

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG121918\
 Data File : BG038733.D
 Acq On : 19 Dec 2018 19:05
 Operator : JU/SJ
 Sample : PB115792BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB115792BL

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	5.138	344	349	357	rVB	16459	26809	1.83%	0.372%
2	5.608	421	429	442	rBV	300313	505831	34.58%	7.020%
3	7.212	692	702	713	rBV	306830	556266	38.02%	7.720%
4	7.582	758	765	778	rBV	297399	530590	36.27%	7.363%
5	8.052	838	845	855	rVB	80612	144904	9.90%	2.011%
6	8.376	892	900	909	rBV	228404	413349	28.25%	5.736%
7	9.233	1037	1046	1058	rBV	174991	335163	22.91%	4.651%
8	10.873	1318	1325	1334	rBV	103360	199132	13.61%	2.763%
9	13.311	1733	1740	1752	rBV	512471	830495	56.77%	11.525%
10	14.692	1969	1975	1986	rBV2	174692	299881	20.50%	4.162%
11	16.184	2222	2229	2246	rBV2	357802	618098	42.25%	8.578%
12	17.441	2436	2443	2455	rBV	248297	395674	27.05%	5.491%
13	20.044	2879	2886	2899	rVB	1058960	1462953	100.00%	20.302%
14	21.719	3164	3171	3182	rBV	274120	465807	31.84%	6.464%
15	25.015	3720	3732	3751	rBV2	137683	420869	28.77%	5.841%

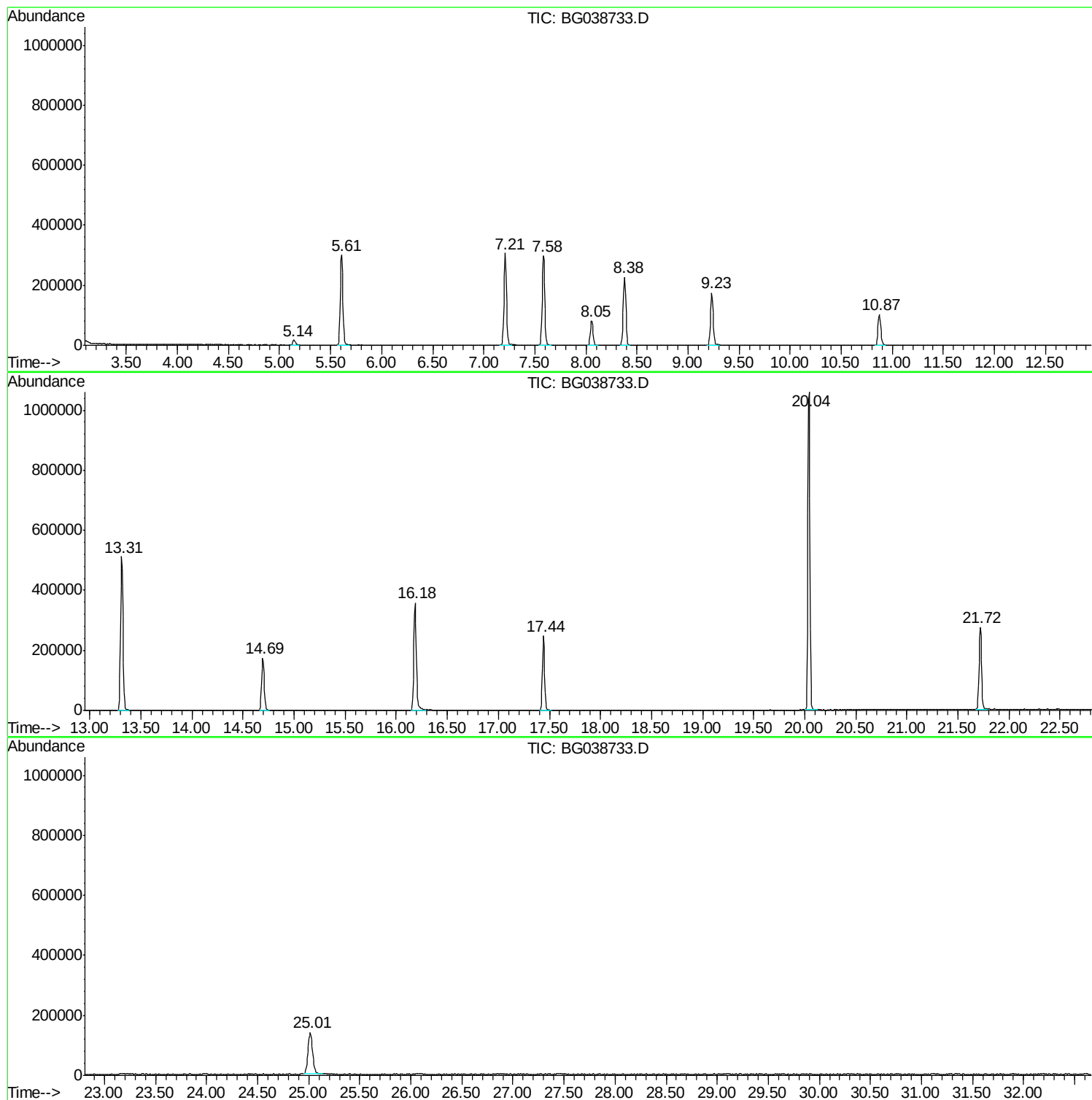
Sum of corrected areas: 7205821

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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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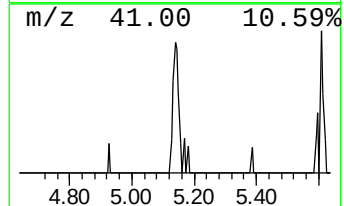
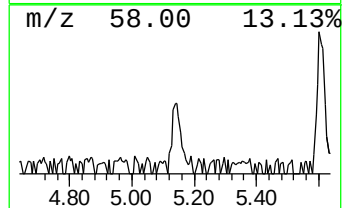
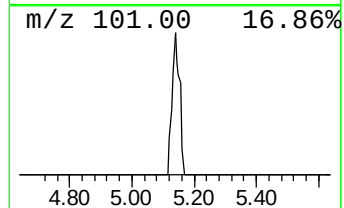
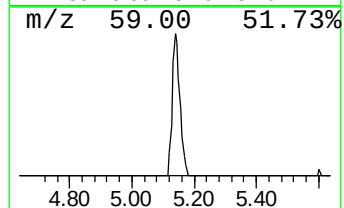
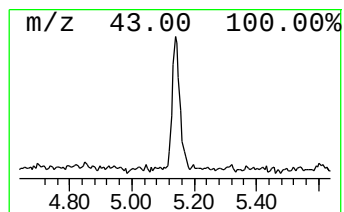
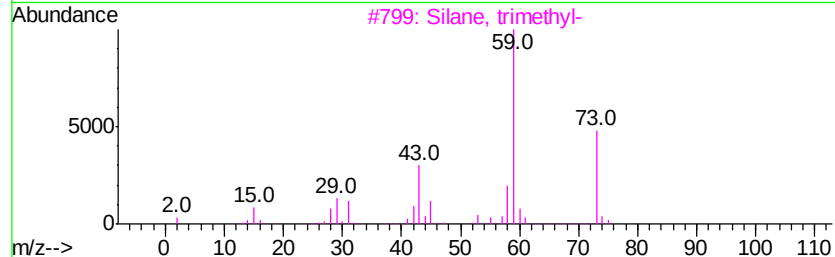
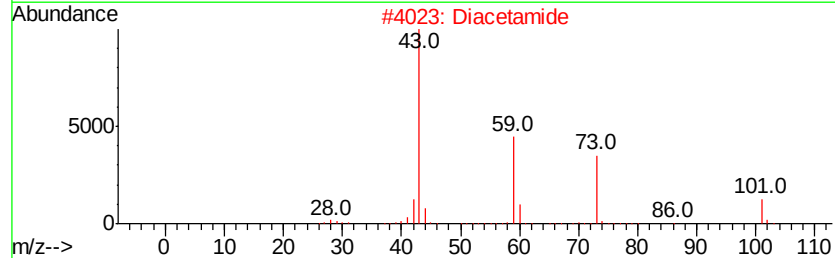
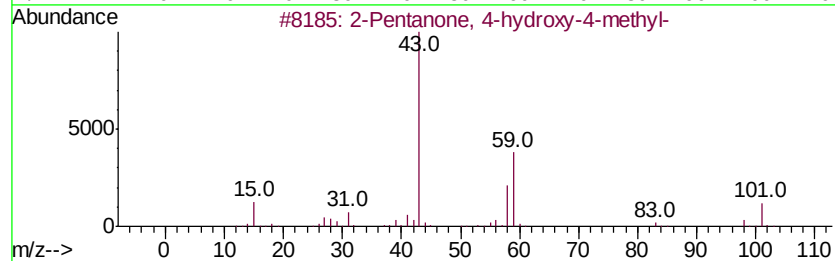
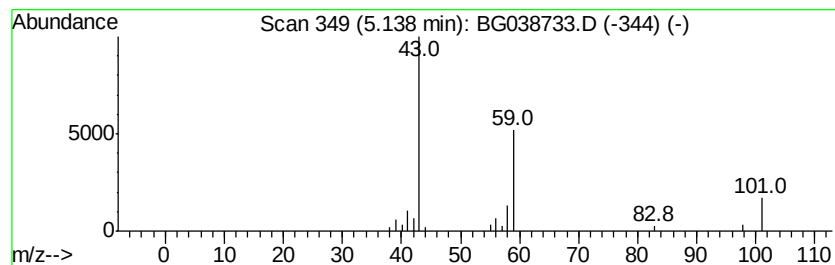
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.14	3.70 ng	26809	1,4-Dichlorobenzene-d4	8.06

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			Diacetamide	101	C4H7NO2	000625-77-4	9
3			Silane, trimethyl-	74	C3H10Si	000993-07-7	9
4			2-Pentanone, 5-(acetyloxy)-	144	C7H12O3	005185-97-7	9
5			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9



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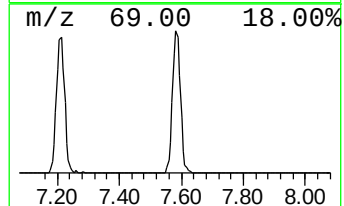
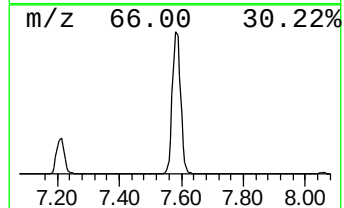
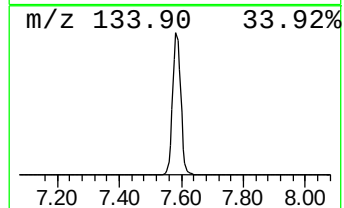
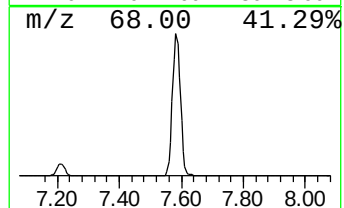
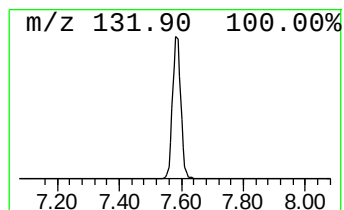
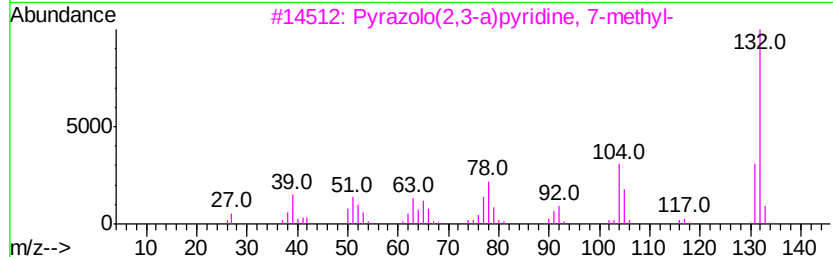
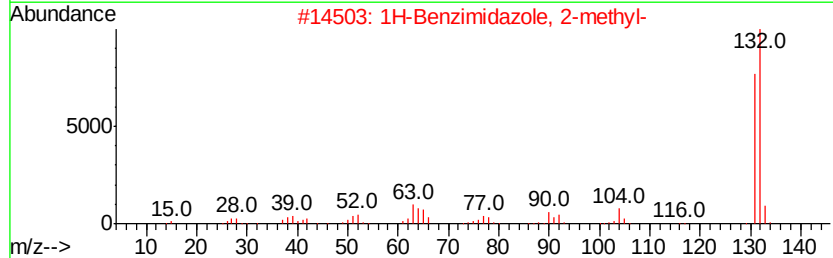
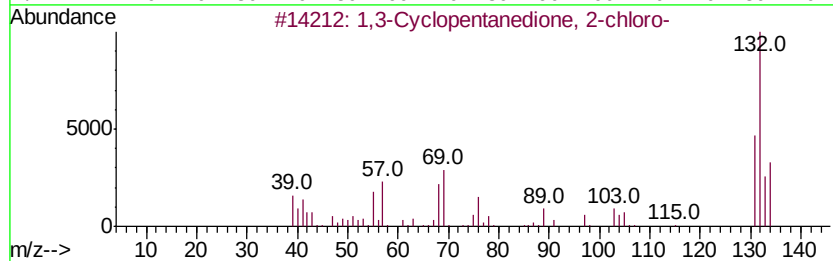
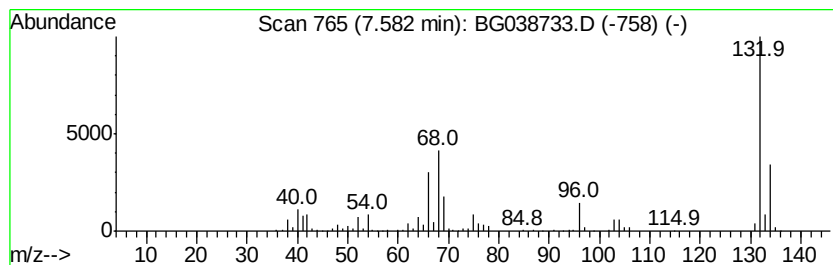
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Peak Number 2 unknown7.58 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	73.23 ng	530590	1,4-Dichlorobenzene-d4	8.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	25
2		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
3		Pyrazolo(2,3-a)pyridine, 7-methyl-	132	C8H8N2	034760-58-2	22
4		5-Aminoindole	132	C8H8N2	005192-03-0	22
5		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.14	3.7	ng	26809	1	8.06	144904	20.0
unknown7.58	7.58	73.2	ng	530590	1	8.06	144904	20.0