

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091922\
 Data File : BP011743.D
 Acq On : 19 Sep 2022 19:49
 Operator : CG/JU
 Sample : N4695-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 INCINERATOR-ASH

Quant Time: Sep 20 04:48:33 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 20 04:42:52 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.940	152	473806	20.000	ng	0.00
21) Naphthalene-d8	10.746	136	1981872	20.000	ng	# 0.00
39) Acenaphthene-d10	14.575	164	1275209	20.000	ng	0.00
64) Phenanthrene-d10	17.328	188	3013361	20.000	ng	0.00
76) Chrysene-d12	21.404	240	3041482	20.000	ng	0.00
86) Perylene-d12	23.851	264	960658	20.000	ng	# 0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.499	112	4306959	165.459	ng	0.00
7) Phenol-d6	7.099	99	5343407	153.994	ng	0.00
23) Nitrobenzene-d5	9.105	82	3317253	99.268	ng	0.00
42) 2,4,6-Tribromophenol	16.069	330	2837114	284.746	ng	0.00
45) 2-Fluorobiphenyl	13.204	172	8299685	101.170	ng	0.00
79) Terphenyl-d14	19.875	244	13308868	95.155	ng	0.00
Target Compounds						
32) Benzoic acid	9.981	122	75125	9.753	ng	Qvalue 88

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091922\
Data File : BP011743.D
Acq On : 19 Sep 2022 19:49
Operator : CG/JU
Sample : N4695-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
INCINERATOR-ASH

Quant Time: Sep 20 04:48:33 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Sep 20 04:42:52 2022
Response via : Initial Calibration

