

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE072115\
 Data File : BE090191.D
 Acq On : 21 Jul 2015 00:58
 Operator : TP/IZ
 Sample : MDL-06-W
 Misc : MDL-WATER-SIM2.2-0.07PPM
 ALS Vial : 14 Sample Multiplier: 1

Quant Time: Jul 21 05:54:03 2015

Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM02.2-EPA-SIM-BE072115.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Tue Jul 21 05:15:45 2015

Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	797	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	3761	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.21	164	2258	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.94	188	5950	0.40	ng/ul	0.00
16) Chrysene-d12	21.12	240	5098	0.40	ng/ul	0.00
20) Perylene-d12	23.43	264	3624	0.40	ng/ul	0.00

System Monitoring Compounds

4) 2-Methylnaphthalene-d10	11.96	152	2305	0.39	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	6686	0.37	ng/ul	0.00

Target Compounds

						Qvalue
3) Naphthalene	10.41	128	558	0.06	ng/ul#	73
5) 2-Methylnaphthalene	12.03	142	364	0.05	ng/ul	94
7) Acenaphthylene	13.93	152	2584	0.05	ng/ul#	84
8) Acenaphthene	14.27	153	1868	0.05	ng/ul	91
9) Fluorene	15.27	166	531	0.05	ng/ul#	93
12) Phenanthrene	16.99	178	3653	0.05	ng/ul#	86
13) Anthracene	17.07	178	3641	0.05	ng/ul#	82
15) Fluoranthene	19.00	202	4521	0.06	ng/ul	86
17) Pyrene	19.36	202	4442	0.07	ng/ul#	89
18) Benzo(a)anthracene	21.10	228	489	0.04	ng/ul#	77
19) Chrysene	21.15	228	1002m	0.06	ng/ul	
21) Benzo(b)fluoranthene	22.72	252	281	0.03	ng/ul#	40
22) Benzo(k)fluoranthene	22.77	252	585m	0.04	ng/ul	
23) Benzo(a)pyrene	23.33	252	321	0.03	ng/ul#	29
24) Indeno(1,2,3-cd)pyrene	25.83	276	1115	0.02	ng/ul	94
25) Dibenzo(a,h)anthracene	25.85	278	159	0.02	ng/ul#	20
26) Benzo(g,h,i)perylene	26.56	276	1292	0.03	ng/ul#	41

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

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