

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM030917\
 Data File : BM009148.D
 Acq On : 09 Mar 2017 16:45
 Operator : SJ/MA
 Sample : I2069-15
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AA0

Manual Integrations
 APPROVED

mohammad
 3/10/2017 6:01:09 PM

Quant Time: Mar 09 17:31:45 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	204	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	652	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.52	164	392	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	907m	0.40	ng/ul	0.00
16) Chrysene-d12	21.44	240	986	0.40	ng/ul	0.00
20) Perylene-d12	23.78	264	1087m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	226	0.21	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	620	0.22	ng/ul	0.00
Target Compounds						
3) Naphthalene	10.73	128	215	0.11	ng/ul#	53
5) 2-Methylnaphthalene	12.34	142	147	0.11	ng/ul	95
12) Phenanthrene	17.30	178	651m	0.23	ng/ul	
15) Fluoranthene	19.32	202	1148	0.31	ng/ul	94
17) Pyrene	19.68	202	1511	0.45	ng/ul	93
18) Benzo(a)anthracene	21.42	228	668	0.20	ng/ul#	88
19) Chrysene	21.47	228	713	0.23	ng/ul#	86
21) Benzo(b)fluoranthene	23.07	252	1120m	0.26	ng/ul	
22) Benzo(k)fluoranthene	23.11	252	280m	0.07	ng/ul	
23) Benzo(a)pyrene	23.67	252	583m	0.15	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.15	276	475m	0.10	ng/ul	
26) Benzo(g,h,i)perylene	26.89	276	461m	0.11	ng/ul	

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

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