

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071522\
 Data File : BM035846.D
 Acq On : 16 Jul 2022 00:52
 Operator : CG/JU
 Sample : N3651-18
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DUP-FILT

Quant Time: Jul 16 03:02:39 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM071322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 14 01:21:21 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.904	152	348362	20.000	ng	0.00
21) Naphthalene-d8	10.710	136	1535313	20.000	ng	0.00
39) Acenaphthene-d10	14.533	164	957310	20.000	ng	0.00
64) Phenanthrene-d10	17.274	188	1920578	20.000	ng	0.00
76) Chrysene-d12	21.445	240	1455513	20.000	ng	-0.01
86) Perylene-d12	23.821	264	1258695	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.463	112	972638	45.350	ng	0.00
7) Phenol-d6	7.063	99	756707	28.207	ng	0.00
23) Nitrobenzene-d5	9.069	82	2142946	79.793	ng	0.00
42) 2,4,6-Tribromophenol	16.021	330	1287468	132.434	ng	0.00
45) 2-Fluorobiphenyl	13.169	172	4926669	71.772	ng	0.00
79) Terphenyl-d14	19.898	244	4689166	62.539	ng	0.00
Target Compounds						
74) Di-n-butylphthalate	18.251	149	13895	2.390	ng	99
84) Bis(2-ethylhexyl)phtha...	21.374	149	66713	4.511	ng	# 98

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

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