

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF042420\
 Data File : BF120199.D
 Acq On : 24 Apr 2020 13:50
 Operator : MA/SJ
 Sample : L2351-10
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-21A

Manual Integrations
 APPROVED

mohammad
 4/27/2020 3:04:21 PM

Quant Time: Apr 24 14:26:13 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 24 10:24:30 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.66	152	177571	20.00	ng	0.00
21) Naphthalene-d8	7.95	136	656148	20.00	ng	0.00
39) Acenaphthene-d10	9.70	164	311306	20.00	ng	0.00
64) Phenanthrene-d10	11.18	188	448550	20.00	ng	0.00
76) Chrysene-d12	13.82	240	321173	20.00	ng	-0.01
87) Perylene-d12	15.22	264	300945	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.27	112	872036	84.87	ng	0.00
7) Phenol-d6	6.32	99	1117146	84.38	ng	0.00
23) Nitrobenzene-d5	7.23	82	689161	54.34	ng	0.00
42) 2,4,6-Tribromophenol	10.49	330	234643	61.56	ng	0.00
45) 2-Fluorobiphenyl	9.03	172	1275980	58.19	ng	0.00
79) Terphenyl-d14	12.77	244	1024643	53.68	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.42	163	199389	8.281	ng	Qvalue 100
75) Fluoranthene	12.39	202	65542	2.545	ng	97
78) Pyrene	12.62	202	82310	3.040	ng	97
81) Benzo(a)anthracene	13.81	228	50408	2.386	ng	# 78
83) Chrysene	13.84	228	54172	2.523	ng	# 93
86) Indeno(1,2,3-cd)pyrene	16.57	276	46423	2.453	ng	95
88) Benzo(b)fluoranthene	14.83	252	96012m	4.435	ng	
90) Benzo(a)pyrene	15.16	252	72896	4.005	ng	# 96
92) Benzo(g,h,i)perylene	16.97	276	46983	2.952	ng	# 90

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

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