

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072516\
 Data File : BG023257.D
 Acq On : 25 Jul 2016 18:38
 Operator : UM/SJ
 Sample : MDL-BLK-S
 Misc : MDL 4PPM
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MDL-BLK-S

Quant Time: Jul 26 03:11:34 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG072016.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jul 21 04:24:30 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.14	152	123957	20.00	ng/ul	0.00
18) Naphthalene-d8	10.98	136	586991	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.81	164	420769	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.57	188	969112	20.00	ng/ul	0.00
75) Chrysene-d12	21.89	240	998915	20.00	ng/ul	0.00
83) Perylene-d12	25.30	264	952650	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.50	96	12658	4.62	ng/uL	0.00
5) Phenol-d5	7.29	99	484745	37.93	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.52	67	384	0.05	ng/ul	0.06
9) 2-Chlorophenol-d4	7.66	132	352745	41.09	ng/ul	0.00
13) 4-Methylphenol-d8	8.85	113	378063	38.02	ng/ul	0.00
19) Nitrobenzene-d5	9.32	128	198059	42.18	ng/ul	0.00
22) 2-Nitrophenol-d4	10.04	143	214160	39.04	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.59	165	368337	35.73	ng/ul	0.00
29) 4-Chloroaniline-d4	11.11	131	515446	48.29	ng/ul	0.00
43) Dimethylphthalate-d6	14.20	166	1372995	39.25	ng/ul	0.00
46) Acenaphthylene-d8	14.50	160	1617371	40.68	ng/ul	0.00
51) 4-Nitrophenol-d4	15.01	143	209672	33.96	ng/ul	0.00
57) Fluorene-d10	15.80	176	1252875	39.60	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.92	200	234648	35.19	ng/ul	0.00
70) Anthracene-d10	17.67	188	1843160	40.63	ng/ul	0.00
76) Pyrene-d10	19.96	212	2062941	40.01	ng/ul	0.00
87) Benzo(a)pyrene-d12	25.07	264	1914470	42.81	ng/ul	0.00

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

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