

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101018\
 Data File : BF109762.D
 Acq On : 11 Oct 2018 00:30
 Operator : JU/SJ
 Sample : J5373-01
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RB37708-7211990-45911RT

Quant Time: Oct 11 07:14:24 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF100818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 08 16:02:11 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.69	152	73688	20.00	ng	0.00
21) Naphthalene-d8	7.97	136	292504	20.00	ng	0.00
38) Acenaphthene-d10	9.72	164	129874	20.00	ng	0.00
63) Phenanthrene-d10	11.20	188	215494	20.00	ng	0.00
75) Chrysene-d12	13.83	240	181973	20.00	ng	0.00
86) Perylene-d12	15.23	264	158968	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	399868	96.32	ng	0.00
7) Phenol-d6	6.35	99	558932	100.79	ng	0.00
23) Nitrobenzene-d5	7.26	82	364764	68.74	ng	-0.01
41) 2,4,6-Tribromophenol	10.51	330	112953	79.82	ng	0.00
44) 2-Fluorobiphenyl	9.05	172	658117	69.79	ng	0.00
78) Terphenyl-d14	12.78	244	487529	51.86	ng	0.00
Target Compounds						
10) Phenol	6.36	94	14017	2.204	ng	87
49) Dimethylphthalate	9.45	163	79109	8.305	ng	99
70) Phenanthrene	11.23	178	22836	2.118	ng	98
74) Fluoranthene	12.41	202	43414	3.854	ng	98
77) Pyrene	12.63	202	40036	3.003	ng	97
80) Benzo(a)anthracene	13.82	228	21707	2.162	ng	96

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

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