

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM033115\
 Data File : BM000782.D
 Acq On : 01 Apr 2015 11:57
 Operator : TP/IZ
 Sample : MDL-02-W
 Misc : MDL-WATER-0.1/0.2PPM
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MDL-02-W

Manual Integrations
 APPROVED

apatel
 4/1/2015 6:18:16 PM

Quant Time: Apr 01 15:30:23 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIMPAH-BM033115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 01 04:17:30 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	36757	5.00	ng	0.00
4) Naphthalene-d8	10.40	136	125867	5.00	ng	0.00
8) Acenaphthene-d10	14.28	164	61097	5.00	ng	0.00
13) Phenanthrene-d10	17.01	188	133351	5.00	ng	0.00
19) Chrysene-d12	21.21	240	111031	5.00	ng	0.00
25) Perylene-d12	23.40	264	103364	5.00	ng	0.00
System Monitoring Compounds						
5) Nitrobenzene-d5	8.77	82	63699	8.24	ng	0.00
9) 2-Fluorobiphenyl	12.89	172	192024	10.13	ng	0.00
21) Terphenyl-d14	19.66	244	145828	10.91	ng	0.00
Target Compounds						Qvalue
6) Naphthalene	10.45	128	2329	0.08	ng	# 93
7) 2-Methylnaphthalene	12.08	142	1372	0.07	ng	96
10) Acenaphthylene	13.98	152	1785	0.07	ng	98
11) Acenaphthene	14.33	154	1418	0.08	ng	99
12) Fluorene	15.32	166	1623	0.08	ng	98
14) Hexachlorobenzene	16.33	284	613	0.08	ng	# 76
16) Phenanthrene	17.05	178	2915m	0.08	ng	
17) Anthracene	17.15	178	2288	0.07	ng	98
18) Fluoranthene	19.09	202	2674	0.07	ng	99
20) Pyrene	19.45	202	2794	0.08	ng	99
22) Benzo(a)anthracene	21.20	228	2677	0.08	ng	97
23) Chrysene	21.24	228	2957	0.09	ng	97
24) Indeno(1,2,3-cd)pyrene	25.58	276	2657	0.08	ng	99
26) Benzo(b)fluoranthene	22.74	252	2494	0.08	ng	# 81
27) Benzo(k)fluoranthene	22.79	252	2511	0.08	ng	# 80
28) Benzo(a)pyrene	23.30	252	2193	0.08	ng	# 73
29) Dibenzo(a,h)anthracene	25.60	278	2156	0.08	ng	# 80
30) Benzo(g,h,i)perylene	26.26	276	2480	0.08	ng	# 92

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

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