

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061121\
 Data File : BM030404.D
 Acq On : 11 Jun 2021 14:49
 Operator : CG/JU
 Sample : M2612-33
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 18-B12(54-58)

Quant Time: Jun 11 15:20:57 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM060921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 11 14:10:50 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.839	152	160287	20.000	ng	0.00
21) Naphthalene-d8	10.628	136	565518	20.000	ng	0.00
39) Acenaphthene-d10	14.457	164	313300	20.000	ng	0.00
64) Phenanthrene-d10	17.198	188	607554	20.000	ng	0.00
76) Chrysene-d12	21.356	240	645662	20.000	ng	# 0.00
86) Perylene-d12	23.633	264	773069	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.422	112	1328413	134.730	ng	0.00
7) Phenol-d6	6.998	99	1451097	101.664	ng	0.00
23) Nitrobenzene-d5	8.992	82	1080681	93.097	ng	0.00
42) 2,4,6-Tribromophenol	15.945	330	526275	133.605	ng	0.00
45) 2-Fluorobiphenyl	13.086	172	2215727	93.004	ng	0.00
79) Terphenyl-d14	19.809	244	3251019	88.440	ng	0.00

Target Compounds	Qvalue

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061121\
Data File : BM030404.D
Acq On : 11 Jun 2021 14:49
Operator : CG/JU
Sample : M2612-33
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
18-B12(54-58)

Quant Time: Jun 11 15:20:57 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM060921.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 11 14:10:50 2021
Response via : Initial Calibration

