

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM042715\
 Data File : BM001121.D
 Acq On : 27 Apr 2015 17:48
 Operator : TP/IZ
 Sample : G1998-03DL 2X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 X1Q38DL

Quant Time: Apr 28 01:52:32 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	1048	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	3516	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	1793	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	4062	0.40	ng/ul	0.00
16) Chrysene-d12	21.16	240	3681	0.40	ng/ul	0.00
20) Perylene-d12	23.33	264	3534	0.40	ng/ul	0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	11.93	152	872	0.19	ng/ul	0.00
14) Fluoranthene-d10	19.00	212	1803	0.19	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.37	128	4332	0.48	ng/ul	99
7) Acenaphthylene	13.91	152	3357	0.41	ng/ul	99
8) Acenaphthene	14.27	153	2605	0.41	ng/ul	98
9) Fluorene	15.26	166	3643	0.50	ng/ul	95
11) Pentachlorophenol	16.63	266	521	0.69	ng/ul	99
12) Phenanthrene	17.00	178	4822	0.41	ng/ul	98
13) Anthracene	17.09	178	2742	0.25	ng/ul	98
17) Pyrene	19.39	202	11479	0.84	ng/ul	98
18) Benzo(a)anthracene	21.14	228	4949	0.42	ng/ul	99
19) Chrysene	21.19	228	7337	0.57	ng/ul	100
22) Benzo(k)fluoranthene	22.71	252	3422	0.26	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.47	276	5489	0.39	ng/ul#	92
26) Benzo(g,h,i)perylene	26.12	276	6096	0.49	ng/ul	99

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

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