

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG100118\
 Data File : BG037090.D
 Acq On : 2 Oct 2018 00:08
 Operator : JU/SJ
 Sample : PB113393BL
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB113393BL

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 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-BG092018.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	5.142	311	315	322	rBV	27134	43737	1.22%	0.258%
2	5.612	389	395	416	rBV	732607	1334676	37.32%	7.869%
3	7.199	659	665	687	rBV	700450	1428757	39.95%	8.424%
4	7.569	721	728	749	rBV	689698	1388688	38.83%	8.188%
5	8.039	802	808	825	rBV	127131	268339	7.50%	1.582%
6	8.362	855	863	876	rBV	565518	1080526	30.21%	6.371%
7	9.202	999	1006	1025	rBV	409276	897875	25.10%	5.294%
8	10.841	1279	1285	1301	rBV	168058	372383	10.41%	2.196%
9	13.286	1694	1701	1726	rBV	1227887	2125788	59.43%	12.534%
10	14.666	1929	1936	1952	rBV2	292142	544425	15.22%	3.210%
11	16.153	2182	2189	2208	rBV	746777	1433394	40.08%	8.451%
12	17.410	2396	2403	2417	rBV2	421114	732365	20.48%	4.318%
13	20.019	2841	2847	2858	rBV	2483676	3576730	100.00%	21.088%
14	21.676	3122	3129	3145	rBV	466960	882688	24.68%	5.204%
15	24.884	3664	3675	3692	rBV2	237213	805952	22.53%	4.752%
16	27.152	4057	4061	4075	rVB9	12021	44497	1.24%	0.262%

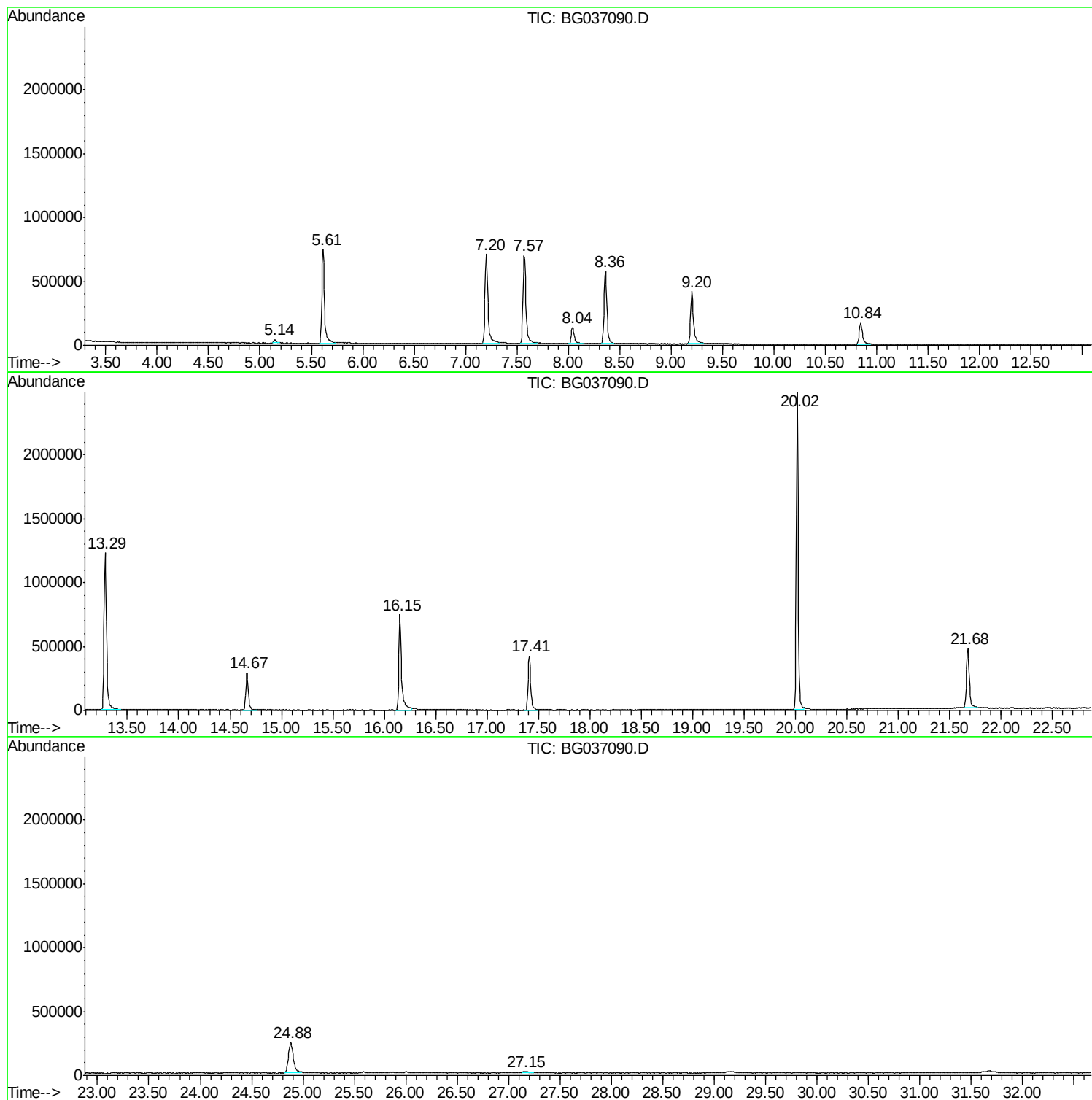
Sum of corrected areas: 16960820

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG092018.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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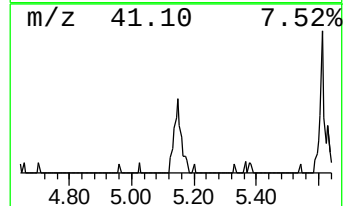
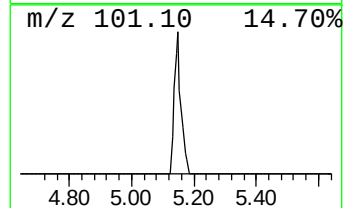
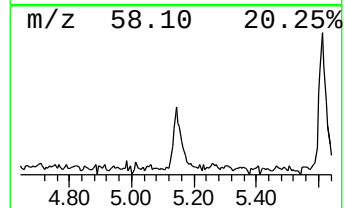
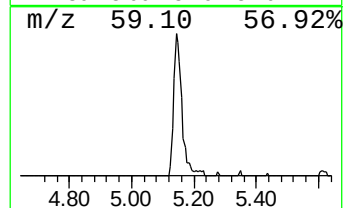
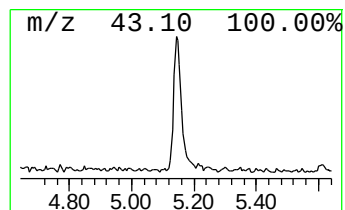
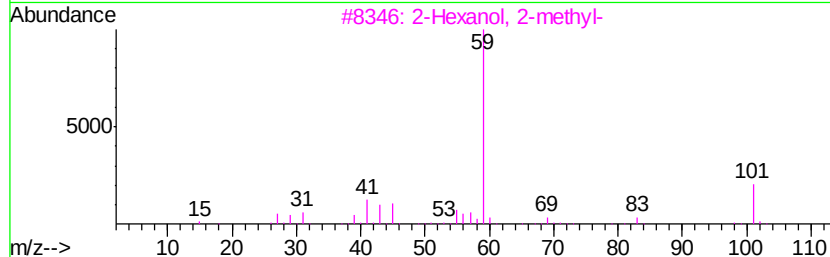
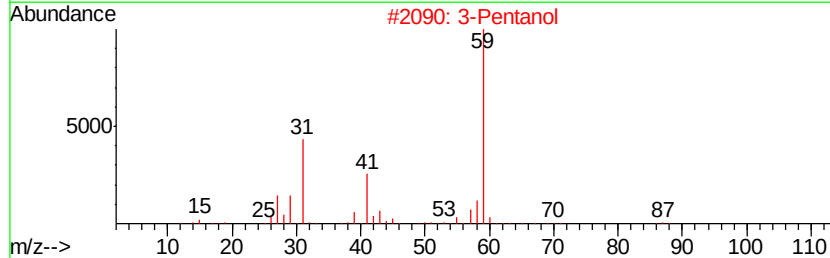
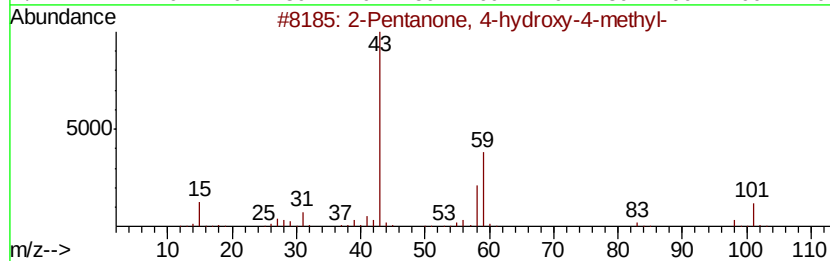
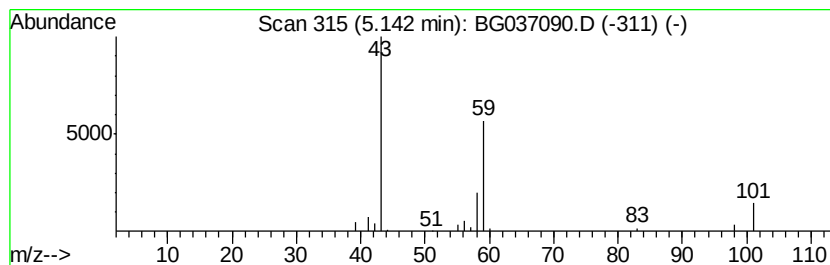
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.14	3.26 ng	43737	1,4-Dichlorobenzene-d4	8.04

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2			3-Pentanol	88	C5H12O	000584-02-1	9
3			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	9
4			Diacetamide	101	C4H7NO2	000625-77-4	7
5			Ethanone, 1-(2-methylcyclopropyl)-	98	C6H10O	000930-56-3	5



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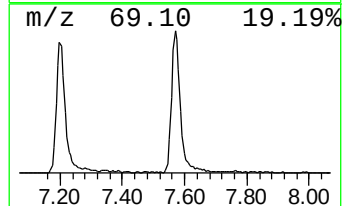
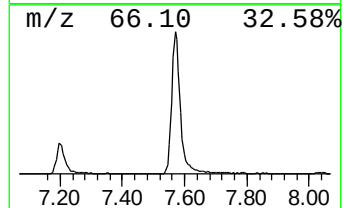
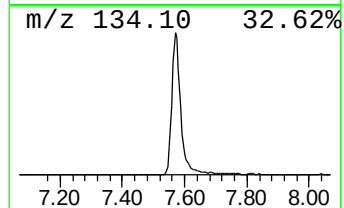
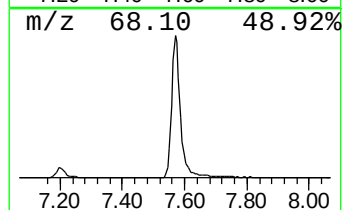
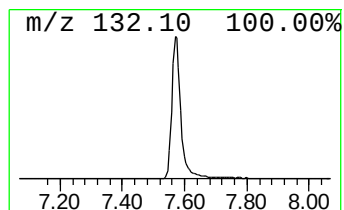
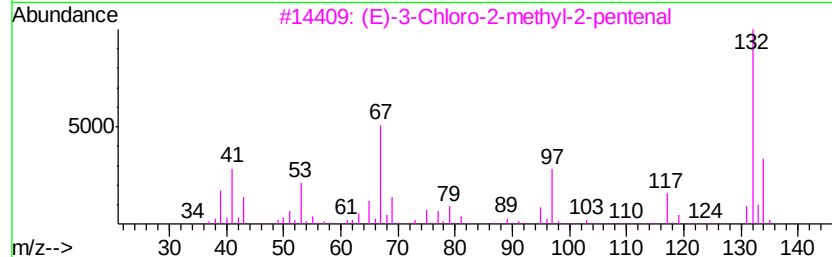
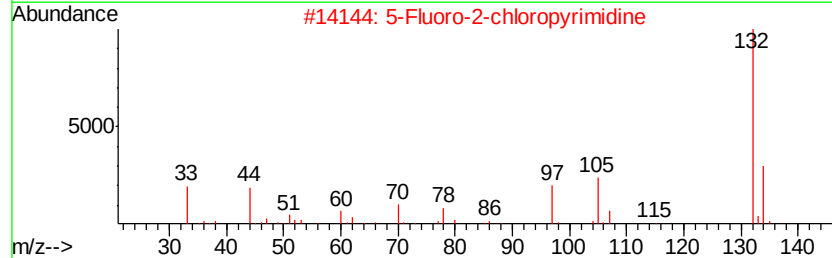
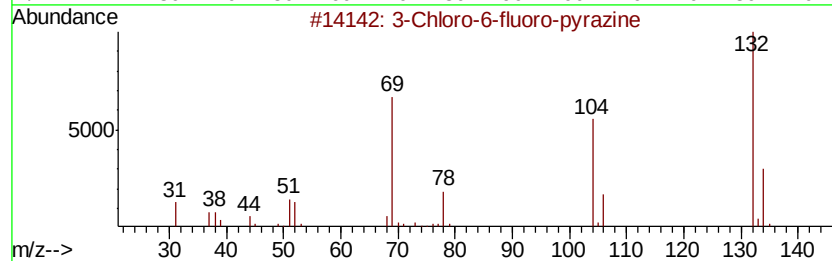
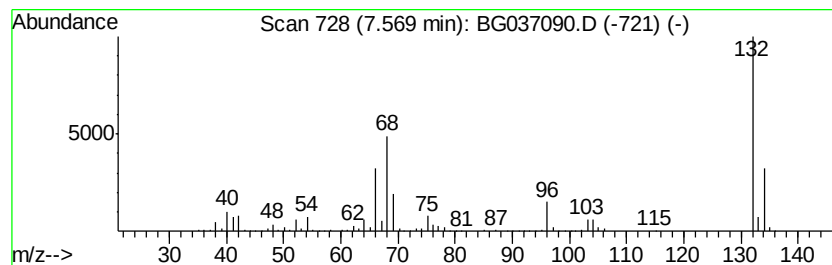
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Peak Number 2 unknown7.57 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.57	103.50 ng	1388690	1,4-Dichlorobenzene-d4	8.04

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		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.14	3.3	ng	43737	1	8.04	268339	20.0
unknown7.57	7.57	103.5	ng	1388690	1	8.04	268339	20.0