

Data Path : Z:\HPCHEM1\BNA G\DATA\BG110415\
 Data File : BG019481.D
 Acq On : 4 Nov 2015 22:08
 Operator : UM/NP
 Sample : G4245-07
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BCY44

Quant Time: Nov 04 23:05:43 2015
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG102815.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 04 22:22:38 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.18	152	28863	20.00	ng/ul	0.00
18) Naphthalene-d8	11.01	136	120778	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.81	164	96018	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.55	188	269607	20.00	ng/ul	0.00
78) Chrysene-d12	21.83	240	360722	20.00	ng/ul	0.00
86) Perylene-d12	25.17	264	342673	20.00	ng/ul	-0.02

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.49	96	675	1.12	ng/uL	0.00
5) Phenol-d5	7.33	99	17811	6.72	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.49	67	66016	39.30	ng/ul	0.00
9) 2-Chlorophenol-d4	7.70	132	47741	28.86	ng/ul	0.00
13) 4-Methylphenol-d8	8.88	113	37481	17.76	ng/ul	0.00
19) Nitrobenzene-d5	9.35	128	32495	36.61	ng/ul	0.00
22) 2-Nitrophenol-d4	10.08	143	38784	36.81	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.62	165	75098	32.85	ng/ul	0.00
29) 4-Chloroaniline-d4	11.14	131	3642	1.57	ng/ul	0.00
44) Dimethylphthalate-d6	14.20	166	306837	38.76	ng/ul	0.00
47) Acenaphthylene-d8	14.51	160	343564	39.19	ng/ul	0.00
52) 4-Nitrophenol-d4	14.99	143	8761	6.97	ng/ul	0.00
58) Fluorene-d10	15.79	176	280579	38.31	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.91	200	63748	36.53	ng/ul	0.00
71) Anthracene-d10	17.64	188	497485	40.48	ng/ul	0.00
79) Pyrene-d10	19.92	212	595236	37.71	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.95	264	671987	39.08	ng/ul	0.00

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

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