

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101118\
 Data File : BF109802.D
 Acq On : 11 Oct 2018 22:27
 Operator : JU/SJ
 Sample : J5394-03
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-GF-02-027-B

Manual Integrations
 APPROVED

Sohil
 10/12/2018 11:01:25 AM

Quant Time: Oct 12 05:00:43 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	6.70	152	73208	20.00	ng	0.00
21) Naphthalene-d8	7.97	136	296131	20.00	ng	0.00
38) Acenaphthene-d10	9.72	164	125657	20.00	ng	0.00
63) Phenanthrene-d10	11.20	188	202304	20.00	ng	0.00
75) Chrysene-d12	13.83	240	199941	20.00	ng	0.00
86) Perylene-d12	15.23	264	227535	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.31	112	399518	96.87	ng	-0.01
7) Phenol-d6	6.34	99	576448	104.63	ng	-0.02
23) Nitrobenzene-d5	7.26	82	374220	69.66	ng	-0.01
41) 2,4,6-Tribromophenol	10.51	330	108233	79.05	ng	0.00
44) 2-Fluorobiphenyl	9.05	172	619503	67.90	ng	-0.01
78) Terphenyl-d14	12.77	244	443428	42.93	ng	-0.01
Target Compounds						
49) Dimethylphthalate	9.45	163	65864	7.147	ng	99
70) Phenanthrene	11.23	178	24303	2.401	ng	97
71) Anthracene	11.29	178	21219	2.028	ng	97
74) Fluoranthene	12.41	202	72507	6.855	ng	99
77) Pyrene	12.63	202	68002	4.642	ng	97
80) Benzo(a)anthracene	13.82	228	51455	4.664	ng	96
82) Chrysene	13.86	228	50906	4.392	ng	96
85) Indeno(1,2,3-cd)pyrene	16.58	276	28358	3.419	ng	# 88
87) Benzo(b)fluoranthene	14.83	252	67464m	5.365	ng	
88) Benzo(k)fluoranthene	14.86	252	32033m	2.536	ng	
89) Benzo(a)pyrene	15.17	252	54659	4.790	ng	96
91) Benzo(a,h,i)perylene	16.99	276	22647	2.420	ng	# 94

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

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