

Data Path : Z:\HPCHEM1\BNA F\DATA\BF082616\
 Data File : BF089922.D
 Acq On : 26 Aug 2016 22:00
 Operator : UM/SJ
 Sample : H4597-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-104

Quant Time: Aug 26 23:26:02 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 26 18:51:25 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.65	152	361200	20.00	ng	0.00
21) Naphthalene-d8	7.93	136	1302077	20.00	ng	-0.01
38) Acenaphthene-d10	9.68	164	657387	20.00	ng	-0.01
63) Phenanthrene-d10	11.15	188	1234759	20.00	ng	0.00
75) Chrysene-d12	13.77	240	1004369	20.00	ng	-0.01
86) Perylene-d12	15.12	264	845375	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.22	112	988996	46.96	ng	0.00
7) Phenol-d6	6.28	99	866162	33.50	ng	-0.01
23) Nitrobenzene-d5	7.22	82	1762057	84.07	ng	0.00
41) 2,4,6-Tribromophenol	10.47	330	600483	96.12	ng	-0.01
44) 2-Fluorobiphenyl	9.02	172	2548118	68.13	ng	0.00
78) Terphenyl-d14	12.74	244	2195099	49.44	ng	0.00
Target Compounds						Qvalue
37) 2-Methylnaphthalene	8.65	142	163152	4.14	ng	99
49) Dimethylphthalate	9.40	163	173863	3.76	ng	99
51) Acenaphthene	9.71	154	100460	2.51	ng	98
54) Dibenzofuran	9.88	168	111555	2.22	ng	# 63
57) Fluorene	10.23	166	194332	4.85	ng	100
70) Phenanthrene	11.18	178	238732	3.83	ng	98

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

Data Path : Z:\HPCHEM1\BNA F\DATA\BF082616\
Data File : BF089922.D
Acq On : 26 Aug 2016 22:00
Operator : UM/SJ
Sample : H4597-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-104

Quant Time: Aug 26 23:26:02 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 26 18:51:25 2016
Response via : Initial Calibration

