

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120519\
 Data File : BF118108.D
 Acq On : 5 Dec 2019 17:37
 Operator : JU
 Sample : K6098-01
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OWL-C1-GW

Quant Time: Dec 06 03:47:26 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	122484	20.00	ng	-0.01
21) Naphthalene-d8	8.16	136	473894	20.00	ng	-0.01
39) Acenaphthene-d10	9.92	164	250561	20.00	ng	-0.02
64) Phenanthrene-d10	11.41	188	480937	20.00	ng	-0.01
76) Chrysene-d12	14.06	240	328885	20.00	ng	-0.01
87) Perylene-d12	15.56	264	298338	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	401011	57.95	ng	0.00
7) Phenol-d6	6.49	99	352185	42.20	ng	-0.02
23) Nitrobenzene-d5	7.43	82	700058	87.82	ng	-0.01
42) 2,4,6-Tribromophenol	10.72	330	251395	94.77	ng	0.00
45) 2-Fluorobiphenyl	9.24	172	1434506	102.71	ng	0.00
79) Terphenyl-d14	13.01	244	1740899	96.74	ng	0.00
Target Compounds						
10) Phenol	6.50	94	23109	2.665	ng	91
50) Dimethylphthalate	9.63	163	314805	17.917	ng	98
52) Acenaphthene	9.95	154	37550	2.629	ng	97
58) Fluorene	10.47	166	62103	4.052	ng	98
71) Phenanthrene	11.43	178	55818	2.308	ng	98

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

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