

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062122\
 Data File : BF128913.D
 Acq On : 21 Jun 2022 22:53
 Operator : CG\JU
 Sample : N3373-09DL 10X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SC-13SDL

Quant Time: Jun 22 01:56:44 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	114941	20.000	ng	0.00
21) Naphthalene-d8	8.039	136	418457	20.000	ng	# 0.00
39) Acenaphthene-d10	9.792	164	202314	20.000	ng	0.00
64) Phenanthrene-d10	11.280	188	308230	20.000	ng	# 0.00
76) Chrysene-d12	13.916	240	243771	20.000	ng	#-0.01
86) Perylene-d12	15.363	264	170333	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.393	112	48966	6.786	ng	0.00
7) Phenol-d6	6.416	99	68334	7.758	ng	0.00
23) Nitrobenzene-d5	7.316	82	48595	5.407	ng	-0.01
42) 2,4,6-Tribromophenol	10.586	330	1712	0.716	ng	0.00
45) 2-Fluorobiphenyl	9.110	172	96490	6.681	ng	0.00
79) Terphenyl-d14	12.857	244	79579	5.673	ng	-0.01
Target Compounds						
						Qvalue
71) Phenanthrene	11.304	178	1117202	72.263	ng	99
75) Fluoranthene	12.492	202	185243	11.463	ng	95

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

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