

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM031017\
 Data File : BM009172.D
 Acq On : 10 Mar 2017 12:47
 Operator : SJ/MA
 Sample : I2107-09
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AJ8

Manual Integrations
 APPROVED

mohammad
 3/13/2017 12:49:11 PM

Quant Time: Mar 10 16:42:21 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM030317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 10 15:19:24 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.88	152	233	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	713	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	430	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	946m	0.40	ng/ul	0.00
16) Chrysene-d12	21.44	240	1026	0.40	ng/ul	0.00
20) Perylene-d12	23.77	264	913m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	209	0.18	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	567	0.19	ng/ul	0.00
Target Compounds				Qvalue		
12) Phenanthrene	17.30	178	320m	0.11	ng/ul	
15) Fluoranthene	19.33	202	932	0.24	ng/ul	88
17) Pyrene	19.69	202	1002	0.28	ng/ul#	91
18) Benzo(a)anthracene	21.42	228	611	0.18	ng/ul#	85
19) Chrysene	21.47	228	708	0.22	ng/ul#	82
21) Benzo(b)fluoranthene	23.07	252	986m	0.27	ng/ul	
22) Benzo(k)fluoranthene	23.11	252	329m	0.10	ng/ul	
23) Benzo(a)pyrene	23.67	252	615m	0.19	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.15	276	412m	0.10	ng/ul	
26) Benzo(g,h,i)perylene	26.89	276	357m	0.10	ng/ul	

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

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