

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012324\
 Data File : BG060411.D
 Acq On : 23 Jan 2024 19:59
 Operator : MA/JU
 Sample : P1221-03
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MH-A

Quant Time: Jan 24 03:18:42 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG010824.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 09 04:07:12 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.927	152	191847	20.000	ng	0.00
21) Naphthalene-d8	10.730	136	937666	20.000	ng	0.00
39) Acenaphthene-d10	14.560	164	658401	20.000	ng	0.00
64) Phenanthrene-d10	17.310	188	1528661	20.000	ng	0.00
76) Chrysene-d12	21.570	240	1347700	20.000	ng	0.00
86) Perylene-d12	24.737	264	1519649	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.500	112	1033957	88.646	ng	0.00
7) Phenol-d6	7.093	99	1442783	83.535	ng	0.00
23) Nitrobenzene-d5	9.084	82	916756	46.167	ng	0.00
42) 2,4,6-Tribromophenol	16.053	330	756031	78.046	ng	0.00
45) 2-Fluorobiphenyl	13.185	172	2488025	52.539	ng	0.00
79) Terphenyl-d14	19.931	244	3374603	55.253	ng	0.00
Target Compounds						
75) Fluoranthene	19.366	202	229641	2.588	ng	Qvalue 89
78) Pyrene	19.731	202	288298	3.350	ng	# 85

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

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