

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG062019\
 Data File : BG041369.D
 Acq On : 20 Jun 2019 22:48
 Operator : HP/JU
 Sample : K3419-03
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TP-B

Quant Time: Jun 21 07:23:25 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG061719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 17 14:58:02 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.08	152	44522	20.00	ng	0.00
21) Naphthalene-d8	10.90	136	172697	20.00	ng	0.00
39) Acenaphthene-d10	14.72	164	117326	20.00	ng	0.00
64) Phenanthrene-d10	17.47	188	254792	20.00	ng	0.00
76) Chrysene-d12	21.74	240	234023	20.00	ng	0.00
87) Perylene-d12	25.07	264	161894	20.00	ng	0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.62	112	321939	132.86	ng	0.00
7) Phenol-d6	7.24	99	482641	137.42	ng	0.00
23) Nitrobenzene-d5	9.26	82	356130	86.73	ng	0.00
42) 2,4,6-Tribromophenol	16.21	330	230040	127.71	ng	0.00
45) 2-Fluorobiphenyl	13.34	172	718934	83.28	ng	0.00
79) Terphenyl-d14	20.06	244	814856	69.01	ng	0.00
Target Compounds						
50) Dimethylphthalate	14.16	163	81433	7.801	ng	98
78) Pyrene	19.87	202	58371	3.698	ng	98
86) Indeno(1,2,3-cd)pyrene	28.89	276	54134	2.994	ng	# 89
88) Benzo(b)fluoranthene	24.02	252	58864m	5.867	ng	
90) Benzo(a)pyrene	24.92	252	48936	5.164	ng	# 90
92) Benzo(g,h,i)perylene	30.09	276	80177	8.505	ng	98

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

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