

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050925\
 Data File : BP024592.D
 Acq On : 09 May 2025 15:46
 Operator : RC/JU
 Sample : Q1968-04
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MH-M

Quant Time: May 09 16:28:09 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP041425.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 09 12:07:46 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.699	152	118424	20.000	ng	0.00
21) Naphthalene-d8	10.469	136	445558	20.000	ng	# 0.00
39) Acenaphthene-d10	14.334	164	260832	20.000	ng	0.00
64) Phenanthrene-d10	17.128	188	542499	20.000	ng	# 0.00
76) Chrysene-d12	21.569	240	704694	20.000	ng	# 0.00
86) Perylene-d12	24.892	264	803227	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.328	112	895054	124.969	ng	0.00
7) Phenol-d6	6.893	99	996480	101.608	ng	0.00
23) Nitrobenzene-d5	8.846	82	742480	95.020	ng	0.00
42) 2,4,6-Tribromophenol	15.845	330	635983	176.234	ng	0.00
45) 2-Fluorobiphenyl	12.940	172	1642675	96.427	ng	0.00
79) Terphenyl-d14	19.845	244	3312385	94.744	ng	0.00

Target Compounds	Qvalue

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

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