

Data Path : Z:\HPCHEM1\BNA G\DATA\BG092115\
 Data File : BG018779.D
 Acq On : 21 Sep 2015 12:56
 Operator : TP
 Sample : G3691-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 H0228

Quant Time: Sep 21 16:14:13 2015
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Sep 21 15:50:58 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.00	152	44052	20.00	ng/ul	0.00
18) Naphthalene-d8	10.80	136	190238	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.63	164	129645	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.37	188	379831	20.00	ng/ul	0.00
78) Chrysene-d12	21.63	240	439619	20.00	ng/ul	-0.01
86) Perylene-d12	24.82	264	434909	20.00	ng/ul	-0.02

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.44	96	1487	1.62	ng/uL	0.00
5) Phenol-d5	7.16	99	26135	7.22	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.32	67	67647	32.27	ng/ul	0.00
9) 2-Chlorophenol-d4	7.53	132	73816	26.94	ng/ul	0.00
13) 4-Methylphenol-d8	8.70	113	49356	16.82	ng/ul	-0.01
19) Nitrobenzene-d5	9.16	128	48177	33.43	ng/ul	0.00
22) 2-Nitrophenol-d4	9.88	143	51941	35.84	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.43	165	89774	29.42	ng/ul	0.00
29) 4-Chloroaniline-d4	10.94	131	3843	1.08	ng/ul	0.00
44) Dimethylphthalate-d6	14.03	166	382493	36.66	ng/ul	0.00
47) Acenaphthylene-d8	14.32	160	415038	33.57	ng/ul	0.00
52) 4-Nitrophenol-d4	14.83	143	13662	6.82	ng/ul	0.00
58) Fluorene-d10	15.62	176	317917	34.85	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.73	200	51309	28.54	ng/ul	-0.01
71) Anthracene-d10	17.47	188	577276	34.52	ng/ul	0.00
79) Pyrene-d10	19.76	212	686266	33.93	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.60	264	707477	33.43	ng/ul	-0.02

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

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