

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061121\
 Data File : BM030402.D
 Acq On : 11 Jun 2021 13:36
 Operator : CG/JU
 Sample : M2612-24
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 18-B12(18-22)

Quant Time: Jun 11 14:12:38 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	166949	20.000	ng	0.00
21) Naphthalene-d8	10.622	136	596251	20.000	ng	# 0.00
39) Acenaphthene-d10	14.457	164	357823	20.000	ng	0.00
64) Phenanthrene-d10	17.192	188	717516	20.000	ng	# 0.00
76) Chrysene-d12	21.357	240	772844	20.000	ng	0.00
86) Perylene-d12	23.639	264	879255	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.422	112	1405241	136.835	ng	0.00
7) Phenol-d6	7.005	99	1557714	104.779	ng	0.00
23) Nitrobenzene-d5	8.993	82	1152157	94.139	ng	0.00
42) 2,4,6-Tribromophenol	15.945	330	654681	145.523	ng	0.00
45) 2-Fluorobiphenyl	13.081	172	2406984	88.461	ng	0.00
79) Terphenyl-d14	19.810	244	3828594	87.012	ng	0.00
Target Compounds						
32) Benzoic acid	9.857	122	5915	5.740	ng	Qvalue 94

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

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