

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF041019\
 Data File : BF113711.D
 Acq On : 10 Apr 2019 22:20
 Operator : JU/SJ
 Sample : K2342-03 5X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-11-S

Manual Integrations
 APPROVED

Sohil
 4/11/2019 2:40:00 PM

Quant Time: Apr 11 07:54:19 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 04 04:05:31 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.67	152	136669	20.00	ng	-0.01
21) Naphthalene-d8	7.96	136	503069	20.00	ng	0.00
39) Acenaphthene-d10	9.71	164	213189	20.00	ng	-0.01
64) Phenanthrene-d10	11.19	188	385661	20.00	ng	-0.01
76) Chrysene-d12	13.83	240	345978	20.00	ng	-0.02
87) Perylene-d12	15.23	264	251945	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	132330	16.51	ng	0.00
7) Phenol-d6	6.34	99	176628	17.88	ng	0.00
23) Nitrobenzene-d5	7.25	82	89454	10.42	ng	-0.01
42) 2,4,6-Tribromophenol	10.50	330	36178	14.13	ng	-0.01
45) 2-Fluorobiphenyl	9.03	172	170422	12.87	ng	-0.02
79) Terphenyl-d14	12.77	244	153674	9.04	ng	-0.02
Target Compounds						
71) Phenanthrene	11.22	178	105362	5.498	ng	97
75) Fluoranthene	12.40	202	203844	9.926	ng	98
78) Pyrene	12.63	202	182603	6.936	ng	98
81) Benzo(a)anthracene	13.82	228	116202	5.822	ng	99
83) Chrysene	13.85	228	108716	5.374	ng	98
86) Indeno(1,2,3-cd)pyrene	16.60	276	57911	3.100	ng	95
88) Benzo(b)fluoranthene	14.84	252	129196m	8.377	ng	
89) Benzo(k)fluoranthene	14.86	252	41296m	2.984	ng	
90) Benzo(a)pyrene	15.18	252	87426	6.380	ng	98
92) Benzo(g,h,i)perylene	17.02	276	55236	4.223	ng	95

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

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