

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063021\
 Data File : BP006114.D
 Acq On : 30 Jun 2021 18:48
 Operator : CG/JU
 Sample : M2875-19
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MW-12

Quant Time: Jul 01 03:07:17 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 30 16:39:37 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.893	152	93977	20.000	ng	0.00
21) Naphthalene-d8	10.693	136	346827	20.000	ng	0.00
39) Acenaphthene-d10	14.522	164	196632	20.000	ng	0.00
64) Phenanthrene-d10	17.275	188	354424	20.000	ng	0.00
76) Chrysene-d12	21.351	240	323982	20.000	ng	0.00
86) Perylene-d12	23.751	264	392811	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.469	112	79928	13.647	ng	0.00
7) Phenol-d6	7.081	99	56687	7.468	ng	0.00
23) Nitrobenzene-d5	9.058	82	306931	43.387	ng	0.00
42) 2,4,6-Tribromophenol	16.016	330	117112	34.706	ng	0.00
45) 2-Fluorobiphenyl	13.151	172	854853	57.088	ng	0.00
79) Terphenyl-d14	19.822	244	1085296	62.726	ng	0.00
Target Compounds						
						Qvalue
27) 2,4-Dimethylphenol	9.916	122	53814	10.306	ng	# 87
31) Naphthalene	10.746	128	1008948	50.278	ng	100
37) 2-Methylnaphthalene	12.351	142	154948	10.845	ng	99
38) 1-Methylnaphthalene	12.569	142	143287	10.647	ng	98

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

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