

Data Path : Z:\HPCHEM1\BNA F\DATA\BF081217\
 Data File : BF097643.D
 Acq On : 12 Aug 2017 19:48
 Operator : SJ/JU
 Sample : I4705-04
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-B-COMP

Manual Integrations
 APPROVED

sohil
 8/14/2017 7:06:16 PM

Quant Time: Aug 14 06:48:01 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 11 15:55:57 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.50	152	114335	20.00	ng	-0.01
21) Naphthalene-d8	9.53	136	481421	20.00	ng	-0.01
38) Acenaphthene-d10	12.35	164	173135	20.00	ng	-0.02
63) Phenanthrene-d10	14.74	188	254589	20.00	ng	-0.02
75) Chrysene-d12	18.44	240	181675	20.00	ng	-0.02
86) Perylene-d12	20.12	264	174680	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	636728	88.52	ng	0.01
7) Phenol-d6	7.06	99	828096	92.80	ng	0.00
23) Nitrobenzene-d5	8.41	82	463596	60.38	ng	-0.01
41) 2,4,6-Tribromophenol	13.64	330	127227	96.67	ng	-0.02
44) 2-Fluorobiphenyl	11.30	172	779662	67.96	ng	-0.02
78) Terphenyl-d14	17.10	244	443094	50.48	ng	-0.02
Target Compounds						
49) Dimethylphthalate	11.99	163	204375	15.317	ng	98
74) Fluoranthene	16.56	202	29651	2.238	ng	95
82) Chrysene	18.46	228	23612	2.237	ng	# 89
85) Indeno(1,2,3-cd)pyrene	21.34	276	24496	2.984	ng	93
87) Benzo(b)fluoranthene	19.71	252	83908m	7.999	ng	
89) Benzo(a)pyrene	20.06	252	46522	4.832	ng	# 96
91) Benzo(g,h,i)perylene	21.69	276	20290	2.683	ng	94

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

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