

Data Path : Z:\HPCHEM1\BNA G\DATA\BG080817\
 Data File : BG028198.D
 Acq On : 8 Aug 2017 20:23
 Operator : SJ/JU
 Sample : I4622-05
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :

Quant Time: Aug 09 02:33:56 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM-EPA-BG080217.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Aug 09 02:30:54 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.36	152	28686	20.00	ng/ul	0.00
18) Naphthalene-d8	11.17	136	118831	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.96	164	86254	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.71	188	262088	20.00	ng/ul	0.00
75) Chrysene-d12	22.04	240	310416	20.00	ng/ul	0.00
83) Perylene-d12	25.55	264	289142	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	4.10	96	64	0.10	ng/uL	0.26
5) Phenol-d5	0.00	99	0	0.00	ng/ul	
7) Bis-(2-Chloroethyl)ether-d	7.53	67	61	0.03	ng/ul	-0.14
9) 2-Chlorophenol-d4	0.00	132	0	0.00	ng/ul	
13) 4-Methylphenol-d8	0.00	113	0	0.00	ng/ul	
19) Nitrobenzene-d5	0.00	128	0	0.00	ng/ul	
22) 2-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
26) 2,4-Dichlorophenol-d3	0.00	165	0	0.00	ng/ul	
29) 4-Chloroaniline-d4	11.41	131	134	0.06	ng/ul	0.10
43) Dimethylphthalate-d6	0.00	166	0	0.00	ng/ul	
46) Acenaphthylene-d8	14.66	160	75	0.01	ng/ul	0.00
51) 4-Nitrophenol-d4	14.90	143	59	0.05	ng/ul	-0.25
57) Fluorene-d10	15.97	176	422	0.06	ng/ul	0.02
62) 4,6-Dinitro-2-methylphenol	16.16	200	73	0.04	ng/ul	0.09
70) Anthracene-d10	17.71	188	262088	20.85	ng/ul	-0.10
76) Pyrene-d10	20.12	212	53	0.00	ng/ul	0.03
87) Benzo(a)pyrene-d12	25.55	264	277276	20.48	ng/ul	0.24

Target Compounds	Ovalue

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

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