

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM042715\
 Data File : BM001119.D
 Acq On : 27 Apr 2015 14:51
 Operator : TP/IZ
 Sample : G1998-03
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 X1Q38

Quant Time: Apr 28 01:44:38 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-SIM-BM042715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 27 14:06:45 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.54	152	1014	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	3480	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	1731	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	3940	0.40	ng/ul	0.00
16) Chrysene-d12	21.16	240	3732	0.40	ng/ul	0.00
20) Perylene-d12	23.32	264	3480	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.93	152	1612	0.36	ng/ul	0.00
14) Fluoranthene-d10	19.00	212	3463	0.39	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.37	128	8293	0.93	ng/ul	98
7) Acenaphthylene	13.91	152	6452	0.81	ng/ul	98
8) Acenaphthene	14.27	153	5020	0.82	ng/ul	99
9) Fluorene	15.26	166	6909	0.98	ng/ul	95
11) Pentachlorophenol	16.61	266	1057	1.44	ng/ul	98
12) Phenanthrene	17.00	178	9031	0.79	ng/ul	98
13) Anthracene	17.09	178	5304	0.51	ng/ul	99
17) Pyrene	19.38	202	22381	1.61	ng/ul	98
18) Benzo(a)anthracene	21.14	228	9615	0.80	ng/ul	97
19) Chrysene	21.19	228	14305	1.09	ng/ul	98
22) Benzo(k)fluoranthene	22.71	252	6573	0.50	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.46	276	10471	0.75	ng/ul#	91
26) Benzo(g,h,i)perylene	26.12	276	11685	0.96	ng/ul	97

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

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