

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE100819\
 Data File : BE100608.D
 Acq On : 8 Oct 2019 14:47
 Operator : JU
 Sample : K5109-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 MW-1

Quant Time: Oct 09 04:08:06 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	4145	0.40	ng	0.00
7) Naphthalene-d8	10.60	136	20184	0.40	ng	0.00
13) Acenaphthene-d10	14.46	164	12530	0.40	ng	0.00
19) Phenanthrene-d10	17.17	188	28161	0.40	ng	0.00
27) Chrysene-d12	21.33	240	33230	0.40	ng	0.00
34) Perylene-d12	23.78	264	37988	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.40	112	3117	0.28	ng	0.00
5) Phenol-d6	6.98	99	3114	0.22	ng	-0.01
8) Nitrobenzene-d5	8.96	82	4206	0.33	ng	0.00
11) 2-Methylnaphthalene-d10	12.19	152	10753	0.36	ng	-0.01
14) 2,4,6-Tribromophenol	15.93	330	1578	0.55	ng	0.00
15) 2-Fluorobiphenyl	13.07	172	16603	0.37	ng	-0.01
25) Fluoranthene-d10	19.20	212	131877	0.38	ng	0.00
29) Terphenyl-d14	19.79	244	31306	0.43	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.29	88	3822399	547.514	ng	Qvalue # 86
22) Pentachlorophenol	16.83	266	108	0.358	ng	86
32) Bis(2-ethylhexyl)phthalate	21.26	149	290128	1.635	ng	100

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

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