

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM031021\
 Data File : BM029076.D
 Acq On : 10 Mar 2021 18:59
 Operator : CG/JU
 Sample : M1603-05
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 32-NORTH

Quant Time: Mar 10 23:40:14 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	7.663	152	7529	20.00	ng	0.00
21) Naphthalene-d8	10.457	136	29854	20.00	ng	# 0.00
39) Acenaphthene-d10	14.333	164	17446	20.00	ng	0.00
64) Phenanthrene-d10	17.080	188	37284	20.00	ng	0.00
76) Chrysene-d12	21.256	240	40410	20.00	ng	# 0.00
86) Perylene-d12	23.415	264	42998	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.240	112	63193	139.12	ng	0.00
7) Phenol-d6	6.863	99	72590	120.99	ng	0.00
23) Nitrobenzene-d5	8.840	82	49478	89.42	ng	0.00
42) 2,4,6-Tribromophenol	15.833	330	31560	152.63	ng	0.00
45) 2-Fluorobiphenyl	12.945	172	114353	87.26	ng	0.00
79) Terphenyl-d14	19.709	244	208091	95.08	ng	0.00
Target Compounds						
10) Phenol	6.893	94	1300	2.18	ng	Qvalue 83
32) Benzoic acid	9.798	122	680	2.22	ng	# 78

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

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