

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF031722\
 Data File : BF127371.D
 Acq On : 17 Mar 2022 12:53
 Operator : CG\JU
 Sample : N1943-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B1

Quant Time: Mar 17 13:46:51 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF031022.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 10 16:51:53 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.863	152	90275	20.000	ng	0.00
21) Naphthalene-d8	8.145	136	334251	20.000	ng	#-0.01
39) Acenaphthene-d10	9.904	164	177641	20.000	ng	-0.01
64) Phenanthrene-d10	11.398	188	287980	20.000	ng	# 0.00
76) Chrysene-d12	14.045	240	151522	20.000	ng	# 0.00
86) Perylene-d12	15.533	264	166755	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.492	112	510687	90.940	ng	0.00
7) Phenol-d6	6.498	99	709819	91.884	ng	0.00
23) Nitrobenzene-d5	7.433	82	480903	59.418	ng	0.00
42) 2,4,6-Tribromophenol	10.698	330	97893	51.191	ng	-0.01
45) 2-Fluorobiphenyl	9.222	172	817229	65.041	ng	-0.01
79) Terphenyl-d14	12.986	244	649595	73.822	ng	0.00

Target Compounds	Qvalue

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

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