

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE121319\
 Data File : BE100886.D
 Acq On : 13 Dec 2019 14:36
 Operator : JU
 Sample : K6185-18DL 2X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 DUP03-20191205DL

Quant Time: Dec 13 15:37:29 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE120419.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 13 10:42:38 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.76	152	4954	0.40	ng	0.00
7) Naphthalene-d8	10.54	136	21830	0.40	ng	0.00
13) Acenaphthene-d10	14.39	164	12042	0.40	ng	0.00
19) Phenanthrene-d10	17.12	188	30320	0.40	ng	0.00
27) Chrysene-d12	21.29	240	25906	0.40	ng	-0.01
34) Perylene-d12	23.74	264	29920	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.37	112	1086	0.09	ng	0.00
5) Phenol-d6	6.93	99	985	0.06	ng	0.00
8) Nitrobenzene-d5	8.90	82	2744	0.26	ng	0.00
11) 2-Methylnaphthalene-d10	12.12	152	6691	0.21	ng	-0.01
14) 2,4,6-Tribromophenol	15.86	330	378	0.28	ng	-0.01
15) 2-Fluorobiphenyl	13.02	172	9347	0.20	ng	0.00
25) Fluoranthene-d10	19.15	212	68791	0.19	ng	0.00
29) Terphenyl-d14	19.75	244	13095	0.21	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.34	88	35276	3.906	ng	Qvalue 90

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

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