

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF031919\
 Data File : BF113134.D
 Acq On : 19 Mar 2019 19:15
 Operator : JU/SJ
 Sample : K1935-10
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-7(6-7)

Manual Integrations
 APPROVED

Sohil
 3/20/2019 6:51:25 PM

Quant Time: Mar 20 02:25:34 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 13 03:42:55 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.72	152	165742	20.00	ng	-0.01
21) Naphthalene-d8	8.00	136	450422	20.00	ng	-0.01
39) Acenaphthene-d10	9.76	164	180311	20.00	ng	0.00
64) Phenanthrene-d10	11.24	188	296021	20.00	ng	0.00
76) Chrysene-d12	13.87	240	242712	20.00	ng	0.00
87) Perylene-d12	15.29	264	206814	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	992076	93.11	ng	0.00
7) Phenol-d6	6.37	99	1287440	96.19	ng	0.00
23) Nitrobenzene-d5	7.29	82	739395	98.23	ng	-0.01
42) 2,4,6-Tribromophenol	10.55	330	220036	114.95	ng	0.00
45) 2-Fluorobiphenyl	9.09	172	1017680	97.16	ng	0.00
79) Terphenyl-d14	12.83	244	767667	61.31	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.47	163	79880	6.008	ng	Qvalue # 92
52) Acenaphthene	9.79	154	53539	4.696	ng	96
71) Phenanthrene	11.27	178	102451	6.956	ng	# 92
72) Anthracene	11.32	178	46087	2.989	ng	# 77
75) Fluoranthene	12.46	202	159571	10.113	ng	95
78) Pyrene	12.69	202	214219	10.470	ng	97
81) Benzo(a)anthracene	13.87	228	134707	8.008	ng	# 89
83) Chrysene	13.90	228	154852	9.398	ng	# 89
86) Indeno(1,2,3-cd)pyrene	16.65	276	58850	4.189	ng	# 92
88) Benzo(b)fluoranthene	14.89	252	138019m	10.317	ng	
89) Benzo(k)fluoranthene	14.91	252	42066m	3.472	ng	
90) Benzo(a)pyrene	15.23	252	112463	9.505	ng	# 86
91) Dibenzo(a,h)anthracene	16.66	278	31557	3.125	ng	# 65
92) Benzo(g,h,i)perylene	17.06	276	92394	9.113	ng	96

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

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