

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011221\
 Data File : BP004511.D
 Acq On : 12 Jan 2021 18:47
 Operator : CG/JU
 Sample : L5221-09
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 124021

Quant Time: Jan 12 23:30:23 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270E-BP011221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 12 15:19:32 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.811	152	232271	20.00	ng	0.00
21) Naphthalene-d8	10.610	136	870100	20.00	ng	# 0.00
39) Acenaphthene-d10	14.463	164	526133	20.00	ng	0.00
64) Phenanthrene-d10	17.222	188	1106594	20.00	ng	0.00
76) Chrysene-d12	21.316	240	1112646	20.00	ng	0.00
86) Perylene-d12	23.669	264	1137530	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.405	112	1546538	116.05	ng	0.00
7) Phenol-d6	6.987	99	1652448	90.53	ng	0.00
23) Nitrobenzene-d5	8.969	82	1216299	79.03	ng	0.00
42) 2,4,6-Tribromophenol	15.963	330	1089933	119.15	ng	0.00
45) 2-Fluorobiphenyl	13.081	172	3240578	80.07	ng	0.00
79) Terphenyl-d14	19.786	244	4661410	76.54	ng	0.00
Target Compounds						
10) Phenol	7.011	94	245389	14.02	ng	Qvalue 95
50) Dimethylphthalate	13.916	163	147766	3.39	ng	98
70) Pentachlorophenol	16.875	266	252147	31.80	ng	97

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011221\
Data File : BP004511.D
Acq On : 12 Jan 2021 18:47
Operator : CG/JU
Sample : L5221-09
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
124021

Quant Time: Jan 12 23:30:23 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270E-BP011221.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jan 12 15:19:32 2021
Response via : Initial Calibration

