

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG041019\
 Data File : BG040417.D
 Acq On : 10 Apr 2019 16:18
 Operator : JU/SJ
 Sample : K2345-04
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TP-2

Manual Integrations
 APPROVED

Sohil
 4/11/2019 2:46:46 PM

Quant Time: Apr 11 05:30:26 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG032719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 28 05:47:00 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.05	152	73223	20.00	ng	-0.02
21) Naphthalene-d8	10.86	136	296785	20.00	ng	-0.02
39) Acenaphthene-d10	14.67	164	170377	20.00	ng	-0.03
64) Phenanthrene-d10	17.42	188	362201	20.00	ng	-0.02
76) Chrysene-d12	21.70	240	358623	20.00	ng	-0.03
87) Perylene-d12	24.97	264	380273	20.00	ng	-0.06
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	615954	139.84	ng	-0.02
7) Phenol-d6	7.21	99	875670	143.85	ng	-0.02
23) Nitrobenzene-d5	9.22	82	539981	96.71	ng	-0.02
42) 2,4,6-Tribromophenol	16.17	330	279596	151.84	ng	-0.02
45) 2-Fluorobiphenyl	13.30	172	1135262	98.73	ng	-0.02
79) Terphenyl-d14	20.03	244	1322509	82.96	ng	-0.02
Target Compounds						
50) Dimethylphthalate	14.12	163	138414	10.380	ng	98
75) Fluoranthene	19.48	202	109487	4.857	ng	92
78) Pyrene	19.84	202	98073	4.159	ng	# 93
81) Benzo(a)anthracene	21.68	228	83943	3.800	ng	96
83) Chrysene	21.74	228	73618m	3.468	ng	
88) Benzo(b)fluoranthene	23.93	252	86239	3.704	ng	# 89
90) Benzo(a)pyrene	24.82	252	57875	2.649	ng	# 96

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

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