

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101218\
 Data File : BG037313.D
 Acq On : 13 Oct 2018 6:06
 Operator : JU/SJ
 Sample : J5383-02
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 100918-SIDI-F-D-M

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	5.197	316	325	334	rBV	49329	85399	1.51%	0.406%
2	5.649	396	402	412	rBV	336024	556313	9.83%	2.643%
3	7.253	666	675	687	rBV	218053	425702	7.52%	2.023%
4	7.641	730	741	751	rBV	629259	1100664	19.45%	5.229%
5	8.117	815	822	833	rVB	198942	355844	6.29%	1.691%
6	8.440	868	877	887	rVB	881944	1598046	28.24%	7.593%
7	9.292	1014	1022	1033	rBV	738044	1352634	23.91%	6.427%
8	10.937	1293	1302	1312	rBV	267138	503573	8.90%	2.393%
9	13.375	1709	1717	1727	rVB	2097684	3343248	59.09%	15.884%
10	14.192	1849	1856	1863	rBV	57520	84633	1.50%	0.402%
11	14.750	1942	1951	1960	rVB2	460891	768257	13.58%	3.650%
12	16.231	2191	2203	2218	rBV2	1095018	1703564	30.11%	8.094%
13	17.494	2410	2418	2427	rBV2	672546	1067515	18.87%	5.072%
14	20.097	2854	2861	2870	rBV	3975882	5658361	100.00%	26.884%
15	21.395	3077	3082	3087	rBV2	45642	70687	1.25%	0.336%
16	21.783	3140	3148	3160	rVB2	693380	1205717	21.31%	5.729%
17	25.126	3706	3717	3732	rVB2	388864	1167455	20.63%	5.547%

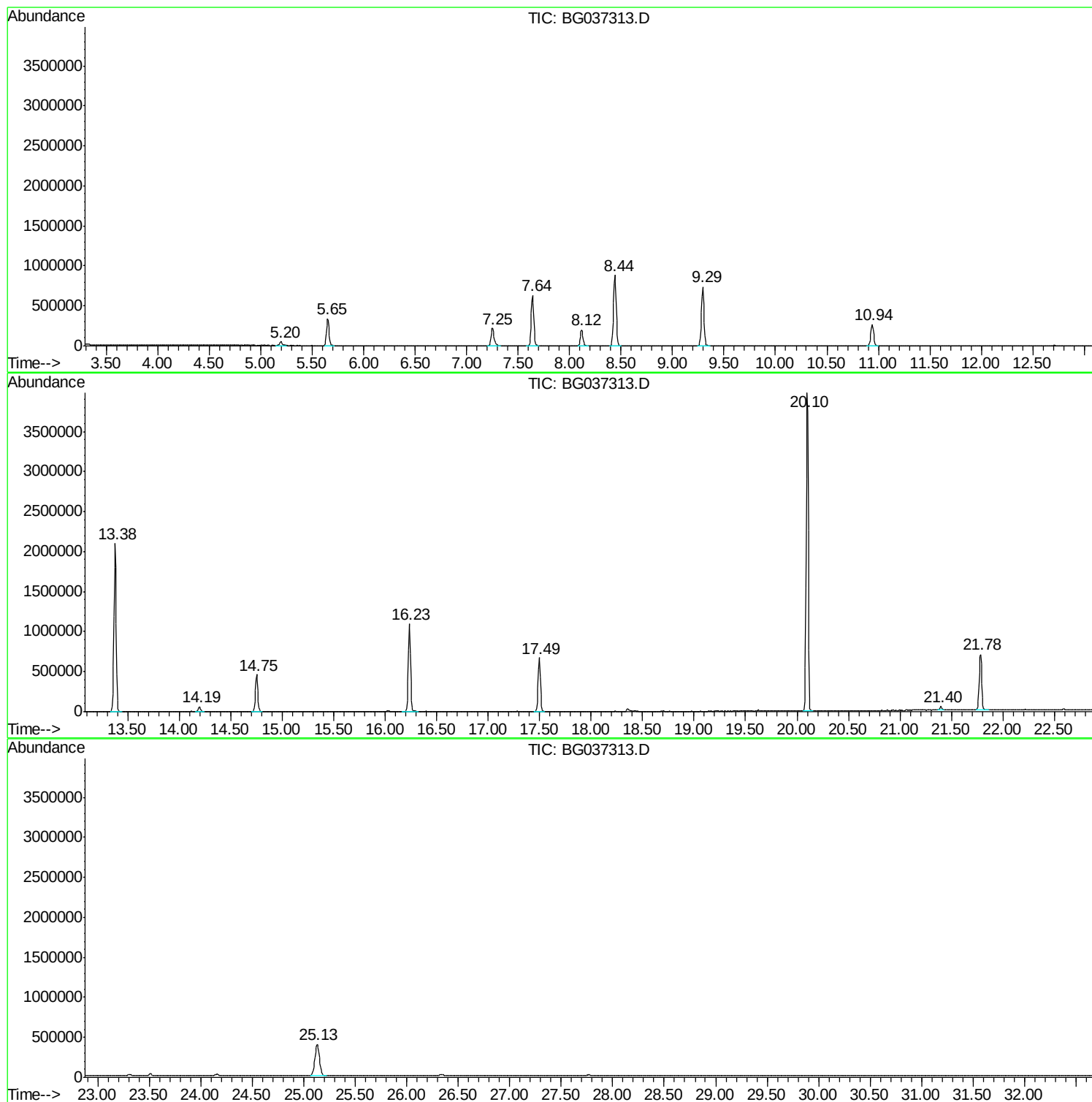
Sum of corrected areas: 21047612

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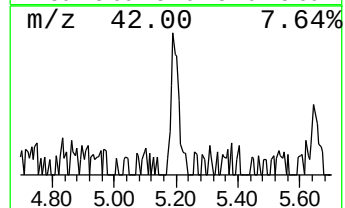
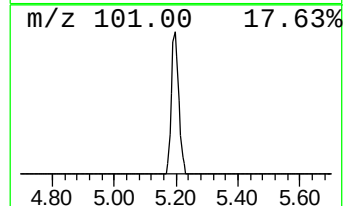
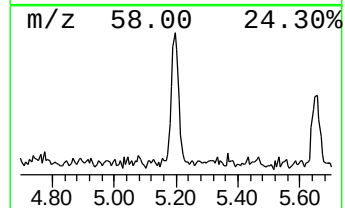
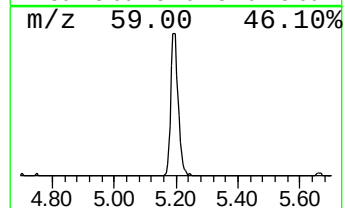
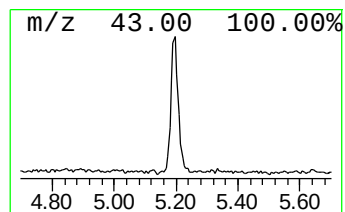
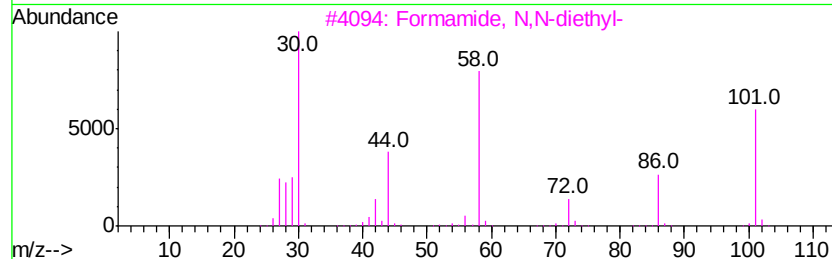
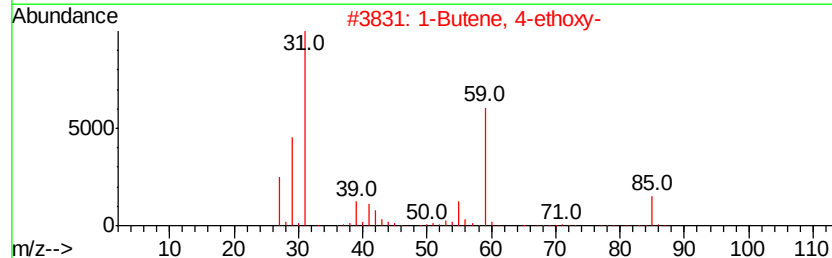
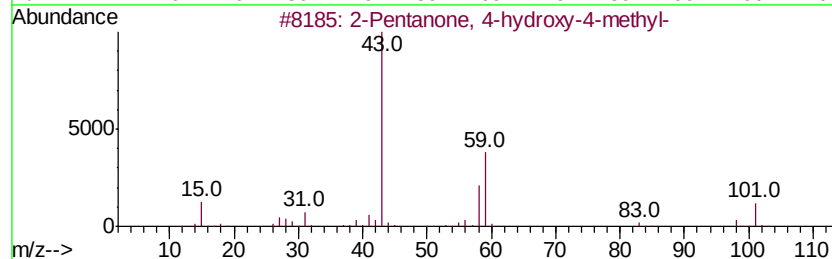
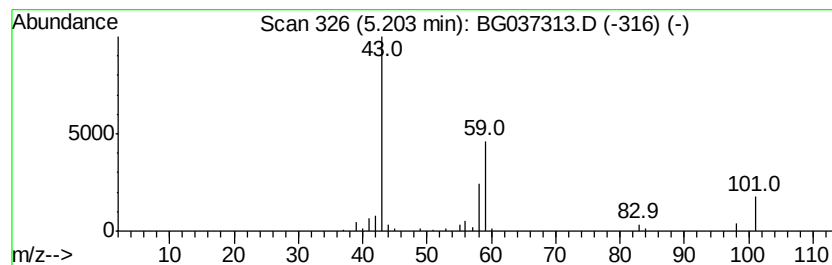
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.20	4.80 ng	85399	1,4-Dichlorobenzene-d4	8.12

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	78
2			1-Butene, 4-ethoxy-	100	C6H12O	044611-46-3	12
3			Formamide, N,N-diethyl-	101	C5H11NO	000617-84-5	9
4			Acetamide, N-(aminocarbonyl)-	102	C3H6N2O2	000591-07-1	9
5			5-Hexen-2-one	98	C6H10O	000109-49-9	9



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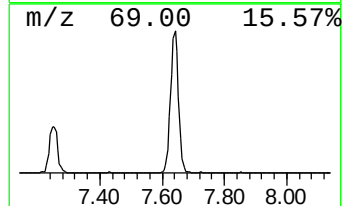
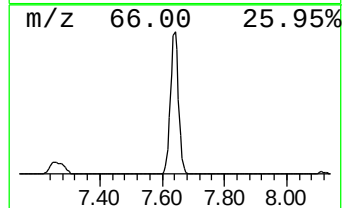
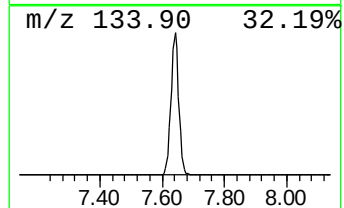
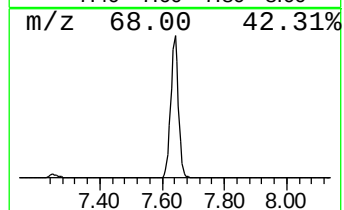
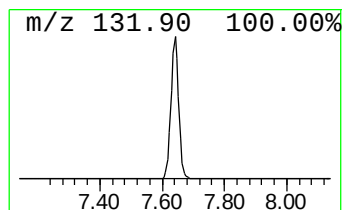
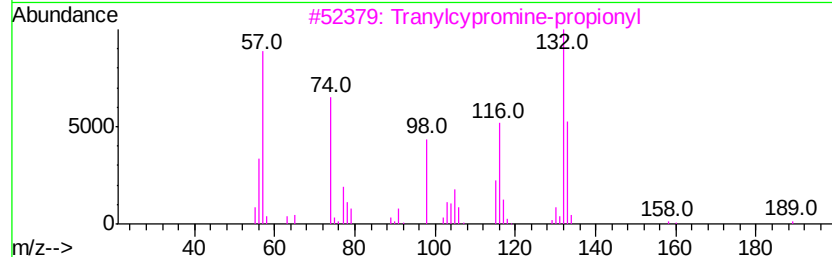
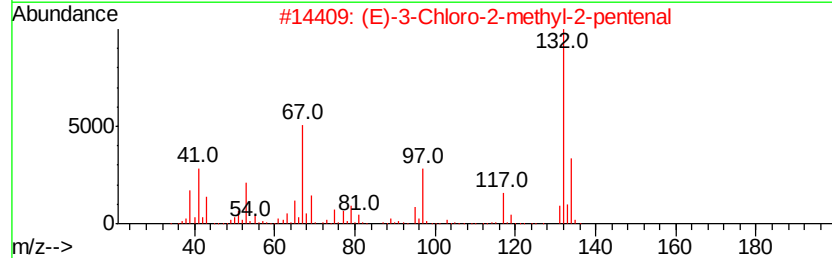
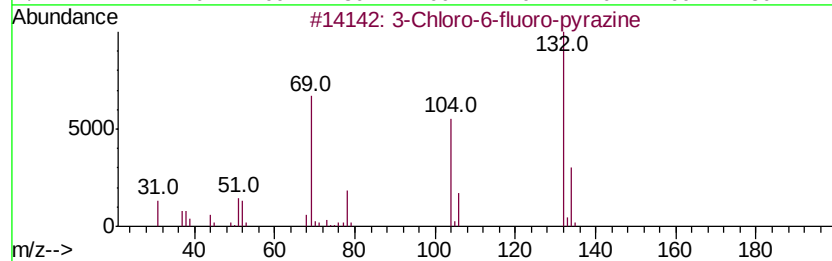
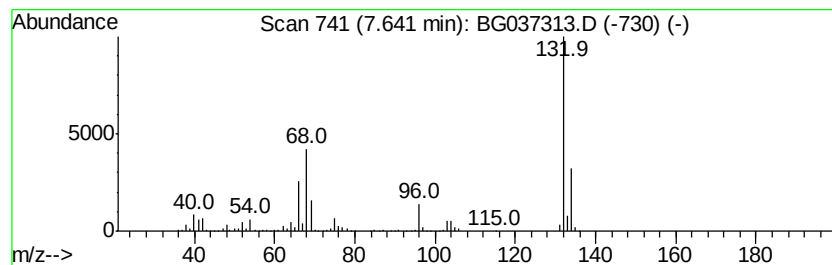
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Peak Number 2 unknown7.64 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.64	61.86 ng	1100660	1,4-Dichlorobenzene-d4	8.12

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	16
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		5-Aminoindole	132	C8H8N2	005192-03-0	10
5		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.20	4.8	ng	85399	1	8.12	355844	20.0
unknown7.64	7.64	61.9	ng	1100660	1	8.12	355844	20.0