

Data Path : Z:\HPCHEM1\BNA G\DATA\BG122717\
 Data File : BG031803.D
 Acq On : 27 Dec 2017 19:14
 Operator : SJ/JU
 Sample : I7058-04
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 122117-TIDO-F-U-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:\HPCHEM1\BNA_G\METHODS\8270-BG122117.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	3.437	19	25	35	rBV2	71394	152827	4.07%	1.098%
2	3.525	35	40	50	rVB	38580	61896	1.65%	0.445%
3	6.222	491	499	510	rBV	155792	258789	6.89%	1.859%
4	7.902	777	785	799	rVB	97423	187445	4.99%	1.346%
5	8.314	846	855	865	rBV	303506	558488	14.86%	4.011%
6	8.807	931	939	948	rBB	121556	218209	5.81%	1.567%
7	9.142	987	996	1009	rBB	463151	872874	23.23%	6.269%
8	10.000	1134	1142	1156	rBV	323321	613072	16.31%	4.403%
9	11.692	1421	1430	1439	rBV	171649	320805	8.54%	2.304%
10	14.066	1826	1834	1844	rBV	1369546	2198321	58.49%	15.787%
11	15.435	2061	2067	2075	rVB2	330795	557475	14.83%	4.004%
12	16.904	2310	2317	2329	rBV	765767	1215054	32.33%	8.726%
13	18.167	2525	2532	2540	rBV	515782	803723	21.39%	5.772%
14	20.699	2956	2963	2970	rBV	2803672	3758171	100.00%	26.989%
15	22.056	3188	3194	3200	rBV6	25960	46884	1.25%	0.337%
16	22.509	3264	3271	3279	rBV2	543142	1027514	27.34%	7.379%
17	26.263	3896	3910	3926	rBV2	320956	1073034	28.55%	7.706%

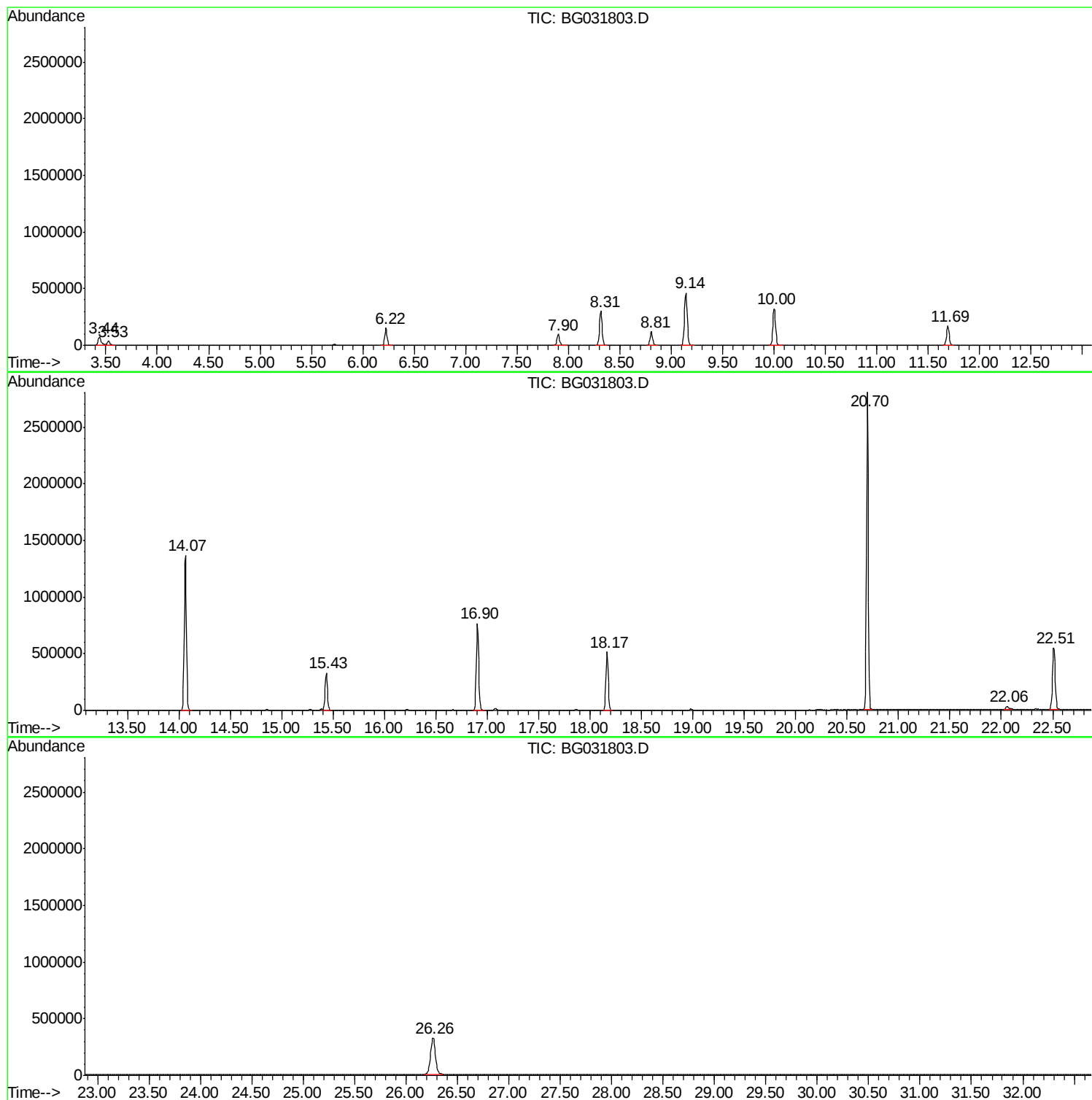
Sum of corrected areas: 13924581

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG122117.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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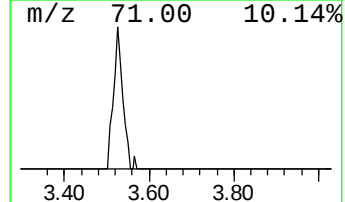
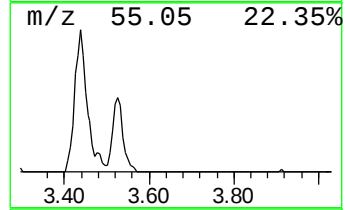
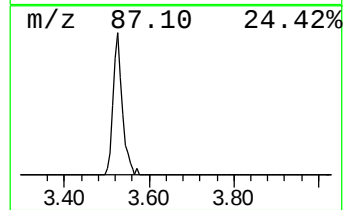
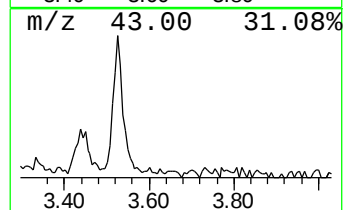
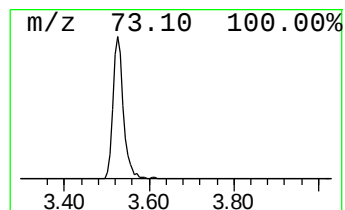
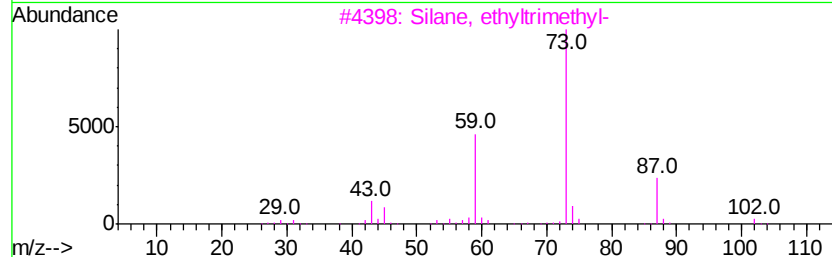
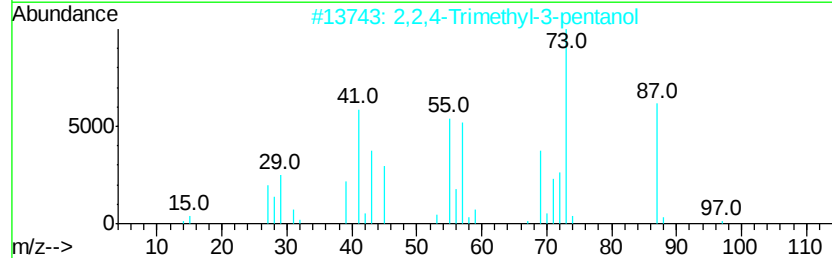
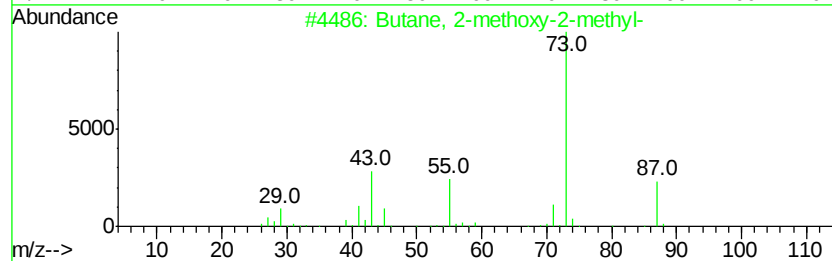
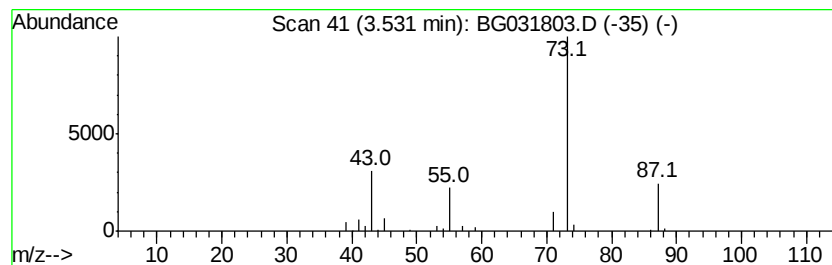
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG122117.M
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Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.53	5.67 ng	61896	1,4-Dichlorobenzene-d4	8.81

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	43
3			Silane, ethyltrimethyl-	102	C5H14Si	003439-38-1	9
4			3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	9
5			3-Hexanol, 2,4-dimethyl-	130	C8H18O	013432-25-2	9



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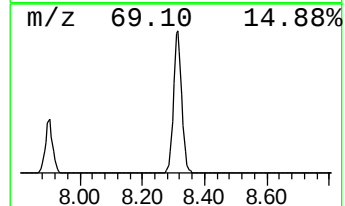
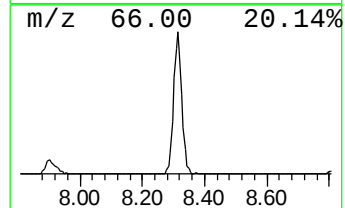
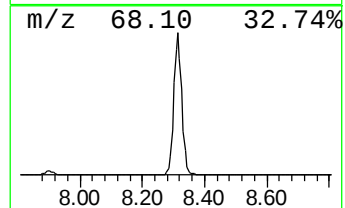
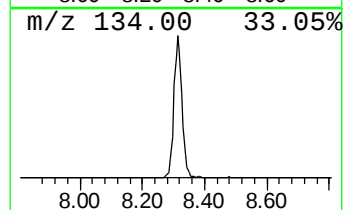
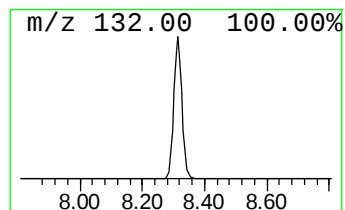
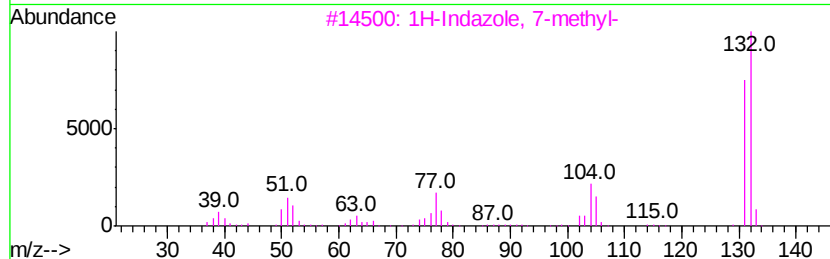
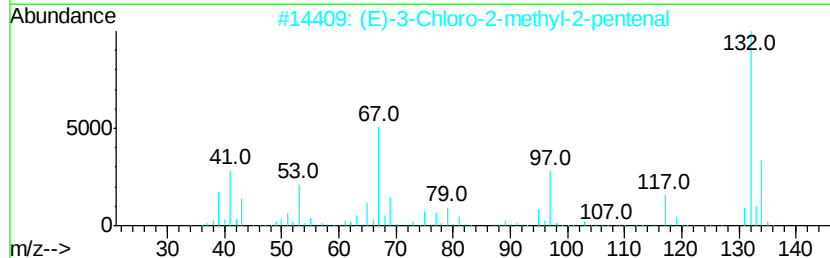
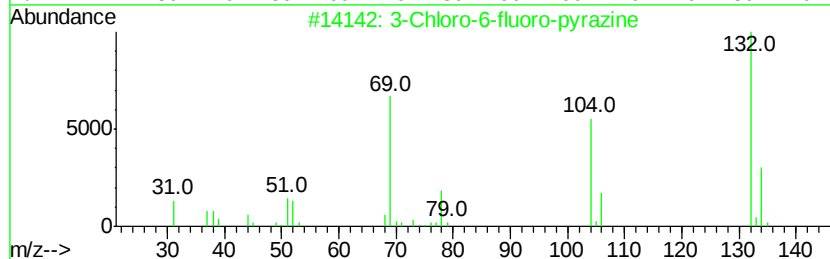
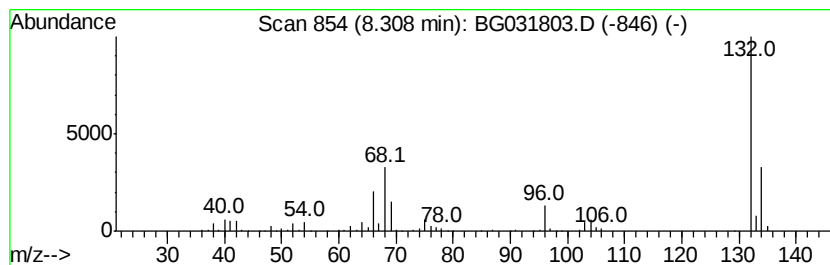
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Peak Number 3 unknown8.31 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.31	51.19 ng	558488	1,4-Dichlorobenzene-d4	8.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	35
3		1H-Indazole, 7-methyl-	132	C8H8N2	1000316-00-7	10
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	3.53	5.7	ng	61896	1	8.81	218209	20.0
unknown8.31	8.31	51.2	ng	558488	1	8.81	218209	20.0