

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012224\
 Data File : BG060391.D
 Acq On : 22 Jan 2024 19:08
 Operator : MA/JU
 Sample : P1207-03
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EC68-WEST

Quant Time: Jan 23 00:39:12 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	7.929	152	191133	20.000	ng	0.00
21) Naphthalene-d8	10.732	136	840264	20.000	ng	0.00
39) Acenaphthene-d10	14.563	164	645429	20.000	ng	0.00
64) Phenanthrene-d10	17.312	188	1609198	20.000	ng	0.00
76) Chrysene-d12	21.572	240	1479090	20.000	ng	0.00
86) Perylene-d12	24.739	264	1697782	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.503	112	1545050	132.959	ng	0.00
7) Phenol-d6	7.095	99	2035926	118.318	ng	0.00
23) Nitrobenzene-d5	9.087	82	1272067	71.486	ng	0.00
42) 2,4,6-Tribromophenol	16.055	330	1213427	127.781	ng	0.00
45) 2-Fluorobiphenyl	13.188	172	3393995	73.110	ng	0.00
79) Terphenyl-d14	19.933	244	5209282	91.338	ng	0.00
Target Compounds						
75) Fluoranthene	19.369	202	229312	2.455	ng	Qvalue 87
78) Pyrene	19.727	202	228499	2.419	ng	# 84

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

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