

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011019\
 Data File : BF111998.D
 Acq On : 10 Jan 2019 20:10
 Operator : JU/SJ
 Sample : K1071-01
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CPSB-17(10-15)

Manual Integrations
 APPROVED

Sohil
 1/11/2019 4:55:04 PM

Quant Time: Jan 11 16:48:35 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF010719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 07 14:33:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	79895	20.00	ng	0.00
21) Naphthalene-d8	8.19	136	196618m	20.00	ng	0.02
39) Acenaphthene-d10	9.96	164	90114m	20.00	ng	0.03
64) Phenanthrene-d10	11.42	188	228023	20.00	ng	0.00
76) Chrysene-d12	14.06	240	204590	20.00	ng	0.00
87) Perylene-d12	15.54	264	174012	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	470021	100.65	ng	-0.02
7) Phenol-d6	6.52	99	583679	96.97	ng	-0.01
23) Nitrobenzene-d5	7.45	82	483332	131.38	ng	-0.01
42) 2,4,6-Tribromophenol	10.73	330	144777	123.52	ng	0.00
45) 2-Fluorobiphenyl	9.27	172	463856	75.06	ng	0.02
79) Terphenyl-d14	13.00	244	698872	64.84	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.53	94	36570	5.590	ng	94
37) 2-Methylnaphthalene	8.93	142	158461m	27.399	ng	
52) Acenaphthene	9.99	154	172353m	30.567	ng	
55) Dibenzofuran	10.16	168	194306m	23.695	ng	
58) Fluorene	10.49	166	238833m	40.026	ng	
71) Phenanthrene	11.44	178	766548	62.877	ng	99
72) Anthracene	11.49	178	130012	10.505	ng	96
75) Fluoranthene	12.64	202	453137	36.206	ng	97
78) Pyrene	12.87	202	424240	30.137	ng	100
81) Benzo(a)anthracene	14.05	228	150759	12.804	ng	# 94
83) Chrysene	14.09	228	161626	14.397	ng	# 94
84) Bis(2-ethylhexyl)phthalate	14.03	149	314294	40.213	ng	96
86) Indeno(1,2,3-cd)pyrene	17.03	276	84807	10.362	ng	98
88) Benzo(b)fluoranthene	15.11	252	204966m	19.691	ng	
89) Benzo(k)fluoranthene	15.13	252	59514m	6.082	ng	
90) Benzo(a)pyrene	15.48	252	121396	13.264	ng	# 70
91) Dibenzo(a,h)anthracene	17.04	278	22885	3.109	ng	# 33
92) Benzo(g,h,i)perylene	17.49	276	92883	12.900	ng	94

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

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