

Data Path : Z:\HPCHEM1\BNA G\DATA\BG081617\
 Data File : BG028377.D
 Acq On : 16 Aug 2017 18:57
 Operator : SJ/JU
 Sample : I4720-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :

Quant Time: Aug 17 06:53:49 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM-EPA-BG080217.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Aug 17 06:52:43 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.34	152	35354	20.00	ng/ul	0.00
18) Naphthalene-d8	11.16	136	169383	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.95	164	129972	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.70	188	340220	20.00	ng/ul	0.00
75) Chrysene-d12	22.03	240	354788	20.00	ng/ul	0.00
83) Perylene-d12	25.53	264	351676	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.82	96	3562	4.69	ng/uL	0.00
5) Phenol-d5	7.50	99	22450	5.78	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.66	67	56755	23.08	ng/ul	0.00
9) 2-Chlorophenol-d4	7.88	132	52372	21.38	ng/ul	0.00
13) 4-Methylphenol-d8	9.04	113	42850	13.20	ng/ul	0.00
19) Nitrobenzene-d5	9.51	128	34438	26.00	ng/ul	0.00
22) 2-Nitrophenol-d4	10.23	143	42247	29.09	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.78	165	76898	25.59	ng/ul	0.00
29) 4-Chloroaniline-d4	11.30	131	7523	2.25	ng/ul	0.00
43) Dimethylphthalate-d6	14.34	166	326963	28.28	ng/ul	0.00
46) Acenaphthylene-d8	14.65	160	349135	26.79	ng/ul	0.00
51) 4-Nitrophenol-d4	15.16	143	9266	5.04	ng/ul	0.00
57) Fluorene-d10	15.94	176	271585	26.19	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	16.06	200	51500	23.30	ng/ul	0.00
70) Anthracene-d10	17.80	188	457825	28.06	ng/ul	0.00
76) Pyrene-d10	20.08	212	501589	27.57	ng/ul	0.00
87) Benzo(a)pyrene-d12	25.29	264	443428	26.93	ng/ul	0.00

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

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