

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110818\
 Data File : BF110603.D
 Acq On : 8 Nov 2018 23:35
 Operator : JU/SJ
 Sample : J5768-01 2X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OK-02-110718-A

Manual Integrations
 APPROVED

Sohil
 11/9/2018 4:26:13 PM

Quant Time: Nov 09 12:46:03 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 08 14:14:47 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	84107	20.00	ng	0.00
21) Naphthalene-d8	8.23	136	351943	20.00	ng	0.00
39) Acenaphthene-d10	9.98	164	157880	20.00	ng	-0.01
64) Phenanthrene-d10	11.47	188	262349	20.00	ng	-0.01
76) Chrysene-d12	14.13	240	175264	20.00	ng	-0.01
87) Perylene-d12	15.65	264	165421	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	217434	41.99	ng	0.00
7) Phenol-d6	6.56	99	281990	42.59	ng	-0.01
23) Nitrobenzene-d5	7.50	82	175914	28.79	ng	-0.01
42) 2,4,6-Tribromophenol	10.77	330	52740	33.15	ng	-0.01
45) 2-Fluorobiphenyl	9.30	172	336122	30.41	ng	0.00
79) Terphenyl-d14	13.06	244	236259	25.25	ng	-0.01
Target Compounds						
10) Phenol	6.57	94	20316	2.646	ng	Qvalue # 69
50) Dimethylphthalate	9.69	163	42808	3.634	ng	99
58) Fluorene	10.53	166	24383	2.247	ng	98
71) Phenanthrene	11.50	178	226514	16.116	ng	98
72) Anthracene	11.56	178	54999	3.879	ng	98
75) Fluoranthene	12.70	202	213768	15.170	ng	97
78) Pyrene	12.93	202	200328	14.845	ng	98
81) Benzo(a)anthracene	14.12	228	83695	7.989	ng	97
83) Chrysene	14.15	228	72909	7.305	ng	97
86) Indeno(1,2,3-cd)pyrene	17.21	276	26532	3.705	ng	# 86
88) Benzo(b)fluoranthene	15.20	252	74625m	7.541	ng	
89) Benzo(k)fluoranthene	15.23	252	28401m	2.904	ng	
90) Benzo(a)pyrene	15.59	252	58053	6.457	ng	98
92) Benzo(a,h,i)perylene	17.69	276	26189	3.469	ng	# 92

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

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