

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120519\
 Data File : BF118109.D
 Acq On : 5 Dec 2019 18:05
 Operator : JU
 Sample : K6098-03
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OWL-C4-GW

Quant Time: Dec 06 06:12:14 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	120173	20.00	ng	-0.01
21) Naphthalene-d8	8.16	136	450938	20.00	ng	-0.01
39) Acenaphthene-d10	9.92	164	222391	20.00	ng	-0.02
64) Phenanthrene-d10	11.41	188	368767	20.00	ng	-0.01
76) Chrysene-d12	14.06	240	191717	20.00	ng	-0.01
87) Perylene-d12	15.56	264	251378	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	740347	109.05	ng	0.00
7) Phenol-d6	6.50	99	652524	79.69	ng	-0.01
23) Nitrobenzene-d5	7.43	82	724750	95.54	ng	-0.01
42) 2,4,6-Tribromophenol	10.72	330	317497	134.85	ng	0.00
45) 2-Fluorobiphenyl	9.24	172	1387501	111.93	ng	0.00
79) Terphenyl-d14	13.00	244	1220594	116.36	ng	-0.01
Target Compounds						
10) Phenol	6.51	94	110061	12.934	ng	98
32) Benzoic acid	7.87	122	15056	3.038	ng	92
50) Dimethylphthalate	9.63	163	290650	18.637	ng	99
52) Acenaphthene	9.95	154	42009	3.314	ng	98
58) Fluorene	10.47	166	62902	4.624	ng	98
71) Phenanthrene	11.43	178	43286	2.335	ng	96

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

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