

Data Path : Z:\HPCHEM1\BNA F\DATA\BF112017\
 Data File : BF100739.D
 Acq On : 20 Nov 2017 19:26
 Operator : SJ/JU
 Sample : I6521-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MH-(0-12)

Manual Integrations
 APPROVED

SOHIL
 11/21/2017 2:15:27 PM

Quant Time: Nov 21 04:37:48 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 17 15:22:08 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	99956	20.00	ng	-0.01
21) Naphthalene-d8	8.15	136	395646	20.00	ng	-0.02
38) Acenaphthene-d10	9.91	164	188130	20.00	ng	-0.01
63) Phenanthrene-d10	11.39	188	376154	20.00	ng	-0.01
75) Chrysene-d12	14.03	240	291399	20.00	ng	-0.02
86) Perylene-d12	15.50	264	206841	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	734918	117.86	ng	0.00
7) Phenol-d6	6.50	99	889782	115.72	ng	0.00
23) Nitrobenzene-d5	7.43	82	549205	91.77	ng	-0.02
41) 2,4,6-Tribromophenol	10.70	330	253829	132.41	ng	-0.01
44) 2-Fluorobiphenyl	9.23	172	1005537	81.39	ng	-0.01
78) Terphenyl-d14	12.98	244	975203	76.90	ng	0.00
Target Compounds						
10) Phenol	6.51	94	28000	3.469	ng	97
49) Dimethylphthalate	9.62	163	93065	6.479	ng	99
74) Fluoranthene	12.60	202	107052	5.100	ng	94
77) Pyrene	12.83	202	89663	4.317	ng	98
80) Benzo(a)anthracene	14.02	228	44071	2.511	ng	98
82) Chrysene	14.06	228	37237	2.191	ng	95
87) Benzo(b)fluoranthene	15.07	252	32319m	2.637	ng	
89) Benzo(a)pyrene	15.43	252	21915	2.017	ng	98

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

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