

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112918\
 Data File : BF111097.D
 Acq On : 29 Nov 2018 18:20
 Operator : JU/SJ
 Sample : J5767-21 2X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

mohammad
 11/30/2018 3:49:44 PM

Quant Time: Nov 30 03:05:25 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	65159	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	270760	20.00	ng	0.00
39) Acenaphthene-d10	9.97	164	120511	20.00	ng	-0.01
64) Phenanthrene-d10	11.47	188	186042	20.00	ng	0.00
76) Chrysene-d12	14.13	240	150031	20.00	ng	0.00
87) Perylene-d12	15.66	264	109088	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.55	112	175623	44.70	ng	0.00
7) Phenol-d6	6.56	99	235418	45.98	ng	-0.01
23) Nitrobenzene-d5	7.50	82	146487	30.96	ng	-0.01
42) 2,4,6-Tribromophenol	10.77	330	44467	37.15	ng	-0.01
45) 2-Fluorobiphenyl	9.29	172	252055	30.30	ng	0.00
79) Terphenyl-d14	13.05	244	185809	24.11	ng	-0.01
Target Compounds						
10) Phenol	6.57	94	18590	3.224	ng	Qvalue 87
50) Dimethylphthalate	9.69	163	25486	2.884	ng	98
71) Phenanthrene	11.49	178	36635	3.718	ng	99
75) Fluoranthene	12.69	202	46933	4.634	ng	98
78) Pyrene	12.93	202	42445	3.873	ng	98
81) Benzo(a)anthracene	14.12	228	21258m	2.400	ng	
83) Chrysene	14.15	228	18952	2.173	ng	98
88) Benzo(b)fluoranthene	15.21	252	18004m	2.809	ng	
90) Benzo(a)pyrene	15.60	252	12874	2.158	ng	# 82

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

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