

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061121\
 Data File : BM030405.D
 Acq On : 11 Jun 2021 15:26
 Operator : CG/JU
 Sample : M2612-36
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 18-B12(66-70)

Quant Time: Jun 11 16:08:15 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	7.840	152	188118	20.000	ng	0.00
21) Naphthalene-d8	10.622	136	669722	20.000	ng	# 0.00
39) Acenaphthene-d10	14.457	164	377616	20.000	ng	0.00
64) Phenanthrene-d10	17.198	188	724192	20.000	ng	0.00
76) Chrysene-d12	21.356	240	695394	20.000	ng	# 0.00
86) Perylene-d12	23.633	264	776890	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.422	112	1487030	128.505	ng	0.00
7) Phenol-d6	6.998	99	1597215	95.346	ng	0.00
23) Nitrobenzene-d5	8.992	82	1212453	88.198	ng	0.00
42) 2,4,6-Tribromophenol	15.945	330	604688	127.365	ng	0.00
45) 2-Fluorobiphenyl	13.086	172	2568909	89.463	ng	0.00
79) Terphenyl-d14	19.809	244	3386364	85.533	ng	0.00

Target Compounds	Qvalue

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

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