

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070618\
 Data File : BF107171.D
 Acq On : 6 Jul 2018 14:32
 Operator : JU/SJ
 Sample : J3826-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OK-03-070218

Manual Integrations
 APPROVED

Sohil
 7/9/2018 1:58:09 PM

Quant Time: Jul 06 16:20:25 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	56979	20.00	ng	-0.02
21) Naphthalene-d8	8.22	136	235156	20.00	ng	-0.02
38) Acenaphthene-d10	9.98	164	121317	20.00	ng	-0.02
63) Phenanthrene-d10	11.48	188	232540	20.00	ng	-0.02
75) Chrysene-d12	14.14	240	170349	20.00	ng	-0.03
86) Perylene-d12	15.67	264	178794	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	306253	91.01	ng	-0.01
7) Phenol-d6	6.59	99	400963	95.42	ng	-0.02
23) Nitrobenzene-d5	7.51	82	247481	58.49	ng	-0.03
41) 2,4,6-Tribromophenol	10.78	330	121024	86.07	ng	-0.02
44) 2-Fluorobiphenyl	9.30	172	479765	63.14	ng	-0.02
78) Terphenyl-d14	13.07	244	498387	70.87	ng	-0.02
Target Compounds						
49) Dimethylphthalate	9.70	163	80325	8.828	ng	99
70) Phenanthrene	11.50	178	46551	4.150	ng	98
74) Fluoranthene	12.70	202	105015	8.495	ng	94
77) Pyrene	12.93	202	135255	12.041	ng	98
80) Benzo(a)anthracene	14.12	228	76840	7.401	ng	94
82) Chrysene	14.16	228	78647	8.234	ng	98
85) Indeno(1,2,3-cd)pyrene	17.24	276	39048	4.817	ng	# 88
87) Benzo(b)fluoranthene	15.21	252	98122m	8.835	ng	
88) Benzo(k)fluoranthene	15.24	252	34655m	3.277	ng	
89) Benzo(a)pyrene	15.60	252	76016	7.699	ng	95
91) Benzo(g,h,i)perylene	17.73	276	36699	4.487	ng	# 89

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

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