

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM020515\
 Data File : BM000430.D
 Acq On : 05 Feb 2015 15:38
 Operator : TP/IZ
 Sample : G1147-03
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 X1G08

Quant Time: Feb 06 08:58:50 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	1943	0.40	ng/ul	0.00
2) Naphthalene-d8	10.51	136	6400	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.38	164	4123	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.12	188	7921	0.40	ng/ul	0.00
16) Chrysene-d12	21.31	240	7192	0.40	ng/ul	0.00
20) Perylene-d12	23.56	264	6230	0.40	ng/ul	0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.11	152	5970	0.70	ng/ul	0.00
14) Fluoranthene-d10	19.15	212	12965	0.75	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.56	128	15780	0.97	ng/ul	99
7) Acenaphthylene	14.10	152	13333	0.78	ng/ul	98
8) Acenaphthene	14.43	153	10240	0.80	ng/ul	99
9) Fluorene	15.43	166	15240	1.00	ng/ul	97
11) Pentachlorophenol	16.78	266	2511	1.36	ng/ul	100
12) Phenanthrene	17.16	178	19301	0.85	ng/ul	99
13) Anthracene	17.26	178	11137	0.53	ng/ul	100
17) Pyrene	19.55	202	46823	1.70	ng/ul	98
18) Benzo(a)anthracene	21.29	228	18828	0.83	ng/ul	99
19) Chrysene	21.35	228	28855	1.18	ng/ul	90
22) Benzo(k)fluoranthene	22.93	252	12340	0.56	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.82	276	18885	0.79	ng/ul#	94
26) Benzo(g,h,i)perylene	26.51	276	21272	1.02	ng/ul	98

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

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