

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG050919\
 Data File : BG040847.D
 Acq On : 10 May 2019 12:11
 Operator : HP/JU
 Sample : K2764-15
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 P026-SS001-1824-01

Manual Integrations
 APPROVED

Sohil
 5/13/2019 8:39:11 PM

Quant Time: May 10 15:35:58 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	8.14	152	61091	20.00	ng	-0.02
21) Naphthalene-d8	10.97	136	263367	20.00	ng	-0.02
39) Acenaphthene-d10	14.77	164	188567	20.00	ng	-0.02
64) Phenanthrene-d10	17.51	188	455077	20.00	ng	-0.02
76) Chrysene-d12	21.80	240	436963	20.00	ng	-0.03
87) Perylene-d12	25.15	264	468434	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.67	112	496655	143.31	ng	0.00
7) Phenol-d6	7.28	99	750336	140.62	ng	-0.01
23) Nitrobenzene-d5	9.32	82	522951	91.75	ng	-0.02
42) 2,4,6-Tribromophenol	16.26	330	365869	141.16	ng	-0.02
45) 2-Fluorobiphenyl	13.40	172	978929	74.97	ng	-0.02
79) Terphenyl-d14	20.11	244	1520777	77.10	ng	-0.02
Target Compounds						
50) Dimethylphthalate	14.21	163	98775	6.261	ng	99
71) Phenanthrene	17.56	178	122202	4.985	ng	96
75) Fluoranthene	19.56	202	299107	9.792	ng	95
78) Pyrene	19.92	202	562836	20.551	ng	99
81) Benzo(a)anthracene	21.78	228	194434	7.041	ng	99
83) Chrysene	21.84	228	241148	9.182	ng	98
86) Indeno(1,2,3-cd)pyrene	28.99	276	111678	3.517	ng	# 87
88) Benzo(b)fluoranthene	24.09	252	232621	8.126	ng	96
89) Benzo(k)fluoranthene	24.15	252	67306m	2.518	ng	
90) Benzo(a)pyrene	24.99	252	142038m	5.242	ng	
92) Benzo(g,h,i)perylene	30.22	276	129744	4.754	ng	97

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

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