

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN041019\
 Data File : BN005135.D
 Acq On : 10 Apr 2019 19:29
 Operator : JU/SJ
 Sample : K2311-02
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 0168-SITE-WATER

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_N\METHODS\8270-BN032619.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	5.240	378	383	390	rBB	70282	97328	23.62%	5.192%
2	6.828	647	653	659	rBB	54164	78054	18.94%	4.164%
3	7.187	708	714	720	rBB	118299	178445	43.31%	9.520%
4	7.652	788	793	798	rBB	21751	32379	7.86%	1.727%
5	7.963	841	846	852	rBB	86556	134351	32.61%	7.168%
6	8.816	985	991	997	rBB	71636	112447	27.29%	5.999%
7	10.434	1260	1266	1272	rBB	31911	48398	11.75%	2.582%
8	12.916	1681	1688	1694	rBB	197158	284513	69.05%	15.179%
9	14.298	1918	1923	1929	rBB2	46452	66356	16.10%	3.540%
10	15.798	2173	2178	2185	rBB	170824	221172	53.68%	11.799%
11	17.051	2386	2391	2397	rBB2	55503	73535	17.85%	3.923%
12	19.686	2834	2839	2844	rBV	351054	412022	100.00%	21.981%
13	21.251	3100	3105	3110	rBV	58692	69055	16.76%	3.684%
14	23.480	3477	3484	3491	rVB2	38699	66372	16.11%	3.541%

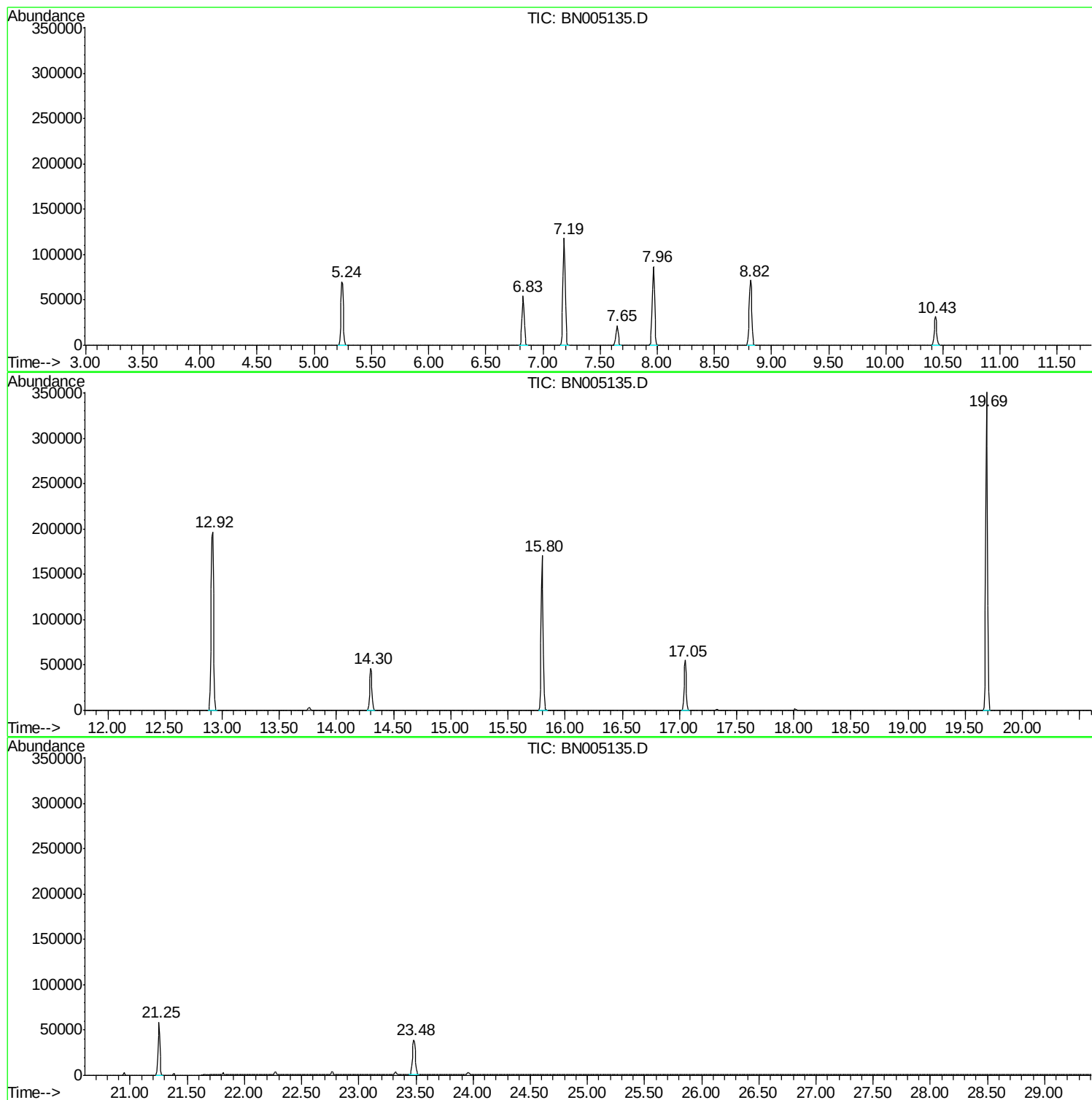
Sum of corrected areas: 1874427

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-BN032619.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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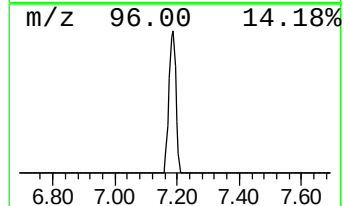
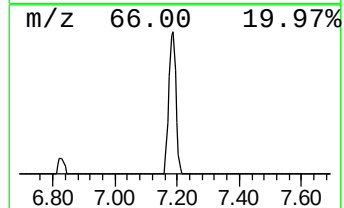
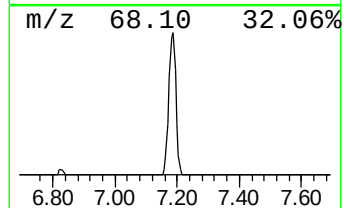
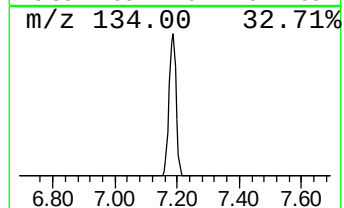
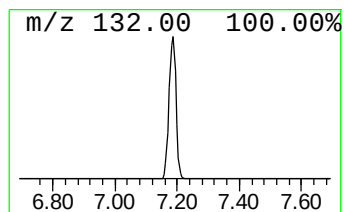
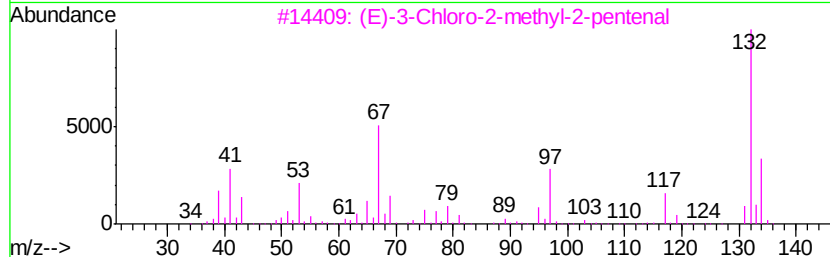
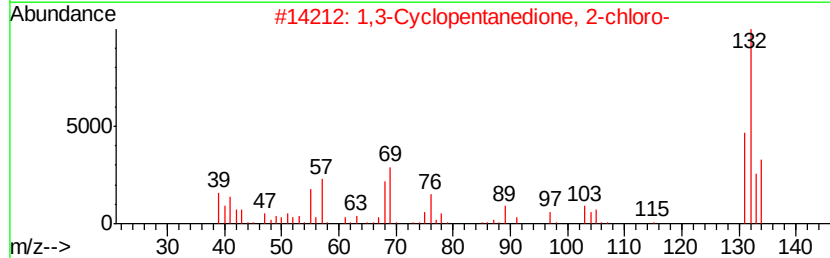
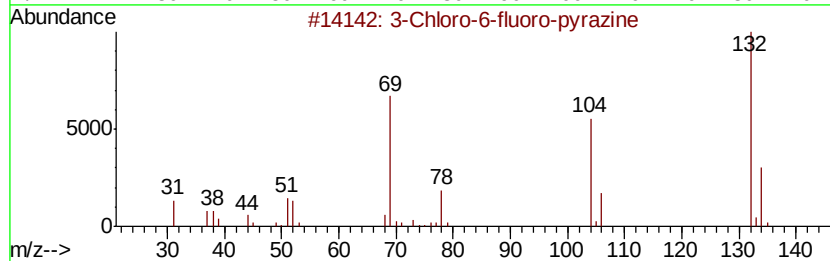
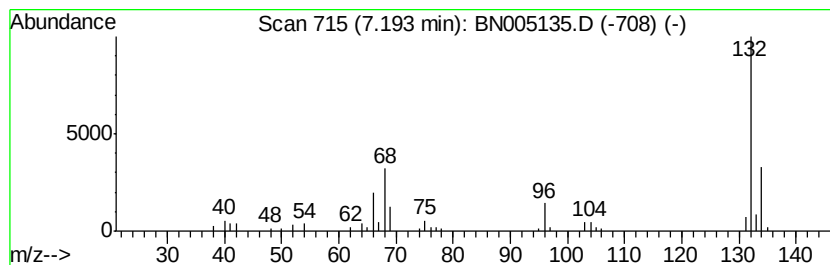
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Peak Number 1 unknown7.19 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.19	110.22 ng	178445	1,4-Dichlorobenzene-d4	7.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
2		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	23
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
4		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
5		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown7.19	7.19	110.2	ng	178445	1	7.65	32379	20.0