

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG122018\
 Data File : BG038767.D
 Acq On : 20 Dec 2018 18:13
 Operator : JU/SJ
 Sample : PB115836BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB115836BL

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	5.139	345	349	359	rBV	19058	30287	1.86%	0.413%
2	5.603	419	428	441	rBV	305063	528721	32.56%	7.218%
3	7.207	694	701	718	rBV	322771	580374	35.74%	7.923%
4	7.583	757	765	776	rBV	307453	548116	33.75%	7.483%
5	8.053	837	845	854	rVB	71003	127989	7.88%	1.747%
6	8.377	891	900	911	rBV	237658	428434	26.38%	5.849%
7	9.234	1038	1046	1057	rBV	186253	347906	21.42%	4.750%
8	10.874	1318	1325	1336	rBV	88644	170013	10.47%	2.321%
9	13.312	1732	1740	1754	rBV	547212	872662	53.74%	11.913%
10	14.693	1968	1975	1986	rBV2	150516	256454	15.79%	3.501%
11	16.185	2222	2229	2248	rBV3	396988	678330	41.77%	9.260%
12	17.442	2436	2443	2453	rBV2	214195	350054	21.56%	4.779%
13	20.039	2879	2885	2895	rBV	1229921	1623998	100.00%	22.171%
14	21.720	3163	3171	3181	rBV	236596	416962	25.68%	5.692%
15	25.010	3721	3731	3746	rBV2	114865	364699	22.46%	4.979%

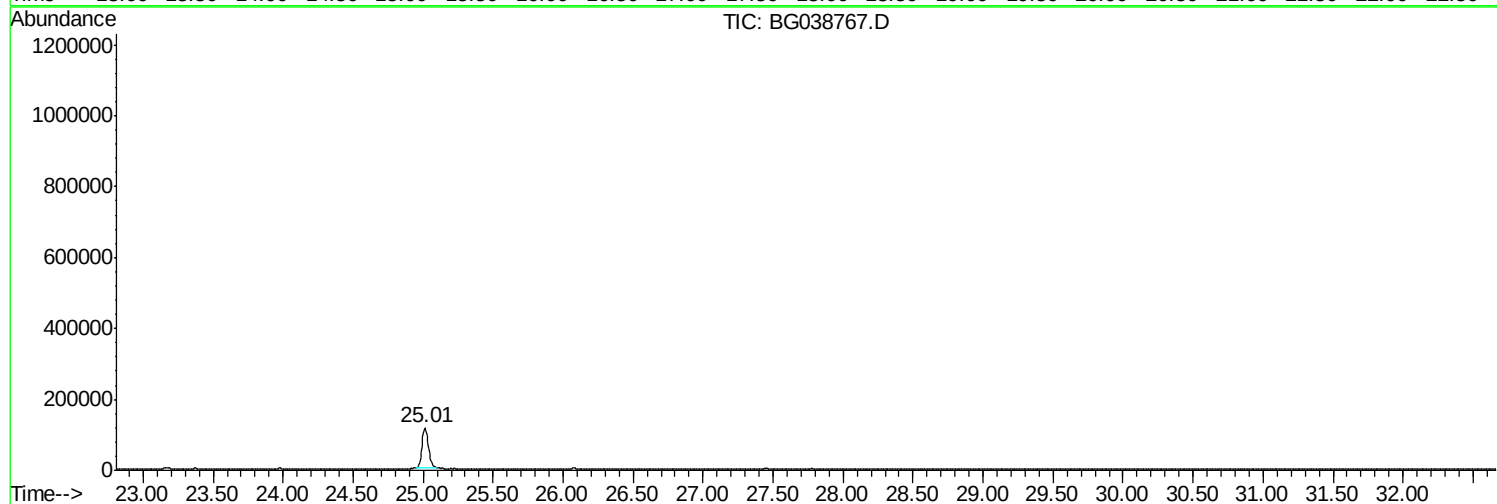
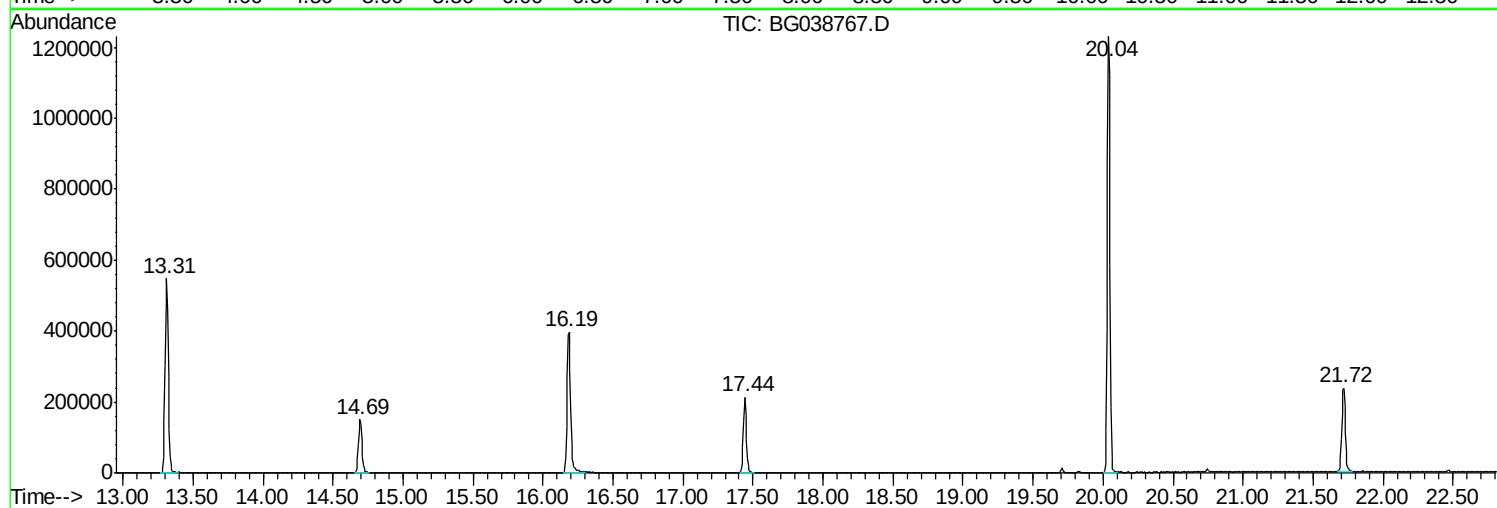
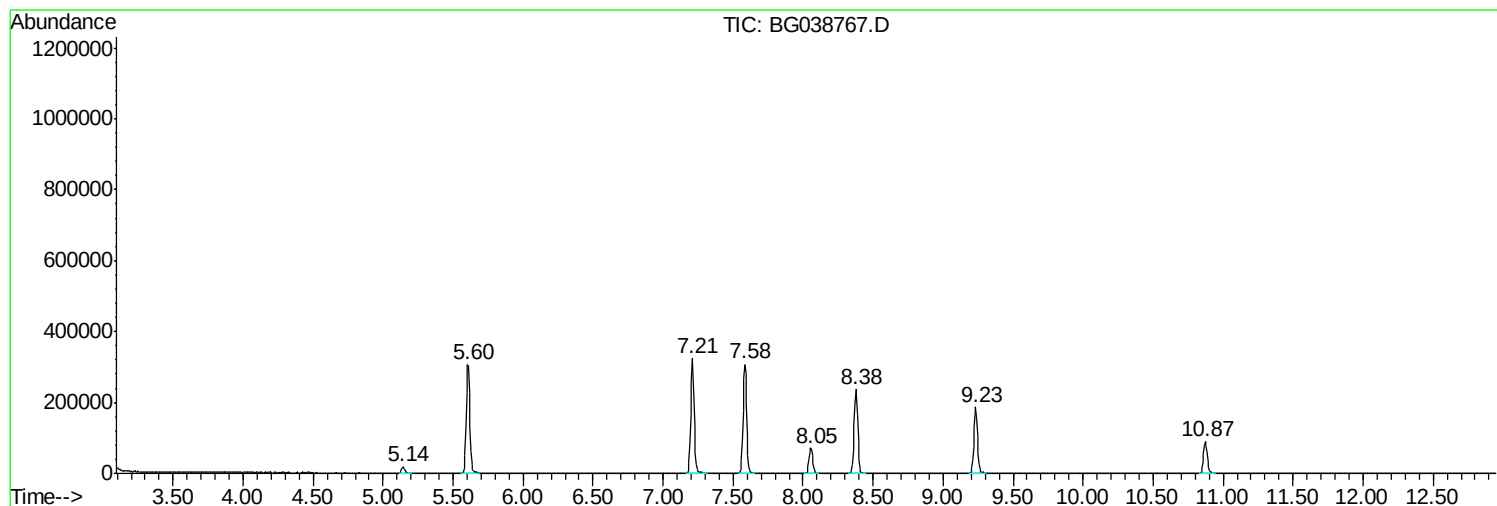
Sum of corrected areas: 7324999

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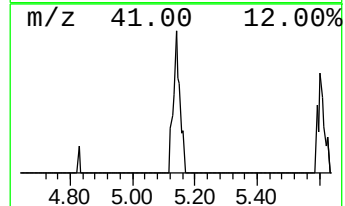
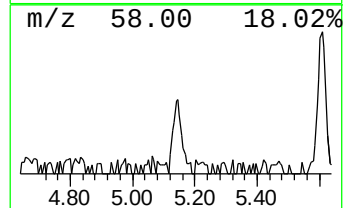
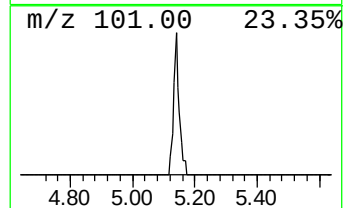
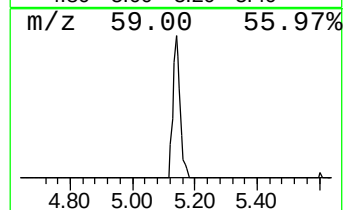
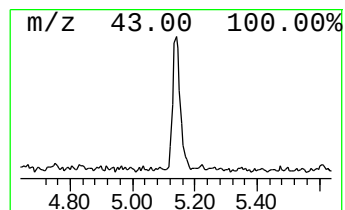
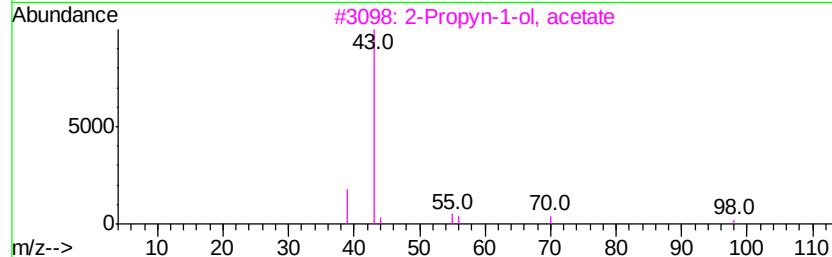
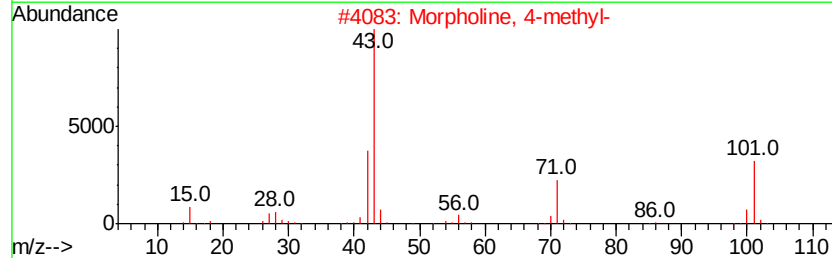
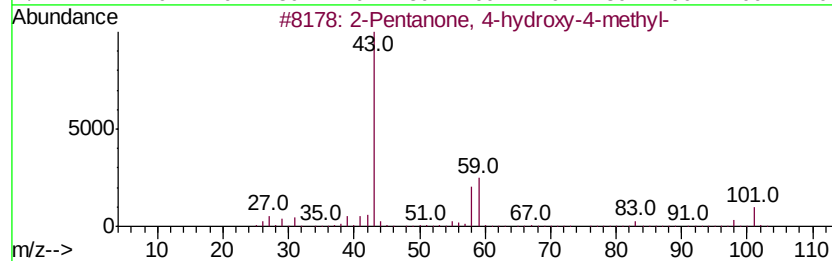
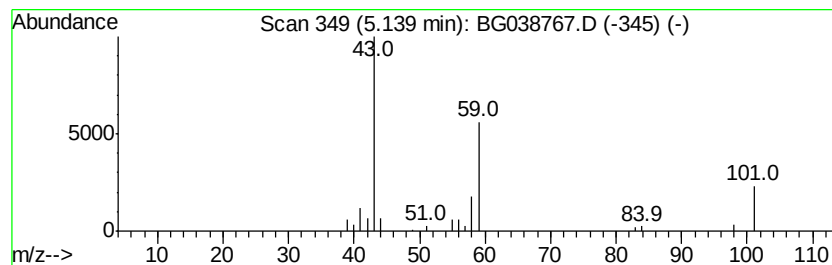
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.14	4.73 ng	30287	1,4-Dichlorobenzene-d4	8.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
3		2-Propyn-1-ol, acetate	98	C5H6O2	000627-09-8	9
4		2-Methyl-2,3-pentanediol	118	C6H14O2	007795-80-4	9
5		1,3,4-Oxadiazole, 2-(acetyloxy)-...	172	C7H12N2O3	066398-57-0	9



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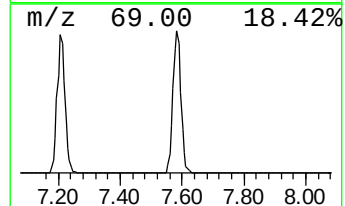
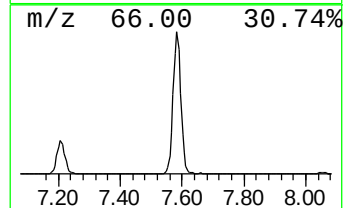
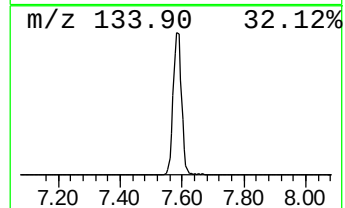
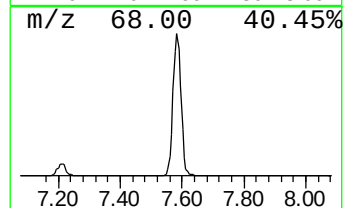
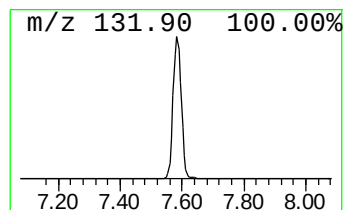
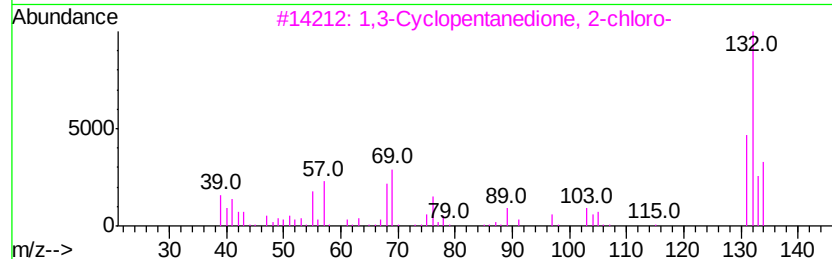
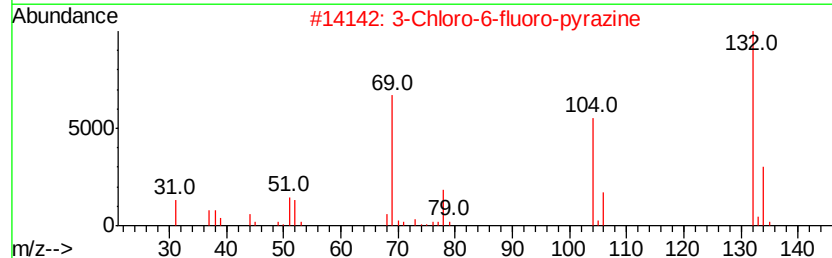
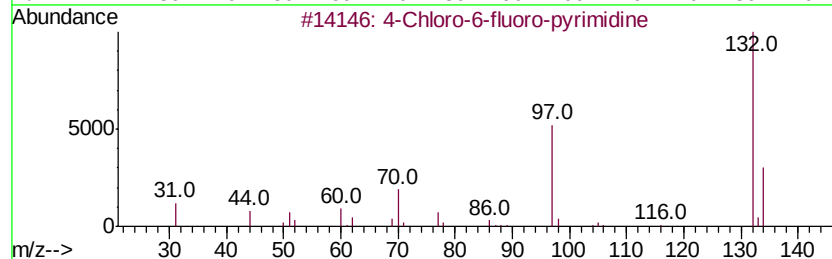
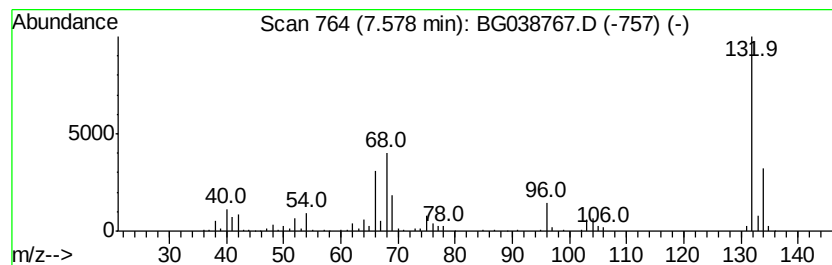
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Peak Number 2 unknown7.58 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	85.65 ng	548116	1,4-Dichlorobenzene-d4	8.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	17
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	12
5		Tranlycypromine-propionyl	189	C12H15NO	1000123-86-3	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.14	4.7	ng	30287	1	8.05	127989	20.0
unknown7.58	7.58	85.7	ng	548116	1	8.05	127989	20.0