

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG111218\
 Data File : BG037947.D
 Acq On : 12 Nov 2018 16:47
 Operator : JU/SJ
 Sample : J5895-03
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-2

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	5.173	316	321	328	rBV	46367	77598	2.25%	0.384%
2	5.632	392	399	410	rBV	982531	1620291	46.92%	8.022%
3	7.236	664	672	685	rBV	1045959	1958183	56.70%	9.695%
4	7.618	729	737	748	rBV	994351	1761852	51.02%	8.723%
5	8.088	811	817	826	rVB	170500	303803	8.80%	1.504%
6	8.411	865	872	881	rBV	728156	1315649	38.10%	6.514%
7	9.269	1010	1018	1032	rBV	648571	1203806	34.86%	5.960%
8	10.908	1288	1297	1306	rBV	252310	479726	13.89%	2.375%
9	13.346	1703	1712	1724	rBV	1772948	2885909	83.57%	14.289%
10	14.169	1846	1852	1864	rBV	115595	173655	5.03%	0.860%
11	14.727	1938	1947	1956	rBV2	424148	713804	20.67%	3.534%
12	16.213	2193	2200	2214	rBV	1188590	1926035	55.77%	9.536%
13	17.471	2407	2414	2426	rBV2	511341	798377	23.12%	3.953%
14	18.329	2554	2560	2570	rBV2	34870	66819	1.93%	0.331%
15	20.074	2850	2857	2877	rBV	2426146	3453283	100.00%	17.098%
16	21.754	3136	3143	3155	rBV	416236	784149	22.71%	3.882%
17	23.429	3421	3428	3435	rBV4	18011	39247	1.14%	0.194%
18	25.080	3694	3709	3730	rBV3	162523	635032	18.39%	3.144%

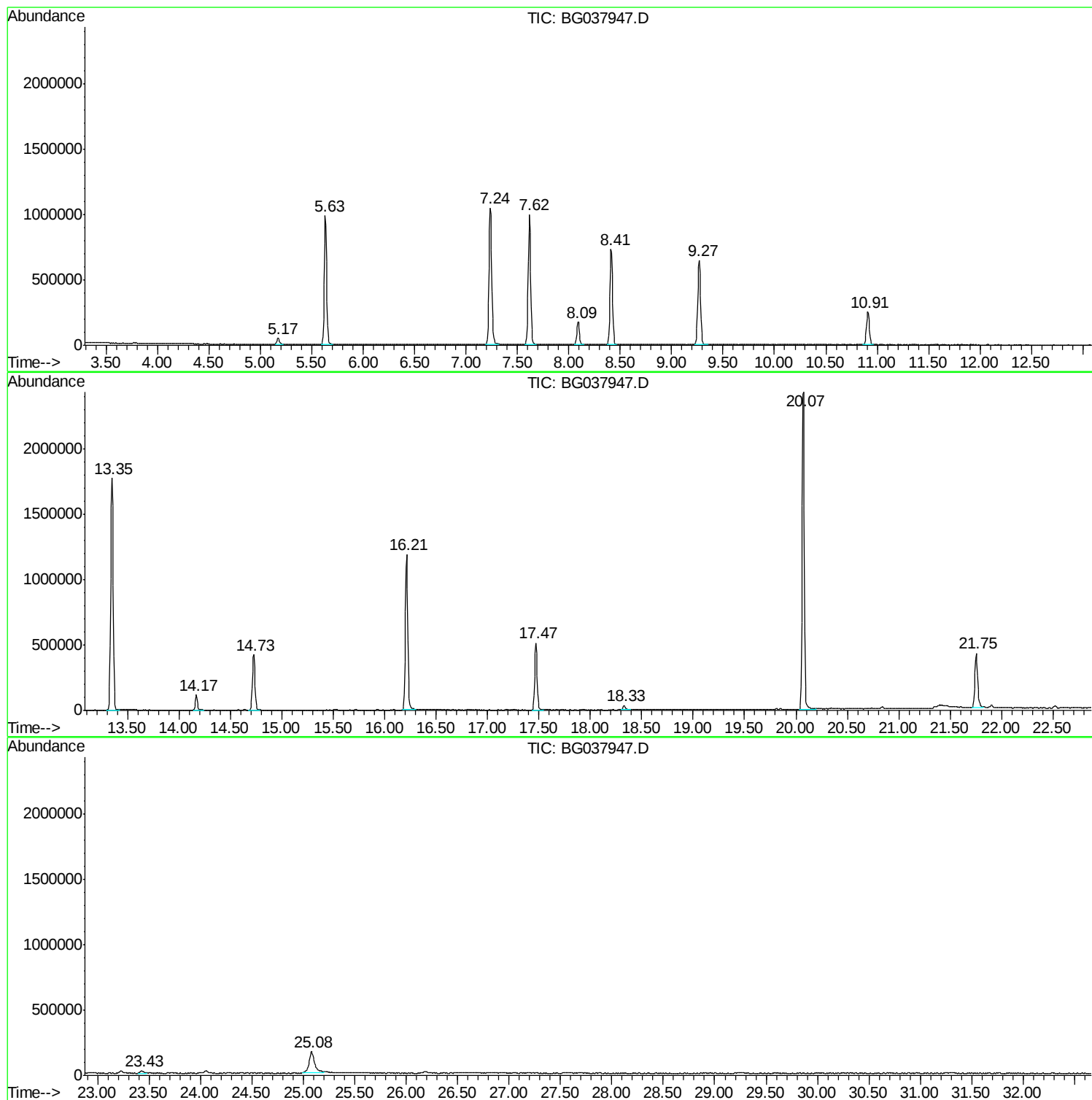
Sum of corrected areas: 20197218

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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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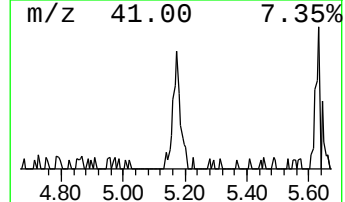
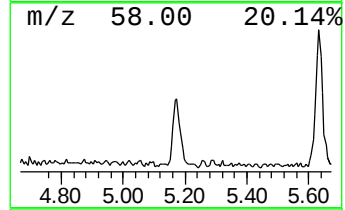
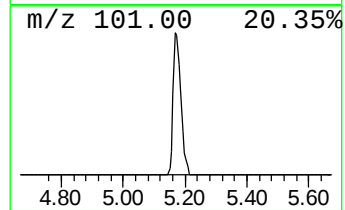
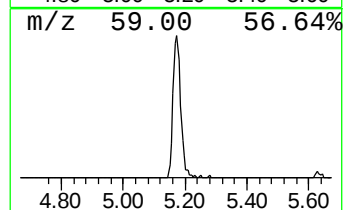
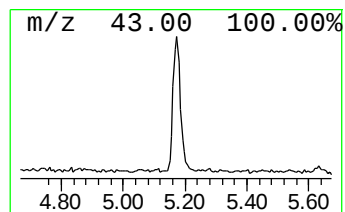
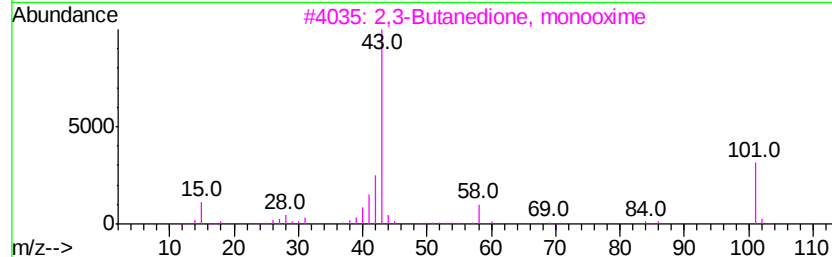
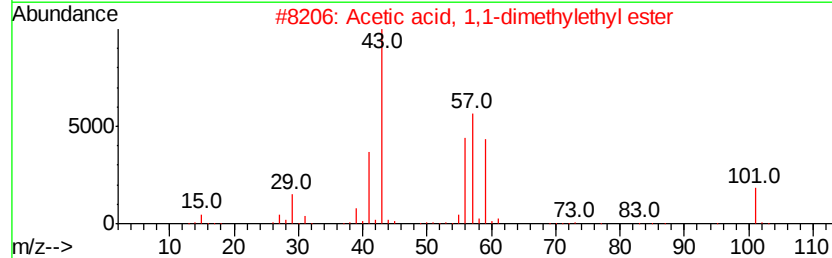
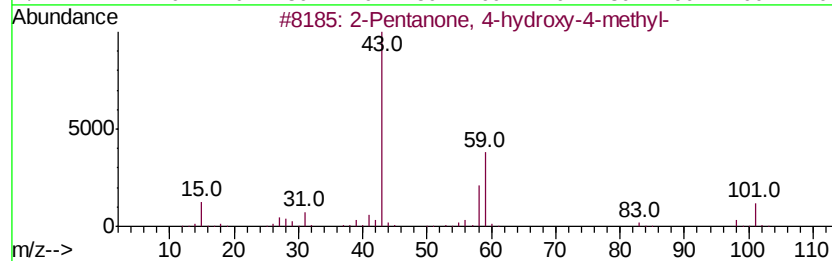
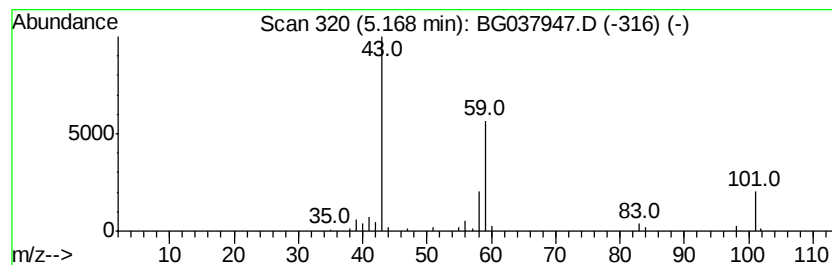
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.17	5.11 ng	77598	1,4-Dichlorobenzene-d4	8.09

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
4			1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	23
5			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	23



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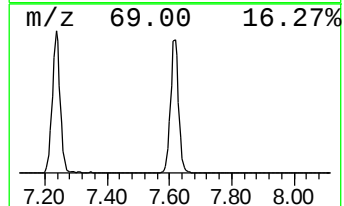
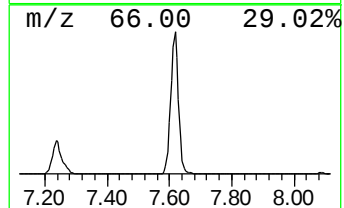
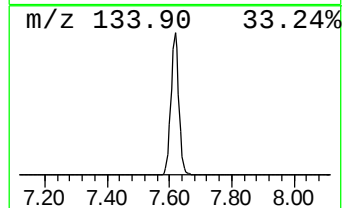
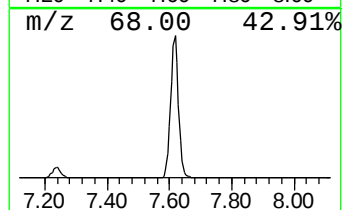
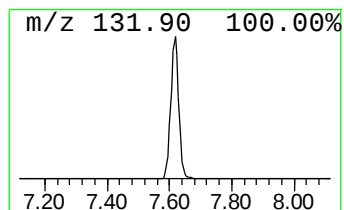
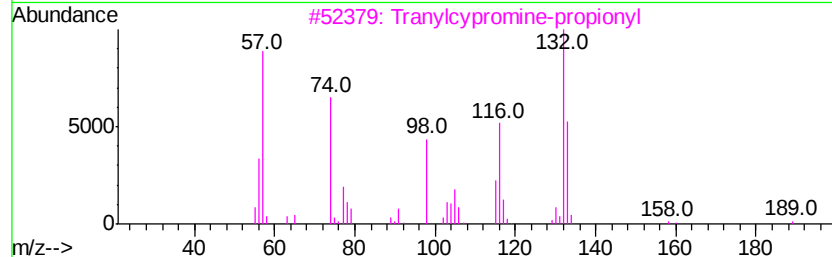
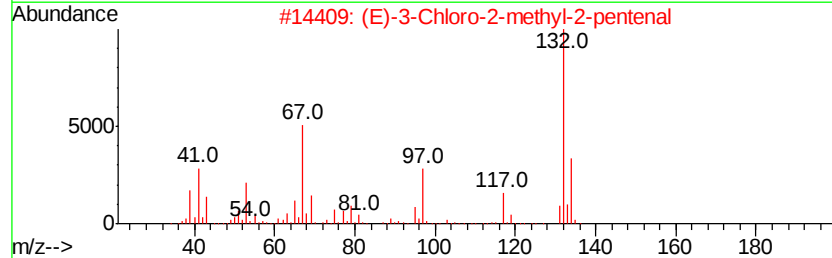
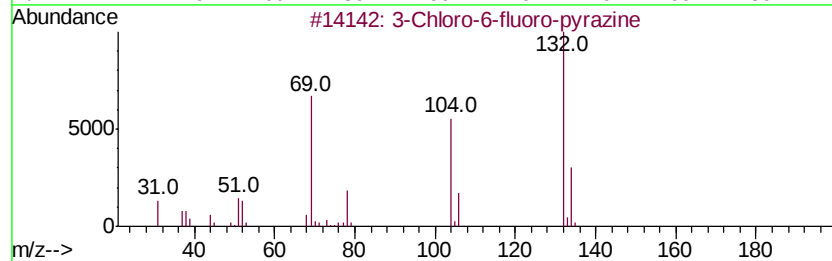
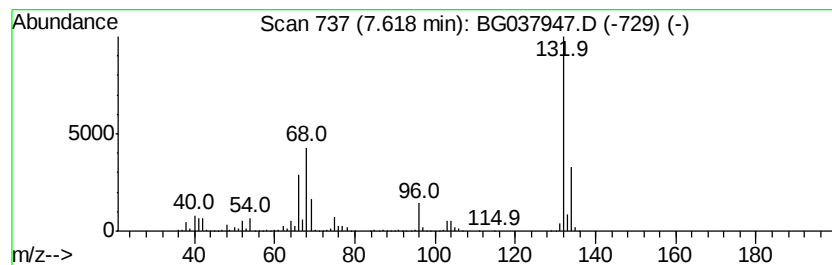
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Peak Number 2 unknown7.62 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.62	115.99 ng	1761850	1,4-Dichlorobenzene-d4	8.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		Amphetaminil	250	C17H18N2	017590-01-1	12
5		Xenon	132	Xe	007440-63-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.17	5.1	ng	77598	1	8.09	303803	20.0
unknown7.62	7.62	116.0	ng	1761850	1	8.09	303803	20.0