

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM033115\
 Data File : BM000781.D
 Acq On : 01 Apr 2015 11:22
 Operator : TP/IZ
 Sample : MDL-01-W
 Misc : MDL-WATER-0.1/0.2PPM
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MDL-01-W

Manual Integrations
 APPROVED

apatel
 4/1/2015 6:18:12 PM

Quant Time: Apr 01 15:29:14 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	35908	5.00	ng	0.00
4) Naphthalene-d8	10.40	136	123883	5.00	ng	0.00
8) Acenaphthene-d10	14.28	164	59669	5.00	ng	0.00
13) Phenanthrene-d10	17.01	188	130495	5.00	ng	0.00
19) Chrysene-d12	21.21	240	107525	5.00	ng	0.00
25) Perylene-d12	23.40	264	100214	5.00	ng	0.00
System Monitoring Compounds						
5) Nitrobenzene-d5	8.77	82	30224	4.07	ng	0.00
9) 2-Fluorobiphenyl	12.89	172	110645	5.98	ng	0.00
21) Terphenyl-d14	19.66	244	137816	10.65	ng	0.00
Target Compounds						Qvalue
6) Naphthalene	10.45	128	1219	0.04	ng	# 84
7) 2-Methylnaphthalene	12.08	142	753	0.04	ng	# 96
10) Acenaphthylene	13.98	152	1136	0.04	ng	98
11) Acenaphthene	14.33	154	910	0.05	ng	96
12) Fluorene	15.32	166	1158	0.06	ng	98
14) Hexachlorobenzene	16.33	284	489	0.06	ng	# 97
16) Phenanthrene	17.05	178	2455m	0.07	ng	
17) Anthracene	17.15	178	1952	0.07	ng	99
18) Fluoranthene	19.09	202	2472	0.07	ng	99
20) Pyrene	19.45	202	2625	0.07	ng	99
22) Benzo(a)anthracene	21.20	228	2507	0.08	ng	96
23) Chrysene	21.24	228	2878	0.09	ng	97
24) Indeno(1,2,3-cd)pyrene	25.59	276	2581	0.08	ng	98
26) Benzo(b)fluoranthene	22.74	252	2383	0.08	ng	# 80
27) Benzo(k)fluoranthene	22.79	252	2490	0.08	ng	# 76
28) Benzo(a)pyrene	23.30	252	2105	0.08	ng	# 71
29) Dibenzo(a,h)anthracene	25.60	278	2046	0.08	ng	# 78
30) Benzo(g,h,i)perylene	26.26	276	2373	0.08	ng	# 91

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

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