

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM020515\
 Data File : BM000433.D
 Acq On : 05 Feb 2015 19:01
 Operator : TP/IZ
 Sample : G1147-03DL 2X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 X1G08DL

Quant Time: Feb 06 06:39:26 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-SIM-BM020515.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 05 14:01:48 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	1988	0.40	ng/ul	0.00
2) Naphthalene-d8	10.51	136	6521	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.38	164	4115	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.12	188	8035	0.40	ng/ul	0.00
16) Chrysene-d12	21.31	240	7215	0.40	ng/ul	0.00
20) Perylene-d12	23.55	264	6563	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.11	152	3043	0.35	ng/ul	0.00
14) Fluoranthene-d10	19.15	212	6445	0.37	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.56	128	8013	0.48	ng/ul	100
7) Acenaphthylene	14.10	152	7013	0.41	ng/ul	99
8) Acenaphthene	14.43	153	5227	0.41	ng/ul	98
9) Fluorene	15.43	166	7520	0.50	ng/ul	99
11) Pentachlorophenol	16.78	266	1040	0.56	ng/ul	96
12) Phenanthrene	17.16	178	9484	0.41	ng/ul	99
13) Anthracene	17.26	178	5591	0.26	ng/ul	99
17) Pyrene	19.55	202	23079	0.83	ng/ul	98
18) Benzo(a)anthracene	21.29	228	9845	0.43	ng/ul	97
19) Chrysene	21.34	228	14703	0.60	ng/ul	98
22) Benzo(k)fluoranthene	22.93	252	6539	0.28	ng/ul#	94
24) Indeno(1,2,3-cd)pyrene	25.82	276	9828	0.39	ng/ul	94
26) Benzo(g,h,i)perylene	26.50	276	10945	0.50	ng/ul	98

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

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