

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE020419\
 Data File : BE098710.D
 Acq On : 4 Feb 2019 16:49
 Operator : JU/SJ
 Sample : PB116795BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PB116795BL

Quant Time: Feb 04 23:58:18 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE012419.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 24 14:44:44 2019
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.815	152	878	0.40	ng	0.00
7) Naphthalene-d8	10.614	136	3513	0.40	ng	# 0.01
13) Acenaphthene-d10	14.485	164	2324	0.40	ng	0.01
19) Phenanthrene-d10	17.229	188	5686	0.40	ng	# 0.00
27) Chrysene-d12	21.414	240	6782	0.40	ng	# 0.00
34) Perylene-d12	23.923	264	6706	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.408	112	701	0.27	ng	0.00
5) Phenol-d6	7.020	99	870	0.23	ng	0.02
8) Nitrobenzene-d5	9.030	82	624	0.22	ng	0.05
11) 2-Methylnaphthalene-d10	12.224	152	1689	0.25	ng	0.01
14) 2,4,6-Tribromophenol	15.997	330	234	0.23	ng	0.02
15) 2-Fluorobiphenyl	13.117	172	2273	0.23	ng	0.01
25) Fluoranthene-d10	19.264	212	23360	0.29	ng	0.00
29) Terphenyl-d14	19.861	244	3957	0.23	ng	0.00

Target Compounds Qvalue

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

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