

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG031721\
 Data File : BG048398.D
 Acq On : 17 Mar 2021 15:07
 Operator : CG/JU
 Sample : M1652-08
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PE-4D-(7.0-7.5)

Quant Time: Mar 17 15:47:44 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG031621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 16 16:05:10 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.115	152	47779	20.00	ng	0.00
21) Naphthalene-d8	10.941	136	166393	20.00	ng	# 0.00
39) Acenaphthene-d10	14.754	164	130632	20.00	ng	0.00
64) Phenanthrene-d10	17.504	188	328111	20.00	ng	# 0.00
76) Chrysene-d12	21.799	240	383828	20.00	ng	# 0.00
86) Perylene-d12	25.142	264	391512	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.659	112	300972	108.20	ng	0.00
7) Phenol-d6	7.263	99	411515	100.42	ng	0.00
23) Nitrobenzene-d5	9.284	82	291533	77.85	ng	0.00
42) 2,4,6-Tribromophenol	16.241	330	209603	125.41	ng	0.00
45) 2-Fluorobiphenyl	13.379	172	573735	61.90	ng	0.00
79) Terphenyl-d14	20.113	244	1222953	61.58	ng	0.00
Target Compounds						
10) Phenol	7.287	94	10870	2.60	ng	# 59

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

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