

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG033116\
 Data File : BG021597.D
 Acq On : 31 Mar 2016 17:26
 Operator : UM/SJ
 Sample : PB89375BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK75

Quant Time: Apr 01 01:29:59 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	7781	20.00	ng/ul	0.00
18) Naphthalene-d8	10.85	136	37886	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.67	164	26873	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.41	188	68937	20.00	ng/ul	0.00
78) Chrysene-d12	21.66	240	80129	20.00	ng/ul	0.00
86) Perylene-d12	24.88	264	76953	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.44	96	998	5.75	ng/uL	0.00
5) Phenol-d5	7.34	99	21612	29.38	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.37	67	14644	33.60	ng/ul	0.00
9) 2-Chlorophenol-d4	7.63	132	19001	34.01	ng/ul	0.00
13) 4-Methylphenol-d8	8.84	113	20036	31.80	ng/ul	0.00
19) Nitrobenzene-d5	9.22	128	10079	35.46	ng/ul	0.00
22) 2-Nitrophenol-d4	9.94	143	11431	35.18	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.56	165	19385	33.38	ng/ul	0.00
29) 4-Chloroaniline-d4	11.02	131	6764	8.91	ng/ul	0.01
44) Dimethylphthalate-d6	14.07	166	79650	35.60	ng/ul	0.00
47) Acenaphthylene-d8	14.37	160	91769	33.27	ng/ul	0.00
52) 4-Nitrophenol-d4	15.10	143	9144	25.57	ng/ul	0.00
58) Fluorene-d10	15.66	176	70337	34.86	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.79	200	12328	27.07	ng/ul	0.00
71) Anthracene-d10	17.51	188	116355	34.44	ng/ul	0.00
79) Pyrene-d10	19.79	212	134747	36.06	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.66	264	135104	38.04	ng/ul	0.00

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

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