

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092515\
 Data File : BG018900.D
 Acq On : 26 Sep 2015 5:02
 Operator : UM/NP
 Sample : G3797-11
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 D9G79

Manual Integrations
 APPROVED

MMDadoda
 9/28/2015 6:14:11 PM

Quant Time: Sep 26 06:55:11 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG091015.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Sep 26 04:18:39 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.99	152	42403	20.00	ng/ul	0.00
18) Naphthalene-d8	10.79	136	175998	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.61	164	120330	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.35	188	366516	20.00	ng/ul	0.00
78) Chrysene-d12	21.60	240	455716	20.00	ng/ul	0.00
86) Perylene-d12	24.78	264	464422	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.43	96	1154	1.31	ng/uL	0.00
5) Phenol-d5	7.16	99	24092	6.92	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.32	67	61715	30.59	ng/ul	0.00
9) 2-Chlorophenol-d4	7.53	132	69446	26.33	ng/ul	0.00
13) 4-Methylphenol-d8	8.70	113	46229	16.36	ng/ul	0.00
19) Nitrobenzene-d5	9.15	128	43248	32.43	ng/ul	0.00
22) 2-Nitrophenol-d4	9.87	143	50357	37.56	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.41	165	88316	31.28	ng/ul	0.00
29) 4-Chloroaniline-d4	10.93	131	1717	0.52	ng/ul	0.00
44) Dimethylphthalate-d6	14.01	166	391333	40.41	ng/ul	0.00
47) Acenaphthylene-d8	14.31	160	317641	27.68	ng/ul	0.00
52) 4-Nitrophenol-d4	14.81	143	13733	7.39	ng/ul	0.00
58) Fluorene-d10	15.60	176	211501	24.98	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.72	200	42160	24.30	ng/ul	0.00
71) Anthracene-d10	17.45	188	194422m	12.05	ng/ul	0.00
79) Pyrene-d10	19.74	212	421790	20.12	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.56	264	325508	14.40	ng/ul	0.00

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

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