

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101220\
 Data File : BG046866.D
 Acq On : 13 Oct 2020 7:10
 Operator : CG/JU
 Sample : L4317-02
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 11991

Quant Time: Oct 13 08:16:01 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG092220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 07 11:38:17 2020
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.103	152	61623	20.00	ng	0.00
21) Naphthalene-d8	10.917	136	236166	20.00	ng	0.00
39) Acenaphthene-d10	14.725	164	171812	20.00	ng	-0.01
64) Phenanthrene-d10	17.468	188	459897	20.00	ng	#-0.01
76) Chrysene-d12	21.740	240	513366	20.00	ng	#-0.02
86) Perylene-d12	25.012	264	503116	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.670	112	517260	134.47	ng	0.00
7) Phenol-d6	7.257	99	657943	121.26	ng	0.00
23) Nitrobenzene-d5	9.266	82	487691	109.86	ng	0.00
42) 2,4,6-Tribromophenol	16.211	330	403851	178.26	ng	0.00
45) 2-Fluorobiphenyl	13.350	172	1128742	93.01	ng	-0.01
79) Terphenyl-d14	20.077	244	2180679	85.13	ng	0.00

Target Compounds	Qvalue

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

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