

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100219\
 Data File : BF117294.D
 Acq On : 2 Oct 2019 12:50
 Operator : JU
 Sample : K4659-04RE 2X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 JC-02-093019-B

Quant Time: Oct 03 01:49:14 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 13 16:23:17 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.81	152	34020	20.00	ng	-0.01
21) Naphthalene-d8	8.09	136	134515	20.00	ng	-0.02
39) Acenaphthene-d10	9.85	164	64101	20.00	ng	-0.01
64) Phenanthrene-d10	11.33	188	96070	20.00	ng	0.00
76) Chrysene-d12	13.97	240	92661	20.00	ng	0.00
87) Perylene-d12	15.43	264	80008	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	80064	36.74	ng	-0.01
7) Phenol-d6	6.45	99	112208	39.85	ng	0.00
23) Nitrobenzene-d5	7.37	82	65637	25.64	ng	-0.02
42) 2,4,6-Tribromophenol	10.63	330	18298	34.15	ng	-0.01
45) 2-Fluorobiphenyl	9.16	172	121720	29.69	ng	-0.01
79) Terphenyl-d14	12.92	244	119283	24.06	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.56	163	25993	5.727	ng	97
71) Phenanthrene	11.36	178	13045	2.391	ng	98
75) Fluoranthene	12.55	202	22474	4.008	ng	92
78) Pyrene	12.78	202	24148	3.106	ng	97
83) Chrysene	14.00	228	12352	2.046	ng	# 89
88) Benzo(b)fluoranthene	15.02	252	12578m	2.595	ng	

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

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