

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100919\
 Data File : BF117356.D
 Acq On : 9 Oct 2019 15:45
 Operator : JU
 Sample : K5281-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-PS-01-062

Quant Time: Oct 09 22:54:50 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	6.80	152	42993	20.00	ng	0.00
21) Naphthalene-d8	8.08	136	175369	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	89274	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	128502	20.00	ng	0.00
76) Chrysene-d12	13.97	240	85389	20.00	ng	0.00
87) Perylene-d12	15.43	264	101842	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	252429	91.67	ng	0.00
7) Phenol-d6	6.45	99	351596	98.79	ng	0.00
23) Nitrobenzene-d5	7.36	82	225263	67.49	ng	0.00
42) 2,4,6-Tribromophenol	10.63	330	66345	88.92	ng	0.00
45) 2-Fluorobiphenyl	9.16	172	393586	68.93	ng	0.00
79) Terphenyl-d14	12.90	244	278657	61.00	ng	0.00
Target Compounds						
10) Phenol	6.46	94	11751	2.995	ng	Qvalue 84
50) Dimethylphthalate	9.55	163	83790	13.256	ng	98
71) Phenanthrene	11.35	178	19028	2.607	ng	97
75) Fluoranthene	12.54	202	43045	5.740	ng	95
78) Pyrene	12.77	202	34350	4.794	ng	99
81) Benzo(a)anthracene	13.96	228	16850	2.954	ng	98
83) Chrysene	13.99	228	17147	3.082	ng	96
86) Indeno(1,2,3-cd)pyrene	16.89	276	8930	2.200	ng	97
88) Benzo(b)fluoranthene	15.00	252	22615m	3.666	ng	
90) Benzo(a)pyrene	15.37	252	15335	2.833	ng	# 82

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

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