

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM030817\
 Data File : BM009123.D
 Acq On : 08 Mar 2017 23:29
 Operator : SJ/MA
 Sample : I2062-18
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AK4

Manual Integrations
 APPROVED

mohammad
 3/9/2017 2:01:43 PM

Quant Time: Mar 09 02:57:17 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM030317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 09 02:44:03 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.88	152	266	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	849	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.52	164	495	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	1140m	0.40	ng/ul	0.00
16) Chrysene-d12	21.45	240	1189	0.40	ng/ul	0.00
20) Perylene-d12	23.78	264	1120m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	235	0.17	ng/ul	0.00
14) Fluoranthene-d10	19.30	212	636	0.18	ng/ul	0.00
Target Compounds						
3) Naphthalene	10.73	128	102	0.04	ng/ul#	10
12) Phenanthrene	17.31	178	491m	0.14	ng/ul	
13) Anthracene	17.40	178	162	0.05	ng/ul#	37
15) Fluoranthene	19.33	202	1581	0.34	ng/ul	96
17) Pyrene	19.69	202	1572	0.38	ng/ul	95
18) Benzo(a)anthracene	21.43	228	1153	0.29	ng/ul#	92
19) Chrysene	21.48	228	1008	0.27	ng/ul#	89
21) Benzo(b)fluoranthene	23.07	252	1551m	0.35	ng/ul	
22) Benzo(k)fluoranthene	23.11	252	584m	0.15	ng/ul	
23) Benzo(a)pyrene	23.68	252	1188m	0.30	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.16	276	649m	0.13	ng/ul	
26) Benzo(g,h,i)perylene	26.89	276	618m	0.14	ng/ul	

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

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