

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF041020\
 Data File : BF119909.D
 Acq On : 10 Apr 2020 21:48
 Operator : MA/SJ
 Sample : L2199-12
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-37-36-COMP

Manual Integrations
 APPROVED

mohammad
 4/14/2020 3:39:31 PM

Quant Time: Apr 11 04:00:23 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 10 13:48:19 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.75	152	174180	20.00	ng	0.00
21) Naphthalene-d8	8.04	136	615848	20.00	ng	0.00
39) Acenaphthene-d10	9.79	164	267785	20.00	ng	0.00
64) Phenanthrene-d10	11.28	188	449019	20.00	ng	0.00
76) Chrysene-d12	13.93	240	346675	20.00	ng	0.00
87) Perylene-d12	15.36	264	280869	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.36	112	904504	89.74	ng	0.00
7) Phenol-d6	6.39	99	1152108	88.71	ng	0.00
23) Nitrobenzene-d5	7.32	82	677561	56.92	ng	0.00
42) 2,4,6-Tribromophenol	10.59	330	230499	70.30	ng	0.00
45) 2-Fluorobiphenyl	9.12	172	1171276	62.10	ng	0.00
79) Terphenyl-d14	12.87	244	1070604	51.96	ng	0.00

Target Compounds				Qvalue		
31) Naphthalene	8.06	128	67087	2.070	ng	98
37) 2-Methylnaphthalene	8.75	142	78540	3.575	ng	98
38) 1-Methylnaphthalene	8.85	142	74346	3.591	ng	99
49) Acenaphthylene	9.66	152	59050	2.230	ng	# 93
50) Dimethylphthalate	9.51	163	181064	8.742	ng	100
52) Acenaphthene	9.83	154	76018	4.304	ng	97
71) Phenanthrene	11.31	178	925245	38.718	ng	99
72) Anthracene	11.36	178	331630	13.584	ng	97
75) Fluoranthene	12.50	202	517557	20.072	ng	97
78) Pyrene	12.73	202	862097	29.498	ng	99
81) Benzo(a)anthracene	13.92	228	337040	14.778	ng	96
83) Chrysene	13.95	228	311096	13.425	ng	98
86) Indeno(1,2,3-cd)pyrene	16.77	276	125676	6.152	ng	98
88) Benzo(b)fluoranthene	14.95	252	283576m	14.035	ng	
89) Benzo(k)fluoranthene	14.97	252	78853m	4.523	ng	
90) Benzo(a)pyrene	15.30	252	279392	16.446	ng	# 94
91) Dibenzo(a,h)anthracene	16.77	278	36558	2.429	ng	# 80
92) Benzo(g,h,i)perylene	17.20	276	152683	10.279	ng	97

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

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