

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050725\
 Data File : BP024563.D
 Acq On : 07 May 2025 14:41
 Operator : RC/JU
 Sample : Q1947-04
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MH-P

Quant Time: May 07 15:43:41 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP041425.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 18 12:04:48 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.675	152	116263	20.000	ng	-0.05
21) Naphthalene-d8	10.457	136	413670	20.000	ng	#-0.04
39) Acenaphthene-d10	14.334	164	114583	20.000	ng	-0.02
64) Phenanthrene-d10	17.151	188	357126	20.000	ng	# 0.00
76) Chrysene-d12	21.604	240	93816	20.000	ng	# 0.01
86) Perylene-d12	24.945	264	2581	20.000	ng	# 0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.299	112	878269	124.905	ng	-0.05
7) Phenol-d6	6.863	99	662147	68.772	ng	-0.05
23) Nitrobenzene-d5	8.822	82	698634	96.301	ng	-0.05
42) 2,4,6-Tribromophenol	15.863	330	481319	303.611	ng	0.00
45) 2-Fluorobiphenyl	12.934	172	1399838	187.053	ng	-0.03
79) Terphenyl-d14	19.875	244	1689854	363.065	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.798	163	18939	2.273	ng	97
78) Pyrene	19.663	202	17444	2.926	ng	98

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

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