

Data Path : Z:\HPCHEM1\BNA M\DATA\BM010715\
 Data File : BM000132.D
 Acq On : 07 Jan 2015 17:03
 Operator : TP/IZ
 Sample : MDL-06-W
 Misc : MDL-WATER-0.07PPM-SIM
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MDL-06-W

Quant Time: Jan 08 11:57:48 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	3011	0.40	ng/ul	0.00
2) Naphthalene-d8	10.64	136	9837	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	5720	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.23	188	10652	0.40	ng/ul	0.02
16) Chrysene-d12	21.41	240	9060	0.40	ng/ul	0.00
20) Perylene-d12	23.70	264	8198	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.22	152	4689	0.36	ng/ul	0.00
14) Fluoranthene-d10	19.25	212	7951	0.35	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
3) Naphthalene	10.68	128	1441	0.05	ng/ul#	88
5) 2-Methylnaphthalene	12.30	142	931	0.05	ng/ul	99
7) Acenaphthylene	14.20	152	1438	0.05	ng/ul#	85
8) Acenaphthene	14.54	153	940	0.05	ng/ul	94
9) Fluorene	15.52	166	1096	0.05	ng/ul#	86
11) Pentachlorophenol	16.89	266	69	0.04	ng/ul#	59
12) Phenanthrene	17.26	178	1747	0.05	ng/ul#	84
13) Anthracene	17.35	178	1609	0.05	ng/ul#	82
15) Fluoranthene	19.28	202	2017	0.06	ng/ul	88
17) Pyrene	19.65	202	2061	0.05	ng/ul#	92
18) Benzo(a)anthracene	21.39	228	1674	0.06	ng/ul#	91
19) Chrysene	21.44	228	1823	0.06	ng/ul	95
21) Benzo(b)fluoranthene	23.01	252	1701	0.05	ng/ul#	62
22) Benzo(k)fluoranthene	23.05	252	1803	0.06	ng/ul#	66
23) Benzo(a)pyrene	23.61	252	1495	0.05	ng/ul#	52
24) Indeno(1,2,3-cd)pyrene	26.05	276	1773	0.05	ng/ul#	78
25) Dibenzo(a,h)anthracene	26.07	278	1474	0.06	ng/ul#	56
26) Benzo(g,h,i)perylene	26.75	276	1633	0.06	ng/ul#	73

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

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