

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111022\
 Data File : BF131120.D
 Acq On : 10 Nov 2022 13:24
 Operator : CG\JU
 Sample : N5378-07
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-15-20221028-12.5-13.5

Quant Time: Nov 11 00:46:58 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 10 04:15:17 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.887	152	74020	20.000	ng	0.00
21) Naphthalene-d8	8.169	136	214877	20.000	ng	-0.01
39) Acenaphthene-d10	9.934	164	116729	20.000	ng	0.00
64) Phenanthrene-d10	11.428	188	195531	20.000	ng	# 0.00
76) Chrysene-d12	14.074	240	112085	20.000	ng	0.00
86) Perylene-d12	15.568	264	111084	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	415827	95.537	ng	0.00
7) Phenol-d6	6.516	99	534027	93.847	ng	0.00
23) Nitrobenzene-d5	7.451	82	325175	68.279	ng	-0.01
42) 2,4,6-Tribromophenol	10.722	330	106257	88.080	ng	-0.01
45) 2-Fluorobiphenyl	9.251	172	478790	57.259	ng	0.00
79) Terphenyl-d14	13.016	244	517791	80.997	ng	0.00
Target Compounds						
87) Indeno(1,2,3-cd)pyrene	17.080	276	4127	4.282	ng	# 91
92) Benzo(g,h,i)perylene	17.539	276	4660	4.908	ng	# 95

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

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