

Data Path : Z:\HPCHEM1\BNA M\DATA\BM121615\
 Data File : BM003575.D
 Acq On : 16 Dec 2015 12:10
 Operator : UM/IZ
 Sample : G4734-09
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4KW7

Quant Time: Dec 17 14:39:36 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-SIM-BM121515.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 17 13:23:22 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.71	152	3599	0.40	ng/ul	0.00
4) Naphthalene-d8	10.50	136	14148	0.40	ng/ul	0.00
8) Acenaphthene-d10	14.37	164	7603	0.40	ng/ul	0.00
12) Phenanthrene-d10	17.11	188	18262	0.40	ng/ul	0.00
18) Chrysene-d12	21.31	240	20389	0.40	ng/ul	0.00
22) Perylene-d12	23.57	264	17406	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	0.00	96	0d	0.00	ng/uL	
6) 2-Methylnaphthalene-d10	12.10	152	6308	0.37	ng/ul	0.00
16) Fluoranthene-d10	19.14	212	16179	0.42	ng/ul	0.00
Target Compounds				Ovalue		
5) Naphthalene	10.55	128	14279	0.45	ng/ul	97
7) 2-Methylnaphthalene	12.16	142	20149	0.94	ng/ul	99
9) Acenaphthylene	14.07	152	16138	0.56	ng/ul	95
10) Acenaphthene	14.42	153	33064	1.41	ng/ul	88
13) Pentachlorophenol	16.78	266	687	0.26	ng/ul	94
14) Phenanthrene	17.14	178	28186	0.60	ng/ul#	92
17) Fluoranthene	19.17	202	55772	1.09	ng/ul	96
19) Pyrene	19.54	202	45691	0.83	ng/ul	92
20) Benzo(a)anthracene	21.29	228	27089	0.54	ng/ul	97
23) Benzo(b)fluoranthene	22.88	252	47276	0.80	ng/ul	97
24) Benzo(k)fluoranthene	22.92	252	54648	1.01	ng/ul	95
25) Benzo(a)pyrene	23.46	252	26980	0.54	ng/ul	96
26) Indeno(1,2,3-cd)pyrene	25.83	276	43658	0.75	ng/ul#	89
27) Dibenzo(a,h)anthracene	25.84	278	43418	0.93	ng/ul	99
28) Benzo(a,h,i)perylene	26.52	276	46974	0.93	ng/ul	99

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

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