

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122223\
 Data File : BG060141.D
 Acq On : 22 Dec 2023 16:51
 Operator : MA/JU
 Sample : 05900-22
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 GHL-SS-64a

Quant Time: Dec 22 17:34:40 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG122123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 22 03:42:04 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.944	152	241968	20.000	ng	# 0.00
21) Naphthalene-d8	10.746	136	1104459	20.000	ng	0.00
39) Acenaphthene-d10	14.577	164	901176	20.000	ng	0.00
64) Phenanthrene-d10	17.321	188	2207674	20.000	ng	0.00
76) Chrysene-d12	21.587	240	1502504	20.000	ng	0.00
86) Perylene-d12	24.765	264	1584769	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.511	112	1488788	96.951	ng	0.00
7) Phenol-d6	7.104	99	2345900	90.773	ng	0.00
23) Nitrobenzene-d5	9.101	82	1329063	58.950	ng	0.00
42) 2,4,6-Tribromophenol	16.070	330	1055745	87.010	ng	0.00
45) 2-Fluorobiphenyl	13.202	172	3412170	56.854	ng	0.00
79) Terphenyl-d14	19.942	244	4483149	53.993	ng	0.00

Target Compounds	Qvalue

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

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