

Data Path : Z:\HPCHEM1\BNA M\DATA\BM102315\
 Data File : BM002976.D
 Acq On : 22 Oct 2015 13:56
 Operator : UM/IZ
 Sample : G4074-09
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 JG9K3

Manual Integrations
 APPROVED

mmdadoda
 10/23/2015 4:43:27 PM

Quant Time: Oct 23 03:06:10 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-SIM-BM102215.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 23 01:01:42 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.64	152	5647	0.40	ng/ul	0.00
4) Naphthalene-d8	11.49	136	24773	0.40	ng/ul	0.00
8) Acenaphthene-d10	15.23	164	13914	0.40	ng/ul	0.00
12) Phenanthrene-d10	17.94	188	32414m	0.40	ng/ul	0.00
18) Chrysene-d12	22.03	240	28370	0.40	ng/ul	0.00
22) Perylene-d12	24.60	264	24114	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.71	96	12038	2.00	ng/uL	0.00
6) 2-Methylnaphthalene-d10	13.04	152	6135	0.15	ng/ul	0.00
16) Fluoranthene-d10	19.93	212	14646	0.17	ng/ul	0.00
Target Compounds						
5) Naphthalene	11.54	128	833	0.01	ng/ul#	58
7) 2-Methylnaphthalene	13.11	142	534m	0.01	ng/ul	
9) Acenaphthylene	14.96	152	2213	0.03	ng/ul#	80
14) Phenanthrene	17.98	178	16370m	0.16	ng/ul	
17) Fluoranthene	19.95	202	77318	0.64	ng/ul	93
19) Pyrene	20.30	202	67291	0.65	ng/ul	97
20) Benzo(a)anthracene	22.01	228	18618	0.20	ng/ul#	83
21) Chrysene	22.07	228	55641	0.58	ng/ul#	90
23) Benzo(b)fluoranthene	23.80	252	59155m	0.69	ng/ul	
26) Indeno(1,2,3-cd)pyrene	27.31	276	17085	0.19	ng/ul#	91
28) Benzo(g,h,i)perylene	28.15	276	18504	0.25	ng/ul#	55

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

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