

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012721\
 Data File : BM028581.D
 Acq On : 27 Jan 2021 22:11
 Operator : CG/JU
 Sample : M1235-03 10X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 FK-SF-SAMPLE-2

Quant Time: Jan 27 23:35:37 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM012021.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 27 13:20:04 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.969	152	31348	20.00	ng	0.00
21) Naphthalene-d8	10.780	136	123561	20.00	ng	# 0.00
39) Acenaphthene-d10	14.615	164	69172	20.00	ng	0.00
64) Phenanthrene-d10	17.345	188	142874	20.00	ng	# 0.00
76) Chrysene-d12	21.480	240	140275	20.00	ng	# 0.00
86) Perylene-d12	23.750	264	145378	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.522	112	5892	2.97	ng	-0.01
7) Phenol-d6	7.169	99	15587	5.77	ng	-0.02
23) Nitrobenzene-d5	9.139	82	10117	3.88	ng	-0.01
42) 2,4,6-Tribromophenol	16.098	330	60	0.07	ng	-0.01
45) 2-Fluorobiphenyl	13.233	172	23810	4.32	ng	0.00
79) Terphenyl-d14	19.939	244	32883	4.23	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	17.386	178	33660	3.83	ng	98
75) Fluoranthene	19.386	202	42643	4.32	ng	96
78) Pyrene	19.745	202	38361	3.81	ng	98
88) Benzo(b)fluoranthene	23.068	252	21714	2.20	ng	# 75

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

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