

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF082517\
 Data File : BF098035.D
 Acq On : 25 Aug 2017 17:28
 Operator : SJ/JU
 Sample : PB101788BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB101788BL

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_F\METHODS\8270-BF081517.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	4.951	482	487	497	rVB	214851	313155	8.75%	1.077%
2	5.357	550	556	559	rBV	3135257	3364581	93.96%	11.566%
3	6.375	723	729	732	rBV	2955561	3472637	96.98%	11.938%
4	6.510	747	752	755	rBV	3181777	3301243	92.19%	11.349%
5	6.739	787	791	794	rBV	786460	715181	19.97%	2.459%
6	6.892	813	817	821	rBV	2662166	2546848	71.13%	8.755%
7	7.304	881	887	890	rBV	1956498	2156852	60.23%	7.415%
8	8.022	1004	1009	1012	rBV	905674	932131	26.03%	3.204%
9	9.098	1186	1192	1195	rBV	3426772	3580749	100.00%	12.310%
10	9.769	1301	1306	1310	rBV	972845	973675	27.19%	3.347%
11	10.563	1434	1441	1444	rBV	1830721	2020943	56.44%	6.947%
12	11.251	1553	1558	1562	rBV	1138239	990622	27.67%	3.405%
13	12.845	1823	1829	1832	rBV	2930187	3252520	90.83%	11.181%
14	13.880	2001	2005	2009	rBV	864749	857731	23.95%	2.949%
15	15.298	2241	2246	2251	rBV	524164	610427	17.05%	2.098%

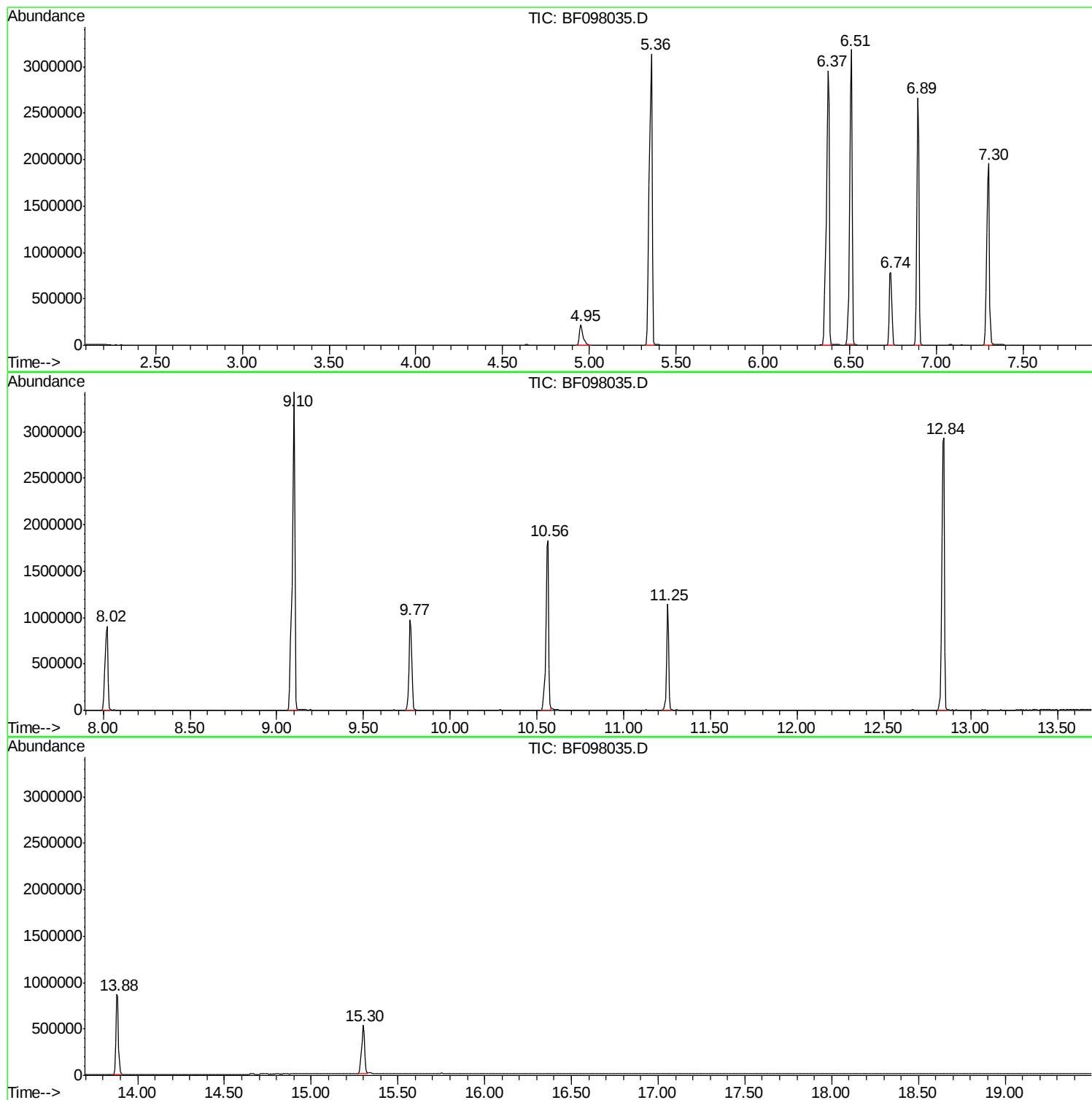
Sum of corrected areas: 29089295

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081517.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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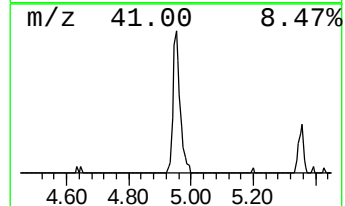
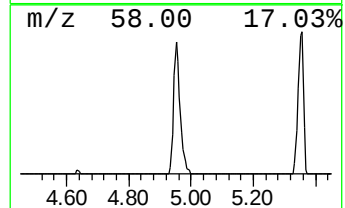
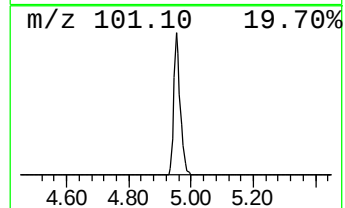
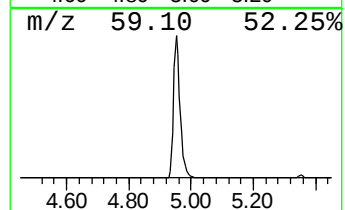
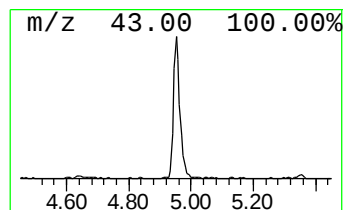
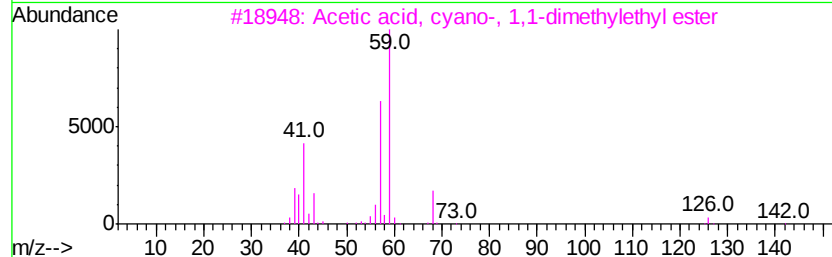
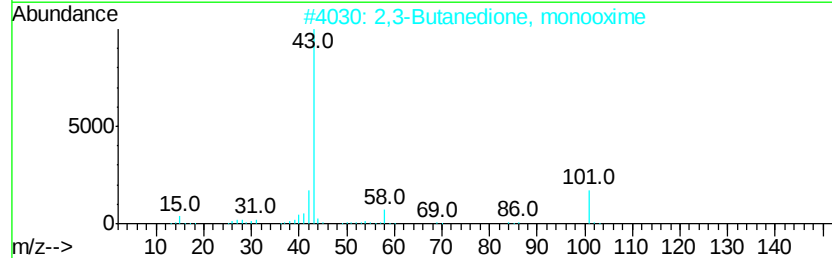
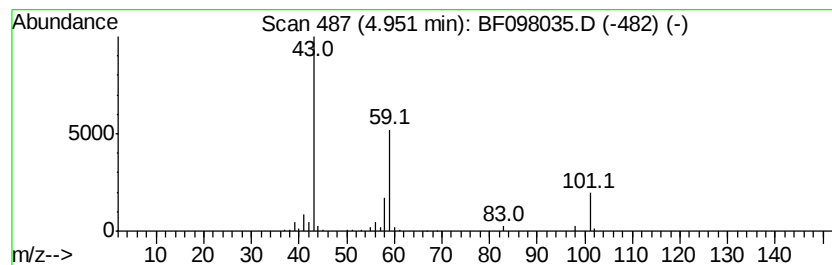
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.95	8.76 ng	313155	1,4-Dichlorobenzene-d4	6.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
3		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
4		Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	17
5		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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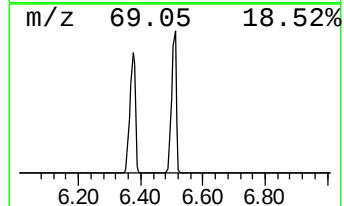
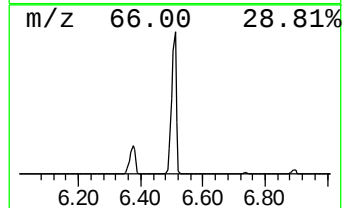
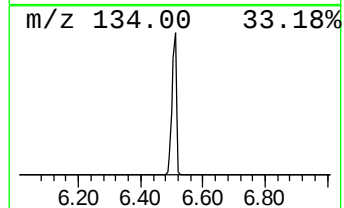
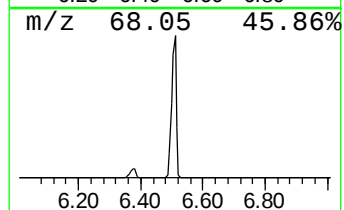
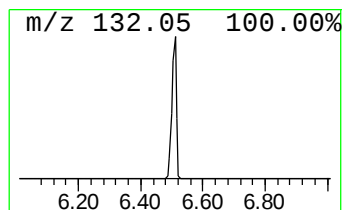
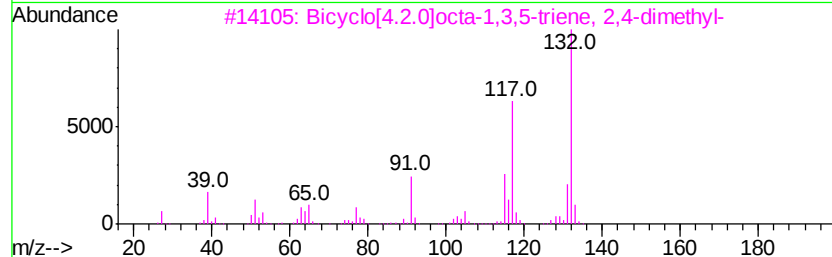
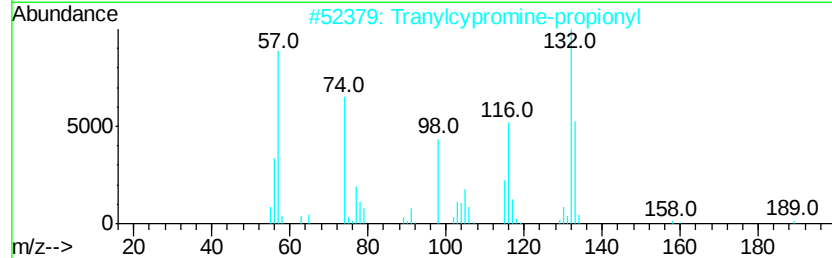
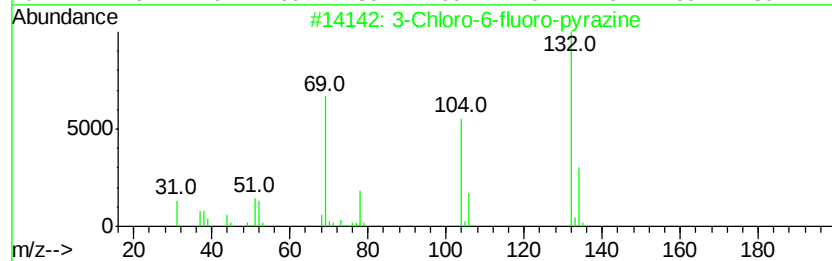
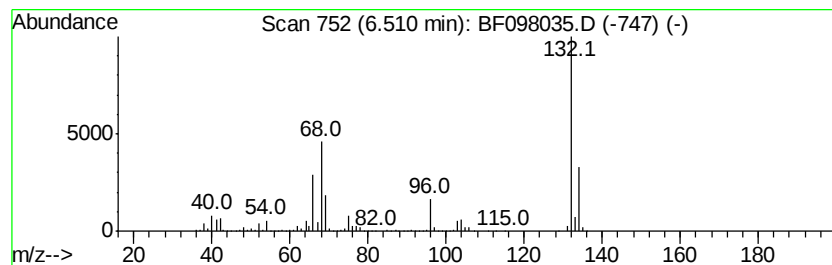
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Peak Number 2 unknown6.51 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.51	92.32 ng	3301240	1,4-Dichlorobenzene-d4	6.74

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		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		Xenon	132	Xe	007440-63-3	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...	4.95	8.8	ng	313155	1	6.74	715181	20.0
unknown6.51	6.51	92.3	ng	3301240	1	6.74	715181	20.0