

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF092518\
 Data File : BF109329.D
 Acq On : 26 Sep 2018 10:52
 Operator : JU/SJ
 Sample : PB113191BL
 Misc :
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB113191BL

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-BF092218.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	4.269	368	371	382	rBV	43416	52491	1.63%	0.244%
2	4.904	474	479	493	rBV	400276	518079	16.11%	2.404%
3	5.322	545	550	553	rBV	2033681	1885261	58.64%	8.749%
4	6.346	718	724	726	rBV	1901927	1962166	61.03%	9.105%
5	6.475	742	746	749	rBV	2290959	1976992	61.49%	9.174%
6	6.704	782	785	789	rBV	877612	698018	21.71%	3.239%
7	6.863	808	812	815	rBV	2589759	2254199	70.11%	10.461%
8	7.269	875	881	884	rBV	1964081	1806639	56.19%	8.384%
9	7.987	999	1003	1007	rBV	1039199	862093	26.81%	4.001%
10	9.063	1181	1186	1190	rBV	3307119	3215100	100.00%	14.920%
11	9.739	1296	1301	1313	rVB	982659	906427	28.19%	4.206%
12	10.522	1430	1434	1450	rBV	1092775	1052860	32.75%	4.886%
13	11.216	1548	1552	1555	rBV	841855	756642	23.53%	3.511%
14	12.798	1816	1821	1824	rBV	2678143	2429337	75.56%	11.273%
15	13.839	1994	1998	2002	rBV	815034	692920	21.55%	3.216%
16	15.239	2231	2236	2241	rBV	411795	480040	14.93%	2.228%

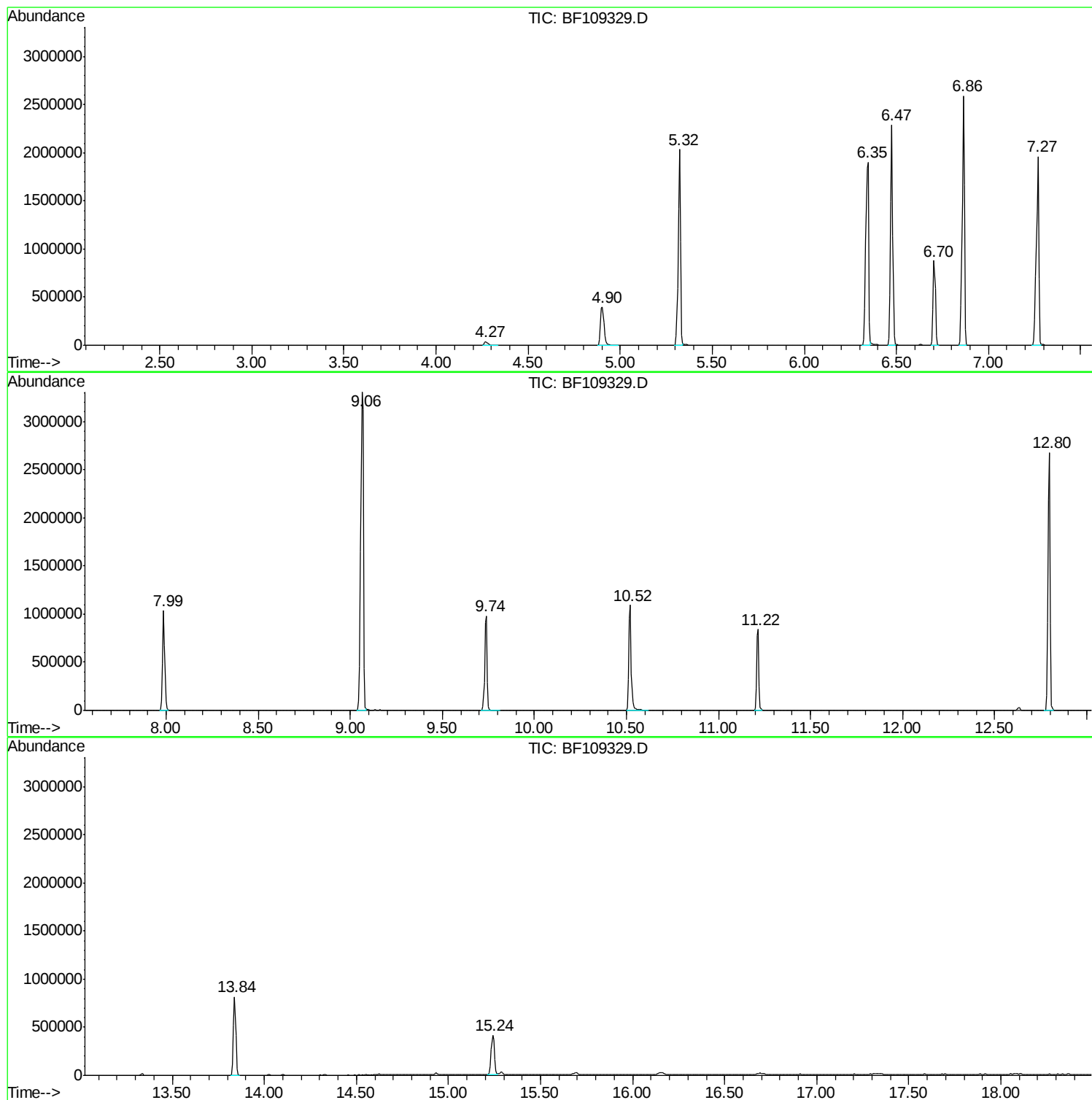
Sum of corrected areas: 21549264

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.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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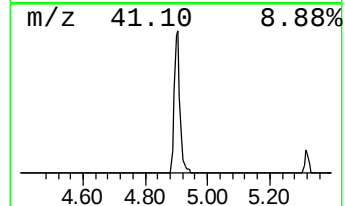
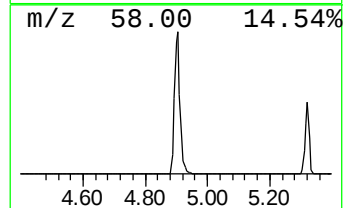
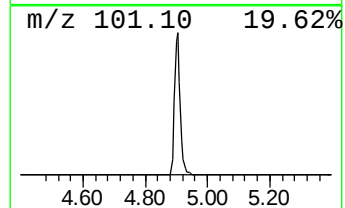
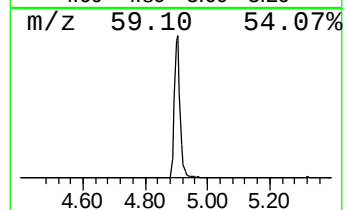
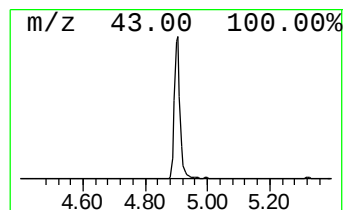
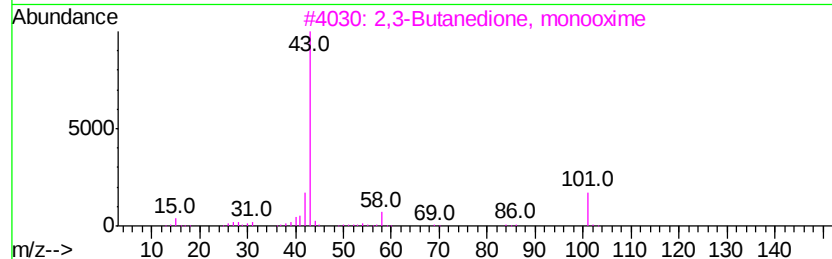
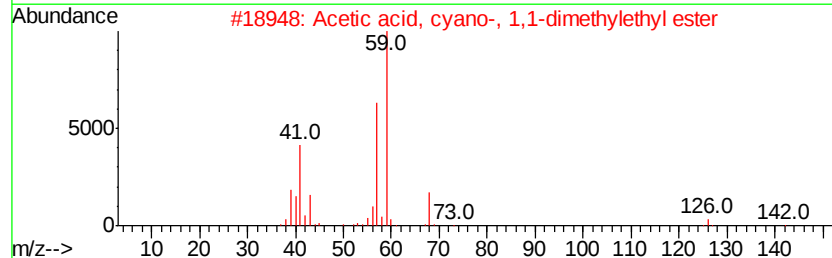
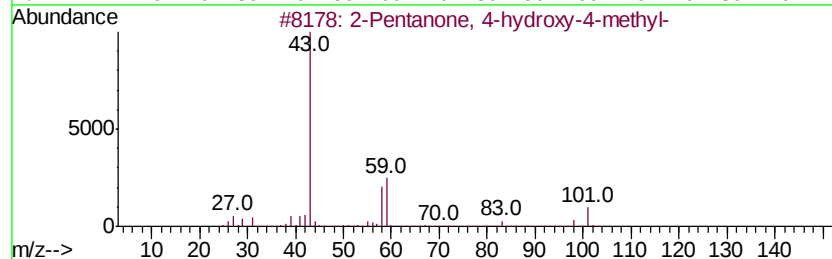
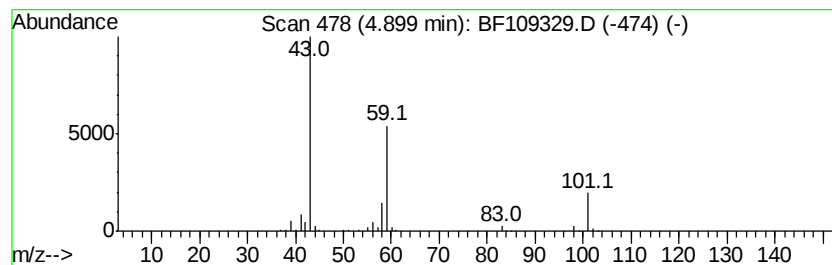
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.90	14.84 ng	518079	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	64
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			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5			Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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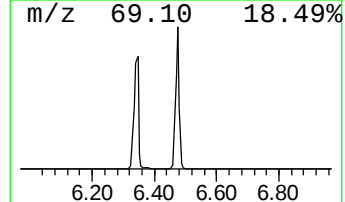
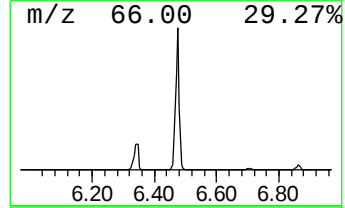
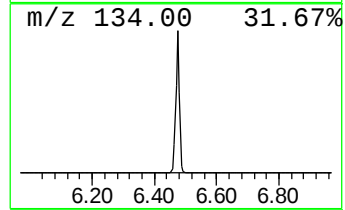
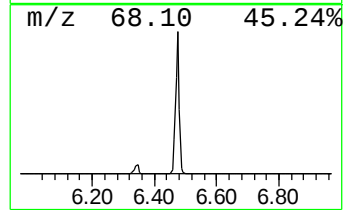
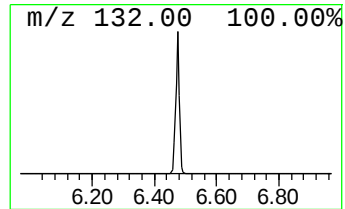
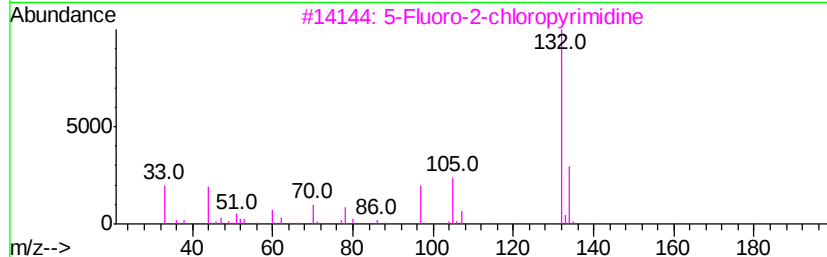
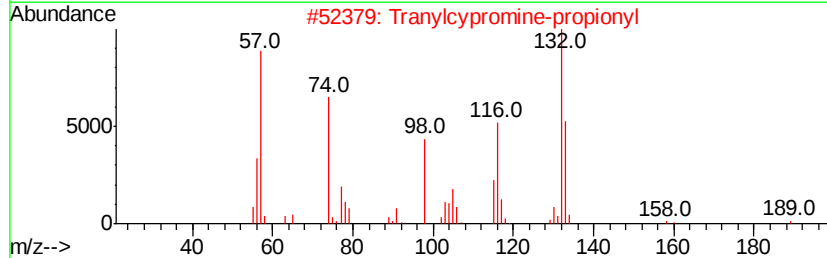
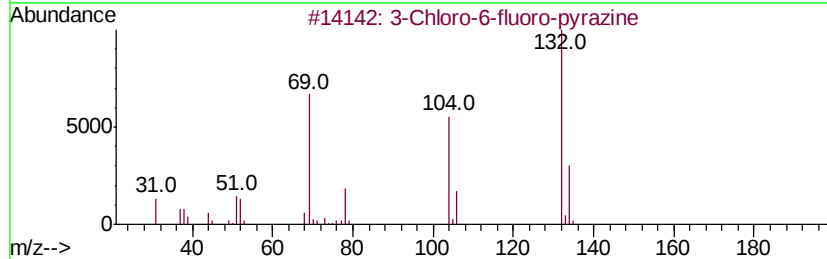
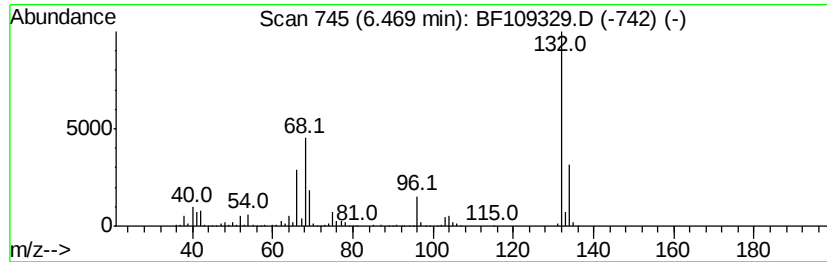
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Peak Number 2 unknown6.47 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.47	56.65 ng	1976990	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		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
4		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9
5		1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.90	14.8	ng	518079	1	6.70	698018	20.0
unknown6.47	6.47	56.6	ng	1976990	1	6.70	698018	20.0