

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100318\
 Data File : BF109551.D
 Acq On : 3 Oct 2018 17:25
 Operator : JU/SJ
 Sample : J5194-05 2X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HR-06-100218-C

Manual Integrations
 APPROVED

Sohil
 10/4/2018 11:44:08 AM

Quant Time: Oct 03 19:10:09 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 22 13:32:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	91085	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	300040	20.00	ng	0.00
38) Acenaphthene-d10	9.73	164	127747	20.00	ng	0.00
63) Phenanthrene-d10	11.21	188	231018	20.00	ng	0.00
75) Chrysene-d12	13.84	240	228363	20.00	ng	0.00
86) Perylene-d12	15.24	264	165690	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	248651	44.96	ng	0.00
7) Phenol-d6	6.35	99	327781	46.45	ng	-0.01
23) Nitrobenzene-d5	7.26	82	194609	38.29	ng	-0.02
41) 2,4,6-Tribromophenol	10.52	330	61352	48.72	ng	-0.01
44) 2-Fluorobiphenyl	9.06	172	317376	37.47	ng	-0.02
78) Terphenyl-d14	12.79	244	303973	32.67	ng	-0.01
Target Compounds						
49) Dimethylphthalate	9.45	163	90361	9.725	ng	# 99
70) Phenanthrene	11.23	178	54841	4.754	ng	98
74) Fluoranthene	12.42	202	148876	12.077	ng	95
77) Pyrene	12.64	202	142958	8.735	ng	99
80) Benzo(a)anthracene	13.83	228	79701	5.794	ng	98
82) Chrysene	13.87	228	71505m	5.245	ng	
85) Indeno(1,2,3-cd)pyrene	16.60	276	24572	2.224	ng	# 89
87) Benzo(b)fluoranthene	14.85	252	66812m	7.338	ng	
88) Benzo(k)fluoranthene	14.86	252	28470m	3.295	ng	
89) Benzo(a)pyrene	15.19	252	45575	5.555	ng	# 87
91) Benzo(g,h,i)perylene	17.00	276	23703	3.583	ng	# 89

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

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