

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062615\
 Data File : BE089958.D
 Acq On : 27 Jun 2015 2:11
 Operator : TP/IZ
 Sample : G2653-06
 Misc : LOQ-SIM-WATER-0.2PPM
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 LOQ-WATER2015-02

Quant Time: Jun 27 03:08:45 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BE062315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 25 16:51:27 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	29301	5.00	ng	0.00
6) Naphthalene-d8	10.39	136	136748	5.00	ng	0.00
12) Acenaphthene-d10	14.24	164	81030	5.00	ng	0.00
18) Phenanthrene-d10	16.98	188	193351	5.00	ng	0.00
25) Chrysene-d12	21.13	240	204381	5.00	ng	0.00
32) Perylene-d12	23.47	264	180359	5.00	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.26	112	55621	10.51	ng	0.00
5) Phenol-d6	6.81	99	89697	11.07	ng	0.00
7) Nitrobenzene-d5	8.77	82	65220	6.41	ng	0.00
13) 2,4,6-Tribromophenol	15.74	330	9669	6.81	ng	0.00
14) 2-Fluorobiphenyl	12.87	172	154210	6.53	ng	0.00
27) Terphenyl-d14	19.60	244	228171	6.47	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.26	88	642	0.17	ng	# 76
3) n-Nitrosodimethylamine	3.55	42	640	0.12	ng	# 80
8) Nitrobenzene	8.81	77	1300	0.12	ng	96
9) Naphthalene	10.44	128	3524	0.12	ng	94
10) Hexachlorobutadiene	10.74	225	554	0.12	ng	98
11) 2-Methylnaphthalene	12.05	142	2352	0.11	ng	98
15) Acenaphthylene	13.95	152	15103	0.11	ng	100
16) Acenaphthene	14.30	154	2358	0.11	ng	96
17) Fluorene	15.29	166	3204	0.12	ng	# 96
19) 4-Bromophenyl-phenylether	16.18	248	867	0.11	ng	93
20) Hexachlorobenzene	16.30	284	4150	0.12	ng	99
21) Pentachlorophenol	16.63	266	293	0.58	ng	# 89
22) Phenanthrene	17.01	178	5475	0.11	ng	97
23) Anthracene	17.12	178	4579	0.10	ng	99
24) Fluoranthene	19.02	202	5328	0.10	ng	100
26) Pyrene	19.38	202	5552	0.11	ng	100
28) Benzo(a)anthracene	21.12	228	5696	0.12	ng	98
29) Chrysene	21.17	228	5803	0.11	ng	98
30) Bis(2-ethylhexyl)phthalate	21.07	149	630	0.15	ng	# 100
31) Indeno(1,2,3-cd)pyrene	25.88	276	3545	0.09	ng	# 100
33) Benzo(b)fluoranthene	22.75	252	3847	0.10	ng	# 95
34) Benzo(k)fluoranthene	22.80	252	4285	0.10	ng	# 95
35) Benzo(a)pyrene	23.36	252	3320	0.10	ng	# 91
36) Dibenzo(a,h)anthracene	25.90	278	2860	0.09	ng	92
37) Benzo(g,h,i)perylene	26.62	276	3339	0.10	ng	95

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

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