

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112018\
 Data File : BF110890.D
 Acq On : 20 Nov 2018 13:23
 Operator : JU/SJ
 Sample : J6010-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FS-RT3786

Manual Integrations
 APPROVED

Sohil
 11/21/2018 3:17:48 PM

Quant Time: Nov 20 14:35:03 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	77879	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	322696	20.00	ng	0.00
39) Acenaphthene-d10	9.99	164	147216	20.00	ng	0.00
64) Phenanthrene-d10	11.49	188	253313	20.00	ng	0.00
76) Chrysene-d12	14.13	240	171910	20.00	ng	0.00
87) Perylene-d12	15.66	264	176059	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.57	112	602888	125.74	ng	0.00
7) Phenol-d6	6.57	99	760192	123.99	ng	0.00
23) Nitrobenzene-d5	7.51	82	496918	88.68	ng	0.00
42) 2,4,6-Tribromophenol	10.79	330	165334	111.46	ng	0.00
45) 2-Fluorobiphenyl	9.30	172	847336	82.20	ng	0.00
79) Terphenyl-d14	13.07	244	637953	69.50	ng	0.00
Target Compounds						
10) Phenol	6.59	94	40800	5.739	ng	# 70
50) Dimethylphthalate	9.69	163	64162	5.841	ng	# 97
75) Fluoranthene	12.70	202	46982	3.453	ng	96
78) Pyrene	12.94	202	47559	3.593	ng	96
81) Benzo(a)anthracene	14.13	228	25553	2.487	ng	95
83) Chrysene	14.16	228	25655	2.621	ng	97
86) Indeno(1,2,3-cd)pyrene	17.26	276	17848	2.541	ng	98
88) Benzo(b)fluoranthene	15.21	252	39931m	3.791	ng	
90) Benzo(a)pyrene	15.60	252	27751	2.900	ng	# 91
92) Benzo(g,h,i)perylene	17.74	276	16344	2.034	ng	95

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

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