

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF031524\
 Data File : BF137620.D
 Acq On : 15 Mar 2024 20:32
 Operator : CG\JU
 Sample : P1769-01DL 50X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-2DL

Quant Time: Mar 15 23:24:53 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF022924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 14 14:45:20 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.728	152	121800	20.000	ng	0.00
21) Naphthalene-d8	8.004	136	475231	20.000	ng	0.00
39) Acenaphthene-d10	9.757	164	263433	20.000	ng	0.00
64) Phenanthrene-d10	11.239	188	467263	20.000	ng	0.00
76) Chrysene-d12	13.874	240	311826	20.000	ng	0.00
86) Perylene-d12	15.298	264	317708	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.357	112	7680	0.955	ng	0.02
7) Phenol-d6	6.363	99	8749	0.754	ng	0.00
23) Nitrobenzene-d5	7.293	82	17815	1.742	ng	0.00
42) 2,4,6-Tribromophenol	10.551	330	5900	2.550	ng	0.00
45) 2-Fluorobiphenyl	9.081	172	39202	2.329	ng	0.00
79) Terphenyl-d14	12.822	244	45956	2.027	ng	0.00
Target Compounds						
10) Phenol	6.375	94	1308268	104.707	ng	Qvalue 96
17) 2-Methylphenol	6.981	107	369689	46.577	ng	99
20) 3+4-Methylphenols	7.134	107	348990	38.392	ng	# 53
27) 2,4-Dimethylphenol	7.663	122	158329	20.773	ng	98

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

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