

Data Path : Z:\HPCHEM1\BNA G\DATA\BG081816\
 Data File : BG023769.D
 Acq On : 18 Aug 2016 22:18
 Operator : UM/SJ
 Sample : H4460-15
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BD2T7

Manual Integrations
 APPROVED

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 8/19/2016 6:37:07 PM

Quant Time: Aug 19 03:02:35 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG081816.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Aug 19 01:43:20 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.10	152	101664	20.00	ng/ul	-0.02
18) Naphthalene-d8	10.93	136	436464	20.00	ng/ul	-0.02
35) Acenaphthene-d10	14.77	164	286545	20.00	ng/ul	-0.01
61) Phenanthrene-d10	17.53	188	678535m	20.00	ng/ul	-0.01
75) Chrysene-d12	21.84	240	782997m	20.00	ng/ul	0.00
83) Perylene-d12	25.22	264	729432	20.00	ng/ul	-0.01

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System Monitoring Compounds

3) 1,4-Dioxane-d8	3.47	96	3033m	1.35	ng/uL	0.00
5) Phenol-d5	7.26	99	65312	6.85	ng/ul	-0.02
7) Bis-(2-Chloroethyl)ether-d	7.41	67	211511	32.38	ng/ul	-0.02
9) 2-Chlorophenol-d4	7.63	132	175141	25.88	ng/ul	-0.02
13) 4-Methylphenol-d8	8.81	113	118072	15.84	ng/ul	-0.02
19) Nitrobenzene-d5	9.27	128	112651	31.95	ng/ul	-0.02
22) 2-Nitrophenol-d4	10.00	143	132388	32.93	ng/ul	-0.01
26) 2,4-Dichlorophenol-d3	10.55	165	208841	28.73	ng/ul	-0.02
29) 4-Chloroaniline-d4	11.08	131	22142	2.49	ng/ul	0.00
43) Dimethylphthalate-d6	14.16	166	814234	34.28	ng/ul	-0.01
46) Acenaphthylene-d8	14.46	160	922011	34.28	ng/ul	-0.02
51) 4-Nitrophenol-d4	15.01	143	24474m	6.37	ng/ul	0.01
57) Fluorene-d10	15.76	176	701635	33.83	ng/ul	-0.01
62) 4,6-Dinitro-2-methylphenol	15.90	200	131874m	30.66	ng/ul	-0.01
70) Anthracene-d10	17.63	188	1069049	34.17	ng/ul	-0.01
76) Pyrene-d10	19.92	212	1241494	33.04	ng/ul	0.00
87) Benzo(a)pyrene-d12	24.99	264	1156868	34.64	ng/ul	-0.01

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

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