

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112918\
 Data File : BF111095.D
 Acq On : 29 Nov 2018 17:25
 Operator : JU/SJ
 Sample : J5767-17 2X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

mohammad
 11/30/2018 3:49:37 PM

Quant Time: Nov 30 02:56:15 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	60747	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	253590	20.00	ng	0.00
39) Acenaphthene-d10	9.97	164	113077	20.00	ng	-0.01
64) Phenanthrene-d10	11.47	188	170347	20.00	ng	0.00
76) Chrysene-d12	14.13	240	135519	20.00	ng	0.00
87) Perylene-d12	15.66	264	118725	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.55	112	164453	44.89	ng	0.00
7) Phenol-d6	6.56	99	222165	46.54	ng	0.00
23) Nitrobenzene-d5	7.50	82	139023	31.37	ng	0.00
42) 2,4,6-Tribromophenol	10.77	330	40884	36.41	ng	0.00
45) 2-Fluorobiphenyl	9.29	172	253288	32.45	ng	0.00
79) Terphenyl-d14	13.04	244	173489	24.92	ng	-0.01
Target Compounds						
10) Phenol	6.57	94	17943	3.338	ng	93
50) Dimethylphthalate	9.69	163	20016	2.414	ng	96
52) Acenaphthene	10.00	154	24296	3.526	ng	96
58) Fluorene	10.53	166	23054	2.948	ng	99
71) Phenanthrene	11.49	178	236211	26.182	ng	99
72) Anthracene	11.55	178	59531	6.479	ng	98
75) Fluoranthene	12.69	202	285045	30.740	ng	97
78) Pyrene	12.92	202	239213	24.168	ng	97
81) Benzo(a)anthracene	14.12	228	127318	15.913	ng	99
83) Chrysene	14.15	228	107411	13.637	ng	99
86) Indeno(1,2,3-cd)pyrene	17.26	276	33176	5.868	ng	98
88) Benzo(b)fluoranthene	15.20	252	117147m	16.796	ng	
89) Benzo(k)fluoranthene	15.23	252	42284m	5.890	ng	
90) Benzo(a)pyrene	15.60	252	78336	12.066	ng	100
92) Benzo(a,h,i)perylene	17.74	276	28741	5.565	ng	96

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

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