

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG021720\
 Data File : BG044459.D
 Acq On : 17 Feb 2020 22:55
 Operator : CG/JU
 Sample : L1547-15
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MARCY-GREEN-BH-03-C

Quant Time: Feb 18 06:35:08 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG013020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 10 16:11:50 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.89	152	76239	20.00	ng	-0.02
21) Naphthalene-d8	10.69	136	307241	20.00	ng	-0.01
39) Acenaphthene-d10	14.53	164	194960	20.00	ng	-0.02
64) Phenanthrene-d10	17.27	188	436521	20.00	ng	-0.02
76) Chrysene-d12	21.51	240	437602	20.00	ng	-0.02
87) Perylene-d12	24.59	264	469870	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	440836	102.05	ng	0.00
7) Phenol-d6	7.06	99	567444	97.82	ng	-0.01
23) Nitrobenzene-d5	9.04	82	344934	63.38	ng	-0.02
42) 2,4,6-Tribromophenol	16.02	330	222808	97.35	ng	-0.01
45) 2-Fluorobiphenyl	13.15	172	808208	69.42	ng	-0.01
79) Terphenyl-d14	19.89	244	1267540	59.58	ng	-0.01
Target Compounds						Qvalue
50) Dimethylphthalate	13.99	163	32855	2.34	ng	100
71) Phenanthrene	17.31	178	112760	5.33	ng	95
75) Fluoranthene	19.33	202	140604	5.75	ng	95
78) Pyrene	19.69	202	137254	4.88	ng	94
81) Benzo(a)anthracene	21.50	228	67316	2.50	ng	91
83) Chrysene	21.56	228	67489	2.59	ng	93
88) Benzo(b)fluoranthene	23.62	252	78523	2.90	ng	97
90) Benzo(a)pyrene	24.44	252	55302	2.16	ng	99

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

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