

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111022\
 Data File : BG055554.D
 Acq On : 11 Nov 2022 1:06
 Operator : CG/JU
 Sample : N5476-01
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RED-006-5

Quant Time: Nov 11 02:04:51 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG110922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 10 03:39:46 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.268	152	45140	20.000	ng	0.00
21) Naphthalene-d8	11.094	136	187333	20.000	ng	0.00
39) Acenaphthene-d10	14.883	164	126922	20.000	ng	0.00
64) Phenanthrene-d10	17.627	188	303446	20.000	ng	0.00
76) Chrysene-d12	21.922	240	309299	20.000	ng	0.00
86) Perylene-d12	25.342	264	324781	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.794	112	184232	77.738	ng	0.00
7) Phenol-d6	7.398	99	147046	44.919	ng	0.00
23) Nitrobenzene-d5	9.437	82	344027	124.496	ng	0.00
42) 2,4,6-Tribromophenol	16.364	330	250799	202.521	ng	0.00
45) 2-Fluorobiphenyl	13.514	172	803445	95.485	ng	0.00
79) Terphenyl-d14	20.212	244	1257456	81.658	ng	0.00

Target Compounds	Qvalue

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

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