

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062122\
 Data File : BF128914.D
 Acq On : 21 Jun 2022 23:22
 Operator : CG\JU
 Sample : N3373-10DL 10X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SC-13EDL

Quant Time: Jun 22 01:57:07 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	110618	20.000	ng	0.00
21) Naphthalene-d8	8.039	136	403353	20.000	ng	# 0.00
39) Acenaphthene-d10	9.792	164	197703	20.000	ng	0.00
64) Phenanthrene-d10	11.274	188	303509	20.000	ng	0.00
76) Chrysene-d12	13.915	240	248483	20.000	ng	#-0.01
86) Perylene-d12	15.362	264	174593	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.393	112	50447	7.264	ng	0.00
7) Phenol-d6	6.416	99	68571	8.089	ng	0.00
23) Nitrobenzene-d5	7.316	82	48306	5.576	ng	-0.01
42) 2,4,6-Tribromophenol	10.580	330	2472	1.058	ng	0.00
45) 2-Fluorobiphenyl	9.110	172	95533	6.769	ng	0.00
79) Terphenyl-d14	12.857	244	82476	5.768	ng	-0.01
Target Compounds						
						Qvalue
71) Phenanthrene	11.304	178	811395	53.299	ng	99
75) Fluoranthene	12.492	202	154213	9.691	ng	96

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

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