

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032222\
 Data File : BG052783.D
 Acq On : 22 Mar 2022 13:21
 Operator : CG/JU
 Sample : N1974-13
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 B-3

Quant Time: Mar 22 16:38:21 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG031522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 15 15:56:10 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.147	152	41471	20.000	ng	-0.01
21) Naphthalene-d8	10.955	136	167689	20.000	ng	#-0.02
39) Acenaphthene-d10	14.767	164	38562	20.000	ng	0.00
64) Phenanthrene-d10	17.505	188	182779	20.000	ng	#-0.01
76) Chrysene-d12	21.793	240	40140	20.000	ng	#-0.02
86) Perylene-d12	25.118	264	66	20.000	ng	#-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.703	112	284955	119.663	ng	0.00
7) Phenol-d6	7.289	99	205523	57.247	ng	0.00
23) Nitrobenzene-d5	9.304	82	320317	95.873	ng	-0.01
42) 2,4,6-Tribromophenol	16.247	330	208316	402.023	ng	-0.01
45) 2-Fluorobiphenyl	13.392	172	676504	258.737	ng	-0.01
79) Terphenyl-d14	20.113	244	730445	323.705	ng	-0.01

Target Compounds	Qvalue

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

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