

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG051524\
 Data File : BG061208.D
 Acq On : 15 May 2024 16:21
 Operator : MA/JU
 Sample : P2469-04
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MH-A

Quant Time: May 15 16:56:54 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG043024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 01 05:11:25 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.929	152	23769	20.000	ng	-0.03
21) Naphthalene-d8	10.720	136	92365	20.000	ng	#-0.04
39) Acenaphthene-d10	14.557	164	71543	20.000	ng	-0.03
64) Phenanthrene-d10	17.306	188	168912	20.000	ng	-0.03
76) Chrysene-d12	21.566	240	177695	20.000	ng	-0.03
86) Perylene-d12	24.692	264	207324	20.000	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.526	112	134209	114.833	ng	-0.02
7) Phenol-d6	7.106	99	168460	97.436	ng	-0.03
23) Nitrobenzene-d5	9.081	82	138656	74.417	ng	-0.04
42) 2,4,6-Tribromophenol	16.049	330	161442	112.895	ng	-0.03
45) 2-Fluorobiphenyl	13.182	172	383157	79.194	ng	-0.03
79) Terphenyl-d14	19.927	244	974928	91.249	ng	-0.03
Target Compounds						
77) Benzidine	19.545	184	28397	10.704	ng	Qvalue 97

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

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