

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF103122\
 Data File : BF130884.D
 Acq On : 31 Oct 2022 17:42
 Operator : CG\JU
 Sample : N5329-03
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CBH-594

Quant Time: Nov 01 02:29:42 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF102722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Oct 29 05:41:34 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.940	152	78111	20.000	ng	0.00
21) Naphthalene-d8	8.228	136	280931	20.000	ng	# 0.00
39) Acenaphthene-d10	9.986	164	131280	20.000	ng	-0.01
64) Phenanthrene-d10	11.480	188	206714	20.000	ng	0.00
76) Chrysene-d12	14.127	240	159910	20.000	ng	0.00
86) Perylene-d12	15.645	264	124363	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.557	112	307226	68.186	ng	0.00
7) Phenol-d6	6.551	99	236823	40.447	ng	-0.01
23) Nitrobenzene-d5	7.504	82	547179	117.446	ng	-0.01
42) 2,4,6-Tribromophenol	10.780	330	172841	156.401	ng	0.00
45) 2-Fluorobiphenyl	9.304	172	903225	104.020	ng	0.00
79) Terphenyl-d14	13.074	244	780815	89.476	ng	0.00
Target Compounds						
						Qvalue
38) 1-Methylnaphthalene	9.039	142	27367	3.050	ng	97
52) Acenaphthene	10.022	154	119044	16.244	ng	99
58) Fluorene	10.533	166	50899	6.085	ng	98
71) Phenanthrene	11.504	178	23341	2.188	ng	98
75) Fluoranthene	12.698	202	27010	2.383	ng	98

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

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