

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102318\
 Data File : BG037488.D
 Acq On : 23 Oct 2018 22:45
 Operator : JU/SJ
 Sample : J5626-06
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EB-2018.10.19-SED

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-BG101818.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	5.189	316	323	333	rBV	38897	69028	1.43%	0.317%
2	5.647	394	401	409	rBV	485825	824465	17.10%	3.789%
3	7.245	665	673	684	rBV	324784	579945	12.03%	2.665%
4	7.633	730	739	753	rVB	865615	1554343	32.25%	7.143%
5	8.109	812	820	831	rVB	218803	389402	8.08%	1.789%
6	8.432	867	875	886	rVB	764875	1415358	29.36%	6.504%
7	9.284	1011	1020	1030	rBV	691374	1275238	26.46%	5.860%
8	10.929	1292	1300	1308	rVB	320258	596826	12.38%	2.743%
9	13.362	1707	1714	1726	rBV	1945759	3119783	64.72%	14.337%
10	14.742	1941	1949	1958	rBV2	549903	927701	19.25%	4.263%
11	16.223	2194	2201	2215	rBV	1536432	2436287	50.54%	11.196%
12	17.486	2409	2416	2427	rBV	779843	1194139	24.77%	5.488%
13	20.083	2852	2858	2872	rBV	3375726	4820159	100.00%	22.150%
14	21.769	3138	3145	3155	rVB2	750672	1257429	26.09%	5.778%
15	24.096	3535	3541	3552	rVB6	19727	53898	1.12%	0.248%
16	25.107	3701	3713	3726	rBV2	398859	1247007	25.87%	5.730%

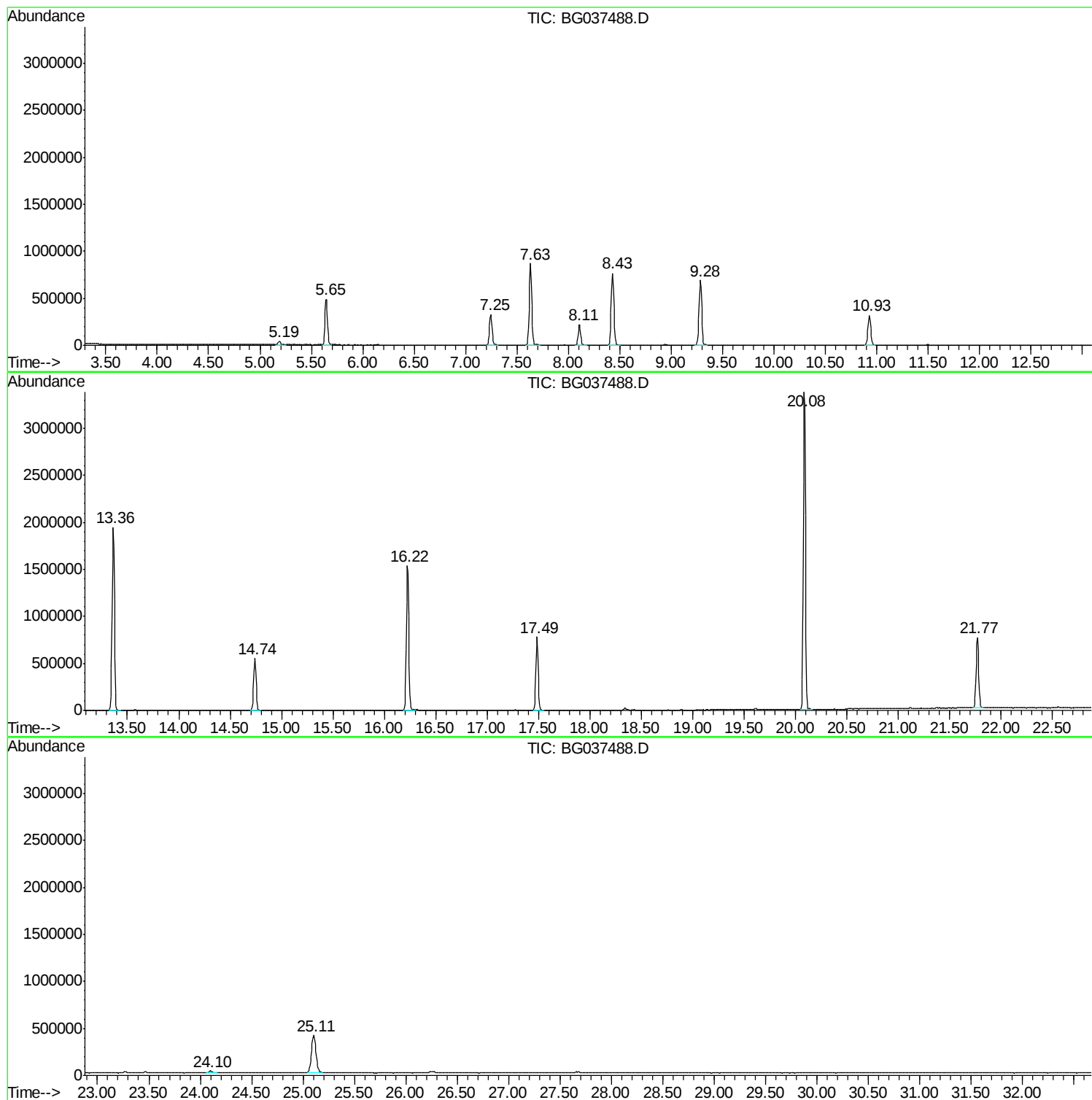
Sum of corrected areas: 21761008

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG101818.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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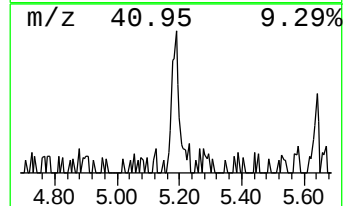
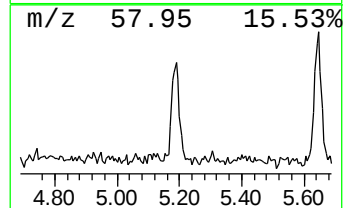
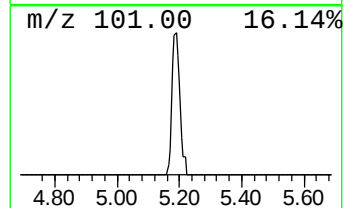
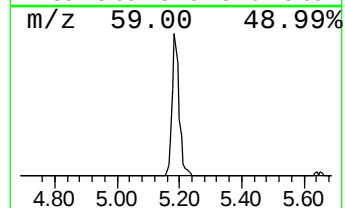
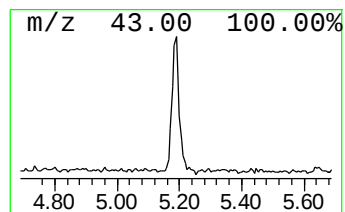
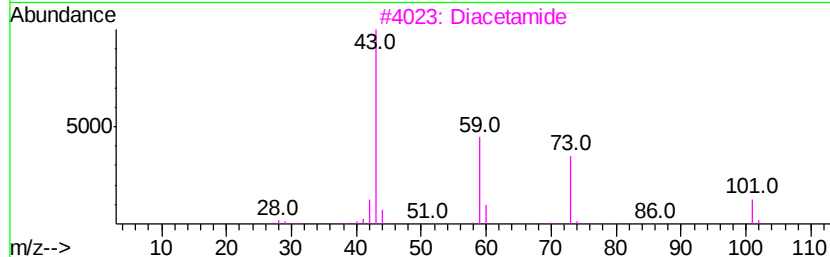
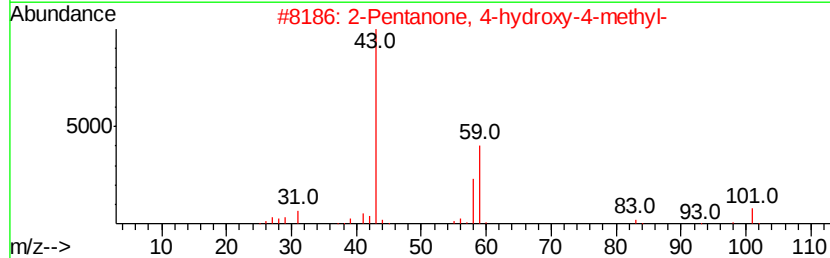
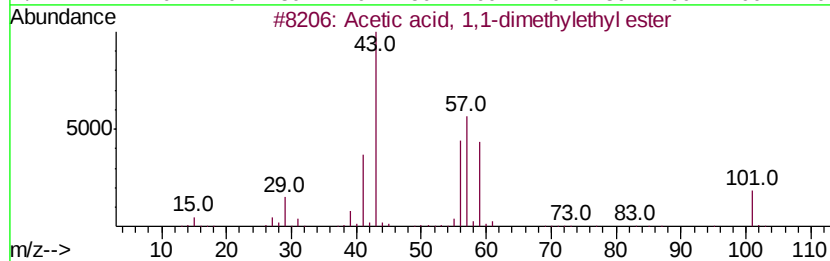
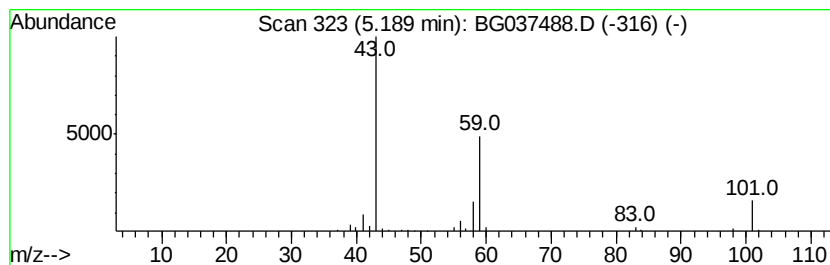
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.19	3.55 ng	69028	1,4-Dichlorobenzene-d4	8.10

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	45
2			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	42
3			Diacetamide	101	C4H7NO2	000625-77-4	9
4			2-Pentanone, 4-mercapto-4-methyl-	132	C6H12OS	019872-52-7	9
5			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9



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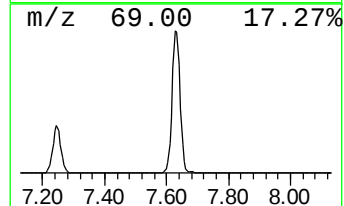
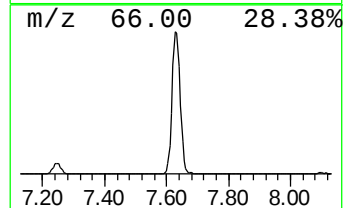
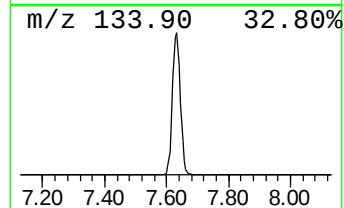
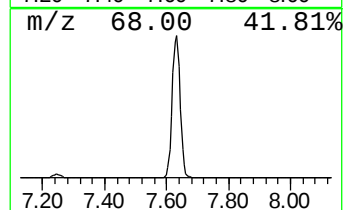
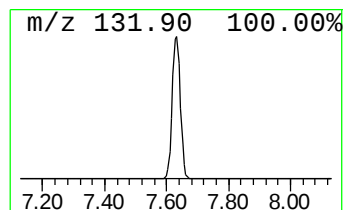
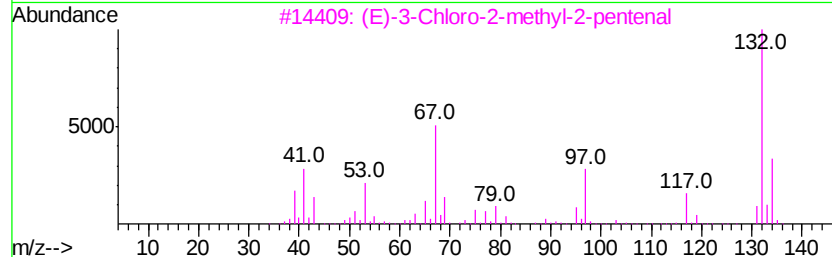
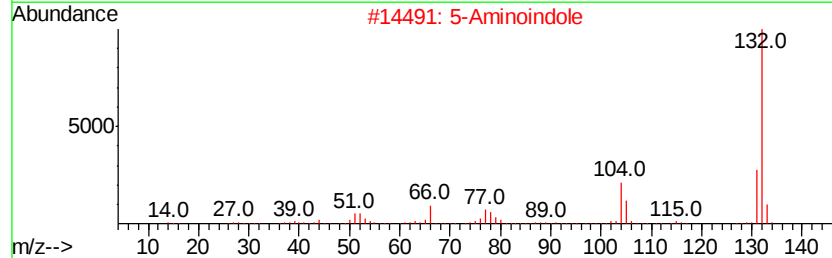
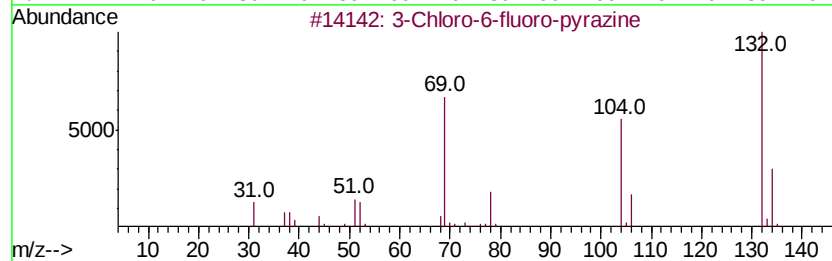
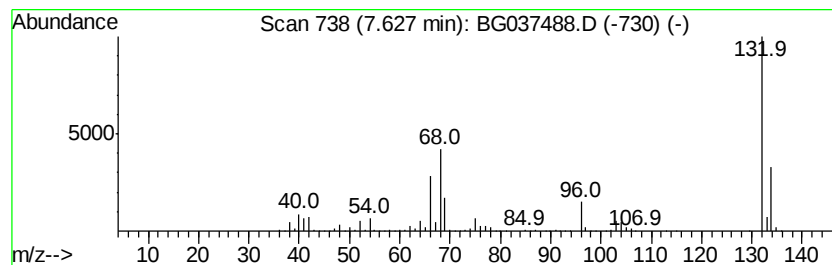
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Peak Number 2 unknown7.63 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.63	79.83 ng	1554340	1,4-Dichlorobenzene-d4	8.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.19	3.5	ng	69028	1	8.10	389402	20.0
unknown7.63	7.63	79.8	ng	1554340	1	8.10	389402	20.0