

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG100119\
 Data File : BG042979.D
 Acq On : 1 Oct 2019 22:50
 Operator : JU
 Sample : K4946-10
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 GP-5 (1-3)

Quant Time: Oct 02 01:58:07 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG092519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 25 15:35:52 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.07	152	75183	20.00	ng	0.00
21) Naphthalene-d8	10.87	136	279571	20.00	ng	0.00
39) Acenaphthene-d10	14.69	164	199475	20.00	ng	0.00
64) Phenanthrene-d10	17.43	188	466015	20.00	ng	0.00
76) Chrysene-d12	21.70	240	496669	20.00	ng	-0.01
87) Perylene-d12	24.92	264	593763	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.65	112	422376	100.26	ng	0.00
7) Phenol-d6	7.24	99	632388	104.24	ng	0.00
23) Nitrobenzene-d5	9.23	82	405539	70.51	ng	0.00
42) 2,4,6-Tribromophenol	16.18	330	365236	106.48	ng	0.00
45) 2-Fluorobiphenyl	13.32	172	995120	72.32	ng	0.00
79) Terphenyl-d14	20.04	244	1702366	73.04	ng	0.00
Target Compounds						
10) Phenol	7.26	94	23028	4.137	ng	89
50) Dimethylphthalate	14.14	163	169700	11.354	ng	98
78) Pyrene	19.84	202	60128	2.028	ng	95

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

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