

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063021\
 Data File : BP006116.D
 Acq On : 30 Jun 2021 19:57
 Operator : CG/JU
 Sample : M2875-17
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MW-2

Manual Integrations
 APPROVED

mohammad
 7/1/2021 2:52:46 PM

Quant Time: Jul 01 03:07:46 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	7.887	152	71171	20.000	ng	0.00
21) Naphthalene-d8	10.693	136	253915	20.000	ng	# 0.00
39) Acenaphthene-d10	14.522	164	143010	20.000	ng	0.00
64) Phenanthrene-d10	17.269	188	286789	20.000	ng	# 0.00
76) Chrysene-d12	21.351	240	355926	20.000	ng	0.00
86) Perylene-d12	23.751	264	437322	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.469	112	262674	59.223	ng	0.00
7) Phenol-d6	7.075	99	175349	30.502	ng	0.00
23) Nitrobenzene-d5	9.063	82	460167	88.851	ng	0.00
42) 2,4,6-Tribromophenol	16.016	330	339040	138.149	ng	0.00
45) 2-Fluorobiphenyl	13.151	172	1002639	92.063	ng	0.00
79) Terphenyl-d14	19.822	244	1593734	83.845	ng	0.00
Target Compounds						
10) Phenol	7.105	94	23026	4.009	ng	# 1
27) 2,4-Dimethylphenol	9.916	122	84311	22.055	ng	# 93
31) Naphthalene	10.746	128	1101807	74.997	ng	100
37) 2-Methylnaphthalene	12.351	142	163799	15.660	ng	99
38) 1-Methylnaphthalene	12.569	142	166806m	16.930	ng	

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

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