

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG071818\
 Data File : BG035733.D
 Acq On : 19 Jul 2018 1:24
 Operator : JU/SJ
 Sample : PB111232BL
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB111232BL

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-BG071218.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	5.150	309	317	324	rBV	117050	185851	4.15%	0.964%
2	5.620	390	397	411	rBV	901688	1501312	33.53%	7.785%
3	7.206	659	667	683	rBV	899574	1610807	35.98%	8.353%
4	7.577	723	730	744	rBV	877277	1537561	34.34%	7.973%
5	8.041	800	809	817	rBV	174557	306731	6.85%	1.591%
6	8.358	852	863	876	rBV	674751	1177576	26.30%	6.107%
7	9.198	998	1006	1019	rBV	509027	949115	21.20%	4.922%
8	10.843	1278	1286	1294	rBV	193150	374095	8.36%	1.940%
9	13.286	1694	1702	1714	rBV	1415639	2298523	51.34%	11.920%
10	14.661	1928	1936	1946	rBV	323028	569643	12.72%	2.954%
11	16.147	2182	2189	2202	rBV	762061	1409059	31.47%	7.307%
12	17.404	2396	2403	2414	rBV	527419	833932	18.63%	4.325%
13	20.013	2840	2847	2858	rBV	3310189	4476891	100.00%	23.216%
14	21.669	3122	3129	3139	rBV	634273	1077602	24.07%	5.588%
15	24.883	3664	3676	3690	rBV2	331316	974863	21.78%	5.055%

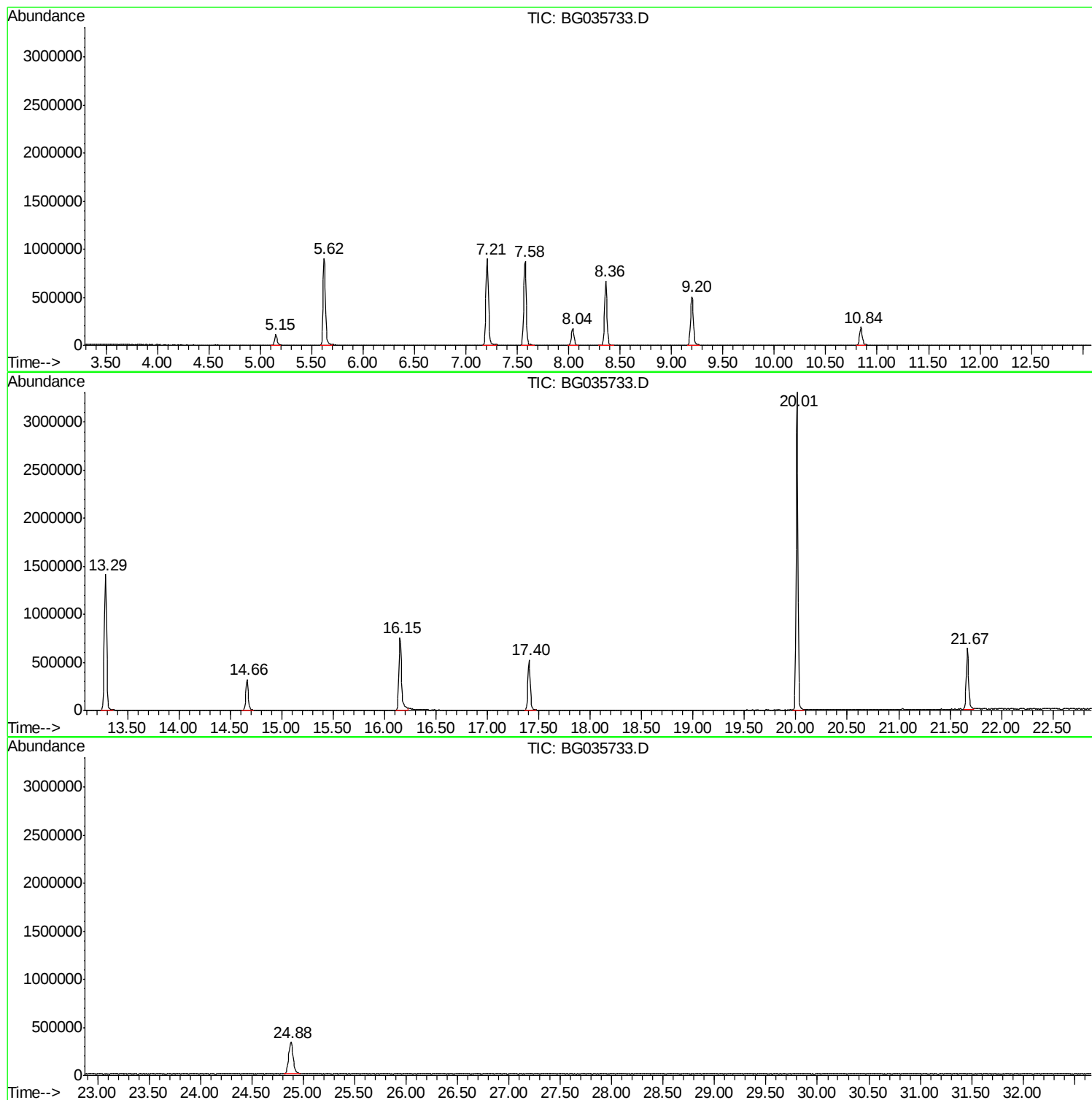
Sum of corrected areas: 19283561

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



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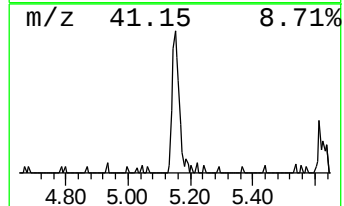
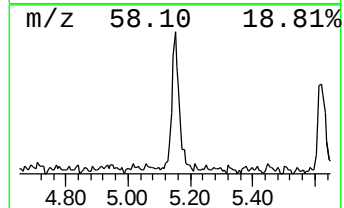
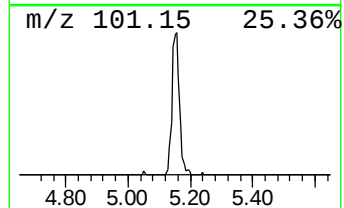
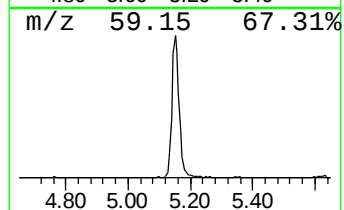
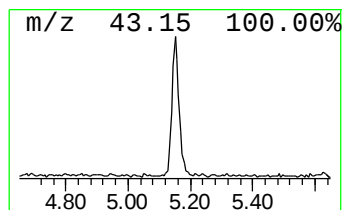
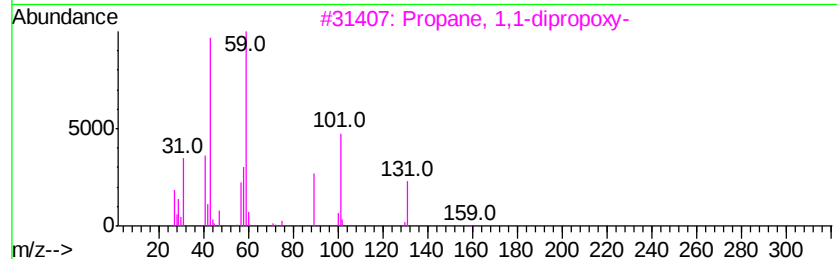
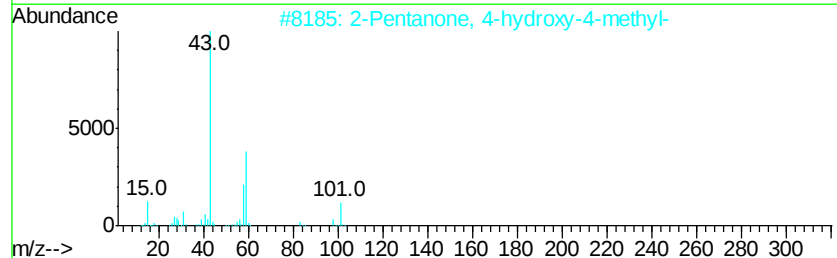
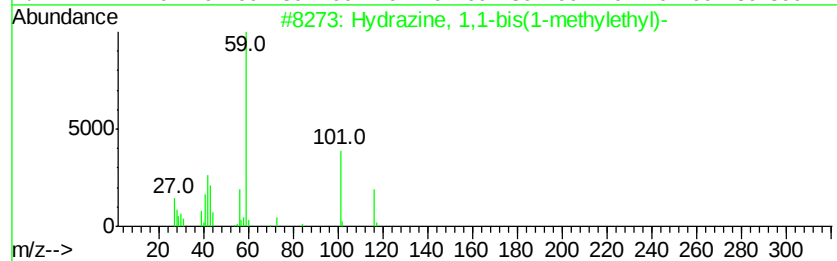
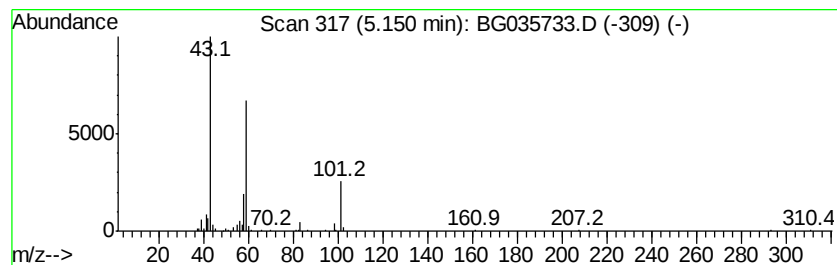
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Peak Number 1 unknown5.15 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.15	12.12 ng	185851	1,4-Dichlorobenzene-d4	8.04

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	47
2			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	47
3			Propane, 1,1-dipropoxy-	160	C9H20O2	004744-11-0	38
4			Dimethadione	129	C5H7N03	000695-53-4	33
5			2,3-Butanedione, monooxime	101	C4H7N02	000057-71-6	25



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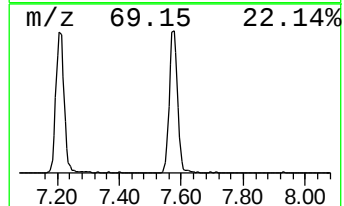
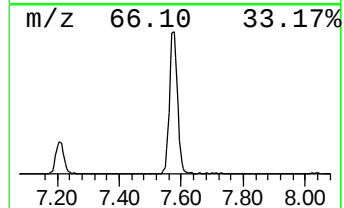
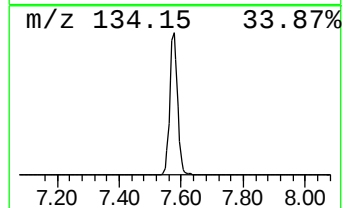
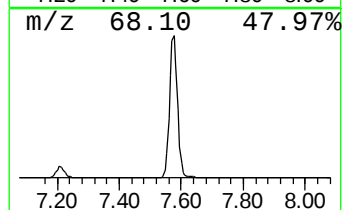
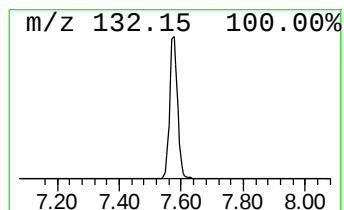
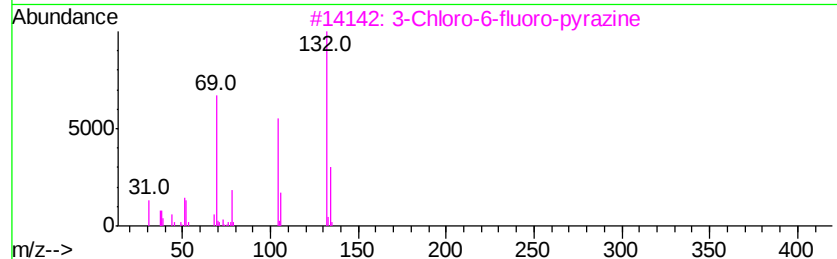
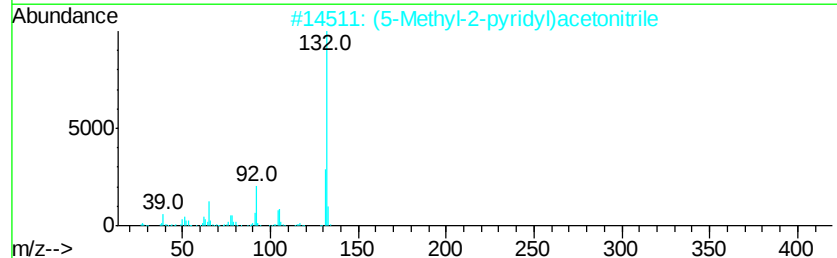
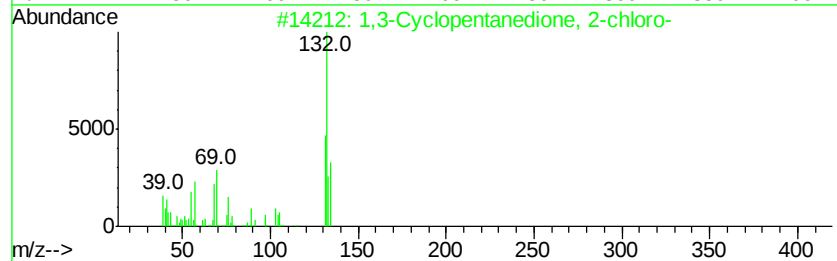
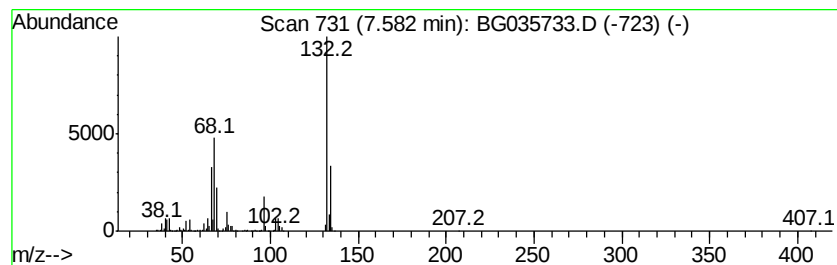
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Peak Number 2 unknown7.58 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	100.26 ng	1537560	1,4-Dichlorobenzene-d4	8.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	25
2		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	22
3		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
5		5-Aminoindole	132	C8H8N2	005192-03-0	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown5.15	5.15	12.1	ng	185851	1	8.04	306731	20.0
unknown7.58	7.58	100.3	ng	1537560	1	8.04	306731	20.0