

Data Path : U:\HPCHEM1\BNA F\DATA\BF011718\
 Data File : BF102187.D
 Acq On : 18 Jan 2018 9:02
 Operator : SJ/JU
 Sample : J1144-03DL 2X
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 IGW-3DL

Quant Time: Jan 18 10:00:18 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF010918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 17 15:28:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.72	152	147101	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	530737	20.00	ng	0.00
38) Acenaphthene-d10	9.76	164	219961	20.00	ng	0.00
63) Phenanthrene-d10	11.24	188	310753	20.00	ng	0.00
75) Chrysene-d12	13.86	240	261679	20.00	ng	-0.01
86) Perylene-d12	15.28	264	253751	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	556732	53.44	ng	0.00
7) Phenol-d6	6.35	99	660659	53.15	ng	0.00
23) Nitrobenzene-d5	7.29	82	390097	43.19	ng	0.00
41) 2,4,6-Tribromophenol	10.54	330	99004	51.58	ng	0.00
44) 2-Fluorobiphenyl	9.08	172	680414	48.33	ng	0.00
78) Terphenyl-d14	12.82	244	438839	34.82	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.36	94	28591	2.036	ng	98
31) Naphthalene	8.03	128	477102	17.990	ng	# 95
37) 2-Methylnaphthalene	8.72	142	766217	43.887	ng	98
49) Dimethylphthalate	9.47	163	72869	4.130	ng	# 90
51) Acenaphthene	9.79	154	29811	2.053	ng	# 67
57) Fluorene	10.30	166	83655	6.072	ng	# 90
70) Phenanthrene	11.26	178	159495	8.787	ng	97

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

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