

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG062918\
 Data File : BG035509.D
 Acq On : 29 Jun 2018 11:42
 Operator : JU/SJ
 Sample : PB110463BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB110463BL

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	5.175	315	321	329	rVB	31835	49058	1.62%	0.405%
2	5.639	392	400	410	rBV	491102	769559	25.38%	6.361%
3	7.225	663	670	680	rBV	499081	868017	28.62%	7.174%
4	7.595	726	733	744	rBV	478975	813430	26.82%	6.723%
5	8.065	807	813	820	rVB	101134	172196	5.68%	1.423%
6	8.383	860	867	878	rVB	339139	599333	19.76%	4.954%
7	9.223	1001	1010	1024	rBV	289449	522932	17.24%	4.322%
8	10.867	1281	1290	1299	rBV	124615	238713	7.87%	1.973%
9	13.305	1696	1705	1717	rBV	910753	1361735	44.90%	11.255%
10	14.680	1932	1939	1948	rBV2	238847	387772	12.79%	3.205%
11	16.166	2185	2192	2205	rBV	810896	1188965	39.21%	9.827%
12	17.423	2400	2406	2415	rBV2	405136	596208	19.66%	4.928%
13	20.032	2843	2850	2858	rBV	2269653	3032674	100.00%	25.065%
14	21.688	3125	3132	3141	rBV	487750	796018	26.25%	6.579%
15	24.913	3669	3681	3694	rBV2	251991	702393	23.16%	5.805%

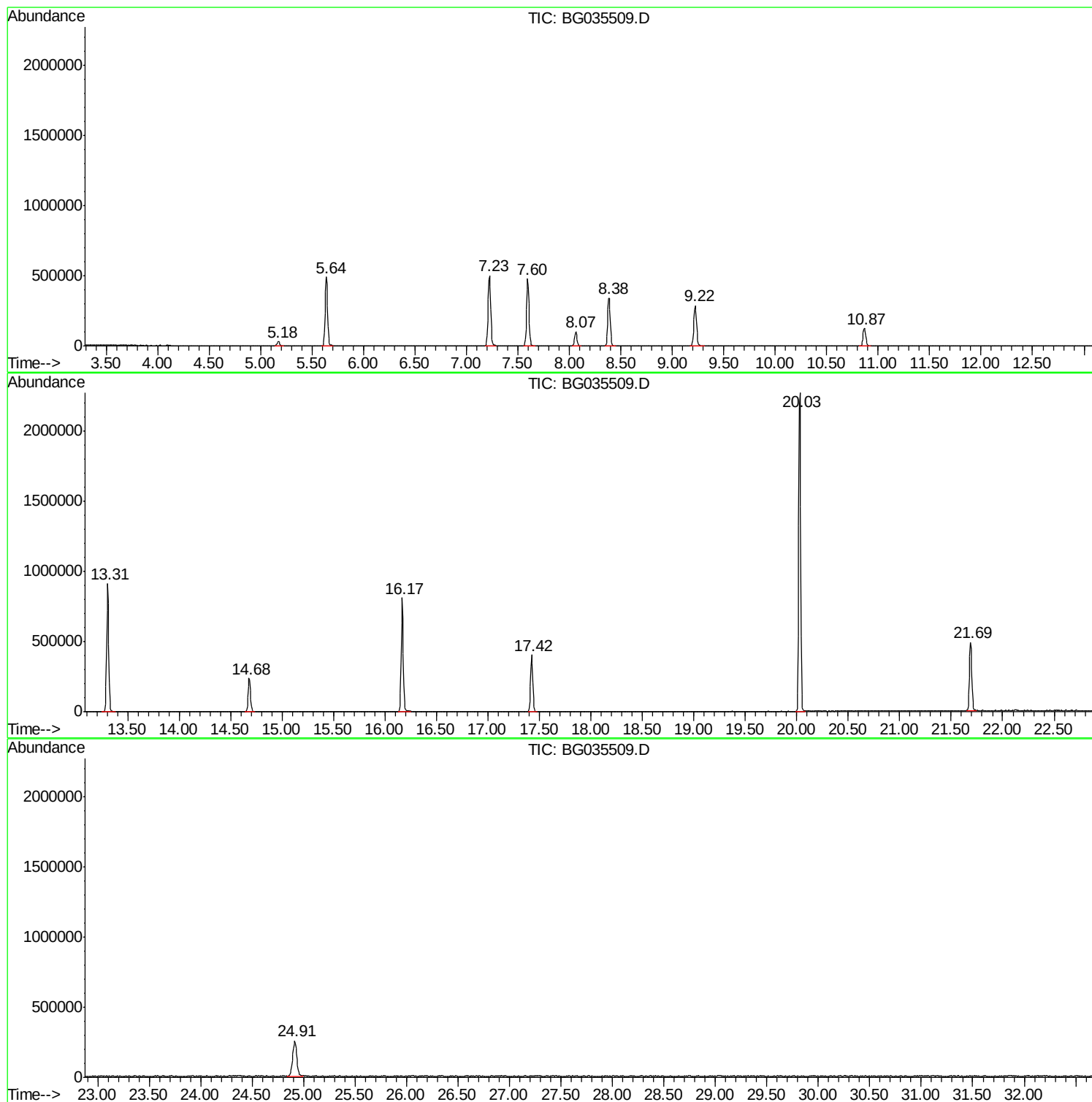
Sum of corrected areas: 12099003

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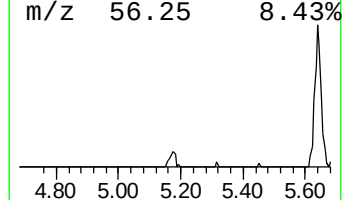
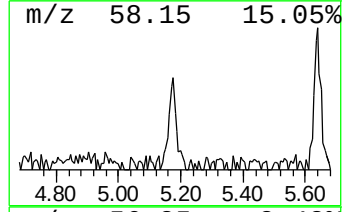
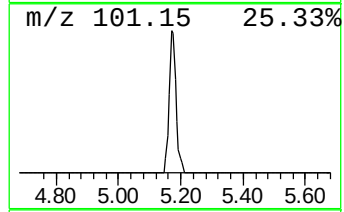
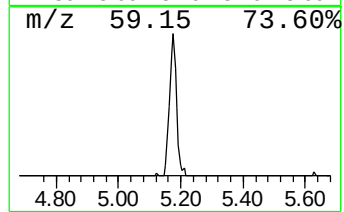
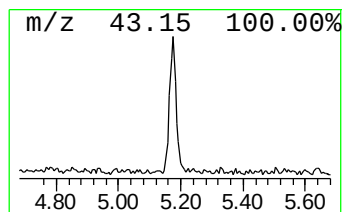
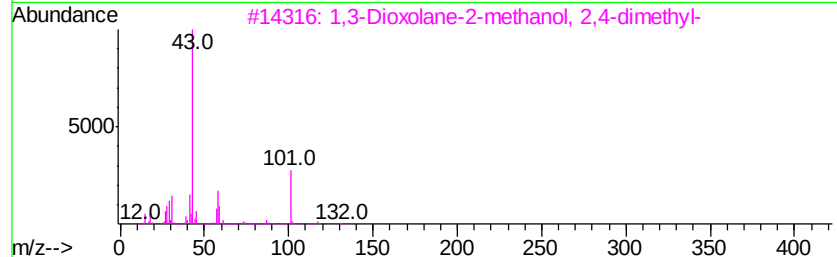
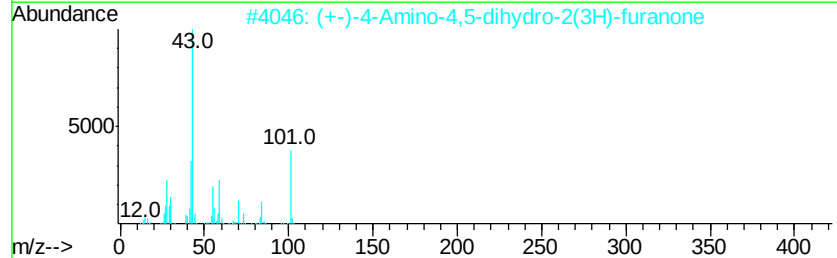
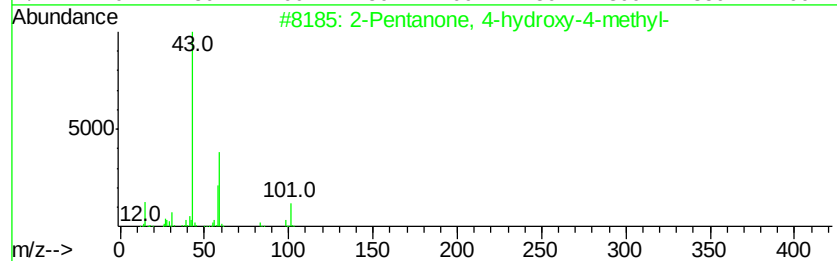
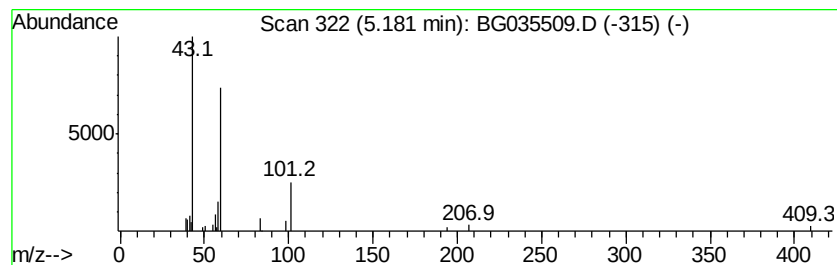
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.18	5.70 ng	49058	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	50
2			(+)-4-Amino-4,5-dihydro-2(3H)-furanone	101	C4H7NO2	016504-58-8	28
3			1,3-Dioxolane-2-methanol, 2,4-dimethyl-	132	C6H12O3	053951-43-2	23
4			2-Propanone, 1-cyclopropyl-	98	C6H10O	004160-75-2	9
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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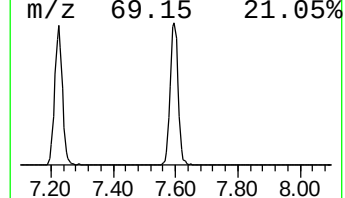
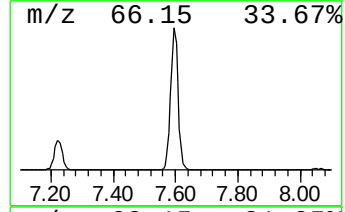
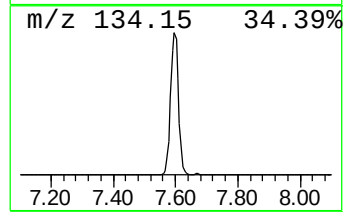
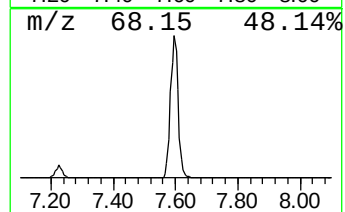
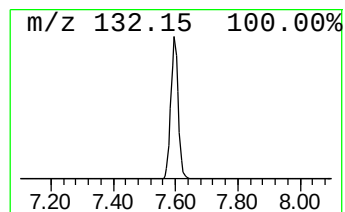
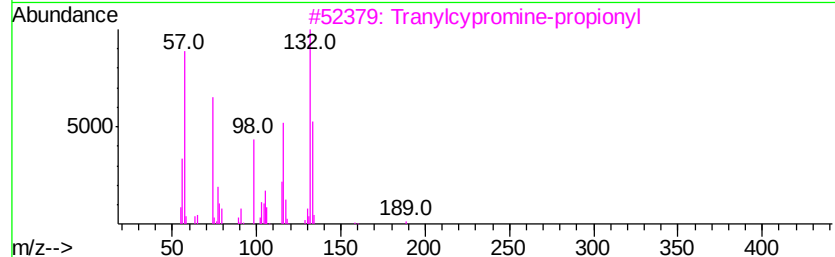
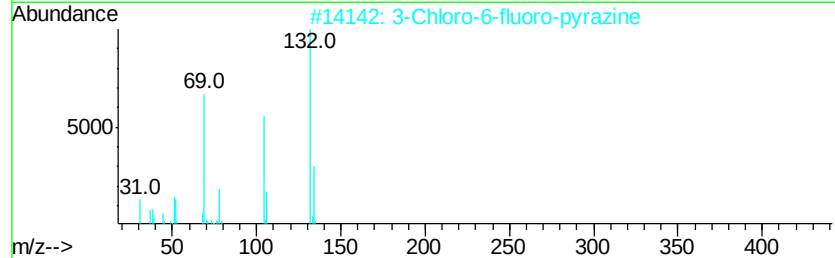
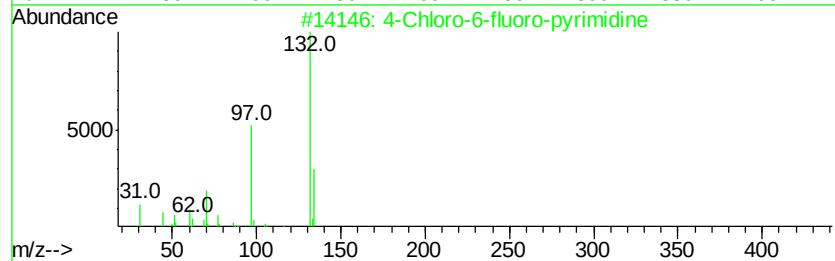
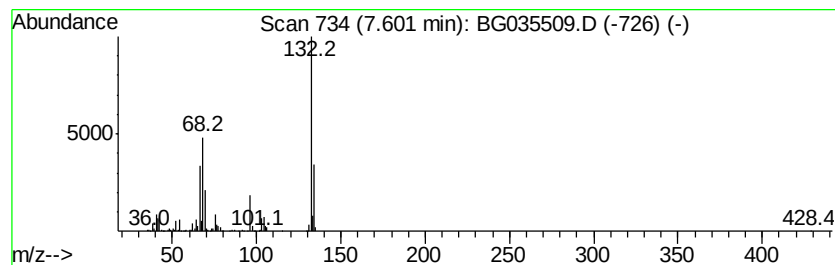
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Peak Number 2 unknown7.60 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.60	94.48 ng	813430	1,4-Dichlorobenzene-d4	8.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		Tranvlcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4		N-(-3,4-Methylenedioxyphenylisop...	267	C17H17NO2	1000378-96-6	10
5		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.18	5.7	ng	49058	1	8.06	172196	20.0
unknown7.60	7.60	94.5	ng	813430	1	8.06	172196	20.0