

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM021315\
 Data File : BM000522.D
 Acq On : 13 Feb 2015 17:30
 Operator : TP/IZ
 Sample : G1249-12DL 2X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4FMDL

Quant Time: Feb 14 00:54:06 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	3762	0.40	ng/ul	0.00
2) Naphthalene-d8	10.49	136	11902	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.35	164	7114	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.09	188	14309	0.40	ng/ul	0.00
16) Chrysene-d12	21.29	240	12249	0.40	ng/ul	0.00
20) Perylene-d12	23.51	264	10576	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.09	152	2657	0.17	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	5636	0.19	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.54	128	13091	0.47	ng/ul	100
7) Acenaphthylene	14.07	152	10563	0.39	ng/ul	99
8) Acenaphthene	14.40	153	8374	0.41	ng/ul#	81
9) Fluorene	15.40	166	11741	0.47	ng/ul	98
11) Pentachlorophenol	16.75	266	509	0.28	ng/ul	93
12) Phenanthrene	17.13	178	15435	0.40	ng/ul	98
13) Anthracene	17.23	178	8775	0.25	ng/ul	98
17) Pyrene	19.51	202	37038	0.84	ng/ul	92
18) Benzo(a)anthracene	21.27	228	14037	0.41	ng/ul	97
19) Chrysene	21.32	228	22912	0.58	ng/ul	98
22) Benzo(k)fluoranthene	22.89	252	9800	0.28	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.74	276	13433	0.36	ng/ul#	93
26) Benzo(g,h,i)perylene	26.43	276	15170	0.46	ng/ul	100

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

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