

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF051520\
 Data File : BF120330.D
 Acq On : 15 May 2020 19:15
 Operator : MA/SJ
 Sample : L2665-02
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 0375-AMENDED-COMP-MP-A

Quant Time: May 16 04:29:33 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF051420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 15 12:31:31 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.62	152	273774	20.00	ng	0.00
21) Naphthalene-d8	7.90	136	925132	20.00	ng	0.00
39) Acenaphthene-d10	9.66	164	479836	20.00	ng	0.00
64) Phenanthrene-d10	11.14	188	814676	20.00	ng	0.00
76) Chrysene-d12	13.77	240	455385	20.00	ng	0.00
87) Perylene-d12	15.17	264	498230	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.22	112	1375099	97.73	ng	0.00
7) Phenol-d6	6.27	99	1731517	97.31	ng	0.00
23) Nitrobenzene-d5	7.18	82	1028803	76.14	ng	-0.01
42) 2,4,6-Tribromophenol	10.45	330	361639	92.44	ng	0.00
45) 2-Fluorobiphenyl	8.98	172	1837607	78.68	ng	0.00
79) Terphenyl-d14	12.73	244	1507243	73.12	ng	0.00

Target Compounds				Qvalue		
31) Naphthalene	7.92	128	249974	6.402	ng	# 90
37) 2-Methylnaphthalene	8.62	142	453524	16.974	ng	97
38) 1-Methylnaphthalene	8.72	142	361288	13.941	ng	# 99
50) Dimethylphthalate	9.38	163	329783	10.872	ng	# 97
52) Acenaphthene	9.69	154	102036	4.419	ng	93
58) Fluorene	10.20	166	197863	7.813	ng	# 94
71) Phenanthrene	11.16	178	1258684	36.768	ng	98
72) Anthracene	11.22	178	323262	8.821	ng	96
73) Carbazole	11.37	167	104076	3.039	ng	# 77
75) Fluoranthene	12.35	202	934347	25.465	ng	99
78) Pyrene	12.58	202	969986	29.878	ng	98
81) Benzo(a)anthracene	13.77	228	431280	17.956	ng	# 95
83) Chrysene	13.80	228	447758	17.361	ng	# 93
86) Indeno(1,2,3-cd)pyrene	16.49	276	144743	6.183	ng	96
88) Benzo(b)fluoranthene	14.78	252	375337m	13.794	ng	
89) Benzo(k)fluoranthene	14.80	252	103413m	4.387	ng	
90) Benzo(a)pyrene	15.11	252	323715	13.266	ng	# 94
91) Dibenzo(a,h)anthracene	16.49	278	50093	2.317	ng	# 73
92) Benzo(a,h,i)perylene	16.89	276	164017	7.524	ng	98

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

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