

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011119\
 Data File : BF112014.D
 Acq On : 11 Jan 2019 14:46
 Operator : JU/SJ
 Sample : K1088-02
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CPSB-06(10-15)

Manual Integrations
 APPROVED

Sohil
 1/14/2019 4:10:13 PM

Quant Time: Jan 11 22:31:45 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF010719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 11 21:37:26 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	118852	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	321757	20.00	ng	0.00
39) Acenaphthene-d10	9.92	164	136723	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	212268	20.00	ng	0.02
76) Chrysene-d12	14.10	240	201378	20.00	ng	0.06
87) Perylene-d12	15.60	264	160583	20.00	ng	0.08
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	697538	100.41	ng	0.00
7) Phenol-d6	6.52	99	890187	99.42	ng	0.00
23) Nitrobenzene-d5	7.43	82	435714	72.37	ng	0.00
42) 2,4,6-Tribromophenol	10.72	330	163611	92.00	ng	0.01
45) 2-Fluorobiphenyl	9.23	172	697330	74.37	ng	0.00
79) Terphenyl-d14	13.03	244	671308	63.28	ng	0.04
Target Compounds						
10) Phenol	6.53	94	44352	4.557	ng	# 75
50) Dimethylphthalate	9.62	163	106459	10.365	ng	# 89
52) Acenaphthene	9.95	154	173823	20.318	ng	# 90
55) Dibenzofuran	10.12	168	89142	7.165	ng	# 84
58) Fluorene	10.47	166	146957	16.233	ng	# 93
71) Phenanthrene	11.44	178	310778	27.384	ng	# 90
72) Anthracene	11.50	178	159857	13.876	ng	# 84
75) Fluoranthene	12.67	202	587954	50.465	ng	# 94
78) Pyrene	12.90	202	559994m	40.415	ng	# 76
81) Benzo(a)anthracene	14.09	228	184288	15.901	ng	# 76
83) Chrysene	14.13	228	175790	15.908	ng	# 86
84) Bis(2-ethylhexyl)phthalate	14.06	149	257755	33.505	ng	# 98
86) Indeno(1,2,3-cd)pyrene	17.09	276	41677	5.173	ng	# 1
88) Benzo(b)fluoranthene	15.16	252	115889m	12.064	ng	# 71
89) Benzo(k)fluoranthene	15.17	252	35568m	3.939	ng	#
90) Benzo(a)pyrene	15.53	252	93571	11.078	ng	#
92) Benzo(a,h,i)perylene	17.53	276	43336	6.522	ng	#

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

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