

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102918\
 Data File : BF110282.D
 Acq On : 30 Oct 2018 00:27
 Operator : JU/SJ
 Sample : J5194-09 5X
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HR-05-102618-B

Manual Integrations
 APPROVED

Sohil
 10/30/2018 4:42:47 PM

Quant Time: Oct 30 12:56:51 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 29 17:29:00 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	118074	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	452094	20.00	ng	0.00
39) Acenaphthene-d10	10.02	164	180127	20.00	ng	0.00
64) Phenanthrene-d10	11.50	188	283795	20.00	ng	0.00
76) Chrysene-d12	14.16	240	196161	20.00	ng	0.00
87) Perylene-d12	15.69	264	137929	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.58	112	114864	15.84	ng	0.00
7) Phenol-d6	6.58	99	145896	15.73	ng	-0.01
23) Nitrobenzene-d5	7.53	82	89929	11.80	ng	0.00
42) 2,4,6-Tribromophenol	10.80	330	20882	11.70	ng	0.00
45) 2-Fluorobiphenyl	9.32	172	187957	16.39	ng	0.00
79) Terphenyl-d14	13.09	244	139124	14.75	ng	0.00
Target Compounds						
71) Phenanthrene	11.53	178	39045	2.629	ng	97
75) Fluoranthene	12.73	202	70537	4.680	ng	99
78) Pyrene	12.96	202	100222	7.127	ng	97
81) Benzo(a)anthracene	14.14	228	45082	3.875	ng	# 91
83) Chrysene	14.18	228	44580	4.035	ng	# 95
88) Benzo(b)fluoranthene	15.23	252	32804m	4.119	ng	
90) Benzo(a)pyrene	15.63	252	27288	3.723	ng	# 84
92) Benzo(g,h,i)perylene	17.74	276	18082	2.902	ng	# 81

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

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