

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101118\
 Data File : BG037282.D
 Acq On : 12 Oct 2018 8:25
 Operator : JU/SJ
 Sample : J5371-03
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 100818-JETT-F-D-S

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	5.200	320	325	330	rBV	38275	63348	1.15%	0.290%
2	5.653	395	402	410	rBV	354288	560351	10.13%	2.568%
3	7.251	668	674	685	rVB	229358	436078	7.88%	1.999%
4	7.645	733	741	749	rBV	624723	1097035	19.83%	5.028%
5	8.121	815	822	830	rBV	218790	388039	7.02%	1.779%
6	8.444	869	877	886	rBV	864463	1539641	27.83%	7.057%
7	9.296	1014	1022	1031	rBV	747417	1362554	24.63%	6.245%
8	10.941	1295	1302	1311	rVB	315081	587962	10.63%	2.695%
9	13.379	1708	1717	1725	rBV	2162416	3499112	63.26%	16.038%
10	14.196	1850	1856	1863	rBV	100905	157030	2.84%	0.720%
11	14.754	1944	1951	1961	rVB2	551171	906535	16.39%	4.155%
12	16.240	2197	2204	2217	rBV3	1052960	1687367	30.50%	7.734%
13	17.498	2410	2418	2427	rBV2	753905	1202836	21.75%	5.513%
14	20.101	2855	2861	2868	rBV	3951764	5531452	100.00%	25.353%
15	21.405	3078	3083	3093	rBV3	47471	82566	1.49%	0.378%
16	21.787	3141	3148	3161	rVB	767027	1308401	23.65%	5.997%
17	24.172	3545	3554	3559	rBV4	25016	61075	1.10%	0.280%
18	25.136	3706	3718	3733	rBV3	420134	1346657	24.35%	6.172%

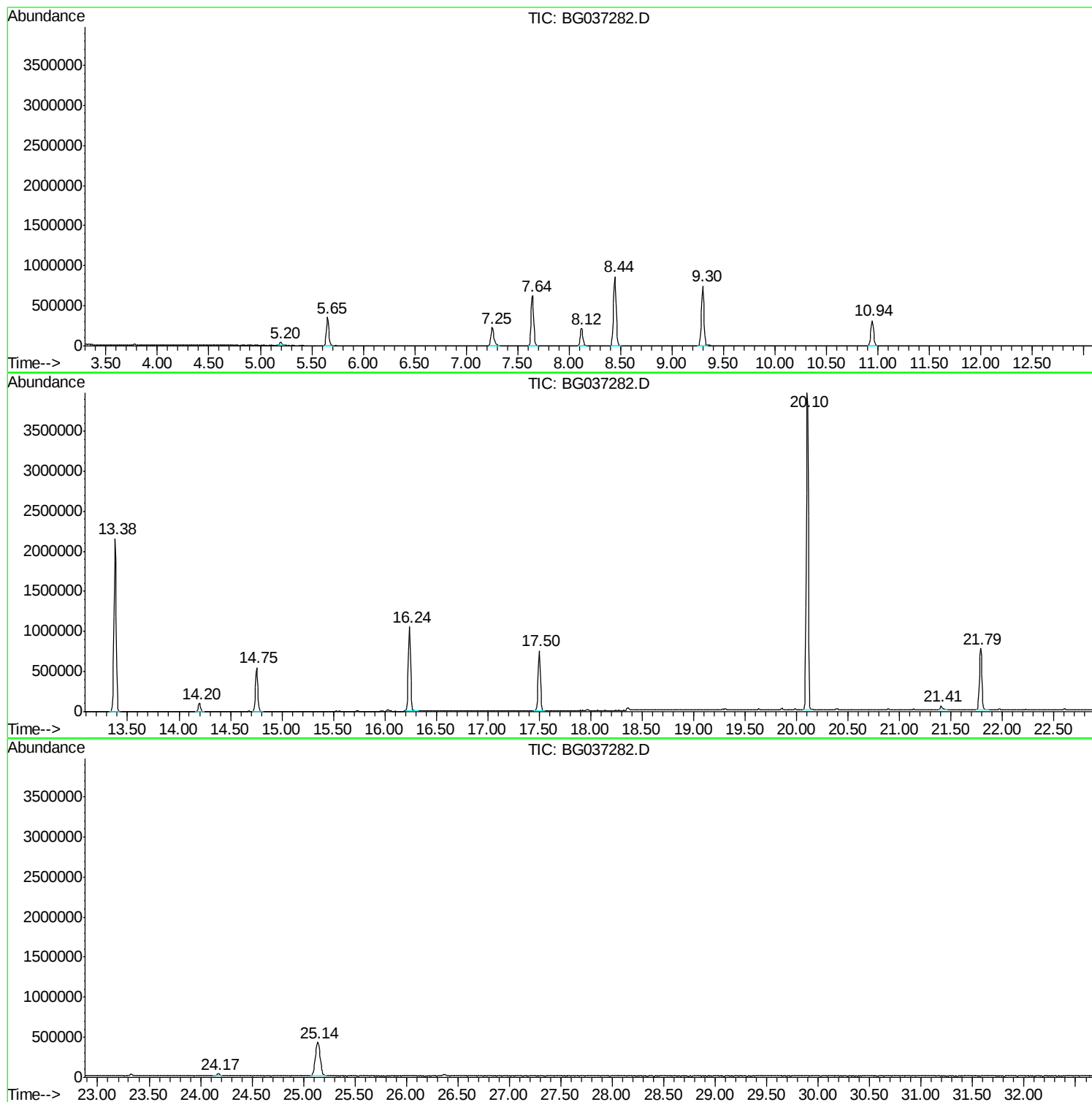
Sum of corrected areas: 21818039

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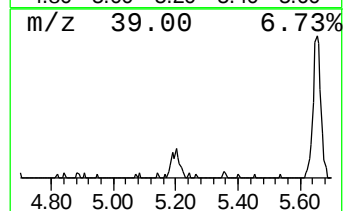
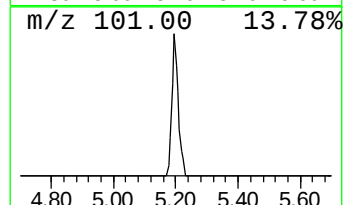
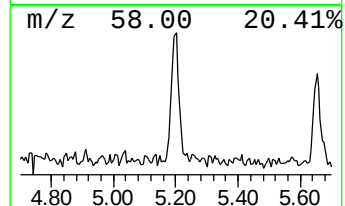
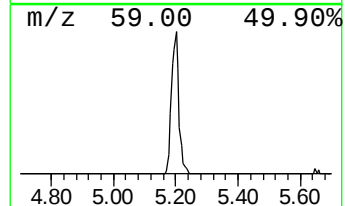
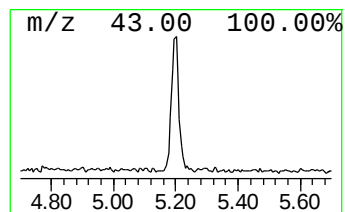
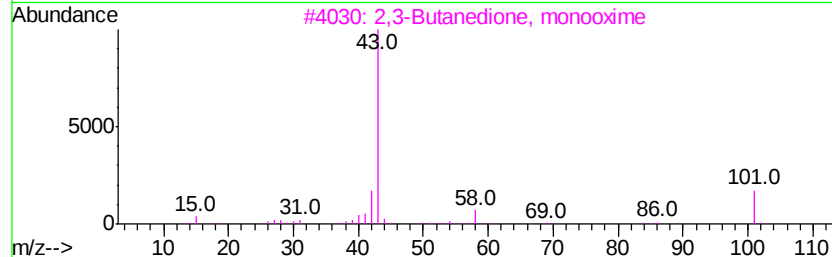
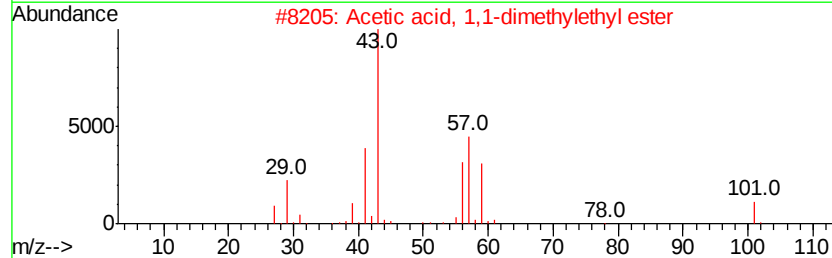
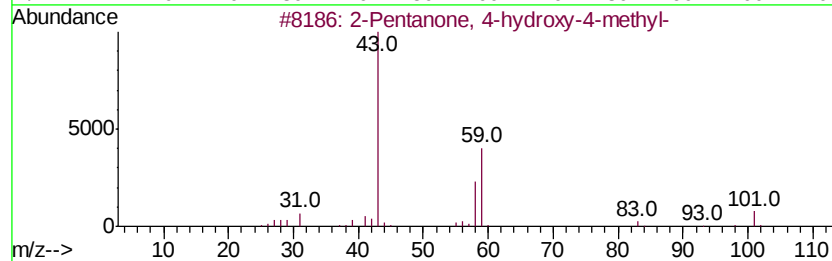
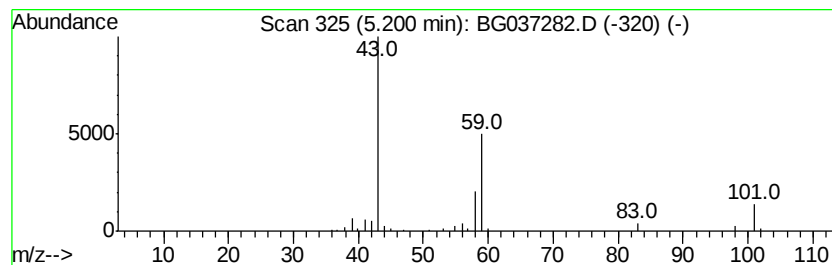
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.20	3.27 ng	63348	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			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	23
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4			5-Hexen-2-one	98	C6H10O	000109-49-9	9
5			4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	9



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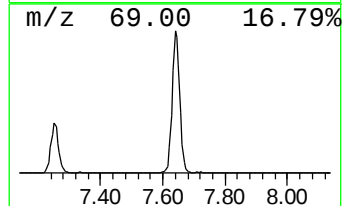
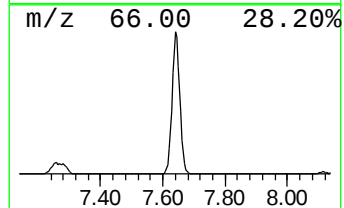
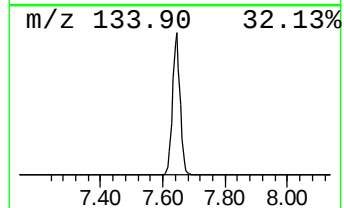
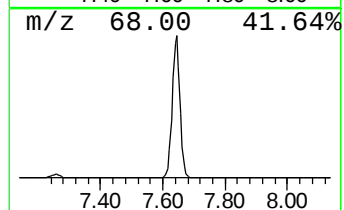
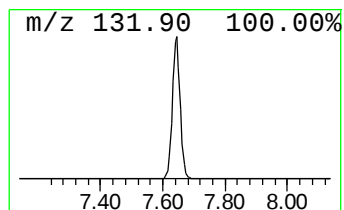
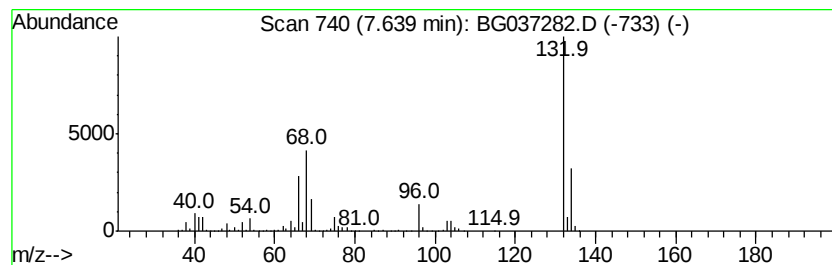
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Peak Number 2 unknown7.64 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.64	56.54 ng	1097040	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	17
3		Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	10
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.20	3.3	ng	63348	1	8.12	388039	20.0
unknown7.64	7.64	56.5	ng	1097040	1	8.12	388039	20.0