

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060322\
 Data File : BP010698.D
 Acq On : 03 Jun 2022 18:13
 Operator : CG/JU
 Sample : N3170-07
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 20220409-FIELD-BLANK-2

Quant Time: Jun 04 02:01:26 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP052822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jun 04 01:56:43 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.852	152	115751	20.000	ng	0.00
21) Naphthalene-d8	10.651	136	456855	20.000	ng	0.00
39) Acenaphthene-d10	14.486	164	302050	20.000	ng	-0.01
64) Phenanthrene-d10	17.245	188	653669	20.000	ng	0.00
76) Chrysene-d12	21.333	240	639012	20.000	ng	#-0.01
86) Perylene-d12	23.739	264	651012	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.440	112	382568	60.295	ng	0.00
7) Phenol-d6	7.034	99	294676	32.714	ng	0.00
23) Nitrobenzene-d5	9.010	82	884019	91.484	ng	-0.01
42) 2,4,6-Tribromophenol	15.986	330	652959	190.976	ng	0.00
45) 2-Fluorobiphenyl	13.110	172	1938151	91.068	ng	0.00
79) Terphenyl-d14	19.804	244	3605800	105.989	ng	0.00

Target Compounds	Qvalue

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

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