

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE040121\
 Data File : BE101173.D
 Acq On : 1 Apr 2021 12:24
 Operator : CG/JU
 Sample : PB135471BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PB135471BL

Quant Time: Apr 01 13:07:50 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_E\Methods\8270-SIM-BE032421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 01 12:49:55 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.842	152	4760	0.40	ng	0.00
7) Naphthalene-d8	10.627	136	19047	0.40	ng	# 0.00
13) Acenaphthene-d10	14.457	164	9945	0.40	ng	0.00
19) Phenanthrene-d10	17.212	188	22276	0.40	ng	# 0.00
29) Chrysene-d12	21.366	240	26631	0.40	ng	# 0.00
36) Perylene-d12	23.849	264	33861	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.435	112	4252	0.38	ng	0.00
5) Phenol-d6	7.001	99	5627	0.32	ng	0.00
8) Nitrobenzene-d5	8.991	82	4166	0.37	ng	0.01
11) 2-Methylnaphthalene-d10	12.209	152	9414	0.30	ng	0.00
14) 2,4,6-Tribromophenol	15.957	330	675	0.38	ng	0.00
15) 2-Fluorobiphenyl	13.088	172	12749	0.34	ng	0.00
27) Fluoranthene-d10	19.225	212	83297	0.30	ng	0.00
31) Terphenyl-d14	19.823	244	19320	0.32	ng	0.00

Target Compounds Qvalue

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

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