

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111821\
 Data File : BP008023.D
 Acq On : 18 Nov 2021 19:01
 Operator : CG/JU
 Sample : M4724-17
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SUB-FLOOR-BRICK

Quant Time: Nov 19 02:24:25 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 15 05:36:49 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.834	152	91072	20.000	ng	0.00
21) Naphthalene-d8	10.628	136	323778	20.000	ng	-0.01
39) Acenaphthene-d10	14.469	164	206435	20.000	ng	-0.01
64) Phenanthrene-d10	17.228	188	450809	20.000	ng	#-0.01
76) Chrysene-d12	21.322	240	602210	20.000	ng	#-0.01
86) Perylene-d12	23.722	264	746927	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.423	112	600254	106.921	ng	0.00
7) Phenol-d6	7.011	99	713037	95.471	ng	-0.01
23) Nitrobenzene-d5	8.987	82	435372	71.958	ng	-0.02
42) 2,4,6-Tribromophenol	15.963	330	348353	107.251	ng	-0.01
45) 2-Fluorobiphenyl	13.093	172	1036809	70.451	ng	-0.01
79) Terphenyl-d14	19.792	244	1934395	65.133	ng	-0.01
Target Compounds						
						Qvalue
74) Di-n-butylphthalate	18.192	149	59643	2.273	ng	98
80) Butylbenzylphthalate	20.475	149	34301	2.307	ng	97
83) Chrysene	21.357	228	86663	2.281	ng	98
84) Bis(2-ethylhexyl)phtha...	21.239	149	51862	2.443	ng	# 98
85) Di-n-octyl phthalate	22.163	149	103376	2.642	ng	# 96
88) Benzo(b)fluoranthene	22.992	252	103793	2.334	ng	98
89) Benzo(k)fluoranthene	23.039	252	96645	2.176	ng	99

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

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