

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF021519\
 Data File : BF112592.D
 Acq On : 15 Feb 2019 21:26
 Operator : JU/SJ
 Sample : K1358-03 2X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TR-04-021419

Manual Integrations
 APPROVED

Sohil
 2/18/2019 12:06:41 PM

Quant Time: Feb 16 00:14:02 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 28 10:38:18 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	28941	20.00	ng	-0.01
21) Naphthalene-d8	8.12	136	120411	20.00	ng	-0.01
39) Acenaphthene-d10	9.87	164	57686	20.00	ng	-0.02
64) Phenanthrene-d10	11.36	188	98985	20.00	ng	0.00
76) Chrysene-d12	14.01	240	74082	20.00	ng	0.00
87) Perylene-d12	15.49	264	56417	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	187543	137.64	ng	-0.01
7) Phenol-d6	6.48	99	233239	123.19	ng	-0.01
23) Nitrobenzene-d5	7.39	82	147725	70.69	ng	-0.02
42) 2,4,6-Tribromophenol	10.67	330	71047	105.81	ng	-0.01
45) 2-Fluorobiphenyl	9.19	172	302335	74.13	ng	-0.02
79) Terphenyl-d14	12.94	244	239759	60.96	ng	0.00
Target Compounds						
10) Phenol	6.50	94	8371	4.030	ng	Qvalue # 76
50) Dimethylphthalate	9.58	163	12433	2.774	ng	97
71) Phenanthrene	11.39	178	23117	4.492	ng	97
75) Fluoranthene	12.59	202	57145	10.353	ng	97
78) Pyrene	12.82	202	56400	10.996	ng	97
81) Benzo(a)anthracene	14.00	228	25859	5.938	ng	97
83) Chrysene	14.04	228	25654m	6.171	ng	
86) Indeno(1,2,3-cd)pyrene	17.00	276	15157	4.601	ng	96
88) Benzo(b)fluoranthene	15.06	252	43475m	12.885	ng	
89) Benzo(k)fluoranthene	15.09	252	12522m	3.875	ng	
90) Benzo(a)pyrene	15.43	252	28950	9.536	ng	# 93
92) Benzo(a,h,i)perylene	17.46	276	15300	6.628	ng	# 90

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

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