

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE121119\
 Data File : BE100862.D
 Acq On : 12 Dec 2019 2:05
 Operator : JU
 Sample : K6185-01DL 2X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 RE103D1-20191204DL

Quant Time: Dec 12 05:42:35 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE120419.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 04 17:18:11 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.76	152	4963	0.40	ng	-0.01
7) Naphthalene-d8	10.54	136	23567	0.40	ng	0.00
13) Acenaphthene-d10	14.39	164	12899	0.40	ng	-0.01
19) Phenanthrene-d10	17.12	188	29190	0.40	ng	-0.01
27) Chrysene-d12	21.29	240	21429	0.40	ng	-0.01
34) Perylene-d12	23.74	264	24343	0.40	ng	-0.02
System Monitoring Compounds						
4) 2-Fluorophenol	5.36	112	1107	0.09	ng	-0.01
5) Phenol-d6	6.93	99	1085	0.06	ng	0.00
8) Nitrobenzene-d5	8.90	82	2666	0.24	ng	0.00
11) 2-Methylnaphthalene-d10	12.14	152	7082	0.21	ng	0.00
14) 2,4,6-Tribromophenol	15.88	330	352	0.25	ng	0.00
15) 2-Fluorobiphenyl	13.02	172	10275	0.20	ng	-0.01
25) Fluoranthene-d10	19.15	212	57865	0.17	ng	-0.01
29) Terphenyl-d14	19.75	244	10937	0.21	ng	-0.01
Target Compounds						
2) 1,4-Dioxane	3.34	88	26444	2.922	ng	Qvalue # 89

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

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