

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012120\
 Data File : BF118650.D
 Acq On : 21 Jan 2020 23:37
 Operator : CG/JU
 Sample : L1210-06
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 Q012-WC1

Manual Integrations
 APPROVED

mohammad
 1/22/2020 1:36:31 PM

Quant Time: Jan 22 05:04:10 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 20 14:45:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	402617	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	1414738	20.00	ng	-0.01
39) Acenaphthene-d10	9.92	164	718696	20.00	ng	-0.01
64) Phenanthrene-d10	11.40	188	1069014	20.00	ng	-0.01
76) Chrysene-d12	14.03	240	909819	20.00	ng	-0.01
87) Perylene-d12	15.51	264	1003822	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	1842592	85.00	ng	0.00
7) Phenol-d6	6.50	99	2340088	83.05	ng	-0.01
23) Nitrobenzene-d5	7.43	82	1495705	59.14	ng	-0.01
42) 2,4,6-Tribromophenol	10.70	330	579763	69.74	ng	-0.01
45) 2-Fluorobiphenyl	9.23	172	2728531	62.40	ng	-0.01
79) Terphenyl-d14	12.98	244	2454815	58.85	ng	-0.01
Target Compounds						
50) Dimethylphthalate	9.62	163	216802	4.630	ng	98
71) Phenanthrene	11.42	178	260894	5.059	ng	93
75) Fluoranthene	12.61	202	329980	6.066	ng	93
78) Pyrene	12.84	202	320677	5.192	ng	92
81) Benzo(a)anthracene	14.02	228	184146	3.365	ng	98
83) Chrysene	14.06	228	203079	3.872	ng	96
88) Benzo(b)fluoranthene	15.07	252	251503m	3.986	ng	
90) Benzo(a)pyrene	15.44	252	149226	2.742	ng	# 88

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

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