

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF110817\
 Data File : BF100296.D
 Acq On : 8 Nov 2017 17:22
 Operator : SJ/JU
 Sample : I6305-01 10X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PL-01-110317

Manual Integrations
 APPROVED

Sohil
 11/9/2017 3:41:36 PM

Quant Time: Nov 08 17:51:25 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 08 14:33:31 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	159283	20.00	ng	0.00
21) Naphthalene-d8	8.18	136	653676	20.00	ng	0.00
38) Acenaphthene-d10	9.93	164	299225	20.00	ng	0.00
63) Phenanthrene-d10	11.42	188	534181	20.00	ng	0.00
75) Chrysene-d12	14.06	240	347325	20.00	ng	-0.01
86) Perylene-d12	15.54	264	344208	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	118297	11.91	ng	-0.01
7) Phenol-d6	6.50	99	148091	12.09	ng	-0.02
23) Nitrobenzene-d5	7.45	82	76995	7.79	ng	-0.02
41) 2,4,6-Tribromophenol	10.72	330	33493	10.98	ng	-0.01
44) 2-Fluorobiphenyl	9.25	172	157091	7.99	ng	-0.01
78) Terphenyl-d14	13.00	244	112129	7.42	ng	-0.01
Target Compounds						
49) Dimethylphthalate	9.65	163	49996	2.188	ng	99
70) Phenanthrene	11.44	178	202982	7.217	ng	97
71) Anthracene	11.49	178	60058	2.105	ng	98
74) Fluoranthene	12.63	202	282396	9.474	ng	95
77) Pyrene	12.86	202	317645	12.830	ng	99
80) Benzo(a)anthracene	14.04	228	123637	5.911	ng	95
82) Chrysene	14.08	228	107329	5.299	ng	97
85) Indeno(1,2,3-cd)pyrene	17.02	276	58577	3.576	ng	# 93
87) Benzo(b)fluoranthene	15.10	252	140527m	6.891	ng	
89) Benzo(a)pyrene	15.47	252	103744	5.738	ng	100
91) Benzo(g,h,i)perylene	17.47	276	61010	4.215	ng	95

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

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