

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040722\
 Data File : BG053073.D
 Acq On : 7 Apr 2022 14:33
 Operator : CG/JU
 Sample : N2207-08RE
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 20220352-AMENDED-098-21-01RE

Quant Time: Apr 08 07:22:50 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG031522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 06 14:35:39 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.122	152	28949	20.000	ng	0.00
21) Naphthalene-d8	10.942	136	111790	20.000	ng	0.00
39) Acenaphthene-d10	14.748	164	81852	20.000	ng	0.00
64) Phenanthrene-d10	17.492	188	207834	20.000	ng	0.00
76) Chrysene-d12	21.774	240	217547	20.000	ng	0.00
86) Perylene-d12	25.081	264	218300	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.690	112	16958	10.202	ng	0.00
7) Phenol-d6	7.282	99	116683	46.560	ng	0.00
23) Nitrobenzene-d5	9.285	82	170326	76.471	ng	0.00
42) 2,4,6-Tribromophenol	16.229	330	9766	8.879	ng	0.00
45) 2-Fluorobiphenyl	13.368	172	356276	64.195	ng	0.00
79) Terphenyl-d14	20.094	244	722433	59.072	ng	0.00

Target Compounds	Qvalue

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

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