

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG041720\
 Data File : BG045038.D
 Acq On : 17 Apr 2020 18:26
 Operator : CG/JU
 Sample : L2300-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 WATER-RECLAIM-SYSTEM-SLUDG

Quant Time: Apr 17 19:15:54 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG041520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 17 11:08:29 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.03	152	56651	20.00	ng	0.00
21) Naphthalene-d8	10.83	136	242730	20.00	ng	0.00
39) Acenaphthene-d10	14.65	164	202482	20.00	ng	0.00
64) Phenanthrene-d10	17.40	188	495626	20.00	ng	0.00
76) Chrysene-d12	21.66	240	484188	20.00	ng	0.00
87) Perylene-d12	24.85	264	513350	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	433578	126.36	ng	0.00
7) Phenol-d6	7.18	99	641772	125.88	ng	0.00
23) Nitrobenzene-d5	9.18	82	491976	92.48	ng	0.00
42) 2,4,6-Tribromophenol	16.15	330	476010	152.17	ng	0.00
45) 2-Fluorobiphenyl	13.28	172	1281065	92.77	ng	0.00
79) Terphenyl-d14	20.01	244	2234774	100.60	ng	0.00
Target Compounds						
10) Phenol	7.21	94	63933	12.532	ng	84
20) 3+4-Methylphenols	8.79	107	251456	45.640	ng	91
32) Benzoic acid	10.07	122	13830	10.357	ng	92

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

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