

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040422\
 Data File : BG053010.D
 Acq On : 4 Apr 2022 16:29
 Operator : CG/JU
 Sample : N2207-09
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

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

Quant Time: Apr 05 01:07:05 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG031522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 24 16:48:57 2022
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Christian Giraldo 04/05/2022
 Supervised By :Jagrut Upadhyay 04/05/2022

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.128	152	36989	20.000	ng	#-0.02
21) Naphthalene-d8	10.942	136	140787	20.000	ng	-0.02
39) Acenaphthene-d10	14.748	164	98252	20.000	ng	-0.02
64) Phenanthrene-d10	17.486	188	231903	20.000	ng	-0.02
76) Chrysene-d12	21.768	240	256686	20.000	ng	-0.03
86) Perylene-d12	25.070	264	260686	20.000	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.696	112	455m	0.214	ng	0.00
7) Phenol-d6	7.276	99	7732	2.415	ng	-0.01
23) Nitrobenzene-d5	9.285	82	275155	98.092	ng	-0.02
42) 2,4,6-Tribromophenol	16.229	330	628	0.476	ng	-0.02
45) 2-Fluorobiphenyl	13.374	172	578367	86.818	ng	-0.02
79) Terphenyl-d14	20.094	244	1236319	85.678	ng	-0.02

Target Compounds	Qvalue

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

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