

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM031017\
 Data File : BM009171.D
 Acq On : 10 Mar 2017 12:11
 Operator : SJ/MA
 Sample : I2107-15
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AL0

Manual Integrations
 APPROVED

mohammad
 3/13/2017 12:49:06 PM

Quant Time: Mar 10 16:37:29 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	228	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	724	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	423	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	928m	0.40	ng/ul	0.00
16) Chrysene-d12	21.44	240	1064	0.40	ng/ul	0.00
20) Perylene-d12	23.77	264	963m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	200	0.17	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	554	0.19	ng/ul	0.00
Target Compounds						
3) Naphthalene	10.73	128	282	0.14	ng/ul#	59
5) 2-Methylnaphthalene	12.34	142	231	0.15	ng/ul	100
12) Phenanthrene	17.30	178	426m	0.15	ng/ul	
15) Fluoranthene	19.33	202	793	0.21	ng/ul	87
17) Pyrene	19.69	202	805	0.22	ng/ul#	85
18) Benzo(a)anthracene	21.42	228	409	0.12	ng/ul#	75
19) Chrysene	21.47	228	528	0.16	ng/ul#	77
21) Benzo(b)fluoranthene	23.07	252	712m	0.19	ng/ul	
22) Benzo(k)fluoranthene	23.11	252	243m	0.07	ng/ul	
23) Benzo(a)pyrene	23.67	252	471m	0.14	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.15	276	328m	0.08	ng/ul	
26) Benzo(g,h,i)perylene	26.88	276	287m	0.08	ng/ul	

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

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