

Data Path : Z:\HPCHEM1\BNA M\DATA\BM112015\
 Data File : BM003221.D
 Acq On : 20 Nov 2015 15:09
 Operator : UM/IZ
 Sample : G4345-03
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4J00

Quant Time: Nov 21 04:15:38 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-SIM-BM111815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Nov 21 03:26:56 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	12515	0.40	ng/ul	0.00
4) Naphthalene-d8	10.53	136	48309	0.40	ng/ul	0.00
8) Acenaphthene-d10	14.38	164	25922	0.40	ng/ul	0.00
12) Phenanthrene-d10	17.12	188	55878	0.40	ng/ul	0.00
18) Chrysene-d12	21.33	240	56308	0.40	ng/ul	0.00
22) Perylene-d12	23.60	264	53201	0.40	ng/ul	0.01
System Monitoring Compounds						
2) 1,4-Dioxane-d8	0.00	96	0d	0.00	ng/uL	
6) 2-Methylnaphthalene-d10	12.12	152	20990	0.32	ng/ul	0.00
16) Fluoranthene-d10	19.17	212	54395	0.40	ng/ul	0.00
Target Compounds				Ovalue		
5) Naphthalene	10.58	128	69639	0.56	ng/ul	99
7) 2-Methylnaphthalene	12.19	142	43565	0.54	ng/ul#	100
11) Fluorene	15.44	166	69253	0.70	ng/ul	99
13) Pentachlorophenol	16.78	266	15538	1.18	ng/ul	96
14) Phenanthrene	17.17	178	124590	0.77	ng/ul	99
17) Fluoranthene	19.19	202	120443	0.67	ng/ul	94
20) Benzo(a)anthracene	21.31	228	129697	0.79	ng/ul	94
24) Benzo(k)fluoranthene	22.95	252	122259	0.64	ng/ul#	92
26) Indeno(1,2,3-cd)pyrene	25.87	276	111193	0.52	ng/ul	96
28) Benzo(g,h,i)perylene	26.58	276	192593	1.02	ng/ul	98

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

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