

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM031121\
 Data File : BM029090.D
 Acq On : 11 Mar 2021 12:01
 Operator : CG/JU
 Sample : M1603-06
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 33-NORTH

Quant Time: Mar 11 12:32:51 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM022321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 09 10:08:02 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.657	152	8289	20.00	ng	-0.01
21) Naphthalene-d8	10.451	136	31263	20.00	ng	#-0.01
39) Acenaphthene-d10	14.327	164	18087	20.00	ng	-0.01
64) Phenanthrene-d10	17.080	188	39271	20.00	ng	0.00
76) Chrysene-d12	21.250	240	40616	20.00	ng	#-0.01
86) Perylene-d12	23.409	264	40062	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.234	112	61875	123.73	ng	-0.01
7) Phenol-d6	6.857	99	65889	99.75	ng	-0.01
23) Nitrobenzene-d5	8.839	82	49957	86.21	ng	0.00
42) 2,4,6-Tribromophenol	15.833	330	30305	141.37	ng	0.00
45) 2-Fluorobiphenyl	12.939	172	110166	81.09	ng	-0.01
79) Terphenyl-d14	19.703	244	200069	90.95	ng	0.00
Target Compounds						
10) Phenol	6.887	94	1335	2.03	ng	Qvalue 94

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

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