

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111021\
 Data File : BN017378.D
 Acq On : 10 Nov 2021 10:04
 Operator : CG/JU
 Sample : PB140437BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB140437BL

Quant Time: Nov 10 11:04:22 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN110921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 10 11:00:44 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.542	152	28909	0.400	ng	0.00
7) Naphthalene-d8	10.307	136	128623	0.400	ng	# 0.00
13) Acenaphthene-d10	14.181	164	66597	0.400	ng	0.00
19) Phenanthrene-d10	16.934	188	121188	0.400	ng	0.00
29) Chrysene-d12	21.144	240	100404	0.400	ng	# 0.01
35) Perylene-d12	23.339	264	90589	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.182	112	21586	0.324	ng	0.02
5) Phenol-d6	6.740	99	19338	0.263	ng	0.00
8) Nitrobenzene-d5	8.684	82	24576	0.318	ng	0.00
11) 2-Methylnaphthalene-d10	11.911	152	63372	0.335	ng	0.00
14) 2,4,6-Tribromophenol	15.691	330	3018	0.121	ng	0.01
15) 2-Fluorobiphenyl	12.798	172	83050	0.328	ng	0.00
27) Fluoranthene-d10	18.975	212	104106	0.337	ng	0.00
31) Terphenyl-d14	19.590	244	88626	0.396	ng	0.00

Target Compounds Qvalue

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111021\
Data File : BN017378.D
Acq On : 10 Nov 2021 10:04
Operator : CG/JU
Sample : PB140437BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
PB140437BL

Quant Time: Nov 10 11:04:22 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN110921.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Nov 10 11:00:44 2021
Response via : Initial Calibration

