

Data Path : Z:\HPCHEM1\BNA G\DATA\BG101716\
 Data File : BG024384.D
 Acq On : 17 Oct 2016 15:46
 Operator : UM/SJ
 Sample : PB94041BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK41

Quant Time: Oct 17 17:20:44 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG0100716.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Oct 15 05:09:35 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.29	152	152224	20.00	ng/ul	0.00
18) Naphthalene-d8	11.11	136	681774	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.90	164	590817	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.64	188	1284206	20.00	ng/ul	0.00
75) Chrysene-d12	21.93	240	1440047	20.00	ng/ul	0.00
83) Perylene-d12	25.38	264	1420975	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.64	96	13447	4.19	ng/uL	0.00
5) Phenol-d5	7.41	99	297200	20.93	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.60	67	209226	21.79	ng/ul	0.00
9) 2-Chlorophenol-d4	7.81	132	209695	21.44	ng/ul	0.00
13) 4-Methylphenol-d8	8.97	113	250692	21.83	ng/ul	0.00
19) Nitrobenzene-d5	9.46	128	113479	23.45	ng/ul	0.00
22) 2-Nitrophenol-d4	10.18	143	136201	25.16	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.72	165	255944	21.71	ng/ul	0.00
29) 4-Chloroaniline-d4	11.24	131	338530	27.33	ng/ul	0.00
43) Dimethylphthalate-d6	14.29	166	931636	22.58	ng/ul	0.00
46) Acenaphthylene-d8	14.60	160	1080355	22.86	ng/ul	0.00
51) 4-Nitrophenol-d4	15.06	143	158295	21.64	ng/ul	0.00
57) Fluorene-d10	15.88	176	876299	22.47	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.99	200	158637	20.55	ng/ul	0.00
70) Anthracene-d10	17.74	188	1316117	22.69	ng/ul	0.00
76) Pyrene-d10	20.01	212	1497365	22.28	ng/ul	0.00
87) Benzo(a)pyrene-d12	25.15	264	1518969	23.76	ng/ul	0.00

Target Compounds Ovalue

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

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