

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112018\
 Data File : BF110898.D
 Acq On : 20 Nov 2018 17:01
 Operator : JU/SJ
 Sample : J5993-12
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

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

Quant Time: Nov 21 04:28:07 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 13 12:43:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.94	152	54400	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	223882	20.00	ng	0.00
39) Acenaphthene-d10	9.98	164	103791	20.00	ng	-0.01
64) Phenanthrene-d10	11.48	188	177611	20.00	ng	0.00
76) Chrysene-d12	14.13	240	125651	20.00	ng	0.00
87) Perylene-d12	15.67	264	109583	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	134942	40.29	ng	0.00
7) Phenol-d6	6.57	99	103431	24.15	ng	0.00
23) Nitrobenzene-d5	7.50	82	284601	73.21	ng	-0.01
42) 2,4,6-Tribromophenol	10.78	330	93583	89.48	ng	0.00
45) 2-Fluorobiphenyl	9.30	172	538106	74.04	ng	0.00
79) Terphenyl-d14	13.06	244	389893	58.11	ng	-0.01
Target Compounds						
10) Phenol	6.58	94	18174	3.660	ng	Qvalue # 77
31) Naphthalene	8.25	128	166134	15.807	ng	99
37) 2-Methylnaphthalene	8.94	142	35875	5.207	ng	100
38) 1-Methylnaphthalene	9.04	142	29463	4.447	ng	# 95
50) Dimethylphthalate	9.69	163	63762	8.233	ng	98
71) Phenanthrene	11.50	178	19577	2.057	ng	97

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

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