

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092322\
 Data File : BG055018.D
 Acq On : 23 Sep 2022 18:50
 Operator : CG/JU
 Sample : N4751-07
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PE-6

Quant Time: Sep 24 02:28:49 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG091922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 19 17:41:40 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.054	152	38618	20.000	ng	0.00
21) Naphthalene-d8	10.880	136	154737	20.000	ng	# 0.00
39) Acenaphthene-d10	14.699	164	109461	20.000	ng	0.00
64) Phenanthrene-d10	17.449	188	254289	20.000	ng	# 0.00
76) Chrysene-d12	21.726	240	253495	20.000	ng	0.00
86) Perylene-d12	25.028	264	265178	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.598	112	178031	90.204	ng	0.00
7) Phenol-d6	7.220	99	236009	89.214	ng	0.00
23) Nitrobenzene-d5	9.235	82	134887	50.526	ng	0.00
42) 2,4,6-Tribromophenol	16.192	330	100819	78.622	ng	0.00
45) 2-Fluorobiphenyl	13.325	172	372985	53.701	ng	0.00
79) Terphenyl-d14	20.046	244	636118	52.492	ng	0.00
Target Compounds						
75) Fluoranthene	19.500	202	39740	2.423	ng	Qvalue 97

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

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