

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092322\
 Data File : BG055016.D
 Acq On : 23 Sep 2022 17:26
 Operator : CG/JU
 Sample : N4751-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PE-1

Quant Time: Sep 24 02:28:22 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	8.059	152	38948	20.000	ng	0.00
21) Naphthalene-d8	10.885	136	153423	20.000	ng	# 0.00
39) Acenaphthene-d10	14.704	164	110315	20.000	ng	0.00
64) Phenanthrene-d10	17.448	188	260264	20.000	ng	# 0.00
76) Chrysene-d12	21.725	240	259981	20.000	ng	0.00
86) Perylene-d12	25.033	264	273641	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.597	112	239358	120.249	ng	0.00
7) Phenol-d6	7.219	99	312909	117.281	ng	0.00
23) Nitrobenzene-d5	9.234	82	178948	67.605	ng	0.00
42) 2,4,6-Tribromophenol	16.190	330	145572	112.642	ng	0.00
45) 2-Fluorobiphenyl	13.323	172	501460	71.639	ng	0.00
79) Terphenyl-d14	20.051	244	927927	74.661	ng	0.00
Target Compounds						
75) Fluoranthene	19.498	202	38578	2.298	ng	95
78) Pyrene	19.863	202	34329	2.009	ng	96

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

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