

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101019\
 Data File : BF117369.D
 Acq On : 10 Oct 2019 19:22
 Operator : JU
 Sample : K5145-01 2X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HD-01-100919

Quant Time: Oct 11 06:54:14 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 09 01:26:00 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	42711	20.00	ng	-0.01
21) Naphthalene-d8	8.07	136	175441	20.00	ng	-0.01
39) Acenaphthene-d10	9.83	164	85186	20.00	ng	-0.01
64) Phenanthrene-d10	11.32	188	120155	20.00	ng	-0.01
76) Chrysene-d12	13.96	240	84398	20.00	ng	-0.01
87) Perylene-d12	15.42	264	96222	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	95384	34.87	ng	0.00
7) Phenol-d6	6.44	99	140240	39.67	ng	-0.01
23) Nitrobenzene-d5	7.36	82	80125	24.00	ng	-0.01
42) 2,4,6-Tribromophenol	10.62	330	22963	32.25	ng	-0.01
45) 2-Fluorobiphenyl	9.15	172	150691	27.66	ng	-0.01
79) Terphenyl-d14	12.90	244	101471	22.47	ng	-0.01
Target Compounds						
50) Dimethylphthalate	9.55	163	15682	2.600	ng	99
75) Fluoranthene	12.53	202	24219	3.454	ng	98
78) Pyrene	12.76	202	38605	5.451	ng	97
81) Benzo(a)anthracene	13.95	228	17681m	3.136	ng	
83) Chrysene	13.99	228	19656m	3.574	ng	
86) Indeno(1,2,3-cd)pyrene	16.87	276	18856	4.700	ng	# 91
88) Benzo(b)fluoranthene	15.00	252	36244m	6.218	ng	
89) Benzo(k)fluoranthene	15.01	252	17605m	3.189	ng	
90) Benzo(a)pyrene	15.36	252	33670	6.583	ng	# 90
92) Benzo(g,h,i)perylene	17.32	276	18808	4.632	ng	96

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

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