

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102122\
 Data File : BG055244.D
 Acq On : 21 Oct 2022 21:53
 Operator : CG/JU
 Sample : N5217-06
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-G1-COMP

Quant Time: Oct 22 01:32:53 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG101922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 20 09:21:42 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.286	152	48860	20.000	ng	0.00
21) Naphthalene-d8	11.112	136	223699	20.000	ng	0.00
39) Acenaphthene-d10	14.895	164	154319	20.000	ng	0.00
64) Phenanthrene-d10	17.628	188	338890	20.000	ng	0.00
76) Chrysene-d12	21.905	240	233532	20.000	ng	-0.01
86) Perylene-d12	25.313	264	241755	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.806	112	321773	118.484	ng	0.00
7) Phenol-d6	7.422	99	453549	109.601	ng	0.00
23) Nitrobenzene-d5	9.455	82	281099	60.666	ng	0.00
42) 2,4,6-Tribromophenol	16.370	330	174071	108.859	ng	0.00
45) 2-Fluorobiphenyl	13.526	172	618036	63.971	ng	0.00
79) Terphenyl-d14	20.201	244	882823	70.438	ng	0.00
Target Compounds						
75) Fluoranthene	19.660	202	42817	2.137	ng	98
78) Pyrene	20.019	202	35534	2.019	ng	99

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

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