

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061221\
 Data File : BM030426.D
 Acq On : 12 Jun 2021 22:00
 Operator : CG/JU
 Sample : M2676-11
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 HEAT-C

Quant Time: Jun 14 01:01:16 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM060921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 11 14:10:50 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.834	152	164480	20.000	ng	0.00
21) Naphthalene-d8	10.622	136	608759	20.000	ng	# 0.00
39) Acenaphthene-d10	14.457	164	354435	20.000	ng	0.00
64) Phenanthrene-d10	17.192	188	733926	20.000	ng	0.00
76) Chrysene-d12	21.350	240	814416	20.000	ng	0.00
86) Perylene-d12	23.627	264	884993	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.416	112	803656	79.430	ng	0.00
7) Phenol-d6	6.993	99	1289840	88.063	ng	0.00
23) Nitrobenzene-d5	8.987	82	896024	71.707	ng	0.00
42) 2,4,6-Tribromophenol	15.939	330	275472	61.817	ng	0.00
45) 2-Fluorobiphenyl	13.080	172	2094787	77.723	ng	0.00
79) Terphenyl-d14	19.803	244	3536063	76.262	ng	0.00

Target Compounds	Qvalue

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

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