

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030921\
 Data File : BM029052.D
 Acq On : 09 Mar 2021 18:52
 Operator : CG/JU
 Sample : M1576-09
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 IDW-GW804-SO-030421

Quant Time: Mar 09 23:39:24 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	7.669	152	8894	20.00	ng	0.00
21) Naphthalene-d8	10.457	136	32923	20.00	ng	# 0.00
39) Acenaphthene-d10	14.333	164	18453	20.00	ng	0.00
64) Phenanthrene-d10	17.080	188	39307	20.00	ng	0.00
76) Chrysene-d12	21.256	240	41040	20.00	ng	# 0.00
86) Perylene-d12	23.415	264	42332	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.246	112	77771	144.94	ng	0.00
7) Phenol-d6	6.869	99	82725	116.72	ng	0.00
23) Nitrobenzene-d5	8.840	82	60149	98.57	ng	0.00
42) 2,4,6-Tribromophenol	15.833	330	34388	157.24	ng	0.00
45) 2-Fluorobiphenyl	12.945	172	131802	95.09	ng	0.00
79) Terphenyl-d14	19.709	244	221550	99.68	ng	0.00
Target Compounds						
10) Phenol	6.893	94	1642	2.33	ng	Qvalue 77

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

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