

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG031716\
 Data File : BG021433.D
 Acq On : 17 Mar 2016 15:25
 Operator : UM/SJ
 Sample : PB88946BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK46

Quant Time: Mar 18 04:04:01 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG031616.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 18 04:03:18 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.08	152	7328	20.00	ng/ul	0.00
18) Naphthalene-d8	10.89	136	35110	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.70	164	24495	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.44	188	62778	20.00	ng/ul	0.00
78) Chrysene-d12	21.70	240	72457	20.00	ng/ul	0.00
86) Perylene-d12	24.93	264	70098	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.48	96	1036	6.34	ng/uL	0.00
5) Phenol-d5	7.33	99	21656	31.26	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.40	67	14524	35.39	ng/ul	0.00
9) 2-Chlorophenol-d4	7.65	132	17824	33.88	ng/ul	0.00
13) 4-Methylphenol-d8	8.85	113	18992	32.01	ng/ul	0.00
19) Nitrobenzene-d5	9.25	128	9075	34.45	ng/ul	0.00
22) 2-Nitrophenol-d4	9.97	143	10476	34.79	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.58	165	17978	33.40	ng/ul	0.00
29) 4-Chloroaniline-d4	11.03	131	28356	40.30	ng/ul	0.00
44) Dimethylphthalate-d6	14.10	166	80400	39.42	ng/ul	0.00
47) Acenaphthylene-d8	14.39	160	88734	35.29	ng/ul	0.00
52) 4-Nitrophenol-d4	15.07	143	10217	31.34	ng/ul	0.00
58) Fluorene-d10	15.69	176	66973	36.42	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.81	200	12270	29.58	ng/ul	0.00
71) Anthracene-d10	17.54	188	117886	38.32	ng/ul	0.00
79) Pyrene-d10	19.81	212	137054	40.56	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.71	264	144266	44.59	ng/ul	0.00

Target Compounds Qvalue

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

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