

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP111119\
 Data File : BP001024.D
 Acq On : 11 Nov 2019 19:11
 Operator : JU
 Sample : K5760-02
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 RBR-50

Quant Time: Nov 12 06:11:54 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP110619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 06 17:07:13 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.38	152	241400	20.00	ng	0.00
21) Naphthalene-d8	10.40	136	888210	20.00	ng	0.00
39) Acenaphthene-d10	13.26	164	500262	20.00	ng	0.00
64) Phenanthrene-d10	15.66	188	954881	20.00	ng	0.00
76) Chrysene-d12	19.29	240	805002	20.00	ng	0.00
87) Perylene-d12	21.08	264	909924	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	6.30	112	1213116	91.47	ng	0.00
7) Phenol-d6	7.82	99	1607469	95.85	ng	0.00
23) Nitrobenzene-d5	9.25	82	991215	61.73	ng	-0.01
42) 2,4,6-Tribromophenol	14.55	330	601562	100.73	ng	0.00
45) 2-Fluorobiphenyl	12.18	172	2054201	61.84	ng	0.00
79) Terphenyl-d14	17.93	244	2394519	59.17	ng	0.00
Target Compounds						
10) Phenol	7.83	94	95919	5.229	ng	97
50) Dimethylphthalate	12.85	163	157985	4.271	ng	99
75) Fluoranthene	17.40	202	174548	3.160	ng	98
78) Pyrene	17.70	202	152919	2.688	ng	97
83) Chrysene	19.32	228	103916	2.183	ng	96
88) Benzo(b)fluoranthene	20.58	252	136063	2.509	ng	99

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

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP111119\
Data File : BP001024.D
Acq On : 11 Nov 2019 19:11
Operator : JU
Sample : K5760-02
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
RBR-50

Quant Time: Nov 12 06:11:54 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP110619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Nov 06 17:07:13 2019
Response via : Initial Calibration

