

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062922\
 Data File : BG054144.D
 Acq On : 29 Jun 2022 14:17
 Operator : CG/JU
 Sample : N3488-10
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-12-17-19-20220623

Quant Time: Jun 29 14:53:30 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG061322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 27 00:27:54 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.071	152	22391	20.000	ng	0.00
21) Naphthalene-d8	10.879	136	87146	20.000	ng	0.00
39) Acenaphthene-d10	14.692	164	67239	20.000	ng	0.00
64) Phenanthrene-d10	17.436	188	176432	20.000	ng	# 0.00
76) Chrysene-d12	21.713	240	188278	20.000	ng	0.00
86) Perylene-d12	24.968	264	190861	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.638	112	139870	119.431	ng	0.00
7) Phenol-d6	7.230	99	191171	120.330	ng	0.00
23) Nitrobenzene-d5	9.228	82	126523	81.588	ng	0.00
42) 2,4,6-Tribromophenol	16.185	330	141488	130.969	ng	0.00
45) 2-Fluorobiphenyl	13.317	172	368336	79.658	ng	0.00
79) Terphenyl-d14	20.045	244	820289	87.749	ng	0.00

Target Compounds	Qvalue

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

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