

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM033115\
 Data File : BM000784.D
 Acq On : 01 Apr 2015 13:09
 Operator : TP/IZ
 Sample : MDL-04-W
 Misc : MDL-WATER-0.1/0.2PPM
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MDL-04-W

Manual Integrations
 APPROVED

apatel
 4/1/2015 6:18:20 PM

Quant Time: Apr 01 15:31:56 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	7.62	152	34136	5.00	ng	0.00
4) Naphthalene-d8	10.40	136	114536	5.00	ng	0.00
8) Acenaphthene-d10	14.28	164	56501	5.00	ng	0.00
13) Phenanthrene-d10	17.01	188	125368	5.00	ng	0.00
19) Chrysene-d12	21.21	240	104082	5.00	ng	0.00
25) Perylene-d12	23.40	264	96525	5.00	ng	0.00
System Monitoring Compounds						
5) Nitrobenzene-d5	8.77	82	56376	8.02	ng	0.00
9) 2-Fluorobiphenyl	12.89	172	171284	9.77	ng	0.00
21) Terphenyl-d14	19.66	244	136205	10.88	ng	0.00
Target Compounds						Qvalue
6) Naphthalene	10.45	128	2100	0.08	ng	# 91
7) 2-Methylnaphthalene	12.08	142	1199	0.07	ng	96
10) Acenaphthylene	13.99	152	1564	0.06	ng	97
11) Acenaphthene	14.33	154	1259	0.07	ng	98
12) Fluorene	15.32	166	1465	0.08	ng	100
14) Hexachlorobenzene	16.33	284	577	0.08	ng	# 70
16) Phenanthrene	17.05	178	2668m	0.08	ng	
17) Anthracene	17.15	178	2155	0.07	ng	99
18) Fluoranthene	19.09	202	2536	0.07	ng	100
20) Pyrene	19.45	202	2616	0.08	ng	99
22) Benzo(a)anthracene	21.20	228	2465	0.08	ng	97
23) Chrysene	21.24	228	2765	0.09	ng	96
24) Indeno(1,2,3-cd)pyrene	25.59	276	2456	0.08	ng	98
26) Benzo(b)fluoranthene	22.74	252	2270	0.07	ng	# 77
27) Benzo(k)fluoranthene	22.79	252	2371	0.08	ng	# 78
28) Benzo(a)pyrene	23.30	252	2015	0.08	ng	# 71
29) Dibenzo(a,h)anthracene	25.60	278	1919	0.07	ng	# 78
30) Benzo(g,h,i)perylene	26.26	276	2271	0.08	ng	# 91

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

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