

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062122\
 Data File : BF128911.D
 Acq On : 21 Jun 2022 21:55
 Operator : CG\JU
 Sample : N3373-11DL 10X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SC-13WDL

Quant Time: Jun 22 01:56:03 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 22 01:02:29 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.751	152	113247	20.000	ng	0.00
21) Naphthalene-d8	8.039	136	405132	20.000	ng	# 0.00
39) Acenaphthene-d10	9.792	164	192483	20.000	ng	0.00
64) Phenanthrene-d10	11.280	188	286627	20.000	ng	# 0.00
76) Chrysene-d12	13.921	240	232140	20.000	ng	# 0.00
86) Perylene-d12	15.362	264	159596	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.392	112	60725	8.541	ng	0.00
7) Phenol-d6	6.416	99	75971	8.754	ng	0.00
23) Nitrobenzene-d5	7.316	82	53790	6.182	ng	-0.01
42) 2,4,6-Tribromophenol	10.580	330	11564	5.085	ng	0.00
45) 2-Fluorobiphenyl	9.110	172	97368	7.086	ng	0.00
79) Terphenyl-d14	12.863	244	81201	6.078	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	11.304	178	757355	52.680	ng	100
75) Fluoranthene	12.492	202	317527	21.129	ng	95

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

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