

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG012020\
 Data File : BG044125.D
 Acq On : 21 Jan 2020 9:32
 Operator : JU
 Sample : PB126194BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB126194BL

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-BG123019.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	5.032	325	331	344	rVB	299191	458224	8.26%	1.575%
2	5.508	405	412	427	rBV	1336435	2189849	39.50%	7.527%
3	7.101	674	683	696	rBV	1429339	2506574	45.21%	8.616%
4	7.465	737	745	756	rBV	1401542	2461772	44.40%	8.462%
5	7.929	817	824	833	rVB	251919	428321	7.73%	1.472%
6	8.252	871	879	888	rBV	1065323	1836530	33.12%	6.313%
7	9.086	1011	1021	1043	rBV	823114	1502498	27.10%	5.164%
8	10.732	1292	1301	1318	rBV	334079	647049	11.67%	2.224%
9	13.188	1712	1719	1736	rBV	2219113	3773597	68.06%	12.971%
10	14.568	1946	1954	1968	rBV2	601974	1014342	18.29%	3.487%
11	16.055	2200	2207	2227	rBV	1786700	3000061	54.11%	10.312%
12	17.312	2413	2421	2434	rBV	720457	1179507	21.27%	4.054%
13	19.927	2859	2866	2879	rBV	4062670	5544455	100.00%	19.057%
14	21.554	3137	3143	3155	rBV	726699	1215180	21.92%	4.177%
15	23.810	3522	3527	3534	rVB6	29160	59875	1.08%	0.206%
16	24.662	3662	3672	3690	rBV2	431548	1275537	23.01%	4.384%

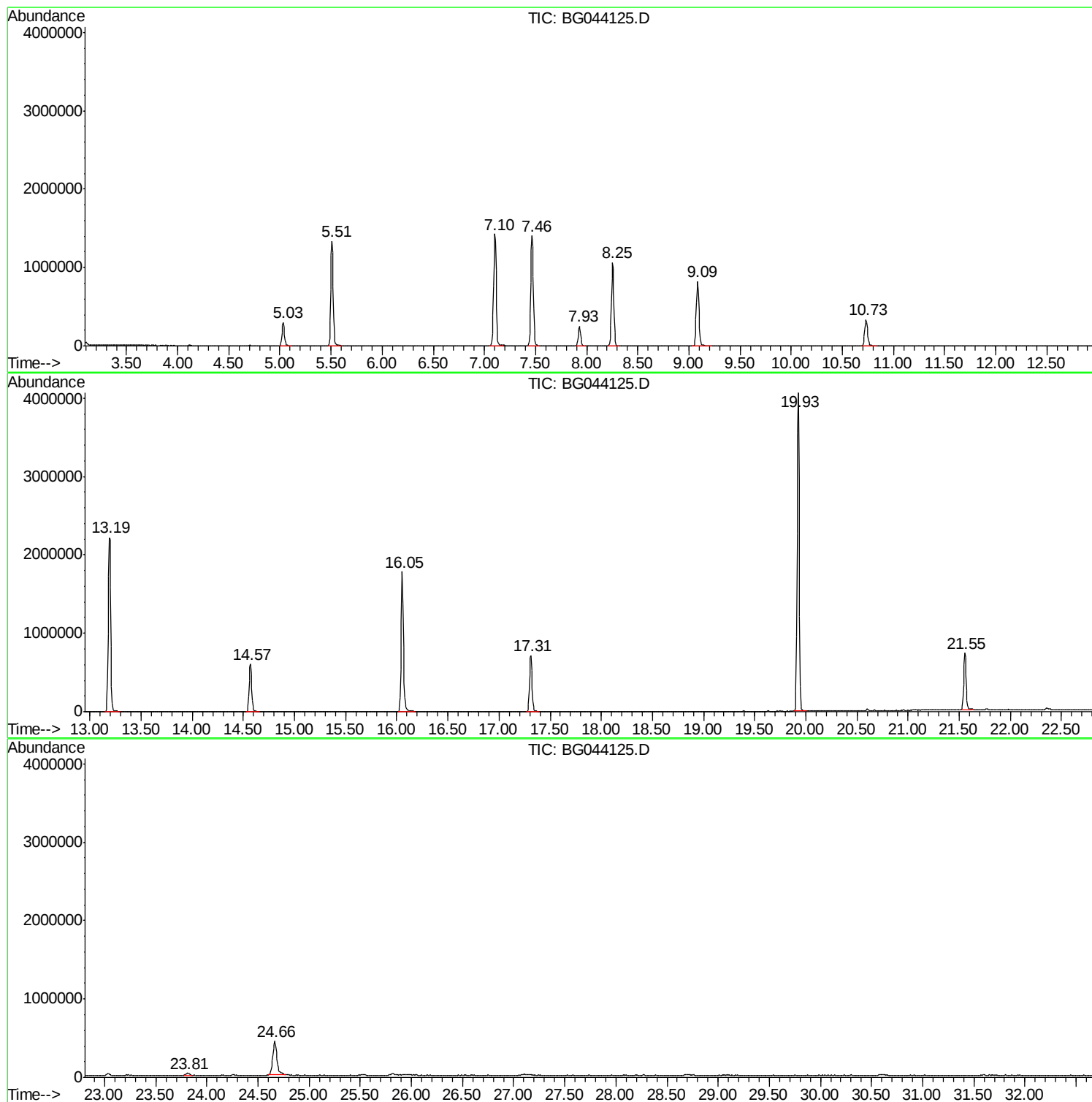
Sum of corrected areas: 29093371

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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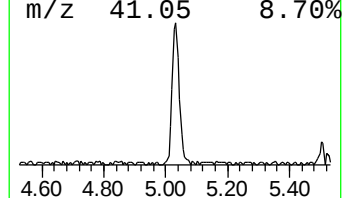
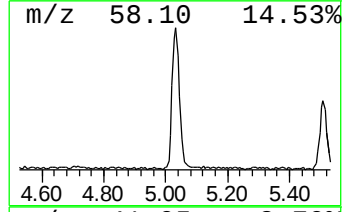
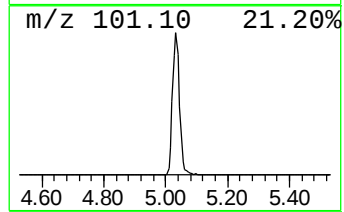
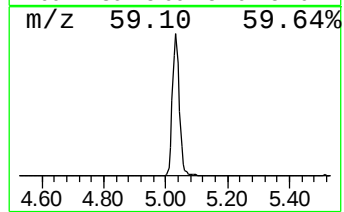
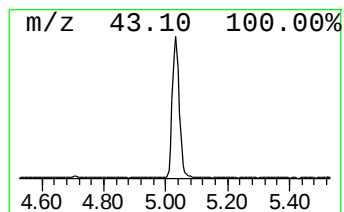
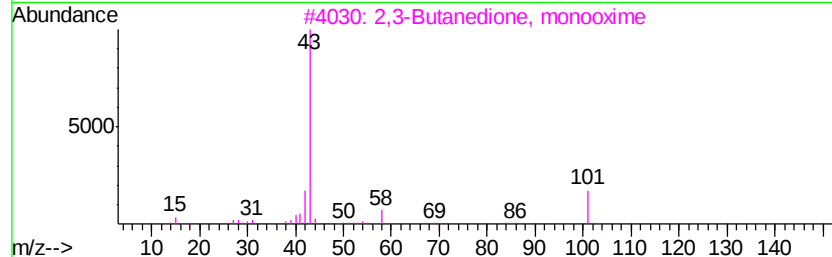
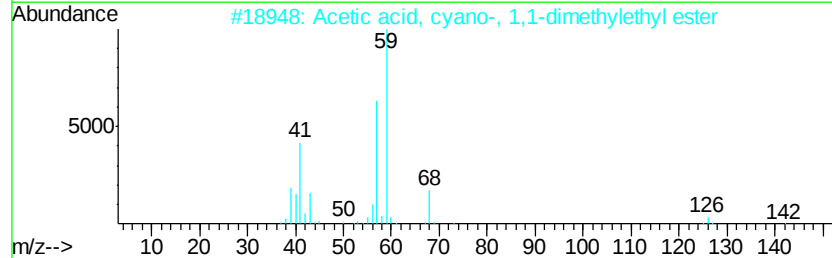
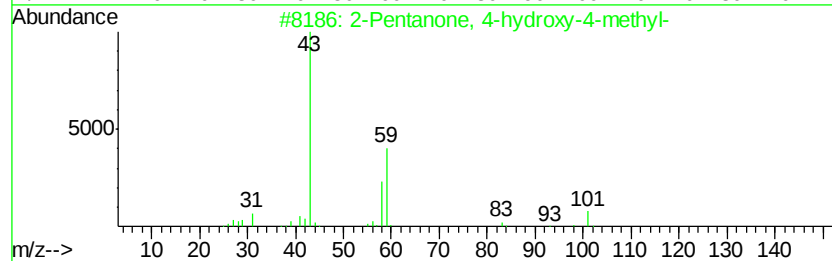
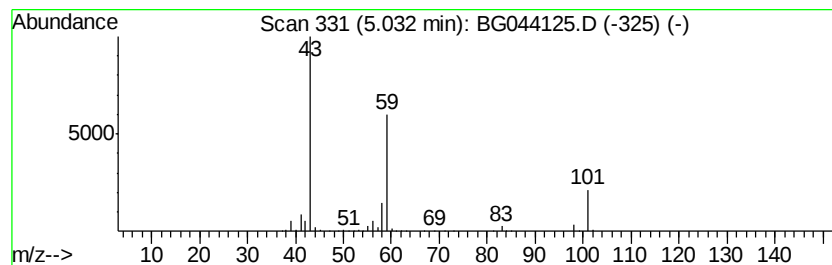
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.03	21.40 ng	458224	1,4-Dichlorobenzene-d4	7.93

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40
2			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4			2-Propanone, 1-(1-methylethoxy)-	116	C6H12O2	042781-12-4	9
5			1,2-Butanediol	90	C4H10O2	000584-03-2	9



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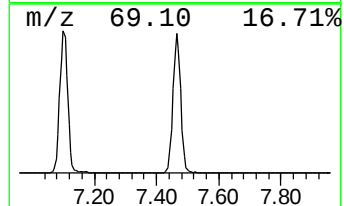
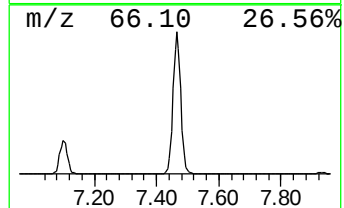
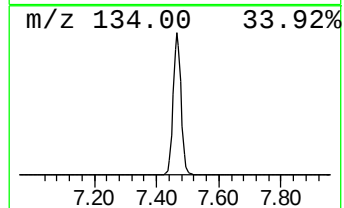
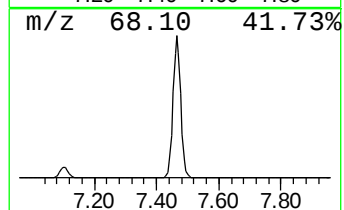
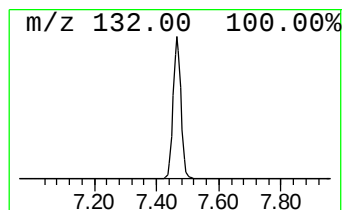
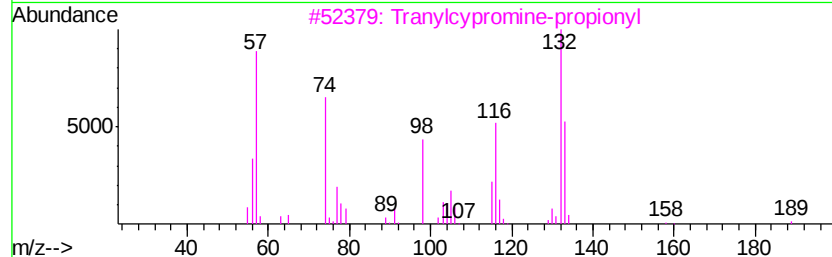
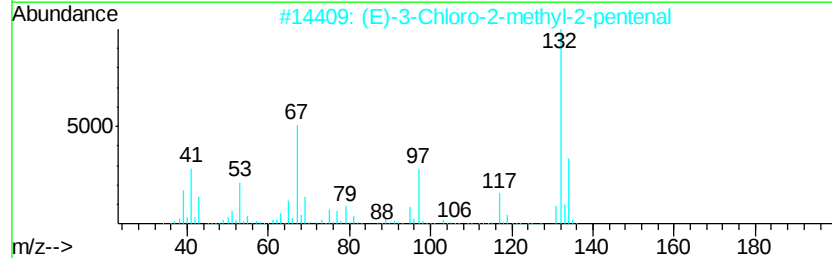
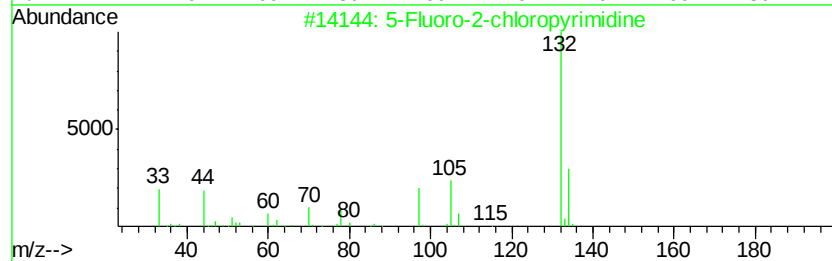
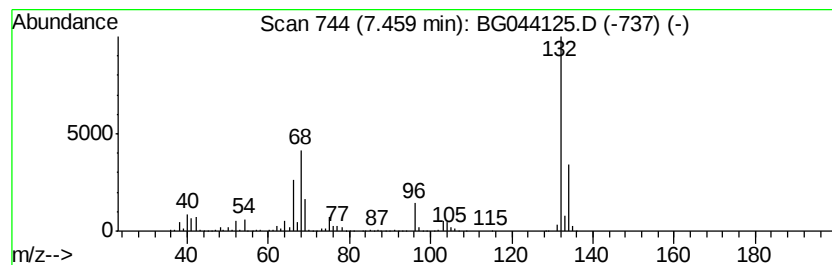
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Peak Number 2 unknown7.46 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.46	114.95 ng	2461770	1,4-Dichlorobenzene-d4	7.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.03	21.4	ng	458224	1	7.93	428321	20.0
unknown7.46	7.46	115.0	ng	2461770	1	7.93	428321	20.0