

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110518\
 Data File : BG037820.D
 Acq On : 6 Nov 2018 2:14
 Operator : JU/SJ
 Sample : PB114415BL
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB114415BL

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-BG101818.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	5.177	315	321	329	rBV	40135	65518	1.73%	0.327%
2	5.635	393	399	415	rBV	920128	1545866	40.74%	7.726%
3	7.239	664	672	683	rBV	983517	1760231	46.39%	8.798%
4	7.621	729	737	749	rBV	929168	1664934	43.88%	8.321%
5	8.097	810	818	825	rBV	194084	342236	9.02%	1.711%
6	8.420	865	873	890	rVB	691917	1263807	33.31%	6.317%
7	9.272	1008	1018	1028	rBV	589919	1103524	29.08%	5.515%
8	10.917	1289	1298	1307	rBV	287677	540415	14.24%	2.701%
9	13.350	1703	1712	1724	rBV	1744814	2826672	74.50%	14.128%
10	14.731	1939	1947	1955	rBV2	477510	811574	21.39%	4.056%
11	16.217	2193	2200	2215	rBV	1012556	1686400	44.45%	8.429%
12	17.474	2406	2414	2424	rBV2	581365	941354	24.81%	4.705%
13	20.077	2850	2857	2875	rBV	2663496	3794254	100.00%	18.964%
14	21.758	3136	3143	3165	rBV	498825	969140	25.54%	4.844%
15	25.095	3699	3711	3726	rBV2	193319	691871	18.23%	3.458%

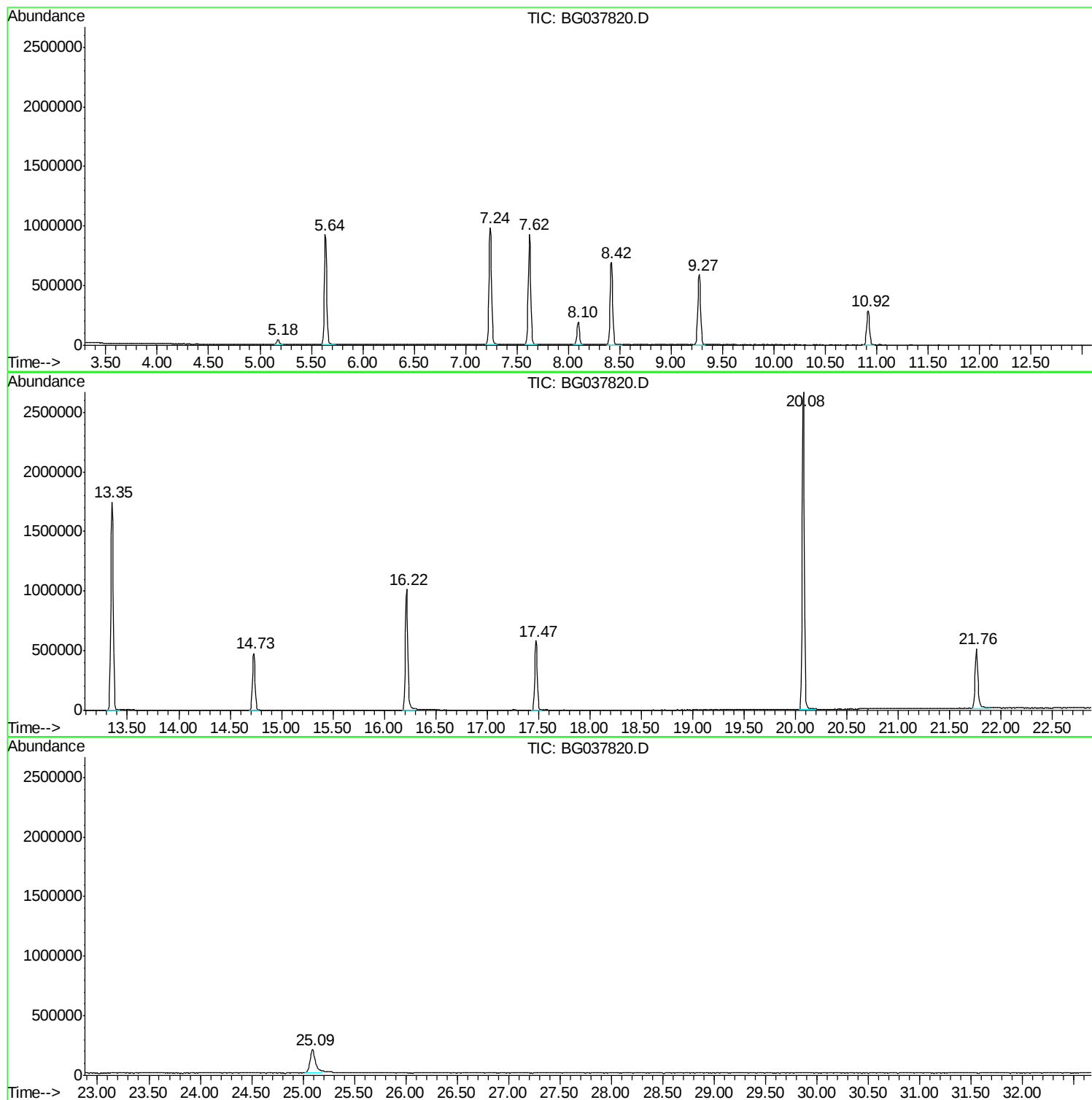
Sum of corrected areas: 20007796

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG101818.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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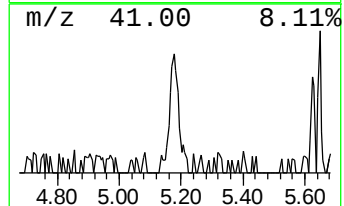
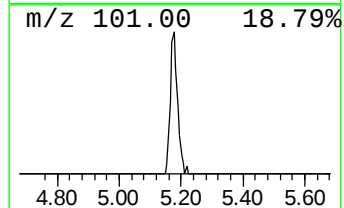
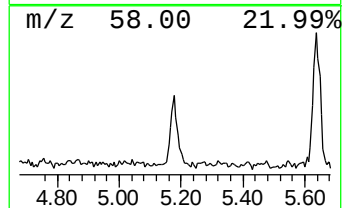
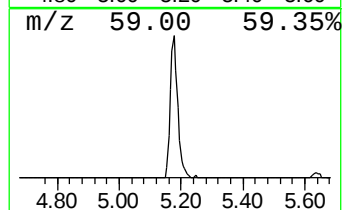
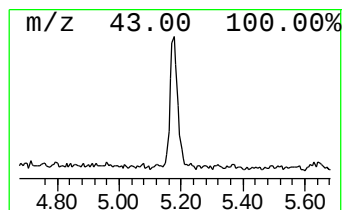
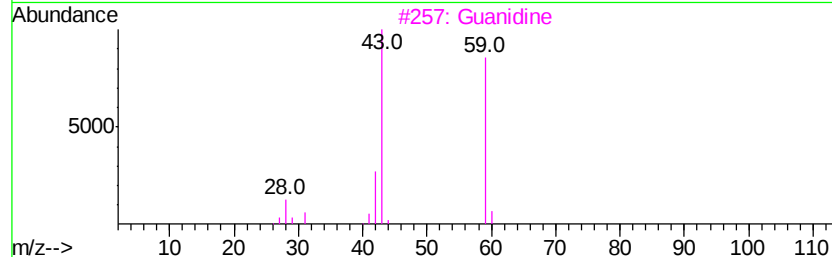
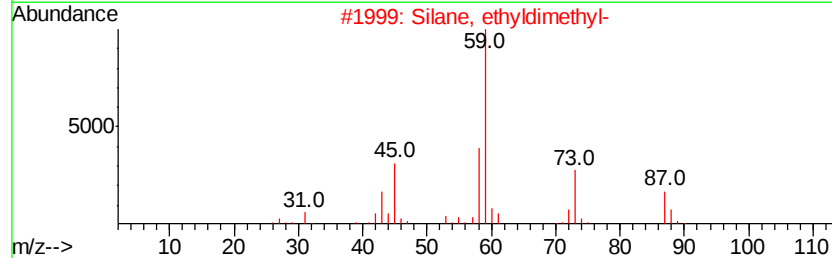
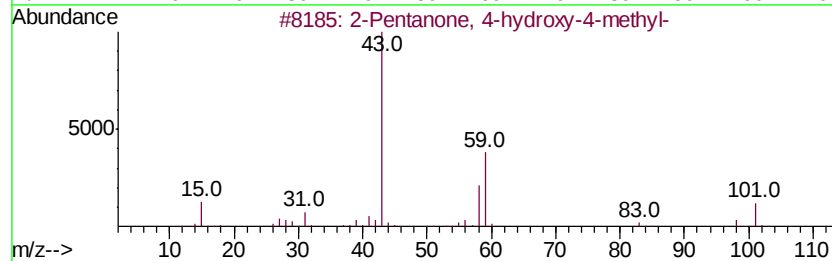
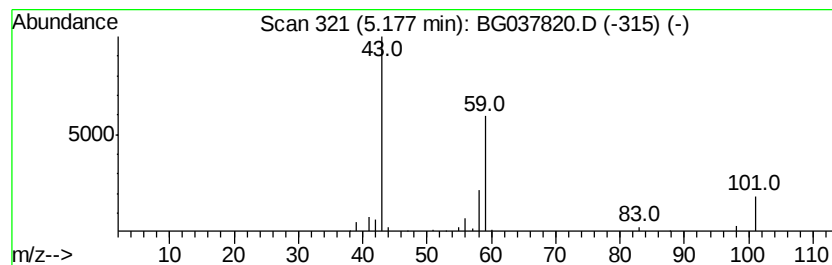
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.18	3.83 ng	65518	1,4-Dichlorobenzene-d4	8.10

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			Silane, ethyldimethyl-	88	C4H12Si	000758-21-4	25
3			Guanidine	59	CH5N3	000113-00-8	9
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5			Acetone	58	C3H6O	000067-64-1	9



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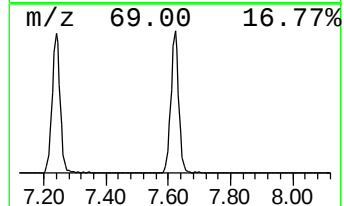
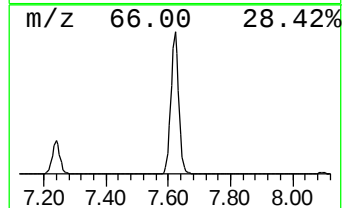
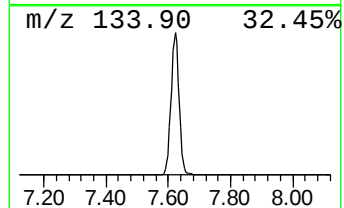
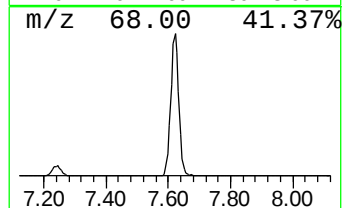
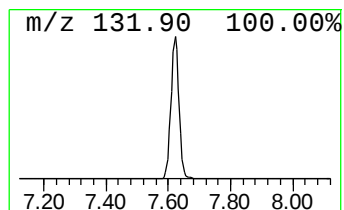
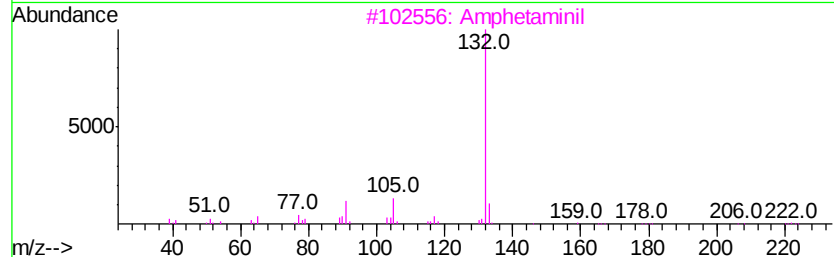
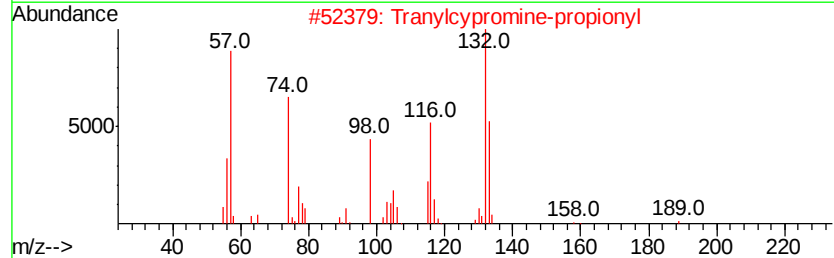
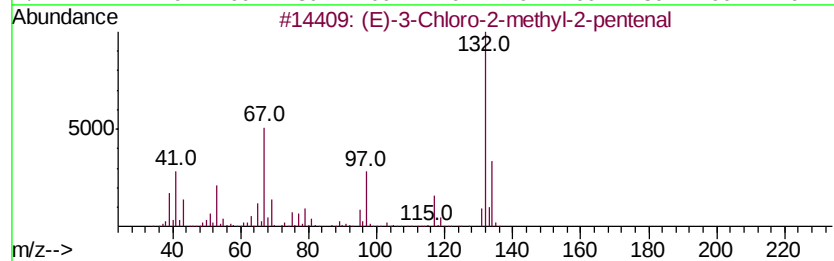
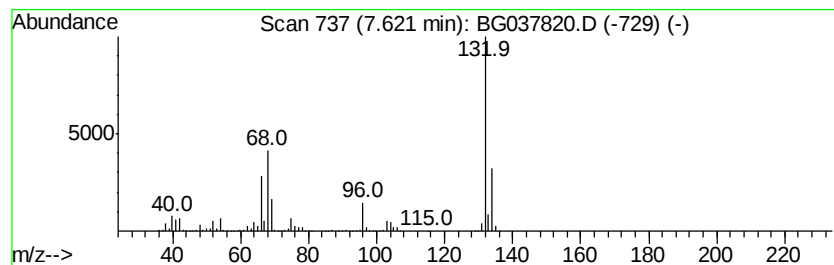
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Peak Number 2 unknown7.62 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.62	97.30 ng	1664930	1,4-Dichlorobenzene-d4	8.10

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
2			Tranlylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
3			Amphetaminil	250	C17H18N2	017590-01-1	12
4			N-(-phenylisopropyl)benzalimine	223	C16H17N	1000379-01-3	10
5			N-(-3,4-Methylenedioxyphenylisop...	267	C17H17NO2	1000378-96-6	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.18	3.8	ng	65518	1	8.10	342236	20.0
unknown7.62	7.62	97.3	ng	1664930	1	8.10	342236	20.0