

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE010715\
 Data File : BE089346.D
 Acq On : 8 Jan 2015 1:25
 Operator : TP/IZ
 Sample : G1001-05
 Misc : LOD-WATER-0.5PPM-SIM
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 LOD-WATER2015-01

Quant Time: Jan 08 15:37:26 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-PAH-BASE-BE010715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 08 15:21:44 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	18164	5.00	ng	0.00
2) Naphthalene-d8	8.51	136	71299	5.00	ng	0.00
6) Acenaphthene-d10	10.65	164	34049	5.00	ng	0.00
11) Phenanthrene-d10	12.72	188	51426	5.00	ng	0.00
15) Chrysene-d12	18.07	240	36782	5.00	ng	0.00
21) Perylene-d12	21.12	264	33107	5.00	ng	0.00

System Monitoring Compounds						
3) Nitrobenzene-d5	7.64	82	35880	7.24	ng	0.00
7) 2-Fluorobiphenyl	9.85	172	59885	6.45	ng	0.00
17) Terphenyl-d14	15.85	244	46138	7.08	ng	0.00

Target Compounds						Qvalue
4) Naphthalene	8.53	128	1218	0.08	ng	# 91
5) 2-Methylnaphthalene	9.39	142	745	0.07	ng	# 95
8) Acenaphthylene	10.48	152	4227	0.07	ng	# 85
9) Acenaphthene	10.69	154	777	0.09	ng	# 89
10) Fluorene	11.34	166	722	0.07	ng	# 99
12) Phenanthrene	12.75	178	4230	0.08	ng	# 90
13) Anthracene	12.84	178	3846	0.07	ng	# 85
14) Fluoranthene	14.96	202	887	0.07	ng	# 85
16) Pyrene	15.41	202	919	0.08	ng	# 87
18) Benzo(a)anthracene	18.05	228	3204	0.09	ng	# 85
19) Chrysene	18.12	228	3006	0.08	ng	# 84
20) Indeno(1,2,3-cd)pyrene	23.24	276	2502	0.07	ng	# 76
22) Benzo(b)fluoranthene	20.38	252	527	0.07	ng	# 50
23) Benzo(k)fluoranthene	20.43	252	735	0.08	ng	# 65
24) Benzo(a)pyrene	21.01	252	532	0.07	ng	# 65
25) Dibenzo(a,h)anthracene	23.31	278	449	0.07	ng	# 46
26) Benzo(g,h,i)perylene	23.82	276	2318	0.07	ng	# 62

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

