

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041921\
 Data File : BM029545.D
 Acq On : 19 Apr 2021 17:55
 Operator : CG/JU
 Sample : M2074-06
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 VNPH2-B-3D

Quant Time: Apr 20 04:15:42 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM040121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 16 11:56:50 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.663	152	111995	20.00	ng	-0.01
21) Naphthalene-d8	10.445	136	466518	20.00	ng	-0.01
39) Acenaphthene-d10	14.304	164	276740	20.00	ng	-0.01
64) Phenanthrene-d10	17.051	188	507939	20.00	ng	0.00
76) Chrysene-d12	21.227	240	359698	20.00	ng	# 0.00
86) Perylene-d12	23.433	264	356843	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.263	112	621763	91.66	ng	0.00
7) Phenol-d6	6.845	99	832704	86.96	ng	0.00
23) Nitrobenzene-d5	8.822	82	572793	56.29	ng	0.00
42) 2,4,6-Tribromophenol	15.798	330	239800	86.69	ng	0.00
45) 2-Fluorobiphenyl	12.927	172	960786	48.37	ng	0.00
79) Terphenyl-d14	19.680	244	1114303	58.84	ng	0.00
Target Compounds						
						Qvalue
10) Phenol	6.869	94	24657	2.62	ng	79
20) 3+4-Methylphenols	8.434	107	28174	3.39	ng	95
31) Naphthalene	10.492	128	58670	2.39	ng	100
78) Pyrene	19.468	202	57697	2.31	ng	98
84) Bis(2-ethylhexyl)phtha...	21.145	149	137680	7.95	ng	# 93

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

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