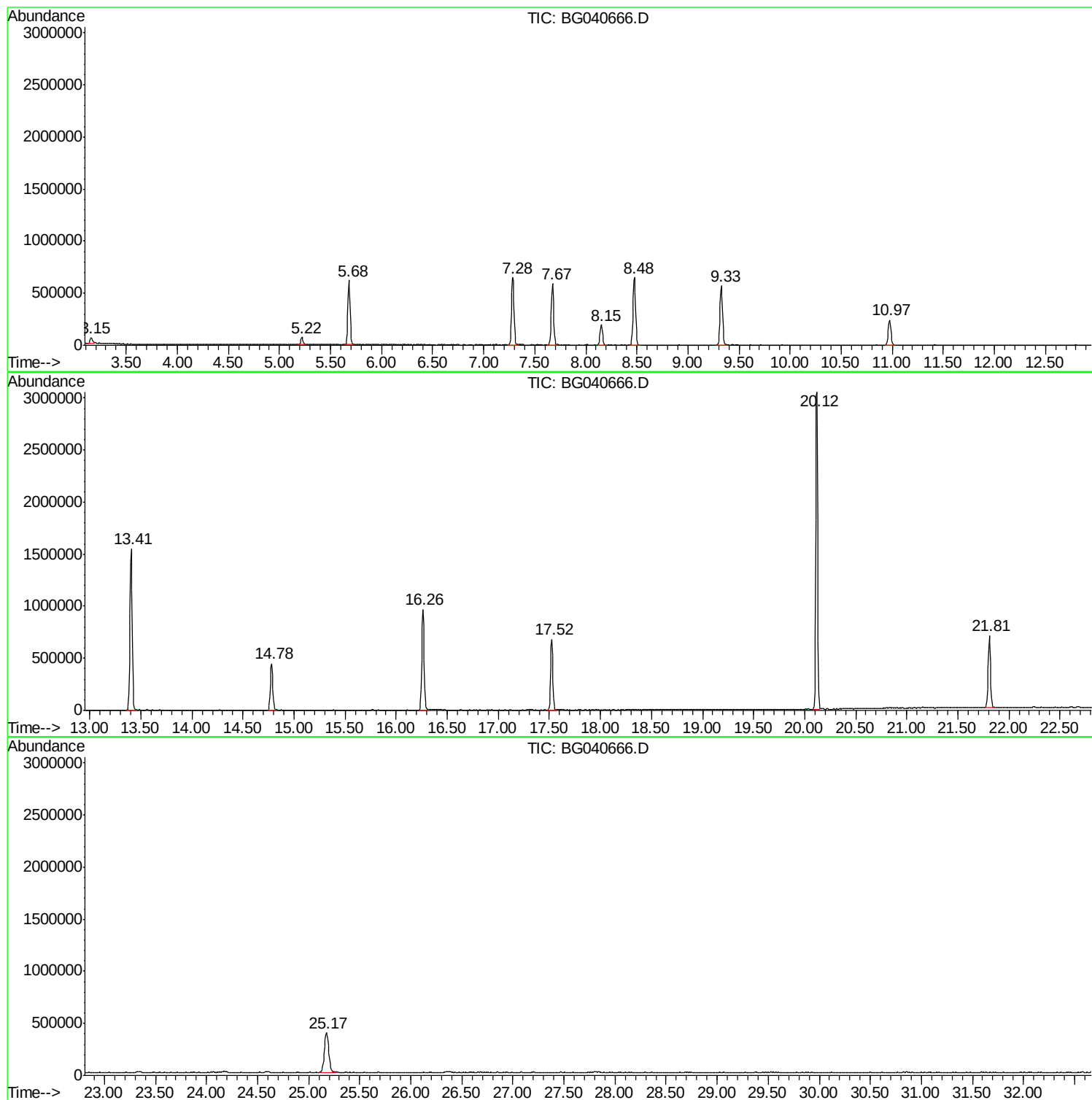
Sum of corrected areas: 17954892

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG042319.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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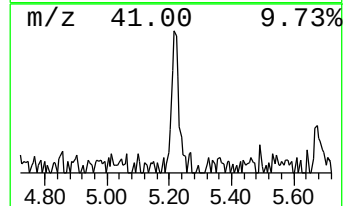
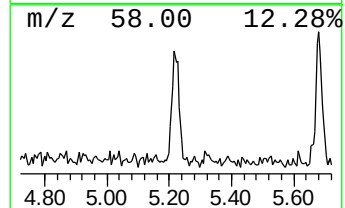
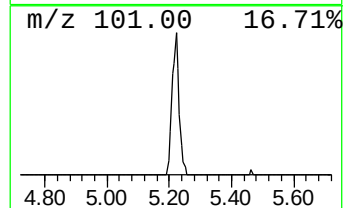
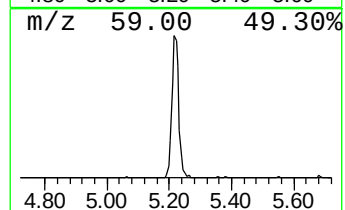
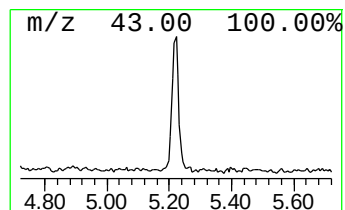
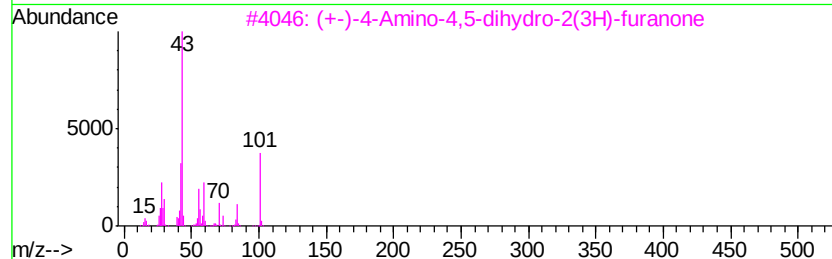
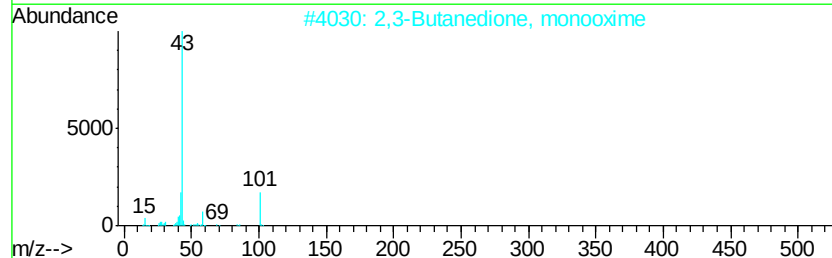
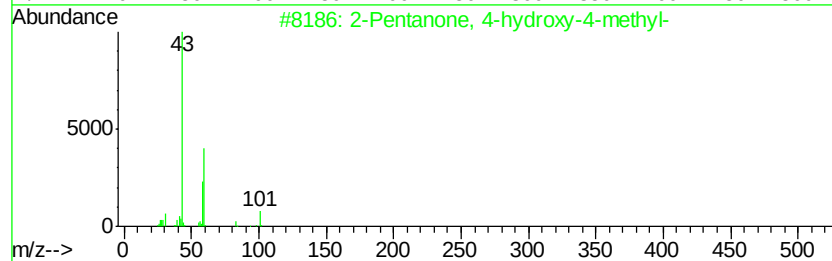
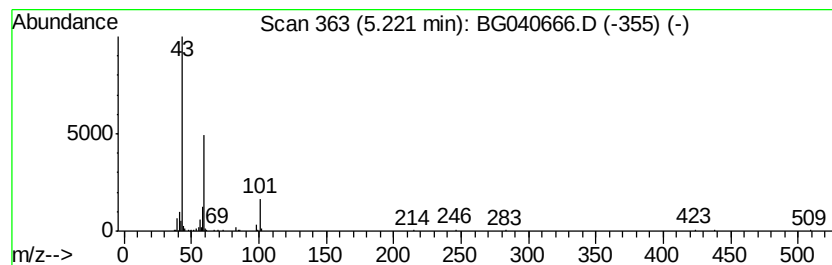
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.22	6.79 ng	112278	1,4-Dichlorobenzene-d4	8.15

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
3			(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9
4			Acetamide, N-(aminocarbonyl)-	102	C3H6N2O2	000591-07-1	9
5			Propanenitrile, 3-ethoxy-	99	C5H9NO	002141-62-0	9



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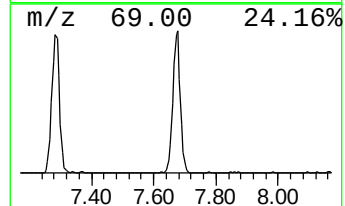
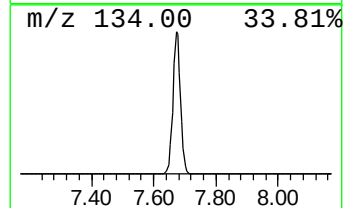
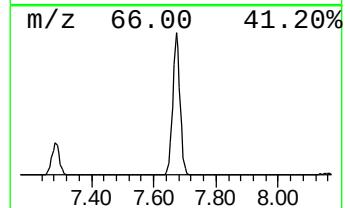
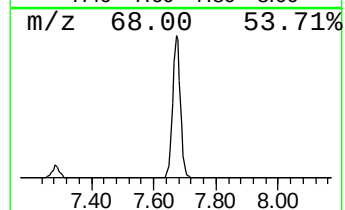
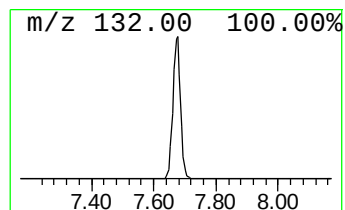
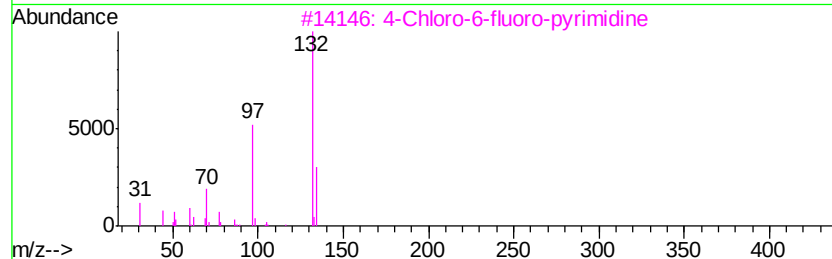
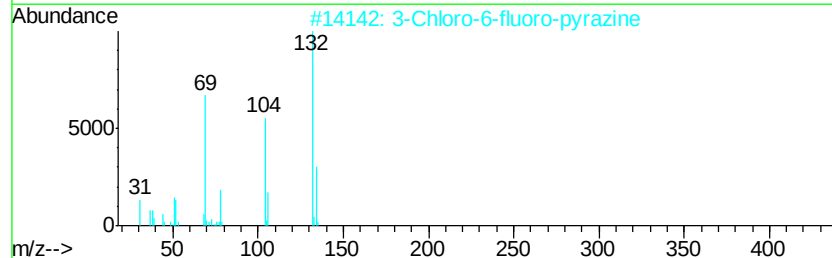
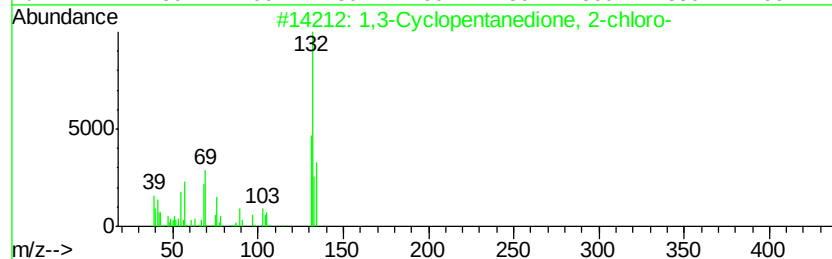
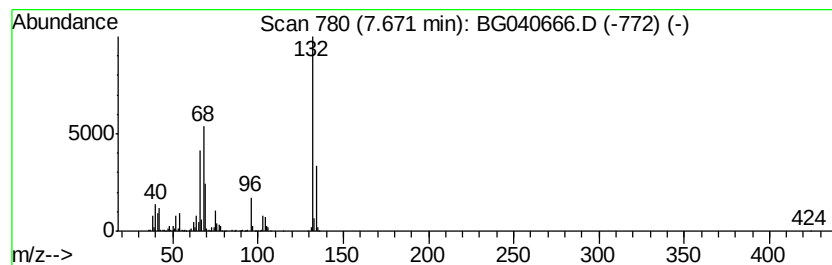
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Peak Number 3 unknown7.67 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.67	61.70 ng	1020590	1,4-Dichlorobenzene-d4	8.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	22
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	14
3		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	14
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.22	6.8	ng	112278	1	8.15	330800	20.0
unknown7.67	7.67	61.7	ng	1020590	1	8.15	330800	20.0