

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072619\
 Data File : BF115804.D
 Acq On : 26 Jul 2019 23:10
 Operator : HP/JU
 Sample : K3626-12
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SU-01-072519-B

Quant Time: Jul 27 02:14:25 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 25 13:16:07 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	155746	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	587987	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	324066	20.00	ng	0.00
64) Phenanthrene-d10	11.39	188	586091	20.00	ng	0.00
76) Chrysene-d12	14.03	240	441629	20.00	ng	0.00
87) Perylene-d12	15.50	264	396287	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	970821	101.96	ng	0.01
7) Phenol-d6	6.50	99	1062936	87.98	ng	0.00
23) Nitrobenzene-d5	7.43	82	985174	97.42	ng	0.00
42) 2,4,6-Tribromophenol	10.70	330	461114	121.90	ng	0.00
45) 2-Fluorobiphenyl	9.23	172	1782726	96.29	ng	0.00
79) Terphenyl-d14	12.97	244	2062385	86.17	ng	0.00
Target Compounds						
31) Naphthalene	8.17	128	94494	3.318	ng	100
37) 2-Methylnaphthalene	8.86	142	56404	2.988	ng	96
38) 1-Methylnaphthalene	8.96	142	43301	2.415	ng	98
71) Phenanthrene	11.41	178	77315	2.574	ng	97
73) Carbazole	11.62	167	64335	2.363	ng	99

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

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072619\
Data File : BF115804.D
Acq On : 26 Jul 2019 23:10
Operator : HP/JU
Sample : K3626-12
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SU-01-072519-B

Quant Time: Jul 27 02:14:25 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jul 25 13:16:07 2019
Response via : Initial Calibration

