

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG070218\
 Data File : BG035566.D
 Acq On : 3 Jul 2018 00:52
 Operator : JU/SJ
 Sample : PB110755BL
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB110755BL

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-BG060718.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	5.173	315	321	329	rBV	80239	123837	3.00%	0.755%
2	5.643	395	401	414	rVB	749063	1201910	29.07%	7.330%
3	7.229	663	671	682	rBV	774593	1336063	32.31%	8.148%
4	7.599	727	734	748	rBV	735708	1259912	30.47%	7.684%
5	8.063	805	813	821	rBV	123954	205490	4.97%	1.253%
6	8.386	861	868	879	rBV	545144	948319	22.94%	5.783%
7	9.226	1003	1011	1018	rBV	441868	793602	19.19%	4.840%
8	10.871	1283	1291	1297	rBV	141286	266102	6.44%	1.623%
9	13.309	1697	1706	1718	rBV	1320574	2054106	49.68%	12.527%
10	14.684	1932	1940	1950	rBV3	248756	414941	10.04%	2.531%
11	16.170	2186	2193	2205	rBV	980109	1581528	38.25%	9.645%
12	17.421	2400	2406	2413	rBV2	383809	600625	14.53%	3.663%
13	20.029	2843	2850	2861	rBV	3120904	4134568	100.00%	25.215%
14	21.692	3125	3133	3143	rBV	468833	775559	18.76%	4.730%
15	24.911	3671	3681	3695	rBV2	232268	700659	16.95%	4.273%

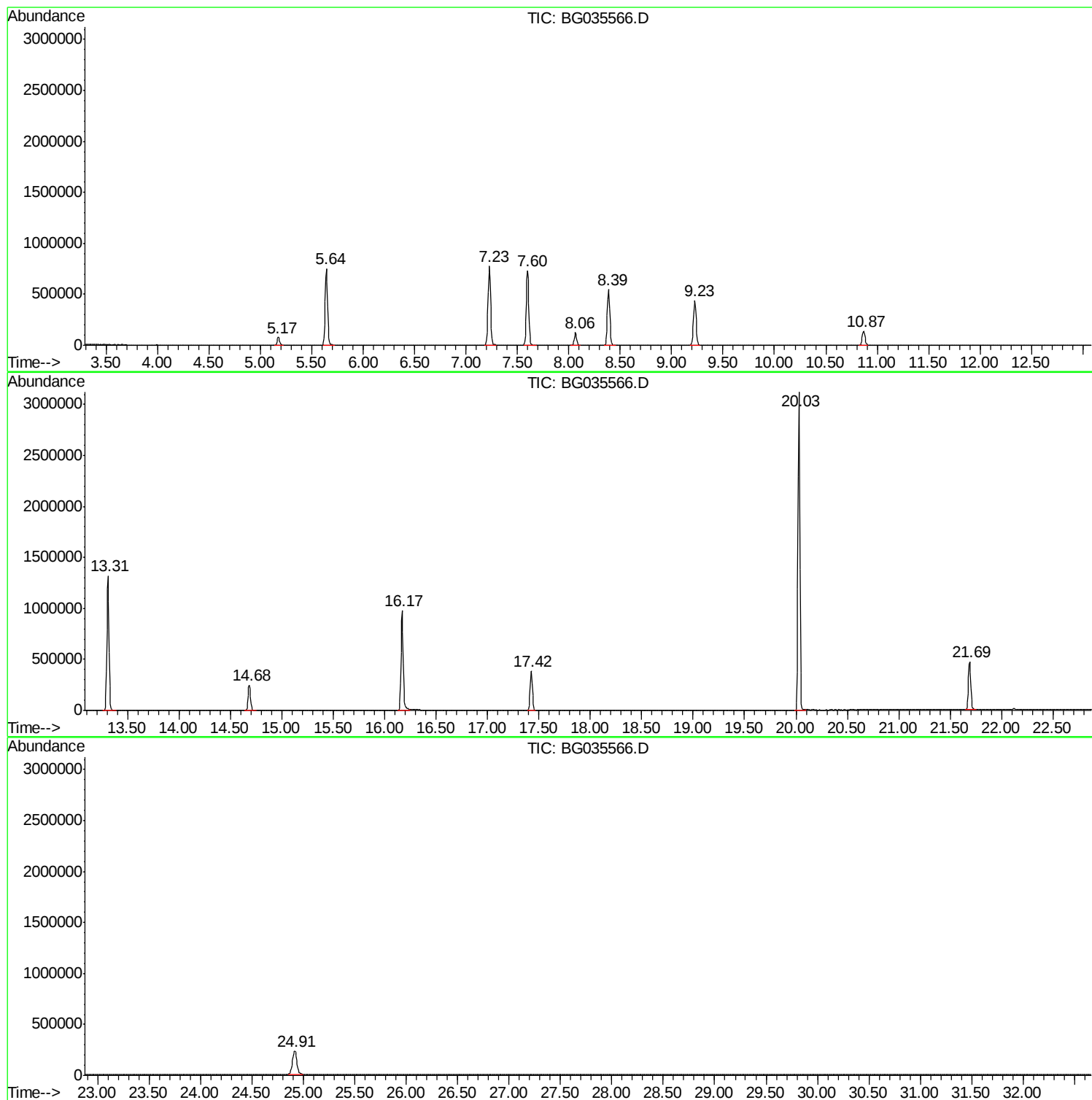
Sum of corrected areas: 16397221

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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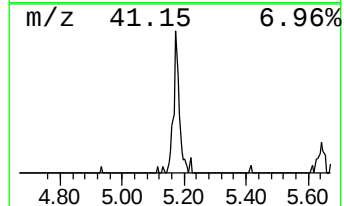
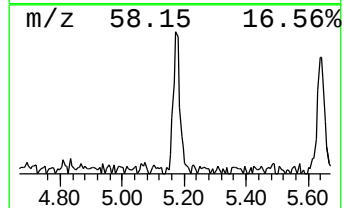
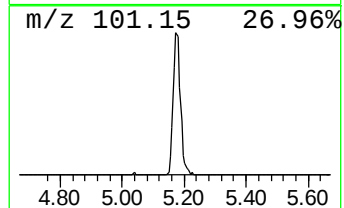
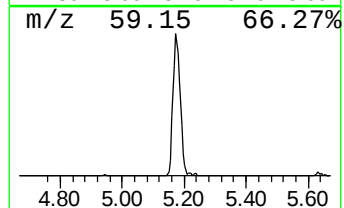
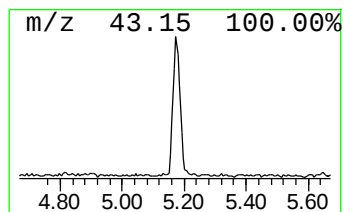
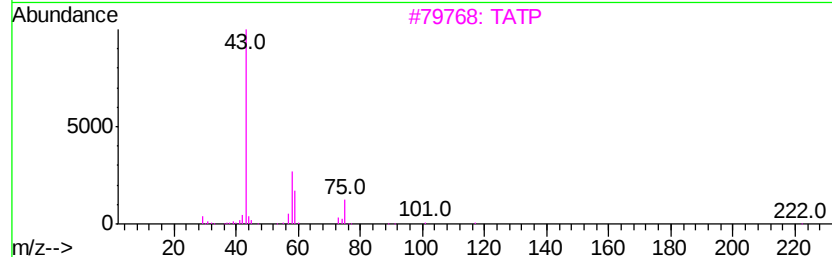
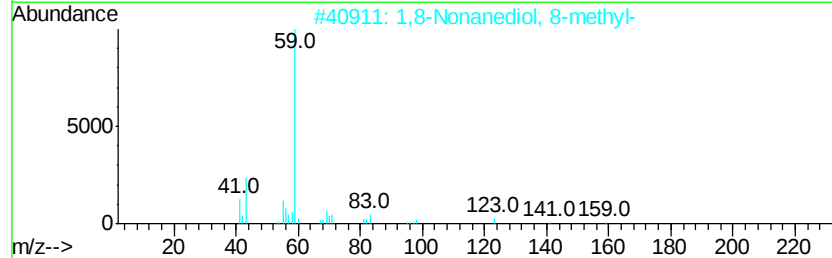
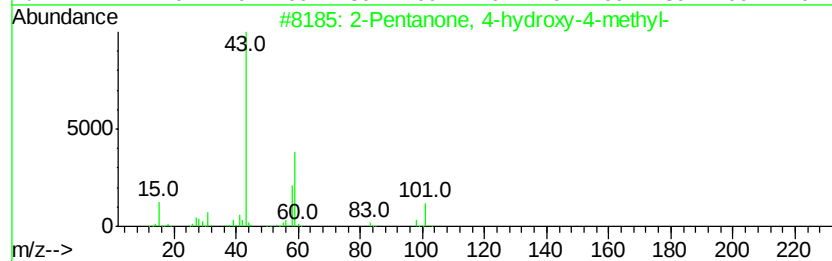
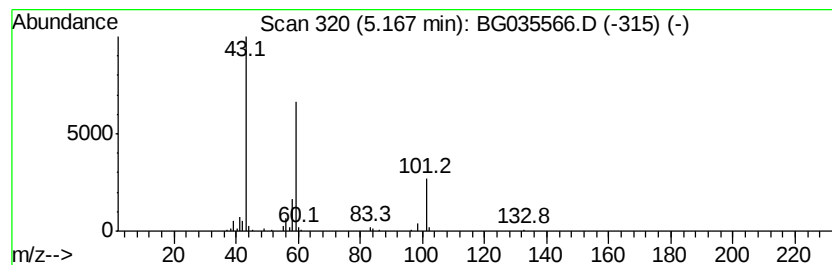
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.17	12.05 ng	123837	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	36
2	1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	23
3	TATP	222	C9H18O6	017088-37-8	12
4	Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	12
5	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10



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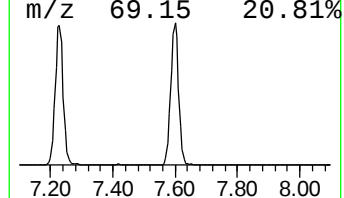
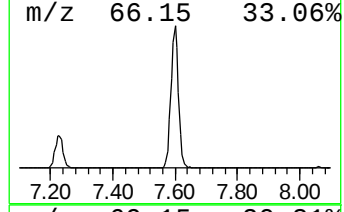
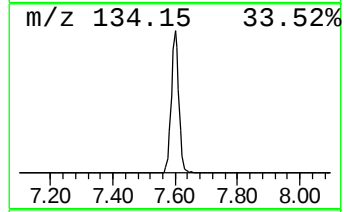
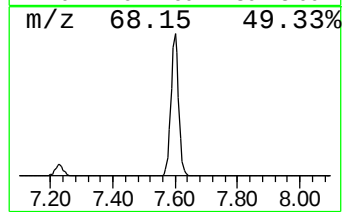
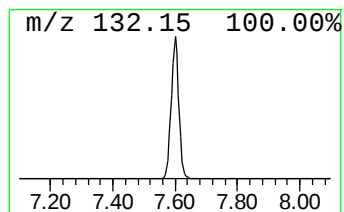
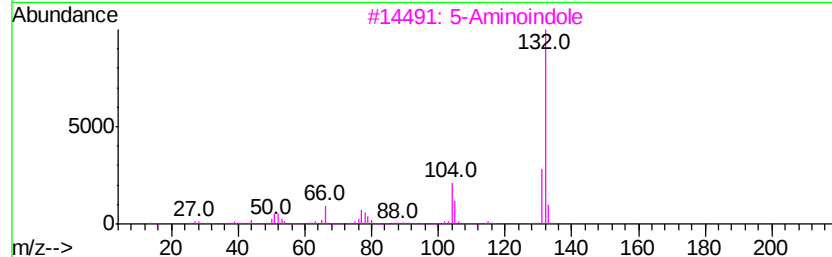
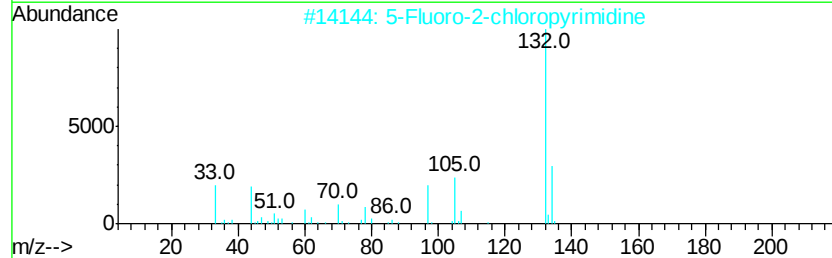
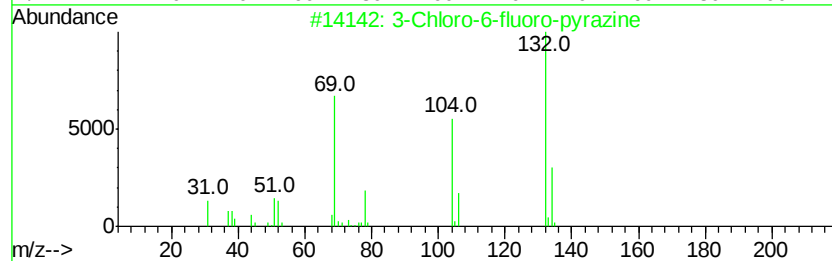
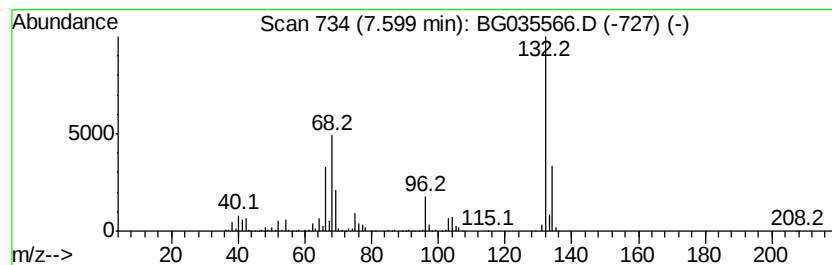
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Peak Number 2 unknown7.60 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.60	122.63 ng	1259910	1,4-Dichlorobenzene-d4	8.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	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...	5.17	12.1	ng	123837	1	8.06	205490	20.0
unknown7.60	7.60	122.6	ng	1259910	1	8.06	205490	20.0