

Data Path : Z:\HPCHEM1\BNA M\DATA\BM010715\
 Data File : BM000128.D
 Acq On : 07 Jan 2015 14:41
 Operator : TP/IZ
 Sample : MDL-02-W
 Misc : MDL-WATER-0.07PPM-SIM
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MDL-02-W

Manual Integrations
 APPROVED

TEJASKUMAR
 1/9/2015 5:44:49 PM

Quant Time: Jan 08 11:53:24 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-SIM-BM010715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 08 10:41:47 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	2889	0.40	ng/ul	0.00
2) Naphthalene-d8	10.64	136	9614	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	5634	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.23	188	10870	0.40	ng/ul	0.02
16) Chrysene-d12	21.41	240	9032	0.40	ng/ul	0.00
20) Perylene-d12	23.70	264	8212	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	4535	0.35	ng/ul	0.00
14) Fluoranthene-d10	19.25	212	8182	0.35	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.68	128	1150	0.04	ng/ul#	85
5) 2-Methylnaphthalene	12.29	142	765	0.04	ng/ul	98
7) Acenaphthylene	14.20	152	1214	0.04	ng/ul#	82
8) Acenaphthene	14.54	153	780	0.04	ng/ul	95
9) Fluorene	15.54	166	946	0.05	ng/ul#	91
11) Pentachlorophenol	16.90	266	41	0.02	ng/ul#	70
12) Phenanthrene	17.26	178	1503	0.04	ng/ul#	82
13) Anthracene	17.35	178	1358	0.04	ng/ul#	81
15) Fluoranthene	19.28	202	1766	0.05	ng/ul	87
17) Pyrene	19.65	202	1822	0.05	ng/ul#	92
18) Benzo(a)anthracene	21.39	228	1445	0.05	ng/ul	91
19) Chrysene	21.44	228	1631m	0.05	ng/ul	
21) Benzo(b)fluoranthene	23.00	252	1488	0.05	ng/ul#	58
22) Benzo(k)fluoranthene	23.05	252	1424	0.05	ng/ul#	57
23) Benzo(a)pyrene	23.61	252	1324	0.05	ng/ul#	44
24) Indeno(1,2,3-cd)pyrene	26.06	276	1532	0.05	ng/ul	98
25) Dibenzo(a,h)anthracene	26.08	278	1270	0.05	ng/ul#	55
26) Benzo(a,h,i)perylene	26.76	276	1387	0.05	ng/ul#	74

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

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