

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040722\
 Data File : BG053072.D
 Acq On : 7 Apr 2022 13:54
 Operator : CG/JU
 Sample : N2207-09RE
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SPLP-LEACHATE-AMENDED-096-21-01RE

Quant Time: Apr 08 07:22:42 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG031522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 06 14:35:39 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.123	152	37073	20.000	ng	0.00
21) Naphthalene-d8	10.937	136	135527	20.000	ng	0.00
39) Acenaphthene-d10	14.749	164	97667	20.000	ng	0.00
64) Phenanthrene-d10	17.493	188	220452	20.000	ng	0.00
76) Chrysene-d12	21.775	240	219182	20.000	ng	0.00
86) Perylene-d12	25.082	264	223554	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.691	112	550	0.258	ng	0.00
7) Phenol-d6	7.283	99	6823	2.126	ng	0.00
23) Nitrobenzene-d5	9.286	82	267578	99.093	ng	0.00
42) 2,4,6-Tribromophenol	16.235	330	650	0.495	ng	0.00
45) 2-Fluorobiphenyl	13.375	172	564174	85.195	ng	0.00
79) Terphenyl-d14	20.095	244	1097671	89.085	ng	0.00

Target Compounds	Qvalue

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

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