

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010419\
 Data File : BG038932.D
 Acq On : 4 Jan 2019 16:32
 Operator : JU/SJ
 Sample : PB116158BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB116158BL

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-BG121218.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	5.129	341	347	354	rBV	20754	34778	1.72%	0.398%
2	5.599	419	427	442	rBV	371923	616759	30.49%	7.052%
3	7.203	693	700	711	rBV	383962	685892	33.91%	7.842%
4	7.573	756	763	773	rBV	365050	650641	32.17%	7.439%
5	8.043	835	843	852	rBV	74180	129190	6.39%	1.477%
6	8.367	890	898	907	rBV	278392	494366	24.44%	5.652%
7	9.224	1034	1044	1054	rBV	212417	410421	20.29%	4.692%
8	10.864	1315	1323	1333	rBV	93093	174313	8.62%	1.993%
9	13.302	1731	1738	1747	rBV	678152	1080757	53.43%	12.357%
10	14.683	1967	1973	1985	rVB2	170446	285118	14.10%	3.260%
11	16.175	2220	2227	2242	rBV2	534497	913854	45.18%	10.448%
12	17.432	2434	2441	2456	rVB	251396	401372	19.84%	4.589%
13	20.035	2878	2884	2893	rVB	1507581	2022587	100.00%	23.125%
14	21.710	3162	3169	3181	rBV	255199	450270	22.26%	5.148%
15	25.006	3720	3730	3745	rVB2	130861	396116	19.58%	4.529%

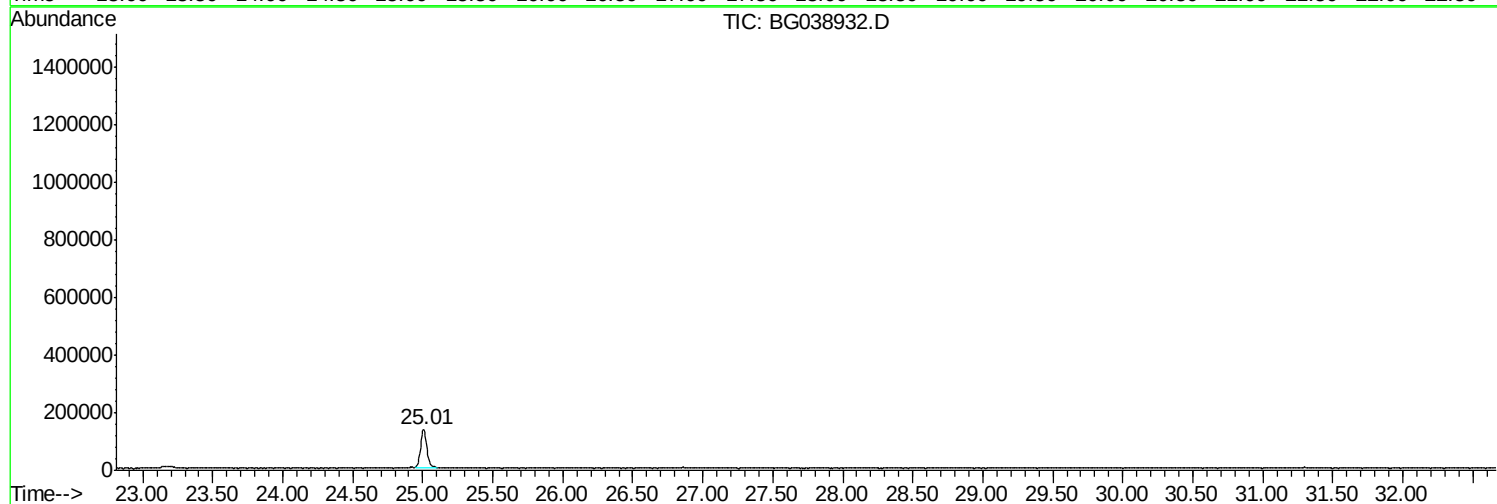
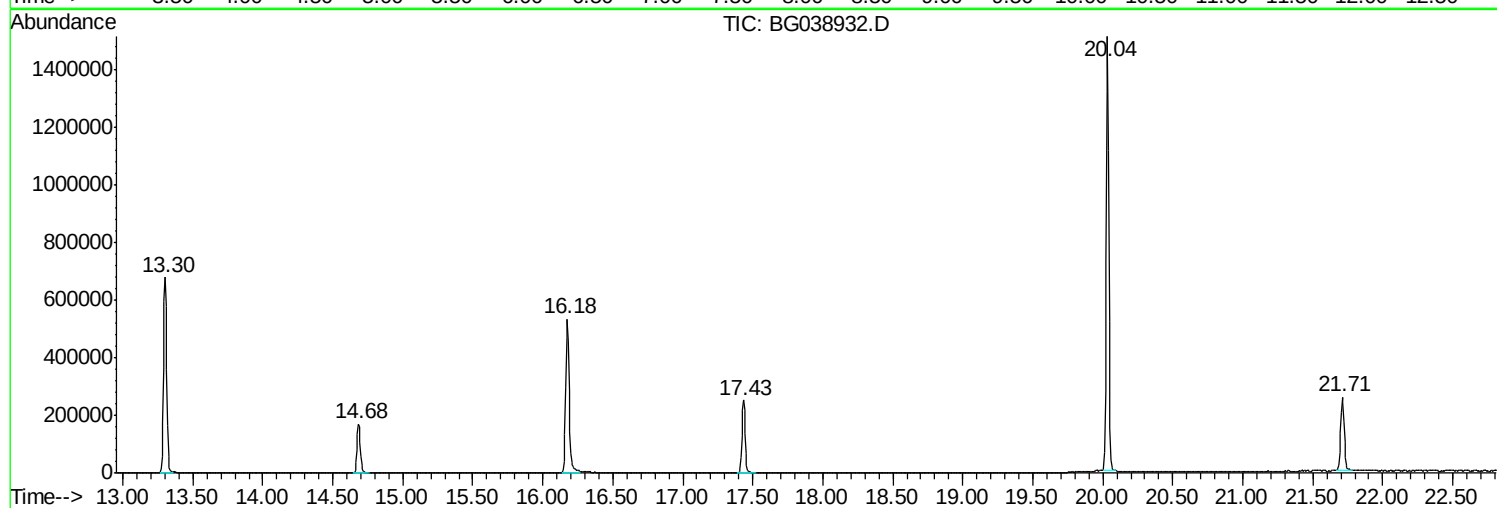
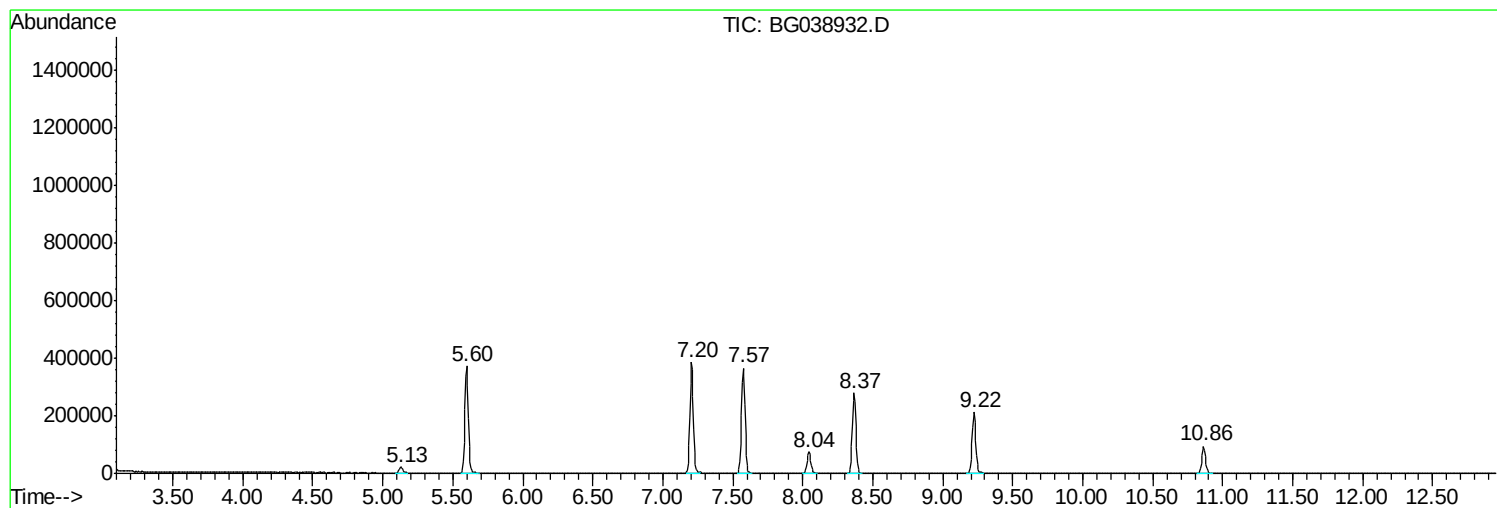
Sum of corrected areas: 8746434

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.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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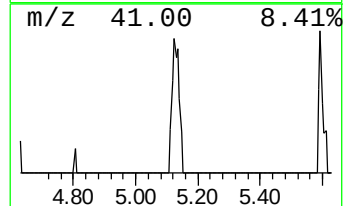
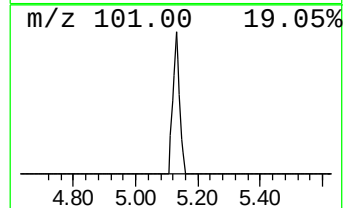
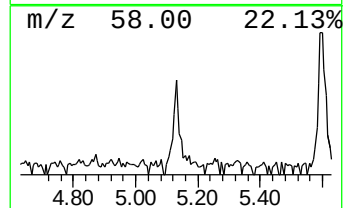
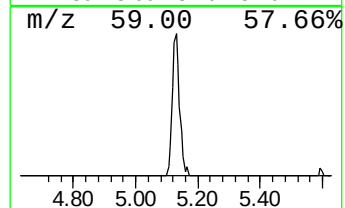
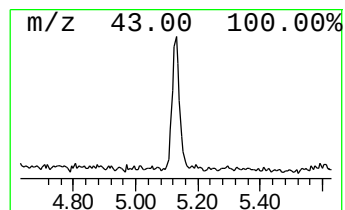
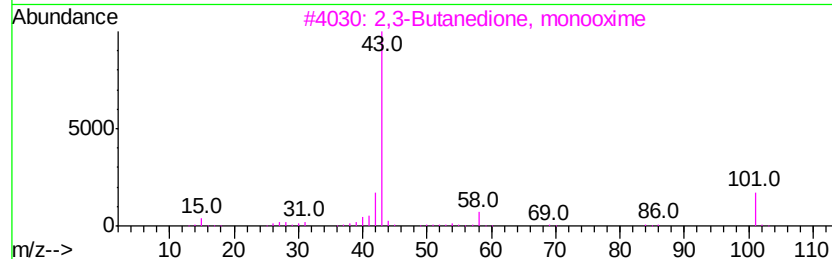
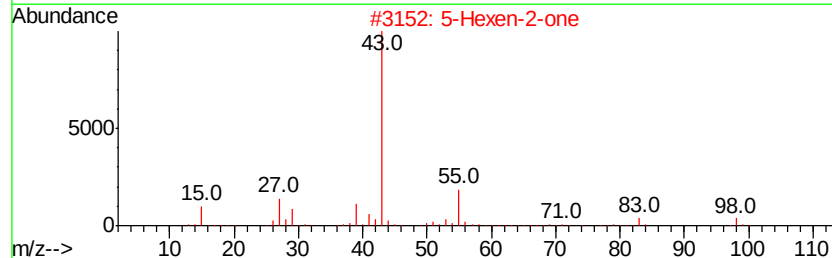
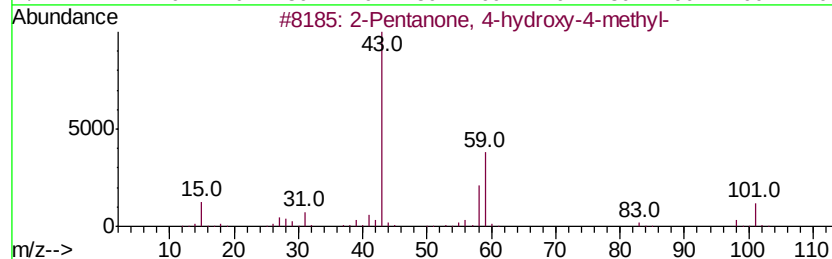
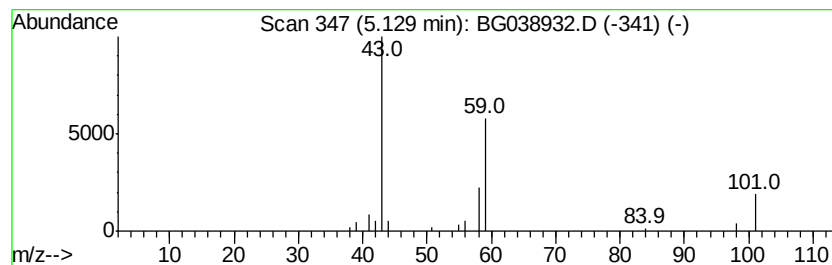
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.13	5.38 ng	34778	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	83
2			5-Hexen-2-one	98	C6H10O	000109-49-9	10
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4			2-Heptanone, 7,7-dichloro-	182	C7H12Cl2O	066241-43-8	9
5			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	9



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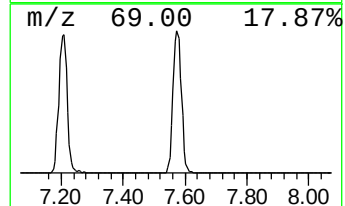
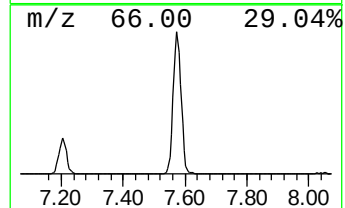
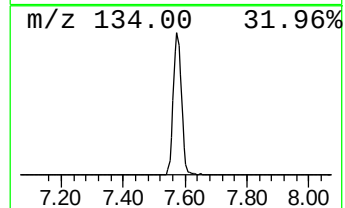
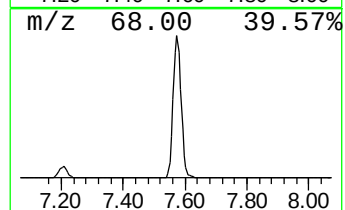
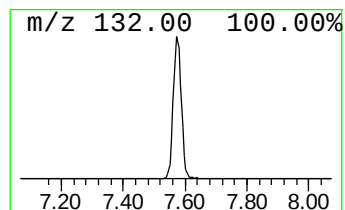
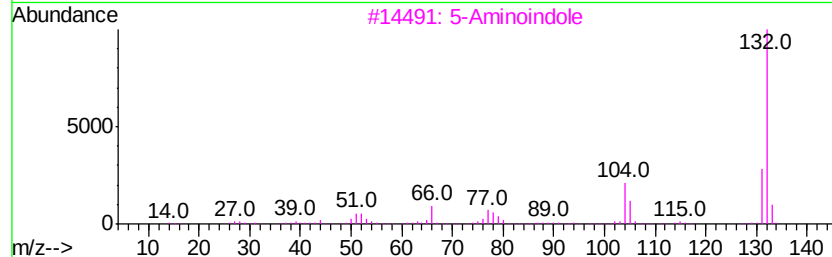
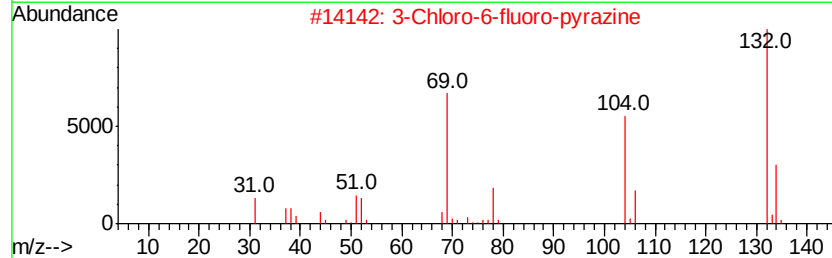
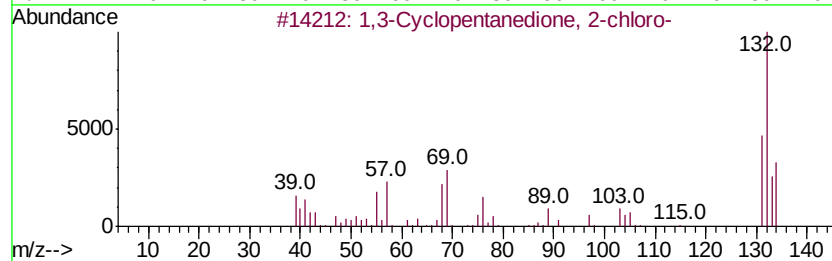
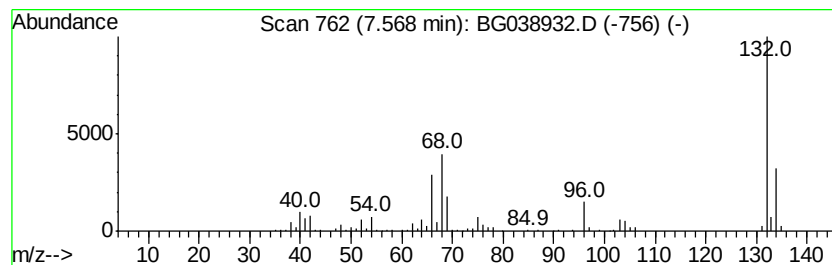
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Peak Number 2 unknown7.57 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.57	100.73 ng	650641	1,4-Dichlorobenzene-d4	8.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	35
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		1,2,4-Trifluorobenzene	132	C6H3F3	000367-23-7	10
5		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.13	5.4	ng	34778	1	8.04	129190	20.0
unknown7.57	7.57	100.7	ng	650641	1	8.04	129190	20.0