

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM021315\
 Data File : BM000521.D
 Acq On : 13 Feb 2015 16:55
 Operator : TP/IZ
 Sample : G1249-12
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4FM0

Quant Time: Feb 14 05:11:12 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	3993	0.40	ng/ul	0.00
2) Naphthalene-d8	10.49	136	13022	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.35	164	7261	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.09	188	15071	0.40	ng/ul	0.00
16) Chrysene-d12	21.29	240	12780	0.40	ng/ul	0.00
20) Perylene-d12	23.51	264	10759	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.09	152	5764	0.35	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	11468	0.36	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.54	128	28046	0.92	ng/ul	98
7) Acenaphthylene	14.07	152	22325	0.82	ng/ul	99
8) Acenaphthene	14.40	153	17653	0.84	ng/ul#	80
9) Fluorene	15.40	166	25234	1.00	ng/ul	95
11) Pentachlorophenol	16.76	266	398	0.21	ng/ul	98
12) Phenanthrene	17.13	178	32168	0.79	ng/ul	98
13) Anthracene	17.23	178	18674	0.51	ng/ul	97
17) Pyrene	19.53	202	77125	1.68	ng/ul	98
18) Benzo(a)anthracene	21.27	228	29377	0.82	ng/ul	99
19) Chrysene	21.32	228	45446	1.09	ng/ul	97
22) Benzo(k)fluoranthene	22.89	252	19966	0.55	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.74	276	27733	0.73	ng/ul	95
26) Benzo(g,h,i)perylene	26.44	276	31404	0.94	ng/ul	98

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

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