

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE040921\
 Data File : BE101267.D
 Acq On : 9 Apr 2021 10:47
 Operator : CG/JU
 Sample : PB135423BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PB135423BL

Quant Time: Apr 09 11:31:18 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	2602	0.40	ng	0.00
7) Naphthalene-d8	10.614	136	11190	0.40	ng	0.00
13) Acenaphthene-d10	14.442	164	7676	0.40	ng	0.00
19) Phenanthrene-d10	17.196	188	24488	0.40	ng	#-0.01
29) Chrysene-d12	21.355	240	21499	0.40	ng	0.00
36) Perylene-d12	23.837	264	22311	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.423	112	3297	0.54	ng	0.00
5) Phenol-d6	7.001	99	4641	0.48	ng	0.01
8) Nitrobenzene-d5	8.977	82	4022	0.48	ng	0.00
11) 2-Methylnaphthalene-d10	12.195	152	7832	0.30	ng	0.00
14) 2,4,6-Tribromophenol	15.956	330	753	0.37	ng	0.00
15) 2-Fluorobiphenyl	13.073	172	11081	0.40	ng	0.00
27) Fluoranthene-d10	19.215	212	140510	0.33	ng	0.00
31) Terphenyl-d14	19.813	244	30480	0.53	ng	0.00

Target Compounds Qvalue

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

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