

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071018\
 Data File : BF107292.D
 Acq On : 11 Jul 2018 00:56
 Operator : JU/SJ
 Sample : J3907-01
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S-1

Quant Time: Jul 11 02:26:15 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 29 14:38:29 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.94	152	32776	20.00	ng	-0.02
21) Naphthalene-d8	8.22	136	116633	20.00	ng	-0.02
38) Acenaphthene-d10	9.98	164	54857	20.00	ng	-0.02
63) Phenanthrene-d10	11.48	188	115509	20.00	ng	-0.02
75) Chrysene-d12	14.14	240	95102	20.00	ng	-0.03
86) Perylene-d12	15.67	264	71826	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.58	112	249981	129.15	ng	-0.02
7) Phenol-d6	6.59	99	294569	121.86	ng	-0.02
23) Nitrobenzene-d5	7.51	82	170712	81.34	ng	-0.03
41) 2,4,6-Tribromophenol	10.78	330	86086	135.40	ng	-0.02
44) 2-Fluorobiphenyl	9.30	172	315420	99.91	ng	-0.03
78) Terphenyl-d14	13.07	244	406818	103.62	ng	-0.02
Target Compounds						
10) Phenol	6.60	94	7140	2.588	ng	91
49) Dimethylphthalate	9.69	163	88372	21.479	ng	100
83) Bis(2-ethylhexyl)phthalate	14.07	149	7184	2.179	ng	97

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

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