

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102125\
 Data File : BP025987.D
 Acq On : 21 Oct 2025 15:35
 Operator : RC/JU
 Sample : Q3369-04
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MH-22

Quant Time: Oct 21 15:57:55 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP101525.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 15 23:48:51 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.678	152	171438	20.00	ng	-0.02
21) Naphthalene-d8	10.448	136	635897	20.00	ng	-0.01
39) Acenaphthene-d10	14.319	164	385704	20.00	ng	0.00
64) Phenanthrene-d10	17.136	188	696834	20.00	ng	0.02
76) Chrysene-d12	21.583	240	587973	20.00	ng	# 0.00
86) Perylene-d12	24.924	264	769574	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.325	112	1341130	128.26	ng	-0.01
7) Phenol-d6	6.896	99	1576808	108.06	ng	-0.01
23) Nitrobenzene-d5	8.831	82	1193018	84.33	ng	-0.02
42) 2,4,6-Tribromophenol	15.848	330	628398	121.02	ng	0.00
45) 2-Fluorobiphenyl	12.919	172	2479566	82.40	ng	-0.01
79) Terphenyl-d14	19.860	244	2878647	78.25	ng	0.00

Target Compounds	Qvalue

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

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