

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021020\
 Data File : BG044377.D
 Acq On : 11 Feb 2020 6:29
 Operator : CG/JU
 Sample : L1449-01
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 S-1-BY-TANK-WEST

Quant Time: Feb 11 08:05:32 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	7.90	152	68187	20.00	ng	0.00
21) Naphthalene-d8	10.70	136	299103	20.00	ng	0.00
39) Acenaphthene-d10	14.54	164	208062	20.00	ng	0.00
64) Phenanthrene-d10	17.28	188	492985	20.00	ng	0.00
76) Chrysene-d12	21.53	240	462278	20.00	ng	0.00
87) Perylene-d12	24.61	264	488417	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.48	112	480805	124.44	ng	0.00
7) Phenol-d6	7.07	99	644184	124.16	ng	0.00
23) Nitrobenzene-d5	9.06	82	388678	73.36	ng	0.00
42) 2,4,6-Tribromophenol	16.03	330	287788	117.82	ng	0.00
45) 2-Fluorobiphenyl	13.17	172	951704	76.60	ng	0.00
79) Terphenyl-d14	19.90	244	1542877	68.65	ng	0.00
Target Compounds						Qvalue
50) Dimethylphthalate	13.99	163	59780	3.997	ng	98
75) Fluoranthene	19.34	202	154181	5.583	ng	96
78) Pyrene	19.70	202	130562	4.398	ng	94
81) Benzo(a)anthracene	21.51	228	95390	3.348	ng	97
83) Chrysene	21.57	228	83606	3.036	ng	97
88) Benzo(b)fluoranthene	23.64	252	119476	4.245	ng	96
90) Benzo(a)pyrene	24.46	252	78635	2.955	ng	99

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

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