

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF111518\
 Data File : BF110799.D
 Acq On : 15 Nov 2018 20:13
 Operator : JU/SJ
 Sample : J5767-15 2X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CL-01-111418-D

Manual Integrations
 APPROVED

Sohil
 11/16/2018 10:07:27 AM

Quant Time: Nov 16 07:30:09 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	60468	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	241737	20.00	ng	0.00
39) Acenaphthene-d10	9.98	164	105635	20.00	ng	-0.01
64) Phenanthrene-d10	11.48	188	174246	20.00	ng	0.00
76) Chrysene-d12	14.13	240	157672	20.00	ng	0.00
87) Perylene-d12	15.66	264	112824	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.55	112	193253	51.91	ng	-0.01
7) Phenol-d6	6.56	99	244770	51.42	ng	-0.01
23) Nitrobenzene-d5	7.50	82	150704	35.90	ng	-0.01
42) 2,4,6-Tribromophenol	10.77	330	49187	46.21	ng	-0.01
45) 2-Fluorobiphenyl	9.29	172	278306	37.63	ng	-0.01
79) Terphenyl-d14	13.06	244	236707	28.12	ng	-0.01
Target Compounds						
10) Phenol	6.57	94	12300	2.228	ng	Qvalue 86
50) Dimethylphthalate	9.69	163	27671	3.511	ng	100
71) Phenanthrene	11.50	178	44457	4.762	ng	97
75) Fluoranthene	12.70	202	50025	5.345	ng	98
78) Pyrene	12.93	202	70621	5.817	ng	97
81) Benzo(a)anthracene	14.12	228	28509m	3.025	ng	
83) Chrysene	14.16	228	32270	3.594	ng	96
88) Benzo(b)fluoranthene	15.20	252	21804m	3.230	ng	
90) Benzo(a)pyrene	15.60	252	15563	2.538	ng	# 82

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

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