

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071119\
 Data File : BF115505.D
 Acq On : 11 Jul 2019 20:15
 Operator : HP/JU
 Sample : K3746-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LID-A

Quant Time: Jul 12 05:23:31 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF070519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 10 16:47:05 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	147889	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	572351	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	295556	20.00	ng	-0.01
64) Phenanthrene-d10	11.42	188	589550	20.00	ng	0.00
76) Chrysene-d12	14.05	240	486990	20.00	ng	-0.02
87) Perylene-d12	15.53	264	467396	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	853790	89.20	ng	0.00
7) Phenol-d6	6.52	99	1129820	92.81	ng	0.00
23) Nitrobenzene-d5	7.45	82	690960	71.33	ng	0.00
42) 2,4,6-Tribromophenol	10.72	330	339335	103.84	ng	-0.01
45) 2-Fluorobiphenyl	9.25	172	1314791	78.06	ng	0.00
79) Terphenyl-d14	13.00	244	1557477	65.48	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.64	163	209359	9.438	ng	100
71) Phenanthrene	11.44	178	100231	3.233	ng	96
75) Fluoranthene	12.63	202	186532	5.714	ng	96
78) Pyrene	12.86	202	125242	3.307	ng	97
83) Chrysene	14.08	228	121897	3.802	ng	98

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

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071119\
Data File : BF115505.D
Acq On : 11 Jul 2019 20:15
Operator : HP/JU
Sample : K3746-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
LID-A

Quant Time: Jul 12 05:23:31 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF070519.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 10 16:47:05 2019
Response via : Initial Calibration

