

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE040821\
 Data File : BE101240.D
 Acq On : 8 Apr 2021 10:04
 Operator : CG/JU
 Sample : PB135597BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PB135597BL

Quant Time: Apr 08 11:35:13 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_E\Methods\8270-SIM-BE040621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 07 10:22:45 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.830	152	4282	0.40	ng	0.00
7) Naphthalene-d8	10.614	136	17256	0.40	ng	0.00
13) Acenaphthene-d10	14.444	164	9916	0.40	ng	0.00
19) Phenanthrene-d10	17.196	188	21135	0.40	ng	#-0.01
29) Chrysene-d12	21.354	240	26332	0.40	ng	# 0.00
36) Perylene-d12	23.833	264	25302	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.423	112	5234	0.52	ng	0.00
5) Phenol-d6	6.989	99	7129	0.45	ng	0.00
8) Nitrobenzene-d5	8.977	82	6007	0.47	ng	0.00
11) 2-Methylnaphthalene-d10	12.196	152	11675	0.29	ng	0.00
14) 2,4,6-Tribromophenol	15.956	330	879	0.33	ng	0.00
15) 2-Fluorobiphenyl	13.075	172	16487	0.46	ng	0.00
27) Fluoranthene-d10	19.215	212	107832	0.29	ng	0.00
31) Terphenyl-d14	19.812	244	26386	0.37	ng	0.00

Target Compounds Qvalue

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

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