

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF022122\
 Data File : BF126985.D
 Acq On : 21 Feb 2022 20:48
 Operator : CG\JU
 Sample : N1602-12
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SPLP-LEACHATE-AMENDED-3-COMP

Quant Time: Feb 22 03:51:48 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 21 13:16:57 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.916	152	91163	20.000	ng	0.00
21) Naphthalene-d8	8.198	136	366390	20.000	ng	# 0.00
39) Acenaphthene-d10	9.957	164	190943	20.000	ng	0.00
64) Phenanthrene-d10	11.445	188	336345	20.000	ng	# 0.00
76) Chrysene-d12	14.086	240	145078	20.000	ng	#-0.01
86) Perylene-d12	15.586	264	161576	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.540	112	459555	77.913	ng	0.00
7) Phenol-d6	6.534	99	427306	53.689	ng	-0.01
23) Nitrobenzene-d5	7.481	82	879384	108.335	ng	-0.01
42) 2,4,6-Tribromophenol	10.751	330	313194	142.461	ng	0.00
45) 2-Fluorobiphenyl	9.275	172	1461321	112.667	ng	0.00
79) Terphenyl-d14	13.033	244	1381613	139.995	ng	0.00

Target Compounds	Qvalue

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

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