

Data Path : Z:\HPCHEM1\BNA_M\Data\BM061815\
 Data File : BM001733.D
 Acq On : 18 Jun 2015 10:30
 Operator : TP/IZ
 Sample : PB84013BL
 Misc : SBLK70
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK70

Quant Time: Jun 18 17:13:13 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-SIM-BM061815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 18 04:33:02 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	1002	0.40	ng/ul	0.00
2) Naphthalene-d8	10.46	136	3531	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.32	164	1645	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.06	188	3561	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	2731	0.40	ng/ul	0.00
20) Perylene-d12	23.47	264	2410	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.06	152	1395	0.32	ng/ul	0.01
14) Fluoranthene-d10	19.09	212	2507	0.33	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.50	128	55	0.01	ng/ul#	1
5) 2-Methylnaphthalene	12.12	142	28	0.00	ng/ul	85
7) Acenaphthylene	14.03	152	29	0.00	ng/ul#	1
8) Acenaphthene	14.37	153	20	0.00	ng/ul#	72
9) Fluorene	15.37	166	29	0.00	ng/ul#	77
11) Pentachlorophenol	16.72	266	15	0.02	ng/ul#	48
12) Phenanthrene	17.11	178	48	0.00	ng/ul#	1
13) Anthracene	17.20	178	30	0.00	ng/ul#	1
15) Fluoranthene	19.13	202	59	0.01	ng/ul#	1
17) Pyrene	19.49	202	60	0.01	ng/ul#	1
18) Benzo(a)anthracene	21.23	228	59	0.01	ng/ul#	11
19) Chrysene	21.28	228	83	0.01	ng/ul#	36
21) Benzo(b)fluoranthene	22.81	252	86	0.01	ng/ul#	1
22) Benzo(k)fluoranthene	22.85	252	73	0.01	ng/ul#	1
23) Benzo(a)pyrene	23.37	252	53	0.01	ng/ul#	1
24) Indeno(1,2,3-cd)pyrene	25.70	276	80	0.01	ng/ul	91
25) Dibenzo(a,h)anthracene	25.73	278	74	0.01	ng/ul#	1
26) Benzo(g,h,i)perylene	26.38	276	76	0.01	ng/ul#	1

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

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