

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112918\
 Data File : BF111098.D
 Acq On : 29 Nov 2018 18:47
 Operator : JU/SJ
 Sample : J5767-23 2X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CL-02-112818-D

Manual Integrations
 APPROVED

mohammad
 11/30/2018 3:49:46 PM

Quant Time: Nov 30 03:12:47 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF112618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 29 10:35:31 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.93	152	58643	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	241307	20.00	ng	0.00
39) Acenaphthene-d10	9.97	164	98752	20.00	ng	-0.01
64) Phenanthrene-d10	11.47	188	156879	20.00	ng	0.00
76) Chrysene-d12	14.13	240	138921	20.00	ng	0.00
87) Perylene-d12	15.67	264	95618	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.55	112	160126	45.28	ng	0.00
7) Phenol-d6	6.56	99	214519	46.55	ng	0.00
23) Nitrobenzene-d5	7.50	82	128777	30.54	ng	0.00
42) 2,4,6-Tribromophenol	10.77	330	38071	38.82	ng	-0.01
45) 2-Fluorobiphenyl	9.29	172	221546	32.50	ng	0.00
79) Terphenyl-d14	13.05	244	180577	25.30	ng	0.00
Target Compounds						
10) Phenol	6.58	94	17080	3.291	ng	83
50) Dimethylphthalate	9.68	163	18742	2.588	ng	99
75) Fluoranthene	12.69	202	28813	3.374	ng	95
78) Pyrene	12.93	202	34850	3.435	ng	98
83) Chrysene	14.16	228	18474	2.288	ng	# 90
88) Benzo(b)fluoranthene	15.21	252	19914m	3.545	ng	
90) Benzo(a)pyrene	15.60	252	12960	2.479	ng	# 1
92) Benzo(g,h,i)perylene	17.76	276	8569	2.060	ng	# 70

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

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