

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP110719\
 Data File : BP000995.D
 Acq On : 07 Nov 2019 10:09
 Operator : JU
 Sample : PB124565BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB124565BL

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_P\METHODS\8270-BP110619.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	5.740	617	621	637	rBV	730385	1152326	8.39%	1.426%
2	6.311	711	718	722	rBV	5402751	6935461	50.48%	8.584%
3	7.822	967	975	978	rBV	5241875	7245536	52.74%	8.968%
4	8.016	1001	1008	1027	rBV	5853043	7550195	54.96%	9.345%
5	8.375	1063	1069	1074	rBV	1357698	1514640	11.02%	1.875%
6	8.616	1104	1110	1115	rBV	5310148	6304533	45.89%	7.803%
7	9.252	1211	1218	1222	rBV	3828087	4596864	33.46%	5.690%
8	10.404	1407	1414	1419	rBV	1673813	1957229	14.25%	2.423%
9	12.181	1709	1716	1721	rBV	8292781	10489427	76.35%	12.983%
10	13.263	1894	1900	1912	rBV	2136445	2545339	18.53%	3.150%
11	14.557	2113	2120	2147	rBV	5874775	7434607	54.11%	9.202%
12	15.657	2302	2307	2312	rBV	2709829	2945650	21.44%	3.646%
13	17.939	2689	2695	2699	rBV	11015338	13738564	100.00%	17.005%
14	19.292	2920	2925	2935	rBV	2912480	3167931	23.06%	3.921%
15	21.080	3223	3229	3246	rVB	2312202	3215199	23.40%	3.980%

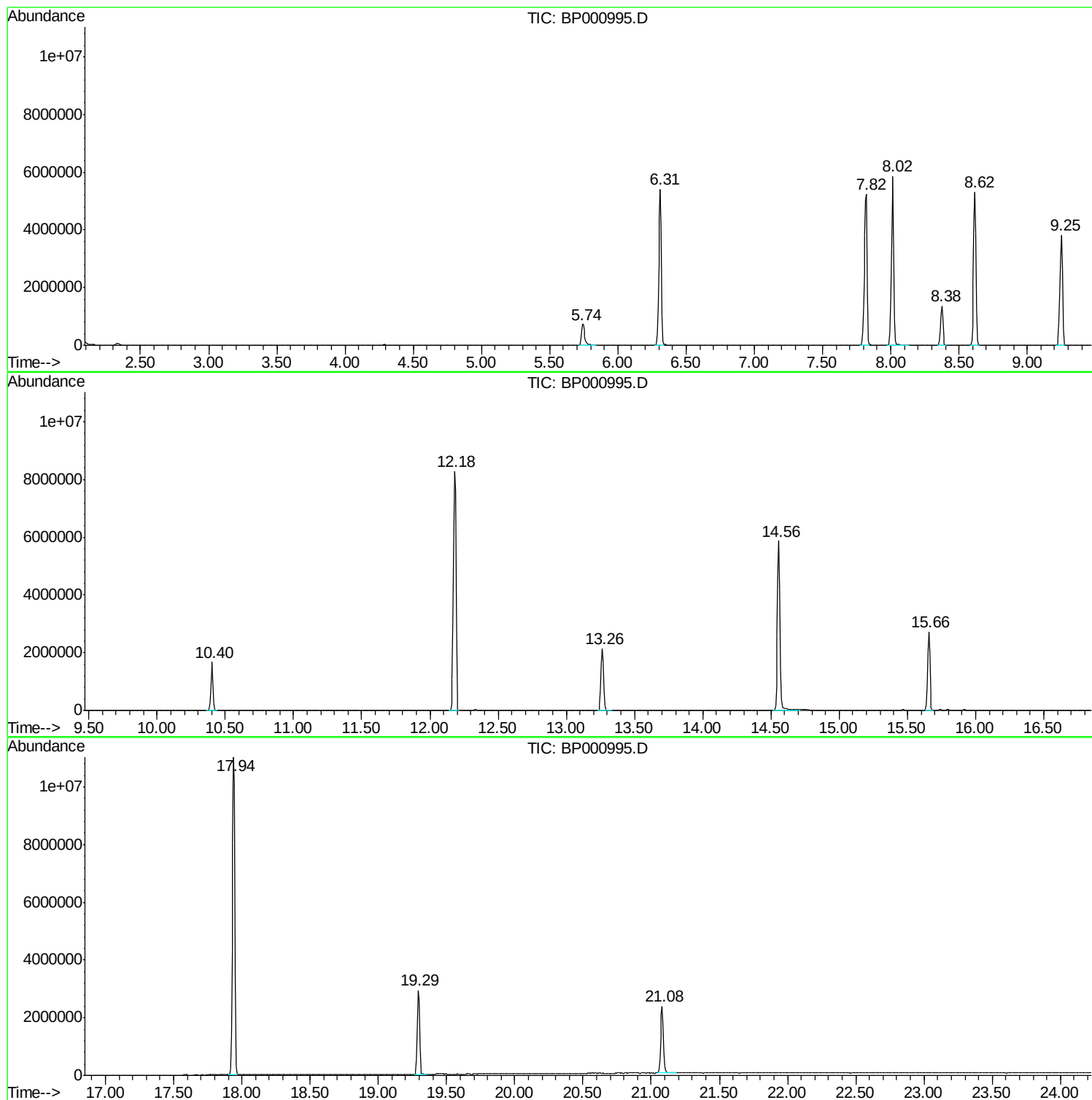
Sum of corrected areas: 80793501

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP110619.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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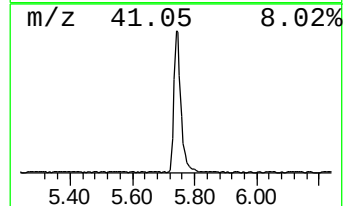
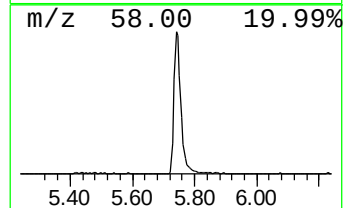
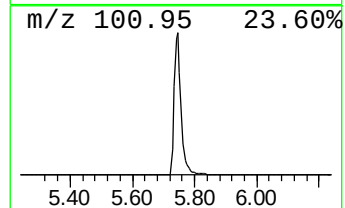
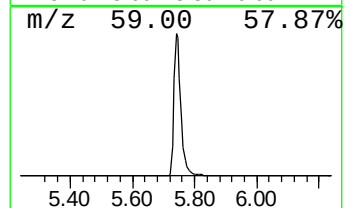
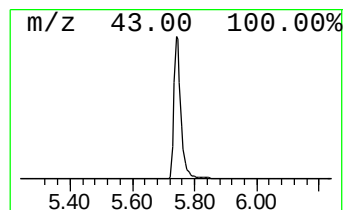
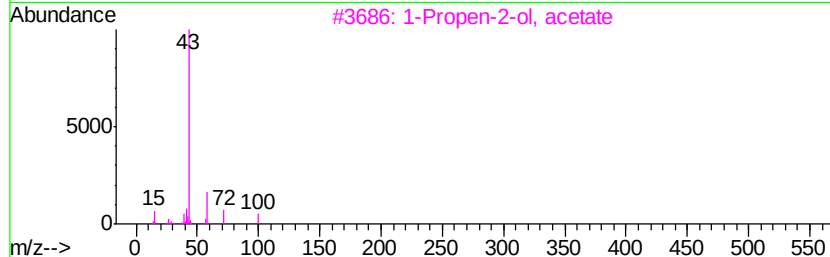
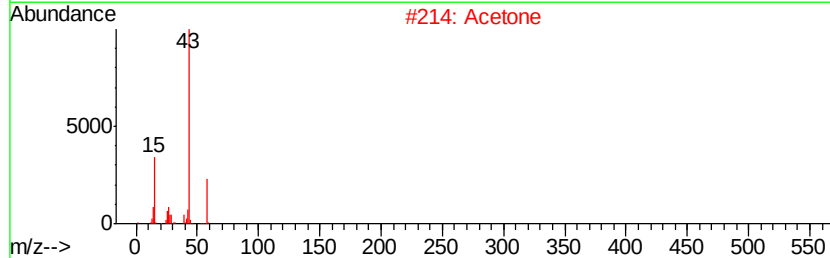
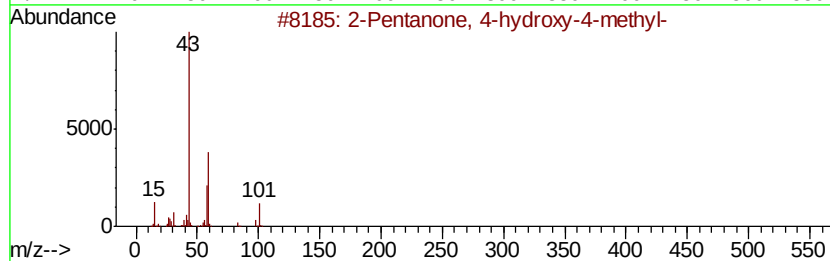
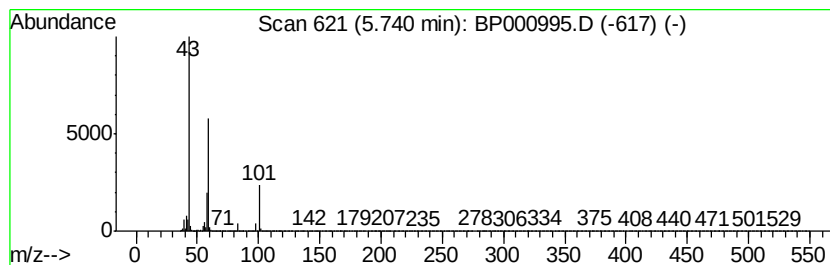
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.74	15.22 ng	1152330	1,4-Dichlorobenzene-d4	8.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2		Acetone	58	C3H6O	000067-64-1	12
3		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
4		5-Hexen-2-one	98	C6H10O	000109-49-9	9
5		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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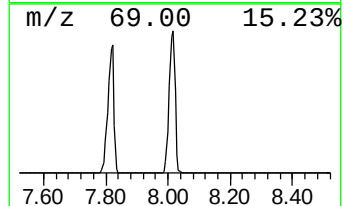
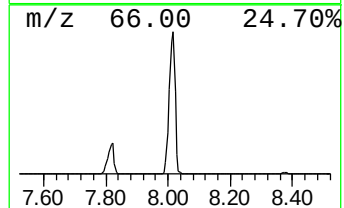
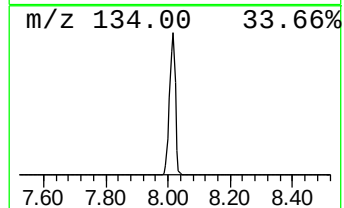
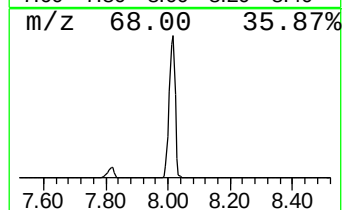
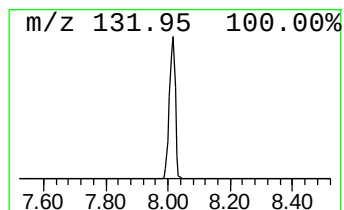
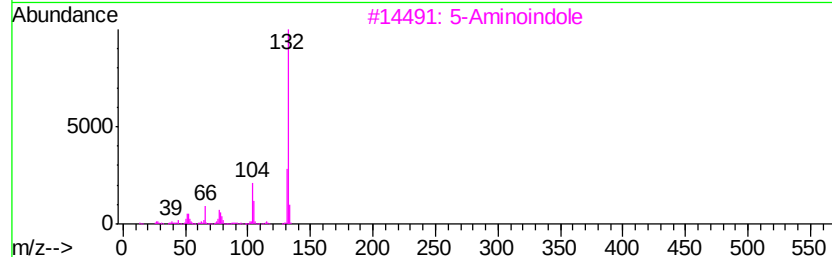
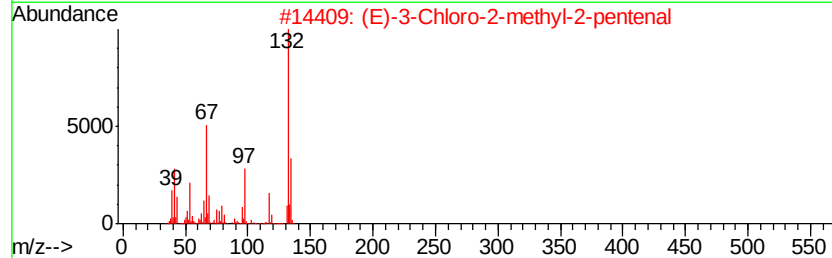
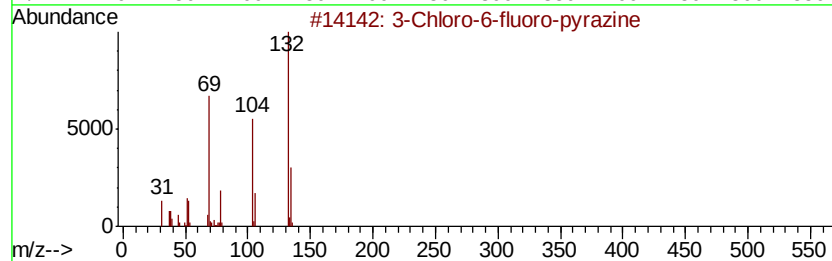
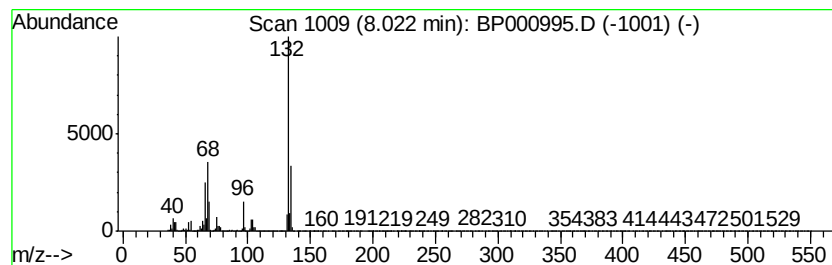
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Peak Number 2 unknown8.02 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.02	99.70 ng	7550200	1,4-Dichlorobenzene-d4	8.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	27
3		5-Aminoindole	132	C8H8N2	005192-03-0	22
4		N-(-3,4-Methylenedioxyphenylisop...	267	C17H17NO2	1000378-96-6	10
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.74	15.2	ng	1152330	1	8.38	1514640	20.0
unknown8.02	8.02	99.7	ng	7550200	1	8.38	1514640	20.0