

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092221\
 Data File : BP007097.D
 Acq On : 22 Sep 2021 14:20
 Operator : CG/JU
 Sample : PB139245BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB139245BL

Quant Time: Sep 22 15:12:22 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 21 16:50:24 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.899	152	266446	20.000	ng	0.00
21) Naphthalene-d8	10.704	136	1075285	20.000	ng	# 0.00
39) Acenaphthene-d10	14.528	164	694619	20.000	ng	0.00
64) Phenanthrene-d10	17.280	188	1543899	20.000	ng	# 0.00
76) Chrysene-d12	21.363	240	1494704	20.000	ng	# 0.00
86) Perylene-d12	23.780	264	1507170	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.475	112	2003725	125.884	ng	0.00
7) Phenol-d6	7.069	99	2450910	112.660	ng	0.00
23) Nitrobenzene-d5	9.063	82	1391081	75.339	ng	0.00
42) 2,4,6-Tribromophenol	16.022	330	1027000	114.043	ng	0.00
45) 2-Fluorobiphenyl	13.157	172	3596706	74.181	ng	0.00
79) Terphenyl-d14	19.839	244	6003982	76.973	ng	0.00

Target Compounds	Qvalue

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092221\
Data File : BP007097.D
Acq On : 22 Sep 2021 14:20
Operator : CG/JU
Sample : PB139245BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
PB139245BL

Quant Time: Sep 22 15:12:22 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Sep 21 16:50:24 2021
Response via : Initial Calibration

