

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111418\
 Data File : BF110768.D
 Acq On : 14 Nov 2018 17:00
 Operator : JU/SJ
 Sample : J5911-07 2X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FDSB-06(10-12)

Manual Integrations
 APPROVED

Sohil
 11/15/2018 10:59:26 AM

Quant Time: Nov 14 17:39:28 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 13 12:43:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	81607	20.00	ng	0.00
21) Naphthalene-d8	8.23	136	336269	20.00	ng	0.00
39) Acenaphthene-d10	9.99	164	149741	20.00	ng	0.00
64) Phenanthrene-d10	11.48	188	235417	20.00	ng	0.00
76) Chrysene-d12	14.13	240	190012	20.00	ng	0.00
87) Perylene-d12	15.66	264	160144	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	280400	55.81	ng	0.00
7) Phenol-d6	6.57	99	362021	56.35	ng	0.00
23) Nitrobenzene-d5	7.51	82	215988	36.99	ng	0.00
42) 2,4,6-Tribromophenol	10.78	330	69113	45.81	ng	0.00
45) 2-Fluorobiphenyl	9.30	172	382697	36.50	ng	0.00
79) Terphenyl-d14	13.06	244	281599	27.75	ng	0.00
Target Compounds						
10) Phenol	6.58	94	19141	2.569	ng	Qvalue # 74
37) 2-Methylnaphthalene	8.94	142	45040	4.352	ng	97
38) 1-Methylnaphthalene	9.04	142	44320m	4.453	ng	
50) Dimethylphthalate	9.69	163	39228	3.511	ng	# 95
52) Acenaphthene	10.02	154	19609	2.141	ng	# 82
55) Dibenzofuran	10.19	168	50110	3.739	ng	98
58) Fluorene	10.53	166	71238	6.921	ng	96

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

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