

Data Path : Z:\HPCHEM1\BNA M\DATA\BM102215\
 Data File : BM002965.D
 Acq On : 22 Oct 2015 01:49
 Operator : UM/IZ
 Sample : G4075-10
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 JG9L1

Manual Integrations
 APPROVED

MMdadoda
 10/22/2015 4:57:01 PM

Quant Time: Oct 22 13:31:07 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 : Wed Oct 21 16:23:48 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.64	152	7637	0.40	ng/ul	-0.02
4) Naphthalene-d8	11.49	136	33502	0.40	ng/ul	-0.01
8) Acenaphthene-d10	15.23	164	18282	0.40	ng/ul	0.00
12) Phenanthrene-d10	17.94	188	38923m	0.40	ng/ul	-0.02
18) Chrysene-d12	22.03	240	28484	0.40	ng/ul	-0.01
22) Perylene-d12	24.60	264	23752	0.40	ng/ul	-0.02
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.71	96	23927	2.93	ng/uL	0.00
6) 2-Methylnaphthalene-d10	13.04	152	11490	0.20	ng/ul	-0.01
16) Fluoranthene-d10	19.93	212	22858	0.22	ng/ul	0.00
Target Compounds						
17) Fluoranthene	19.95	202	8332	0.06	ng/ul	83
19) Pyrene	20.30	202	11331	0.11	ng/ul#	86
20) Benzo(a)anthracene	22.01	228	4173m	0.04	ng/ul	
21) Chrysene	22.07	228	8935m	0.09	ng/ul	
26) Indeno(1,2,3-cd)pyrene	27.32	276	3965	0.05	ng/ul#	59
28) Benzo(a,h,i)perylene	28.16	276	6042	0.08	ng/ul#	26

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

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