

Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\
 Data File : BG018975.D
 Acq On : 30 Sep 2015 16:38
 Operator : UM/NP
 Sample : G3803-03
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 E43J3

Manual Integrations
 APPROVED

MMDadoda
 10/1/2015 6:19:29 PM

Quant Time: Oct 01 05:26:23 2015
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 01 04:12:22 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.99	152	50189	20.00	ng/ul	0.00
18) Naphthalene-d8	10.79	136	227132	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.61	164	166651	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.36	188	492160	20.00	ng/ul	0.00
78) Chrysene-d12	21.61	240	565691	20.00	ng/ul	0.00
86) Perylene-d12	24.79	264	569677	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.43	96	1391m	1.33	ng/uL	0.00
5) Phenol-d5	7.16	99	28919	7.02	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.31	67	47324	19.81	ng/ul	0.00
9) 2-Chlorophenol-d4	7.52	132	62159	19.91	ng/ul	0.00
13) 4-Methylphenol-d8	8.70	113	49222	14.72	ng/ul	0.00
19) Nitrobenzene-d5	9.14	128	37691	21.90	ng/ul	0.00
22) 2-Nitrophenol-d4	9.87	143	41962	24.25	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.41	165	81299	22.31	ng/ul	0.00
29) 4-Chloroaniline-d4	10.93	131	6538	1.54	ng/ul	0.00
44) Dimethylphthalate-d6	14.01	166	305870	22.81	ng/ul	0.00
47) Acenaphthylene-d8	14.31	160	355163	22.35	ng/ul	0.00
52) 4-Nitrophenol-d4	14.83	143	17366	6.75	ng/ul	0.00
58) Fluorene-d10	15.60	176	281246	23.98	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.72	200	48631	20.87	ng/ul	0.00
71) Anthracene-d10	17.45	188	484899	22.38	ng/ul	0.00
79) Pyrene-d10	19.74	212	595670	22.89	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.57	264	628926	22.68	ng/ul	0.00

Target Compounds Ovalue

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

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