

Data Path : Z:\HPCHEM1\BNA M\DATA\BM010715\
 Data File : BM000127.D
 Acq On : 07 Jan 2015 14:05
 Operator : TP/IZ
 Sample : MDL-01-W
 Misc : MDL-WATER-0.07PPM-SIM
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MDL-01-W

Manual Integrations
 APPROVED

TEJASKUMAR
 1/9/2015 5:44:51 PM

Quant Time: Jan 08 11:52:02 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	3303	0.40	ng/ul	0.00
2) Naphthalene-d8	10.64	136	11043	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	6694	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.23	188	11478	0.40	ng/ul	0.02
16) Chrysene-d12	21.41	240	8297	0.40	ng/ul	0.00
20) Perylene-d12	23.71	264	7393	0.40	ng/ul	0.01

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	5421	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.25	212	8069	0.33	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.68	128	1446	0.05	ng/ul#	89
5) 2-Methylnaphthalene	12.29	142	946	0.05	ng/ul	99
7) Acenaphthylene	14.20	152	1527	0.05	ng/ul#	89
8) Acenaphthene	14.54	153	986	0.05	ng/ul	94
9) Fluorene	15.54	166	1135	0.05	ng/ul#	92
11) Pentachlorophenol	16.88	266	32	0.02	ng/ul#	84
12) Phenanthrene	17.26	178	1675	0.05	ng/ul#	86
13) Anthracene	17.35	178	1544	0.05	ng/ul#	82
15) Fluoranthene	19.28	202	1857	0.05	ng/ul	89
17) Pyrene	19.65	202	1871	0.05	ng/ul#	93
18) Benzo(a)anthracene	21.39	228	1359	0.05	ng/ul#	89
19) Chrysene	21.44	228	1572m	0.05	ng/ul	
21) Benzo(b)fluoranthene	23.01	252	1364	0.05	ng/ul#	56
22) Benzo(k)fluoranthene	23.05	252	1428	0.05	ng/ul#	56
23) Benzo(a)pyrene	23.61	252	1445	0.05	ng/ul#	44
24) Indeno(1,2,3-cd)pyrene	26.04	276	1547	0.05	ng/ul#	78
25) Dibenzo(a,h)anthracene	26.08	278	1242	0.05	ng/ul#	40
26) Benzo(g,h,i)perylene	26.75	276	1377	0.05	ng/ul#	73

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

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