

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE040721\
 Data File : BE101222.D
 Acq On : 7 Apr 2021 11:11
 Operator : CG/JU
 Sample : PB135530BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PB135530BL

Quant Time: Apr 07 11:54:46 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	7.839	152	3984	0.40	ng	0.00
7) Naphthalene-d8	10.615	136	16068	0.40	ng	0.00
13) Acenaphthene-d10	14.443	164	8792	0.40	ng	0.00
19) Phenanthrene-d10	17.195	188	21614	0.40	ng	#-0.01
29) Chrysene-d12	21.354	240	22048	0.40	ng	# 0.00
36) Perylene-d12	23.831	264	20795	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.421	112	4256	0.45	ng	0.00
5) Phenol-d6	6.987	99	5684	0.39	ng	0.00
8) Nitrobenzene-d5	8.978	82	4629	0.39	ng	0.00
11) 2-Methylnaphthalene-d10	12.196	152	10518	0.28	ng	0.00
14) 2,4,6-Tribromophenol	15.956	330	572	0.24	ng	0.00
15) 2-Fluorobiphenyl	13.075	172	15050	0.48	ng	0.00
27) Fluoranthene-d10	19.215	212	106079	0.28	ng	0.00
31) Terphenyl-d14	19.812	244	25197	0.43	ng	0.00

Target Compounds Qvalue

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

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