

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE021015\
 Data File : BE089563.D
 Acq On : 11 Feb 2015 00:18
 Operator : TP/IZ
 Sample : MDL-02-W
 Misc : MDL-WATER-0.1/0.2
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 MDL-02-W

Quant Time: Feb 11 02:41:44 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	65186	5.00	ng	0.00
2) Naphthalene-d8	12.32	136	249461	5.00	ng	0.00
6) Acenaphthene-d10	16.50	164	115859	5.00	ng	0.00
11) Phenanthrene-d10	19.84	188	205299	5.00	ng	0.00
16) Chrysene-d12	23.72	240	238427	5.00	ng	0.00
22) Perylene-d12	25.40	264	237733	5.00	ng	0.00

System Monitoring Compounds						
3) Nitrobenzene-d5	10.68	82	136575	8.54	ng	0.00
7) 2-Fluorobiphenyl	14.96	172	208099	6.63	ng	0.00
18) Terphenyl-d14	22.36	244	253850	8.55	ng	0.00

Target Compounds						Qvalue
4) Naphthalene	12.37	128	4045	0.07	ng	# 94
5) 2-Methylnaphthalene	14.02	142	2079	0.07	ng	97
8) Acenaphthylene	16.15	152	10511	0.05	ng	99
9) Acenaphthene	16.58	154	1852	0.06	ng	93
10) Fluorene	17.83	166	1994	0.06	ng	98
12) Pentachlorophenol	19.52	266	45	0.34	ng	# 73
13) Phenanthrene	19.88	178	3559	0.07	ng	98
14) Anthracene	19.98	178	2593	0.06	ng	99
15) Fluoranthene	21.78	202	2809	0.06	ng	97
17) Pyrene	22.09	202	3064	0.06	ng	99
19) Benzo(a)anthracene	23.70	228	13446	0.07	ng	94
20) Chrysene	23.74	228	16638	0.07	ng	98
21) Indeno(1,2,3-cd)pyrene	26.83	276	5410	0.06	ng	# 85
23) Benzo(b)fluoranthene	24.97	252	1275	0.26	ng	# 88
24) Benzo(k)fluoranthene	25.00	252	2932m	0.05	ng	
25) Benzo(a)pyrene	25.33	252	1455	0.22	ng	# 82
26) Dibenzo(a,h)anthracene	26.86	278	960	0.14	ng	# 68
27) Benzo(g,h,i)perylene	27.26	276	7957	0.04	ng	# 80

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

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