

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042922\
 Data File : BG053383.D
 Acq On : 29 Apr 2022 18:58
 Operator : CG/JU
 Sample : N2616-11
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MH-20D

Quant Time: May 02 01:03:41 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG040822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Apr 24 01:13:30 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.099	152	19487	20.000	ng	#-0.01
21) Naphthalene-d8	10.913	136	75137	20.000	ng	0.00
39) Acenaphthene-d10	14.726	164	58161	20.000	ng	0.00
64) Phenanthrene-d10	17.469	188	169906	20.000	ng	-0.01
76) Chrysene-d12	21.757	240	208824	20.000	ng	0.00
86) Perylene-d12	25.053	264	210760	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.667	112	169609	149.197	ng	0.00
7) Phenol-d6	7.265	99	218341	124.547	ng	0.00
23) Nitrobenzene-d5	9.262	82	188393	101.806	ng	-0.01
42) 2,4,6-Tribromophenol	16.212	330	159781	172.677	ng	-0.01
45) 2-Fluorobiphenyl	13.345	172	414432	98.255	ng	-0.01
79) Terphenyl-d14	20.077	244	1165612	94.271	ng	0.00

Target Compounds	Qvalue

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

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