

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112118\
 Data File : BF110928.D
 Acq On : 21 Nov 2018 15:10
 Operator : JU/SJ
 Sample : J5768-05
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OK-03-112018

Manual Integrations
 APPROVED

mohammad
 11/23/2018 11:49:00 AM

Quant Time: Nov 21 22:35:32 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	61631	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	258434	20.00	ng	0.00
39) Acenaphthene-d10	9.98	164	117544	20.00	ng	-0.01
64) Phenanthrene-d10	11.48	188	198371	20.00	ng	0.00
76) Chrysene-d12	14.13	240	137606	20.00	ng	0.00
87) Perylene-d12	15.67	264	113443	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	375183	98.88	ng	0.00
7) Phenol-d6	6.57	99	475819	98.07	ng	0.00
23) Nitrobenzene-d5	7.50	82	309424	68.95	ng	-0.01
42) 2,4,6-Tribromophenol	10.78	330	101120	85.38	ng	0.00
45) 2-Fluorobiphenyl	9.29	172	542529	65.92	ng	-0.01
79) Terphenyl-d14	13.06	244	398722	54.26	ng	-0.01
Target Compounds						
10) Phenol	6.58	94	24436	4.343	ng	Qvalue # 76
50) Dimethylphthalate	9.69	163	52892	6.030	ng	100
71) Phenanthrene	11.50	178	52958	4.983	ng	98
75) Fluoranthene	12.70	202	124896	11.722	ng	98
78) Pyrene	12.93	202	119960	11.322	ng	98
81) Benzo(a)anthracene	14.12	228	67676	8.228	ng	98
83) Chrysene	14.16	228	56448	7.204	ng	97
86) Indeno(1,2,3-cd)pyrene	17.26	276	21820	3.881	ng	94
88) Benzo(b)fluoranthene	15.22	252	69638m	10.261	ng	
89) Benzo(k)fluoranthene	15.24	252	20984m	3.129	ng	
90) Benzo(a)pyrene	15.60	252	45727	7.416	ng	98
92) Benzo(a,h,i)perylene	17.75	276	19478	3.762	ng	# 94

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

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