

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF103020\
 Data File : BF122213.D
 Acq On : 31 Oct 2020 10:40
 Operator : JU/CG
 Sample : L4568-02
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RECLAMATION-PIT

Quant Time: Nov 01 23:51:41 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 30 10:16:54 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.710	152	88298	20.00	ng	0.00
21) Naphthalene-d8	7.992	136	322821	20.00	ng	0.00
39) Acenaphthene-d10	9.745	164	142942	20.00	ng	0.00
64) Phenanthrene-d10	11.233	188	217942	20.00	ng	0.00
76) Chrysene-d12	13.880	240	195475	20.00	ng	0.00
86) Perylene-d12	15.315	264	186014	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.351	112	682826	114.14	ng	0.02
7) Phenol-d6	6.375	99	750415	97.44	ng	0.00
23) Nitrobenzene-d5	7.281	82	607539	96.52	ng	0.00
42) 2,4,6-Tribromophenol	10.539	330	202916	114.76	ng	0.00
45) 2-Fluorobiphenyl	9.069	172	952808	98.08	ng	0.00
79) Terphenyl-d14	12.816	244	777621	75.81	ng	0.00
Target Compounds						
10) Phenol	6.387	94	56206	7.07	ng	Qvalue 88
20) 3+4-Methylphenols	7.157	107	115026	18.45	ng	# 51
50) Dimethylphthalate	9.469	163	72281	6.92	ng	99

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

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