

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042222\
 Data File : BG053269.D
 Acq On : 22 Apr 2022 20:17
 Operator : CG/JU
 Sample : N2494-02
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MH-1S

Quant Time: Apr 24 01:16:09 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	8.107	152	27773	20.000	ng	0.00
21) Naphthalene-d8	10.921	136	108230	20.000	ng	0.00
39) Acenaphthene-d10	14.733	164	80055	20.000	ng	0.00
64) Phenanthrene-d10	17.477	188	207669	20.000	ng	0.00
76) Chrysene-d12	21.759	240	215679	20.000	ng	0.00
86) Perylene-d12	25.061	264	215033	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.675	112	210809	130.114	ng	0.00
7) Phenol-d6	7.273	99	271086	108.500	ng	0.00
23) Nitrobenzene-d5	9.270	82	241292	90.523	ng	0.00
42) 2,4,6-Tribromophenol	16.220	330	186005	146.042	ng	0.00
45) 2-Fluorobiphenyl	13.359	172	528931	91.106	ng	0.00
79) Terphenyl-d14	20.079	244	1159987	90.834	ng	0.00

Target Compounds	Qvalue

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

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