

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120518\
 Data File : BF111118.D
 Acq On : 5 Dec 2018 20:21
 Operator : JU/SJ
 Sample : J5764-27 2X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 JC-02-113018-D

Quant Time: Dec 06 01:27:23 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 05 17:30:55 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	550670	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	2143298	20.00	ng	0.00
39) Acenaphthene-d10	9.84	164	959344	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	1420660	20.00	ng	0.00
76) Chrysene-d12	13.96	240	1027199	20.00	ng	0.00
87) Perylene-d12	15.39	264	723932	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.41	112	1818002	51.52	ng	0.00
7) Phenol-d6	6.43	99	2334926	51.36	ng	0.00
23) Nitrobenzene-d5	7.36	82	1393093	38.48	ng	0.00
42) 2,4,6-Tribromophenol	10.62	330	358381	47.00	ng	0.00
45) 2-Fluorobiphenyl	9.16	172	2401663	38.47	ng	0.00
79) Terphenyl-d14	12.90	244	1663680	33.17	ng	0.00
Target Compounds						
10) Phenol	6.44	94	148755	3.001	ng	97
50) Dimethylphthalate	9.55	163	163871	2.454	ng	99
71) Phenanthrene	11.35	178	143990	2.107	ng	96
75) Fluoranthene	12.53	202	177199	2.650	ng	98
78) Pyrene	12.76	202	167324	2.091	ng	93
88) Benzo(b)fluoranthene	14.97	252	102843	2.507	ng	# 84
89) Benzo(k)fluoranthene	14.97	252	102843	2.593	ng	# 84

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

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