

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG111419\
 Data File : BG043631.D
 Acq On : 14 Nov 2019 13:22
 Operator : JU
 Sample : PB124780BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB124780BL

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-BG102119.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	5.099	336	342	355	rBV	110163	172956	5.21%	1.294%
2	5.592	419	426	440	rBV	542011	932945	28.09%	6.982%
3	7.197	690	699	716	rBV	550958	1044331	31.44%	7.815%
4	7.543	750	758	774	rBV	584000	1047509	31.54%	7.839%
5	7.996	828	835	844	rBV	88019	152037	4.58%	1.138%
6	8.319	882	890	904	rBV	452792	807993	24.33%	6.047%
7	9.159	1025	1033	1052	rBV	302815	618221	18.61%	4.626%
8	10.804	1305	1313	1324	rBV	84901	190022	5.72%	1.422%
9	13.248	1722	1729	1742	rBV	986634	1642810	49.46%	12.294%
10	14.629	1958	1964	1979	rBV	184020	320915	9.66%	2.402%
11	16.121	2211	2218	2239	rBV	856207	1522919	45.85%	11.397%
12	17.373	2425	2431	2447	rBV	228730	424370	12.78%	3.176%
13	19.987	2869	2876	2888	rBV	2357651	3321321	100.00%	24.855%
14	21.638	3149	3157	3176	rBV	280022	574074	17.28%	4.296%
15	24.823	3689	3699	3716	rBV2	182052	590202	17.77%	4.417%

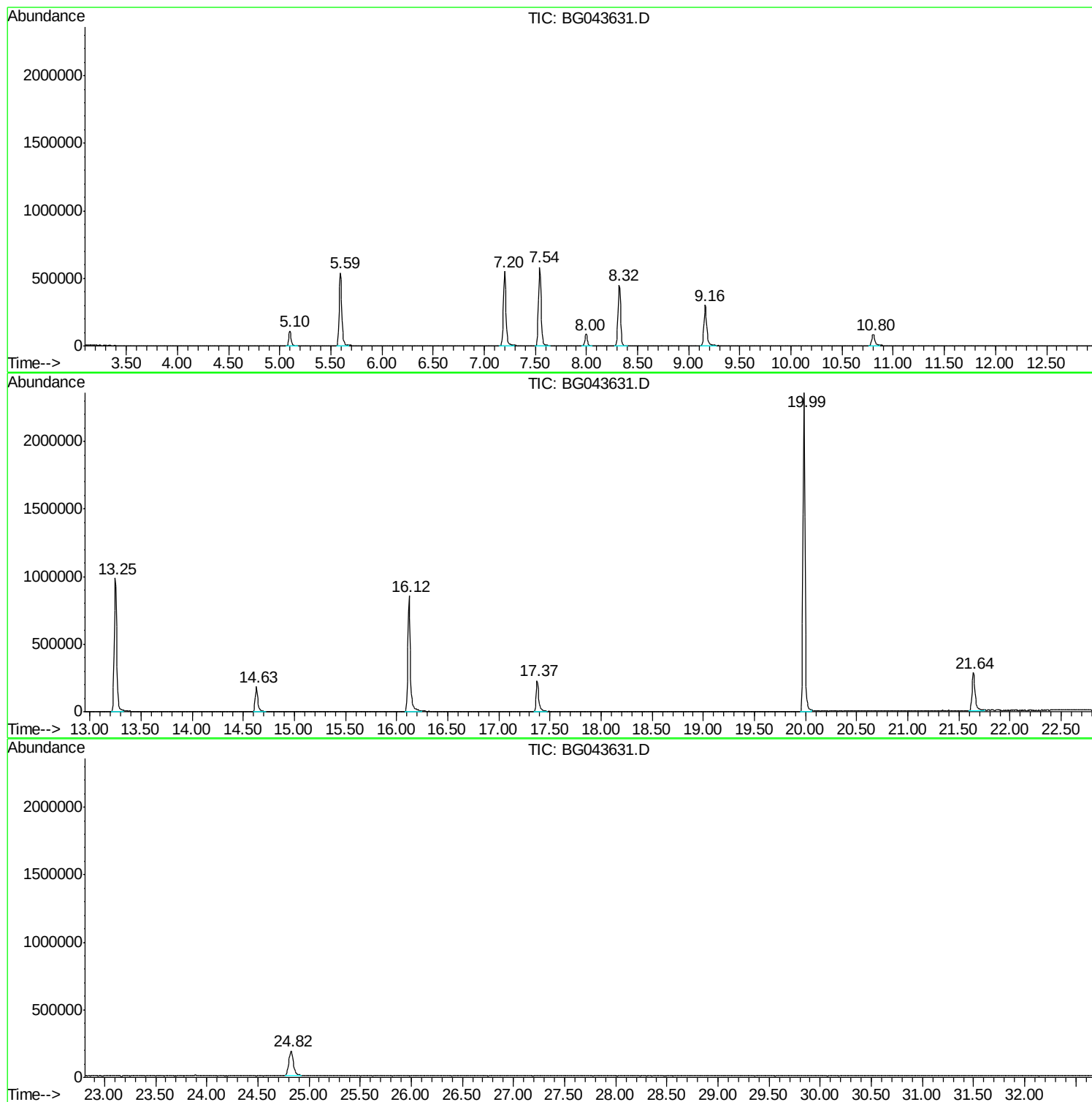
Sum of corrected areas: 13362625

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



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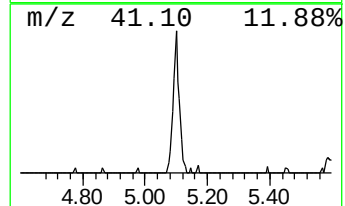
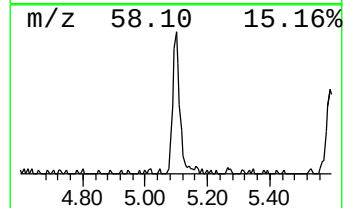
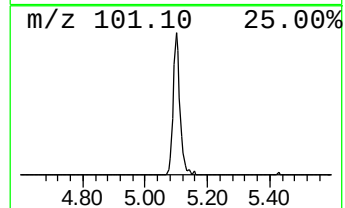
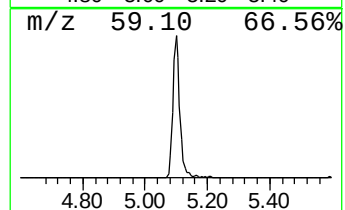
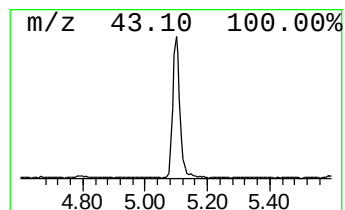
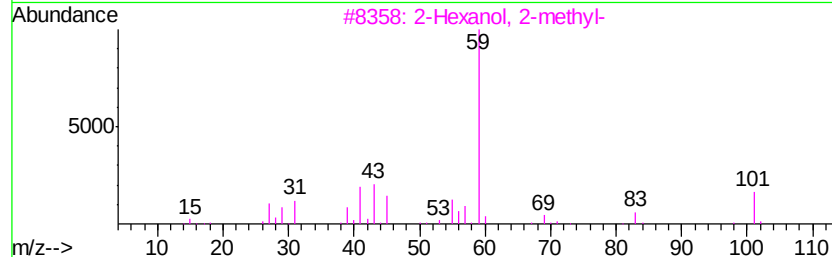
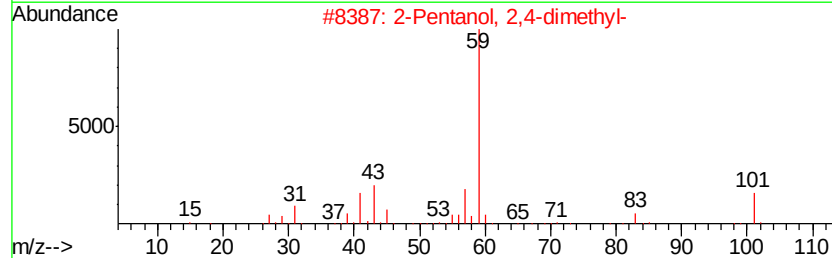
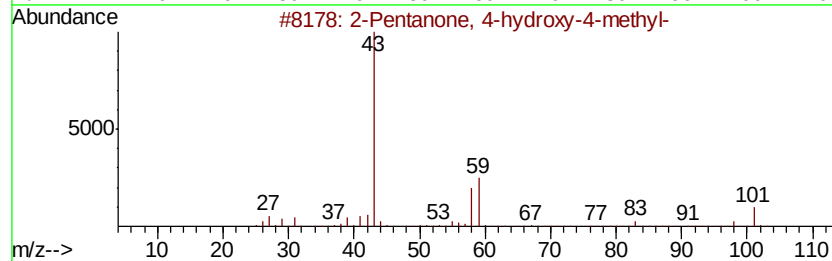
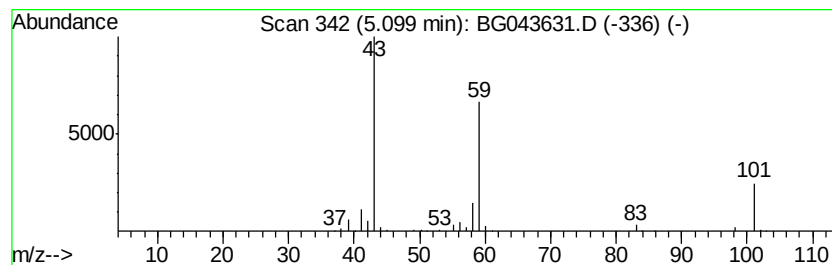
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.10	22.75 ng	172956	1,4-Dichlorobenzene-d4	8.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	42
3		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	40
4		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
5		3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	16



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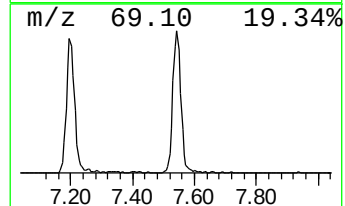
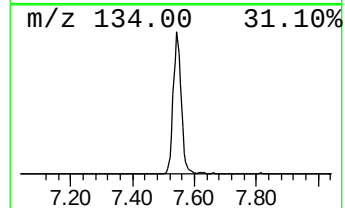
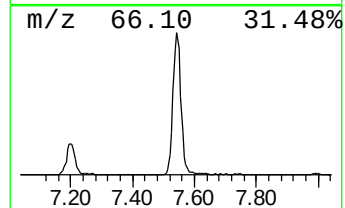
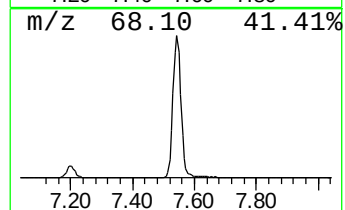
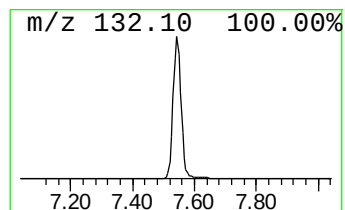
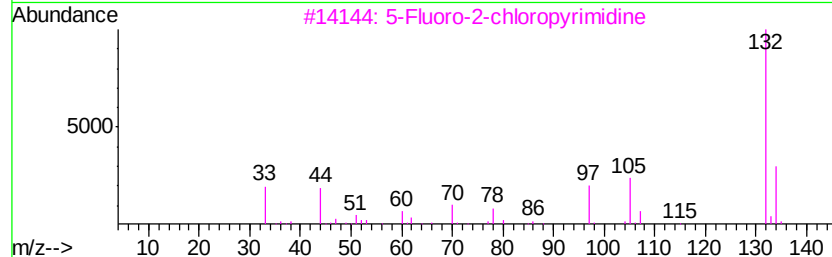
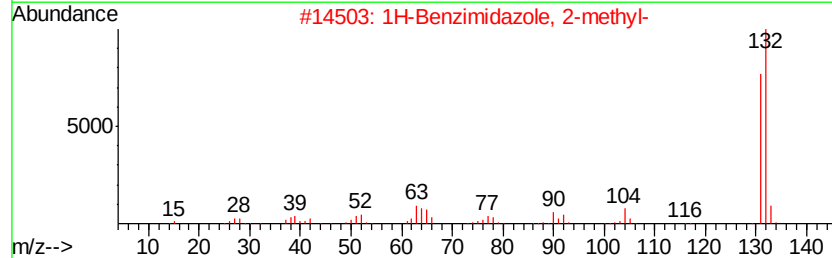
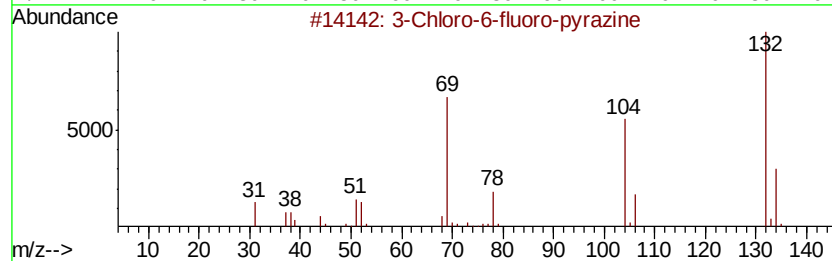
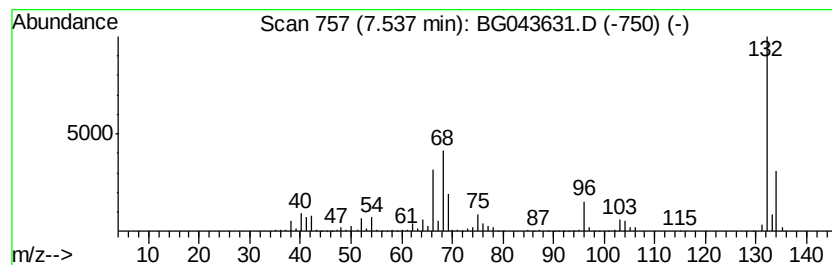
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Peak Number 2 unknown7.54 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.54	137.80 ng	1047510	1,4-Dichlorobenzene-d4	8.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.10	22.8	ng	172956	1	8.00	152037	20.0
unknown7.54	7.54	137.8	ng	1047510	1	8.00	152037	20.0