

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE081718\
 Data File : BE097301.D
 Acq On : 17 Aug 2018 17:49
 Operator : SJ/JU
 Sample : J4424-07
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A3YR7

Manual Integrations
 APPROVED

Sohil
 8/20/2018 4:15:31 PM

Quant Time: Aug 17 23:47:57 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE081718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 17 23:33:44 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	1037	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	5276	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.01	164	2929	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	6832m	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	5710m	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	4377m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.72	152	1657	0.20	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	4002m	0.22	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	16.81	178	678m	0.038	ng/ul	
15) Fluoranthene	18.84	202	873m	0.047	ng/ul	
17) Pyrene	19.20	202	967	0.056	ng/ul	93
18) Benzo(a)anthracene	20.96	228	396	0.028	ng/ul#	78
19) Chrysene	21.01	228	422	0.025	ng/ul#	72
21) Benzo(b)fluoranthene	22.54	252	354m	0.030	ng/ul	
23) Benzo(a)pyrene	23.11	252	286m	0.026	ng/ul	

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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