

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM033115\
 Data File : BM000786.D
 Acq On : 01 Apr 2015 14:20
 Operator : TP/IZ
 Sample : MDL-06-W
 Misc : MDL-WATER-0.1/0.2PPM
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MDL-06-W

Manual Integrations
 APPROVED

apatel
 4/1/2015 6:18:31 PM

Quant Time: Apr 01 15:33:29 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	35264	5.00	ng	0.00
4) Naphthalene-d8	10.40	136	118796	5.00	ng	0.00
8) Acenaphthene-d10	14.28	164	57352	5.00	ng	0.00
13) Phenanthrene-d10	17.01	188	128908	5.00	ng	0.00
19) Chrysene-d12	21.21	240	109617	5.00	ng	0.00
25) Perylene-d12	23.40	264	100767	5.00	ng	0.00
System Monitoring Compounds						
5) Nitrobenzene-d5	8.77	82	55495	7.62	ng	0.00
9) 2-Fluorobiphenyl	12.89	172	167626	9.42	ng	0.00
21) Terphenyl-d14	19.66	244	133184	10.10	ng	0.00
Target Compounds						Qvalue
6) Naphthalene	10.45	128	2302	0.08	ng	# 92
7) 2-Methylnaphthalene	12.08	142	1382	0.08	ng	95
10) Acenaphthylene	13.98	152	1791	0.07	ng	100
11) Acenaphthene	14.33	154	1358	0.08	ng	97
12) Fluorene	15.32	166	1610	0.08	ng	98
14) Hexachlorobenzene	16.33	284	625	0.08	ng	# 99
16) Phenanthrene	17.05	178	2865m	0.08	ng	
17) Anthracene	17.15	178	2372	0.08	ng	98
18) Fluoranthene	19.09	202	2786	0.08	ng	98
20) Pyrene	19.45	202	2885	0.08	ng	98
22) Benzo(a)anthracene	21.20	228	2715	0.09	ng	97
23) Chrysene	21.24	228	3064	0.09	ng	98
24) Indeno(1,2,3-cd)pyrene	25.59	276	2718	0.08	ng	100
26) Benzo(b)fluoranthene	22.74	252	2512	0.08	ng	# 80
27) Benzo(k)fluoranthene	22.79	252	2659	0.09	ng	# 79
28) Benzo(a)pyrene	23.30	252	2229	0.08	ng	# 75
29) Dibenzo(a,h)anthracene	25.60	278	2143	0.08	ng	# 80
30) Benzo(g,h,i)perylene	26.26	276	2539	0.09	ng	# 91

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

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