

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE073015\
 Data File : BE090299.D
 Acq On : 30 Jul 2015 15:48
 Operator : TP/IZ
 Sample : MDL-07-W
 Misc : MDL-SIM2.2-WATER-0.15PPM
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 MDL-07-W

Manual Integrations
 APPROVED

MMDadoda
 7/31/2015 3:07:46 PM

Quant Time: Jul 31 06:42:50 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM02.2-EPA-SIM-BE072815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 31 06:10:41 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	920	0.40	ng/ul	0.00
2) Naphthalene-d8	10.36	136	3704	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.21	164	2174	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.94	188	4771	0.40	ng/ul	0.00
16) Chrysene-d12	21.12	240	2791	0.40	ng/ul	0.00
20) Perylene-d12	23.43	264	1648	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.95	152	2470	0.43	ng/ul	0.01
14) Fluoranthene-d10	18.97	212	5319	0.40	ng/ul	0.00
Target Compounds						
					Qvalue	
3) Naphthalene	10.41	128	1211	0.12	ng/ul#	86
5) 2-Methylnaphthalene	12.03	142	743	0.12	ng/ul	96
7) Acenaphthylene	13.93	152	5623	0.12	ng/ul#	92
8) Acenaphthene	14.27	153	4053	0.12	ng/ul	98
9) Fluorene	15.27	166	1043	0.12	ng/ul#	96
11) Pentachlorophenol	16.62	266	49	0.16	ng/ul#	89
12) Phenanthrene	16.99	178	6002	0.12	ng/ul#	92
13) Anthracene	17.07	178	6224	0.12	ng/ul#	87
15) Fluoranthene	19.00	202	7218	0.12	ng/ul	95
17) Pyrene	19.36	202	7383	0.14	ng/ul#	91
18) Benzo(a)anthracene	21.10	228	444	0.11	ng/ul#	78
19) Chrysene	21.15	228	1647m	0.14	ng/ul	
21) Benzo(b)fluoranthene	22.73	252	229	0.09	ng/ul#	54
22) Benzo(k)fluoranthene	22.77	252	917m	0.12	ng/ul	
23) Benzo(a)pyrene	23.33	252	441m	0.13	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.81	276	674	0.06	ng/ul#	81
25) Dibenzo(a,h)anthracene	25.85	278	133m	0.11	ng/ul	
26) Benzo(g,h,i)perylene	26.56	276	1405	0.10	ng/ul#	50

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

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