

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF100322\
 Data File : BF130501.D
 Acq On : 03 Oct 2022 16:28
 Operator : CG\JU
 Sample : N4888-06
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 REGULATORY-FOUNDATION-1

Quant Time: Oct 04 03:16:33 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 28 02:59:31 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.633	152	84115	20.000	ng	-0.01
21) Naphthalene-d8	7.922	136	355975	20.000	ng	0.00
39) Acenaphthene-d10	9.669	164	118739	20.000	ng	-0.01
64) Phenanthrene-d10	11.163	188	163891	20.000	ng	# 0.00
76) Chrysene-d12	13.786	240	147501	20.000	ng	# 0.00
86) Perylene-d12	15.174	264	155827	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.239	112	368963	76.081	ng	0.00
7) Phenol-d6	6.275	99	600450	98.954	ng	-0.02
23) Nitrobenzene-d5	7.204	82	440402	63.205	ng	-0.01
42) 2,4,6-Tribromophenol	10.463	330	56619	49.764	ng	0.00
45) 2-Fluorobiphenyl	8.998	172	600120	73.597	ng	-0.01
79) Terphenyl-d14	12.745	244	448437	50.361	ng	0.00

Target Compounds	Qvalue

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

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