

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120519\
 Data File : BF118110.D
 Acq On : 5 Dec 2019 18:33
 Operator : JU
 Sample : K6098-04
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OWL-C5-GW

Quant Time: Dec 06 04:06:38 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 03 15:16:53 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	116416	20.00	ng	-0.01
21) Naphthalene-d8	8.16	136	447012	20.00	ng	-0.01
39) Acenaphthene-d10	9.92	164	236048	20.00	ng	-0.02
64) Phenanthrene-d10	11.41	188	432891	20.00	ng	-0.01
76) Chrysene-d12	14.06	240	265221	20.00	ng	-0.01
87) Perylene-d12	15.56	264	248977	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	659809	100.32	ng	0.00
7) Phenol-d6	6.50	99	561923	70.84	ng	-0.01
23) Nitrobenzene-d5	7.43	82	677548	90.10	ng	-0.01
42) 2,4,6-Tribromophenol	10.72	330	355819	142.38	ng	0.00
45) 2-Fluorobiphenyl	9.24	172	1326227	100.80	ng	0.00
79) Terphenyl-d14	13.00	244	1617481	111.46	ng	-0.01
Target Compounds						
10) Phenol	6.51	94	61831	7.501	ng	97
50) Dimethylphthalate	9.63	163	295134	17.830	ng	99
52) Acenaphthene	9.95	154	28386	2.110	ng	# 94
58) Fluorene	10.47	166	57346	3.972	ng	# 93

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

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