

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF021918\
 Data File : BF103099.D
 Acq On : 19 Feb 2018 23:36
 Operator : SJ/JU
 Sample : J1600-06
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 A2 SB44 BOT-N 7

Manual Integrations
 APPROVED

Sohil
 2/20/2018 10:20:57 AM

Quant Time: Feb 20 01:31:47 2018
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 16 14:34:49 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	182262	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	753044	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	322514	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	490755	20.00	ng	0.00
75) Chrysene-d12	14.04	240	321793	20.00	ng	-0.01
86) Perylene-d12	15.54	264	310600	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	1367037	113.91	ng	0.01
7) Phenol-d6	6.52	99	1637477	113.32	ng	0.00
23) Nitrobenzene-d5	7.45	82	1064197	98.03	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	307576	118.49	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	1236172	63.60	ng	0.00
78) Terphenyl-d14	12.99	244	787739	57.96	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.62	163	158472	6.302	ng	Qvalue 100
70) Phenanthrene	11.43	178	79173	2.897	ng	98
74) Fluoranthene	12.62	202	194660	7.079	ng	98
77) Pyrene	12.84	202	192649	7.635	ng	99
80) Benzo(a)anthracene	14.03	228	144987	7.164	ng	98
82) Chrysene	14.07	228	144644	7.090	ng	98
85) Indeno(1,2,3-cd)pyrene	17.04	276	88326	5.220	ng	98
87) Benzo(b)fluoranthene	15.10	252	221237m	10.986	ng	
88) Benzo(k)fluoranthene	15.12	252	72306m	3.758	ng	
89) Benzo(a)pyrene	15.47	252	158011	8.717	ng	98
91) Benzo(g,h,i)perylene	17.50	276	84290	5.853	ng	100

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

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