

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102225\
 Data File : BP026011.D
 Acq On : 22 Oct 2025 23:22
 Operator : RC/JU
 Sample : Q3402-09
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 41-LB

Quant Time: Oct 23 10:42:40 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	7.702	152	165237	20.000	ng	0.00
21) Naphthalene-d8	10.466	136	671609	20.000	ng	0.00
39) Acenaphthene-d10	14.325	164	486656	20.000	ng	0.00
64) Phenanthrene-d10	17.137	188	975190	20.000	ng	0.02
76) Chrysene-d12	21.595	240	646839	20.000	ng	0.01
86) Perylene-d12	24.936	264	704631	20.000	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.337	112	1382070	137.133	ng	0.00
7) Phenol-d6	6.908	99	1669600	118.709	ng	0.00
23) Nitrobenzene-d5	8.849	82	1269953	84.992	ng	0.00
42) 2,4,6-Tribromophenol	15.848	330	791149	120.758	ng	0.00
45) 2-Fluorobiphenyl	12.925	172	2936341	77.336	ng	0.00
79) Terphenyl-d14	19.860	244	3771580	93.187	ng	0.00
Target Compounds						
77) Benzidine	19.501	184	226518	10.929	ng	Qvalue 98

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

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