

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG072618\
 Data File : BG035927.D
 Acq On : 27 Jul 2018 3:01
 Operator : JU/SJ
 Sample : PB111478BL
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB111478BL

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	5.135	309	314	324	rVB	56263	86253	3.06%	0.739%
2	5.605	387	394	408	rBV	429649	726209	25.78%	6.225%
3	7.191	657	664	672	rBV	434145	777938	27.62%	6.668%
4	7.561	719	727	739	rBV	443570	775449	27.53%	6.647%
5	8.025	796	806	815	rBV	105663	184153	6.54%	1.578%
6	8.349	853	861	870	rBV	333206	583945	20.73%	5.005%
7	9.189	995	1004	1018	rBV	232551	476651	16.92%	4.085%
8	10.833	1275	1284	1301	rBV	110892	243650	8.65%	2.088%
9	13.271	1692	1699	1711	rBV	781985	1253621	44.51%	10.745%
10	14.652	1927	1934	1949	rBV3	223297	404827	14.37%	3.470%
11	16.138	2180	2187	2204	rBV	660865	1124486	39.93%	9.638%
12	17.395	2393	2401	2418	rBV	395980	673644	23.92%	5.774%
13	19.998	2838	2844	2854	rBV	2132500	2816427	100.00%	24.140%
14	21.654	3119	3126	3139	rBV	450705	805282	28.59%	6.902%
15	24.856	3660	3671	3684	rBV3	237837	734369	26.07%	6.294%

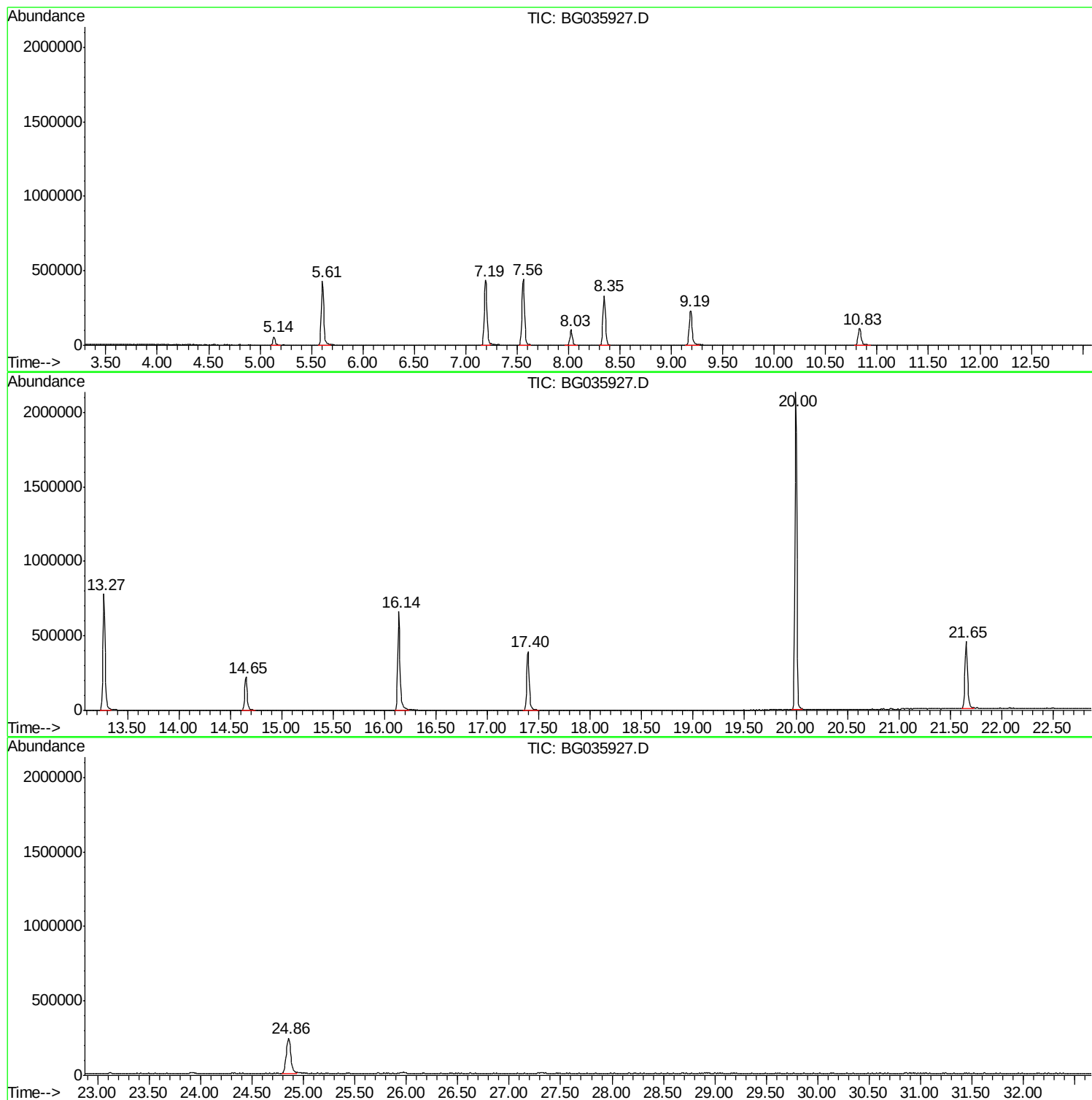
Sum of corrected areas: 11666904

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG071218.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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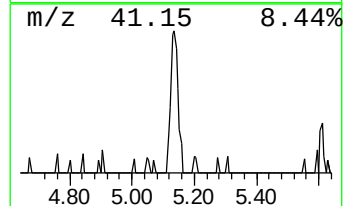
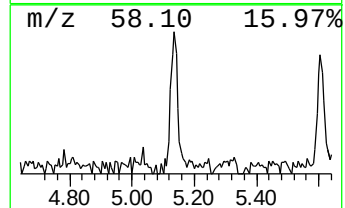
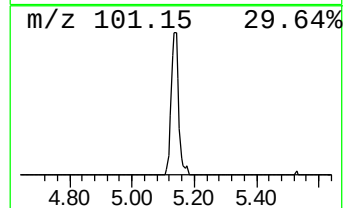
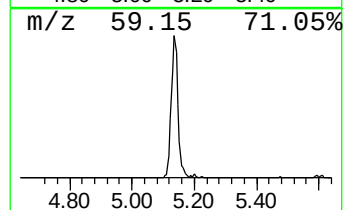
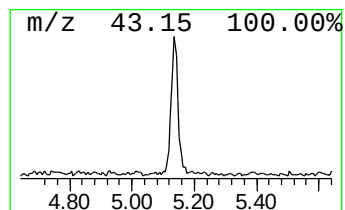
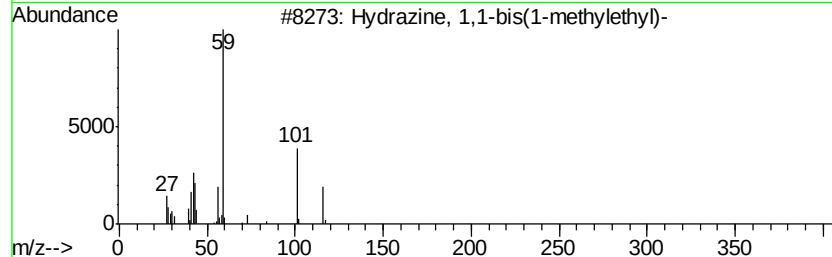
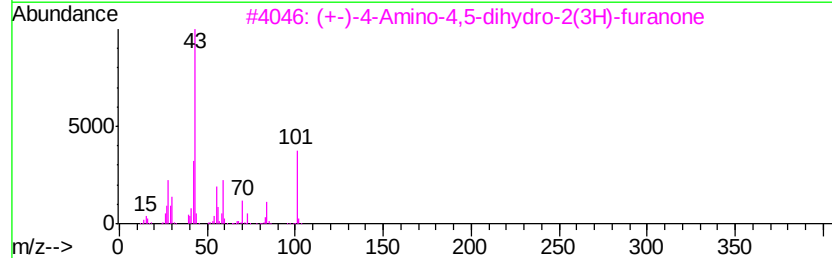
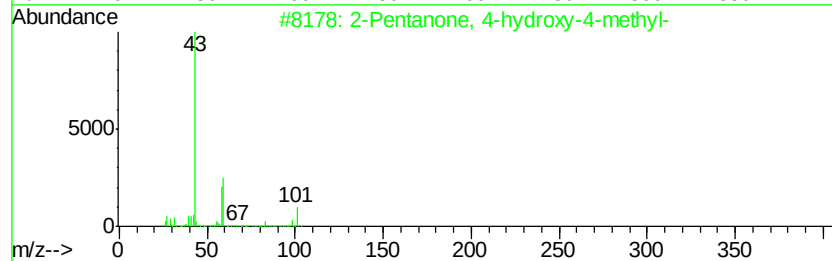
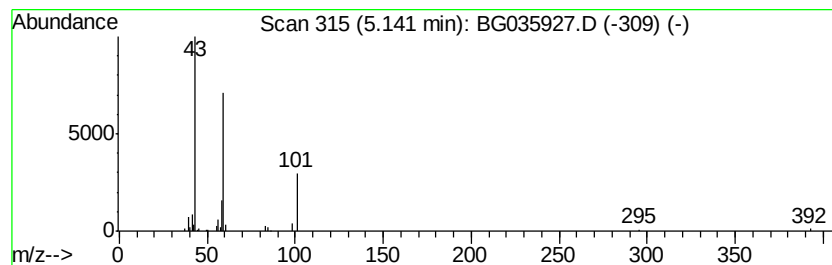
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.14	9.37 ng	86253	1,4-Dichlorobenzene-d4	8.03

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2			(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	37
3			Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	33
4			1,2-Butanediol	90	C4H10O2	000584-03-2	32
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	32



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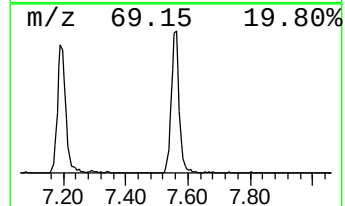
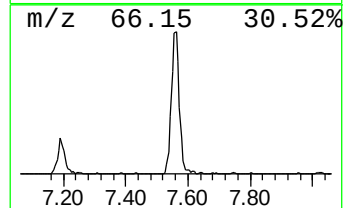
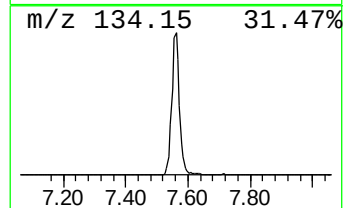
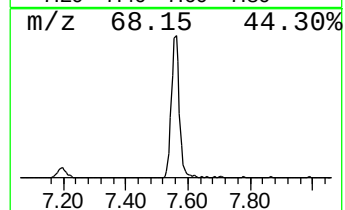
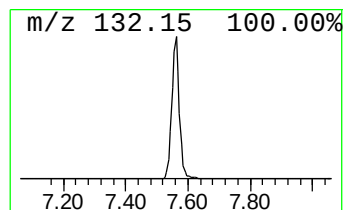
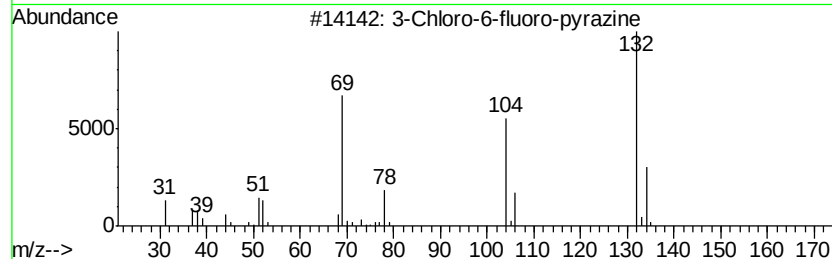
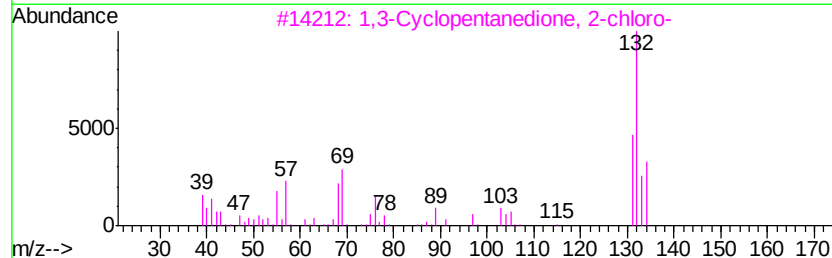
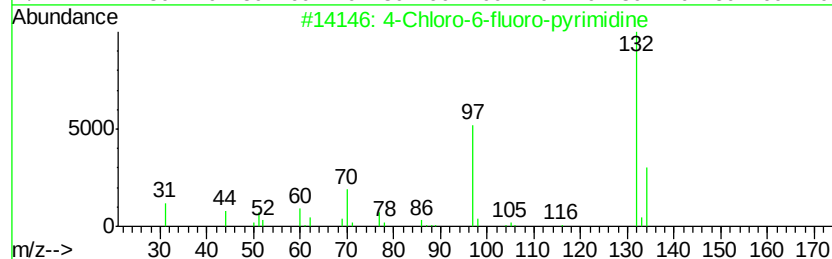
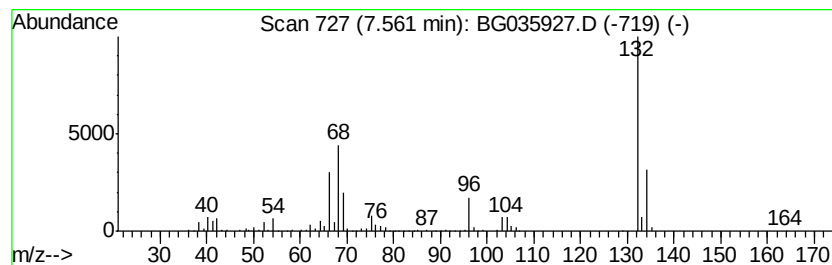
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Peak Number 2 unknown7.56 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.56	84.22 ng	775449	1,4-Dichlorobenzene-d4	8.03

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	17
3		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
5		Tranlylcypromine-propionyl	189	C12H15NO	1000123-86-3	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.14	9.4	ng	86253	1	8.03	184153	20.0
unknown7.56	7.56	84.2	ng	775449	1	8.03	184153	20.0