

Data Path : Z:\HPCHEM1\BNA G\DATA\BG030316\
 Data File : BG021268.D
 Acq On : 4 Mar 2016 13:19
 Operator : UM/SJ
 Sample : H1583-11
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 D9R54

Quant Time: Mar 04 23:51:55 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG030316.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 04 23:46:39 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.13	152	10278	20.00	ng/ul	0.00
18) Naphthalene-d8	10.93	136	51291	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.74	164	30427	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.47	188	67358	20.00	ng/ul	0.00
78) Chrysene-d12	21.73	240	73578	20.00	ng/ul	0.00
86) Perylene-d12	24.99	264	67981	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.54	96	204	0.81	ng/uL	0.00
5) Phenol-d5	7.28	99	5058	4.65	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.45	67	20985	28.91	ng/ul	0.00
9) 2-Chlorophenol-d4	7.66	132	16792	22.11	ng/ul	0.00
13) 4-Methylphenol-d8	8.82	113	10526	12.25	ng/ul	0.00
19) Nitrobenzene-d5	9.30	128	11970	28.81	ng/ul	0.00
22) 2-Nitrophenol-d4	10.01	143	12291	27.75	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.55	165	19062	24.08	ng/ul	0.00
29) 4-Chloroaniline-d4	0.00	131	0	0.00	ng/ul	0.00
44) Dimethylphthalate-d6	14.14	166	76255	29.91	ng/ul	0.00
47) Acenaphthylene-d8	14.44	160	95403	30.07	ng/ul	0.00
52) 4-Nitrophenol-d4	14.92	143	1892	3.81	ng/ul	0.00
58) Fluorene-d10	15.72	176	67980	30.36	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.83	200	12826	26.53	ng/ul	0.00
71) Anthracene-d10	17.57	188	105555	31.51	ng/ul	0.00
79) Pyrene-d10	19.85	212	119002	31.95	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.77	264	114136	31.10	ng/ul	0.00

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

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