

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060717\
 Data File : BF095746.D
 Acq On : 7 Jun 2017 15:15
 Operator : SJ/MA
 Sample : PB99520BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB99520BL

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-BF060517.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	4.598	508	512	530	rBV	57955	168844	6.15%	0.909%
2	5.281	623	628	649	rBV	1099041	1835957	66.86%	9.883%
3	6.939	905	910	922	rBV	1157521	1788594	65.13%	9.629%
4	7.039	922	927	942	rVB	1487430	2061451	75.07%	11.097%
5	7.381	981	985	997	rBV	290453	365586	13.31%	1.968%
6	7.622	1020	1026	1044	rBV	1132209	1533195	55.83%	8.254%
7	8.298	1135	1141	1154	rBV	818016	1133273	41.27%	6.101%
8	9.410	1326	1330	1344	rBV	387741	521995	19.01%	2.810%
9	11.198	1626	1634	1649	rBV	1824713	2523603	91.90%	13.585%
10	12.233	1805	1810	1823	rBV2	473091	621017	22.62%	3.343%
11	13.533	2025	2031	2051	rBV	1050279	1447967	52.73%	7.795%
12	14.627	2212	2217	2234	rBV2	509593	658917	24.00%	3.547%
13	17.009	2615	2622	2626	rBV2	2035169	2746004	100.00%	14.783%
14	18.345	2844	2849	2858	rBV	520024	643419	23.43%	3.464%
15	20.009	3128	3132	3143	rBV	413776	526205	19.16%	2.833%

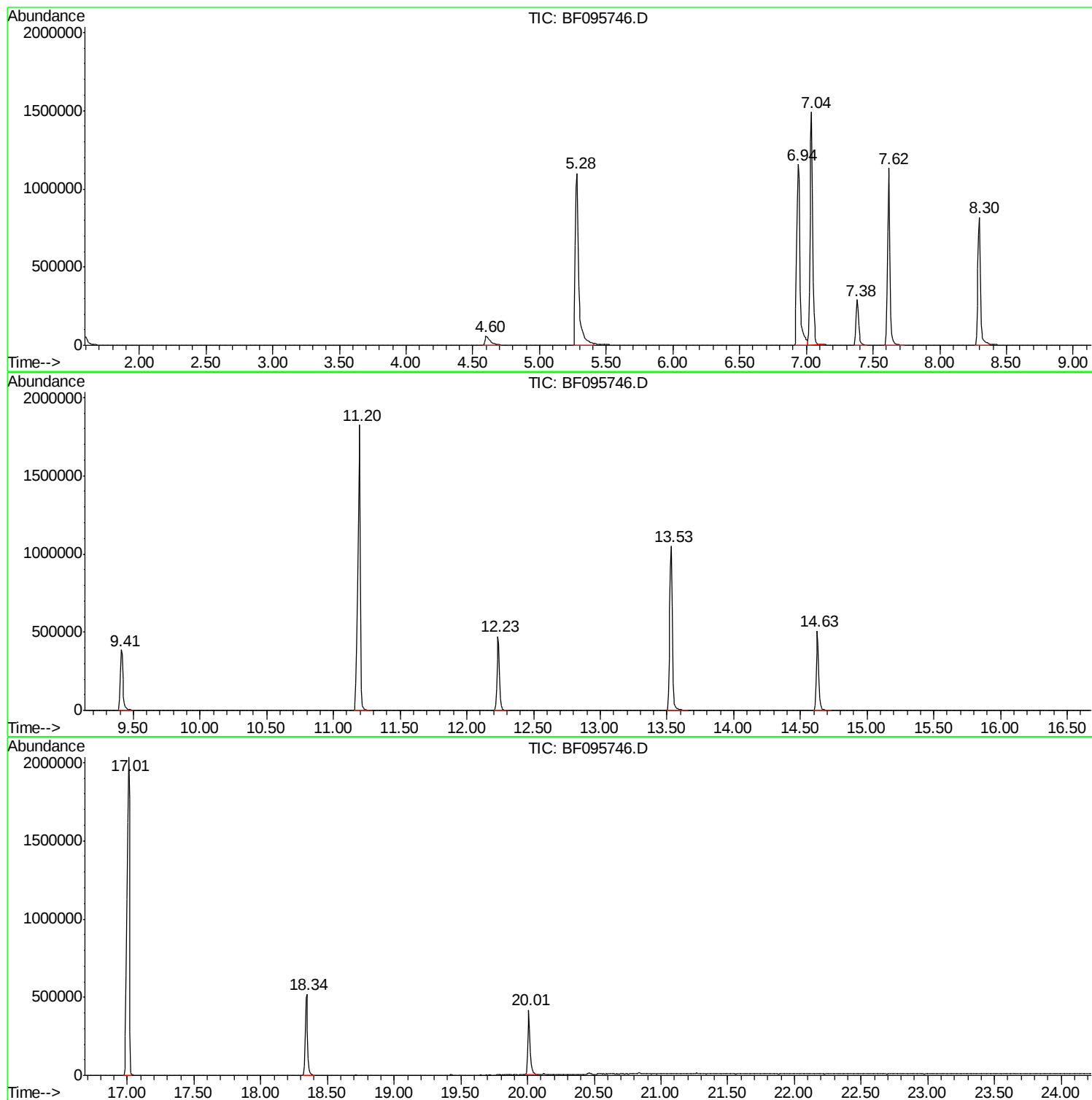
Sum of corrected areas: 18576027

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060517.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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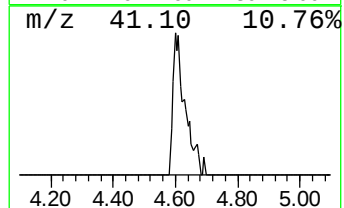
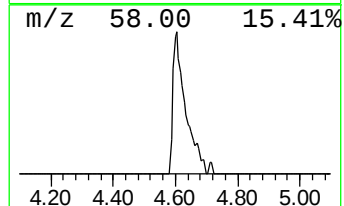
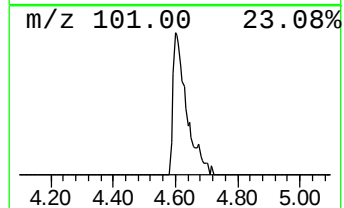
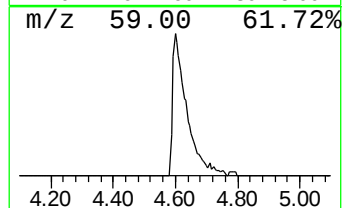
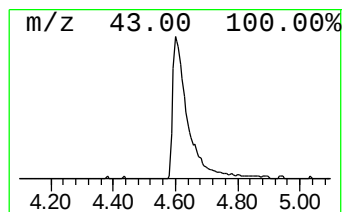
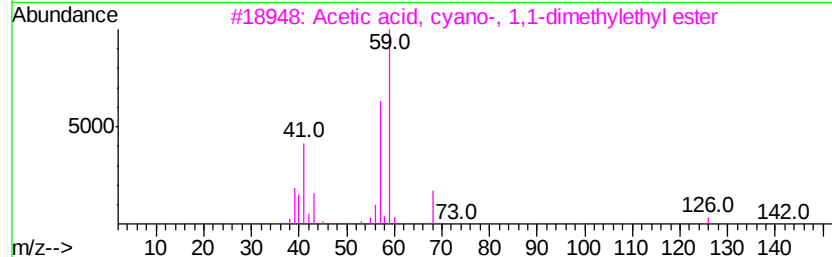
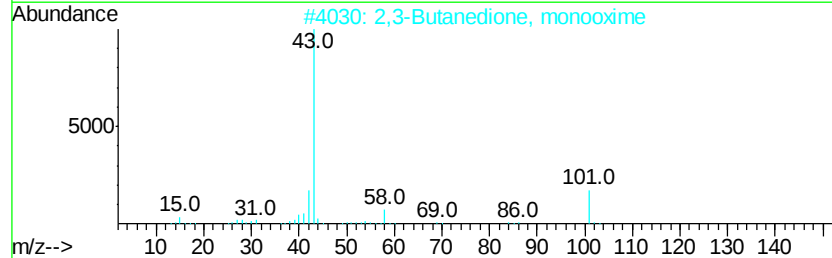
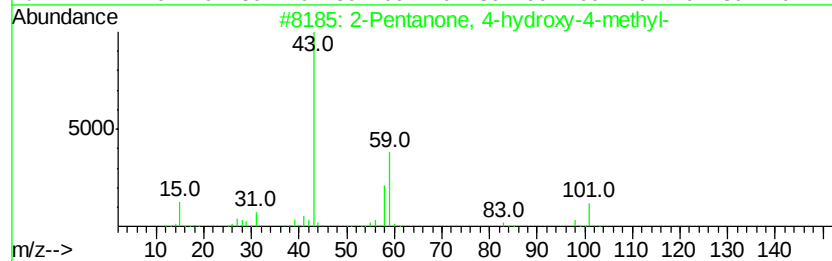
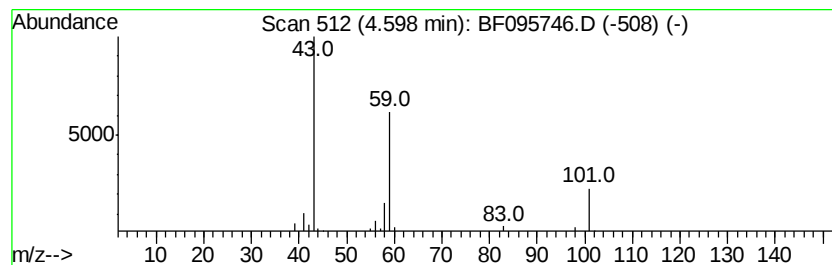
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.60	9.24 ng	168844	1,4-Dichlorobenzene-d4	7.38

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56	
2	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25	
3	Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17	
4	3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	17	
5	5-Hexen-2-one	98	C6H10O	000109-49-9	9	



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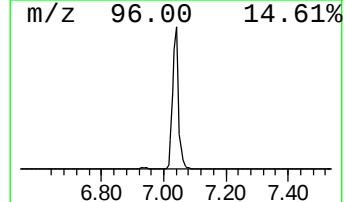
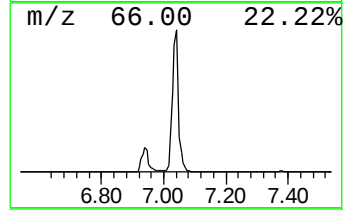
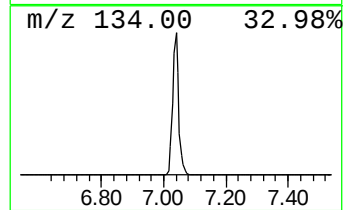
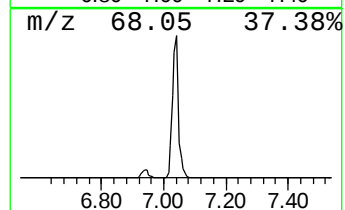
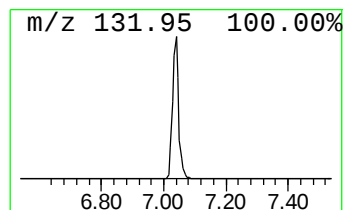
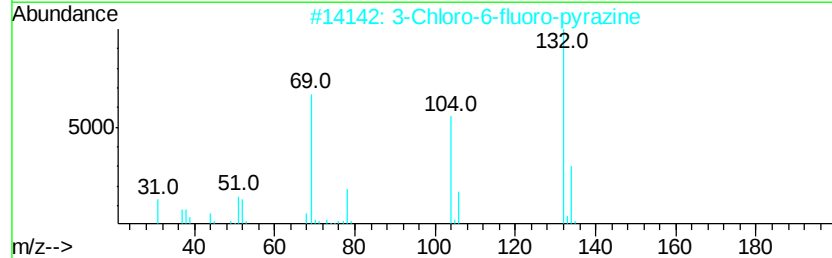
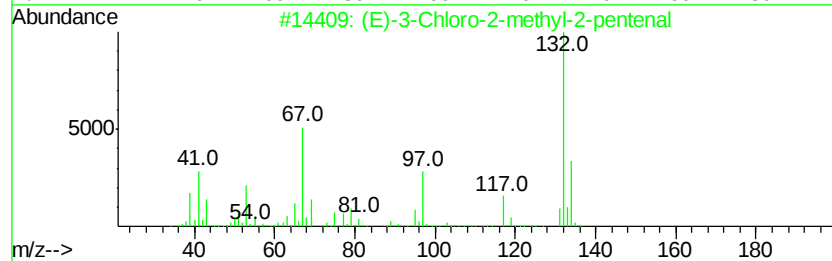
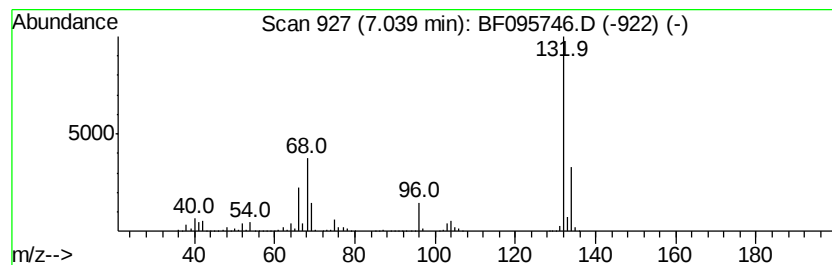
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Peak Number 2 unknown7.04 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.04	112.78 ng	2061450	1,4-Dichlorobenzene-d4	7.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	35
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.60	9.2	ng	168844	1	7.38	365586	20.0
unknown7.04	7.04	112.8	ng	2061450	1	7.38	365586	20.0