

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG111618\
 Data File : BG038037.D
 Acq On : 16 Nov 2018 17:21
 Operator : JU/SJ
 Sample : J5965-07
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MW-7

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-BG111318.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	5.165	314	319	326	rBV	36417	59753	1.99%	0.377%
2	5.629	390	398	408	rBV	764175	1266489	42.08%	7.994%
3	7.233	662	671	680	rBV	786281	1388802	46.14%	8.766%
4	7.615	728	736	749	rVB	734060	1329254	44.16%	8.390%
5	8.085	809	816	824	rVB	149945	261119	8.68%	1.648%
6	8.408	863	871	881	rVB	558920	991214	32.93%	6.256%
7	9.266	1009	1017	1027	rBV	463847	862595	28.66%	5.444%
8	10.905	1288	1296	1305	rVB	204530	384379	12.77%	2.426%
9	13.343	1703	1711	1718	rBV	1260830	2036956	67.68%	12.857%
10	14.718	1938	1945	1953	rBV3	327356	546948	18.17%	3.452%
11	16.210	2191	2199	2215	rBV3	933983	1496234	49.71%	9.444%
12	17.468	2407	2413	2421	rBV2	436942	689784	22.92%	4.354%
13	20.071	2849	2856	2867	rBV	2151884	3009801	100.00%	18.997%
14	21.751	3135	3142	3153	rVB2	449644	770173	25.59%	4.861%
15	25.071	3692	3707	3719	rBV2	242770	750199	24.93%	4.735%

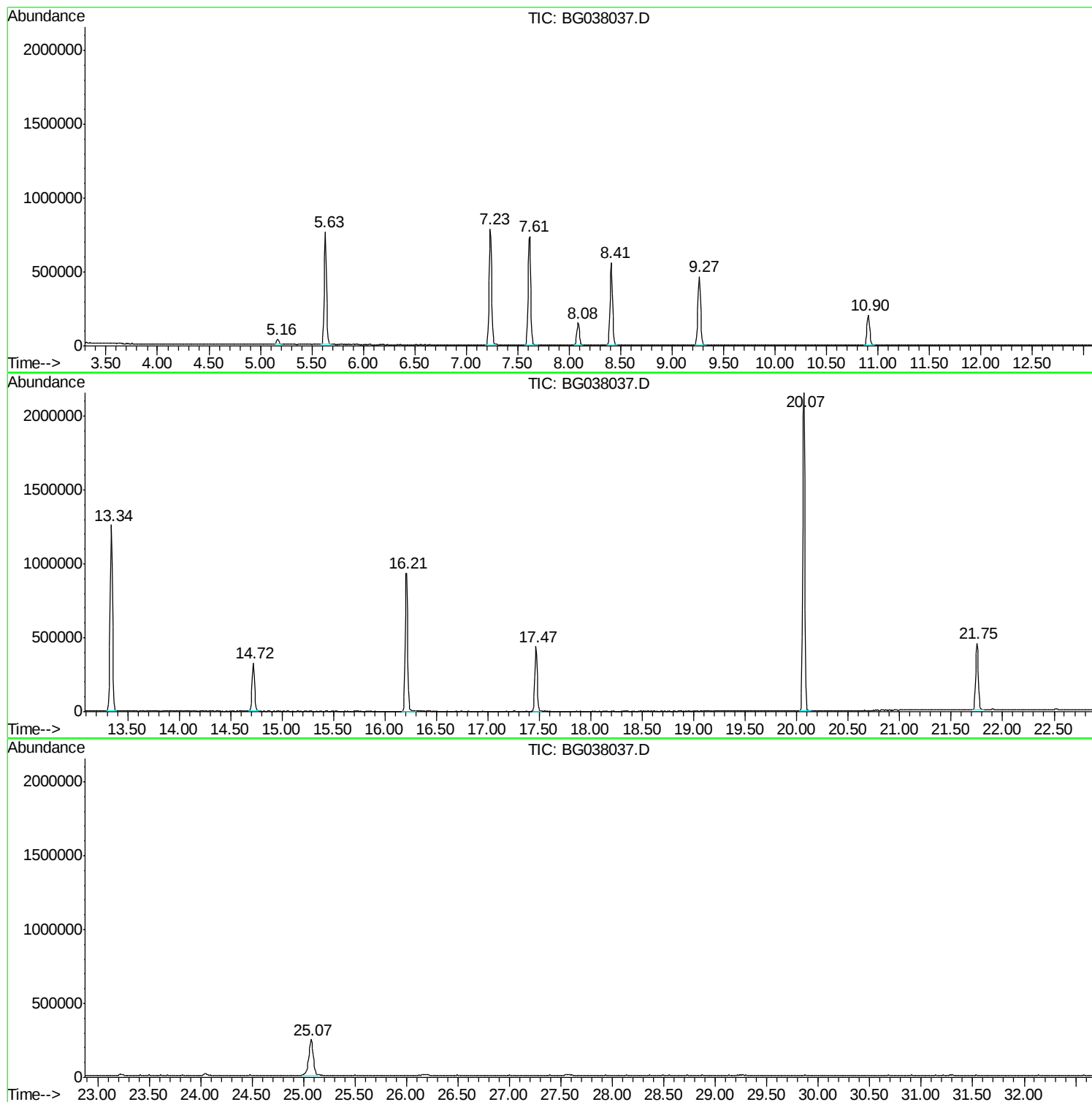
Sum of corrected areas: 15843700

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG111318.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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ClientSampled :
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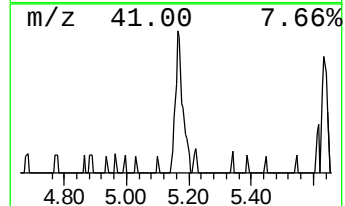
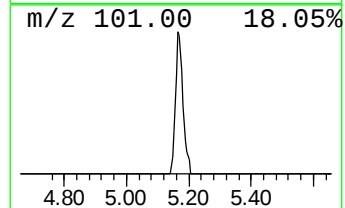
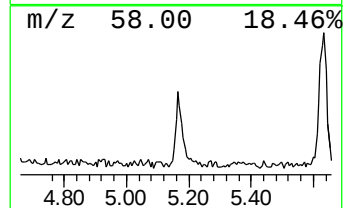
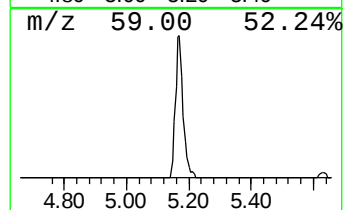
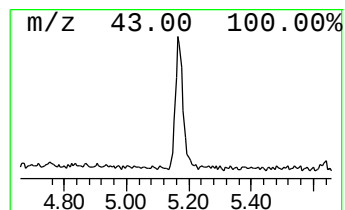
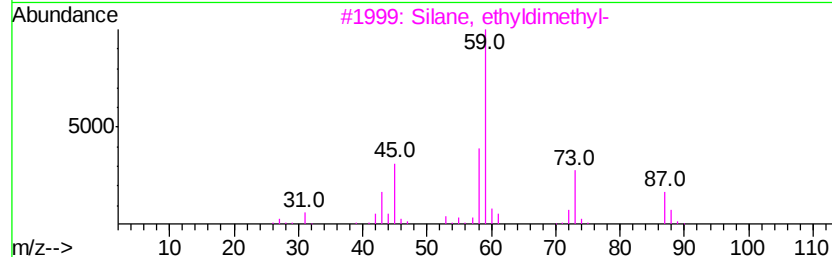
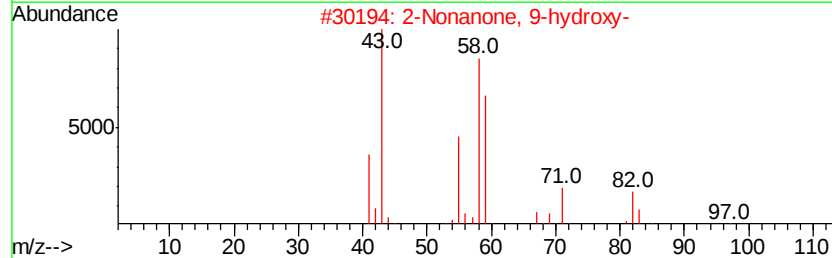
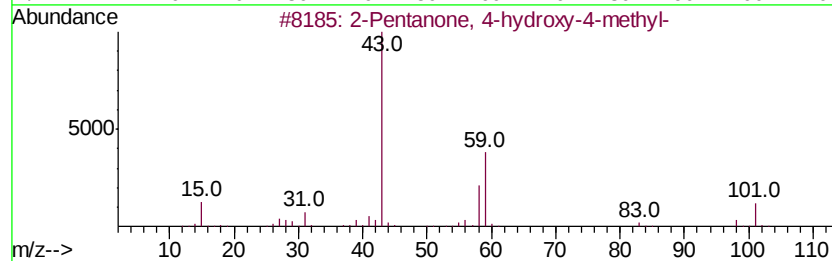
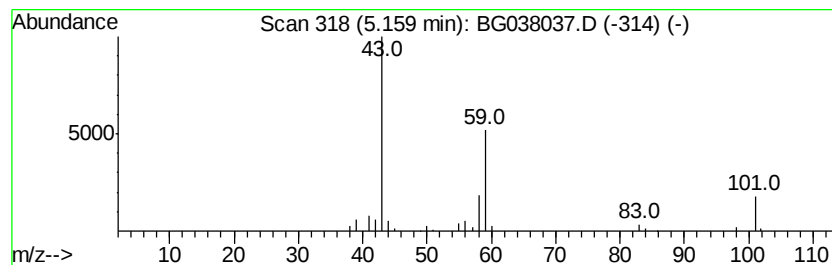
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.16	4.58 ng	59753	1,4-Dichlorobenzene-d4	8.08

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	78
2			2-Nonanone, 9-hydroxy-	158	C9H18O2	025368-56-3	28
3			Silane, ethyldimethyl-	88	C4H12Si	000758-21-4	25
4			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	12
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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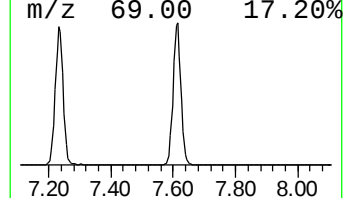
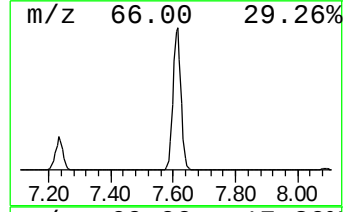
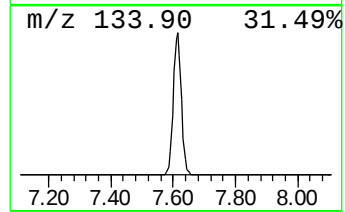
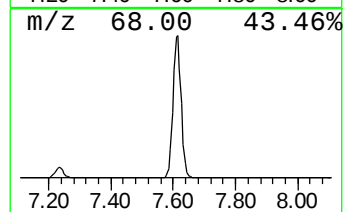
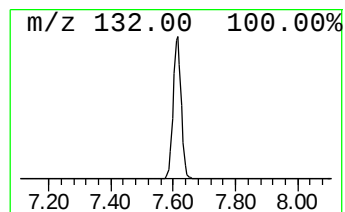
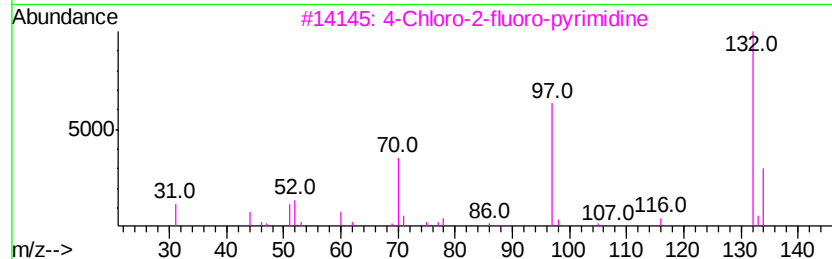
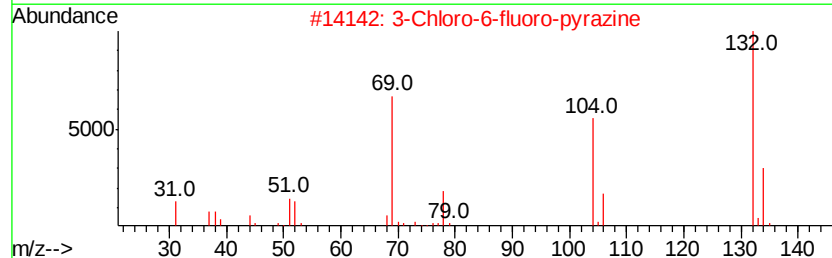
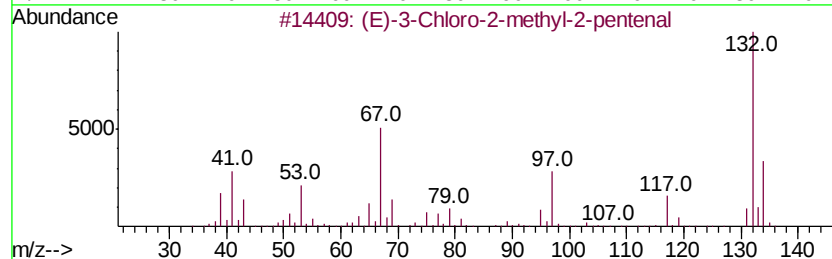
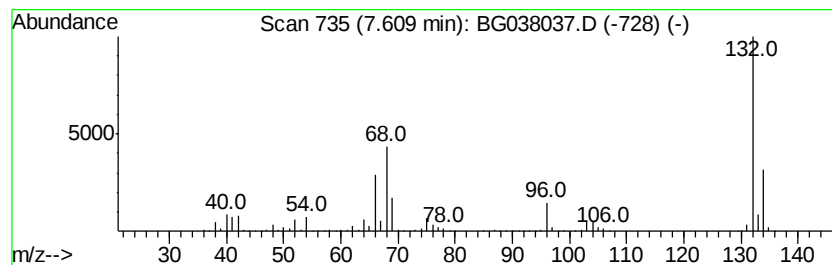
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Peak Number 2 unknown7.61 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.61	101.81 ng	1329250	1,4-Dichlorobenzene-d4	8.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	30
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	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...	5.16	4.6	ng	59753	1	8.08	261119	20.0
unknown7.61	7.61	101.8	ng	1329250	1	8.08	261119	20.0