

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111022\
 Data File : BF131133.D
 Acq On : 10 Nov 2022 20:39
 Operator : CG\JU
 Sample : N5520-05
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 124042

Quant Time: Nov 11 00:51:11 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	77348	20.000	ng	0.00
21) Naphthalene-d8	8.169	136	271240	20.000	ng	-0.01
39) Acenaphthene-d10	9.933	164	136623	20.000	ng	0.00
64) Phenanthrene-d10	11.422	188	202481	20.000	ng	0.00
76) Chrysene-d12	14.074	240	161994	20.000	ng	0.00
86) Perylene-d12	15.568	264	162858	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	432861	95.171	ng	0.00
7) Phenol-d6	6.516	99	563808	94.818	ng	0.00
23) Nitrobenzene-d5	7.451	82	353641	58.826	ng	-0.01
42) 2,4,6-Tribromophenol	10.722	330	113323	80.259	ng	-0.01
45) 2-Fluorobiphenyl	9.251	172	568062	58.043	ng	0.00
79) Terphenyl-d14	13.010	244	401699	43.477	ng	0.00
Target Compounds						
70) Pentachlorophenol	11.233	266	2193	7.098	ng	# 81
84) Bis(2-ethylhexyl)phtha...	14.045	149	14599	2.718	ng	97
87) Indeno(1,2,3-cd)pyrene	17.080	276	5213	4.225	ng	# 95
92) Benzo(g,h,i)perylene	17.545	276	4998	4.763	ng	# 90

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

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