

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF032119\
 Data File : BF113220.D
 Acq On : 21 Mar 2019 20:44
 Operator : JU/SJ
 Sample : K2004-09 20X
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-22(7-8)

Quant Time: Mar 22 02:01:30 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 13 03:42:55 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.72	152	173879	20.00	ng	-0.01
21) Naphthalene-d8	8.00	136	674195	20.00	ng	-0.01
39) Acenaphthene-d10	9.76	164	311124	20.00	ng	-0.01
64) Phenanthrene-d10	11.24	188	532786	20.00	ng	0.00
76) Chrysene-d12	13.87	240	384964	20.00	ng	0.00
87) Perylene-d12	15.30	264	325504	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	23072	2.06	ng	-0.01
7) Phenol-d6	6.37	99	42448	3.02	ng	0.00
23) Nitrobenzene-d5	7.28	82	24263	2.15	ng	-0.02
42) 2,4,6-Tribromophenol	10.55	330	7758	2.35	ng	0.00
45) 2-Fluorobiphenyl	9.08	172	52549	0.76	ng	-0.01
79) Terphenyl-d14	12.82	244	40710	2.05	ng	0.00
Target Compounds						
78) Pyrene	12.67	202	165984	5.115	ng	Qvalue 95
81) Benzo(a)anthracene	13.86	228	179756	6.737	ng	# 79
83) Chrysene	13.90	228	179290	6.860	ng	# 80
90) Benzo(a)pyrene	15.23	252	120994	6.497	ng	# 75
92) Benzo(g,h,i)perylene	17.07	276	79681	4.993	ng	# 87

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

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