

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG032018\
 Data File : BG033289.D
 Acq On : 21 Mar 2018 5:50
 Operator : SJ/JU
 Sample : PB107504BL
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB107504BL

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-BG031918.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	5.796	420	427	433	rBV	32958	57366	1.41%	0.331%
2	6.307	507	514	535	rBV	649971	1320450	32.44%	7.625%
3	7.994	795	801	827	rBV	660398	1479270	36.34%	8.543%
4	8.399	864	870	895	rBV	616547	1359866	33.41%	7.853%
5	8.893	947	954	966	rBV2	86691	211196	5.19%	1.220%
6	9.228	1001	1011	1030	rBV	471622	1004341	24.67%	5.800%
7	10.091	1151	1158	1187	rBV	358475	870205	21.38%	5.025%
8	11.777	1438	1445	1461	rBV2	135330	315230	7.74%	1.820%
9	14.139	1840	1847	1863	rBV	1235466	2211518	54.33%	12.771%
10	15.508	2072	2080	2092	rBV3	293051	515626	12.67%	2.978%
11	16.983	2324	2331	2352	rBV	757371	1424522	35.00%	8.226%
12	18.240	2539	2545	2559	rBV	440309	747193	18.36%	4.315%
13	20.761	2968	2974	2987	rBV	3065825	4070340	100.00%	23.506%
14	22.594	3279	3286	3297	rBV2	438855	884376	21.73%	5.107%
15	26.396	3921	3933	3950	rBV2	239570	844957	20.76%	4.880%

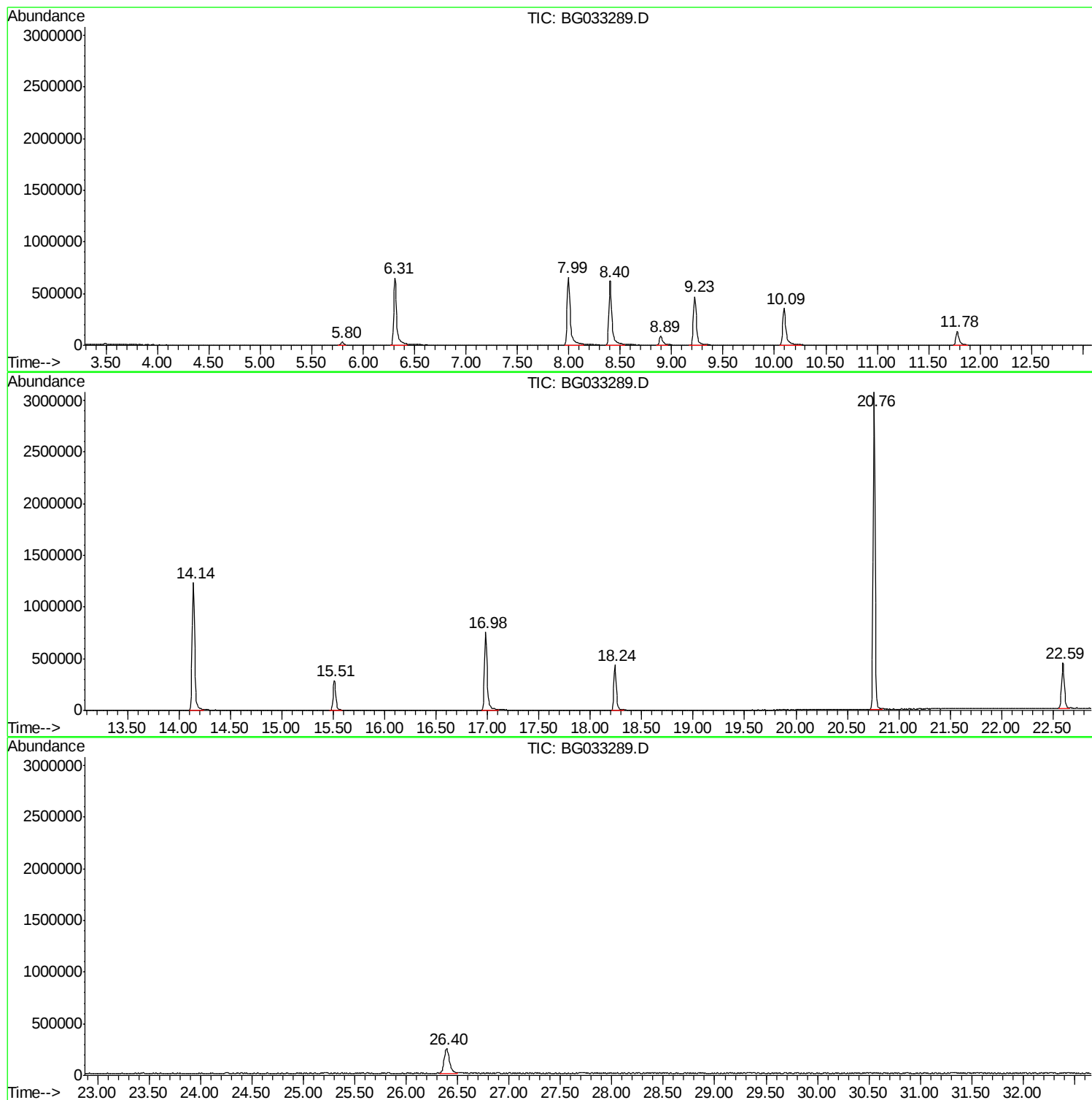
Sum of corrected areas: 17316456

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG031918.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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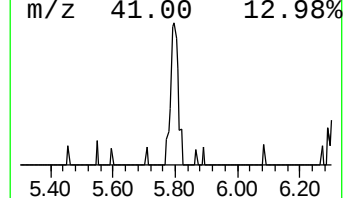
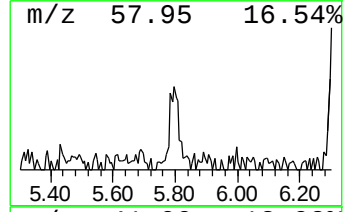
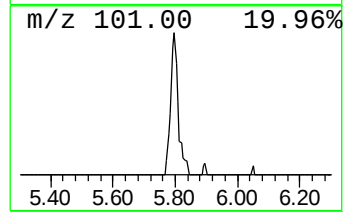
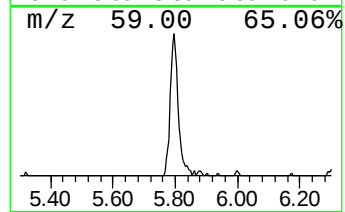
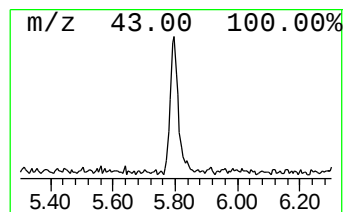
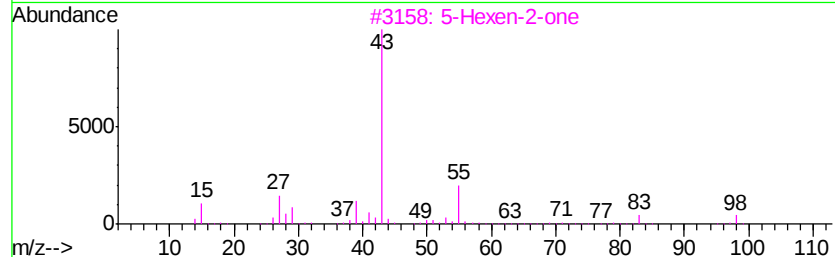
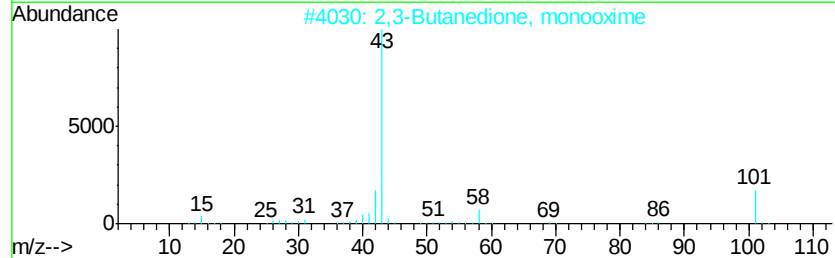
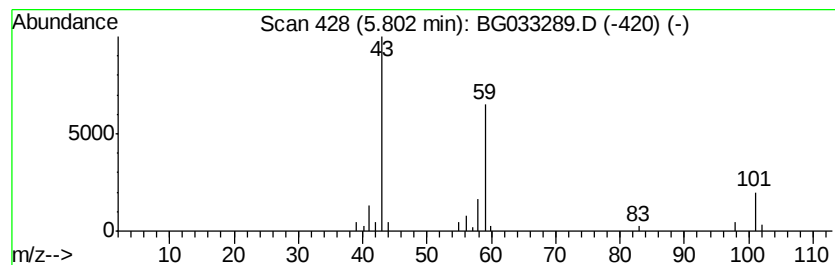
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.80	5.43 ng	57366	1,4-Dichlorobenzene-d4	8.89

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
3			5-Hexen-2-one	98	C6H10O	000109-49-9	9
4			1,3,4-Oxadiazole, 2-(acetyloxy)-...	172	C7H12N2O3	066398-57-0	9
5			3-Pentanol	88	C5H12O	000584-02-1	9



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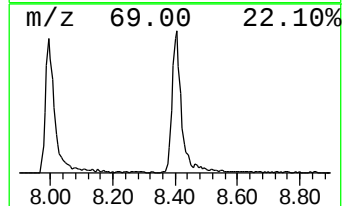
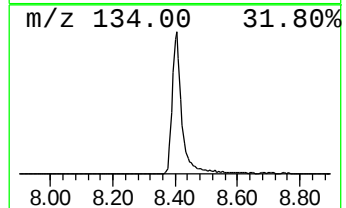
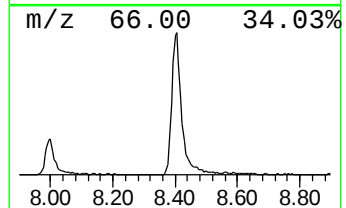
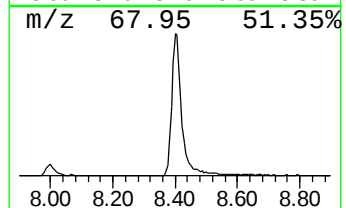
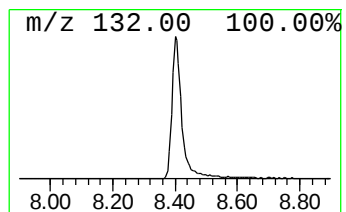
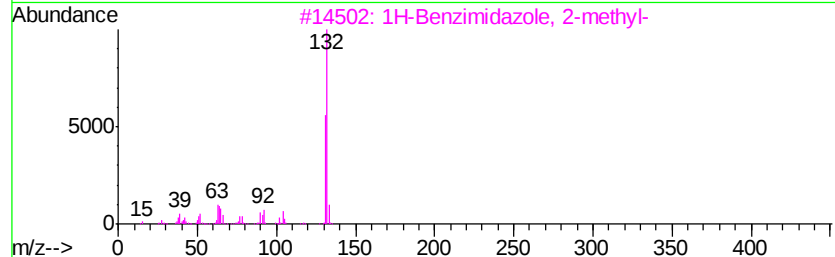
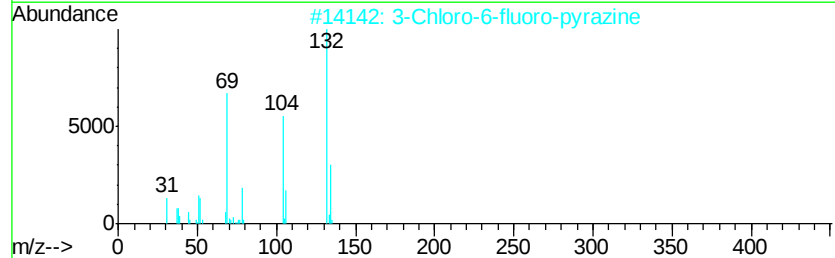
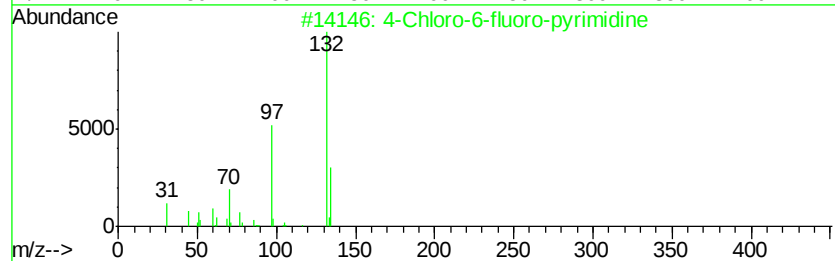
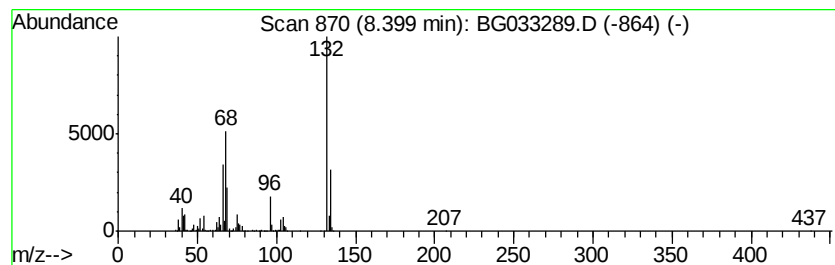
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Peak Number 2 unknown8.40 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.40	128.78 ng	1359870	1,4-Dichlorobenzene-d4	8.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4		Imidazole, 4,5-dicvno-1-methyl-	132	C6H4N4	019485-35-9	12
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.80	5.4	ng	57366	1	8.89	211196	20.0
unknown8.40	8.40	128.8	ng	1359870	1	8.89	211196	20.0