

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG123022\
 Data File : BG056171.D
 Acq On : 30 Dec 2022 15:06
 Operator : CG/JU
 Sample : PB149967TB
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB149967TB

Quant Time: Dec 31 00:17:49 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG122722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 28 05:20:14 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.341	152	49546	20.000	ng	0.00
21) Naphthalene-d8	11.167	136	169411	20.000	ng	# 0.00
39) Acenaphthene-d10	14.945	164	136370	20.000	ng	0.00
64) Phenanthrene-d10	17.677	188	369551	20.000	ng	0.00
76) Chrysene-d12	21.972	240	446099	20.000	ng	0.00
86) Perylene-d12	25.426	264	432861	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.855	112	377600	136.076	ng	0.00
7) Phenol-d6	7.471	99	503567	118.167	ng	0.00
23) Nitrobenzene-d5	9.510	82	261132	78.846	ng	0.00
42) 2,4,6-Tribromophenol	16.419	330	221514	128.257	ng	0.00
45) 2-Fluorobiphenyl	13.576	172	826918	83.735	ng	0.00
79) Terphenyl-d14	20.256	244	1998211	80.227	ng	0.00

Target Compounds	Qvalue

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

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