

Data Path : Z:\HPCHEM1\BNA F\DATA\BF121616\
 Data File : BF091696.D
 Acq On : 17 Dec 2016 2:03
 Operator : UM/SJ
 Sample : H6099-14 10X
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP12-2.5-3FT

Quant Time: Dec 17 04:42:58 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 16 00:35:41 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.77	152	216615	20.00	ng	-0.01
21) Naphthalene-d8	8.06	136	884112	20.00	ng	-0.01
38) Acenaphthene-d10	9.81	164	383538	20.00	ng	-0.01
63) Phenanthrene-d10	11.30	188	527577	20.00	ng	-0.01
75) Chrysene-d12	13.93	240	451575	20.00	ng	-0.01
86) Perylene-d12	15.33	264	339866	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.37	112	168649	13.12	ng	-0.01
7) Phenol-d6	6.42	99	214168	13.36	ng	-0.01
23) Nitrobenzene-d5	7.33	82	130759	8.25	ng	-0.02
41) 2,4,6-Tribromophenol	10.60	330	35310	11.08	ng	-0.01
44) 2-Fluorobiphenyl	9.14	172	237411	10.70	ng	-0.01
78) Terphenyl-d14	12.88	244	154658	7.77	ng	-0.02
Target Compounds						Qvalue
70) Phenanthrene	11.32	178	151393	5.79	ng	96
74) Fluoranthene	12.50	202	240981	8.81	ng	95
77) Pyrene	12.73	202	212480	5.93	ng	98
80) Benzo(a)anthracene	13.92	228	101010	3.85	ng	# 93
82) Chrysene	13.95	228	109354m	3.99	ng	
87) Benzo(b)fluoranthene	14.93	252	110771m	5.47	ng	
88) Benzo(k)fluoranthene	14.95	252	47313m	2.66	ng	
89) Benzo(a)pyrene	15.28	252	77031	4.22	ng	# 92
91) Benzo(g,h,i)perylene	17.08	276	49043	2.95	ng	# 85

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

