

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092022\
 Data File : BF130352.D
 Acq On : 20 Sep 2022 19:30
 Operator : CG\JU
 Sample : N4714-01 5X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OK-01-091922

Quant Time: Sep 21 05:17:54 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 19 15:06:14 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.663	152	116603	20.000	ng	0.00
21) Naphthalene-d8	7.945	136	407461	20.000	ng	# 0.00
39) Acenaphthene-d10	9.692	164	186087	20.000	ng	0.00
64) Phenanthrene-d10	11.175	188	266920	20.000	ng	0.00
76) Chrysene-d12	13.810	240	191857	20.000	ng	# 0.00
86) Perylene-d12	15.204	264	142573	20.000	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.263	112	138767	20.642	ng	0.00
7) Phenol-d6	6.304	99	161320	19.178	ng	0.00
23) Nitrobenzene-d5	7.222	82	100196	12.563	ng	-0.01
42) 2,4,6-Tribromophenol	10.480	330	32989	18.501	ng	0.00
45) 2-Fluorobiphenyl	9.022	172	183265	14.341	ng	0.00
79) Terphenyl-d14	12.757	244	154072	13.303	ng	0.00
Target Compounds						
88) Benzo(b)fluoranthene	14.810	252	27278	2.931	ng	# 71
90) Benzo(a)pyrene	15.145	252	19274	2.614	ng	# 62

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

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