

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE021015\
 Data File : BE089564.D
 Acq On : 11 Feb 2015 00:57
 Operator : TP/IZ
 Sample : MDL-03-W
 Misc : MDL-WATER-0.1/0.2
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 MDL-03-W

Quant Time: Feb 11 02:43:36 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIMPAH-BE021015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 11 01:27:13 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.42	152	54653	5.00	ng	0.00
2) Naphthalene-d8	12.32	136	211632	5.00	ng	0.00
6) Acenaphthene-d10	16.50	164	100680	5.00	ng	0.00
11) Phenanthrene-d10	19.84	188	180393	5.00	ng	0.00
16) Chrysene-d12	23.71	240	203528	5.00	ng	0.00
22) Perylene-d12	25.40	264	198216	5.00	ng	0.00

System Monitoring Compounds						
3) Nitrobenzene-d5	10.68	82	107210	7.91	ng	0.00
7) 2-Fluorobiphenyl	14.95	172	166047	6.09	ng	0.00
18) Terphenyl-d14	22.36	244	206481	8.14	ng	0.00

Target Compounds						Qvalue
4) Naphthalene	12.37	128	3003	0.06	ng	# 96
5) 2-Methylnaphthalene	14.03	142	1585	0.06	ng	94
8) Acenaphthylene	16.15	152	7901	0.05	ng	98
9) Acenaphthene	16.58	154	1449	0.05	ng	91
10) Fluorene	17.83	166	1533	0.05	ng	# 96
12) Pentachlorophenol	19.52	266	34	0.34	ng	# 76
13) Phenanthrene	19.88	178	2832	0.06	ng	97
14) Anthracene	19.97	178	2058	0.06	ng	96
15) Fluoranthene	21.78	202	2158	0.05	ng	99
17) Pyrene	22.10	202	2308	0.06	ng	98
19) Benzo(a)anthracene	23.70	228	10521	0.06	ng	93
20) Chrysene	23.74	228	12870	0.07	ng	97
21) Indeno(1,2,3-cd)pyrene	26.83	276	5025	0.06	ng	# 90
23) Benzo(b)fluoranthene	24.97	252	861	0.25	ng	# 86
24) Benzo(k)fluoranthene	25.00	252	1793m	0.04	ng	
25) Benzo(a)pyrene	25.34	252	981	0.21	ng	# 81
26) Dibenzo(a,h)anthracene	26.87	278	742	0.14	ng	# 67
27) Benzo(g,h,i)perylene	27.26	276	4983	0.03	ng	# 77

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

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