

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101618\
 Data File : BG037387.D
 Acq On : 17 Oct 2018 12:08
 Operator : JU/SJ
 Sample : PB113906BL
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB113906BL

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-BG101118.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	5.186	318	323	330	rBV	38386	64129	1.46%	0.290%
2	5.650	394	402	412	rBV	1004977	1646131	37.45%	7.456%
3	7.254	666	675	683	rBV	1011326	1815789	41.31%	8.225%
4	7.636	731	740	752	rBV	946689	1687306	38.38%	7.643%
5	8.112	814	821	831	rVB	213719	383215	8.72%	1.736%
6	8.435	867	876	885	rVB	686554	1242681	28.27%	5.629%
7	9.287	1012	1021	1033	rBV	597090	1071336	24.37%	4.853%
8	10.932	1293	1301	1310	rVB	311162	576373	13.11%	2.611%
9	13.365	1708	1715	1725	rVB	1642535	2690102	61.19%	12.185%
10	14.746	1940	1950	1960	rBV2	542739	902441	20.53%	4.088%
11	16.226	2195	2202	2221	rBV2	1080085	1798642	40.92%	8.147%
12	17.489	2410	2417	2427	rBV2	792858	1227529	27.92%	5.560%
13	20.092	2853	2860	2868	rBV	3134378	4396017	100.00%	19.912%
14	21.773	3139	3146	3160	rVB	800682	1334945	30.37%	6.047%
15	25.116	3704	3715	3730	rBV2	406138	1240080	28.21%	5.617%

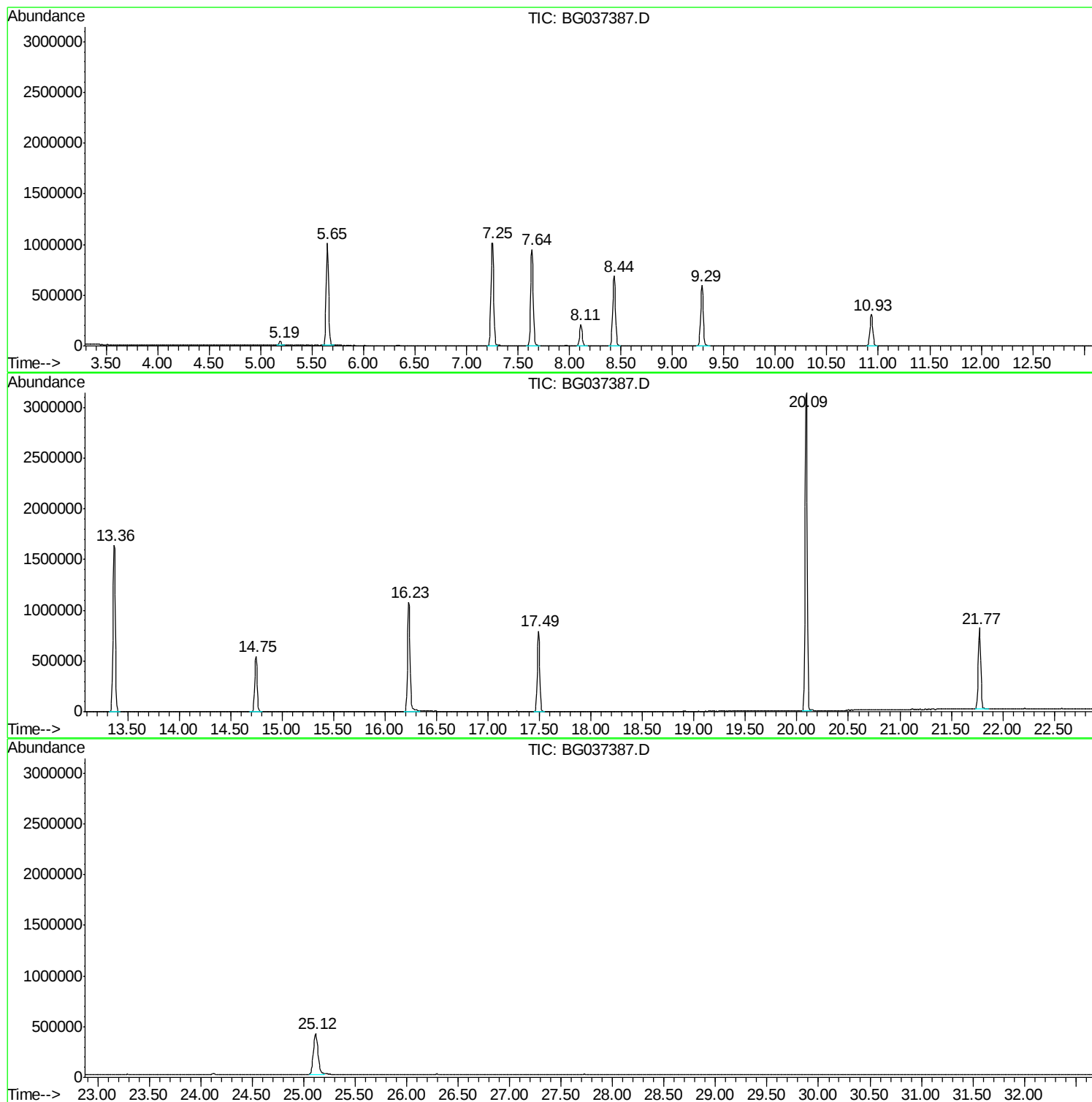
Sum of corrected areas: 22076716

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG101118.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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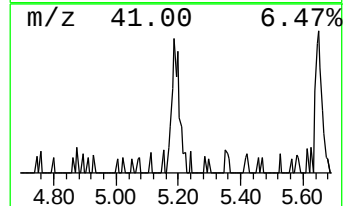
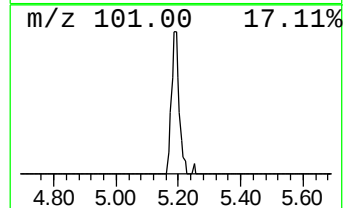
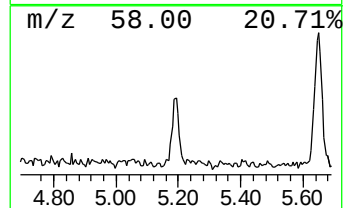
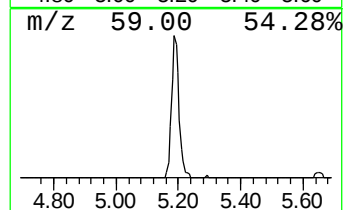
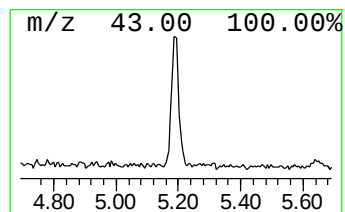
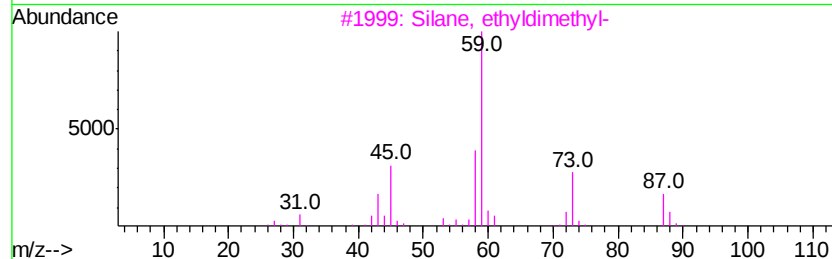
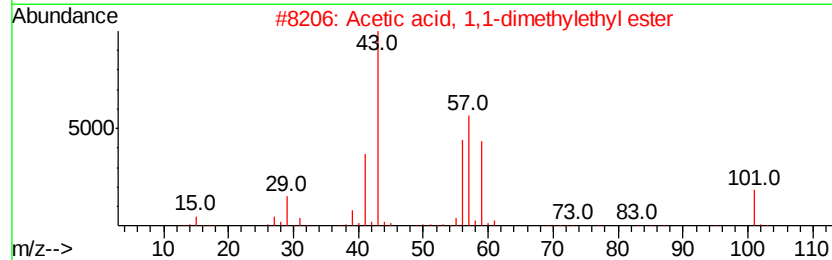
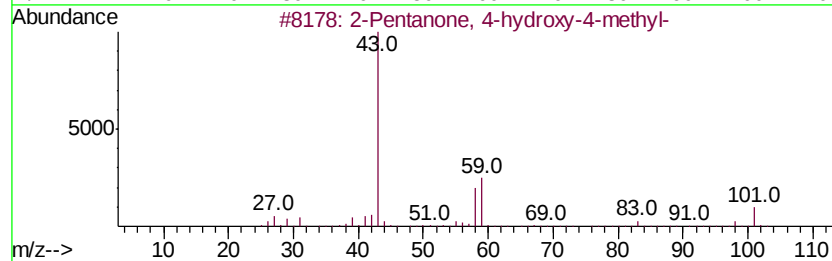
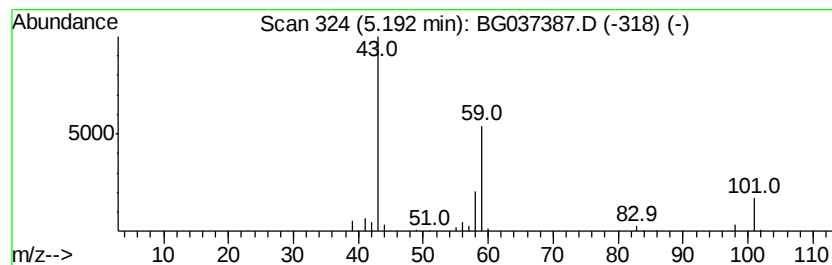
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.19	3.35 ng	64129	1,4-Dichlorobenzene-d4	8.11

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	42
3			Silane, ethyldimethyl-	88	C4H12Si	000758-21-4	28
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5			1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	9



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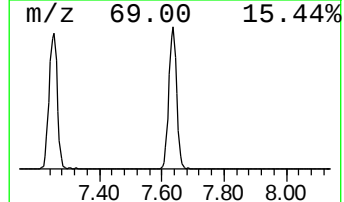
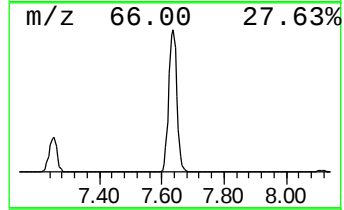
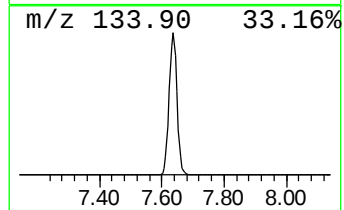
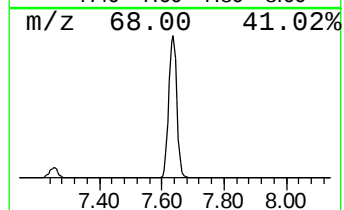
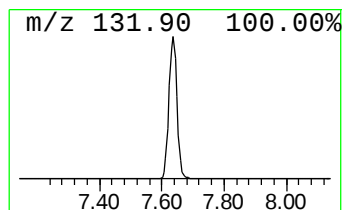
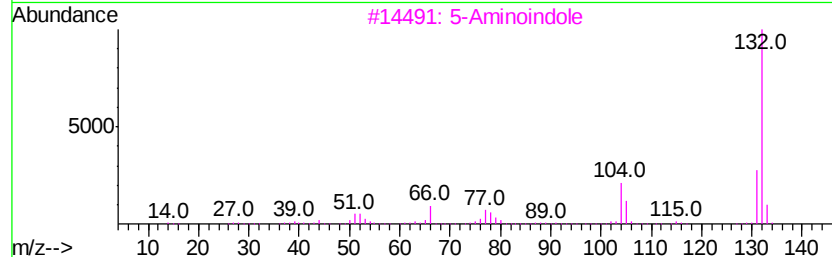
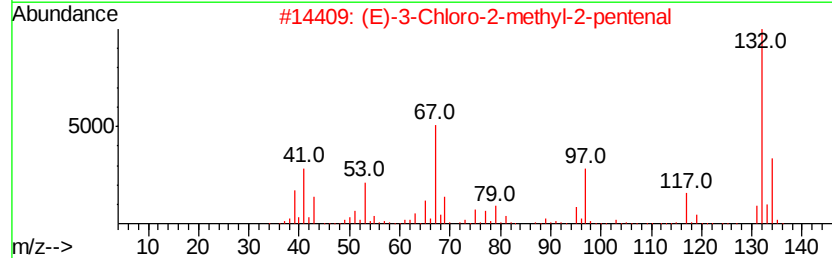
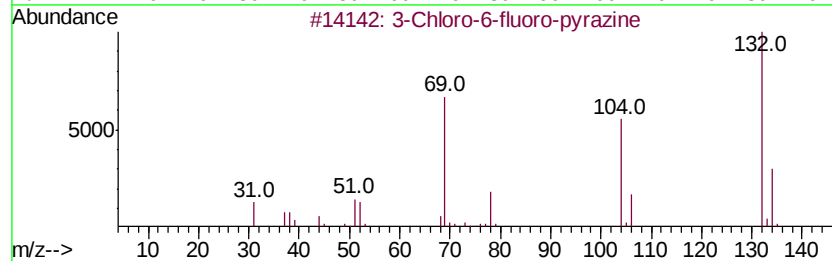
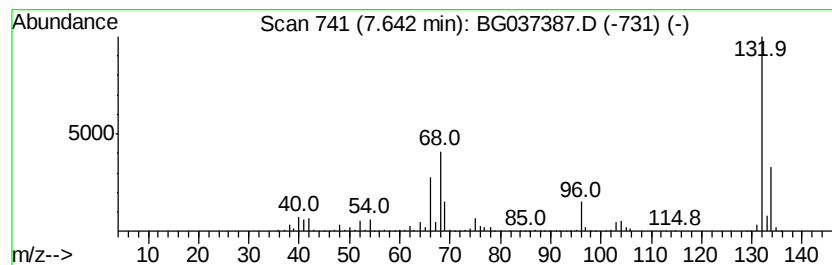
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Peak Number 2 unknown7.64 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.64	88.06 ng	1687310	1,4-Dichlorobenzene-d4	8.11

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3			5-Aminoindole	132	C8H8N2	005192-03-0	10
4			Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5			5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.19	3.4	ng	64129	1	8.11	383215	20.0
unknown7.64	7.64	88.1	ng	1687310	1	8.11	383215	20.0