

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE040721\
 Data File : BE101236.D
 Acq On : 7 Apr 2021 19:50
 Operator : CG/JU
 Sample : M1951-04
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SW-2

Quant Time: Apr 08 03:23:54 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	7.838	152	3502	0.40	ng	0.00
7) Naphthalene-d8	10.615	136	16641	0.40	ng	0.00
13) Acenaphthene-d10	14.443	164	13021	0.40	ng	0.00
19) Phenanthrene-d10	17.197	188	40431	0.40	ng	#-0.01
29) Chrysene-d12	21.354	240	39156	0.40	ng	# 0.00
36) Perylene-d12	23.834	264	30198	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.420	112	1970	0.24	ng	0.00
5) Phenol-d6	6.998	99	2140	0.17	ng	0.00
8) Nitrobenzene-d5	8.978	82	4529	0.37	ng	0.00
11) 2-Methylnaphthalene-d10	12.196	152	9662	0.25	ng	0.00
14) 2,4,6-Tribromophenol	15.957	330	2169	0.62	ng	0.00
15) 2-Fluorobiphenyl	13.074	172	14980	0.32	ng	0.00
27) Fluoranthene-d10	19.214	212	231261	0.33	ng	0.00
31) Terphenyl-d14	19.817	244	55557	0.53	ng	0.00
Target Compounds						Qvalue
9) Naphthalene	10.667	128	5057	0.03	ng	# 95
25) Phenanthrene	17.242	178	4911	0.04	ng	# 88
28) Fluoranthene	19.243	202	6910	0.04	ng	# 93
30) Pyrene	19.605	202	6264	0.04	ng	# 92
37) Benzo(b)fluoranthene	23.075	252	2368	0.02	ng	# 48

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

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