

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE072115\
 Data File : BE090199.D
 Acq On : 21 Jul 2015 7:05
 Operator : TP/IZ
 Sample : MDL-06-W
 Misc : MDL-WATER-SIM2.2-0.15PPM
 ALS Vial : 22 Sample Multiplier: 1

Quant Time: Jul 21 09:07:07 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	800	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	3746	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.22	164	2212	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.94	188	5797	0.40	ng/ul	0.00
16) Chrysene-d12	21.12	240	4113	0.40	ng/ul	0.00
20) Perylene-d12	23.43	264	2931	0.40	ng/ul	0.00

System Monitoring Compounds

4) 2-Methylnaphthalene-d10	11.96	152	2280	0.38	ng/ul	0.00
14) Fluoranthene-d10	18.97	212	6378	0.36	ng/ul	0.00

Target Compounds

						Qvalue
3) Naphthalene	10.42	128	1077	0.11	ng/ul#	87
5) 2-Methylnaphthalene	12.03	142	723	0.11	ng/ul	97
7) Acenaphthylene	13.93	152	4989	0.11	ng/ul	94
8) Acenaphthene	14.27	153	3654	0.11	ng/ul	98
9) Fluorene	15.27	166	1047	0.11	ng/ul#	97
12) Phenanthrene	16.99	178	7084	0.11	ng/ul	93
13) Anthracene	17.07	178	7044	0.11	ng/ul	93
15) Fluoranthene	19.00	202	8570	0.11	ng/ul	93
17) Pyrene	19.36	202	8538	0.16	ng/ul	96
18) Benzo(a)anthracene	21.10	228	656	0.06	ng/ul#	83
19) Chrysene	21.15	228	2038m	0.16	ng/ul	
21) Benzo(b)fluoranthene	22.72	252	371	0.05	ng/ul#	48
22) Benzo(k)fluoranthene	22.77	252	889m	0.08	ng/ul	
23) Benzo(a)pyrene	23.32	252	428	0.05	ng/ul#	50
24) Indeno(1,2,3-cd)pyrene	25.82	276	1408	0.04	ng/ul	99
25) Dibenzo(a,h)anthracene	25.84	278	205	0.03	ng/ul#	16
26) Benzo(g,h,i)perylene	26.56	276	1999	0.06	ng/ul#	52

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

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