

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100721\
 Data File : BG050493.D
 Acq On : 7 Oct 2021 21:50
 Operator : CG/JU
 Sample : M4088-02
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 WC-20211005-WATER

Quant Time: Oct 08 06:06:20 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 06 15:51:01 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.259	152	47269	20.00	ng	0.00
21) Naphthalene-d8	11.085	136	195480	20.00	ng	# 0.00
39) Acenaphthene-d10	14.875	164	132731	20.00	ng	0.00
64) Phenanthrene-d10	17.618	188	303577	20.00	ng	# 0.00
76) Chrysene-d12	21.913	240	270432	20.00	ng	# 0.00
86) Perylene-d12	25.327	264	266999	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.791	112	350017	110.74	ng	0.00
7) Phenol-d6	7.395	99	436112	95.30	ng	0.00
23) Nitrobenzene-d5	9.428	82	379978	80.47	ng	0.00
42) 2,4,6-Tribromophenol	16.355	330	165459	130.69	ng	0.00
45) 2-Fluorobiphenyl	13.500	172	697402	79.08	ng	0.00
79) Terphenyl-d14	20.210	244	1121092	80.79	ng	0.00

Target Compounds	Qvalue

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

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