

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE040821\
 Data File : BE101243.D
 Acq On : 8 Apr 2021 11:57
 Operator : CG/JU
 Sample : M1951-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SW-1

Quant Time: Apr 08 13:02:24 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	7.829	152	4027	0.40	ng	# 0.00
7) Naphthalene-d8	10.615	136	19358	0.40	ng	0.00
13) Acenaphthene-d10	14.443	164	14868	0.40	ng	0.00
19) Phenanthrene-d10	17.197	188	42614	0.40	ng	#-0.01
29) Chrysene-d12	21.355	240	32419	0.40	ng	# 0.00
36) Perylene-d12	23.831	264	31361	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.422	112	1877	0.20	ng	0.00
5) Phenol-d6	6.988	99	2181	0.15	ng	0.00
8) Nitrobenzene-d5	8.978	82	4867	0.34	ng	0.00
11) 2-Methylnaphthalene-d10	12.196	152	10540	0.23	ng	0.00
14) 2,4,6-Tribromophenol	15.942	330	2143	0.54	ng	-0.01
15) 2-Fluorobiphenyl	13.074	172	16326	0.31	ng	0.00
27) Fluoranthene-d10	19.214	212	189697	0.26	ng	0.00
31) Terphenyl-d14	19.812	244	49148	0.57	ng	0.00
Target Compounds						
9) Naphthalene	10.667	128	6055	0.03	ng	# 95
25) Phenanthrene	17.242	178	5689	0.05	ng	# 91
28) Fluoranthene	19.243	202	6753	0.04	ng	# 89
30) Pyrene	19.605	202	5286	0.05	ng	# 92
37) Benzo(b)fluoranthene	23.075	252	2319	0.02	ng	# 51

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

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