

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_G\DATA\BG011818\
 Data File : BG032145.D
 Acq On : 18 Jan 2018 20:47
 Operator : SJ/JU
 Sample : J1187-02
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EPE-106-4A(34-36)

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:\HPCHEM1\BNA_G\METHODS\8270-BG011218.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	5.661	396	404	413	rBV	23994	40534	1.32%	0.297%
2	6.166	481	490	501	rBV	422067	730897	23.89%	5.349%
3	7.847	768	776	787	rBV	446833	833050	27.23%	6.097%
4	8.252	837	845	861	rBB	463091	842482	27.53%	6.166%
5	8.740	921	928	939	rBV2	103694	188005	6.14%	1.376%
6	9.075	978	985	1000	rVB	334137	629198	20.56%	4.605%
7	9.938	1123	1132	1144	rBV	261677	489345	15.99%	3.581%
8	11.624	1411	1419	1427	rBV	146350	277410	9.07%	2.030%
9	14.004	1817	1824	1835	rVB	938890	1483699	48.49%	10.859%
10	14.803	1953	1960	1966	rBV	96401	145668	4.76%	1.066%
11	15.379	2051	2058	2067	rBV2	287740	478794	15.65%	3.504%
12	16.854	2302	2309	2318	rBV	1030935	1588989	51.93%	11.630%
13	18.111	2517	2523	2531	rBV2	469491	733564	23.97%	5.369%
14	20.302	2892	2896	2904	rBV2	29297	40463	1.32%	0.296%
15	20.649	2949	2955	2963	rBV	2304025	3059788	100.00%	22.394%
16	21.989	3178	3183	3191	rVB9	20355	34632	1.13%	0.253%
17	22.447	3253	3261	3271	rVB	560430	1048919	34.28%	7.677%
18	26.149	3880	3891	3905	rBV2	317405	1017875	33.27%	7.450%

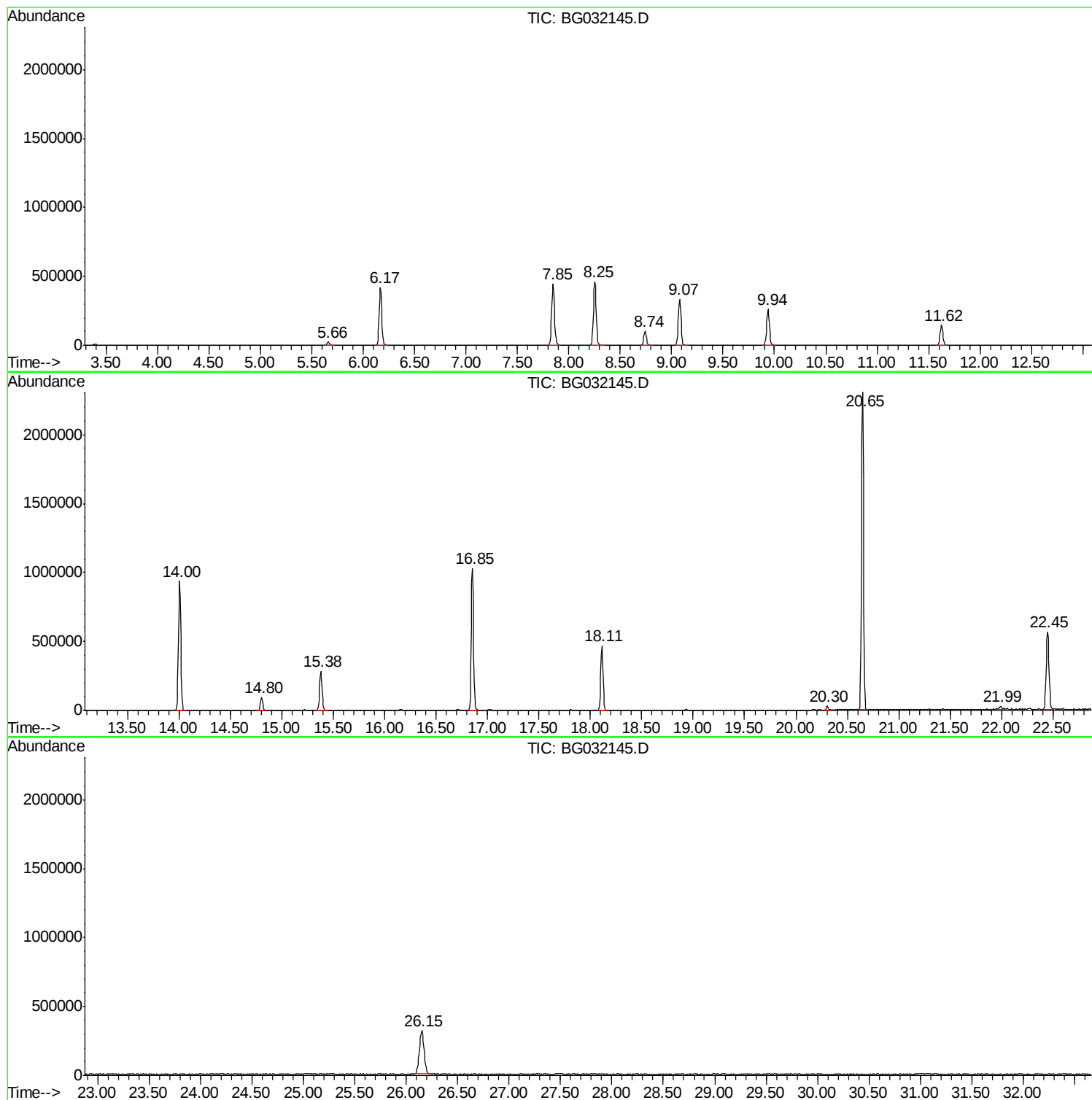
Sum of corrected areas: 13663312

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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG011218.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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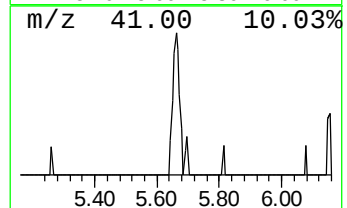
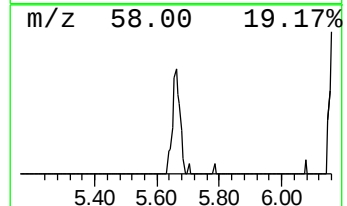
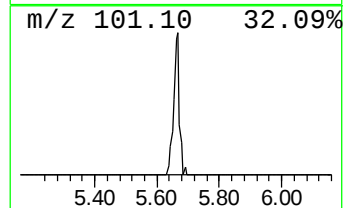
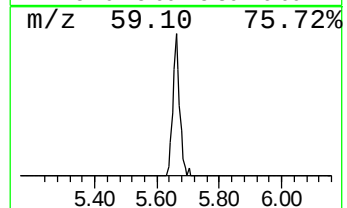
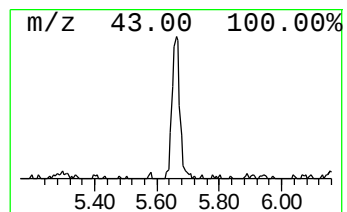
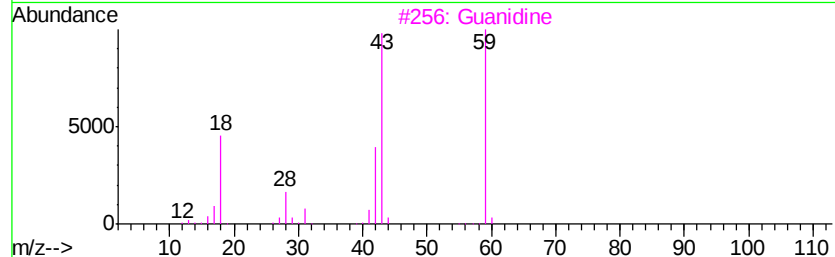
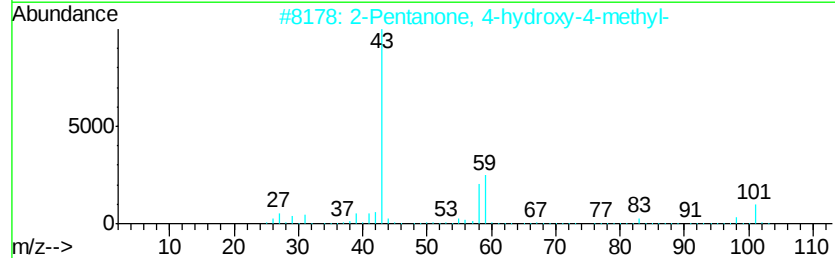
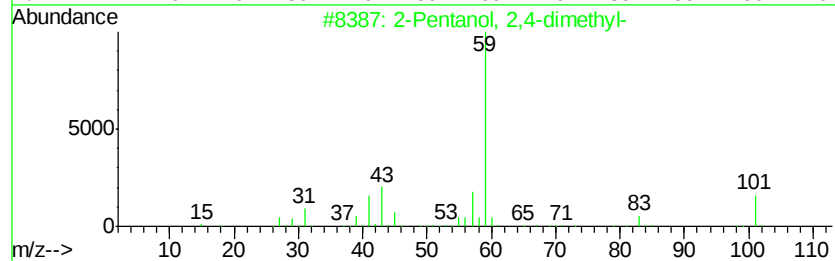
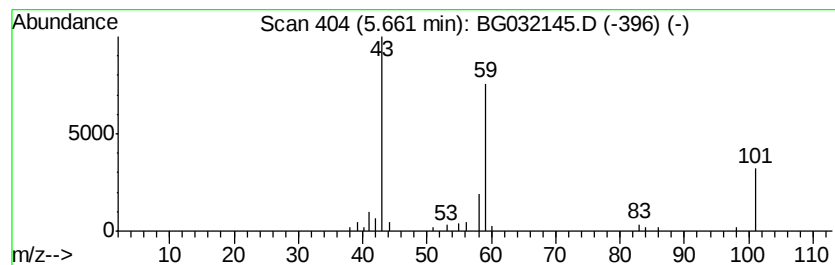
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TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanol, 2,4-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.66	4.31 ng	40534	1,4-Dichlorobenzene-d4	8.75

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	59
2			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
3			Guanidine	59	CH5N3	000113-00-8	50
4			3-Heptanol, 4-methyl-	130	C8H18O	014979-39-6	36
5			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33



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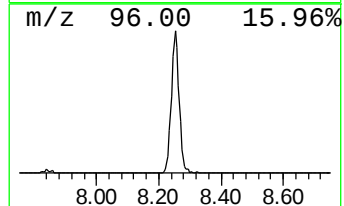
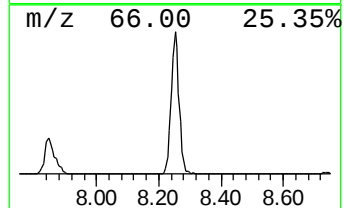
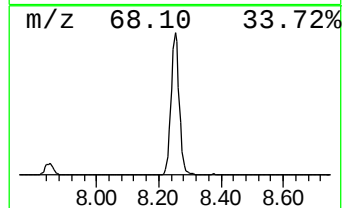
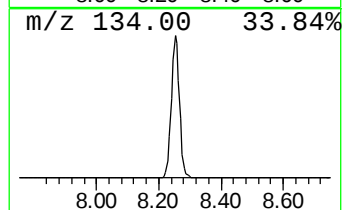
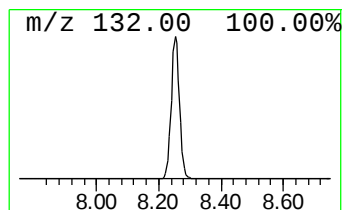
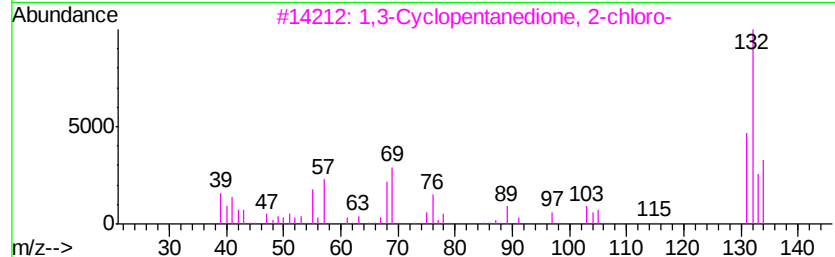
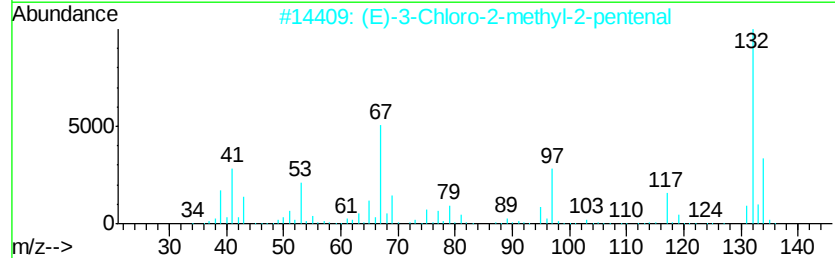
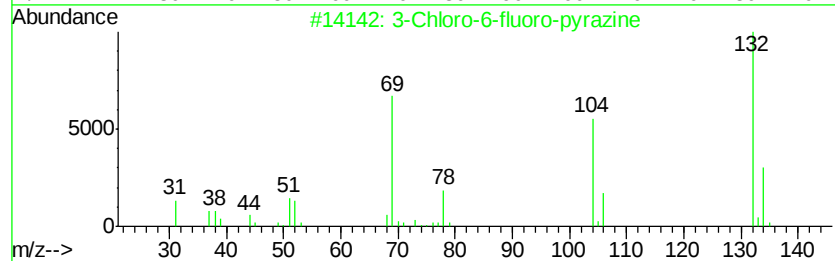
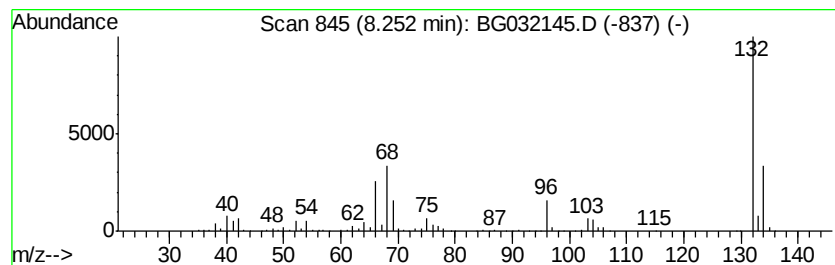
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Peak Number 2 unknown8.25 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.25	89.62 ng	842482	1,4-Dichlorobenzene-d4	8.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
3		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	23
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
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-Pentanol, 2,4-d...	5.66	4.3	ng	40534	1	8.75	188005	20.0
unknown8.25	8.25	89.6	ng	842482	1	8.75	188005	20.0