

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071719\
 Data File : BF115639.D
 Acq On : 17 Jul 2019 22:59
 Operator : HP/JU
 Sample : K3840-03 5X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-SF-02-005-B

Manual Integrations
 APPROVED

mohammad
 7/19/2019 2:26:16 PM

Quant Time: Jul 18 06:50:17 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 12 14:03:52 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	133240	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	489732	20.00	ng	0.00
39) Acenaphthene-d10	9.92	164	183074	20.00	ng	-0.01
64) Phenanthrene-d10	11.40	188	292751	20.00	ng	-0.01
76) Chrysene-d12	14.03	240	217356	20.00	ng	-0.02
87) Perylene-d12	15.50	264	177147	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	101799	12.50	ng	0.00
7) Phenol-d6	6.50	99	198982	19.25	ng	-0.01
23) Nitrobenzene-d5	7.43	82	120077	14.26	ng	-0.02
42) 2,4,6-Tribromophenol	10.70	330	11101	5.19	ng	-0.02
45) 2-Fluorobiphenyl	9.23	172	199643	19.09	ng	-0.02
79) Terphenyl-d14	12.98	244	147927	12.56	ng	-0.02
Target Compounds						
50) Dimethylphthalate	9.62	163	28039	2.036	ng	99
71) Phenanthrene	11.42	178	48463	3.230	ng	96
75) Fluoranthene	12.61	202	76728	4.839	ng	98
78) Pyrene	12.84	202	72193	3.920	ng	98
81) Benzo(a)anthracene	14.02	228	28886	2.017	ng	# 90
83) Chrysene	14.06	228	29863	2.077	ng	# 91
88) Benzo(b)fluoranthene	15.07	252	25632m	2.247	ng	

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

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