

Data Path : Z:\HPCHEM1\BNA G\DATA\BG081816\
 Data File : BG023793.D
 Acq On : 19 Aug 2016 13:29
 Operator : UM/SJ
 Sample : H4492-15
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 JHPW0

Quant Time: Aug 19 16:53:43 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG081816.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Aug 19 04:06:00 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.09	152	123611	20.00	ng/ul	0.00
18) Naphthalene-d8	10.93	136	541341	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.77	164	366034	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.53	188	812245	20.00	ng/ul	0.00
75) Chrysene-d12	21.85	240	849371	20.00	ng/ul	0.00
83) Perylene-d12	25.23	264	831443	20.00	ng/ul	0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.46	96	14212	5.22	ng/uL	0.00
5) Phenol-d5	7.25	99	366366	31.61	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.41	67	267158	33.64	ng/ul	0.00
9) 2-Chlorophenol-d4	7.62	132	271015	32.94	ng/ul	0.00
13) 4-Methylphenol-d8	8.81	113	309868	34.19	ng/ul	0.00
19) Nitrobenzene-d5	9.27	128	146934	33.60	ng/ul	0.00
22) 2-Nitrophenol-d4	10.00	143	187833	37.67	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.55	165	304793	33.81	ng/ul	0.00
29) 4-Chloroaniline-d4	11.07	131	405415	36.75	ng/ul	0.00
43) Dimethylphthalate-d6	14.17	166	1017090	33.52	ng/ul	0.00
46) Acenaphthylene-d8	14.46	160	1139333	33.16	ng/ul	0.00
51) 4-Nitrophenol-d4	14.99	143	121807	24.84	ng/ul	0.00
57) Fluorene-d10	15.77	176	884532	33.39	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.90	200	69413	13.48	ng/ul	0.00
70) Anthracene-d10	17.63	188	1287716	34.39	ng/ul	0.00
76) Pyrene-d10	19.92	212	1410696	34.61	ng/ul	0.00
87) Benzo(a)pyrene-d12	25.00	264	1382907	36.33	ng/ul	0.01

Target Compounds

					Ovalue	
6) Phenol	7.28	94	33872	2.85	ng/ul#	75
44) Dimethylphthalate	14.21	163	663662	21.87	ng/ul	98

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

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