

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF021819\
 Data File : BF112609.D
 Acq On : 18 Feb 2019 19:00
 Operator : JU/SJ
 Sample : K1518-13
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CABLE-BRIDGE-FOUNDATION-D-7

Manual Integrations
 APPROVED

Sohil
 2/19/2019 11:57:32 AM

Quant Time: Feb 19 10:42:42 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	6.82	152	59984	20.00	ng	-0.02
21) Naphthalene-d8	8.10	136	239801	20.00	ng	-0.02
39) Acenaphthene-d10	9.86	164	119258	20.00	ng	-0.03
64) Phenanthrene-d10	11.35	188	228590	20.00	ng	-0.02
76) Chrysene-d12	14.00	240	158813	20.00	ng	-0.01
87) Perylene-d12	15.47	264	148006	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.45	112	360916	127.80	ng	-0.03
7) Phenol-d6	6.47	99	466755	118.95	ng	-0.02
23) Nitrobenzene-d5	7.39	82	299498	71.96	ng	-0.02
42) 2,4,6-Tribromophenol	10.66	330	160106	115.34	ng	-0.02
45) 2-Fluorobiphenyl	9.18	172	585540	69.44	ng	-0.02
79) Terphenyl-d14	12.93	244	609159	72.24	ng	-0.02

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
10) Phenol	6.49	94	9971	2.316	ng	# 64
49) Acenaphthylene	9.72	152	30363	2.562	ng	99
50) Dimethylphthalate	9.57	163	50222	5.420	ng	99
71) Phenanthrene	11.37	178	107077	9.010	ng	97
72) Anthracene	11.43	178	40629	3.432	ng	98
73) Carbazole	11.59	167	26891	2.303	ng	96
75) Fluoranthene	12.57	202	205951	16.158	ng	97
78) Pyrene	12.80	202	199634	18.156	ng	99
81) Benzo(a)anthracene	13.99	228	85441	9.152	ng	98
83) Chrysene	14.02	228	93143	10.452	ng	97
86) Indeno(1,2,3-cd)pyrene	16.97	276	60383m	8.551	ng	
88) Benzo(b)fluoranthene	15.04	252	158666m	17.925	ng	
89) Benzo(k)fluoranthene	15.07	252	44779m	5.282	ng	
90) Benzo(a)pyrene	15.42	252	73801	9.266	ng	97
91) Dibenzo(a,h)anthracene	16.96	278	14759	2.299	ng	# 79
92) Benzo(g,h,i)perylene	17.42	276	60941	10.063	ng	97

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

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