

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102418\
 Data File : BF110113.D
 Acq On : 24 Oct 2018 14:25
 Operator : JU/SJ
 Sample : J5610-07 5X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-4(0-2)

Manual Integrations
 APPROVED

Sohil
 10/25/2018 12:41:44 PM

Quant Time: Oct 25 09:49:42 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 23 15:06:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	145161	20.00	ng	-0.01
21) Naphthalene-d8	8.25	136	549457	20.00	ng	-0.02
39) Acenaphthene-d10	10.01	164	216085	20.00	ng	-0.02
64) Phenanthrene-d10	11.50	188	367437	20.00	ng	-0.01
76) Chrysene-d12	14.14	240	319441	20.00	ng	-0.02
87) Perylene-d12	15.67	264	233932	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.57	112	186977	20.98	ng	-0.02
7) Phenol-d6	6.57	99	232910	20.43	ng	-0.03
23) Nitrobenzene-d5	7.52	82	124634	13.45	ng	-0.03
42) 2,4,6-Tribromophenol	10.80	330	36722	17.16	ng	-0.02
45) 2-Fluorobiphenyl	9.32	172	224993	16.35	ng	-0.02
79) Terphenyl-d14	13.08	244	189812	12.36	ng	-0.02
Target Compounds						
50) Dimethylphthalate	9.72	163	34473	2.161	ng	99
75) Fluoranthene	12.72	202	101090	5.181	ng	98
78) Pyrene	12.94	202	107816	4.708	ng	96
81) Benzo(a)anthracene	14.14	228	55519	2.931	ng	# 93
83) Chrysene	14.17	228	52041m	2.892	ng	
88) Benzo(b)fluoranthene	15.22	252	58103m	4.301	ng	
90) Benzo(a)pyrene	15.60	252	34564	2.781	ng	# 84

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

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