

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081621\
 Data File : BM031774.D
 Acq On : 17 Aug 2021 05:30
 Operator : CG/JU
 Sample : M3425-18
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 23N

Quant Time: Aug 17 07:29:10 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM081621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 16 16:42:38 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.551	152	42131	20.000	ng	# 0.00
21) Naphthalene-d8	10.310	136	177941	20.000	ng	# 0.00
39) Acenaphthene-d10	14.186	164	115329	20.000	ng	0.00
64) Phenanthrene-d10	16.939	188	253280	20.000	ng	0.00
76) Chrysene-d12	21.139	240	238257	20.000	ng	0.00
86) Perylene-d12	23.285	264	211477	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.187	112	368070	140.218	ng	0.00
7) Phenol-d6	6.745	99	441851	116.164	ng	0.00
23) Nitrobenzene-d5	8.704	82	359556	84.607	ng	0.00
42) 2,4,6-Tribromophenol	15.686	330	197598	139.111	ng	0.00
45) 2-Fluorobiphenyl	12.798	172	740722	85.487	ng	0.00
79) Terphenyl-d14	19.586	244	1424075	100.705	ng	0.00

Target Compounds	Qvalue

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

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