

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM041715\
 Data File : BM001037.D
 Acq On : 17 Apr 2015 18:24
 Operator : TP/IZ
 Sample : G1749-03
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 X1N62

Quant Time: Apr 18 00:27:33 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-SIM-BM041715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 17 15:43:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	1553	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	5451	0.40	ng/ul	0.01
6) Acenaphthene-d10	14.24	164	2775	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.98	188	5997	0.40	ng/ul	0.00
16) Chrysene-d12	21.19	240	5205	0.40	ng/ul	0.00
20) Perylene-d12	23.37	264	4678	0.40	ng/ul	0.01

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.96	152	2180	0.31	ng/ul	0.00
14) Fluoranthene-d10	19.03	212	4475	0.34	ng/ul	0.01

Target Compounds				Qvalue		
3) Naphthalene	10.41	128	3401	0.24	ng/ul#	96
5) 2-Methylnaphthalene	12.03	142	2444	0.26	ng/ul	98
7) Acenaphthylene	13.94	152	9340	0.76	ng/ul	98
8) Acenaphthene	14.30	153	3142	0.33	ng/ul	96
9) Fluorene	15.30	166	4072	0.36	ng/ul	89
12) Phenanthrene	17.03	178	12541	0.72	ng/ul	94
13) Anthracene	17.12	178	3928	0.25	ng/ul	95
15) Fluoranthene	19.06	202	8328	0.43	ng/ul	91
17) Pyrene	19.42	202	9299	0.47	ng/ul	96
18) Benzo(a)anthracene	21.17	228	13114	0.80	ng/ul	95
19) Chrysene	21.22	228	8160	0.45	ng/ul	94
21) Benzo(b)fluoranthene	22.71	252	6141	0.36	ng/ul	99
23) Benzo(a)pyrene	23.26	252	8768	0.54	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.53	276	7536	0.42	ng/ul	99
25) Dibenzo(a,h)anthracene	25.54	278	6444	0.44	ng/ul	98
26) Benzo(g,h,i)perylene	26.19	276	7924	0.50	ng/ul	96

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

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