

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF121917\
 Data File : BF101535.D
 Acq On : 19 Dec 2017 21:03
 Operator : SJ/JU
 Sample : I6981-01 10X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PHMHA-WCS

Manual Integrations
 APPROVED

Sohil
 12/20/2017 3:48:02 PM

Quant Time: Dec 20 07:52:47 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 18 15:38:59 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	155269	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	622683	20.00	ng	0.00
38) Acenaphthene-d10	9.85	164	246031	20.00	ng	0.00
63) Phenanthrene-d10	11.34	188	390039	20.00	ng	0.00
75) Chrysene-d12	13.99	240	314702	20.00	ng	0.00
86) Perylene-d12	15.46	264	319504	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.44	112	96586	9.27	ng	0.02
7) Phenol-d6	6.46	99	123159	9.49	ng	0.00
23) Nitrobenzene-d5	7.39	82	75239	6.73	ng	0.00
41) 2,4,6-Tribromophenol	10.65	330	9417	3.97	ng	0.00
44) 2-Fluorobiphenyl	9.16	172	147112	9.28	ng	0.00
78) Terphenyl-d14	12.92	244	98906	7.58	ng	0.00
Target Compounds						
70) Phenanthrene	11.36	178	63076	2.908	ng	Qvalue 97
74) Fluoranthene	12.56	202	131879	5.792	ng	96
77) Pyrene	12.79	202	112069	4.745	ng	99
80) Benzo(a)anthracene	13.97	228	52982	2.550	ng	96
82) Chrysene	14.02	228	49451	2.520	ng	94
87) Benzo(b)fluoranthene	15.03	252	59276m	3.044	ng	
89) Benzo(a)pyrene	15.40	252	39466	2.198	ng	# 90

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

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