

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_G\DATA\BG010318\
 Data File : BG031887.D
 Acq On : 3 Jan 2018 12:51
 Operator : SJ/JU
 Sample : I7106-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EPE-106-2(25-27)

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	3.423	17	23	34	rBV2	38146	76244	2.19%	0.478%
2	5.702	402	411	422	rBV	25557	47190	1.36%	0.296%
3	6.207	490	497	510	rVB	497754	865583	24.90%	5.428%
4	7.888	774	783	794	rBV	519917	984617	28.33%	6.174%
5	8.299	844	853	870	rBV	526539	979328	28.17%	6.141%
6	8.787	928	936	945	rBB2	127151	230108	6.62%	1.443%
7	9.122	984	993	1004	rBB	421543	779662	22.43%	4.889%
8	9.980	1130	1139	1149	rBV	290626	557592	16.04%	3.496%
9	11.666	1417	1426	1434	rBV	184398	350122	10.07%	2.195%
10	14.045	1823	1831	1840	rBV	1141674	1848153	53.17%	11.589%
11	14.839	1960	1966	1973	rBV	69626	104601	3.01%	0.656%
12	15.414	2054	2064	2073	rVB2	356854	596323	17.15%	3.739%
13	16.889	2308	2315	2327	rBV	1122193	1782902	51.29%	11.180%
14	18.146	2521	2529	2539	rBV2	565881	896761	25.80%	5.623%
15	20.679	2954	2960	2969	rBV	2634878	3476101	100.00%	21.797%
16	22.312	3233	3238	3248	rVB2	57730	93663	2.69%	0.587%
17	22.488	3260	3268	3277	rBV	589884	1117159	32.14%	7.005%
18	23.793	3483	3490	3497	rBV4	20588	43489	1.25%	0.273%
19	26.213	3891	3902	3917	rBV3	329273	1117740	32.15%	7.009%

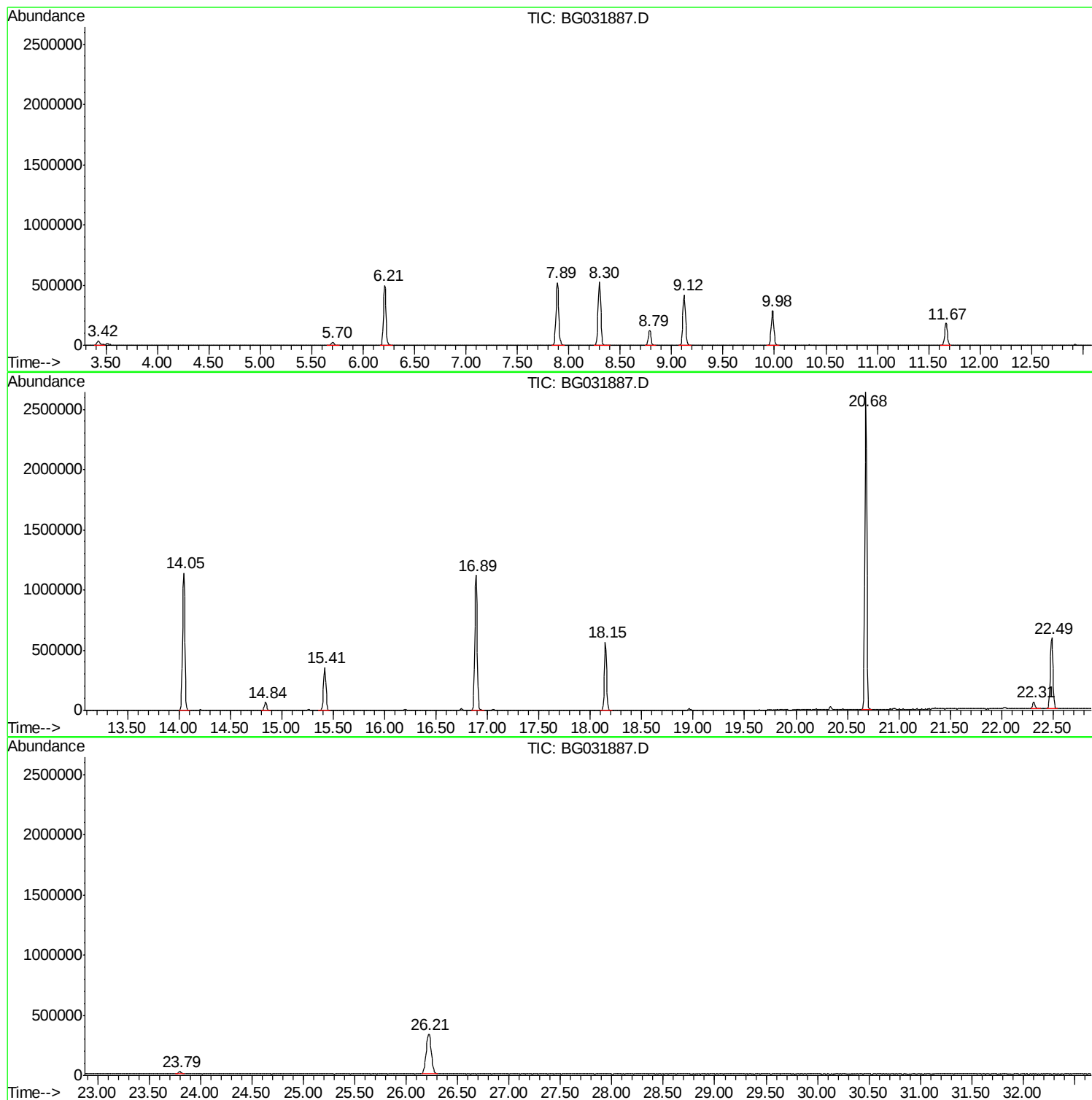
Sum of corrected areas: 15947338

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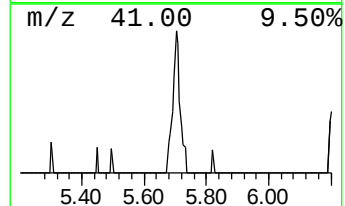
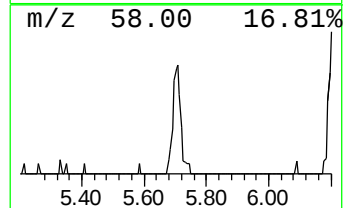
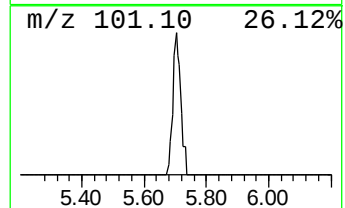
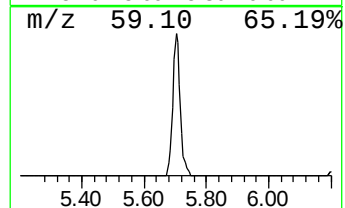
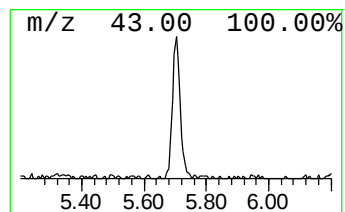
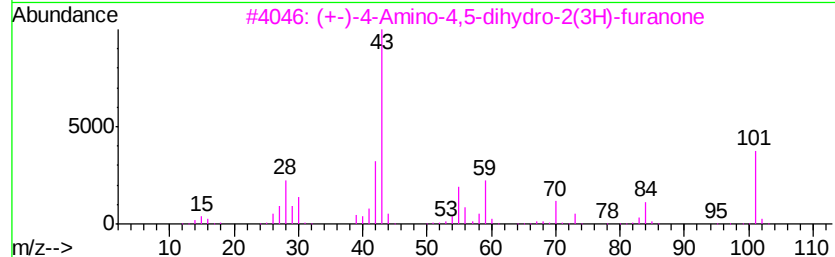
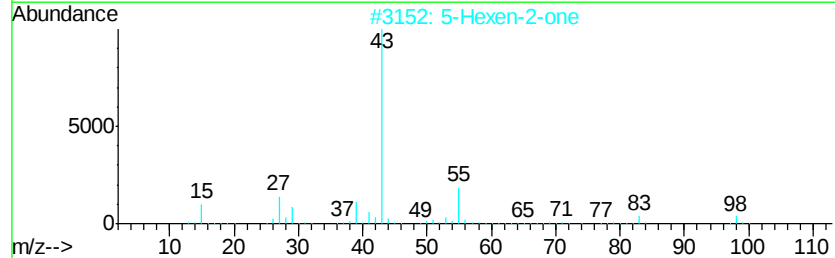
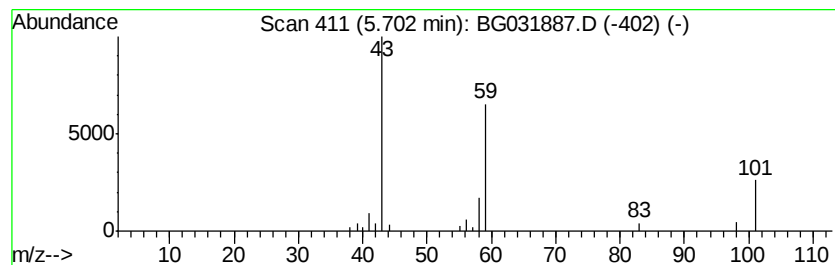
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.70	4.10 ng	47190	1,4-Dichlorobenzene-d4	8.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40
2		5-Hexen-2-one	98	C6H10O	000109-49-9	10
3		(+)-4-Amino-4,5-dihydro-2(3H)-furanone	101	C4H7NO2	016504-58-8	9
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5		Acetone	58	C3H6O	000067-64-1	9



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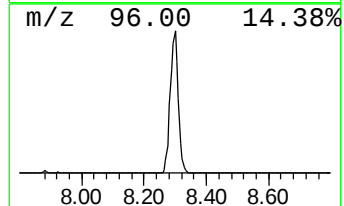
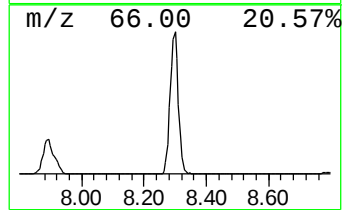
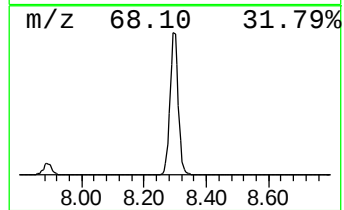
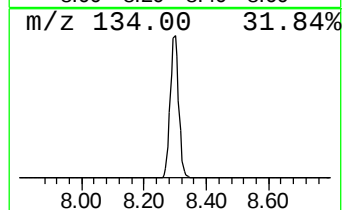
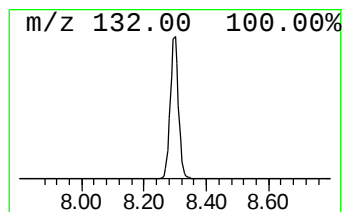
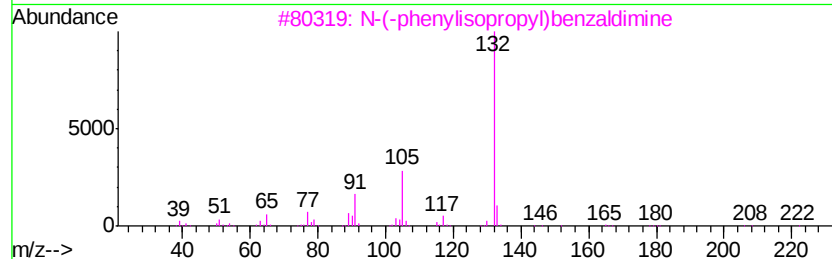
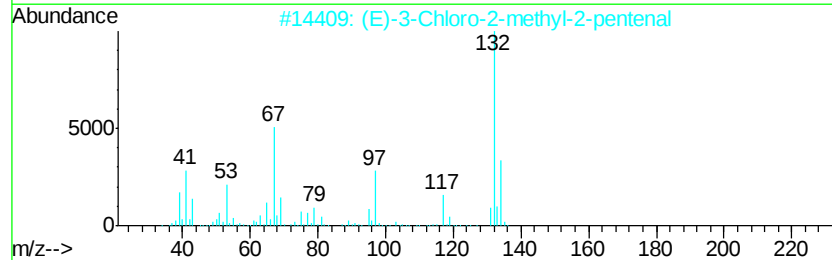
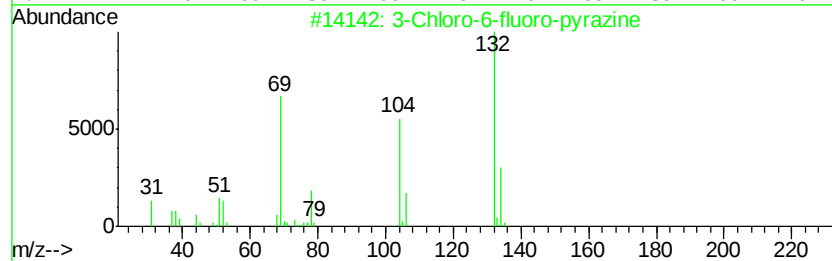
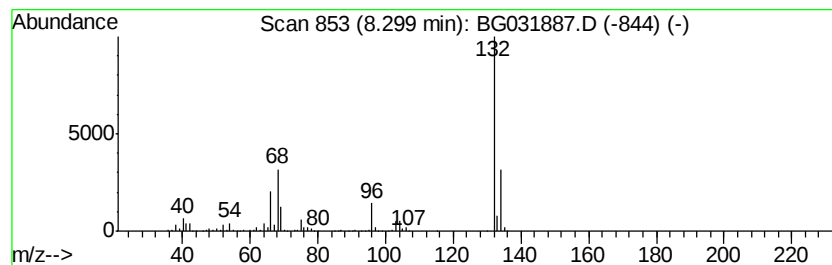
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Peak Number 3 unknown8.30 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.30	85.12 ng	979328	1,4-Dichlorobenzene-d4	8.79

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	25
3		N-(-phenylisopropyl)benzalimine	223	C16H17N	1000379-01-3	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.70	4.1	ng	47190	1	8.79	230108	20.0
unknown8.30	8.30	85.1	ng	979328	1	8.79	230108	20.0