

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE040821\
 Data File : BE101247.D
 Acq On : 8 Apr 2021 14:33
 Operator : CG/JU
 Sample : SP5532
 Misc : 8270 SIM SURROGATE
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SP5532

Quant Time: Apr 08 15:23:58 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	5980	0.40	ng	0.00
7) Naphthalene-d8	10.614	136	25008	0.40	ng	0.00
13) Acenaphthene-d10	14.443	164	13858	0.40	ng	0.00
19) Phenanthrene-d10	17.196	188	24692	0.40	ng	#-0.01
29) Chrysene-d12	21.354	240	26702	0.40	ng	# 0.00
36) Perylene-d12	23.830	264	29316	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.421	112	6920	0.49	ng	0.00
5) Phenol-d6	6.987	99	9244	0.42	ng	0.00
8) Nitrobenzene-d5	8.977	82	8419	0.45	ng	0.00
11) 2-Methylnaphthalene-d10	12.195	152	13663	0.23	ng	0.00
14) 2,4,6-Tribromophenol	15.956	330	996	0.27	ng	0.00
15) 2-Fluorobiphenyl	13.074	172	19600	0.39	ng	0.00
27) Fluoranthene-d10	19.213	212	74351	0.17	ng	0.00
31) Terphenyl-d14	19.811	244	18000	0.25	ng	0.00

Target Compounds	Qvalue

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

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