

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP100719\
 Data File : BP000557.D
 Acq On : 07 Oct 2019 20:57
 Operator : JU
 Sample : K5213-09
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 W-1

Quant Time: Oct 08 10:55:42 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP100119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 01 18:03:42 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.76	152	266368	20.00	ng	0.00
21) Naphthalene-d8	8.05	136	933811	20.00	ng	0.00
39) Acenaphthene-d10	9.82	164	487439	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	854957	20.00	ng	0.00
76) Chrysene-d12	13.99	240	760125	20.00	ng	0.01
87) Perylene-d12	15.47	264	800531	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.39	112	1373401	82.88	ng	0.00
7) Phenol-d6	6.41	99	1242940	62.24	ng	0.00
23) Nitrobenzene-d5	7.33	82	1227802	80.30	ng	0.00
42) 2,4,6-Tribromophenol	10.62	330	638288	122.88	ng	0.00
45) 2-Fluorobiphenyl	9.13	172	2428475	80.61	ng	0.00
79) Terphenyl-d14	12.93	244	2811452	74.89	ng	0.01
Target Compounds						
10) Phenol	6.42	94	143252	7.007	ng	92
38) 1-Methylnaphthalene	8.86	142	181366	6.556	ng	# 97
50) Dimethylphthalate	9.53	163	553471	16.341	ng	# 89
58) Fluorene	10.37	166	60370	2.176	ng	# 83
78) Pyrene	12.77	202	156560	2.986	ng	# 93

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

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP100719\
Data File : BP000557.D
Acq On : 07 Oct 2019 20:57
Operator : JU
Sample : K5213-09
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
W-1

Quant Time: Oct 08 10:55:42 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP100119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Oct 01 18:03:42 2019
Response via : Initial Calibration

