

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM021315\
 Data File : BM000518.D
 Acq On : 13 Feb 2015 10:56
 Operator : TP/IZ
 Sample : G1227-03
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 X1H34

Quant Time: Feb 13 12:05:04 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-SIM-BM021215.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 12 16:45:55 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.70	152	3992	0.40	ng/ul	0.00
2) Naphthalene-d8	10.49	136	13221	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.35	164	7837	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.09	188	15828	0.40	ng/ul	0.00
16) Chrysene-d12	21.29	240	13948	0.40	ng/ul	0.00
20) Perylene-d12	23.51	264	11739	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.09	152	5570	0.33	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	11619	0.35	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.54	128	26496	0.85	ng/ul	99
7) Acenaphthylene	14.08	152	21594	0.73	ng/ul	99
8) Acenaphthene	14.40	153	16673	0.74	ng/ul#	79
9) Fluorene	15.40	166	24466	0.89	ng/ul	97
12) Phenanthrene	17.13	178	31364	0.73	ng/ul	99
13) Anthracene	17.23	178	18296	0.47	ng/ul	98
17) Pyrene	19.53	202	75193	1.50	ng/ul	98
18) Benzo(a)anthracene	21.27	228	29252	0.74	ng/ul	99
19) Chrysene	21.32	228	44100	0.97	ng/ul	98
22) Benzo(k)fluoranthene	22.89	252	18580	0.47	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.75	276	27114	0.66	ng/ul#	94
26) Benzo(g,h,i)perylene	26.44	276	30568	0.84	ng/ul	98

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

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