

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM033115\
 Data File : BM000787.D
 Acq On : 01 Apr 2015 14:55
 Operator : TP/IZ
 Sample : MDL-07-W
 Misc : MDL-WATER-0.1/0.2PPM
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MDL-07-W

Manual Integrations
 APPROVED

apatel
 4/1/2015 6:18:35 PM

Quant Time: Apr 01 15:49:36 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	37572	5.00	ng	0.00
4) Naphthalene-d8	10.40	136	125635	5.00	ng	0.00
8) Acenaphthene-d10	14.28	164	60572	5.00	ng	0.00
13) Phenanthrene-d10	17.01	188	136670	5.00	ng	0.00
19) Chrysene-d12	21.21	240	113674	5.00	ng	0.00
25) Perylene-d12	23.40	264	105211	5.00	ng	0.00
System Monitoring Compounds						
5) Nitrobenzene-d5	8.77	82	58319	7.57	ng	0.00
9) 2-Fluorobiphenyl	12.89	172	174360	9.28	ng	0.00
21) Terphenyl-d14	19.66	244	139870	10.23	ng	0.00
Target Compounds						Qvalue
6) Naphthalene	10.45	128	2129	0.07	ng	# 92
7) 2-Methylnaphthalene	12.08	142	1242	0.07	ng	96
10) Acenaphthylene	13.99	152	1617	0.06	ng	99
11) Acenaphthene	14.33	154	1263	0.07	ng	97
12) Fluorene	15.32	166	1471	0.07	ng	97
14) Hexachlorobenzene	16.33	284	587	0.07	ng	# 96
16) Phenanthrene	17.05	178	2681m	0.07	ng	
17) Anthracene	17.15	178	2169	0.07	ng	99
18) Fluoranthene	19.09	202	2576	0.07	ng	99
20) Pyrene	19.45	202	2657	0.07	ng	99
22) Benzo(a)anthracene	21.19	228	2588	0.08	ng	# 91
23) Chrysene	21.24	228	2806	0.08	ng	96
24) Indeno(1,2,3-cd)pyrene	25.59	276	2519	0.07	ng	99
26) Benzo(b)fluoranthene	22.74	252	2332	0.07	ng	# 79
27) Benzo(k)fluoranthene	22.79	252	2406	0.08	ng	# 78
28) Benzo(a)pyrene	23.30	252	2087	0.07	ng	# 71
29) Dibenzo(a,h)anthracene	25.61	278	1998	0.07	ng	# 83
30) Benzo(g,h,i)perylene	26.26	276	2339	0.08	ng	# 91

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

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