

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF111017\
 Data File : BF100385.D
 Acq On : 10 Nov 2017 12:39
 Operator : SJ/JU
 Sample : I6384-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EAST-RUTHERFORD-ROLLOFFS

Manual Integrations
 APPROVED

SOHIL
 11/13/2017 4:43:11 PM

Quant Time: Nov 10 16:00:33 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 10 15:44:40 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	186141	20.00	ng	0.00
21) Naphthalene-d8	8.18	136	751331	20.00	ng	0.00
38) Acenaphthene-d10	9.94	164	338008	20.00	ng	0.00
63) Phenanthrene-d10	11.43	188	547672	20.00	ng	0.00
75) Chrysene-d12	14.06	240	429930	20.00	ng	0.00
86) Perylene-d12	15.54	264	412438	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	1487125	128.07	ng	0.01
7) Phenol-d6	6.53	99	1849502	129.16	ng	0.00
23) Nitrobenzene-d5	7.47	82	1208136	106.31	ng	0.00
41) 2,4,6-Tribromophenol	10.73	330	434585	126.18	ng	0.00
44) 2-Fluorobiphenyl	9.26	172	1835910	82.71	ng	0.00
78) Terphenyl-d14	13.02	244	1448924	77.44	ng	0.00
Target Compounds						
10) Phenol	6.54	94	75645	5.032	ng	94
49) Dimethylphthalate	9.65	163	131208	5.084	ng	# 98
70) Phenanthrene	11.45	178	169265	5.870	ng	# 96
74) Fluoranthene	12.64	202	163265	5.342	ng	97
77) Pyrene	12.87	202	135948	4.436	ng	98
82) Chrysene	14.09	228	50296	2.006	ng	95
87) Benzo(b)fluoranthene	15.10	252	61287m	2.508	ng	

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

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