

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF042120\
 Data File : BF120114.D
 Acq On : 21 Apr 2020 20:41
 Operator : MA/SJ
 Sample : L2330-04
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WB-10-2

Manual Integrations
 APPROVED

mohammad
 4/24/2020 4:17:17 AM

Quant Time: Apr 22 03:51:03 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 21 13:48:22 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.69	152	143130	20.00	ng	0.00
21) Naphthalene-d8	7.98	136	527026	20.00	ng	0.00
39) Acenaphthene-d10	9.73	164	243201	20.00	ng	0.00
64) Phenanthrene-d10	11.22	188	356705	20.00	ng	0.00
76) Chrysene-d12	13.87	240	283319	20.00	ng	0.00
87) Perylene-d12	15.29	264	249936	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.29	112	588933	71.11	ng	0.00
7) Phenol-d6	6.33	99	750449	70.32	ng	0.00
23) Nitrobenzene-d5	7.26	82	469363	46.07	ng	0.00
42) 2,4,6-Tribromophenol	10.53	330	138870	46.64	ng	0.00
45) 2-Fluorobiphenyl	9.06	172	864690	50.48	ng	0.00
79) Terphenyl-d14	12.81	244	658813	39.13	ng	0.00
Target Compounds						
22) Acetophenone	7.10	105	37548	3.042	ng	Qvalue # 87
31) Naphthalene	8.00	128	1220783	44.027	ng	99
37) 2-Methylnaphthalene	8.69	142	269547	14.338	ng	99
38) 1-Methylnaphthalene	8.79	142	257744	14.546	ng	98
46) 1,1'-Biphenyl	9.16	154	73561	3.584	ng	100
49) Acenaphthylene	9.59	152	71218	2.961	ng	# 89
50) Dimethylphthalate	9.46	163	117477	6.245	ng	99
52) Acenaphthene	9.77	154	155951	9.721	ng	95
55) Dibenzofuran	9.94	168	105531	4.794	ng	# 87
58) Fluorene	10.29	166	209187	11.882	ng	98
71) Phenanthrene	11.25	178	815768	42.971	ng	99
72) Anthracene	11.30	178	186189	9.601	ng	98
73) Carbazole	11.46	167	37868	2.065	ng	# 82
75) Fluoranthene	12.44	202	558701	27.275	ng	98
78) Pyrene	12.67	202	644332	26.977	ng	99
81) Benzo(a)anthracene	13.86	228	240068	12.880	ng	# 91
83) Chrysene	13.89	228	243587m	12.862	ng	
86) Indeno(1,2,3-cd)pyrene	16.66	276	89620	5.368	ng	99
88) Benzo(b)fluoranthene	14.89	252	217738m	12.110	ng	
89) Benzo(k)fluoranthene	14.91	252	61812m	3.984	ng	
90) Benzo(a)pyrene	15.23	252	166214	10.995	ng	# 96
92) Benzo(g,h,i)perylene	17.07	276	96954	7.335	ng	96

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

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