

Data Path : V:\HPCHEM1\BNA G\DATA\BG033116\
 Data File : BG021595.D
 Acq On : 31 Mar 2016 13:12
 Operator : UM/SJ
 Sample : PB89256BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :

Quant Time: Apr 01 01:29:48 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG031616.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Apr 01 01:28:23 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.05	152	9419	20.00	ng/ul	0.00
18) Naphthalene-d8	10.85	136	42638	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.67	164	28057	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.41	188	70503	20.00	ng/ul	0.00
78) Chrysene-d12	21.67	240	78864	20.00	ng/ul	0.00
86) Perylene-d12	24.88	264	74842	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.45	96	1303	6.21	ng/uL	0.00
5) Phenol-d5	7.34	99	24938	28.00	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.38	67	16938	32.11	ng/ul	0.00
9) 2-Chlorophenol-d4	7.63	132	22233	32.88	ng/ul	0.00
13) 4-Methylphenol-d8	8.85	113	23171	30.38	ng/ul	0.00
19) Nitrobenzene-d5	9.22	128	10883	34.02	ng/ul	0.00
22) 2-Nitrophenol-d4	9.94	143	12432	34.00	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.56	165	22094	33.80	ng/ul	0.00
29) 4-Chloroaniline-d4	11.01	131	4565	5.34	ng/ul	0.00
44) Dimethylphthalate-d6	14.07	166	84015	35.96	ng/ul	0.00
47) Acenaphthylene-d8	14.37	160	101340	35.19	ng/ul	0.00
52) 4-Nitrophenol-d4	15.10	143	9707	26.00	ng/ul	0.00
58) Fluorene-d10	15.66	176	75201	35.70	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.79	200	12759	27.39	ng/ul	0.00
71) Anthracene-d10	17.51	188	125901	36.44	ng/ul	0.00
79) Pyrene-d10	19.79	212	145214	39.49	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.66	264	148315	42.94	ng/ul	0.00

Target Compounds Ovalue

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

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