

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101024\
 Data File : BF139744.D
 Acq On : 10 Oct 2024 14:08
 Operator : RC/JU
 Sample : P4340-04
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 1B-7

Quant Time: Oct 10 14:45:51 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF100124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 02 04:58:47 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.869	152	79835	20.000	ng	0.00
21) Naphthalene-d8	8.145	136	298273	20.000	ng	0.00
39) Acenaphthene-d10	9.904	164	165062	20.000	ng	0.00
64) Phenanthrene-d10	11.386	188	286061	20.000	ng	0.00
76) Chrysene-d12	14.027	240	167873	20.000	ng	0.00
86) Perylene-d12	15.504	264	172959	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.499	112	589448	127.971	ng	0.01
7) Phenol-d6	6.504	99	702672	113.440	ng	0.00
23) Nitrobenzene-d5	7.428	82	559843	95.030	ng	0.00
42) 2,4,6-Tribromophenol	10.692	330	212038	133.072	ng	0.00
45) 2-Fluorobiphenyl	9.222	172	1012439	94.163	ng	0.00
79) Terphenyl-d14	12.975	244	896713	87.647	ng	0.00
Target Compounds						
20) 3+4-Methylphenols	7.281	107	16169	3.125	ng	Qvalue # 48
27) 2,4-Dimethylphenol	7.804	122	8372	2.608	ng	95

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

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