

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112018\
 Data File : BF110902.D
 Acq On : 20 Nov 2018 18:50
 Operator : JU/SJ
 Sample : J5763-03 2X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HR-06-111918-B

Manual Integrations
 APPROVED

Sohil
 11/21/2018 3:18:03 PM

Quant Time: Nov 21 04:43:22 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	6.94	152	64615	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	264382	20.00	ng	0.00
39) Acenaphthene-d10	9.98	164	108021	20.00	ng	-0.01
64) Phenanthrene-d10	11.48	188	165817	20.00	ng	0.00
76) Chrysene-d12	14.13	240	136110	20.00	ng	0.00
87) Perylene-d12	15.67	264	91515	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	234180	58.87	ng	0.00
7) Phenol-d6	6.56	99	296889	58.36	ng	-0.01
23) Nitrobenzene-d5	7.50	82	185060	40.31	ng	-0.01
42) 2,4,6-Tribromophenol	10.78	330	50715	46.59	ng	0.00
45) 2-Fluorobiphenyl	9.29	172	303008	40.06	ng	-0.01
79) Terphenyl-d14	13.06	244	221924	30.53	ng	0.00
Target Compounds						
10) Phenol	6.58	94	14495	2.457	ng	85
50) Dimethylphthalate	9.69	163	28479	3.533	ng	99
71) Phenanthrene	11.50	178	35676	4.016	ng	98
75) Fluoranthene	12.70	202	43518	4.886	ng	98
78) Pyrene	12.94	202	38846	3.707	ng	98
81) Benzo(a)anthracene	14.12	228	17693	2.175	ng	# 91
88) Benzo(b)fluoranthene	15.22	252	12136m	2.217	ng	

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

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