

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG051119\
 Data File : BG040908.D
 Acq On : 12 May 2019 14:10
 Operator : HP/JU
 Sample : K2766-02
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 P026-SS006-0612-01

Manual Integrations
 APPROVED

Sohil
 5/14/2019 8:01:54 PM

Quant Time: May 13 05:13:58 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	54061	20.00	ng	-0.03
21) Naphthalene-d8	10.96	136	233221	20.00	ng	-0.03
39) Acenaphthene-d10	14.76	164	170788	20.00	ng	-0.03
64) Phenanthrene-d10	17.51	188	427967	20.00	ng	-0.03
76) Chrysene-d12	21.79	240	425678	20.00	ng	-0.03
87) Perylene-d12	25.14	264	467390	20.00	ng	-0.06
System Monitoring Compounds						
5) 2-Fluorophenol	5.67	112	427152	139.28	ng	-0.02
7) Phenol-d6	7.28	99	638632	135.24	ng	-0.02
23) Nitrobenzene-d5	9.31	82	441169	87.41	ng	-0.03
42) 2,4,6-Tribromophenol	16.25	330	344083	146.57	ng	-0.02
45) 2-Fluorobiphenyl	13.39	172	988452	83.58	ng	-0.03
79) Terphenyl-d14	20.10	244	1557869	81.08	ng	-0.03
Target Compounds						
50) Dimethylphthalate	14.20	163	135261	9.466	ng	98
75) Fluoranthene	19.55	202	135164	4.705	ng	97
78) Pyrene	19.92	202	270988	10.157	ng	98
81) Benzo(a)anthracene	21.77	228	112796	4.193	ng	96
83) Chrysene	21.84	228	151398	5.917	ng	99
86) Indeno(1,2,3-cd)pyrene	28.97	276	125463	4.056	ng	96
88) Benzo(b)fluoranthene	24.06	252	184782	6.470	ng	98
89) Benzo(k)fluoranthene	24.12	252	59257m	2.222	ng	
90) Benzo(a)pyrene	24.97	252	126351	4.673	ng	96
92) Benzo(g,h,i)perylene	30.17	276	155816	5.722	ng	98

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

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