

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111218\
 Data File : BF110690.D
 Acq On : 12 Nov 2018 22:26
 Operator : JU/SJ
 Sample : J5901-01
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S-N

Manual Integrations
 APPROVED

Sohil
 11/13/2018 11:16:33 AM

Quant Time: Nov 13 06:33:40 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 09 15:13:13 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	86170	20.00	ng	0.00
21) Naphthalene-d8	8.23	136	375070	20.00	ng	0.00
39) Acenaphthene-d10	9.99	164	172797	20.00	ng	0.00
64) Phenanthrene-d10	11.48	188	310778	20.00	ng	0.00
76) Chrysene-d12	14.13	240	177778	20.00	ng	-0.01
87) Perylene-d12	15.65	264	165580	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	522978	98.58	ng	0.00
7) Phenol-d6	6.57	99	694080	102.31	ng	0.00
23) Nitrobenzene-d5	7.51	82	435449	66.86	ng	0.00
42) 2,4,6-Tribromophenol	10.78	330	129353	74.29	ng	0.00
45) 2-Fluorobiphenyl	9.30	172	725806	59.99	ng	0.00
79) Terphenyl-d14	13.06	244	558209	58.80	ng	-0.01
Target Compounds						
10) Phenol	6.58	94	42847	5.447	ng	Qvalue # 69
50) Dimethylphthalate	9.69	163	128326	9.953	ng	99
71) Phenanthrene	11.50	178	78129	4.692	ng	99
75) Fluoranthene	12.70	202	174132	10.432	ng	98
78) Pyrene	12.93	202	145923	10.660	ng	97
81) Benzo(a)anthracene	14.12	228	72259	6.800	ng	97
83) Chrysene	14.15	228	60921	6.018	ng	98
86) Indeno(1,2,3-cd)pyrene	17.22	276	34532	4.754	ng	96
88) Benzo(b)fluoranthene	15.20	252	74262m	7.497	ng	
89) Benzo(k)fluoranthene	15.23	252	28586m	2.921	ng	
90) Benzo(a)pyrene	15.59	252	57650	6.406	ng	100
92) Benzo(a,h,i)perylene	17.69	276	31195m	4.128	ng	

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

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