

Data Path : Z:\HPCHEM1\BNA_M\Data\BM030917\
 Data File : BM009151.D
 Acq On : 09 Mar 2017 18:33
 Operator : SJ/MA
 Sample : I2069-21
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AJ6

Manual Integrations
 APPROVED

mohammad
 3/10/2017 6:01:15 PM

Quant Time: Mar 20 09:27:10 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 03:16:46 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.88	152	180	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	579	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	362	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	833m	0.40	ng/ul	0.00
16) Chrysene-d12	21.44	240	967	0.40	ng/ul	0.00
20) Perylene-d12	23.78	264	833m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	284	0.30	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	859	0.33	ng/ul	0.00
Target Compounds				Qvalue		
12) Phenanthrene	17.30	178	545m	0.21	ng/ul	
13) Anthracene	17.40	178	157	0.06	ng/ul#	39
15) Fluoranthene	19.33	202	767	0.22	ng/ul	88
17) Pyrene	19.69	202	871	0.26	ng/ul#	89
18) Benzo(a)anthracene	21.42	228	456	0.14	ng/ul#	75
19) Chrysene	21.47	228	408	0.13	ng/ul#	71
21) Benzo(b)fluoranthene	23.07	252	637m	0.19	ng/ul	
22) Benzo(k)fluoranthene	23.11	252	223m	0.08	ng/ul	
23) Benzo(a)pyrene	23.67	252	407m	0.14	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.15	276	315m	0.08	ng/ul	
26) Benzo(g,h,i)perylene	26.89	276	291m	0.09	ng/ul	

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

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