

Data Path : Z:\HPCHEM1\BNA F\DATA\BF022718\
 Data File : BF103288.D
 Acq On : 28 Feb 2018 3:21
 Operator : SJ/JU
 Sample : J1683-03
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RFK-RMAB-MW10-GW

Quant Time: Feb 28 04:29:16 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 27 17:05:20 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	142475	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	534393	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	222973	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	318332	20.00	ng	0.00
75) Chrysene-d12	14.05	240	243111	20.00	ng	0.00
86) Perylene-d12	15.56	264	186837	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	432884	45.95	ng	0.00
7) Phenol-d6	6.50	99	354478	30.75	ng	0.00
23) Nitrobenzene-d5	7.45	82	571907	61.12	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	164718	87.28	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	966449	71.96	ng	0.00
78) Terphenyl-d14	12.99	244	681583	67.39	ng	0.00
Target Compounds						
10) Phenol	6.52	94	53547	4.002	ng	92
22) Acetophenone	7.26	105	304935	24.446	ng	# 58
31) Naphthalene	8.18	128	421825	16.123	ng	97
49) Dimethylphthalate	9.63	163	51990	2.907	ng	# 57

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

Data Path : Z:\HPCHEM1\BNA F\DATA\BF022718\
Data File : BF103288.D
Acq On : 28 Feb 2018 3:21
Operator : SJ/JU
Sample : J1683-03
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
RFK-RMAB-MW10-GW

Quant Time: Feb 28 04:29:16 2018
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Feb 27 17:05:20 2018
Response via : Initial Calibration

