

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG050319\
 Data File : BG040733.D
 Acq On : 3 May 2019 20:23
 Operator : JU/SJ
 Sample : K2667-03
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SP-1-B

Quant Time: May 04 07:16:01 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	8.15	152	42745	20.00	ng	-0.01
21) Naphthalene-d8	10.97	136	185912	20.00	ng	-0.02
39) Acenaphthene-d10	14.77	164	142606	20.00	ng	-0.02
64) Phenanthrene-d10	17.52	188	389601	20.00	ng	-0.02
76) Chrysene-d12	21.80	240	377430	20.00	ng	-0.02
87) Perylene-d12	25.16	264	404553	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.68	112	383555	158.17	ng	0.00
7) Phenol-d6	7.28	99	586629	157.12	ng	-0.01
23) Nitrobenzene-d5	9.32	82	406228	100.96	ng	-0.02
42) 2,4,6-Tribromophenol	16.26	330	331672	169.21	ng	-0.01
45) 2-Fluorobiphenyl	13.40	172	876813	88.80	ng	-0.02
79) Terphenyl-d14	20.11	244	1528501	89.72	ng	-0.02
Target Compounds						
50) Dimethylphthalate	14.22	163	124989	10.476	ng	Qvalue # 95
71) Phenanthrene	17.56	178	108894	5.189	ng	97
75) Fluoranthene	19.56	202	139874	5.349	ng	97
78) Pyrene	19.92	202	235803	9.968	ng	97
81) Benzo(a)anthracene	21.78	228	100527	4.214	ng	98
83) Chrysene	21.85	228	107642	4.745	ng	95
88) Benzo(b)fluoranthene	24.09	252	79975	3.235	ng	99
90) Benzo(a)pyrene	24.99	252	80745	3.450	ng	98

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

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