

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070318\
 Data File : BF107094.D
 Acq On : 3 Jul 2018 20:46
 Operator : JU/SJ
 Sample : J3840-12
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-4-S

Manual Integrations
 APPROVED

Sohil
 7/5/2018 4:29:39 PM

Quant Time: Jul 05 04:08:04 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	6.95	152	48501	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	166372	20.00	ng	-0.01
38) Acenaphthene-d10	10.00	164	73033	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	153254	20.00	ng	0.00
75) Chrysene-d12	14.15	240	128124	20.00	ng	-0.02
86) Perylene-d12	15.69	264	97148	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	284309	99.26	ng	0.00
7) Phenol-d6	6.60	99	340546	95.21	ng	-0.01
23) Nitrobenzene-d5	7.52	82	186572	62.32	ng	-0.02
41) 2,4,6-Tribromophenol	10.80	330	88497	104.55	ng	0.00
44) 2-Fluorobiphenyl	9.31	172	340986	77.77	ng	-0.01
78) Terphenyl-d14	13.08	244	436237	82.47	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.71	163	62999	11.501	ng	99
70) Phenanthrene	11.52	178	117478	15.893	ng	99
71) Anthracene	11.57	178	31065	4.117	ng	98
74) Fluoranthene	12.71	202	249654	30.643	ng	97
77) Pyrene	12.95	202	236860	28.034	ng	98
80) Benzo(a)anthracene	14.14	228	108965	13.954	ng	99
82) Chrysene	14.17	228	90164	12.551	ng	99
85) Indeno(1,2,3-cd)pyrene	17.28	276	35467	5.818	ng	# 88
87) Benzo(b)fluoranthene	15.23	252	87259m	14.459	ng	
88) Benzo(k)fluoranthene	15.25	252	28188m	4.905	ng	
89) Benzo(a)pyrene	15.62	252	63036	11.750	ng	# 94
90) Dibenzo(a,h)anthracene	17.27	278	9648m	2.122	ng	
91) Benzo(a,h,i)perylene	17.77	276	36333	8.176	ng	# 90

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

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