

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF101118\
 Data File : BF109824.D
 Acq On : 12 Oct 2018 9:28
 Operator : JU/SJ
 Sample : J5296-16
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 100418-DBRI-E-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:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF100818.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	4.875	471	474	483	rBV	41790	57916	2.56%	0.440%
2	5.310	545	548	572	rBV	234378	434142	19.16%	3.296%
3	6.345	720	724	741	rBV	121123	319157	14.08%	2.423%
4	6.463	741	744	770	rVB	945579	1111814	49.07%	8.441%
5	6.692	780	783	806	rBV	294707	445884	19.68%	3.385%
6	6.851	806	810	838	rVV	1577247	1615025	71.27%	12.261%
7	7.257	876	879	914	rBV	904184	1320844	58.29%	10.028%
8	7.975	998	1001	1018	rBV	380990	495411	21.86%	3.761%
9	9.051	1180	1184	1203	rBV	1817179	2155568	95.13%	16.365%
10	9.445	1247	1251	1268	rVB	75814	95857	4.23%	0.728%
11	9.722	1294	1298	1316	rBV	468390	493869	21.79%	3.749%
12	10.510	1428	1432	1451	rBV	481054	704282	31.08%	5.347%
13	11.198	1546	1549	1569	rBV	315950	470928	20.78%	3.575%
14	12.780	1814	1818	1838	rBV	2223394	2265997	100.00%	17.203%
15	13.686	1969	1972	1984	rBV6	12771	33465	1.48%	0.254%
16	13.827	1992	1996	2009	rBV	483292	607854	26.83%	4.615%
17	14.904	2172	2179	2186	rBV	27985	34977	1.54%	0.266%
18	15.227	2229	2234	2250	rBV	251263	439934	19.41%	3.340%
19	15.645	2300	2305	2312	rVB2	25224	36941	1.63%	0.280%
20	16.098	2378	2382	2390	rVB2	18766	32057	1.41%	0.243%

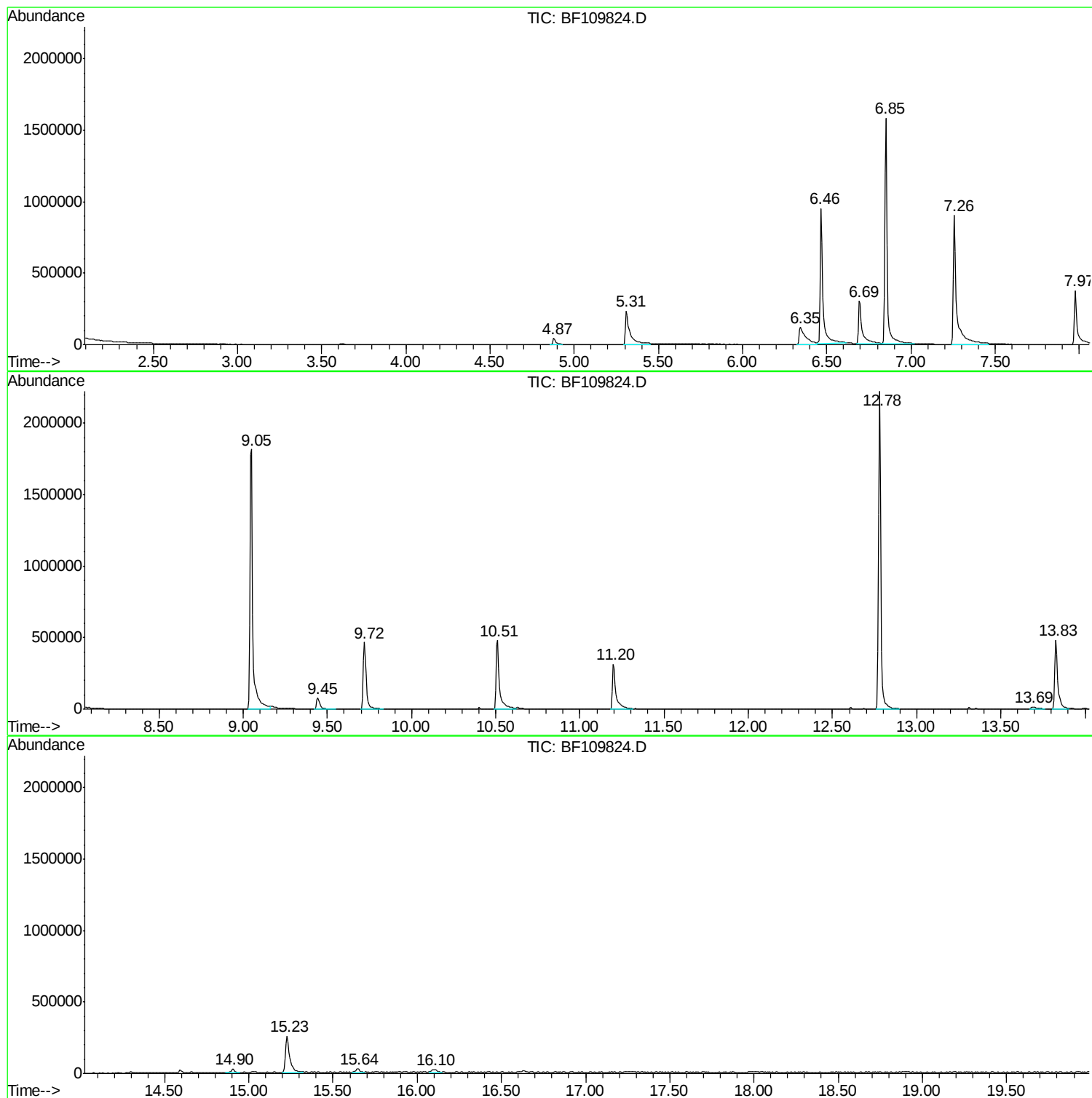
Sum of corrected areas: 13171922

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF100818.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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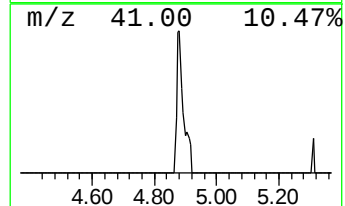
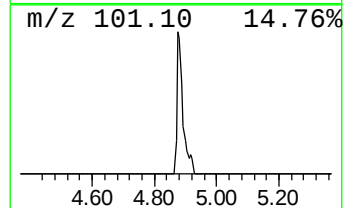
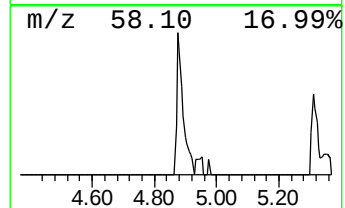
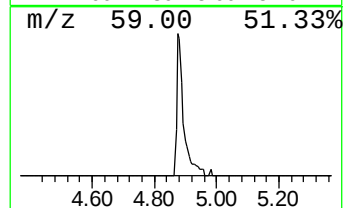
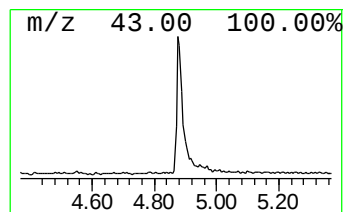
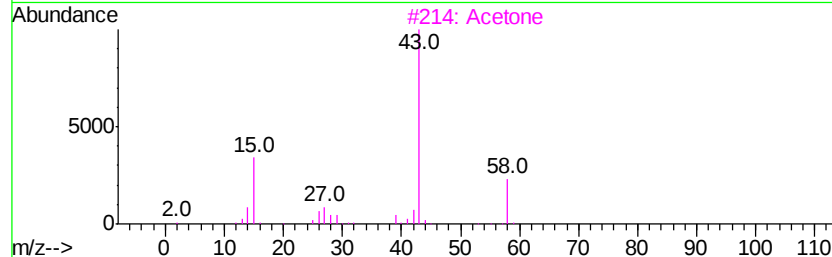
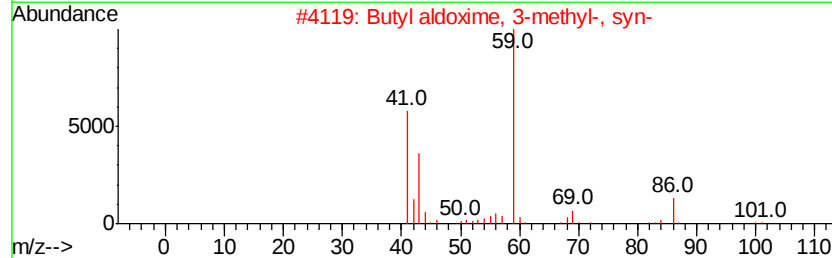
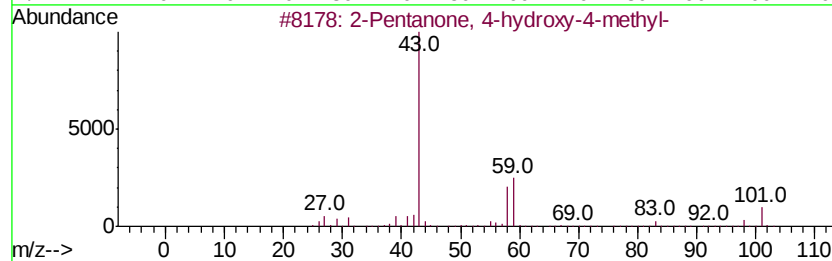
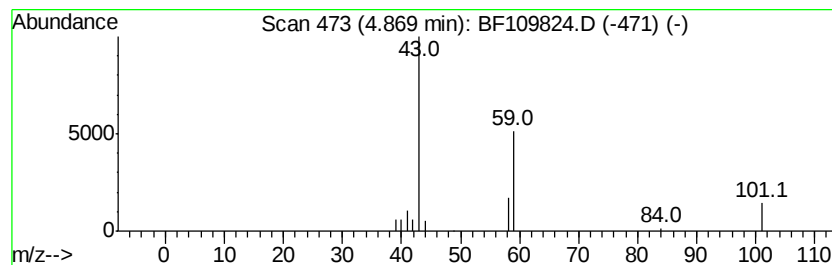
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.87	2.60 ng	57916	1,4-Dichlorobenzene-d4	6.70

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	17
3			Acetone	58	C3H6O	000067-64-1	9
4			4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	9
5			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	9



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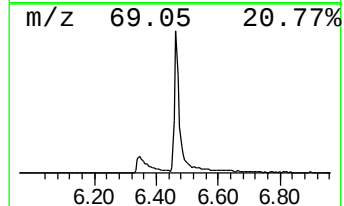
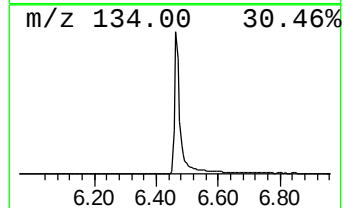
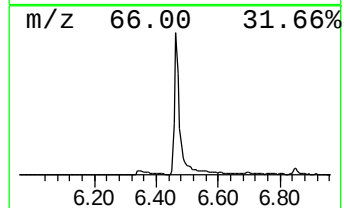
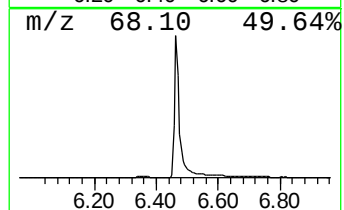
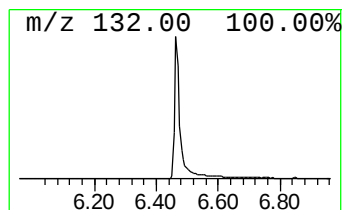
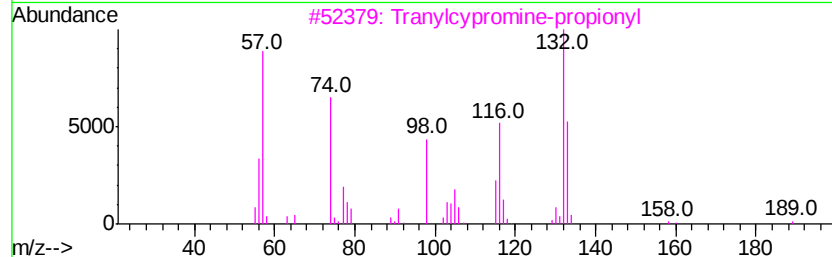
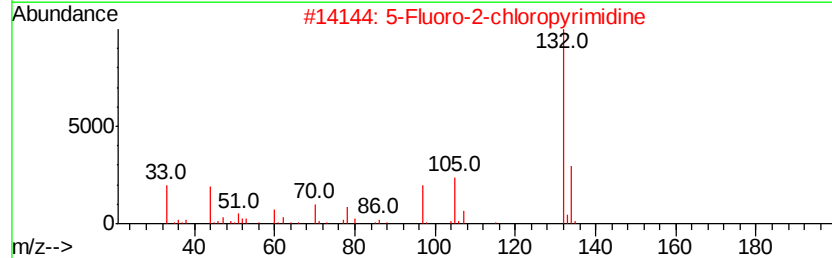
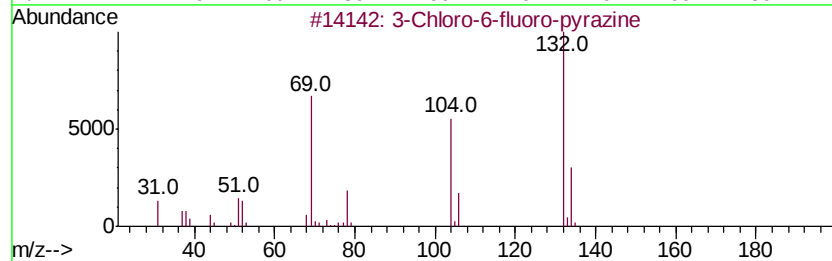
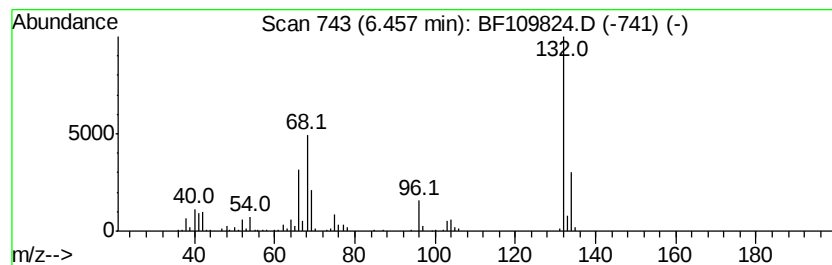
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Peak Number 2 unknown6.46 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.46	49.87 ng	1111810	1,4-Dichlorobenzene-d4	6.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	10
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.87	2.6	ng	57916	1	6.70	445884	20.0
unknown6.46	6.46	49.9	ng	1111810	1	6.70	445884	20.0