

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042717\
 Data File : BF094670.D
 Acq On : 27 Apr 2017 18:19
 Operator : SJ/MA
 Sample : I2855-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OWS-2

Manual Integrations
 APPROVED

mohammad
 4/28/2017 3:00:40 PM

Quant Time: Apr 28 04:43:48 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF041717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 17 14:59:23 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	5.78	152	119361	20.00	ng	0.01
21) Naphthalene-d8	7.27	136	478168	20.00	ng	0.00
38) Acenaphthene-d10	9.40	164	154157	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	378937	20.00	ng	0.00
75) Chrysene-d12	14.47	240	240685	20.00	ng	0.00
86) Perylene-d12	16.09	264	27585	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	4.34	112	895280	124.44	ng	0.04
7) Phenol-d6	5.40	99	837359	98.73	ng	0.00
23) Nitrobenzene-d5	6.42	82	732834	94.63	ng	0.00
41) 2,4,6-Tribromophenol	10.38	330	263855	218.82	ng	0.00
44) 2-Fluorobiphenyl	8.59	172	1368814	130.65	ng	0.00
78) Terphenyl-d14	13.20	244	1053499	117.00	ng	0.00
Target Compounds						
14) 1,2-Dichlorobenzene	5.97	146	18483	2.20	ng	Qvalue 95
20) 3+4-Methylphenols	6.28	107	24706	2.82	ng	# 54
32) Benzoic acid	7.08	122	96093m	25.54	ng	

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

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042717\
Data File : BF094670.D
Acq On : 27 Apr 2017 18:19
Operator : SJ/MA
Sample : I2855-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
OWS-2

Manual Integrations
APPROVED

mohammad
4/28/2017 3:00:40 PM

Quant Time: Apr 28 04:43:48 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF041717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Apr 17 14:59:23 2017
Response via : Initial Calibration

