

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080221\
 Data File : BF124901.D
 Acq On : 2 Aug 2021 11:27
 Operator : JU/CG
 Sample : M3216-11
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SPLP-LEACHATE-AMENDED-COM-

Quant Time: Aug 02 11:59:33 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF072221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 02 11:15:51 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.851	152	5170	20.000	ng	0.00
21) Naphthalene-d8	8.134	136	19988	20.000	ng	0.00
39) Acenaphthene-d10	9.892	164	11134	20.000	ng	0.00
64) Phenanthrene-d10	11.375	188	20832	20.000	ng	0.00
76) Chrysene-d12	14.010	240	14488	20.000	ng	-0.01
86) Perylene-d12	15.480	264	11150	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.463	112	18426	60.345	ng	-0.01
7) Phenol-d6	6.475	99	13820	36.103	ng	-0.01
23) Nitrobenzene-d5	7.422	82	37077	110.403	ng	0.00
42) 2,4,6-Tribromophenol	10.680	330	14957	156.460	ng	0.00
45) 2-Fluorobiphenyl	9.216	172	73089	96.416	ng	0.00
79) Terphenyl-d14	12.963	244	88732	104.983	ng	0.00

Target Compounds	Qvalue

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

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