

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE021015\
 Data File : BE089562.D
 Acq On : 10 Feb 2015 23:39
 Operator : TP/IZ
 Sample : MDL-01-W
 Misc : MDL-WATER-0.1/0.2
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 MDL-01-W

Quant Time: Feb 11 02:40:11 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	9.41	152	65062	5.00	ng	0.00
2) Naphthalene-d8	12.32	136	247801	5.00	ng	0.00
6) Acenaphthene-d10	16.50	164	117158	5.00	ng	0.00
11) Phenanthrene-d10	19.84	188	207530	5.00	ng	0.00
16) Chrysene-d12	23.71	240	239868	5.00	ng	0.00
22) Perylene-d12	25.40	264	240250	5.00	ng	0.00

System Monitoring Compounds						
3) Nitrobenzene-d5	10.68	82	153143	9.64	ng	0.00
7) 2-Fluorobiphenyl	14.95	172	231379	7.29	ng	0.00
18) Terphenyl-d14	22.36	244	281101	9.41	ng	0.00

Target Compounds						Qvalue
4) Naphthalene	12.37	128	4185	0.08	ng	# 95
5) 2-Methylnaphthalene	14.03	142	2180	0.07	ng	99
8) Acenaphthylene	16.15	152	11006	0.06	ng	99
9) Acenaphthene	16.58	154	1961	0.06	ng	96
10) Fluorene	17.83	166	2141	0.06	ng	98
12) Pentachlorophenol	19.51	266	57	0.35	ng	# 83
13) Phenanthrene	19.88	178	3630	0.07	ng	99
14) Anthracene	19.97	178	2881	0.07	ng	99
15) Fluoranthene	21.78	202	3017	0.06	ng	98
17) Pyrene	22.10	202	3306	0.07	ng	99
19) Benzo(a)anthracene	23.70	228	14781m	0.07	ng	
20) Chrysene	23.74	228	17123	0.08	ng	99
21) Indeno(1,2,3-cd)pyrene	26.83	276	7229	0.06	ng	96
23) Benzo(b)fluoranthene	24.97	252	1376	0.26	ng	# 89
24) Benzo(k)fluoranthene	25.00	252	3477m	0.06	ng	
25) Benzo(a)pyrene	25.34	252	1657	0.22	ng	# 89
26) Dibenzo(a,h)anthracene	26.87	278	1311	0.15	ng	# 79
27) Benzo(g,h,i)perylene	27.26	276	8429	0.04	ng	# 82

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

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