

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM033115\
 Data File : BM000783.D
 Acq On : 01 Apr 2015 12:33
 Operator : TP/IZ
 Sample : MDL-03-W
 Misc : MDL-WATER-0.1/0.2PPM
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MDL-03-W

Quant Time: Apr 01 15:31:01 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	35890	5.00	ng	0.00
4) Naphthalene-d8	10.40	136	122517	5.00	ng	0.00
8) Acenaphthene-d10	14.28	164	58865	5.00	ng	0.00
13) Phenanthrene-d10	17.01	188	128903	5.00	ng	0.00
19) Chrysene-d12	21.22	240	104628	5.00	ng	0.01
25) Perylene-d12	23.40	264	97772	5.00	ng	0.00

System Monitoring Compounds						
5) Nitrobenzene-d5	8.77	82	58593	7.80	ng	0.00
9) 2-Fluorobiphenyl	12.89	172	177047	9.70	ng	0.00
21) Terphenyl-d14	19.66	244	134292	10.67	ng	0.00

Target Compounds						Qvalue
6) Naphthalene	10.45	128	2116	0.07	ng	# 91
7) 2-Methylnaphthalene	12.08	142	1252	0.07	ng	96
10) Acenaphthylene	13.98	152	1619	0.06	ng	99
11) Acenaphthene	14.33	154	1273	0.07	ng	98
12) Fluorene	15.32	166	1472	0.07	ng	99
14) Hexachlorobenzene	16.33	284	587	0.08	ng	# 95
16) Phenanthrene	17.05	178	2673	0.08	ng	96
17) Anthracene	17.15	178	2084	0.07	ng	98
18) Fluoranthene	19.09	202	2411	0.07	ng	99
20) Pyrene	19.45	202	2517	0.07	ng	98
22) Benzo(a)anthracene	21.20	228	2353	0.08	ng	96
23) Chrysene	21.25	228	2657	0.08	ng	# 90
24) Indeno(1,2,3-cd)pyrene	25.59	276	2414	0.07	ng	97
26) Benzo(b)fluoranthene	22.74	252	2149	0.07	ng	# 77
27) Benzo(k)fluoranthene	22.79	252	2413	0.08	ng	# 75
28) Benzo(a)pyrene	23.31	252	1939	0.07	ng	# 70
29) Dibenzo(a,h)anthracene	25.61	278	1893	0.07	ng	# 79
30) Benzo(g,h,i)perylene	26.26	276	2232	0.08	ng	# 87

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

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