

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041221\
 Data File : BP005270.D
 Acq On : 12 Apr 2021 17:25
 Operator : CG/JU
 Sample : M1967-10
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SB1-B

Quant Time: Apr 12 18:37:54 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP040921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 09 18:44:11 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.675	152	33515	20.00	ng	# 0.00
21) Naphthalene-d8	10.469	136	116287	20.00	ng	# 0.00
39) Acenaphthene-d10	14.340	164	71650	20.00	ng	0.00
64) Phenanthrene-d10	17.104	188	169810	20.00	ng	# 0.00
76) Chrysene-d12	21.204	240	188505	20.00	ng	0.00
86) Perylene-d12	23.474	264	186389	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.281	112	268479	84.07	ng	0.00
7) Phenol-d6	6.863	99	341154	93.63	ng	0.00
23) Nitrobenzene-d5	8.840	82	231380	62.75	ng	0.00
42) 2,4,6-Tribromophenol	15.839	330	29732	32.87	ng	0.00
45) 2-Fluorobiphenyl	12.951	172	299124	56.36	ng	0.00
79) Terphenyl-d14	19.680	244	484613	50.50	ng	0.00
Target Compounds						
						Qvalue
10) Phenol	6.893	94	228030	61.10	ng	75
20) 3+4-Methylphenols	8.463	107	9062	3.38	ng	83
50) Dimethylphthalate	13.804	163	12858	2.44	ng	# 96
75) Fluoranthene	19.128	202	34962	2.97	ng	96
78) Pyrene	19.480	202	38329	3.27	ng	97
84) Bis(2-ethylhexyl)phtha...	21.116	149	58917	8.57	ng	99
88) Benzo(b)fluoranthene	22.786	252	32186	2.64	ng	# 82

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

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