

Data Path : Z:\HPCHEM1\BNA G\DATA\BG032717\
 Data File : BG026446.D
 Acq On : 27 Mar 2017 23:50
 Operator : SJ/MA
 Sample : I2358-16
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 COBMO

Quant Time: Mar 28 11:04:44 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM-EPA-BG032717MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Mar 28 01:23:51 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.06	152	189585	20.00	ng/ul	0.01
18) Naphthalene-d8	10.85	136	772014	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.65	164	450152	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.38	188	1057939	20.00	ng/ul	0.00
77) Chrysene-d12	21.63	240	1228099	20.00	ng/ul	0.00
85) Perylene-d12	24.81	264	1210026	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.47	96	20434	5.16	ng/uL	0.00
5) Phenol-d5	7.21	99	92527	6.63	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.37	67	258138	32.99	ng/ul	0.00
9) 2-Chlorophenol-d4	7.58	132	330199	26.38	ng/ul	0.00
13) 4-Methylphenol-d8	8.75	113	180281	16.03	ng/ul	0.00
19) Nitrobenzene-d5	9.21	128	201011	36.43	ng/ul	0.00
22) 2-Nitrophenol-d4	9.92	143	236668	36.60	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.47	165	362328	30.53	ng/ul	0.00
29) 4-Chloroaniline-d4	10.98	131	6474	0.55	ng/ul	0.00
43) Dimethylphthalate-d6	14.05	166	1217018	34.91	ng/ul	0.00
46) Acenaphthylene-d8	14.35	160	1451354	33.42	ng/ul	0.00
51) 4-Nitrophenol-d4	14.85	143	43660	6.44	ng/ul	0.01
57) Fluorene-d10	15.63	176	1102240	34.81	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.75	200	234380	34.45	ng/ul	0.00
70) Anthracene-d10	17.48	188	1711577	35.63	ng/ul	0.00
78) Pyrene-d10	19.75	212	1942401	35.46	ng/ul	0.00
89) Benzo(a)pyrene-d12	24.59	264	1918409	36.90	ng/ul	0.00

Target Compounds

						Ovalue
6) Phenol	7.24	94	45007	3.09	ng/ul#	90
10) 2-Chlorophenol	7.61	128	97274	8.03	ng/ul	99
20) Nitrobenzene	9.25	77	2367086	215.95	ng/ul	88
27) 2,4-Dichlorophenol	10.49	162	27539	2.35	ng/ul#	85
36) 1,2,4,5-Tetrachlorobenzene	12.85	216	149306	10.53	ng/ul#	95
72) 1,2,3,4-Tetrachlorobenzene	13.46	216	392667	28.18	ng/uL	98

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

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