

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092220\
 Data File : BP003351.D
 Acq On : 22 Sep 2020 22:01
 Operator : CG/JU
 Sample : L4103-11 10X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 T-1

Quant Time: Sep 23 03:35:36 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP091520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 22 12:48:08 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.569	152	149908	20.00	ng	0.00
21) Naphthalene-d8	10.352	136	621710	20.00	ng	# 0.00
39) Acenaphthene-d10	14.222	164	382707	20.00	ng	0.00
64) Phenanthrene-d10	16.981	188	713754	20.00	ng	# 0.00
76) Chrysene-d12	21.075	240	795285	20.00	ng	0.00
86) Perylene-d12	23.269	264	748803	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.187	112	11216	1.31	ng	0.00
7) Phenol-d6	6.781	99	12698	1.11	ng	0.01
23) Nitrobenzene-d5	8.740	82	7620	0.72	ng	0.02
42) 2,4,6-Tribromophenol	15.728	330	5604	0.96	ng	0.00
45) 2-Fluorobiphenyl	12.840	172	24895	0.95	ng	0.00
79) Terphenyl-d14	19.557	244	38173	1.00	ng	0.00
Target Compounds						
71) Phenanthrene	17.022	178	134805	3.47	ng	# 77

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092220\
Data File : BP003351.D
Acq On : 22 Sep 2020 22:01
Operator : CG/JU
Sample : L4103-11 10X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
T-1

Quant Time: Sep 23 03:35:36 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP091520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Sep 22 12:48:08 2020
Response via : Initial Calibration

