

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG051119\
 Data File : BG040903.D
 Acq On : 12 May 2019 11:00
 Operator : HP/JU
 Sample : K2768-12
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 P026-SS015-1218-01

Manual Integrations
 APPROVED

Sohil
 5/14/2019 8:01:50 PM

Quant Time: May 13 04:58:27 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG042319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 24 05:35:33 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.13	152	50353	20.00	ng	-0.03
21) Naphthalene-d8	10.96	136	219597	20.00	ng	-0.03
39) Acenaphthene-d10	14.76	164	160149	20.00	ng	-0.03
64) Phenanthrene-d10	17.51	188	413701	20.00	ng	-0.03
76) Chrysene-d12	21.79	240	417731	20.00	ng	-0.03
87) Perylene-d12	25.13	264	449854	20.00	ng	-0.07
System Monitoring Compounds						
5) 2-Fluorophenol	5.67	112	431183	150.95	ng	-0.02
7) Phenol-d6	7.28	99	651386	148.10	ng	-0.02
23) Nitrobenzene-d5	9.31	82	445484	93.74	ng	-0.03
42) 2,4,6-Tribromophenol	16.25	330	350525	159.24	ng	-0.02
45) 2-Fluorobiphenyl	13.39	172	946641	85.37	ng	-0.03
79) Terphenyl-d14	20.10	244	1569261	83.23	ng	-0.03
Target Compounds						
50) Dimethylphthalate	14.20	163	171799	12.821	ng	97
75) Fluoranthene	19.55	202	70802	2.550	ng	94
78) Pyrene	19.91	202	138845	5.303	ng	96
83) Chrysene	21.84	228	60879	2.425	ng	97
88) Benzo(b)fluoranthene	24.06	252	64912m	2.361	ng	

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

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