

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081122\
 Data File : BG054607.D
 Acq On : 11 Aug 2022 16:29
 Operator : CG/JU
 Sample : N4017-02
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB09

Quant Time: Aug 12 06:04:15 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 08 14:48:51 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.970	152	15519	20.000	ng	0.00
21) Naphthalene-d8	10.779	136	60384	20.000	ng	# 0.00
39) Acenaphthene-d10	14.610	164	48253	20.000	ng	0.00
64) Phenanthrene-d10	17.359	188	128266	20.000	ng	# 0.00
76) Chrysene-d12	21.625	240	148704	20.000	ng	# 0.00
86) Perylene-d12	24.827	264	152921	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.556	112	88468	119.022	ng	0.00
7) Phenol-d6	7.171	99	124947	122.679	ng	0.00
23) Nitrobenzene-d5	9.140	82	84640	76.332	ng	0.00
42) 2,4,6-Tribromophenol	16.108	330	105460	104.093	ng	0.00
45) 2-Fluorobiphenyl	13.229	172	246449	72.050	ng	0.00
79) Terphenyl-d14	19.968	244	541005	72.173	ng	0.00

Target Compounds	Qvalue

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

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