

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101218\
 Data File : BG037320.D
 Acq On : 13 Oct 2018 10:40
 Operator : JU/SJ
 Sample : J5383-09
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 100918-TIPC-F-U-B

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 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	5.193	319	324	333	rBV	36750	64338	1.26%	0.325%
2	5.651	395	402	411	rVB	330734	538600	10.51%	2.720%
3	7.249	667	674	685	rBV	213412	395575	7.72%	1.998%
4	7.637	733	740	749	rBV	573167	1024289	19.98%	5.173%
5	8.119	814	822	833	rBV	192517	342877	6.69%	1.732%
6	8.442	869	877	885	rVB	796925	1442181	28.14%	7.284%
7	9.294	1014	1022	1031	rBV	660473	1222380	23.85%	6.173%
8	10.939	1292	1302	1311	rBV	270543	493922	9.64%	2.494%
9	13.371	1709	1716	1724	rBV	1853362	3038302	59.28%	15.344%
10	14.194	1849	1856	1862	rBV	62654	91974	1.79%	0.465%
11	14.746	1942	1950	1959	rVB2	453706	772538	15.07%	3.902%
12	16.233	2196	2203	2213	rBV	1042858	1578304	30.79%	7.971%
13	17.496	2411	2418	2427	rBV2	668644	1059191	20.66%	5.349%
14	18.354	2559	2564	2571	rBV2	41913	63818	1.25%	0.322%
15	20.099	2854	2861	2868	rBV	3716046	5125743	100.00%	25.887%
16	21.391	3076	3081	3090	rBV3	59842	99976	1.95%	0.505%
17	21.779	3140	3147	3154	rBV	681460	1177279	22.97%	5.946%
18	23.506	3435	3441	3448	rVB	51993	101552	1.98%	0.513%
19	25.128	3705	3717	3729	rBV3	383971	1167804	22.78%	5.898%

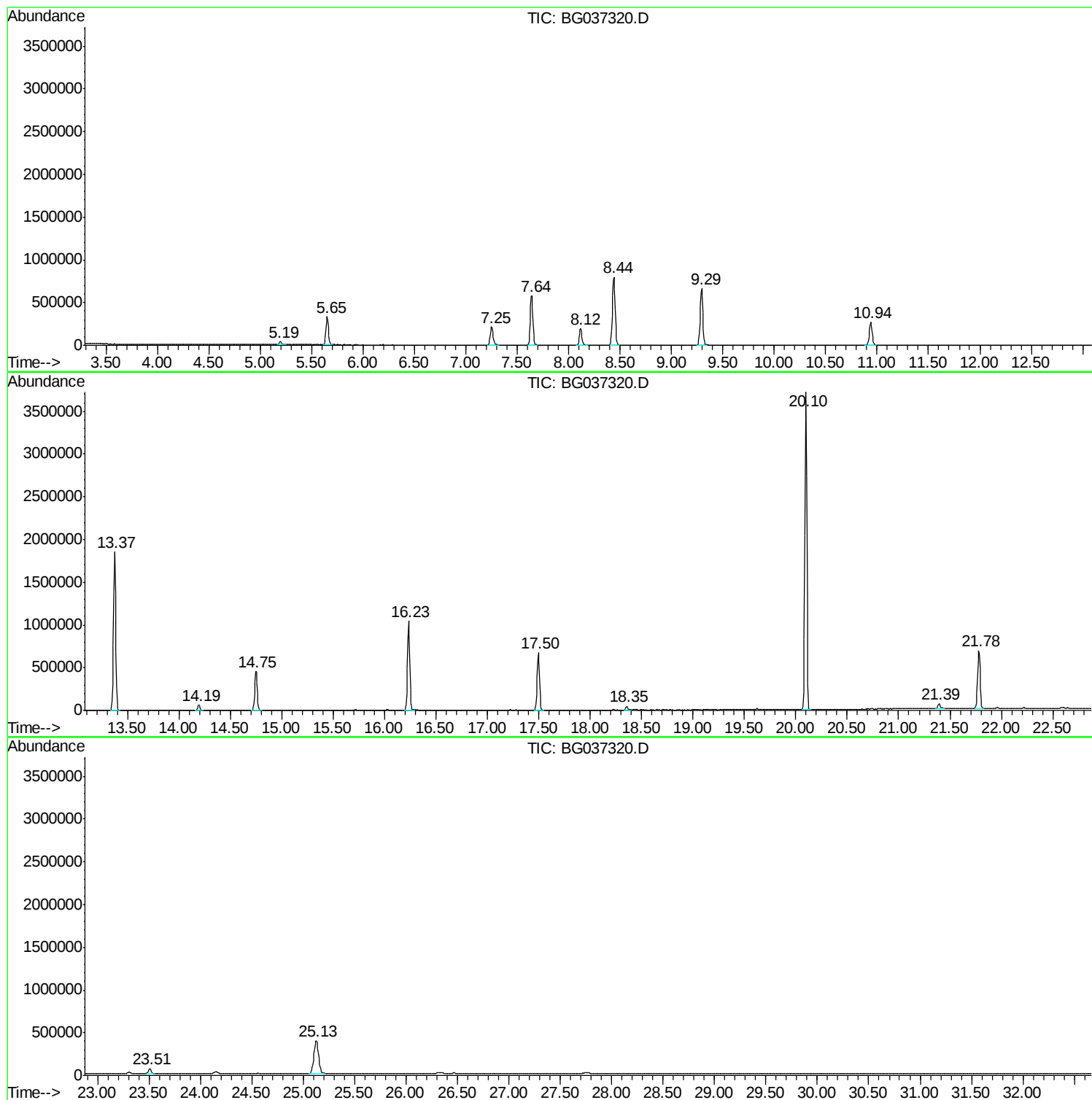
Sum of corrected areas: 19800643

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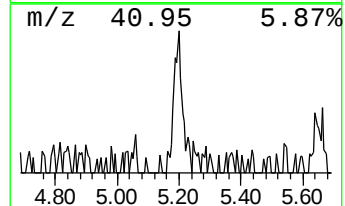
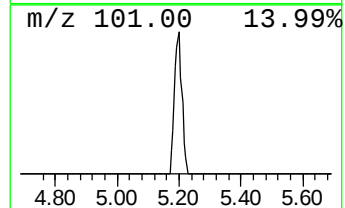
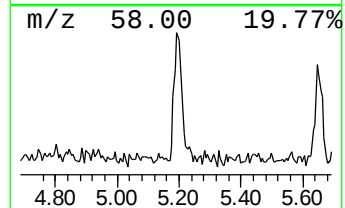
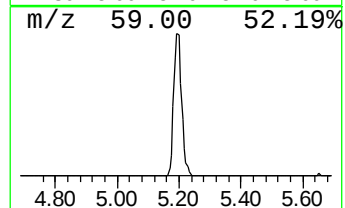
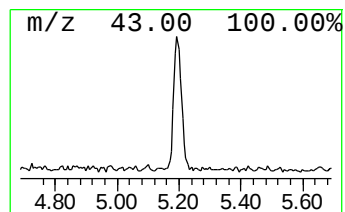
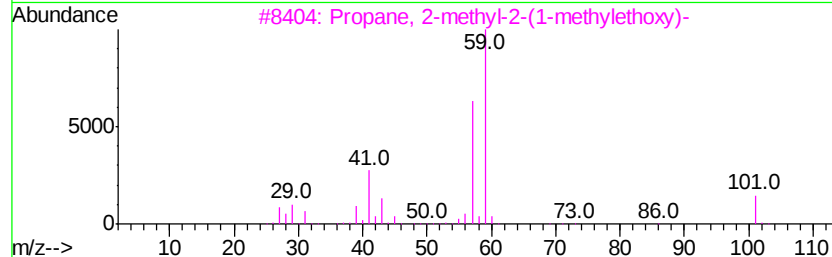
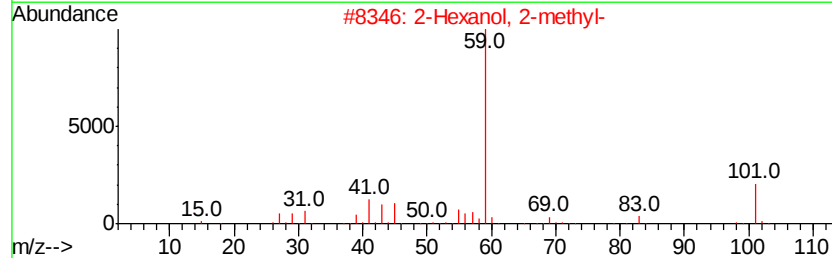
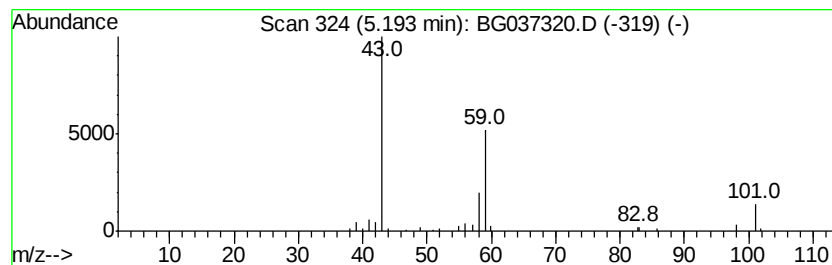
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.19	3.75 ng	64338	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	53
2			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	37
3			Propane, 2-methyl-2-(1-methylethoxy)-	116	C7H16O	017348-59-3	25
4			2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	25
5			3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	17



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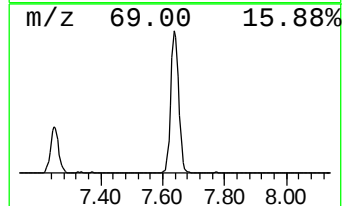
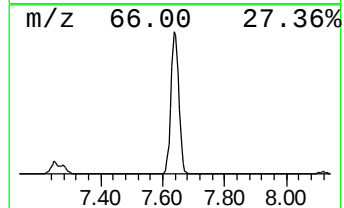
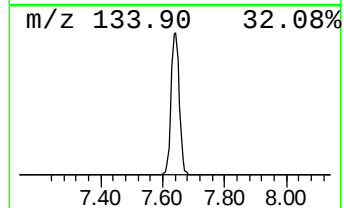
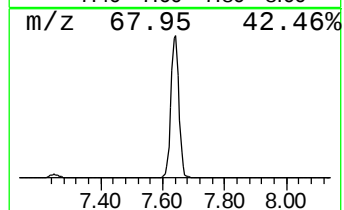
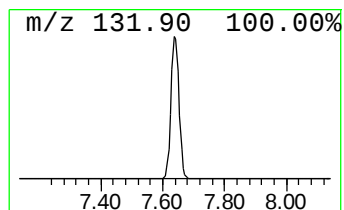
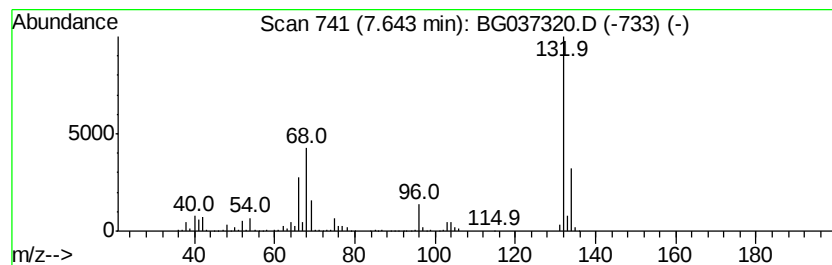
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Peak Number 2 unknown7.64 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.64	59.75 ng	1024290	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		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	10
4		1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	9
5		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.19	3.8	ng	64338	1	8.12	342877	20.0
unknown7.64	7.64	59.8	ng	1024290	1	8.12	342877	20.0