

Data Path : Z:\HPCHEM1\BNA F\DATA\BF112917\
 Data File : BF100999.D
 Acq On : 30 Nov 2017 2:08
 Operator : SJ/JU
 Sample : I6578-05DL 5X
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BBR-2-E(4-5)DL

Manual Integrations
 APPROVED

Sohil
 11/30/2017 4:32:27 PM

Quant Time: Nov 30 07:36:32 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 28 15:21:02 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	51989	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	188736	20.00	ng	0.00
38) Acenaphthene-d10	9.89	164	73126	20.00	ng	0.00
63) Phenanthrene-d10	11.38	188	129435	20.00	ng	0.00
75) Chrysene-d12	14.03	240	117962	20.00	ng	0.00
86) Perylene-d12	15.50	264	91944	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	68033	20.98	ng	-0.01
7) Phenol-d6	6.47	99	81140	20.29	ng	-0.02
23) Nitrobenzene-d5	7.41	82	44098	15.45	ng	-0.02
41) 2,4,6-Tribromophenol	10.68	330	15910	21.35	ng	0.00
44) 2-Fluorobiphenyl	9.21	172	99383	20.69	ng	0.00
78) Terphenyl-d14	12.96	244	89031	17.34	ng	0.00
Target Compounds						
70) Phenanthrene	11.40	178	89369	13.114	ng	98
71) Anthracene	11.46	178	18878	2.730	ng	97
74) Fluoranthene	12.60	202	216804	30.018	ng	93
77) Pyrene	12.83	202	219653	26.122	ng	98
80) Benzo(a)anthracene	14.02	228	109126	15.362	ng	98
82) Chrysene	14.05	228	105114	15.281	ng	97
85) Indeno(1,2,3-cd)pyrene	16.97	276	40600	7.297	ng	# 89
87) Benzo(b)fluoranthene	15.07	252	107260m	19.691	ng	
88) Benzo(k)fluoranthene	15.09	252	27349m	5.314	ng	
89) Benzo(a)pyrene	15.43	252	73500	15.220	ng	95
90) Dibenzo(a,h)anthracene	16.99	278	9274m	2.348	ng	
91) Benzo(a,h,i)perylene	17.43	276	37899	9.803	ng	# 91

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

