

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE100819\
 Data File : BE100611.D
 Acq On : 8 Oct 2019 16:37
 Operator : JU
 Sample : K5109-04
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 MW-4

Quant Time: Oct 09 04:11:52 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE100319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 09 01:42:20 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.82	152	4025	0.40	ng	0.00
7) Naphthalene-d8	10.60	136	19261	0.40	ng	0.00
13) Acenaphthene-d10	14.46	164	11627	0.40	ng	0.00
19) Phenanthrene-d10	17.17	188	26026	0.40	ng	0.00
27) Chrysene-d12	21.33	240	32422	0.40	ng	0.00
34) Perylene-d12	23.79	264	36994	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.40	112	3145	0.29	ng	0.00
5) Phenol-d6	6.98	99	3232	0.24	ng	-0.01
8) Nitrobenzene-d5	8.97	82	4379	0.36	ng	0.00
11) 2-Methylnaphthalene-d10	12.20	152	10447	0.36	ng	-0.01
14) 2,4,6-Tribromophenol	15.93	330	1285	0.48	ng	0.00
15) 2-Fluorobiphenyl	13.07	172	13954	0.33	ng	-0.01
25) Fluoranthene-d10	19.20	212	134714	0.42	ng	0.00
29) Terphenyl-d14	19.79	244	30827	0.43	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.28	88	56699	8.364	ng	# 87
23) Phenanthrene	17.21	178	1531	0.021	ng	# 86
32) Bis(2-ethylhexyl)phthalate	21.26	149	39520	0.228	ng	100

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

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