

Data Path : Z:\HPCHEM1\BNA G\DATA\BG092115\
 Data File : BG018796.D
 Acq On : 21 Sep 2015 23:42
 Operator : TP
 Sample : G3696-17
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0AP1

Manual Integrations
 APPROVED

MMDadoda
 9/22/2015 2:14:45 PM

Quant Time: Sep 22 02:50:00 2015
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 22 02:06:40 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.00	152	49808	20.00	ng/ul	0.00
18) Naphthalene-d8	10.80	136	215654	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.62	164	146367	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.37	188	394332	20.00	ng/ul	0.00
78) Chrysene-d12	21.63	240	448688	20.00	ng/ul	0.00
86) Perylene-d12	24.81	264	450743	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.44	96	1307	1.26	ng/uL	0.00
5) Phenol-d5	7.16	99	21691	5.30	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.32	67	69065	29.14	ng/ul	0.00
9) 2-Chlorophenol-d4	7.53	132	72186	23.30	ng/ul	0.00
13) 4-Methylphenol-d8	8.70	113	45349	13.66	ng/ul	0.00
19) Nitrobenzene-d5	9.16	128	48670m	29.79	ng/ul	0.00
22) 2-Nitrophenol-d4	9.88	143	51479	31.33	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.42	165	88038	25.45	ng/ul	0.00
29) 4-Chloroaniline-d4	0.00	131	0d	0.00	ng/ul	
44) Dimethylphthalate-d6	14.03	166	377438	32.04	ng/ul	0.00
47) Acenaphthylene-d8	14.32	160	420396	30.12	ng/ul	0.00
52) 4-Nitrophenol-d4	14.83	143	9930	4.39	ng/ul	0.00
58) Fluorene-d10	15.62	176	326253	31.67	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.73	200	51294	27.48	ng/ul	0.00
71) Anthracene-d10	17.47	188	566204	32.61	ng/ul	0.00
79) Pyrene-d10	19.75	212	667731	32.35	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.59	264	609338	27.78	ng/ul	0.00

Target Compounds Ovalue

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

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