

Data Path : Z:\HPCHEM1\BNA G\DATA\BG121615\
 Data File : BG020213.D
 Acq On : 16 Dec 2015 6:31
 Operator : UM/SJ
 Sample : G4786-23
 Misc :
 ALS Vial : 59 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 G9C97

Quant Time: Dec 16 07:36:10 2015
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG121415.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 16 03:45:51 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.10	152	20383	20.00	ng/ul	0.00
18) Naphthalene-d8	10.91	136	97542	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.72	164	74215	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.47	188	190944	20.00	ng/ul	0.00
78) Chrysene-d12	21.74	240	203464	20.00	ng/ul	0.00
86) Perylene-d12	25.01	264	190168	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.47	96	456	1.24	ng/uL	0.00
5) Phenol-d5	7.25	99	14035	7.81	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.42	67	34491	32.16	ng/ul	0.00
9) 2-Chlorophenol-d4	7.63	132	35877	27.44	ng/ul	0.00
13) 4-Methylphenol-d8	8.80	113	28119	18.29	ng/ul	0.00
19) Nitrobenzene-d5	9.27	128	24190	31.81	ng/ul	0.00
22) 2-Nitrophenol-d4	9.99	143	28847	34.09	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.53	165	51828	30.24	ng/ul	0.00
29) 4-Chloroaniline-d4	11.05	131	1873	0.97	ng/ul	0.01
44) Dimethylphthalate-d6	14.13	166	196483	32.73	ng/ul	0.00
47) Acenaphthylene-d8	14.42	160	228729	32.75	ng/ul	0.00
52) 4-Nitrophenol-d4	14.91	143	6609	6.14	ng/ul	0.00
58) Fluorene-d10	15.71	176	176490	32.41	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.82	200	35244	31.14	ng/ul	0.00
71) Anthracene-d10	17.57	188	279630	31.78	ng/ul	0.00
79) Pyrene-d10	19.85	212	309638	32.65	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.78	264	312278	32.92	ng/ul	0.00

Target Compounds Ovalue

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

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