

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF031122\
 Data File : BF127302.D
 Acq On : 11 Mar 2022 16:22
 Operator : CG\JU
 Sample : PB143171BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB143171BL

Quant Time: Mar 11 17:26:34 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF031022.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 11 12:42:13 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.869	152	123546	20.000	ng	0.00
21) Naphthalene-d8	8.151	136	457106	20.000	ng	0.00
39) Acenaphthene-d10	9.910	164	267089	20.000	ng	0.00
64) Phenanthrene-d10	11.398	188	480028	20.000	ng	# 0.00
76) Chrysene-d12	14.045	240	381516	20.000	ng	# 0.00
86) Perylene-d12	15.533	264	286002	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	917975	119.445	ng	0.02
7) Phenol-d6	6.504	99	1279306	121.005	ng	0.00
23) Nitrobenzene-d5	7.434	82	913203	82.505	ng	0.00
42) 2,4,6-Tribromophenol	10.704	330	363693	126.493	ng	0.00
45) 2-Fluorobiphenyl	9.228	172	1550794	82.089	ng	0.00
79) Terphenyl-d14	12.992	244	1816767	81.999	ng	0.00

Target Compounds	Qvalue

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

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