

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090922\
 Data File : BG054907.D
 Acq On : 9 Sep 2022 21:57
 Operator : CG/JU
 Sample : N4538-07 5X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MBD-SB-1-2-0-5

Quant Time: Sep 10 02:03:31 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG082522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 10 01:53:42 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.140	152	48585	20.000	ng	0.00
21) Naphthalene-d8	10.972	136	200749	20.000	ng	0.00
39) Acenaphthene-d10	14.780	164	130066	20.000	ng	0.00
64) Phenanthrene-d10	17.524	188	263938	20.000	ng	0.00
76) Chrysene-d12	21.819	240	247792	20.000	ng	0.00
86) Perylene-d12	25.197	264	261988	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.685	112	55766	19.987	ng	0.00
7) Phenol-d6	7.300	99	78164	20.485	ng	0.00
23) Nitrobenzene-d5	9.316	82	45253	11.834	ng	0.00
42) 2,4,6-Tribromophenol	16.266	330	29034	17.864	ng	0.00
45) 2-Fluorobiphenyl	13.399	172	125119	14.458	ng	0.00
79) Terphenyl-d14	20.121	244	182320	14.579	ng	0.00

Target Compounds	Qvalue

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

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