

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092822\
 Data File : BP011833.D
 Acq On : 28 Sep 2022 16:54
 Operator : CG/JU
 Sample : PB147943TB
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB147943TB

Quant Time: Sep 29 02:28:02 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 28 02:03:33 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.905	152	340288	20.000	ng	0.00
21) Naphthalene-d8	10.716	136	1429416	20.000	ng	# 0.00
39) Acenaphthene-d10	14.551	164	858360	20.000	ng	0.00
64) Phenanthrene-d10	17.304	188	1775751	20.000	ng	0.00
76) Chrysene-d12	21.386	240	1462389	20.000	ng	-0.01
86) Perylene-d12	23.821	264	1220442	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.469	112	2987145	154.024	ng	0.00
7) Phenol-d6	7.069	99	3851340	134.305	ng	0.00
23) Nitrobenzene-d5	9.075	82	2116759	77.509	ng	0.00
42) 2,4,6-Tribromophenol	16.039	330	1046725	181.102	ng	0.00
45) 2-Fluorobiphenyl	13.175	172	4776338	87.996	ng	0.00
79) Terphenyl-d14	19.857	244	6582590	103.499	ng	-0.01

Target Compounds	Qvalue

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092822\
Data File : BP011833.D
Acq On : 28 Sep 2022 16:54
Operator : CG/JU
Sample : PB147943TB
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
PB147943TB

Quant Time: Sep 29 02:28:02 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092122.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Sep 28 02:03:33 2022
Response via : Initial Calibration

